



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 2, 2026 – 04:06 pm BST

PDB ID : 9RK9 / pdb_00009rk9
Title : Fervidolysin B (FerB) from Fervidobacterium pennivorans
Authors : Opperman, D.J.; Tolmie, C.; Kruger, M.W.
Deposited on : 2025-06-13
Resolution : 1.49 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	:	1.8.4, CSD as541be (2020)
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

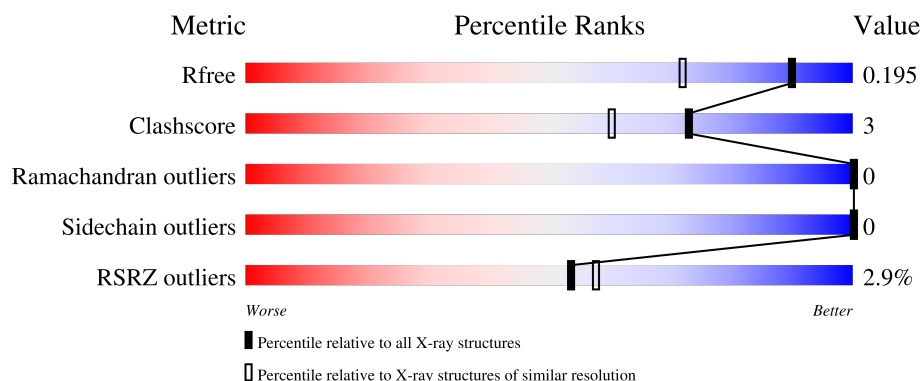
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

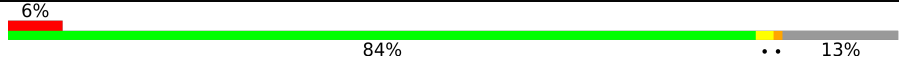
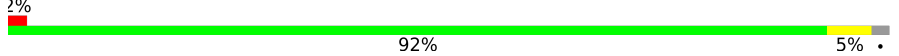
The reported resolution of this entry is 1.49 Å.

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	4037 (1.50-1.50)
Clashscore	190562	4235 (1.50-1.50)
Ramachandran outliers	187476	4153 (1.50-1.50)
Sidechain outliers	187428	4150 (1.50-1.50)
RSRZ outliers	180081	4039 (1.50-1.50)

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	97	
2	B	332	

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 3558 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 Subtilisin-like serine protease.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
1	A	84	Total	C	N	O	0	0	0
			656	416	105	135			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	initiating methionine	UNP H9UCM0

- Molecule 2 is a protein called Subtilisin-like serine protease.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	324	Total	C	N	O	S	0	2	0
			2500	1590	417	484	9			

There are 10 discrepancies between the modelled and reference sequences:

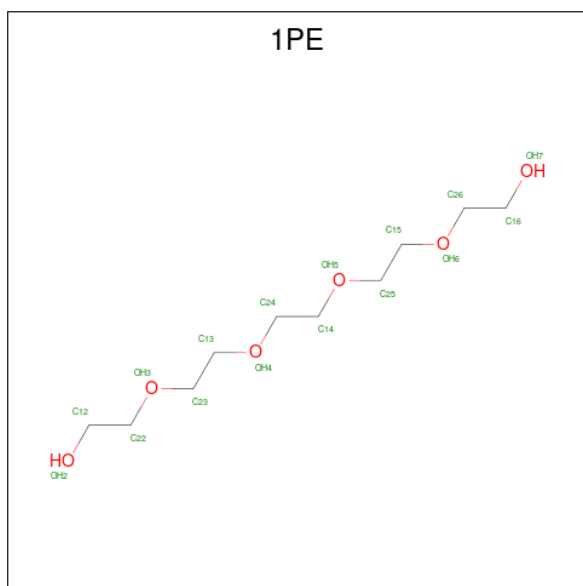
Chain	Residue	Modelled	Actual	Comment	Reference
B	420	SER	-	expression tag	UNP H9UCM0
B	421	ALA	-	expression tag	UNP H9UCM0
B	422	HIS	-	expression tag	UNP H9UCM0
B	423	HIS	-	expression tag	UNP H9UCM0
B	424	HIS	-	expression tag	UNP H9UCM0
B	425	HIS	-	expression tag	UNP H9UCM0
B	426	HIS	-	expression tag	UNP H9UCM0
B	427	HIS	-	expression tag	UNP H9UCM0
B	428	HIS	-	expression tag	UNP H9UCM0
B	429	HIS	-	expression tag	UNP H9UCM0

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



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

- Molecule 4 is PENTAETHYLENE GLYCOL (CCD ID: 1PE) (formula: $C_{10}H_{22}O_6$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			16	10	6		
4	B	1	Total	C	O	0	0
			13	8	5		

- Molecule 5 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$).

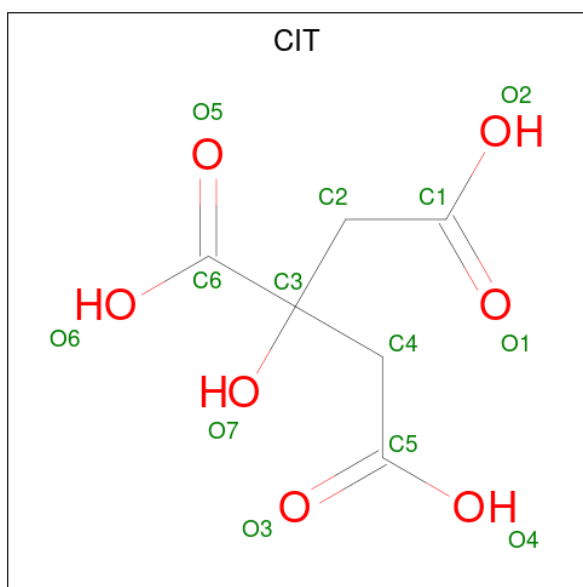


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			7	4	3		
5	B	1	Total	C	O	0	0
			7	4	3		

- Molecule 6 is CALCIUM ION (CCD ID: CA) (formula: Ca) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	2	Total	Ca	0	0
			2	2		

- Molecule 7 is CITRIC ACID (CCD ID: CIT) (formula: C₆H₈O₇).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	B	1	Total	C	O	0	0
			13	6	7		

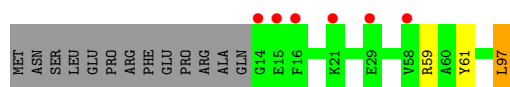
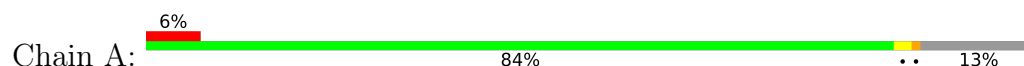
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	70	Total	O	0	0
			70	70		
8	B	257	Total	O	0	11
			268	268		

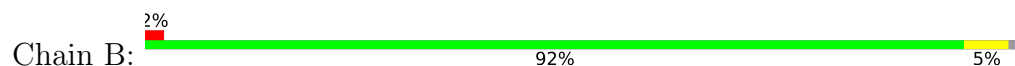
3 Residue-property plots [i](#)

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: Subtilisin-like serine protease



- Molecule 2: Subtilisin-like serine protease



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	53.57Å 71.83Å 141.07Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	64.01 – 1.49 64.01 – 1.49	Depositor EDS
% Data completeness (in resolution range)	99.9 (64.01-1.49) 99.9 (64.01-1.49)	Depositor EDS
R_{merge}	0.18	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.00 (at 1.49Å)	Xtriage
Refinement program	REFMAC 5.8.0258	Depositor
R, R_{free}	0.174 , 0.195 0.174 , 0.195	Depositor DCC
R_{free} test set	1970 reflections (2.20%)	wwPDB-VP
Wilson B-factor (Å ²)	19.7	Xtriage
Anisotropy	0.183	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 36.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	3558	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.02% 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: GOL, PEG, CA, CIT, 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	1.14	1/667 (0.1%)	1.21	0/903
2	B	1.11	3/2568 (0.1%)	1.17	1/3513 (0.0%)
All	All	1.12	4/3235 (0.1%)	1.18	1/4416 (0.0%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	97	LEU	C-O	6.61	1.36	1.23
2	B	147	ILE	C-O	6.19	1.32	1.24
2	B	173	HIS	CE1-NE2	6.17	1.38	1.32
2	B	390	TYR	C-O	5.22	1.30	1.24

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	215	ASP	CA-CB-CG	5.96	118.56	112.60

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	656	0	644	6	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	2500	0	2382	16	0
3	A	6	0	8	1	0
4	A	16	0	22	5	0
4	B	13	0	17	0	0
5	A	7	0	10	0	0
5	B	7	0	10	0	0
6	B	2	0	0	0	0
7	B	13	0	5	1	0
8	A	70	0	0	3	0
8	B	268	0	0	0	0
All	All	3558	0	3098	20	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 (20) 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:97:LEU:C	2:B:357[B]:SER:OG	2.07	0.98
1:A:59:ARG:HD3	8:A:752:HOH:O	1.86	0.74
1:A:61:TYR:HA	4:A:602:1PE:H131	1.75	0.69
2:B:235:SER:HB3	2:B:357[A]:SER:HB2	1.81	0.62
4:A:602:1PE:H162	2:B:223:TRP:HE1	1.64	0.61
2:B:235:SER:HB3	2:B:357[A]:SER:CB	2.35	0.56
4:A:602:1PE:C16	2:B:223:TRP:HE1	2.21	0.54
1:A:97:LEU:CA	2:B:357[B]:SER:OG	2.59	0.50
2:B:404:LYS:NZ	7:B:505:CIT:H22	2.27	0.50
2:B:336:GLY:HA3	2:B:352:PHE:HB3	1.98	0.46
2:B:173:HIS:HA	2:B:340:THR:O	2.17	0.45
1:A:97:LEU:HD13	2:B:263:GLY:HA2	1.99	0.44
3:A:601:GOL:H32	2:B:306:TYR:CE1	2.52	0.43
1:A:97:LEU:OXT	2:B:357[B]:SER:OG	2.36	0.43
4:A:602:1PE:H152	8:A:764:HOH:O	2.19	0.42
2:B:182:THR:HG21	2:B:193:TRP:CD1	2.55	0.41
2:B:402:TRP:CH2	2:B:404:LYS:HA	2.55	0.41
2:B:110:HIS:CD2	2:B:110:HIS:H	2.37	0.40
2:B:173:HIS:CD2	2:B:339:SER:HB3	2.57	0.40
4:A:602:1PE:C15	8:A:764:HOH:O	2.70	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	82/97 (84%)	82 (100%)	0	0	100	100
2	B	324/332 (98%)	311 (96%)	13 (4%)	0	100	100
All	All	406/429 (95%)	393 (97%)	13 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	73/85 (86%)	73 (100%)	0	100	100
2	B	256/262 (98%)	256 (100%)	0	100	100
All	All	329/347 (95%)	329 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	113	ASN
2	B	127	ASN
2	B	216	ASN
2	B	289	ASN
2	B	328	GLN
2	B	348	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 8 ligands modelled in this entry, 2 are monoatomic - leaving 6 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	1PE	B	503	-	12,12,15	0.32	0	11,11,14	0.35	0
5	PEG	B	504	-	6,6,6	0.29	0	5,5,5	0.08	0
4	1PE	A	602	-	15,15,15	0.36	0	14,14,14	0.54	0
3	GOL	A	601	-	5,5,5	0.12	0	5,5,5	0.39	0
7	CIT	B	505	-	12,12,12	1.38	1 (8%)	17,17,17	1.36	2 (11%)
5	PEG	A	603	-	6,6,6	0.11	0	5,5,5	0.16	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	1PE	B	503	-	-	6/10/10/13	-
5	PEG	B	504	-	-	0/4/4/4	-
4	1PE	A	602	-	-	7/13/13/13	-

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GOL	A	601	-	-	2/4/4/4	-
7	CIT	B	505	-	-	6/16/16/16	-
5	PEG	A	603	-	-	3/4/4/4	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	B	505	CIT	O1-C1	3.03	1.32	1.22

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	B	505	CIT	O5-C6-C3	-3.82	116.85	122.25
7	B	505	CIT	O6-C6-C3	2.24	116.95	113.05

There are no chirality outliers.

All (24) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	601	GOL	O1-C1-C2-C3
7	B	505	CIT	C1-C2-C3-C4
7	B	505	CIT	C1-C2-C3-C6
7	B	505	CIT	C2-C3-C4-C5
7	B	505	CIT	O7-C3-C4-C5
7	B	505	CIT	C6-C3-C4-C5
4	A	602	1PE	C15-C25-OH5-C14
4	A	602	1PE	OH5-C14-C24-OH4
4	A	602	1PE	OH4-C13-C23-OH3
5	A	603	PEG	O1-C1-C2-O2
4	B	503	1PE	C24-C14-OH5-C25
7	B	505	CIT	C1-C2-C3-O7
4	A	602	1PE	OH7-C16-C26-OH6
5	A	603	PEG	O2-C3-C4-O4
3	A	601	GOL	O1-C1-C2-O2
4	A	602	1PE	C13-C23-OH3-C22
5	A	603	PEG	C4-C3-O2-C2
4	B	503	1PE	C23-C13-OH4-C24
4	B	503	1PE	C15-C25-OH5-C14
4	B	503	1PE	OH5-C14-C24-OH4
4	B	503	1PE	OH4-C13-C23-OH3

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms
4	A	602	1PE	C23-C13-OH4-C24
4	A	602	1PE	OH6-C15-C25-OH5
4	B	503	1PE	C12-C22-OH3-C23

There are no ring outliers.

3 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	602	1PE	5	0
3	A	601	GOL	1	0
7	B	505	CIT	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 [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	84/97 (86%)	0.54	6 (7%)	22 24	16, 25, 44, 53	0
2	B	324/332 (97%)	-0.04	6 (1%)	66 70	8, 19, 31, 68	2 (0%)
All	All	408/429 (95%)	0.08	12 (2%)	53 58	8, 20, 39, 68	2 (0%)

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	421	ALA	6.3
2	B	346	TYR	3.7
1	A	14	GLY	3.6
2	B	308	LYS	2.6
1	A	16	PHE	2.5
2	B	420	SER	2.5
2	B	99	THR	2.3
1	A	58	VAL	2.2
1	A	15	GLU	2.2
2	B	167	ASP	2.2
1	A	29	GLU	2.2
1	A	21	LYS	2.2

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

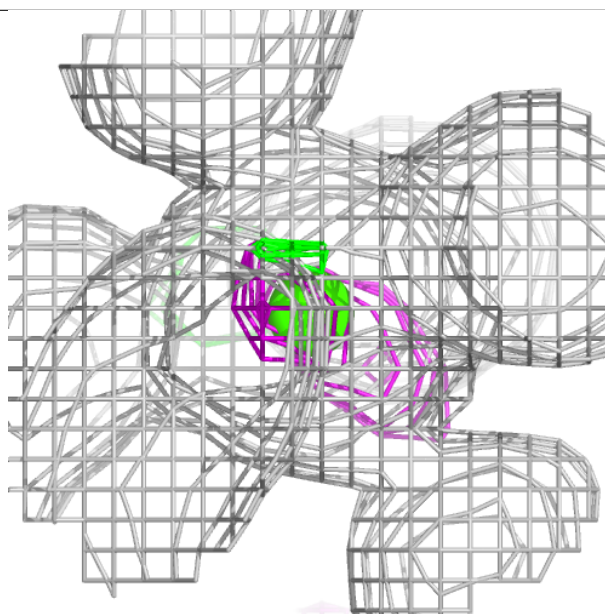
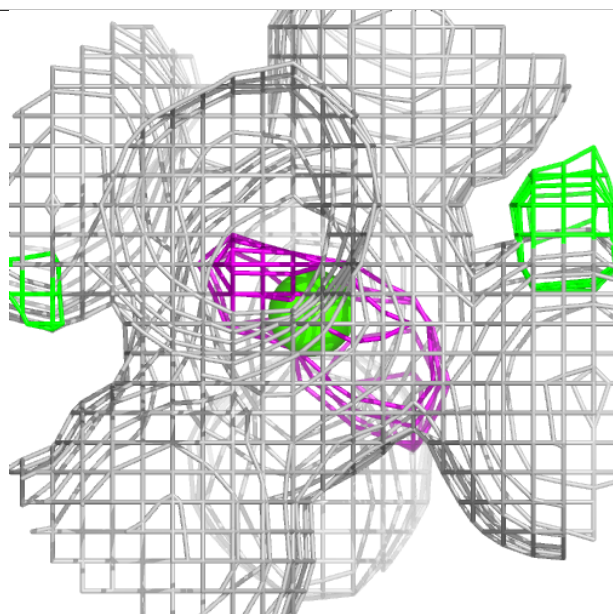
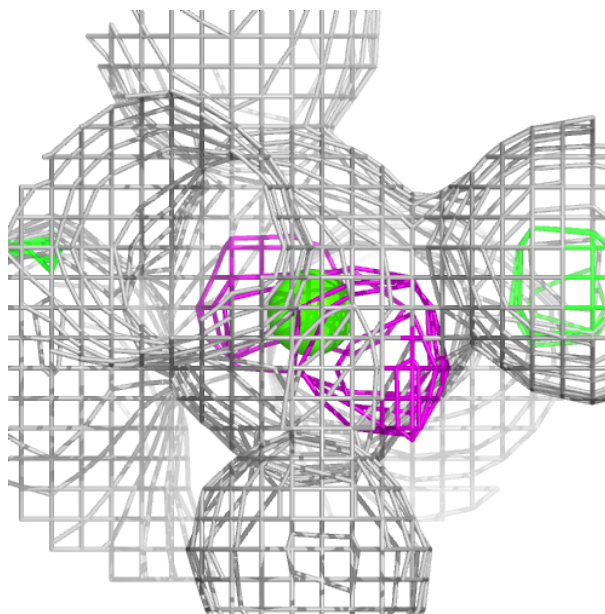
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	1PE	B	503	13/16	0.83	0.19	49,53,61,63	0
5	PEG	A	603	7/7	0.86	0.19	49,52,60,60	0
7	CIT	B	505	13/13	0.86	0.15	34,56,65,68	0
4	1PE	A	602	16/16	0.88	0.15	33,39,49,54	0
3	GOL	A	601	6/6	0.91	0.12	49,52,55,55	0
5	PEG	B	504	7/7	0.94	0.10	32,36,43,47	0
6	CA	B	502	1/1	0.99	0.06	21,21,21,21	0
6	CA	B	501	1/1	0.99	0.03	20,20,20,20	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.

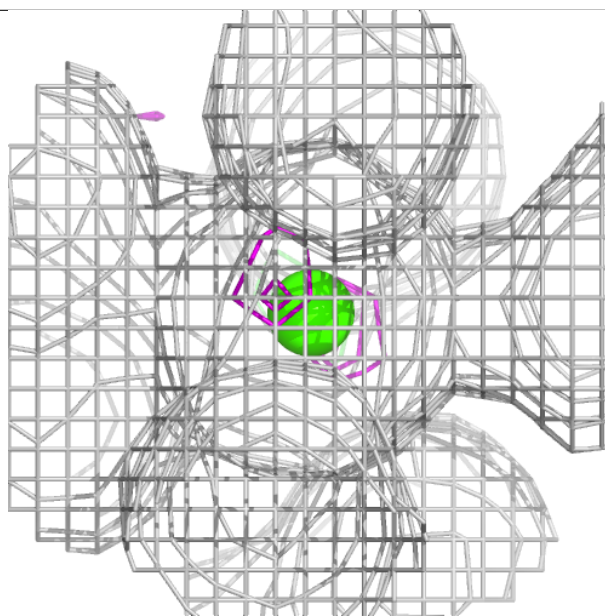
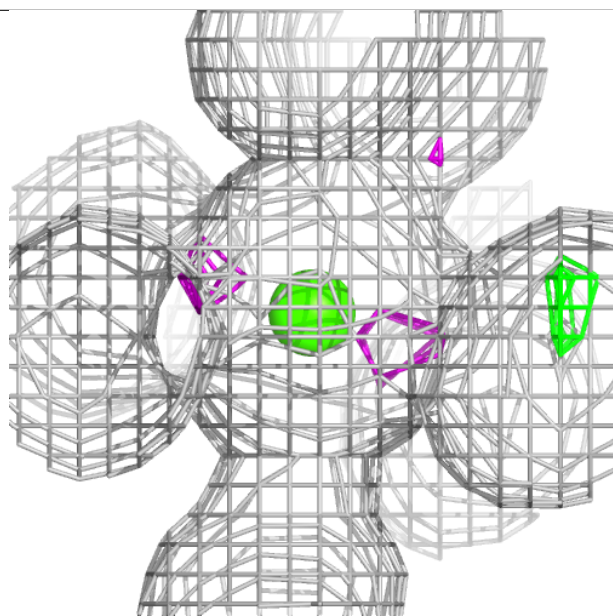
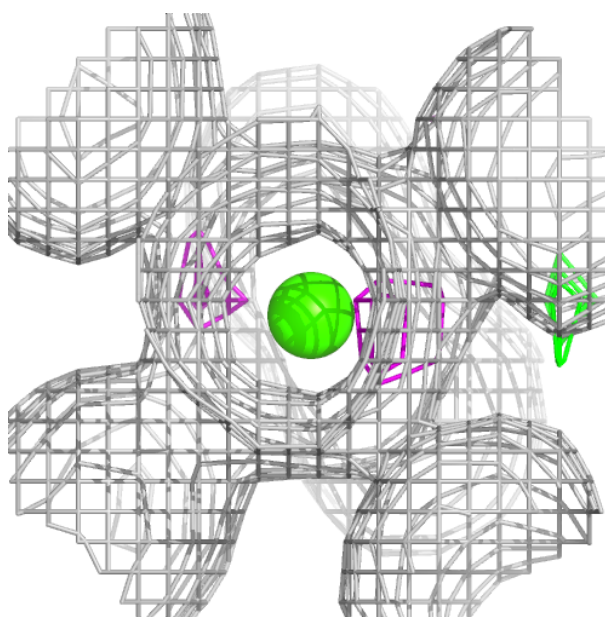
Electron density around CA B 502:

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



Electron density around CA B 501:

$2mF_o-DF_c$ (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.