



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 9, 2026 – 11:02 AM UTC

PDB ID : 4EWT / pdb_00004ewt
Title : The crystal structure of a putative aminohydrolase from methicillin resistant Staphylococcus aureus
Authors : Girish, T.S.; Vivek, B.; Colaco, M.; Misquith, S.; Gopal, B.
Deposited on : 2012-04-27
Resolution : 2.10 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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

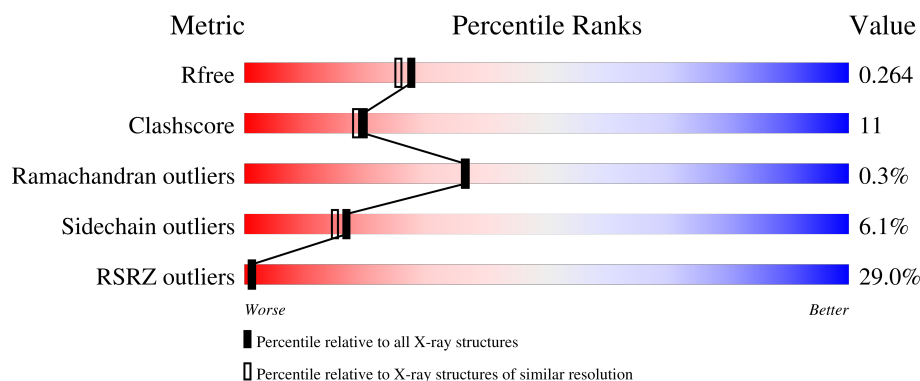
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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	392	12% 77% 20% ..
1	B	392	42% 67% 23% 5% .
1	C	392	9% 79% 17% ..
1	D	392	51% 84% 12% ..

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 12344 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.

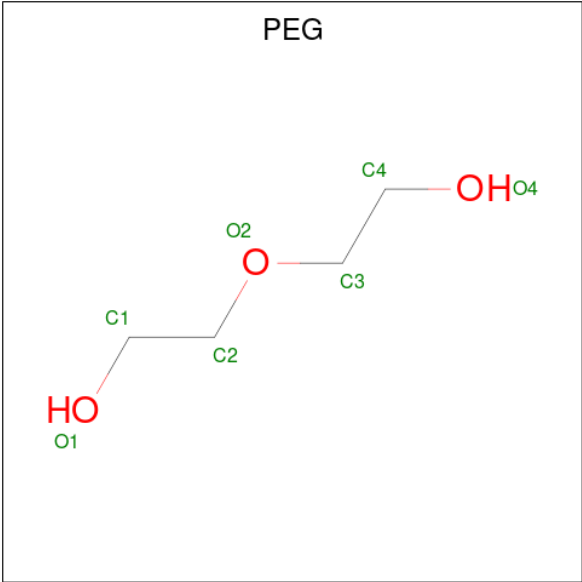
- Molecule 1 is a protein called Peptidase, M20/M25/M40 family.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	389	Total	C	N	O	S	0	1	0
			2985	1896	499	577	13			
1	C	389	Total	C	N	O	S	0	1	0
			2996	1907	496	580	13			
1	D	389	Total	C	N	O	S	0	2	0
			2980	1892	496	578	14			
1	B	375	Total	C	N	O	S	0	1	0
			2827	1795	473	546	13			

- Molecule 2 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn).

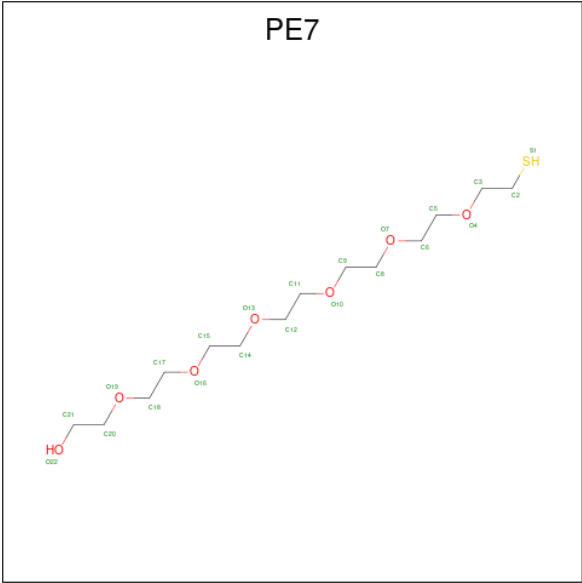
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Mn	0	0
			2	2		
2	C	2	Total	Mn	0	0
			2	2		
2	D	2	Total	Mn	0	0
			2	2		
2	B	2	Total	Mn	0	0
			2	2		

- Molecule 3 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: C₄H₁₀O₃).



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

- Molecule 4 is 1-DEOXY-1-THIO-HEPTAETHYLENE GLYCOL (CCD ID: PE7) (formula: $C_{14}H_{30}O_7S$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	D	1	Total	C	O	S	0	0
			22	14	7	1		
4	B	1	Total	C	O	S	0	0
			22	14	7	1		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	B	1	Total	C	O	S	0	0
			22	14	7	1		

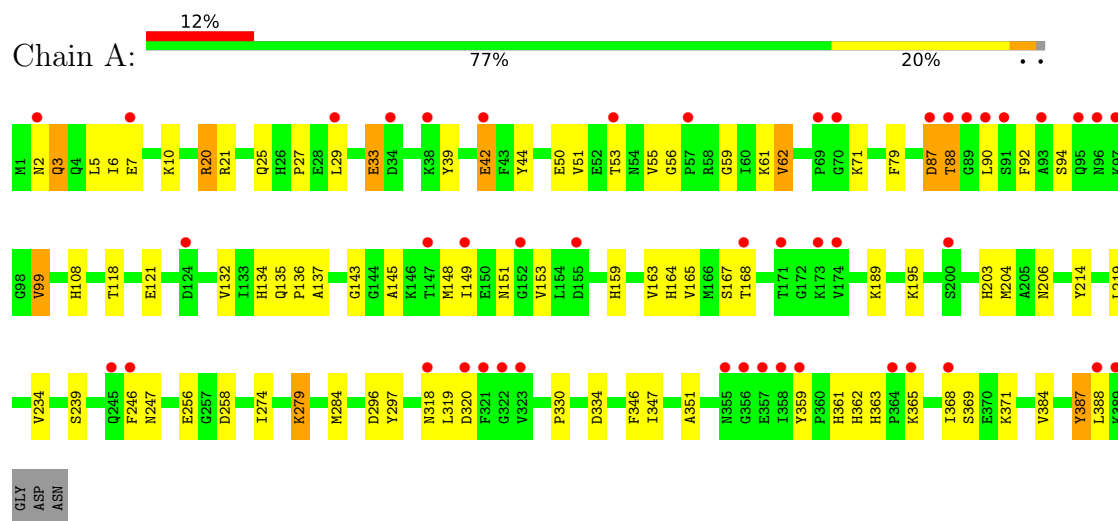
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	125	Total	O	0	0
			125	125		
5	C	127	Total	O	0	0
			127	127		
5	D	110	Total	O	0	0
			110	110		
5	B	113	Total	O	0	0
			113	113		

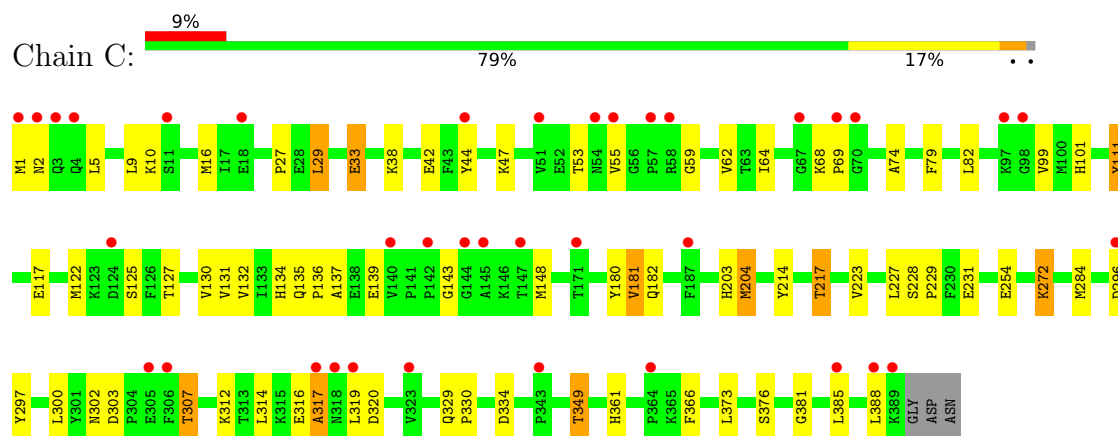
3 Residue-property plots [i](#)

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.

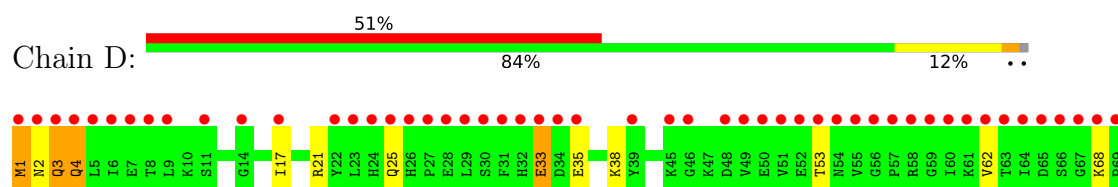
- Molecule 1: Peptidase, M20/M25/M40 family

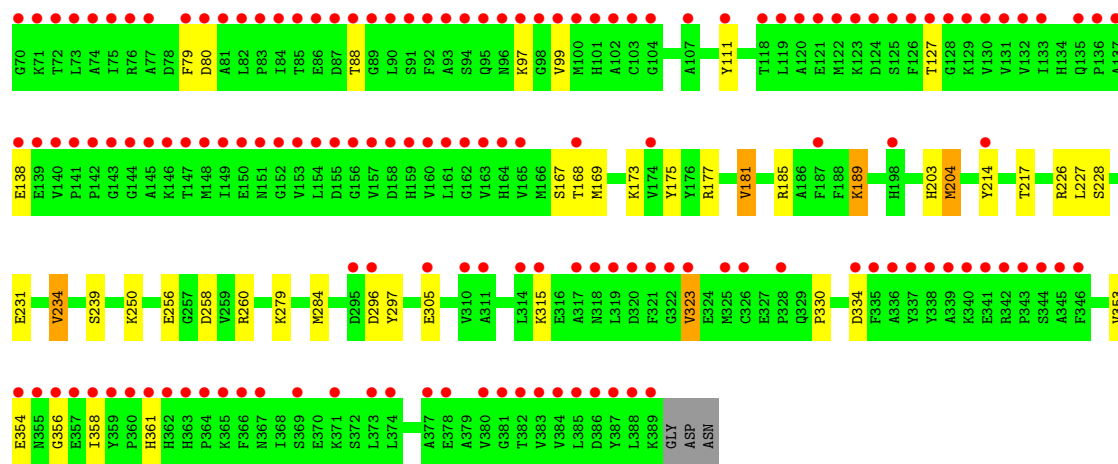


- Molecule 1: Peptidase, M20/M25/M40 family

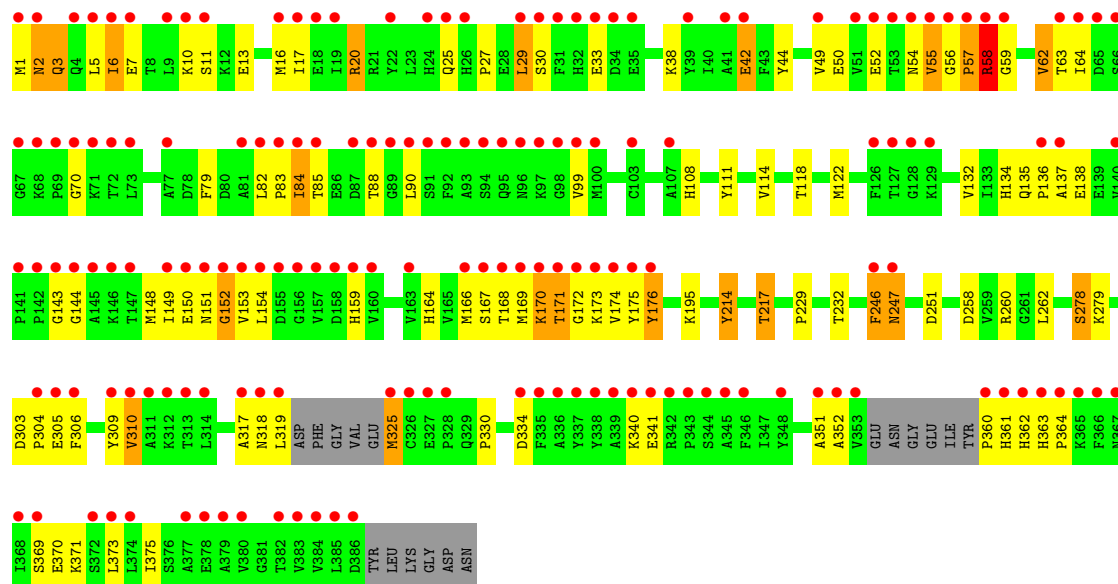
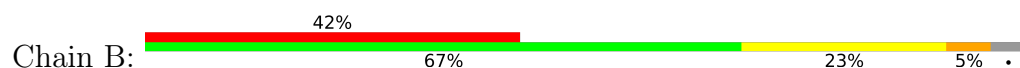


- Molecule 1: Peptidase, M20/M25/M40 family





● Molecule 1: Peptidase, M20/M25/M40 family



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	44.62Å 120.11Å 132.41Å 115.40° 94.64° 96.55°	Depositor
Resolution (Å)	39.32 – 2.10 39.32 – 2.10	Depositor EDS
% Data completeness (in resolution range)	93.8 (39.32-2.10) 93.8 (39.32-2.10)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.79 (at 2.10Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.199 , 0.229 0.241 , 0.264	Depositor DCC
R_{free} test set	7153 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	26.4	Xtriage
Anisotropy	0.057	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 36.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.016 for -h,-k,h+k+l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	12344	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.65% 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: PE7, MN, PEG

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	0.76	0/3052	0.94	8/4134 (0.2%)
1	B	1.01	4/2890 (0.1%)	1.14	19/3917 (0.5%)
1	C	0.75	4/3065 (0.1%)	0.88	7/4152 (0.2%)
1	D	0.81	0/3048	0.90	7/4130 (0.2%)
All	All	0.84	8/12055 (0.1%)	0.97	41/16333 (0.3%)

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

The worst 5 of 8 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	6	ILE	CA-CB	-7.73	1.45	1.54
1	C	223	VAL	C-O	-7.02	1.15	1.24
1	B	251	ASP	C-O	-5.38	1.17	1.24
1	C	272	LYS	CD-CE	-5.21	1.36	1.52
1	C	204	MET	C-O	-5.16	1.17	1.24

The worst 5 of 41 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	152	GLY	N-CA-C	13.38	131.12	113.99
1	B	151	ASN	N-CA-C	-11.94	93.28	110.59
1	B	58	ARG	CA-C-N	-11.62	110.30	121.57
1	B	58	ARG	C-N-CA	-11.62	110.30	121.57

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Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	88	THR	N-CA-C	-11.00	94.73	110.24

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	152	GLY	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	A	2985	0	2920	71	0
1	B	2827	0	2709	102	0
1	C	2996	0	2932	58	0
1	D	2980	0	2895	52	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
2	C	2	0	0	0	0
2	D	2	0	0	0	0
3	A	7	0	10	0	0
4	B	44	0	60	1	0
4	D	22	0	30	0	0
5	A	125	0	0	7	0
5	B	113	0	0	2	0
5	C	127	0	0	1	0
5	D	110	0	0	3	0
All	All	12344	0	11556	269	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 11.

The worst 5 of 269 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:1:MET:SD	1:D:3:GLN:HG3	1.71	1.31

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:371:LYS:O	1:B:375:ILE:HD12	1.34	1.22
1:B:2:ASN:O	1:B:6:ILE:HG12	1.37	1.21
1:A:88:THR:CG2	1:A:90:LEU:HG	1.78	1.14
1:C:82:LEU:HD21	1:C:139:GLU:HG2	1.18	1.14

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	388/392 (99%)	377 (97%)	11 (3%)	0	100	100
1	B	370/392 (94%)	343 (93%)	22 (6%)	5 (1%)	9	5
1	C	388/392 (99%)	379 (98%)	9 (2%)	0	100	100
1	D	389/392 (99%)	379 (97%)	10 (3%)	0	100	100
All	All	1535/1568 (98%)	1478 (96%)	52 (3%)	5 (0%)	36	36

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	170	LYS
1	B	171	THR
1	B	167	SER
1	B	70	GLY
1	B	172	GLY

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	316/326 (97%)	297 (94%)	19 (6%)	17	15
1	B	290/326 (89%)	273 (94%)	17 (6%)	18	16
1	C	318/326 (98%)	297 (93%)	21 (7%)	15	13
1	D	314/326 (96%)	296 (94%)	18 (6%)	18	17
All	All	1238/1304 (95%)	1163 (94%)	75 (6%)	17	15

5 of 75 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	354	GLU
1	B	325	MET
1	B	11	SER
1	B	84	ILE
1	C	55	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 23 such sidechains are listed below:

Mol	Chain	Res	Type
1	D	193	GLN
1	D	362	HIS
1	D	361	HIS
1	B	96	ASN
1	A	203	HIS

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 12 ligands modelled in this entry, 8 are monoatomic - leaving 4 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	$\# Z > 2$
4	PE7	B	404	-	21,21,21	0.52	0	19,20,20	0.54	0
4	PE7	D	403	-	21,21,21	0.48	0	19,20,20	0.77	0
4	PE7	B	403	-	21,21,21	0.52	0	19,20,20	0.67	0
3	PEG	A	403	-	6,6,6	0.42	0	5,5,5	0.31	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
4	PE7	B	404	-	-	7/19/19/19	-
4	PE7	D	403	-	-	6/19/19/19	-
4	PE7	B	403	-	-	12/19/19/19	-
3	PEG	A	403	-	-	1/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

5 of 26 torsion outliers are listed below:

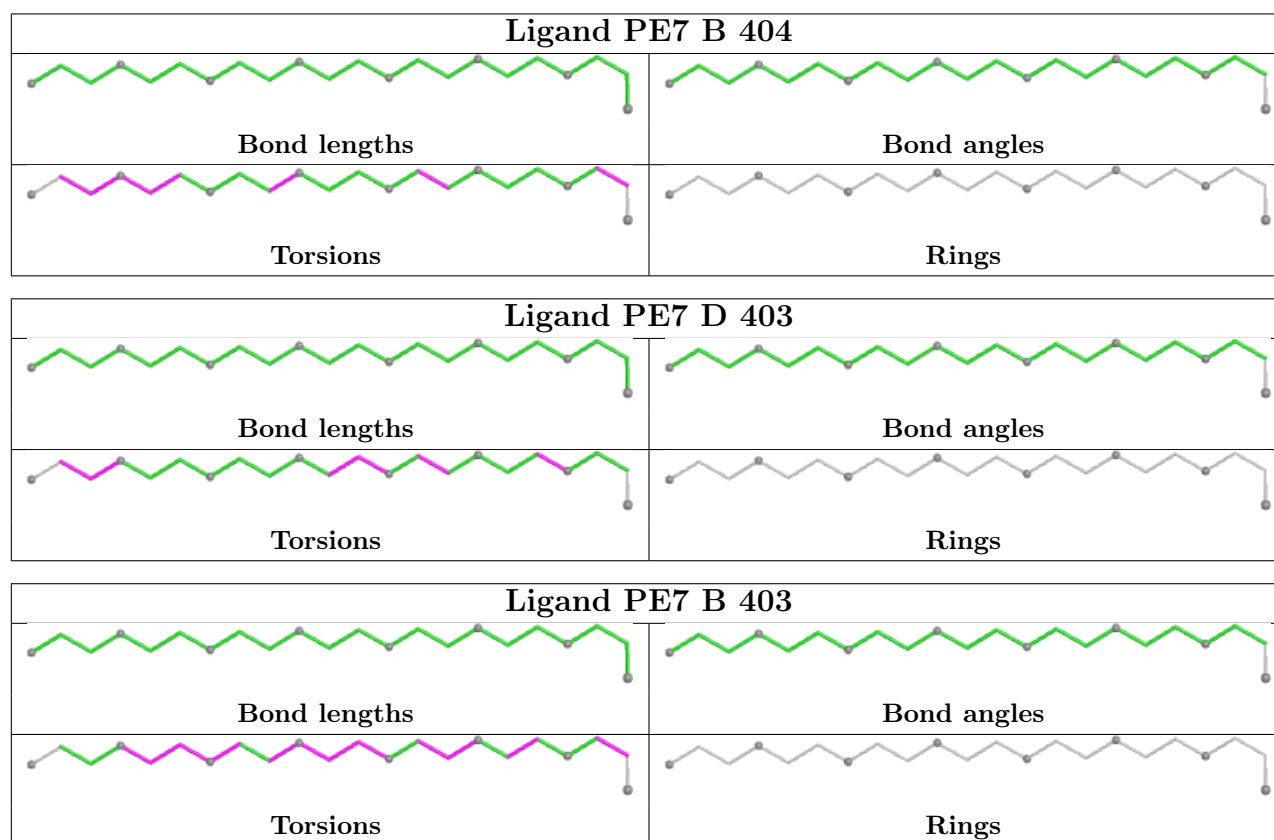
Mol	Chain	Res	Type	Atoms
4	B	403	PE7	S1-C2-C3-O4
4	B	403	PE7	O16-C17-C18-O19
4	D	403	PE7	O10-C11-C12-O13
3	A	403	PEG	O2-C3-C4-O4
4	B	403	PE7	O4-C5-C6-O7

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	403	PE7	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 [i](#)

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	389/392 (99%)	0.74	46 (11%) 9 9	10, 38, 67, 92	1 (0%)
1	B	375/392 (95%)	1.72	165 (44%) 0 1	8, 37, 64, 76	1 (0%)
1	C	389/392 (99%)	0.75	37 (9%) 14 14	8, 36, 57, 70	1 (0%)
1	D	389/392 (99%)	2.34	199 (51%) 0 0	7, 44, 91, 130	2 (0%)
All	All	1542/1568 (98%)	1.38	447 (28%) 1 1	7, 38, 71, 130	5 (0%)

The worst 5 of 447 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	389	LYS	12.3
1	D	98	GLY	9.6
1	D	343	PRO	9.0
1	D	337	TYR	8.3
1	D	156	GLY	7.7

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](#)

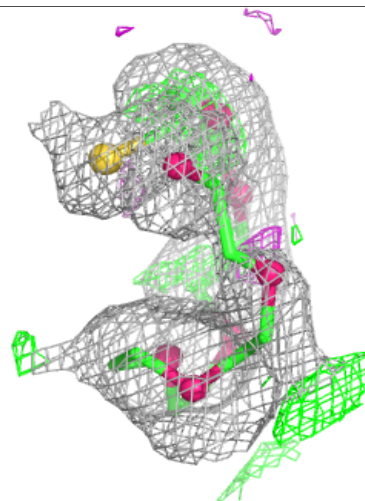
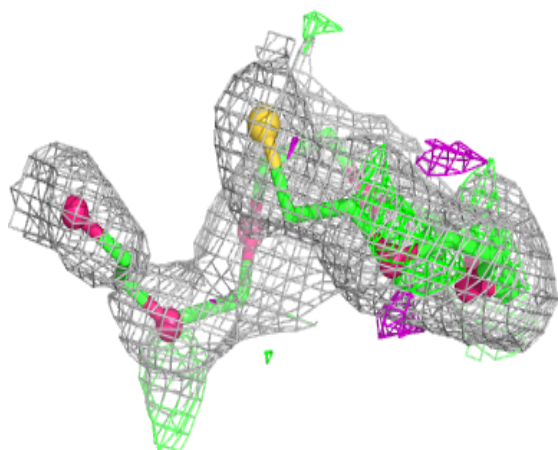
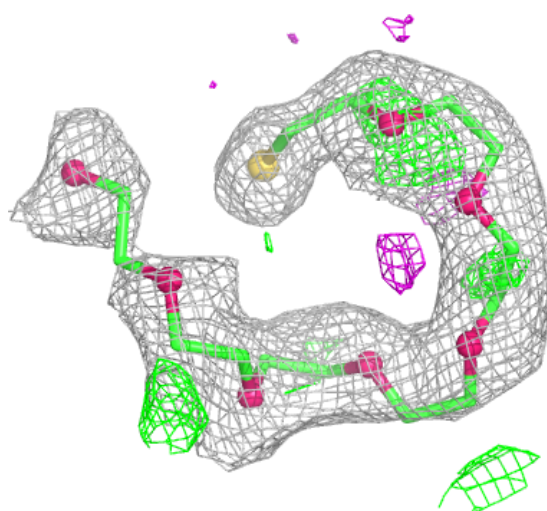
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q < 0.9’ lists 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
4	PE7	B	403	22/22	0.83	0.23	54,62,67,67	0
4	PE7	B	404	22/22	0.87	0.18	40,48,56,61	0
4	PE7	D	403	22/22	0.89	0.16	35,38,52,58	0
3	PEG	A	403	7/7	0.91	0.15	36,36,39,41	0
2	MN	B	401	1/1	0.91	0.13	69,69,69,69	0
2	MN	D	401	1/1	0.94	0.09	57,57,57,57	0
2	MN	B	402	1/1	0.95	0.06	48,48,48,48	0
2	MN	A	401	1/1	0.96	0.07	55,55,55,55	0
2	MN	C	402	1/1	0.97	0.05	35,35,35,35	0
2	MN	C	401	1/1	0.97	0.06	49,49,49,49	0
2	MN	A	402	1/1	0.98	0.06	35,35,35,35	0
2	MN	D	402	1/1	0.98	0.06	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.

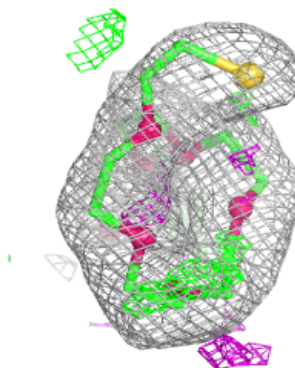
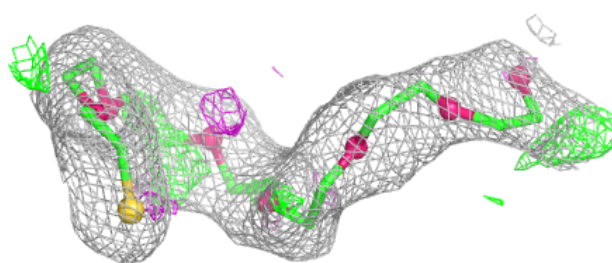
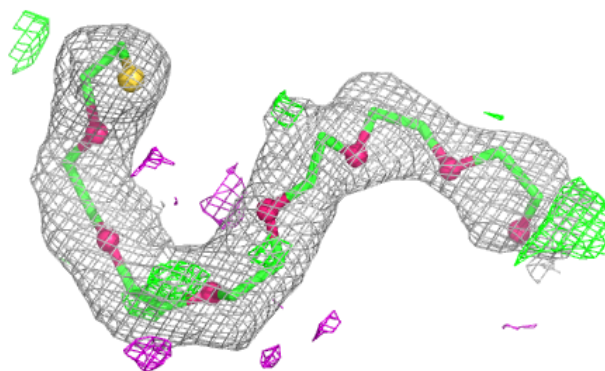
Electron density around PE7 B 403:

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

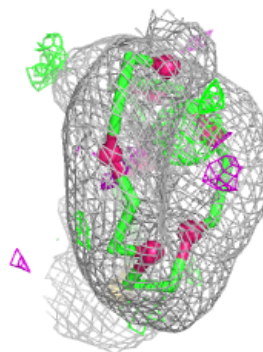
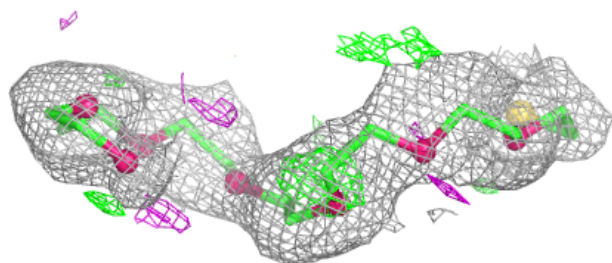
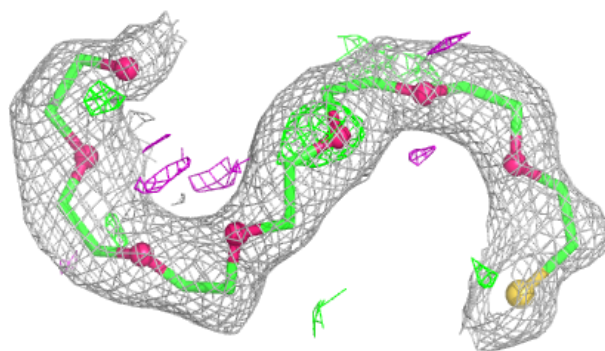


Electron density around PE7 B 404:

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

**Electron density around PE7 D 403:**

$2mF_o-DF_c$ (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.