Full wwPDB X-ray Structure Validation Report ⓘ

Aug 22, 2020 – 04:55 AM BST

PDB ID	:	5FL1
Title	:	Structure of a hydrolase with an inhibitor
Authors	:	Cekic, N.; Heinonen, J.E.; Stubbs, K.A.; Roth, C.; McEachern, E.J.; Davies, G.J.; Vocadlo, D.J.
Deposited on	:	2015-10-20
Resolution	:	1.95 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 ⓘ symbol.

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

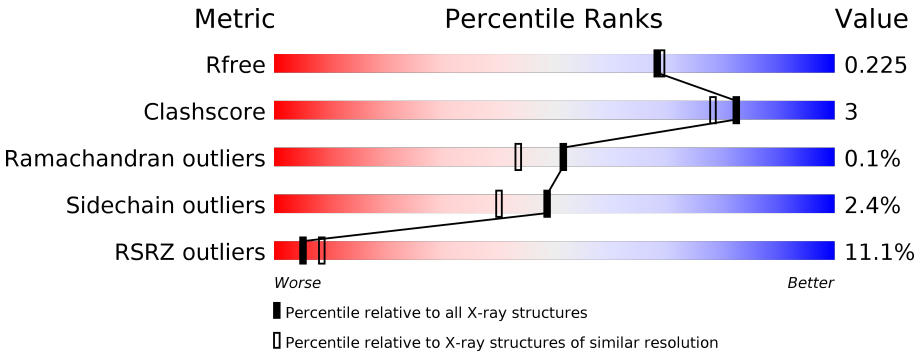
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.13.1
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.13.1

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 1.95 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))
R_{free}	130704	2580 (1.96-1.96)
Clashscore	141614	2705 (1.96-1.96)
Ramachandran outliers	138981	2678 (1.96-1.96)
Sidechain outliers	138945	2678 (1.96-1.96)
RSRZ outliers	127900	2539 (1.96-1.96)

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% The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	716	12% 91% 6% •
1	B	716	10% 86% 6% • 7%

2 Entry composition

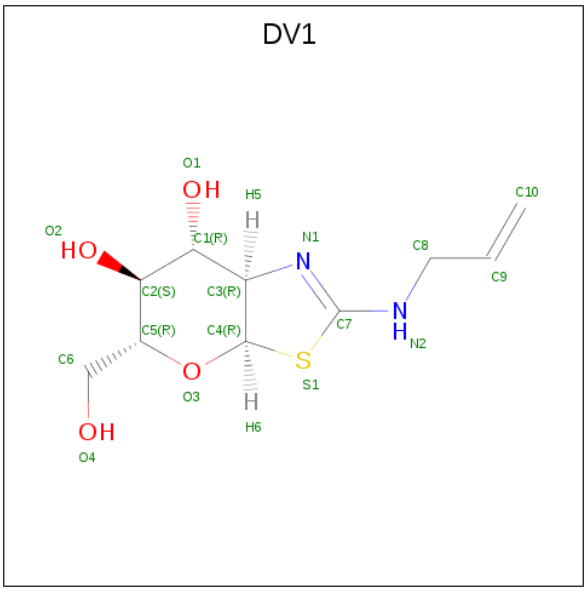
There are 5 unique types of molecules in this entry. The entry contains 22762 atoms, of which 11013 are hydrogens and 0 are deuteriums.

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 O-GLCNACASE BT_4395.

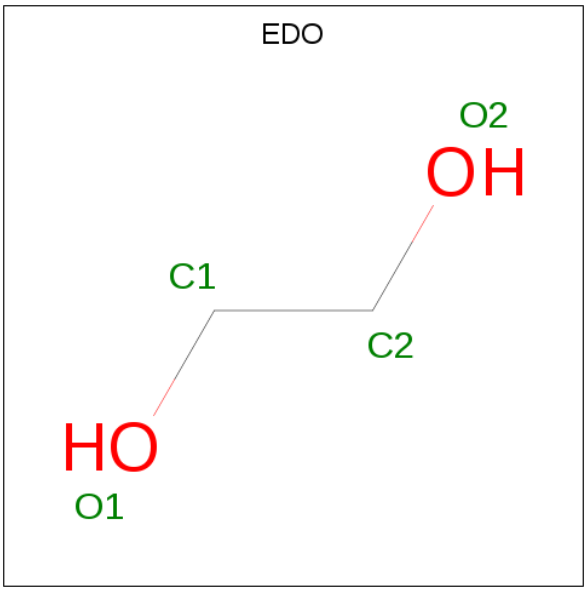
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	703	Total	C	H	N	O	S	0	3	0
			11360	3662	5636	968	1075	19			
1	B	665	Total	C	H	N	O	S	0	0	0
			10725	3464	5321	913	1009	18			

- Molecule 2 is (3 {a} {R},5 {R},6 {S},7 {R},7 {a} {R})-5-(hydroxymethyl)-2-(prop-2-enylamino)-5,6,7,7 {a}-tetrahydro-3 {a} {H}-pyrano[3,2-d][1,3]thiazole-6,7-diol (three-letter code: DV1) (formula: $C_{10}H_{16}N_2O_4S$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total	C	H	N	O	S	0	0
			33	10	16	2	4	1		
2	B	1	Total	C	H	N	O	S	0	0
			33	10	16	2	4	1		

- Molecule 3 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $C_2H_6O_2$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C H O 10 2 6 2	0	0
3	A	1	Total C H O 10 2 6 2	0	0
3	A	1	Total C H O 10 2 6 2	0	0
3	B	1	Total C H O 10 2 6 2	0	0

- Molecule 4 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	1	Total Ca 1 1	0	0

- Molecule 5 is water.

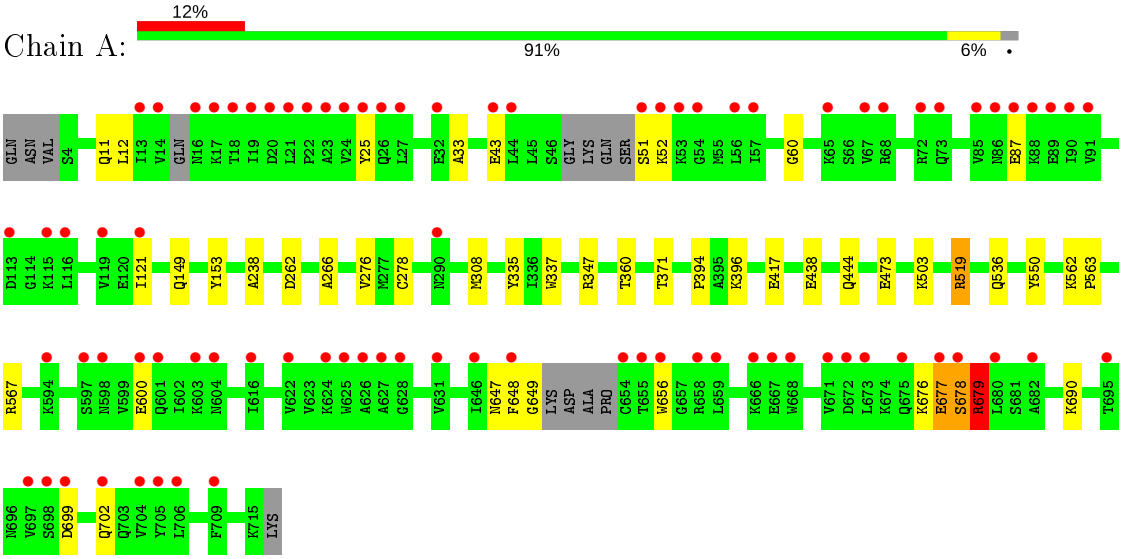
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	305	Total O 305 305	0	0
5	B	265	Total O 265 265	0	0

3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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 and electron density. 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. A red dot above a residue indicates a poor fit to the electron density (RSRZ > 2). 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.

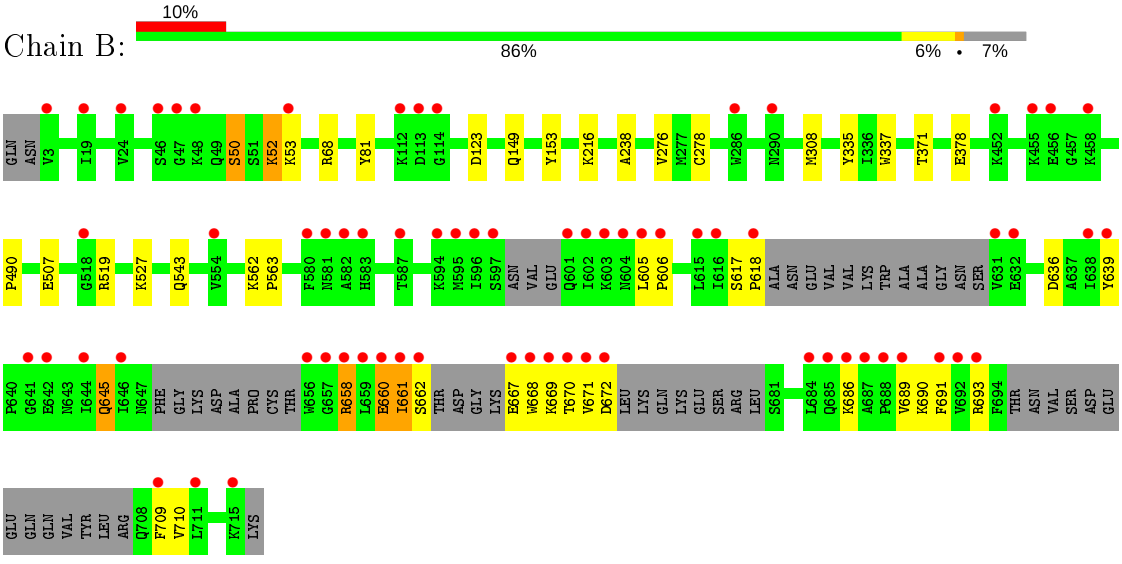
- Molecule 1: O-GLCNACASE BT_4395

Chain A:



- Molecule 1: O-GLCNACASE BT_4395

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 2 21 21	Depositor
Cell constants a, b, c, α, β, γ	51.50Å 162.11Å 223.21Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.08 – 1.95 49.08 – 1.95	Depositor EDS
% Data completeness (in resolution range)	92.8 (49.08-1.95) 93.6 (49.08-1.95)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.37 (at 1.95Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, R_{free}	0.193 , 0.222 0.197 , 0.225	Depositor DCC
R_{free} test set	6442 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å^2)	27.7	Xtriage
Anisotropy	1.208	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.39 , 43.7	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	22762	wwPDB-VP
Average B, all atoms (Å^2)	53.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 19.28% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CA, DV1, EDO

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	0.42	0/5873	0.55	0/7958
1	B	0.43	0/5537	0.56	0/7499
All	All	0.42	0/11410	0.55	0/15457

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand.A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	600	GLU	Peptide

5.2 Too-close contacts

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	5724	5636	5640	29	0
1	B	5404	5321	5321	29	0
2	A	17	16	0	0	0
2	B	17	16	0	0	0
3	A	12	18	18	2	0
3	B	4	6	6	0	0
4	B	1	0	0	0	0
5	A	305	0	0	4	0
5	B	265	0	0	1	0
All	All	11749	11013	10985	56	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 3.

All (56) 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:A:649:GLY:O	1:A:679:ARG:NH2	2.04	0.89
1:B:662:SER:OG	1:B:691:PHE:N	2.21	0.73
1:A:647:ASN:OD1	1:A:679:ARG:NH1	2.23	0.71
1:B:662:SER:CB	1:B:690:LYS:H	2.03	0.71
1:A:648:PHE:N	1:A:679:ARG:HH12	1.87	0.70
1:A:536:GLN:NE2	5:A:2253:HOH:O	2.32	0.63
1:B:543:GLN:OE1	5:B:2253:HOH:O	2.16	0.58
1:A:25:TYR:O	1:A:51:SER:N	2.36	0.58
1:B:617:SER:HB2	1:B:618:PRO:HD3	1.87	0.57
1:B:50:SER:OG	1:B:52:LYS:O	2.24	0.56
1:B:661:ILE:HG22	1:B:668:TRP:HA	1.88	0.55
1:B:662:SER:CB	1:B:691:PHE:H	2.20	0.54
1:B:662:SER:OG	1:B:690:LYS:N	2.42	0.53
1:B:658:ARG:HD3	1:B:670:THR:HG21	1.91	0.53
1:B:661:ILE:HG22	1:B:669:LYS:N	2.25	0.52
1:A:567:ARG:NH1	5:A:2268:HOH:O	2.40	0.50
1:B:660:GLU:HG3	1:B:693:ARG:HB3	1.94	0.50
1:A:676:LYS:HG3	1:A:677:GLU:N	2.28	0.49
1:A:417:GLU:OE1	1:A:417:GLU:N	2.43	0.47
1:A:503:LYS:NZ	1:B:507:GLU:OE2	2.48	0.47
1:A:308:MET:HA	1:A:335:TYR:O	2.16	0.46
1:A:519:ARG:H	1:A:519:ARG:HD2	1.80	0.46
1:A:648:PHE:H	1:A:679:ARG:HH12	1.60	0.46
1:B:662:SER:CB	1:B:690:LYS:N	2.76	0.46
1:B:639:TYR:O	1:B:689:VAL:N	2.48	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:667:GLU:C	1:B:668:TRP:HD1	2.19	0.46
1:B:660:GLU:HB2	1:B:693:ARG:H	1.81	0.46
1:A:438[A]:GLU:OE1	5:A:2191:HOH:O	2.21	0.45
1:B:661:ILE:HG22	1:B:669:LYS:H	1.81	0.44
1:A:562[A]:LYS:HB3	1:A:563:PRO:HD3	1.99	0.44
1:A:690:LYS:NZ	5:A:2290:HOH:O	2.43	0.44
1:B:308:MET:HA	1:B:335:TYR:O	2.18	0.44
1:B:661:ILE:HG13	1:B:662:SER:N	2.32	0.44
1:B:671:VAL:HG12	1:B:672:ASP:H	1.82	0.44
1:A:262:ASP:HA	1:A:266:ALA:HB3	2.00	0.44
1:A:562[B]:LYS:HB3	1:A:563:PRO:HD3	2.00	0.43
1:A:473:GLU:OE2	1:B:527:LYS:NZ	2.37	0.43
1:A:12:LEU:HD13	1:A:121:ILE:HG12	2.00	0.43
1:B:668:TRP:HH2	1:B:693:ARG:NH1	2.16	0.43
1:A:648:PHE:N	1:A:679:ARG:NH1	2.62	0.43
1:B:562:LYS:HB3	1:B:563:PRO:HD3	2.00	0.42
1:A:550:TYR:CG	3:A:1719:EDO:H22	2.55	0.42
1:B:238:ALA:HA	1:B:276:VAL:O	2.18	0.42
1:B:605:LEU:HB2	1:B:606:PRO:CD	2.49	0.42
1:B:378:GLU:HG3	1:B:490:PRO:HB2	2.02	0.42
1:A:238:ALA:HA	1:A:276:VAL:O	2.20	0.42
1:A:678:SER:O	1:A:679:ARG:HB2	2.18	0.42
1:B:81:TYR:CZ	1:B:123:ASP:HB3	2.55	0.41
1:A:656:TRP:CE2	1:A:702:GLN:HG2	2.55	0.41
1:B:149:GLN:HB3	1:B:153:TYR:CZ	2.56	0.41
1:A:347:ARG:O	3:A:1717:EDO:O2	2.29	0.41
1:A:149:GLN:HB3	1:A:153:TYR:CZ	2.56	0.41
1:A:360:THR:OG1	1:A:394:PRO:O	2.39	0.41
1:B:645:GLN:HG3	1:B:710:VAL:CG1	2.50	0.41
1:A:648:PHE:H	1:A:679:ARG:NH1	2.19	0.40
1:A:33:ALA:HB2	1:A:60:GLY:HA2	2.04	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	698/716 (98%)	673 (96%)	24 (3%)	1 (0%)	51	43
1	B	651/716 (91%)	617 (95%)	33 (5%)	1 (0%)	47	38
All	All	1349/1432 (94%)	1290 (96%)	57 (4%)	2 (0%)	51	43

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	679	ARG
1	B	519	ARG

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	622/630 (99%)	607 (98%)	15 (2%)	49	40
1	B	586/630 (93%)	571 (97%)	15 (3%)	46	36
All	All	1208/1260 (96%)	1178 (98%)	30 (2%)	49	38

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11[A]	GLN
1	A	11[B]	GLN
1	A	43	GLU
1	A	52	LYS
1	A	87	GLU
1	A	278	CYS
1	A	337	TRP
1	A	371	THR
1	A	396	LYS
1	A	444	GLN
1	A	519	ARG

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Mol	Chain	Res	Type
1	A	677	GLU
1	A	678	SER
1	A	679	ARG
1	A	699	ASP
1	B	50	SER
1	B	52	LYS
1	B	53	LYS
1	B	68	ARG
1	B	216	LYS
1	B	278	CYS
1	B	337	TRP
1	B	371	THR
1	B	636	ASP
1	B	645	GLN
1	B	658	ARG
1	B	660	GLU
1	B	661	ILE
1	B	686	LYS
1	B	709	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 7 ligands modelled in this entry, 1 is monoatomic - leaving 6 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	$\#\|Z\| > 2$
2	DV1	A	1716	-	16,18,18	1.01	1 (6%)	15,25,25	0.87	0
3	EDO	B	1717	-	3,3,3	0.37	0	2,2,2	0.21	0
3	EDO	A	1718	-	3,3,3	0.54	0	2,2,2	0.47	0
3	EDO	A	1719	-	3,3,3	0.51	0	2,2,2	0.41	0
2	DV1	B	1716	-	16,18,18	1.01	1 (6%)	15,25,25	0.98	0
3	EDO	A	1717	-	3,3,3	0.42	0	2,2,2	0.22	0

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
2	DV1	A	1716	-	-	1/6/34/34	0/2/2/2
3	EDO	B	1717	-	-	0/1/1/1	-
3	EDO	A	1718	-	-	0/1/1/1	-
3	EDO	A	1719	-	-	0/1/1/1	-
2	DV1	B	1716	-	-	1/6/34/34	0/2/2/2
3	EDO	A	1717	-	-	0/1/1/1	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1716	DV1	C7-S1	-3.31	1.67	1.76
2	A	1716	DV1	C7-S1	-3.22	1.67	1.76

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1716	DV1	N2-C8-C9-C10
2	B	1716	DV1	N2-C8-C9-C10

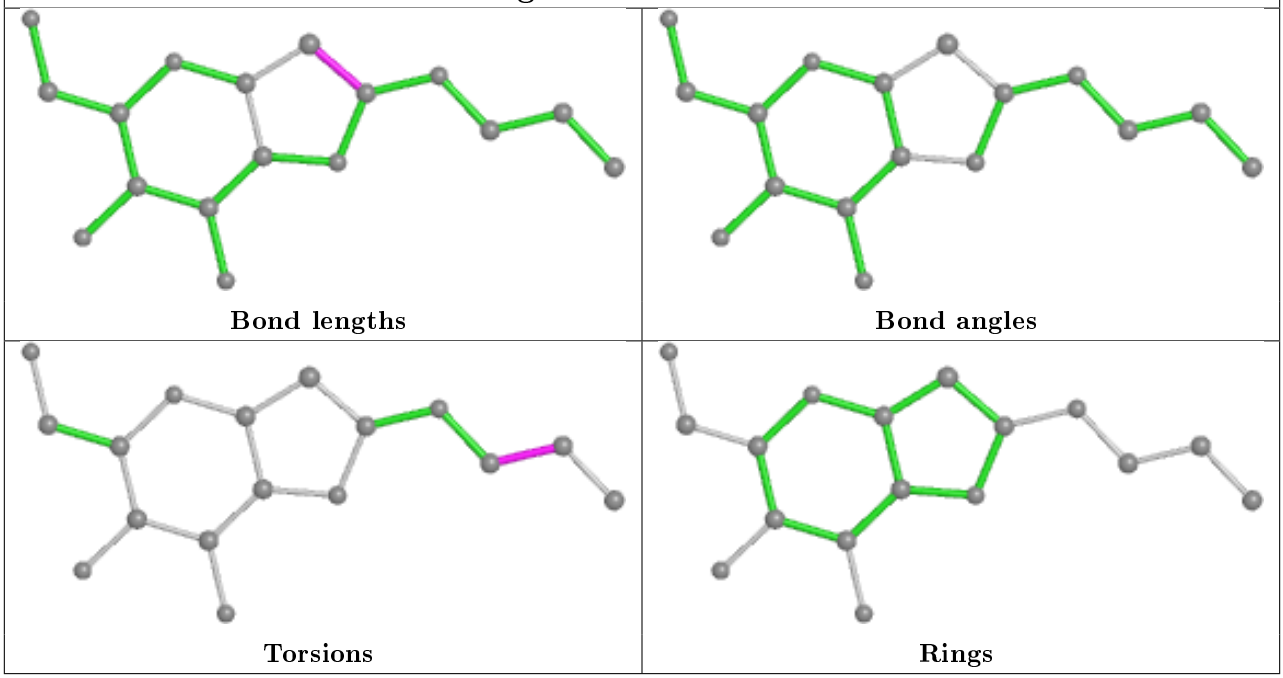
There are no ring outliers.

2 monomers are involved in 2 short contacts:

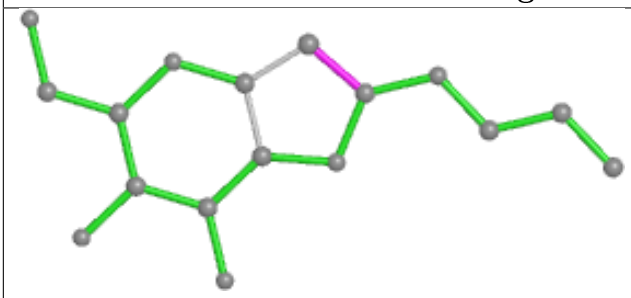
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1719	EDO	1	0
3	A	1717	EDO	1	0

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.

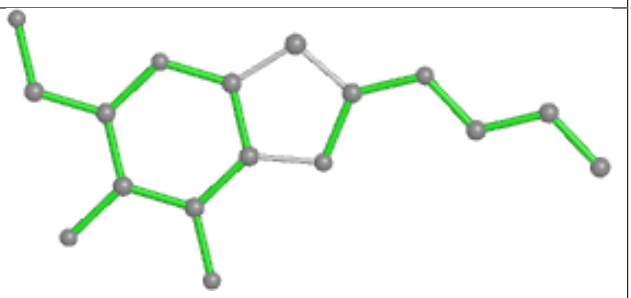
Ligand DV1 A 1716



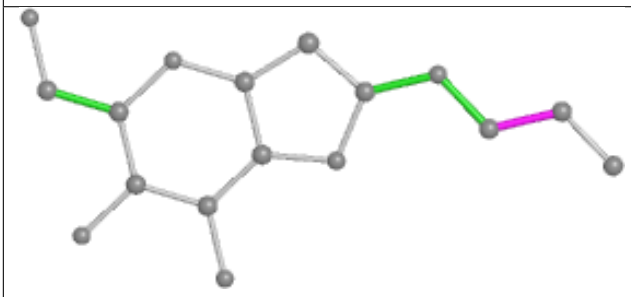
Ligand DV1 B 1716



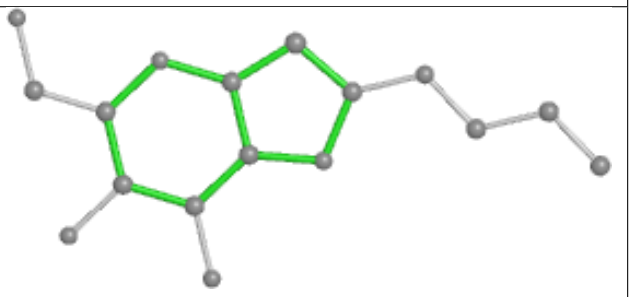
Bond lengths



Bond angles



Torsions



Rings

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled '#RSRZ> 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95^{th} percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled 'Q< 0.9' lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å^2)	Q<0.9
1	A	703/716 (98%)	0.75	83 (11%) 4 7	24, 44, 98, 137	0
1	B	665/716 (92%)	0.66	69 (10%) 6 10	21, 38, 91, 139	0
All	All	1368/1432 (95%)	0.70	152 (11%) 5 8	21, 41, 95, 139	0

All (152) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	602	ILE	10.5
1	B	656	TRP	8.1
1	A	53	LYS	7.5
1	A	601	GLN	7.4
1	A	24	VAL	7.2
1	A	88	LYS	6.9
1	A	627	ALA	6.7
1	B	661	ILE	6.4
1	B	596	ILE	6.1
1	A	13	ILE	5.6
1	A	655	THR	5.6
1	B	646	ILE	5.4
1	B	659	LEU	5.4
1	B	668	TRP	5.4
1	A	14	VAL	5.1
1	B	290	ASN	5.0
1	A	56	LEU	4.9
1	B	603	LYS	4.7
1	B	601	GLN	4.5
1	A	54	GLY	4.5
1	A	119	VAL	4.5
1	B	687	ALA	4.4
1	B	597	SER	4.4
1	B	671	VAL	4.4

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Mol	Chain	Res	Type	RSRZ
1	A	52	LYS	4.3
1	B	670	THR	4.3
1	B	48	LYS	4.3
1	A	673	LEU	4.2
1	A	675	GLN	4.2
1	B	604	ASN	4.2
1	B	685	GLN	4.2
1	A	628	GLY	4.2
1	B	518	GLY	4.2
1	B	662	SER	4.2
1	B	581	ASN	4.1
1	B	692	VAL	4.1
1	A	648	PHE	4.1
1	A	115	LYS	4.0
1	B	583	HIS	4.0
1	B	582	ALA	3.9
1	B	644	ILE	3.9
1	B	684	LEU	3.8
1	A	25	TYR	3.7
1	A	85	VAL	3.7
1	B	631	VAL	3.7
1	B	53	LYS	3.7
1	A	697	VAL	3.7
1	A	654	CYS	3.7
1	A	21	LEU	3.6
1	A	27	LEU	3.6
1	A	116	LEU	3.6
1	A	51	SER	3.6
1	B	667	GLU	3.6
1	A	22	PRO	3.6
1	B	47	GLY	3.5
1	B	688	PRO	3.5
1	A	604	ASN	3.5
1	A	87	GLU	3.5
1	B	3	VAL	3.4
1	A	18	THR	3.3
1	A	626	ALA	3.3
1	A	65	LYS	3.3
1	A	598	ASN	3.3
1	A	625	TRP	3.3
1	B	657	GLY	3.2
1	B	605	LEU	3.2

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Mol	Chain	Res	Type	RSRZ
1	A	91	VAL	3.2
1	A	646	ILE	3.2
1	A	622	VAL	3.1
1	B	638	ILE	3.1
1	B	458	LYS	3.1
1	A	666	LYS	3.1
1	A	19	ILE	3.1
1	B	455	LYS	3.1
1	A	680	LEU	3.0
1	A	23	ALA	3.0
1	B	709	PHE	3.0
1	A	671	VAL	3.0
1	A	600	GLU	2.9
1	B	618	PRO	2.9
1	A	699	ASP	2.9
1	A	603	LYS	2.9
1	A	659	LEU	2.9
1	B	689	VAL	2.8
1	A	698	SER	2.8
1	A	677	GLU	2.8
1	A	17	LYS	2.8
1	A	678	SER	2.8
1	A	16	ASN	2.8
1	A	44	LEU	2.8
1	B	711	LEU	2.8
1	A	667	GLU	2.8
1	B	456	GLU	2.8
1	B	660	GLU	2.8
1	B	693	ARG	2.7
1	A	57	ILE	2.7
1	A	90	ILE	2.7
1	B	46	SER	2.7
1	B	114	GLY	2.7
1	A	67	VAL	2.7
1	B	595	MET	2.7
1	A	68	ARG	2.7
1	A	72	ARG	2.7
1	B	669	LYS	2.7
1	A	121	ILE	2.6
1	B	19	ILE	2.6
1	A	668	TRP	2.6
1	A	624	LYS	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	20	ASP	2.5
1	A	597	SER	2.5
1	B	658	ARG	2.5
1	B	672	ASP	2.5
1	A	89	GLU	2.5
1	B	452	LYS	2.5
1	A	706	LEU	2.5
1	B	715	LYS	2.4
1	A	656	TRP	2.4
1	A	705	TYR	2.4
1	B	580	PHE	2.4
1	B	615	LEU	2.4
1	B	113	ASP	2.4
1	A	658	ARG	2.4
1	A	616	ILE	2.4
1	B	616	ILE	2.3
1	A	682	ALA	2.3
1	B	641	GLY	2.3
1	A	73	GLN	2.3
1	A	113	ASP	2.3
1	A	290	ASN	2.2
1	A	26	GLN	2.2
1	A	86	ASN	2.2
1	A	709	PHE	2.2
1	A	32	GLU	2.2
1	A	695	THR	2.2
1	B	554	VAL	2.2
1	B	686	LYS	2.2
1	A	704	VAL	2.1
1	B	24	VAL	2.1
1	B	632	GLU	2.1
1	B	286	TRP	2.1
1	B	594	LYS	2.1
1	B	606	PRO	2.1
1	A	702	GLN	2.1
1	A	631	VAL	2.1
1	A	672	ASP	2.1
1	A	594	LYS	2.1
1	A	43	GLU	2.0
1	B	691	PHE	2.0
1	B	587	THR	2.0
1	B	112	LYS	2.0

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Mol	Chain	Res	Type	RSRZ
1	B	639	TYR	2.0
1	B	642	GLU	2.0

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

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

6.3 Carbohydrates (i)

There are no monosaccharides in this entry.

6.4 Ligands (i)

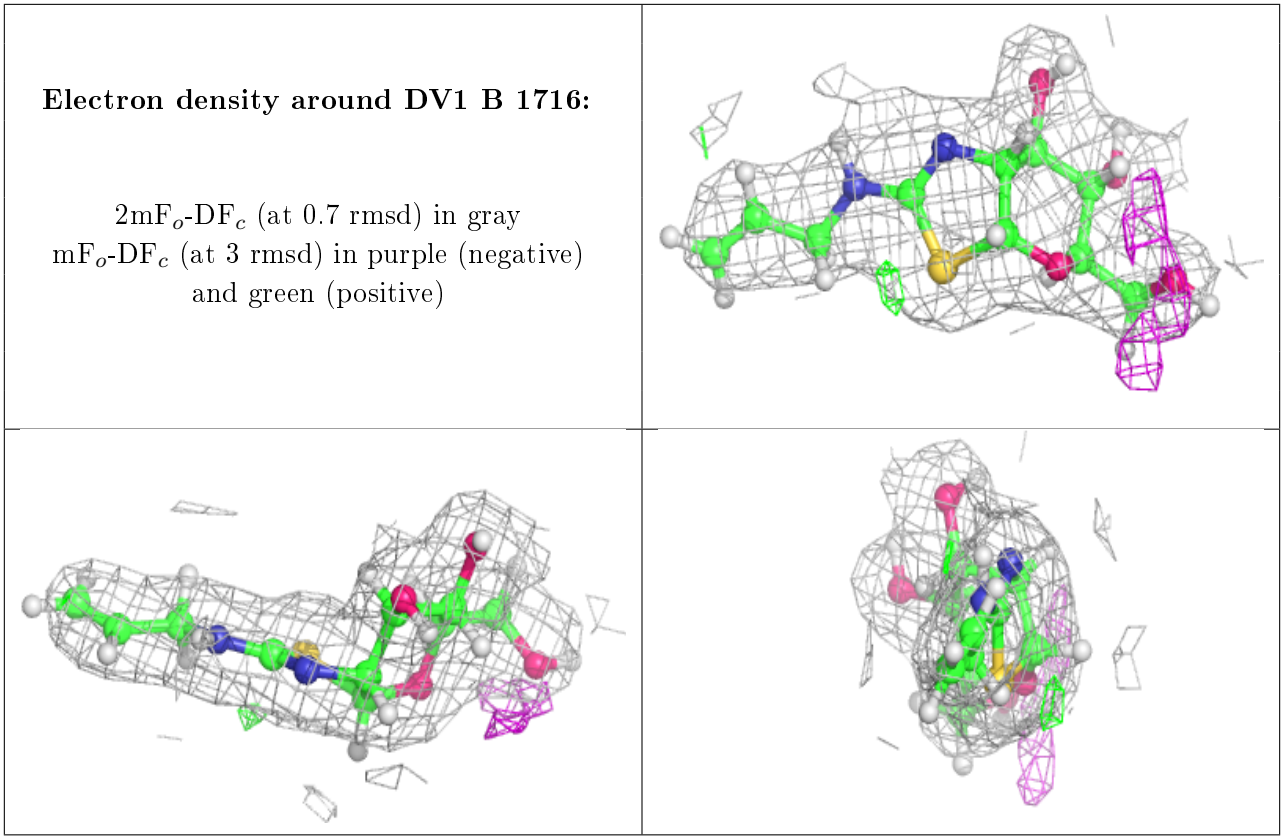
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	EDO	A	1718	4/4	0.90	0.23	42,51,53,54	0
3	EDO	A	1719	4/4	0.90	0.17	33,40,44,45	0
3	EDO	B	1717	4/4	0.95	0.20	43,52,56,58	0
3	EDO	A	1717	4/4	0.96	0.17	33,40,42,43	0
2	DV1	B	1716	17/17	0.97	0.14	25,31,44,45	0
2	DV1	A	1716	17/17	0.97	0.14	23,31,42,42	0
4	CA	B	1718	1/1	0.97	0.14	33,33,33,33	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

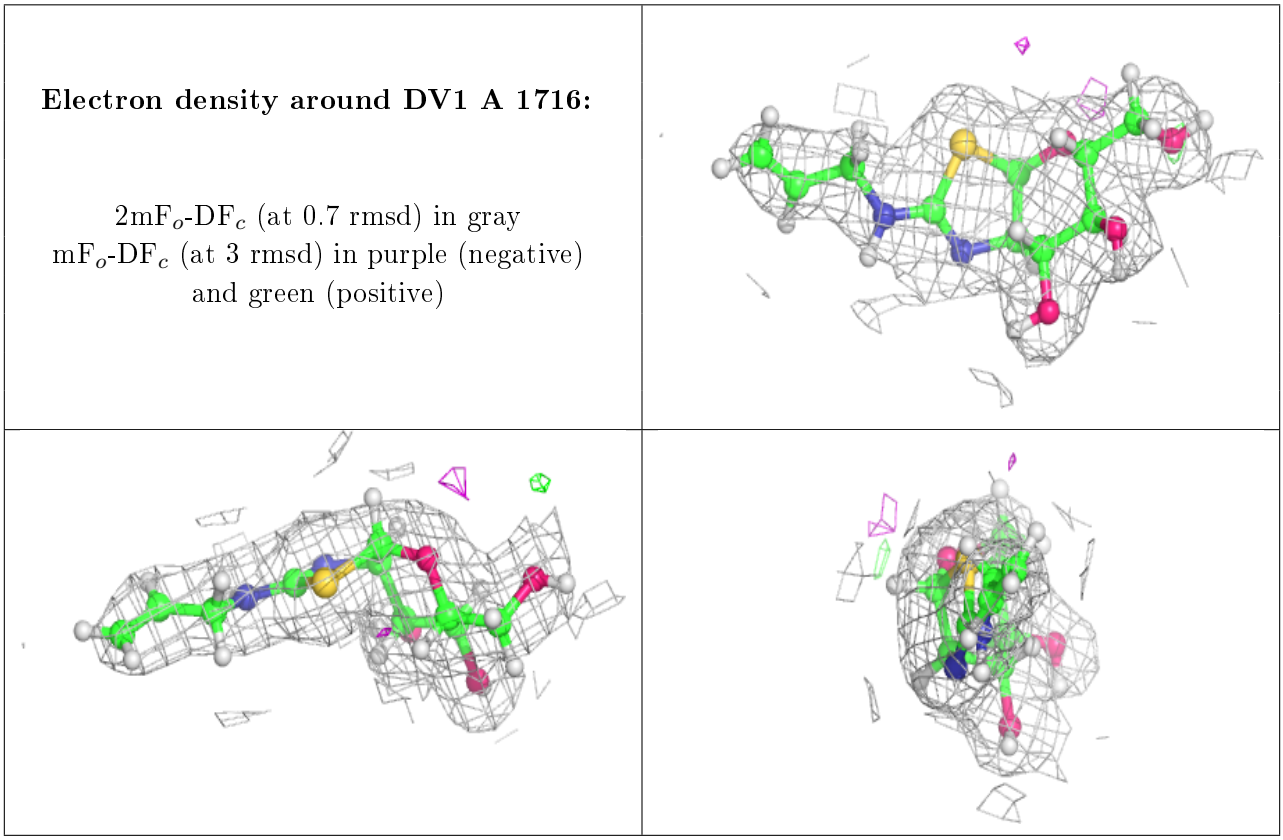
Electron density around DV1 B 1716:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)



Electron density around DV1 A 1716:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)



6.5 Other polymers (i)

There are no such residues in this entry.