



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 03:22 PM UTC

PDB ID : 3AEL / pdb_00003ael
Title : Reaction intermediate structure of Entamoeba histolytica methionine gamma-lyase 1 containing methionine imine-pyridoxamine-5'-phosphate and alpha-amino-alpha, beta-butenic acid-pyridoxal-5'-phosphate
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Deposited on : 2010-02-10
Resolution : 2.00 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

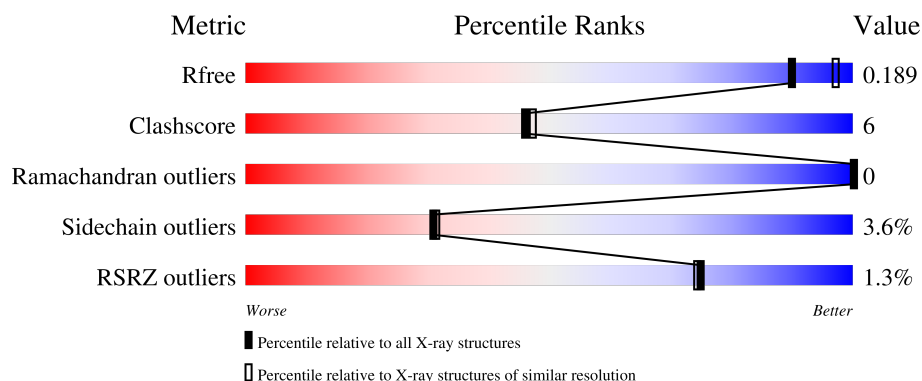
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	389	2% 86% 12% ..
1	B	389	% 87% 11% ...
1	C	389	2% 86% 13% .
1	D	389	% 84% 13% ..

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	4LM	A	2001[B]	-	-	X	-

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 12836 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.

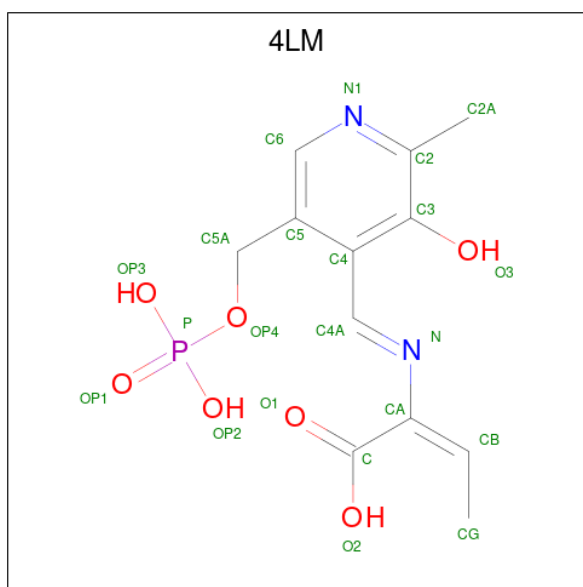
- Molecule 1 is a protein called Methionine gamma-lyase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	387	Total	C	N	O	S	0	1	0
			2952	1878	497	553	24			
1	B	386	Total	C	N	O	S	0	1	0
			2937	1867	495	551	24			
1	C	387	Total	C	N	O	S	0	1	0
			2952	1878	497	553	24			
1	D	384	Total	C	N	O	S	0	1	0
			2923	1859	492	548	24			

There are 4 discrepancies between the modelled and reference sequences:

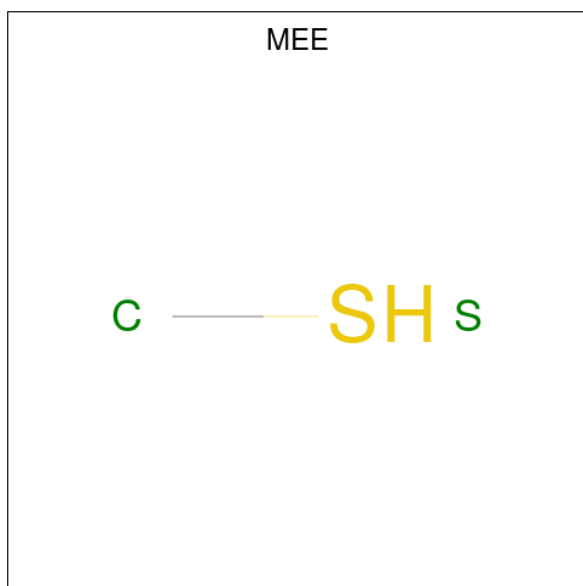
Chain	Residue	Modelled	Actual	Comment	Reference
A	308	LEU	SER	SEE REMARK 999	UNP Q86D28
B	808	LEU	SER	SEE REMARK 999	UNP Q86D28
C	1308	LEU	SER	SEE REMARK 999	UNP Q86D28
D	1808	LEU	SER	SEE REMARK 999	UNP Q86D28

- Molecule 2 is (2E)-2-{{[(1E)-{3-hydroxy-2-methyl-5-[(phosphonoxy)methyl]pyridin-4-yl}methylidene]amino}but-2-enoic acid (CCD ID: 4LM) (formula: C₁₂H₁₅N₂O₇P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	1
			39	23	3	11	2		

- Molecule 3 is METHANETHIOL (CCD ID: MEE) (formula: CH₄S).



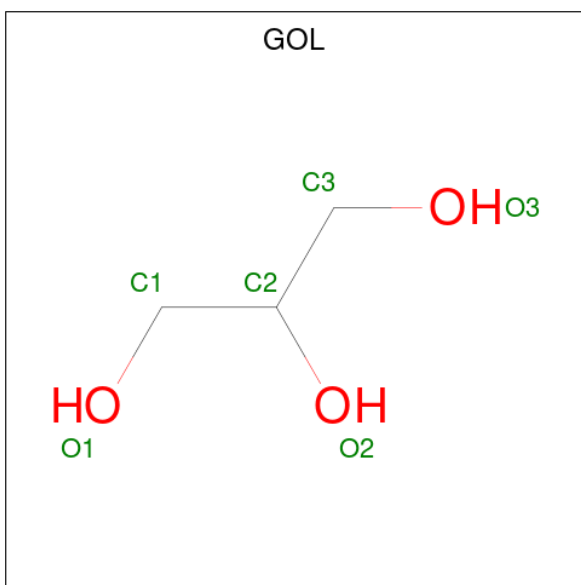
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	S	0	0
			2	1	1		

- Molecule 4 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



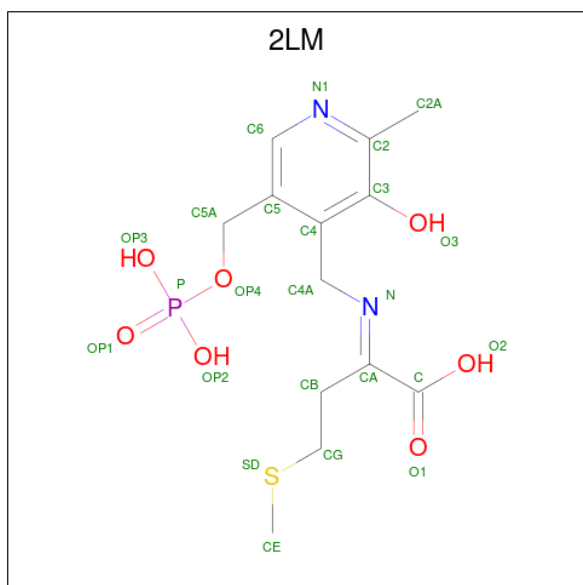
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		
4	C	1	Total	O	S	0	0
			5	4	1		
4	D	1	Total	O	S	0	0
			5	4	1		

- Molecule 5 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



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

- Molecule 6 is (2E)-2-[(3-hydroxy-2-methyl-5-[(phosphonooxy)methyl]pyridin-4-yl)methyl]imino-4-(methylsulfanyl)butanoic acid (CCD ID: 2LM) (formula: C₁₃H₁₉N₂O₇PS).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
6	B	1	Total	C	N	O	P	S	0	0
			24	13	2	7	1	1		
6	C	1	Total	C	N	O	P	S	0	0
			24	13	2	7	1	1		
6	D	1	Total	C	N	O	P	S	0	0
			24	13	2	7	1	1		

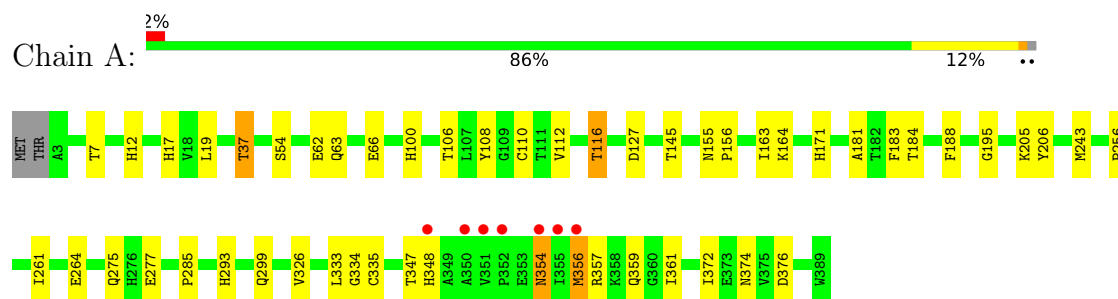
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	291	Total	O	0	0
			291	291		
7	B	185	Total	O	0	0
			185	185		
7	C	300	Total	O	0	0
			300	300		
7	D	151	Total	O	0	0
			151	151		

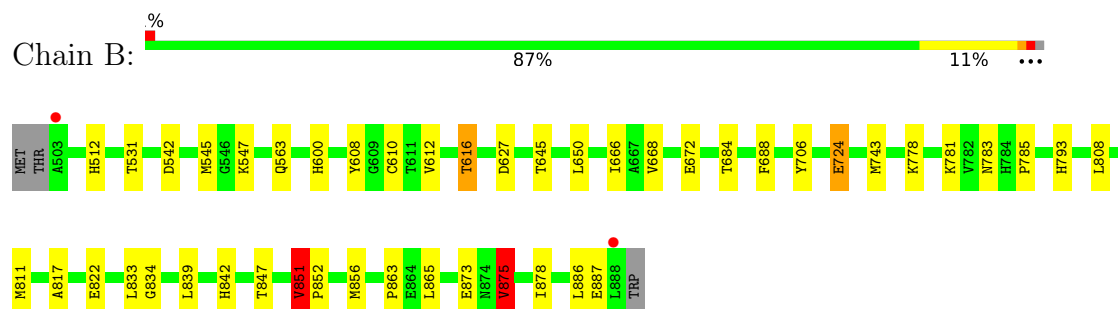
3 Residue-property plots [i](#)

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.

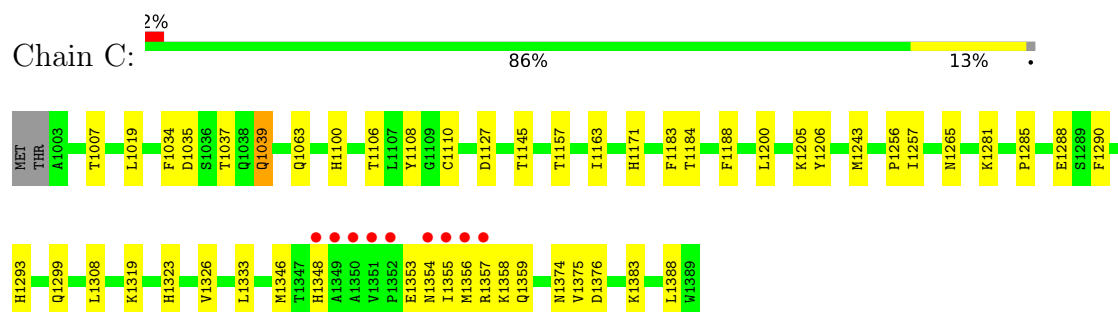
- Molecule 1: Methionine gamma-lyase



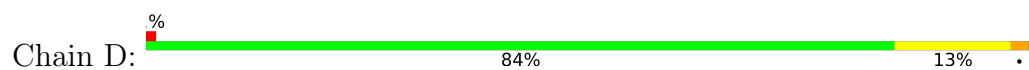
- Molecule 1: Methionine gamma-lyase

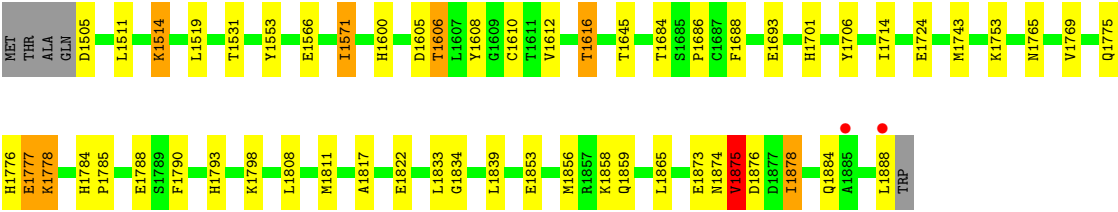


- Molecule 1: Methionine gamma-lyase



- Molecule 1: Methionine gamma-lyase





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	98.97Å 85.22Å 114.25Å 90.00° 101.98° 90.00°	Depositor
Resolution (Å)	34.30 – 2.00 34.30 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.6 (34.30-2.00) 99.7 (34.30-2.00)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.54 (at 2.00Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.153 , 0.188 0.154 , 0.189	Depositor DCC
R_{free} test set	6294 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	23.5	Xtriage
Anisotropy	0.242	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 47.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	12836	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.60% 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: 2LM, MEE, 4LM, GOL, SO4

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	0.95	1/3018 (0.0%)	0.93	2/4080 (0.0%)
1	B	0.82	0/3001	0.92	3/4057 (0.1%)
1	C	0.99	1/3018 (0.0%)	0.91	1/4080 (0.0%)
1	D	0.76	0/2987	0.86	3/4038 (0.1%)
All	All	0.89	2/12024 (0.0%)	0.90	9/16255 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	1346	MET	CA-C	8.35	1.56	1.52
1	A	181	ALA	CA-CB	5.55	1.61	1.53

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	875	VAL	CB-CA-C	-7.02	101.80	112.05
1	D	1875	VAL	CB-CA-C	-6.24	103.72	112.14
1	B	851	VAL	CB-CA-C	5.78	117.13	110.89
1	C	1206	TYR	N-CA-C	5.75	117.55	111.28
1	A	372	ILE	N-CA-C	5.18	117.16	112.29
1	D	1784	HIS	CA-C-N	5.13	124.79	119.56
1	D	1784	HIS	C-N-CA	5.13	124.79	119.56
1	A	326	VAL	N-CA-C	-5.12	106.81	111.67
1	B	851	VAL	N-CA-CB	5.03	114.76	110.08

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	2952	0	2957	36	0
1	B	2937	0	2947	32	0
1	C	2952	0	2957	38	0
1	D	2923	0	2934	40	0
2	A	39	0	22	10	0
3	A	2	0	0	1	0
4	A	5	0	0	0	0
4	B	5	0	0	0	0
4	C	5	0	0	0	0
4	D	5	0	0	0	0
5	A	6	0	8	0	0
5	C	6	0	8	0	0
6	B	24	0	15	2	0
6	C	24	0	16	4	0
6	D	24	0	15	0	0
7	A	291	0	0	6	0
7	B	185	0	0	3	0
7	C	300	0	0	8	0
7	D	151	0	0	5	0
All	All	12836	0	11879	139	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 6.

All (139) 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:D:1505:ASP:HB2	7:D:3136:HOH:O	1.65	0.96
1:D:1765:ASN:ND2	1:D:1875:VAL:HG22	1.83	0.93
2:A:2001[B]:4LM:CG	2:A:2001[B]:4LM:H4A	2.03	0.89
1:B:778:LYS:HE3	1:B:886:LEU:O	1.76	0.85
1:B:822:GLU:OE2	1:C:1037:THR:HG22	1.77	0.84
1:A:293:HIS:HE1	7:A:3362:HOH:O	1.62	0.82
1:B:612:VAL:O	1:B:616:THR:HB	1.84	0.77
1:C:1357:ARG:HA	7:C:3273:HOH:O	1.86	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:2001[B]:4LM:H4A	2:A:2001[B]:4LM:HGB	1.68	0.74
1:D:1874:ASN:HD21	1:D:1876:ASP:HB2	1.54	0.73
1:D:1612:VAL:O	1:D:1616:THR:HB	1.90	0.72
1:A:100:HIS:HD2	1:A:145:THR:OG1	1.73	0.72
1:A:112:VAL:O	1:A:116:THR:HB	1.91	0.70
1:A:205:LYS:HZ3	2:A:2001[B]:4LM:HGB	1.58	0.69
1:C:1100:HIS:HD2	1:C:1145:THR:OG1	1.76	0.68
1:D:1600:HIS:HD2	1:D:1645:THR:OG1	1.77	0.68
1:B:842:HIS:HE1	1:B:863:PRO:O	1.76	0.68
1:B:600:HIS:HD2	1:B:645:THR:OG1	1.75	0.68
1:C:1293:HIS:HE1	7:C:3475:HOH:O	1.76	0.67
1:D:1606:THR:HB	1:D:1859:GLN:HG2	1.77	0.67
2:A:2001[B]:4LM:CG	2:A:2001[B]:4LM:C4A	2.71	0.67
1:D:1514:LYS:CD	1:D:1514:LYS:H	2.08	0.66
1:C:1319:LYS:O	1:C:1323:HIS:HD2	1.79	0.66
1:D:1606:THR:HG21	7:D:3669:HOH:O	1.96	0.65
2:A:2001[B]:4LM:H4A	2:A:2001[B]:4LM:HGA	1.78	0.64
1:C:1205:LYS:HE2	6:C:2004:2LM:H4A	1.78	0.64
2:A:2001[B]:4LM:HG	1:D:1553:TYR:CZ	2.33	0.63
1:B:873:GLU:HB2	1:B:878:ILE:HD11	1.80	0.63
1:B:781:LYS:HD3	1:B:783:ASN:HD21	1.62	0.63
1:A:261:ILE:HD11	1:C:1257:ILE:HG12	1.81	0.62
1:B:512:HIS:HE1	1:D:1873:GLU:OE1	1.83	0.62
1:C:1205:LYS:CE	6:C:2004:2LM:H4A	2.30	0.61
2:A:2001[B]:4LM:HG	1:D:1553:TYR:CE2	2.36	0.61
1:B:600:HIS:HE1	1:B:627:ASP:OD2	1.84	0.60
1:A:243:MET:HE1	1:B:743:MET:CE	2.32	0.60
1:C:1281:LYS:HE2	7:C:3161:HOH:O	2.02	0.59
1:A:285:PRO:O	1:A:293:HIS:HD2	1.85	0.59
1:A:356:MET:HE2	1:A:357:ARG:HA	1.84	0.59
1:A:354:ASN:H	1:A:354:ASN:HD22	1.50	0.58
1:A:37:THR:HG23	1:D:1822:GLU:OE1	2.04	0.58
1:D:1514:LYS:NZ	1:D:1566:GLU:HB3	2.19	0.57
1:C:1354:ASN:HD22	1:C:1355:ILE:HG13	1.69	0.57
1:D:1808:LEU:HD21	1:D:1865:LEU:HD13	1.86	0.57
1:A:264:GLU:HG3	7:A:3574:HOH:O	2.05	0.56
1:B:531:THR:HG22	7:C:3077:HOH:O	2.05	0.56
1:D:1514:LYS:H	1:D:1514:LYS:HD3	1.71	0.55
1:D:1875:VAL:HG23	7:D:3342:HOH:O	2.06	0.55
2:A:2001[A]:4LM:HGB	3:A:2002:MEE:S	2.47	0.54
1:C:1100:HIS:HE1	1:C:1127:ASP:OD2	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:1874:ASN:ND2	1:D:1876:ASP:HB2	2.21	0.54
1:A:62:GLU:O	1:A:66:GLU:HG3	2.08	0.54
1:A:205:LYS:NZ	2:A:2001[B]:4LM:HGB	2.20	0.53
1:A:243:MET:CE	1:B:743:MET:HE1	2.38	0.53
1:D:1769:VAL:HG21	1:D:1878:ILE:HD11	1.88	0.53
1:A:63:GLN:HG2	7:A:3795:HOH:O	2.08	0.53
1:D:1608:TYR:CE2	1:D:1610:CYS:HB2	2.42	0.53
1:D:1684:THR:HG22	1:D:1688:PHE:HB2	1.91	0.53
6:B:2003:2LM:O1	6:B:2003:2LM:HGA	2.08	0.52
1:B:668:VAL:O	1:B:672:GLU:HG3	2.10	0.52
1:D:1765:ASN:ND2	1:D:1875:VAL:CG2	2.67	0.51
1:C:1374:ASN:HD21	1:C:1376:ASP:HB2	1.75	0.51
1:B:684:THR:HG22	1:B:688:PHE:HB2	1.93	0.51
1:C:1035:ASP:H	1:C:1039:GLN:NE2	2.09	0.51
1:C:1243:MET:HE1	1:D:1743:MET:CE	2.41	0.51
1:D:1706:TYR:CE2	1:D:1834:GLY:HA2	2.47	0.50
1:C:1184:THR:HG22	1:C:1188:PHE:HB2	1.93	0.50
1:A:243:MET:CE	1:B:743:MET:CE	2.88	0.50
7:B:3192:HOH:O	1:C:1037:THR:HG22	2.12	0.49
1:A:184:THR:HG22	1:A:188:PHE:HB2	1.95	0.49
1:A:54:SER:HB2	1:D:1714:ILE:HD12	1.95	0.49
1:D:1571:ILE:HD12	1:D:1686:PRO:HB2	1.94	0.48
1:C:1243:MET:CE	1:D:1743:MET:CE	2.92	0.48
1:A:206:TYR:CE2	1:A:334:GLY:HA2	2.48	0.48
1:A:243:MET:HE1	1:B:743:MET:HE1	1.96	0.47
1:A:7:THR:HG22	1:A:256:PRO:HB2	1.97	0.47
1:C:1353:GLU:O	1:C:1357:ARG:HG3	2.14	0.47
1:C:1285:PRO:O	1:C:1293:HIS:HD2	1.98	0.47
1:A:374:ASN:HD21	1:A:376:ASP:HB2	1.80	0.47
1:C:1034:PHE:HA	1:C:1039:GLN:HE22	1.80	0.47
1:B:873:GLU:HG3	1:D:1511:LEU:HD12	1.97	0.46
1:A:163:ILE:H	1:A:299:GLN:HE22	1.62	0.46
7:A:3125:HOH:O	1:D:1531:THR:HG22	2.16	0.46
1:C:1108:TYR:CE1	6:C:2004:2LM:N	2.83	0.46
1:B:808:LEU:HD21	1:B:865:LEU:HD13	1.98	0.46
1:A:156:PRO:HD3	1:A:348:HIS:CE1	2.50	0.46
1:A:171:HIS:HE1	1:A:195:GLY:O	1.99	0.46
1:C:1171:HIS:HD2	7:C:3007:HOH:O	1.99	0.46
1:C:1163:ILE:H	1:C:1299:GLN:HE22	1.64	0.46
1:A:100:HIS:HE1	1:A:127:ASP:OD2	1.99	0.46
1:D:1884:GLN:O	1:D:1888:LEU:HG	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:1765:ASN:HD21	1:D:1875:VAL:HG22	1.72	0.45
1:C:1007:THR:HG22	1:C:1256:PRO:HB2	1.97	0.45
1:A:347:THR:HB	7:A:3309:HOH:O	2.17	0.45
1:B:542:ASP:HA	1:B:545:MET:HE3	1.97	0.45
1:A:335[A]:CYS:SG	7:D:3905:HOH:O	2.60	0.45
1:B:724:GLU:H	1:B:724:GLU:CD	2.25	0.45
1:D:1811:MET:HG3	1:D:1817:ALA:HA	1.99	0.45
1:D:1605:ASP:OD2	1:D:1606:THR:HG22	2.17	0.45
1:B:706:TYR:CE1	1:B:834:GLY:HA2	2.52	0.44
1:C:1348:HIS:CD2	7:C:3489:HOH:O	2.70	0.44
1:D:1875:VAL:CG2	7:D:3342:HOH:O	2.64	0.44
6:C:2004:2LM:O1	6:C:2004:2LM:HGA	2.18	0.44
1:C:1383:LYS:HE3	7:C:3589:HOH:O	2.18	0.44
1:B:600:HIS:CE1	1:B:627:ASP:OD2	2.69	0.43
1:B:811:MET:HG3	1:B:817:ALA:HA	2.00	0.43
1:D:1785:PRO:O	1:D:1793:HIS:HD2	2.01	0.43
1:A:17:HIS:HD2	7:A:3344:HOH:O	2.00	0.43
1:B:851:VAL:HA	1:B:852:PRO:HD3	1.94	0.43
1:C:1319:LYS:O	1:C:1323:HIS:CD2	2.67	0.43
1:B:781:LYS:HD3	1:B:783:ASN:ND2	2.32	0.43
1:A:261:ILE:CD1	1:C:1257:ILE:HG12	2.47	0.43
1:A:106:THR:HA	1:A:359:GLN:HG2	2.01	0.42
1:A:356:MET:CE	1:A:357:ARG:HA	2.48	0.42
1:B:650:LEU:HA	7:B:3571:HOH:O	2.18	0.42
1:C:1265:ASN:ND2	1:C:1375:VAL:HB	2.34	0.42
1:B:847:THR:HA	6:B:2003:2LM:HEA	2.01	0.42
1:C:1157:THR:HA	1:C:1308:LEU:HD13	2.01	0.42
1:C:1108:TYR:CE2	1:C:1110:CYS:HB2	2.54	0.42
1:D:1511:LEU:HD23	1:D:1753:LYS:HD2	2.00	0.42
1:B:547:LYS:HB2	1:B:547:LYS:HE2	1.85	0.42
1:C:1358:LYS:HE2	7:C:3686:HOH:O	2.20	0.42
1:D:1776:HIS:CE1	1:D:1778:LYS:HB2	2.55	0.42
1:C:1243:MET:CE	1:D:1743:MET:HE1	2.49	0.42
1:C:1374:ASN:ND2	1:C:1376:ASP:HB2	2.33	0.42
1:D:1686:PRO:HD3	1:D:1701:HIS:CE1	2.55	0.42
1:D:1777:GLU:HG2	1:D:1778:LYS:N	2.34	0.42
1:C:1106:THR:HA	1:C:1359:GLN:HG2	2.02	0.41
1:C:1285:PRO:HA	1:C:1290:PHE:CD2	2.55	0.41
1:A:12:HIS:CE1	1:C:1326:VAL:O	2.73	0.41
1:B:785:PRO:O	1:B:793:HIS:HD2	2.04	0.41
1:A:155:ASN:HA	1:A:156:PRO:HA	1.83	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:1200:LEU:HD12	1:C:1200:LEU:C	2.46	0.41
1:B:608:TYR:CE2	1:B:610:CYS:HB2	2.56	0.41
1:B:650:LEU:HD21	1:B:666:ILE:HD13	2.03	0.41
1:B:875:VAL:HG22	7:B:3864:HOH:O	2.20	0.41
2:A:2001[B]:4LM:C4A	2:A:2001[B]:4LM:HGA	2.45	0.40
1:A:108:TYR:CE2	1:A:110:CYS:HB2	2.56	0.40
1:A:356:MET:HE3	1:A:361:ILE:O	2.22	0.40
1:D:1785:PRO:HA	1:D:1790:PHE:CD2	2.57	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	386/389 (99%)	376 (97%)	10 (3%)	0	100	100
1	B	385/389 (99%)	377 (98%)	8 (2%)	0	100	100
1	C	386/389 (99%)	376 (97%)	10 (3%)	0	100	100
1	D	383/389 (98%)	375 (98%)	8 (2%)	0	100	100
All	All	1540/1556 (99%)	1504 (98%)	36 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	321/322 (100%)	311 (97%)	10 (3%)	35	37
1	B	320/322 (99%)	311 (97%)	9 (3%)	38	41
1	C	321/322 (100%)	313 (98%)	8 (2%)	42	45
1	D	319/322 (99%)	300 (94%)	19 (6%)	17	14
All	All	1281/1288 (100%)	1235 (96%)	46 (4%)	31	31

All (46) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	19	LEU
1	A	37	THR
1	A	116	THR
1	A	164	LYS
1	A	183	PHE
1	A	275	GLN
1	A	277	GLU
1	A	333	LEU
1	A	354	ASN
1	A	356	MET
1	B	563	GLN
1	B	616	THR
1	B	724	GLU
1	B	833	LEU
1	B	839	LEU
1	B	851	VAL
1	B	856	MET
1	B	875	VAL
1	B	887	GLU
1	C	1019	LEU
1	C	1039	GLN
1	C	1063	GLN
1	C	1183	PHE
1	C	1288	GLU
1	C	1333	LEU
1	C	1356	MET
1	C	1388	LEU
1	D	1514	LYS
1	D	1519	LEU
1	D	1571	ILE
1	D	1606	THR
1	D	1616	THR
1	D	1693	GLU

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Mol	Chain	Res	Type
1	D	1724	GLU
1	D	1775	GLN
1	D	1777	GLU
1	D	1778	LYS
1	D	1788	GLU
1	D	1798	LYS
1	D	1833	LEU
1	D	1839	LEU
1	D	1853	GLU
1	D	1856	MET
1	D	1858	LYS
1	D	1875	VAL
1	D	1878	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (37) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	17	HIS
1	A	100	HIS
1	A	117	HIS
1	A	171	HIS
1	A	265	ASN
1	A	293	HIS
1	A	299	GLN
1	A	348	HIS
1	A	354	ASN
1	A	374	ASN
1	B	512	HIS
1	B	563	GLN
1	B	600	HIS
1	B	765	ASN
1	B	783	ASN
1	B	793	HIS
1	B	799	GLN
1	B	842	HIS
1	B	884	GLN
1	C	1039	GLN
1	C	1100	HIS
1	C	1117	HIS
1	C	1265	ASN
1	C	1283	ASN
1	C	1293	HIS

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Mol	Chain	Res	Type
1	C	1299	GLN
1	C	1323	HIS
1	C	1354	ASN
1	C	1359	GLN
1	C	1374	ASN
1	D	1600	HIS
1	D	1617	HIS
1	D	1765	ASN
1	D	1775	GLN
1	D	1783	ASN
1	D	1793	HIS
1	D	1874	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](#)

12 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	$\# Z > 2$
4	SO4	C	2008	-	4,4,4	0.20	0	6,6,6	0.24	0
2	4LM	A	2001[A]	-	21,22,22	1.31	1 (4%)	26,31,31	1.75	5 (19%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	2LM	C	2004	-	23,24,24	2.86	4 (17%)	24,33,33	1.84	7 (29%)
6	2LM	B	2003	-	23,24,24	3.15	4 (17%)	24,33,33	1.64	7 (29%)
4	SO4	D	2009	-	4,4,4	0.23	0	6,6,6	0.13	0
5	GOL	C	2011	-	5,5,5	0.35	0	5,5,5	0.28	0
4	SO4	A	2006	-	4,4,4	0.19	0	6,6,6	0.24	0
6	2LM	D	2005	-	23,24,24	2.98	4 (17%)	24,33,33	1.42	3 (12%)
5	GOL	A	2010	-	5,5,5	0.53	0	5,5,5	0.39	0
4	SO4	B	2007	-	4,4,4	0.22	0	6,6,6	0.12	0
3	MEE	A	2002	-	0,1,1	-	-	-	-	-
2	4LM	A	2001[B]	-	21,22,22	1.28	1 (4%)	26,31,31	1.81	5 (19%)

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	4LM	A	2001[A]	-	-	3/15/17/17	0/1/1/1
6	2LM	B	2003	-	-	9/18/19/19	0/1/1/1
5	GOL	C	2011	-	-	0/4/4/4	-
6	2LM	D	2005	-	-	9/18/19/19	0/1/1/1
5	GOL	A	2010	-	-	0/4/4/4	-
6	2LM	C	2004	-	-	9/18/19/19	0/1/1/1
2	4LM	A	2001[B]	-	-	2/15/17/17	0/1/1/1

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	B	2003	2LM	C4A-N	-11.97	1.29	1.46
6	D	2005	2LM	C4A-N	-11.59	1.30	1.46
6	C	2004	2LM	CA-N	9.61	1.41	1.28
6	C	2004	2LM	C4A-N	-8.32	1.34	1.46
6	B	2003	2LM	CA-N	6.70	1.37	1.28
6	D	2005	2LM	CA-N	6.35	1.36	1.28
2	A	2001[A]	4LM	CA-N	-3.24	1.37	1.41
6	B	2003	2LM	C2A-C2	2.78	1.54	1.50
2	A	2001[B]	4LM	CA-N	-2.63	1.38	1.41
6	C	2004	2LM	O2-C	-2.57	1.23	1.30
6	D	2005	2LM	C2A-C2	2.39	1.54	1.50
6	B	2003	2LM	P-OP4	2.24	1.67	1.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	C	2004	2LM	CB-CA	2.16	1.56	1.50
6	D	2005	2LM	P-OP2	-2.15	1.46	1.54

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	2001[A]	4LM	O2-C-CA	5.18	121.71	113.40
2	A	2001[B]	4LM	O2-C-CA	5.03	121.47	113.40
6	C	2004	2LM	O2-C-O1	-4.36	113.52	123.90
6	B	2003	2LM	CE-SD-CG	4.02	121.16	100.32
6	C	2004	2LM	CE-SD-CG	3.60	118.95	100.32
2	A	2001[B]	4LM	C4-C3-C2	3.50	122.11	120.14
2	A	2001[A]	4LM	C4-C3-C2	3.41	122.06	120.14
6	D	2005	2LM	CE-SD-CG	3.37	117.76	100.32
6	C	2004	2LM	C3-C4-C5	-2.92	116.08	118.73
2	A	2001[B]	4LM	C4A-N-CA	2.90	129.92	122.63
6	C	2004	2LM	C6-C5-C4	2.86	120.22	118.06
6	B	2003	2LM	C3-C2-N1	-2.67	117.59	120.96
2	A	2001[B]	4LM	C3-C4-C5	-2.54	116.24	118.28
2	A	2001[A]	4LM	C4A-N-CA	2.53	128.99	122.63
2	A	2001[A]	4LM	C3-C4-C5	-2.45	116.31	118.28
6	C	2004	2LM	C4A-C4-C5	2.42	122.39	119.76
6	D	2005	2LM	C6-C5-C4	2.40	119.88	118.06
6	B	2003	2LM	OP4-P-OP1	-2.35	100.10	106.44
6	B	2003	2LM	C2A-C2-N1	2.34	122.05	117.64
6	B	2003	2LM	OP3-P-OP1	2.33	119.93	110.83
6	C	2004	2LM	C5A-C5-C6	-2.24	115.71	119.36
6	B	2003	2LM	O2-C-O1	-2.15	118.78	123.90
2	A	2001[B]	4LM	O2-C-O1	-2.14	118.80	123.90
6	D	2005	2LM	O2-C-O1	-2.10	118.91	123.90
2	A	2001[A]	4LM	O2-C-O1	-2.09	118.94	123.90
6	C	2004	2LM	C4-C4A-N	2.04	124.76	114.56
6	B	2003	2LM	C5-C6-N1	-2.01	120.56	123.83

There are no chirality outliers.

All (32) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	B	2003	2LM	O1-C-CA-N
6	B	2003	2LM	O2-C-CA-N
6	B	2003	2LM	O2-C-CA-CB
6	B	2003	2LM	C3-C4-C4A-N

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Mol	Chain	Res	Type	Atoms
6	B	2003	2LM	C5-C4-C4A-N
6	B	2003	2LM	C-CA-CB-CG
6	C	2004	2LM	O1-C-CA-N
6	C	2004	2LM	O2-C-CA-N
6	C	2004	2LM	O2-C-CA-CB
6	C	2004	2LM	C3-C4-C4A-N
6	C	2004	2LM	C5-C4-C4A-N
6	C	2004	2LM	C-CA-CB-CG
6	D	2005	2LM	O1-C-CA-N
6	D	2005	2LM	O2-C-CA-N
6	D	2005	2LM	O2-C-CA-CB
6	D	2005	2LM	C3-C4-C4A-N
6	D	2005	2LM	C5-C4-C4A-N
6	D	2005	2LM	C-CA-CB-CG
6	C	2004	2LM	O1-C-CA-CB
6	B	2003	2LM	CA-CB-CG-SD
6	C	2004	2LM	CA-CB-CG-SD
6	D	2005	2LM	CA-CB-CG-SD
6	D	2005	2LM	CB-CG-SD-CE
2	A	2001[A]	4LM	O1-C-CA-CB
2	A	2001[B]	4LM	O1-C-CA-CB
6	B	2003	2LM	N-CA-CB-CG
6	C	2004	2LM	N-CA-CB-CG
6	D	2005	2LM	N-CA-CB-CG
6	B	2003	2LM	C6-C5-C5A-OP4
2	A	2001[A]	4LM	O2-C-CA-CB
2	A	2001[B]	4LM	O2-C-CA-CB
2	A	2001[A]	4LM	N-CA-CB-CG

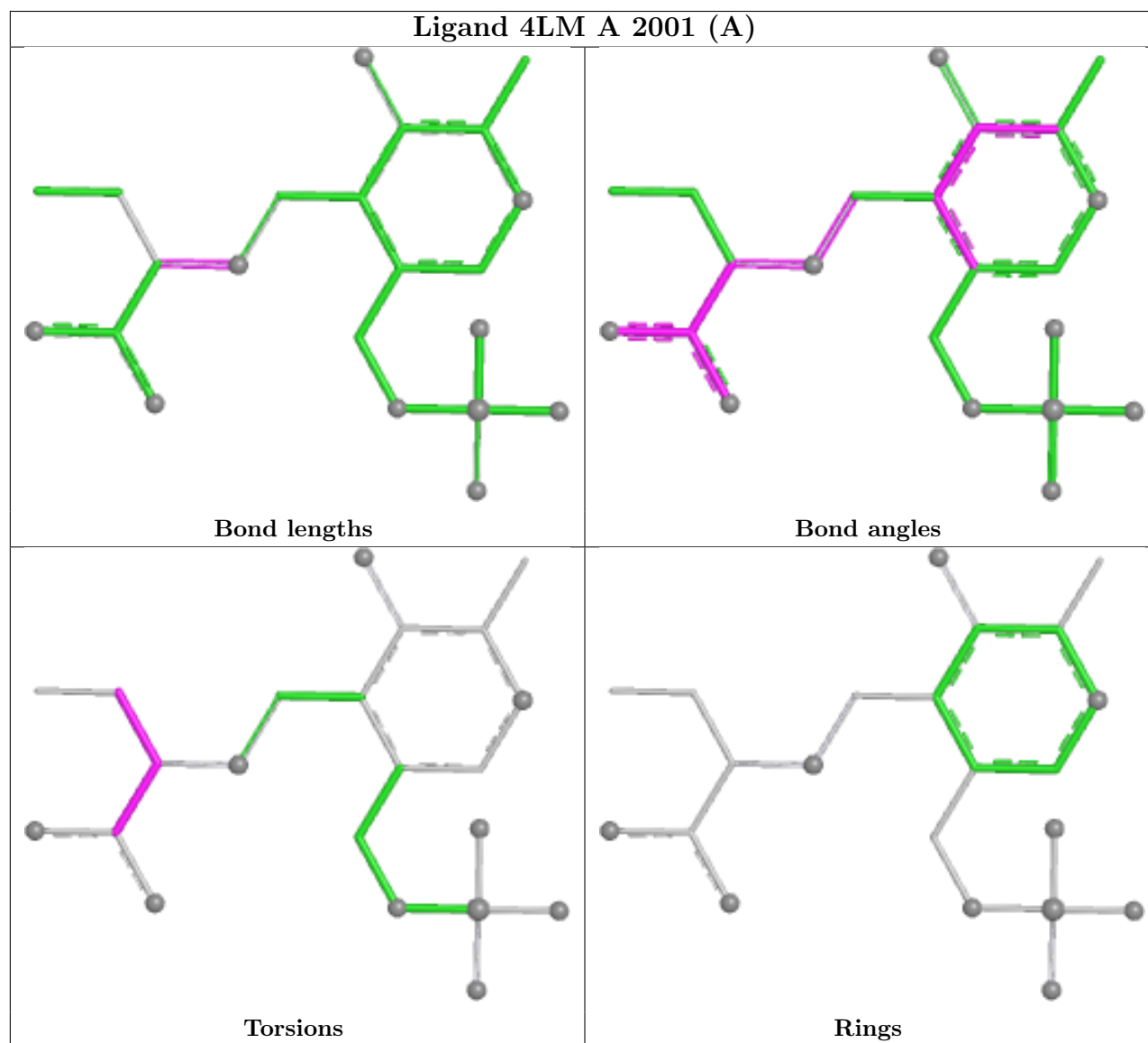
There are no ring outliers.

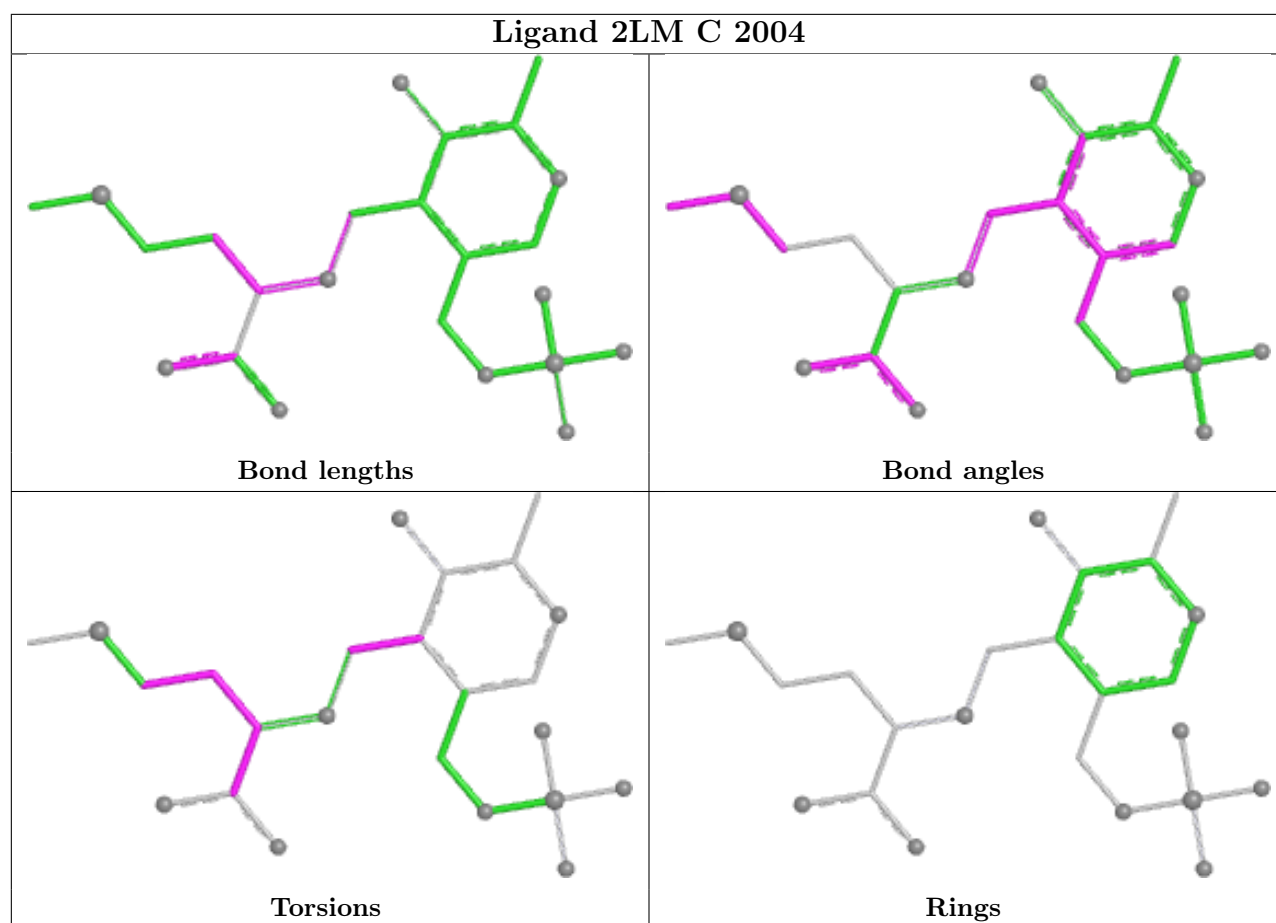
5 monomers are involved in 16 short contacts:

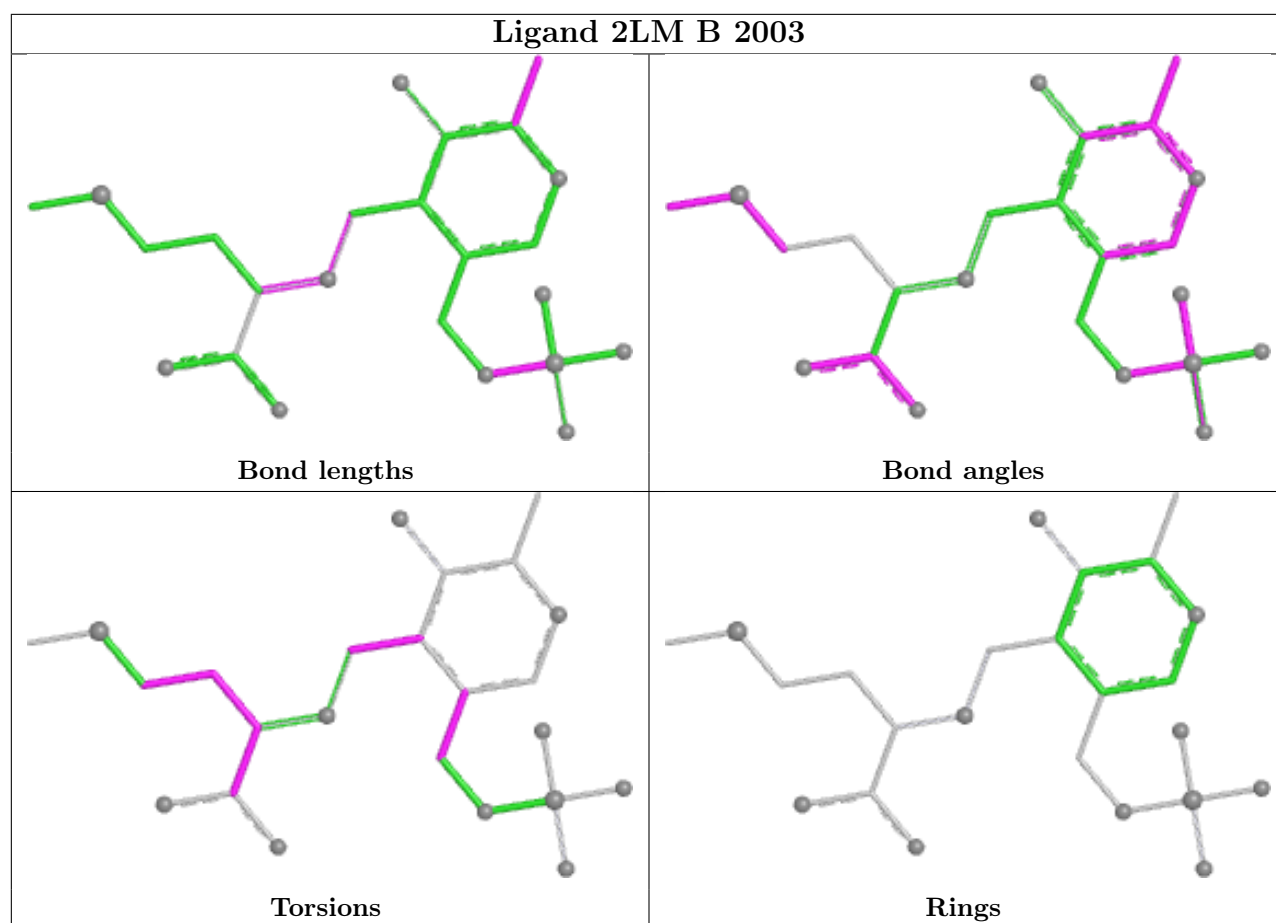
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	2001[A]	4LM	1	0
6	C	2004	2LM	4	0
6	B	2003	2LM	2	0
3	A	2002	MEE	1	0
2	A	2001[B]	4LM	9	0

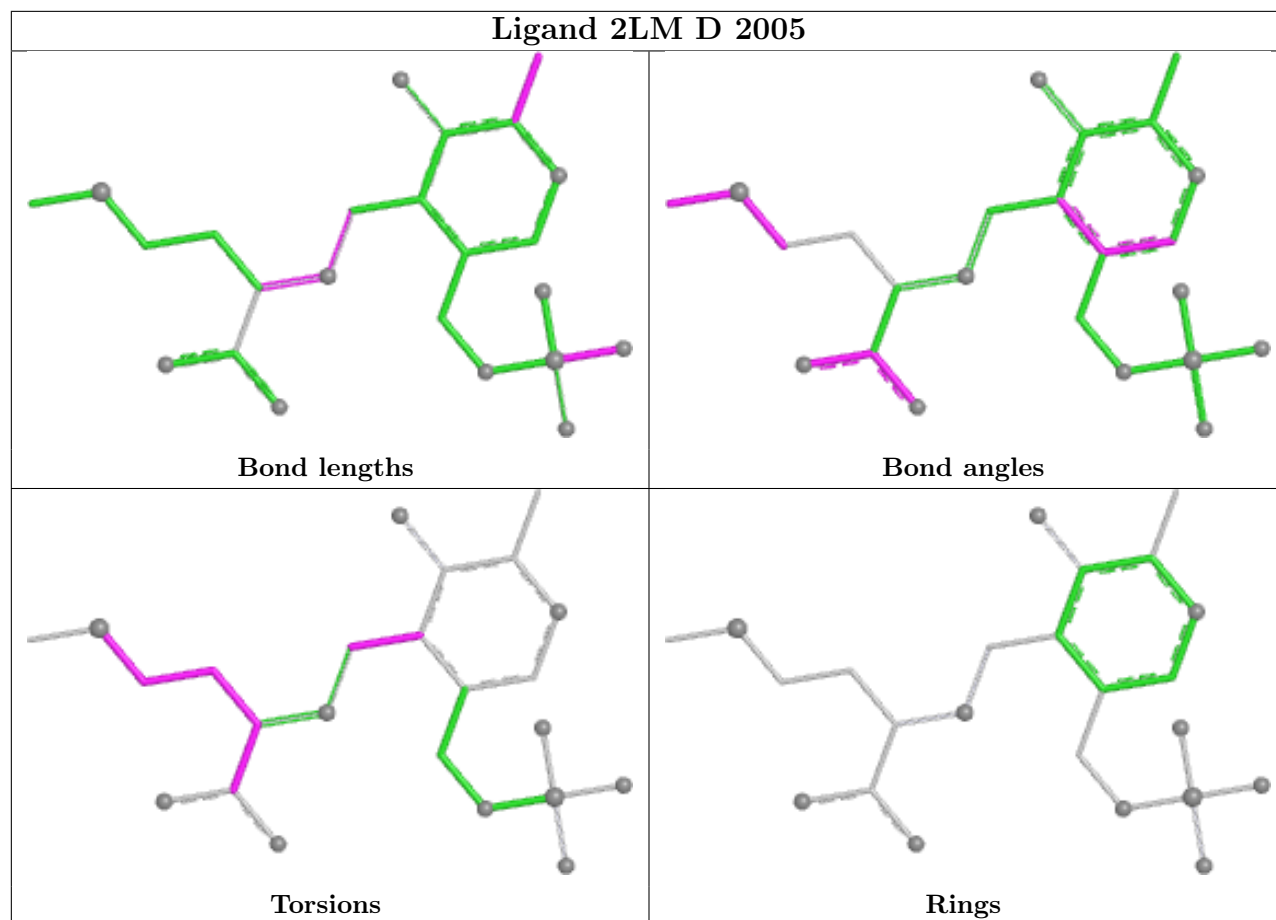
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will

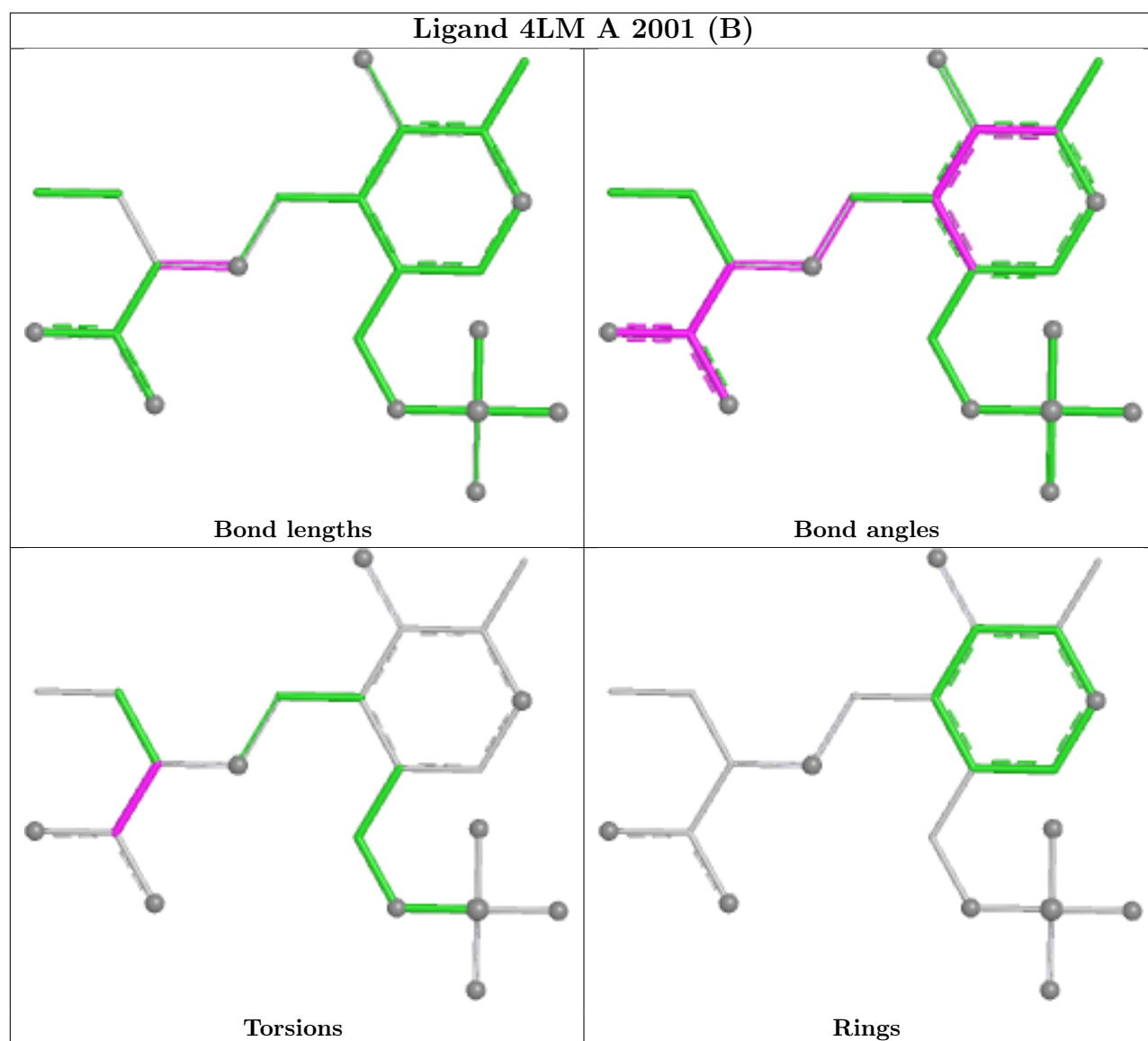
also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted 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	387/389 (99%)	-0.65	7 (1%) 67 67	13, 21, 40, 61	2 (0%)
1	B	386/389 (99%)	-0.39	2 (0%) 87 87	17, 33, 56, 73	2 (0%)
1	C	387/389 (99%)	-0.63	9 (2%) 61 60	11, 20, 39, 62	2 (0%)
1	D	384/389 (98%)	-0.30	2 (0%) 87 87	16, 35, 59, 75	1 (0%)
All	All	1544/1556 (99%)	-0.49	20 (1%) 75 74	11, 25, 54, 75	7 (0%)

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	1351	VAL	4.5
1	B	503	ALA	3.6
1	C	1355	ILE	3.4
1	A	356	MET	3.3
1	A	351	VAL	3.1
1	C	1350	ALA	3.0
1	A	355	ILE	3.0
1	D	1888	LEU	2.9
1	A	352	PRO	2.9
1	C	1356	MET	2.8
1	B	888	LEU	2.8
1	C	1348	HIS	2.6
1	C	1352	PRO	2.4
1	A	348	HIS	2.4
1	C	1354	ASN	2.3
1	C	1349	ALA	2.3
1	C	1357	ARG	2.3
1	D	1885	ALA	2.3
1	A	354	ASN	2.2
1	A	350	ALA	2.1

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](#)

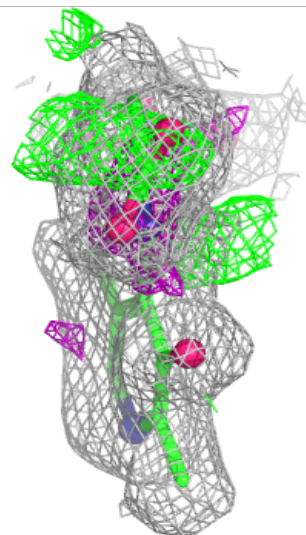
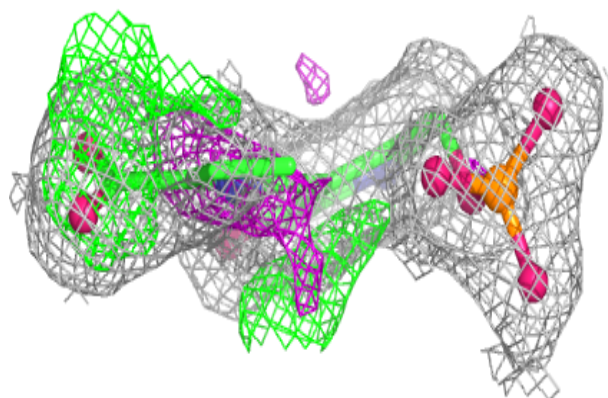
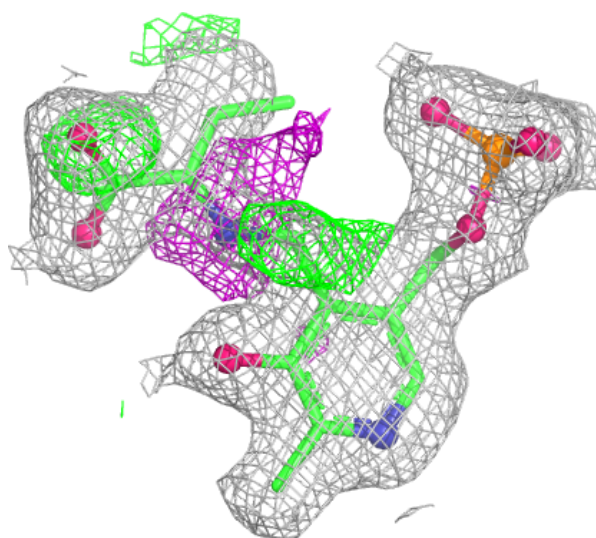
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
3	MEE	A	2002	2/2	0.74	0.28	68,68,68,68	0
5	GOL	A	2010	6/6	0.88	0.11	39,44,46,47	0
4	SO4	D	2009	5/5	0.89	0.07	71,72,72,73	0
4	SO4	B	2007	5/5	0.89	0.08	79,79,80,80	0
4	SO4	C	2008	5/5	0.91	0.08	58,58,59,59	0
5	GOL	C	2011	6/6	0.92	0.11	44,47,48,50	0
4	SO4	A	2006	5/5	0.93	0.07	62,62,62,62	0
2	4LM	A	2001[B]	22/22	0.93	0.10	13,18,27,28	17
2	4LM	A	2001[A]	22/22	0.93	0.10	13,18,27,27	17
6	2LM	B	2003	24/24	0.94	0.10	17,29,40,53	0
6	2LM	C	2004	24/24	0.94	0.10	16,20,37,51	0
6	2LM	D	2005	24/24	0.95	0.09	21,30,40,55	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.

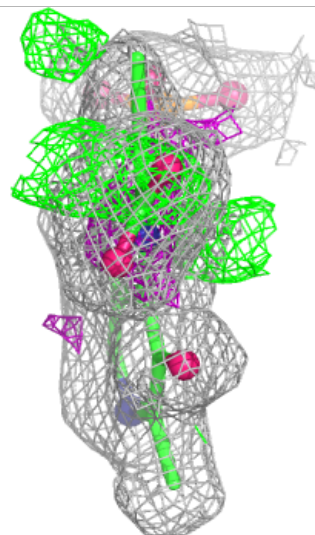
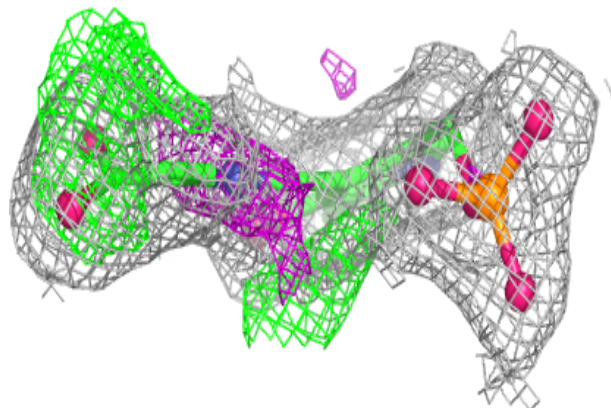
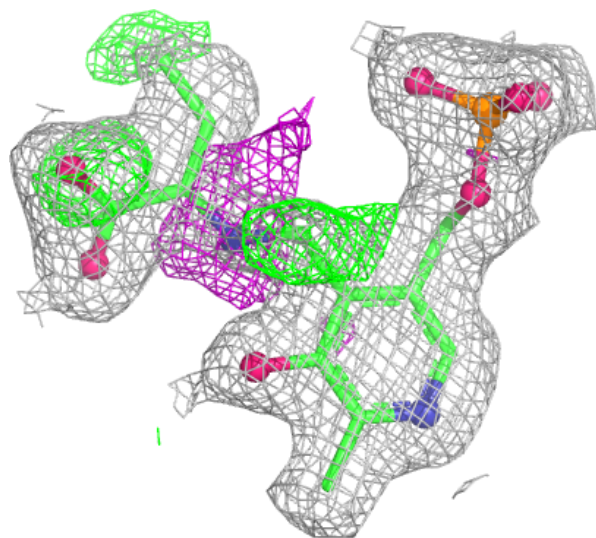
Electron density around 4LM A 2001 (B):

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



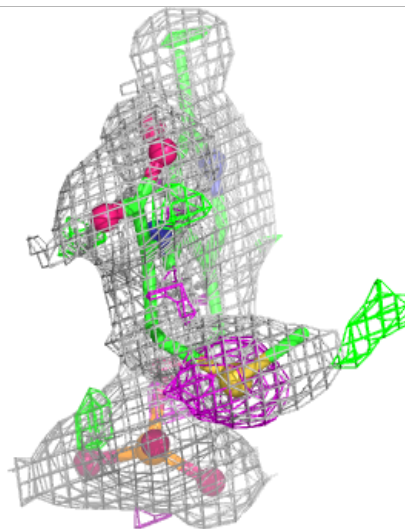
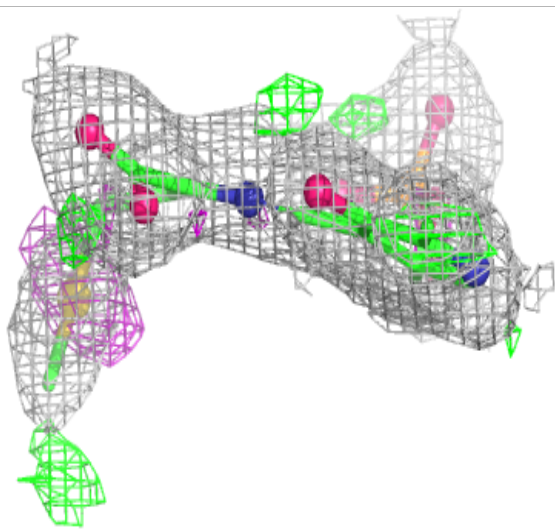
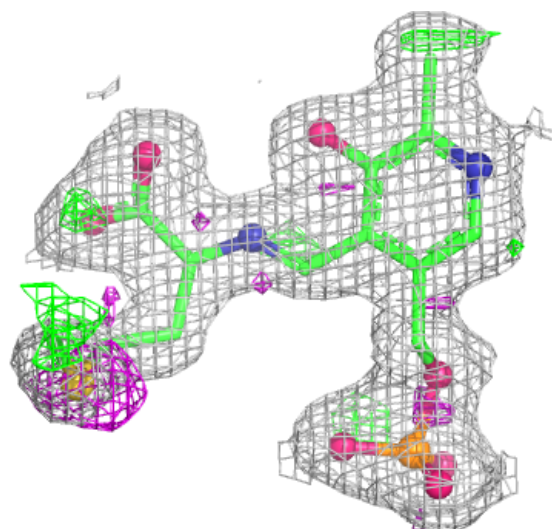
Electron density around 4LM A 2001 (A):

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



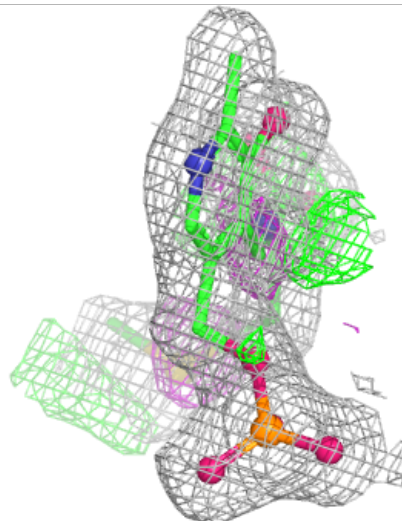
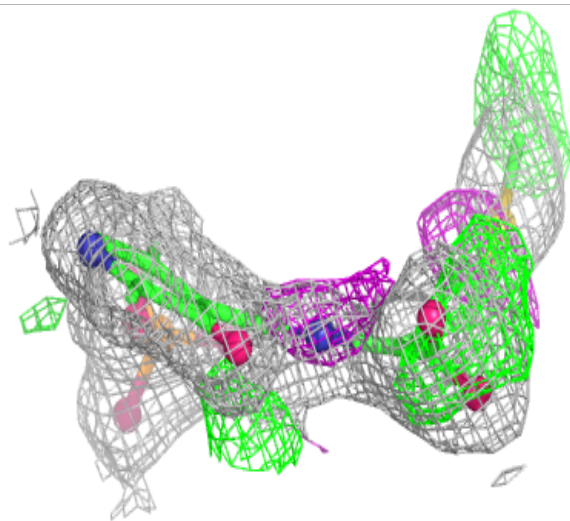
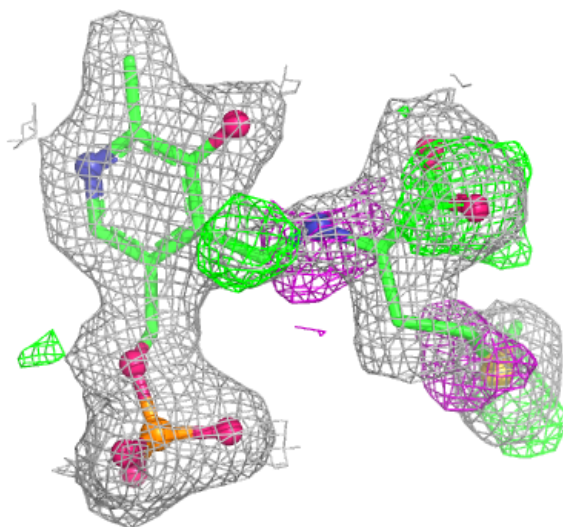
Electron density around 2LM B 2003:

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



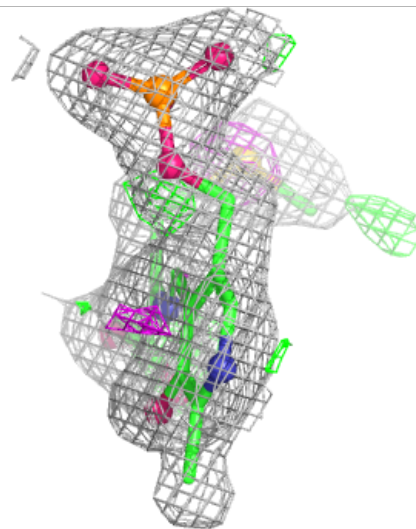
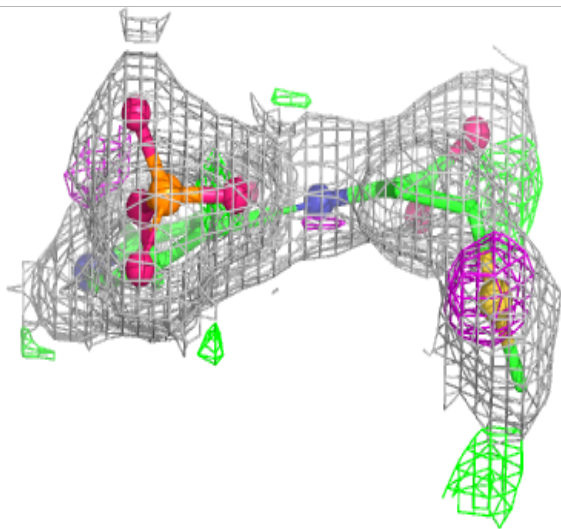
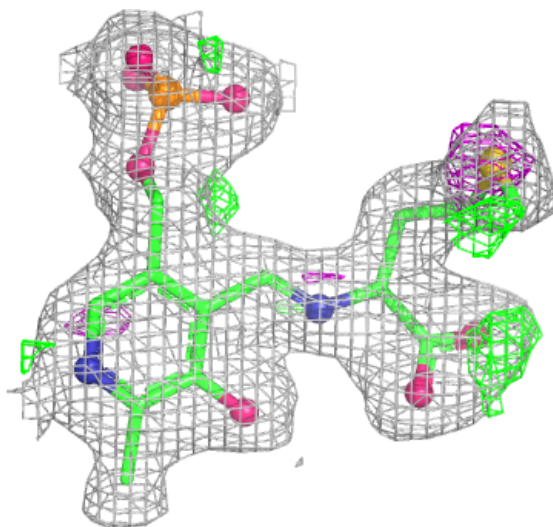
Electron density around 2LM C 2004:

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



Electron density around 2LM D 2005:

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



6.5 Other polymers ⓘ

There are no such residues in this entry.