



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 04:59 PM UTC

PDB ID : 2IRV / pdb_00002irv
Title : Crystal structure of GlpG, a rhomboid intramembrane serine protease
Authors : Bibi, E.; Fass, D.; Ben-Shem, A.
Deposited on : 2006-10-16
Resolution : 2.30 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 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

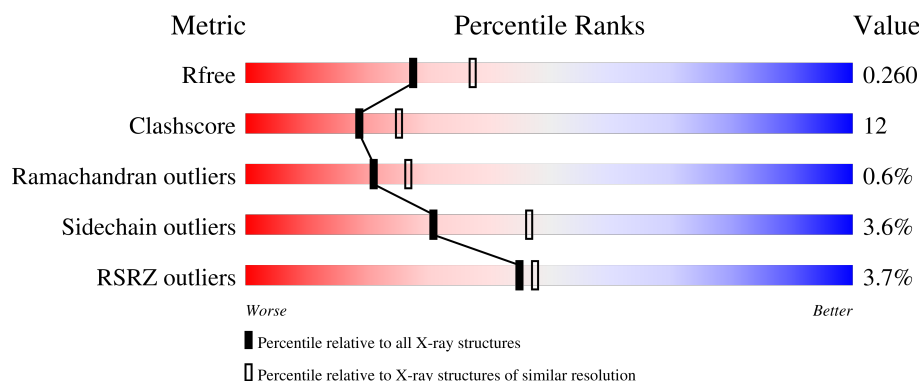
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	182	5% 65% 29% ...
1	B	182	2% 71% 26% ...

2 Entry composition [i](#)

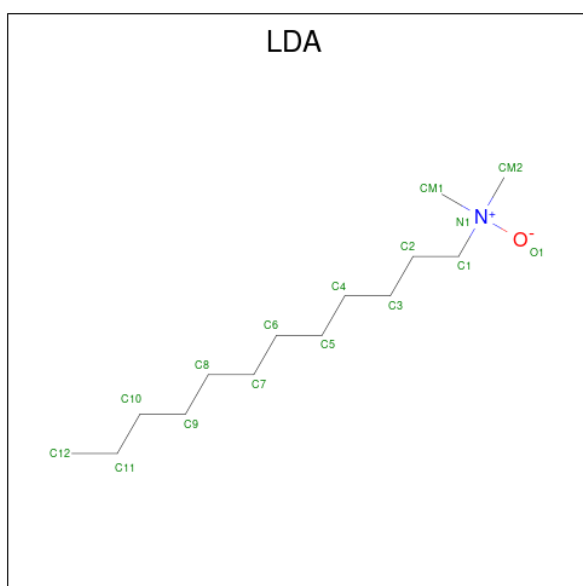
There are 6 unique types of molecules in this entry. The entry contains 3196 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Protein glpG.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	176	Total	C	N	O	S	0	0	0
			1399	945	225	220	9			
1	B	179	Total	C	N	O	S	0	0	0
			1426	964	226	225	11			

- Molecule 2 is LAURYL DIMETHYLAMINE-N-OXIDE (CCD ID: LDA) (formula: $C_{14}H_{31}NO$).



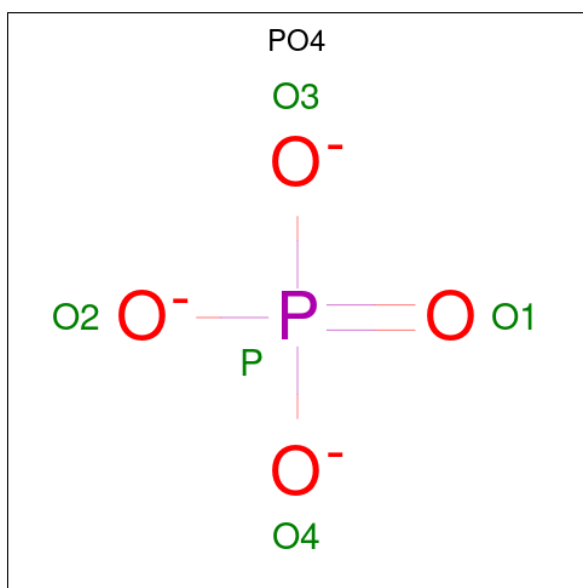
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			10	8	1	1		
2	A	1	Total	C	N	O	0	0
			13	11	1	1		
2	A	1	Total	C	N	O	0	0
			14	12	1	1		
2	B	1	Total	C	N	O	0	0
			16	14	1	1		

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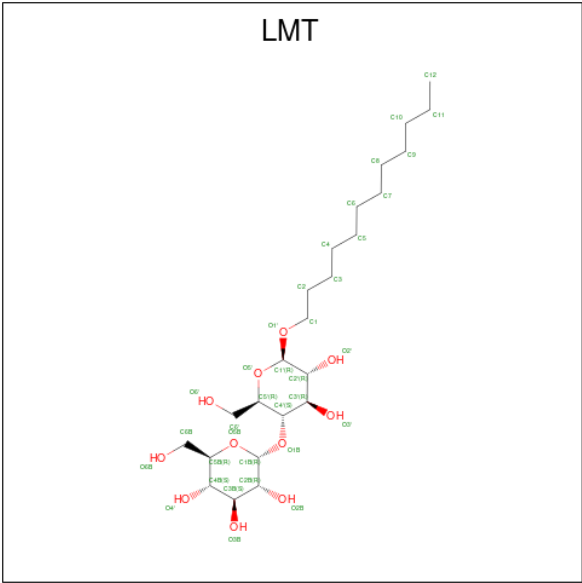
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	B	1	Total	C	N	O	0	0
			16	14	1	1		
2	B	1	Total	C	N	O	0	0
			16	14	1	1		
2	B	1	Total	C	N	O	0	0
			13	11	1	1		
2	B	1	Total	C	N	O	0	0
			12	10	1	1		
2	B	1	Total	C	N	O	0	0
			9	7	1	1		
2	B	1	Total	C	N	O	0	0
			10	8	1	1		
2	B	1	Total	C	N	O	0	0
			11	9	1	1		
2	B	1	Total	C	N	O	0	0
			13	11	1	1		
2	B	1	Total	C	N	O	0	0
			14	12	1	1		
2	B	1	Total	C	N	O	0	0
			11	9	1	1		
2	B	1	Total	C	N	O	0	0
			10	8	1	1		
2	B	1	Total	C	N	O	0	0
			13	11	1	1		

- Molecule 3 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



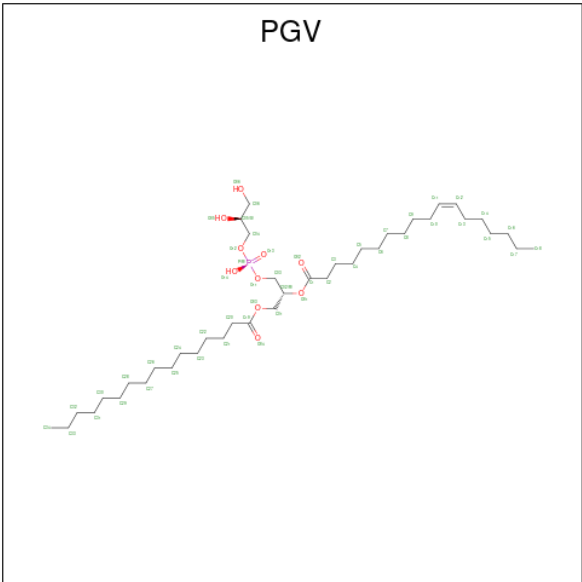
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	O	P	0	0
			5	4	1		

- Molecule 4 is DODECYL-BETA-D-MALTOSE (CCD ID: LMT) (formula: C₂₄H₄₆O₁₁).



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

- Molecule 5 is (1R)-2-[[[(2S)-2,3-DIHYDROXYPROPYL]OXY}(HYDROXY)PHOSPHORYL]OXY}-1-[(PALMITOYLOXY)METHYL]ETHYL (11E)-OCTADEC-11-ENOATE (CCD ID: PGV) (formula: C₄₀H₇₇O₁₀P).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	B	1	Total C O P 26 15 10 1	0	0
5	B	1	Total C 8 8	0	0
5	B	1	Total C 8 8	0	0
5	B	1	Total C 7 7	0	0

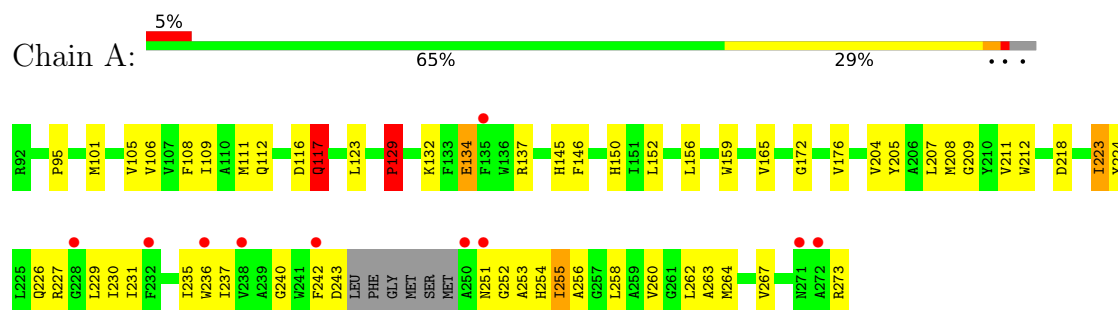
- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	25	Total O 25 25	0	0
6	B	56	Total O 56 56	0	0

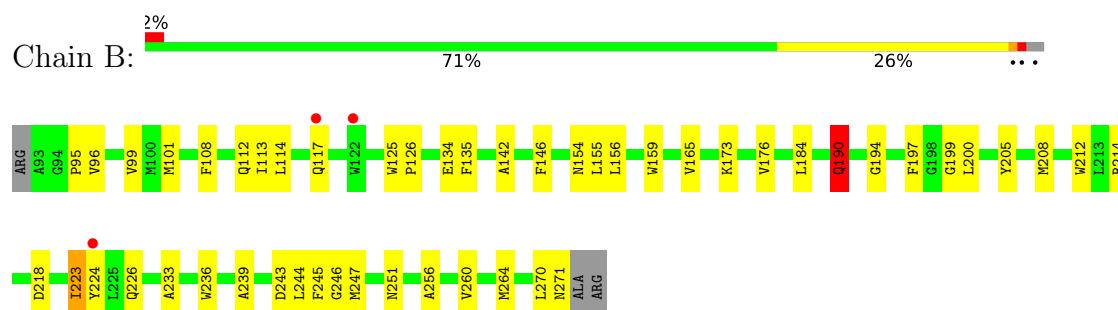
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: Protein glpG



• Molecule 1: Protein glpG



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	49.82Å 69.35Å 67.58Å 90.00° 101.03° 90.00°	Depositor
Resolution (Å)	50.00 – 2.30 50.00 – 2.30	Depositor EDS
% Data completeness (in resolution range)	(Not available) (50.00-2.30) 84.3 (50.00-2.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.15 (at 2.31Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.220 , 0.263 0.236 , 0.260	Depositor DCC
R_{free} test set	1175 reflections (6.90%)	wwPDB-VP
Wilson B-factor (Å ²)	24.2	Xtriage
Anisotropy	0.087	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 52.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	3196	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.13% 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: PO4, PGV, LDA, LMT

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.58	4/1446 (0.3%)	0.97	8/1969 (0.4%)
1	B	0.53	2/1475 (0.1%)	0.90	3/2008 (0.1%)
All	All	0.55	6/2921 (0.2%)	0.94	11/3977 (0.3%)

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	117	GLN	CD-OE1	5.36	1.33	1.23
1	B	226	GLN	CD-OE1	5.35	1.33	1.23
1	B	190	GLN	CD-OE1	5.34	1.33	1.23
1	A	254	HIS	ND1-CE1	5.24	1.37	1.32
1	A	150	HIS	ND1-CE1	5.12	1.37	1.32
1	A	145	HIS	ND1-CE1	5.11	1.37	1.32

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	145	HIS	CB-CG-CD2	-6.62	122.60	131.20
1	A	150	HIS	CB-CG-CD2	-6.43	122.84	131.20
1	A	254	HIS	CB-CG-CD2	-6.42	122.86	131.20
1	A	123	LEU	N-CA-C	6.11	120.90	113.38
1	B	134	GLU	N-CA-C	-5.93	100.73	109.62
1	B	146	PHE	N-CA-C	5.84	119.72	112.59
1	A	146	PHE	N-CA-C	5.81	119.67	112.59
1	A	254	HIS	CB-CG-ND1	5.61	131.11	122.70
1	A	150	HIS	CB-CG-ND1	5.50	130.96	122.70
1	A	145	HIS	CB-CG-ND1	5.48	130.92	122.70
1	B	135	PHE	N-CA-C	5.04	118.89	112.34

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	1399	0	1400	37	0
1	B	1426	0	1431	39	0
2	A	37	0	62	2	0
2	B	164	0	285	11	0
3	B	5	0	0	0	0
4	B	35	0	46	1	0
5	B	49	0	56	4	0
6	A	25	0	0	1	0
6	B	56	0	0	2	0
All	All	3196	0	3280	76	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 12.

All (76) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:165:VAL:HA	1:A:223:ILE:HD11	1.55	0.88
1:B:165:VAL:HA	1:B:223:ILE:HD11	1.56	0.88
1:B:156:LEU:HD23	2:B:3837:LDA:HM21	1.58	0.86
1:A:211:VAL:HG13	1:A:223:ILE:HG23	1.59	0.84
1:B:154:ASN:HB3	1:B:200:LEU:HD12	1.60	0.81
1:B:114:LEU:HD21	2:B:3834:LDA:H42	1.63	0.78
1:B:101:MET:HE3	1:B:155:LEU:HD13	1.69	0.74
1:B:251:ASN:HB3	6:B:19:HOH:O	1.89	0.71
1:A:117:GLN:H	1:A:117:GLN:CD	2.03	0.66
1:A:152:LEU:O	1:A:156:LEU:HG	1.95	0.66
1:B:108:PHE:O	1:B:112:GLN:HG2	1.96	0.66
1:A:251:ASN:C	1:A:253:ALA:H	2.02	0.66
1:A:101:MET:SD	1:A:159:TRP:HE3	2.19	0.65
1:A:218:ASP:OD1	1:A:273:ARG:HD2	1.97	0.65
1:A:263:ALA:O	1:A:267:VAL:HG23	1.98	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:95:PRO:HA	2:A:3842:LDA:H11	1.81	0.62
1:B:184:LEU:HD21	4:B:3888:LMT:H92	1.82	0.60
1:A:106:VAL:HG22	2:B:3835:LDA:H122	1.82	0.60
1:B:239:ALA:HB1	1:B:247:MET:CE	2.32	0.60
1:B:239:ALA:HB1	1:B:247:MET:HE1	1.84	0.60
1:A:159:TRP:CD1	2:A:3850:LDA:HM21	2.37	0.59
1:B:256:ALA:O	1:B:260:VAL:HG23	2.04	0.58
1:A:111:MET:HE2	1:A:116:ASP:OD1	2.03	0.57
1:A:226:GLN:HB2	1:A:229:LEU:HG	1.87	0.57
1:B:184:LEU:HD23	1:B:260:VAL:CG2	2.35	0.56
1:B:245:PHE:HD2	1:B:247:MET:HE2	1.71	0.56
1:B:95:PRO:HA	2:B:3836:LDA:H31	1.89	0.55
1:A:165:VAL:CA	1:A:223:ILE:HD11	2.32	0.55
1:B:159:TRP:HB2	2:B:3837:LDA:HM12	1.88	0.55
1:A:134:GLU:HB2	1:A:137:ARG:HD2	1.90	0.54
1:A:251:ASN:C	1:A:253:ALA:N	2.66	0.54
1:B:95:PRO:O	1:B:99:VAL:HG23	2.08	0.53
1:B:205:TYR:HE1	5:B:5136:PGV:H02	1.73	0.53
1:B:96:VAL:HG22	2:B:3838:LDA:H72	1.90	0.52
1:B:142:ALA:HA	1:B:199:GLY:O	2.09	0.52
1:A:212:TRP:CD1	1:A:230:ILE:HD12	2.44	0.52
1:A:251:ASN:O	1:A:253:ALA:N	2.43	0.52
1:A:231:ILE:O	1:A:235:ILE:HG13	2.09	0.52
1:A:129:PRO:O	1:A:132:LYS:HB2	2.10	0.52
1:A:204:VAL:O	1:A:208:MET:HG3	2.11	0.51
1:B:205:TYR:CE1	5:B:5136:PGV:H02	2.45	0.51
1:B:95:PRO:HD3	2:B:3836:LDA:HM22	1.92	0.51
1:A:176:VAL:HG12	1:A:264:MET:HE1	1.93	0.51
1:A:101:MET:SD	1:A:159:TRP:CE3	3.02	0.50
1:A:108:PHE:O	1:A:112:GLN:HG2	2.11	0.50
1:B:125:TRP:O	1:B:126:PRO:C	2.56	0.49
1:A:205:TYR:CE2	1:A:237:ILE:HD11	2.48	0.48
1:B:190:GLN:HA	1:B:194:GLY:O	2.13	0.48
1:A:256:ALA:O	1:A:260:VAL:HG23	2.14	0.48
1:A:105:VAL:O	1:A:109:ILE:HG13	2.14	0.48
1:B:159:TRP:CB	2:B:3837:LDA:HM12	2.43	0.47
1:B:236:TRP:HZ2	5:B:5136:PGV:H221	1.79	0.47
1:B:243:ASP:OD1	1:B:246:GLY:HA2	2.15	0.46
1:A:227:ARG:O	1:A:230:ILE:HG22	2.15	0.46
1:B:176:VAL:HG12	1:B:264:MET:HE1	1.98	0.46
1:A:258:LEU:O	1:A:262:LEU:HG	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:212:TRP:CZ2	2:B:3847:LDA:H12	2.51	0.45
1:A:209:GLY:O	1:A:262:LEU:HD23	2.17	0.45
1:B:117:GLN:HG3	6:B:1:HOH:O	2.16	0.45
1:A:255:ILE:HG22	1:A:256:ALA:N	2.31	0.45
1:A:172:GLY:O	1:A:176:VAL:HG23	2.16	0.44
1:B:165:VAL:HA	1:B:223:ILE:CD1	2.38	0.44
1:B:270:LEU:O	1:B:271:ASN:CB	2.66	0.44
1:A:242:PHE:O	1:A:243:ASP:C	2.62	0.43
1:B:244:LEU:O	1:B:247:MET:HG3	2.18	0.43
1:A:101:MET:HE2	1:A:101:MET:HB3	1.85	0.42
1:B:113:ILE:HD12	2:B:3834:LDA:H71	2.01	0.42
1:B:154:ASN:HB3	1:B:200:LEU:CD1	2.42	0.42
1:B:99:VAL:HG22	2:B:3836:LDA:H71	2.02	0.42
1:B:236:TRP:CZ2	5:B:5136:PGV:H221	2.55	0.42
1:B:197:PHE:CD1	1:B:197:PHE:C	2.97	0.41
1:B:214:ARG:O	1:B:218:ASP:N	2.52	0.41
1:A:240:GLY:HA3	6:A:46:HOH:O	2.19	0.41
1:A:207:LEU:HD23	1:A:207:LEU:HA	1.96	0.40
1:B:208:MET:CE	1:B:233:ALA:HB2	2.51	0.40
1:A:236:TRP:CE3	1:A:236:TRP:HA	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	172/182 (94%)	157 (91%)	13 (8%)	2 (1%)	10	12
1	B	177/182 (97%)	169 (96%)	8 (4%)	0	100	100
All	All	349/364 (96%)	326 (93%)	21 (6%)	2 (1%)	21	27

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	252	GLY
1	A	129	PRO

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	138/145 (95%)	132 (96%)	6 (4%)	26	39
1	B	143/145 (99%)	139 (97%)	4 (3%)	38	56
All	All	281/290 (97%)	271 (96%)	10 (4%)	31	47

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	117	GLN
1	A	129	PRO
1	A	134	GLU
1	A	223	ILE
1	A	224	TYR
1	A	255	ILE
1	B	173	LYS
1	B	190	GLN
1	B	223	ILE
1	B	224	TYR

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

22 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	LDA	A	3850	-	11,13,15	0.89	1 (9%)	12,15,17	0.91	0
2	LDA	B	3836	-	13,15,15	1.03	2 (15%)	14,17,17	0.89	0
2	LDA	B	3847	-	7,9,15	1.45	1 (14%)	8,11,17	0.53	0
2	LDA	A	3845	-	10,12,15	0.95	1 (10%)	11,14,17	0.68	0
2	LDA	B	3839	-	6,8,15	1.08	1 (16%)	7,10,17	0.48	0
2	LDA	B	3837	-	10,12,15	0.96	1 (10%)	11,14,17	0.69	0
5	PGV	B	5139	-	6,6,50	0.35	0	5,5,56	0.76	0
5	PGV	B	5136	-	25,25,50	1.27	3 (12%)	28,31,56	1.69	3 (10%)
2	LDA	B	3841	-	8,10,15	1.17	2 (25%)	9,12,17	0.62	0
2	LDA	A	3842	-	7,9,15	1.16	1 (14%)	8,11,17	0.56	0
4	LMT	B	3888	-	36,36,36	0.91	2 (5%)	47,47,47	1.05	1 (2%)
3	PO4	B	5156	-	4,4,4	1.64	0	6,6,6	0.48	0
2	LDA	B	3843	-	10,12,15	0.93	1 (10%)	11,14,17	0.73	0
2	LDA	B	3834	-	13,15,15	0.99	2 (15%)	14,17,17	1.00	1 (7%)
2	LDA	B	3848	-	10,12,15	1.00	2 (20%)	11,14,17	0.71	0
2	LDA	B	3844	-	11,13,15	1.19	2 (18%)	12,15,17	0.84	0
2	LDA	B	3835	-	13,15,15	0.88	1 (7%)	14,17,17	0.93	0
2	LDA	B	3838	-	9,11,15	1.03	1 (11%)	10,13,17	0.62	0
2	LDA	B	3840	-	7,9,15	1.14	1 (14%)	8,11,17	0.49	0
5	PGV	B	5138	-	7,7,50	0.35	0	6,6,56	0.77	0
5	PGV	B	5137	-	7,7,50	0.40	0	6,6,56	0.80	0
2	LDA	B	3846	-	8,10,15	1.04	1 (12%)	9,12,17	0.69	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
2	LDA	A	3850	-	-	0/11/11/13	-
2	LDA	B	3836	-	-	0/13/13/13	-
2	LDA	B	3847	-	-	0/7/7/13	-
2	LDA	A	3845	-	-	0/10/10/13	-
2	LDA	B	3839	-	-	0/6/6/13	-
2	LDA	B	3837	-	-	0/10/10/13	-
5	PGV	B	5139	-	-	0/4/4/55	-
5	PGV	B	5136	-	-	7/29/29/55	-
2	LDA	B	3841	-	-	0/8/8/13	-
2	LDA	A	3842	-	-	0/7/7/13	-
4	LMT	B	3888	-	-	0/21/61/61	0/2/2/2
2	LDA	B	3843	-	-	0/10/10/13	-
2	LDA	B	3834	-	-	0/13/13/13	-
2	LDA	B	3848	-	-	0/10/10/13	-
2	LDA	B	3844	-	-	0/11/11/13	-
2	LDA	B	3835	-	-	0/13/13/13	-
2	LDA	B	3838	-	-	0/9/9/13	-
2	LDA	B	3840	-	-	0/7/7/13	-
5	PGV	B	5138	-	-	0/5/5/55	-
5	PGV	B	5137	-	-	0/5/5/55	-
2	LDA	B	3846	-	-	0/8/8/13	-

All (26) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	5136	PGV	O01-C02	3.38	1.54	1.46
2	B	3844	LDA	C1-N1	-3.19	1.48	1.51
2	B	3847	LDA	C1-N1	-3.10	1.48	1.51
5	B	5136	PGV	O01-C1	2.80	1.41	1.35
4	B	3888	LMT	O1'-C1'	2.68	1.44	1.40
2	A	3850	LDA	O1-N1	2.59	1.48	1.42
5	B	5136	PGV	C03-C02	2.50	1.58	1.50
2	B	3834	LDA	C1-N1	-2.50	1.48	1.51
2	B	3836	LDA	C1-N1	-2.47	1.49	1.51
4	B	3888	LMT	O5'-C1'	2.45	1.48	1.41
2	B	3846	LDA	O1-N1	2.38	1.47	1.42
2	B	3838	LDA	O1-N1	2.36	1.47	1.42
2	B	3835	LDA	O1-N1	2.35	1.47	1.42
2	B	3836	LDA	O1-N1	2.35	1.47	1.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	3834	LDA	O1-N1	2.33	1.47	1.42
2	B	3841	LDA	O1-N1	2.24	1.47	1.42
2	B	3840	LDA	O1-N1	2.19	1.47	1.42
2	B	3843	LDA	O1-N1	2.16	1.47	1.42
2	B	3841	LDA	C1-N1	-2.16	1.49	1.51
2	A	3845	LDA	O1-N1	2.15	1.47	1.42
2	B	3848	LDA	O1-N1	2.14	1.47	1.42
2	B	3839	LDA	O1-N1	2.12	1.47	1.42
2	B	3844	LDA	O1-N1	2.11	1.47	1.42
2	A	3842	LDA	O1-N1	2.11	1.47	1.42
2	B	3837	LDA	O1-N1	2.07	1.47	1.42
2	B	3848	LDA	C1-N1	-2.00	1.49	1.51

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	5136	PGV	O01-C1-C2	5.59	121.06	111.09
5	B	5136	PGV	C02-O01-C1	4.83	126.38	117.85
4	B	3888	LMT	C1-O1'-C1'	-3.57	107.58	113.68
5	B	5136	PGV	C01-O03-C19	2.46	126.10	117.12
2	B	3834	LDA	C7-C6-C5	-2.01	104.19	114.37

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	B	5136	PGV	O03-C01-C02-O01
5	B	5136	PGV	O03-C01-C02-C03
5	B	5136	PGV	C03-O11-P-O13
5	B	5136	PGV	C04-O12-P-O13
5	B	5136	PGV	C01-C02-O01-C1
5	B	5136	PGV	O01-C02-C03-O11
5	B	5136	PGV	C01-C02-C03-O11

There are no ring outliers.

10 monomers are involved in 18 short contacts:

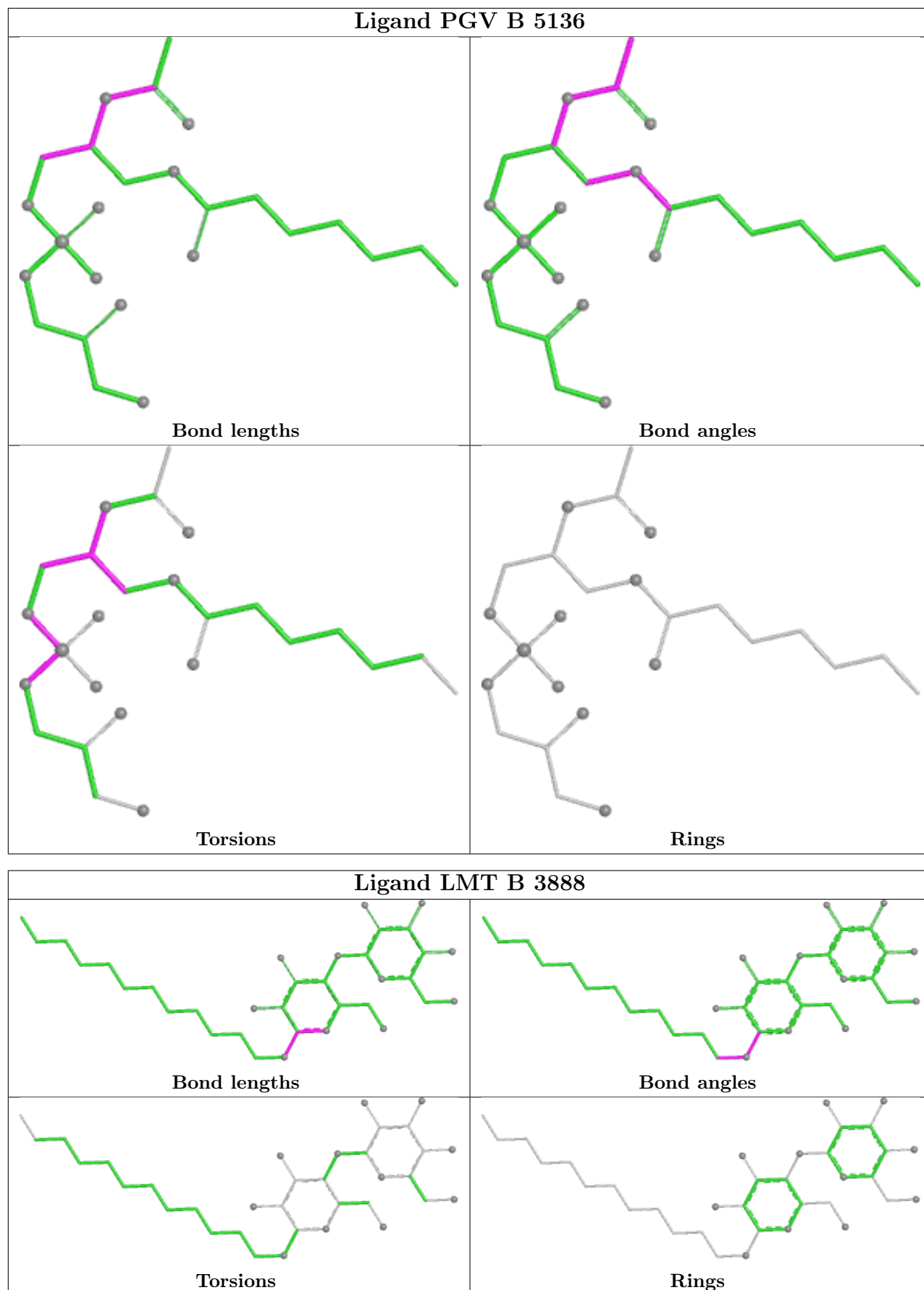
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	3850	LDA	1	0
2	B	3836	LDA	3	0
2	B	3847	LDA	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	3837	LDA	3	0
5	B	5136	PGV	4	0
2	A	3842	LDA	1	0
4	B	3888	LMT	1	0
2	B	3834	LDA	2	0
2	B	3835	LDA	1	0
2	B	3838	LDA	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	176/182 (96%)	0.31	10 (5%) 29 31	19, 28, 51, 63	0
1	B	179/182 (98%)	-0.14	3 (1%) 69 70	13, 22, 31, 41	0
All	All	355/364 (97%)	0.08	13 (3%) 45 48	13, 24, 46, 63	0

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	251	ASN	3.3
1	A	271	ASN	3.0
1	A	135	PHE	2.9
1	A	236	TRP	2.8
1	B	117	GLN	2.4
1	B	224	TYR	2.3
1	A	242	PHE	2.3
1	A	238	VAL	2.2
1	A	232	PHE	2.2
1	B	122	TRP	2.1
1	A	250	ALA	2.1
1	A	272	ALA	2.1
1	A	228	GLY	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 ⓘ

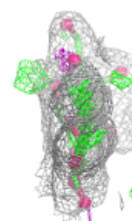
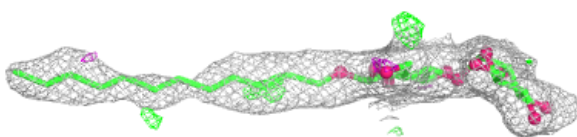
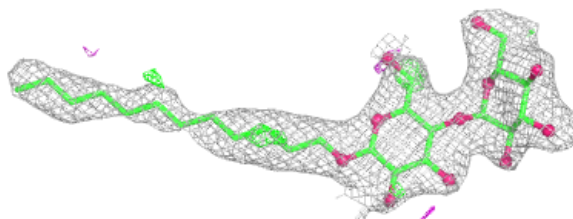
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	LDA	B	3838	12/16	0.65	0.23	42,44,52,52	0
2	LDA	B	3846	11/16	0.68	0.17	38,39,43,44	0
2	LDA	A	3845	13/16	0.71	0.19	56,58,62,62	0
2	LDA	A	3850	14/16	0.72	0.22	41,52,61,61	0
2	LDA	B	3839	9/16	0.73	0.19	55,56,57,58	0
2	LDA	B	3840	10/16	0.74	0.19	49,50,51,51	0
2	LDA	B	3848	13/16	0.75	0.16	56,58,59,60	0
2	LDA	B	3837	13/16	0.76	0.19	53,54,55,56	0
2	LDA	B	3847	10/16	0.77	0.18	33,34,36,37	0
2	LDA	B	3843	13/16	0.78	0.17	42,43,47,47	0
2	LDA	A	3842	10/16	0.78	0.21	52,53,55,56	0
2	LDA	B	3834	16/16	0.79	0.17	32,45,62,62	0
2	LDA	B	3836	16/16	0.81	0.17	39,45,52,53	0
4	LMT	B	3888	35/35	0.82	0.15	26,33,41,43	0
5	PGV	B	5136	26/51	0.82	0.17	40,48,52,53	0
2	LDA	B	3835	16/16	0.83	0.16	35,39,50,50	0
2	LDA	B	3841	11/16	0.84	0.13	39,41,44,44	0
5	PGV	B	5137	8/51	0.84	0.12	35,36,37,37	0
2	LDA	B	3844	14/16	0.86	0.13	36,37,41,41	0
5	PGV	B	5138	8/51	0.88	0.12	32,33,34,34	0
5	PGV	B	5139	7/51	0.90	0.14	38,39,39,40	0
3	PO4	B	5156	5/5	0.95	0.07	27,27,29,30	0

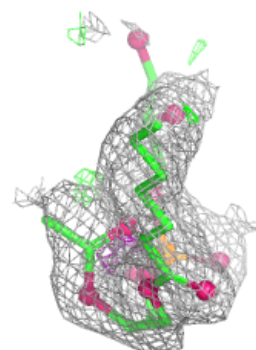
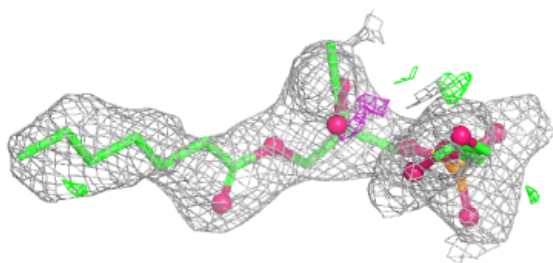
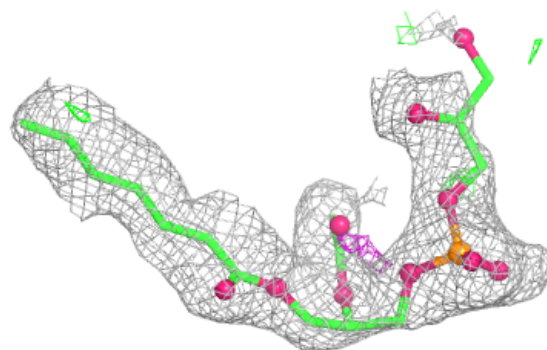
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around LMT B 3888:

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

**Electron density around PGV B 5136:**

$2mF_o-DF_c$ (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.