Full wwPDB X-ray Structure Validation Report ⓘ

Sep 19, 2023 – 09:07 PM EDT

PDB ID	:	5KNK
Title	:	Lipid A secondary acyltransferase LpxM from Acinetobacter baumannii with catalytic residue substitution (E127A)
Authors	:	Dovala, D.L.; Hu, Q.; Metzger IV, L.E.
Deposited on	:	2016-06-28
Resolution	:	1.90 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:

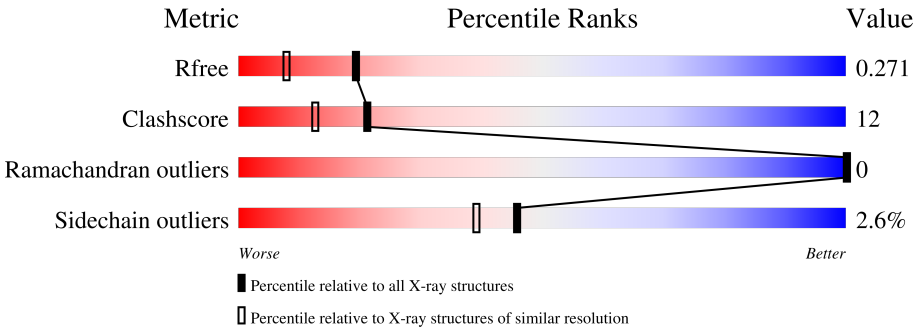
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35.1
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35.1

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	6207 (1.90-1.90)
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)

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 <=5%

Mol	Chain	Length	Quality of chain
1	B	333	77% 19%

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
5	11A	B	410	-	-	X	-

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 2846 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.

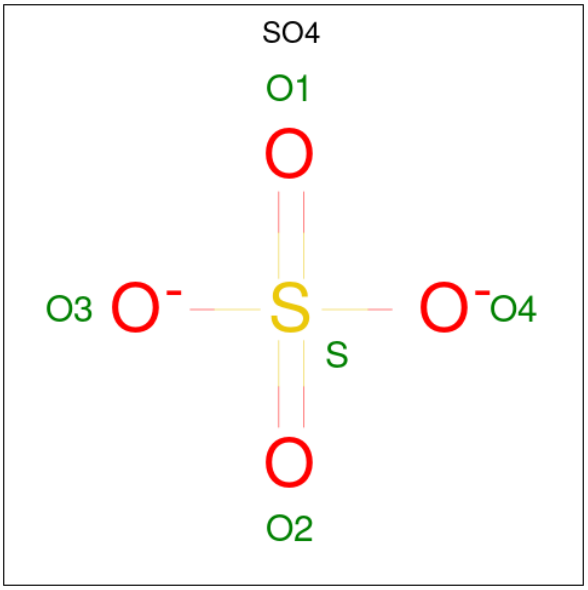
- Molecule 1 is a protein called Lipid A biosynthesis lauroyl acyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	B	324	2620	1685	452	476	7	0	9	0

There are 8 discrepancies between the modelled and reference sequences:

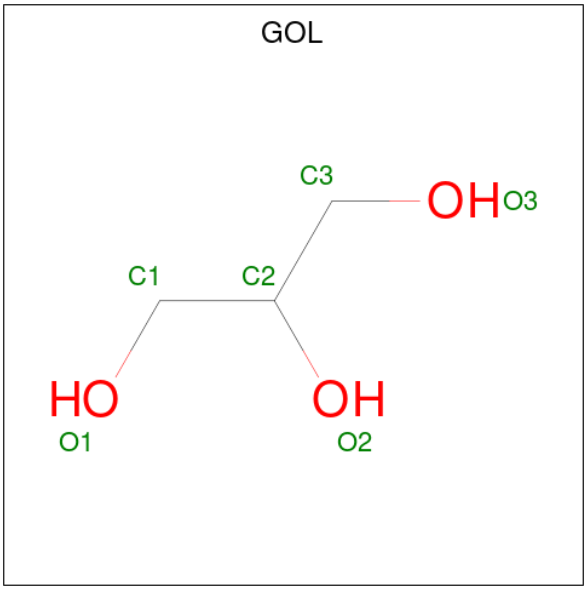
Chain	Residue	Modelled	Actual	Comment	Reference
B	-5	MET	-	initiating methionine	UNP S3TFW2
B	-4	HIS	-	expression tag	UNP S3TFW2
B	-3	HIS	-	expression tag	UNP S3TFW2
B	-2	HIS	-	expression tag	UNP S3TFW2
B	-1	HIS	-	expression tag	UNP S3TFW2
B	0	HIS	-	expression tag	UNP S3TFW2
B	1	HIS	-	expression tag	UNP S3TFW2
B	127	ALA	GLU	engineered mutation	UNP S3TFW2

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



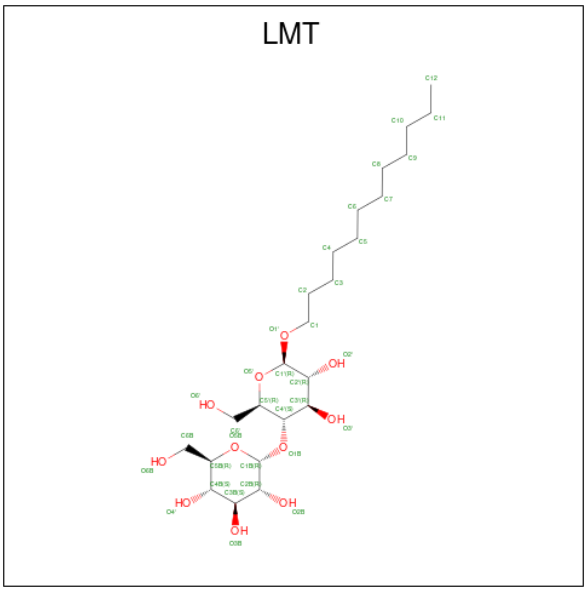
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	1	Total O S 5 4 1	0	0
2	B	1	Total O S 5 4 1	0	0
2	B	1	Total O S 5 4 1	0	0
2	B	1	Total O S 5 4 1	0	0

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



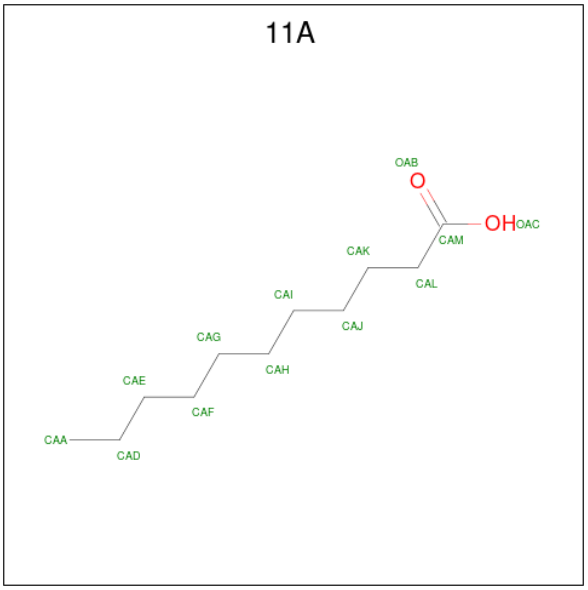
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	1	Total C O 6 3 3	0	0
3	B	1	Total C O 6 3 3	0	0
3	B	1	Total C O 6 3 3	0	0

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



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	1	Total C O 35 24 11	0	0
4	B	1	Total C O 35 24 11	0	0

- Molecule 5 is UNDECANOIC ACID (three-letter code: 11A) (formula: $C_{11}H_{22}O_2$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	B	1	Total C O 13 11 2	0	0

- Molecule 6 is water.

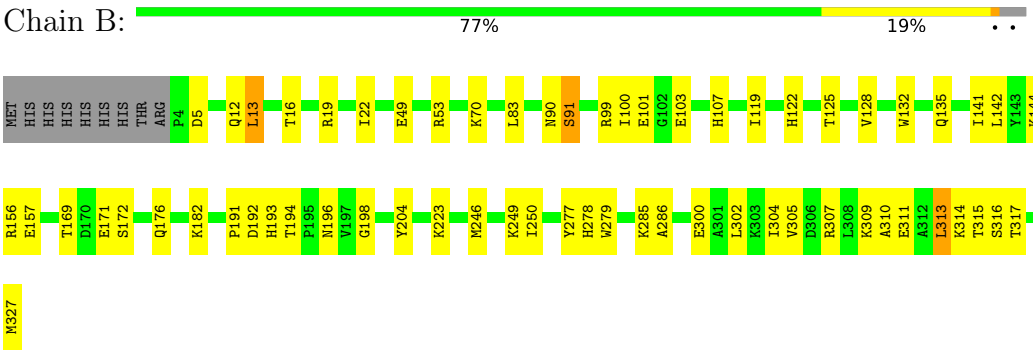
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	105	Total	O	0	0
			105	105		

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. 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. 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: Lipid A biosynthesis lauroyl acyltransferase

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α, β, γ	59.30Å 145.47Å 87.61Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	72.73 – 1.90 72.73 – 1.44	Depositor EDS
% Data completeness (in resolution range)	99.9 (72.73-1.90) 76.6 (72.73-1.44)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	0.11 (at 1.43Å)	Xtriage
Refinement program	PHENIX 1.10_2155	Depositor
R, R_{free}	0.225 , 0.271 0.225 , 0.271	Depositor DCC
R_{free} test set	2000 reflections (2.88%)	wwPDB-VP
Wilson B-factor (Å^2)	14.7	Xtriage
Anisotropy	0.510	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.35 , 42.1	EDS
L-test for twinning[2]	$< \|L\| > = 0.47$, $< L^2 > = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.93	EDS
Total number of atoms	2846	wwPDB-VP
Average B, all atoms (Å^2)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.26% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: SO4, LMT, GOL, 11A

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	B	0.42	1/2701 (0.0%)	0.60	3/3663 (0.1%)

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	B	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	320	GLU	C-N	5.09	1.44	1.34

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	313	LEU	CA-CB-CG	6.03	129.17	115.30
1	B	313	LEU	CB-CG-CD2	5.21	119.86	111.00
1	B	13	LEU	CA-CB-CG	5.13	127.10	115.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	5	ASP	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	B	2620	0	2638	64	0
2	B	20	0	0	0	0
3	B	18	0	24	2	0
4	B	70	0	92	13	0
5	B	13	0	21	16	0
6	B	105	0	0	5	1
All	All	2846	0	2775	67	1

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 12.

All (67) 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:83:LEU:HD11	4:B:408:LMT:H91	1.47	0.96
1:B:310:ALA:O	1:B:313:LEU:HD22	1.75	0.86
1:B:19:ARG:NH1	6:B:502:HOH:O	2.15	0.79
1:B:286:ALA:HA	4:B:409:LMT:H71	1.66	0.77
1:B:171:GLU:OE1	6:B:501:HOH:O	2.02	0.76
1:B:194:THR:HG21	4:B:409:LMT:H52	1.68	0.73
1:B:122:HIS:HB2	5:B:410:11A:HAI2	1.75	0.67
1:B:285:LYS:HB3	4:B:409:LMT:H42	1.76	0.67
1:B:193:HIS:NE2	5:B:410:11A:HAK1	2.09	0.67
1:B:277:TYR:CE1	5:B:410:11A:HAH1	2.32	0.65
1:B:144[B]:LYS:HE3	1:B:169[B]:THR:HG21	1.79	0.63
1:B:278:HIS:HB2	5:B:410:11A:HAJ1	1.80	0.63
1:B:313:LEU:HD23	1:B:314:LYS:H	1.64	0.63
1:B:277:TYR:OH	5:B:410:11A:HAF2	1.99	0.61
1:B:132:TRP:O	1:B:135:GLN:HG2	2.00	0.61
1:B:310:ALA:O	1:B:313:LEU:CD2	2.48	0.61
1:B:286:ALA:HB2	4:B:409:LMT:H41	1.83	0.60
1:B:153:ARG:C	1:B:153:ARG:HD3	2.23	0.58
1:B:198:GLY:O	4:B:409:LMT:H112	2.04	0.57
1:B:194:THR:HB	4:B:409:LMT:H22	1.86	0.57
1:B:12:GLN:O	1:B:16:THR:HG23	2.06	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:70:LYS:NZ	6:B:505:HOH:O	2.38	0.55
1:B:153:ARG:NH1	1:B:157:GLU:HB2	2.21	0.55
1:B:193:HIS:NE2	5:B:410:11A:HAI1	2.22	0.54
1:B:279:TRP:CZ3	5:B:410:11A:HAA1	2.43	0.54
1:B:119:ILE:HG23	1:B:191:PRO:HG3	1.90	0.54
1:B:313:LEU:O	1:B:316:SER:OG	2.23	0.53
1:B:193:HIS:HD1	4:B:409:LMT:C1B	2.19	0.53
1:B:194:THR:OG1	4:B:409:LMT:H31	2.09	0.51
1:B:53:ARG:HB3	1:B:302:LEU:HD11	1.92	0.51
1:B:193:HIS:CD2	5:B:410:11A:HAI1	2.46	0.51
1:B:313:LEU:HD23	1:B:314:LYS:N	2.27	0.50
1:B:49:GLU:OE2	1:B:53:ARG:NH2	2.45	0.50
1:B:70:LYS:HG3	6:B:585:HOH:O	2.10	0.49
1:B:90:ASN:HB2	3:B:405:GOL:H2	1.95	0.48
1:B:277:TYR:CE1	5:B:410:11A:HAF2	2.49	0.48
1:B:277:TYR:HE1	5:B:410:11A:HAD1	1.77	0.48
4:B:409:LMT:H52	4:B:409:LMT:H82	1.57	0.47
1:B:307:ARG:O	1:B:311:GLU:HG3	2.14	0.47
1:B:194:THR:HG21	4:B:409:LMT:H82	1.96	0.47
1:B:300:GLU:O	1:B:304:ILE:HG13	2.15	0.47
1:B:192:ASP:OD1	1:B:192:ASP:N	2.48	0.46
1:B:246:MET:HE2	1:B:250:ILE:HD12	1.98	0.46
1:B:100:ILE:HD13	1:B:132:TRP:CZ2	2.51	0.46
1:B:16:THR:O	1:B:19:ARG:HG2	2.16	0.46
1:B:277:TYR:CZ	5:B:410:11A:HAH1	2.51	0.46
1:B:152:ASP:O	1:B:156:ARG:HG3	2.17	0.45
1:B:103:GLU:HG2	1:B:107:HIS:CD2	2.51	0.45
1:B:204:TYR:CE2	5:B:410:11A:HAA3	2.52	0.45
1:B:99:ARG:NH1	1:B:101:GLU:OE1	2.46	0.45
1:B:204:TYR:OH	5:B:410:11A:HAA2	2.17	0.45
1:B:305:VAL:HG12	1:B:309:LYS:HD2	1.99	0.45
1:B:125:THR:O	1:B:128:VAL:HG12	2.18	0.44
1:B:249:LYS:HE3	1:B:249:LYS:HB3	1.86	0.44
1:B:204:TYR:HE2	5:B:410:11A:HAA3	1.82	0.44
4:B:409:LMT:H21	4:B:409:LMT:H51	1.22	0.43
1:B:142:LEU:HB3	1:B:169[A]:THR:HG22	2.01	0.43
1:B:182:LYS:HG2	6:B:525:HOH:O	2.19	0.43
1:B:172:SER:O	1:B:176:GLN:HG3	2.19	0.43
1:B:141:ILE:HG22	1:B:142:LEU:O	2.19	0.42
1:B:277:TYR:CZ	5:B:410:11A:HAF2	2.55	0.42
1:B:277:TYR:CE1	5:B:410:11A:HAD1	2.53	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:285:LYS:HB3	1:B:285:LYS:HE3	1.77	0.42
1:B:91:SER:H	3:B:405:GOL:H12	1.84	0.41
1:B:144[B]:LYS:HE2	1:B:196:ASN:ND2	2.36	0.41
1:B:315:THR:OG1	1:B:323[B]:GLN:NE2	2.53	0.41

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:B:502:HOH:O	6:B:502:HOH:O[4_555]	2.07	0.13

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	B	331/333 (99%)	319 (96%)	12 (4%)	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	B	281/291 (97%)	273 (97%)	8 (3%)	43 36

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	13	LEU
1	B	22	ILE
1	B	91	SER
1	B	153	ARG
1	B	223[A]	LYS
1	B	223[B]	LYS
1	B	317	THR
1	B	327	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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)

10 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	SO4	B	404	-	4,4,4	0.18	0	6,6,6	0.21	0
5	11A	B	410	-	12,12,12	0.78	0	12,12,12	0.86	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
3	GOL	B	407	-	5,5,5	0.50	0	5,5,5	0.36	0
2	SO4	B	403	-	4,4,4	0.11	0	6,6,6	0.26	0
4	LMT	B	408	-	36,36,36	1.11	3 (8%)	47,47,47	1.02	3 (6%)
2	SO4	B	401	-	4,4,4	0.18	0	6,6,6	0.24	0
2	SO4	B	402	-	4,4,4	0.15	0	6,6,6	0.19	0
3	GOL	B	406	-	5,5,5	0.58	0	5,5,5	0.25	0
4	LMT	B	409	-	36,36,36	1.20	3 (8%)	47,47,47	1.76	8 (17%)
3	GOL	B	405	-	5,5,5	0.56	0	5,5,5	0.42	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
5	11A	B	410	-	-	8/10/10/10	-
3	GOL	B	407	-	-	4/4/4/4	-
4	LMT	B	408	-	-	7/21/61/61	0/2/2/2
3	GOL	B	406	-	-	2/4/4/4	-
4	LMT	B	409	-	-	11/21/61/61	0/2/2/2
3	GOL	B	405	-	-	2/4/4/4	-

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	409	LMT	O5B-C1B	3.70	1.51	1.41
4	B	409	LMT	O5'-C1'	3.70	1.51	1.41
4	B	408	LMT	O5B-C1B	3.43	1.50	1.41
4	B	408	LMT	O5'-C1'	2.92	1.49	1.41
4	B	409	LMT	O5'-C5'	2.15	1.49	1.44
4	B	408	LMT	O5B-C5B	2.00	1.49	1.44

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	B	409	LMT	C1B-O1B-C4'	-5.83	103.54	117.96
4	B	409	LMT	O5'-C5'-C6'	3.58	115.34	106.44
4	B	409	LMT	C1-O1'-C1'	3.49	119.63	113.84
4	B	409	LMT	C1'-O5'-C5'	-3.29	107.24	113.69
4	B	409	LMT	O1B-C1B-C2B	3.17	116.30	108.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	B	409	LMT	C3B-C4B-C5B	2.87	115.36	110.24
4	B	408	LMT	C1B-O1B-C4'	-2.33	112.19	117.96
4	B	409	LMT	C4B-C3B-C2B	2.24	114.74	110.82
4	B	409	LMT	O5B-C5B-C4B	2.05	113.42	109.69
4	B	408	LMT	C1'-O5'-C5'	-2.03	109.70	113.69
4	B	408	LMT	C6B-C5B-C4B	-2.02	108.27	113.00

There are no chirality outliers.

All (34) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	405	GOL	C1-C2-C3-O3
3	B	406	GOL	C1-C2-C3-O3
3	B	407	GOL	C1-C2-C3-O3
4	B	409	LMT	C2-C1-O1'-C1'
4	B	409	LMT	C2-C3-C4-C5
4	B	409	LMT	C5-C6-C7-C8
5	B	410	11A	CAJ-CAK-CAL-CAM
4	B	408	LMT	C11-C10-C9-C8
5	B	410	11A	CAD-CAE-CAF-CAG
4	B	409	LMT	C4-C5-C6-C7
3	B	406	GOL	O2-C2-C3-O3
3	B	407	GOL	O2-C2-C3-O3
4	B	409	LMT	C7-C8-C9-C10
4	B	408	LMT	C3-C4-C5-C6
5	B	410	11A	CAE-CAF-CAG-CAH
4	B	408	LMT	C4B-C5B-C6B-O6B
4	B	408	LMT	C7-C8-C9-C10
4	B	409	LMT	C2'-C1'-O1'-C1
5	B	410	11A	CAF-CAG-CAH-CAI
4	B	409	LMT	C11-C10-C9-C8
5	B	410	11A	CAG-CAH-CAI-CAJ
4	B	409	LMT	C1-C2-C3-C4
3	B	407	GOL	O1-C1-C2-O2
3	B	407	GOL	O1-C1-C2-C3
4	B	408	LMT	O5B-C5B-C6B-O6B
4	B	408	LMT	C5-C6-C7-C8
3	B	405	GOL	O2-C2-C3-O3
4	B	409	LMT	O5B-C5B-C6B-O6B
5	B	410	11A	CAK-CAL-CAM-OAC
5	B	410	11A	CAK-CAL-CAM-OAB
4	B	409	LMT	C6-C7-C8-C9

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Mol	Chain	Res	Type	Atoms
5	B	410	11A	CAH-CAI-CAJ-CAK
4	B	408	LMT	C1-C2-C3-C4
4	B	409	LMT	C4B-C5B-C6B-O6B

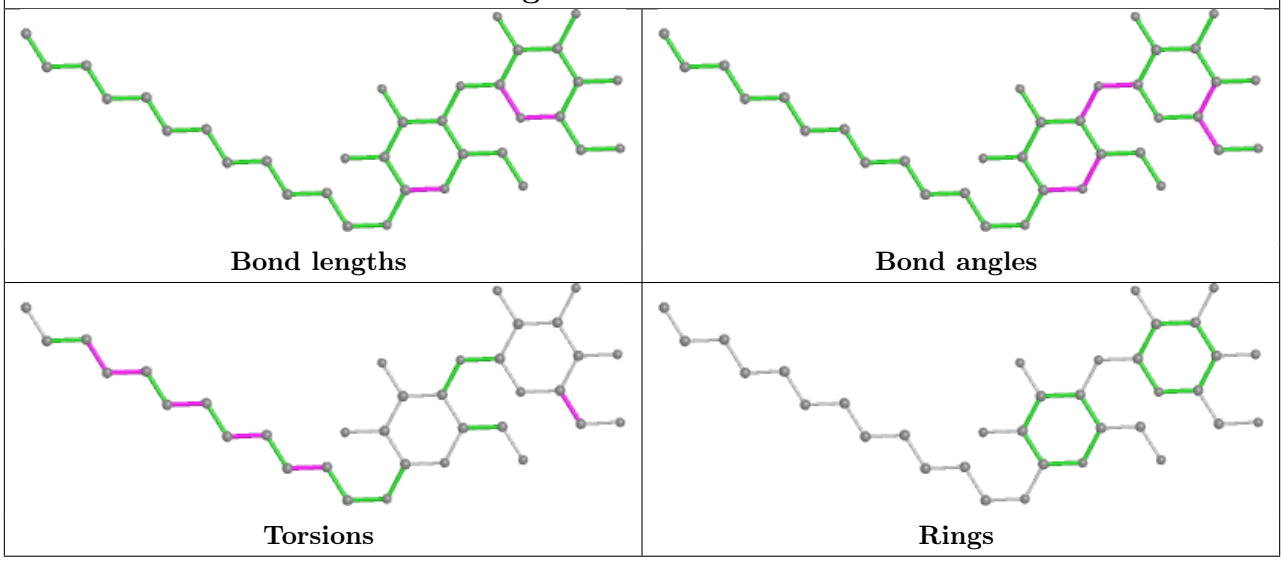
There are no ring outliers.

4 monomers are involved in 31 short contacts:

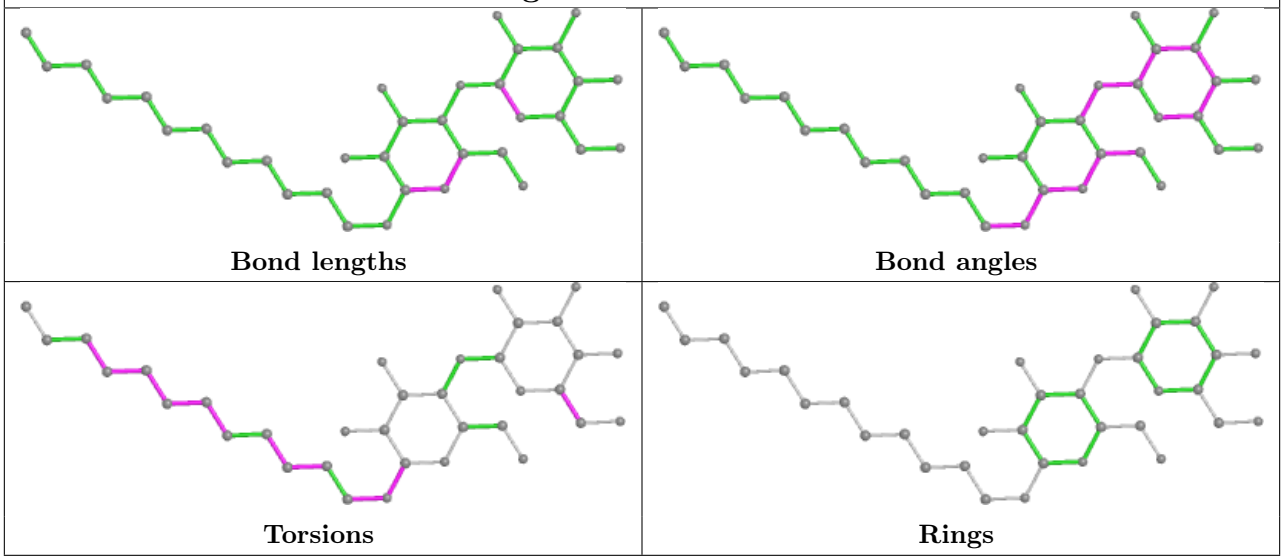
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	410	11A	16	0
4	B	408	LMT	1	0
4	B	409	LMT	12	0
3	B	405	GOL	2	0

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 then 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.

Ligand LMT B 408



Ligand LMT B 409



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

Unable to reproduce the depositors R factor - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

Unable to reproduce the depositors R factor - this section is therefore empty.

6.3 Carbohydrates

Unable to reproduce the depositors R factor - this section is therefore empty.

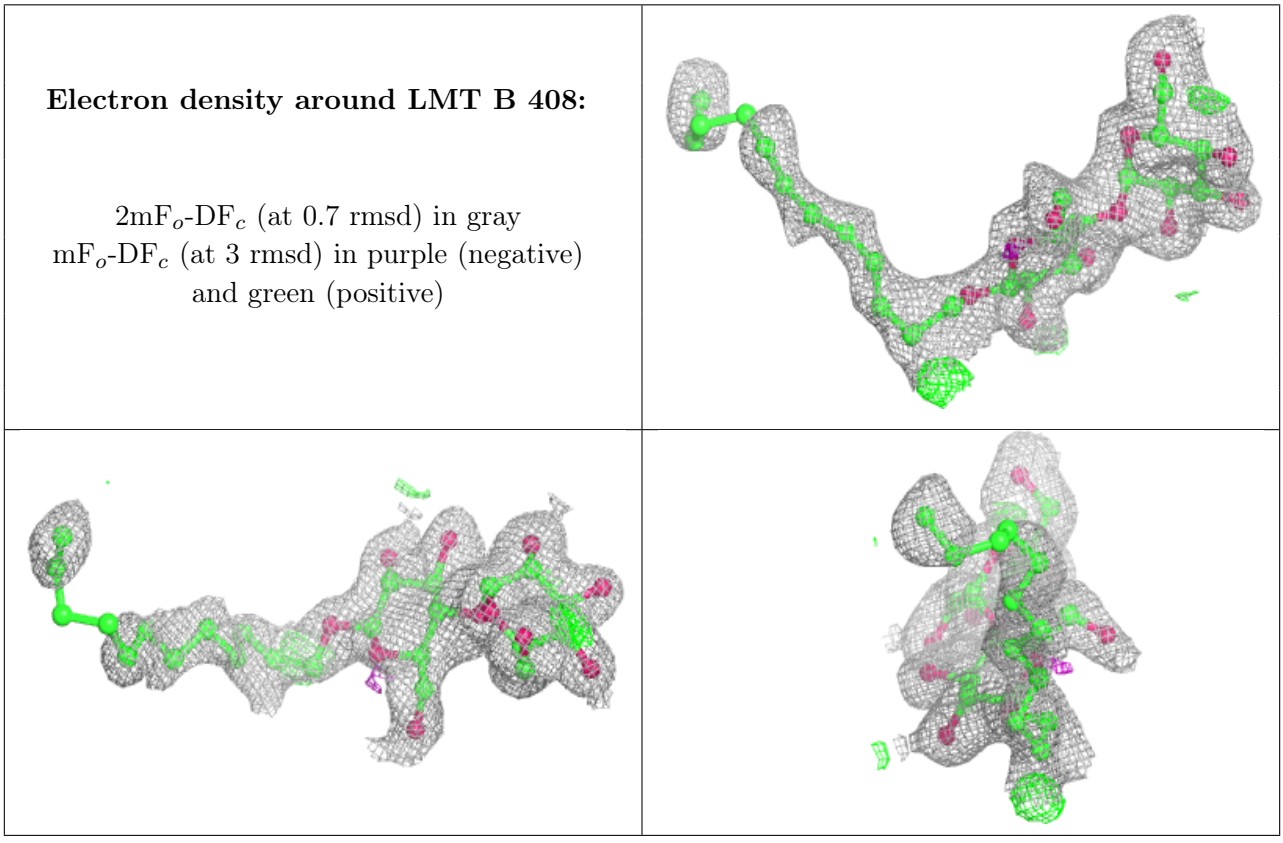
6.4 Ligands

Unable to reproduce the depositors R factor - this section is therefore empty.

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.

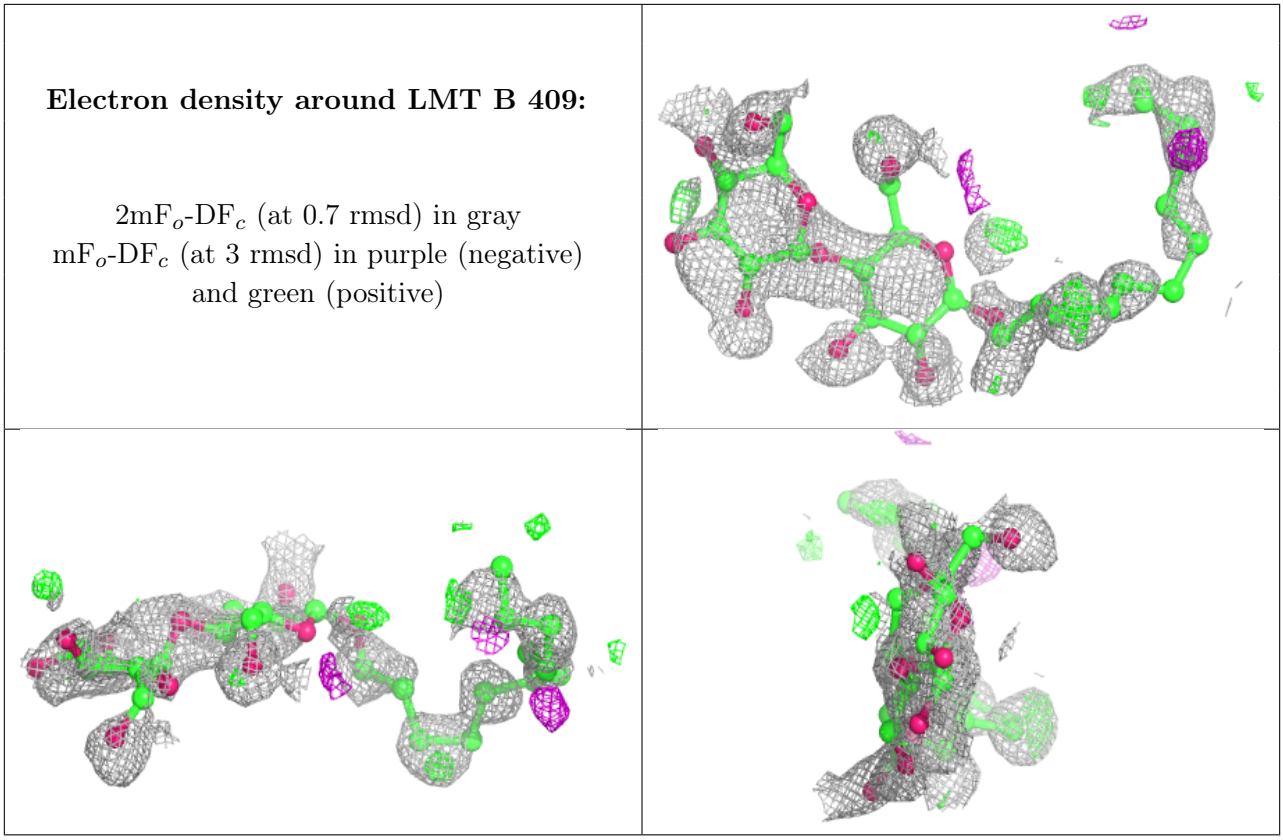
Electron density around LMT B 408:

$2mF_o$-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 409:

$2mF_o$-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)

Unable to reproduce the depositors R factor - this section is therefore empty.