



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 19, 2017 – 05:23 PM EDT

PDB ID : 5VC2
Title : Crystal structure of a serine hydroxymethyltransferase from *Helicobacter pylori*
Authors : Seattle Structural Genomics Center for Infectious Disease (SSGCID)
Deposited on : 2017-03-30
Resolution : 1.90 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

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20029077
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	rb-20029077

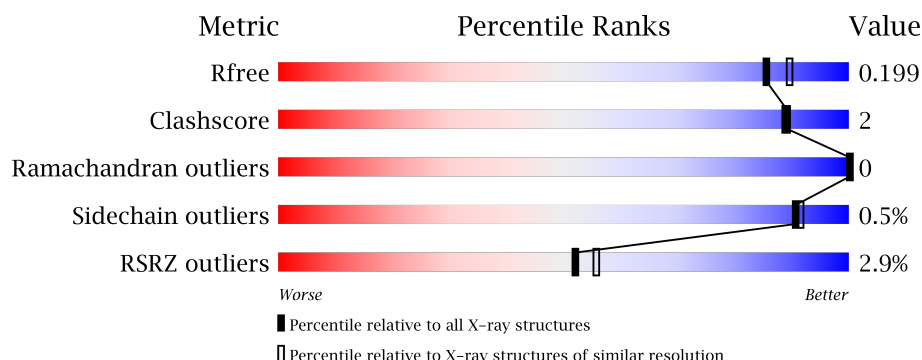
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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}	100719	5047 (1.90-1.90)
Clashscore	112137	5731 (1.90-1.90)
Ramachandran outliers	110173	5669 (1.90-1.90)
Sidechain outliers	110143	5670 (1.90-1.90)
RSRZ outliers	101464	5100 (1.90-1.90)

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. 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	424	<div> <div>3%</div> <div> <div></div> <div>87%</div> <div>9%</div> </div> </div>
1	B	424	<div> <div>3%</div> <div> <div></div> <div>83%</div> <div>7%</div> <div>10%</div> </div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6479 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 Serine hydroxymethyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	386	Total	C	N	O	S	0	4	0
			2973	1897	503	559	14			
1	B	380	Total	C	N	O	S	0	5	0
			2911	1863	493	542	13			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-7	MET	-	initiating methionine	UNP B5Z9V7
A	-6	ALA	-	expression tag	UNP B5Z9V7
A	-5	HIS	-	expression tag	UNP B5Z9V7
A	-4	HIS	-	expression tag	UNP B5Z9V7
A	-3	HIS	-	expression tag	UNP B5Z9V7
A	-2	HIS	-	expression tag	UNP B5Z9V7
A	-1	HIS	-	expression tag	UNP B5Z9V7
A	0	HIS	-	expression tag	UNP B5Z9V7
B	-7	MET	-	initiating methionine	UNP B5Z9V7
B	-6	ALA	-	expression tag	UNP B5Z9V7
B	-5	HIS	-	expression tag	UNP B5Z9V7
B	-4	HIS	-	expression tag	UNP B5Z9V7
B	-3	HIS	-	expression tag	UNP B5Z9V7
B	-2	HIS	-	expression tag	UNP B5Z9V7
B	-1	HIS	-	expression tag	UNP B5Z9V7
B	0	HIS	-	expression tag	UNP B5Z9V7

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	331	Total	O	0	1
			332	332		
2	B	262	Total	O	0	1
			263	263		

- Molecule 1: Serine hydroxymethyltransferase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	57.97Å 91.02Å 161.32Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	41.81 – 1.90 41.81 – 1.90	Depositor EDS
% Data completeness (in resolution range)	99.0 (41.81-1.90) 99.0 (41.81-1.90)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.36 (at 1.89Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.165 , 0.199 0.164 , 0.199	Depositor DCC
R_{free} test set	1954 reflections (2.90%)	DCC
Wilson B-factor (Å ²)	29.4	Xtriage
Anisotropy	0.405	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 55.5	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	6479	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.21% 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

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	0.37	0/3040	0.52	0/4103
1	B	0.35	0/2982	0.50	0/4031
All	All	0.36	0/6022	0.51	0/8134

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

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	2973	0	2945	11	0
1	B	2911	0	2868	17	0
2	A	332	0	0	4	0
2	B	263	0	0	1	0
All	All	6479	0	5813	26	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 2.

All (26) 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)
2:A:689:HOH:O	1:B:44:MET:HG3	2.00	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:232:ARG:HE	1:B:260:LEU:HD23	1.69	0.58
1:B:403:ALA:O	1:B:407[B]:GLN:HG2	2.04	0.57
1:B:32:SER:HB2	1:B:226:LYS:HE3	1.88	0.56
1:A:284:LYS:NZ	2:A:504:HOH:O	2.38	0.55
1:A:32:SER:HB2	1:A:226:LYS:HE3	1.87	0.55
1:B:236:ILE:HD12	1:B:251:ILE:HD11	1.88	0.54
1:A:112[B]:LYS:NZ	2:A:503:HOH:O	2.38	0.53
1:B:30:ILE:HG13	1:B:338:THR:HG22	1.89	0.53
1:B:11:GLU:OE2	2:B:501:HOH:O	2.18	0.51
1:B:26[A]:HIS:HD2	1:B:338:THR:OG1	1.93	0.51
1:A:232:ARG:NH2	1:B:52:TYR:O	2.41	0.46
1:B:305:VAL:HB	1:B:315:LEU:HB2	1.99	0.45
1:A:115:GLY:O	1:A:141:SER:HA	2.16	0.45
1:B:115:GLY:O	1:B:141:SER:HA	2.17	0.44
1:A:135:LYS:HE3	1:A:137:TYR:OH	2.18	0.43
1:B:2:ALA:N	1:B:415:ILE:HD12	2.33	0.43
1:B:38:PRO:HB3	1:B:416:PHE:CE2	2.54	0.43
1:B:290[B]:MET:HG2	1:B:361:SER:OG	2.17	0.43
1:A:51:LYS:HE2	1:A:52:TYR:CZ	2.54	0.42
1:B:26[A]:HIS:CD2	1:B:338:THR:OG1	2.72	0.42
1:A:91:GLN:HB2	1:A:251:ILE:HG21	2.01	0.42
1:A:118:LEU:HD11	2:A:642:HOH:O	2.20	0.41
1:B:172:SER:HB3	1:B:200:HIS:CE1	2.56	0.41
1:B:40:VAL:HG13	1:B:266:ALA:HB1	2.02	0.40
1:A:146:LEU:HD23	1:A:146:LEU:HA	1.84	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	382/424 (90%)	371 (97%)	11 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	377/424 (89%)	365 (97%)	12 (3%)	0	100	100
All	All	759/848 (90%)	736 (97%)	23 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	314/350 (90%)	313 (100%)	1 (0%)	94	94
1	B	302/350 (86%)	300 (99%)	2 (1%)	87	87
All	All	616/700 (88%)	613 (100%)	3 (0%)	91	91

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	342	ASN
1	B	162	VAL
1	B	342	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA [i](#)

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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	386/424 (91%)	-0.05	11 (2%) 53 57	19, 31, 59, 100	0
1	B	380/424 (89%)	-0.04	11 (2%) 52 56	20, 36, 62, 93	0
All	All	766/848 (90%)	-0.05	22 (2%) 52 56	19, 34, 61, 100	0

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	3	TYR	5.6
1	B	258	GLY	5.3
1	A	131	SER	4.7
1	A	130	VAL	4.1
1	B	146	LEU	3.8
1	A	118	LEU	3.3
1	B	323	TYR	3.2
1	A	67	VAL	3.1
1	A	120	CYS	2.9
1	A	119	SER	2.6
1	A	3	TYR	2.6
1	B	251	ILE	2.6
1	A	146	LEU	2.5
1	A	258	GLY	2.5
1	B	322	PRO	2.4
1	B	348	THR	2.4
1	B	2	ALA	2.3
1	B	325	GLY	2.2
1	A	352	PHE	2.2
1	A	133	THR	2.2
1	B	320[A]	ASP	2.0
1	B	416	PHE	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 carbohydrates in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.