



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 10:22 AM UTC

PDB ID : 6IFQ / pdb_00006ifq
Title : Crystal structure of a standalone versatile EAL protein from *Vibrio cholerae* O395 - Apo form
Authors : Yadav, M.; Pal, K.; Sen, U.
Deposited on : 2018-09-20
Resolution : 1.95 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
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

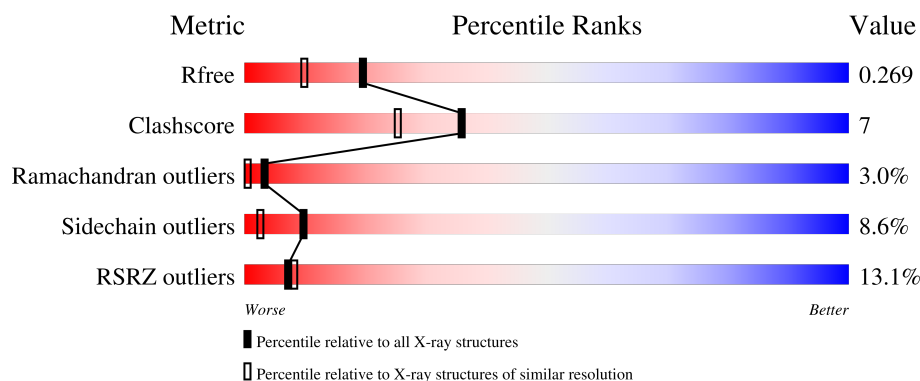
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.95 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	3494 (1.96-1.96)
Clashscore	190562	3612 (1.96-1.96)
Ramachandran outliers	187476	3587 (1.96-1.96)
Sidechain outliers	187428	3587 (1.96-1.96)
RSRZ outliers	180081	3495 (1.96-1.96)

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	257	14% 69% 18% 5% 8%
1	B	257	11% 73% 15% 8%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3844 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 cyclic di nucleotide phosphodiesterase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	237	Total	C	N	O	S	0	0	0
			1895	1215	316	354	10			
1	A	237	Total	C	N	O	S	0	0	0
			1895	1215	316	354	10			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	15	SER	CYS	engineered mutation	UNP A0A0H3AJ04
A	15	SER	CYS	engineered mutation	UNP A0A0H3AJ04

- Molecule 2 is CALCIUM ION (CCD ID: CA) (formula: Ca) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Ca	0	0
			1	1		
2	A	1	Total	Ca	0	0
			1	1		

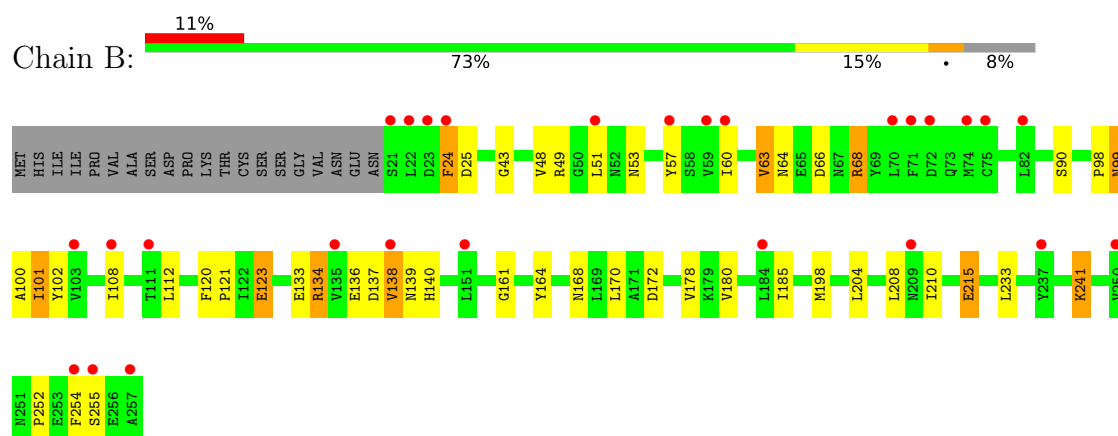
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	29	Total	O	0	0
			29	29		
3	A	23	Total	O	0	0
			23	23		

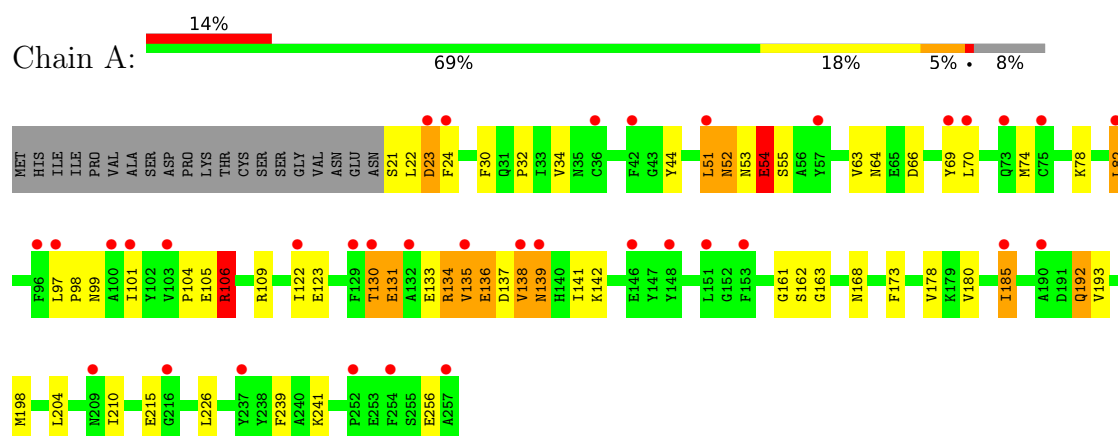
3 Residue-property plots [i](#)

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: cyclic di nucleotide phosphodiesterase



- Molecule 1: cyclic di nucleotide phosphodiesterase



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	157.70Å 42.81Å 73.57Å 90.00° 94.61° 90.00°	Depositor
Resolution (Å)	51.59 – 1.95 51.59 – 1.95	Depositor EDS
% Data completeness (in resolution range)	97.8 (51.59-1.95) 98.4 (51.59-1.95)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.26 (at 1.95Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.207 , 0.258 (Not available) , 0.269	Depositor DCC
R_{free} test set	1838 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å ²)	39.4	Xtriage
Anisotropy	0.875	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 54.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	3844	wwPDB-VP
Average B, all atoms (Å ²)	83.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.63% 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: CA

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	0.44	0/1934	0.76	0/2613
1	B	0.45	0/1934	0.80	0/2613
All	All	0.44	0/3868	0.78	0/5226

There are no bond length outliers.

There are no bond angle outliers.

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	1895	0	1871	32	0
1	B	1895	0	1871	27	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	23	0	0	0	0
3	B	29	0	0	0	0
All	All	3844	0	3742	55	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 7.

All (55) 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:137:ASP:O	1:A:139:ASN:N	2.23	0.71
1:B:180:VAL:HG11	1:B:198:MET:HE1	1.73	0.70
1:A:130:THR:OG1	1:A:131:GLU:N	2.23	0.70
1:A:135:VAL:O	1:A:137:ASP:N	2.24	0.70
1:A:180:VAL:HG11	1:A:198:MET:HE1	1.74	0.70
1:A:97:LEU:O	1:A:99:ASN:N	2.20	0.69
1:B:178:VAL:HG23	1:B:210:ILE:HG21	1.78	0.65
1:A:52:ASN:O	1:A:54:GLU:N	2.29	0.64
1:B:68:ARG:HD3	1:B:99:ASN:HD22	1.63	0.64
1:B:136:GLU:HB3	1:B:140:HIS:HB3	1.78	0.63
1:B:101:ILE:HG13	1:B:102:TYR:H	1.64	0.62
1:B:25:ASP:HB3	1:B:51:LEU:HD11	1.84	0.60
1:A:139:ASN:HA	1:A:142:LYS:HB3	1.83	0.60
1:A:52:ASN:OD1	1:A:52:ASN:N	2.20	0.60
1:A:162:SER:OG	1:A:163:GLY:N	2.33	0.59
1:A:104:PRO:O	1:A:106:ARG:N	2.35	0.57
1:A:64:ASN:HD22	1:A:66:ASP:H	1.52	0.55
1:B:137:ASP:O	1:B:139:ASN:N	2.34	0.55
1:A:74:MET:HG3	1:A:78:LYS:HE2	1.91	0.52
1:B:64:ASN:HD22	1:B:66:ASP:H	1.56	0.52
1:A:178:VAL:HG23	1:A:210:ILE:HG21	1.93	0.49
1:B:134:ARG:HB2	1:B:164:TYR:CE1	2.48	0.49
1:A:130:THR:HG1	1:A:131:GLU:H	1.60	0.49
1:A:133:GLU:O	1:A:136:GLU:HG2	2.13	0.48
1:A:30:PHE:HE2	1:A:82:LEU:HD13	1.79	0.47
1:A:134:ARG:HG3	1:A:135:VAL:N	2.30	0.47
1:A:192:GLN:CD	1:A:192:GLN:H	2.23	0.46
1:B:57:TYR:HA	1:B:60:ILE:HG22	1.97	0.46
1:B:123:GLU:H	1:B:123:GLU:CD	2.24	0.46
1:B:134:ARG:NH1	1:A:131:GLU:OE2	2.44	0.46
1:B:138:VAL:HG11	1:B:172:ASP:HB2	1.98	0.45
1:B:180:VAL:O	1:B:215:GLU:HG3	2.16	0.45
1:A:34:VAL:HG21	1:A:239:PHE:CE1	2.51	0.45
1:A:32:PRO:HB3	1:A:44:TYR:CE2	2.51	0.45
1:B:161:GLY:O	1:A:168:ASN:HB3	2.17	0.45
1:A:51:LEU:H	1:A:51:LEU:HD22	1.81	0.45
1:B:241:LYS:H	1:B:241:LYS:HD2	1.82	0.44
1:A:185:ILE:HG13	1:A:226:LEU:HD11	1.97	0.44
1:A:22:LEU:HD12	1:A:70:LEU:HD22	1.98	0.44
1:B:48:VAL:O	1:B:49:ARG:HD2	2.17	0.44
1:B:168:ASN:HB3	1:A:161:GLY:O	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:108:ILE:O	1:B:112:LEU:HG	2.18	0.44
1:A:22:LEU:HB3	1:A:23:ASP:H	1.58	0.43
1:B:208:LEU:HD23	1:B:208:LEU:HA	1.87	0.43
1:B:24:PHE:HZ	1:B:63:VAL:HG12	1.83	0.42
1:A:69:TYR:HB2	1:A:101:ILE:HG23	2.00	0.42
1:B:252:PRO:HA	1:B:255:SER:OG	2.19	0.42
1:A:52:ASN:HB2	1:A:55:SER:OG	2.20	0.42
1:B:98:PRO:O	1:B:100:ALA:N	2.53	0.42
1:A:142:LYS:HG3	1:A:173:PHE:HB3	2.01	0.41
1:B:43:GLY:HA2	1:B:90:SER:OG	2.20	0.41
1:A:82:LEU:HD23	1:A:82:LEU:HA	1.86	0.41
1:B:241:LYS:H	1:B:241:LYS:CD	2.34	0.41
1:B:208:LEU:HD21	1:A:193:VAL:HG13	2.01	0.40
1:B:120:PHE:HA	1:B:121:PRO:HD3	1.87	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	235/257 (91%)	209 (89%)	15 (6%)	11 (5%)	2	0
1	B	235/257 (91%)	218 (93%)	14 (6%)	3 (1%)	9	3
All	All	470/514 (91%)	427 (91%)	29 (6%)	14 (3%)	3	0

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	101	ILE
1	A	53	ASN
1	A	54	GLU
1	A	136	GLU

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Mol	Chain	Res	Type
1	A	138	VAL
1	A	105	GLU
1	A	131	GLU
1	A	98	PRO
1	A	106	ARG
1	B	99	ASN
1	A	23	ASP
1	A	24	PHE
1	A	135	VAL
1	B	138	VAL

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	204/222 (92%)	183 (90%)	21 (10%)	7	1
1	B	204/222 (92%)	190 (93%)	14 (7%)	14	5
All	All	408/444 (92%)	373 (91%)	35 (9%)	10	2

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	24	PHE
1	B	53	ASN
1	B	63	VAL
1	B	68	ARG
1	B	123	GLU
1	B	133	GLU
1	B	134	ARG
1	B	170	LEU
1	B	185	ILE
1	B	204	LEU
1	B	215	GLU
1	B	233	LEU
1	B	241	LYS

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Mol	Chain	Res	Type
1	B	254	PHE
1	A	21	SER
1	A	51	LEU
1	A	52	ASN
1	A	54	GLU
1	A	63	VAL
1	A	82	LEU
1	A	106	ARG
1	A	109	ARG
1	A	122	ILE
1	A	123	GLU
1	A	130	THR
1	A	134	ARG
1	A	138	VAL
1	A	139	ASN
1	A	141	ILE
1	A	185	ILE
1	A	192	GLN
1	A	204	LEU
1	A	215	GLU
1	A	241	LYS
1	A	256	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	53	ASN
1	B	64	ASN
1	B	99	ASN
1	B	154	GLN
1	B	174	GLN
1	B	251	ASN
1	A	64	ASN
1	A	154	GLN
1	A	192	GLN

5.3.3 RNA ⓘ

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](#)

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	237/257 (92%)	0.95	35 (14%) 5 6	41, 70, 157, 175	0
1	B	237/257 (92%)	0.91	27 (11%) 10 11	40, 68, 152, 180	0
All	All	474/514 (92%)	0.93	62 (13%) 7 8	40, 70, 155, 180	0

All (62) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	254	PHE	5.4
1	B	257	ALA	4.5
1	A	237	TYR	4.3
1	A	252	PRO	4.1
1	A	257	ALA	4.0
1	A	132	ALA	4.0
1	B	70	LEU	3.9
1	A	96	PHE	3.8
1	B	184	LEU	3.7
1	A	103	VAL	3.7
1	A	24	PHE	3.7
1	B	138	VAL	3.5
1	A	254	PHE	3.5
1	B	59	VAL	3.5
1	B	237	TYR	3.4
1	B	23	ASP	3.3
1	A	209	ASN	3.0
1	A	153	PHE	2.9
1	B	135	VAL	2.9
1	A	148	TYR	2.9
1	A	51	LEU	2.9
1	A	185	ILE	2.9
1	B	108	ILE	2.8
1	B	60	ILE	2.7

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Mol	Chain	Res	Type	RSRZ
1	B	250	VAL	2.6
1	A	138	VAL	2.6
1	B	22	LEU	2.6
1	A	129	PHE	2.5
1	B	255	SER	2.5
1	A	139	ASN	2.5
1	A	97	LEU	2.4
1	A	216	GLY	2.4
1	B	51	LEU	2.4
1	B	82	LEU	2.4
1	B	24	PHE	2.3
1	A	130	THR	2.3
1	A	146	GLU	2.3
1	A	135	VAL	2.3
1	A	101	ILE	2.3
1	B	72	ASP	2.3
1	B	103	VAL	2.3
1	A	122	ILE	2.3
1	B	74	MET	2.3
1	B	151	LEU	2.3
1	A	190	ALA	2.2
1	B	209	ASN	2.2
1	A	23	ASP	2.2
1	B	71	PHE	2.2
1	B	21	SER	2.2
1	A	69	TYR	2.1
1	B	111	THR	2.1
1	A	75	CYS	2.1
1	A	70	LEU	2.1
1	A	82	LEU	2.1
1	A	100	ALA	2.1
1	A	42	PHE	2.1
1	B	57	TYR	2.0
1	A	73	GLN	2.0
1	B	75	CYS	2.0
1	A	36	CYS	2.0
1	A	151	LEU	2.0
1	A	57	TYR	2.0

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

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](#)

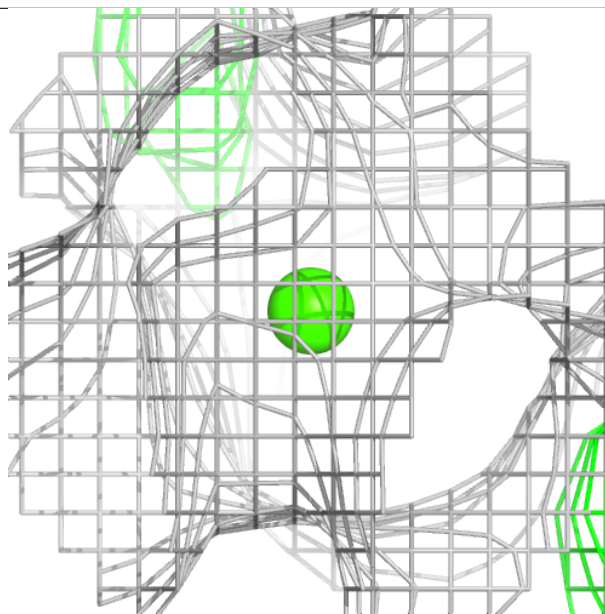
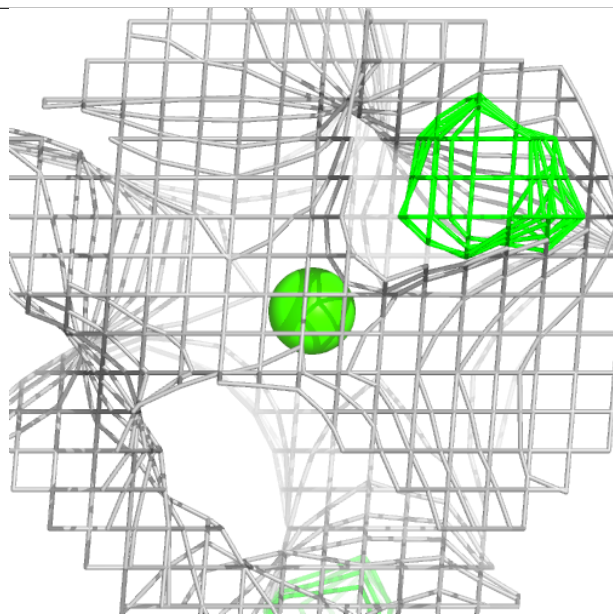
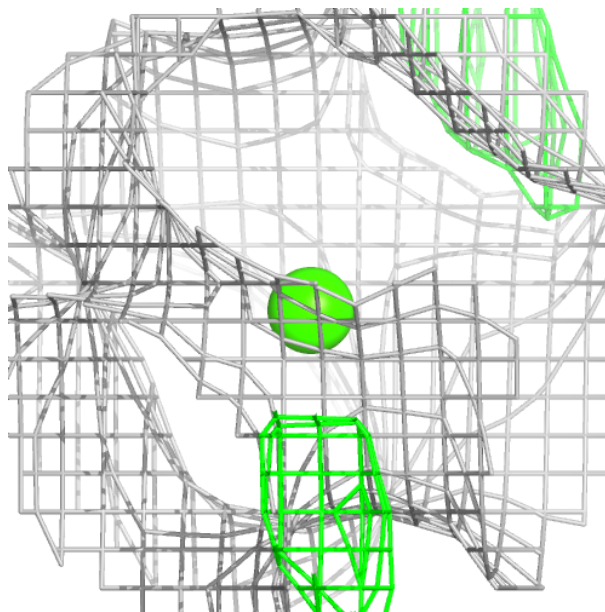
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
2	CA	B	301	1/1	0.97	0.05	59,59,59,59	0
2	CA	A	501	1/1	0.97	0.13	64,64,64,64	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.

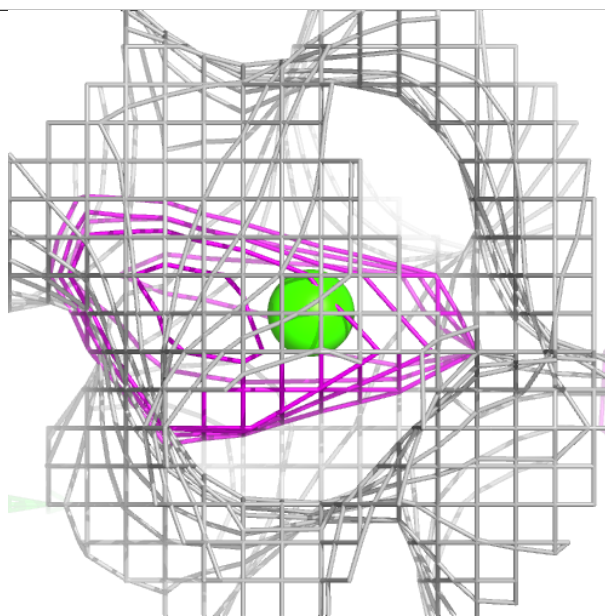
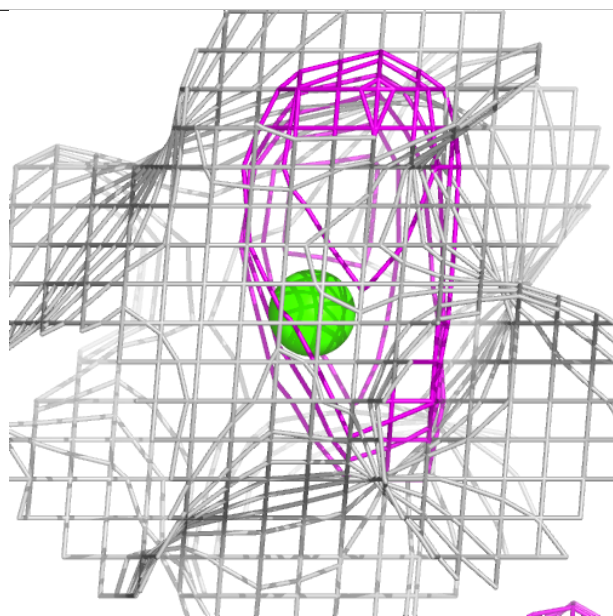
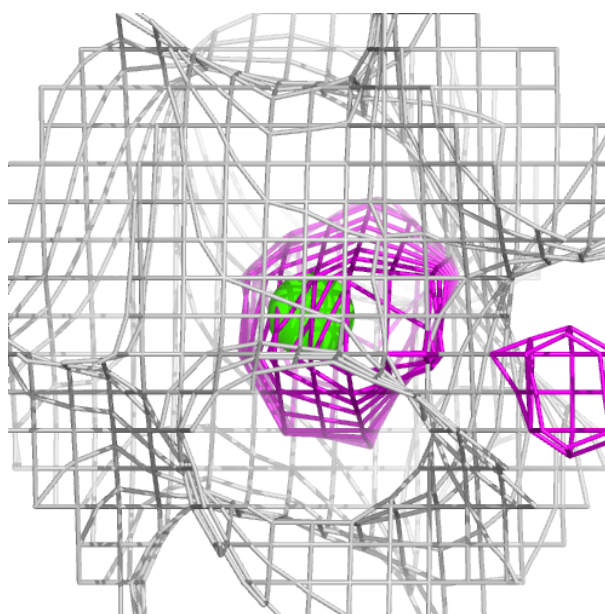
Electron density around CA B 301:

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



Electron density around CA A 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.