Full wwPDB X-ray Structure Validation Report (i)

May 26, 2020 – 06:16 pm BST

PDB ID	:	1MY3
Title	:	crystal structure of glutamate receptor ligand-binding core in complex with bromo-willardiine in the Zn crystal form
Authors	:	Jin, R.; Gouaux, E.
Deposited on	:	2002-10-03
Resolution	:	1.75 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

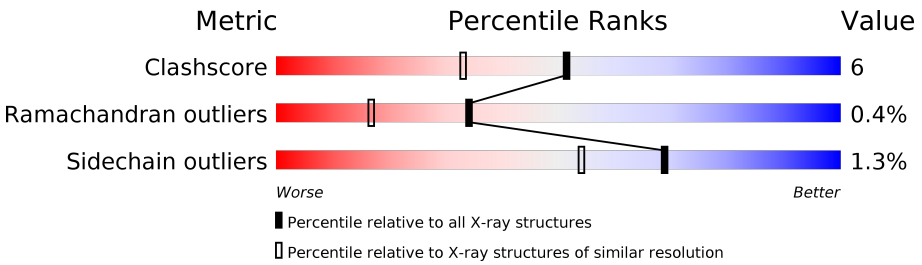
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.75 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	141614	2466 (1.76-1.76)
Ramachandran outliers	138981	2437 (1.76-1.76)
Sidechain outliers	138945	2437 (1.76-1.76)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=5%

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	263	86% 11% • •
1	B	263	87% 10% • •
1	C	263	83% 14% • •

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 6576 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called GLUTAMATE RECEPTOR 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	258	1959	1250	325	370	14	0	0	0
1	B	259	1966	1255	325	372	14	0	0	0
1	C	258	1956	1249	322	372	13	0	0	0

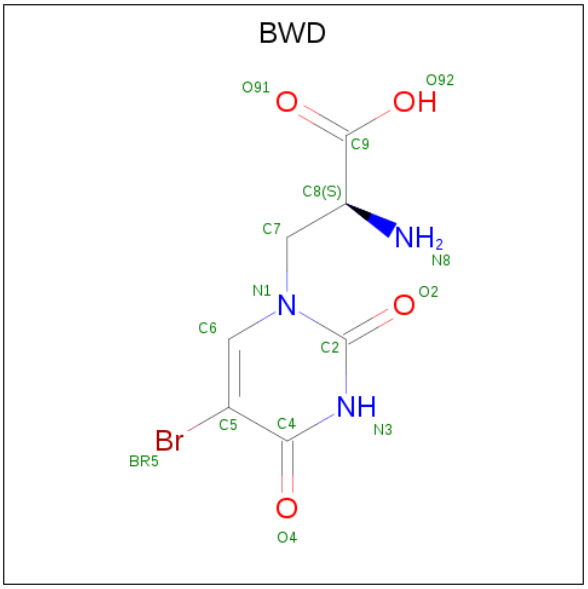
There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	GLY	-	CLONING ARTIFACT	UNP P19491
A	2	ALA	-	CLONING ARTIFACT	UNP P19491
A	118	GLY	-	LINKER	UNP P19491
A	119	THR	-	LINKER	UNP P19491
B	1	GLY	-	CLONING ARTIFACT	UNP P19491
B	2	ALA	-	CLONING ARTIFACT	UNP P19491
B	118	GLY	-	LINKER	UNP P19491
B	119	THR	-	LINKER	UNP P19491
C	1	GLY	-	CLONING ARTIFACT	UNP P19491
C	2	ALA	-	CLONING ARTIFACT	UNP P19491
C	118	GLY	-	LINKER	UNP P19491
C	119	THR	-	LINKER	UNP P19491

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Zn		
2	B	2	2	2	0	0
2	A	1	1	1	0	0
2	C	2	2	2	0	0

- Molecule 3 is 2-AMINO-3-(5-BROMO-2,4-DIOXO-3,4-DIHYDRO-2H-PYRIMIDIN-1-YL)-PROPIONIC ACID (three-letter code: BWD) (formula: $C_7H_8BrN_3O_4$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	Br	C	N	O		
3	A	1	15	1	7	3	4	0	0
			Total	Br	C	N	O		
3	B	1	15	1	7	3	4	0	0
			Total	Br	C	N	O		
3	C	1	15	1	7	3	4	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	A	241	241	241	0	0
			Total	O		
4	B	254	254	254	0	0
			Total	O		
4	C	150	150	150	0	0

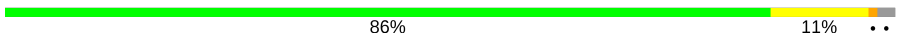
3 Residue-property plots (i)

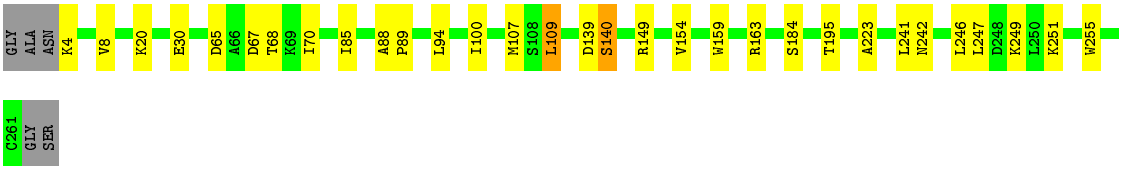
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

Note EDS was not executed.

- Molecule 1: GLUTAMATE RECEPTOR 2

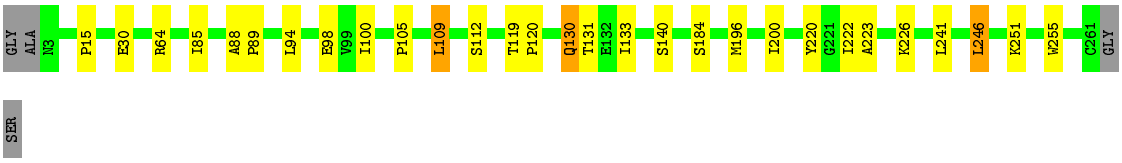
Chain A:





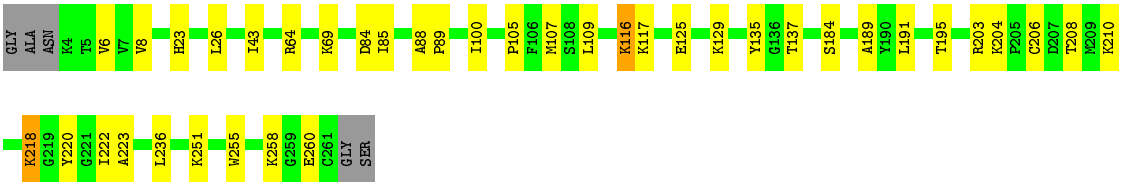
- Molecule 1: GLUTAMATE RECEPTOR 2

Chain B:



- Molecule 1: GLUTAMATE RECEPTOR 2

Chain C:



4 Data and refinement statistics (i)

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α, β, γ	114.06Å 163.41Å 47.85Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.29 – 1.75	Depositor
% Data completeness (in resolution range)	94.8 (29.29-1.75)	Depositor
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR	Depositor
R, R_{free}	0.211 , 0.255	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	6576	wwPDB-VP
Average B, all atoms (Å^2)	22.0	wwPDB-VP

5 Model quality (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, BWD

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\#\|Z\| >5$	RMSZ	$\#\|Z\| >5$
1	A	0.57	1/1995 (0.1%)	0.70	3/2688 (0.1%)
1	B	0.53	0/2002	0.70	2/2703 (0.1%)
1	C	0.50	0/1992	0.68	0/2685
All	All	0.53	1/5989 (0.0%)	0.69	5/8076 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	149	ARG	NE-CZ	13.34	1.50	1.33

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	149	ARG	CD-NE-CZ	-7.77	112.72	123.60
1	B	109	LEU	CA-CB-CG	5.84	128.74	115.30
1	A	109	LEU	CA-CB-CG	5.38	127.67	115.30
1	A	140	SER	N-CA-C	5.14	124.88	111.00
1	B	140	SER	N-CA-C	5.07	124.70	111.00

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1959	0	1941	20	0
1	B	1966	0	1939	19	0
1	C	1956	0	1939	27	0
2	A	1	0	0	0	0
2	B	2	0	0	0	0
2	C	2	0	0	0	0
3	A	15	0	7	0	0
3	B	15	0	7	0	0
3	C	15	0	7	0	0
4	A	241	0	0	5	0
4	B	254	0	0	4	0
4	C	150	0	0	5	0
All	All	6576	0	5840	65	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 6.

All (65) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:130:GLN:HE22	1:B:133:ILE:H	1.20	0.90
1:A:68:THR:HG23	1:A:70:ILE:H	1.48	0.78
1:A:65:ASP:HB3	1:A:68:THR:HG22	1.65	0.77
1:B:226:LYS:HE2	4:B:954:HOH:O	1.88	0.73
1:C:64:ARG:HH21	1:C:69:LYS:HD3	1.55	0.71
1:A:68:THR:HG21	4:A:865:HOH:O	1.95	0.66
1:C:100:ILE:HD12	1:C:223:ALA:HB1	1.82	0.61
1:C:218:LYS:HE2	4:C:830:HOH:O	2.01	0.60
1:A:65:ASP:HB3	1:A:68:THR:CG2	2.31	0.59
1:C:204:LYS:HA	1:C:206:CYS:N	2.17	0.59
1:A:154:VAL:HG23	4:A:973:HOH:O	2.03	0.58
1:C:125:GLU:O	1:C:129:LYS:HG2	2.05	0.57
1:C:8:VAL:HG22	1:C:85:ILE:CG2	2.33	0.57
1:A:241:LEU:HD22	1:A:246:LEU:HD22	1.86	0.57
1:C:204:LYS:HA	1:C:206:CYS:H	1.70	0.57
1:B:184:SER:HB2	4:B:939:HOH:O	2.06	0.56
1:B:130:GLN:NE2	1:B:133:ILE:H	1.97	0.55
1:C:116:LYS:O	1:C:208:THR:HA	2.07	0.55
1:C:107:MET:SD	1:C:195:THR:HG21	2.47	0.55
1:C:210:LYS:HE2	4:C:951:HOH:O	2.07	0.55
1:A:4:LYS:N	4:A:839:HOH:O	2.40	0.54
1:B:130:GLN:HE22	1:B:133:ILE:N	1.98	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:43:ILE:HD11	1:C:222:ILE:HD11	1.90	0.54
1:A:88:ALA:HB1	1:A:89:PRO:HD2	1.90	0.53
1:A:159:TRP:O	1:A:163:ARG:HG2	2.11	0.51
1:C:43:ILE:HD13	1:C:222:ILE:HD13	1.93	0.50
1:A:241:LEU:HD22	1:A:246:LEU:CD2	2.41	0.50
1:C:23:HIS:HA	1:C:26:LEU:HD12	1.95	0.49
1:C:218:LYS:HE3	4:C:811:HOH:O	2.11	0.49
1:A:251:LYS:O	1:A:255:TRP:HB2	2.11	0.49
1:B:112:SER:HB2	4:B:881:HOH:O	2.12	0.49
1:C:203:ARG:HH22	1:C:258:LYS:HB3	1.77	0.48
1:C:88:ALA:HB1	1:C:89:PRO:HD2	1.95	0.48
1:C:204:LYS:HG2	1:C:260:GLU:OE2	2.14	0.48
1:C:137:THR:HG22	1:C:191:LEU:HB2	1.95	0.48
1:A:8:VAL:HG22	1:A:85:ILE:CG2	2.45	0.47
1:B:88:ALA:HB1	1:B:89:PRO:HD2	1.95	0.47
1:B:131:THR:HG21	4:C:941:HOH:O	2.15	0.47
1:B:131:THR:HG22	1:B:131:THR:O	2.17	0.45
1:B:94:LEU:O	1:B:98:GLU:HG3	2.17	0.45
1:C:184:SER:HB2	4:C:891:HOH:O	2.17	0.45
1:A:139:ASP:O	1:A:140:SER:HB2	2.17	0.44
1:A:20:LYS:HD2	1:A:30:GLU:O	2.17	0.44
1:B:241:LEU:HD22	1:B:246:LEU:HD22	2.00	0.44
1:A:94:LEU:HD12	1:C:236:LEU:HD22	2.00	0.43
1:C:251:LYS:O	1:C:255:TRP:HB2	2.19	0.43
1:A:100:ILE:HD12	1:A:223:ALA:HB1	2.00	0.43
1:A:184:SER:HB3	4:A:868:HOH:O	2.18	0.43
1:A:249:LYS:HD2	4:A:1016:HOH:O	2.18	0.43
1:B:119:THR:HA	1:B:120:PRO:HD3	1.93	0.43
1:C:64:ARG:HH21	1:C:69:LYS:CD	2.27	0.43
1:B:100:ILE:HD12	1:B:223:ALA:HB1	2.01	0.42
1:B:64:ARG:HB3	1:B:64:ARG:HE	1.66	0.42
1:C:6:VAL:HG13	1:C:84:ASP:HB2	2.02	0.42
1:C:135:TYR:HA	1:C:189:ALA:O	2.20	0.42
1:B:105:PRO:HA	1:B:220:TYR:O	2.20	0.42
1:B:196:MET:O	1:B:200:ILE:HG12	2.19	0.42
1:C:203:ARG:NH2	1:C:258:LYS:HB3	2.35	0.41
1:A:242:ASN:HB2	1:A:247:LEU:HD12	2.02	0.41
1:C:105:PRO:HA	1:C:220:TYR:O	2.21	0.41
1:B:251:LYS:HE2	4:B:1000:HOH:O	2.19	0.41
1:A:107:MET:SD	1:A:195:THR:HG21	2.61	0.41
1:B:251:LYS:O	1:B:255:TRP:HB2	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:85:ILE:HD11	1:B:222:ILE:HG21	2.03	0.41
1:C:43:ILE:HD11	1:C:222:ILE:CD1	2.50	0.40

There are no symmetry-related clashes.

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	256/263 (97%)	249 (97%)	6 (2%)	1 (0%)	34 17
1	B	257/263 (98%)	253 (98%)	4 (2%)	0	100 100
1	C	256/263 (97%)	248 (97%)	6 (2%)	2 (1%)	19 6
All	All	769/789 (98%)	750 (98%)	16 (2%)	3 (0%)	34 17

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	67	ASP
1	C	116	LYS
1	C	117	LYS

5.3.2 Protein sidechains (i)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	203/219 (93%)	202 (100%)	1 (0%)	88	83
1	B	202/219 (92%)	197 (98%)	5 (2%)	47	25
1	C	203/219 (93%)	201 (99%)	2 (1%)	76	63
All	All	608/657 (92%)	600 (99%)	8 (1%)	69	54

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	109	LEU
1	B	15	PRO
1	B	30	GLU
1	B	109	LEU
1	B	130	GLN
1	B	246	LEU
1	C	109	LEU
1	C	218	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	130	GLN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains (i)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates (i)

There are no carbohydrates in this entry.

5.6 Ligand geometry (i)

Of 8 ligands modelled in this entry, 5 are monoatomic - leaving 3 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
3	BWD	B	801	-	10,15,15	4.06	3 (30%)	11,21,21	3.00	4 (36%)
3	BWD	A	800	-	10,15,15	3.94	2 (20%)	11,21,21	3.03	4 (36%)
3	BWD	C	802	-	10,15,15	4.09	2 (20%)	11,21,21	2.98	4 (36%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	BWD	B	801	-	-	0/4/8/8	0/1/1/1
3	BWD	A	800	-	-	0/4/8/8	0/1/1/1
3	BWD	C	802	-	-	0/4/8/8	0/1/1/1

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	802	BWD	C6-N1	-11.97	1.32	1.46
3	B	801	BWD	C6-N1	-11.63	1.33	1.46
3	A	800	BWD	C6-N1	-11.51	1.33	1.46
3	B	801	BWD	BR5-C5	-3.57	1.87	1.97
3	C	802	BWD	BR5-C5	-3.28	1.88	1.97
3	A	800	BWD	BR5-C5	-3.03	1.89	1.97
3	B	801	BWD	C7-C8	-2.31	1.50	1.53

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	A	800	BWD	C5-C6-N1	8.36	123.51	109.53
3	C	802	BWD	C5-C6-N1	8.15	123.16	109.53
3	B	801	BWD	C5-C6-N1	8.12	123.12	109.53
3	B	801	BWD	BR5-C5-C6	4.18	121.64	110.56
3	C	802	BWD	BR5-C5-C6	4.18	121.63	110.56
3	A	800	BWD	BR5-C5-C6	3.91	120.93	110.56

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	B	801	BWD	N3-C2-N1	-2.28	114.13	116.59
3	A	800	BWD	N3-C2-N1	-2.26	114.15	116.59
3	C	802	BWD	N3-C2-N1	-2.23	114.18	116.59
3	B	801	BWD	C5-C4-N3	-2.20	113.22	116.20
3	A	800	BWD	C5-C4-N3	-2.18	113.25	116.20
3	C	802	BWD	C5-C4-N3	-2.03	113.45	116.20

There are no chirality outliers.

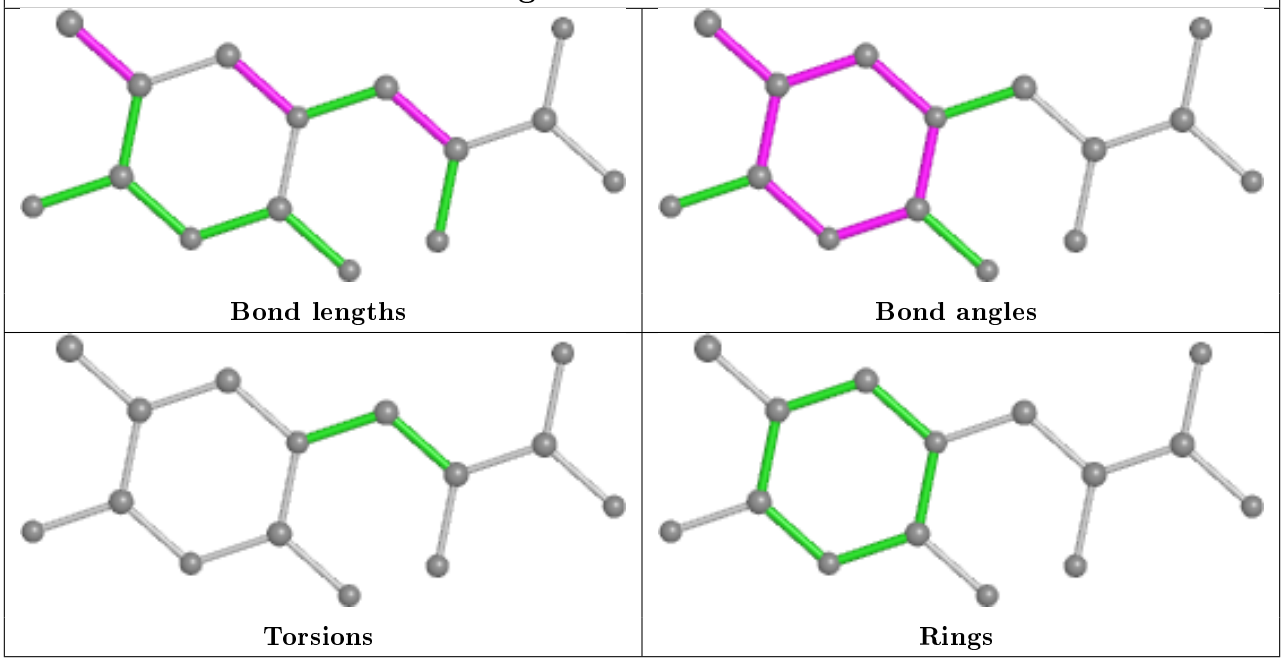
There are no torsion outliers.

There are no ring outliers.

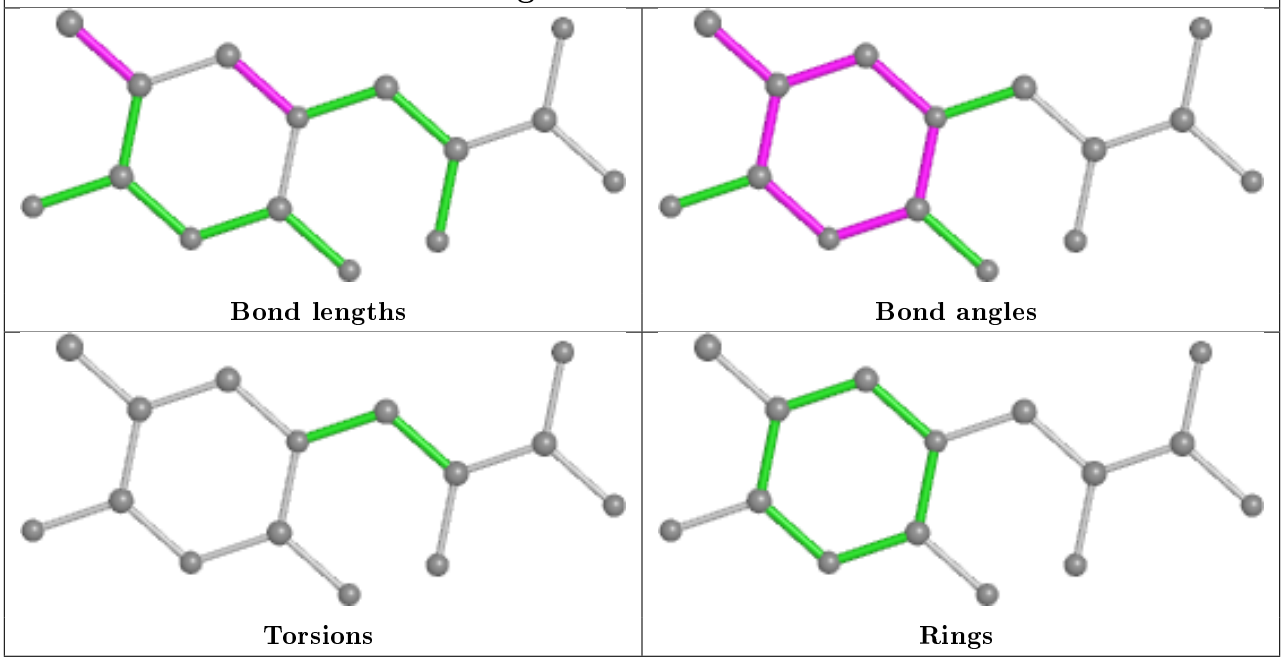
No monomer is involved in short contacts.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less then 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

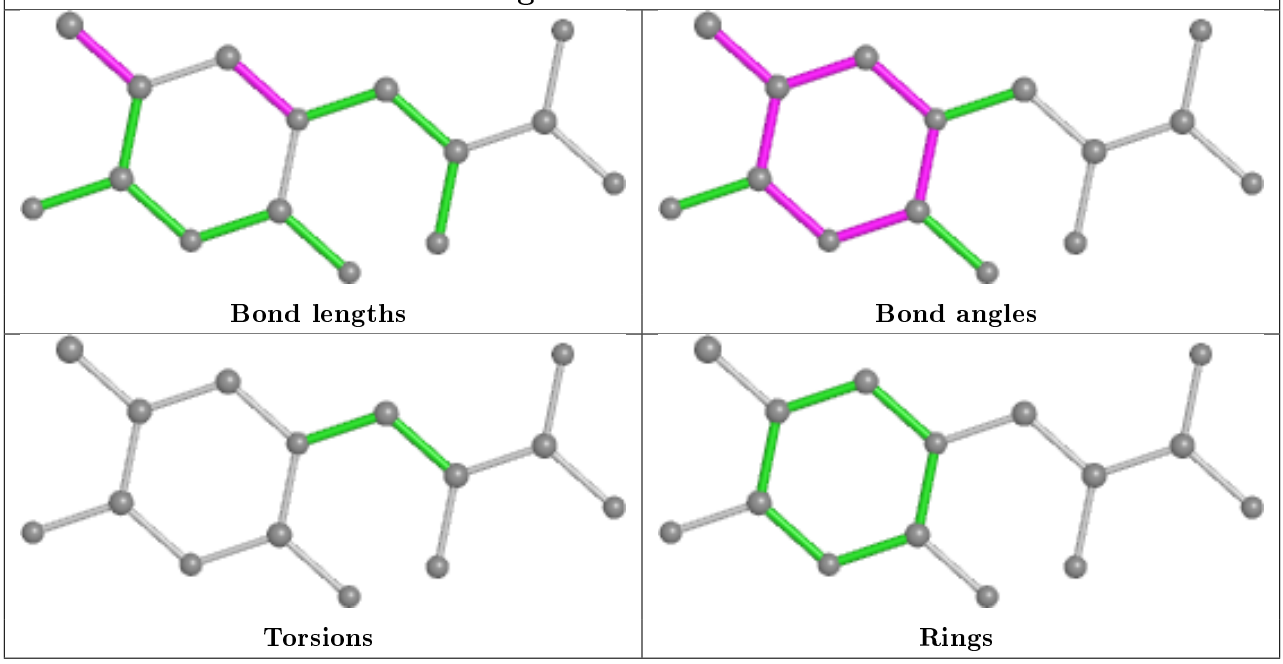
Ligand BWD B 801



Ligand BWD A 800



Ligand BWD C 802



5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data (i)

6.1 Protein, DNA and RNA chains (i)

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains (i)

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates (i)

EDS was not executed - this section is therefore empty.

6.4 Ligands (i)

EDS was not executed - this section is therefore empty.

6.5 Other polymers (i)

EDS was not executed - this section is therefore empty.