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Open Notebook Science Challenge

Solubilities

of

Organic Compounds

in

Organic Solvents

Compiled and Measured by:
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Rajarshi Guha, Research Scientist at the NIH Chemical Genomics Center
Antony Williams, Vice President of Strategic Development, ChemSpider at the Royal Society of Chemistry
Bill Hooker, Postdoctoral Researcher in Molecular Biology
Andrew Lang, Professor of Mathematics at Oral Roberts University
Brent Friesen, Associate Professor of Chemistry at Dominican University
and
Tim Bohinski, David Bulger, Matthew Federici, Jenny Hale, Jenna Mancinelli, Khalid Mirza, Marshall Moritz, Daniel Rein, Cedric Tchakounte, and Hai Truong

Version: February 11, 2010

Preface

The Open Notebook Science Solubility Challenge

Solubility is an important consideration for many chemistry applications. Synthetic chemists usually use a solvent to perform reactions and knowledge of the solubility of the starting materials or products can be very useful to pick an appropriate solvent. Analytical chemists can use solubility to design separation techniques and factor in dynamic range considerations. Physical chemists can create and evaluate their models of how molecules interact in the solubilization and precipitation processes.

Solubility data can be obtained from a variety of online and offline sources. As with all chemical data, it can be a challenge to evaluate reported measurements. Some databases offer no references while others provide citations to peer reviewed journal articles. Given the choice, more weight is generally given to the latter. This is reasonable in most cases because more information about the purity of compounds and the methods used are available in peer-reviewed articles.

However, the information for how a specific measurement was obtained within a journal article is not generally provided. General methods are provided but the raw data for a specific measurement are typically not published. Peer review is not intended to validate individual measurements - its function is to ensure that the authors made appropriate conclusions based on their processed datasets and the state of knowledge in the field.

The Open Notebook Science Challenge was initiated in the fall of 2008 as the result of a discussion on a train in the UK between Jean-Claude Bradley and Cameron Neylon.[1,2] The concept was very simple: create a crowdsourcing opportunity for the chemistry community to contribute solubility measurements under Open Notebook Science conditions. This method of publication entails providing immediate public access to the chemist's laboratory notebook, as well as all raw data used to compute the measurements.[3,4]

On Sept 3, 2008 the first ONSC measurements were recorded by Bradley and Neylon at the University of Southampton in Neylon's laboratory.[5] The project was soon sponsored by Submeta, offering ten $500 awards for students in the US or the UK who best recorded how they performed their experiments.[6] Furthermore, the first 3 winners also received one year subscriptions to Nature magazine, thanks to a sponsorship from the Nature Publishing Group.[7] Sigma-Aldrich supported the contest by donating chemicals upon request.[8]

Students were evaluated by a group of judges who convened once a month to deliberate the next award. Judges also provided feedback to the students by commenting on their lab notebook pages directly on the wiki. Their expertise ranged from chemistry to mathematics, spectroscopy and molecular biology.

Techniques

Participants in the ONS Challenge were not required to use a specific method to measure solubility - although they were required to properly document their experiments and analyses. Due to its simplicity, most measurements in the past year were made using the SAMS NMR technique, requiring no volume measurement or calibration curves.[9] Two assumptions are made with this method. The first is that the volume of solute and solvent are additive, with the error becoming negligible at low solubility values. The second is that NMR integration values are proportional to the amount of solvent and solute. Some deviations from this have been observed for default NMR parameters and in later experiments long relaxation times are introduced into the protocol (D1 = 50s).[10]

Data Curation

Since an Open Notebook approach is used in this work, those interested in the validity of the measurements can assess the methods used - both for the preparation of saturated solutions and the raw data from the measurements. Over time, values in the database are likely to improve and possibly some errors may be uncovered and corrected. However, on the whole, we feel that the values provided in this work should be of use to chemists trying to gain an appreciation of solubility for most applications. This is especially the case for values that are not obtainable from any other source.

When clearly erroneous data points are discovered, they are flagged in the database as "DONOTUSE". This way interfaces with the dataset can ignore these values while allowing anyone to investigate why the data points were flagged. This might happen when early experiments did not allow for sufficient mixing or NMR D1 relaxation times were long enough to fully integrate peaks of interest. Out of 681 reported measurements, 51 are currently marked in this way. A shared Google Spreadsheet is used to collect and curate the dataset. This allows easy data entry while providing a simple way to interrogate the database for visualization applications via the Google API.[11]

Literature data and format conversions

An additional 400 solubility measurements from the literature are included in the database. These generally correspond to compounds that are structurally identical or similar to the compounds measured by the ONS Challenge participants. These values are averaged in with the values from the participants, with appropriate references provided. In order to compare values, conversions from molar fraction or g solute/100g solvent to molarity were made by assuming that the volumes are additive and obtaining the density of the solutes in most cases from the predicted values in ChemSpider.[12]

For the convenience of chemists with diverse applications, all three formats are provided. For the cases where solutes are miscible with the solvent, the molarity reported is simply the solute's density. The practical interpretation of this is that solutions of any molarity below the solute's density can be prepared.

In the process of converting units and averaging heterogeneous data sources, no attempt has been made to track significant figures.

Those interested in any information about the precision of measurements should consult each individual data source. This may not be an easy task for measurements only carried out once and where factors such as the quality of spectral peaks and baselines are not optimal.

This collection will be most valuable for those who do not require highly precise measurements for their applications. For example, synthetic chemists can easily use rough estimates of solubility to select appropriate solvents for a reaction. In any case, one would be wise to consider all measurements as provisional, regardless of the source. As more data are collected, subsequent editions of this book will adjust values accordingly.

Predicted Solubilities

Predicted solubility values for selected solutes in a range of solvents are provided in a section. Specifically, solutes are included when measurements from at least 5 different solvents are available. A method using Abraham descriptors depends on the experimental solubility measurements from several solvents to make predictions, which is detailed in that section of the book. For this reason, some aqueous solubility measurements are included, which are generally available from the literature. Predicted aqueous solubilities are used when experimental values are not available. The focus of this collection remains on non-aqueous solubility.

Consistent with how the experimental measurements are made available, the predicted solubility values are provided as a work in progress. The purpose in providing them is to suggest solvents of interest for various applications. The boiling point of each solvent is also listed in the table to allow a convenient selection. When available, experimental measurements are listed next to the predicted values. This information can be helpful to gauge the usefulness of the model to some extent but does not guarantee its reliability for the other solvents. As more measurements are collected the reliability of the predictions is likely to increase and this will be reflected in future editions of this book.

Searching the database

The values in this database can be accessed and filtered in various ways. More information is available at the ONS Challenge wiki[13] and Chapter 16 of the book "Beautiful Data".[14]

Database version

Archived as an Excel Spreadsheet by WebCite on February 11, 2010.[15]

Notebook archive

This is the first edition to include a full archive of the ONS Challenge notebook. A space export from Wikispaces provides an initial version of all the HTML pages in the notebook with local hyperlinks to copies of all images and files uploaded onto the wiki. All of the Google Spreadsheets are automatically downloaded as Excel spreadsheets and placed in the same "files" folder as the images. NMR spectra, stored as JCAMP-DX files, are placed in the "spectra" folder. All of the HTML pages are reformatted to provide local references to both Excel spreadsheets and the JCAMP-DX files.

The notebook archive is meant to represent a snapshot of the state of all source documents at the time of the publication of an edition of this book. When used from a server with web services running, clicking on links to the spectra will allow interaction via a browser interface, including zooming in or out and integration of the NMR spectrum. When accessed in stand-alone mode after downloading or directly from a CD, everything will work the same, except that JCAMP-DX files must be open from JSpecView running on the desktop. Excel files will retain any calculations in the cells of the original Google Spreadsheets but dynamic values generated from calling web services - such the script that automatically integrates NMR spectra - will be frozen as simple values. However the link to the web service used will be stored in the cell as a comment. Links to external websites are not crawled and embedded Google Spreadsheets or videos are not copied. These will work but will reflect live data on the web.

The February 11, 2010 version of the notebook archive is available on a hosted site[16], on a CD[17] or by download[18].

References

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[17] Notebook Archive CD. **http://www.lulu.com/content/compact-disc/onschallenge-archive/8337307**
[18] Notebook Archive ZIP download. **http://showme.physics.drexel.edu/onsc/archives/onschallenge2010-02-11/onschallenge2010-02-11.zip**

Contents

Judges

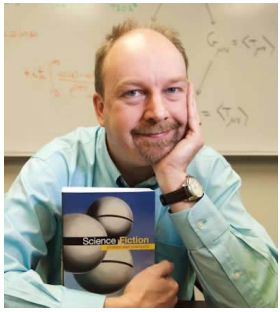
Jean-Claude Bradley is an Associate Professor of Chemistry and E-Learning Coordinator for the College of Arts and Sciences at Drexel University. He leads the UsefulChem project, an initiative started in the summer of 2005 to make the scientific process as transparent as possible by publishing all research work in real time to a collection of public blogs, wikis and other web pages. Jean-Claude coined the term Open Notebook Science to distinguish this approach from other more restricted forms of Open Science. Jean-Claude teaches undergraduate organic chemistry courses with most content freely available on public blogs, wikis, games, Second Life and audio and video podcasts. He has a Ph.D. in organic chemistry and has published articles and obtained patents in the areas of synthetic and mechanistic chemistry, gene therapy, nanotechnology and scientific knowledge management. **http://usefulchem.blogspot.com**

Cameron Neylon is a biophysicist who has always worked in interdisciplinary areas and is an advocate of open research practice and improved data management. He currently works as Senior Scientist in Biomolecular Sciences at the ISIS Neutron Scattering facility at the Science and Technology Facilities Council (STFC). Along with his work in structural biology and biophysics his research and writing focuses on the interface of web technology with science and the successful (and unsuccessful) application of generic and specially designed tools in the academic research environment. He is a founder member of the Open Knowledge Foundation Science Working Group and writes regularly on Open Research at his blog, Science in the Open. **http://blog.openwetware.org/scienceintheopen**

Rajarshi Guha is a Research Scientist at the NIH Chemical Genomics Center (NCGC). At the NCGC he is involved in cheminformatics methodology and software development for various aspects of high throughput screening for small molecules and RNAi. His research makes extensive use of statistical methods and has been applied to a variety of biological systems. Along with algorithm development, he is extensively involved in cheminformatics software development, including development of toolkits, web services and integration of these into distributed infrastructures. As a believer in Open Source and Open Data, much of his research and software is available under Open Source licenses. **http://blog.rguha.net**

Antony Williams is the host of ChemSpider, the online chemistry portal leading the charge towards open and collaborative chemistry. ChemSpider provides access to over 23 million unique chemical entities sourced from almost 300 data sources and provides a platform for community based depositions and curation to clean up internet-based chemistry. Antony spent over a decade in the commercial scientific software business as Chief Science Officer for a Cheminformatics software company and during his tenure oversaw their product development, marketing and sales teams. He is an accomplished NMR spectroscopist with over 100 peer-reviewed publications. During his career he was the NMR Technology Leader for the Eastman-Kodak company and has worked in both academia and national government research institutions. He has recently taken his passion for providing access to chemistry related information and software services to the masses by hosting the ChemSpider service. **http://www.chemspider.com/blog/**

Bill Hooker is a molecular biologist by trade and has worked on G protein signaling, PCR diagnostics, anti-schistosome vaccines, HIV replication and molecular mechanisms of cancer. He serves as an associate editor on the innovative Open Access journal BioMed Central Research Notes and as a member of the advisory board of the Berglund Center for Internet Studies. He is currently the R&D team leader at a small biotech company in Portland, OR. He has never had an idea that couldn't be improved by sharing it with as many people as possible, and doesn't believe that anyone else has, either. He is particularly interested in extending the "open" ethos of Open Source software and Open Access publishing to all aspects of science, from raw data to hypothesis testing by way of distributed, collaborative effort. **http://www.sennoma.net**

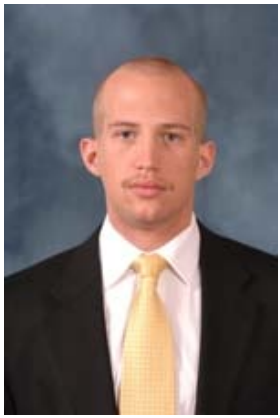


Andrew Lang is a Professor of Mathematics at Oral Roberts University. His PhD training is in the area of quantum field theory in curved spacetime. While remaining active in this area he has always enjoyed working collaboratively on interdisciplinary projects ranging from modeling basketball free throws to the stability of spinning spacecraft under thrust. His interests include the relationship between science and science fiction and the epistemological differences between teleology and metaphysical naturalism. As a result of being a judge for the Open Notebook Science Challenge he has developed an appreciation for openess in science, multi-dimensional data visualization and cheminformatics. **http://webapps.oru.edu/new_php/blog/index.php?user_id=23**

Educational Partners

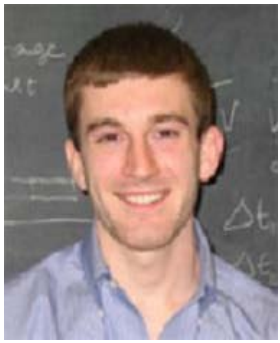
J. Brent Friesen from Dominican University incorporated the Open Notebook Science Solubility Challenge into his Sophomore Organic Chemistry laboratory course in January 2009. He has been a professor in the natural science department at Dominican University since 1999. He regularly teaches Organic Chemistry (lecture and laboratory), Biochemistry, Introduction to Organic Chemistry, Chemistry of Natural Products, and Forensic Chemistry. His research focuses on the automation of countercurrent separation technology, anti-tuberculosis compounds, and interactions with Reichardt's Dye. Friesen has also done research with undergraduates on projects such as "Biorenewable Solvents for Countercurrent Chromatography" and "Used Coffee Grounds: A New Source of Biofuels." Friesen is a regular volunteer with the Oak Park Education Foundation's Global Village, a program matching scientists with elementary school classrooms. His recent articles include "Saying What You Mean: Teaching Mechanisms in Organic Chemistry" (Journal of Chemical Education) and "Countercurrent Separation of Natural Products" (Journal of Natural Products).

Students



Tim Bohinski (April 2009 Submeta Award winner) is a senior at Drexel University majoring in chemistry, with a strong interest in organic and theoretical chemistry. He will be graduating with a B.S. in June 2010. He is currently performing undergraduate research in the theoretical/physical field, studying nanoparticle optics, specifically in the area of plasmoms. He plans to continue his studies at the doctorate level. Tim is also a member of the Drexel American Chemical Society and was a member of the Drexel Crew Team. In my spare time he enjoys the outdoors, staying fit, and homebrewing. LinkedIn profile: **http://www.linkedin.com/pub/timothy-bohinski/18/110/a7 Supervisor: Jean-Claude Bradley, Drexel**

David Bulger (February 2009 Submeta Award winner) is a junior Chemistry major at Oral Roberts University in Tulsa, OK. Since freshman year, he has been involved in research ranging from water purity, red tide in Second Life, collaboration between water utilities and academia, NMR solubility determination and Sortase A inhibition. After his undergraduate degree, he plans to pursue an M.D. or M.D./Ph.D. **http://biolab.isis.rl.ac.uk/david_bulger. Supervisor: Robert Stewart, Oral Roberts University**



Matthew Federici (June 2009 Submeta Award winner) graduated from the University of Virginia in 2008. After receiving the Naval ROTC Tweeddale Scholarship Award in 2004, Matthew studied naval nuclear engineering and interned on the USS Alexandria. As a varsity student athlete, he doubled majored in mechanical engineering as well as aerospace engineering. Matthew spent his senior year at UVa studying biomechanics under Dr. Pradip Sheth. His undergraduate academic career finished by winning the Cyrus Society Award, given to a select few of student athletes with outstanding academic performance and community service. Matthew is pursuing his PhD in Mechanical Engineering under Dr. Minjun Kim at Drexel University. As a GAANN fellow, he is currently investigating biomimetics and drug delivery systems. Matthew plans to attend medical school following the completion of his dissertation. Website: **http://microfluidics.tripod.com/ Supervisor: Jean-Claude Bradley, Drexel**

Jennifer (Jenny) Hale (December 2008 Submeta Award winner) gained a BSc in Chemistry with Pharmacology from the University of Southampton, UK, in 2004. Following her degree she spent 15 months working on the synthesis of phosphoramidites before taking up a PhD in the department of chemical biology in 2006. Jenny's area of research for her PhD has been the directed evolution of a beta-glucuronidase into a beta-galactosidase via neutral drift, a phenomenon found in natural evolution that has only recently begun to be explored in laboratory evolution. Additionally, her research has been used in the development and testing of an electronic laboratory notebook in the form of a blog: LaBLog **(http://blogs.chem.soton.ac.uk/neutral_drift)** as well as wider investigations into Open Notebook Science. Outside of the laboratory, Jenny is a keen ballroom and latin dancer and spent a year as president of the Southampton University Ballroom and Latin Dance Society, as well as competing in numerous university dancesport competitions. **Supervisor: Cameron Neylon, Southampton**

Jenna Mancinelli (September 2009 Submeta Award winner) is currently a junior at Drexel University, studying Biology. She is particularly interested in human physiology and its relation to clinical medicine. She hopes to attend medical school following her graduation from Drexel University. In her spare time, Jenna enjoys going to concerts, studying Italian and experimenting with her culinary skills. **Supervisor: Jean-Claude Bradley, Drexel**

Khalid Baig Mirza (January 2009 Submeta Award winner) obtained a BSc in Chemistry and Biology from Osmania University, India. He earned an MS in Chemistry from Fairleigh Dickinson University at Madison, NJ in 2002. He is currently working towards a PhD at Drexel University under the guidance of Prof. Jean-Claude Bradley. His research focuses on the synthesis of potentially active anti-malarial diamides obtained from a Ugi reaction. His research is performed in the open as 'Open Notebook Science'. He enjoys gardening, cricket and keeping up to date on world affairs. **Supervisor: Jean-Claude Bradley, Drexel**

Marshall Moritz (July 2009 Submeta Award winner) is a chemistry and mathematics double major at Syracuse University in Syracuse, NY with an anticipated graduation in May 2011. After studying abroad in Hong Kong, he is looking forward to continuing his studies at SU and pursuing further research not only in solubility determination by NMR but also in other areas of organic chemistry. Following graduation, he hopes to continue his education in graduate school with a masters degree. He is an avid cyclist, cycling cross country in 2006, and is also a traveling and music enthusiast. **Supervisor: Jean-Claude Bradley, Drexel**

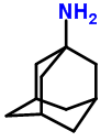
Daniel Rein (August 2009 Submeta Award winner) is currently a student studying chemistry and biology at Drexel University. He plays guitar in Jazz Band and is passionate about music. For freshman year he studied physics at Northeastern. **http://www.linkedin.com/pub/daniel-rein/18/31a/72b Supervisor: Jean-Claude Bradley, Drexel**

Cedric Tchakounte (March 2009 Submeta Award winner) is a senior majoring in Biological Sciences & Biotechnology at Drexel University. He is a distinguished student, having received multiple Dean's List honors and the Who's Who Among Students in American Universities & Colleges award. He has also shown his dedication to community and service by being a mentor for Guided Youths. He enjoys performing research, because he believes it to be a cornerstone to learning true science. With his academic chapter as an undergraduate coming to an end, Cedric looks forward to what awaits. He plans on pursuing a career in medicine. **Supervisor: Jean-Claude Bradley, Drexel**

Hai Truong (December 2009 Submeta Award winner) is pursuing a BS in Chemistry at Drexel University from 2009. After graduating with honors from a high school for the Gifted in Vietnam, he came to the United States for college. In 2008, he attended Collin College in McKinney,TX. At Collin, he was a member of Phi Theta Kappa, an international honor society for two-year colleges. In 2009, he transfered to Drexel University with a Dean's Scholarship. Beside academic studies, Hai enjoys swimming and playing tennis. **Supervisor: Jean-Claude Bradley, Drexel**

Solubilities

1-adamantylamine $C_{10}H_{17}N$



Compound Data

Molecular weight	151.249	**H bond acceptors**	1	**Rule of 5 violations**	0
Compound type	amine	**H bond donors**	2	**ACD/ALogP**	2.22
Phase 25°C	solid	**Rotatable bonds**	1	**Predicted density**	1.066 g/cm^3
SMILES	C1C2CC3CC1CC(C2)(C3)N				
InChIKey	DKNWSYNQZKUICI-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
methanol	0.36	0.02	7.62

1-octadecylamine $C_{18}H_{39}N$[69, 908]

Compound Data

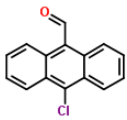
Molecular weight	269.509	**H bond acceptors**	1	**Rule of 5 violations**	1
Compound type	amine	**H bond donors**	2	**ACD/ALogP**	8.37
Phase 25°C	solid	**Rotatable bonds**	17	**Predicted density**	0.818 g/cm^3
SMILES	NCCCCCCCCCCCCCCCCCC				

InChIKey	REYJJPSVUYRZGE-UHFFFAOYSA-N

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
acetonitrile	0.08	0.00	2.96
ethanol	1.77	0.20	146.72
hexane	0.13	0.02	5.42
methanol	1.68	0.14	134.67
THF	1.68	0.23	112.18
toluene	1.34	0.20	74.24

10-chloro-9-anthraldehyde $C_{15}H_9ClO$[57]



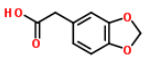
Compound Data

Molecular weight	240.684	**H bond acceptors**	1	**Rule of 5 violations**	0
Compound type	aldehyde	**H bond donors**	0	**ACD/ALogP**	4.7
Phase 25°C	solid	**Rotatable bonds**	1	**Predicted density**	1.327 g/cm^3
SMILES	O=Cc2c1c(cccc1)c(Cl)c3c2cccc3				
InChIKey	SHYBXXMECBHHFH-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
ethanol	0.00	0.00	0.00
methanol	0.00	0.00	0.00
THF	0.08	0.01	2.16

2-(1,3-benzodioxol-5-yl)acetic acid $C_9H_8O_4$[98, 82]



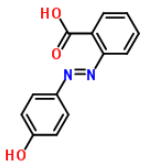
Compound Data

Molecular weight	180.157	**H bond acceptors**	4	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	1.37
Phase 25°C	solid	**Rotatable bonds**	2	**Predicted density**	1.406 g/cm^3
SMILES	O=C(O)Cc1ccc2OCOc2c1				
InChIKey	ODVLMCWNGKLROU-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
acetonitrile	0.67	0.04	17.68
DMSO	5.37	0.55	282.22
THF	2.73	0.25	83.68

2-(4-hydroxyphenyl-azo)benzoic acid $C_{13}H_{10}N_2O_3$[3, 32]



Compound Data

Molecular weight	242.23	**H bond acceptors**	5	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	2	**ACD/ALogP**	3.72
Phase 25°C	solid	**Rotatable bonds**	4	**Predicted density**	1.3 g/cm^3
SMILES	Oc2ccc(N=Nc1ccccc1C(O)=O)cc2				
InChIKey	DWQOTEPNRWVUDA-CCEZHUSRSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
ethanol	0.09	0.01	2.68
methanol	0.06	0.00	1.95
THF	0.55	0.05	16.42

2-chloro-5-nitrobenzaldehyde $C_7H_4ClNO_3$[105, 69, 208, 33, 205]

Compound Data

Molecular weight	185.565	**H bond acceptors**	4	**Rule of 5 violations**	0
Compound type	aldehyde	**H bond donors**	0	**ACD/ALogP**	2.5
Phase 25°C	solid	**Rotatable bonds**	2	**Predicted density**	1.485 g/cm^3
SMILES	O=Cc1cc(ccc1Cl)[N+]([O-])=O				
InChIKey	VFVHWCKUHAEDMY-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
acetonitrile	2.31	0.15	80.47
chloroform	2.90	0.27	56.27
DMSO	2.93	0.25	78.05
ethanol	*	*	*
methanol	*	*	*
THF	2.79	0.25	87.80
toluene	1.74	0.19	47.44

* This aldehyde reacts with alcohols to form a hemiacetal.

2-chloro-5-nitrobenzoic acid $C_7H_4ClNO_4$[905]

Compound Data

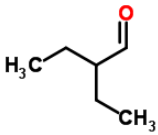
Molecular weight	201.564	**H bond acceptors**	5	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	2.02
Phase 25°C	solid	**Rotatable bonds**	2	**Predicted density**	1.602 g/cm^3
SMILES	O=C(O)c1cc(ccc1Cl)[N+]([O-])=O				
InChIKey	QUEKGYQTRJVEQC-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
1-butanol	0.78	0.07	21.66
1-decanol	0.39	0.07	9.98
1-heptanol	0.54	0.08	14.24
1-hexanol	0.64	0.08	17.19
1-octanol	0.48	0.07	12.51
1-pentanol	0.74	0.08	20.28
1-propanol	0.95	0.08	27.36
2-butanol	0.91	0.09	25.86
2-methyl-1-propanol	0.63	0.06	17.22
2-methyl-2-propanol	1.21	0.12	35.78
2-pentanol	0.76	0.08	20.94
2-propanol	1.13	0.09	33.57
3-methyl-1-butanol	0.68	0.07	18.53
butyl acetate	0.46	0.06	11.11
dibutyl ether	0.10	0.02	2.62
diethyl ether	0.57	0.06	16.86
diisopropyl ether	0.19	0.03	5.18
ethanol	1.40	0.09	43.91
ethyl acetate	0.75	0.08	18.59
methyl acetate	0.89	0.08	22.25

pentyl acetate	0.43	0.06	10.39
THF	2.97	0.27	105.73

2-ethylbutyraldehyde $C_6H_{12}O$[22]



Compound Data

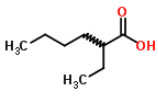
Molecular weight	100.159	**H bond acceptors**	1	**Rule of 5 violations**	0
Compound type	aldehyde	**H bond donors**	0	**ACD/ALogP**	1.79
Phase 25°C	liquid	**Rotatable bonds**	3	**Predicted density**	0.799 g/cm^3
SMILES	O=CC(CC)CC				
InChIKey	UNNGUFMVYQJGTD-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
methanol	8.13	¤	¤

¤ Solute is very soluble/miscible, conversion fail.

2-ethylhexanoic acid $C_8H_{16}O_2$[22]



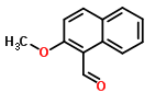
Compound Data

Molecular weight	144.211	**H bond acceptors**	2	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	2.72
Phase 25°C	liquid	**Rotatable bonds**	5	**Predicted density**	0.926 g/cm^3
SMILES	O=C(O)C(CC)CCCC				
InChIKey	OBETXYAYXDNJHR-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
methanol	6.26	0.91	4777.15

2-methoxy-1-naphthaldehyde $C_{12}H_{10}O_2$[205]



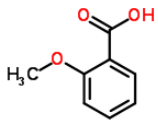
Compound Data

Molecular weight	186.207	**H bond acceptors**	2	**Rule of 5 violations**	0
Compound type	aldehyde	**H bond donors**	0	**ACD/ALogP**	2.95
Phase 25°C	solid	**Rotatable bonds**	2	**Predicted density**	1.169 g/cm^3
SMILES	O=Cc1c2c(ccc1OC)cccc2				
InChIKey	YIQGLTKAOHRZOL-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
methanol	1.26	0.06	38.98

2-methoxybenzoic acid $C_8H_8O_3$[905, 34, 907]



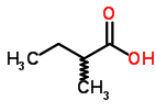
Compound Data

Molecular weight	152.147	**H bond acceptors**	3	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	1.5
Phase 25°C	solid	**Rotatable bonds**	2	**Predicted density**	1.207 g/cm^3
SMILES	O=C(O)c1ccccc1OC				
InChIKey	ILUJQPXNXACGAN-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
1-butanol	0.57	0.05	11.61
1-decanol	0.22	0.04	4.16
1-heptanol	0.34	0.05	6.59
1-hexanol	0.37	0.05	7.24
1-octanol	0.30	0.05	5.76
1-pentanol	0.47	0.05	9.37
1-propanol	0.76	0.06	16.09
1,4-dioxane	1.53	0.14	28.99
2-butanol	0.50	0.05	10.14
2-ethyl-1-hexanol	0.25	0.04	4.78
2-methyl-1-butanol	0.39	0.04	7.71
2-methyl-1-pentanol	0.35	0.04	6.84
2-methyl-1-propanol	0.41	0.04	8.21
2-methyl-2-propanol	0.53	0.05	10.75
2-pentanol	0.39	0.04	7.71
2-propanol	0.59	0.05	12.26
3-methyl-1-butanol	0.36	0.04	7.09
4-methyl-2-pentanol	0.32	0.04	6.26
butyl acetate	0.34	0.04	6.10
chloroform	0.46	0.04	5.01
dibutyl ether	0.04	0.01	0.78
diethyl ether	0.24	0.02	5.13
diisopropyl ether	0.06	0.01	1.21
ethanol	1.19	0.08	27.31
ethyl acetate	5.63	0.66	328.57
methanol	1.82	0.09	47.72
methyl acetate	0.86	0.07	16.16
pentyl acetate	0.26	0.04	4.64
propylene carbonate	0.99	0.13	14.28
THF	2.02	0.18	45.61
water	0.03	0.00	0.46

2-methyl butyric acid $C_5H_{10}O_2$[22]



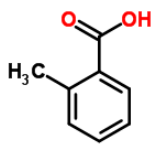
Compound Data

Molecular weight	102.132	**H bond acceptors**	2	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	1.13
Phase 25°C	liquid	**Rotatable bonds**	2	**Predicted density**	0.962 g/cm^3
SMILES	O=C(O)C(C)CC				
InChIKey	WLAMNBDJUVNPJU-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
methanol	9.16	0.93	4515.12

2-methylbenzoic acid $C_8H_8O_2$[905, 907]



Compound Data

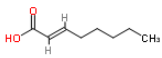
Molecular weight	136.148	**H bond acceptors**	2	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	2.36
Phase 25°C	solid	**Rotatable bonds**	1	**Predicted density**	1.151 g/cm^3
SMILES	O=C(O)c1ccccc1C				
InChIKey	ZWLPBLYKEWSWPD-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)

1-butanol	1.61	0.15	33.64
1-decanol	1.04	0.18	19.50
1-heptanol	1.28	0.18	25.04
1-hexanol	1.38	0.17	27.52
1-octanol	1.16	0.18	22.24
1-pentanol	1.49	0.16	30.37
1-propanol	1.77	0.14	38.34
1,4-dioxane	2.64	0.25	52.53
2-butanol	1.75	0.17	37.51
2-methyl-1-butanol	1.25	0.14	24.69
2-methyl-1-propanol	1.28	0.12	25.64
2-methyl-2-propanol	2.19	0.21	50.05
2-pentanol	1.68	0.19	35.28
2-propanol	1.95	0.16	43.63
3-methyl-1-butanol	1.34	0.15	26.80
4-methyl-2-pentanol	1.39	0.17	27.93
butyl acetate	1.17	0.15	20.87
dibutyl ether	0.58	0.09	10.87
diethyl ether	1.65	0.17	38.03
diisopropyl ether	0.86	0.11	17.20
ethanol	2.08	0.14	48.15
ethyl acetate	1.49	0.15	27.42
methyl acetate	1.51	0.13	27.56
pentyl acetate	1.04	0.15	18.31
propylene carbonate	0.46	0.06	5.50
THF	2.92	0.26	67.18
water	0.01	0.00	0.14

2-octenoic acid $C_8H_{14}O_2$[22]



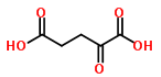
Compound Data

Molecular weight	142.196	**H bond acceptors**	2	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	2.93
Phase 25°C	liquid	**Rotatable bonds**	5	**Predicted density**	0.955 g/cm^3
SMILES	O=C(O)/C=C/CCCCC				
InChIKey	CWMPPVPFLSZGCY-VOTSOKGWSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
methanol	6.64	0.96	11068.64

2-oxopentanedioic acid $C_5H_6O_5$[25]



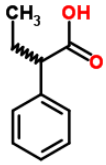
Compound Data

Molecular weight	146.098	**H bond acceptors**	5	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	2	**ACD/ALogP**	-1.43
Phase 25°C	solid	**Rotatable bonds**	4	**Predicted density**	1.499 g/cm^3
SMILES	O=C(O)C(=O)CCC(=O)O				
InChIKey	KPGXRSRHYNQIFN-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
ethanol	2.99	0.20	79.04
methanol	5.02	0.29	190.70
THF	3.12	0.26	72.46

2-phenylbutyric acid $C_{10}H_{12}O_2$[79, 83]



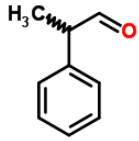
Compound Data

Molecular weight	164.201	**H bond acceptors**	2	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	2.38
Phase 25°C	solid	**Rotatable bonds**	3	**Predicted density**	1.09 g/cm^3
SMILES	O=C(O)C(c1ccccc1)CC				
InChIKey	OFJWFSNDPCAWDK-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
acetonitrile	6.18	0.83	1968.04
DMSO	5.97	0.81	886.12
ethanol	6.20	0.85	1977.18
THF	5.96	0.82	1059.60
toluene	5.35	0.74	519.73

2-phenylpropanal $C_9H_{10}O$[21]



Compound Data

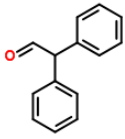
Molecular weight	134.175	**H bond acceptors**	1	**Rule of 5 violations**	0
Compound type	aldehyde	**H bond donors**	0	**ACD/ALogP**	2.13
Phase 25°C	liquid	**Rotatable bonds**	2	**Predicted density**	0.98 g/cm^3
SMILES	c1ccccc1C(C)C=O				
InChIKey	IQVAERDLDAZARL-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
methanol	7.53	¤	¤

¤ Solute is very soluble/miscible, conversion fail.

2,2-diphenylacetaldehyde $C_{14}H_{12}O$[21]



Compound Data

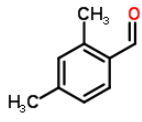
Molecular weight	196.245	**H bond acceptors**	1	**Rule of 5 violations**	0
Compound type	aldehyde	**H bond donors**	0	**ACD/ALogP**	3.67
Phase 25°C	liquid	**Rotatable bonds**	3	**Predicted density**	1.069 g/cm^3
SMILES	c1ccccc1C(c1ccccc1)C(=O)				
InChIKey	HLLGFGBLKOIZOM-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
methanol	5.64	¤	¤

¤ Solute is very soluble/miscible, conversion fail.

2,4-dimethylbenzaldehyde $C_9H_{10}O$[21]



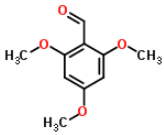
Compound Data

Molecular weight	134.175	**H bond acceptors**	1	**Rule of 5 violations**	0
Compound type	aldehyde	**H bond donors**	0	**ACD/ALogP**	2.56
Phase 25°C	liquid	**Rotatable bonds**	1	**Predicted density**	1.003 g/cm^3
SMILES	O=Cc1ccc(C)cc1C				
InChIKey	GISVICWQYMUPJF-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
methanol	7.17	0.88	3128.10

2,4,6-trimethoxybenzaldehyde $C_{10}H_{12}O_4$[82]



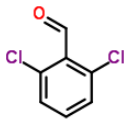
Compound Data

Molecular weight	196.2	**H bond acceptors**	4	**Rule of 5 violations**	0
Compound type	aldehyde	**H bond donors**	0	**ACD/ALogP**	1.49
Phase 25°C	solid	**Rotatable bonds**	4	**Predicted density**	1.133 g/cm^3
SMILES	O=Cc1c(OC)cc(OC)cc1OC				
InChIKey	CRBZVDLXAIFERF-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
THF	0.14	0.01	3.11

2,6-dichlorobenzaldehyde $C_7H_4Cl_2O$[208, 205]



Compound Data

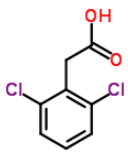
Molecular weight	175.012	**H bond acceptors**	1	**Rule of 5 violations**	0
Compound type	aldehyde	**H bond donors**	0	**ACD/ALogP**	3.03
Phase 25°C	solid	**Rotatable bonds**	1	**Predicted density**	1.4 g/cm^3
SMILES	O=Cc1c(Cl)cccc1Cl				
InChIKey	DMIYKWPEFRFTPY-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
acetonitrile	1.35	0.08	38.05
chloroform	3.41	0.32	69.35
ethanol	*	*	*
methanol	*	*	*
THF	2.48	0.22	69.58
toluene	1.74	0.19	44.68

* This aldehyde reacts with alcohols to form a hemiacetal.

2,6-dichlorophenylacetic acid $C_8H_6Cl_2O_2$[85, 82]



Compound Data

Molecular weight	205.038	**H bond acceptors**	2	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	2.71
Phase 25°C	solid	**Rotatable bonds**	2	**Predicted density**	1.456 g/cm^3
SMILES	Clc1cccc(Cl)c1CC(=O)O				
InChIKey	SFAILOOQFZNOAU-UHFFFAOYSA-N				

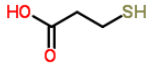
Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
THF	3.03	0.30	120.05

3-mercaptopropionic acid $C_3H_6O_2S$[22]

Compound Data

Molecular weight	106.144	**H bond acceptors**	2	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	0.43



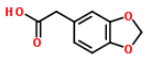
Phase 25°C	liquid	**Rotatable bonds**	3	**Predicted density**	1.223 g/cm^3
SMILES	C(CS)C(=O)O				
InChIKey	DKIDEFUBRARXTE-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
methanol	11.48	0.99	44306.20

3,4-(methylenedioxy)phenylacetic acid $C_9H_8O_4$[138]

Compound Data



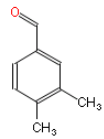
Molecular weight	180.157	**H bond acceptors**	4	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	1.37
Phase 25°C	solid	**Rotatable bonds**	2	**Predicted density**	1.406 g/cm^3
SMILES	O=C(O)Cc1ccc2OCOc2c1				
InChIKey	ODVLMCWNGKLROU-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
acetonitrile	0.10	0.01	2.44

3,4-dimethylbenzaldehyde $C_9H_{10}O$[59]

Compound Data



Molecular weight	134.175	**H bond acceptors**	1	**Rule of 5 violations**	0
Compound type	aldehyde	**H bond donors**	0	**ACD/ALogP**	2.56
Phase 25°C	liquid	**Rotatable bonds**	1	**Predicted density**	1.003 g/cm^3
SMILES	Cc1ccc(cc1C)C=O				
InChIKey	POQJHLBMLVTHAU-UHFFFAOYSA-N				

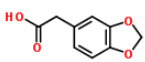
Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
ethanol	7.54	¤	¤
methanol	7.54	¤	¤
THF	7.54	¤	¤

¤ Solute is very soluble/miscible, conversion fail.

3,4-methy lendioxyp henylacet ic acid $C_9H_8O_4$[129, 132, 131]

Compound Data



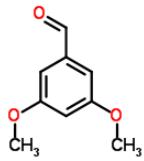
Molecular weight	180.157	**H bond acceptors**	4	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	1.37
Phase 25°C	solid	**Rotatable bonds**	2	**Predicted density**	1.406 g/cm^3
SMILES	O=C(O)Cc1ccc2OCOc2c1				
InChIKey	ODVLMCWNGKLROU-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
ethanol	0.55	0.03	13.67
methanol	0.94	0.04	25.57

THF	1.95	0.17	51.81

3,5-dimethoxybenzaldehyde $C_9H_{10}O_3$[205, 208]



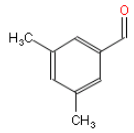
Compound Data

Molecular weight	166.174	**H bond acceptors**	3	**Rule of 5 violations**	0
Compound type	aldehyde	**H bond donors**	0	**ACD/ALogP**	1.53
Phase 25°C	solid	**Rotatable bonds**	3	**Predicted density**	1.114 g/cm^3
SMILES	O=Cc1cc(OC)cc(OC)c1				
InChIKey	VFZRZRDOXPRTSC-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
methanol	1.71	0.09	50.66
THF	4.03	0.45	185.73

3,5-dimethylbenzaldehyde $C_9H_{10}O$[63]



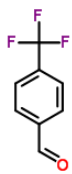
Compound Data

Molecular weight	134.175	**H bond acceptors**	1	**Rule of 5 violations**	0
Compound type	aldehyde	**H bond donors**	0	**ACD/ALogP**	2.56
Phase 25°C	liquid	**Rotatable bonds**	1	**Predicted density**	1.003 g/cm^3
SMILES	Cc1cc(C=O)cc(C)c1				
InChIKey	NBEFMISJJNGCIZ-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
ethanol	7.44	0.99	27092.92
methanol	7.44	0.99	28064.38
THF	7.44	0.99	23376.63

4-(trifluoromethyl)benzaldehyde $C_8H_5F_3O$[21]



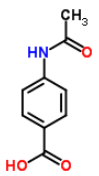
Compound Data

Molecular weight	174.12	**H bond acceptors**	1	**Rule of 5 violations**	0
Compound type	aldehyde	**H bond donors**	0	**ACD/ALogP**	2.61
Phase 25°C	liquid	**Rotatable bonds**	1	**Predicted density**	1.293 g/cm^3
SMILES	c1cc(C=O)ccc1C(F)(F)F				
InChIKey	BEOBZEOPTQQELP-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
methanol	7.32	0.96	11867.65

4-acetamidobenzoic acid $C_9H_9NO_3$[137]



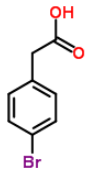
Compound Data

Molecular weight	179.173	**H bond acceptors**	4	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	2	**ACD/ALogP**	1.31
Phase 25°C	solid	**Rotatable bonds**	2	**Predicted density**	1.326 g/cm^3
SMILES	O=C(Nc1ccc(cc1)C(=O)O)C				
InChIKey	QCXJEYYXVJIFCE-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
acetonitrile	0.01	0.00	0.24
benzene	0.00	0.00	0.00
ethanol	0.23	0.01	5.45
methanol	0.19	0.01	4.64
THF	0.17	0.01	3.45

4-bromophenylacetic acid $C_8H_7BrO_2$[85]



Compound Data

Molecular weight	215.044	**H bond acceptors**	2	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	2.28
Phase 25°C	solid	**Rotatable bonds**	2	**Predicted density**	1.616 g/cm^3
SMILES	Brc1ccc(cc1)CC(=O)O				
InChIKey	QOWSWEBLNVACCL-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
THF	3.68	0.37	171.85

4-chloro-3-nitrobenzoic acid $C_7H_4ClNO_4$[905, 907]

Compound Data

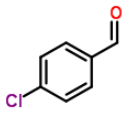
Molecular weight	201.564	**H bond acceptors**	5	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	2.37
Phase 25°C	solid	**Rotatable bonds**	2	**Predicted density**	1.602 g/cm^3
SMILES	O=[N+]([O-])c1cc(ccc1Cl)C(=O)O				
InChIKey	DFXQXFGFOLXAPO-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
1-butanol	0.31	0.03	8.08
1-decanol	0.19	0.04	4.74
1-heptanol	0.25	0.04	6.34
1-hexanol	0.27	0.03	6.90
1-octanol	0.23	0.04	5.80
1-pentanol	0.27	0.03	6.95
1-propanol	0.37	0.03	9.84
1,4-dioxane	1.43	0.13	35.32
2-butanol	0.32	0.03	8.39
2-ethyl-1-hexanol	0.15	0.02	3.75
2-methyl-1-pentanol	0.19	0.02	4.82
2-methyl-1-propanol	0.19	0.02	4.90
2-methyl-2-propanol	0.38	0.04	10.01
2-pentanol	0.30	0.03	7.77
2-propanol	0.39	0.03	10.45
3-methyl-1-butanol	0.24	0.03	6.17
4-methyl-2-pentanol	0.23	0.03	5.89
butyl acetate	0.23	0.03	5.39
dibutyl ether	0.05	0.01	1.30
diethyl ether	0.24	0.02	6.80
diisopropyl ether	0.10	0.01	2.69
ethanol	0.55	0.03	15.27
ethyl acetate	0.36	0.04	8.46
methyl acetate	0.44	0.04	10.34
methyl butyrate	0.21	0.02	4.88

pentyl acetate	0.16	0.02	3.73
propyl acetate	0.26	0.03	6.08
propylene carbonate	0.15	0.02	2.56
THF	1.95	0.17	57.61
water	0.00	0.00	0.00

4-chlorobenzaldehyde C_7H_5ClO[7, 209]



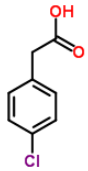
Compound Data

Molecular weight	140.567	**H bond acceptors**	1	**Rule of 5 violations**	0
Compound type	aldehyde	**H bond donors**	0	**ACD/ALogP**	2.21
Phase 25°C	solid	**Rotatable bonds**	1	**Predicted density**	1.243 g/cm^3
SMILES	O=Cc1ccc(Cl)cc1				
InChIKey	AVPYQKSLYISFPO-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
acetonitrile	5.58	0.45	285.27
chloroform	3.61	0.33	57.17
ethanol	2.81	0.20	74.23
methanol	3.55	0.20	110.72
THF	5.29	0.51	204.25

4-chlorophenylacetic acid $C_8H_7ClO_2$[4, 75, 73]



Compound Data

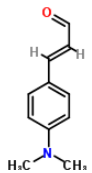
Molecular weight	170.593	**H bond acceptors**	2	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	2.1
Phase 25°C	solid	**Rotatable bonds**	2	**Predicted density**	1.324 g/cm^3
SMILES	Clc1ccc(cc1)CC(=O)O				
InChIKey	CDPKJZJVTHSESZ-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
acetonitrile	1.67	0.10	48.72
DMSO	5.74	0.61	342.14
methanol	7.76	¤	¤
THF	4.45	0.45	197.18
toluene	0.89	0.10	19.69

¤ Solute is very soluble/miscible, conversion fail.

4-dimethylamino cinnamaldehyde $C_{11}H_{13}NO$[3, 32]



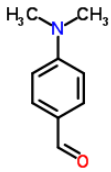
Compound Data

Molecular weight	175.227	**H bond acceptors**	2	**Rule of 5 violations**	0
Compound type	aldehyde	**H bond donors**	0	**ACD/ALogP**	2.63
Phase 25°C	solid	**Rotatable bonds**	3	**Predicted density**	1.057 g/cm^3
SMILES	O=CC=Cc1ccc(N(C)C)cc1				
InChIKey	RUKJCCIJLIMGEP-ONEGZZNKSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
ethanol	0.06	0.00	1.25
methanol	0.07	0.00	1.65
THF	0.42	0.03	8.75

4-dimethylaminobenzaldehyde $C_9H_{11}NO$[208, 205]



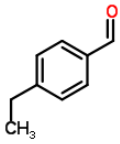
Compound Data

Molecular weight	149.19	**H bond acceptors**	2	**Rule of 5 violations**	0
Compound type	aldehyde	**H bond donors**	0	**ACD/ALogP**	1.81
Phase 25°C	solid	**Rotatable bonds**	2	**Predicted density**	1.072 g/cm^3
SMILES	O=Cc1ccc(N(C)C)cc1				
InChIKey	BGNGWHSBYQYVRX-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
acetonitrile	2.60	0.18	81.37
chloroform	4.02	0.42	90.76
ethanol	0.63	0.04	13.21
methanol	1.36	0.07	33.24
THF	2.50	0.23	63.27
toluene	1.54	0.17	33.57

4-ethylbenzaldehyde $C_9H_{10}O$[63]



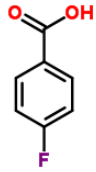
Compound Data

Molecular weight	134.175	**H bond acceptors**	1	**Rule of 5 violations**	0
Compound type	aldehyde	**H bond donors**	0	**ACD/ALogP**	2.63
Phase 25°C	liquid	**Rotatable bonds**	2	**Predicted density**	1.001 g/cm^3
SMILES	CCc1ccc(C=O)cc1				
InChIKey	QNGNSVIICDLXHT-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
ethanol	7.30	0.95	5840.38
methanol	7.30	0.94	6049.80
THF	7.30	0.96	5039.27

4-fluorobenzoic acid $C_7H_5FO_2$[106, 121, 907]



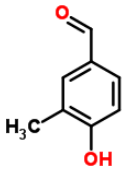
Compound Data

Molecular weight	140.112	**H bond acceptors**	2	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	2.07
Phase 25°C	solid	**Rotatable bonds**	1	**Predicted density**	1.319 g/cm^3
SMILES	O=C(O)c1ccc(F)cc1				
InChIKey	BBYDXOIZLAWGSL-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
acetonitrile	0.39	0.02	7.63
DMSO	4.19	0.35	96.26
ethanol	0.80	0.05	15.71
methanol	0.76	0.03	15.38
THF	2.49	0.21	52.47
toluene	0.00	0.00	0.00
water	0.01	0.00	0.14

4-hydroxy-3-methylbenzaldehyde $C_8H_8O_2$[57]



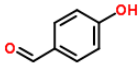
Compound Data

Molecular weight	136.148	**H bond acceptors**	2	**Rule of 5 violations**	0
Compound type	aldehyde	**H bond donors**	1	**ACD/ALogP**	1.85
Phase 25°C	solid	**Rotatable bonds**	2	**Predicted density**	1.175 g/cm^3
SMILES	O=Cc1ccc(O)c(c1)C				
InChIKey	BAKYASSDAXQKKY-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
ethanol	2.15	0.14	49.98
methanol	2.78	0.15	74.15
THF	3.05	0.27	71.04

4-hydroxybenzaldehyde $C_7H_6O_2$[208, 57, 205, 58, 88, 85, 909]



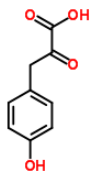
Compound Data

Molecular weight	122.121	**H bond acceptors**	2	**Rule of 5 violations**	0
Compound type	aldehyde	**H bond donors**	1	**ACD/ALogP**	1.39
Phase 25°C	solid	**Rotatable bonds**	2	**Predicted density**	1.226 g/cm^3
SMILES	O=Cc1ccc(O)cc1				
InChIKey	RGHHSNMVTDWUBI-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
acetonitrile	1.46	0.09	27.93
chloroform	1.75	0.14	17.26
ethanol	2.52	0.17	52.82
methanol	3.39	0.18	83.19
THF	3.70	0.32	79.10
toluene	0.02	0.00	0.28
water	0.11	0.00	1.36

4-hydroxyphenylpyruvic acid $C_9H_8O_4$[15]



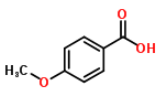
Compound Data

Molecular weight	180.157	**H bond acceptors**	4	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	2	**ACD/ALogP**	-0.2
Phase 25°C	solid	**Rotatable bonds**	4	**Predicted density**	1.397 g/cm^3
SMILES	c1cc(O)ccc1CC(=O)C(=O)O				
InChIKey	KKADPXVIOXHVKN-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
ethanol	0.71	0.04	18.05
methanol	0.91	0.04	24.67

4-methoxybenzoic acid $C_8H_8O_3$[905]



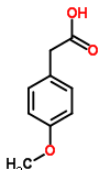
Compound Data

Molecular weight	152.147	**H bond acceptors**	3	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	1.96
Phase 25°C	solid	**Rotatable bonds**	2	**Predicted density**	1.207 g/cm^3
SMILES	COc1ccc(cc1)C(O)=O				
InChIKey	ZEYHEAKUIGZSGI-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
1-butanol	0.12	0.01	2.30
1-decanol	0.07	0.01	1.30
1-heptanol	0.09	0.01	1.69
1-hexanol	0.10	0.01	1.89
1-octanol	0.08	0.01	1.49
1-pentanol	0.10	0.01	1.90
1-propanol	0.13	0.01	2.53
1,4-dioxane	0.39	0.04	6.27
2-butanol	0.13	0.01	2.51
2-ethyl-1-hexanol	0.07	0.01	1.31
2-methyl-1-propanol	0.08	0.01	1.54
2-methyl-2-propanol	0.17	0.02	3.29
2-propanol	0.14	0.01	2.74
3-methyl-1-butanol	0.08	0.01	1.52
butyl acetate	0.08	0.01	1.39
dibutyl ether	0.02	0.00	0.39
diethyl ether	0.09	0.01	1.89
diisopropyl ether	0.03	0.00	0.60
ethanol	0.20	0.01	4.00
ethyl acetate	0.13	0.01	2.24
pentyl acetate	0.06	0.01	1.04
THF	0.70	0.06	12.92

4-methoxyphenylacetic acid $C_9H_{10}O_3$[89, 136, 92, 910]



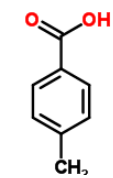
Compound Data

Molecular weight	166.174	**H bond acceptors**	3	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	1.42
Phase 25°C	solid	**Rotatable bonds**	3	**Predicted density**	1.179 g/cm^3
SMILES	O=C(O)Cc1ccc(OC)cc1				
InChIKey	NRPFNQUDKRYCNX-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
acetonitrile	2.23	0.15	72.11
DMSO	4.50	0.47	186.41
methanol	2.50	0.14	85.19
THF	3.31	0.33	114.38
toluene	0.37	0.04	7.45
water	0.04	0.00	0.67

4-methylbenzoic acid $C_8H_8O_2$[2, 25]



Compound Data

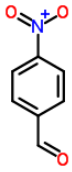
Molecular weight	136.148	**H bond acceptors**	2	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	2.36
Phase 25°C	solid	**Rotatable bonds**	1	**Predicted density**	1.151 g/cm^3
SMILES	O=C(O)c1ccc(cc1)C				
InChIKey	LPNBBFKOUUSUDB-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
ethanol	0.78	0.05	14.89
methanol	0.78	0.04	15.54

THF	1.60	0.14	29.84

4-nitrobenzaldehyde $C_7H_5NO_3$[904, 111, 98, 212, 33, 205, 82, 122]



Compound Data

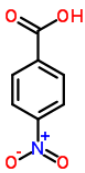
Molecular weight	151.12	**H bond acceptors**	4	**Rule of 5 violations**	0
Compound type	aldehyde	**H bond donors**	0	**ACD/ALogP**	1.56
Phase 25°C	solid	**Rotatable bonds**	2	**Predicted density**	1.338 g/cm^3
SMILES	c1cc(C=O)ccc1[N+](=O)[O-]				
InChIKey	BXRFQSNOROATLV-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
1-butanol	*	*	*
1-propanol	*	*	*
2-propanol	*	*	*
acetonitrile	1.13	0.07	26.16
benzene	0.46	0.04	8.40
carbon tetrachloride	0.07	0.01	0.63
chloroform	1.02	0.08	11.57
cyclohexane	0.01	0.00	0.19
dichloromethane	1.29	0.09	18.23
DMF	1.68	0.15	36.02
DMSO	1.24	0.09	19.83
ethanol	*	*	*
ethylene glycol	*	*	*
methanol	*	*	*
THF	0.88	0.07	16.33
toluene	0.46	0.05	8.42

* This aldehyde reacts with alcohols to form a hemiacetal.

4-nitrobenzoic acid $C_7H_5NO_4$[905]



Compound Data

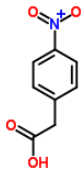
Molecular weight	167.119	**H bond acceptors**	5	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	1.89
Phase 25°C	solid	**Rotatable bonds**	2	**Predicted density**	1.468 g/cm^3
SMILES	O=[N+]([O-])c1ccc(C(=O)O)cc1				
InChIKey	OTLNPYWUJOZPPA-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
1-butanol	0.09	0.01	1.89
1-decanol	0.05	0.01	1.01
1-heptanol	0.08	0.01	1.65
1-hexanol	0.08	0.01	1.65
1-octanol	0.06	0.01	1.23
1-pentanol	0.09	0.01	1.87
1-propanol	0.11	0.01	2.34
1,4-dioxane	0.47	0.04	8.34
2-butanol	0.10	0.01	2.11
2-methyl-1-butanol	0.05	0.01	1.04
2-methyl-1-pentanol	0.05	0.01	1.03
2-methyl-1-propanol	0.07	0.01	1.47
2-methyl-2-butanol	0.17	0.02	3.57
2-methyl-2-propanol	0.15	0.01	3.17
2-pentanol	0.08	0.01	1.67
2-propanol	0.11	0.01	2.35

3-methyl-1-butanol	0.07	0.01	1.46
4-methyl-2-pentanol	0.07	0.01	1.45
butyl acetate	0.08	0.01	1.52
dibutyl ether	0.02	0.00	0.43
diethyl ether	0.10	0.01	2.30
diisopropyl ether	0.03	0.00	0.66
ethanol	0.14	0.01	3.05
ethyl acetate	0.13	0.01	2.46
methanol	0.18	0.01	4.08
methyl acetate	0.14	0.01	2.62
pentyl acetate	0.05	0.01	0.95
propylene carbonate	0.06	0.01	0.84
THF	0.78	0.06	15.82

4-nitrophenylacetic acid $C_8H_7NO_4$[15]



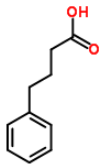
Compound Data

Molecular weight	181.145	**H bond acceptors**	5	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	1.24
Phase 25°C	solid	**Rotatable bonds**	3	**Predicted density**	1.407 g/cm^3
SMILES	c1cc(CC(=O)O)ccc1[N+](=O)[O-]				
InChIKey	YBADLXQNJCMBKR-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
ethanol	0.71	0.04	18.15
methanol	1.29	0.06	37.21

4-phenylbutyric acid $C_{10}H_{12}O_2$[72]



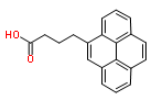
Compound Data

Molecular weight	164.201	**H bond acceptors**	2	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	2.42
Phase 25°C	solid	**Rotatable bonds**	4	**Predicted density**	1.095 g/cm^3
SMILES	O=C(O)CCCc1ccccc1				
InChIKey	OBKXEAXTFZPCHS-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
1-hexanol	0.19	0.02	3.94
chloroform	3.99	0.44	108.74
diethyl ether	4.02	0.51	226.42
methanol	4.54	0.38	310.15
toluene	4.06	0.52	195.66

4-pyrenebutanoic acid $C_{20}H_{16}O_2$[61, 99, 51, 30, 24, 91, 84]



Compound Data

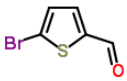
Molecular weight	288.34	**H bond acceptors**	2	**Rule of 5 violations**	1
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	5.37
Phase 25°C	solid	**Rotatable bonds**	4	**Predicted density**	1.296 g/cm^3
SMILES	O=C(O)CCCc4cc2cccc1ccc3c(c12)c4ccc3				
InChIKey	GSBSDUZVSCTUKA-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration	Mole Fraction (X)	pph (g/100g)

	(M)		
2-propanol	0.06	0.00	2.22
acetonitrile	0.00	0.00	0.00
benzene	0.00	0.00	0.00
carbon tetrachloride	0.00	0.00	0.00
chloroform	0.03	0.00	0.58
cyclohexane	0.00	0.00	0.00
cyclopentane	0.00	0.00	0.00
dichloromethane	0.07	0.00	1.64
diethyl ether	0.02	0.00	0.79
DMF	1.88	0.21	107.11
DMSO	2.07	0.21	100.37
ethanol	0.06	0.00	2.25
hexane	0.00	0.00	0.00
methanol	0.01	0.00	0.51
THF	0.55	0.05	19.99
toluene	0.00	0.00	0.00

5-bromothiophene-2-carbaldehyde C_5H_3BrOS[21]



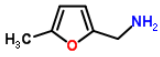
Compound Data

Molecular weight	191.046	**H bond acceptors**	1	**Rule of 5 violations**	0
Compound type	aldehyde	**H bond donors**	0	**ACD/ALogP**	1.97
Phase 25°C	liquid	**Rotatable bonds**	1	**Predicted density**	1.789 g/cm^3
SMILES	C1=C(SC(=C1)Br)C=O				
InChIKey	GFBVUFQNHLUCPX-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
methanol	9.22	0.96	15186.93

5-methylfurfurylamine C_6H_9NO[26]



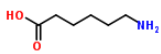
Compound Data

Molecular weight	111.142	**H bond acceptors**	2	**Rule of 5 violations**	0
Compound type	amine	**H bond donors**	2	**ACD/ALogP**	0.71
Phase 25°C	liquid	**Rotatable bonds**	2	**Predicted density**	1.024 g/cm^3
SMILES	o1c(ccc1CN)C				
InChIKey	YSEAGSCGERFGBL-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
methanol	8.97	0.94	5010.81

6-aminocaproic acid $C_6H_{13}NO_2$[14]



Compound Data

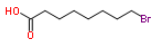
Molecular weight	131.173	**H bond acceptors**	3	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	3	**ACD/ALogP**	-0.11
Phase 25°C	solid	**Rotatable bonds**	6	**Predicted density**	1.042 g/cm^3
SMILES	C(CCC(=O)O)CCN				
InChIKey	SLXKOJJOQWFEFD-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)

ethanol	0.00	0.00	0.00
methanol	0.06	0.00	1.05

8-bromooctanoic acid $C_8H_{15}BrO_2$[3, 66]



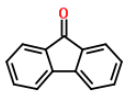
Compound Data

Molecular weight	223.107	**H bond acceptors**	2	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	2.76
Phase 25°C	solid	**Rotatable bonds**	7	**Predicted density**	1.324 g/cm^3
SMILES	BrCCCCCCCC(=O)O				
InChIKey	BKJFDZSBZWHRNH-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
ethanol	4.07	0.43	370.56
methanol	3.28	0.24	217.27
THF	3.56	0.42	219.59

9-fluorenone $C_{13}H_8O$[905]



Compound Data

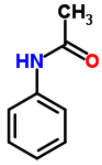
Molecular weight	180.202	**H bond acceptors**	1	**Rule of 5 violations**	0
Compound type	non-Ugi related	**H bond donors**	0	**ACD/ALogP**	3.58
Phase 25°C	solid	**Rotatable bonds**	0	**Predicted density**	1.244 g/cm^3
SMILES	O=C3c1ccccc1c2c3cccc2				
InChIKey	YLQWCDOCJODRMT-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
1-butanol	0.37	0.03	8.75
1-decanol	0.42	0.08	9.73
1-heptanol	0.43	0.06	10.08
1-hexanol	0.41	0.05	9.63
1-octanol	0.43	0.07	10.04
1-pentanol	0.41	0.05	9.69
1-propanol	0.38	0.03	9.12
2-butanol	0.33	0.03	7.80
2-ethyl-1-hexanol	0.35	0.06	8.09
2-methyl-1-pentanol	0.36	0.05	8.41
2-methyl-1-propanol	0.25	0.02	5.84
2-methyl-2-butanol	0.40	0.04	9.43
2-methyl-2-propanol	0.31	0.03	7.27
2-pentanol	0.35	0.04	8.21
2-propanol	0.29	0.02	6.90
2,2,4-trimethylpentane	0.07	0.01	1.80
3-methyl-1-butanol	0.31	0.03	7.23
4-methyl-2-pentanol	0.29	0.04	6.73
acetonitrile	1.57	0.10	49.02
benzene	2.61	0.27	86.63
butyronitrile	1.93	0.19	61.50
carbon tetrachloride	1.83	0.18	26.44
cyclohexane	0.15	0.02	3.50
cyclooctane	0.17	0.02	3.98
cyclopentanol	0.82	0.07	16.70
decane	0.12	0.02	3.00
dibutyl ether	0.37	0.06	9.03
dichloromethane	4.03	0.40	139.36
diethyl ether	0.88	0.09	24.76

diisopropyl ether	0.37	0.05	9.29
ethanol	3.08	0.25	128.48
heptane	0.11	0.02	2.90
hexadecane	0.12	0.03	2.85
hexane	0.10	0.01	2.71
methanol	0.37	0.02	9.36
methyl tert-butyl ether	0.76	0.09	20.52
methylcyclohexane	0.14	0.02	3.32
nonane	0.12	0.02	3.04
octane	0.11	0.02	2.83
propionitrile	1.99	0.17	65.35
tert-butylcyclohexane	0.12	0.02	2.71
toluene	2.17	0.25	65.48
undecane	0.12	0.03	2.96

acetanilide C_8H_9NO[35]



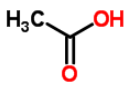
Compound Data

Molecular weight	135.163	**H bond acceptors**	2	**Rule of 5 violations**	0
Compound type	amide	**H bond donors**	1	**ACD/ALogP**	1.08
Phase 25°C	solid	**Rotatable bonds**	1	**Predicted density**	1.103 g/cm^3
SMILES	O=C(Nc1ccccc1)C				
InChIKey	FZERHIULMFGESH-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
acetone	1.42	0.11	30.10

acetic acid $C_2H_4O_2$[22]



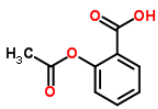
Compound Data

Molecular weight	60.052	**H bond acceptors**	2	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	-0.29
Phase 25°C	liquid	**Rotatable bonds**	0	**Predicted density**	1.068 g/cm^3
SMILES	O=C(O)C				
InChIKey	QTBSBXVTEAMEQO-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
methanol	17.47	0.98	7876.42

acetylsalicylic acid $C_9H_8O_4$[72, 34, 903, 907]



Compound Data

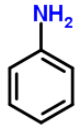
Molecular weight	180.157	**H bond acceptors**	4	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	1.19
Phase 25°C	solid	**Rotatable bonds**	3	**Predicted density**	1.29 g/cm^3
SMILES	O=C(Oc1ccccc1C(=O)O)C				
InChIKey	BSYNRYMUTXBXSQ-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
1-hexanol	0.00	0.00	0.00
chloroform	0.18	0.01	2.24
diethyl ether	0.17	0.02	4.27

methanol	1.30	0.06	38.00
toluene	0.00	0.00	0.00
water	0.03	0.00	0.45

aniline C_6H_7N[29]



Compound Data

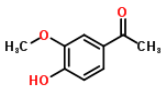
Molecular weight	93.1265	**H bond acceptors**	1	**Rule of 5 violations**	0
Compound type	amine	**H bond donors**	2	**ACD/ALogP**	0.94
Phase 25°C	liquid	**Rotatable bonds**	1	**Predicted density**	1.015 g/cm^3
SMILES	Nc1ccccc1				
InChIKey	PAYRUJLWNCNPSJ-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
methanol	10.97	¤	¤

¤ Solute is very soluble/miscible, conversion fail.

apocynin $C_9H_{10}O_3$[906]



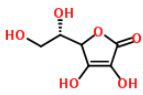
Compound Data

Molecular weight	166.174	**H bond acceptors**	3	**Rule of 5 violations**	0
Compound type	non-Ugi related	**H bond donors**	1	**ACD/ALogP**	1.39
Phase 25°C	solid	**Rotatable bonds**	3	**Predicted density**	1.158 g/cm^3
SMILES	Oc1ccc(cc1OC)C(C)=O				
InChIKey	DFYRUELUNQRZTB-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
1-propanol	0.32	0.02	7.01
1,2-dichloroethane	0.41	0.04	6.17
butanone	0.66	0.06	15.41
toluene	0.06	0.01	1.15

ascorbic acid $C_9H_8O_4$[72]



Compound Data

Molecular weight	180.157	**H bond acceptors**	4	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	1.19
Phase 25°C	solid	**Rotatable bonds**	3	**Predicted density**	1.29 g/cm^3
SMILES	O\C1=C(/O)C(=O)O[C@@H]1[C@@H](O)CO				
InChIKey	CIWBSHSKHKDKBQ-JLAZNSOCSA-N				

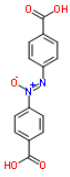
Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
1-hexanol	0.00	0.00	0.00
chloroform	0.00	0.00	0.00
diethyl ether	0.00	0.00	0.00
toluene	0.00	0.00	0.00

azoxydibenzoic acid $C_{14}H_{10}N_2O_5$[72]

Compound Data

Molecular weight	286.24	**H bond acceptors**	7	**Rule of 5 violations**	0

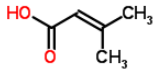


Compound type	carboxylic acid	**H bond donors**	2	**ACD/ALogP**	4.05
Phase 25°C	solid	**Rotatable bonds**	4	**Predicted density**	1.41 g/cm^3
SMILES	O=C(O)c2ccc([N+](/[O-])=N/c1ccc(cc1)C(=O)O)cc2				
InChIKey	ZYVHVGJGLOEEKD-NXVVXOECSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
1-hexanol	0.00	0.00	0.00
chloroform	0.00	0.00	0.00
diethyl ether	0.00	0.00	0.00
toluene	0.00	0.00	0.00

B,B-dimethylacrylic acid $C_5H_8O_2$[136]



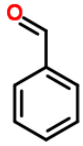
Compound Data

Molecular weight	100.116	**H bond acceptors**	2	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	1.35
Phase 25°C	solid	**Rotatable bonds**	1	**Predicted density**	1.01 g/cm^3
SMILES	O=C(O)\C=C(/C)C				
InChIKey	YYPNJNDODFVZLE-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
methanol	4.73	0.27	118.40

benzaldehyde C_7H_6O[21]



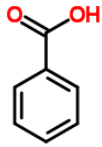
Compound Data

Molecular weight	106.122	**H bond acceptors**	1	**Rule of 5 violations**	0
Compound type	aldehyde	**H bond donors**	0	**ACD/ALogP**	1.64
Phase 25°C	liquid	**Rotatable bonds**	1	**Predicted density**	1.049 g/cm^3
SMILES	c1ccccc1C=O				
InChIKey	HUMNYLRZRPPJDN-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
methanol	9.85	0.99	39374.95

benzoic acid $C_7H_6O_2$[155, 37, 902, 58, 160, 43, 136, 9, 907, 900]



Compound Data

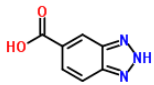
Molecular weight	122.121	**H bond acceptors**	2	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	1.9
Phase 25°C	solid	**Rotatable bonds**	1	**Predicted density**	1.197 g/cm^3
SMILES	O=C(O)c1ccccc1				
InChIKey	WPYMKLBDIGXBTP-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
1-octanol	1.48	0.22	25.97
acetone	1.35	0.11	24.77

acetonitrile	0.75	0.04	13.37
benzene	0.48	0.04	7.06
chloroform	1.80	0.15	17.95
ethanol	2.55	0.17	54.02
methanol	2.84	0.15	64.81
THF	3.37	0.29	69.27
toluene	0.65	0.07	9.81
water	0.03	0.00	0.36

benzotriazole-5-carboxylic acid $C_7H_5N_3O_2$[3, 32]



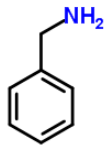
Compound Data

Molecular weight	163.133	**H bond acceptors**	5	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	2	**ACD/ALogP**	0.82
Phase 25°C	solid	**Rotatable bonds**	1	**Predicted density**	1.617 g/cm^3
SMILES	O=C(O)c1ccc2nnnc2c1				
InChIKey					

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
ethanol	0.04	0.00	0.73
methanol	0.04	0.00	0.76
THF	0.02	0.00	0.36

benzylamine C_7H_9N[29]



Compound Data

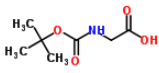
Molecular weight	107.153	**H bond acceptors**	1	**Rule of 5 violations**	0
Compound type	amine	**H bond donors**	2	**ACD/ALogP**	1.09
Phase 25°C	liquid	**Rotatable bonds**	2	**Predicted density**	0.979 g/cm^3
SMILES	NCc1ccccc1				
InChIKey	WGQKYBSKWIADBV-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
methanol	9.16	¤	¤

¤ Solute is very soluble/miscible, conversion fail.

boc-glycine $C_7H_{13}NO_4$[143, 71, 207, 135, 133, 116]



Compound Data

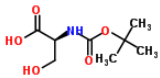
Molecular weight	175.182	**H bond acceptors**	5	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	2	**ACD/ALogP**	0.75
Phase 25°C	solid	**Rotatable bonds**	4	**Predicted density**	1.159 g/cm^3
SMILES	O=C(OC(C)(C)C)NCC(=O)O				
InChIKey	VRPJIFMKZZEXLR-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
2-propanol	3.63	0.38	177.58
acetonitrile	2.46	0.18	91.84
benzene	0.25	0.02	5.32
chloroform	3.85	0.42	107.55
dichloromethane	4.14	0.43	154.79
diethyl ether	2.65	0.31	105.51

DMSO	4.78	0.55	274.57
ethanol	4.11	0.39	243.70
hexane	0.25	0.03	6.74
methanol	4.72	0.41	384.60
THF	3.72	0.40	164.41
toluene	0.19	0.02	3.93

boc-serine $C_8H_{15}NO_5$[98, 85]



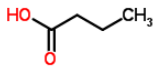
Compound Data

Molecular weight	205.208	**H bond acceptors**	6	**Rule of 5 violations**	0
Compound type	boc serine	**H bond donors**	3	**ACD/ALogP**	0.22
Phase 25°C	solid	**Rotatable bonds**	6	**Predicted density**	1.24 g/cm^3
SMILES	O=C(OC(C)(C)C)N[C@H](C(=O)O)CO				
InChIKey	FHOAKXBXYSJBGX-YFKPBYRVSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
methanol	3.62	0.28	246.06
THF	3.11	0.34	145.78

butyric acid $C_4H_8O_2$[22]



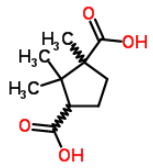
Compound Data

Molecular weight	88.1051	**H bond acceptors**	2	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	0.78
Phase 25°C	liquid	**Rotatable bonds**	2	**Predicted density**	0.987 g/cm^3
SMILES	CCCC(=O)O				
InChIKey	FERIUCNNQQJTOY-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
methanol	10.94	0.95	5462.12

camphoric acid $C_{10}H_{16}O_4$[902]



Compound Data

Molecular weight	200.232	**H bond acceptors**	4	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	2	**ACD/ALogP**	1.47
Phase 25°C	solid	**Rotatable bonds**	2	**Predicted density**	1.177 g/cm^3
SMILES	O=C(O)C1CCC(C(=O)O)(C)C1(C)C				
InChIKey	LSPHULWDVZXLIL-UHFFFAOYSA-N				

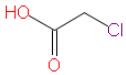
Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
1-propanol	1.72	0.16	61.24
methanol	2.75	0.18	137.41

chloroacetic acid $C_2H_3ClO_2$[95]

Compound Data

Molecular weight	94.497	**H bond acceptors**	2	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	0.18
Phase 25°C	solid	**Rotatable bonds**	1	**Predicted density**	1.398 g/cm^3

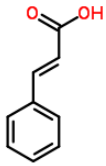


SMILES	ClCC(O)=O
InChIKey	FOCAUTSVDIKZOP-UHFFFAOYSA-N

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
acetonitrile	12.45	0.81	993.98
DMSO	11.20	0.77	396.40
methanol	9.45	0.53	328.30
THF	9.63	0.69	288.10
toluene	1.41	0.14	16.95

cinnamic acid $C_9H_8O_2$[72, 902, 48, 907]



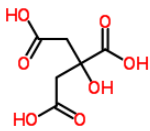
Compound Data

Molecular weight	148.159	**H bond acceptors**	2	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	2.38
Phase 25°C	solid	**Rotatable bonds**	2	**Predicted density**	1.184 g/cm^3
SMILES	O=C(O)\C=C\c1ccccc1				
InChIKey	WBYWAXJHAXSJNI-VOTSOKGWSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
1-hexanol	0.00	0.00	0.00
chloroform	0.93	0.08	10.40
diethyl ether	0.43	0.04	9.17
ethanol	0.86	0.05	18.31
methanol	1.10	0.05	25.19
toluene	0.30	0.03	5.30
water	0.00	0.00	0.00

citric acid anhydrous $C_6H_8O_7$[902, 25]



Compound Data

Molecular weight	192.124	**H bond acceptors**	7	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	4	**ACD/ALogP**	-1.33
Phase 25°C	solid	**Rotatable bonds**	6	**Predicted density**	1.762 g/cm^3
SMILES	O=C(O)CC(O)(C(=O)O)CC(=O)O				
InChIKey	KRKNYBCHXYNGOX-UHFFFAOYSA-N				

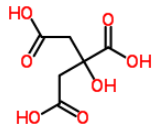
Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
chloroform	0.00	0.00	0.00
ethanol	1.60	0.10	47.74
methanol	3.08	0.16	118.32
THF	1.80	0.15	47.60

citric acid monohydrate $C_6H_8O_7$[2]

Compound Data

Molecular weight	192.124	**H bond acceptors**	7	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	4	**ACD/ALogP**	-1.33



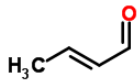
Phase 25°C	solid	**Rotatable bonds**	6	**Predicted density**	1.762 g/cm^3
SMILES	O=C(O)CC(O)(C(=O)O)CC(=O)O				
InChIKey	KRKNYBCHXYNGOX-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
ethanol	1.78	0.12	54.40
methanol	2.27	0.11	76.97
THF	1.52	0.13	38.72

crotonaldehyde C_4H_6O[21]

Compound Data



Molecular weight	70.0898	**H bond acceptors**	1	**Rule of 5 violations**	0
Compound type	aldehyde	**H bond donors**	0	**ACD/ALogP**	0.88
Phase 25°C	solid	**Rotatable bonds**	1	**Predicted density**	0.819 g/cm^3
SMILES	CC=CC=O				
InChIKey	MLUCVPSAIODCQM-NSCUHMNNSA-N				

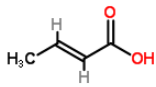
Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
methanol	12.17	¤	¤

¤ Solute is very soluble/miscible, conversion fail.

crotonic acid $C_4H_6O_2$[64, 62, 907]

Compound Data



Molecular weight	86.0892	**H bond acceptors**	2	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	0.94
Phase 25°C	solid	**Rotatable bonds**	1	**Predicted density**	1.039 g/cm^3
SMILES	O=C(O)/C=C/C				
InChIKey	LDHQCZJRKDOVOX-NSCUHMNNSA-N				

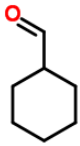
Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
acetonitrile	3.59	0.22	58.89
benzene	4.45	0.39	69.51
chloroform	7.34	0.60	107.51
DMSO	7.76	0.61	170.26
ethanol	6.15	0.43	138.41
hexane	1.07	0.13	14.97
methanol	7.62	0.47	236.33
THF	8.09	0.66	233.69
toluene	3.52	0.34	49.12
water	0.89	0.02	8.29

cyclohexanecarbaldehyde $C_7H_{12}O$[21]

Compound Data

Molecular weight	112.17	**H bond acceptors**	1	**Rule of 5 violations**	0
Compound type	aldehyde	**H bond donors**	0	**ACD/ALogP**	2.02

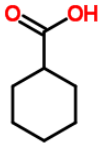


Phase 25°C	liquid	**Rotatable bonds**	1	**Predicted density**	0.992 g/cm
SMILES	C1CCC(CC1)C=O				
InChIKey	KVFDZFBHBWTVID-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
methanol	8.29	0.85	1972.34

cyclohexanecarboxylic acid $C_7H_{12}O_2$[22]



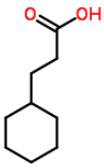
Compound Data

Molecular weight	128.169	**H bond acceptors**	2	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	1.88
Phase 25°C	liquid	**Rotatable bonds**	1	**Predicted density**	1.079 g/cm^3
SMILES	C1CCC(CC1)C(=O)O				
InChIKey	NZNMSOFKMUBTKW-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
methanol	8.06	0.89	3220.96

cyclohexanepropionic acid $C_9H_{16}O_2$[22]



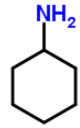
Compound Data

Molecular weight	156.222	**H bond acceptors**	2	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	3.09
Phase 25°C	liquid	**Rotatable bonds**	3	**Predicted density**	1.006 g/cm^3
SMILES	C1CCC(CC1)CCC(=O)O				
InChIKey	HJZLEGIHUQOJBA-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
methanol	5.84	0.73	1301.33

cyclohexylamine $C_6H_{13}N$[29]



Compound Data

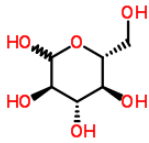
Molecular weight	99.1741	**H bond acceptors**	1	**Rule of 5 violations**	0
Compound type	amine	**H bond donors**	2	**ACD/ALogP**	1.3
Phase 25°C	liquid	**Rotatable bonds**	1	**Predicted density**	0.869 g/cm^3
SMILES	NC1CCCCC1				
InChIKey	PAFZNILMFXTMIY-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
methanol	8.74	0.99	45092.19

D-glucose $C_6H_{12}O_6$[1, 207]

Compound Data

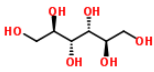


Molecular weight	180.156	**H bond acceptors**	6	**Rule of 5 violations**	1
Compound type	non-Ugi related	**H bond donors**	5	**ACD/ALogP**	-2.57
Phase 25°C	solid	**Rotatable bonds**	6	**Predicted density**	1.732 g/cm^3
SMILES	O[C@H]1[C@H](O)[C@H](OC(O)[C@@H]1O)CO				
InChIKey	WQZGKKKJIJFFOK-GASJEMHNSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
ethanol	0.01	0.00	0.12
methanol	0.04	0.00	0.84
THF	0.02	0.00	0.30

D-mannitol $C_6H_{14}O_6$[1, 207]



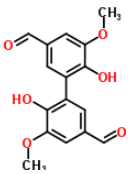
Compound Data

Molecular weight	182.172	**H bond acceptors**	6	**Rule of 5 violations**	1
Compound type	non-Ugi related	**H bond donors**	6	**ACD/ALogP**	-2.68
Phase 25°C	solid	**Rotatable bonds**	11	**Predicted density**	1.596 g/cm^3
SMILES	O[C@H]([C@H](O)CO)[C@H](O)[C@H](O)CO				
InChIKey	FBPFZTCFMRRESA-KVTDHHQDSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
ethanol	0.01	0.00	0.12
methanol	0.01	0.00	0.24
THF	0.02	0.00	0.30

dehydrodivanillin $C_{16}H_{14}O_6$[906]



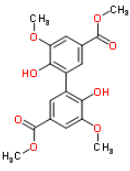
Compound Data

Molecular weight	302.279	**H bond acceptors**	6	**Rule of 5 violations**	0
Compound type	non-Ugi related	**H bond donors**	2	**ACD/ALogP**	2.7
Phase 25°C	solid	**Rotatable bonds**	7	**Predicted density**	1.356 g/cm^3
SMILES	O=Cc2cc(c1cc(cc(OC)c1O)C=O)c(O)c(OC)c2				
InChIKey	NSTQUZVZBUTVPY-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
1-propanol	0.00	0.00	0.00
1,2-dichloroethane	0.00	0.00	0.00
butanone	0.00	0.00	0.00
toluene	0.00	0.00	0.00

di(methyl vanillate) $C_{18}H_{18}O_8$[906]



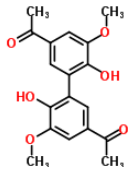
Compound Data

Molecular weight	362.331	**H bond acceptors**	8	**Rule of 5 violations**	0
Compound type	non-Ugi related	**H bond donors**	2	**ACD/ALogP**	2.28
Phase 25°C	solid	**Rotatable bonds**	9	**Predicted density**	1.315 g/cm^3
SMILES	O=C(OC)c2cc(c1cc(cc(OC)c1O)C(=O)OC)c(O)c(OC)c2				
InChIKey	DBUKZZKLTUXIQM-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
1-propanol	0.00	0.00	0.00
1,2-dichloroethane	0.01	0.00	0.31
butanone	0.00	0.00	0.00
toluene	0.00	0.00	0.00

diapocynin $C_{18}H_{18}O_6$[906]



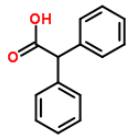
Compound Data

Molecular weight	330.332	**H bond acceptors**	6	**Rule of 5 violations**	0
Compound type	non-Ugi related	**H bond donors**	2	**ACD/ALogP**	2.84
Phase 25°C	solid	**Rotatable bonds**	7	**Predicted density**	1.257 g/cm^3
SMILES	O=C(c2cc(c1cc(cc(OC)c1O)C(=O)C)c(O)c(OC)c2)C				
InChIKey	HLNDPICGHQGWSU-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
1-propanol	0.00	0.00	0.00
1,2-dichloroethane	0.00	0.00	0.00
butanone	0.00	0.00	0.00
toluene	0.00	0.00	0.00

diphenylacetic acid $C_{14}H_{12}O_2$[81, 77, 907]



Compound Data

Molecular weight	212.244	**H bond acceptors**	2	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	2.79
Phase 25°C	solid	**Rotatable bonds**	3	**Predicted density**	1.174 g/cm^3
SMILES	O=C(O)C(c1ccccc1)c2ccccc2				
InChIKey	PYHXGXCGESYPCW-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
acetonitrile	0.46	0.03	14.26
DMSO	2.56	0.25	92.04
ethanol	1.15	0.08	39.51
methanol	1.27	0.07	46.47
THF	2.91	0.33	144.17
toluene	0.17	0.02	4.27
water	0.00	0.00	0.00

formic acid CH_2O_2[22]

Compound Data

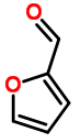
Molecular weight	46.0254	**H bond acceptors**	2	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	-0.47
Phase 25°C	liquid	**Rotatable bonds**	0	**Predicted density**	1.154 g/cm^3
SMILES	O=CO				
InChIKey	BDAGIHXWWSANSR-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
methanol	26.50	¤	¤

¤ Solute is very soluble/miscible, conversion fail.

furfuraldehyde $C_5H_4O_2$[20]



Compound Data

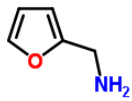
Molecular weight	96.0841	**H bond acceptors**	2	**Rule of 5 violations**	0
Compound type	aldehyde	**H bond donors**	0	**ACD/ALogP**	0.43
Phase 25°C	liquid	**Rotatable bonds**	1	**Predicted density**	1.145 g/cm^3
SMILES	O=Cc1ccco1				
InChIKey	HYBBIBNJHNGZAN-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
methanol	12.02	¤	¤

¤ Solute is very soluble/miscible, conversion fail.

furfurylamine C_5H_7NO[26]



Compound Data

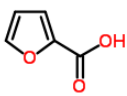
Molecular weight	97.1152	**H bond acceptors**	2	**Rule of 5 violations**	0
Compound type	amine	**H bond donors**	2	**ACD/ALogP**	-0.43
Phase 25°C	liquid	**Rotatable bonds**	2	**Predicted density**	1.053 g/cm^3
SMILES	o1c(ccc1)CN				
InChIKey	DDRPCXLAQZKBJP-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
methanol	11.32	¤	¤

¤ Solute is very soluble/miscible, conversion fail.

furoic acid $C_5H_4O_3$[104, 911]



Compound Data

Molecular weight	112.084	**H bond acceptors**	3	**Rule of 5 violations**	0
Compound type	carboxlyic acid	**H bond donors**	1	**ACD/ALogP**	0.98
Phase 25°C	solid	**Rotatable bonds**	1	**Predicted density**	1.322 g/cm^3
SMILES	OC(=O)c1ccco1				
InChIKey	SMNDYUVBFMFKNZ-UHFFFAOYSA-N				

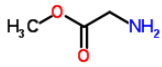
Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
acetonitrile	1.10	0.06	18.14
DMSO	6.64	0.52	154.95
methanol	4.12	0.21	94.36
THF	4.07	0.33	77.05
toluene	0.06	0.01	0.78
water	0.32	0.03	4.23

glycine methyl ester $C_3H_7NO_2$[207]

Compound Data

Molecular weight	89.0932	**H bond acceptors**	3	**Rule of 5 violations**	0
Compound type	amine	**H bond donors**	2	**ACD/ALogP**	-1.37
Phase 25°C	solid	**Rotatable bonds**	3	**Predicted density**	1.045 g/cm^3

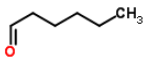


SMILES	O=C(OC)CN
InChIKey	KQSSATDQUYCRGS-UHFFFAOYSA-N

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
ethanol	0.19	0.01	2.21
methanol	1.32	0.06	17.60
THF	0.04	0.00	0.40

hexanaldehyde $C_6H_{12}O$[21]



Compound Data

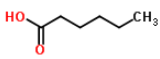
Molecular weight	100.159	**H bond acceptors**	1	**Rule of 5 violations**	0
Compound type	aldehyde	**H bond donors**	0	**ACD/ALogP**	2.37
Phase 25°C	liquid	**Rotatable bonds**	4	**Predicted density**	0.801 g/cm^3
SMILES	CCCCCC=O				
InChIKey	JARKCYVAAOWBJS-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
methanol	8.33	¤	¤

¤ Solute is very soluble/miscible, conversion fail.

hexanoic acid $C_6H_{12}O_2$[22]



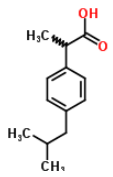
Compound Data

Molecular weight	116.158	**H bond acceptors**	2	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	1.88
Phase 25°C	liquid	**Rotatable bonds**	4	**Predicted density**	0.95 g/cm^3
SMILES	CCCCCC(=O)O				
InChIKey	FUZZWVXGSFPDMH-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
methanol	7.98	0.93	5071.51

ibuprofen $C_{13}H_{18}O_2$[905, 907]



Compound Data

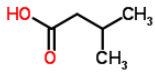
Molecular weight	206.281	**H bond acceptors**	2	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	3.5
Phase 25°C	solid	**Rotatable bonds**	4	**Predicted density**	1.029 g/cm^3
SMILES	O=C(O)C(c1ccc(cc1)CC(C)C)C				
InChIKey	HEFNNWSXXWATRW-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
1-decanol	1.12	0.22	35.98
1-octanol	1.20	0.20	39.60

1-pentanol	1.46	0.18	52.50
1-propanol	1.52	0.14	56.72
2-butanol	1.78	0.20	71.27
2-methyl-1-propanol	1.76	0.20	70.04
2-propanol	2.22	0.23	104.32
ethanol	1.18	0.08	40.88
methanol	1.16	0.06	41.41
water	0.00	0.00	0.00

isovaleric acid $C_5H_{10}O_2$[22]



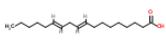
Compound Data

Molecular weight	102.132	**H bond acceptors**	2	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	1.26
Phase 25°C	liquid	**Rotatable bonds**	2	**Predicted density**	0.962 g/cm^3
SMILES	CC(C)CC(=O)O				
InChIKey	GWYFCOCPABKNJV-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
methanol	9.06	0.91	3222.50

linoleic acid $C_{18}H_{32}O_2$[22]



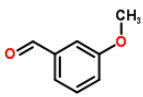
Compound Data

Molecular weight	280.445	**H bond acceptors**	2	**Rule of 5 violations**	1
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	7.06
Phase 25°C	liquid	**Rotatable bonds**	14	**Predicted density**	0.911 g/cm^3
SMILES	CCCCCC=CCC=CCCCCCCCC(=O)O				
InChIKey	OYHQOLUKZRVURQ-AVQMFFATSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
methanol	3.22	0.94	13712.82

m-anisaldehyde $C_8H_8O_2$[21]



Compound Data

Molecular weight	136.148	**H bond acceptors**	2	**Rule of 5 violations**	0
Compound type	aldehyde	**H bond donors**	0	**ACD/ALogP**	1.62
Phase 25°C	liquid	**Rotatable bonds**	2	**Predicted density**	1.088 g/cm^3
SMILES	COc1cc(C=O)ccc1				
InChIKey	WMPDAIZRQDCGFH-UHFFFAOYSA-N				

Solubility Data

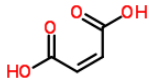
Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
methanol	8.20	¤	¤

¤ Solute is very soluble/miscible, conversion fail.

maleic acid $C_4H_4O_4$[39, 40]

Compound Data

Molecular weight	116.072	**H bond acceptors**	4	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	2	**ACD/ALogP**	0.21

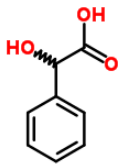


Phase 25°C	solid	**Rotatable bonds**	2	**Predicted density**	1.499 g/cm^3
SMILES	C(=CC(=O)O)C(=O)O				
InChIKey	VZCYOOQTPOCHFL-OWOJBTEDSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
ethanol	3.41	0.21	68.95
methanol	3.82	0.19	83.62

mandelic acid $C_8H_8O_3$[72, 71]



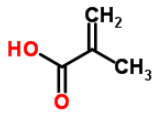
Compound Data

Molecular weight	152.147	**H bond acceptors**	3	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	2	**ACD/ALogP**	0.66
Phase 25°C	solid	**Rotatable bonds**	3	**Predicted density**	1.321 g/cm^3
SMILES	O=C(O)C(O)c1ccccc1				
InChIKey	IWYDHOAUDWTVEP-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
1-hexanol	0.00	0.00	0.00
chloroform	0.12	0.01	1.23
diethyl ether	0.77	0.08	17.51
methanol	3.54	0.20	120.77
toluene	0.00	0.00	0.00

methacrylic acid $C_4H_6O_2$[22]



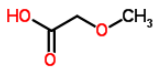
Compound Data

Molecular weight	86.0892	**H bond acceptors**	2	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	0.63
Phase 25°C	liquid	**Rotatable bonds**	1	**Predicted density**	1.023 g/cm^3
SMILES	CC(=C)C(=O)O				
InChIKey	CERQOIWHTDAKMF-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
methanol	11.79	0.98	17218.73

methoxyacetic acid $C_3H_6O_3$[22]



Compound Data

Molecular weight	90.0779	**H bond acceptors**	3	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	-0.49
Phase 25°C	liquid	**Rotatable bonds**	2	**Predicted density**	1.139 g/cm^3
SMILES	COCC(=O)O				
InChIKey	RMIODHQZRUFFFF-UHFFFAOYSA-N				

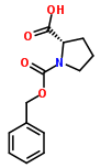
Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)

methanol	13.03	¤	¤

¤ Solute is very soluble/miscible, conversion fail.

N-Cbz-L-p roline $C_{13}H_{15}NO_4$[134]



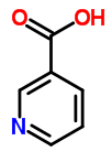
Compound Data

Molecular weight	249.262	**H bond acceptors**	5	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	1.21
Phase 25°C	solid	**Rotatable bonds**	4	**Predicted density**	1.309 g/cm^3
SMILES	O=C(O)[C@H]2N(C(=O)OCc1ccccc1)CCC2				
InChIKey	JXGVXCZADZNAMJ-NSHDSACASA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
methanol	4.26	0.49	746.90

nicotinic acid $C_6H_5NO_2$[100, 137, 907]



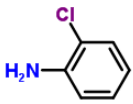
Compound Data

Molecular weight	123.109	**H bond acceptors**	3	**Rule of 5 violations**	0
Compound type	carboxlyic acid	**H bond donors**	1	**ACD/ALogP**	0.29
Phase 25°C	solid	**Rotatable bonds**	1	**Predicted density**	1.293 g/cm^3
SMILES	O=C(O)c1cccnc1				
InChIKey	PVNIIMVLHYAWGP-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
acetonitrile	0.00	0.00	0.00
benzene	0.00	0.00	0.00
DMSO	0.57	0.04	6.71
ethanol	0.09	0.01	1.43
methanol	0.06	0.00	1.03
THF	0.06	0.00	0.75
toluene	0.00	0.00	0.00
water	0.37	0.01	4.73

o-chloroaniline C_6H_6ClN[29]



Compound Data

Molecular weight	127.572	**H bond acceptors**	1	**Rule of 5 violations**	0
Compound type	amine	**H bond donors**	2	**ACD/ALogP**	1.93
Phase 25°C	liquid	**Rotatable bonds**	1	**Predicted density**	1.23 g/cm^3
SMILES	Nc1ccccc1Cl				
InChIKey	AKCRQHGQIJBRMN-UHFFFAOYSA-N				

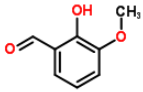
Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
methanol	9.51	0.97	11802.88

o-vanillin $C_8H_8O_3$[208, 205]

Compound Data

Molecular weight	152.147	**H bond acceptors**	3	**Rule of 5 violations**	0
Compound type	aldehyde	**H bond donors**	1	**ACD/ALogP**	1.31



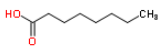
Phase 25°C	solid	**Rotatable bonds**	3	**Predicted density**	1.231 g/cm^3
SMILES	Oc1c(cccc1OC)C=O				
InChIKey	JJVNINGBHGBWJH-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
acetonitrile	2.56	0.17	76.28
ethanol	3.04	0.22	94.99
methanol	2.27	0.12	63.75
THF	5.37	0.56	268.76
toluene	3.82	0.43	126.41

octanoic acid $C_8H_{16}O_2$[22]

Compound Data



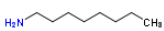
Molecular weight	144.211	**H bond acceptors**	2	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	2.92
Phase 25°C	liquid	**Rotatable bonds**	6	**Predicted density**	0.929 g/cm^3
SMILES	CCCCCCCC(=O)O				
InChIKey	WWZKQHOCKIZLMA-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
methanol	6.31	0.93	5899.86

octylamine $C_8H_{19}N$[29]

Compound Data



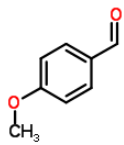
Molecular weight	129.243	**H bond acceptors**	1	**Rule of 5 violations**	0
Compound type	amine	**H bond donors**	2	**ACD/ALogP**	3.24
Phase 25°C	liquid	**Rotatable bonds**	7	**Predicted density**	0.786 g/cm^3
SMILES	NCCCCCCCC				
InChIKey	IOQPZZOEVPZRBK-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
methanol	6.05	0.98	20005.33

p-anisaldehyde $C_8H_8O_2$[21]

Compound Data



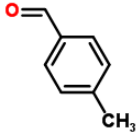
Molecular weight	136.148	**H bond acceptors**	2	**Rule of 5 violations**	0
Compound type	aldehyde	**H bond donors**	0	**ACD/ALogP**	1.64
Phase 25°C	liquid	**Rotatable bonds**	2	**Predicted density**	1.088 g/cm^3
SMILES	COc1ccc(C=O)cc1				
InChIKey	ZRSNZINYAWTAHE-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
methanol	8.22	¤	¤

¤ Solute is very soluble/miscible, conversion fail.

p-tolualdehyde C_8H_8O[21]



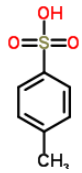
Compound Data

Molecular weight	120.148	**H bond acceptors**	1	**Rule of 5 violations**	0
Compound type	aldehyde	**H bond donors**	0	**ACD/ALogP**	2.01
Phase 25°C	liquid	**Rotatable bonds**	1	**Predicted density**	1.023 g/cm^3
SMILES	O=Cc1ccc(C)cc1				
InChIKey	FXLOVSHXALFLKQ-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
methanol	8.47	0.99	25859.25

p-toluenesulfonic acid $C_7H_8O_3S$[107, 136]



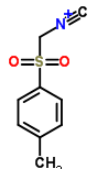
Compound Data

Molecular weight	172.202	**H bond acceptors**	3	**Rule of 5 violations**	0
Compound type	non-Ugi related	**H bond donors**	1	**ACD/ALogP**	-0.88
Phase 25°C	solid	**Rotatable bonds**	1	**Predicted density**	1.34 g/cm^3
SMILES	O=S(=O)(O)c1ccc(cc1)C				
InChIKey	JOXIMZWYDAKGHI-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
acetonitrile	0.11	0.01	2.69
ethanol	4.69	0.41	260.62
toluene	0.00	0.00	0.00

p-toluenesulfonylmethyl isocyanide $C_9H_9NO_2S$[208]



Compound Data

Molecular weight	195.238	**H bond acceptors**	3	**Rule of 5 violations**	0
Compound type	isonitrile	**H bond donors**	0	**ACD/ALogP**	1.38
Phase 25°C	solid	**Rotatable bonds**	2	**Predicted density**	‡
SMILES	O=S(=O)(c1ccc(cc1)C)C[N+]#[C-]				
InChIKey	CFOAUYCPAUGDFF-UHFFFAOYSA-N				

Solubility Data

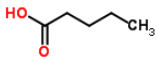
Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
acetonitrile	1.88	‡	‡
chloroform	0.21	‡	‡
ethanol	0.18	‡	‡
THF	1.86	‡	‡
toluene	0.14	‡	‡

‡ Solute density prediction fail.

pentanoic acid $C_5H_{10}O_2$[22]

Compound Data

Molecular weight	102.132	**H bond acceptors**	2	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	1.34
Phase 25°C	liquid	**Rotatable bonds**	3	**Predicted density**	0.966 g/cm^3

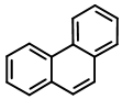


SMILES	CCCCC(=O)O
InChIKey	NQPDZGIKBAWPEJ-UHFFFAOYSA-N

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
methanol	9.19	0.93	4393.39

phenanthrene $C_{14}H_{10}$[902]



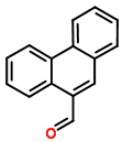
Compound Data

Molecular weight	178.229	**H bond acceptors**	0	**Rule of 5 violations**	0
Compound type	non-Ugi related	**H bond donors**	0	**ACD/ALogP**	4.55
Phase 25°C	solid	**Rotatable bonds**	0	**Predicted density**	1.13 g/cm^3
SMILES	c3cc2ccc1ccccc1c2cc3				
InChIKey	YNPNZTXNASCQKK-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
benzene	2.02	0.21	60.52
carbon disulfide	3.05	0.26	83.20
carbon tetrachloride	1.73	0.18	24.99
diethyl ether	1.36	0.15	42.04
ethanol	0.21	0.01	4.96
hexane	0.32	0.04	8.90

phenanthrene-9-carboxaldehyde $C_{15}H_{10}O$[30, 60, 136, 4, 242, 51, 24]



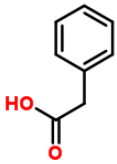
Compound Data

Molecular weight	206.239	**H bond acceptors**	1	**Rule of 5 violations**	0
Compound type	aldehyde	**H bond donors**	0	**ACD/ALogP**	4.28
Phase 25°C	solid	**Rotatable bonds**	1	**Predicted density**	1.217 g/cm^3
SMILES	O=Cc2cc3c(c1c2cccc1)cccc3				
InChIKey	QECIGCMPORCORE-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
1,1,2-trichlorotrifluoroethane	0.02	0.00	0.25
2-propanol	0.07	0.01	1.85
acetonitrile	0.16	0.01	4.39
benzene	0.66	0.06	17.56
chloroform	0.04	0.00	0.55
cyclohexane	0.07	0.01	1.85
cyclopentane	0.03	0.00	0.79
dichloromethane	0.00	0.00	0.00
diethyl ether	0.10	0.01	2.86
DMF	1.25	0.12	37.60
DMSO	0.77	0.06	16.62
ethanol	0.10	0.01	2.69
hexane	0.07	0.01	2.16
methanol	0.11	0.00	2.93
THF	2.32	0.23	87.22
toluene	0.14	0.01	3.40

phenylacetic acid $C_8H_8O_2$[86, 123, 907]



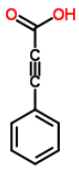
Compound Data

Molecular weight	136.148	**H bond acceptors**	2	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	1.72
Phase 25°C	solid	**Rotatable bonds**	2	**Predicted density**	1.164 g/cm^3
SMILES	O=C(O)Cc1ccccc1				
InChIKey	WLJVXDMOQOGPHL-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
acetonitrile	3.87	0.28	128.87
benzene	3.35	0.33	85.91
DMF	3.76	0.36	105.03
DMSO	6.35	0.64	305.77
THF	5.62	0.57	247.02
toluene	1.94	0.21	39.23
water	0.13	0.00	1.80

phenylpropynoic acid $C_9H_6O_2$[98, 82, 85]



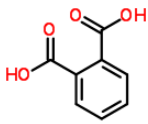
Compound Data

Molecular weight	146.143	**H bond acceptors**	2	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	1.7
Phase 25°C	solid	**Rotatable bonds**	1	**Predicted density**	1.24 g/cm^3
SMILES	O=C(C#Cc1ccccc1)O				
InChIKey	XNERWVPQCYSMLC-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
acetonitrile	1.38	0.08	32.24
DMSO	3.33	0.28	72.89
methanol	2.89	0.16	85.06
THF	4.34	0.42	143.85

phthalic acid $C_8H_6O_4$[138, 47]



Compound Data

Molecular weight	166.131	**H bond acceptors**	4	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	2	**ACD/ALogP**	1.22
Phase 25°C	solid	**Rotatable bonds**	2	**Predicted density**	1.451 g/cm^3
SMILES	O=C(O)c1ccccc1C(=O)O				
InChIKey	XNGIFLGASWRNHJ-UHFFFAOYSA-N				

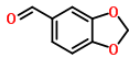
Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
chloroform	0.51	0.04	6.00
methanol	0.43	0.02	9.98

piperonal $C_8H_6O_3$[71, 30, 907]

Compound Data

Molecular weight	150.131	**H bond acceptors**	3	**Rule of 5 violations**	0
Compound type	aldehyde	**H bond donors**	0	**ACD/ALogP**	0.79

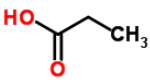


Phase 25°C	solid	**Rotatable bonds**	1	**Predicted density**	1.337 g/cm^3
SMILES	O=Cc1ccc2OCOc2c1				
InChIKey	SATCULPHIDQDRE-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
acetonitrile	7.11	0.66	708.73
chloroform	7.95	0.86	741.57
dichloromethane	7.59	0.78	616.11
diethyl ether	6.38	0.69	460.15
DMSO	6.05	0.57	257.75
ethanol	4.99	0.40	218.45
hexane	0.15	0.02	3.39
methanol	7.29	0.63	798.18
THF	7.44	0.78	750.82
toluene	6.40	0.71	392.09
water	0.02	0.00	0.30

propanoic acid $C_3H_6O_2$[22]



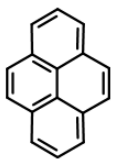
Compound Data

Molecular weight	74.0785	**H bond acceptors**	2	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	0.31
Phase 25°C	liquid	**Rotatable bonds**	1	**Predicted density**	1.019 g/cm^3
SMILES	CCC(=O)O				
InChIKey	XBDQKXXYIPTUBI-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
methanol	13.41	0.96	5249.72

pyrene $C_{16}H_{10}$[108, 109]



Compound Data

Molecular weight	202.251	**H bond acceptors**	0	**Rule of 5 violations**	1
Compound type	non-Ugi related	**H bond donors**	0	**ACD/ALogP**	5.19
Phase 25°C	solid	**Rotatable bonds**	0	**Predicted density**	1.248 g/cm^3
SMILES	c3ccc2ccc1cccc4c1c2c3cc4				
InChIKey	BBEAQIROQSPTKN-UHFFFAOYSA-N				

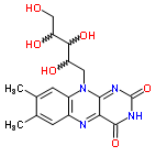
Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
acetonitrile	0.10	0.01	2.75

riboflavin $C_{17}H_{20}N_4O_6$[46]

Compound Data

Molecular weight	376.364	**H bond acceptors**	10	**Rule of 5 violations**	2
Compound type	non-Ugi related	**H bond donors**	5	**ACD/ALogP**	-2.02
Phase 25°C	solid	**Rotatable bonds**	9	**Predicted density**	1.65 g/cm^3
SMILES	O=C2/N=C\1/N(c3cc(c(cc3/N=C/1C(=O)N2)C)C)CC(O)C(O)C(O)CO				

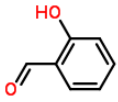


InChIKey	AUNGANRZJHBGPY-UHFFFAOYSA-N

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
methanol	0.00	0.00	0.00

salicylaldehyde $C_7H_6O_2$[21]



Compound Data

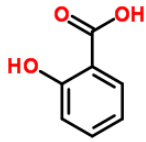
Molecular weight	122.121	**H bond acceptors**	2	**Rule of 5 violations**	0
Compound type	aldehyde	**H bond donors**	1	**ACD/ALogP**	1.61
Phase 25°C	liquid	**Rotatable bonds**	2	**Predicted density**	1.226 g/cm^3
SMILES	c1cccc(C=O)c1O				
InChIKey	SMQUZDBALVYZAC-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
methanol	10.04	¤	¤

¤ Solute is very soluble/miscible, conversion fail.

salicylic acid $C_7H_6O_3$[34, 902, 901, 138, 52, 907]



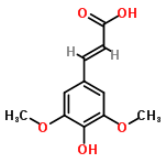
Compound Data

Molecular weight	138.121	**H bond acceptors**	3	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	2	**ACD/ALogP**	2.06
Phase 25°C	solid	**Rotatable bonds**	2	**Predicted density**	1.375 g/cm^3
SMILES	O=C(O)c1ccccc1O				
InChIKey	YGSDEFSMJLZEOE-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
chloroform	0.19	0.01	1.74
dichloromethane	0.15	0.01	1.68
ethanol	2.34	0.15	54.17
methanol	2.65	0.13	66.24
toluene	0.11	0.01	1.76
water	0.01	0.00	0.14

sinapic acid $C_{11}H_{12}O_5$[25, 2]



Compound Data

Molecular weight	224.21	**H bond acceptors**	5	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	2	**ACD/ALogP**	1.29
Phase 25°C	solid	**Rotatable bonds**	5	**Predicted density**	1.307 g/cm^3
SMILES	O=C(O)\C=C\c1cc(OC)c(O)c(OC)c1				
InChIKey	PCMORTLOPMLEFB-ONEGZZNKSA-N				

Solubility Data

Solvent	Concentration	Mole Fraction (X)	pph (g/100g)

	(M)		
ethanol	0.79	0.05	26.27
methanol	1.04	0.05	37.69
THF	0.45	0.04	11.95

sodium chloride $ClNa$[1, 207, 900]

Compound Data

Molecular weight	58.4428	**H bond acceptors**	0	**Rule of 5 violations**	0
Compound type	non-Ugi related	**H bond donors**	0	**ACD/ALogP**	†
Phase 25°C	solid	**Rotatable bonds**	0	**Predicted density**	‡
SMILES	Cl[Na]				
InChIKey	FAPWRFPIFSIZLT-UHFFFAOYSA-M				

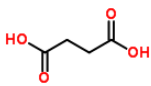
Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
ethanol	0.00	0.00	0.00
methanol	0.21	‡	‡
THF	0.04	‡	‡

† ACD/ALogP prediction fail.

‡ Solute density prediction fail.

succinic acid $C_4H_6O_4$[45, 44, 42]



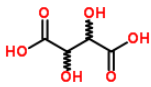
Compound Data

Molecular weight	118.088	**H bond acceptors**	4	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	2	**ACD/ALogP**	-0.59
Phase 25°C	solid	**Rotatable bonds**	3	**Predicted density**	1.408 g/cm^3
SMILES	OC(CCC(O)=O)=O				
InChIKey	KDYFGRWQOYBRFD-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
2-propanol	0.34	0.03	5.22
ethanol	0.57	0.03	9.06

tartaric acid $C_4H_6O_6$[38]



Compound Data

Molecular weight	150.087	**H bond acceptors**	6	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	4	**ACD/ALogP**	-1.43
Phase 25°C	solid	**Rotatable bonds**	5	**Predicted density**	1.886 g/cm^3
SMILES	C(C(C(=O)O)O)(C(=O)O)O				
InChIKey	FEWJPZIEWOKRBE-UHFFFAOYSA-N				

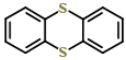
Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
acetone	0.79	0.06	16.39

thianthrene $C_{12}H_8S_2$[905]

Compound Data

Molecular weight	216.322	**H bond acceptors**	0	**Rule of 5 violations**	0
Compound type	non-Ugi related	**H bond donors**	0	**ACD/ALogP**	4.57



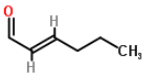
Phase 25°C	solid	**Rotatable bonds**	0	**Predicted density**	1.31 g/cm^3
SMILES	S1c3c(Sc2c1cccc2)cccc3				
InChIKey	GVIJJXMXTUZIOD-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
1-decanol	0.04	0.01	1.05
1,4-dioxane	0.27	0.02	6.14
acetonitrile	0.02	0.00	0.58
decane	0.03	0.01	0.89
diethyl ether	0.08	0.01	2.39
diisopropyl ether	0.05	0.01	1.44
ethylene glycol	0.02	0.00	0.40
hexadecane	0.03	0.01	0.84
nonane	0.03	0.01	0.90
undecane	0.03	0.01	0.88

trans-2-hexen-1-al $C_6H_{10}O$[22]

Compound Data



Molecular weight	98.143	**H bond acceptors**	1	**Rule of 5 violations**	0
Compound type	aldehyde	**H bond donors**	0	**ACD/ALogP**	1.58
Phase 25°C	liquid	**Rotatable bonds**	3	**Predicted density**	0.828 g/cm^3
SMILES	CCCC=CC=O				
InChIKey	MBDOYVRWFFCFHM-SNAWJCMRSA-N				

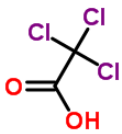
Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
methanol	8.62	¤	¤

¤ Solute is very soluble/miscible, conversion fail.

trichloroacetic acid $C_2HCl_3O_2$[22]

Compound Data



Molecular weight	163.387	**H bond acceptors**	2	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	1.67
Phase 25°C	liquid	**Rotatable bonds**	0	**Predicted density**	1.807 g/cm^3
SMILES	C(=O)(C(Cl)(Cl)Cl)O				
InChIKey	YNJBWRMUSHSURL-UHFFFAOYSA-N				

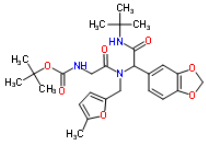
Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
methanol	9.91	0.80	2068.61

UCExp216-3A $C_{26}H_{35}N_3O_7$[113]

Compound Data

Molecular weight	501.572	**H bond acceptors**	10	**Rule of 5 violations**	2
Compound type	Ugi Product	**H bond donors**	2	**ACD/ALogP**	4.06
Phase 25°C	solid	**Rotatable bonds**	10	**Predicted density**	1.209 g/cm^3

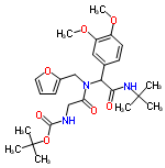


SMILES	CC(C)(C)NC(=O)C(c1ccc2OCOc2c1)N(Cc3ccc(C)o3)C(=O)CNC(=O)OC(C)(C)C
InChIKey	MTSJLRXXXJHLDI-UHFFFAOYSA-N

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
acetonitrile	0.04	0.00	2.73
DMSO	0.28	0.02	14.46
ethanol	0.03	0.00	1.95
methanol	0.06	0.00	4.10
toluene	0.01	0.00	0.58

Ugi product 104C (UC) $C_{26}H_{37}N_3O_7$[65]



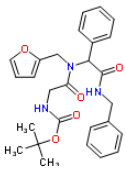
Compound Data

Molecular weight	503.588	**H bond acceptors**	10	**Rule of 5 violations**	2
Compound type	Ugi Product	**H bond donors**	2	**ACD/ALogP**	3.48
Phase 25°C	solid	**Rotatable bonds**	12	**Predicted density**	1.156 g/cm^3
SMILES	O=C(CNC(=O)OC(C)(C)C)N(Cc1ccco1)C(c2ccc(OC)c(OC)c2)C(=O)NC(C)(C)C				
InChIKey	RXLIYJRDQONCRV-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
methanol	0.29	0.01	22.20

Ugi product 108C (UC) $C_{27}H_{31}N_3O_5$[65]



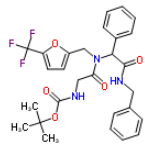
Compound Data

Molecular weight	477.552	**H bond acceptors**	8	**Rule of 5 violations**	0
Compound type	Ugi Product	**H bond donors**	2	**ACD/ALogP**	4.43
Phase 25°C	solid	**Rotatable bonds**	11	**Predicted density**	1.198 g/cm^3
SMILES	O=C(CNC(=O)OC(C)(C)C)N(Cc1ccco1)C(c2ccccc2)C(=O)NCc3ccccc3				
InChIKey	COBIEQVPXUIPRH-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
methanol	0.36	0.02	26.66

Ugi product 109C(UC) $C_{28}H_{30}F_3N_3O_5$[65]



Compound Data

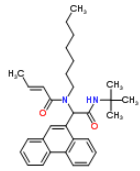
Molecular weight	545.55	**H bond acceptors**	8	**Rule of 5 violations**	1
Compound type	Ugi Product	**H bond donors**	2	**ACD/ALogP**	5
Phase 25°C	solid	**Rotatable bonds**	11	**Predicted density**	1.263 g/cm^3
SMILES	O=C(CNC(=O)OC(C)(C)C)N(Cc1ccc(o1)C(F)(F)F)C(c2ccccc2)C(=O)NCc3ccccc3				
InChIKey	XRHVBZVCUGMHPN-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration	Mole Fraction (X)	pph (g/100g)

	(M)		
methanol	0.15	0.01	11.62

Ugi product 148B $C_{31}H_{40}N_2O_2$[145, 96]



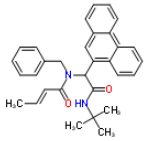
Compound Data

Molecular weight	472.661	**H bond acceptors**	4	**Rule of 5 violations**	1
Compound type	Ugi Product	**H bond donors**	1	**ACD/ALogP**	7.47
Phase 25°C	solid	**Rotatable bonds**	11	**Predicted density**	1.067 g/cm^3
SMILES	CC(C)(C)NC(=O)C(N(CCCCCCC)C(=O)\C=C\C)c2cc3ccccc3c1ccccc12				
InChIKey	HOKCUULFCASAQT-FRKPEAEDSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
acetonitrile	0.02	0.00	1.28
benzene	0.05	0.00	2.77
ethanol	0.00	0.00	0.00
methanol	0.03	0.00	1.91
THF	0.40	0.04	25.03
toluene	0.14	0.02	8.10

Ugi product 150D(UC) $C_{31}H_{32}N_2O_2$[78, 65]



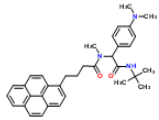
Compound Data

Molecular weight	464.598	**H bond acceptors**	4	**Rule of 5 violations**	1
Compound type	Ugi Product	**H bond donors**	1	**ACD/ALogP**	6.18
Phase 25°C	solid	**Rotatable bonds**	7	**Predicted density**	1.148 g/cm^3
SMILES	CC(C)(C)NC(=O)C(N(Cc1ccccc1)C(=O)\C=C\C)c3cc4ccccc4c2ccccc23				
InChIKey	PBZQTKRWYXTXIS-WLRTZDKTSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
acetonitrile	0.04	0.00	2.53
DMSO	0.18	0.01	8.21
ethanol	0.04	0.00	2.42
methanol	0.07	0.00	4.12
THF	0.49	0.05	31.41
toluene	0.04	0.00	2.17

Ugi product 171K (UC) $C_{35}H_{39}N_3O_2$[65]



Compound Data

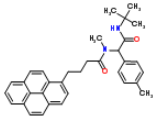
Molecular weight	533.703	**H bond acceptors**	5	**Rule of 5 violations**	2
Compound type	Ugi Product	**H bond donors**	1	**ACD/ALogP**	6.83
Phase 25°C	solid	**Rotatable bonds**	9	**Predicted density**	1.177 g/cm^3
SMILES	CN(C)c1ccc(cc1)C(C(=O)NC(C)(C)C)N(C)C(=O)CCCc5ccc4ccc3cccc2ccc5c4c23				
InChIKey	XWJCEFGKUMNBLK-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
methanol	0.02	0.00	1.43

Ugi product 173B (UC) $C_{34}H_{36}N_2O_2$[65]

Compound Data

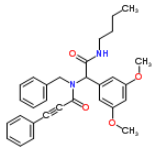


Molecular weight	504.662	**H bond acceptors**	4	**Rule of 5 violations**	2
Compound type	Ugi Product	**H bond donors**	1	**ACD/ALogP**	7.18
Phase 25°C	solid	**Rotatable bonds**	8	**Predicted density**	1.169 g/cm^3
SMILES	CC(C)(C)NC(=O)C(N(C)C(=O)CCCc4ccc3ccc2cccc1ccc4c3c12)c5ccc(C)cc5				
InChIKey	KNHNINYGWDFMQZ-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
methanol	0.08	0.00	5.55

Ugi product 173G (UC) $C_{30}H_{32}N_2O_4$[65]



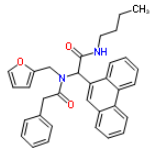
Compound Data

Molecular weight	484.586	**H bond acceptors**	6	**Rule of 5 violations**	1
Compound type	Ugi Product	**H bond donors**	1	**ACD/ALogP**	6.39
Phase 25°C	solid	**Rotatable bonds**	11	**Predicted density**	1.19 g/cm^3
SMILES	COc1cc(cc(OC)c1)C(C(=O)NCCCC)N(Cc2ccccc2)C(=O)C#Cc3ccccc3				
InChIKey	VOGZWAORWZGOTE-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
methanol	0.18	0.01	12.50

Ugi product 176C (UC) $C_{33}H_{32}N_2O_3$[155, 127, 144]



Compound Data

Molecular weight	504.619	**H bond acceptors**	5	**Rule of 5 violations**	2
Compound type	Ugi Product	**H bond donors**	1	**ACD/ALogP**	6.9
Phase 25°C	solid	**Rotatable bonds**	10	**Predicted density**	1.197 g/cm^3
SMILES	O=C(Cc1ccccc1)N(Cc2ccco2)C(c4cc5ccccc5c3ccccc34)C(=O)NCCCC				
InChIKey	RGGYLRDGJAJICY-UHFFFAOYSA-N				

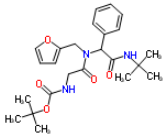
Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
1-octanol	0.01	0.00	0.62
1,4-dioxane	0.19	0.02	10.47
acetonitrile	0.02	0.00	1.36
benzene	0.03	0.00	1.76
diethyl ether	0.02	0.00	1.39
DMF	0.27	0.02	17.67
DMSO	0.19	0.01	9.48
ethanol	0.01	0.00	0.65
methanol	0.02	0.00	1.35
THF	0.28	0.02	17.72
toluene	0.03	0.00	1.76

Ugi product 206B(UC) $C_{24}H_{33}N_3O_5$[56]

Compound Data

Molecular weight	443.536	**H bond acceptors**	8	**Rule of 5 violations**	0
Compound type	Ugi Product	**H bond donors**	2	**ACD/ALogP**	3.74
Phase 25°C	solid	**Rotatable bonds**	10	**Predicted density**	1.145 g/cm^3
SMILES	O=C(CNC(=O)OC(C)(C)C)N(Cc1ccco1)C(c2ccccc2)C(=O)NC(C)(C)C				



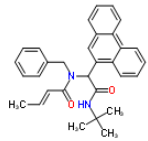
InChIKey	BXAOUWIVXLQYOZ-UHFFFAOYSA-N

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
2-propanol	0.00	0.00	0.00
acetonitrile	0.02	0.00	1.20
benzene	0.01	0.00	0.51
carbon tetrachloride	0.00	0.00	0.00
chloroform	0.57	0.06	21.63
cyclohexane	0.00	0.00	0.00
cyclopentane	0.00	0.00	0.00
dichloromethane	0.24	0.02	9.37
diethyl ether	0.00	0.00	0.00
DMF	0.24	0.02	13.49
DMSO	0.23	0.02	10.19
ethanol	0.02	0.00	1.15
hexane	0.00	0.00	0.00
methanol	0.05	0.00	3.00
THF	0.26	0.02	14.19
toluene	0.01	0.00	0.51

Ugi product 214C(UC) $C_{31}H_{32}N_2O_2$[27]

Compound Data



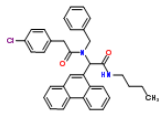
Molecular weight	464.598	**H bond acceptors**	4	**Rule of 5 violations**	1
Compound type	Ugi Product	**H bond donors**	1	**ACD/ALogP**	6.18
Phase 25°C	solid	**Rotatable bonds**	7	**Predicted density**	1.148 g/cm^3
SMILES	CC(C)(C)NC(=O)C(N(Cc1ccccc1)C(=O)\C=C\C)c3cc4ccccc4c2ccccc23				
InChIKey	PBZQTKRWYXTXIS-WLRTZDKTSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
methanol	0.01	0.00	0.62
THF	0.35	0.03	20.96

Ugi product 215F(UC) $C_{35}H_{33}ClN_2O_2$[71]

Compound Data



Molecular weight	549.102	**H bond acceptors**	4	**Rule of 5 violations**	2
Compound type	Ugi Product	**H bond donors**	1	**ACD/ALogP**	8.33
Phase 25°C	solid	**Rotatable bonds**	10	**Predicted density**	1.219 g/cm^3
SMILES	Clc5ccc(CC(=O)N(Cc1ccccc1)C(c3cc4ccccc4c2ccccc23)C(=O)NCCCC)cc5				
InChIKey	NCOSMKWGNCJPCR-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
acetonitrile	0.04	0.00	2.99
DMSO	0.38	0.03	22.91
ethanol	0.06	0.00	4.34
methanol	0.12	0.01	9.25
THF	0.61	0.06	51.09

toluene	0.07	0.01	4.56

Ugi Product 216-4A $C_{28}H_{40}N_2O_7$[97, 93]

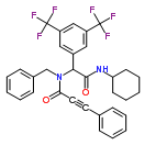
Compound Data

Molecular weight	516.626	**H bond acceptors**	9	**Rule of 5 violations**	1
Compound type	Ugi Product	**H bond donors**	2	**ACD/ALogP**	4.08
Phase 25°C	solid	**Rotatable bonds**	13	**Predicted density**	1.12 g/cm^3
SMILES	Cc1ccc(o1)CC(C(c2ccc(c(c2)OC)OC)C(=O)NC(C)(C)C)C(=O)CNC(=O)OC(C)(C)C				
InChIKey					

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
acetonitrile	0.13	0.01	9.17
DMSO	0.06	0.00	2.90
ethanol	0.07	0.00	4.79
methanol	0.18	0.01	13.47
toluene	0.20	0.02	13.07

Ugi product 234 C32H28F6N2O2[65]



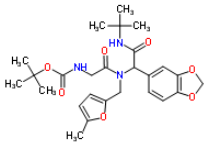
Compound Data

Molecular weight	586.57	**H bond acceptors**	4	**Rule of 5 violations**	2
Compound type	Ugi Product	**H bond donors**	1	**ACD/ALogP**	8.56
Phase 25°C	solid	**Rotatable bonds**	7	**Predicted density**	1.33 g/cm^3
SMILES	c1ccc(cc1)CN(C(c2cc(cc(c2)C(F)(F)F)C(F)(F)F)C(=O)NC3CCCCC3)C(=O)C#Cc4ccccc4				
InChIKey					

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
ethanol	0.23	0.01	19.25

Ugi product 62E (UC) $C_{26}H_{35}N_3O_7$[65]



Compound Data

Molecular weight	501.572	**H bond acceptors**	10	**Rule of 5 violations**	2
Compound type	Ugi Product	**H bond donors**	2	**ACD/ALogP**	4.06
Phase 25°C	solid	**Rotatable bonds**	10	**Predicted density**	1.209 g/cm^3
SMILES	CC(C)(C)NC(=O)C(c1ccc2OCOc2c1)N(Cc3ccc(C)o3)C(=O)CNC(=O)OC(C)(C)C				
InChIKey	MTSJLRXXXJHLDI-UHFFFAOYSA-N				

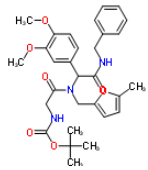
Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
methanol	0.13	0.01	9.15

Ugi product 64C (UC) $C_{30}H_{37}N_3O_7$[65]

Compound Data

Molecular weight	551.631	**H bond acceptors**	10	**Rule of 5 violations**	2
Compound type	Ugi Product	**H bond donors**	2	**ACD/ALogP**	4.63

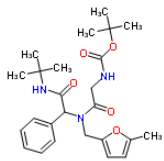


Phase 25°C	solid	**Rotatable bonds**	13	**Predicted density**	1.192 g/cm^3
SMILES	CC(C)(C)OC(=O)NCC(=O)N(Cc1ccc(C)o1)C(C(=O)NCc2ccccc2)c3ccc(OC)c(OC)c3				
InChIKey	VMMFSZQZVZNXSK-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
methanol	0.03	0.00	2.23

Ugi product 86B (UC) $C_{25}H_{35}N_3O_5$[65]



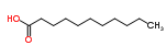
Compound Data

Molecular weight	457.562	**H bond acceptors**	8	**Rule of 5 violations**	0
Compound type	Ugi Product	**H bond donors**	2	**ACD/ALogP**	4.2
Phase 25°C	solid	**Rotatable bonds**	10	**Predicted density**	1.133 g/cm^3
SMILES	O=C(CNC(=O)OC(C)(C)C)N(Cc1ccc(C)o1)C(c2ccccc2)C(=O)NC(C)(C)C				
InChIKey	WDAGUORBKBNETC-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
methanol	0.17	0.01	11.09

undecanoic acid $C_{11}H_{22}O_2$[58]



Compound Data

Molecular weight	186.291	**H bond acceptors**	2	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	4.5
Phase 25°C	solid	**Rotatable bonds**	9	**Predicted density**	0.909 g/cm^3
SMILES	O=C(O)CCCCCCCCCC				
InChIKey	ZDPHROOEEOARMN-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
ethanol	3.94	0.55	488.75
methanol	4.00	0.49	549.05

urea CH_4N_2O[53]

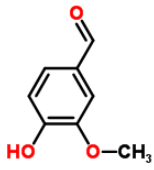
Compound Data

Molecular weight	60.0553	**H bond acceptors**	3	**Rule of 5 violations**	0
Compound type	amide	**H bond donors**	4	**ACD/ALogP**	-2.11
Phase 25°C	solid	**Rotatable bonds**	0	**Predicted density**	1.212 g/cm^3
SMILES	NC(=O)N				
InChIKey	XSQUKJJJFZCRTK-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
methanol	2.66	0.12	24.44

vanillin $C_8H_8O_3$[906, 4, 55, 11, 10, 207, 19]



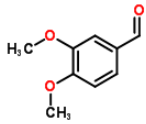
Compound Data

Molecular weight	152.147	**H bond acceptors**	3	**Rule of 5 violations**	0
Compound type	aldehyde	**H bond donors**	1	**ACD/ALogP**	1.19
Phase 25°C	solid	**Rotatable bonds**	3	**Predicted density**	1.231 g/cm^3
SMILES	Oc1ccc(cc1OC)C=O				
InChIKey	MWOOGOJBHIARFG-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
1-propanol	1.82	0.15	44.94
1,2-dichloroethane	1.17	0.10	17.74
acetonitrile	2.36	0.15	67.86
butanone	2.14	0.21	56.32
ethanol	2.47	0.17	69.35
methanol	4.16	0.27	173.01
THF	3.59	0.34	108.89
toluene	0.30	0.03	5.44

veratraldehyde $C_9H_{10}O_3$[41, 71, 208]



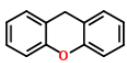
Compound Data

Molecular weight	166.174	**H bond acceptors**	3	**Rule of 5 violations**	0
Compound type	aldehyde	**H bond donors**	0	**ACD/ALogP**	1.61
Phase 25°C	solid	**Rotatable bonds**	3	**Predicted density**	1.114 g/cm^3
SMILES	COc1cc(ccc1OC)C=O				
InChIKey	WJUFSDZVCOTFON-UHFFFAOYSA-N				

Solubility Data

Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
acetone	3.03	0.29	119.01
acetonitrile	5.76	0.69	904.54
chloroform	5.70	0.75	421.71
dichloromethane	6.06	0.81	837.51
diethyl ether	3.55	0.43	170.84
DMSO	5.31	0.64	386.17
ethanol	5.49	0.64	645.97
hexane	0.07	0.01	1.74
methanol	5.87	0.67	1041.50
THF	5.33	0.68	480.29
toluene	4.54	0.60	267.44

xanthene $C_{13}H_{10}O$[905]



Compound Data

Molecular weight	182.218	**H bond acceptors**	1	**Rule of 5 violations**	0
Compound type	non-Ugi related	**H bond donors**	0	**ACD/ALogP**	3.93
Phase 25°C	solid	**Rotatable bonds**	0	**Predicted density**	1.159 g/cm^3
SMILES	O2c1ccccc1Cc3c2cccc3				
InChIKey	GJCOSYZMQJWQCA-UHFFFAOYSA-N				

Solubility Data

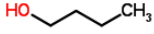
Solvent	Concentration (M)	Mole Fraction (X)	pph (g/100g)
1-butanol	0.19	0.02	4.43
1-decanol	0.24	0.05	5.49
1-heptanol	0.23	0.03	5.30
1-hexanol	0.22	0.03	5.09

1-octanol	0.24	0.04	5.52
1-pentanol	0.20	0.02	4.64
1-propanol	0.15	0.01	3.52
1,2-dichloroethane	1.62	0.15	33.77
2-butanol	0.13	0.01	3.02
2-methyl-1-pentanol	0.16	0.02	3.67
2-methyl-1-propanol	0.12	0.01	2.78
2-methyl-2-butanol	0.18	0.02	4.16
2-methyl-2-propanol	0.12	0.01	2.77
2-pentanol	0.16	0.02	3.70
2-propanol	0.11	0.01	2.58
2,2,4-trimethylpentane	0.15	0.02	3.95
3-methyl-1-butanol	0.15	0.02	3.46
4-methyl-2-pentanol	0.14	0.02	3.22
acetonitrile	0.35	0.02	9.03
carbon tetrachloride	1.25	0.12	16.71
cyclohexane	0.39	0.04	9.58
cyclooctane	0.38	0.05	9.32
cyclopentanol	0.33	0.03	6.32
decane	0.24	0.05	6.19
dibutyl ether	0.50	0.08	12.68
diethyl ether	0.79	0.08	22.39
diisopropyl ether	0.41	0.06	10.54
ethanol	0.10	0.01	2.37
heptane	0.24	0.03	6.54
hexadecane	0.24	0.07	5.88
hexane	0.23	0.03	6.44
methanol	0.10	0.00	2.46
methyl tert-butyl ether	0.65	0.08	17.59
methylcyclohexane	0.33	0.04	8.17
nonane	0.24	0.04	6.28
octane	0.25	0.04	6.67
undecane	0.23	0.05	5.85

Solvents

Below we present data for the solvents used in the Open Notebook Science Challenge. We present various properties such as density and boiling point together with important solvation related parameters such as dielectric constants and dipole moments [1000-1100].

1-butanol



Compound Data

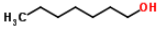
Molecular weight	74.1216
Predicted density	0.805 g/cm^3
Boiling point	117.7 °C
Dipole moment	1.66
Dielectric constant	17.8
SMILES	CCCCO
InChIKey	LRHPLDYGYMQRHN-UHFFFAOYSA-N

1-decanol

Compound Data

Molecular weight	158.2811
Predicted density	0.828 g/cm^3
Boiling point	227.8 °C
SMILES	OCCCCCCCCCC
InChIKey	MWKFXSUHUHTGQN-UHFFFAOYSA-N

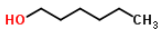
1-heptanol



Compound Data

Molecular weight	116.2013
Predicted density	0.82 g/cm^3
Boiling point	176.9 °C
SMILES	OCCCCCCC
InChIKey	BBMCTIGTTCKYKF-UHFFFAOYSA-N

1-hexanol



Compound Data

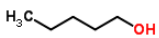
Molecular weight	102.1748
Predicted density	0.816 g/cm^3
Boiling point	158.2 °C
Dipole moment	1.55
Dielectric constant	13.3
SMILES	OCCCCCC
InChIKey	ZSIAUFGUXNUGDI-UHFFFAOYSA-N

1-octanol

Compound Data

Molecular weight	130.2279
Predicted density	0.823 g/cm^3
Boiling point	194.7 °C
SMILES	OCCCCCCCC
InChIKey	KBPLFHHGFOOTCA-UHFFFAOYSA-N

1-pentanol

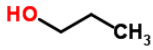


Compound Data

Molecular weight	88.1482
Predicted density	0.811 g/cm^3
Boiling point	138.5 °C
SMILES	OCCCCC
InChIKey	AMQJEAYHLZJPGS-UHFFFAOYSA-N

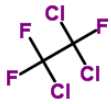
1-propanol

Compound Data



Molecular weight	60.095
Predicted density	0.795 g/cm^3
Boiling point	95.8 °C
Dipole moment	1.68
Dielectric constant	20.1
SMILES	CCCO
InChIKey	BDERNNFJNOPAEC-UHFFFAOYSA-N

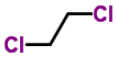
1,1,2-trichlorotrifluoroethane



Compound Data

Molecular weight	187.3756
Predicted density	1.67 g/cm^3
Boiling point	50.9 °C
SMILES	ClC(F)(F)C(Cl)(Cl)F
InChIKey	AJDIZQLSFPQPEY-UHFFFAOYSA-N

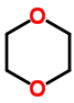
1,2-dichloroethane



Compound Data

Molecular weight	98.9592
Predicted density	1.173 g/cm^3
Boiling point	83.5 °C
SMILES	ClCCCl
InChIKey	WSLDOOZREJYCGB-UHFFFAOYSA-N

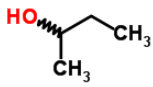
1,4-dioxane



Compound Data

Molecular weight	88.1051
Predicted density	0.995 g/cm^3
Boiling point	102.9 °C
SMILES	O1CCOCC1
InChIKey	RYHBNJHYFVUHQT-UHFFFAOYSA-N

2-butanol



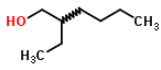
Compound Data

Molecular weight	74.1216
Predicted density	0.801 g/cm^3
Boiling point	96.6 °C
SMILES	OC(C)CC
InChIKey	BTANRVKWQNVYAZ-UHFFFAOYSA-N

2-ethyl-1-hexanol

Compound Data

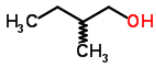
Molecular weight	130.2279
Predicted density	0.821 g/cm^3



Boiling point	184.6 °C
SMILES	OCC(CC)CCCC
InChIKey	YIWUKEYIRIRTPP-UHFFFAOYSA-N

2-methyl-1-butanol

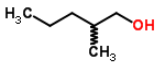
Compound Data



Molecular weight	88.1482
Predicted density	0.809 g/cm^3
Boiling point	128.7 °C
SMILES	OCC(C)CC
InChIKey	QPRQEDXDYOZYLA-UHFFFAOYSA-N

2-methyl-1-pentanol

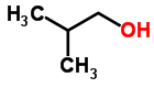
Compound Data



Molecular weight	102.1748
Predicted density	0.814 g/cm^3
Boiling point	148 °C
SMILES	OCC(C)CCC
InChIKey	PFNHSEQQEPMLNI-UHFFFAOYSA-N

2-methyl-1-propanol

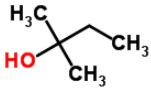
Compound Data



Molecular weight	74.1216
Predicted density	0.801 g/cm^3
Boiling point	105 °C
SMILES	OCC(C)C
InChIKey	ZXEKIIBDNHEJCQ-UHFFFAOYSA-N

2-methyl-2-butanol

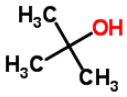
Compound Data



Molecular weight	88.1482
Predicted density	0.811 g/cm^3
Boiling point	102 °C
SMILES	OC(C)(C)CC
InChIKey	MSXVEPNJUHWQHW-UHFFFAOYSA-N

2-methyl-2-propanol

Compound Data

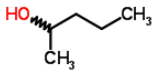


Molecular weight	74.1216
Predicted density	0.804 g/cm^3
Boiling point	84.6 °C
SMILES	OC(C)(C)C

InChIKey	DKGAVHZHDRPRBM-UHFFFAOYSA-N

2-pentanol

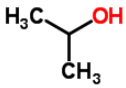
Compound Data



Molecular weight	88.1482
Predicted density	0.809 g/cm^3
Boiling point	118.8 °C
SMILES	OC(C)CCC
InChIKey	JYVLIDXNZAXMDK-UHFFFAOYSA-N

2-propanol

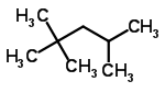
Compound Data



Molecular weight	60.095
Predicted density	0.791 g/cm^3
Boiling point	73 °C
Dipole moment	1.66
Dielectric constant	1.89
SMILES	CC(O)C
InChIKey	KFZMGEQAYNKOFK-UHFFFAOYSA-N

2,2,4-trimethylpentane

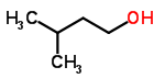
Compound Data



Molecular weight	114.2285
Predicted density	0.709 g/cm^3
Boiling point	98.8 °C
SMILES	CC(C)CC(C)(C)C
InChIKey	NHTMVDHEPJAVLT-UHFFFAOYSA-N

3-methyl-1-butanol

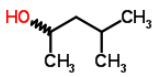
Compound Data



Molecular weight	88.1482
Predicted density	0.809 g/cm^3
Boiling point	131.2 °C
SMILES	OCCC(C)C
InChIKey	PHTQWCKDNZKARW-UHFFFAOYSA-N

4-methyl-2-pentanol

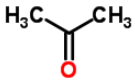
Compound Data



Molecular weight	102.1748
Predicted density	0.811 g/cm^3
Boiling point	133.5 °C
SMILES	OC(C)CC(C)C
InChIKey	WVYWICLMDOOCFB-UHFFFAOYSA-N

acetone

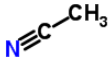
Compound Data



Molecular weight	58.0791
Predicted density	0.772 g/cm^3
Boiling point	46.5 °C
Dipole moment	2.88
Dielectric constant	20.7
SMILES	CC(=O)C
InChIKey	CSCPPACGZOOCGX-UHFFFAOYSA-N

acetonitrile

Compound Data



Molecular weight	41.0519
Predicted density	0.747 g/cm^3
Boiling point	63.5 °C
Dipole moment	3.92
Dielectric constant	36.6
SMILES	CCN
InChIKey	QUSNBJAOOMFDIB-UHFFFAOYSA-N

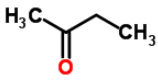
benzene

Compound Data

Molecular weight	78.1118
Predicted density	0.873 g/cm^3
Boiling point	78.8 °C
Dipole moment	0
Dielectric constant	2.28
SMILES	c1ccccc1
InChIKey	UHOVQNZJYSORNB-UHFFFAOYSA-N

butanone

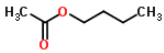
Compound Data



Molecular weight	72.1057
Predicted density	0.786 g/cm^3
Boiling point	75.6 °C
SMILES	O=C(C)CC
InChIKey	ZWEHNKRNPOVVGH-UHFFFAOYSA-N

butyl acetate

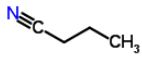
Compound Data



Molecular weight	116.1583
Predicted density	0.886 g/cm^3
Boiling point	126.6 °C
SMILES	O=C(OCCCC)C
InChIKey	DKPFZGUDAPQIHT-UHFFFAOYSA-N

butyronitrile

Compound Data



Molecular weight	69.1051
Predicted density	0.785 g/cm^3
Boiling point	117.3 °C
SMILES	N#CCCC
InChIKey	KVNRLNFWIYMESJ-UHFFFAOYSA-N

carbon disulfide

Compound Data

Molecular weight	76.1407
Predicted density	1.259 g/cm^3
Boiling point	46.2 °C
SMILES	S=C=S
InChIKey	QGJOPFRUJISHPQ-UHFFFAOYSA-N

carbon tetrachloride

Compound Data

Molecular weight	153.8227
Predicted density	1.697 g/cm^3
Boiling point	76 °C
Dipole moment	0
Dielectric constant	2.24
SMILES	[C](Cl)(Cl)(Cl)Cl
InChIKey	VZGDMQKNWNREIO-UHFFFAOYSA-N

chloroform

Compound Data

Molecular weight	119.3776
Predicted density	1.5 g/cm^3
Boiling point	61.2 °C
Dipole moment	1.15
Dielectric constant	5.5
SMILES	C(Cl)(Cl)Cl
InChIKey	HEDRZPFGACZZDS-UHFFFAOYSA-N

cyclohexane

Compound Data

Molecular weight	84.1595
Predicted density	0.79 g/cm^3
Boiling point	80.7 °C
Dipole moment	0
Dielectric constant	18.5
SMILES	C1CCCCC1
InChIKey	XDTMQSROBMDMFD-UHFFFAOYSA-N

cyclooctane

Compound Data

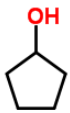
Molecular weight	112.2126
Predicted density	0.79 g/cm^3
Boiling point	152 °C
SMILES	C1CCCCCCC1
InChIKey	WJTCGQSWYFHTAC-UHFFFAOYSA-N

cyclopentane

Compound Data

Molecular weight	70.1329
Predicted density	0.79 g/cm^3
Boiling point	49.2 °C
SMILES	C1CCCC1
InChIKey	RGSFGYAAUTVSQA-UHFFFAOYSA-N

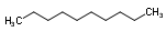
cyclopentanol



Compound Data

Molecular weight	86.1323
Predicted density	1.004 g/cm^3
Boiling point	140.8 °C
SMILES	OC1CCCC1
InChIKey	XCIXKGXIYUWCLL-UHFFFAOYSA-N

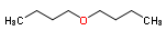
decane



Compound Data

Molecular weight	142.2817
Predicted density	0.734 g/cm^3
Boiling point	174.9 °C
SMILES	C(CCCCCCCC)C
InChIKey	DIOQZVSQGTUSAI-UHFFFAOYSA-N

dibutyl ether



Compound Data

Molecular weight	130.2279
Predicted density	0.78 g/cm^3
Boiling point	142.1 °C
SMILES	O(CCCC)CCCC
InChIKey	DURPTKYDGMDSBL-UHFFFAOYSA-N

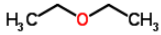
dichloromethane

Compound Data

Molecular weight	84.9326
Predicted density	1.252 g/cm^3
Boiling point	39.6 °C

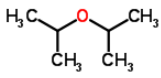
Dipole moment	1.6
Dielectric constant	9.08
SMILES	C(Cl)Cl
InChIKey	YMWUJEATGCHHMB-UHFFFAOYSA-N

diethyl ether



Compound Data	
Molecular weight	74.1216
Predicted density	0.734 g/cm^3
Boiling point	33.2 °C
Dipole moment	1.15
Dielectric constant	4.34
SMILES	CCOCC
InChIKey	RTZKZFJDLAIYFH-UHFFFAOYSA-N

diisopropyl ether

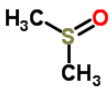


Compound Data	
Molecular weight	102.1748
Predicted density	0.758 g/cm^3
Boiling point	68.3 °C
SMILES	O(C(C)C)C(C)C
InChIKey	ZAFNJMIOTHYJRJ-UHFFFAOYSA-N

DMF

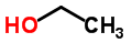
Compound Data	
Molecular weight	73.0938
Predicted density	0.87 g/cm^3
Boiling point	153 °C
Dipole moment	3.82
Dielectric constant	38.3
SMILES	O=C(NC)C
InChIKey	OHLUUHNLEMFGTQ-UHFFFAOYSA-N

DMSO



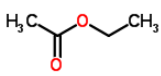
Compound Data	
Molecular weight	78.1334
Predicted density	1.099 g/cm^3
Boiling point	189 °C
Dipole moment	3.96
Dielectric constant	47.2
SMILES	CS(=O)C
InChIKey	IAZDPXIOMUYVGZ-UHFFFAOYSA-N

ethanol



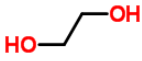
Compound Data	
Molecular weight	46.0684
Predicted density	0.78 g/cm^3
Boiling point	72.6 °C
Dipole moment	1.69
Dielectric constant	24.3
SMILES	CCO
InChIKey	LFQSCWFLJHTTHZ-UHFFFAOYSA-N

ethyl acetate



Compound Data	
Molecular weight	88.1051
Predicted density	0.898 g/cm^3
Boiling point	73.9 °C
Dipole moment	1.78
Dielectric constant	6.02
SMILES	CCOC(=O)C1=NN(C(=N1)C(Cl)(Cl)Cl)C2=C(C=C(C=C2)Cl)Cl
InChIKey	GMBRUAIJEFRHFQ-UHFFFAOYSA-N

ethylene glycol

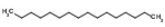


Compound Data	
Molecular weight	62.0678
Predicted density	1.097 g/cm^3
Boiling point	197.5 °C
SMILES	OCCO
InChIKey	LYCAIKOWRPUZTN-UHFFFAOYSA-N

heptane

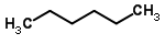
Compound Data	
Molecular weight	100.2019
Predicted density	0.695 g/cm^3
Boiling point	98.8 °C
SMILES	CCCCCCC
InChIKey	IMNFDUFMRHMDMM-UHFFFAOYSA-N

hexadecane



Compound Data	
Molecular weight	226.4412
Predicted density	0.773 g/cm^3
Boiling point	286.6 °C
SMILES	CCCCCCCCCCCCCCCC
InChIKey	DCAYPVUWAIABOU-UHFFFAOYSA-N

hexane

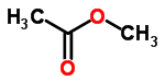


Compound Data	
Molecular weight	86.1754
Predicted density	0.675 g/cm^3
Boiling point	68.5 °C
Dipole moment	0.08
Dielectric constant	2.02
SMILES	CCCCCC
InChIKey	VLKZOEOYAKHREP-UHFFFAOYSA-N

methanol

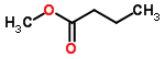
Compound Data	
Molecular weight	32.0419
Predicted density	0.753 g/cm^3
Boiling point	48.1 °C
Dipole moment	1.7
Dielectric constant	33
SMILES	CO
InChIKey	OKKJLVBELUTLKV-UHFFFAOYSA-N

methyl acetate



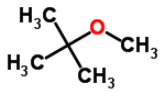
Compound Data	
Molecular weight	74.0785
Predicted density	0.908 g/cm^3
Boiling point	44 °C
SMILES	O=C(OC)C
InChIKey	KXKVLQRXCPHEJC-UHFFFAOYSA-N

methyl butyrate



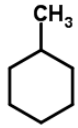
Compound Data	
Molecular weight	102.1317
Predicted density	0.891 g/cm^3
Boiling point	104.2 °C
SMILES	O=C(OC)CCC
InChIKey	UUIQMZJEGPQKFD-UHFFFAOYSA-N

methyl tert-butyl ether



Compound Data	
Molecular weight	88.1482
Predicted density	0.75 g/cm^3
Boiling point	55.2 °C
SMILES	O(C(C)(C)C)C
InChIKey	BZLVMXJERCGZMT-UHFFFAOYSA-N

methylcyclohexane



Compound Data

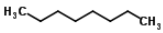
Molecular weight	98.1861
Predicted density	0.776 g/cm^3
Boiling point	101.1 °C
SMILES	CC1CCCCC1
InChIKey	UAEPNZWRGJTJPN-UHFFFAOYSA-N

nonane

Compound Data

Molecular weight	128.2551
Predicted density	0.724 g/cm^3
Boiling point	151.7 °C
SMILES	CCCCCCCCC
InChIKey	BKIMMITUMNQMOS-UHFFFAOYSA-N

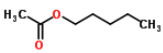
octane



Compound Data

Molecular weight	114.2285
Predicted density	0.711 g/cm^3
Boiling point	126.4 °C
SMILES	CCCCCCCC
InChIKey	TVMXDCGIABBOFY-UHFFFAOYSA-N

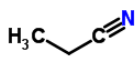
pentyl acetate



Compound Data

Molecular weight	130.1849
Predicted density	0.882 g/cm^3
Boiling point	149.9 °C
SMILES	O=C(OCCCCC)C
InChIKey	PGMYKACGEOXYJE-UHFFFAOYSA-N

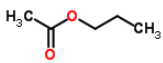
propionitrile



Compound Data

Molecular weight	55.0785
Predicted density	0.771 g/cm^3
Boiling point	91.3 °C
SMILES	N#CCC
InChIKey	FVSKHRXBFJPNKK-UHFFFAOYSA-N

propyl acetate

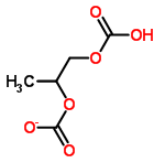


Compound Data

Molecular weight	102.1317
Predicted density	0.891 g/cm^3
Boiling point	101.4 °C
SMILES	O=C(OCCC)C

InChIKey	YKYONYBAUNKHLG-UHFFFAOYSA-N

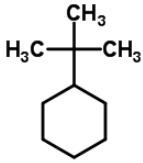
propylene carbonate



Compound Data

Molecular weight	163.106
Predicted density	1.205 g/cm^3
Boiling point	332.3 °C
SMILES	OC(=O)OCC(C)OC(=O)[O-]
InChIKey	ZEBXBLIKXVICMJ-UHFFFAOYSA-M

tert-butylcyclohexane



Compound Data

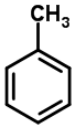
Molecular weight	140.2658
Predicted density	0.812 g/cm^3
Boiling point	168.3 °C
SMILES	CC(C)(C)C1CCCCC1
InChIKey	XTVMZZBLCLWBPM-UHFFFAOYSA-N

THF

Compound Data

Molecular weight	72.1057
Predicted density	0.904 g/cm^3
Boiling point	68.3 °C
Dipole moment	1.63
Dielectric constant	7.52
SMILES	C1CCOC1
InChIKey	WYURNTSHIVDZCO-UHFFFAOYSA-N

toluene



Compound Data

Molecular weight	92.1384
Predicted density	0.871 g/cm^3
Boiling point	110.6 °C
Dipole moment	0.36
Dielectric constant	2.38
SMILES	c1ccccc1C
InChIKey	YXFVVABEGXRONW-UHFFFAOYSA-N

undecane

Compound Data

Molecular weight	156.3083
Predicted density	0.743 g/cm^3
Boiling point	196.3 °C
SMILES	C(CCCCCCCC)CC
InChIKey	RSJKGSCJYJTIGS-UHFFFAOYSA-N

water

Compound Data

Molecular weight	18.0153
Predicted density	0.998 g/cm^3
Boiling point	100 °C
SMILES	O
InChIKey	XLYOFNOQVPJJNP-UHFFFAOYSA-N

Table 1. Abraham solubility coefficients

Solvent	c	e	s	a	b	v
1-butanol	0.152	0.438	-1.177	0.096	-3.919	4.122
1-decanol	-0.062	0.754	-1.461	0.063	-4.053	4.293
1-heptanol	-0.026	0.491	-1.258	0.035	-4.155	4.415
1-hexanol	0.044	0.47	-1.153	0.083	-4.057	4.249
1-octanol	-0.034	0.489	-1.044	-0.024	-4.235	4.218
1-pentanol	0.08	0.521	-1.294	0.208	-3.908	4.208
1-propanol	0.148	0.436	-1.098	0.389	-3.893	4.036
1,2-dichloroethane	0.183	0.294	-0.134	-2.801	-4.291	4.18
1,4-dioxane	0.098	0.35	-0.083	-0.556	-4.826	4.172
2-butanol	0.194	0.383	-0.956	0.134	-3.606	3.829
2-methyl-1-propanol	0.161	0.31	-1.069	0.183	-3.774	4.04
2-methyl-2-propanol	0.197	0.136	-0.916	0.318	-4.031	4.112
2-pentanol	0.115	0.455	-1.331	0.206	-3.745	4.201
2-propanol	0.102	0.315	-1.02	0.532	-3.865	4.023
2,2,4-trimethylpentane	0.318	0.555	-1.737	-3.677	-4.864	4.417
3-methyl-1-butanol	0.123	0.37	-1.243	0.074	-3.781	4.208
acetone	0.313	0.312	-0.121	-0.608	-4.753	3.942
acetonitrile	0.413	0.077	0.326	-1.566	-4.391	3.364
benzene	0.142	0.464	-0.588	-3.009	-4.625	4.491
butanone	0.246	0.256	-0.08	-0.767	-4.855	4.148
butyl acetate	0.248	0.356	-0.501	-0.867	-4.973	4.281
carbon disulfide	0.047	0.686	-0.943	-3.603	-5.818	4.921
carbon tetrachloride	0.199	0.523	-1.159	-3.56	-4.594	4.618
chloroform	0.191	0.105	-0.403	-3.112	-3.514	4.395
cyclohexane	0.159	0.784	-1.678	-3.74	-4.929	4.577
decane	0.16	0.585	-1.734	-3.435	-5.078	4.582
dibutyl ether	0.203	0.369	-0.954	-1.488	-5.426	4.508
dichloromethane	0.319	0.102	-0.187	-3.058	-4.09	4.324
diethyl ether	0.33	0.401	-0.814	-0.457	-4.959	4.32
DMF	-0.305	-0.058	0.343	0.358	-4.865	4.486
DMSO	-0.194	0.327	0.791	1.26	-4.54	3.361
ethanol	0.208	0.409	-0.959	0.186	-3.645	3.928
ethyl acetate	0.328	0.369	-0.446	-0.7	-4.904	4.15
ethylene glycol	-0.243	0.695	-0.67	0.726	-2.399	2.67
heptane	0.325	0.67	-2.061	-3.317	-4.733	4.543
hexadecane	0.087	0.667	-1.617	-3.587	-4.869	4.433
hexane	0.361	0.579	-1.723	-3.599	-4.764	4.344
methanol	0.329	0.299	-0.671	0.08	-3.389	3.512
methyl acetate	0.351	0.223	-0.15	-1.035	-4.527	3.972
methyl tert-butyl ether	0.376	0.264	-0.788	-1.078	-5.03	4.41
methylcyclohexane	0.246	0.782	-1.982	-3.517	-4.293	4.528
nonane	0.24	0.619	-1.713	-3.532	-4.921	4.482
octane	0.223	0.642	-1.647	-3.48	-5.067	4.526
THF	0.207	0.372	-0.392	-0.236	-4.934	4.447
toluene	0.143	0.527	-0.72	-3.01	-4.824	4.545
undecane	0.058	0.603	-1.661	-3.421	-5.12	4.619

water	0	0	0	0	0	0

Table 2. Hansen solubility parameters

Solvent	δ_d	δ_p	δ_h
1-butanol	16	5.7	15.8
1-decanol	17.5	2.6	10
1-heptanol	15.1	8	13
1-hexanol	14.1	8.6	12.7
1-octanol	17	3.3	11.9
1-pentanol	15.9	5.9	13.9
1-propanol	16	6.8	17.4
1,4-dioxane	19	1.8	7.4
2-butanol	15.8	5.7	14.5
2-methyl-1-butanol	16	5.1	14.3
2-methyl-1-propanol	15.1	5.7	15.9
2-methyl-2-propanol	15.2	5.1	14.7
2-pentanol	15.6	6.4	13.3
2-propanol	15.8	6.1	16.4
2,2,4-trimethylpentane	14.1	0	0
3-methyl-1-butanol	15.8	5.2	13.3
acetone	15.5	10.4	7
acetonitrile	15.3	18	6.1
benzene	18.4	0	2
butyl acetate	15.8	3.7	6.3
carbon tetrachloride	17.8	0	0.6
chloroform	17.8	3.1	5.7
cyclohexane	16.8	0	0.2
cyclooctane	17.5	0	0
cyclopentane	16.4	0	1.8
decane	15.7	0	0
dibutyl ether	15.7	4.5	4.2
dichloromethane	18.2	6.3	6.1
diethyl ether	14.5	2.9	5.1
DMF	17.4	16.7	11.3
DMSO	18.4	16.4	10.2
ethanol	15.8	8.8	19.4
ethyl acetate	15.8	5.3	7.2
ethylene glycol	17	11	26
heptane	15.3	0	0
hexadecane	16.3	0	0
hexane	14.9	0	0
methanol	15.1	12.3	22.3
methyl acetate	15.5	7.2	7.6
nonane	15.7	0	0
octane	15.5	0	0
propylene carbonate	20	18	4.1
THF	16.8	5.7	8
toluene	18	1.4	2
undecane	16	0	0

Solubility Modeling - Abraham General Solvation Model

Abraham Descriptors

Abraham *et al.*, have developed equations, the *Abraham solvation equations,* that can be used to predict the solubility of organic compounds in 84 different organic solvents [1]. The Abraham method assumes that the partition coefficient between water and a solvent, P_s, is given by the ratio of solubilities of a solute in the solvent, S_s, and in water, S_w,

P_s = S_s / S_w. (1)

If this assumption is reasonable for the solute and solvent in question, then the solubility of the solute in the solvent can be calculated from the predicted partition coefficient,

log S_s = log S_w + c + e E + s S + a A + b B + v V, (2)

where the continually refined (as more experimental data becomes available) coefficients c, e, s, b, and v vary by solvent, see Table 1, and E, S, A, B, and V are solute descriptors, described as follows:

E is the solute ***excess molar refractivity*** *in units of (cubic cm per mol)/10.*
S is the solute ***dipolarity/polarizability.***
A is the ***overall (summation) hydrogen bond acidity*** .
B is the ***overall (summation) hydrogen bond basicity*** .
V is the ***McGowan characteristic volume*** *in units of (cubic cm per mol)/100.*

Descriptor Calculation Methods

Various methods are used to determine the solute descriptors either from first principles or by working backwards from known logP and solubility values. Abraham *et al.* [2] have published a review of methods. The following sections detail the method used in this book to calculate the Abraham solute descriptors.

log S_w - The logarithm of the molar solubility in water

If the water solubility is known, then log S_w is simply the logarithm of this value. If the water solubility is not known, it can be approximated using the free web-service provided by VCCLab [http://www.vcclab.org/web/alogps/]. If the web service fails, the value can be approximated using the 1-octanol solubility via equation (1). If neither the water or the 1-octanol solubility is known, then log S_w is calculated using linear regression from all known solubility data, see S, A, and B below. If log S_w is predicted rather than measured, it will be italicized when presented.

V - The McGowan characteristic volume

V is the McGowan approximation for the molecular volume. It can be calculated from fragment contributions from the solute's molecular formula and the number of bonds. We calculated the McGowan volume from predicted values for the molar volume found on ChemSpider by using

V (intrinsic) = 0.597 + 0.6823 V. *(3)* [3]

E - The excess molar refractivity

E is the molar refraction, MR, of the compound calculated using McGowan's volume less the molar refraction of an alkane with the same McGowan volume,

E = (MR) − 2.83195 V + 0.52553, *(4)*

where

MR = 10 (η^2 −1)/(η^2 + 2) V where η is the refractive index. *(5)*

We calculated the E from predicted values for MR (intrinsic) found on ChemSpider and by using

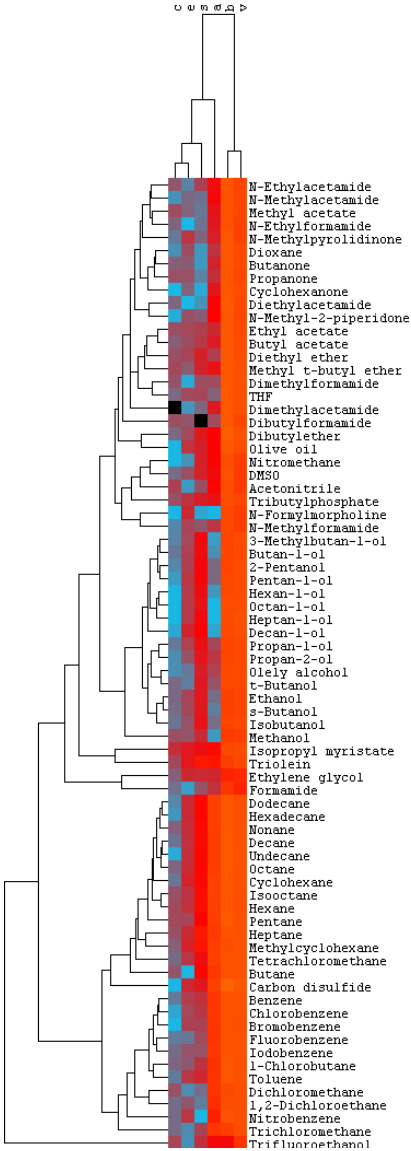
MR (intrinsic) = V (intrinsic) * MR / V *(6)*.

S, A, and B

S, A, and B can be calculated from GLC [2], RP-HPLC [4], or simple linear regression from known solubility data or known logP values. We caluclated S, A, and B by performing linear regression on the solutes with at least five solubility values in the Open Notebook Science Challenge database. When using the descriptors to predict the solubility of the solute in other organic solvents it is important to note that the more solvents used in the regression analysis, the more accurate the descriptors are likely to be. Solubility values for carboxylic acids in toluene or benzene were not used in the regression analysis since they are known to dimerize. Similarly, care must be taken when using Abraham descriptors to predict the solubility of carboxylic acids in toluene and benzene [5].

Clustering

By giving each solvent descriptors (c, e, s, a, b, v) it becomes possible to determine which solvents are similar by comparing their descriptors - much like as is done with Hansen solubility parameters . Using recently updated values [1], we used HappieClust to perform a cluster analysis (agglomerative hierarchical clustering) to see which solvents are considered similar in the Abraham general solvation method. The results are depicted in the figure below.



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Summary of Results

A summary of the results of the regression analyses giving the Abraham solute descriptors are presented below in Table 3.

Table 3. Abraham solute descriptors

Solute	E	S	A	B	V
1-octadecylamine	0.011	1.290	0.034	1.979	3.951
2-chloro-5-nitrobenzaldehyde	1.221	1.359	-0.054	0.182	0.954
2-chloro-5-nitrobenzoic acid	1.220	1.109	0.773	0.178	0.967
2-methoxybenzoic acid	0.850	1.328	0.422	0.477	0.970
2-methylbenzoic acid	0.854	0.689	0.368	0.284	0.857
2-phenylbutyric acid	0.878	1.077	0.009	0.583	1.332
4-acetamidobenzoic acid	1.279	1.112	0.793	0.734	1.104
4-chloro-3-nitrobenzoic acid	1.220	1.317	0.583	0.371	0.967
4-chlorobenzaldehyde	0.933	0.716	0.080	0.087	0.781
4-chlorophenylacetic acid	0.980	0.565	0.265	0.257	1.013
4-dimethylaminobenzaldehyde	1.190	1.602	0.079	0.756	1.162
4-fluorobenzoic acid	0.726	0.720	0.400	0.168	0.680
4-hydroxybenzaldehyde	0.918	1.003	0.439	0.225	0.583
4-methoxybenzoic acid	0.850	1.290	0.404	0.552	0.970
4-methoxyphenylacetic acid	0.877	1.137	0.101	0.678	1.189
4-nitrobenzaldehyde	1.049	1.634	-0.031	0.289	0.780
4-nitrobenzoic acid	1.048	1.301	0.518	0.349	0.793
4-phenylbutyric acid	0.896	3.980	0.540	-0.076	1.321
4-pyrenebutanoic acid	3.727	2.030	0.411	0.855	2.385
9-fluorenone	1.638	1.392	0.097	0.233	1.247
acetylsalicylic acid	0.946	5.432	0.815	-0.028	1.170
benzoic acid	0.788	0.144	0.100	0.239	0.618
boc-glycine	0.402	0.823	0.255	0.784	1.338
chloroacetic acid	0.502	0.372	0.096	0.042	0.114
cinnamic acid	1.161	4.115	0.445	-0.194	0.957
crotonic acid	0.474	0.405	0.130	0.155	0.339
diphenylacetic acid	1.565	1.499	0.147	1.083	1.773
furoic acid	0.590	0.351	0.479	0.125	0.366
ibuprofen	0.942	1.122	0.444	1.364	2.061
mandelic acid	0.970	6.447	1.008	-0.527	0.812
nicotinic acid	0.762	0.738	0.461	0.681	0.519
o-vanillin	1.024	0.802	0.046	0.448	0.935
phenanthrene	2.100	0.840	0.057	0.198	1.435
phenanthrene-9-carboxaldehyde	2.494	1.166	0.327	0.423	1.606
phenylacetic acid	0.832	0.917	0.082	0.455	0.837
piperonal	1.032	0.928	0.118	0.175	0.769
salicylic acid	0.920	0.110	0.410	0.111	0.595
thianthrene	2.281	1.312	-0.101	0.786	1.543
UCExp216-3A	2.332	4.791	-0.007	3.981	5.203
Ugi product 148B	3.440	5.683	0.240	3.436	5.615
Ugi product 150D(UC)	4.342	4.362	0.051	3.236	5.052
Ugi product 176C (UC)	4.713	4.490	0.151	3.935	5.298
Ugi product 206B(UC)	1.915	2.463	0.301	3.663	4.800
Ugi product 215F(UC)	5.307	4.975	0.047	3.742	5.726
Ugi Product 216-4A	1.716	4.742	-0.317	4.359	5.883
vanillin	1.024	1.086	0.365	0.411	0.935
veratraldehyde	0.889	1.091	0.204	0.572	1.310
xanthene	1.517	1.171	-0.085	0.454	1.428

Measured and predicted values

1-octadecylamine $C_{18}H_{39}N$ - 6 solvents [69, 908]

Compound Data

Molecular weight	269.509	**H bond acceptors**	1	**Rule of 5 violations**	1
Compound type	amine	**H bond donors**	2	**ACD/ALogP**	8.37
Phase 25°C	solid	**Rotatable bonds**	17	**Predicted density**	0.818 g/cm^3
SMILES	NCCCCCCCCCCCCCCCCCC				
InChIKey	REYJJPSVUYRZGE-UHFFFAOYSA-N				

Abraham Solute Descriptors

log S_w	E	S	A	B	V
-6.910	0.011	1.290	0.034	1.979	3.951

Solubility Data

	Concentration (M)		
Solvent	**Measured**	**Predicted**	**Solvent bp °C**
acetonitrile	0.08	0.03	63.5
ethanol	1.77	2.39	72.6
hexane	0.13	0.07	68.5
methanol	1.68	0.54	48.1
THF	1.68	3.91	68.3
toluene	1.34	4.14	110.6
1-butanol	-	1.83	117.7
1-decanol	-	1.24	227.8
1-heptanol	-	4.66	176.9
1-hexanol	-	2.59	158.2
1-octanol	-	1.00	194.7
1-pentanol	-	2.55	138.5
1-propanol	-	1.21	95.8
1,2-dichloroethane	-	> 9	83.5
1,4-dioxane	-	1.00	102.9
2-butanol	-	1.13	96.6
2-methyl-1-propanol	-	2.37	105.0
2-methyl-2-propanol	-	2.44	84.6
2-pentanol	-	4.87	118.8
2-propanol	-	1.39	73.0
2,2,4-trimethylpentane	-	0.07	98.8
3-methyl-1-butanol	-	5.75	131.2
acetone	-	0.25	46.5
benzene	-	> 9	78.8
butanone	-	0.98	75.6
butyl acetate	-	0.55	126.6
carbon disulfide	-	0.54	46.2
carbon tetrachloride	-	6.78	76.0
chloroform	-	> 9	61.2
cyclohexane	-	0.20	80.7
decane	-	0.09	174.9
dibutyl ether	-	0.12	142.1
dichloromethane	-	> 9	39.6
diethyl ether	-	0.41	33.2
DMF	-	> 9	153.0
DMSO	-	0.02	189.0
ethyl acetate	-	0.33	73.9
ethylene glycol	-	0.01	197.5
heptane	-	0.17	98.8
hexadecane	-	0.07	286.6
methyl acetate	-	0.89	44.0
methyl tert-butyl ether	-	0.77	55.2
methylcyclohexane	-	1.15	101.1
nonane	-	0.09	151.7
octane	-	0.09	126.4

undecane	-	0.10	196.3
water	-	0.00	100.0

2-chloro-5-nitrobenzaldehyde $C_7H_4ClNO_3$ - 5 solvents [105, 69, 208]

Compound Data

Molecular weight	185.565	**H bond acceptors**	4	**Rule of 5 violations**	0
Compound type	aldehyde	**H bond donors**	0	**ACD/ALogP**	2.5
Phase 25°C	solid	**Rotatable bonds**	2	**Predicted density**	1.485 g/cm^3
SMILES	O=Cc1cc(ccc1Cl)[N+]([O-])=O				
InChIKey	VFVHWCKUHAEDMY-UHFFFAOYSA-N				

Abraham Solute Descriptors

log S_w	E	S	A	B	V
-3.100	1.221	1.359	-0.054	0.182	0.954

Solubility Data

	Concentration (M)		
Solvent	**Measured**	**Predicted**	**Solvent bp °C**
acetonitrile	2.31	2.23	63.5
chloroform	2.90	2.49	61.2
DMSO	2.93	3.12	189
THF	2.79	2.45	68.3
toluene	1.74	2.15	110.6
1-butanol	-	0.16	117.7
1-decanol	-	0.13	227.8
1-heptanol	-	0.17	176.9
1-hexanol	-	0.18	158.2
1-octanol	-	0.20	194.7
1-pentanol	-	0.14	138.5
1-propanol	-	0.16	95.8
1,2-dichloroethane	-	4.18	83.5
1,4-dioxane	-	2.80	102.9
2-butanol	-	0.18	96.6
2-methyl-1-propanol	-	0.14	105.0
2-methyl-2-propanol	-	0.16	84.6
2-pentanol	-	0.12	118.8
2-propanol	-	0.13	73.0
2,2,4-trimethylpentane	-	0.12	98.8
3-methyl-1-butanol	-	0.13	131.2
acetone	-	2.29	46.5
benzene	-	2.62	78.8
butanone	-	2.94	75.6
butyl acetate	-	1.35	126.6
carbon disulfide	-	2.18	46.2
carbon tetrachloride	-	0.85	76.0
cyclohexane	-	0.26	80.7
decane	-	0.11	174.9
dibutyl ether	-	0.45	142.1
dichloromethane	-	4.35	39.6
diethyl ether	-	0.72	33.2
DMF	-	2.33	153.0
ethanol	-	0.24	72.6
ethyl acetate	-	1.51	73.9
ethylene glycol	-	0.05	197.5
heptane	-	0.08	98.8
hexadecane	-	0.14	286.6
hexane	-	0.13	68.5
methanol	-	0.26	48.1
methyl acetate	-	2.21	44.0
methyl tert-butyl ether	-	0.76	55.2
methylcyclohexane	-	0.14	101.1

nonane	-	0.14	151.7
octane	-	0.18	126.4
undecane	-	0.13	196.3
water	-	0.00	100.0

2-chloro-5-nitrobenzoic acid $C_7H_4ClNO_4$ - 22 solvents [905]

Compound Data

Molecular weight	201.564	**H bond acceptors**	5	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	2.02
Phase 25°C	solid	**Rotatable bonds**	2	**Predicted density**	1.602 g/cm^3
SMILES	O=C(O)c1cc(ccc1Cl)[N+]([O-])=O				
InChIKey	QUEKGYQTRJVEQC-UHFFFAOYSA-N				

Abraham Solute Descriptors

log S_w	E	S	A	B	V
-2.940	1.220	1.109	0.773	0.178	0.967

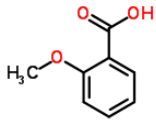
Solubility Data

	Concentration (M)		
Solvent	**Measured**	**Predicted**	**Solvent bp °C**
1-butanol	0.78	0.64	117.7
1-decanol	0.39	0.60	227.8
1-heptanol	0.54	0.63	176.9
1-hexanol	0.64	0.71	158.2
1-octanol	0.48	0.59	194.7
1-pentanol	0.74	0.76	138.5
1-propanol	0.95	1.09	95.8
2-butanol	0.91	0.67	96.6
2-methyl-1-propanol	0.63	0.62	105
2-methyl-2-propanol	1.21	0.82	84.6
2-pentanol	0.76	0.65	118.8
2-propanol	1.13	1.08	73
3-methyl-1-butanol	0.68	0.52	131.2
butyl acetate	0.46	0.59	126.6
dibutyl ether	0.10	0.08	142.1
diethyl ether	0.57	0.83	33.2
diisopropyl ether	0.19	*	68.3
ethanol	1.40	1.00	72.6
ethyl acetate	0.75	0.88	73.9
methyl acetate	0.89	0.57	44
pentyl acetate	0.43	*	149.9
THF	2.97	3.38	68.3
1,2-dichloroethane	-	0.04	83.5
1,4-dioxane	-	1.74	102.9
2,2,4-trimethylpentane	-	0.00	98.8
acetone	-	1.31	46.5
acetonitrile	-	0.16	63.5
benzene	-	0.02	78.8
butanone	-	1.22	75.6
carbon disulfide	-	0.01	46.2
carbon tetrachloride	-	0.00	76.0
chloroform	-	0.01	61.2
cyclohexane	-	0.00	80.7
decane	-	0.00	174.9
dichloromethane	-	0.02	39.6
DMF	-	6.55	153.0
DMSO	-	> 9	189.0
ethylene glycol	-	0.44	197.5
heptane	-	0.00	98.8
hexadecane	-	0.00	286.6

hexane	-	0.00	68.5
methanol	-	0.74	48.1
methyl tert-butyl ether	-	0.27	55.2
methylcyclohexane	-	0.00	101.1
nonane	-	0.00	151.7
octane	-	0.00	126.4
toluene	-	0.02	110.6
undecane	-	0.00	196.3
water	-	0.00	100.0

* This solvent doesn't have recorded Abraham coefficients.

2-methoxybenzoic acid $C_8H_8O_3$ - 31 solvents [905, 34, 907]



Compound Data

Molecular weight	152.147	**H bond acceptors**	3	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	1.5
Phase 25°C	solid	**Rotatable bonds**	2	**Predicted density**	1.207 g/cm^3
SMILES	O=C(O)c1ccccc1OC				
InChIKey	ILUJQPXNXACGAN-UHFFFAOYSA-N				

Abraham Solute Descriptors

log S_w	E	S	A	B	V
-1.523	0.850	1.328	0.422	0.477	0.970

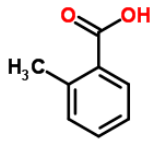
Solubility Data

	Concentration (M)		
Solvent	**Measured**	**Predicted**	**Solvent bp °C**
1-butanol	0.57	0.40	117.7
1-decanol	0.22	0.24	227.8
1-heptanol	0.34	0.33	176.9
1-hexanol	0.37	0.41	158.2
1-octanol	0.30	0.34	194.7
1-pentanol	0.47	0.39	138.5
1-propanol	0.76	0.57	95.8
1,4-dioxane	1.53	1.87	102.9
2-butanol	0.50	0.60	96.6
2-ethyl-1-hexanol	0.25	*	184.6
2-methyl-1-butanol	0.39	*	128.7
2-methyl-1-pentanol	0.35	*	148
2-methyl-1-propanol	0.41	0.48	105
2-methyl-2-propanol	0.53	0.59	84.6
2-pentanol	0.39	0.39	118.8
2-propanol	0.59	0.60	73
3-methyl-1-butanol	0.36	0.37	131.2
4-methyl-2-pentanol	0.32	*	133.5
butyl acetate	0.34	0.60	126.6
chloroform	0.46	0.31	61.2
dibutyl ether	0.04	0.08	142.1
diethyl ether	0.24	0.50	33.2
diisopropyl ether	0.06	*	68.3
ethanol	1.19	0.81	72.6
ethyl acetate	5.63	0.82	73.9
methanol	1.82	0.98	48.1
methyl acetate	0.86	1.19	44
pentyl acetate	0.26	*	149.9
propylene carbonate	0.99	*	332.3
THF	2.02	2.18	68.3
water	0.03	0.03	100
1,2-dichloroethane	-	0.36	83.5
2,2,4-trimethylpentane	-	0.00	98.8
acetone	-	1.56	46.5

acetonitrile	-	0.79	63.5
benzene	-	0.13	78.8
butanone	-	1.65	75.6
carbon disulfide	-	0.02	46.2
carbon tetrachloride	-	0.02	76.0
cyclohexane	-	0.00	80.7
decane	-	0.00	174.9
dichloromethane	-	0.39	39.6
DMF	-	5.74	153.0
DMSO	-	> 9	189.0
ethylene glycol	-	0.49	197.5
heptane	-	0.00	98.8
hexadecane	-	0.00	286.6
hexane	-	0.00	68.5
methyl tert-butyl ether	-	0.28	55.2
methylcyclohexane	-	0.00	101.1
nonane	-	0.00	151.7
octane	-	0.00	126.4
toluene	-	0.09	110.6
undecane	-	0.00	196.3

* This solvent doesn't have recorded Abraham coefficients.

2-methylbenzoic acid $C_8H_8O_2$ - 27 solvents [905, 907]



Compound Data

Molecular weight	136.148	**H bond acceptors**	2	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	2.36
Phase 25°C	solid	**Rotatable bonds**	1	**Predicted density**	1.151 g/cm^3
SMILES	O=C(O)c1ccccc1C				
InChIKey	ZWLPBLYKEWSWPD-UHFFFAOYSA-N				

Abraham Solute Descriptors

log S_w	E	S	A	B	V
-2.000	0.854	0.689	0.368	0.284	0.857

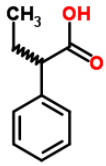
Solubility Data

	Concentration (M)		
Solvent	**Measured**	**Predicted**	**Solvent bp °C**
1-butanol	1.61	1.48	117.7
1-decanol	1.04	1.34	227.8
1-heptanol	1.28	1.39	176.9
1-hexanol	1.38	1.49	158.2
1-octanol	1.16	1.17	194.7
1-pentanol	1.49	1.61	138.5
1-propanol	1.77	1.82	95.8
1,4-dioxane	2.64	2.19	102.9
2-butanol	1.75	1.48	96.6
2-methyl-1-butanol	1.25	*	128.7
2-methyl-1-propanol	1.28	1.41	105
2-methyl-2-propanol	2.19	1.51	84.6
2-pentanol	1.68	1.59	118.8
2-propanol	1.95	1.64	73
3-methyl-1-butanol	1.34	1.39	131.2
4-methyl-2-pentanol	1.39	*	133.5
butyl acetate	1.17	1.40	126.6
dibutyl ether	0.58	0.43	142.1
diethyl ether	1.65	1.73	33.2
diisopropyl ether	0.86	*	68.3
ethanol	2.08	1.98	72.6
ethyl acetate	1.49	1.75	73.9
methyl acetate	1.51	1.50	44

pentyl acetate	1.04	*	149.9
propylene carbonate	0.46	*	332.3
THF	2.92	3.79	68.3
water	0.01	0.01	100
1,2-dichloroethane	-	0.47	83.5
2,2,4-trimethylpentane	-	0.04	98.8
acetone	-	2.00	46.5
acetonitrile	-	0.58	63.5
benzene	-	0.37	78.8
butanone	-	2.01	75.6
carbon disulfide	-	0.17	46.2
carbon tetrachloride	-	0.16	76.0
chloroform	-	0.43	61.2
cyclohexane	-	0.07	80.7
decane	-	0.05	174.9
dichloromethane	-	0.50	39.6
DMF	-	3.00	153.0
DMSO	-	4.83	189.0
ethylene glycol	-	0.58	197.5
heptane	-	0.06	98.8
hexadecane	-	0.04	286.6
hexane	-	0.05	68.5
methanol	-	1.58	48.1
methyl tert-butyl ether	-	1.03	55.2
methylcyclohexane	-	0.08	101.1
nonane	-	0.05	151.7
octane	-	0.06	126.4
toluene	-	0.33	110.6
undecane	-	0.05	196.3

* This solvent doesn't have recorded Abraham coefficients.

2-phenylbutyric acid $C_{10}H_{12}O_2$ - 5 solvents [79, 83]



Compound Data

Molecular weight	164.201	**H bond acceptors**	2	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	2.38
Phase 25°C	solid	**Rotatable bonds**	3	**Predicted density**	1.09 g/cm^3
SMILES	O=C(O)C(c1ccccc1)CC				
InChIKey	OFJWFSNDPCAWDK-UHFFFAOYSA-N				

Abraham Solute Descriptors

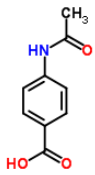
log S_w	E	S	A	B	V
-2.070	0.878	1.077	0.009	0.583	1.332

Solubility Data

	Concentration (M)		
Solvent	**Measured**	**Predicted**	**Solvent bp °C**
acetonitrile	6.18	4.67	63.5
DMSO	5.97	5.21	189
ethanol	6.20	3.75	72.6
THF	5.96	12.21	68.3
toluene	5.35	9.46	110.6
1-butanol	-	2.55	117.7
1-decanol	-	2.06	227.8
1-heptanol	-	2.75	176.9
1-hexanol	-	2.76	158.2
1-octanol	-	2.24	194.7
1-pentanol	-	2.53	138.5
1-propanol	-	2.45	95.8
1,2-dichloroethane	-	> 9	83.5
1,4-dioxane	-	> 9	102.9

2-butanol	-	2.69	96.6
2-methyl-1-propanol	-	2.48	105.0
2-methyl-2-propanol	-	2.46	84.6
2-pentanol	-	2.67	118.8
2-propanol	-	2.09	73.0
2,2,4-trimethylpentane	-	0.76	98.8
3-methyl-1-butanol	-	2.76	131.2
acetone	-	7.27	46.5
benzene	-	> 9	78.8
butanone	-	> 9	75.6
butyl acetate	-	5.59	126.6
carbon disulfide	-	4.96	46.2
carbon tetrachloride	-	6.06	76.0
chloroform	-	> 9	61.2
cyclohexane	-	1.45	80.7
decane	-	0.71	174.9
dibutyl ether	-	1.81	142.1
dichloromethane	-	> 9	39.6
diethyl ether	-	3.94	33.2
DMF	-	> 9	153.0
ethyl acetate	-	5.82	73.9
ethylene glycol	-	0.55	197.5
heptane	-	0.77	98.8
hexadecane	-	0.79	286.6
hexane	-	0.83	68.5
methanol	-	3.18	48.1
methyl acetate	-	> 9	44.0
methyl tert-butyl ether	-	4.19	55.2
methylcyclohexane	-	1.68	101.1
nonane	-	0.87	151.7
octane	-	0.97	126.4
undecane	-	0.74	196.3
water	-	0.01	100.0

4-acetamidobenzoic acid $C_9H_9NO_3$ - 5 solvents [137]



Compound Data

Molecular weight	179.173	**H bond acceptors**	4	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	2	**ACD/ALogP**	1.31
Phase 25°C	solid	**Rotatable bonds**	2	**Predicted density**	1.326 g/cm^3
SMILES	O=C(Nc1ccc(cc1)C(=O)O)C				
InChIKey	QCXJEYYXVJIFCE-UHFFFAOYSA-N				

Abraham Solute Descriptors

log S_w	E	S	A	B	V
-2.120	1.279	1.112	0.793	0.734	1.104

Solubility Data

	Concentration (M)		
Solvent	**Measured**	**Predicted**	**Solvent bp °C**
acetonitrile	0.01	0.01	63.5
benzene	0.00	0.00	78.8
ethanol	0.23	0.22	72.6
methanol	0.19	0.20	48.1
THF	0.17	0.17	68.3
1-butanol	-	0.11	117.7
1-decanol	-	0.09	227.8
1-heptanol	-	0.09	176.9
1-hexanol	-	0.10	158.2
1-octanol	-	0.07	194.7
1-pentanol	-	0.13	138.5
1-propanol	-	0.19	95.8

1,2-dichloroethane	-	0.00	83.5
1,4-dioxane	-	0.09	102.9
2-butanol	-	0.15	96.6
2-methyl-1-propanol	-	0.12	105.0
2-methyl-2-propanol	-	0.12	84.6
2-pentanol	-	0.14	118.8
2-propanol	-	0.19	73.0
2,2,4-trimethylpentane	-	0.00	98.8
3-methyl-1-butanol	-	0.10	131.2
acetone	-	0.07	46.5
butanone	-	0.06	75.6
butyl acetate	-	0.03	126.6
carbon disulfide	-	0.00	46.2
carbon tetrachloride	-	0.00	76.0
chloroform	-	0.00	61.2
cyclohexane	-	0.00	80.7
decane	-	0.00	174.9
dibutyl ether	-	0.00	142.1
dichloromethane	-	0.00	39.6
diethyl ether	-	0.04	33.2
DMF	-	0.35	153.0
DMSO	-	2.29	189.0
ethyl acetate	-	0.04	73.9
ethylene glycol	-	0.35	197.5
heptane	-	0.00	98.8
hexadecane	-	0.00	286.6
hexane	-	0.00	68.5
methyl acetate	-	0.04	44.0
methyl tert-butyl ether	-	0.01	55.2
methylcyclohexane	-	0.00	101.1
nonane	-	0.00	151.7
octane	-	0.00	126.4
toluene	-	0.00	110.6
undecane	-	0.00	196.3
water	-	0.01	100.0

4-chloro-3-nitrobenzoic acid $C_7H_4ClNO_4$ - 30 solvents [905, 907]

Compound Data

Molecular weight	201.564	**H bond acceptors**	5	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	2.37
Phase 25°C	solid	**Rotatable bonds**	2	**Predicted density**	1.602 g/cm^3
SMILES	O=[N+]([O-])c1cc(ccc1Cl)C(=O)O				
InChIKey	DFXQXFGFOLXAPO-UHFFFAOYSA-N				

Abraham Solute Descriptors

log S_w	E	S	A	B	V
-2.310	1.220	1.317	0.583	0.371	0.967

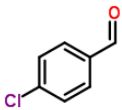
Solubility Data

	Concentration (M)		
Solvent	**Measured**	**Predicted**	**Solvent bp °C**
1-butanol	0.31	0.26	117.7
1-decanol	0.19	0.20	227.8
1-heptanol	0.25	0.23	176.9
1-hexanol	0.27	0.28	158.2
1-octanol	0.23	0.24	194.7
1-pentanol	0.27	0.28	138.5
1-propanol	0.37	0.41	95.8
1,4-dioxane	1.43	1.07	102.9
2-butanol	0.32	0.34	96.6

2-ethyl-1-hexanol	0.15	*	184.6
2-methyl-1-pentanol	0.19	*	148
2-methyl-1-propanol	0.19	0.27	105
2-methyl-2-propanol	0.38	0.33	84.6
2-pentanol	0.30	0.25	118.8
2-propanol	0.39	0.40	73
3-methyl-1-butanol	0.24	0.22	131.2
4-methyl-2-pentanol	0.23	*	133.5
butyl acetate	0.23	0.32	126.6
dibutyl ether	0.05	0.04	142.1
diethyl ether	0.24	0.32	33.2
diisopropyl ether	0.10	*	68.3
ethanol	0.55	0.49	72.6
ethyl acetate	0.36	0.47	73.9
methyl acetate	0.44	0.47	44
methyl butyrate	0.21	*	104.2
pentyl acetate	0.16	*	149.9
propyl acetate	0.26	*	101.4
propylene carbonate	0.15	*	332.3
THF	1.95	1.48	68.3
water	0.00	0.00	100
1,2-dichloroethane	-	0.08	83.5
2,2,4-trimethylpentane	-	0.00	98.8
acetone	-	0.83	46.5
acetonitrile	-	0.22	63.5
benzene	-	0.03	78.8
butanone	-	0.81	75.6
carbon disulfide	-	0.01	46.2
carbon tetrachloride	-	0.00	76.0
chloroform	-	0.04	61.2
cyclohexane	-	0.00	80.7
decane	-	0.00	174.9
dichloromethane	-	0.06	39.6
DMF	-	3.24	153.0
DMSO	-	> 9	189.0
ethylene glycol	-	0.34	197.5
heptane	-	0.00	98.8
hexadecane	-	0.00	286.6
hexane	-	0.00	68.5
methanol	-	0.49	48.1
methyl tert-butyl ether	-	0.13	55.2
methylcyclohexane	-	0.00	101.1
nonane	-	0.00	151.7
octane	-	0.00	126.4
toluene	-	0.02	110.6
undecane	-	0.00	196.3

* This solvent doesn't have recorded Abraham coefficients.

4-chlorobenzaldehyde C_7H_5ClO - 5 solvents [7, 209]



Compound Data

Molecular weight	140.567	**H bond acceptors**	1	**Rule of 5 violations**	0
Compound type	aldehyde	**H bond donors**	0	**ACD/ALogP**	2.21
Phase 25°C	solid	**Rotatable bonds**	1	**Predicted density**	1.243 g/cm^3
SMILES	O=Cc1ccc(Cl)cc1				
InChIKey	AVPYQKSLYISFPO-UHFFFAOYSA-N				

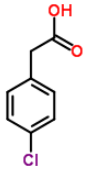
Abraham Solute Descriptors

log S_w	E	S	A	B	V
-2.270	0.933	0.716	0.080	0.087	0.781

Solubility Data

Solvent	Concentration (M) Measured	Predicted	Solvent bp °C
acetonitrile	5.58	3.70	63.5
chloroform	3.61	4.06	61.2
ethanol	2.81	2.50	72.6
methanol	3.55	2.05	48.1
THF	5.29	10.68	68.3
1-butanol	-	2.16	117.7
1-decanol	-	2.14	227.8
1-heptanol	-	2.24	176.9
1-hexanol	-	2.29	158.2
1-octanol	-	2.13	194.7
1-pentanol	-	2.15	138.5
1-propanol	-	2.20	95.8
1,2-dichloroethane	-	5.74	83.5
1,4-dioxane	-	7.75	102.9
2-butanol	-	1.92	96.6
2-methyl-1-propanol	-	1.80	105.0
2-methyl-2-propanol	-	1.92	84.6
2-pentanol	-	1.94	118.8
2-propanol	-	1.75	73.0
2,2,4-trimethylpentane	-	1.13	98.8
3-methyl-1-butanol	-	1.87	131.2
acetone	-	7.31	46.5
benzene	-	5.60	78.8
butanone	-	8.19	75.6
butyl acetate	-	6.21	126.6
carbon disulfide	-	6.18	46.2
carbon tetrachloride	-	3.23	76.0
cyclohexane	-	1.84	80.7
decane	-	1.14	174.9
dibutyl ether	-	3.34	142.1
dichloromethane	-	6.12	39.6
diethyl ether	-	5.72	33.2
DMF	-	5.32	153.0
DMSO	-	5.47	189.0
ethyl acetate	-	6.94	73.9
ethylene glycol	-	0.39	197.5
heptane	-	1.19	98.8
hexadecane	-	1.08	286.6
hexane	-	1.22	68.5
methyl acetate	-	6.42	44.0
methyl tert-butyl ether	-	5.11	55.2
methylcyclohexane	-	1.47	101.1
nonane	-	1.29	151.7
octane	-	1.54	126.4
toluene	-	5.48	110.6
undecane	-	1.12	196.3
water	-	0.01	100.0

4-chlorophenylacetic acid $C_8H_7ClO_2$ - 5 solvents [4, 75, 73]



Compound Data

Molecular weight	170.593	**H bond acceptors**	2	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	2.1
Phase 25°C	solid	**Rotatable bonds**	2	**Predicted density**	1.324 g/cm^3
SMILES	Clc1ccc(cc1)CC(=O)O				
InChIKey	CDPKJZJVTHSESZ-UHFFFAOYSA-N				

Abraham Solute Descriptors

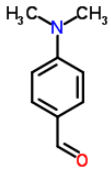
log S	E	S	A	B	V

	w				
-2.450	0.980	0.565	0.265	0.257	1.013

Solubility Data

	Concentration (M)		
Solvent	**Measured**	**Predicted**	**Solvent bp °C**
acetonitrile	1.67	1.22	63.5
DMSO	5.74	4.95	189
methanol	7.76	3.16	48.1
THF	4.45	11.84	68.3
toluene	0.89	2.33	110.6
1-butanol	-	4.56	117.7
1-decanol	-	5.30	227.8
1-heptanol	-	5.09	176.9
1-hexanol	-	4.85	158.2
1-octanol	-	3.82	194.7
1-pentanol	-	5.27	138.5
1-propanol	-	4.95	95.8
1,2-dichloroethane	-	2.16	83.5
1,4-dioxane	-	6.05	102.9
2-butanol	-	3.68	96.6
2-methyl-1-propanol	-	3.81	105.0
2-methyl-2-propanol	-	3.76	84.6
2-pentanol	-	5.07	118.8
2-propanol	-	4.04	73.0
2,2,4-trimethylpentane	-	0.48	98.8
3-methyl-1-butanol	-	4.39	131.2
acetone	-	5.13	46.5
benzene	-	2.38	78.8
butanone	-	5.65	75.6
butyl acetate	-	4.92	126.6
carbon disulfide	-	1.86	46.2
carbon tetrachloride	-	1.44	76.0
chloroform	-	2.18	61.2
cyclohexane	-	0.81	80.7
decane	-	0.53	174.9
dibutyl ether	-	2.25	142.1
dichloromethane	-	2.40	39.6
diethyl ether	-	6.20	33.2
DMF	-	5.89	153.0
ethyl acetate	-	5.56	73.9
ethylene glycol	-	0.78	197.5
heptane	-	0.74	98.8
hexadecane	-	0.46	286.6
hexane	-	0.53	68.5
methyl acetate	-	4.17	44.0
methyl tert-butyl ether	-	4.24	55.2
methylcyclohexane	-	0.98	101.1
nonane	-	0.58	151.7
octane	-	0.68	126.4
undecane	-	0.52	196.3
water	-	0.00	100.0

4-dimethylaminobenzaldehyde $C_9H_{11}NO$ - 6 solvents [208, 205]



Compound Data

Molecular weight	149.19	**H bond acceptors**	2	**Rule of 5 violations**	0
Compound type	aldehyde	**H bond donors**	0	**ACD/ALogP**	1.81
Phase 25°C	solid	**Rotatable bonds**	2	**Predicted density**	1.072 g/cm^3
SMILES	O=Cc1ccc(N(C)C)cc1				
InChIKey	BGNGWHSBYQYVRX-UHFFFAOYSA-N				

Abraham Solute Descriptors

log S_w	E	S	A	B	V
-1.050	1.190	1.602	0.079	0.756	1.162

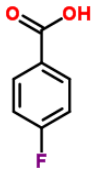
Solubility Data

	Concentration (M)		
Solvent	**Measured**	**Predicted**	**Solvent bp °C**
acetonitrile	2.60	2.77	63.5
chloroform	4.02	6.68	61.2
ethanol	0.63	0.86	72.6
methanol	1.36	1.22	48.1
THF	2.50	2.46	68.3
toluene	1.54	0.92	110.6
1-butanol	-	0.37	117.7
1-decanol	-	0.24	227.8
1-heptanol	-	0.31	176.9
1-hexanol	-	0.38	158.2
1-octanol	-	0.33	194.7
1-pentanol	-	0.34	138.5
1-propanol	-	0.43	95.8
1,2-dichloroethane	-	4.57	83.5
1,4-dioxane	-	3.07	102.9
2-butanol	-	0.63	96.6
2-methyl-1-propanol	-	0.42	105.0
2-methyl-2-propanol	-	0.40	84.6
2-pentanol	-	0.35	118.8
2-propanol	-	0.39	73.0
2,2,4-trimethylpentane	-	0.02	98.8
3-methyl-1-butanol	-	0.36	131.2
acetone	-	2.40	46.5
benzene	-	1.54	78.8
butanone	-	2.89	75.6
butyl acetate	-	0.92	126.6
carbon disulfide	-	0.22	46.2
carbon tetrachloride	-	0.34	76.0
cyclohexane	-	0.04	80.7
decane	-	0.02	174.9
dibutyl ether	-	0.12	142.1
dichloromethane	-	6.04	39.6
diethyl ether	-	0.49	33.2
DMF	-	4.88	153.0
DMSO	-	> 9	189.0
ethyl acetate	-	1.15	73.9
ethylene glycol	-	0.64	197.5
heptane	-	0.02	98.8
hexadecane	-	0.03	286.6
hexane	-	0.03	68.5
methyl acetate	-	2.74	44.0
methyl tert-butyl ether	-	0.41	55.2
methylcyclohexane	-	0.05	101.1
nonane	-	0.02	151.7
octane	-	0.03	126.4
undecane	-	0.02	196.3
water	-	0.09	100.0

4-fluorobenzoic acid $C_7H_5FO_2$ - 7 solvents [106, 121, 907]

Compound Data

Molecular weight	140.112	**H bond acceptors**	2	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	2.07
Phase 25°C	solid	**Rotatable bonds**	1	**Predicted density**	1.319 g/cm^3



SMILES O=C(O)c1ccc(F)cc1

InChIKey BBYDXOIZLAWGSL-UHFFFAOYSA-N

Abraham Solute Descriptors

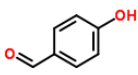
log S_w	E	S	A	B	V
-2.000	0.726	0.720	0.400	0.168	0.680

Solubility Data

	Concentration (M)		
Solvent	**Measured**	**Predicted**	**Solvent bp °C**
acetonitrile	0.39	0.42	63.5
DMSO	4.19	4.36	189
ethanol	0.80	0.89	72.6
methanol	0.76	0.82	48.1
THF	2.49	1.98	68.3
toluene	0.00	0.12	110.6
water	0.01	0.01	100
1-butanol	-	0.64	117.7
1-decanol	-	0.50	227.8
1-heptanol	-	0.55	176.9
1-hexanol	-	0.63	158.2
1-octanol	-	0.52	194.7
1-pentanol	-	0.65	138.5
1-propanol	-	0.83	95.8
1,2-dichloroethane	-	0.20	83.5
1,4-dioxane	-	1.25	102.9
2-butanol	-	0.69	96.6
2-methyl-1-propanol	-	0.64	105.0
2-methyl-2-propanol	-	0.76	84.6
2-pentanol	-	0.63	118.8
2-propanol	-	0.79	73.0
2,2,4-trimethylpentane	-	0.02	98.8
3-methyl-1-butanol	-	0.57	131.2
acetone	-	1.24	46.5
benzene	-	0.13	78.8
butanone	-	1.18	75.6
butyl acetate	-	0.75	126.6
carbon disulfide	-	0.06	46.2
carbon tetrachloride	-	0.05	76.0
chloroform	-	0.14	61.2
cyclohexane	-	0.02	80.7
decane	-	0.02	174.9
dibutyl ether	-	0.22	142.1
dichloromethane	-	0.19	39.6
diethyl ether	-	0.91	33.2
DMF	-	1.89	153.0
ethyl acetate	-	0.99	73.9
ethylene glycol	-	0.30	197.5
heptane	-	0.02	98.8
hexadecane	-	0.01	286.6
hexane	-	0.02	68.5
methyl acetate	-	0.86	44.0
methyl tert-butyl ether	-	0.53	55.2
methylcyclohexane	-	0.02	101.1
nonane	-	0.02	151.7
octane	-	0.02	126.4
undecane	-	0.02	196.3

4-hydroxybenzaldehyde $C_7H_6O_2$ - 7 solvents [208, 57, 205, 58, 88, 85, 909]

Compound Data



Molecular weight	122.121	**H bond acceptors**	2	**Rule of 5 violations**	0
Compound type	aldehyde	**H bond donors**	1	**ACD/ALogP**	1.39
Phase 25°C	solid	**Rotatable bonds**	2	**Predicted density**	1.226 g/cm^3
SMILES	O=Cc1ccc(O)cc1				
InChIKey	RGHHSNMVTDWUBI-UHFFFAOYSA-N				

Abraham Solute Descriptors

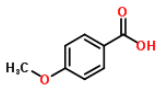
log S_w	E	S	A	B	V
-0.959	0.918	1.003	0.439	0.225	0.583

Solubility Data

	Concentration (M)		
Solvent	**Measured**	**Predicted**	**Solvent bp °C**
acetonitrile	1.46	1.38	63.5
chloroform	1.75	0.22	61.2
ethanol	2.52	1.65	72.6
methanol	3.39	1.97	48.1
THF	3.70	3.79	68.3
toluene	0.02	0.16	110.6
water	0.11	0.11	100
1-butanol	-	0.96	117.7
1-decanol	-	0.67	227.8
1-heptanol	-	0.73	176.9
1-hexanol	-	0.92	158.2
1-octanol	-	0.81	194.7
1-pentanol	-	0.93	138.5
1-propanol	-	1.38	95.8
1,2-dichloroethane	-	0.40	83.5
1,4-dioxane	-	3.04	102.9
2-butanol	-	1.29	96.6
2-methyl-1-propanol	-	1.01	105.0
2-methyl-2-propanol	-	1.19	84.6
2-pentanol	-	0.87	118.8
2-propanol	-	1.32	73.0
2,2,4-trimethylpentane	-	0.01	98.8
3-methyl-1-butanol	-	0.79	131.2
acetone	-	3.05	46.5
benzene	-	0.19	78.8
butanone	-	2.72	75.6
butyl acetate	-	1.30	126.6
carbon disulfide	-	0.06	46.2
carbon tetrachloride	-	0.05	76.0
cyclohexane	-	0.01	80.7
decane	-	0.01	174.9
dibutyl ether	-	0.24	142.1
dichloromethane	-	0.34	39.6
diethyl ether	-	1.34	33.2
DMF	-	5.10	153.0
DMSO	-	> 9	189.0
ethyl acetate	-	1.87	73.9
ethylene glycol	-	1.26	197.5
heptane	-	0.01	98.8
hexadecane	-	0.01	286.6
hexane	-	0.01	68.5
methyl acetate	-	1.96	44.0
methyl tert-butyl ether	-	0.69	55.2
methylcyclohexane	-	0.01	101.1
nonane	-	0.01	151.7

octane	-	0.02	126.4
undecane	-	0.01	196.3

4-methoxybenzoic acid $C_8H_8O_3$ - 22 solvents [905]



Compound Data

Molecular weight	152.147	**H bond acceptors**	3	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	1.96
Phase 25°C	solid	**Rotatable bonds**	2	**Predicted density**	1.207 g/cm^3
SMILES	COc1ccc(cc1)C(O)=O				
InChIKey	ZEYHEAKUIGZSGI-UHFFFAOYSA-N				

Abraham Solute Descriptors

log S_w	E	S	A	B	V
-1.880	0.850	1.290	0.404	0.552	0.970

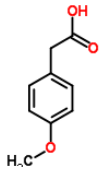
Solubility Data

	Concentration (M)		
Solvent	**Measured**	**Predicted**	**Solvent bp °C**
1-butanol	0.12	0.10	117.7
1-decanol	0.07	0.06	227.8
1-heptanol	0.09	0.08	176.9
1-hexanol	0.10	0.10	158.2
1-octanol	0.08	0.08	194.7
1-pentanol	0.10	0.10	138.5
1-propanol	0.13	0.14	95.8
1,4-dioxane	0.39	0.37	102.9
2-butanol	0.13	0.15	96.6
2-ethyl-1-hexanol	0.07	*	184.6
2-methyl-1-propanol	0.08	0.12	105
2-methyl-2-propanol	0.17	0.14	84.6
2-propanol	0.14	0.14	73
3-methyl-1-butanol	0.08	0.10	131.2
butyl acetate	0.08	0.12	126.6
dibutyl ether	0.02	0.02	142.1
diethyl ether	0.09	0.10	33.2
diisopropyl ether	0.03	*	68.3
ethanol	0.20	0.21	72.6
ethyl acetate	0.13	0.17	73.9
pentyl acetate	0.06	*	149.9
THF	0.70	0.43	68.3
1,2-dichloroethane	-	0.09	83.5
2-pentanol	-	0.10	118.8
2,2,4-trimethylpentane	-	0.00	98.8
acetone	-	0.32	46.5
acetonitrile	-	0.17	63.5
benzene	-	0.03	78.8
butanone	-	0.33	75.6
carbon disulfide	-	0.00	46.2
carbon tetrachloride	-	0.01	76.0
chloroform	-	0.09	61.2
cyclohexane	-	0.00	80.7
decane	-	0.00	174.9
dichloromethane	-	0.10	39.6
DMF	-	1.05	153.0
DMSO	-	3.09	189.0
ethylene glycol	-	0.15	197.5
heptane	-	0.00	98.8
hexadecane	-	0.00	286.6
hexane	-	0.00	68.5
methanol	-	0.26	48.1
methyl acetate	-	0.25	44.0

methyl tert-butyl ether	-	0.06	55.2
methylcyclohexane	-	0.00	101.1
nonane	-	0.00	151.7
octane	-	0.00	126.4
toluene	-	0.02	110.6
undecane	-	0.00	196.3
water	-	0.01	100.0

* This solvent doesn't have recorded Abraham coefficients.

4-methoxyphenylacetic acid $C_9H_{10}O_3$ - 6 solvents [89, 136, 92, 910]



Compound Data

Molecular weight	166.174	**H bond acceptors**	3	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	1.42
Phase 25°C	solid	**Rotatable bonds**	3	**Predicted density**	1.179 g/cm^3
SMILES	O=C(O)Cc1ccc(OC)cc1				
InChIKey	NRPFNQUDKRYCNX-UHFFFAOYSA-N				

Abraham Solute Descriptors

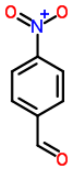
log S_w	E	S	A	B	V
-1.398	0.877	1.137	0.101	0.678	1.189

Solubility Data

	Concentration (M)		
Solvent	**Measured**	**Predicted**	**Solvent bp °C**
acetonitrile	2.23	2.08	63.5
DMSO	4.50	4.37	189
methanol	2.50	2.07	48.1
THF	3.31	4.06	68.3
toluene	0.37	1.66	110.6
water	0.04	0.04	100
1-butanol	-	1.13	117.7
1-decanol	-	0.80	227.8
1-heptanol	-	1.03	176.9
1-hexanol	-	1.14	158.2
1-octanol	-	0.90	194.7
1-pentanol	-	1.10	138.5
1-propanol	-	1.21	95.8
1,2-dichloroethane	-	4.67	83.5
1,4-dioxane	-	3.51	102.9
2-butanol	-	1.47	96.6
2-methyl-1-propanol	-	1.21	105.0
2-methyl-2-propanol	-	1.16	84.6
2-pentanol	-	1.20	118.8
2-propanol	-	1.09	73.0
2,2,4-trimethylpentane	-	0.10	98.8
3-methyl-1-butanol	-	1.21	131.2
acetone	-	2.85	46.5
benzene	-	2.42	78.8
butanone	-	3.51	75.6
butyl acetate	-	1.68	126.6
carbon disulfide	-	0.53	46.2
carbon tetrachloride	-	0.91	76.0
chloroform	-	> 9	61.2
cyclohexane	-	0.18	80.7
decane	-	0.09	174.9
dibutyl ether	-	0.38	142.1
dichloromethane	-	7.21	39.6
diethyl ether	-	1.22	33.2
DMF	-	5.11	153.0
ethyl acetate	-	1.93	73.9

ethylene glycol	-	0.67	197.5
heptane	-	0.11	98.8
hexadecane	-	0.11	286.6
hexane	-	0.12	68.5
methyl acetate	-	3.37	44.0
methyl tert-butyl ether	-	1.09	55.2
methylcyclohexane	-	0.25	101.1
nonane	-	0.12	151.7
octane	-	0.13	126.4
undecane	-	0.09	196.3

4-nitrobenzaldehyde $C_7H_5NO_3$ - 10 solvents [111, 98, 904, 212, 82, 122]



Compound Data

Molecular weight	151.12	**H bond acceptors**	4	**Rule of 5 violations**	0
Compound type	aldehyde	**H bond donors**	0	**ACD/ALogP**	1.56
Phase 25°C	solid	**Rotatable bonds**	2	**Predicted density**	1.338 g/cm^3
SMILES	c1cc(C=O)ccc1[N+](=O)[O-]				
InChIKey	BXRFQSNOROATLV-UHFFFAOYSA-N				

Abraham Solute Descriptors

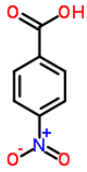
log S_w	E	S	A	B	V
-2.280	1.049	1.634	-0.031	0.289	0.780

Solubility Data

	Concentration (M)		
Solvent	**Measured**	**Predicted**	**Solvent bp °C**
acetonitrile	1.13	1.41	63.5
benzene	0.46	0.44	78.8
carbon tetrachloride	0.07	0.09	76
chloroform	1.02	0.74	61.2
cyclohexane	0.01	0.02	80.7
dichloromethane	1.29	1.33	39.6
DMF	1.68	0.99	153
DMSO	1.24	2.70	189
THF	0.88	0.53	68.3
toluene	0.46	0.30	110.6
1-butanol	-	0.03	117.7
1-decanol	-	0.02	227.8
1-heptanol	-	0.02	176.9
1-hexanol	-	0.03	158.2
1-octanol	-	0.04	194.7
1-pentanol	-	0.02	138.5
1-propanol	-	0.03	95.8
1,2-dichloroethane	-	1.25	83.5
1,4-dioxane	-	0.84	102.9
2-butanol	-	0.05	96.6
2-methyl-1-propanol	-	0.03	105.0
2-methyl-2-propanol	-	0.04	84.6
2-pentanol	-	0.02	118.8
2-propanol	-	0.03	73.0
2,2,4-trimethylpentane	-	0.01	98.8
3-methyl-1-butanol	-	0.02	131.2
acetone	-	0.76	46.5
butanone	-	0.91	75.6
butyl acetate	-	0.28	126.6
carbon disulfide	-	0.16	46.2
decane	-	0.01	174.9
dibutyl ether	-	0.06	142.1
diethyl ether	-	0.12	33.2
ethanol	-	0.06	72.6
ethyl acetate	-	0.35	73.9

ethylene glycol	-	0.03	197.5
heptane	-	0.00	98.8
hexadecane	-	0.01	286.6
methanol	-	0.11	48.1
methyl acetate	-	0.76	44.0
methyl tert-butyl ether	-	0.13	55.2
methylcyclohexane	-	0.01	101.1
nonane	-	0.01	151.7
octane	-	0.01	126.4
undecane	-	0.01	196.3
water	-	0.01	100.0

4-nitrobenzoic acid $C_7H_5NO_4$ - 29 solvents [905]



Compound Data

Molecular weight	167.119	**H bond acceptors**	5	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	1.89
Phase 25°C	solid	**Rotatable bonds**	2	**Predicted density**	1.468 g/cm^3
SMILES	O=[N+]([O-])c1ccc(C(=O)O)cc1				
InChIKey	OTLNPYWUJOZPPA-UHFFFAOYSA-N				

Abraham Solute Descriptors

log S_w	E	S	A	B	V
-2.120	1.048	1.301	0.518	0.349	0.793

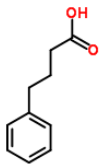
Solubility Data

	Concentration (M)		
Solvent	**Measured**	**Predicted**	**Solvent bp °C**
1-butanol	0.09	0.08	117.7
1-decanol	0.05	0.05	227.8
1-heptanol	0.08	0.06	176.9
1-hexanol	0.08	0.08	158.2
1-octanol	0.06	0.07	194.7
1-pentanol	0.09	0.08	138.5
1-propanol	0.11	0.13	95.8
1,4-dioxane	0.47	0.38	102.9
2-butanol	0.10	0.12	96.6
2-methyl-1-butanol	0.05	*	128.7
2-methyl-1-pentanol	0.05	*	148
2-methyl-1-propanol	0.07	0.09	105
2-methyl-2-butanol	0.17	*	102
2-methyl-2-propanol	0.15	0.11	84.6
2-pentanol	0.08	0.07	118.8
2-propanol	0.11	0.13	73
3-methyl-1-butanol	0.07	0.07	131.2
4-methyl-2-pentanol	0.07	*	133.5
butyl acetate	0.08	0.12	126.6
dibutyl ether	0.02	0.01	142.1
diethyl ether	0.10	0.11	33.2
diisopropyl ether	0.03	*	68.3
ethanol	0.14	0.16	72.6
ethyl acetate	0.13	0.17	73.9
methanol	0.18	0.20	48.1
methyl acetate	0.14	0.20	44
pentyl acetate	0.05	*	149.9
propylene carbonate	0.06	*	332.3
THF	0.78	0.45	68.3
1,2-dichloroethane	-	0.04	83.5
2,2,4-trimethylpentane	-	0.00	98.8
acetone	-	0.33	46.5
acetonitrile	-	0.13	63.5

benzene	-	0.01	78.8
butanone	-	0.31	75.6
carbon disulfide	-	0.00	46.2
carbon tetrachloride	-	0.00	76.0
chloroform	-	0.02	61.2
cyclohexane	-	0.00	80.7
decane	-	0.00	174.9
dichloromethane	-	0.03	39.6
DMF	-	1.02	153.0
DMSO	-	6.21	189.0
ethylene glycol	-	0.14	197.5
heptane	-	0.00	98.8
hexadecane	-	0.00	286.6
hexane	-	0.00	68.5
methyl tert-butyl ether	-	0.05	55.2
methylcyclohexane	-	0.00	101.1
nonane	-	0.00	151.7
octane	-	0.00	126.4
toluene	-	0.01	110.6
undecane	-	0.00	196.3
water	-	0.01	100.0

* This solvent doesn't have recorded Abraham coefficients.

4-phenylbutyric acid $C_{10}H_{12}O_2$ - 5 solvents [72]



Compound Data

Molecular weight	164.201	**H bond acceptors**	2	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	2.42
Phase 25°C	solid	**Rotatable bonds**	4	**Predicted density**	1.095 g/cm^3
SMILES	O=C(O)CCCc1ccccc1				
InChIKey	OBKXEAXTFZPCHS-UHFFFAOYSA-N				

Abraham Solute Descriptors

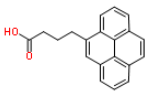
log S_w	E	S	A	B	V
-2.510	0.896	3.980	0.540	-0.076	1.321

Solubility Data

	Concentration (M)		
Solvent	**Measured**	**Predicted**	**Solvent bp °C**
1-hexanol	0.19	0.21	158.2
chloroform	3.99	3.67	61.2
diethyl ether	4.02	5.98	33.2
methanol	4.54	2.27	48.1
toluene	4.06	0.97	110.6
1-butanol	-	0.14	117.7
1-decanol	-	0.02	227.8
1-heptanol	-	0.12	176.9
1-octanol	-	0.42	194.7
1-pentanol	-	0.07	138.5
1-propanol	-	0.31	95.8
1,2-dichloroethane	-	> 9	83.5
1,4-dioxane	-	> 9	102.9
2-butanol	-	0.42	96.6
2-methyl-1-propanol	-	0.25	105.0
2-methyl-2-propanol	-	1.18	84.6
2-pentanol	-	0.05	118.8
2-propanol	-	0.51	73.0
2,2,4-trimethylpentane	-	0.00	98.8
3-methyl-1-butanol	-	0.08	131.2
acetone	-	> 9	46.5
acetonitrile	-	> 9	63.5

benzene	-	2.33	78.8
butanone	-	> 9	75.6
butyl acetate	-	> 9	126.6
carbon disulfide	-	0.25	46.2
carbon tetrachloride	-	0.01	76.0
cyclohexane	-	0.00	80.7
decane	-	0.00	174.9
dibutyl ether	-	0.62	142.1
dichloromethane	-	> 9	39.6
DMF	-	> 9	153.0
DMSO	-	> 9	189.0
ethyl acetate	-	> 9	73.9
ethylene glycol	-	0.20	197.5
heptane	-	0.00	98.8
hexadecane	-	0.00	286.6
hexane	-	0.00	68.5
methyl acetate	-	> 9	44.0
methyl tert-butyl ether	-	3.92	55.2
methylcyclohexane	-	0.00	101.1
nonane	-	0.00	151.7
octane	-	0.00	126.4
THF	-	> 9	68.3
undecane	-	0.00	196.3
water	-	0.00	100.0

4-pyrenebutanoic acid $C_{20}H_{16}O_2$ - 16 solvents [61, 99, 51, 30, 24, 91, 84]



Compound Data

Molecular weight	288.34	**H bond acceptors**	2	**Rule of 5 violations**	1
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	5.37
Phase 25°C	solid	**Rotatable bonds**	4	**Predicted density**	1.296 g/cm^3
SMILES	O=C(O)CCCc4cc2cccc1ccc3c(c12)c4ccc3				
InChIKey	GSBSDUZVSCTUKA-UHFFFAOYSA-N				

Abraham Solute Descriptors

log S_w	E	S	A	B	V
-7.150	3.727	2.030	0.411	0.855	2.385

Solubility Data

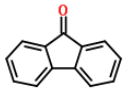
	Concentration (M)		
Solvent	**Measured**	**Predicted**	**Solvent bp °C**
2-propanol	0.06	0.04	73.0
acetonitrile	0.00	0.01	63.5
benzene	0.00	0.11	78.8
carbon tetrachloride	0.00	0.02	76
chloroform	0.03	0.06	61.2
cyclohexane	0.00	0.00	80.7
cyclopentane	0.00	*	49.2
dichloromethane	0.07	0.05	39.6
diethyl ether	0.02	0.08	33.2
DMF	1.88	0.51	153
DMSO	2.07	1.36	189
ethanol	0.06	0.09	72.6
hexane	0.00	0.00	68.5
methanol	0.01	0.03	48.1
THF	0.55	0.87	68.3
toluene	0.00	0.09	110.6
1-butanol	-	0.06	117.7
1-decanol	-	0.27	227.8
1-heptanol	-	0.12	176.9
1-hexanol	-	0.10	158.2
1-octanol	-	0.09	194.7

1-pentanol	-	0.11	138.5
1-propanol	-	0.07	95.8
1,2-dichloroethane	-	0.10	83.5
1,4-dioxane	-	0.48	102.9
2-butanol	-	0.04	96.6
2-methyl-1-propanol	-	0.03	105.0
2-methyl-2-propanol	-	0.02	84.6
2-pentanol	-	0.07	118.8
2,2,4-trimethylpentane	-	0.00	98.8
3-methyl-1-butanol	-	0.05	131.2
acetone	-	0.15	46.5
butanone	-	0.21	75.6
butyl acetate	-	0.10	126.6
carbon disulfide	-	0.07	46.2
decane	-	0.00	174.9
dibutyl ether	-	0.01	142.1
ethyl acetate	-	0.12	73.9
ethylene glycol	-	0.03	197.5
heptane	-	0.00	98.8
hexadecane	-	0.00	286.6
methyl acetate	-	0.08	44.0
methyl tert-butyl ether	-	0.02	55.2
methylcyclohexane	-	0.00	101.1
nonane	-	0.00	151.7
octane	-	0.00	126.4
undecane	-	0.00	196.3
water	-	0.00	100.0

* This solvent doesn't have recorded Abraham coefficients.

9-fluorenone $C_{13}H_8O$ - 43 solvents [905]

Compound Data



Molecular weight	180.202	**H bond acceptors**	1	**Rule of 5 violations**	0
Compound type	non-Ugi related	**H bond donors**	0	**ACD/ALogP**	3.58
Phase 25°C	solid	**Rotatable bonds**	0	**Predicted density**	1.244 g/cm^3
SMILES	O=C3c1ccccc1c2c3cccc2				
InChIKey	YLQWCDOCJODRMT-UHFFFAOYSA-N				

Abraham Solute Descriptors

log S_w	E	S	A	B	V
-3.910	1.638	1.392	0.097	0.233	1.247

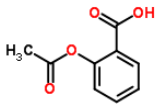
Solubility Data

	Concentration (M)		
Solvent	**Measured**	**Predicted**	**Solvent bp °C**
1-butanol	0.37	0.36	117.7
1-decanol	0.42	0.44	227.8
1-heptanol	0.43	0.46	176.9
1-hexanol	0.41	0.46	158.2
1-octanol	0.43	0.47	194.7
1-pentanol	0.41	0.38	138.5
1-propanol	0.38	0.39	95.8
2-butanol	0.33	0.34	96.6
2-ethyl-1-hexanol	0.35	*	184.6
2-methyl-1-pentanol	0.36	*	148
2-methyl-1-propanol	0.25	0.28	105
2-methyl-2-butanol	0.40	*	102
2-methyl-2-propanol	0.31	0.29	84.6
2-pentanol	0.35	0.31	118.8
2-propanol	0.29	0.29	73
2,2,4-trimethylpentane	0.07	0.08	98.8

3-methyl-1-butanol	0.31	0.29	131.2
4-methyl-2-pentanol	0.29	*	133.5
acetonitrile	1.57	1.27	63.5
benzene	2.61	2.55	78.8
butyronitrile	1.93	*	117.3
carbon tetrachloride	1.83	0.75	76
cyclohexane	0.15	0.25	80.7
cyclooctane	0.17	*	152
cyclopentanol	0.82	*	140.8
decane	0.12	0.10	174.9
dibutyl ether	0.37	0.61	142.1
dichloromethane	4.03	2.89	39.6
diethyl ether	0.88	1.36	33.2
diisopropyl ether	0.37	*	68.3
ethanol	3.08	0.50	72.6
heptane	0.11	0.08	98.8
hexadecane	0.12	0.12	286.6
hexane	0.10	0.09	68.5
methanol	0.37	0.37	48.1
methyl tert-butyl ether	0.76	1.06	55.2
methylcyclohexane	0.14	0.15	101.1
nonane	0.12	0.12	151.7
octane	0.11	0.16	126.4
propionitrile	1.99	*	91.3
tert-butylcyclohexane	0.12	*	168.3
toluene	2.17	2.23	110.6
undecane	0.12	0.12	196.3
1,2-dichloroethane	-	3.24	83.5
1,4-dioxane	-	4.70	102.9
acetone	-	3.14	46.5
butanone	-	4.10	75.6
butyl acetate	-	2.10	126.6
carbon disulfide	-	2.41	46.2
chloroform	-	1.79	61.2
DMF	-	4.62	153.0
DMSO	-	6.16	189.0
ethyl acetate	-	2.33	73.9
ethylene glycol	-	0.08	197.5
methyl acetate	-	2.50	44.0
THF	-	5.44	68.3
water	-	0.00	100.0

* This solvent doesn't have recorded Abraham coefficients.

acetylsalicylic acid $C_9H_8O_4$ - 6 solvents [72, 34, 903, 907]



Compound Data

Molecular weight	180.157	**H bond acceptors**	4	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	1.19
Phase 25°C	solid	**Rotatable bonds**	3	**Predicted density**	1.29 g/cm^3
SMILES	O=C(Oc1ccccc1C(=O)O)C				
InChIKey	BSYNRYMUTXBXSQ-UHFFFAOYSA-N				

Abraham Solute Descriptors

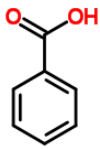
log S_w	E	S	A	B	V
-1.602	0.946	5.432	0.815	-0.028	1.170

Solubility Data

	Concentration (M)		
Solvent	**Measured**	**Predicted**	**Solvent bp °C**
1-hexanol	0.00	0.01	158.2
chloroform	0.18	0.16	61.2

diethyl ether	0.17	0.32	33.2
methanol	1.30	0.43	48.1
toluene	0.00	0.01	110.6
water	0.03	0.03	100
1-butanol	-	0.00	117.7
1-decanol	-	0.00	227.8
1-heptanol	-	0.00	176.9
1-octanol	-	0.02	194.7
1-pentanol	-	0.00	138.5
1-propanol	-	0.01	95.8
1,2-dichloroethane	-	7.20	83.5
1,4-dioxane	-	> 9	102.9
2-butanol	-	0.03	96.6
2-methyl-1-propanol	-	0.01	105.0
2-methyl-2-propanol	-	0.09	84.6
2-pentanol	-	0.00	118.8
2-propanol	-	0.03	73.0
2,2,4-trimethylpentane	-	0.00	98.8
3-methyl-1-butanol	-	0.00	131.2
acetone	-	> 9	46.5
acetonitrile	-	> 9	63.5
benzene	-	0.05	78.8
butanone	-	> 9	75.6
butyl acetate	-	5.04	126.6
carbon disulfide	-	0.00	46.2
carbon tetrachloride	-	0.00	76.0
cyclohexane	-	0.00	80.7
decane	-	0.00	174.9
dibutyl ether	-	0.01	142.1
dichloromethane	-	3.00	39.6
DMF	-	> 9	153.0
DMSO	-	> 9	189.0
ethyl acetate	-	> 9	73.9
ethylene glycol	-	0.09	197.5
heptane	-	0.00	98.8
hexadecane	-	0.00	286.6
hexane	-	0.00	68.5
methyl acetate	-	> 9	44.0
methyl tert-butyl ether	-	0.15	55.2
methylcyclohexane	-	0.00	101.1
nonane	-	0.00	151.7
octane	-	0.00	126.4
THF	-	> 9	68.3
undecane	-	0.00	196.3

benzoic acid $C_7H_6O_2$ - 10 solvents [155, 37, 902, 58, 160, 43, 136, 9, 907, 900]



Compound Data

Molecular weight	122.121	**H bond acceptors**	2	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	1.9
Phase 25°C	solid	**Rotatable bonds**	1	**Predicted density**	1.197 g/cm^3
SMILES	O=C(O)c1ccccc1				
InChIKey	WPYMKLBDIGXBTP-UHFFFAOYSA-N				

Abraham Solute Descriptors

log S_w	E	S	A	B	V
-1.538	0.788	0.144	0.100	0.239	0.618

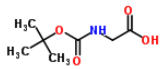
Solubility Data

Concentration (M)

Solvent	Measured	Predicted	Solvent bp °C
1-octanol	1.48	1.81	194.7

acetone	1.35	1.75	46.5
acetonitrile	0.75	0.72	63.5
benzene	0.48	1.80	78.8
chloroform	1.80	1.75	61.2
ethanol	2.55	2.69	72.6
methanol	2.84	1.99	48.1
THF	3.37	2.83	68.3
toluene	0.65	1.87	110.6
water	0.03	0.03	100
1-butanol	-	2.58	117.7
1-decanol	-	2.99	227.8
1-heptanol	-	2.41	176.9
1-hexanol	-	2.38	158.2
1-pentanol	-	2.85	138.5
1-propanol	-	2.51	95.8
1,2-dichloroethane	-	1.37	83.5
1,4-dioxane	-	1.56	102.9
2-butanol	-	2.18	96.6
2-methyl-1-propanol	-	2.13	105.0
2-methyl-2-propanol	-	1.76	84.6
2-pentanol	-	2.93	118.8
2-propanol	-	1.92	73.0
2,2,4-trimethylpentane	-	1.47	98.8
3-methyl-1-butanol	-	2.53	131.2
butanone	-	1.68	75.6
butyl acetate	-	1.95	126.6
carbon disulfide	-	1.61	46.2
carbon tetrachloride	-	2.03	76.0
cyclohexane	-	1.89	80.7
decane	-	1.29	174.9
dibutyl ether	-	1.45	142.1
dichloromethane	-	1.68	39.6
diethyl ether	-	2.70	33.2
DMF	-	0.64	153.0
DMSO	-	0.57	189.0
ethyl acetate	-	2.19	73.9
ethylene glycol	-	0.66	197.5
heptane	-	2.32	98.8
hexadecane	-	1.15	286.6
hexane	-	1.66	68.5
methyl acetate	-	1.73	44.0
methyl tert-butyl ether	-	2.23	55.2
methylcyclohexane	-	2.89	101.1
nonane	-	1.53	151.7
octane	-	1.56	126.4
undecane	-	1.11	196.3

boc-glycine $C_7H_{13}NO_4$ - 12 solvents [143, 71, 207, 135, 133, 116]



Compound Data

Molecular weight	175.182	**H bond acceptors**	5	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	2	**ACD/ALogP**	0.75
Phase 25°C	solid	**Rotatable bonds**	4	**Predicted density**	1.159 g/cm^3
SMILES	O=C(OC(C)(C)C)NCC(=O)O				
InChIKey	VRPJIFMKZZEXLR-UHFFFAOYSA-N				

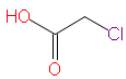
Abraham Solute Descriptors

log S_w	E	S	A	B	V
-1.260	0.402	0.823	0.255	0.784	1.338

Solubility Data

	Concentration (M)		
Solvent	**Measured**	**Predicted**	**Solvent bp °C**
2-propanol	3.63	4.15	73.0
acetonitrile	2.46	1.29	63.5
benzene	0.25	1.59	78.8
chloroform	3.85	9.41	61.2
dichloromethane	4.14	5.56	39.6
diethyl ether	2.65	2.18	33.2
DMSO	4.78	3.87	189
ethanol	4.11	5.88	72.6
hexane	0.25	0.12	68.5
methanol	4.72	5.01	48.1
THF	3.72	6.26	68.3
toluene	0.19	1.08	110.6
1-butanol	-	3.70	117.7
1-decanol	-	2.30	227.8
1-heptanol	-	3.43	176.9
1-hexanol	-	3.55	158.2
1-octanol	-	2.30	194.7
1-pentanol	-	3.85	138.5
1-propanol	-	4.06	95.8
1,2-dichloroethane	-	2.79	83.5
1,4-dioxane	-	3.70	102.9
2-butanol	-	4.29	96.6
2-methyl-1-propanol	-	4.37	105.0
2-methyl-2-propanol	-	4.59	84.6
2-pentanol	-	4.80	118.8
2,2,4-trimethylpentane	-	0.10	98.8
3-methyl-1-butanol	-	4.73	131.2
acetone	-	2.97	46.5
butanone	-	3.73	75.6
butyl acetate	-	2.13	126.6
carbon disulfide	-	0.25	46.2
carbon tetrachloride	-	0.73	76.0
cyclohexane	-	0.14	80.7
decane	-	0.10	174.9
dibutyl ether	-	0.51	142.1
DMF	-	> 9	153.0
ethyl acetate	-	2.40	73.9
ethylene glycol	-	1.27	197.5
heptane	-	0.14	98.8
hexadecane	-	0.09	286.6
methyl acetate	-	3.62	44.0
methyl tert-butyl ether	-	1.80	55.2
methylcyclohexane	-	0.29	101.1
nonane	-	0.11	151.7
octane	-	0.12	126.4
undecane	-	0.09	196.3
water	-	0.05	100.0

chloroacetic acid $C_2H_3ClO_2$ - 5 solvents [95]



Compound Data

Molecular weight	94.497	**H bond acceptors**	2	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	0.18
Phase 25°C	solid	**Rotatable bonds**	1	**Predicted density**	1.398 g/cm^3
SMILES	ClCC(O)=O				
InChIKey	FOCAUTSVDIKZOP-UHFFFAOYSA-N				

Abraham Solute Descriptors

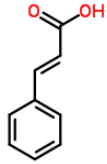
log S_w	E	S	A	B	V

0.470	0.502	0.372	0.096	0.042	0.114

Solubility Data

	Concentration (M)		
Solvent	**Measured**	**Predicted**	**Solvent bp °C**
acetonitrile	12.45	12.35	63.5
DMSO	11.20	11.16	189
methanol	9.45	9.23	48.1
THF	9.63	9.88	68.3
toluene	1.41	4.33	110.6
1-butanol	-	5.23	117.7
1-decanol	-	3.70	227.8
1-heptanol	-	3.58	176.9
1-hexanol	-	4.39	158.2
1-octanol	-	3.92	194.7
1-pentanol	-	4.63	138.5
1-propanol	-	5.78	95.8
1,2-dichloroethane	-	6.00	83.5
1,4-dioxane	-	8.55	102.9
2-butanol	-	6.29	96.6
2-methyl-1-propanol	-	5.11	105.0
2-methyl-2-propanol	-	5.30	84.6
2-pentanol	-	4.57	118.8
2-propanol	-	4.99	73.0
2,2,4-trimethylpentane	-	2.33	98.8
3-methyl-1-butanol	-	4.41	131.2
acetone	-	> 9	46.5
benzene	-	4.52	78.8
butanone	-	> 9	75.6
butyl acetate	-	8.05	126.6
carbon disulfide	-	3.03	46.2
carbon tetrachloride	-	3.11	76.0
chloroform	-	4.16	61.2
cyclohexane	-	2.26	80.7
decane	-	1.81	174.9
dibutyl ether	-	4.43	142.1
dichloromethane	-	6.29	39.6
diethyl ether	-	8.68	33.2
DMF	-	4.03	153.0
ethyl acetate	-	> 9	73.9
ethylene glycol	-	3.98	197.5
heptane	-	2.32	98.8
hexadecane	-	1.76	286.6
hexane	-	2.69	68.5
methyl acetate	-	> 9	44.0
methyl tert-butyl ether	-	7.47	55.2
methylcyclohexane	-	2.34	101.1
nonane	-	2.23	151.7
octane	-	2.35	126.4
undecane	-	1.57	196.3
water	-	2.95	100.0

cinnamic acid $C_9H_8O_2$ - 7 solvents [72, 902, 48, 907]



Compound Data

Molecular weight	148.159	**H bond acceptors**	2	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	2.38
Phase 25°C	solid	**Rotatable bonds**	2	**Predicted density**	1.184 g/cm^3
SMILES	O=C(O)\C=C\c1ccccc1				
InChIKey	WBYWAXJHAXSJNI-VOTSOKGWSA-N				

Abraham Solute Descriptors

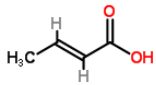
log S_w	E	S	A	B	V
-2.310	1.161	4.115	0.445	-0.194	0.957

Solubility Data

	Concentration (M)		
Solvent	**Measured**	**Predicted**	**Solvent bp °C**
1-hexanol	0.00	0.03	158.2
chloroform	0.93	0.71	61.2
diethyl ether	0.43	1.07	33.2
ethanol	0.86	0.09	72.6
methanol	1.10	0.46	48.1
toluene	0.30	0.27	110.6
water	0.00	0.00	100
1-butanol	-	0.02	117.7
1-decanol	-	0.00	227.8
1-heptanol	-	0.01	176.9
1-octanol	-	0.06	194.7
1-pentanol	-	0.01	138.5
1-propanol	-	0.04	95.8
1,2-dichloroethane	-	> 9	83.5
1,4-dioxane	-	> 9	102.9
2-butanol	-	0.07	96.6
2-methyl-1-propanol	-	0.03	105.0
2-methyl-2-propanol	-	0.14	84.6
2-pentanol	-	0.00	118.8
2-propanol	-	0.06	73.0
2,2,4-trimethylpentane	-	0.00	98.8
3-methyl-1-butanol	-	0.01	131.2
acetone	-	> 9	46.5
acetonitrile	-	> 9	63.5
benzene	-	0.64	78.8
butanone	-	> 9	75.6
butyl acetate	-	> 9	126.6
carbon disulfide	-	0.08	46.2
carbon tetrachloride	-	0.00	76.0
cyclohexane	-	0.00	80.7
decane	-	0.00	174.9
dibutyl ether	-	0.13	142.1
dichloromethane	-	8.52	39.6
DMF	-	> 9	153.0
DMSO	-	> 9	189.0
ethyl acetate	-	> 9	73.9
ethylene glycol	-	0.07	197.5
heptane	-	0.00	98.8
hexadecane	-	0.00	286.6
hexane	-	0.00	68.5
methyl acetate	-	> 9	44.0
methyl tert-butyl ether	-	0.70	55.2
methylcyclohexane	-	0.00	101.1
nonane	-	0.00	151.7
octane	-	0.00	126.4
THF	-	> 9	68.3
undecane	-	0.00	196.3

crotonic acid $C_4H_6O_2$ - 10 solvents [64, 62, 907]

Compound Data



Molecular weight	86.0892	**H bond acceptors**	2	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	0.94
Phase 25°C	solid	**Rotatable bonds**	1	**Predicted density**	1.039 g/cm^3
SMILES	O=C(O)/C=C/C				

InChIKey LDHQCZJRKDOVOX-NSCUHMNNSA-N

Abraham Solute Descriptors

log S_w	E	S	A	B	V
-0.051	0.474	0.405	0.130	0.155	0.339

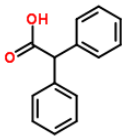
Solubility Data

	Concentration (M)		
Solvent	**Measured**	**Predicted**	**Solvent bp °C**
acetonitrile	3.59	6.14	63.5
benzene	4.45	3.08	78.8
chloroform	7.34	3.70	61.2
DMSO	7.76	6.77	189
ethanol	6.15	5.68	72.6
hexane	1.07	1.43	68.5
methanol	7.62	6.67	48.1
THF	8.09	7.70	68.3
toluene	3.52	2.84	110.6
water	0.89	0.89	100
1-butanol	-	4.32	117.7
1-decanol	-	3.09	227.8
1-heptanol	-	3.20	176.9
1-hexanol	-	3.73	158.2
1-octanol	-	3.13	194.7
1-pentanol	-	3.99	138.5
1-propanol	-	4.74	95.8
1,2-dichloroethane	-	4.04	83.5
1,4-dioxane	-	5.95	102.9
2-butanol	-	4.95	96.6
2-methyl-1-propanol	-	4.30	105.0
2-methyl-2-propanol	-	4.48	84.6
2-pentanol	-	4.09	118.8
2-propanol	-	4.19	73.0
2,2,4-trimethylpentane	-	1.24	98.8
3-methyl-1-butanol	-	3.94	131.2
acetone	-	7.63	46.5
butanone	-	6.90	75.6
butyl acetate	-	5.39	126.6
carbon disulfide	-	1.74	46.2
carbon tetrachloride	-	2.08	76.0
cyclohexane	-	1.27	80.7
decane	-	1.01	174.9
dibutyl ether	-	2.73	142.1
dichloromethane	-	4.74	39.6
diethyl ether	-	5.99	33.2
DMF	-	3.71	153.0
ethyl acetate	-	6.73	73.9
ethylene glycol	-	2.47	197.5
heptane	-	1.36	98.8
hexadecane	-	0.96	286.6
methyl acetate	-	7.18	44.0
methyl tert-butyl ether	-	5.10	55.2
methylcyclohexane	-	1.50	101.1
nonane	-	1.23	151.7
octane	-	1.28	126.4
undecane	-	0.89	196.3

diphenylacetic acid $C_{14}H_{12}O_2$ - 7 solvents [81, 77, 907]

Compound Data

Molecular weight	212.244	**H bond acceptors**	2	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	2.79



Phase 25°C	solid	**Rotatable bonds**	3	**Predicted density**	1.174 g/cm^3
SMILES	O=C(O)C(c1ccccc1)c2ccccc2				
InChIKey	PYHXGXCGESYPCW-UHFFFAOYSA-N				

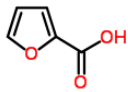
Abraham Solute Descriptors

log S_w	E	S	A	B	V
-2.310	1.565	1.499	0.147	1.083	1.773

Solubility Data

	Concentration (M)		
Solvent	**Measured**	**Predicted**	**Solvent bp °C**
acetonitrile	0.46	0.49	63.5
DMSO	2.56	2.65	189
ethanol	1.15	1.40	72.6
methanol	1.27	1.12	48.1
THF	2.91	2.51	68.3
toluene	0.17	0.93	110.6
water	0.00	0.00	100
1-butanol	-	0.70	117.7
1-decanol	-	0.71	227.8
1-heptanol	-	0.76	176.9
1-hexanol	-	0.79	158.2
1-octanol	-	0.56	194.7
1-pentanol	-	0.81	138.5
1-propanol	-	0.75	95.8
1,2-dichloroethane	-	3.06	83.5
1,4-dioxane	-	2.00	102.9
2-butanol	-	0.90	96.6
2-methyl-1-propanol	-	0.69	105.0
2-methyl-2-propanol	-	0.50	84.6
2-pentanol	-	0.88	118.8
2-propanol	-	0.61	73.0
2,2,4-trimethylpentane	-	0.02	98.8
3-methyl-1-butanol	-	0.81	131.2
acetone	-	1.16	46.5
benzene	-	1.54	78.8
butanone	-	1.59	75.6
butyl acetate	-	0.66	126.6
carbon disulfide	-	0.20	46.2
carbon tetrachloride	-	0.46	76.0
chloroform	-	> 9	61.2
cyclohexane	-	0.06	80.7
decane	-	0.02	174.9
dibutyl ether	-	0.09	142.1
dichloromethane	-	4.75	39.6
diethyl ether	-	0.45	33.2
DMF	-	3.53	153.0
ethyl acetate	-	0.75	73.9
ethylene glycol	-	0.59	197.5
heptane	-	0.03	98.8
hexadecane	-	0.03	286.6
hexane	-	0.02	68.5
methyl acetate	-	1.43	44.0
methyl tert-butyl ether	-	0.33	55.2
methylcyclohexane	-	0.11	101.1
nonane	-	0.03	151.7
octane	-	0.03	126.4
undecane	-	0.02	196.3

furoic acid $C_5H_4O_3$ - 6 solvents [104, 911]



Compound Data

Molecular weight	112.084	**H bond acceptors**	3	**Rule of 5 violations**	0
Compound type	carboxlyic acid	**H bond donors**	1	**ACD/ALogP**	0.98
Phase 25°C	solid	**Rotatable bonds**	1	**Predicted density**	1.322 g/cm^3
SMILES	OC(=O)c1ccco1				
InChIKey	SMNDYUVBFMFKNZ-UHFFFAOYSA-N				

Abraham Solute Descriptors

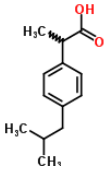
log S_w	E	S	A	B	V
-0.620	0.590	0.351	0.479	0.125	0.366

Solubility Data

	Concentration (M)		
Solvent	**Measured**	**Predicted**	**Solvent bp °C**
acetonitrile	1.10	0.77	63.5
DMSO	6.64	8.41	189
methanol	4.12	3.57	48.1
THF	4.07	3.71	68.3
toluene	0.06	0.16	110.6
water	0.32	0.16	110.6
1-butanol	-	2.79	117.7
1-decanol	-	2.23	227.8
1-heptanol	-	2.08	176.9
1-hexanol	-	2.44	158.2
1-octanol	-	1.88	194.7
1-pentanol	-	2.93	138.5
1-propanol	-	3.80	95.8
1,2-dichloroethane	-	0.22	83.5
1,4-dioxane	-	2.07	102.9
2-butanol	-	3.03	96.6
2-methyl-1-propanol	-	2.79	105.0
2-methyl-2-propanol	-	3.10	84.6
2-pentanol	-	2.93	118.8
2-propanol	-	3.60	73.0
2,2,4-trimethylpentane	-	0.05	98.8
3-methyl-1-butanol	-	2.46	131.2
acetone	-	2.48	46.5
benzene	-	0.16	78.8
butanone	-	1.97	75.6
butyl acetate	-	1.57	126.6
carbon disulfide	-	0.07	46.2
carbon tetrachloride	-	0.08	76.0
chloroform	-	0.15	61.2
cyclohexane	-	0.05	80.7
decane	-	0.05	174.9
dibutyl ether	-	0.53	142.1
dichloromethane	-	0.20	39.6
diethyl ether	-	2.55	33.2
DMF	-	2.34	153.0
ethyl acetate	-	2.20	73.9
ethylene glycol	-	2.18	197.5
heptane	-	0.07	98.8
hexadecane	-	0.04	286.6
hexane	-	0.06	68.5
methyl acetate	-	1.60	44.0
methyl tert-butyl ether	-	1.28	55.2
methylcyclohexane	-	0.07	101.1
nonane	-	0.05	151.7
octane	-	0.06	126.4

undecane	-	0.04	196.3
water	-	0.24	100.0

ibuprofen $C_{13}H_{18}O_2$ - 10 solvents [905, 907]



Compound Data

Molecular weight	206.281	**H bond acceptors**	2	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	3.5
Phase 25°C	solid	**Rotatable bonds**	4	**Predicted density**	1.029 g/cm^3
SMILES	O=C(O)C(c1ccc(cc1)CC(C)C)C				
InChIKey	HEFNNWSXXWATRW-UHFFFAOYSA-N				

Abraham Solute Descriptors

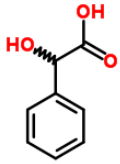
log S_w	E	S	A	B	V
-2.310	0.942	1.122	0.444	1.364	2.061

Solubility Data

	Concentration (M)		
Solvent	**Measured**	**Predicted**	**Solvent bp °C**
1-decanol	1.12	1.10	227.8
1-octanol	1.20	0.70	194.7
1-pentanol	1.46	1.74	138.5
1-propanol	1.52	1.56	95.8
2-butanol	1.78	1.59	96.6
2-methyl-1-propanol	1.76	1.59	105
2-propanol	2.22	1.57	73
ethanol	1.18	2.57	72.6
methanol	1.16	1.57	48.1
water	0.00	0.00	100
1-butanol	-	1.33	117.7
1-heptanol	-	1.44	176.9
1-hexanol	-	1.38	158.2
1,2-dichloroethane	-	0.33	83.5
1,4-dioxane	-	0.62	102.9
2-methyl-2-propanol	-	1.26	84.6
2-pentanol	-	2.39	118.8
2,2,4-trimethylpentane	-	0.00	98.8
3-methyl-1-butanol	-	2.05	131.2
acetone	-	0.34	46.5
acetonitrile	-	0.06	63.5
benzene	-	0.17	78.8
butanone	-	0.47	75.6
butyl acetate	-	0.23	126.6
carbon disulfide	-	0.01	46.2
carbon tetrachloride	-	0.06	76.0
chloroform	-	2.56	61.2
cyclohexane	-	0.01	80.7
decane	-	0.00	174.9
dibutyl ether	-	0.02	142.1
dichloromethane	-	0.74	39.6
diethyl ether	-	0.26	33.2
DMF	-	3.02	153.0
DMSO	-	0.96	189.0
ethyl acetate	-	0.26	73.9
ethylene glycol	-	0.80	197.5
heptane	-	0.01	98.8
hexadecane	-	0.00	286.6
hexane	-	0.00	68.5
methyl acetate	-	0.43	44.0
methyl tert-butyl ether	-	0.15	55.2
methylcyclohexane	-	0.02	101.1
nonane	-	0.00	151.7

octane	-	0.00	126.4
THF	-	1.36	68.3
toluene	-	0.09	110.6
undecane	-	0.00	196.3

mandelic acid $C_8H_8O_3$ - 5 solvents [72, 71]



Compound Data

Molecular weight	152.147	**H bond acceptors**	3	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	2	**ACD/ALogP**	0.66
Phase 25°C	solid	**Rotatable bonds**	3	**Predicted density**	1.321 g/cm^3
SMILES	O=C(O)C(O)c1ccccc1				
InChIKey	IWYDHOAUDWTVEP-UHFFFAOYSA-N				

Abraham Solute Descriptors

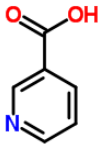
log S_w	E	S	A	B	V
-0.960	0.970	6.447	1.008	-0.527	0.812

Solubility Data

	Concentration (M)		
Solvent	**Measured**	**Predicted**	**Solvent bp °C**
1-hexanol	0.00	0.01	158.2
chloroform	0.12	0.10	61.2
diethyl ether	0.77	1.49	33.2
methanol	3.54	1.13	48.1
toluene	0.00	0.02	110.6
1-butanol	-	0.00	117.7
1-decanol	-	0.00	227.8
1-heptanol	-	0.00	176.9
1-octanol	-	0.02	194.7
1-pentanol	-	0.00	138.5
1-propanol	-	0.02	95.8
1,2-dichloroethane	-	> 9	83.5
1,4-dioxane	-	> 9	102.9
2-butanol	-	0.04	96.6
2-methyl-1-propanol	-	0.01	105.0
2-methyl-2-propanol	-	0.18	84.6
2-pentanol	-	0.00	118.8
2-propanol	-	0.05	73.0
2,2,4-trimethylpentane	-	0.00	98.8
3-methyl-1-butanol	-	0.00	131.2
acetone	-	> 9	46.5
acetonitrile	-	> 9	63.5
benzene	-	0.08	78.8
butanone	-	> 9	75.6
butyl acetate	-	> 9	126.6
carbon disulfide	-	0.00	46.2
carbon tetrachloride	-	0.00	76.0
cyclohexane	-	0.00	80.7
decane	-	0.00	174.9
dibutyl ether	-	0.03	142.1
dichloromethane	-	6.86	39.6
DMF	-	> 9	153.0
DMSO	-	> 9	189.0
ethyl acetate	-	> 9	73.9
ethylene glycol	-	0.21	197.5
heptane	-	0.00	98.8
hexadecane	-	0.00	286.6
hexane	-	0.00	68.5
methyl acetate	-	> 9	44.0
methyl tert-butyl ether	-	0.55	55.2

methylcyclohexane	-	0.00	101.1
nonane	-	0.00	151.7
octane	-	0.00	126.4
THF	-	> 9	68.3
undecane	-	0.00	196.3
water	-	0.11	100.0

nicotinic acid $C_6H_5NO_2$ - 8 solvents [100, 137, 907]



Compound Data

Molecular weight	123.109	**H bond acceptors**	3	**Rule of 5 violations**	0
Compound type	carboxlyic acid	**H bond donors**	1	**ACD/ALogP**	0.29
Phase 25°C	solid	**Rotatable bonds**	1	**Predicted density**	1.293 g/cm^3
SMILES	O=C(O)c1cccnc1				
InChIKey	PVNIIMVLHYAWGP-UHFFFAOYSA-N				

Abraham Solute Descriptors

log S_w	E	S	A	B	V
-0.432	0.762	0.738	0.461	0.681	0.519

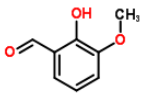
Solubility Data

	Concentration (M)		
Solvent	**Measured**	**Predicted**	**Solvent bp °C**
acetonitrile	0.00	0.02	63.5
benzene	0.00	0.00	78.8
DMSO	0.57	0.28	189
ethanol	0.09	0.11	72.6
methanol	0.06	0.15	48.1
THF	0.06	0.04	68.3
toluene	0.00	0.00	110.6
water	0.37	0.37	100
1-butanol	-	0.05	117.7
1-decanol	-	0.03	227.8
1-heptanol	-	0.03	176.9
1-hexanol	-	0.04	158.2
1-octanol	-	0.03	194.7
1-pentanol	-	0.05	138.5
1-propanol	-	0.07	95.8
1,2-dichloroethane	-	0.01	83.5
1,4-dioxane	-	0.03	102.9
2-butanol	-	0.09	96.6
2-methyl-1-propanol	-	0.06	105.0
2-methyl-2-propanol	-	0.05	84.6
2-pentanol	-	0.06	118.8
2-propanol	-	0.07	73.0
2,2,4-trimethylpentane	-	0.00	98.8
3-methyl-1-butanol	-	0.05	131.2
acetone	-	0.04	46.5
butanone	-	0.03	75.6
butyl acetate	-	0.01	126.6
carbon disulfide	-	0.00	46.2
carbon tetrachloride	-	0.00	76.0
chloroform	-	0.01	61.2
cyclohexane	-	0.00	80.7
decane	-	0.00	174.9
dibutyl ether	-	0.00	142.1
dichloromethane	-	0.01	39.6
diethyl ether	-	0.02	33.2
DMF	-	0.05	153.0
ethyl acetate	-	0.02	73.9
ethylene glycol	-	0.28	197.5
heptane	-	0.00	98.8

hexadecane	-	0.00	286.6
hexane	-	0.00	68.5
methyl acetate	-	0.03	44.0
methyl tert-butyl ether	-	0.01	55.2
methylcyclohexane	-	0.00	101.1
nonane	-	0.00	151.7
octane	-	0.00	126.4
undecane	-	0.00	196.3

o-vanillin $C_8H_8O_3$ - 5 solvents [208, 205]

Compound Data



Molecular weight	152.147	**H bond acceptors**	3	**Rule of 5 violations**	0
Compound type	aldehyde	**H bond donors**	1	**ACD/ALogP**	1.31
Phase 25°C	solid	**Rotatable bonds**	3	**Predicted density**	1.231 g/cm^3
SMILES	Oc1c(cccc1OC)C=O				
InChIKey	JJVNINGBHGBWJH-UHFFFAOYSA-N				

Abraham Solute Descriptors

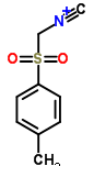
log S_w	E	S	A	B	V
-1.470	1.024	0.802	0.046	0.448	0.935

Solubility Data

	Concentration (M)		
Solvent	**Measured**	**Predicted**	**Solvent bp °C**
acetonitrile	2.56	2.47	63.5
ethanol	3.04	2.74	72.6
methanol	2.27	2.50	48.1
THF	5.37	5.53	68.3
toluene	3.82	3.87	110.6
1-butanol	-	1.96	117.7
1-decanol	-	1.87	227.8
1-heptanol	-	1.86	176.9
1-hexanol	-	1.96	158.2
1-octanol	-	1.61	194.7
1-pentanol	-	2.00	138.5
1-propanol	-	1.97	95.8
1,2-dichloroethane	-	5.84	83.5
1,4-dioxane	-	4.32	102.9
2-butanol	-	2.10	96.6
2-methyl-1-propanol	-	1.77	105.0
2-methyl-2-propanol	-	1.54	84.6
2-pentanol	-	2.02	118.8
2-propanol	-	1.56	73.0
2,2,4-trimethylpentane	-	0.64	98.8
3-methyl-1-butanol	-	1.91	131.2
acetone	-	3.95	46.5
benzene	-	4.64	78.8
butanone	-	4.41	75.6
butyl acetate	-	3.01	126.6
carbon disulfide	-	2.27	46.2
carbon tetrachloride	-	2.71	76.0
chloroform	-	7.93	61.2
cyclohexane	-	1.12	80.7
decane	-	0.57	174.9
dibutyl ether	-	1.16	142.1
dichloromethane	-	7.51	39.6
diethyl ether	-	2.61	33.2
DMF	-	2.98	153.0
DMSO	-	2.98	189.0
ethyl acetate	-	3.40	73.9

ethylene glycol	-	0.83	197.5
heptane	-	0.73	98.8
hexadecane	-	0.64	286.6
hexane	-	0.74	68.5
methyl acetate	-	4.26	44.0
methyl tert-butyl ether	-	2.33	55.2
methylcyclohexane	-	1.38	101.1
nonane	-	0.72	151.7
octane	-	0.78	126.4
undecane	-	0.55	196.3
water	-	0.03	100.0

p-toluenesulfonylmethyl isocyanide $C_9H_9NO_2S$ - 5 solvents [208]



Compound Data

Molecular weight	195.238	**H bond acceptors**	3	**Rule of 5 violations**	0
Compound type	isonitrile	**H bond donors**	0	**ACD/ALogP**	1.38
Phase 25°C	solid	**Rotatable bonds**	2	**Predicted density**	‡
SMILES	O=S(=O)(c1ccc(cc1)C)C[N+]#[C-]				
InChIKey	CFOAUYCPAUGDFF-UHFFFAOYSA-N				

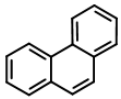
Solubility Data

	Concentration (M)		
Solvent	**Measured**	**Predicted**	**Solvent bp °C**
acetonitrile	1.88	0.03	63.5
chloroform	0.21	0.03	61.2
ethanol	0.18	0.03	72.6
THF	1.86	0.03	68.3
toluene	0.14	0.03	110.6
1-butanol	-	1.96	117.7
1-decanol	-	1.87	227.8
1-heptanol	-	1.86	176.9
1-hexanol	-	1.96	158.2
1-octanol	-	1.61	194.7
1-pentanol	-	2.00	138.5
1-propanol	-	1.97	95.8
1,2-dichloroethane	-	5.84	83.5
1,4-dioxane	-	4.32	102.9
2-butanol	-	2.10	96.6
2-methyl-1-propanol	-	1.77	105.0
2-methyl-2-propanol	-	1.54	84.6
2-pentanol	-	2.02	118.8
2-propanol	-	1.56	73.0
2,2,4-trimethylpentane	-	0.64	98.8
3-methyl-1-butanol	-	1.91	131.2
acetone	-	3.95	46.5
benzene	-	4.64	78.8
butanone	-	4.41	75.6
butyl acetate	-	3.01	126.6
carbon disulfide	-	2.27	46.2
carbon tetrachloride	-	2.71	76.0
cyclohexane	-	1.12	80.7
decane	-	0.57	174.9
dibutyl ether	-	1.16	142.1
dichloromethane	-	7.51	39.6
diethyl ether	-	2.61	33.2
DMF	-	2.98	153.0
DMSO	-	2.98	189.0
ethyl acetate	-	3.40	73.9
ethylene glycol	-	0.83	197.5
heptane	-	0.73	98.8
hexadecane	-	0.64	286.6

hexane	-	0.74	68.5
methanol	-	2.50	48.1
methyl acetate	-	4.26	44.0
methyl tert-butyl ether	-	2.33	55.2
methylcyclohexane	-	1.38	101.1
nonane	-	0.72	151.7
octane	-	0.78	126.4
undecane	-	0.55	196.3
water	-	0.03	100.0

‡ Solute density prediction fail.

phenanthrene $C_{14}H_{10}$ - 6 solvents [902]



Compound Data

Molecular weight	178.229	**H bond acceptors**	0	**Rule of 5 violations**	0
Compound type	non-Ugi related	**H bond donors**	0	**ACD/ALogP**	4.55
Phase 25°C	solid	**Rotatable bonds**	0	**Predicted density**	1.13 g/cm^3
SMILES	c3cc2ccc1ccccc1c2cc3				
InChIKey	YNPNZTXNASCQKK-UHFFFAOYSA-N				

Abraham Solute Descriptors

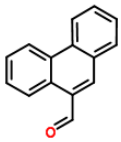
log S_w	E	S	A	B	V
-5.720	2.100	0.840	0.057	0.198	1.435

Solubility Data

	Concentration (M)		
Solvent	**Measured**	**Predicted**	**Solvent bp °C**
benzene	2.02	1.82	78.8
carbon disulfide	3.05	4.78	46.2
carbon tetrachloride	1.73	1.31	76
diethyl ether	1.36	0.91	33.2
ethanol	0.21	0.29	72.6
hexane	0.32	0.31	68.5
1-butanol	-	0.32	117.7
1-decanol	-	0.86	227.8
1-heptanol	-	0.55	176.9
1-hexanol	-	0.44	158.2
1-octanol	-	0.41	194.7
1-pentanol	-	0.44	138.5
1-propanol	-	0.29	95.8
1,2-dichloroethane	-	0.90	83.5
1,4-dioxane	-	1.10	102.9
2-butanol	-	0.18	96.6
2-methyl-1-propanol	-	0.18	105.0
2-methyl-2-propanol	-	0.13	84.6
2-pentanol	-	0.34	118.8
2-propanol	-	0.17	73.0
2,2,4-trimethylpentane	-	0.30	98.8
3-methyl-1-butanol	-	0.27	131.2
acetone	-	0.67	46.5
acetonitrile	-	0.10	63.5
butanone	-	0.88	75.6
butyl acetate	-	0.92	126.6
chloroform	-	0.61	61.2
cyclohexane	-	1.13	80.7
decane	-	0.39	174.9
dibutyl ether	-	0.58	142.1
dichloromethane	-	0.75	39.6
DMF	-	0.43	153.0
DMSO	-	0.27	189.0
ethyl acetate	-	0.90	73.9

ethylene glycol	-	0.02	197.5
heptane	-	0.47	98.8
hexadecane	-	0.40	286.6
methanol	-	0.11	48.1
methyl acetate	-	0.52	44.0
methyl tert-butyl ether	-	0.66	55.2
methylcyclohexane	-	0.89	101.1
nonane	-	0.43	151.7
octane	-	0.58	126.4
THF	-	2.14	68.3
toluene	-	2.09	110.6
undecane	-	0.43	196.3
water	-	0.00	100.0

phenanthrene-9-carboxaldehyde $C_{15}H_{10}O$ - 16 solvents [30, 60, 136, 4, 242, 51, 24]



Compound Data

Molecular weight	206.239	**H bond acceptors**	1	**Rule of 5 violations**	0
Compound type	aldehyde	**H bond donors**	0	**ACD/ALogP**	4.28
Phase 25°C	solid	**Rotatable bonds**	1	**Predicted density**	1.217 g/cm^3
SMILES	O=Cc2cc3c(c1c2cccc1)cccc3				
InChIKey	QECIGCMPORCORE-UHFFFAOYSA-N				

Abraham Solute Descriptors

log S_w	E	S	A	B	V
-5.520	2.494	1.166	0.327	0.423	1.606

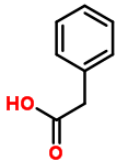
Solubility Data

	Concentration (M)		
Solvent	**Measured**	**Predicted**	**Solvent bp °C**
1,1,2-trichlorotrifluoroethane	0.02	*	50.9
2-propanol	0.07	0.15	73
acetonitrile	0.16	0.03	63.5
benzene	0.66	0.23	78.8
chloroform	0.04	0.10	61.2
cyclohexane	0.07	0.05	80.7
cyclopentane	0.03	*	49.2
dichloromethane	0.00	0.11	39.6
diethyl ether	0.10	0.36	33.2
DMF	1.25	0.50	153
DMSO	0.77	0.83	189
ethanol	0.10	0.26	72.6
hexane	0.07	0.01	68.5
methanol	0.11	0.10	48.1
THF	2.32	1.38	68.3
toluene	0.14	0.24	110.6
1-butanol	-	0.22	117.7
1-decanol	-	0.63	227.8
1-heptanol	-	0.36	176.9
1-hexanol	-	0.31	158.2
1-octanol	-	0.27	194.7
1-pentanol	-	0.34	138.5
1-propanol	-	0.25	95.8
1,2-dichloroethane	-	0.17	83.5
1,4-dioxane	-	0.68	102.9
2-butanol	-	0.15	96.6
2-methyl-1-propanol	-	0.13	105.0
2-methyl-2-propanol	-	0.09	84.6
2-pentanol	-	0.26	118.8
2,2,4-trimethylpentane	-	0.01	98.8
3-methyl-1-butanol	-	0.18	131.2
acetone	-	0.36	46.5

butanone	-	0.43	75.6
butyl acetate	-	0.33	126.6
carbon disulfide	-	0.25	46.2
carbon tetrachloride	-	0.09	76.0
decane	-	0.01	174.9
dibutyl ether	-	0.09	142.1
ethyl acetate	-	0.37	73.9
ethylene glycol	-	0.05	197.5
heptane	-	0.02	98.8
hexadecane	-	0.02	286.6
methyl acetate	-	0.22	44.0
methyl tert-butyl ether	-	0.16	55.2
methylcyclohexane	-	0.05	101.1
nonane	-	0.02	151.7
octane	-	0.02	126.4
undecane	-	0.02	196.3
water	-	0.00	100.0

* This solvent doesn't have recorded Abraham coefficients.

phenylacetic acid $C_8H_8O_2$ - 7 solvents [86, 123, 907]



Compound Data

Molecular weight	136.148	**H bond acceptors**	2	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	1	**ACD/ALogP**	1.72
Phase 25°C	solid	**Rotatable bonds**	2	**Predicted density**	1.164 g/cm^3
SMILES	O=C(O)Cc1ccccc1				
InChIKey	WLJVXDMOQOGPHL-UHFFFAOYSA-N				

Abraham Solute Descriptors

log S_w	E	S	A	B	V
-0.886	0.832	0.917	0.082	0.455	0.837

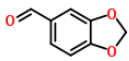
Solubility Data

	Concentration (M)		
Solvent	**Measured**	**Predicted**	**Solvent bp °C**
acetonitrile	3.87	3.78	63.5
benzene	3.35	3.22	78.8
DMF	3.76	4.39	153
DMSO	6.35	5.83	189
THF	5.62	5.32	68.3
toluene	1.94	2.48	110.6
water	0.13	0.13	100
1-butanol	-	1.67	117.7
1-decanol	-	1.23	227.8
1-heptanol	-	1.41	176.9
1-hexanol	-	1.61	158.2
1-octanol	-	1.34	194.7
1-pentanol	-	1.58	138.5
1-propanol	-	1.79	95.8
1,2-dichloroethane	-	5.41	83.5
1,4-dioxane	-	4.73	102.9
2-butanol	-	2.10	96.6
2-methyl-1-propanol	-	1.70	105.0
2-methyl-2-propanol	-	1.64	84.6
2-pentanol	-	1.64	118.8
2-propanol	-	1.56	73.0
2,2,4-trimethylpentane	-	0.30	98.8
3-methyl-1-butanol	-	1.62	131.2
acetone	-	4.57	46.5
butanone	-	4.98	75.6
butyl acetate	-	2.79	126.6

carbon disulfide	-	1.10	46.2
carbon tetrachloride	-	1.46	76.0
chloroform	-	7.00	61.2
cyclohexane	-	0.46	80.7
decane	-	0.26	174.9
dibutyl ether	-	0.85	142.1
dichloromethane	-	7.10	39.6
diethyl ether	-	2.24	33.2
ethanol	-	2.66	72.6
ethyl acetate	-	3.33	73.9
ethylene glycol	-	1.09	197.5
heptane	-	0.30	98.8
hexadecane	-	0.30	286.6
hexane	-	0.35	68.5
methanol	-	3.01	48.1
methyl acetate	-	4.90	44.0
methyl tert-butyl ether	-	1.99	55.2
methylcyclohexane	-	0.55	101.1
nonane	-	0.33	151.7
octane	-	0.36	126.4
undecane	-	0.25	196.3

piperonal $C_8H_6O_3$ - 11 solvents [71, 30, 907]

Compound Data



Molecular weight	150.131	**H bond acceptors**	3	**Rule of 5 violations**	0
Compound type	aldehyde	**H bond donors**	0	**ACD/ALogP**	0.79
Phase 25°C	solid	**Rotatable bonds**	1	**Predicted density**	1.337 g/cm^3
SMILES	O=Cc1ccc2OCOc2c1				
InChIKey	SATCULPHIDQDRE-UHFFFAOYSA-N				

Abraham Solute Descriptors

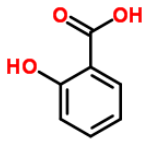
log S_w	E	S	A	B	V
-1.699	1.032	0.928	0.118	0.175	0.769

Solubility Data

	Concentration (M)		
Solvent	**Measured**	**Predicted**	**Solvent bp °C**
acetonitrile	7.11	5.37	63.5
chloroform	7.95	4.22	61.2
dichloromethane	7.59	6.33	39.6
diethyl ether	6.38	4.90	33.2
DMSO	6.05	13.10	189
ethanol	4.99	2.80	72.6
hexane	0.15	0.55	68.5
methanol	7.29	2.71	48.1
THF	7.44	11.41	68.3
toluene	6.40	4.13	110.6
water	0.02	0.02	100
1-butanol	-	2.04	117.7
1-decanol	-	1.83	227.8
1-heptanol	-	1.94	176.9
1-hexanol	-	2.13	158.2
1-octanol	-	2.01	194.7
1-pentanol	-	1.97	138.5
1-propanol	-	2.23	95.8
1,2-dichloroethane	-	6.26	83.5
1,4-dioxane	-	> 9	102.9
2-butanol	-	2.15	96.6
2-methyl-1-propanol	-	1.81	105.0
2-methyl-2-propanol	-	1.92	84.6
2-pentanol	-	1.78	118.8

2-propanol	-	1.83	73.0
2,2,4-trimethylpentane	-	0.49	98.8
3-methyl-1-butanol	-	1.72	131.2
acetone	-	8.95	46.5
benzene	-	4.64	78.8
butanone	-	> 9	75.6
butyl acetate	-	5.91	126.6
carbon disulfide	-	3.33	46.2
carbon tetrachloride	-	1.96	76.0
cyclohexane	-	0.85	80.7
decane	-	0.49	174.9
dibutyl ether	-	2.20	142.1
DMF	-	7.86	153.0
ethyl acetate	-	7.03	73.9
ethylene glycol	-	0.75	197.5
heptane	-	0.48	98.8
hexadecane	-	0.51	286.6
methyl acetate	-	7.65	44.0
methyl tert-butyl ether	-	4.01	55.2
methylcyclohexane	-	0.68	101.1
nonane	-	0.58	151.7
octane	-	0.70	126.4
undecane	-	0.49	196.3

salicylic acid $C_7H_6O_3$ - 6 solvents [34, 902, 901, 138, 52, 907]



Compound Data

Molecular weight	138.121	**H bond acceptors**	3	**Rule of 5 violations**	0
Compound type	carboxylic acid	**H bond donors**	2	**ACD/ALogP**	2.06
Phase 25°C	solid	**Rotatable bonds**	2	**Predicted density**	1.375 g/cm^3
SMILES	O=C(O)c1ccccc1O				
InChIKey	YGSDEFSMJLZEOE-UHFFFAOYSA-N				

Abraham Solute Descriptors

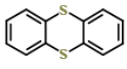
log S_w	E	S	A	B	V
-2.000	0.920	0.110	0.410	0.111	0.595

Solubility Data

	Concentration (M)		
Solvent	**Measured**	**Predicted**	**Solvent bp °C**
chloroform	0.19	0.16	61.2
dichloromethane	0.15	0.18	39.6
ethanol	2.34	3.07	72.6
methanol	2.65	1.89	48.1
toluene	0.11	0.30	110.6
water	0.01	0.01	100
1-butanol	-	3.03	117.7
1-decanol	-	3.98	227.8
1-heptanol	-	2.93	176.9
1-hexanol	-	2.89	158.2
1-octanol	-	2.14	194.7
1-pentanol	-	3.73	138.5
1-propanol	-	3.60	95.8
1,2-dichloroethane	-	0.20	83.5
1,4-dioxane	-	1.34	102.9
2-butanol	-	2.36	96.6
2-methyl-1-propanol	-	2.44	105.0
2-methyl-2-propanol	-	2.24	84.6
2-pentanol	-	3.59	118.8
2-propanol	-	2.89	73.0
2,2,4-trimethylpentane	-	0.16	98.8

3-methyl-1-butanol	-	2.76	131.2
acetone	-	1.43	46.5
acetonitrile	-	0.25	63.5
benzene	-	0.27	78.8
butanone	-	1.22	75.6
butyl acetate	-	1.44	126.6
carbon disulfide	-	0.24	46.2
carbon tetrachloride	-	0.21	76.0
cyclohexane	-	0.22	80.7
decane	-	0.18	174.9
dibutyl ether	-	0.81	142.1
diethyl ether	-	2.76	33.2
DMF	-	0.90	153.0
DMSO	-	1.60	189.0
ethyl acetate	-	1.80	73.9
ethylene glycol	-	0.88	197.5
heptane	-	0.34	98.8
hexadecane	-	0.14	286.6
hexane	-	0.19	68.5
methyl acetate	-	0.94	44.0
methyl tert-butyl ether	-	1.43	55.2
methylcyclohexane	-	0.33	101.1
nonane	-	0.20	151.7
octane	-	0.22	126.4
THF	-	3.21	68.3
undecane	-	0.16	196.3

thianthrene $C_{12}H_8S_2$ - 10 solvents [905]



Compound Data

Molecular weight	216.322	**H bond acceptors**	0	**Rule of 5 violations**	0
Compound type	non-Ugi related	**H bond donors**	0	**ACD/ALogP**	4.57
Phase 25°C	solid	**Rotatable bonds**	0	**Predicted density**	1.31 g/cm^3
SMILES	S1c3c(Sc2c1cccc2)cccc3				
InChIKey	GVIJJXMXTUZIOD-UHFFFAOYSA-N				

Abraham Solute Descriptors

log S_w	E	S	A	B	V
-4.290	2.281	1.312	-0.101	0.786	1.543

Solubility Data

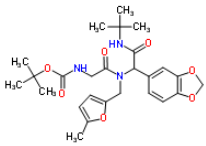
	Concentration (M)		
Solvent	**Measured**	**Predicted**	**Solvent bp °C**
1-decanol	0.04	0.08	227.8
1,4-dioxane	0.27	0.16	102.9
acetonitrile	0.02	0.04	63.5
decane	0.03	0.02	174.9
diethyl ether	0.08	0.05	33.2
diisopropyl ether	0.05	*	68.3
ethylene glycol	0.02	0.02	197.5
hexadecane	0.03	0.04	286.6
nonane	0.03	0.03	151.7
undecane	0.03	0.03	196.3
1-butanol	-	0.04	117.7
1-heptanol	-	0.05	176.9
1-hexanol	-	0.05	158.2
1-octanol	-	0.04	194.7
1-pentanol	-	0.05	138.5
1-propanol	-	0.03	95.8
1,2-dichloroethane	-	0.56	83.5
2-butanol	-	0.04	96.6
2-methyl-1-propanol	-	0.03	105.0

2-methyl-2-propanol	-	0.01	84.6
2-pentanol	-	0.04	118.8
2-propanol	-	0.02	73.0
2,2,4-trimethylpentane	-	0.02	98.8
3-methyl-1-butanol	-	0.04	131.2
acetone	-	0.10	46.5
benzene	-	0.55	78.8
butanone	-	0.12	75.6
butyl acetate	-	0.08	126.6
carbon disulfide	-	0.29	46.2
carbon tetrachloride	-	0.29	76.0
chloroform	-	0.88	61.2
cyclohexane	-	0.10	80.7
dibutyl ether	-	0.02	142.1
dichloromethane	-	0.60	39.6
DMF	-	0.06	153.0
DMSO	-	0.06	189.0
ethanol	-	0.06	72.6
ethyl acetate	-	0.08	73.9
heptane	-	0.03	98.8
hexane	-	0.03	68.5
methanol	-	0.04	48.1
methyl acetate	-	0.11	44.0
methyl tert-butyl ether	-	0.04	55.2
methylcyclohexane	-	0.13	101.1
octane	-	0.04	126.4
THF	-	0.18	68.3
toluene	-	0.43	110.6
water	-	0.00	100.0

* This solvent doesn't have recorded Abraham coefficients.

UCExp216-3A $C_{26}H_{35}N_3O_7$ - 5 solvents [113]

Compound Data



Molecular weight	501.572	**H bond acceptors**	10	**Rule of 5 violations**	2
Compound type	Ugi Product	**H bond donors**	2	**ACD/ALogP**	4.06
Phase 25°C	solid	**Rotatable bonds**	10	**Predicted density**	1.209 g/cm^3
SMILES	CC(C)(C)NC(=O)C(c1ccc2OCOc2c1)N(Cc3ccc(C)o3)C(=O)CNC(=O)OC(C)(C)C				
InChIKey	MTSJLRXXXJHLDI-UHFFFAOYSA-N				

Abraham Solute Descriptors

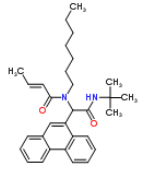
log S_w	E	S	A	B	V
-4.050	2.332	4.791	-0.007	3.981	5.203

Solubility Data

	Concentration (M)		
Solvent	**Measured**	**Predicted**	**Solvent bp °C**
acetonitrile	0.04	0.01	63.5
DMSO	0.28	0.52	189
ethanol	0.03	0.03	72.6
methanol	0.06	0.03	48.1
toluene	0.01	0.02	110.6
1-butanol	-	0.00	117.7
1-decanol	-	0.00	227.8
1-heptanol	-	0.00	176.9
1-hexanol	-	0.00	158.2
1-octanol	-	0.00	194.7
1-pentanol	-	0.00	138.5
1-propanol	-	0.00	95.8

1,2-dichloroethane	-	7.32	83.5
1,4-dioxane	-	0.09	102.9
2-butanol	-	0.01	96.6
2-methyl-1-propanol	-	0.01	105.0
2-methyl-2-propanol	-	0.00	84.6
2-pentanol	-	0.00	118.8
2-propanol	-	0.00	73.0
2,2,4-trimethylpentane	-	0.00	98.8
3-methyl-1-butanol	-	0.01	131.2
acetone	-	0.01	46.5
benzene	-	0.22	78.8
butanone	-	0.05	75.6
butyl acetate	-	0.00	126.6
carbon disulfide	-	0.00	46.2
carbon tetrachloride	-	0.00	76.0
chloroform	-	> 9	61.2
cyclohexane	-	0.00	80.7
decane	-	0.00	174.9
dibutyl ether	-	0.00	142.1
dichloromethane	-	> 9	39.6
diethyl ether	-	0.00	33.2
DMF	-	> 9	153.0
ethyl acetate	-	0.00	73.9
ethylene glycol	-	0.03	197.5
heptane	-	0.00	98.8
hexadecane	-	0.00	286.6
hexane	-	0.00	68.5
methyl acetate	-	0.06	44.0
methyl tert-butyl ether	-	0.00	55.2
methylcyclohexane	-	0.00	101.1
nonane	-	0.00	151.7
octane	-	0.00	126.4
THF	-	0.04	68.3
undecane	-	0.00	196.3
water	-	0.00	100.0

Ugi product 148B $C_{31}H_{40}N_2O_2$ - 6 solvents [145, 96]



Compound Data

Molecular weight	472.661	**H bond acceptors**	4	**Rule of 5 violations**	1
Compound type	Ugi Product	**H bond donors**	1	**ACD/ALogP**	7.47
Phase 25°C	solid	**Rotatable bonds**	11	**Predicted density**	1.067 g/cm^3
SMILES	CC(C)(C)NC(=O)C(N(CCCCCCC)C(=O)\C=C\C)c2cc3ccccc3c1ccccc12				
InChIKey	HOKCUULFCASAQT-FRKPEAEDSA-N				

Abraham Solute Descriptors

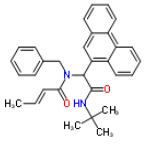
log S_w	E	S	A	B	V
-7.620	3.440	5.683	0.240	3.436	5.615

Solubility Data

	Concentration (M)		
Solvent	**Measured**	**Predicted**	**Solvent bp °C**
acetonitrile	0.02	0.02	63.5
benzene	0.05	0.24	78.8
ethanol	0.00	0.01	72.6
methanol	0.03	0.01	48.1
THF	0.40	0.39	68.3
toluene	0.14	0.03	110.6
1-butanol	-	0.00	117.7
1-decanol	-	0.00	227.8
1-heptanol	-	0.00	176.9
1-hexanol	-	0.00	158.2

1-octanol	-	0.00	194.7
1-pentanol	-	0.00	138.5
1-propanol	-	0.00	95.8
1,2-dichloroethane	-	7.34	83.5
1,4-dioxane	-	0.83	102.9
2-butanol	-	0.00	96.6
2-methyl-1-propanol	-	0.00	105.0
2-methyl-2-propanol	-	0.00	84.6
2-pentanol	-	0.00	118.8
2-propanol	-	0.00	73.0
2,2,4-trimethylpentane	-	0.00	98.8
3-methyl-1-butanol	-	0.00	131.2
acetone	-	0.05	46.5
butanone	-	0.30	75.6
butyl acetate	-	0.01	126.6
carbon disulfide	-	0.00	46.2
carbon tetrachloride	-	0.00	76.0
chloroform	-	> 9	61.2
cyclohexane	-	0.00	80.7
decane	-	0.00	174.9
dibutyl ether	-	0.00	142.1
dichloromethane	-	> 9	39.6
diethyl ether	-	0.00	33.2
DMF	-	> 9	153.0
DMSO	-	> 9	189.0
ethyl acetate	-	0.01	73.9
ethylene glycol	-	0.00	197.5
heptane	-	0.00	98.8
hexadecane	-	0.00	286.6
hexane	-	0.00	68.5
methyl acetate	-	0.14	44.0
methyl tert-butyl ether	-	0.00	55.2
methylcyclohexane	-	0.00	101.1
nonane	-	0.00	151.7
octane	-	0.00	126.4
undecane	-	0.00	196.3
water	-	0.00	100.0

Ugi product 150D(UC) $C_{31}H_{32}N_2O_2$ - 6 solvents [78, 65]



Compound Data

Molecular weight	464.598	**H bond acceptors**	4	**Rule of 5 violations**	1
Compound type	Ugi Product	**H bond donors**	1	**ACD/ALogP**	6.18
Phase 25°C	solid	**Rotatable bonds**	7	**Predicted density**	1.148 g/cm^3
SMILES	CC(C)(C)NC(=O)C(N(Cc1ccccc1)C(=O)\C=C\C)c3cc4ccccc4c2ccccc23				
InChIKey	PBZQTKRWYXTXIS-WLRTZDKTSA-N				

Abraham Solute Descriptors

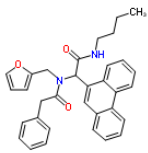
log S_w	E	S	A	B	V
-7.160	4.342	4.362	0.051	3.236	5.052

Solubility Data

	Concentration (M)		
Solvent	**Measured**	**Predicted**	**Solvent bp °C**
acetonitrile	0.04	0.01	63.5
DMSO	0.18	0.74	189
ethanol	0.04	0.05	72.6
methanol	0.07	0.02	48.1
THF	0.49	0.27	68.3
toluene	0.04	0.21	110.6
1-butanol	-	0.01	117.7

1-decanol	-	0.02	227.8
1-heptanol	-	0.02	176.9
1-hexanol	-	0.02	158.2
1-octanol	-	0.01	194.7
1-pentanol	-	0.01	138.5
1-propanol	-	0.01	95.8
1,2-dichloroethane	-	6.33	83.5
1,4-dioxane	-	0.34	102.9
2-butanol	-	0.02	96.6
2-methyl-1-propanol	-	0.01	105.0
2-methyl-2-propanol	-	0.00	84.6
2-pentanol	-	0.02	118.8
2-propanol	-	0.01	73.0
2,2,4-trimethylpentane	-	0.00	98.8
3-methyl-1-butanol	-	0.01	131.2
acetone	-	0.03	46.5
benzene	-	0.99	78.8
butanone	-	0.11	75.6
butyl acetate	-	0.01	126.6
carbon disulfide	-	0.00	46.2
carbon tetrachloride	-	0.03	76.0
chloroform	-	> 9	61.2
cyclohexane	-	0.00	80.7
decane	-	0.00	174.9
dibutyl ether	-	0.00	142.1
dichloromethane	-	> 9	39.6
diethyl ether	-	0.00	33.2
DMF	-	5.21	153.0
ethyl acetate	-	0.01	73.9
ethylene glycol	-	0.03	197.5
heptane	-	0.00	98.8
hexadecane	-	0.00	286.6
hexane	-	0.00	68.5
methyl acetate	-	0.07	44.0
methyl tert-butyl ether	-	0.00	55.2
methylcyclohexane	-	0.00	101.1
nonane	-	0.00	151.7
octane	-	0.00	126.4
undecane	-	0.00	196.3
water	-	0.00	100.0

Ugi product 176C (UC) $C_{33}H_{32}N_2O_3$ - 11 solvents [155, 127, 144]



Compound Data

Molecular weight	504.619	**H bond acceptors**	5	**Rule of 5 violations**	2
Compound type	Ugi Product	**H bond donors**	1	**ACD/ALogP**	6.9
Phase 25°C	solid	**Rotatable bonds**	10	**Predicted density**	1.197 g/cm^3
SMILES	O=C(Cc1ccccc1)N(Cc2ccco2)C(c4cc5ccccc5c3ccccc34)C(=O)NCCCC				
InChIKey	RGGYLRDGJAJICY-UHFFFAOYSA-N				

Abraham Solute Descriptors

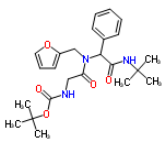
log S_w	E	S	A	B	V
-5.310	4.713	4.490	0.151	3.935	5.298

Solubility Data

	Concentration (M)		
Solvent	**Measured**	**Predicted**	**Solvent bp °C**
1-octanol	0.01	0.01	194.7
1,4-dioxane	0.19	0.12	102.9
acetonitrile	0.02	0.00	63.5
benzene	0.03	0.33	78.8
diethyl ether	0.02	0.00	33.2

DMF	0.27	2.14	153
DMSO	0.19	0.53	189
ethanol	0.01	0.10	72.6
methanol	0.02	0.05	48.1
THF	0.28	0.10	68.3
toluene	0.03	0.05	110.6
1-butanol	-	0.01	117.7
1-decanol	-	0.03	227.8
1-heptanol	-	0.02	176.9
1-hexanol	-	0.02	158.2
1-pentanol	-	0.02	138.5
1-propanol	-	0.01	95.8
1,2-dichloroethane	-	3.14	83.5
2-butanol	-	0.03	96.6
2-methyl-1-propanol	-	0.01	105.0
2-methyl-2-propanol	-	0.00	84.6
2-pentanol	-	0.03	118.8
2-propanol	-	0.01	73.0
2,2,4-trimethylpentane	-	0.00	98.8
3-methyl-1-butanol	-	0.03	131.2
acetone	-	0.01	46.5
butanone	-	0.03	75.6
butyl acetate	-	0.00	126.6
carbon disulfide	-	0.00	46.2
carbon tetrachloride	-	0.01	76.0
chloroform	-	> 9	61.2
cyclohexane	-	0.00	80.7
decane	-	0.00	174.9
dibutyl ether	-	0.00	142.1
dichloromethane	-	> 9	39.6
ethyl acetate	-	0.00	73.9
ethylene glycol	-	0.34	197.5
heptane	-	0.00	98.8
hexadecane	-	0.00	286.6
hexane	-	0.00	68.5
methyl acetate	-	0.03	44.0
methyl tert-butyl ether	-	0.00	55.2
methylcyclohexane	-	0.00	101.1
nonane	-	0.00	151.7
octane	-	0.00	126.4
undecane	-	0.00	196.3
water	-	0.00	100.0

Ugi product 206B(UC) $C_{24}H_{33}N_3O_5$ - 16 solvents [56]



Compound Data

Molecular weight	443.536	**H bond acceptors**	8	**Rule of 5 violations**	0
Compound type	Ugi Product	**H bond donors**	2	**ACD/ALogP**	3.74
Phase 25°C	solid	**Rotatable bonds**	10	**Predicted density**	1.145 g/cm^3
SMILES	O=C(CNC(=O)OC(C)(C)C)N(Cc1ccco1)C(c2ccccc2)C(=O)NC(C)(C)C				
InChIKey	BXAOUWIVXLQYOZ-UHFFFAOYSA-N				

Abraham Solute Descriptors

log S_w	E	S	A	B	V
-4.240	1.915	2.463	0.301	3.663	4.800

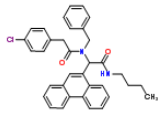
Solubility Data

	Concentration (M)		
Solvent	**Measured**	**Predicted**	**Solvent bp °C**
2-propanol	0.00	0.18	73.0
acetonitrile	0.02	0.00	63.5

benzene	0.01	0.11	78.8
carbon tetrachloride	0.00	0.02	76
chloroform	0.57	279.75	61.2
cyclohexane	0.00	0.00	80.7
cyclopentane	0.00	*	49.2
dichloromethane	0.24	4.64	39.6
diethyl ether	0.00	0.00	33.2
DMF	0.24	1.01	153
DMSO	0.23	0.01	189
ethanol	0.02	0.88	72.6
hexane	0.00	0.00	68.5
methanol	0.05	0.30	48.1
THF	0.26	0.08	68.3
toluene	0.01	0.02	110.6
1-butanol	-	0.20	117.7
1-decanol	-	0.21	227.8
1-heptanol	-	0.36	176.9
1-hexanol	-	0.26	158.2
1-octanol	-	0.07	194.7
1-pentanol	-	0.39	138.5
1-propanol	-	0.18	95.8
1,2-dichloroethane	-	0.48	83.5
1,4-dioxane	-	0.03	102.9
2-butanol	-	0.35	96.6
2-methyl-1-propanol	-	0.32	105.0
2-methyl-2-propanol	-	0.11	84.6
2-pentanol	-	0.94	118.8
2,2,4-trimethylpentane	-	0.00	98.8
3-methyl-1-butanol	-	0.79	131.2
acetone	-	0.00	46.5
butanone	-	0.02	75.6
butyl acetate	-	0.00	126.6
carbon disulfide	-	0.00	46.2
decane	-	0.00	174.9
dibutyl ether	-	0.00	142.1
ethyl acetate	-	0.00	73.9
ethylene glycol	-	0.28	197.5
heptane	-	0.00	98.8
hexadecane	-	0.00	286.6
methyl acetate	-	0.02	44.0
methyl tert-butyl ether	-	0.00	55.2
methylcyclohexane	-	0.00	101.1
nonane	-	0.00	151.7
octane	-	0.00	126.4
undecane	-	0.00	196.3
water	-	0.00	100.0

* This solvent doesn't have recorded Abraham coefficients.

Ugi product 215F(UC) $C_{35}H_{33}ClN_2O_2$ - 6 solvents [71]



Compound Data

Molecular weight	549.102	**H bond acceptors**	4	**Rule of 5 violations**	2
Compound type	Ugi Product	**H bond donors**	1	**ACD/ALogP**	8.33
Phase 25°C	solid	**Rotatable bonds**	10	**Predicted density**	1.219 g/cm^3
SMILES	Clc5ccc(CC(=O)N(Cc1ccccc1)C(c3cc4ccccc4c2ccccc23)C(=O)NCCCC)cc5				
InChIKey	NCOSMKWGNCJPCR-UHFFFAOYSA-N				

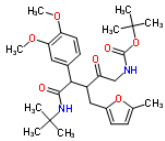
Abraham Solute Descriptors

log S_W	E	S	A	B	V
-7.580	5.307	4.975	0.047	3.742	5.726

Solubility Data

	Concentration (M)		
Solvent	**Measured**	**Predicted**	**Solvent bp °C**
acetonitrile	0.04	0.00	63.5
DMSO	0.38	1.63	189
ethanol	0.06	0.08	72.6
methanol	0.12	0.03	48.1
THF	0.61	0.44	68.3
toluene	0.07	0.41	110.6
1-butanol	-	0.01	117.7
1-decanol	-	0.03	227.8
1-heptanol	-	0.03	176.9
1-hexanol	-	0.02	158.2
1-octanol	-	0.01	194.7
1-pentanol	-	0.02	138.5
1-propanol	-	0.01	95.8
1,2-dichloroethane	-	> 9	83.5
1,4-dioxane	-	0.58	102.9
2-butanol	-	0.02	96.6
2-methyl-1-propanol	-	0.01	105.0
2-methyl-2-propanol	-	0.00	84.6
2-pentanol	-	0.02	118.8
2-propanol	-	0.01	73.0
2,2,4-trimethylpentane	-	0.00	98.8
3-methyl-1-butanol	-	0.02	131.2
acetone	-	0.04	46.5
benzene	-	2.33	78.8
butanone	-	0.15	75.6
butyl acetate	-	0.01	126.6
carbon disulfide	-	0.00	46.2
carbon tetrachloride	-	0.05	76.0
chloroform	-	> 9	61.2
cyclohexane	-	0.00	80.7
decane	-	0.00	174.9
dibutyl ether	-	0.00	142.1
dichloromethane	-	> 9	39.6
diethyl ether	-	0.00	33.2
DMF	-	> 9	153.0
ethyl acetate	-	0.01	73.9
ethylene glycol	-	0.08	197.5
heptane	-	0.00	98.8
hexadecane	-	0.00	286.6
hexane	-	0.00	68.5
methyl acetate	-	0.09	44.0
methyl tert-butyl ether	-	0.00	55.2
methylcyclohexane	-	0.00	101.1
nonane	-	0.00	151.7
octane	-	0.00	126.4
undecane	-	0.00	196.3
water	-	0.00	100.0

Ugi Product 216-4A $C_{28}H_{40}N_2O_7$ - 5 solvents [97, 93]



Compound Data

Molecular weight	516.626	**H bond acceptors**	9	**Rule of 5 violations**	1
Compound type	Ugi Product	**H bond donors**	2	**ACD/ALogP**	4.08
Phase 25°C	solid	**Rotatable bonds**	13	**Predicted density**	1.12 g/cm^3
SMILES	Cc1ccc(o1)CC(C(c2ccc(c(c2)OC)OC)C(=O)NC(C)(C)C)C(=O)CNC(=O)OC(C)(C)C				
InChIKey					

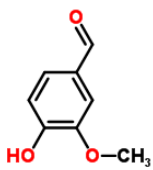
Abraham Solute Descriptors

log S_w	E	S	A	B	V
-4.630	1.716	4.742	-0.317	4.359	5.883

Solubility Data

	Concentration (M)		
Solvent	**Measured**	**Predicted**	**Solvent bp °C**
acetonitrile	0.13	0.04	63.5
DMSO	0.06	0.12	189
ethanol	0.07	0.08	72.6
methanol	0.18	0.08	48.1
toluene	0.20	0.46	110.6
1-butanol	-	0.01	117.7
1-decanol	-	0.00	227.8
1-heptanol	-	0.01	176.9
1-hexanol	-	0.01	158.2
1-octanol	-	0.00	194.7
1-pentanol	-	0.01	138.5
1-propanol	-	0.01	95.8
1,2-dichloroethane	-	> 9	83.5
1,4-dioxane	-	0.23	102.9
2-butanol	-	0.03	96.6
2-methyl-1-propanol	-	0.02	105.0
2-methyl-2-propanol	-	0.01	84.6
2-pentanol	-	0.02	118.8
2-propanol	-	0.01	73.0
2,2,4-trimethylpentane	-	0.00	98.8
3-methyl-1-butanol	-	0.03	131.2
acetone	-	0.02	46.5
benzene	-	5.40	78.8
butanone	-	0.14	75.6
butyl acetate	-	0.00	126.6
carbon disulfide	-	0.00	46.2
carbon tetrachloride	-	0.17	76.0
chloroform	-	> 9	61.2
cyclohexane	-	0.00	80.7
decane	-	0.00	174.9
dibutyl ether	-	0.00	142.1
dichloromethane	-	> 9	39.6
diethyl ether	-	0.00	33.2
DMF	-	> 9	153.0
ethyl acetate	-	0.00	73.9
ethylene glycol	-	0.01	197.5
heptane	-	0.00	98.8
hexadecane	-	0.00	286.6
hexane	-	0.00	68.5
methyl acetate	-	0.23	44.0
methyl tert-butyl ether	-	0.00	55.2
methylcyclohexane	-	0.00	101.1
nonane	-	0.00	151.7
octane	-	0.00	126.4
THF	-	0.12	68.3
undecane	-	0.00	196.3
water	-	0.00	100.0

vanillin $C_8H_8O_3$ - 8 solvents [906, 4, 55, 11, 10, 207, 19]



Compound Data

Molecular weight	152.147	**H bond acceptors**	3	**Rule of 5 violations**	0
Compound type	aldehyde	**H bond donors**	1	**ACD/ALogP**	1.19
Phase 25°C	solid	**Rotatable bonds**	3	**Predicted density**	1.231 g/cm^3
SMILES	Oc1ccc(cc1OC)C=O				

InChIKey MWOOGOJBHIARFG-UHFFFAOYSA-N

Abraham Solute Descriptors

log S_w	E	S	A	B	V
-1.480	1.024	1.086	0.365	0.411	0.935

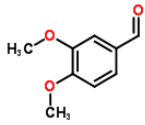
Solubility Data

	Concentration (M)		
Solvent	**Measured**	**Predicted**	**Solvent bp °C**
1-propanol	1.82	1.73	95.8
1,2-dichloroethane	1.17	0.95	83.5
acetonitrile	2.36	1.36	63.5
butanone	2.14	3.49	75.6
ethanol	2.47	2.22	72.6
methanol	4.16	2.22	48.1
THF	3.59	5.31	68.3
toluene	0.30	0.39	110.6
1-butanol	-	1.32	117.7
1-decanol	-	1.03	227.8
1-heptanol	-	1.16	176.9
1-hexanol	-	1.34	158.2
1-octanol	-	1.11	194.7
1-pentanol	-	1.35	138.5
1,4-dioxane	-	3.97	102.9
2-butanol	-	1.64	96.6
2-methyl-1-propanol	-	1.35	105.0
2-methyl-2-propanol	-	1.46	84.6
2-pentanol	-	1.31	118.8
2-propanol	-	1.60	73.0
2,2,4-trimethylpentane	-	0.02	98.8
3-methyl-1-butanol	-	1.20	131.2
acetone	-	3.39	46.5
benzene	-	0.50	78.8
butyl acetate	-	1.70	126.6
carbon disulfide	-	0.14	46.2
carbon tetrachloride	-	0.13	76.0
chloroform	-	0.81	61.2
cyclohexane	-	0.04	80.7
decane	-	0.02	174.9
dibutyl ether	-	0.32	142.1
dichloromethane	-	0.97	39.6
diethyl ether	-	1.62	33.2
DMF	-	7.11	153.0
DMSO	-	> 9	189.0
ethyl acetate	-	2.24	73.9
ethylene glycol	-	1.09	197.5
heptane	-	0.02	98.8
hexadecane	-	0.02	286.6
hexane	-	0.02	68.5
methyl acetate	-	2.57	44.0
methyl tert-butyl ether	-	0.94	55.2
methylcyclohexane	-	0.04	101.1
nonane	-	0.03	151.7
octane	-	0.03	126.4
undecane	-	0.02	196.3
water	-	0.03	100.0

veratraldehyde $C_9H_{10}O_3$ - 11 solvents [41, 71, 208]

Compound Data

Molecular weight	166.174	**H bond acceptors**	3	**Rule of 5 violations**	0
Compound type	aldehyde	**H bond donors**	0	**ACD/ALogP**	1.61



Phase 25°C	solid	**Rotatable bonds**	3	**Predicted density**	1.114 g/cm^3
SMILES	COc1cc(ccc1OC)C=O				
InChIKey	WJUFSDZVCOTFON-UHFFFAOYSA-N				

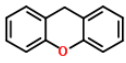
Abraham Solute Descriptors

log S_w	E	S	A	B	V
-2.080	0.889	1.091	0.204	0.572	1.310

Solubility Data

	Concentration (M)		
Solvent	**Measured**	**Predicted**	**Solvent bp °C**
acetone	3.03	5.03	46.5
acetonitrile	5.76	2.16	63.5
chloroform	5.70	7.59	61.2
dichloromethane	6.06	6.76	39.6
diethyl ether	3.55	2.82	33.2
DMSO	5.31	8.78	189
ethanol	5.49	3.51	72.6
hexane	0.07	0.14	68.5
methanol	5.87	2.90	48.1
THF	5.33	9.71	68.3
toluene	4.54	2.14	110.6
1-butanol	-	2.27	117.7
1-decanol	-	1.80	227.8
1-heptanol	-	2.37	176.9
1-hexanol	-	2.44	158.2
1-octanol	-	1.91	194.7
1-pentanol	-	2.36	138.5
1-propanol	-	2.51	95.8
1,2-dichloroethane	-	4.69	83.5
1,4-dioxane	-	6.79	102.9
2-butanol	-	2.48	96.6
2-methyl-1-propanol	-	2.31	105.0
2-methyl-2-propanol	-	2.43	84.6
2-pentanol	-	2.47	118.8
2-propanol	-	2.28	73.0
2,2,4-trimethylpentane	-	0.12	98.8
3-methyl-1-butanol	-	2.41	131.2
benzene	-	2.89	78.8
butanone	-	6.44	75.6
butyl acetate	-	3.36	126.6
carbon disulfide	-	0.86	46.2
carbon tetrachloride	-	1.04	76.0
cyclohexane	-	0.23	80.7
decane	-	0.13	174.9
dibutyl ether	-	0.82	142.1
DMF	-	> 9	153.0
ethyl acetate	-	3.80	73.9
ethylene glycol	-	0.69	197.5
heptane	-	0.15	98.8
hexadecane	-	0.13	286.6
methyl acetate	-	5.14	44.0
methyl tert-butyl ether	-	2.26	55.2
methylcyclohexane	-	0.29	101.1
nonane	-	0.15	151.7
octane	-	0.17	126.4
undecane	-	0.13	196.3
water	-	0.01	100.0

xanthene $C_{13}H_{10}O$ - 37 solvents [905]



Compound Data

Molecular weight	182.218	**H bond acceptors**	1	**Rule of 5 violations**	0
Compound type	non-Ugi related	**H bond donors**	0	**ACD/ALogP**	3.93
Phase 25°C	solid	**Rotatable bonds**	0	**Predicted density**	1.159 g/cm^3
SMILES	O2c1ccccc1Cc3c2cccc3				
InChIKey	GJCOSYZMQJWQCA-UHFFFAOYSA-N				

Abraham Solute Descriptors

log S_w	E	S	A	B	V
-4.270	1.517	1.171	-0.085	0.454	1.428

Solubility Data

	Concentration (M)		
Solvent	**Measured**	**Predicted**	**Solvent bp °C**
1-butanol	0.19	0.18	117.7
1-decanol	0.24	0.24	227.8
1-heptanol	0.23	0.24	176.9
1-hexanol	0.22	0.22	158.2
1-octanol	0.24	0.21	194.7
1-pentanol	0.20	0.20	138.5
1-propanol	0.15	0.16	95.8
1,2-dichloroethane	1.62	2.87	83.5
2-butanol	0.13	0.16	96.6
2-methyl-1-pentanol	0.16	*	148
2-methyl-1-propanol	0.12	0.14	105
2-methyl-2-butanol	0.18	*	102
2-methyl-2-propanol	0.12	0.12	84.6
2-pentanol	0.16	0.18	118.8
2-propanol	0.11	0.11	73
2,2,4-trimethylpentane	0.15	0.18	98.8
3-methyl-1-butanol	0.15	0.17	131.2
4-methyl-2-pentanol	0.14	*	133.5
acetonitrile	0.35	0.38	63.5
carbon tetrachloride	1.25	1.49	76
cyclohexane	0.39	0.53	80.7
cyclooctane	0.38	*	152
cyclopentanol	0.33	*	140.8
decane	0.24	0.19	174.9
dibutyl ether	0.50	0.30	142.1
diethyl ether	0.79	0.47	33.2
diisopropyl ether	0.41	*	68.3
ethanol	0.10	0.23	72.6
heptane	0.24	0.19	98.8
hexadecane	0.24	0.23	286.6
hexane	0.23	0.20	68.5
methanol	0.10	0.16	48.1
methyl tert-butyl ether	0.65	0.48	55.2
methylcyclohexane	0.33	0.45	101.1
nonane	0.24	0.23	151.7
octane	0.25	0.28	126.4
undecane	0.23	0.21	196.3

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