



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 10:52 PM UTC

PDB ID : 8H09 / pdb_00008h09
Title : Structure of the thermolabile hemolysin from *Vibrio alginolyticus* (apo form)
Authors : Ma, Q.; Wang, C.
Deposited on : 2022-09-28
Resolution : 1.81 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
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

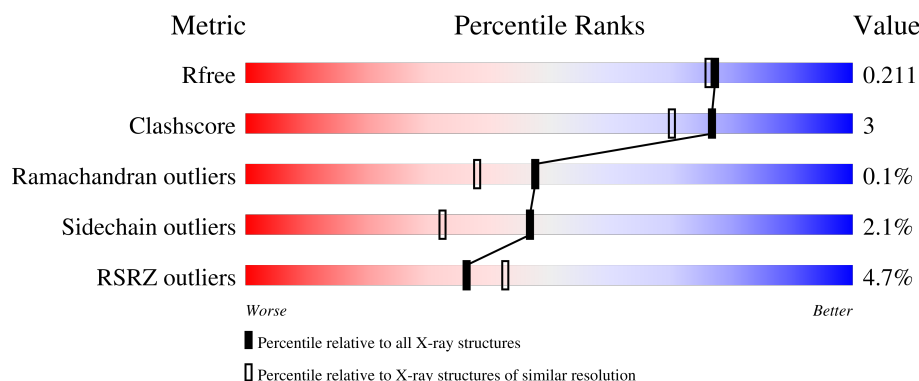
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.81 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	1112 (1.82-1.82)
Clashscore	190562	1148 (1.82-1.82)
Ramachandran outliers	187476	1140 (1.82-1.82)
Sidechain outliers	187428	1140 (1.82-1.82)
RSRZ outliers	180081	1112 (1.82-1.82)

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	426	
1	B	426	

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 6631 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.

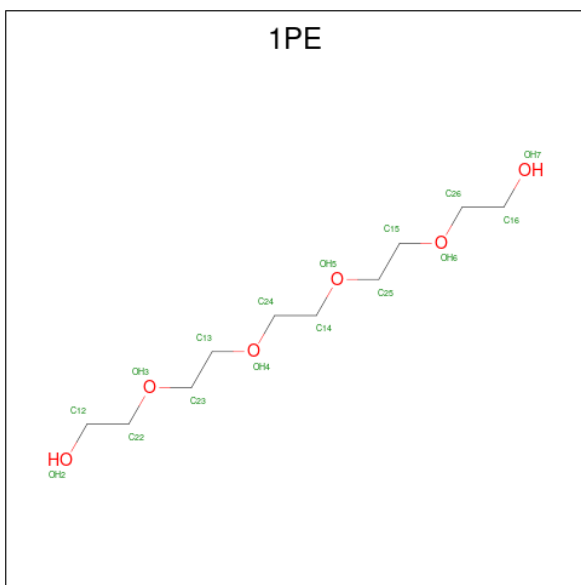
- Molecule 1 is a protein called SGNH/GDSL hydrolase family protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	378	Total	C	N	O	S	0	1	0
			3056	1941	508	594	13			
1	B	386	Total	C	N	O	S	0	0	0
			3099	1966	515	605	13			

There are 16 discrepancies between the modelled and reference sequences:

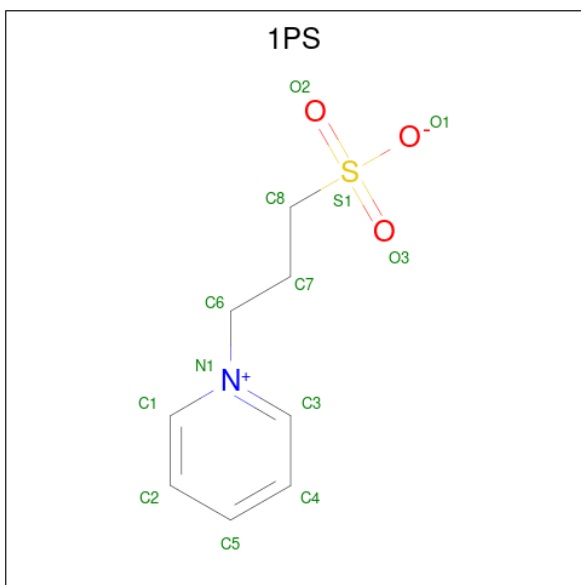
Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	MET	-	initiating methionine	UNP A0A7Y4B3E8
A	0	GLY	-	expression tag	UNP A0A7Y4B3E8
A	419	HIS	-	expression tag	UNP A0A7Y4B3E8
A	420	HIS	-	expression tag	UNP A0A7Y4B3E8
A	421	HIS	-	expression tag	UNP A0A7Y4B3E8
A	422	HIS	-	expression tag	UNP A0A7Y4B3E8
A	423	HIS	-	expression tag	UNP A0A7Y4B3E8
A	424	HIS	-	expression tag	UNP A0A7Y4B3E8
B	-1	MET	-	initiating methionine	UNP A0A7Y4B3E8
B	0	GLY	-	expression tag	UNP A0A7Y4B3E8
B	419	HIS	-	expression tag	UNP A0A7Y4B3E8
B	420	HIS	-	expression tag	UNP A0A7Y4B3E8
B	421	HIS	-	expression tag	UNP A0A7Y4B3E8
B	422	HIS	-	expression tag	UNP A0A7Y4B3E8
B	423	HIS	-	expression tag	UNP A0A7Y4B3E8
B	424	HIS	-	expression tag	UNP A0A7Y4B3E8

- Molecule 2 is PENTAETHYLENE GLYCOL (CCD ID: 1PE) (formula: C₁₀H₂₂O₆).



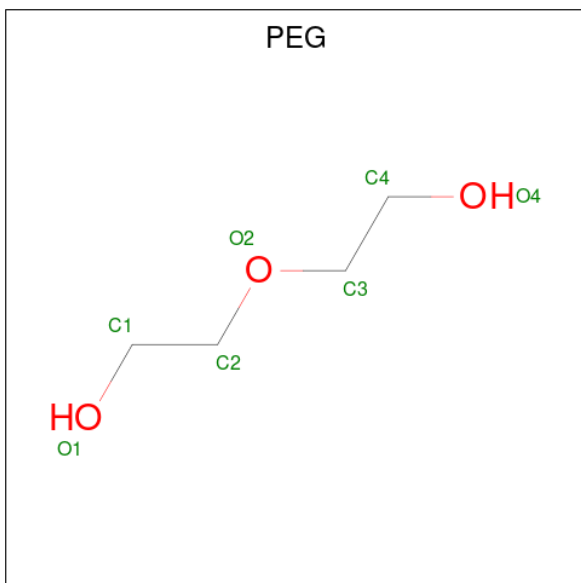
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			16	10	6		
2	A	1	Total	C	O	0	0
			16	10	6		
2	B	1	Total	C	O	0	0
			16	10	6		
2	B	1	Total	C	O	0	0
			16	10	6		
2	B	1	Total	C	O	0	0
			16	10	6		

- Molecule 3 is 3-PYRIDINIUM-1-YLPROPANE-1-SULFONATE (CCD ID: 1PS) (formula: $C_8H_{11}NO_3S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	S	0	0
			13	8	1	3	1		
3	A	1	Total	C	N	O	S	0	0
			13	8	1	3	1		
3	B	1	Total	C	N	O	S	0	0
			13	8	1	3	1		
3	B	1	Total	C	N	O	S	0	0
			13	8	1	3	1		

- Molecule 4 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C O 7 4 3	0	0
4	A	1	Total C O 7 4 3	0	0
4	A	1	Total C O 7 4 3	0	0
4	A	1	Total C O 7 4 3	0	0
4	B	1	Total C O 7 4 3	0	0

- Molecule 5 is MAGNESIUM ION (CCD ID: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total Mg 1 1	0	0
5	B	1	Total Mg 1 1	0	0

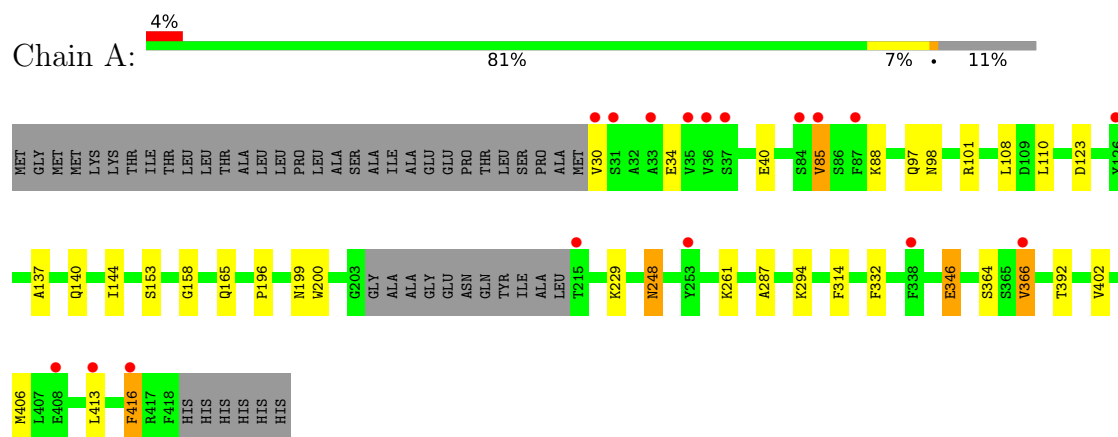
- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	168	Total O 168 168	0	0
6	B	139	Total O 139 139	0	0

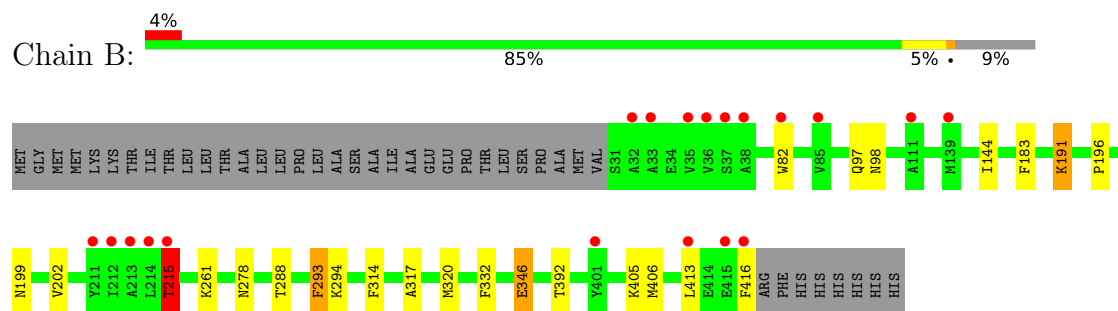
3 Residue-property plots [i](#)

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: SGNH/GDSL hydrolase family protein



- Molecule 1: SGNH/GDSL hydrolase family protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	65.93Å 71.24Å 83.64Å 90.00° 101.32° 90.00°	Depositor
Resolution (Å)	53.78 – 1.81 53.78 – 1.81	Depositor EDS
% Data completeness (in resolution range)	98.8 (53.78-1.81) 98.8 (53.78-1.81)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.17 (at 1.81Å)	Xtriage
Refinement program	BUSTER 2.10.2	Depositor
R, R_{free}	0.180 , 0.204 (Not available) , 0.211	Depositor DCC
R_{free} test set	3353 reflections (4.83%)	wwPDB-VP
Wilson B-factor (Å ²)	27.1	Xtriage
Anisotropy	0.497	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 40.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	6631	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.34% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: MG, PEG, 1PS, 1PE

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	0.82	0/3142	1.03	3/4273 (0.1%)
1	B	0.80	0/3183	1.04	4/4331 (0.1%)
All	All	0.81	0/6325	1.04	7/8604 (0.1%)

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	293	PHE	CA-CB-CG	-5.62	108.18	113.80
1	A	416	PHE	CA-CB-CG	5.19	118.99	113.80
1	B	294	LYS	N-CA-C	-5.17	105.77	111.71
1	A	199	ASN	N-CA-C	5.16	117.14	108.73
1	A	294	LYS	N-CA-C	-5.07	105.88	111.71
1	B	199	ASN	N-CA-C	5.07	116.99	108.73
1	B	215	THR	N-CA-C	5.02	116.38	109.15

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3056	0	2856	21	0
1	B	3099	0	2894	14	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	32	0	44	5	0
2	B	48	0	66	2	0
3	A	26	0	22	1	0
3	B	26	0	22	0	0
4	A	28	0	40	3	0
4	B	7	0	10	1	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
6	A	168	0	0	0	0
6	B	139	0	0	0	0
All	All	6631	0	5954	35	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 3.

All (35) 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:248:ASN:ND2	2:A:501:1PE:H132	1.95	0.80
1:B:97:GLN:HG3	1:B:196:PRO:HB3	1.71	0.71
1:A:97:GLN:HG3	1:A:196:PRO:HB3	1.72	0.71
1:A:248:ASN:HD22	2:A:501:1PE:H132	1.53	0.71
1:B:144:ILE:HD11	1:B:278:ASN:ND2	2.12	0.65
1:A:85:VAL:HG13	1:A:88:LYS:HG3	1.79	0.63
1:B:392:THR:HG21	2:B:501:1PE:H132	1.82	0.60
1:A:346:GLU:H	1:A:346:GLU:CD	2.14	0.56
1:B:215:THR:HG21	2:B:502:1PE:H122	1.86	0.56
1:A:200:TRP:CE2	4:A:505:PEG:H31	2.43	0.53
1:A:261:LYS:HE2	1:A:314:PHE:HB2	1.90	0.52
1:B:183:PHE:CE2	1:B:191:LYS:HE2	2.45	0.50
1:A:392:THR:HG21	2:A:501:1PE:H232	1.92	0.50
1:B:261:LYS:HE2	1:B:314:PHE:HB2	1.92	0.50
1:A:287:ALA:HA	2:A:501:1PE:H261	1.95	0.48
1:B:202:VAL:HG11	1:B:215:THR:HG23	1.97	0.47
1:B:346:GLU:H	1:B:346:GLU:CD	2.24	0.46
1:A:137:ALA:HA	4:A:508:PEG:H42	1.98	0.45
1:A:364:SER:OG	1:A:366:VAL:HG13	2.17	0.45
1:A:108:LEU:HB3	1:A:110:LEU:HG	1.98	0.44
1:B:191:LYS:HA	4:B:506:PEG:H21	1.97	0.44
1:A:402:VAL:O	1:A:406:MET:HG3	2.17	0.44
1:A:144:ILE:HD12	1:A:416:PHE:HB3	2.00	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:153:SER:OG	2:A:501:1PE:H131	2.18	0.43
1:A:101:ARG:NH2	4:A:507:PEG:H21	2.34	0.42
1:A:332:PHE:HB3	1:A:406:MET:HE3	2.02	0.42
1:B:413:LEU:HG	1:B:416:PHE:CD2	2.55	0.42
1:A:413:LEU:HB3	1:A:416:PHE:CD2	2.55	0.42
1:A:123:ASP:O	3:A:504:1PS:H1	2.19	0.41
1:A:30:VAL:HG12	1:A:229:LYS:HD2	2.01	0.41
1:B:288:THR:HA	1:B:293:PHE:CD1	2.55	0.41
1:B:317:ALA:HA	1:B:320:MET:HE3	2.03	0.41
1:B:413:LEU:HG	1:B:416:PHE:HD2	1.86	0.41
1:B:332:PHE:CE1	1:B:405:LYS:HB3	2.56	0.40
1:A:137:ALA:O	1:A:140:GLN:HG3	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	375/426 (88%)	362 (96%)	12 (3%)	1 (0%)	36	26
1	B	384/426 (90%)	375 (98%)	9 (2%)	0	100	100
All	All	759/852 (89%)	737 (97%)	21 (3%)	1 (0%)	48	38

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	158	GLY

5.3.2 Protein sidechains ⓘ

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	326/362 (90%)	318 (98%)	8 (2%)	42	25
1	B	328/362 (91%)	322 (98%)	6 (2%)	51	39
All	All	654/724 (90%)	640 (98%)	14 (2%)	47	32

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	34	GLU
1	A	40	GLU
1	A	85	VAL
1	A	98	ASN
1	A	165	GLN
1	A	248	ASN
1	A	346	GLU
1	A	366	VAL
1	B	82	TRP
1	B	98	ASN
1	B	191	LYS
1	B	215	THR
1	B	346	GLU
1	B	406	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	233	ASN
1	A	248	ASN
1	A	318	GLN
1	A	325	GLN
1	A	328	ASN
1	B	318	GLN
1	B	328	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 16 ligands modelled in this entry, 2 are monoatomic - leaving 14 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	1PS	B	504	-	13,13,13	0.27	0	17,17,17	0.28	0
4	PEG	A	506	-	6,6,6	0.11	0	5,5,5	0.12	0
2	1PE	A	501	-	15,15,15	0.12	0	14,14,14	0.15	0
3	1PS	A	504	-	13,13,13	0.35	0	17,17,17	0.40	0
4	PEG	A	507	-	6,6,6	0.13	0	5,5,5	0.22	0
2	1PE	B	501	-	15,15,15	0.15	0	14,14,14	0.19	0
2	1PE	A	502	-	15,15,15	0.16	0	14,14,14	0.08	0
3	1PS	B	505	-	13,13,13	0.30	0	17,17,17	0.35	0
4	PEG	A	508	-	6,6,6	0.16	0	5,5,5	0.03	0
3	1PS	A	503	-	13,13,13	0.33	0	17,17,17	0.25	0
4	PEG	B	506	-	6,6,6	0.10	0	5,5,5	0.11	0
2	1PE	B	502	-	15,15,15	0.17	0	14,14,14	0.15	0
4	PEG	A	505	-	6,6,6	0.12	0	5,5,5	0.15	0
2	1PE	B	503	-	15,15,15	0.19	0	14,14,14	0.13	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
3	1PS	B	504	-	-	0/7/7/7	0/1/1/1
4	PEG	A	506	-	-	3/4/4/4	-
2	1PE	A	501	-	-	5/13/13/13	-
3	1PS	A	504	-	-	1/7/7/7	0/1/1/1
4	PEG	A	507	-	-	1/4/4/4	-
2	1PE	B	501	-	-	5/13/13/13	-
2	1PE	A	502	-	-	1/13/13/13	-
3	1PS	B	505	-	-	3/7/7/7	0/1/1/1
4	PEG	A	508	-	-	1/4/4/4	-
3	1PS	A	503	-	-	2/7/7/7	0/1/1/1
4	PEG	B	506	-	-	3/4/4/4	-
2	1PE	B	502	-	-	1/13/13/13	-
4	PEG	A	505	-	-	3/4/4/4	-
2	1PE	B	503	-	-	3/13/13/13	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (32) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	505	1PS	C7-C8-S1-O2
3	B	505	1PS	C7-C8-S1-O3
2	A	501	1PE	OH5-C14-C24-OH4
3	A	503	1PS	C7-C8-S1-O1
4	A	507	PEG	O2-C3-C4-O4
2	A	501	1PE	OH6-C15-C25-OH5
3	A	503	1PS	C7-C8-S1-O2
3	B	505	1PS	C7-C8-S1-O1
4	A	506	PEG	C1-C2-O2-C3
2	B	501	1PE	C14-C24-OH4-C13
2	B	502	1PE	C15-C25-OH5-C14
4	B	506	PEG	C4-C3-O2-C2
4	A	505	PEG	C1-C2-O2-C3
4	B	506	PEG	C1-C2-O2-C3
2	A	501	1PE	C15-C25-OH5-C14
4	A	506	PEG	C4-C3-O2-C2

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms
2	B	501	1PE	C16-C26-OH6-C15
2	A	502	1PE	C15-C25-OH5-C14
4	A	508	PEG	C4-C3-O2-C2
2	A	501	1PE	C14-C24-OH4-C13
2	B	503	1PE	C13-C23-OH3-C22
4	B	506	PEG	O2-C3-C4-O4
2	B	501	1PE	OH4-C13-C23-OH3
4	A	505	PEG	C4-C3-O2-C2
4	A	506	PEG	O2-C3-C4-O4
2	B	503	1PE	OH6-C15-C25-OH5
4	A	505	PEG	O2-C3-C4-O4
3	A	504	1PS	C7-C6-N1-C3
2	B	503	1PE	C25-C15-OH6-C26
2	B	501	1PE	OH6-C15-C25-OH5
2	B	501	1PE	C12-C22-OH3-C23
2	A	501	1PE	OH4-C13-C23-OH3

There are no ring outliers.

8 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	501	1PE	5	0
3	A	504	1PS	1	0
4	A	507	PEG	1	0
2	B	501	1PE	1	0
4	A	508	PEG	1	0
4	B	506	PEG	1	0
2	B	502	1PE	1	0
4	A	505	PEG	1	0

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	378/426 (88%)	0.11	17 (4%) 38 44	16, 31, 56, 78	1 (0%)
1	B	386/426 (90%)	0.22	19 (4%) 35 41	20, 34, 66, 80	0
All	All	764/852 (89%)	0.16	36 (4%) 36 43	16, 32, 62, 80	1 (0%)

All (36) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	416	PHE	5.1
1	B	214	LEU	4.8
1	B	36	VAL	4.7
1	B	212	ILE	4.3
1	B	33	ALA	3.9
1	B	213	ALA	3.9
1	B	211	TYR	3.8
1	A	33	ALA	3.7
1	B	37	SER	3.3
1	B	415	GLU	3.2
1	A	215	THR	3.2
1	A	416	PHE	3.2
1	A	338	PHE	3.1
1	B	35	VAL	3.0
1	A	87	PHE	3.0
1	B	139	MET	3.0
1	A	85	VAL	2.9
1	A	253	TYR	2.9
1	B	32	ALA	2.8
1	B	215	THR	2.8
1	A	84	SER	2.8
1	B	413	LEU	2.7
1	B	82	TRP	2.7
1	A	31	SER	2.7

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	111	ALA	2.6
1	A	30	VAL	2.5
1	B	85	VAL	2.5
1	A	413	LEU	2.4
1	A	35	VAL	2.2
1	A	366	VAL	2.2
1	A	408	GLU	2.1
1	B	401	TYR	2.1
1	A	37	SER	2.1
1	B	38	ALA	2.0
1	A	126	TYR	2.0
1	A	36	VAL	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(Å ²)	Q<0.9
3	1PS	A	503	13/13	0.77	0.15	43,49,80,82	0
4	PEG	A	508	7/7	0.77	0.21	65,66,67,67	0
3	1PS	B	505	13/13	0.78	0.16	41,44,76,76	0
4	PEG	B	506	7/7	0.79	0.16	58,65,70,70	0
3	1PS	A	504	13/13	0.80	0.15	44,49,76,77	0
2	1PE	B	503	16/16	0.80	0.17	52,62,67,67	0
4	PEG	A	507	7/7	0.83	0.16	50,52,56,61	0
2	1PE	A	502	16/16	0.84	0.14	47,55,63,63	0
4	PEG	A	506	7/7	0.85	0.17	59,59,63,65	0
2	1PE	A	501	16/16	0.87	0.15	55,59,62,62	0
5	MG	A	509	1/1	0.88	0.10	47,47,47,47	0

Continued on next page...

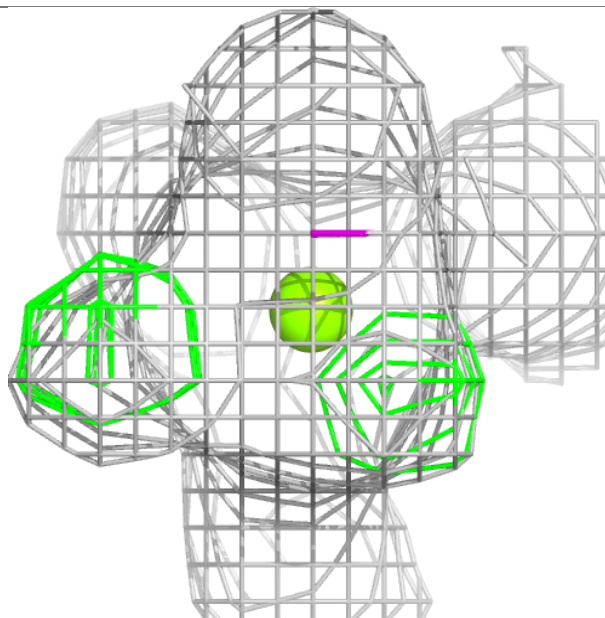
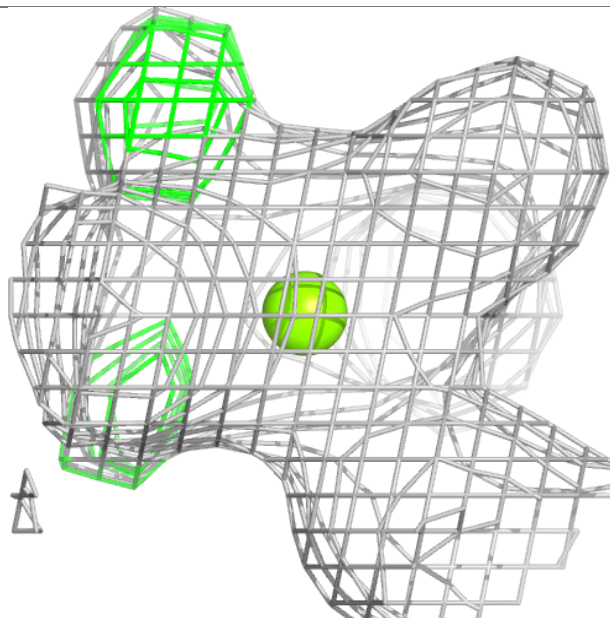
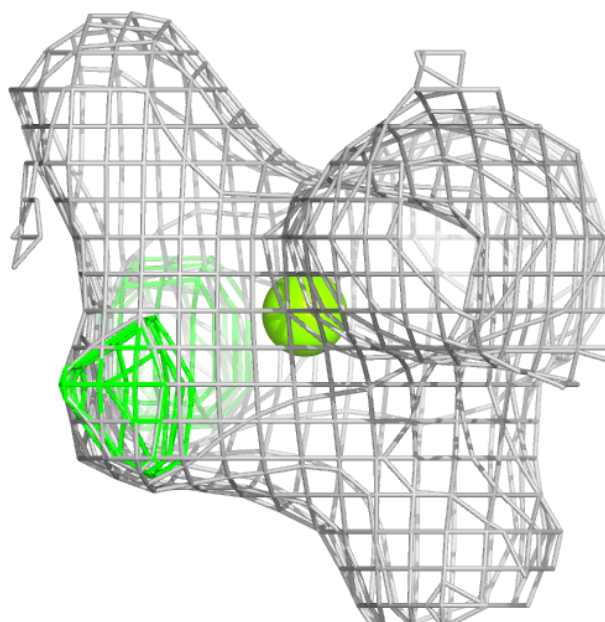
Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	PEG	A	505	7/7	0.89	0.15	53,54,60,61	0
2	1PE	B	501	16/16	0.90	0.12	42,44,51,52	0
3	1PS	B	504	13/13	0.90	0.12	44,51,65,66	0
2	1PE	B	502	16/16	0.90	0.12	40,48,56,57	0
5	MG	B	507	1/1	0.99	0.04	35,35,35,35	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.

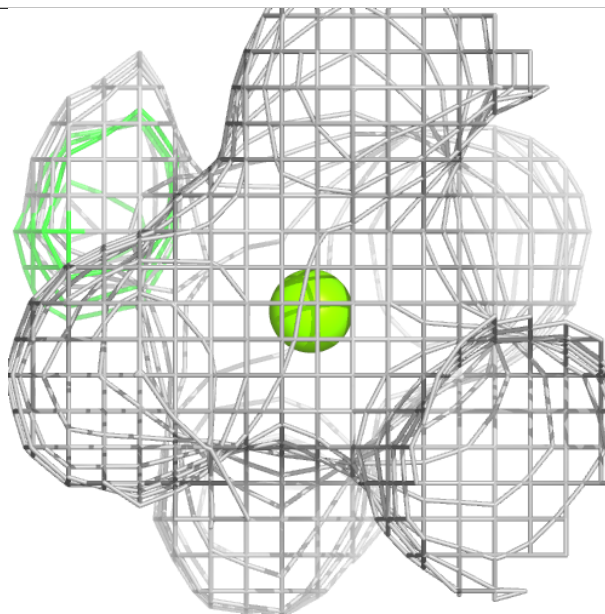
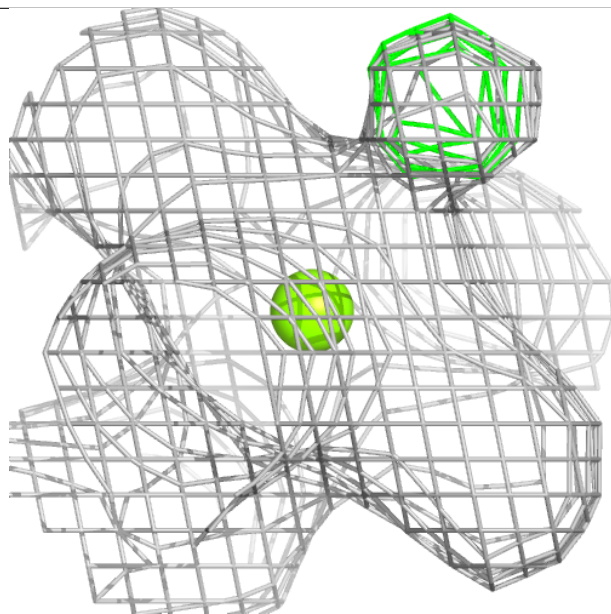
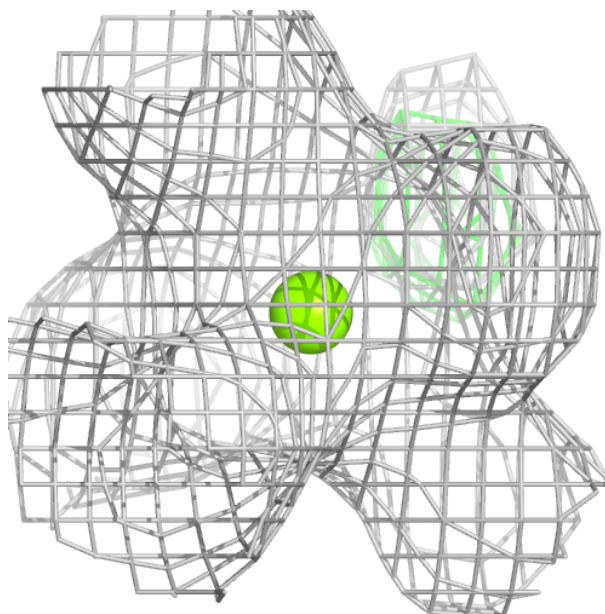
Electron density around MG A 509:

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 MG B 507:

$2mF_o-DF_c$ (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.