



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 11:58 AM UTC

PDB ID : 7C2N / pdb_00007c2n
Title : Crystal structure of mycolic acid transporter MmpL3 from Mycobacterium smegmatis complexed with SPIRO
Authors : Zhang, B.; Yang, X.; Hu, T.; Rao, Z.
Deposited on : 2020-05-08
Resolution : 2.82 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

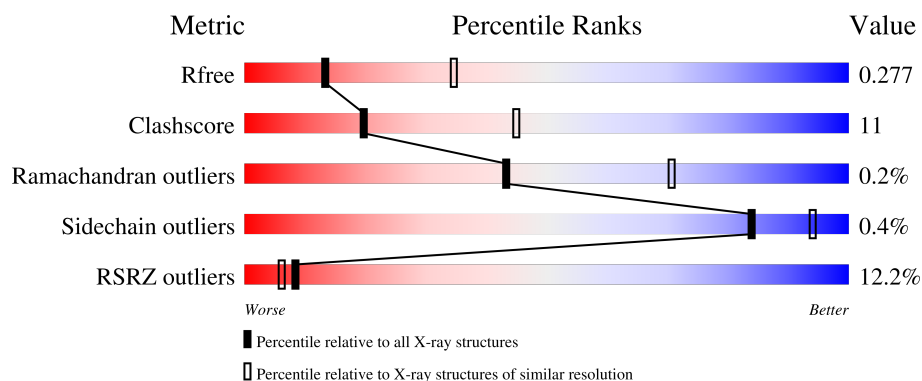
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.82 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	4591 (2.84-2.80)
Clashscore	190562	5010 (2.84-2.80)
Ramachandran outliers	187476	4916 (2.84-2.80)
Sidechain outliers	187428	4918 (2.84-2.80)
RSRZ outliers	180081	4594 (2.84-2.80)

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	934	12% 75% 20% ...

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7178 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 Drug exporters of the RND superfamily-like protein, Endolysin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	900	Total	C	N	O	S	0	0	0
			6952	4482	1182	1256	32			

There are 28 discrepancies between the modelled and reference sequences:

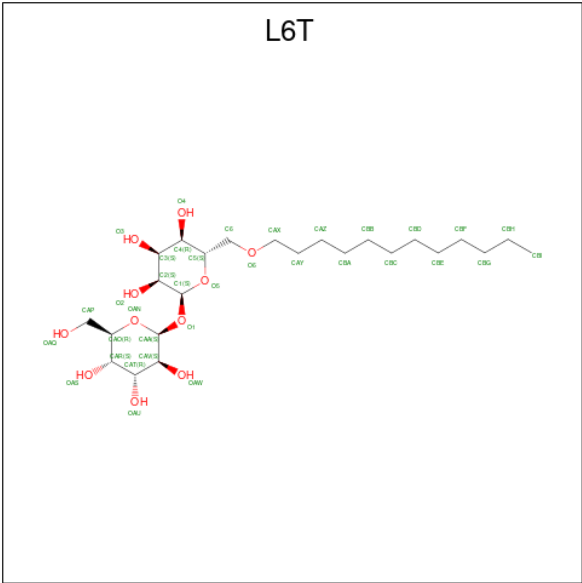
Chain	Residue	Modelled	Actual	Comment	Reference
A	-4	PHE	-	expression tag	UNP I7G2R2
A	-3	GLN	-	expression tag	UNP I7G2R2
A	-2	SER	-	expression tag	UNP I7G2R2
A	-1	ASN	-	expression tag	UNP I7G2R2
A	0	ALA	-	expression tag	UNP I7G2R2
A	749	GLU	-	linker	UNP I7G2R2
A	750	PHE	-	linker	UNP I7G2R2
A	803	THR	CYS	engineered mutation	UNP A0A097J809
A	846	ALA	CYS	engineered mutation	UNP A0A097J809
A	911	GLU	-	expression tag	UNP A0A097J809
A	912	PHE	-	expression tag	UNP A0A097J809
A	913	HIS	-	expression tag	UNP A0A097J809
A	914	LEU	-	expression tag	UNP A0A097J809
A	915	GLY	-	expression tag	UNP A0A097J809
A	916	GLY	-	expression tag	UNP A0A097J809
A	917	ILE	-	expression tag	UNP A0A097J809
A	918	LYS	-	expression tag	UNP A0A097J809
A	919	ALA	-	expression tag	UNP A0A097J809
A	920	PHE	-	expression tag	UNP A0A097J809
A	921	HIS	-	expression tag	UNP A0A097J809
A	922	HIS	-	expression tag	UNP A0A097J809
A	923	HIS	-	expression tag	UNP A0A097J809
A	924	HIS	-	expression tag	UNP A0A097J809
A	925	HIS	-	expression tag	UNP A0A097J809
A	926	HIS	-	expression tag	UNP A0A097J809
A	927	HIS	-	expression tag	UNP A0A097J809
A	928	HIS	-	expression tag	UNP A0A097J809

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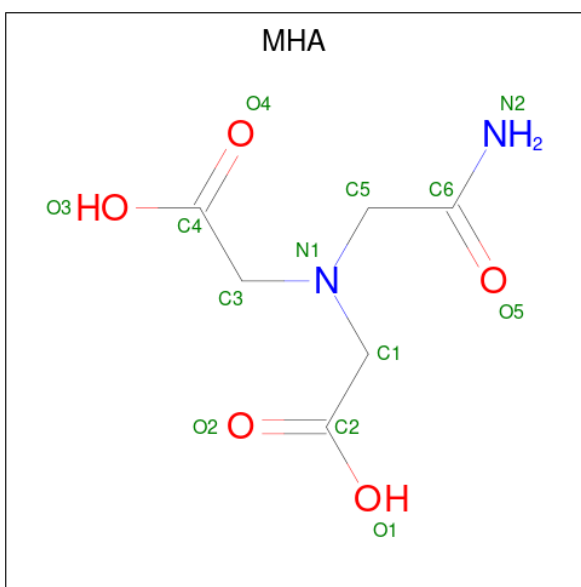
Chain	Residue	Modelled	Actual	Comment	Reference
A	929	HIS	-	expression tag	UNP A0A097J809

- Molecule 2 is alpha-D-glucopyranosyl 6-O-dodecyl-alpha-D-glucopyranoside (CCD ID: L6T) (formula: C₂₄H₄₆O₁₁).



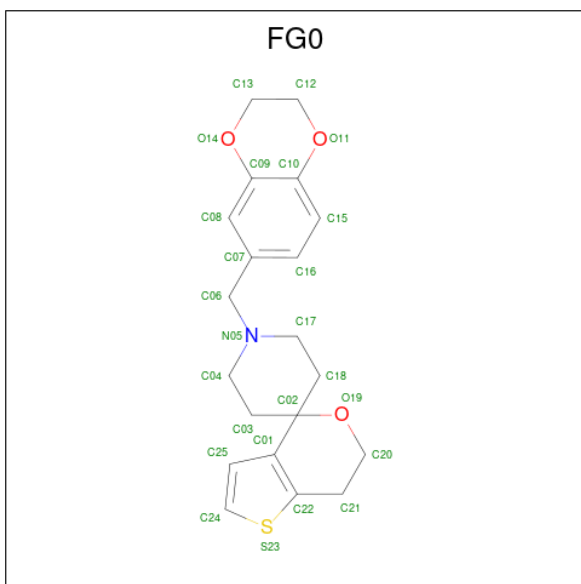
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			35	24	11		
2	A	1	Total	C	O	0	0
			35	24	11		
2	A	1	Total	C	O	0	0
			35	24	11		
2	A	1	Total	C	O	0	0
			35	24	11		
2	A	1	Total	C	O	0	0
			35	24	11		

- Molecule 3 is (CARBAMOYLMETHYL-CARBOXYMETHYL-AMINO)-ACETIC ACID (CCD ID: MHA) (formula: C₆H₁₀N₂O₅).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			13	6	2	5		
3	A	1	Total	C	N	O	0	0
			13	6	2	5		

- Molecule 4 is 1'-(2,3-dihydro-1,4-benzodioxin-6-ylmethyl)spiro[6,7-dihydrothieno[3,2-c]pyran-4,4'-piperidine] (CCD ID: FG0) (formula: C₂₀H₂₃NO₃S).

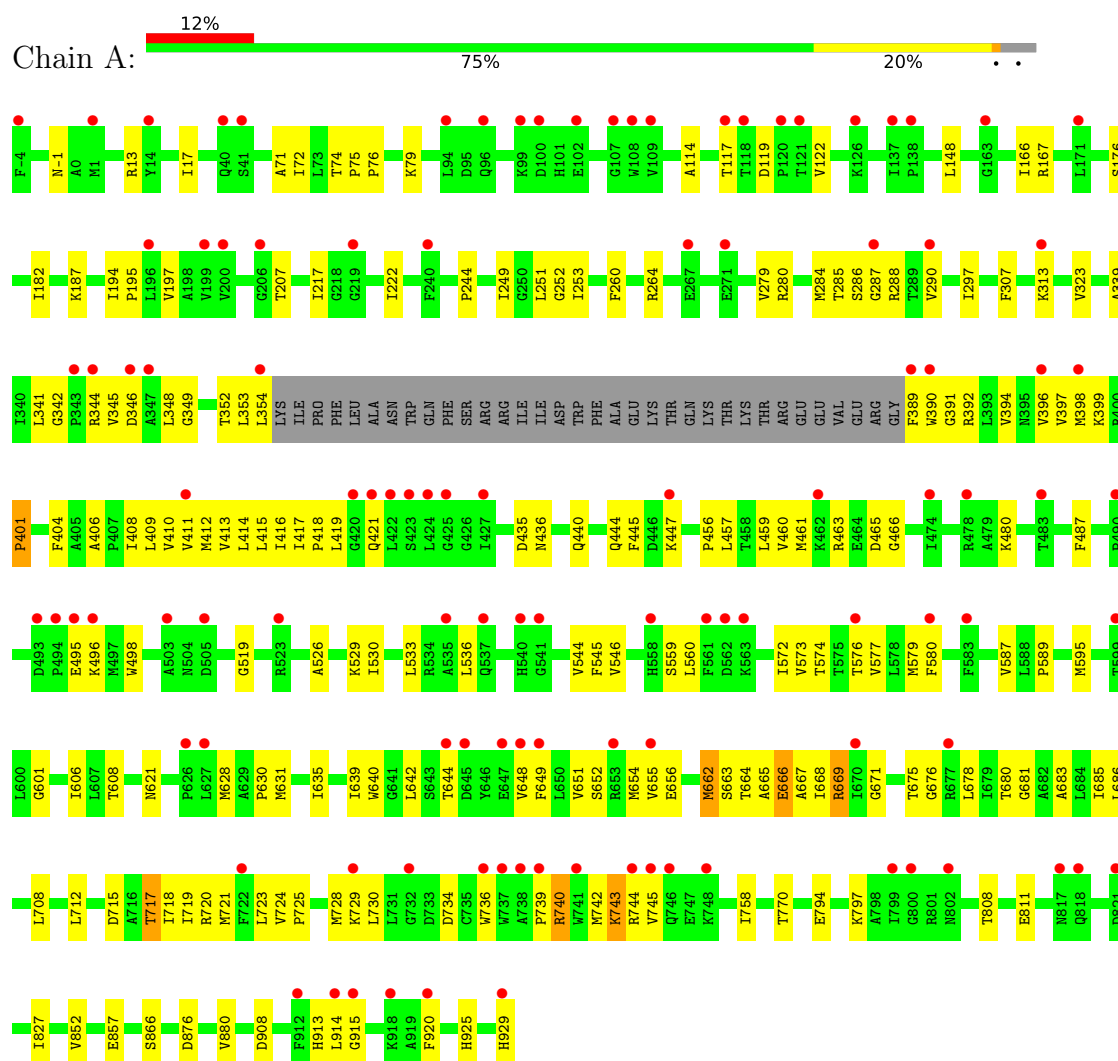


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	S	0	0
			25	20	1	3	1		

3 Residue-property plots

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: Drug exporters of the RND superfamily-like protein, Endolysin



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	86.49Å 140.39Å 145.23Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	44.54 – 2.82 44.54 – 2.82	Depositor EDS
% Data completeness (in resolution range)	99.4 (44.54-2.82) 99.6 (44.54-2.82)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.37 (at 2.81Å)	Xtriage
Refinement program	PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.232 , 0.273 0.238 , 0.277	Depositor DCC
R_{free} test set	2002 reflections (4.60%)	wwPDB-VP
Wilson B-factor (Å ²)	80.6	Xtriage
Anisotropy	0.392	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 42.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.002 for -h,l,k	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	7178	wwPDB-VP
Average B, all atoms (Å ²)	88.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.76% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: FG0, MHA, L6T

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	0.22	0/7097	0.48	3/9642 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	666	GLU	N-CA-C	-6.09	105.70	113.01
1	A	740	ARG	N-CA-C	5.58	116.26	108.00
1	A	717	THR	N-CA-C	5.19	121.86	110.80

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	6952	0	7111	163	0
2	A	175	0	0	2	0
3	A	26	0	16	2	0
4	A	25	0	0	1	0
All	All	7178	0	7127	164	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 11.

All (164) 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:662:MET:HE3	1:A:666:GLU:CB	1.98	0.93
1:A:662:MET:HE3	1:A:666:GLU:HB3	1.53	0.88
1:A:743:LYS:HZ2	1:A:745:VAL:HG13	1.44	0.81
1:A:207:THR:HG21	1:A:346:ASP:HA	1.64	0.78
1:A:743:LYS:NZ	1:A:745:VAL:HG13	1.98	0.78
1:A:654:MET:HG2	1:A:671:GLY:HA3	1.66	0.76
1:A:743:LYS:NZ	1:A:745:VAL:CG1	2.48	0.76
1:A:680:THR:HG22	1:A:720:ARG:HH12	1.53	0.74
1:A:914:LEU:HD23	1:A:914:LEU:H	1.52	0.74
1:A:628:MET:HG3	1:A:630:PRO:HD2	1.74	0.70
1:A:663:SER:O	1:A:667:ALA:HB2	1.92	0.68
1:A:410:VAL:O	1:A:414:LEU:HG	1.92	0.68
1:A:680:THR:CG2	1:A:720:ARG:HH12	2.07	0.68
1:A:297:ILE:HG21	1:A:323:VAL:HG21	1.77	0.67
1:A:421:GLN:O	1:A:608:THR:HG21	1.95	0.66
1:A:409:LEU:O	1:A:413:VAL:HG23	1.96	0.65
1:A:739:PRO:O	1:A:742:MET:HG2	1.98	0.64
1:A:676:GLY:O	1:A:680:THR:HG23	1.97	0.64
1:A:676:GLY:HA2	1:A:720:ARG:HH22	1.63	0.64
1:A:652:SER:O	1:A:656:GLU:HG3	1.98	0.64
1:A:680:THR:HG22	1:A:720:ARG:NH1	2.13	0.62
1:A:348:LEU:HD12	1:A:352:THR:HG21	1.81	0.62
1:A:461:MET:HG2	1:A:544:VAL:HG12	1.82	0.61
1:A:194:ILE:HA	1:A:197:VAL:HG22	1.83	0.60
1:A:406:ALA:O	1:A:410:VAL:HG22	2.02	0.60
1:A:662:MET:CE	1:A:666:GLU:HB3	2.28	0.60
1:A:579:MET:HE3	1:A:648:VAL:HG23	1.84	0.60
1:A:717:THR:HB	1:A:721:MET:HE2	1.85	0.58
1:A:635:ILE:HG13	1:A:639:ILE:HD12	1.86	0.57
1:A:876:ASP:O	1:A:880:VAL:HG23	2.04	0.57
1:A:664:THR:HG1	1:A:736:TRP:HZ3	1.53	0.57
1:A:580:PHE:HD1	1:A:589:PRO:HG2	1.70	0.57
1:A:654:MET:HE1	1:A:728:MET:SD	2.44	0.57
1:A:413:VAL:HA	1:A:416:ILE:HD12	1.87	0.56
1:A:572:ILE:O	1:A:576:THR:HG23	2.06	0.56
1:A:743:LYS:NZ	1:A:745:VAL:HG12	2.21	0.56
1:A:671:GLY:O	1:A:675:THR:OG1	2.20	0.56
1:A:465:ASP:OD1	1:A:466:GLY:N	2.36	0.56
1:A:743:LYS:HZ2	1:A:745:VAL:CG1	2.11	0.56
1:A:929:HIS:ND1	1:A:929:HIS:O	2.38	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:640:TRP:O	1:A:644:THR:HG23	2.06	0.56
1:A:389:PHE:HD2	1:A:392:ARG:HD2	1.71	0.56
1:A:743:LYS:O	1:A:743:LYS:HD3	2.05	0.56
1:A:398:MET:O	1:A:729:LYS:HD2	2.06	0.55
1:A:642:LEU:HA	4:A:1008:FG0:C08	2.36	0.55
1:A:664:THR:O	1:A:667:ALA:HB3	2.07	0.55
1:A:676:GLY:HA2	1:A:720:ARG:NH2	2.21	0.55
1:A:739:PRO:O	1:A:740:ARG:HG2	2.07	0.54
1:A:409:LEU:HD21	1:A:723:LEU:HD11	1.89	0.54
1:A:683:ALA:HB1	1:A:712:LEU:HD11	1.90	0.54
1:A:260:PHE:HA	1:A:678:LEU:HD21	1.91	0.53
1:A:286:SER:O	1:A:290:VAL:HG13	2.08	0.53
1:A:652:SER:O	1:A:655:VAL:HG12	2.09	0.53
1:A:392:ARG:O	1:A:396:VAL:HG23	2.09	0.52
1:A:628:MET:HB3	1:A:631:MET:HG2	1.91	0.52
1:A:-1:ASN:HB3	1:A:288:ARG:NH1	2.25	0.52
3:A:1006:MHA:HC12	3:A:1006:MHA:N2	2.25	0.52
1:A:71:ALA:HB1	1:A:166:ILE:HD11	1.90	0.52
1:A:573:VAL:O	1:A:577:VAL:HG12	2.10	0.52
1:A:681:GLY:O	1:A:685:ILE:HG13	2.10	0.51
1:A:914:LEU:HD23	1:A:914:LEU:N	2.24	0.51
1:A:435:ASP:HA	1:A:440:GLN:HE21	1.73	0.51
1:A:743:LYS:HZ1	1:A:745:VAL:CG1	2.24	0.51
1:A:114:ALA:O	1:A:117:THR:OG1	2.27	0.51
1:A:739:PRO:O	1:A:740:ARG:CG	2.59	0.51
1:A:264:ARG:HD2	1:A:286:SER:OG	2.11	0.51
1:A:401:PRO:CB	1:A:730:LEU:HD11	2.40	0.51
1:A:342:GLY:O	1:A:345:VAL:HG23	2.10	0.51
1:A:394:VAL:O	1:A:398:MET:HG3	2.11	0.51
1:A:353:LEU:C	1:A:354:LEU:HD12	2.35	0.51
1:A:487:PHE:CZ	1:A:533:LEU:HD11	2.45	0.51
1:A:715:ASP:HA	1:A:719:ILE:HG12	1.93	0.51
1:A:715:ASP:O	1:A:720:ARG:HB2	2.11	0.50
1:A:914:LEU:H	1:A:914:LEU:CD2	2.23	0.50
1:A:284:MET:O	1:A:288:ARG:HD2	2.11	0.50
1:A:119:ASP:HB3	1:A:122:VAL:HB	1.93	0.50
1:A:408:ILE:O	1:A:411:VAL:HG22	2.12	0.50
1:A:857:GLU:OE1	1:A:925:HIS:NE2	2.37	0.50
1:A:415:LEU:HD13	1:A:606:ILE:HD11	1.93	0.50
1:A:285:THR:O	1:A:288:ARG:HG2	2.12	0.49
1:A:394:VAL:HG23	1:A:669:ARG:HG3	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:397:VAL:HG12	1:A:404:PHE:CE2	2.48	0.49
1:A:394:VAL:CG2	1:A:669:ARG:HG3	2.43	0.49
1:A:662:MET:HE3	1:A:666:GLU:CG	2.43	0.49
1:A:435:ASP:HA	1:A:440:GLN:NE2	2.28	0.48
1:A:76:PRO:O	1:A:79:LYS:HG3	2.12	0.48
1:A:526:ALA:O	1:A:530:ILE:HD12	2.13	0.48
1:A:665:ALA:C	1:A:667:ALA:N	2.71	0.48
1:A:739:PRO:C	1:A:740:ARG:HG2	2.39	0.48
1:A:580:PHE:CD1	1:A:589:PRO:HG2	2.49	0.47
1:A:663:SER:O	1:A:667:ALA:CB	2.61	0.47
1:A:717:THR:OG1	1:A:718:ILE:N	2.48	0.47
1:A:13:ARG:O	1:A:17:ILE:HG22	2.14	0.47
1:A:222:ILE:HD11	1:A:251:LEU:HB2	1.97	0.47
1:A:417:ILE:HG13	1:A:418:PRO:HD3	1.97	0.47
1:A:459:LEU:HD23	1:A:498:TRP:CZ3	2.49	0.47
1:A:530:ILE:HG23	1:A:546:VAL:HG11	1.95	0.47
1:A:197:VAL:HB	1:A:685:ILE:HG23	1.97	0.47
1:A:770:THR:HG23	1:A:920:PHE:O	2.15	0.47
1:A:794:GLU:HG2	1:A:797:LYS:HE2	1.97	0.47
1:A:724:VAL:O	1:A:728:MET:HE2	2.16	0.46
1:A:398:MET:HE1	1:A:665:ALA:HA	1.97	0.46
1:A:389:PHE:CD2	1:A:392:ARG:HD2	2.49	0.46
1:A:307:PHE:O	1:A:313:LYS:HE3	2.16	0.46
1:A:341:LEU:O	1:A:344:ARG:HD3	2.15	0.46
1:A:463:ARG:HG2	1:A:465:ASP:OD1	2.16	0.46
1:A:399:LYS:C	1:A:401:PRO:HD3	2.40	0.46
1:A:436:ASN:O	1:A:440:GLN:HG3	2.15	0.46
1:A:719:ILE:O	1:A:724:VAL:HG13	2.15	0.46
1:A:416:ILE:HG12	1:A:601:GLY:C	2.40	0.46
1:A:389:PHE:HB3	1:A:392:ARG:HD2	1.97	0.46
1:A:456:PRO:HA	1:A:519:GLY:HA2	1.98	0.46
1:A:655:VAL:HG23	1:A:736:TRP:NE1	2.31	0.46
1:A:391:GLY:HA2	1:A:394:VAL:HG22	1.98	0.46
1:A:487:PHE:HZ	1:A:533:LEU:HD11	1.80	0.46
1:A:587:VAL:HG11	1:A:734:ASP:O	2.16	0.46
1:A:866:SER:OG	3:A:1007:MHA:N2	2.49	0.45
1:A:-1:ASN:HB3	1:A:288:ARG:HH12	1.80	0.45
1:A:194:ILE:N	1:A:195:PRO:HD2	2.32	0.45
1:A:287:GLY:O	1:A:290:VAL:HG22	2.16	0.45
1:A:148:LEU:HD11	1:A:176:SER:HB3	1.99	0.45
1:A:649:PHE:O	1:A:652:SER:OG	2.34	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:167:ARG:NH1	1:A:444:GLN:OE1	2.49	0.44
1:A:252:GLY:HA3	1:A:686:LEU:HD12	1.99	0.44
1:A:412:MET:O	1:A:416:ILE:HG13	2.17	0.44
1:A:621:ASN:OD1	2:A:1005:L6T:OAQ	2.35	0.44
1:A:280:ARG:NH1	1:A:908:ASP:OD1	2.51	0.44
1:A:187:LYS:HB3	1:A:187:LYS:HE3	1.74	0.44
1:A:559:SER:OG	1:A:560:LEU:N	2.51	0.44
1:A:914:LEU:HG	1:A:914:LEU:O	2.16	0.44
1:A:415:LEU:O	1:A:418:PRO:HD2	2.17	0.43
1:A:398:MET:HE3	1:A:668:ILE:HD12	2.00	0.43
1:A:249:ILE:O	1:A:253:ILE:HG13	2.17	0.43
1:A:279:VAL:HG21	1:A:339:ALA:HB2	2.00	0.43
1:A:495:GLU:HB3	1:A:496:LYS:HD3	2.01	0.43
1:A:595:MET:HE2	1:A:595:MET:HB3	1.77	0.43
1:A:480:LYS:O	1:A:536:LEU:HD21	2.19	0.43
1:A:651:VAL:HB	1:A:654:MET:HE2	2.00	0.43
1:A:758:ILE:HG23	1:A:913:HIS:CE1	2.54	0.43
1:A:529:LYS:HD3	1:A:529:LYS:HA	1.73	0.43
1:A:72:ILE:HD11	1:A:445:PHE:HB2	2.01	0.43
1:A:664:THR:OG1	1:A:736:TRP:HZ3	2.02	0.42
1:A:390:TRP:O	1:A:394:VAL:HG22	2.18	0.42
1:A:457:LEU:HD13	1:A:530:ILE:HG13	2.01	0.42
1:A:574:THR:HA	1:A:577:VAL:HG12	2.01	0.42
1:A:495:GLU:HB3	1:A:496:LYS:CD	2.50	0.42
1:A:182:ILE:HG23	1:A:244:PRO:HD3	2.02	0.42
1:A:744:ARG:NH2	2:A:1003:L6T:O2	2.53	0.42
1:A:447:LYS:HE3	1:A:447:LYS:HB2	1.78	0.42
1:A:595:MET:HE3	1:A:723:LEU:HD23	2.02	0.42
1:A:719:ILE:HG13	1:A:720:ARG:N	2.35	0.41
1:A:708:LEU:HD23	1:A:708:LEU:HA	1.79	0.41
1:A:414:LEU:O	1:A:417:ILE:HG12	2.19	0.41
1:A:460:VAL:CG1	1:A:545:PHE:HB2	2.51	0.41
1:A:416:ILE:O	1:A:419:LEU:HB2	2.19	0.41
1:A:715:ASP:HA	1:A:719:ILE:CG1	2.50	0.41
1:A:349:GLY:O	1:A:352:THR:OG1	2.35	0.41
1:A:808:THR:HG23	1:A:811:GLU:H	1.85	0.41
1:A:74:THR:HA	1:A:75:PRO:HD3	1.96	0.41
1:A:827:ILE:HD11	1:A:852:VAL:HG21	2.03	0.41
1:A:217:ILE:HD13	1:A:217:ILE:HA	1.93	0.41
1:A:723:LEU:HD12	1:A:723:LEU:HA	1.75	0.40
1:A:724:VAL:HG22	1:A:725:PRO:HD3	2.04	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:461:MET:HG2	1:A:544:VAL:CG1	2.51	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	896/934 (96%)	868 (97%)	26 (3%)	2 (0%)	43 71

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	915	GLY
1	A	401	PRO

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	738/769 (96%)	735 (100%)	3 (0%)	84 94

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	662	MET

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Mol	Chain	Res	Type
1	A	669	ARG
1	A	743	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	-3	GLN
1	A	-1	ASN
1	A	87	GLN
1	A	159	GLN
1	A	395	ASN
1	A	554	GLN
1	A	621	ASN

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

8 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	L6T	A	1002	-	36,36,36	1.46	7 (19%)	47,47,47	0.92	1 (2%)
3	MHA	A	1006	-	12,12,12	1.55	1 (8%)	15,15,15	1.40	2 (13%)
2	L6T	A	1001	-	36,36,36	1.51	7 (19%)	47,47,47	0.79	1 (2%)
3	MHA	A	1007	-	12,12,12	1.53	1 (8%)	15,15,15	1.44	2 (13%)
2	L6T	A	1005	-	36,36,36	1.46	7 (19%)	47,47,47	1.18	6 (12%)
4	FG0	A	1008	-	27,29,29	3.87	10 (37%)	35,42,42	2.20	11 (31%)
2	L6T	A	1004	-	36,36,36	1.44	7 (19%)	47,47,47	1.18	2 (4%)
2	L6T	A	1003	-	36,36,36	1.49	7 (19%)	47,47,47	1.07	3 (6%)

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	L6T	A	1002	-	-	14/20/60/60	0/2/2/2
3	MHA	A	1006	-	-	9/12/12/12	-
2	L6T	A	1001	-	-	9/20/60/60	0/2/2/2
3	MHA	A	1007	-	-	2/12/12/12	-
2	L6T	A	1005	-	-	12/20/60/60	0/2/2/2
4	FG0	A	1008	-	-	1/4/36/36	0/5/5/5
2	L6T	A	1004	-	-	8/20/60/60	0/2/2/2
2	L6T	A	1003	-	-	9/20/60/60	0/2/2/2

All (47) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1008	FG0	C06-N05	-14.28	1.20	1.47
4	A	1008	FG0	C18-C02	-6.28	1.45	1.53
4	A	1008	FG0	C25-C01	6.21	1.55	1.43
4	A	1008	FG0	C03-C02	-5.93	1.45	1.53
4	A	1008	FG0	O14-C09	5.40	1.43	1.37
2	A	1003	L6T	O5-C5	4.05	1.54	1.44
3	A	1006	MHA	C6-N2	4.01	1.45	1.32
2	A	1001	L6T	O5-C5	3.99	1.54	1.44
3	A	1007	MHA	C6-N2	3.97	1.45	1.32
2	A	1005	L6T	O5-C5	3.93	1.54	1.44
2	A	1002	L6T	O5-C5	3.88	1.53	1.44
2	A	1004	L6T	O5-C5	3.59	1.53	1.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1008	FG0	C22-S23	-3.42	1.63	1.72
2	A	1001	L6T	OAN-CAA	3.32	1.50	1.41
2	A	1004	L6T	OAN-CAA	3.18	1.50	1.41
2	A	1005	L6T	OAN-CAA	3.08	1.49	1.41
4	A	1008	FG0	C08-C09	3.03	1.44	1.38
2	A	1002	L6T	OAN-CAA	3.02	1.49	1.41
2	A	1003	L6T	OAN-CAA	3.00	1.49	1.41
2	A	1004	L6T	C6-C5	-2.96	1.42	1.51
4	A	1008	FG0	O11-C10	2.91	1.40	1.37
2	A	1001	L6T	C6-C5	-2.83	1.43	1.51
2	A	1003	L6T	C6-C5	-2.71	1.43	1.51
2	A	1002	L6T	C6-C5	-2.71	1.43	1.51
2	A	1005	L6T	C6-C5	-2.67	1.43	1.51
4	A	1008	FG0	C24-S23	-2.59	1.58	1.70
2	A	1003	L6T	O5-C1	2.51	1.48	1.41
2	A	1002	L6T	O5-C1	2.44	1.48	1.41
2	A	1003	L6T	C3-C2	-2.42	1.46	1.52
2	A	1002	L6T	C3-C2	-2.40	1.46	1.52
2	A	1001	L6T	O5-C1	2.39	1.48	1.41
2	A	1001	L6T	C3-C2	-2.39	1.46	1.52
2	A	1001	L6T	O3-C3	2.37	1.48	1.43
2	A	1005	L6T	C3-C2	-2.36	1.46	1.52
2	A	1004	L6T	O5-C1	2.30	1.47	1.41
2	A	1004	L6T	O3-C3	2.29	1.48	1.43
2	A	1005	L6T	O3-C3	2.28	1.48	1.43
2	A	1003	L6T	O3-C3	2.27	1.48	1.43
2	A	1005	L6T	O5-C1	2.24	1.47	1.41
2	A	1004	L6T	C3-C2	-2.23	1.46	1.52
4	A	1008	FG0	C09-C10	-2.22	1.36	1.40
2	A	1003	L6T	O2-C2	2.21	1.48	1.43
2	A	1002	L6T	O3-C3	2.20	1.48	1.43
2	A	1004	L6T	O2-C2	2.16	1.48	1.43
2	A	1005	L6T	O2-C2	2.15	1.48	1.43
2	A	1001	L6T	O2-C2	2.11	1.48	1.43
2	A	1002	L6T	O2-C2	2.08	1.48	1.43

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1008	FG0	C24-S23-C22	8.68	101.45	92.08
4	A	1008	FG0	C04-C03-C02	3.59	114.61	111.81
4	A	1008	FG0	O19-C02-C01	3.37	115.04	108.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1004	L6T	C1-C2-C3	3.06	116.44	110.01
2	A	1005	L6T	C1-O5-C5	-3.02	107.81	113.72
2	A	1004	L6T	C4-C3-C2	2.88	115.89	110.83
2	A	1003	L6T	CAA-O1-C1	-2.78	109.34	114.33
2	A	1003	L6T	CAA-CAV-CAT	2.70	115.69	110.01
2	A	1005	L6T	O5-C5-C6	2.64	111.93	106.69
4	A	1008	FG0	C18-C17-N05	2.60	113.81	111.16
4	A	1008	FG0	C21-C22-C01	-2.58	120.22	125.30
3	A	1007	MHA	C6-C5-N1	-2.55	109.75	114.42
3	A	1006	MHA	C5-C6-N2	2.51	120.67	115.61
2	A	1001	L6T	OAN-CAO-CAR	2.48	114.17	109.70
3	A	1006	MHA	C3-N1-C1	2.44	117.39	111.66
2	A	1005	L6T	CAT-CAR-CAO	2.44	114.66	110.23
2	A	1005	L6T	O1-C1-O5	-2.40	104.38	110.69
2	A	1003	L6T	CAA-OAN-CAO	-2.38	109.07	113.72
4	A	1008	FG0	C21-C22-S23	2.38	129.81	123.66
4	A	1008	FG0	C17-C18-C02	2.38	113.67	111.81
2	A	1005	L6T	O1-CAA-CAV	2.38	113.94	108.09
4	A	1008	FG0	C03-C04-N05	2.27	113.47	111.16
3	A	1007	MHA	C5-C6-N2	2.27	120.18	115.61
2	A	1002	L6T	OAN-CAO-CAR	2.18	113.63	109.70
4	A	1008	FG0	C25-C24-S23	-2.17	108.98	112.86
4	A	1008	FG0	C06-N05-C17	-2.11	106.52	111.07
4	A	1008	FG0	C12-O11-C10	2.06	117.30	113.67
2	A	1005	L6T	OAN-CAO-CAR	2.05	113.39	109.70

There are no chirality outliers.

All (64) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1002	L6T	C4-C5-C6-O6
2	A	1002	L6T	OAN-CAA-O1-C1
2	A	1005	L6T	CAV-CAA-O1-C1
3	A	1006	MHA	C4-C3-N1-C1
3	A	1006	MHA	C6-C5-N1-C1
3	A	1006	MHA	N1-C3-C4-O3
3	A	1006	MHA	N1-C3-C4-O4
3	A	1006	MHA	N1-C5-C6-O5
3	A	1006	MHA	N1-C5-C6-N2
2	A	1002	L6T	OAN-CAO-CAP-OAQ
2	A	1001	L6T	CAR-CAO-CAP-OAQ
2	A	1005	L6T	CAR-CAO-CAP-OAQ

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Mol	Chain	Res	Type	Atoms
2	A	1002	L6T	O5-C5-C6-O6
2	A	1005	L6T	OAN-CAO-CAP-OAQ
2	A	1003	L6T	OAN-CAO-CAP-OAQ
2	A	1001	L6T	OAN-CAO-CAP-OAQ
2	A	1002	L6T	CAR-CAO-CAP-OAQ
2	A	1005	L6T	CBA-CBB-CBC-CBD
2	A	1004	L6T	CAX-CAY-CAZ-CBA
2	A	1005	L6T	O5-C1-O1-CAA
2	A	1005	L6T	OAN-CAA-O1-C1
2	A	1004	L6T	OAN-CAO-CAP-OAQ
2	A	1003	L6T	O6-CAX-CAY-CAZ
2	A	1002	L6T	CAV-CAA-O1-C1
2	A	1005	L6T	C2-C1-O1-CAA
2	A	1004	L6T	CBB-CBC-CBD-CBE
2	A	1001	L6T	C4-C5-C6-O6
2	A	1003	L6T	CAR-CAO-CAP-OAQ
2	A	1005	L6T	CBB-CBC-CBD-CBE
2	A	1001	L6T	O5-C5-C6-O6
2	A	1003	L6T	CAX-CAY-CAZ-CBA
2	A	1002	L6T	CBC-CBD-CBE-CBF
2	A	1003	L6T	CAZ-CBA-CBB-CBC
3	A	1007	MHA	N1-C3-C4-O4
2	A	1005	L6T	C5-C6-O6-CAX
2	A	1001	L6T	CBA-CBB-CBC-CBD
2	A	1001	L6T	CBF-CBG-CBH-CBI
2	A	1002	L6T	CBA-CBB-CBC-CBD
2	A	1003	L6T	O5-C5-C6-O6
2	A	1004	L6T	CBE-CBF-CBG-CBH
2	A	1003	L6T	C4-C5-C6-O6
2	A	1003	L6T	CBA-CBB-CBC-CBD
2	A	1002	L6T	CBB-CBC-CBD-CBE
2	A	1002	L6T	CBD-CBE-CBF-CBG
3	A	1007	MHA	N1-C3-C4-O3
3	A	1006	MHA	C2-C1-N1-C3
2	A	1003	L6T	C5-C6-O6-CAX
4	A	1008	FG0	C07-C06-N05-C04
2	A	1005	L6T	CBC-CBD-CBE-CBF
2	A	1005	L6T	CAY-CAX-O6-C6
2	A	1002	L6T	CAY-CAZ-CBA-CBB
2	A	1002	L6T	CAY-CAX-O6-C6
2	A	1004	L6T	CBD-CBE-CBF-CBG
2	A	1002	L6T	C5-C6-O6-CAX

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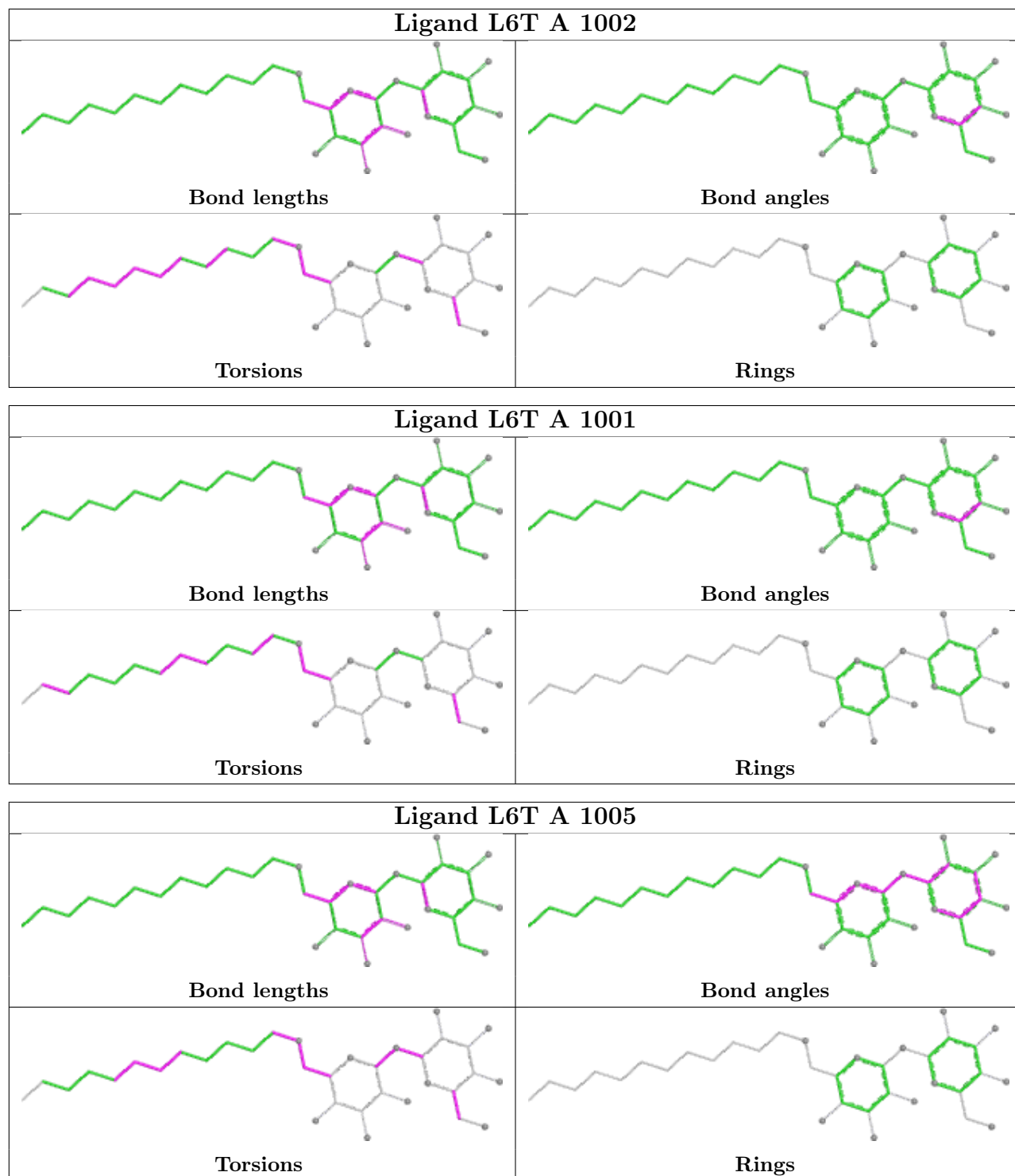
Mol	Chain	Res	Type	Atoms
2	A	1005	L6T	C4-C5-C6-O6
2	A	1001	L6T	C5-C6-O6-CAX
3	A	1006	MHA	C4-C3-N1-C5
2	A	1004	L6T	CAY-CAX-O6-C6
2	A	1001	L6T	O6-CAX-CAY-CAZ
2	A	1004	L6T	C5-C6-O6-CAX
2	A	1001	L6T	CAZ-CBA-CBB-CBC
2	A	1004	L6T	O5-C5-C6-O6
3	A	1006	MHA	N1-C1-C2-O1
2	A	1002	L6T	CBE-CBF-CBG-CBH

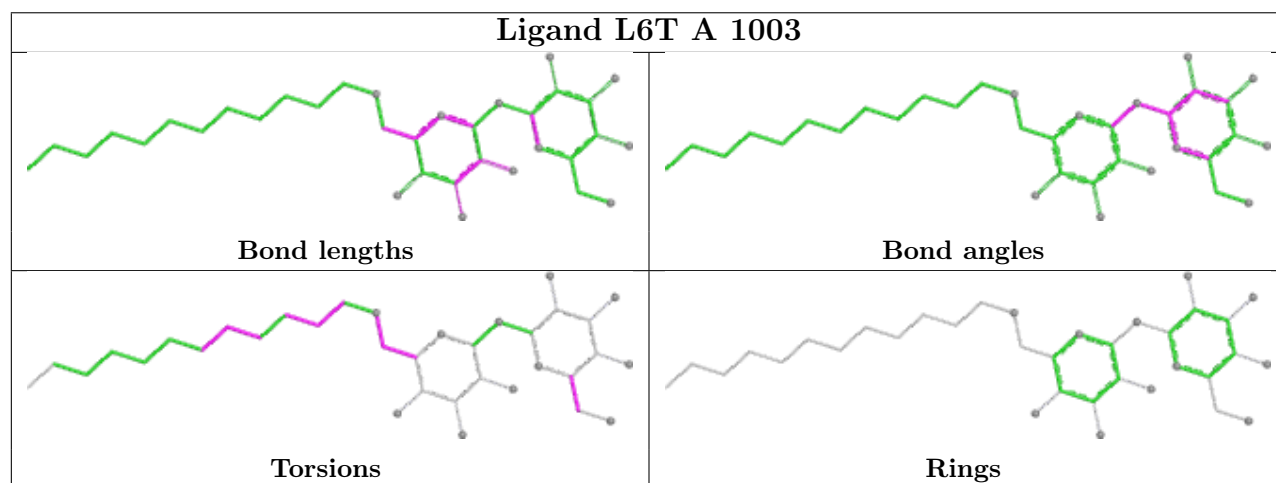
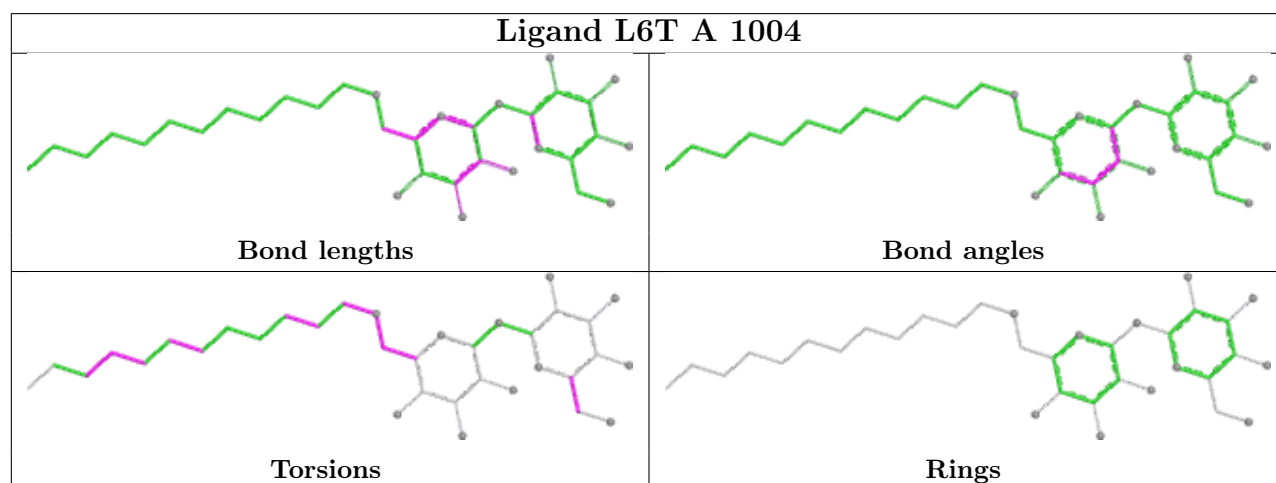
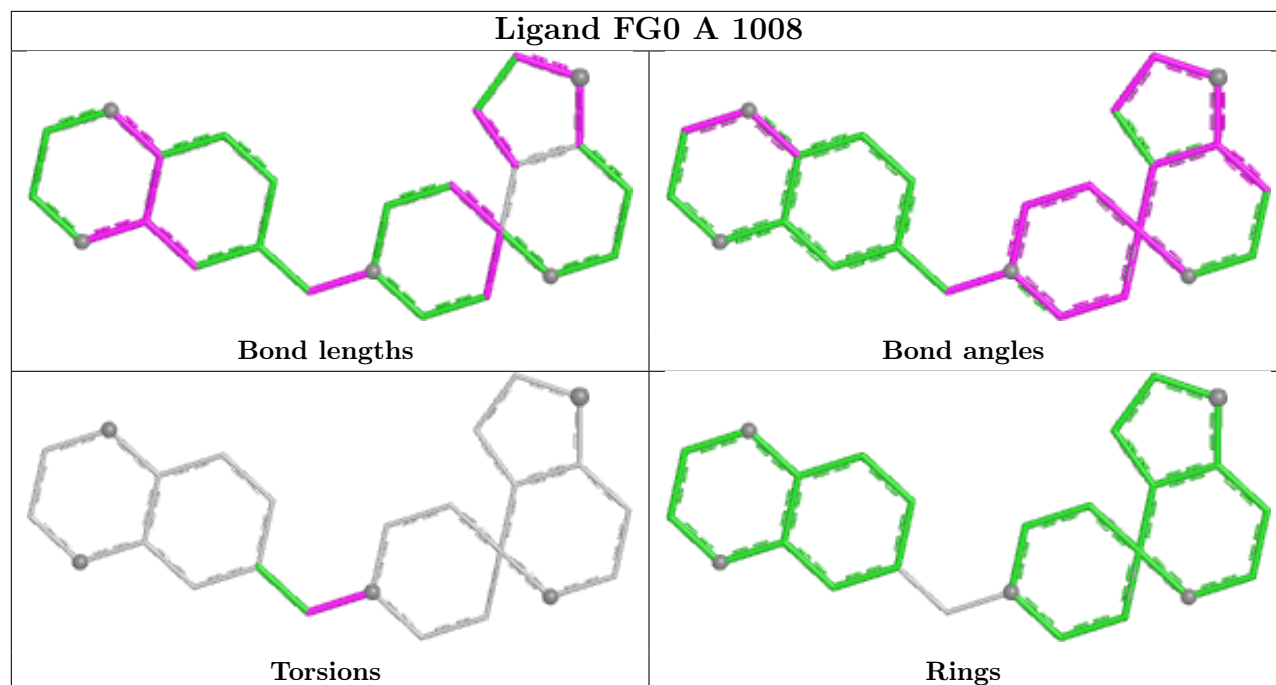
There are no ring outliers.

5 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1006	MHA	1	0
3	A	1007	MHA	1	0
2	A	1005	L6T	1	0
4	A	1008	FG0	1	0
2	A	1003	L6T	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

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, 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	900/934 (96%)	0.76	110 (12%) 8 6	55, 81, 132, 196	0

All (110) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	495	GLU	9.8
1	A	423	SER	7.6
1	A	424	LEU	6.5
1	A	561	PHE	5.7
1	A	494	PRO	5.6
1	A	287	GLY	5.0
1	A	96	GLN	5.0
1	A	914	LEU	4.6
1	A	421	GLN	4.6
1	A	99	LYS	4.6
1	A	493	ASP	4.5
1	A	1	MET	4.4
1	A	163	GLY	4.3
1	A	802	ASN	4.3
1	A	737	TRP	4.3
1	A	108	TRP	4.2
1	A	-4	PHE	4.2
1	A	389	PHE	4.1
1	A	425	GLY	4.1
1	A	627	LEU	4.0
1	A	390	TRP	4.0
1	A	420	GLY	4.0
1	A	496	LYS	4.0
1	A	107	GLY	3.9
1	A	271	GLU	3.9
1	A	347	ALA	3.9
1	A	505	ASP	3.7

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Mol	Chain	Res	Type	RSRZ
1	A	648	VAL	3.5
1	A	653	ARG	3.5
1	A	290	VAL	3.4
1	A	14	TYR	3.4
1	A	137	ILE	3.4
1	A	41	SER	3.3
1	A	344	ARG	3.3
1	A	583	PHE	3.3
1	A	736	TRP	3.3
1	A	818	GLN	3.3
1	A	817	ASN	3.2
1	A	346	ASP	3.2
1	A	929	HIS	3.2
1	A	821	ASP	3.2
1	A	535	ALA	3.1
1	A	267	GLU	3.1
1	A	343	PRO	3.1
1	A	200	VAL	3.0
1	A	745	VAL	3.0
1	A	576	THR	3.0
1	A	918	LYS	3.0
1	A	647	GLU	2.9
1	A	427	ILE	2.9
1	A	40	GLN	2.9
1	A	422	LEU	2.8
1	A	219	GLY	2.8
1	A	599	THR	2.8
1	A	109	VAL	2.7
1	A	563	LYS	2.7
1	A	738	ALA	2.7
1	A	626	PRO	2.6
1	A	354	LEU	2.6
1	A	729	LYS	2.6
1	A	474	ILE	2.6
1	A	799	ILE	2.6
1	A	912	PHE	2.6
1	A	196	LEU	2.5
1	A	411	VAL	2.5
1	A	240	PHE	2.5
1	A	920	PHE	2.5
1	A	120	PRO	2.5
1	A	915	GLY	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	645	ASP	2.5
1	A	199	VAL	2.5
1	A	396	VAL	2.5
1	A	580	PHE	2.5
1	A	746	GLN	2.5
1	A	171	LEU	2.4
1	A	138	PRO	2.4
1	A	126	LYS	2.4
1	A	121	THR	2.4
1	A	655	VAL	2.4
1	A	102	GLU	2.4
1	A	117	THR	2.4
1	A	649	PHE	2.3
1	A	398	MET	2.3
1	A	558	HIS	2.3
1	A	741	TRP	2.3
1	A	483	THR	2.3
1	A	462	LYS	2.3
1	A	677	ARG	2.3
1	A	732	GLY	2.3
1	A	94	LEU	2.2
1	A	100	ASP	2.2
1	A	478	ARG	2.2
1	A	523	ARG	2.2
1	A	206	GLY	2.2
1	A	722	PHE	2.2
1	A	503	ALA	2.2
1	A	562	ASP	2.1
1	A	644	THR	2.1
1	A	313	LYS	2.1
1	A	748	LYS	2.1
1	A	800	GLY	2.1
1	A	540	HIS	2.1
1	A	490	PRO	2.1
1	A	744	ARG	2.1
1	A	537	GLN	2.0
1	A	739	PRO	2.0
1	A	541	GLY	2.0
1	A	670	ILE	2.0
1	A	447	LYS	2.0
1	A	118	THR	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

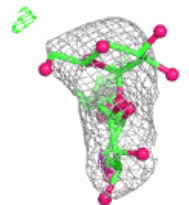
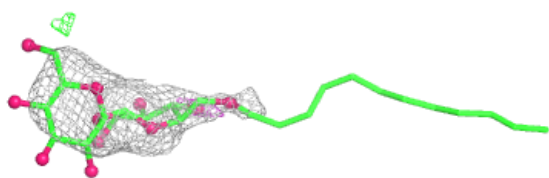
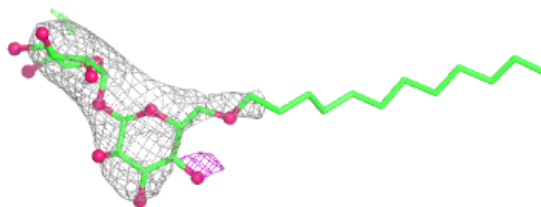
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	L6T	A	1003	35/35	0.77	0.28	92,143,155,165	0
3	MHA	A	1007	13/13	0.81	0.16	70,77,83,90	0
2	L6T	A	1005	35/35	0.82	0.18	66,101,129,131	0
2	L6T	A	1002	35/35	0.85	0.15	57,135,161,168	0
3	MHA	A	1006	13/13	0.87	0.13	66,77,89,92	0
4	FG0	A	1008	25/25	0.89	0.14	82,94,106,108	0
2	L6T	A	1004	35/35	0.90	0.14	75,88,104,120	0
2	L6T	A	1001	35/35	0.91	0.14	59,76,99,101	0

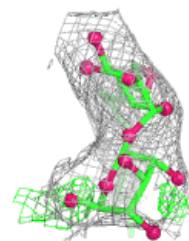
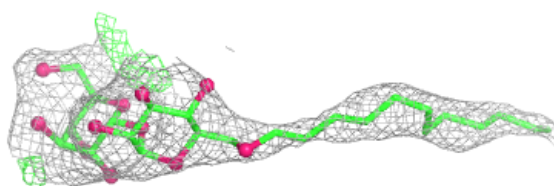
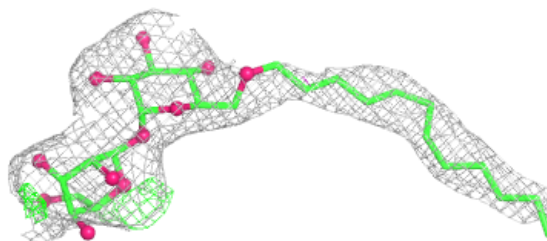
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around L6T A 1003:

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

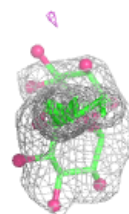
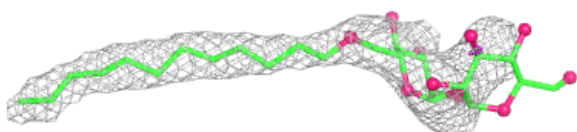
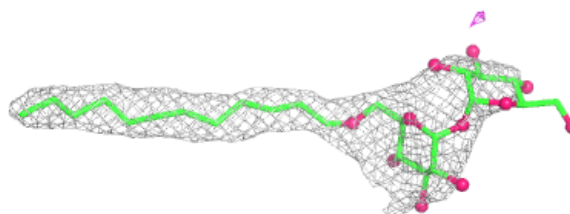
**Electron density around L6T A 1005:**

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

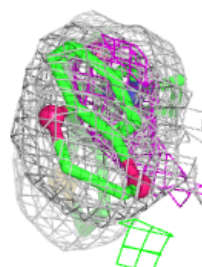
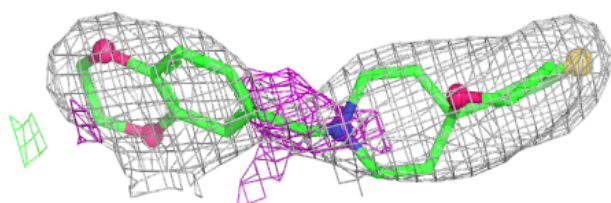
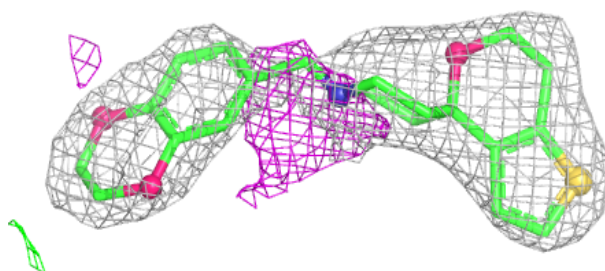


Electron density around L6T A 1002:

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

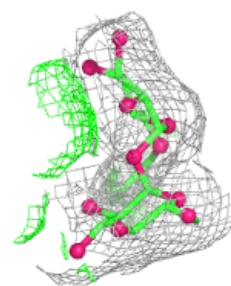
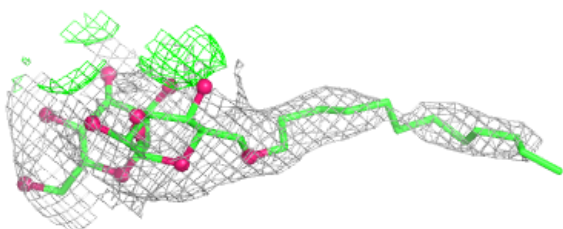
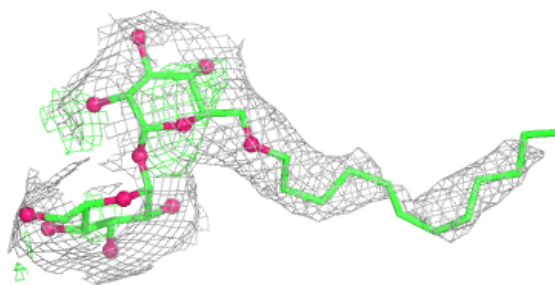
**Electron density around FG0 A 1008:**

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

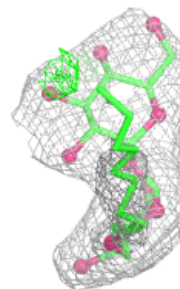
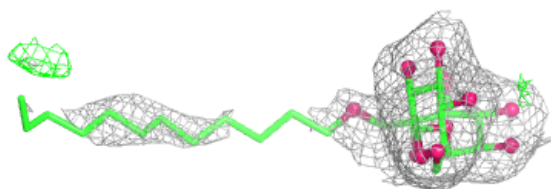
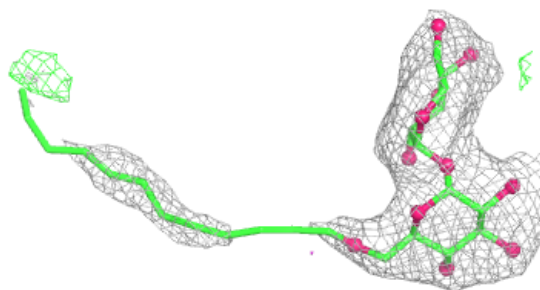


Electron density around L6T A 1004:

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

**Electron density around L6T A 1001:**

$2mF_o-DF_c$ (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.