



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 08:26 PM UTC

PDB ID : 2X0H / pdb_00002x0h
Title : BtGH84 Michaelis complex
Authors : He, Y.; Davies, G.J.
Deposited on : 2009-12-08
Resolution : 2.21 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

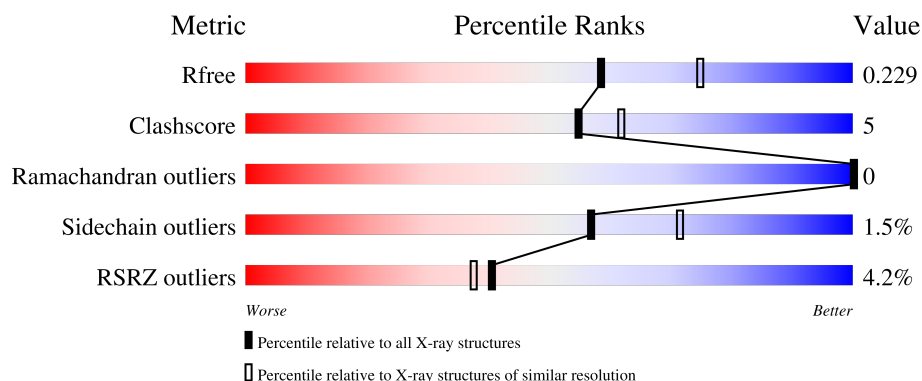
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.21 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	7682 (2.24-2.20)
Clashscore	190562	8402 (2.24-2.20)
Ramachandran outliers	187476	8303 (2.24-2.20)
Sidechain outliers	187428	8304 (2.24-2.20)
RSRZ outliers	180081	7683 (2.24-2.20)

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	737	3% 74% 11% 15%
1	B	737	4% 76% 11% 13%

2 Entry composition [i](#)

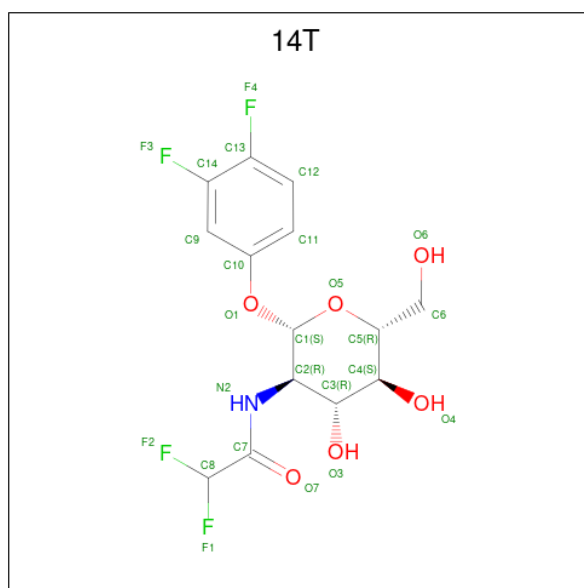
There are 5 unique types of molecules in this entry. The entry contains 11257 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 O-GLCNACASE BT_4395.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	630	Total	C	N	O	S	0	3	0
			5168	3323	868	958	19			
1	B	642	Total	C	N	O	S	0	6	0
			5269	3373	896	981	19			

- Molecule 2 is 3,4-difluorophenyl 2-deoxy-2-[(difluoroacetyl)amino]-beta-D-glucopyranoside (CCD ID: 14T) (formula: C₁₄H₁₅F₄NO₆).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	F	N	O	0	0
			25	14	4	1	6		
2	B	1	Total	C	F	N	O	0	0
			25	14	4	1	6		

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



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

- Molecule 4 is CALCIUM ION (CCD ID: CA) (formula: Ca).

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

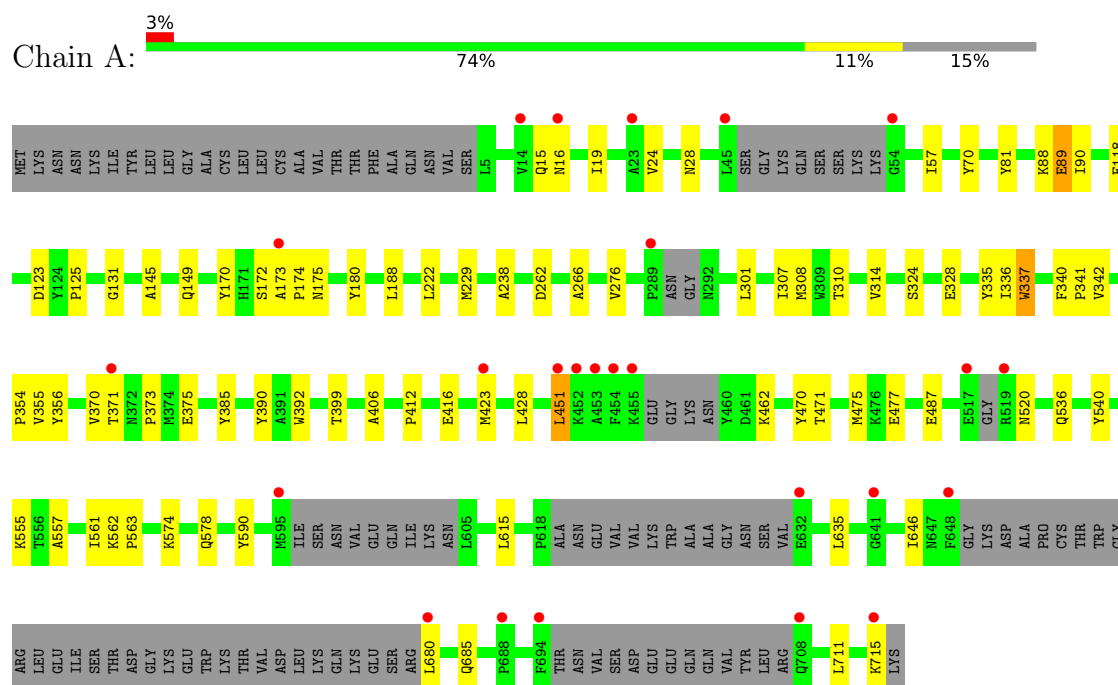
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	325	Total	O	0	0
			325	325		
5	B	408	Total	O	0	0
			408	408		

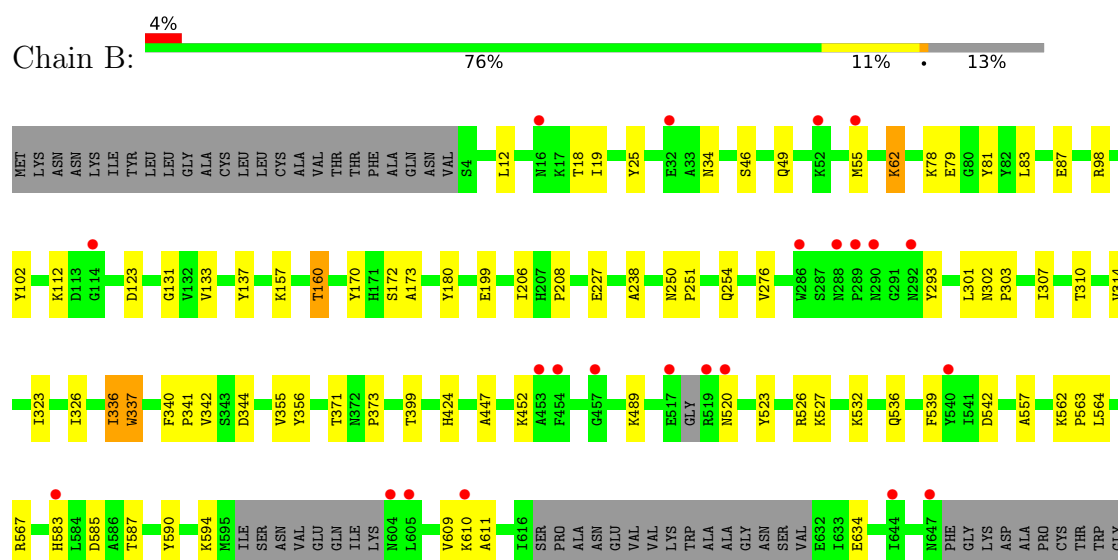
3 Residue-property plots [i](#)

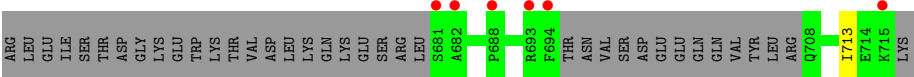
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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



• Molecule 1: O-GLCNACASE BT_4395





4 Data and refinement statistics

Property	Value	Source
Space group	P 2 ₁ 2 ₁ 2 ₁	Depositor
Cell constants a, b, c, α , β , γ	51.63Å 163.15Å 224.55Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.08 – 2.21 48.08 – 2.21	Depositor EDS
% Data completeness (in resolution range)	99.2 (48.08-2.21) 99.2 (48.08-2.21)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.96 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.194 , 0.226 0.201 , 0.229	Depositor DCC
R_{free} test set	4769 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	30.1	Xtriage
Anisotropy	0.152	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 44.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.36$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	11257	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 22.53 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 5.5926e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: CA, 14T, GOL

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	0.90	3/5297 (0.1%)	0.93	2/7175 (0.0%)
1	B	0.97	4/5402 (0.1%)	0.97	6/7313 (0.1%)
All	All	0.94	7/10699 (0.1%)	0.95	8/14488 (0.1%)

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	18	THR	C-N	-7.23	1.25	1.33
1	B	172	SER	C-O	-5.61	1.18	1.24
1	A	390	TYR	C-O	-5.37	1.18	1.24
1	B	539	PHE	C-O	-5.09	1.18	1.24
1	A	172	SER	C-O	-5.07	1.17	1.24
1	A	336	ILE	C-O	-5.05	1.18	1.24
1	B	542	ASP	C-O	-5.03	1.17	1.24

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	326	ILE	N-CA-C	5.87	116.05	110.53
1	A	336	ILE	N-CA-C	5.77	116.52	109.30
1	B	336	ILE	N-CA-C	5.71	116.43	108.89
1	B	611	ALA	CB-CA-C	-5.36	109.94	117.23
1	A	175	ASN	N-CA-C	5.19	118.64	112.72
1	B	160	THR	N-CA-C	5.19	116.87	108.26
1	B	489	LYS	CA-C-N	-5.09	113.40	119.05
1	B	489	LYS	C-N-CA	-5.09	113.40	119.05

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	5168	0	5077	52	0
1	B	5269	0	5179	50	0
2	A	25	0	7	1	0
2	B	25	0	7	1	0
3	A	12	0	16	1	0
3	B	24	0	32	2	0
4	B	1	0	0	0	0
5	A	325	0	0	8	0
5	B	408	0	0	7	0
All	All	11257	0	10318	101	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 5.

All (101) 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:19:ILE:HD12	1:B:55:MET:HE1	1.47	0.96
1:A:355:VAL:O	1:A:399:THR:HG23	1.76	0.85
1:B:536:GLN:HG2	1:B:590:TYR:CD1	2.17	0.79
1:B:355:VAL:O	1:B:399:THR:HG23	1.84	0.76
1:A:324:SER:O	1:A:328:GLU:HG3	1.91	0.70
1:A:375:GLU:CD	5:A:2197:HOH:O	2.34	0.69
1:B:563:PRO:O	1:B:567[B]:ARG:HG3	1.94	0.68
1:A:15:GLN:O	1:A:16:ASN:HB2	1.93	0.68
1:B:585:ASP:OD1	1:B:587:THR:HG23	1.97	0.65
1:B:46:SER:CA	5:B:2040:HOH:O	2.43	0.65
1:A:536:GLN:HG2	1:A:590:TYR:CD1	2.34	0.62
1:B:250:ASN:O	1:B:254:GLN:HG3	2.00	0.61
1:A:356:TYR:HB3	1:A:399:THR:HG21	1.82	0.61
1:B:583[B]:HIS:O	1:B:583[B]:HIS:CD2	2.54	0.61
1:B:532:LYS:O	1:B:536:GLN:HG3	2.01	0.60
1:A:354:PRO:HB2	1:A:399:THR:HG22	1.84	0.60
1:A:562:LYS:HB3	1:A:563:PRO:HD3	1.85	0.57
1:A:646:ILE:O	1:A:680:LEU:HB2	2.03	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:520:ASN:OD1	1:B:523:TYR:N	2.36	0.57
1:A:555:LYS:HE3	5:A:2289:HOH:O	2.04	0.56
1:A:536:GLN:O	1:A:540[A]:TYR:HD2	1.89	0.54
3:A:1716:GOL:H2	5:A:2324:HOH:O	2.08	0.54
1:A:412:PRO:HD2	1:A:487:GLU:OE1	2.08	0.53
1:A:423[B]:MET:CE	1:A:423[B]:MET:HA	2.39	0.53
1:B:340:PHE:O	1:B:371:THR:OG1	2.25	0.53
1:B:19:ILE:CD1	1:B:55:MET:HE1	2.31	0.52
1:B:447:ALA:HB1	1:B:564:LEU:CD1	2.39	0.52
1:A:342:VAL:HG23	5:A:2196:HOH:O	2.10	0.52
1:A:462:LYS:NZ	5:A:2252:HOH:O	2.43	0.52
1:B:83:LEU:HD23	1:B:83:LEU:C	2.36	0.51
1:A:188:LEU:HD13	1:A:229:MET:HE1	1.91	0.51
1:B:78:LYS:O	1:B:79:GLU:HB2	2.11	0.51
1:A:557:ALA:HB1	1:A:561:ILE:HB	1.92	0.50
1:B:356:TYR:HB3	1:B:399:THR:HG21	1.94	0.49
1:B:323:ILE:HD12	1:B:336:ILE:HD11	1.94	0.49
1:B:137:TYR:HB2	3:B:1718:GOL:H32	1.94	0.49
1:A:170:TYR:HB2	1:A:180:TYR:CE1	2.48	0.49
1:B:344:ASP:O	3:B:1718:GOL:H31	2.13	0.48
1:A:423[B]:MET:HA	1:A:423[B]:MET:HE3	1.95	0.48
1:B:81:TYR:CE2	1:B:123:ASP:HB3	2.49	0.48
1:B:310:THR:HA	1:B:337:TRP:O	2.13	0.48
1:A:477:GLU:OE1	1:B:527:LYS:NZ	2.40	0.47
1:A:314:VAL:HG11	2:A:1000:14T:H9	1.96	0.47
1:B:238:ALA:HA	1:B:276:VAL:O	2.14	0.47
1:A:471:THR:O	1:A:475:MET:HG3	2.14	0.47
1:A:262:ASP:HA	1:A:266:ALA:HB3	1.98	0.46
1:A:15:GLN:HB2	1:A:118:GLU:HB3	1.96	0.46
1:A:145:ALA:O	1:A:149:GLN:HG2	2.15	0.46
1:B:81:TYR:CZ	1:B:123:ASP:HB3	2.50	0.46
1:A:173:ALA:HA	1:A:174:PRO:HA	1.65	0.46
1:A:340:PHE:O	1:A:371:THR:OG1	2.34	0.46
1:B:34:ASN:OD1	1:B:34:ASN:C	2.59	0.46
1:A:81:TYR:CZ	1:A:123:ASP:HB3	2.51	0.46
1:A:125:PRO:HB3	1:A:392:TRP:CE3	2.51	0.46
1:A:635:LEU:HD11	1:A:711:LEU:CD2	2.47	0.45
1:A:222:LEU:C	1:A:222:LEU:HD23	2.41	0.45
1:A:301:LEU:HD23	1:A:307:ILE:HD11	1.99	0.45
1:B:314:VAL:HG11	2:B:1000:14T:H9	1.98	0.45
1:A:308:MET:HA	1:A:335:TYR:O	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:562:LYS:NZ	5:A:2306:HOH:O	2.49	0.45
1:B:206:ILE:HG23	1:B:208:PRO:HD3	1.99	0.44
1:B:424:HIS:O	1:B:557:ALA:HA	2.17	0.44
1:B:562:LYS:N	1:B:563:PRO:CD	2.80	0.44
1:B:609:VAL:O	1:B:610:LYS:HD3	2.17	0.44
1:A:310:THR:HA	1:A:337:TRP:O	2.18	0.44
1:A:354:PRO:HD2	5:A:2244:HOH:O	2.17	0.44
1:B:567[B]:ARG:NH2	5:B:2377:HOH:O	2.50	0.44
1:B:594:LYS:HB2	1:B:634:GLU:HB3	1.99	0.44
1:B:173:ALA:HB2	5:B:2186:HOH:O	2.17	0.44
1:A:385:TYR:CD2	1:A:406:ALA:HB2	2.52	0.44
1:B:25:TYR:CE2	1:B:49:GLN:HG3	2.53	0.43
1:A:520:ASN:OD1	1:A:520:ASN:C	2.61	0.43
1:A:341:PRO:HD2	1:A:373:PRO:HA	2.01	0.43
1:A:238:ALA:HA	1:A:276:VAL:O	2.18	0.43
1:A:70:TYR:OH	1:A:89:GLU:OE2	2.37	0.43
1:A:131:GLY:O	1:A:370:VAL:HA	2.19	0.43
1:B:251:PRO:HB3	1:B:293:TYR:CD1	2.54	0.43
1:A:555:LYS:CE	5:A:2289:HOH:O	2.66	0.43
1:A:371:THR:O	1:A:371:THR:HG23	2.19	0.42
1:A:574:LYS:O	1:A:578:GLN:HG3	2.19	0.42
1:B:131:GLY:HA3	1:B:160:THR:O	2.19	0.42
1:B:562:LYS:HB3	1:B:563:PRO:HD3	2.01	0.42
1:B:341:PRO:HD2	1:B:373:PRO:HA	2.01	0.42
1:B:102:TYR:CE1	1:B:157:LYS:HA	2.55	0.42
1:B:447:ALA:HB1	1:B:564:LEU:HD12	2.02	0.41
1:A:81:TYR:CE2	1:A:123:ASP:HB3	2.56	0.41
1:A:470:TYR:HE1	1:B:526:ARG:NH1	2.18	0.41
1:A:536:GLN:HG2	1:A:590:TYR:CG	2.55	0.41
1:B:170:TYR:HB2	1:B:180:TYR:CE1	2.56	0.41
1:B:227:GLU:HB2	5:B:2196:HOH:O	2.21	0.41
1:B:98:ARG:HG3	1:B:199:GLU:HB3	2.03	0.41
1:A:428:LEU:HD12	1:A:428:LEU:N	2.36	0.41
1:B:87:GLU:HG2	5:B:2099:HOH:O	2.20	0.40
1:B:342:VAL:HG23	5:B:2278:HOH:O	2.21	0.40
1:A:57:ILE:HG12	1:A:90:ILE:HB	2.04	0.40
1:A:451:LEU:HD12	1:A:451:LEU:HA	1.83	0.40
1:B:133:VAL:HG13	1:B:133:VAL:O	2.22	0.40
1:B:301:LEU:HD12	1:B:307:ILE:HD11	2.02	0.40
1:A:341:PRO:O	1:A:342:VAL:C	2.64	0.40
1:B:62:LYS:HD2	5:B:2082:HOH:O	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:302:ASN:HA	1:B:303:PRO:HD3	1.95	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	615/737 (83%)	596 (97%)	19 (3%)	0	100	100
1	B	636/737 (86%)	616 (97%)	20 (3%)	0	100	100
All	All	1251/1474 (85%)	1212 (97%)	39 (3%)	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	559/647 (86%)	548 (98%)	11 (2%)	48	63
1	B	570/647 (88%)	564 (99%)	6 (1%)	65	78
All	All	1129/1294 (87%)	1112 (98%)	17 (2%)	57	71

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	19	ILE
1	A	24	VAL
1	A	28	ASN
1	A	88	LYS
1	A	89	GLU
1	A	337	TRP
1	A	416	GLU
1	A	451	LEU
1	A	615	LEU
1	A	685	GLN
1	A	715	LYS
1	B	12	LEU
1	B	62	LYS
1	B	112	LYS
1	B	337	TRP
1	B	452	LYS
1	B	713	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	10	GLN
1	A	11	GLN
1	A	73	GLN
1	A	252	GLN
1	A	254	GLN
1	A	267	GLN
1	A	273	ASN
1	A	274	GLN
1	A	529	ASN
1	A	547	GLN
1	A	583	HIS
1	A	608	GLN
1	A	645	GLN
1	B	15	GLN
1	B	73	GLN
1	B	144	GLN
1	B	349	HIS
1	B	433	HIS
1	B	459	ASN
1	B	529	ASN
1	B	581	ASN

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 ⓘ

Of 9 ligands modelled in this entry, 1 is monoatomic - leaving 8 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	GOL	B	1718	-	5,5,5	0.46	0	5,5,5	0.74	0
3	GOL	A	1717	-	5,5,5	0.36	0	5,5,5	0.41	0
3	GOL	B	1719	-	5,5,5	0.44	0	5,5,5	0.79	0
3	GOL	A	1716	-	5,5,5	0.46	0	5,5,5	1.15	0
2	14T	A	1000	-	25,26,26	1.11	2 (8%)	35,37,37	2.30	8 (22%)
3	GOL	B	1717	-	5,5,5	0.38	0	5,5,5	0.17	0
3	GOL	B	1720	-	5,5,5	0.47	0	5,5,5	0.73	0
2	14T	B	1000	-	25,26,26	1.02	2 (8%)	35,37,37	1.91	5 (14%)

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	GOL	B	1718	-	-	4/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GOL	A	1717	-	-	4/4/4/4	-
3	GOL	B	1719	-	-	2/4/4/4	-
3	GOL	A	1716	-	-	2/4/4/4	-
2	14T	A	1000	-	-	5/14/34/34	0/2/2/2
3	GOL	B	1717	-	-	4/4/4/4	-
3	GOL	B	1720	-	-	1/4/4/4	-
2	14T	B	1000	-	-	3/14/34/34	0/2/2/2

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1000	14T	C1-C2	3.54	1.58	1.53
2	B	1000	14T	C1-C2	2.79	1.57	1.53
2	B	1000	14T	O1-C1	2.39	1.44	1.41
2	A	1000	14T	O1-C1	2.35	1.44	1.41

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1000	14T	C10-O1-C1	7.83	128.85	117.82
2	B	1000	14T	C10-O1-C1	7.48	128.35	117.82
2	A	1000	14T	C1-O5-C5	4.90	123.28	113.72
2	A	1000	14T	C1-C2-N2	4.48	118.43	110.92
2	A	1000	14T	O1-C1-C2	3.88	112.35	107.12
2	B	1000	14T	C1-C2-N2	3.88	117.43	110.92
2	B	1000	14T	C1-O5-C5	3.78	121.10	113.72
2	A	1000	14T	C3-C2-N2	-3.59	104.01	110.62
2	A	1000	14T	O5-C1-O1	-3.34	99.88	108.42
2	A	1000	14T	C2-N2-C7	3.26	128.16	123.20
2	B	1000	14T	O5-C1-C2	2.77	115.82	110.59
2	B	1000	14T	C2-N2-C7	2.50	127.00	123.20
2	A	1000	14T	C1-C2-C3	2.09	115.95	110.19

There are no chirality outliers.

All (25) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1000	14T	N2-C7-C8-F2
2	B	1000	14T	N2-C7-C8-F2
3	A	1716	GOL	C1-C2-C3-O3

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Mol	Chain	Res	Type	Atoms
3	A	1716	GOL	O2-C2-C3-O3
3	A	1717	GOL	O1-C1-C2-C3
3	A	1717	GOL	C1-C2-C3-O3
3	A	1717	GOL	O2-C2-C3-O3
3	B	1717	GOL	O1-C1-C2-C3
3	B	1718	GOL	C1-C2-C3-O3
3	B	1717	GOL	C1-C2-C3-O3
3	B	1718	GOL	O1-C1-C2-C3
3	B	1719	GOL	O1-C1-C2-C3
3	B	1720	GOL	C1-C2-C3-O3
3	B	1717	GOL	O1-C1-C2-O2
3	B	1718	GOL	O2-C2-C3-O3
3	A	1717	GOL	O1-C1-C2-O2
3	B	1719	GOL	O1-C1-C2-O2
2	B	1000	14T	C1-C2-N2-C7
2	A	1000	14T	C3-C2-N2-C7
3	B	1718	GOL	O1-C1-C2-O2
2	B	1000	14T	C3-C2-N2-C7
2	A	1000	14T	C1-C2-N2-C7
3	B	1717	GOL	O2-C2-C3-O3
2	A	1000	14T	C11-C10-O1-C1
2	A	1000	14T	C9-C10-O1-C1

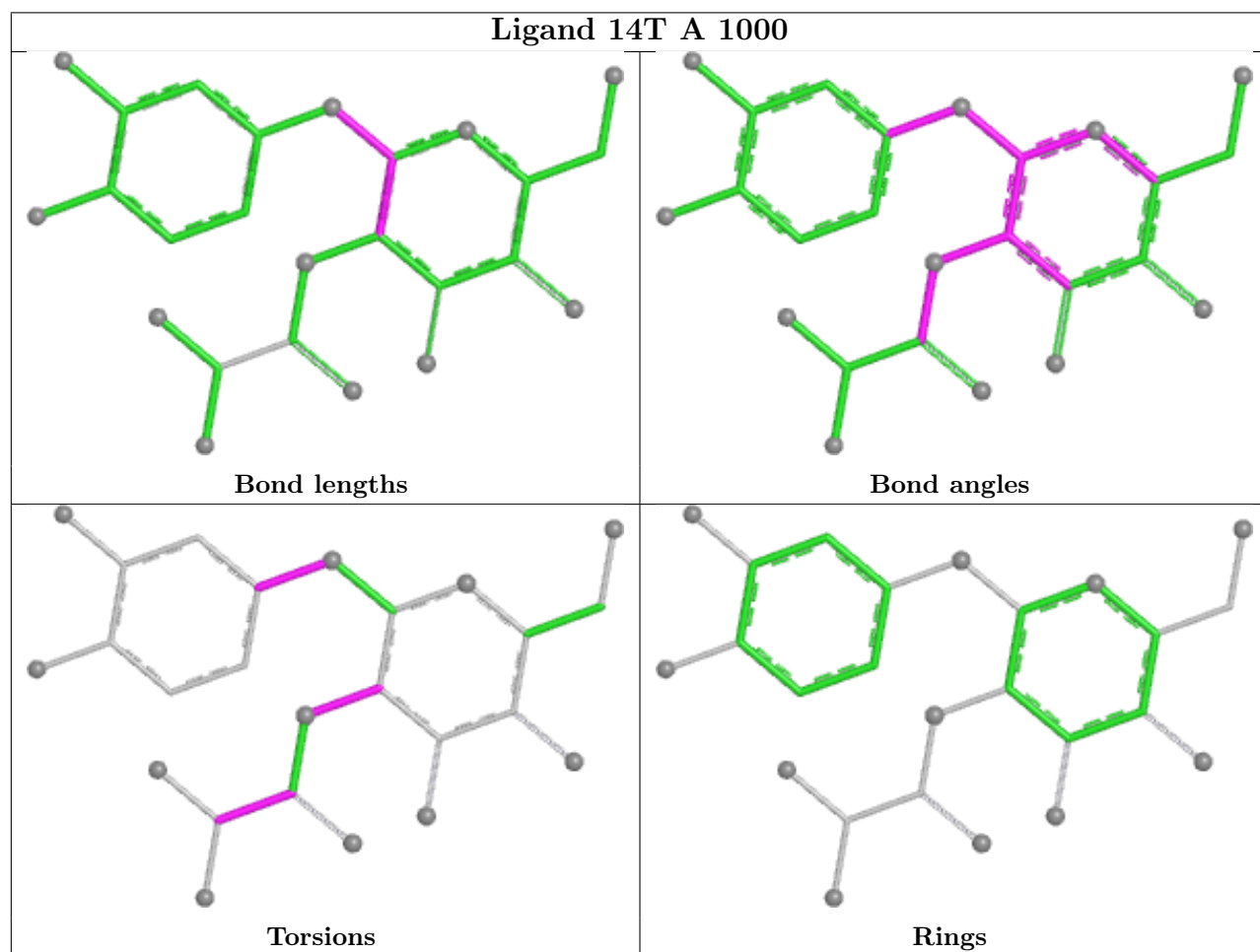
There are no ring outliers.

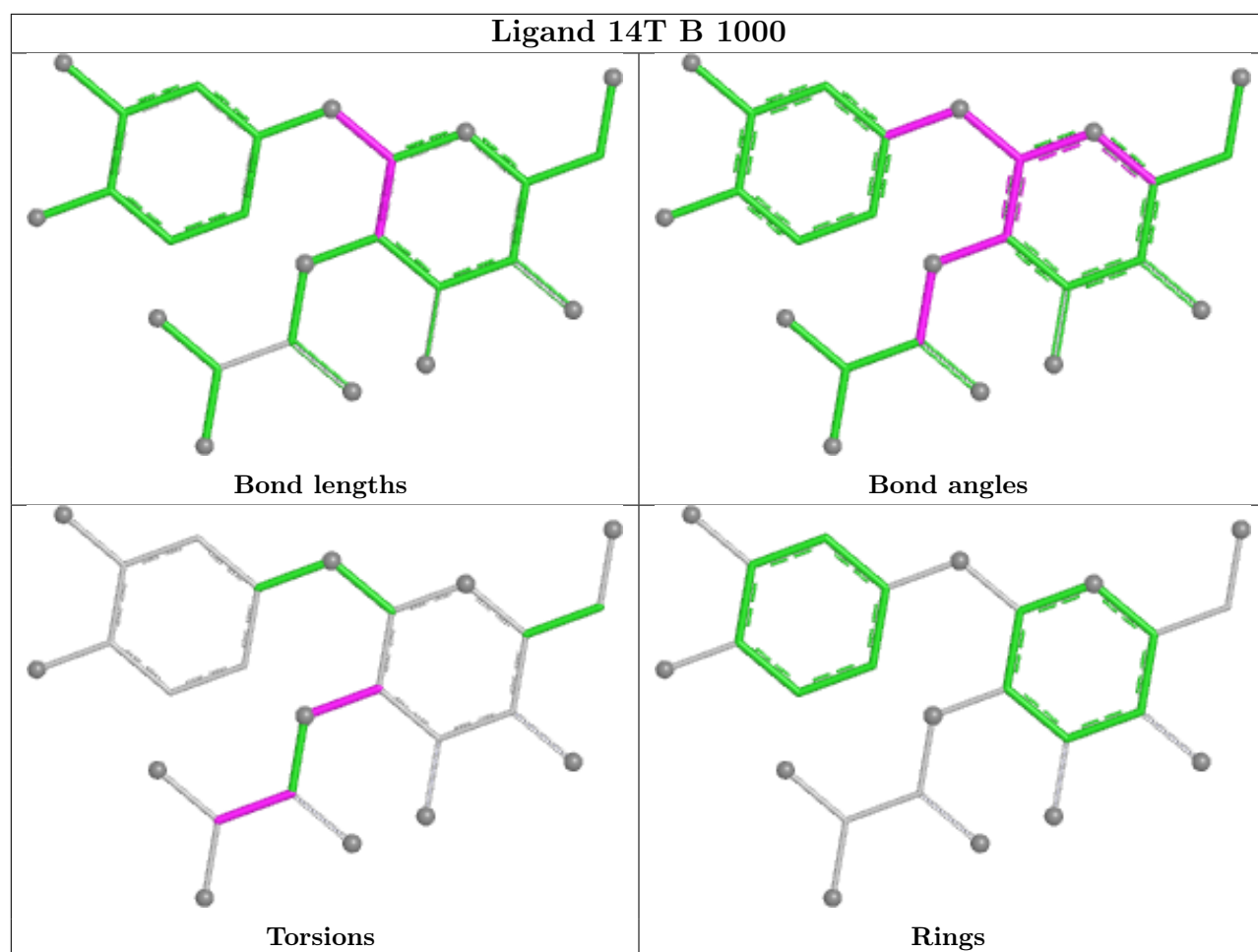
4 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	1718	GOL	2	0
3	A	1716	GOL	1	0
2	A	1000	14T	1	0
2	B	1000	14T	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	630/737 (85%)	0.02	25 (3%)	42 39	16, 37, 74, 94	3 (0%)
1	B	642/737 (87%)	0.02	29 (4%)	38 35	15, 35, 77, 113	6 (0%)
All	All	1272/1474 (86%)	0.02	54 (4%)	40 37	15, 36, 76, 113	9 (0%)

All (54) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	519	ARG	4.9
1	A	54	GLY	4.9
1	B	540	TYR	4.9
1	B	583[A]	HIS	4.8
1	A	519	ARG	4.2
1	A	648	PHE	4.1
1	A	454	PHE	4.1
1	B	647	ASN	4.1
1	B	605	LEU	4.0
1	A	680	LEU	4.0
1	A	694	PHE	4.0
1	B	694	PHE	3.6
1	B	52	LYS	3.5
1	B	457	GLY	3.4
1	B	610	LYS	3.3
1	A	16	ASN	3.3
1	B	286	TRP	3.2
1	B	693	ARG	3.2
1	A	453	ALA	3.2
1	B	16	ASN	3.1
1	B	604	ASN	3.1
1	B	292	ASN	3.0
1	B	454	PHE	2.9
1	A	641	GLY	2.9

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Mol	Chain	Res	Type	RSRZ
1	B	290	ASN	2.9
1	A	14	VAL	2.7
1	B	55	MET	2.6
1	B	288	ASN	2.6
1	B	681	SER	2.6
1	B	517	GLU	2.6
1	B	682	ALA	2.6
1	A	423[A]	MET	2.5
1	A	455	LYS	2.5
1	B	114	GLY	2.5
1	B	644	ILE	2.4
1	B	715	LYS	2.4
1	B	520	ASN	2.4
1	A	715	LYS	2.4
1	A	632	GLU	2.4
1	B	688	PRO	2.4
1	B	453	ALA	2.3
1	A	452	LYS	2.3
1	A	23	ALA	2.3
1	A	708	GLN	2.3
1	A	517	GLU	2.3
1	B	32	GLU	2.3
1	A	371	THR	2.2
1	A	688	PRO	2.2
1	A	173	ALA	2.2
1	A	451	LEU	2.1
1	A	595	MET	2.1
1	A	45	LEU	2.1
1	A	289	PRO	2.1
1	B	289	PRO	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

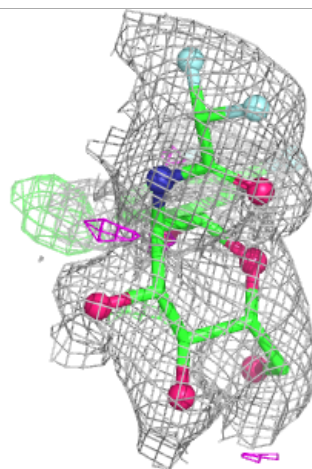
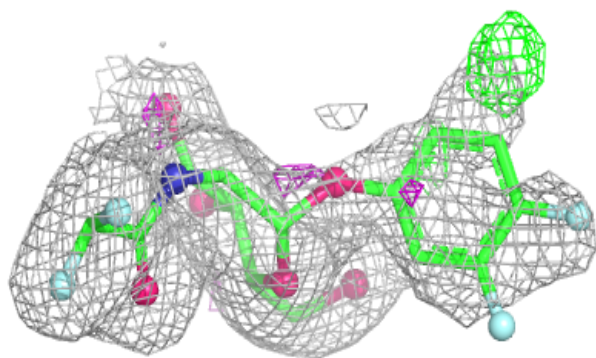
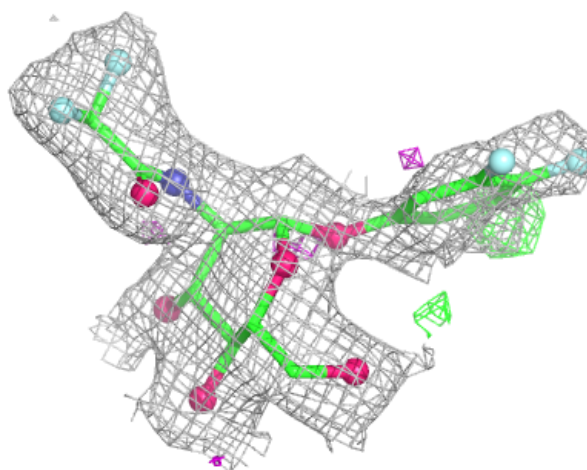
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	GOL	B	1717	6/6	0.81	0.15	63,66,66,67	0
3	GOL	A	1717	6/6	0.86	0.15	52,57,58,60	0
3	GOL	B	1720	6/6	0.86	0.17	47,49,50,52	0
3	GOL	B	1718	6/6	0.88	0.17	43,45,49,50	0
3	GOL	A	1716	6/6	0.88	0.13	40,47,49,52	0
3	GOL	B	1719	6/6	0.91	0.18	47,53,53,57	0
2	14T	B	1000	25/25	0.93	0.11	28,34,57,61	9
2	14T	A	1000	25/25	0.94	0.10	29,37,57,59	9
4	CA	B	1716	1/1	0.99	0.03	30,30,30,30	1

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.

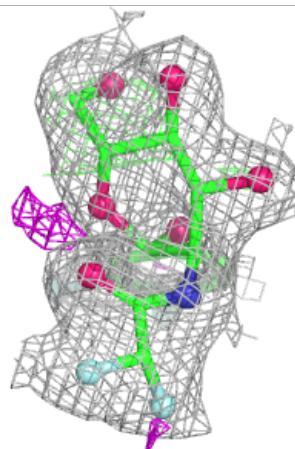
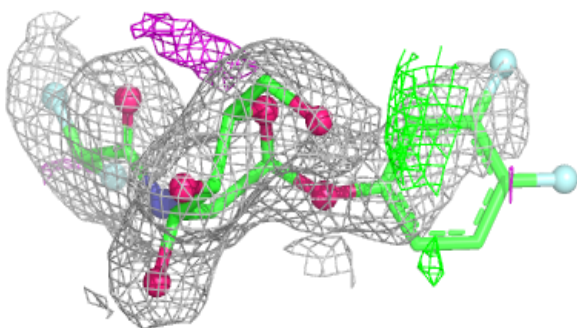
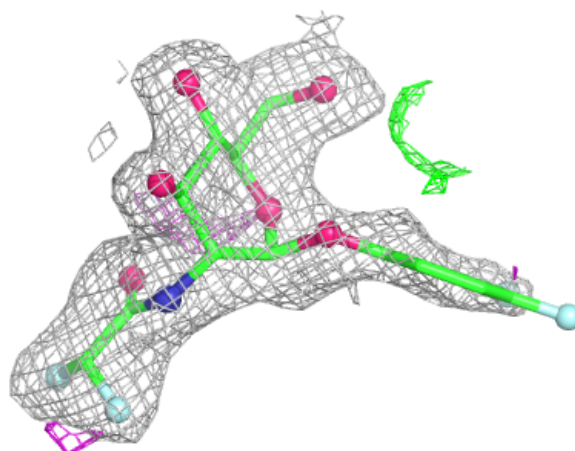
Electron density around 14T B 1000:

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



Electron density around 14T A 1000:

$2mF_o-DF_c$ (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.