



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 04:11 AM UTC

PDB ID : 6HCL / pdb_00006hcl
Title : Crystal structure of a MFS transporter with Ligand at 2.69 Angstroem resolution
Authors : Kalbermatter, D.; Bosshart, P.; Bonetti, S.; Fotiadis, D.
Deposited on : 2018-08-15
Resolution : 2.50 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

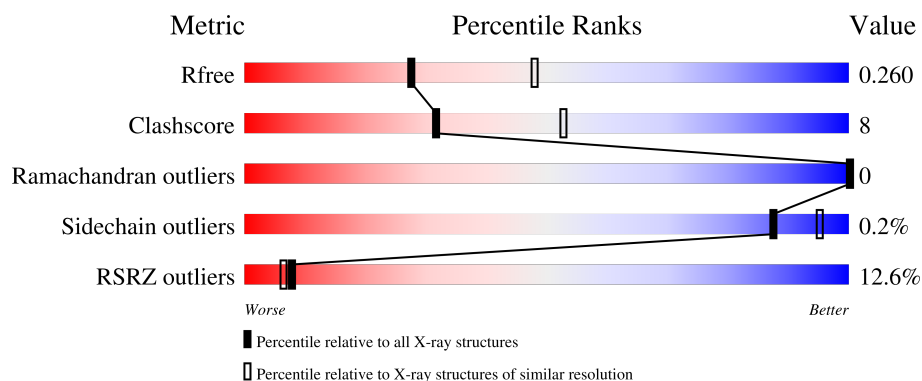
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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}	180053	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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 of 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 $\leq 5\%$. 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	420	
1	B	420	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	2OP	B	502	-	-	X	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5771 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.

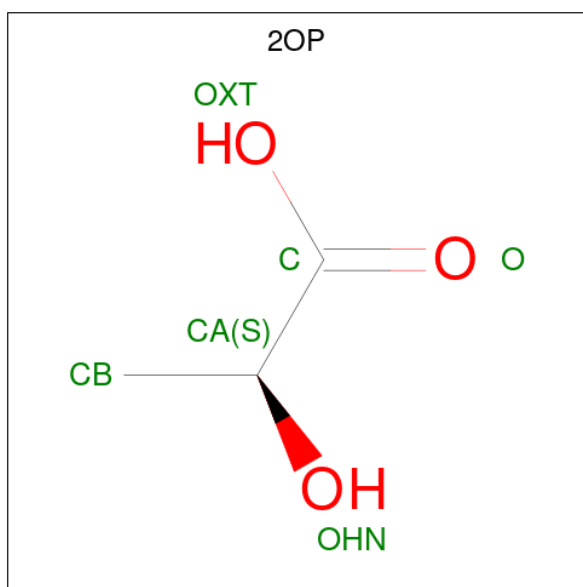
- Molecule 1 is a protein called Major facilitator superfamily MFS_1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	395	Total	C	N	O	S	0	0	0
			2980	1992	470	494	24			
1	B	366	Total	C	N	O	S	0	0	0
			2752	1844	430	455	23			

There are 16 discrepancies between the modelled and reference sequences:

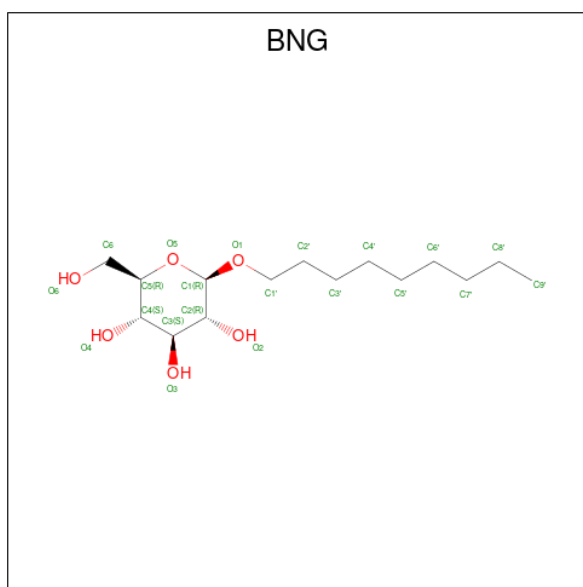
Chain	Residue	Modelled	Actual	Comment	Reference
A	413	LEU	-	expression tag	UNP A0LNN5
A	414	GLU	-	expression tag	UNP A0LNN5
A	415	LEU	-	expression tag	UNP A0LNN5
A	416	GLU	-	expression tag	UNP A0LNN5
A	417	VAL	-	expression tag	UNP A0LNN5
A	418	LEU	-	expression tag	UNP A0LNN5
A	419	PHE	-	expression tag	UNP A0LNN5
A	420	GLN	-	expression tag	UNP A0LNN5
B	413	LEU	-	expression tag	UNP A0LNN5
B	414	GLU	-	expression tag	UNP A0LNN5
B	415	LEU	-	expression tag	UNP A0LNN5
B	416	GLU	-	expression tag	UNP A0LNN5
B	417	VAL	-	expression tag	UNP A0LNN5
B	418	LEU	-	expression tag	UNP A0LNN5
B	419	PHE	-	expression tag	UNP A0LNN5
B	420	GLN	-	expression tag	UNP A0LNN5

- Molecule 2 is (2S)-2-HYDROXYPROPANOIC ACID (CCD ID: 2OP) (formula: C₃H₆O₃).



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

- Molecule 3 is nonyl beta-D-glucopyranoside (CCD ID: BNG) (formula: $C_{15}H_{30}O_6$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			21	15	6		

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	103.42Å 200.02Å 63.32Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	24.83 – 2.50 24.83 – 2.50	Depositor EDS
% Data completeness (in resolution range)	66.7 (24.83-2.50) 66.6 (24.83-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.21 (at 2.50Å)	Xtriage
Refinement program	PHENIX (1.13_2998: ???)	Depositor
R, R_{free}	0.239 , 0.259 0.240 , 0.260	Depositor DCC
R_{free} test set	1548 reflections (3.34%)	wwPDB-VP
Wilson B-factor (Å ²)	94.8	Xtriage
Anisotropy	0.067	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 69.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	5771	wwPDB-VP
Average B, all atoms (Å ²)	115.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.51% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ 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: 2OP, BNG

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	0.32	0/3070	0.51	0/4177
1	B	0.27	0/2834	0.50	0/3855
All	All	0.29	0/5904	0.50	0/8032

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	2980	0	3023	45	0
1	B	2752	0	2784	50	0
2	A	6	0	5	1	0
2	B	12	0	10	4	0
3	A	21	0	30	0	0
All	All	5771	0	5852	94	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 8.

All (94) 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:174:CYS:HB3	1:A:178:MET:HE3	1.51	0.90
1:B:27:LEU:HD22	1:B:87:LEU:HD11	1.62	0.81
1:B:145:LEU:HD22	2:B:502:2OP:HA	1.65	0.77
1:B:174:CYS:HB3	1:B:178:MET:HE3	1.68	0.75
1:B:26:MET:HE2	1:B:178:MET:HE2	1.74	0.70
1:B:159:ILE:HB	1:B:170:VAL:HG22	1.76	0.68
1:A:97:ILE:HD13	1:A:103:LEU:HD13	1.77	0.65
1:A:26:MET:HB3	1:A:178:MET:HE2	1.79	0.63
1:B:95:GLY:O	1:B:168:ARG:NH2	2.34	0.61
1:A:145:LEU:HD22	2:A:501:2OP:HA	1.82	0.61
1:B:97:ILE:HD13	1:B:103:LEU:HD13	1.82	0.60
1:A:170:VAL:O	1:A:174:CYS:HB2	2.02	0.59
1:B:41:GLU:CD	1:B:167:TRP:HE1	2.12	0.58
1:A:396:LEU:HD11	1:B:308:ILE:HD11	1.86	0.58
1:A:31:TRP:O	1:A:35:ILE:HG13	2.04	0.57
1:B:87:LEU:HG	1:B:91:PHE:CE2	2.40	0.57
1:B:84:GLY:HA2	1:B:118:ILE:HD11	1.87	0.56
1:B:348:PRO:HA	1:B:351:GLN:HB2	1.88	0.56
1:A:23:THR:HG23	1:A:178:MET:HB3	1.87	0.55
1:A:245:LEU:HD13	1:A:366:GLY:HA2	1.86	0.55
1:B:245:LEU:HD13	1:B:366:GLY:HA2	1.87	0.55
1:B:16:LEU:HD21	1:B:188:PHE:HB2	1.91	0.53
1:A:97:ILE:HG22	1:A:102:GLN:HG2	1.91	0.52
1:B:382:THR:HG23	1:B:384:TYR:H	1.73	0.52
1:A:107:TYR:O	1:A:111:ALA:HB3	2.09	0.52
1:A:303:GLN:OE1	1:A:327:ILE:HA	2.11	0.51
1:B:139:GLY:O	1:B:143:VAL:HB	2.10	0.50
1:B:247:VAL:HG21	1:B:307:MET:SD	2.51	0.50
1:A:65:PHE:HB3	1:A:66:PRO:HD3	1.93	0.50
1:A:86:LEU:HD13	1:A:113:PHE:CD1	2.47	0.50
1:A:418:LEU:C	1:A:420:GLN:H	2.19	0.50
1:A:95:GLY:HA3	1:A:172:TRP:NE1	2.28	0.49
1:A:295:TYR:CE2	1:A:299:LEU:HD11	2.47	0.49
1:B:70:LEU:HD22	1:B:74:MET:HE3	1.95	0.49
1:A:132:ASP:OD1	1:A:132:ASP:N	2.36	0.48
1:B:107:TYR:O	1:B:111:ALA:HB3	2.13	0.48
1:B:303:GLN:O	1:B:307:MET:HG2	2.14	0.47
1:B:335:PHE:CE2	2:B:502:2OP:HB3	2.50	0.47
1:B:132:ASP:N	1:B:132:ASP:OD1	2.39	0.47
1:B:355:TYR:CE1	1:B:358:LEU:HD23	2.50	0.47
1:A:91:PHE:CD1	1:A:171:PHE:HD1	2.32	0.47
1:A:243:ALA:HB1	1:A:390:ALA:HB1	1.95	0.47

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:303:GLN:O	1:A:307:MET:HG2	2.14	0.47
1:B:26:MET:HB3	1:B:178:MET:HE2	1.97	0.47
1:A:119:TYR:CE1	1:A:123:ILE:HD11	2.50	0.47
1:B:36:LYS:HB3	1:B:37:PRO:HD3	1.97	0.47
1:B:23:THR:HG23	1:B:178:MET:HB3	1.97	0.46
1:A:16:LEU:HD22	1:A:185:ALA:HA	1.98	0.46
1:B:87:LEU:HG	1:B:91:PHE:CZ	2.49	0.46
1:A:409:LYS:HA	1:A:409:LYS:HD2	1.78	0.46
1:B:292:ILE:HG21	1:B:341:THR:OG1	2.15	0.46
1:B:310:ILE:HB	1:B:323:VAL:HG11	1.98	0.46
1:B:30:ALA:O	1:B:33:VAL:HG22	2.16	0.45
1:B:11:TRP:HB3	1:B:130:TRP:HZ2	1.79	0.45
1:B:295:TYR:CZ	1:B:299:LEU:HD11	2.51	0.45
1:B:335:PHE:HE2	1:B:362:CYS:SG	2.39	0.45
1:A:376:LEU:HD22	1:A:385:LEU:HD23	1.99	0.45
1:A:221:TYR:O	1:A:225:LYS:HB3	2.17	0.45
1:A:87:LEU:HA	1:A:111:ALA:HA	1.98	0.45
1:A:295:TYR:CZ	1:A:299:LEU:HD11	2.51	0.45
1:B:58:LEU:O	1:B:62:LEU:HG	2.17	0.45
1:A:86:LEU:O	1:A:110:ILE:HG22	2.17	0.44
1:A:113:PHE:O	1:A:117:MET:HG3	2.18	0.44
1:B:218:ASP:OD1	1:B:348:PRO:HD2	2.18	0.44
1:B:65:PHE:HB3	1:B:66:PRO:HD3	1.99	0.44
1:B:303:GLN:OE1	1:B:327:ILE:HA	2.18	0.44
1:A:217:ARG:O	1:A:409:LYS:HE3	2.18	0.44
1:B:360:THR:O	1:B:364:LEU:HG	2.18	0.43
1:A:27:LEU:HD22	1:A:87:LEU:HD13	2.00	0.43
1:A:218:ASP:OD2	1:A:348:PRO:HD2	2.18	0.43
1:B:167:TRP:HA	1:B:170:VAL:HG23	2.00	0.43
1:A:91:PHE:HB3	1:A:175:GLY:HA3	2.01	0.43
1:A:271:SER:HA	1:A:274:PHE:HD2	1.83	0.43
1:A:150:PHE:HE2	1:A:274:PHE:CE1	2.37	0.43
1:B:42:PHE:HD2	1:B:100:LYS:NZ	2.17	0.43
1:B:58:LEU:HD22	1:B:367:PHE:CG	2.53	0.43
1:B:230:PHE:N	1:B:346:TYR:OH	2.52	0.42
1:A:322:ILE:O	1:A:326:VAL:HG23	2.19	0.42
1:B:280:ARG:NH1	2:B:502:2OP:OXT	2.51	0.42
1:A:109:VAL:O	1:A:113:PHE:HB2	2.19	0.42
1:A:174:CYS:O	1:A:178:MET:HG3	2.20	0.42
1:B:303:GLN:HG3	1:B:307:MET:HE2	2.01	0.42
1:A:384:TYR:O	1:A:388:LEU:HG	2.20	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:218:ASP:CG	1:B:348:PRO:HD2	2.44	0.42
1:A:234:TYR:CE1	1:A:361:ALA:HB2	2.55	0.41
1:A:412:ALA:O	1:A:416:GLU:N	2.52	0.41
1:B:234:TYR:HE2	1:B:357:LEU:O	2.02	0.41
1:A:126:ALA:N	1:A:127:PRO:HD2	2.35	0.41
1:A:377:LYS:O	1:A:381:GLY:N	2.51	0.41
1:B:117:MET:HE2	1:B:117:MET:HB3	1.91	0.41
1:B:280:ARG:HH22	2:B:502:2OP:C	2.33	0.41
1:A:337:LEU:HD23	1:A:337:LEU:HA	1.77	0.40
1:B:64:THR:HG21	1:B:359:PHE:CD2	2.56	0.40
1:B:252:ALA:HA	1:B:269:VAL:HG21	2.02	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	387/420 (92%)	380 (98%)	7 (2%)	0	100	100
1	B	358/420 (85%)	354 (99%)	4 (1%)	0	100	100
All	All	745/840 (89%)	734 (98%)	11 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	291/307 (95%)	290 (100%)	1 (0%)	86	94
1	B	267/307 (87%)	267 (100%)	0	100	100
All	All	558/614 (91%)	557 (100%)	1 (0%)	87	95

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	290	ILE

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

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

4 ligands are modelled in this entry.

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	2OP	B	502	-	4,5,5	1.14	0	2,6,6	1.18	0
2	2OP	A	501	-	4,5,5	1.13	0	2,6,6	0.63	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	BNG	A	502	-	21,21,21	1.10	2 (9%)	26,26,26	1.15	3 (11%)
2	2OP	B	501	-	4,5,5	1.06	0	2,6,6	1.48	1 (50%)

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	2OP	B	502	-	-	2/4/4/4	-
2	2OP	A	501	-	-	2/4/4/4	-
3	BNG	A	502	-	-	3/12/32/32	0/1/1/1
2	2OP	B	501	-	-	4/4/4/4	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	502	BNG	O5-C1	3.03	1.49	1.41
3	A	502	BNG	O1-C1	-2.03	1.36	1.40

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	502	BNG	O1-C1-C2	2.61	112.23	108.27
3	A	502	BNG	C4-C3-C2	2.41	115.06	110.83
3	A	502	BNG	C3-C4-C5	2.20	114.22	110.23
2	B	501	2OP	OXT-C-O	-2.08	119.35	124.08

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	501	2OP	O-C-CA-CB
2	B	501	2OP	OXT-C-CA-CB
3	A	502	BNG	O5-C5-C6-O6
2	B	501	2OP	OXT-C-CA-OHN
3	A	502	BNG	C2'-C1'-O1-C1
3	A	502	BNG	C4-C5-C6-O6
2	A	501	2OP	OXT-C-CA-CB
2	B	502	2OP	OXT-C-CA-CB

Continued on next page...

Continued from previous page...

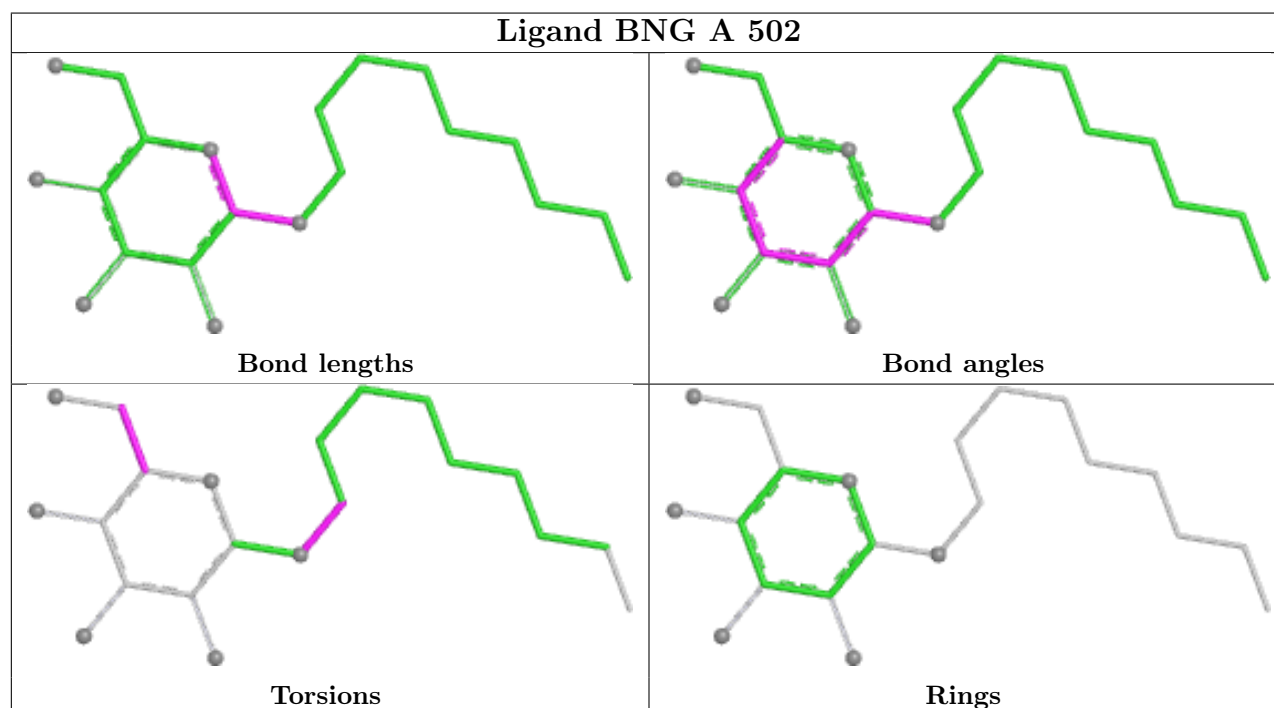
Mol	Chain	Res	Type	Atoms
2	B	501	2OP	O-C-CA-OHN
2	A	501	2OP	O-C-CA-CB
2	B	502	2OP	O-C-CA-CB

There are no ring outliers.

2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	502	2OP	4	0
2	A	501	2OP	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 than 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.



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

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, 95th 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	395/420 (94%)	0.62	43 (10%) 10 9	54, 94, 165, 209	0
1	B	366/420 (87%)	0.80	53 (14%) 6 5	61, 127, 188, 217	0
All	All	761/840 (90%)	0.71	96 (12%) 8 6	54, 107, 180, 217	0

All (96) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	151	LEU	7.8
1	B	125	THR	6.7
1	B	29	TYR	5.4
1	A	420	GLN	5.4
1	B	33	VAL	5.2
1	A	167	TRP	4.8
1	A	38	LEU	4.6
1	B	119	TYR	4.5
1	A	150	PHE	4.4
1	B	91	PHE	4.3
1	B	123	ILE	4.3
1	B	111	ALA	4.3
1	A	29	TYR	4.1
1	B	124	ALA	4.1
1	A	155	LEU	4.1
1	A	73	LYS	3.9
1	B	121	PRO	3.9
1	B	30	ALA	3.9
1	B	129	TRP	3.8
1	A	34	PHE	3.7
1	A	31	TRP	3.6
1	A	196	TRP	3.6
1	B	110	ILE	3.5
1	B	191	PRO	3.4

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	88	ALA	3.4
1	B	92	ILE	3.4
1	B	46	ARG	3.3
1	B	157	THR	3.3
1	A	30	ALA	3.2
1	B	241	SER	3.2
1	A	35	ILE	3.2
1	A	342	CYS	3.1
1	B	34	PHE	3.1
1	A	100	LYS	3.0
1	B	89	ILE	2.9
1	A	159	ILE	2.9
1	B	42	PHE	2.9
1	B	225	LYS	2.9
1	B	258	ALA	2.9
1	A	25	GLY	2.8
1	B	189	LEU	2.8
1	A	156	ALA	2.8
1	B	74	MET	2.8
1	A	169	TYR	2.7
1	B	154	PRO	2.7
1	B	70	LEU	2.7
1	B	407	PRO	2.7
1	A	413	LEU	2.7
1	A	98	GLN	2.7
1	A	28	LEU	2.6
1	A	160	ILE	2.5
1	A	199	ALA	2.5
1	A	172	TRP	2.5
1	B	188	PHE	2.5
1	A	198	PRO	2.5
1	A	88	ALA	2.5
1	B	174	CYS	2.4
1	A	41	GLU	2.4
1	A	87	LEU	2.4
1	A	363	GLY	2.4
1	A	93	LEU	2.4
1	A	212	ALA	2.4
1	B	224	ALA	2.4
1	A	414	GLU	2.4
1	B	134	ARG	2.3
1	B	335	PHE	2.3

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	417	VAL	2.3
1	B	38	LEU	2.3
1	B	151	LEU	2.3
1	B	118	ILE	2.3
1	A	362	CYS	2.3
1	A	158	TYR	2.3
1	B	82	THR	2.3
1	B	98	GLN	2.3
1	A	149	SER	2.3
1	A	345	PHE	2.3
1	A	170	VAL	2.2
1	B	167	TRP	2.2
1	B	231	TRP	2.2
1	A	7	THR	2.2
1	B	19	LEU	2.2
1	B	364	LEU	2.2
1	A	191	PRO	2.2
1	B	171	PHE	2.2
1	B	126	ALA	2.2
1	B	166	GLY	2.2
1	B	97	ILE	2.1
1	B	351	GLN	2.1
1	B	217	ARG	2.1
1	B	156	ALA	2.1
1	A	174	CYS	2.1
1	B	348	PRO	2.1
1	B	85	VAL	2.0
1	B	112	GLY	2.0
1	B	128	LYS	2.0
1	A	128	LYS	2.0

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

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

6.3 Carbohydrates ⓘ

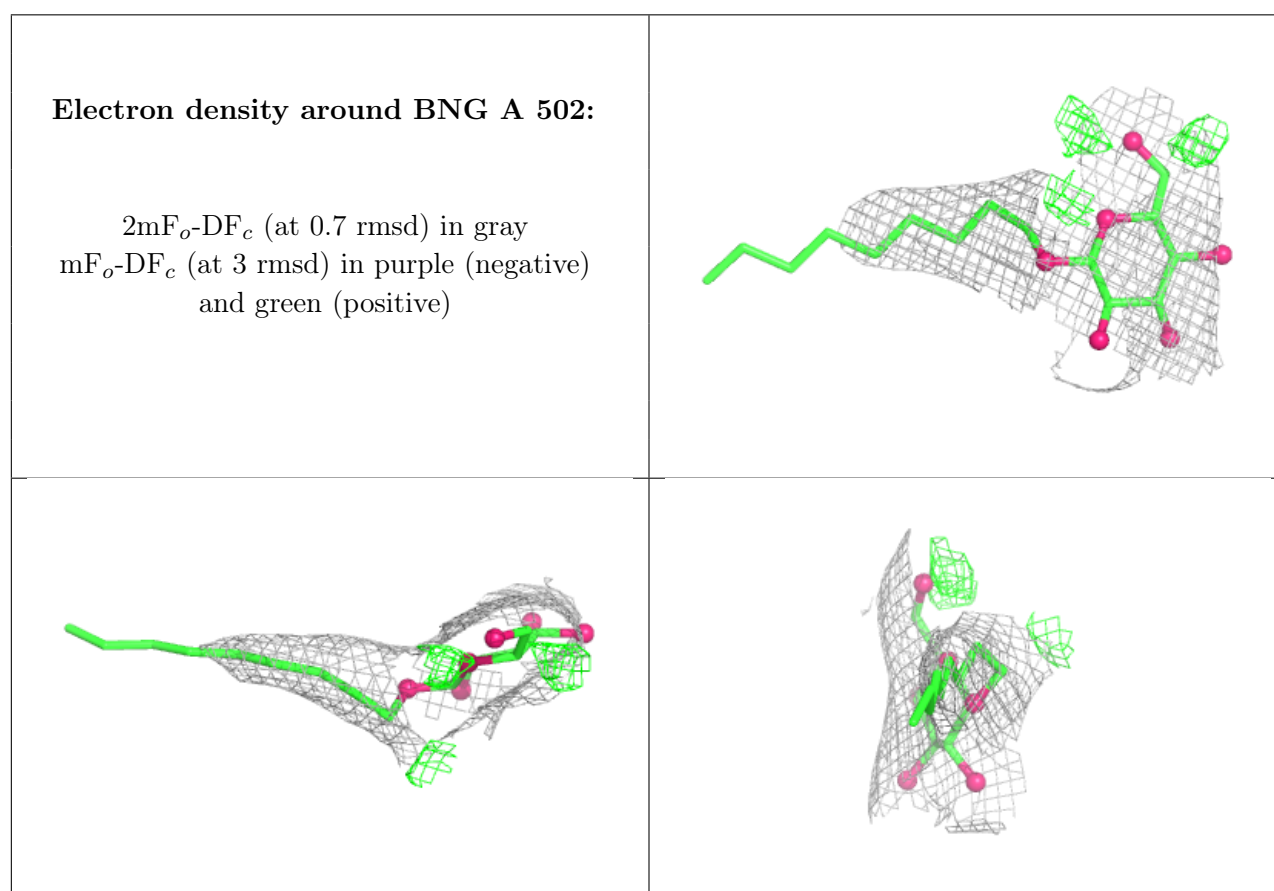
There are no oligosaccharides in this entry.

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, 95th 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
2	2OP	B	501	6/6	0.72	0.25	106,123,128,128	0
2	2OP	B	502	6/6	0.86	0.18	106,115,119,121	0
3	BNG	A	502	21/21	0.86	0.16	101,125,151,164	0
2	2OP	A	501	6/6	0.90	0.20	81,110,110,119	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.



6.5 Other polymers [i](#)

There are no such residues in this entry.