



Full wwPDB X-ray Structure Validation Report ⓘ

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PDB ID : 3R5S
Title : Crystal structure of apo-ViuP
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Deposited on : 2011-03-19
Resolution : 1.79 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

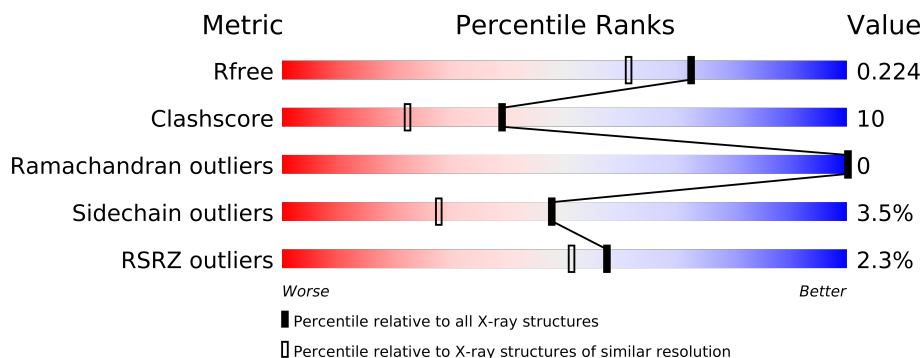
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 1.79 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}	66092	3513 (1.80-1.80)
Clashscore	79885	4461 (1.80-1.80)
Ramachandran outliers	78287	4404 (1.80-1.80)
Sidechain outliers	78261	4403 (1.80-1.80)
RSRZ outliers	66119	3515 (1.80-1.80)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	305	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2713 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Ferric vibriobactin ABC transporter, periplasmic ferric vibriobactin-binding protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	301	Total	C	N	O	Se	0	0	0
			2369	1503	418	445	3			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	326	LEU	-	EXPRESSION TAG	UNP Q9RCF6
A	327	GLU	-	EXPRESSION TAG	UNP Q9RCF6
A	328	HIS	-	EXPRESSION TAG	UNP Q9RCF6
A	329	HIS	-	EXPRESSION TAG	UNP Q9RCF6
A	330	HIS	-	EXPRESSION TAG	UNP Q9RCF6
A	331	HIS	-	EXPRESSION TAG	UNP Q9RCF6
A	332	HIS	-	EXPRESSION TAG	UNP Q9RCF6
A	333	HIS	-	EXPRESSION TAG	UNP Q9RCF6

- Molecule 2 is water.

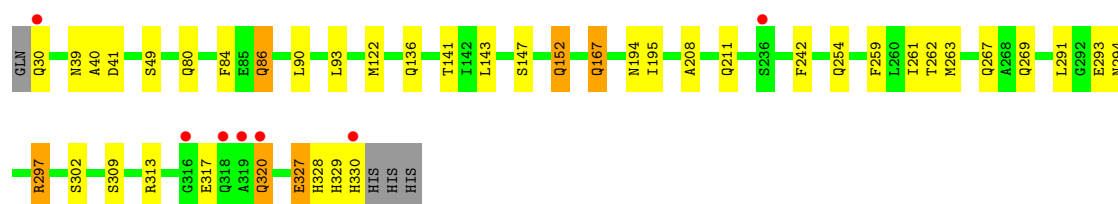
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	344	Total	O	0	0
			344	344		

3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: Ferric vibriobactin ABC transporter, periplasmic ferric vibriobactin-binding protein

Chain A: 



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	76.92Å 61.07Å 66.85Å 90.00° 120.56° 90.00°	Depositor
Resolution (Å)	33.12 – 1.79 33.12 – 1.79	Depositor EDS
% Data completeness (in resolution range)	94.8 (33.12-1.79) 94.6 (33.12-1.79)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.42 (at 1.79Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.6.2_432)	Depositor
R, R_{free}	0.177 , 0.213 0.190 , 0.224	Depositor DCC
R_{free} test set	1910 reflections (7.98%)	DCC
Wilson B-factor (Å ²)	13.0	Xtriage
Anisotropy	0.199	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 29.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 25159 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	2713	wwPDB-VP
Average B, all atoms (Å ²)	15.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.44% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

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	0.39	0/2417	0.52	0/3291

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2369	0	2366	46	0
2	A	344	0	0	8	0
All	All	2713	0	2366	46	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 10.

All (46) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:167:GLN:HE21	1:A:167:GLN:H	1.17	0.91
1:A:86:GLN:HG2	1:A:302:SER:HB3	1.56	0.88
1:A:152:GLN:HE21	1:A:152:GLN:H	1.27	0.82
1:A:297:ARG:HG3	1:A:297:ARG:HH21	1.45	0.80

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:328:HIS:ND1	1:A:329:HIS:HD2	1.87	0.73
1:A:122:MSE:HE2	1:A:143:LEU:HD21	1.72	0.71
1:A:211:GLN:NE2	2:A:479:HOH:O	2.25	0.69
1:A:122:MSE:HE2	1:A:143:LEU:CD2	2.23	0.69
1:A:167:GLN:NE2	1:A:167:GLN:H	1.88	0.68
1:A:313:ARG:HD3	2:A:348:HOH:O	1.96	0.66
1:A:86:GLN:H	1:A:86:GLN:NE2	1.94	0.66
1:A:41:ASP:N	1:A:122:MSE:HE1	2.11	0.65
1:A:86:GLN:HG3	1:A:294:ASN:HB3	1.77	0.65
1:A:40:ALA:HB3	1:A:122:MSE:HE3	1.76	0.64
1:A:40:ALA:HB3	1:A:122:MSE:CE	2.28	0.63
1:A:297:ARG:CG	1:A:297:ARG:HH21	2.12	0.63
1:A:147:SER:HA	1:A:297:ARG:NH1	2.14	0.62
1:A:80:GLN:HG3	2:A:486:HOH:O	2.02	0.59
1:A:86:GLN:H	1:A:86:GLN:HE21	1.49	0.57
1:A:262:THR:O	1:A:262:THR:HG22	2.07	0.55
1:A:269:GLN:HG3	2:A:435:HOH:O	2.06	0.55
1:A:86:GLN:HG3	1:A:294:ASN:CB	2.39	0.52
1:A:147:SER:HA	1:A:297:ARG:HH11	1.75	0.52
1:A:136:GLN:HE22	1:A:141:THR:H	1.57	0.51
1:A:267:GLN:NE2	1:A:293:GLU:OE2	2.40	0.50
1:A:167:GLN:HE21	1:A:167:GLN:N	1.98	0.50
1:A:84:PHE:HB3	1:A:86:GLN:HE22	1.78	0.48
1:A:254:GLN:HE21	1:A:254:GLN:HA	1.78	0.48
1:A:330:HIS:HA	2:A:522:HOH:O	2.13	0.48
1:A:259:PHE:HB3	1:A:291:LEU:HG	1.95	0.48
1:A:136:GLN:NE2	1:A:141:THR:H	2.13	0.47
1:A:254:GLN:NE2	1:A:254:GLN:HA	2.30	0.46
1:A:40:ALA:C	1:A:122:MSE:HE1	2.36	0.46
1:A:261:ILE:HD13	1:A:291:LEU:HD12	1.97	0.45
1:A:263:MSE:HA	1:A:267:GLN:NE2	2.31	0.45
1:A:122:MSE:CE	1:A:143:LEU:HG	2.47	0.45
1:A:49:SER:HB2	2:A:610:HOH:O	2.17	0.45
1:A:297:ARG:HD2	2:A:458:HOH:O	2.18	0.44
1:A:122:MSE:HE3	1:A:143:LEU:HG	1.98	0.44
1:A:90:LEU:HD13	1:A:327:GLU:HG2	2.00	0.44
1:A:195:ILE:HB	1:A:208:ALA:HB3	2.00	0.43
1:A:242:PHE:CZ	1:A:261:ILE:HG13	2.55	0.41
1:A:39:ASN:HB3	1:A:143:LEU:HB3	2.02	0.41
1:A:93:LEU:HG	1:A:330:HIS:CE1	2.55	0.41
1:A:320:GLN:N	2:A:647:HOH:O	2.49	0.41
1:A:309:SER:O	1:A:313:ARG:HG3	2.21	0.41

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	299/305 (98%)	293 (98%)	6 (2%)	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	257/258 (100%)	248 (96%)	9 (4%)	48	28

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	30	GLN
1	A	86	GLN
1	A	152	GLN
1	A	167	GLN
1	A	194	ASN
1	A	297	ARG
1	A	317	GLU
1	A	320	GLN
1	A	327	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	86	GLN
1	A	136	GLN
1	A	152	GLN
1	A	167	GLN
1	A	194	ASN
1	A	211	GLN
1	A	252	GLN
1	A	254	GLN
1	A	267	GLN
1	A	287	GLN
1	A	329	HIS

5.3.3 RNA ⓘ

There are no RNA chains 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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

There are no ligands in this entry.

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 ⓘ

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	301/305 (98%)	-0.11	7 (2%) 57 51	6, 12, 24, 38	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	319	ALA	4.7
1	A	316	GLY	3.3
1	A	30	GLN	2.9
1	A	236	SER	2.4
1	A	318	GLN	2.3
1	A	330	HIS	2.3
1	A	320	GLN	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 ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

There are no ligands in this entry.

6.5 Other polymers ⓘ

There are no such residues in this entry.