



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 09:24 PM UTC

PDB ID : 5I6C / pdb_00005i6c
Title : The structure of the eukaryotic purine/H⁺ symporter, UapA, in complex with Xanthine
Authors : Alguel, Y.; Amillis, S.; Leung, J.; Lambrinidis, G.; Capaldi, S.; Scull, N.J.; Craven, G.; Iwata, S.; Armstrong, A.; Mikros, E.; Diallinas, G.; Cameron, A.D.; Byrne, B.
Deposited on : 2016-02-16
Resolution : 3.70 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](#) ①) 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

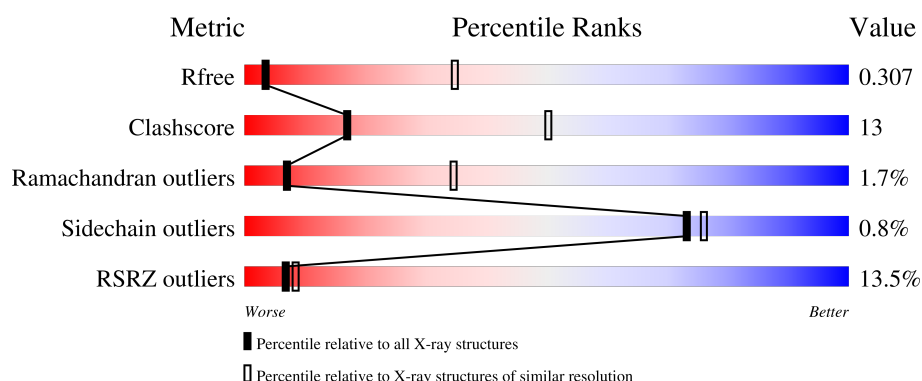
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.70 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	1131 (3.80-3.60)
Clashscore	190562	1171 (3.80-3.60)
Ramachandran outliers	187476	1129 (3.80-3.60)
Sidechain outliers	187428	1126 (3.80-3.60)
RSRZ outliers	180081	1130 (3.80-3.60)

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	574	10% 58% 24% • 16%
1	B	574	12% 59% 21% • 18%

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 7127 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.

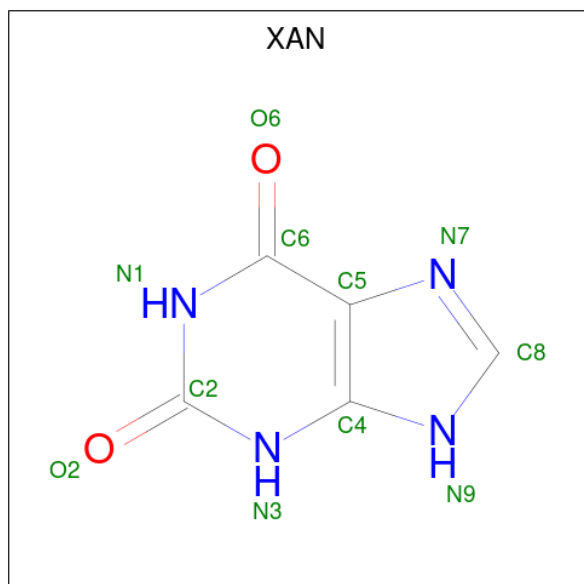
- Molecule 1 is a protein called Uric acid-xanthine permease.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	480	Total	C	N	O	S	0	0	0
			3545	2319	565	626	35			
1	B	468	Total	C	N	O	S	0	0	0
			3434	2242	547	610	35			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	411	VAL	GLY	engineered mutation	UNP Q07307
B	411	VAL	GLY	engineered mutation	UNP Q07307

- Molecule 2 is XANTHINE (CCD ID: XAN) (formula: C₅H₄N₄O₂).



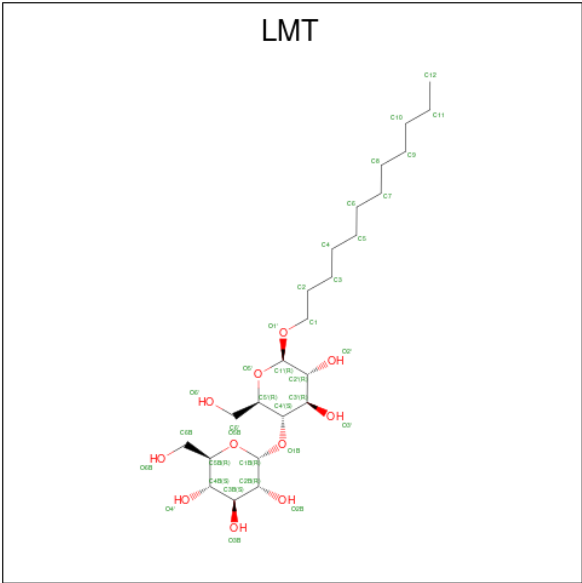
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			11	5	4	2		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	B	1	Total	C	N	O	0	0
			11	5	4	2		

- Molecule 3 is DODECYL-BETA-D-MALTOSIDE (CCD ID: LMT) (formula: C₂₄H₄₆O₁₁).

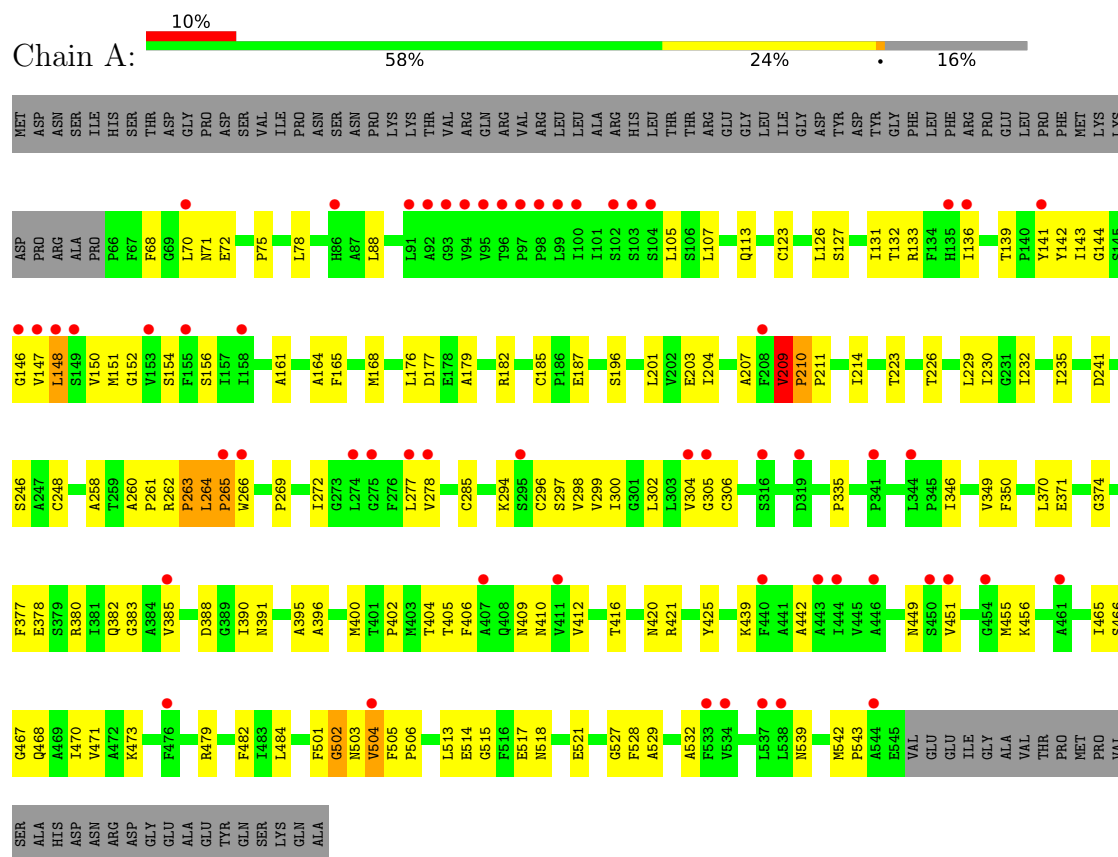


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			35	24	11		
3	A	1	Total	C	O	0	0
			28	17	11		
3	B	1	Total	C	O	0	0
			35	24	11		
3	B	1	Total	C	O	0	0
			28	17	11		

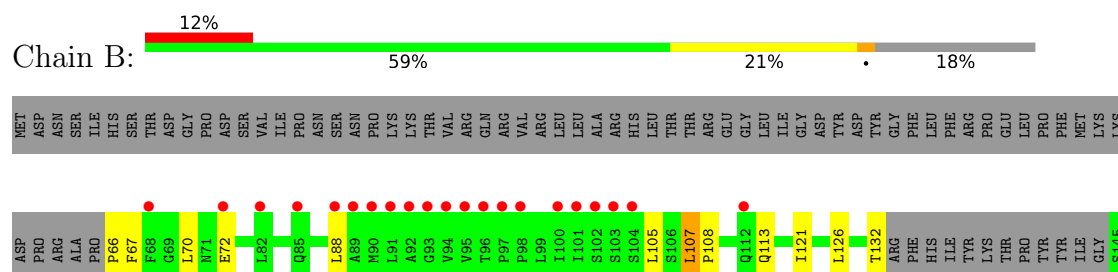
3 Residue-property plots [i](#)

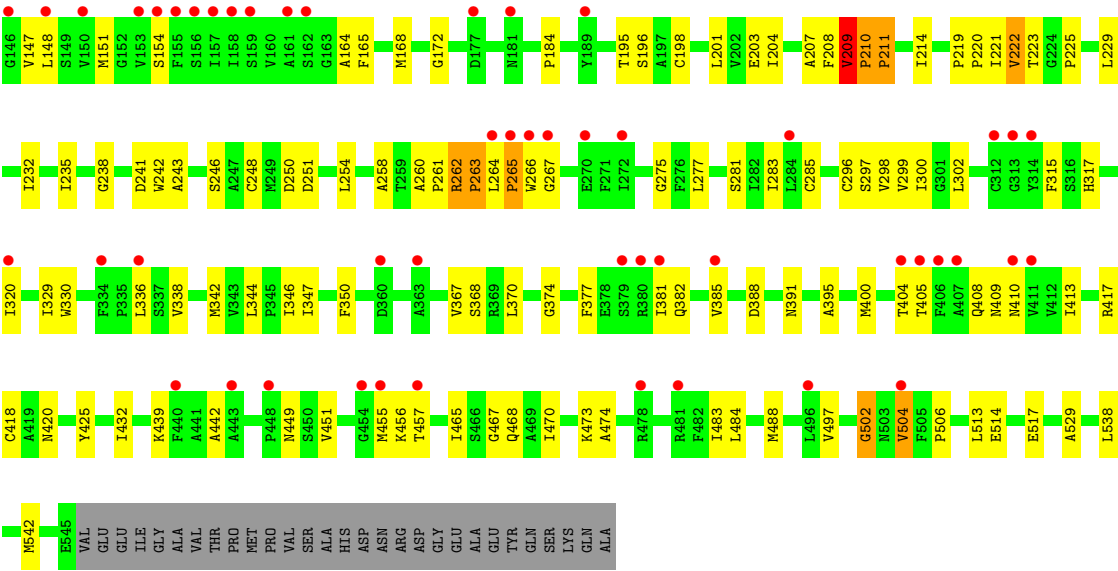
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density.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



• Molecule 1: Uric acid-xanthine permease





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	75.09Å 173.83Å 82.39Å 90.00° 111.11° 90.00°	Depositor
Resolution (Å)	76.86 – 3.70 76.86 – 3.70	Depositor EDS
% Data completeness (in resolution range)	97.7 (76.86-3.70) 97.7 (76.86-3.70)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.50 (at 3.67Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.296 , 0.327 0.303 , 0.307	Depositor DCC
R_{free} test set	1049 reflections (4.30%)	wwPDB-VP
Wilson B-factor (Å ²)	128.3	Xtriage
Anisotropy	0.600	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 177.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.42$, $\langle L^2 \rangle = 0.25$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.86	EDS
Total number of atoms	7127	wwPDB-VP
Average B, all atoms (Å ²)	164.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.83% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: XAN, 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	0.47	0/3627	0.95	10/4949 (0.2%)
1	B	0.46	0/3509	0.95	10/4787 (0.2%)
All	All	0.46	0/7136	0.95	20/9736 (0.2%)

There are no bond length outliers.

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	209	VAL	CA-C-N	-6.73	113.45	120.38
1	A	209	VAL	C-N-CA	-6.73	113.45	120.38
1	B	209	VAL	CA-C-N	-6.59	113.59	120.38
1	B	209	VAL	C-N-CA	-6.59	113.59	120.38
1	A	264	LEU	CA-C-N	-6.14	112.16	119.84
1	A	264	LEU	C-N-CA	-6.14	112.16	119.84
1	A	142	TYR	N-CA-C	6.05	118.66	109.95
1	A	505	PHE	CA-C-N	-5.92	113.66	119.64
1	A	505	PHE	C-N-CA	-5.92	113.66	119.64
1	B	542	MET	CA-C-N	5.65	125.56	119.85
1	B	542	MET	C-N-CA	5.65	125.56	119.85
1	A	262	ARG	CA-C-N	-5.59	112.86	119.84
1	A	262	ARG	C-N-CA	-5.59	112.86	119.84
1	B	262	ARG	CA-C-N	-5.48	112.99	119.84
1	B	262	ARG	C-N-CA	-5.48	112.99	119.84
1	B	439	LYS	N-CA-C	-5.41	106.81	113.41
1	B	222	VAL	N-CA-C	-5.26	107.28	111.81
1	A	439	LYS	N-CA-C	-5.23	107.03	113.41
1	B	497	VAL	CA-C-N	5.17	126.30	119.84
1	B	497	VAL	C-N-CA	5.17	126.30	119.84

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	3545	0	3646	97	0
1	B	3434	0	3537	91	0
2	A	11	0	4	3	0
2	B	11	0	4	2	0
3	A	63	0	72	5	0
3	B	63	0	73	4	0
All	All	7127	0	7336	184	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 13.

All (184) 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:223:THR:HG23	3:B:1002:LMT:H12	1.65	0.78
1:B:67:PHE:HA	1:B:417:ARG:HB3	1.67	0.76
1:A:223:THR:HG23	3:A:1002:LMT:H12	1.66	0.76
1:A:302:LEU:HD12	1:A:451:VAL:HG12	1.72	0.71
1:A:302:LEU:HB2	1:A:455:MET:HG3	1.74	0.70
1:A:502:GLY:HA2	1:A:506:PRO:HB3	1.74	0.70
1:A:151:MET:HE1	1:A:388:ASP:HA	1.73	0.69
1:B:151:MET:HE1	1:B:388:ASP:HA	1.73	0.69
1:A:260:ALA:HB3	1:A:263:PRO:HG3	1.74	0.69
1:A:529:ALA:HB1	3:A:1002:LMT:H122	1.74	0.69
1:B:502:GLY:HA2	1:B:506:PRO:HB3	1.74	0.69
1:A:72:GLU:O	1:A:420:ASN:ND2	2.25	0.68
1:B:201:LEU:HA	1:B:204:ILE:HG12	1.78	0.65
1:A:196:SER:HA	1:A:346:ILE:HG21	1.79	0.64
1:B:261:PRO:O	1:B:262:ARG:HG2	1.98	0.64
1:A:467:GLY:HA2	1:A:470:ILE:HD12	1.80	0.64
1:B:395:ALA:HB1	1:B:400:MET:HB2	1.81	0.63
1:B:260:ALA:HB3	1:B:263:PRO:HG3	1.81	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:504:VAL:HG11	1:A:513:LEU:HD23	1.82	0.62
1:B:121:ILE:HG23	1:B:432:ILE:HG23	1.82	0.61
1:B:113:GLN:HG2	1:B:442:ALA:HA	1.82	0.61
1:A:465:ILE:HD11	1:B:488:MET:HG3	1.84	0.60
1:B:126:LEU:HD13	1:B:391:ASN:HB3	1.81	0.60
1:B:88:LEU:O	1:B:299:VAL:HG21	2.02	0.60
1:B:105:LEU:HD22	1:B:165:PHE:CZ	2.37	0.59
1:A:187:GLU:HA	1:A:335:PRO:HG2	1.85	0.59
1:B:529:ALA:HB1	3:B:1002:LMT:H122	1.84	0.59
1:A:176:LEU:HD13	1:A:182:ARG:HA	1.84	0.59
1:B:164:ALA:O	1:B:168:MET:HG3	2.02	0.58
1:A:269:PRO:HA	1:A:272:ILE:HD12	1.85	0.58
1:B:66:PRO:N	1:B:417:ARG:HH11	2.01	0.58
1:B:504:VAL:HG11	1:B:513:LEU:HD23	1.85	0.58
1:B:302:LEU:HB2	1:B:455:MET:HG3	1.86	0.58
1:A:409:ASN:OD1	1:A:410:ASN:N	2.37	0.57
1:A:263:PRO:HB2	1:A:266:TRP:HB2	1.85	0.57
1:A:154:SER:HA	2:A:1001:XAN:C8	2.35	0.57
1:B:342:MET:O	1:B:346:ILE:HG12	2.05	0.57
1:A:164:ALA:O	1:A:168:MET:HG3	2.03	0.57
1:A:201:LEU:HA	1:A:204:ILE:HG12	1.86	0.57
1:B:473:LYS:NZ	3:B:1002:LMT:O2B	2.39	0.56
1:A:203:GLU:HB2	1:A:350:PHE:HB3	1.88	0.56
1:B:514:GLU:HA	1:B:517:GLU:HB3	1.88	0.56
1:B:302:LEU:HD12	1:B:451:VAL:HG12	1.87	0.55
1:A:302:LEU:HD22	1:A:455:MET:HE2	1.89	0.55
1:B:196:SER:HA	1:B:346:ILE:HG21	1.87	0.55
1:B:70:LEU:HD22	1:B:147:VAL:HG23	1.88	0.55
1:A:70:LEU:HD22	1:A:147:VAL:HG23	1.87	0.55
1:B:132:THR:HB	1:B:425:TYR:OH	2.07	0.54
1:B:121:ILE:HA	1:B:432:ILE:HA	1.90	0.54
1:B:220:PRO:HG2	1:B:474:ALA:HB2	1.88	0.54
1:B:248:CYS:HB3	1:B:258:ALA:O	2.08	0.53
1:B:203:GLU:HB2	1:B:350:PHE:HB3	1.91	0.53
1:A:473:LYS:NZ	3:A:1002:LMT:O2B	2.41	0.53
1:A:88:LEU:O	1:A:299:VAL:HG21	2.08	0.53
1:B:207:ALA:O	1:B:382:GLN:NE2	2.29	0.53
1:A:468:GLN:NE2	1:B:468:GLN:HG3	2.24	0.52
1:B:222:VAL:O	1:B:225:PRO:HD2	2.09	0.51
1:A:229:LEU:HD12	1:A:232:ILE:HB	1.92	0.51
1:A:246:SER:O	1:A:264:LEU:HA	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:285:CYS:SG	1:A:297:SER:HA	2.50	0.51
1:B:72:GLU:O	1:B:420:ASN:ND2	2.43	0.51
1:A:278:VAL:HG21	1:A:451:VAL:HA	1.93	0.51
1:A:294:LYS:C	1:A:296:CYS:H	2.19	0.51
1:B:238:GLY:HA2	1:B:456:LYS:HE3	1.93	0.50
1:B:263:PRO:HB2	1:B:266:TRP:HB2	1.93	0.50
1:B:298:VAL:HG21	3:B:1003:LMT:H6'1	1.94	0.50
1:A:70:LEU:HD21	1:A:370:LEU:HD11	1.94	0.49
1:B:241:ASP:HB3	1:B:449:ASN:O	2.11	0.49
1:B:408:GLN:NE2	2:B:1001:XAN:O2	2.45	0.49
1:A:156:SER:OG	1:A:349:VAL:HG13	2.12	0.49
1:A:514:GLU:HA	1:A:517:GLU:HB3	1.94	0.49
1:B:484:LEU:O	1:B:488:MET:HG2	2.13	0.49
1:A:395:ALA:HB1	1:A:400:MET:HB2	1.94	0.48
1:A:371:GLU:HB2	1:A:380:ARG:HH22	1.78	0.48
1:B:243:ALA:O	1:B:267:GLY:HA3	2.14	0.48
1:A:264:LEU:HB2	1:A:265:PRO:HD3	1.95	0.48
1:A:68:PHE:HB3	1:A:420:ASN:OD1	2.13	0.48
1:A:136:ILE:HG13	1:A:143:ILE:HB	1.96	0.48
1:B:219:PRO:HG2	1:B:221:ILE:HG22	1.95	0.48
1:A:126:LEU:HB3	1:A:391:ASN:CG	2.39	0.47
1:A:517:GLU:O	1:A:521:GLU:HG2	2.14	0.47
1:B:67:PHE:HZ	1:B:368:SER:HA	1.80	0.47
1:A:248:CYS:HB3	1:A:258:ALA:O	2.14	0.47
1:A:144:GLY:O	1:A:383:GLY:HA3	2.15	0.47
1:A:468:GLN:HG3	1:B:468:GLN:NE2	2.29	0.47
1:A:298:VAL:HG21	3:A:1003:LMT:H6'1	1.97	0.47
3:A:1002:LMT:O2B	3:A:1002:LMT:H6E	2.13	0.47
1:B:261:PRO:C	1:B:263:PRO:HD3	2.40	0.47
1:B:220:PRO:HA	1:B:223:THR:HB	1.96	0.47
1:A:210:PRO:O	1:A:214:ILE:HG13	2.14	0.47
1:A:154:SER:HA	2:A:1001:XAN:H8	1.97	0.47
1:B:70:LEU:HD21	1:B:370:LEU:HD11	1.96	0.47
1:B:296:CYS:O	1:B:300:ILE:N	2.35	0.47
1:B:329:ILE:HB	1:B:330:TRP:CE3	2.50	0.47
1:A:207:ALA:O	1:A:382:GLN:NE2	2.46	0.46
1:B:105:LEU:O	1:B:184:PRO:HA	2.14	0.46
1:B:107:LEU:H	1:B:108:PRO:CD	2.28	0.46
1:B:483:ILE:HG12	1:B:538:LEU:HB3	1.97	0.46
1:A:210:PRO:HG2	1:A:214:ILE:HD11	1.97	0.46
1:A:143:ILE:HG21	1:A:390:ILE:HD12	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:242:TRP:HZ3	1:B:457:THR:HG21	1.80	0.46
1:A:226:THR:HG23	1:A:528:PHE:HB2	1.97	0.46
1:A:176:LEU:HD12	1:A:177:ASP:H	1.80	0.46
1:B:246:SER:O	1:B:264:LEU:HA	2.16	0.46
1:A:132:THR:HB	1:A:425:TYR:OH	2.15	0.46
1:B:70:LEU:HD11	1:B:370:LEU:HD11	1.97	0.46
1:B:223:THR:HG22	1:B:470:ILE:HG21	1.97	0.46
1:B:413:ILE:HG23	1:B:418:CYS:O	2.16	0.46
1:A:306:CYS:HB2	1:A:451:VAL:HG11	1.97	0.45
1:A:501:PHE:O	1:A:503:ASN:N	2.50	0.45
1:B:467:GLY:HA2	1:B:470:ILE:HD12	1.97	0.45
1:B:242:TRP:CD1	1:B:275:GLY:HA3	2.51	0.45
1:B:126:LEU:HD22	1:B:391:ASN:OD1	2.16	0.45
1:B:385:VAL:O	1:B:388:ASP:HB3	2.17	0.45
1:A:230:ILE:HG12	1:A:528:PHE:CD2	2.52	0.45
1:A:277:LEU:HA	1:A:277:LEU:HD23	1.65	0.45
1:A:482:PHE:HZ	1:B:283:ILE:HA	1.82	0.45
1:B:409:ASN:OD1	1:B:410:ASN:N	2.50	0.44
1:B:168:MET:HA	1:B:172:GLY:HA3	1.99	0.44
1:B:248:CYS:SG	1:B:263:PRO:HG2	2.57	0.44
1:A:113:GLN:HG2	1:A:442:ALA:HA	2.00	0.44
1:A:154:SER:HA	2:A:1001:XAN:N9	2.33	0.44
1:A:378:GLU:O	1:A:382:GLN:HG2	2.18	0.44
1:A:241:ASP:HB3	1:A:449:ASN:O	2.17	0.44
1:A:412:VAL:O	1:A:416:THR:N	2.51	0.44
1:A:229:LEU:HD23	1:A:527:GLY:C	2.42	0.44
1:B:154:SER:HB2	1:B:404:THR:HG22	1.99	0.44
1:A:241:ASP:OD2	1:A:456:LYS:NZ	2.32	0.43
1:A:404:THR:OG1	1:A:405:THR:N	2.51	0.43
1:A:154:SER:HB2	1:A:404:THR:HG22	1.99	0.43
1:A:196:SER:HB2	1:A:402:PRO:HG3	1.98	0.43
1:B:285:CYS:SG	1:B:297:SER:HA	2.57	0.43
1:A:229:LEU:HD12	1:A:229:LEU:HA	1.85	0.43
1:A:133:ARG:NE	1:A:146:GLY:HA3	2.34	0.43
1:A:223:THR:HG22	1:A:470:ILE:HG21	2.00	0.43
1:A:261:PRO:C	1:A:263:PRO:HD3	2.44	0.43
1:B:277:LEU:O	1:B:281:SER:OG	2.19	0.43
1:A:278:VAL:HG22	1:A:305:GLY:HA3	1.99	0.43
1:B:264:LEU:HB2	1:B:265:PRO:HD3	2.01	0.43
1:B:277:LEU:HD23	1:B:277:LEU:HA	1.81	0.43
1:A:161:ALA:HA	1:A:164:ALA:HB3	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:300:ILE:O	1:A:304:VAL:HG23	2.19	0.43
1:A:75:PRO:O	1:A:78:LEU:N	2.52	0.43
1:A:148:LEU:HD23	1:A:148:LEU:HA	1.89	0.43
1:B:210:PRO:HG2	1:B:214:ILE:HD11	2.00	0.43
1:A:479:ARG:NH2	1:A:539:ASN:O	2.51	0.42
1:B:208:PHE:C	1:B:210:PRO:HD3	2.44	0.42
1:B:229:LEU:HD12	1:B:229:LEU:HA	1.91	0.42
1:B:254:LEU:HD12	1:B:315:PHE:O	2.19	0.42
1:B:232:ILE:O	1:B:235:ILE:HG22	2.19	0.42
1:A:226:THR:OG1	1:A:532:ALA:HB2	2.20	0.42
1:B:210:PRO:N	1:B:211:PRO:HD3	2.35	0.42
1:B:381:ILE:O	1:B:385:VAL:HG23	2.19	0.42
1:A:105:LEU:HD11	1:A:185:CYS:HB3	2.02	0.42
1:A:127:SER:O	1:A:131:ILE:HG13	2.19	0.42
1:B:209:VAL:N	1:B:210:PRO:HD3	2.35	0.42
1:A:105:LEU:HD22	1:A:165:PHE:CZ	2.55	0.42
1:B:317:HIS:HA	1:B:320:ILE:HD12	2.02	0.42
1:A:71:ASN:CG	1:A:421:ARG:HH12	2.28	0.42
1:A:385:VAL:O	1:A:388:ASP:HB3	2.19	0.42
1:B:344:LEU:HA	1:B:347:ILE:HD12	2.00	0.42
1:B:221:ILE:HD12	1:B:221:ILE:HA	1.88	0.41
1:B:250:ASP:OD1	1:B:251:ASP:N	2.44	0.41
1:A:542:MET:HE3	1:A:543:PRO:HD2	2.02	0.41
1:B:456:LYS:HE3	1:B:456:LYS:HB2	1.81	0.41
1:A:150:VAL:HG12	1:A:406:PHE:HB2	2.02	0.41
1:A:396:ALA:HA	1:A:400:MET:O	2.21	0.41
1:A:467:GLY:O	1:A:471:VAL:HG23	2.20	0.41
1:B:210:PRO:O	1:B:214:ILE:HG13	2.21	0.41
1:A:374:GLY:O	1:A:377:PHE:HB3	2.21	0.41
1:B:154:SER:HA	2:B:1001:XAN:C8	2.51	0.41
1:B:195:THR:O	1:B:198:CYS:HB3	2.21	0.41
1:A:123:CYS:SG	1:A:152:GLY:HA2	2.60	0.41
1:A:515:GLY:O	1:A:518:ASN:HB2	2.21	0.41
1:B:336:LEU:O	1:B:338:VAL:HG23	2.21	0.41
1:A:377:PHE:O	1:A:380:ARG:HB2	2.21	0.40
1:A:468:GLN:O	1:A:471:VAL:HB	2.21	0.40
1:B:367:VAL:HG21	1:B:410:ASN:HB3	2.03	0.40
1:B:374:GLY:O	1:B:377:PHE:HB3	2.21	0.40
1:A:232:ILE:O	1:A:235:ILE:HG22	2.21	0.40
1:A:484:LEU:HB3	1:B:465:ILE:HD12	2.03	0.40
1:A:209:VAL:H	1:A:382:GLN:NE2	2.19	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:518:ASN:O	1:A:521:GLU:HB2	2.21	0.40
1:B:404:THR:OG1	1:B:405:THR:N	2.54	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	478/574 (83%)	444 (93%)	25 (5%)	9 (2%)	6	33
1	B	464/574 (81%)	437 (94%)	20 (4%)	7 (2%)	8	36
All	All	942/1148 (82%)	881 (94%)	45 (5%)	16 (2%)	7	34

All (16) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	210	PRO
1	B	210	PRO
1	B	148	LEU
1	A	139	THR
1	A	148	LEU
1	A	211	PRO
1	A	263	PRO
1	A	265	PRO
1	A	502	GLY
1	B	211	PRO
1	B	263	PRO
1	B	265	PRO
1	A	179	ALA
1	B	107	LEU
1	A	107	LEU
1	B	502	GLY

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	382/464 (82%)	378 (99%)	4 (1%)	68	74
1	B	371/464 (80%)	369 (100%)	2 (0%)	81	80
All	All	753/928 (81%)	747 (99%)	6 (1%)	73	76

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	141	TYR
1	A	209	VAL
1	A	466	SER
1	A	504	VAL
1	B	209	VAL
1	B	504	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	382	GLN

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

6 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	XAN	A	1001	-	12,12,12	1.08	1 (8%)	14,17,17	2.56	6 (42%)
3	LMT	B	1003	-	29,29,36	1.27	5 (17%)	40,40,47	1.12	2 (5%)
2	XAN	B	1001	-	12,12,12	1.11	1 (8%)	14,17,17	2.29	6 (42%)
3	LMT	A	1002	-	36,36,36	1.17	5 (13%)	47,47,47	1.02	4 (8%)
3	LMT	A	1003	-	29,29,36	1.26	4 (13%)	40,40,47	0.96	2 (5%)
3	LMT	B	1002	-	36,36,36	1.05	2 (5%)	47,47,47	1.08	4 (8%)

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	XAN	A	1001	-	-	-	0/2/2/2
3	LMT	B	1003	-	-	8/14/54/61	0/2/2/2
2	XAN	B	1001	-	-	-	0/2/2/2
3	LMT	A	1002	-	-	11/21/61/61	0/2/2/2
3	LMT	A	1003	-	-	9/14/54/61	0/2/2/2
3	LMT	B	1002	-	-	12/21/61/61	0/2/2/2

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1003	LMT	O3'-C3'	-2.63	1.36	1.43
3	A	1002	LMT	O3'-C3'	-2.58	1.36	1.43
3	A	1003	LMT	O2'-C2'	-2.58	1.36	1.43
3	B	1002	LMT	O3'-C3'	-2.49	1.36	1.43
3	B	1003	LMT	O3'-C3'	-2.48	1.36	1.43
3	A	1002	LMT	O3B-C3B	-2.25	1.37	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1002	LMT	O2'-C2'	-2.22	1.37	1.43
3	B	1003	LMT	O2'-C2'	-2.18	1.37	1.43
2	A	1001	XAN	C6-N1	-2.16	1.34	1.38
3	A	1002	LMT	O4'-C4B	-2.15	1.37	1.43
2	B	1001	XAN	C6-N1	-2.11	1.34	1.38
3	B	1003	LMT	O4'-C4B	-2.10	1.37	1.43
3	A	1003	LMT	O1'-C1'	-2.09	1.36	1.40
3	A	1002	LMT	O2B-C2B	-2.07	1.37	1.43
3	B	1003	LMT	O3B-C3B	-2.02	1.37	1.43
3	B	1002	LMT	O2'-C2'	-2.02	1.38	1.43
3	B	1003	LMT	O2B-C2B	-2.02	1.38	1.43
3	A	1003	LMT	O2B-C2B	-2.01	1.38	1.43

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1001	XAN	N9-C8-N7	-5.26	106.88	112.98
2	A	1001	XAN	C6-N1-C2	-4.73	119.85	126.37
2	B	1001	XAN	C6-N1-C2	-4.55	120.10	126.37
2	B	1001	XAN	N9-C8-N7	-4.24	108.06	112.98
2	A	1001	XAN	C8-N7-C5	3.57	109.20	104.38
3	B	1003	LMT	C1'-O5'-C5'	-3.21	107.46	113.72
3	B	1002	LMT	O5B-C5B-C4B	3.16	115.40	109.70
2	A	1001	XAN	N1-C2-N3	2.98	120.42	115.74
3	B	1002	LMT	C3'-C4'-C5'	-2.94	104.42	110.93
2	B	1001	XAN	C5-C6-N1	2.77	120.30	113.25
2	B	1001	XAN	C8-N7-C5	2.71	108.04	104.38
2	B	1001	XAN	O6-C6-C5	-2.70	119.41	126.53
3	A	1003	LMT	C3'-C4'-C5'	-2.68	104.98	110.93
3	A	1002	LMT	O5B-C5B-C4B	2.55	114.30	109.70
2	A	1001	XAN	C5-C6-N1	2.48	119.56	113.25
3	B	1003	LMT	C3'-C4'-C5'	-2.44	105.52	110.93
2	B	1001	XAN	N1-C2-N3	2.40	119.52	115.74
3	A	1002	LMT	C3'-C4'-C5'	-2.32	105.79	110.93
2	A	1001	XAN	O6-C6-C5	-2.25	120.59	126.53
3	A	1002	LMT	C1'-O5'-C5'	-2.25	109.33	113.72
3	B	1002	LMT	C1'-O5'-C5'	-2.23	109.36	113.72
3	B	1002	LMT	O1'-C1'-C2'	2.21	111.64	108.27
3	A	1002	LMT	C6B-C5B-C4B	-2.20	107.62	113.02
3	A	1003	LMT	O5'-C1'-O1'	-2.03	105.24	110.04

There are no chirality outliers.

All (40) torsion outliers are listed below:

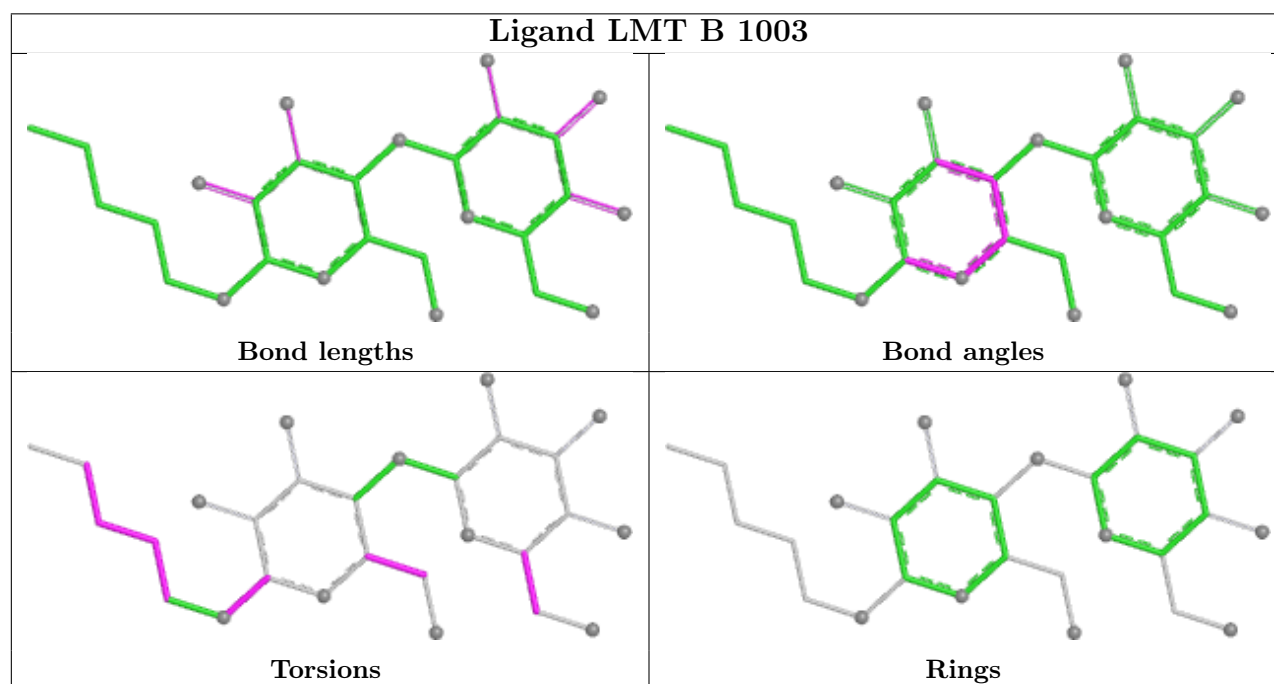
Mol	Chain	Res	Type	Atoms
3	B	1002	LMT	C3'-C4'-O1B-C1B
3	A	1003	LMT	C4'-C5'-C6'-O6'
3	A	1002	LMT	C3'-C4'-O1B-C1B
3	A	1003	LMT	O5'-C5'-C6'-O6'
3	A	1003	LMT	O5B-C5B-C6B-O6B
3	B	1003	LMT	O5'-C5'-C6'-O6'
3	A	1002	LMT	C4B-C5B-C6B-O6B
3	B	1002	LMT	C4B-C5B-C6B-O6B
3	A	1002	LMT	O5'-C1'-O1'-C1
3	B	1002	LMT	O5'-C1'-O1'-C1
3	A	1002	LMT	C2'-C1'-O1'-C1
3	B	1002	LMT	C2'-C1'-O1'-C1
3	A	1003	LMT	C4B-C5B-C6B-O6B
3	A	1002	LMT	O5B-C5B-C6B-O6B
3	B	1003	LMT	C4'-C5'-C6'-O6'
3	B	1002	LMT	O5B-C5B-C6B-O6B
3	B	1002	LMT	O1'-C1-C2-C3
3	A	1002	LMT	O1'-C1-C2-C3
3	B	1003	LMT	C2'-C1'-O1'-C1
3	B	1003	LMT	O1'-C1-C2-C3
3	B	1003	LMT	C1-C2-C3-C4
3	B	1003	LMT	O5'-C1'-O1'-C1
3	A	1002	LMT	C5'-C4'-O1B-C1B
3	A	1003	LMT	O1'-C1-C2-C3
3	A	1002	LMT	C1-C2-C3-C4
3	B	1002	LMT	C1-C2-C3-C4
3	A	1002	LMT	C2-C3-C4-C5
3	A	1003	LMT	O5'-C1'-O1'-C1
3	B	1002	LMT	C2-C3-C4-C5
3	A	1002	LMT	C11-C10-C9-C8
3	B	1002	LMT	C6-C7-C8-C9
3	A	1003	LMT	C1-C2-C3-C4
3	B	1003	LMT	O5B-C5B-C6B-O6B
3	A	1003	LMT	C2-C3-C4-C5
3	B	1002	LMT	C5'-C4'-O1B-C1B
3	A	1002	LMT	C4-C5-C6-C7
3	A	1003	LMT	C2'-C1'-O1'-C1
3	B	1003	LMT	C2-C3-C4-C5
3	B	1002	LMT	O5'-C5'-C6'-O6'
3	B	1002	LMT	C3-C4-C5-C6

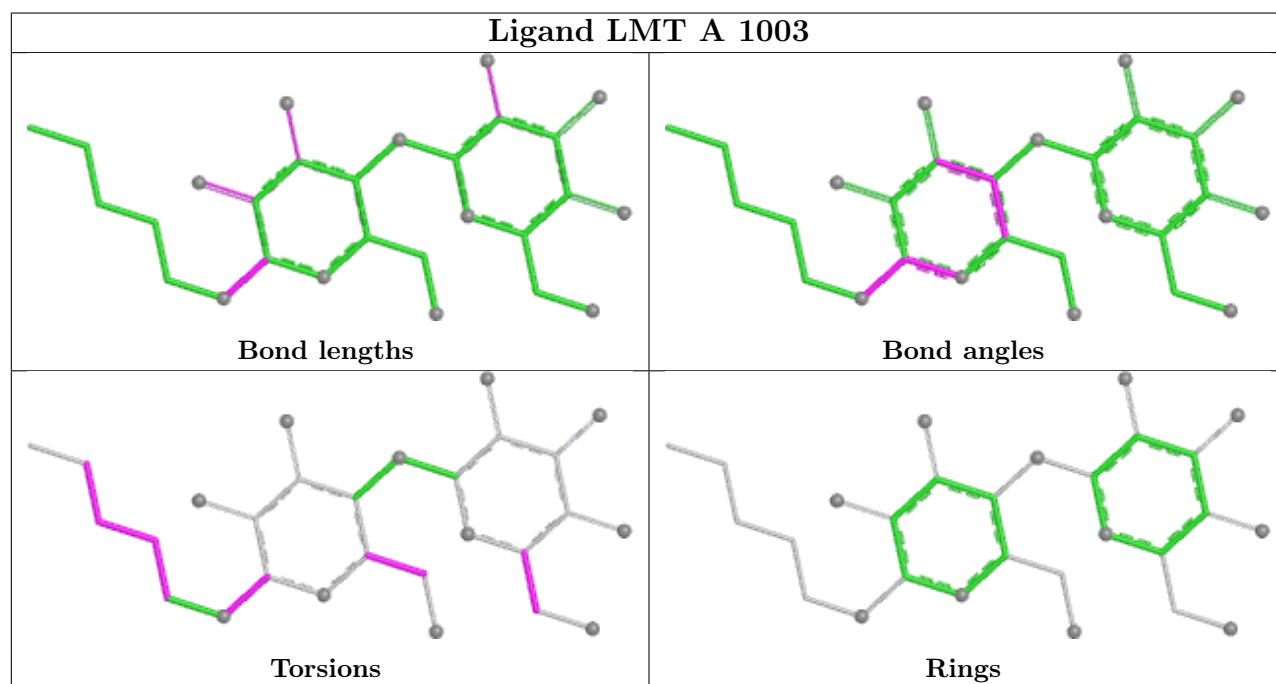
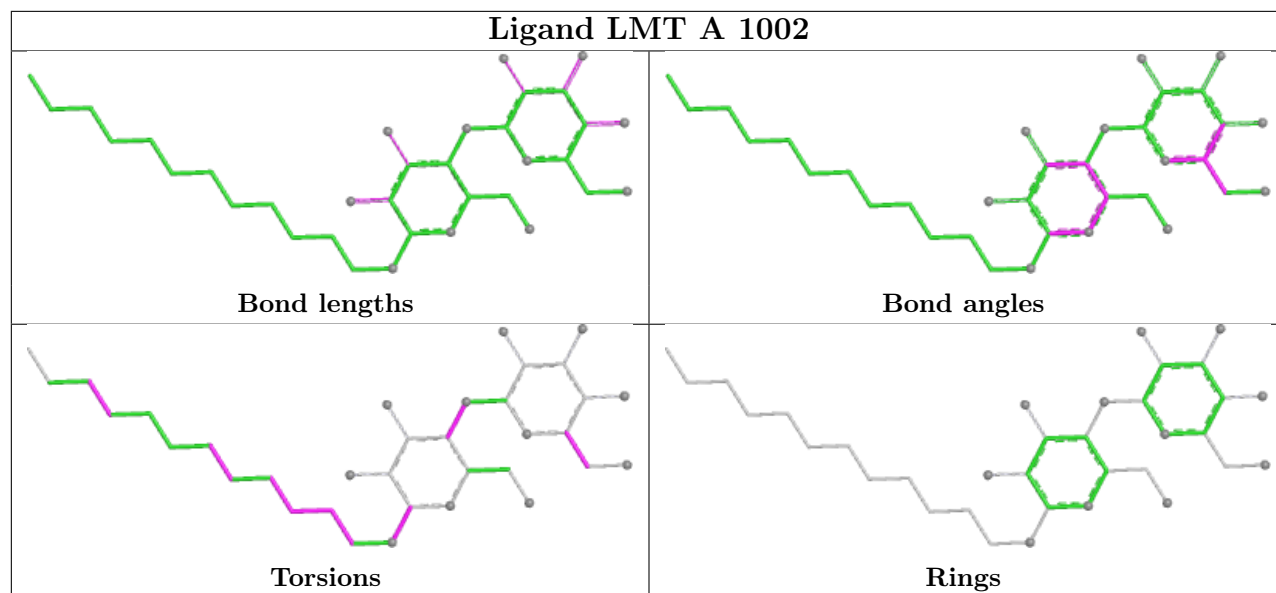
There are no ring outliers.

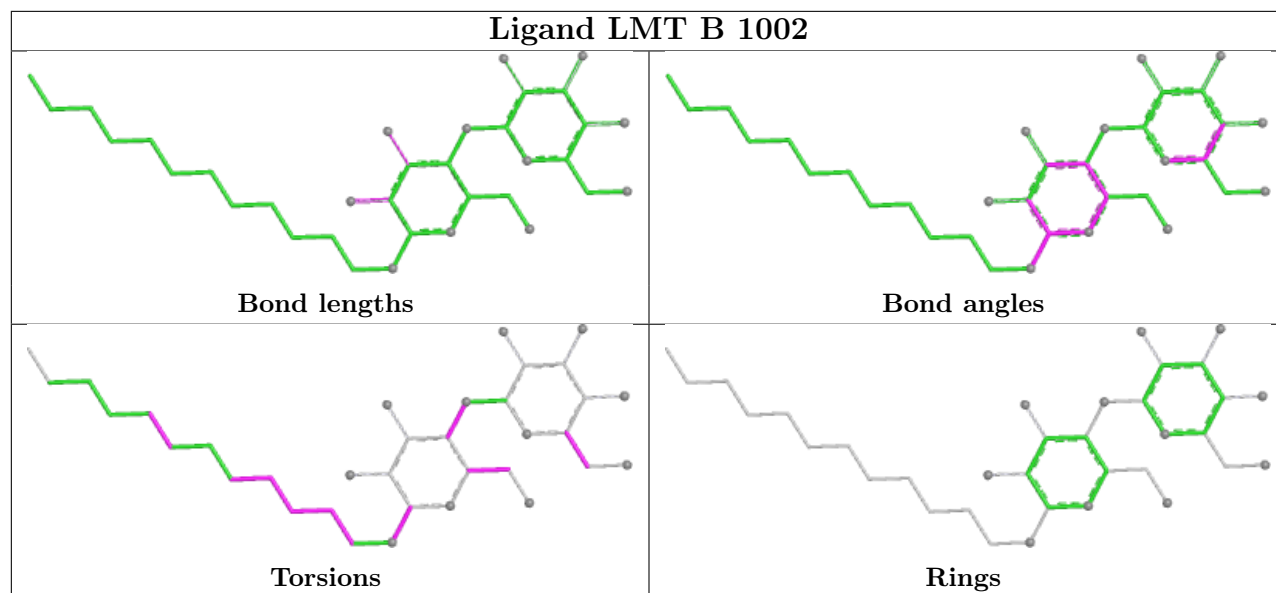
6 monomers are involved in 14 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1001	XAN	3	0
3	B	1003	LMT	1	0
2	B	1001	XAN	2	0
3	A	1002	LMT	4	0
3	A	1003	LMT	1	0
3	B	1002	LMT	3	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	480/574 (83%)	0.61	57 (11%) 9 10	115, 146, 187, 212	0
1	B	468/574 (81%)	0.92	71 (15%) 5 7	127, 177, 223, 257	0
All	All	948/1148 (82%)	0.76	128 (13%) 7 8	115, 158, 212, 257	0

All (128) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	158	ILE	10.1
1	B	97	PRO	9.3
1	A	93	GLY	9.1
1	B	406	PHE	9.0
1	B	266	TRP	8.7
1	A	149	SER	7.9
1	B	478	ARG	7.9
1	B	443	ALA	7.7
1	B	91	LEU	7.7
1	B	100	ILE	7.6
1	B	98	PRO	7.5
1	B	157	ILE	7.4
1	B	156	SER	6.9
1	A	91	LEU	6.9
1	A	94	VAL	6.8
1	B	101	ILE	6.7
1	B	94	VAL	6.5
1	A	95	VAL	6.4
1	A	96	THR	6.3
1	B	155	PHE	6.1
1	B	88	LEU	5.8
1	B	102	SER	5.8
1	B	96	THR	5.6
1	A	92	ALA	5.6

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Mol	Chain	Res	Type	RSRZ
1	A	208	PHE	5.5
1	B	154	SER	5.5
1	B	93	GLY	5.5
1	B	95	VAL	5.5
1	A	265	PRO	5.5
1	A	97	PRO	5.2
1	A	146	GLY	5.2
1	B	312	CYS	5.1
1	B	267	GLY	5.0
1	A	136	ILE	4.7
1	A	504	VAL	4.7
1	A	533	PHE	4.4
1	B	104	SER	4.3
1	B	407	ALA	4.3
1	B	265	PRO	4.2
1	B	360	ASP	4.1
1	A	305	GLY	4.1
1	B	411	VAL	4.0
1	A	440	PHE	3.9
1	B	112	GLN	3.9
1	B	504	VAL	3.9
1	B	481	ARG	3.9
1	A	141	TYR	3.9
1	A	443	ALA	3.8
1	B	153	VAL	3.8
1	A	537	LEU	3.7
1	A	103	SER	3.7
1	B	92	ALA	3.6
1	B	181	ASN	3.6
1	A	135	HIS	3.6
1	A	304	VAL	3.5
1	B	162	SER	3.5
1	B	90	MET	3.5
1	B	89	ALA	3.5
1	B	314	TYR	3.5
1	A	100	ILE	3.5
1	B	177	ASP	3.4
1	B	68	PHE	3.4
1	B	313	GLY	3.3
1	B	85	GLN	3.3
1	A	385	VAL	3.3
1	B	334	PHE	3.3

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Mol	Chain	Res	Type	RSRZ
1	A	70	LEU	3.3
1	B	159	SER	3.3
1	B	440	PHE	3.3
1	A	461	ALA	3.3
1	A	344	LEU	3.2
1	B	448	PRO	3.2
1	A	538	LEU	3.2
1	A	444	ILE	3.1
1	A	277	LEU	3.1
1	B	161	ALA	3.1
1	B	380	ARG	3.1
1	A	148	LEU	3.1
1	B	496	LEU	3.0
1	A	98	PRO	3.0
1	A	153	VAL	3.0
1	B	385	VAL	3.0
1	B	189	TYR	2.9
1	B	284	LEU	2.9
1	A	104	SER	2.9
1	A	451	VAL	2.8
1	A	266	TRP	2.8
1	B	146	GLY	2.8
1	B	103	SER	2.8
1	B	72	GLU	2.8
1	A	411	VAL	2.7
1	A	295	SER	2.7
1	A	102	SER	2.6
1	A	158	ILE	2.6
1	B	320	ILE	2.6
1	A	341	PRO	2.6
1	A	544	ALA	2.6
1	A	476	PHE	2.6
1	B	363	ALA	2.6
1	A	407	ALA	2.5
1	B	404	THR	2.5
1	A	147	VAL	2.5
1	A	86	HIS	2.4
1	A	275	GLY	2.4
1	B	272	ILE	2.4
1	A	446	ALA	2.4
1	B	336	LEU	2.4
1	A	316	SER	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	410	ASN	2.4
1	B	270	GLU	2.4
1	A	534	VAL	2.4
1	B	82	LEU	2.4
1	B	264	LEU	2.4
1	A	155	PHE	2.3
1	B	379	SER	2.3
1	A	274	LEU	2.2
1	B	148	LEU	2.2
1	A	450	SER	2.2
1	A	99	LEU	2.2
1	B	381	ILE	2.2
1	A	454	GLY	2.2
1	B	455	MET	2.2
1	A	278	VAL	2.1
1	B	454	GLY	2.1
1	B	405	THR	2.1
1	B	150	VAL	2.1
1	A	319	ASP	2.1
1	B	457	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	XAN	A	1001	11/11	0.78	0.13	119,119,129,129	0
3	LMT	A	1003	28/35	0.88	0.19	115,142,162,169	0
3	LMT	B	1002	35/35	0.88	0.21	160,195,206,208	0

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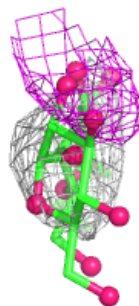
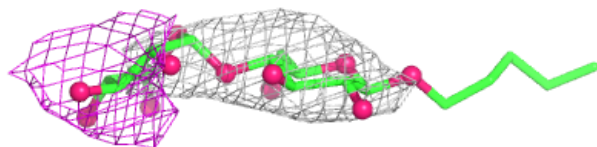
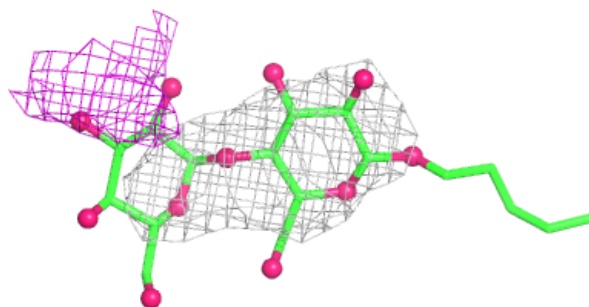
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	XAN	B	1001	11/11	0.92	0.10	158,160,161,162	0
3	LMT	A	1002	35/35	0.95	0.13	122,149,162,162	0
3	LMT	B	1003	28/35	0.95	0.11	136,151,167,169	0

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.

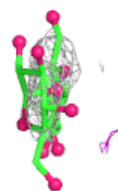
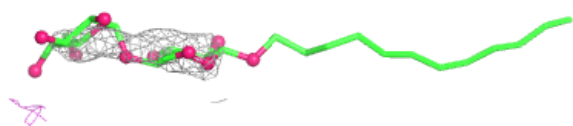
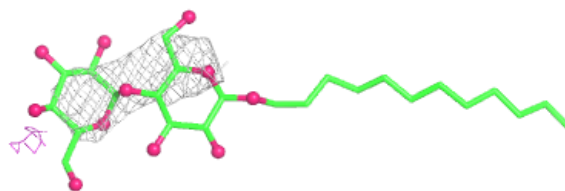
Electron density around LMT A 1003:

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

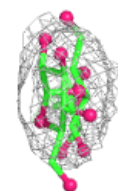
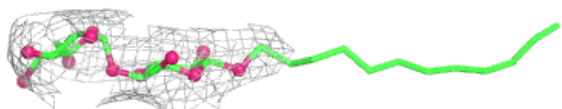
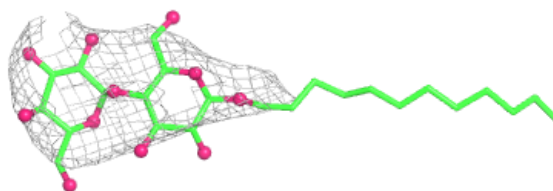


Electron density around LMT B 1002:

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

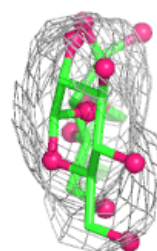
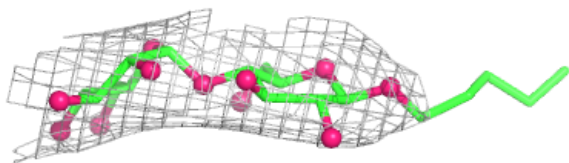
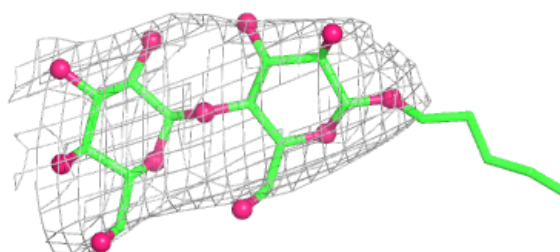
**Electron density around LMT A 1002:**

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



Electron density around LMT B 1003:

$2mF_o-DF_c$ (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.