



Full wwPDB X-ray Structure Validation Report ⓘ

Dec 12, 2023 – 09:25 pm GMT

PDB ID : 4C48
Title : Crystal structure of AcrB-AcrZ complex
Authors : Du, D.; James, N.; Klimont, E.; Luisi, B.F.
Deposited on : 2013-09-02
Resolution : 3.30 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 : 4.02b-467
Mogul : 1.8.4, CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.36
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) : 2.36

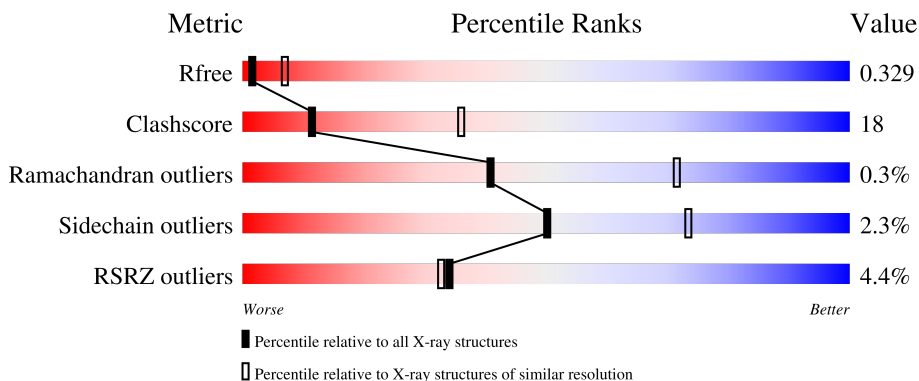
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

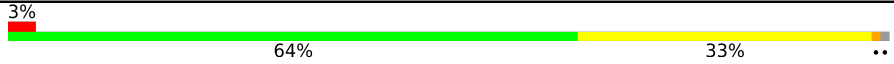
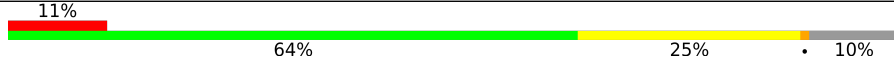

The reported resolution of this entry is 3.30 Å.

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	1149 (3.34-3.26)
Clashscore	141614	1205 (3.34-3.26)
Ramachandran outliers	138981	1183 (3.34-3.26)
Sidechain outliers	138945	1182 (3.34-3.26)
RSRZ outliers	127900	1115 (3.34-3.26)

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	1047	
2	B	169	
3	C	49	

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
4	LMT	A	2042	-	-	-	X
4	LMT	A	2043	-	-	-	X
4	LMT	A	2044	-	-	-	X
4	LMT	A	2046	-	-	-	X
4	LMT	A	2047	-	-	-	X

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 9390 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 ACRIFLAVINE RESISTANCE PROTEIN B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	1032	7788	5013	1285	1446	44	0	0	0

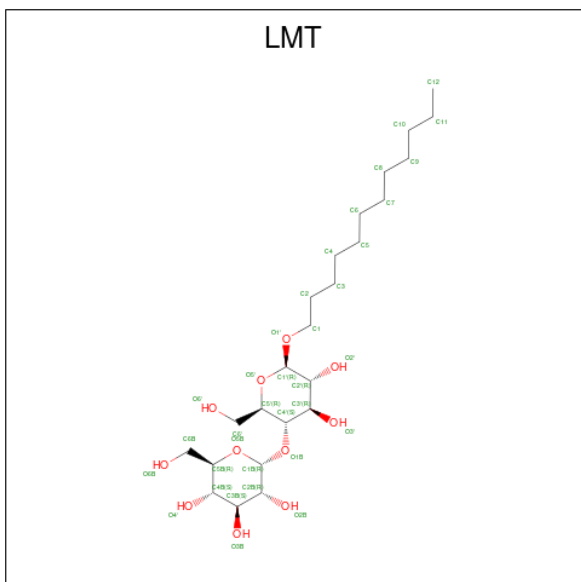
- Molecule 2 is a protein called DARPIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	152	1145	723	199	222	1	0	0	0

- Molecule 3 is a protein called UNCHARACTERIZED MEMBRANE PROTEIN YBHT.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	C	46	326	219	48	56	3	0	0	0

- Molecule 4 is DODECYL-BETA-D-MALTOSE (three-letter code: LMT) (formula: $C_{24}H_{46}O_{11}$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C 10 10	0	0
4	A	1	Total C 6 6	0	0
4	A	1	Total C 8 8	0	0
4	A	1	Total C 12 12	0	0
4	A	1	Total C O 35 24 11	0	0
4	A	1	Total C 8 8	0	0
4	A	1	Total C 4 4	0	0
4	A	1	Total C 6 6	0	0
4	A	1	Total C 6 6	0	0

- Molecule 5 is NICKEL (II) ION (three-letter code: NI) (formula: Ni).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total Ni 1 1	0	0

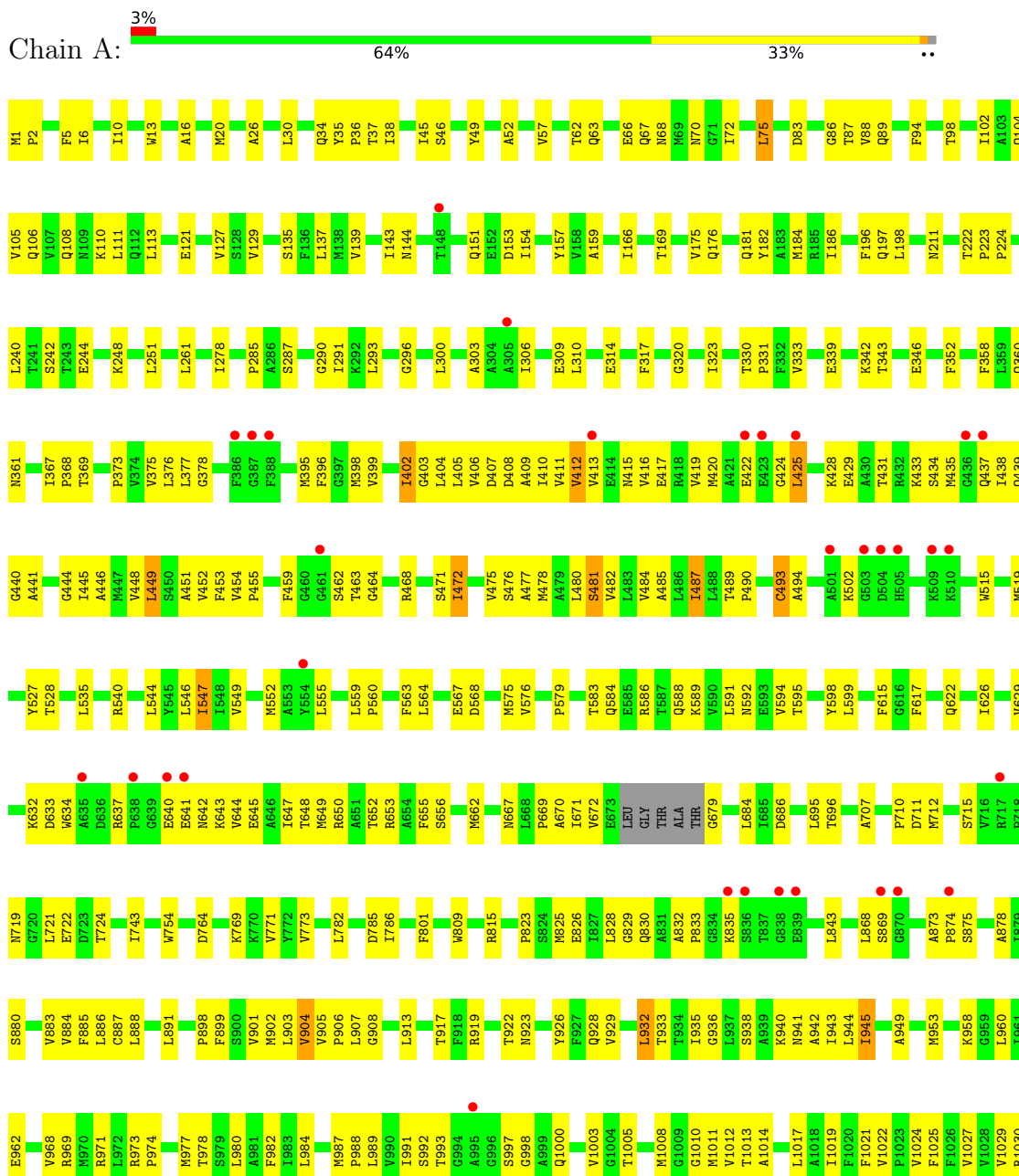
- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	28	Total O 28 28	0	0
6	B	7	Total O 7 7	0	0

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: ACRIFLAVINE RESISTANCE PROTEIN B





- Molecule 2: DARPIN



- Molecule 3: UNCHARACTERIZED MEMBRANE PROTEIN YBHT



4 Data and refinement statistics

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, α , β , γ	145.01Å 145.01Å 538.34Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	29.67 – 3.30 29.66 – 3.30	Depositor EDS
% Data completeness (in resolution range)	99.5 (29.67-3.30) 92.7 (29.66-3.30)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.38 (at 3.31Å)	Xtrriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.281 , 0.322 0.283 , 0.329	Depositor DCC
R_{free} test set	1656 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	80.7	Xtrriage
Anisotropy	0.249	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.26 , 61.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.86	EDS
Total number of atoms	9390	wwPDB-VP
Average B, all atoms (Å ²)	97.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.82% 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: NI, LMT

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	0.58	0/7936	0.70	1/10780 (0.0%)
2	B	0.44	0/1164	0.59	0/1584
3	C	0.50	0/330	0.63	0/448
All	All	0.56	0/9430	0.68	1/12812 (0.0%)

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	2

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	449	LEU	CA-CB-CG	5.27	127.41	115.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	222	THR	Mainchain,Peptide

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	7788	0	7885	278	0
2	B	1145	0	1125	29	0
3	C	326	0	346	23	0
4	A	95	0	137	10	0
5	A	1	0	0	0	0
6	A	28	0	0	1	0
6	B	7	0	0	0	0
All	All	9390	0	9493	333	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 18.

All (333) 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)
3:C:6:LYS:HG3	3:C:10:PHE:CE2	1.51	1.46
3:C:6:LYS:CG	3:C:10:PHE:HE2	1.45	1.28
1:A:575:MET:HE2	1:A:662:MET:CE	1.64	1.24
1:A:575:MET:CE	1:A:662:MET:HE2	1.74	1.16
3:C:6:LYS:CG	3:C:10:PHE:CE2	2.28	1.02
1:A:575:MET:CE	1:A:662:MET:CE	2.37	0.98
2:B:60:LEU:HD21	2:B:94:GLU:HB3	1.48	0.93
4:A:2043:LMT:H3O2	4:A:2043:LMT:H2O1	1.19	0.87
3:C:6:LYS:O	3:C:10:PHE:HD2	1.56	0.87
1:A:70:ASN:O	1:A:110:LYS:NZ	2.08	0.86
4:A:2043:LMT:O2B	4:A:2043:LMT:H4'	1.76	0.86
1:A:960:LEU:HD21	1:A:1030:ARG:HB2	1.57	0.85
1:A:555:LEU:HD22	1:A:913:LEU:HB3	1.62	0.82
3:C:6:LYS:O	3:C:10:PHE:CD2	2.35	0.79
3:C:6:LYS:HG3	3:C:10:PHE:HE2	0.64	0.79
2:B:26:ARG:NH1	2:B:29:GLU:OE1	2.19	0.76
1:A:291:ILE:HD13	1:A:306:ILE:HD13	1.68	0.75
1:A:938:SER:HB3	1:A:1014:ALA:HB1	1.69	0.74
1:A:360:GLN:O	1:A:361:ASN:ND2	2.20	0.73
1:A:901:VAL:HG21	1:A:943:ILE:HG13	1.70	0.72
1:A:575:MET:HE2	1:A:662:MET:HE2	0.81	0.72
1:A:568:ASP:OD1	1:A:634:TRP:NE1	2.23	0.72
2:B:41:ASN:HD21	2:B:72:ASP:H	1.34	0.71
1:A:632:LYS:O	1:A:637:ARG:NH1	2.23	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:422:GLU:OE1	1:A:422:GLU:N	2.24	0.70
1:A:463:THR:HG21	6:A:2019:HOH:O	1.91	0.69
4:A:2043:LMT:O2B	4:A:2043:LMT:O3'	2.01	0.69
1:A:684:LEU:HD22	1:A:695:LEU:HD11	1.75	0.69
1:A:137:LEU:HD21	1:A:303:ALA:HB2	1.75	0.69
1:A:633:ASP:OD1	1:A:634:TRP:N	2.26	0.69
4:A:2043:LMT:O2B	4:A:2043:LMT:C4'	2.41	0.68
1:A:57:VAL:HG21	1:A:86:GLY:HA2	1.76	0.68
1:A:375:VAL:HG11	1:A:405:LEU:HD22	1.75	0.68
1:A:454:VAL:HG12	1:A:475:VAL:HG21	1.76	0.68
4:A:2043:LMT:H1B	4:A:2043:LMT:O6B	1.93	0.67
1:A:721:LEU:HD13	1:A:815:ARG:HD3	1.77	0.67
2:B:89:HIS:HB2	2:B:119:LEU:HD22	1.77	0.66
1:A:712:MET:SD	1:A:835:LYS:HE2	2.36	0.66
1:A:984:LEU:HA	1:A:987:MET:HG2	1.78	0.65
1:A:885:PHE:HA	1:A:902:MET:HE3	1.79	0.64
1:A:404:LEU:HD11	1:A:449:LEU:HD21	1.79	0.64
1:A:453:PHE:HB3	1:A:932:LEU:HD22	1.79	0.63
1:A:598:TYR:HE2	1:A:629:VAL:HG21	1.61	0.63
1:A:899:PHE:O	1:A:903:LEU:HD12	1.97	0.63
1:A:1030:ARG:O	1:A:1034:SER:OG	2.15	0.63
1:A:6:ILE:O	1:A:428:LYS:NZ	2.32	0.63
1:A:424:GLY:HA3	1:A:502:LYS:HB3	1.81	0.63
1:A:898:PRO:O	1:A:901:VAL:HG12	1.99	0.63
1:A:38:ILE:HG22	1:A:462:SER:HB2	1.81	0.62
1:A:248:LYS:HA	1:A:261:LEU:HD23	1.80	0.62
1:A:544:LEU:HD12	1:A:547:ILE:HG12	1.80	0.62
1:A:546:LEU:O	1:A:549:VAL:HG12	1.99	0.62
1:A:898:PRO:O	1:A:902:MET:HG2	2.00	0.62
1:A:446:ALA:HB2	1:A:482:VAL:HG21	1.82	0.62
1:A:485:ALA:O	1:A:490:PRO:HD3	2.00	0.61
2:B:63:VAL:HG11	2:B:99:LEU:HD11	1.82	0.61
2:B:59:HIS:O	2:B:63:VAL:HG23	2.00	0.61
1:A:563:PHE:O	1:A:564:LEU:HD23	2.00	0.61
1:A:451:ALA:HB1	1:A:883:VAL:HG13	1.82	0.60
1:A:453:PHE:CE1	1:A:475:VAL:HG22	2.36	0.60
1:A:20:MET:HA	1:A:377:LEU:HD11	1.84	0.60
1:A:110:LYS:HA	1:A:113:LEU:HG	1.82	0.60
1:A:375:VAL:CG1	1:A:405:LEU:HD22	2.32	0.60
1:A:402:ILE:O	1:A:406:VAL:HG13	2.00	0.60
1:A:153:ASP:OD1	1:A:182:TYR:OH	2.20	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:926:TYR:HB3	1:A:1003:VAL:HG23	1.82	0.59
1:A:884:VAL:HG12	1:A:902:MET:CE	2.32	0.59
2:B:61:GLU:O	2:B:65:VAL:HG23	2.03	0.59
1:A:104:GLN:O	1:A:108:GLN:HG2	2.02	0.59
1:A:139:VAL:HG22	1:A:290:GLY:HA2	1.83	0.59
1:A:102:ILE:HA	1:A:105:VAL:HG22	1.85	0.58
1:A:903:LEU:O	1:A:906:PRO:HD2	2.02	0.58
2:B:82:THR:HG22	2:B:84:LEU:H	1.69	0.57
1:A:407:ASP:HA	1:A:410:ILE:HG12	1.86	0.57
1:A:901:VAL:HG23	1:A:942:ALA:HB3	1.86	0.57
1:A:984:LEU:HD23	1:A:987:MET:HG3	1.85	0.57
1:A:35:TYR:HE2	1:A:672:VAL:H	1.53	0.57
1:A:637:ARG:HB2	1:A:642:ASN:HB3	1.85	0.57
1:A:1021:PHE:HB3	1:A:1025:PHE:CZ	2.39	0.57
4:A:2043:LMT:O6B	4:A:2043:LMT:C1B	2.53	0.57
1:A:454:VAL:N	1:A:455:PRO:HD2	2.20	0.57
2:B:109:LYS:HD2	2:B:113:GLY:HA2	1.85	0.57
1:A:169:THR:OG1	1:A:309:GLU:OE2	2.22	0.57
1:A:330:THR:OG1	1:A:331:PRO:HD3	2.05	0.57
1:A:644:VAL:O	1:A:648:THR:HG23	2.05	0.57
1:A:94:PHE:HB3	1:A:98:THR:HG21	1.87	0.56
1:A:137:LEU:HD13	1:A:293:LEU:HD21	1.88	0.56
1:A:944:LEU:HD12	1:A:971:ARG:HH21	1.71	0.56
1:A:110:LYS:HB2	1:A:113:LEU:HD12	1.88	0.56
1:A:973:ARG:NH1	1:A:977:MET:SD	2.79	0.56
1:A:339:GLU:O	1:A:343:THR:HG23	2.05	0.56
1:A:434:SER:HA	1:A:437:GLN:OE1	2.05	0.56
1:A:884:VAL:HG12	1:A:902:MET:HE1	1.86	0.56
1:A:448:VAL:CG2	1:A:884:VAL:HG13	2.36	0.55
1:A:358:PHE:CD2	1:A:977:MET:HG2	2.41	0.55
1:A:444:GLY:HA3	1:A:891:LEU:HD12	1.89	0.55
1:A:425:LEU:HD12	1:A:429:GLU:HB3	1.87	0.55
1:A:36:PRO:O	1:A:38:ILE:HG13	2.07	0.55
1:A:396:PHE:O	1:A:399:VAL:HG12	2.07	0.55
1:A:721:LEU:HD13	1:A:815:ARG:CD	2.36	0.55
1:A:83:ASP:OD1	1:A:87:THR:OG1	2.25	0.54
1:A:310:LEU:HD21	1:A:323:ILE:HD12	1.89	0.54
1:A:37:THR:HG21	1:A:296:GLY:HA2	1.89	0.54
1:A:415:ASN:ND2	1:A:437:GLN:HE21	2.05	0.54
1:A:653:ARG:O	1:A:656:SER:OG	2.25	0.54
1:A:707:ALA:O	1:A:710:PRO:HD3	2.08	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:2043:LMT:O6'	4:A:2043:LMT:H2B	2.07	0.54
1:A:974:PRO:O	1:A:978:THR:HG23	2.07	0.54
1:A:477:ALA:O	1:A:481:SER:HB3	2.07	0.54
1:A:719:ASN:HB2	1:A:828:LEU:HD13	1.90	0.54
3:C:6:LYS:CD	3:C:10:PHE:HE2	2.18	0.54
1:A:1021:PHE:O	1:A:1024:VAL:HG22	2.08	0.54
1:A:989:LEU:HG	1:A:1000:GLN:HB2	1.90	0.53
1:A:476:SER:O	1:A:480:LEU:HG	2.08	0.53
1:A:2:PRO:O	1:A:6:ILE:HD12	2.07	0.53
1:A:412:VAL:O	1:A:416:VAL:HG23	2.09	0.53
3:C:13:ILE:O	3:C:17:VAL:HG23	2.08	0.53
1:A:62:THR:O	1:A:66:GLU:HG3	2.07	0.53
1:A:453:PHE:HE1	1:A:475:VAL:HG22	1.73	0.53
1:A:588:GLN:O	1:A:592:ASN:ND2	2.41	0.53
1:A:997:SER:HA	1:A:1000:GLN:HE22	1.74	0.53
1:A:20:MET:HA	1:A:377:LEU:CD1	2.38	0.53
1:A:449:LEU:HD13	1:A:936:GLY:HA3	1.90	0.53
1:A:971:ARG:NH1	1:A:974:PRO:HB2	2.24	0.53
1:A:144:ASN:HA	1:A:320:GLY:O	2.08	0.53
1:A:919:ARG:HH11	1:A:1005:THR:HG21	1.74	0.53
2:B:28:ASP:O	2:B:32:ILE:HG13	2.09	0.53
1:A:540:ARG:HD2	3:C:38:GLY:HA3	1.90	0.53
1:A:885:PHE:HA	1:A:902:MET:CE	2.39	0.53
2:B:40:VAL:HG21	2:B:69:ASN:OD1	2.09	0.52
1:A:10:ILE:O	1:A:13:TRP:N	2.42	0.52
1:A:431:THR:O	1:A:435:MET:HG2	2.09	0.52
1:A:375:VAL:HG13	1:A:480:LEU:HD12	1.92	0.52
1:A:932:LEU:O	1:A:935:ILE:HG13	2.09	0.52
1:A:559:LEU:HD23	1:A:923:ASN:HB2	1.92	0.52
1:A:560:PRO:HG2	1:A:922:THR:HG22	1.92	0.52
1:A:528:THR:HG21	1:A:969:ARG:HD3	1.92	0.52
4:A:2043:LMT:H11	4:A:2043:LMT:O2'	2.10	0.52
1:A:34:GLN:HB3	1:A:333:VAL:HG11	1.93	0.51
3:C:6:LYS:CD	3:C:10:PHE:CE2	2.91	0.51
1:A:712:MET:HG2	1:A:843:LEU:HD22	1.92	0.51
1:A:568:ASP:OD2	1:A:644:VAL:HG23	2.10	0.51
1:A:111:LEU:HD12	1:A:129:VAL:HG23	1.92	0.51
1:A:982:PHE:CD2	1:A:1011:MET:HG3	2.46	0.51
1:A:352:PHE:HD1	1:A:369:THR:HG21	1.75	0.51
1:A:679:GLY:HA2	1:A:830:GLN:HB3	1.91	0.51
1:A:1013:THR:O	1:A:1017:LEU:HB3	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:398:MET:O	1:A:402:ILE:HG23	2.10	0.51
1:A:949:ALA:O	1:A:953:MET:HG2	2.11	0.51
1:A:615:PHE:O	1:A:615:PHE:CD1	2.63	0.51
1:A:26:ALA:O	1:A:30:LEU:HB2	2.11	0.51
1:A:83:ASP:HA	1:A:815:ARG:HA	1.93	0.51
1:A:377:LEU:HD12	1:A:378:GLY:N	2.26	0.50
1:A:413:VAL:HA	1:A:493:CYS:SG	2.52	0.50
1:A:868:LEU:O	1:A:869:SER:OG	2.26	0.50
1:A:903:LEU:HD22	1:A:1025:PHE:CD1	2.46	0.50
1:A:410:ILE:HD11	1:A:978:THR:HA	1.92	0.50
1:A:437:GLN:HG2	1:A:438:ILE:HG23	1.93	0.50
1:A:464:GLY:O	1:A:468:ARG:HG2	2.11	0.50
1:A:102:ILE:O	1:A:106:GLN:HG3	2.12	0.50
1:A:417:GLU:O	1:A:420:MET:HG2	2.11	0.50
1:A:449:LEU:O	1:A:452:VAL:HG22	2.12	0.50
1:A:2:PRO:HB2	1:A:435:MET:HG3	1.93	0.50
1:A:143:ILE:HD13	1:A:285:PRO:O	2.12	0.50
1:A:584:GLN:HB2	1:A:622:GLN:HG2	1.93	0.50
1:A:595:THR:HG22	1:A:599:LEU:HD12	1.93	0.50
1:A:143:ILE:HD12	1:A:144:ASN:H	1.76	0.49
1:A:431:THR:HG21	1:A:494:ALA:HB2	1.94	0.49
3:C:18:VAL:HA	3:C:21:ILE:HD12	1.94	0.49
1:A:1027:VAL:O	1:A:1031:ARG:N	2.45	0.49
1:A:904:VAL:HG11	1:A:1022:VAL:HG22	1.93	0.49
1:A:905:VAL:HB	1:A:906:PRO:HD3	1.92	0.49
2:B:78:THR:O	2:B:79:LEU:HB2	2.12	0.49
1:A:449:LEU:CD1	1:A:936:GLY:HA3	2.42	0.49
1:A:715:SER:O	1:A:829:GLY:HA2	2.13	0.49
1:A:409:ALA:O	1:A:413:VAL:HG13	2.13	0.49
1:A:527:TYR:OH	1:A:1019:ILE:O	2.30	0.49
2:B:89:HIS:HB2	2:B:119:LEU:CD2	2.43	0.49
1:A:63:GLN:O	1:A:67:GLN:HG3	2.13	0.49
1:A:641:GLU:O	1:A:650:ARG:NH2	2.38	0.49
1:A:880:SER:O	1:A:884:VAL:HG23	2.12	0.49
1:A:45:ILE:HG12	1:A:129:VAL:HG22	1.95	0.49
1:A:913:LEU:O	1:A:917:THR:HG23	2.13	0.49
4:A:2043:LMT:H6'2	4:A:2043:LMT:O3B	2.13	0.49
1:A:428:LYS:HZ2	1:A:494:ALA:HB1	1.78	0.48
1:A:184:MET:HB3	1:A:771:VAL:HG22	1.96	0.48
1:A:369:THR:O	1:A:373:PRO:HD2	2.14	0.48
1:A:764:ASP:OD2	1:A:769:LYS:NZ	2.38	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:157:TYR:HH	1:A:317:PHE:HD1	1.61	0.48
1:A:438:ILE:O	1:A:441:ALA:N	2.41	0.48
1:A:719:ASN:ND2	1:A:826:GLU:OE2	2.47	0.48
1:A:686:ASP:HB2	1:A:695:LEU:HD22	1.95	0.48
1:A:645:GLU:O	1:A:649:MET:HG2	2.13	0.48
1:A:211:ASN:HA	1:A:240:LEU:HD13	1.96	0.48
1:A:940:LYS:HE3	1:A:941:ASN:OD1	2.13	0.48
1:A:49:TYR:CD2	1:A:52:ALA:HB2	2.49	0.47
1:A:445:ILE:HG22	1:A:943:ILE:HD13	1.96	0.47
2:B:34:MET:SD	2:B:40:VAL:HG22	2.54	0.47
1:A:449:LEU:HD23	1:A:478:MET:SD	2.55	0.47
1:A:901:VAL:HG23	1:A:942:ALA:CB	2.44	0.47
2:B:63:VAL:HG11	2:B:99:LEU:CD1	2.45	0.47
1:A:314:GLU:N	1:A:314:GLU:OE1	2.47	0.47
1:A:809:TRP:CE2	2:B:46:VAL:HG23	2.50	0.47
2:B:56:TYR:HD1	2:B:57:TRP:CD1	2.32	0.47
1:A:111:LEU:HD11	1:A:127:VAL:HG12	1.97	0.47
1:A:448:VAL:O	1:A:452:VAL:HG13	2.14	0.47
1:A:960:LEU:HD21	1:A:1030:ARG:CB	2.39	0.47
1:A:1:MET:HB3	1:A:2:PRO:HD3	1.97	0.46
1:A:72:ILE:HG22	1:A:106:GLN:HB3	1.97	0.46
1:A:198:LEU:HD13	1:A:251:LEU:HD13	1.96	0.46
1:A:242:SER:OG	1:A:244:GLU:OE1	2.33	0.46
1:A:643:LYS:O	1:A:647:ILE:HG13	2.15	0.46
1:A:412:VAL:HB	1:A:438:ILE:HD11	1.95	0.46
1:A:34:GLN:HB3	1:A:333:VAL:CG1	2.45	0.46
1:A:358:PHE:CG	1:A:977:MET:HG2	2.50	0.46
1:A:35:TYR:OH	1:A:670:ALA:O	2.24	0.46
1:A:151:GLN:HE22	1:A:278:ILE:HG23	1.80	0.46
1:A:415:ASN:O	1:A:419:VAL:HG23	2.16	0.46
1:A:468:ARG:O	1:A:472:ILE:HG23	2.15	0.46
1:A:154:ILE:HG22	1:A:287:SER:HB3	1.98	0.46
2:B:113:GLY:O	2:B:143:ASP:HA	2.16	0.46
1:A:135:SER:OG	1:A:672:VAL:O	2.17	0.46
1:A:480:LEU:O	1:A:484:VAL:HG13	2.16	0.46
1:A:873:ALA:N	1:A:874:PRO:HD2	2.30	0.45
1:A:399:VAL:O	1:A:402:ILE:HG13	2.16	0.45
1:A:782:LEU:O	1:A:785:ASP:HB2	2.15	0.45
1:A:711:ASP:N	1:A:711:ASP:OD1	2.49	0.45
1:A:72:ILE:HG13	1:A:75:LEU:HD22	1.97	0.45
1:A:471:SER:O	1:A:475:VAL:HG23	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:586:ARG:O	1:A:589:LYS:HB3	2.16	0.45
1:A:962:GLU:OE1	1:A:962:GLU:N	2.44	0.45
1:A:722:GLU:O	1:A:724:THR:HG23	2.17	0.45
1:A:992:SER:CB	1:A:1000:GLN:HE21	2.30	0.45
1:A:367:ILE:HB	1:A:368:PRO:HD3	1.97	0.45
1:A:549:VAL:O	1:A:552:MET:HB3	2.16	0.45
1:A:440:GLY:O	1:A:891:LEU:HD13	2.17	0.45
1:A:137:LEU:CD2	1:A:303:ALA:HB2	2.44	0.45
1:A:567:GLU:HG2	1:A:998:GLY:HA3	1.99	0.45
1:A:696:THR:HA	1:A:825:MET:HE1	1.98	0.45
1:A:540:ARG:HD2	3:C:38:GLY:CA	2.48	0.44
1:A:594:VAL:HG13	1:A:655:PHE:CZ	2.52	0.44
1:A:888:LEU:HD13	1:A:901:VAL:HG11	1.99	0.44
2:B:30:VAL:O	2:B:34:MET:HG2	2.17	0.44
1:A:402:ILE:HG13	1:A:403:GLY:N	2.32	0.44
1:A:754:TRP:CZ2	1:A:786:ILE:HD13	2.52	0.44
1:A:373:PRO:O	1:A:377:LEU:HG	2.17	0.44
1:A:411:VAL:O	1:A:415:ASN:HB2	2.17	0.44
2:B:48:TRP:NE1	2:B:77:ASP:OD2	2.51	0.44
2:B:51:LEU:HD23	2:B:83:PRO:HG2	1.99	0.44
1:A:433:LYS:O	1:A:437:GLN:OE1	2.36	0.44
1:A:980:LEU:HD21	3:C:19:MET:HB2	1.99	0.43
2:B:67:LEU:HD22	2:B:73:VAL:HG22	2.00	0.43
1:A:176:GLN:N	1:A:290:GLY:O	2.50	0.43
1:A:1032:ARG:NH2	3:C:42:LYS:O	2.51	0.43
1:A:352:PHE:CD1	1:A:369:THR:HG21	2.52	0.43
1:A:583:THR:HG22	1:A:586:ARG:HD3	2.00	0.43
3:C:29:LEU:HA	3:C:32:VAL:HG22	1.99	0.43
1:A:110:LYS:HE2	1:A:110:LYS:HB3	1.52	0.43
1:A:888:LEU:HD12	1:A:902:MET:HE1	1.99	0.43
1:A:644:VAL:HG11	1:A:667:ASN:ND2	2.33	0.43
2:B:25:GLY:O	2:B:62:ILE:HD11	2.19	0.43
1:A:721:LEU:CD1	1:A:815:ARG:HD3	2.48	0.43
4:A:2043:LMT:O2B	4:A:2043:LMT:C3'	2.66	0.43
1:A:997:SER:HA	1:A:1000:GLN:NE2	2.33	0.43
1:A:1024:VAL:HA	1:A:1027:VAL:HG22	2.00	0.43
1:A:342:LYS:O	1:A:346:GLU:HG2	2.19	0.43
1:A:988:PRO:O	1:A:991:ILE:HG13	2.18	0.43
1:A:1025:PHE:O	1:A:1029:VAL:HG23	2.19	0.43
2:B:52:HIS:CE1	2:B:83:PRO:HD3	2.54	0.43
1:A:407:ASP:OD1	1:A:408:ASP:N	2.52	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:575:MET:HE1	1:A:662:MET:CE	2.39	0.43
1:A:459:PHE:O	1:A:464:GLY:HA3	2.19	0.42
1:A:809:TRP:CD2	2:B:46:VAL:HG23	2.54	0.42
1:A:535:LEU:HD12	1:A:535:LEU:HA	1.83	0.42
1:A:669:PRO:O	1:A:671:ILE:N	2.52	0.42
3:C:31:GLU:O	3:C:34:ASN:HB2	2.19	0.42
1:A:559:LEU:HD12	1:A:560:PRO:HD2	2.01	0.42
1:A:16:ALA:O	1:A:20:MET:HG3	2.19	0.42
1:A:102:ILE:O	1:A:105:VAL:HG22	2.18	0.42
1:A:576:VAL:HG11	1:A:591:LEU:HD12	2.00	0.42
1:A:875:SER:O	1:A:878:ALA:HB3	2.19	0.42
1:A:904:VAL:HA	1:A:907:LEU:HG	2.01	0.42
3:C:6:LYS:CG	3:C:10:PHE:CD2	2.97	0.42
1:A:422:GLU:O	1:A:502:LYS:HB2	2.19	0.42
2:B:51:LEU:HD11	2:B:63:VAL:HG13	2.00	0.42
3:C:1:MET:O	3:C:4:LEU:HB2	2.19	0.42
1:A:686:ASP:HB3	1:A:823:PRO:HB2	2.01	0.42
1:A:1019:ILE:HD11	3:C:26:ILE:HG21	2.01	0.42
3:C:6:LYS:HG2	3:C:10:PHE:CE2	2.41	0.42
1:A:832:ALA:HB1	1:A:833:PRO:HD2	2.02	0.42
1:A:928:GLN:O	1:A:932:LEU:HD12	2.19	0.42
3:C:6:LYS:HG3	3:C:10:PHE:CD2	2.33	0.42
1:A:376:LEU:HD12	1:A:377:LEU:N	2.35	0.42
1:A:449:LEU:HG	1:A:453:PHE:CE2	2.54	0.42
1:A:1008:MET:HB2	1:A:1008:MET:HE3	1.83	0.42
1:A:929:VAL:O	1:A:933:THR:HG23	2.20	0.41
1:A:454:VAL:N	1:A:455:PRO:CD	2.83	0.41
1:A:395:MET:O	1:A:398:MET:HG2	2.20	0.41
1:A:515:TRP:NE1	1:A:519:MET:SD	2.82	0.41
1:A:617:PHE:HD2	1:A:626:ILE:HD11	1.84	0.41
1:A:958:LYS:HE3	1:A:962:GLU:HG2	2.02	0.41
2:B:143:ASP:OD1	2:B:146:GLY:N	2.53	0.41
1:A:186:ILE:HB	1:A:773:VAL:HG23	2.02	0.41
1:A:743:ILE:H	1:A:743:ILE:HD12	1.85	0.41
1:A:166:ILE:HG22	1:A:175:VAL:HG21	2.02	0.41
1:A:438:ILE:O	1:A:439:GLN:C	2.59	0.41
1:A:908:GLY:O	1:A:1010:GLY:HA2	2.20	0.41
3:C:19:MET:O	3:C:22:ILE:HG13	2.20	0.41
2:B:109:LYS:HG2	2:B:110:ASP:O	2.20	0.41
1:A:159:ALA:HB1	1:A:181:GLN:HB2	2.03	0.41
1:A:886:LEU:HD12	1:A:887:CYS:N	2.36	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:987:MET:HB2	1:A:988:PRO:HD3	2.01	0.41
1:A:992:SER:HB3	1:A:1000:GLN:HE21	1.86	0.41
3:C:23:LEU:HD23	3:C:26:ILE:HD11	2.02	0.41
1:A:88:VAL:HG12	1:A:89:GLN:N	2.36	0.41
1:A:330:THR:O	1:A:333:VAL:N	2.53	0.41
1:A:984:LEU:HA	1:A:987:MET:CG	2.47	0.41
1:A:1012:VAL:HG23	1:A:1013:THR:N	2.36	0.41
1:A:1017:LEU:O	1:A:1021:PHE:HD2	2.04	0.41
1:A:5:PHE:CD1	1:A:487:ILE:HG23	2.56	0.41
1:A:49:TYR:CE1	1:A:121:GLU:HB2	2.56	0.41
1:A:300:LEU:O	1:A:300:LEU:HD23	2.21	0.41
1:A:403:GLY:O	1:A:406:VAL:HG22	2.20	0.41
1:A:415:ASN:ND2	1:A:437:GLN:NE2	2.69	0.41
1:A:448:VAL:HG11	1:A:888:LEU:HD21	2.01	0.41
1:A:649:MET:O	1:A:652:THR:OG1	2.26	0.41
1:A:453:PHE:CD1	1:A:475:VAL:HG22	2.57	0.40
1:A:945:ILE:HD13	1:A:968:VAL:HG23	2.02	0.40
2:B:128:ILE:O	2:B:131:VAL:HG12	2.21	0.40
1:A:110:LYS:O	1:A:113:LEU:HB2	2.21	0.40
1:A:166:ILE:HA	1:A:309:GLU:HG2	2.03	0.40
1:A:196:PHE:O	1:A:197:GLN:HB2	2.22	0.40
1:A:407:ASP:O	1:A:411:VAL:HG22	2.20	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	1028/1047 (98%)	959 (93%)	66 (6%)	3 (0%)	41 71
2	B	150/169 (89%)	143 (95%)	6 (4%)	1 (1%)	22 54
3	C	44/49 (90%)	43 (98%)	1 (2%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	1222/1265 (97%)	1145 (94%)	73 (6%)	4 (0%)	41 71

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	223	PRO
1	A	224	PRO
2	B	139	VAL
1	A	579	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	823/853 (96%)	804 (98%)	19 (2%)	50 73
2	B	116/132 (88%)	113 (97%)	3 (3%)	46 71
3	C	34/41 (83%)	34 (100%)	0	100 100
All	All	973/1026 (95%)	951 (98%)	22 (2%)	50 73

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	46	SER
1	A	68	ASN
1	A	75	LEU
1	A	402	ILE
1	A	412	VAL
1	A	425	LEU
1	A	472	ILE
1	A	481	SER
1	A	487	ILE
1	A	489	THR
1	A	493	CYS
1	A	547	ILE

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Mol	Chain	Res	Type
1	A	640	GLU
1	A	801	PHE
1	A	904	VAL
1	A	932	LEU
1	A	945	ILE
1	A	993	THR
1	A	1034	SER
2	B	60	LEU
2	B	119	LEU
2	B	133	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	360	GLN
1	A	361	ASN
1	A	415	ASN
1	A	517	ASN
1	A	592	ASN
2	B	41	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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 10 ligands modelled in this entry, 1 is monoatomic - leaving 9 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
4	LMT	A	2043	-	36,36,36	0.46	0	47,47,47	0.87	1 (2%)
4	LMT	A	2038	-	9,9,36	0.27	0	8,8,47	0.48	0
4	LMT	A	2039	-	5,5,36	0.25	0	4,4,47	0.43	0
4	LMT	A	2041	-	7,7,36	0.25	0	6,6,47	0.49	0
4	LMT	A	2044	-	7,7,36	0.31	0	6,6,47	0.44	0
4	LMT	A	2046	-	5,5,36	0.30	0	4,4,47	0.29	0
4	LMT	A	2045	-	3,3,36	0.38	0	2,2,47	0.61	0
4	LMT	A	2042	-	11,11,36	0.21	0	10,10,47	0.61	0
4	LMT	A	2047	-	5,5,36	0.32	0	4,4,47	0.28	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
4	LMT	A	2043	-	-	12/21/61/61	0/2/2/2
4	LMT	A	2038	-	-	4/7/7/61	-
4	LMT	A	2039	-	-	3/3/3/61	-
4	LMT	A	2041	-	-	1/5/5/61	-
4	LMT	A	2044	-	-	3/5/5/61	-
4	LMT	A	2046	-	-	1/3/3/61	-
4	LMT	A	2045	-	-	1/1/1/61	-
4	LMT	A	2042	-	-	4/9/9/61	-
4	LMT	A	2047	-	-	1/3/3/61	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	2043	LMT	O5'-C5'-C4'	2.29	114.59	109.75

There are no chirality outliers.

All (30) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	2043	LMT	C2B-C1B-O1B-C4'
4	A	2043	LMT	O5'-C1'-O1'-C1
4	A	2045	LMT	C1-C2-C3-C4
4	A	2043	LMT	O5'-C5'-C6'-O6'
4	A	2043	LMT	O5B-C5B-C6B-O6B
4	A	2043	LMT	C4'-C5'-C6'-O6'
4	A	2043	LMT	C4B-C5B-C6B-O6B
4	A	2043	LMT	C2'-C1'-O1'-C1
4	A	2038	LMT	C2-C3-C4-C5
4	A	2043	LMT	C2-C3-C4-C5
4	A	2041	LMT	C4-C5-C6-C7
4	A	2042	LMT	C4-C5-C6-C7
4	A	2042	LMT	C2-C3-C4-C5
4	A	2038	LMT	C3-C4-C5-C6
4	A	2043	LMT	C5-C6-C7-C8
4	A	2039	LMT	C3-C4-C5-C6
4	A	2047	LMT	C2-C3-C4-C5
4	A	2043	LMT	C1-C2-C3-C4
4	A	2038	LMT	C1-C2-C3-C4
4	A	2044	LMT	C1-C2-C3-C4
4	A	2042	LMT	C6-C7-C8-C9
4	A	2039	LMT	C1-C2-C3-C4
4	A	2046	LMT	C1-C2-C3-C4
4	A	2043	LMT	C9-C10-C11-C12
4	A	2042	LMT	C3-C4-C5-C6
4	A	2044	LMT	C4-C5-C6-C7
4	A	2044	LMT	C2-C3-C4-C5
4	A	2039	LMT	C2-C3-C4-C5
4	A	2038	LMT	C4-C5-C6-C7
4	A	2043	LMT	C6-C7-C8-C9

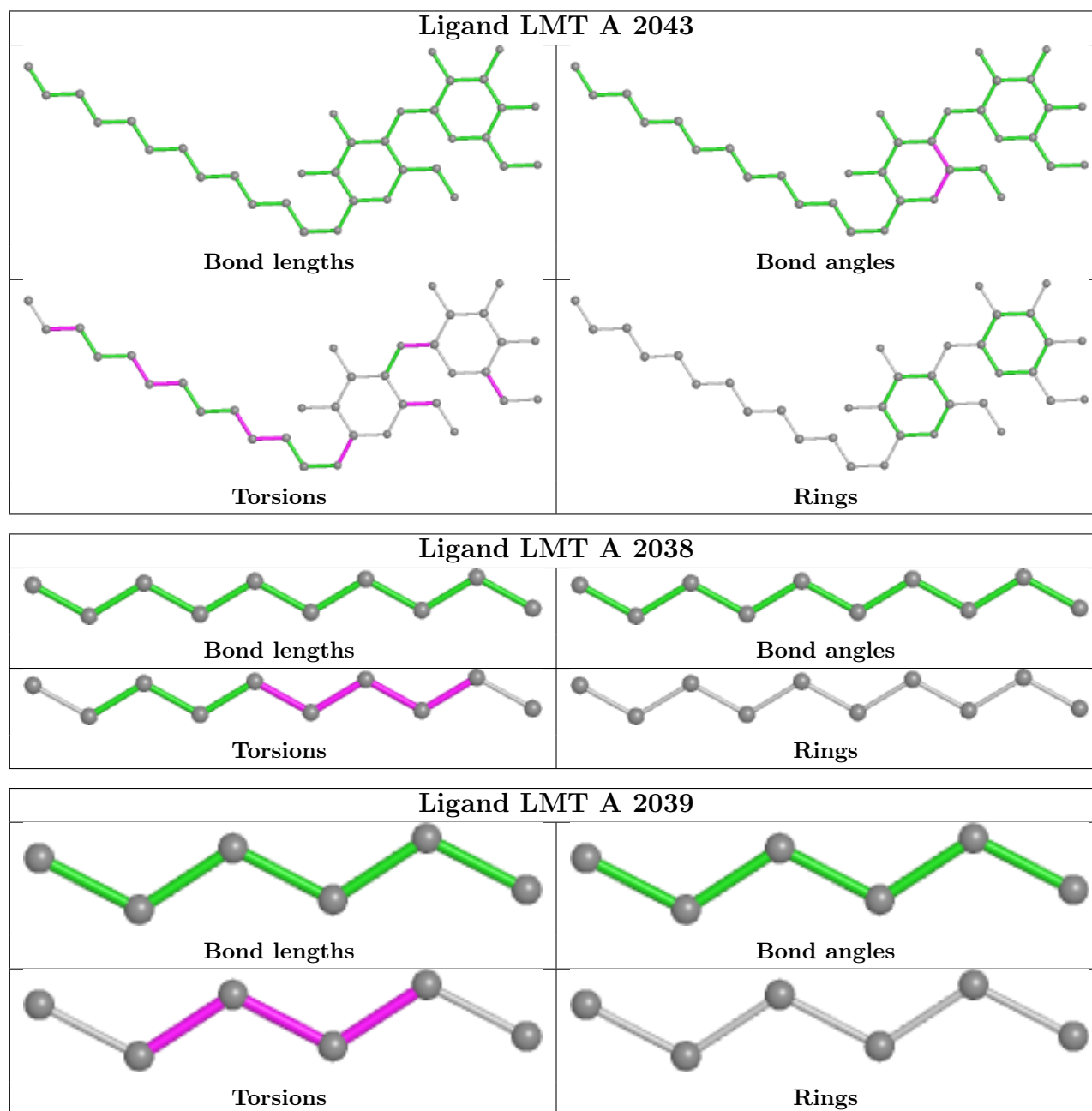
There are no ring outliers.

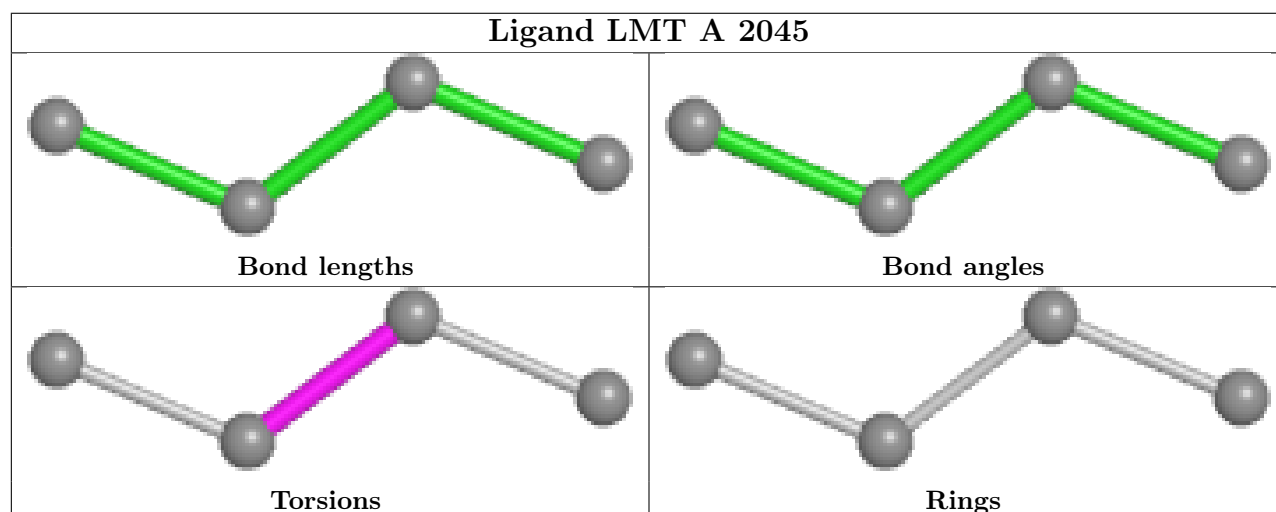
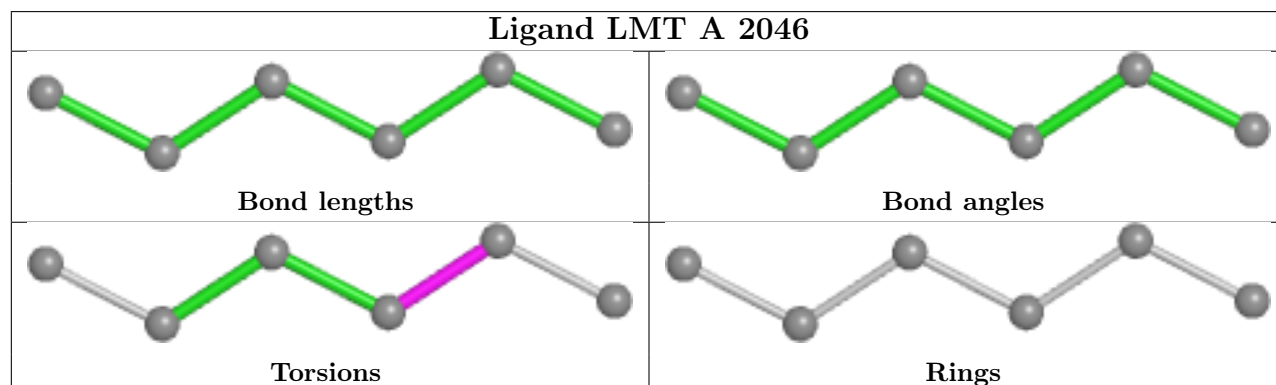
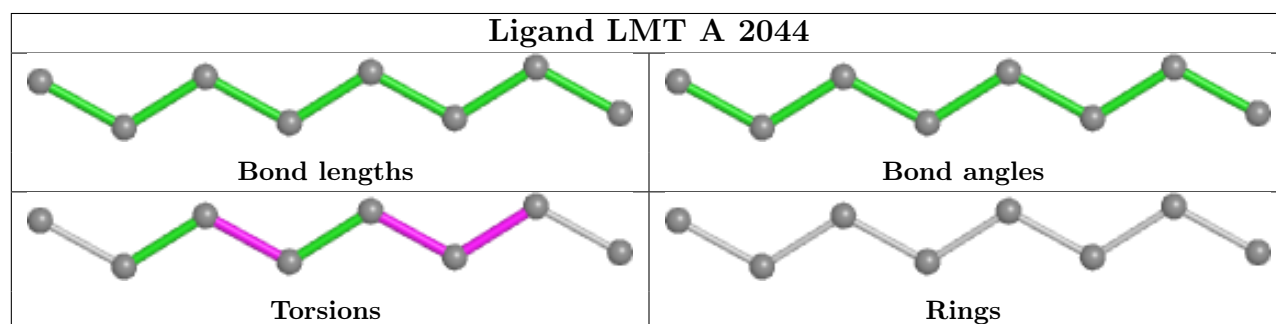
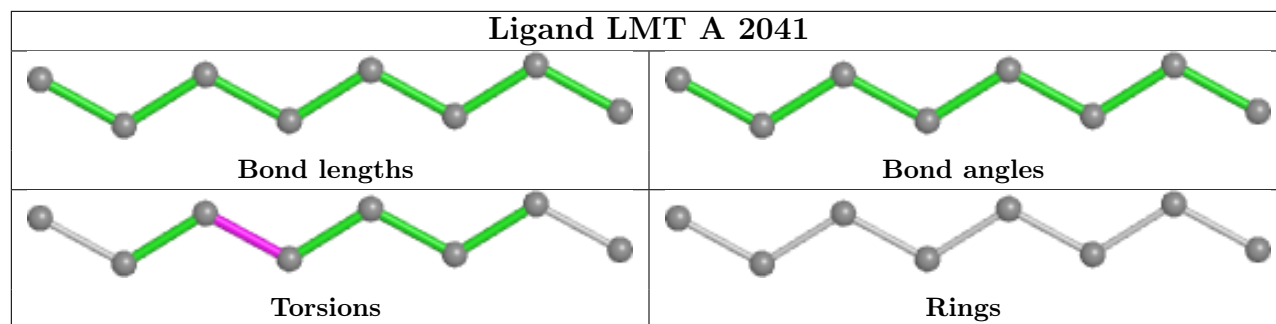
1 monomer is involved in 10 short contacts:

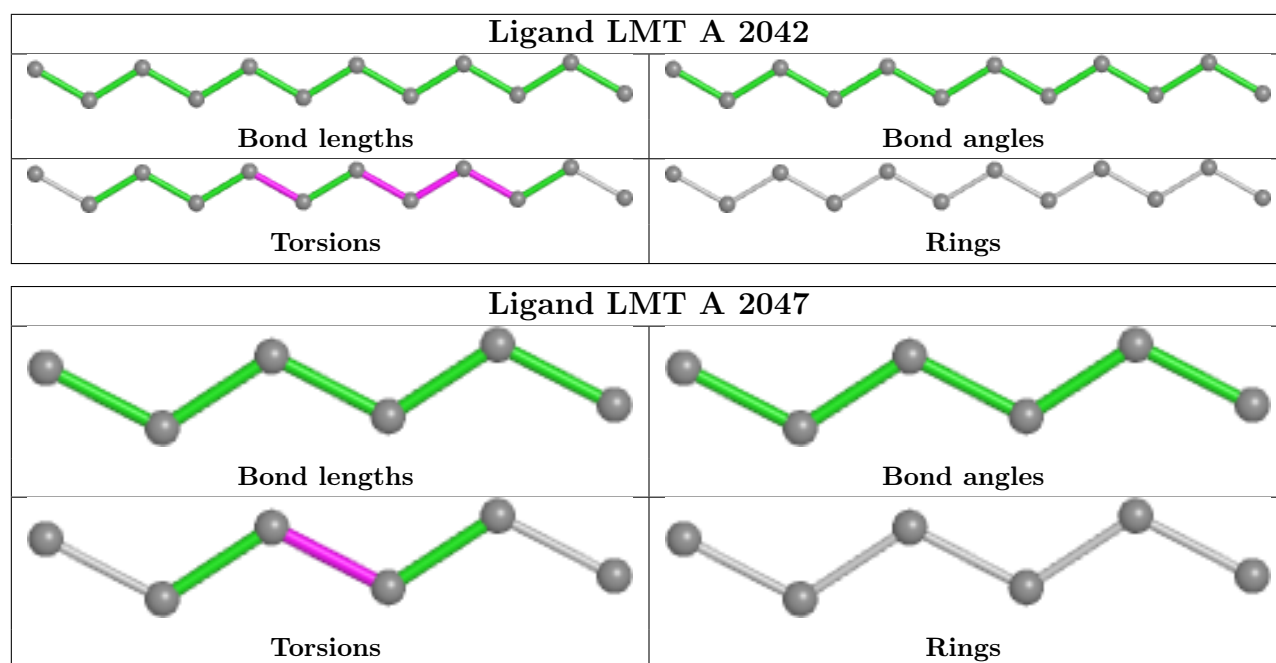
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	2043	LMT	10	0

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	1032/1047 (98%)	-0.04	34 (3%) 46 44	41, 97, 120, 120	0
2	B	152/169 (89%)	0.52	18 (11%) 4 4	92, 119, 120, 120	0
3	C	46/49 (93%)	-0.05	2 (4%) 35 34	83, 118, 120, 120	0
All	All	1230/1265 (97%)	0.03	54 (4%) 34 33	41, 104, 120, 120	0

All (54) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	836	SER	5.5
2	B	35	ALA	5.4
2	B	76	TYR	4.4
1	A	870	GLY	4.2
2	B	77	ASP	3.9
1	A	869	SER	3.8
1	A	386	PHE	3.7
1	A	425	LEU	3.7
1	A	388	PHE	3.5
2	B	36	ASN	3.4
1	A	387	GLY	3.4
1	A	874	PRO	3.4
2	B	84	LEU	3.3
1	A	422	GLU	3.2
1	A	835	LYS	3.2
1	A	839	GLU	3.2
2	B	80	GLY	3.1
2	B	103	GLY	3.1
1	A	501	ALA	3.0
2	B	102	ASN	3.0
3	C	4	LEU	2.9
1	A	305	ALA	2.9
1	A	1034	SER	2.8

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Mol	Chain	Res	Type	RSRZ
2	B	78	THR	2.8
1	A	505	HIS	2.7
1	A	503	GLY	2.7
1	A	509	LYS	2.7
2	B	141	ALA	2.6
1	A	504	ASP	2.6
2	B	162	ALA	2.6
2	B	96	VAL	2.5
1	A	436	GLY	2.5
1	A	461	GLY	2.5
1	A	635	ALA	2.4
3	C	1	MET	2.4
1	A	1035	ARG	2.4
1	A	554	TYR	2.4
1	A	510	LYS	2.4
1	A	838	GLY	2.4
2	B	150	PHE	2.4
1	A	638	PRO	2.3
1	A	413	VAL	2.3
1	A	717	ARG	2.3
1	A	640	GLU	2.3
2	B	157	GLY	2.3
1	A	641	GLU	2.2
2	B	109	LYS	2.2
2	B	158	ASN	2.2
1	A	423	GLU	2.1
1	A	148	THR	2.0
2	B	107	ASN	2.0
2	B	140	ASN	2.0
1	A	995	ALA	2.0
1	A	437	GLN	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](#)

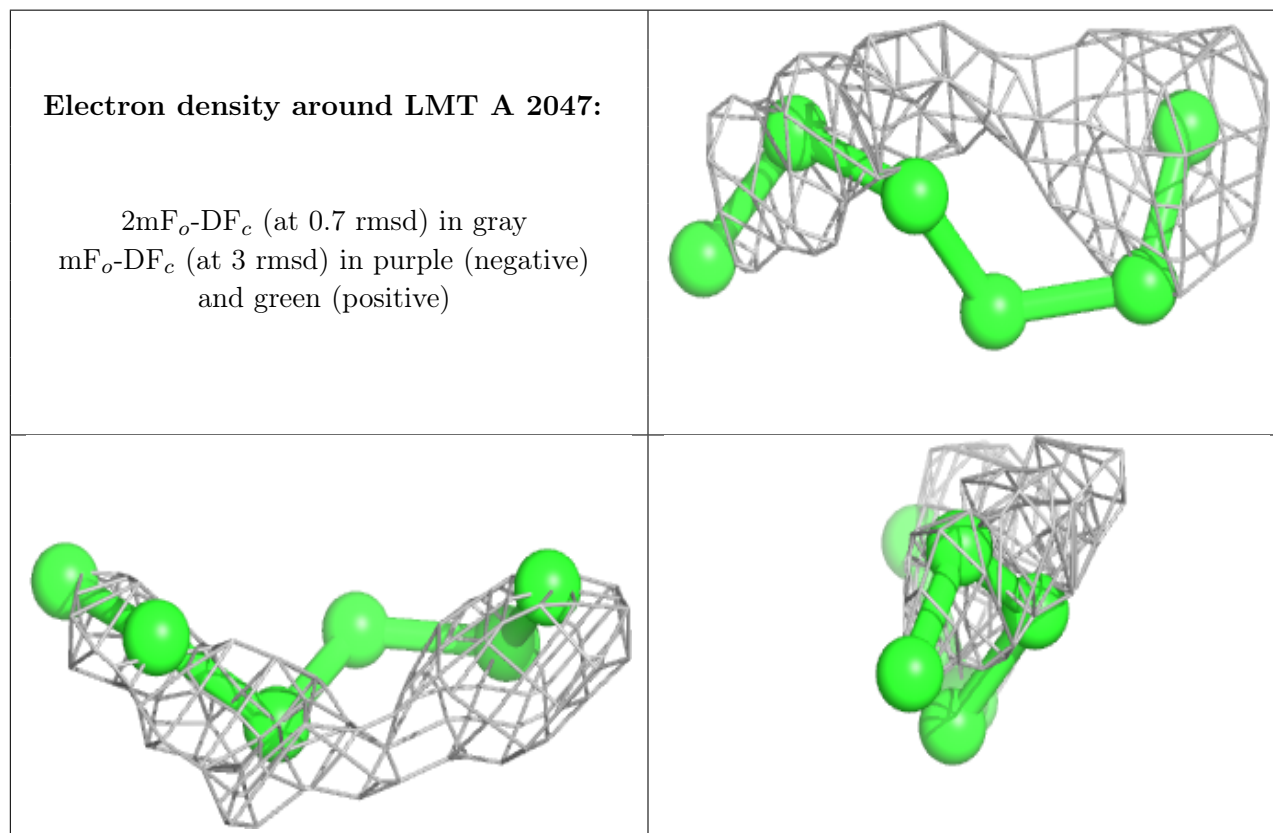
There are no monosaccharides in this entry.

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, 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.

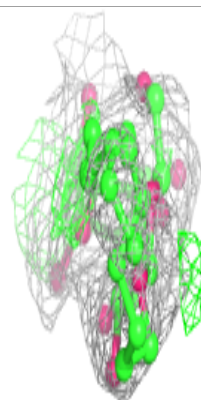
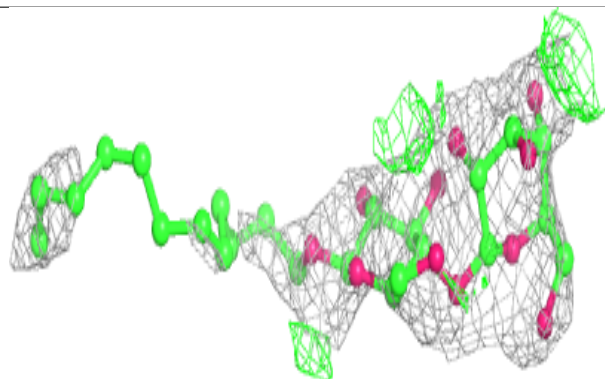
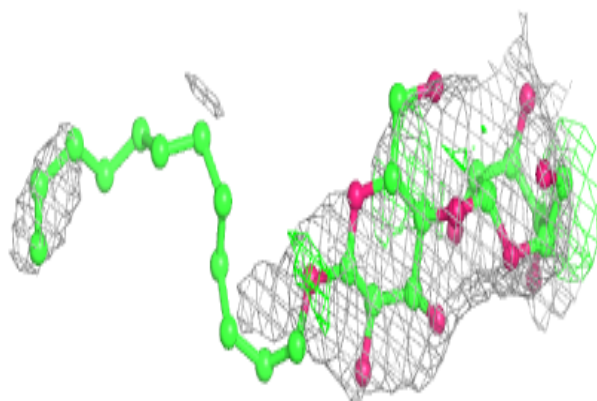
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	LMT	A	2047	6/35	0.53	0.42	81,88,89,89	6
4	LMT	A	2043	35/35	0.67	0.48	68,89,98,100	35
4	LMT	A	2046	6/35	0.71	0.46	97,105,113,114	0
4	LMT	A	2038	10/35	0.71	0.37	111,120,120,120	0
4	LMT	A	2039	6/35	0.73	0.39	57,60,62,63	6
4	LMT	A	2042	12/35	0.76	0.50	59,66,69,70	12
4	LMT	A	2044	8/35	0.76	0.43	88,95,99,100	0
4	LMT	A	2045	4/35	0.79	0.20	58,58,60,60	4
4	LMT	A	2041	8/35	0.83	0.40	95,101,105,107	0
5	NI	A	2040	1/1	0.97	0.04	120,120,120,120	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.

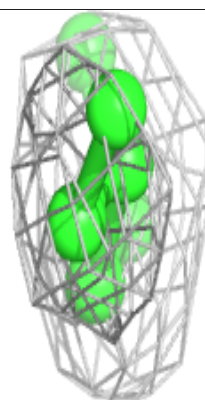
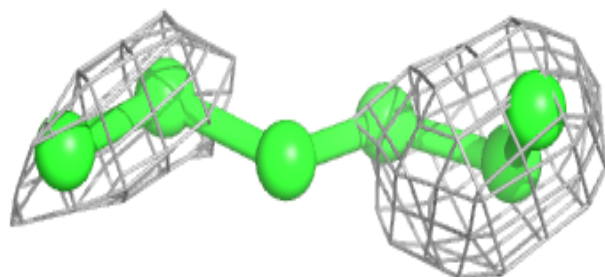
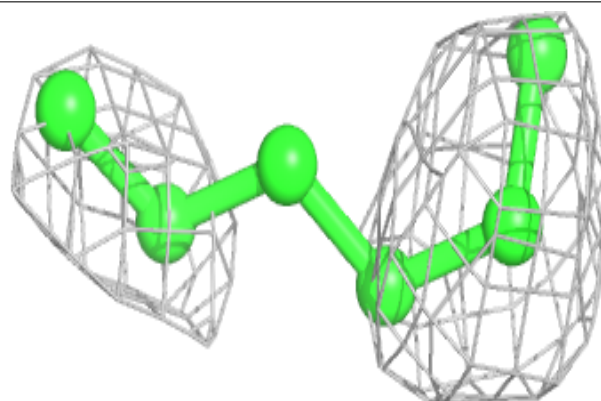


Electron density around LMT A 2043:

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

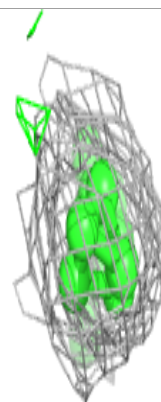
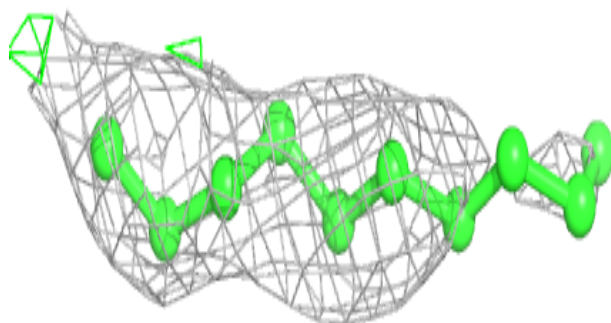
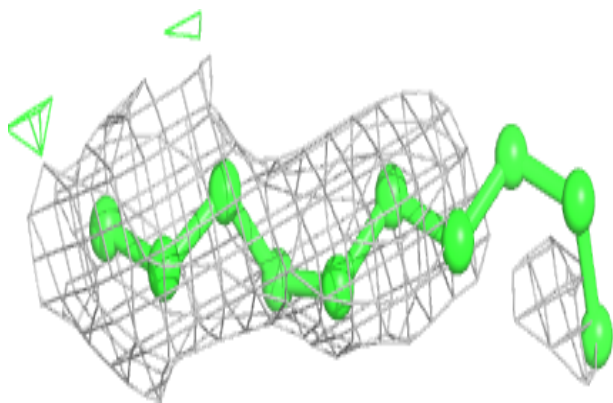
**Electron density around LMT A 2046:**

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

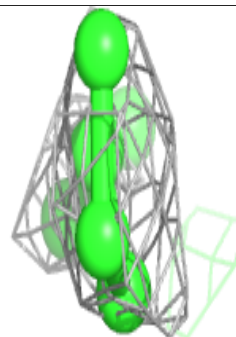
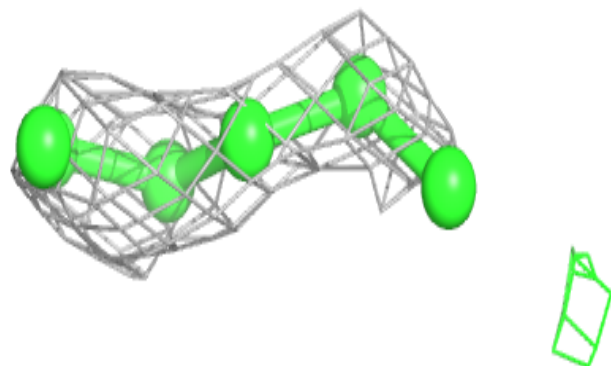
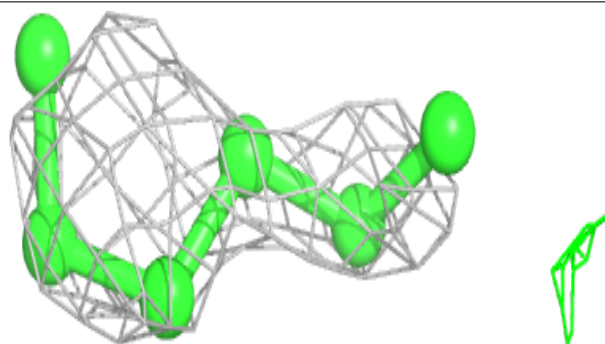


Electron density around LMT A 2038:

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

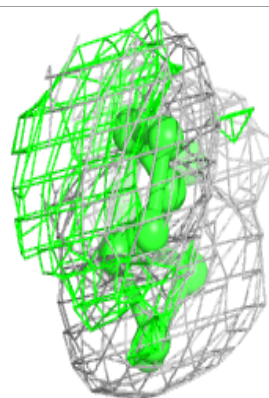
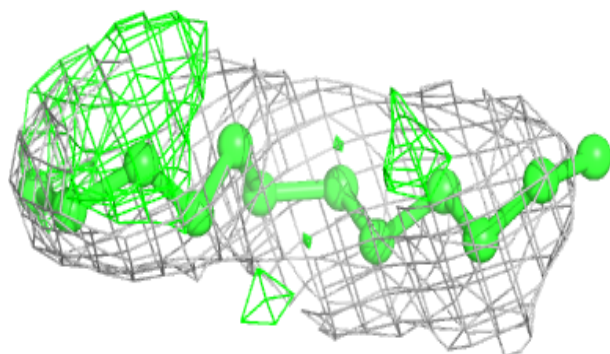
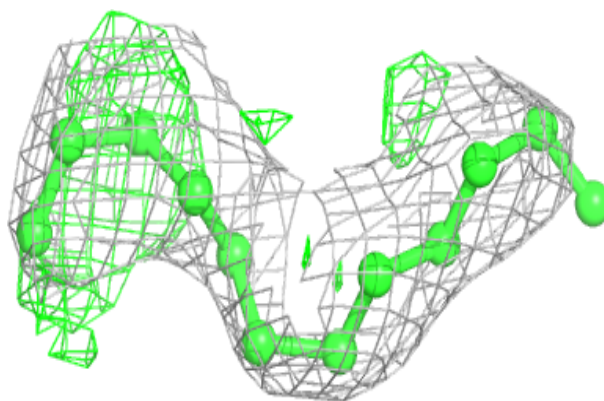
**Electron density around LMT A 2039:**

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

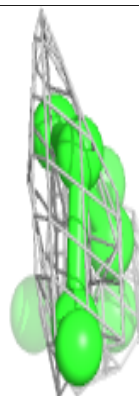
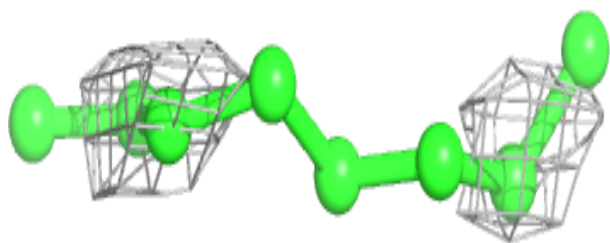
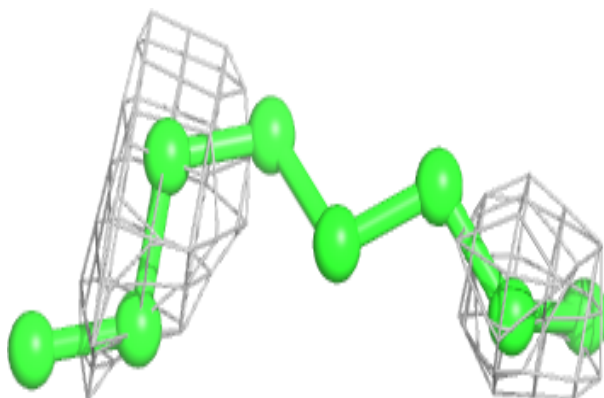


Electron density around LMT A 2042:

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

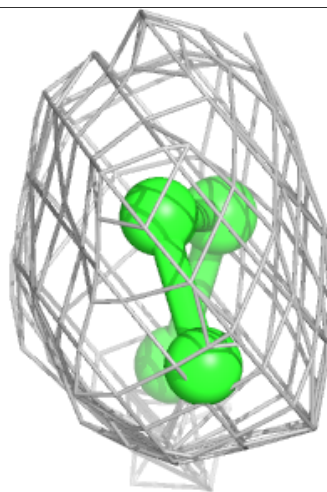
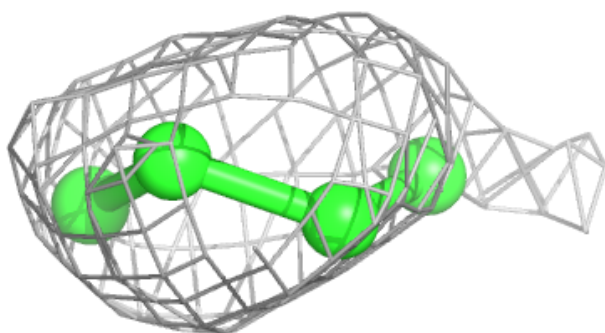
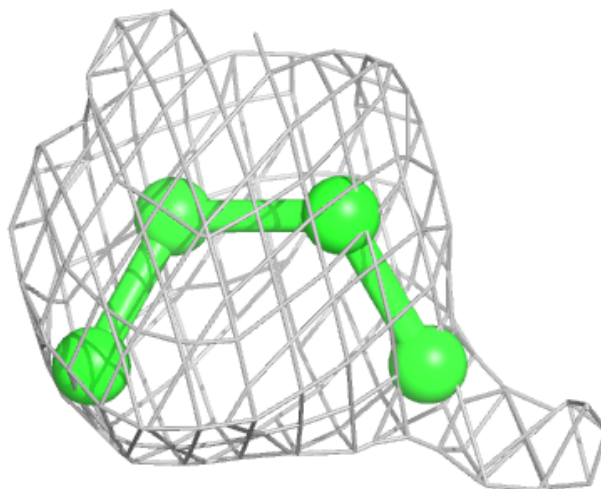
**Electron density around LMT A 2044:**

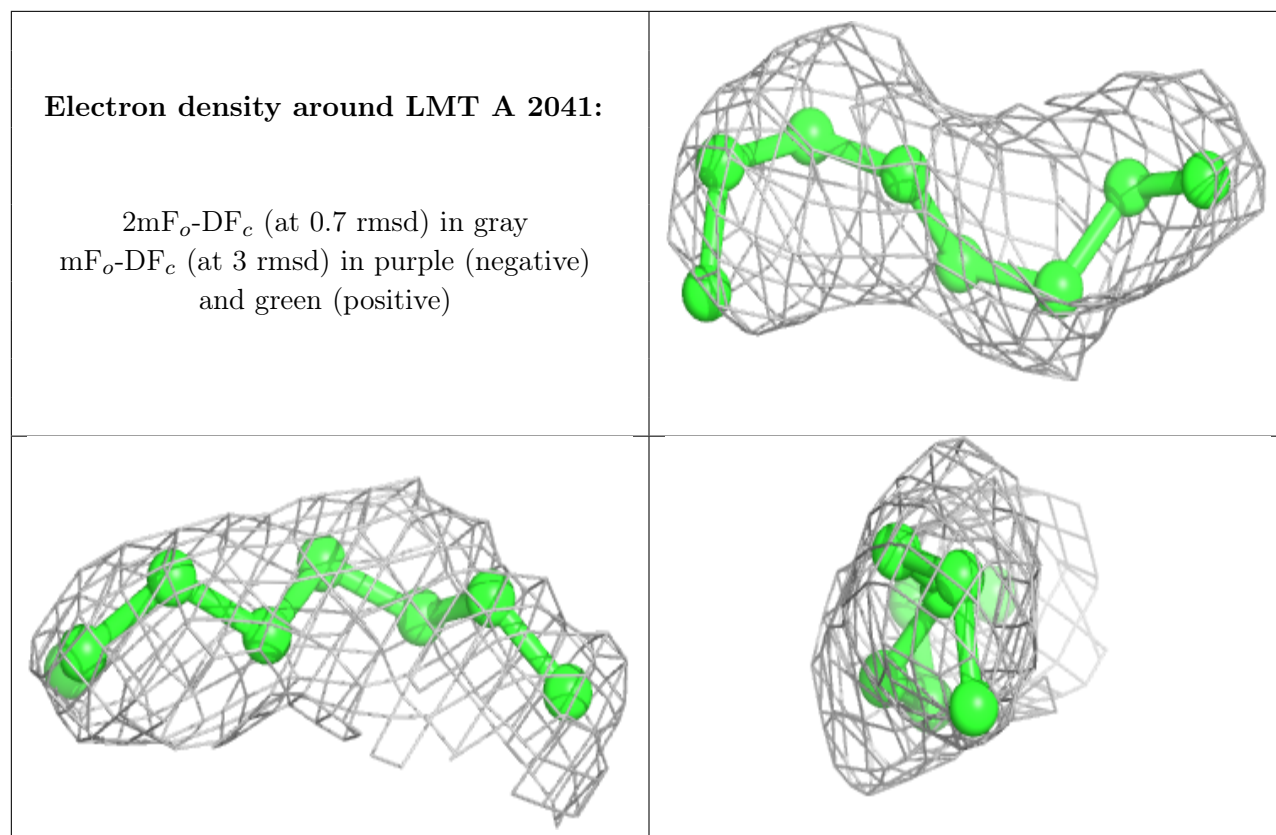
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around LMT A 2045:

$2mF_o-DF_c$ (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.