

Full wwPDB X-ray Structure Validation Report (i)

May 16, 2020 – 05:23 pm BST

PDB ID	:	3C2Q
Title	:	Crystal structure of conserved putative LOR/SDH protein from Methanococ-
		cus maripaludis S2
Authors	:	Duke, N.; Gu, M.; Mulligan, R.; Conrad, B.; Joachimiak, A.; Midwest Center
		for Structural Genomics (MCSG)
Deposited on		
Resolution	:	2.00 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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

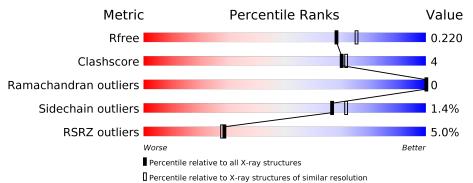
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	$7.0.044 (\mathrm{Gargrove})$
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	$egin{array}{c} { m Whole \ archive} \ (\#{ m Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries},{ m resolution\ range}({ m \AA}))$
R_{free}	130704	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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 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 <=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	А	345	90%	5%	5%
1	В	345	90%	6%	•••

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:



Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	IMD	В	502	-	-	Х	-



3C2Q

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 5386 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 Uncharacterized conserved protein.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace		
1	А	328	Total 2487	C 1578	N 425	O 470			0	0	0
1	В	336	Total 2558	C 1623	N 438	O 482	S 2	Se 13	0	0	0

There are 6 discrepancies between the modelled and reference sequences:

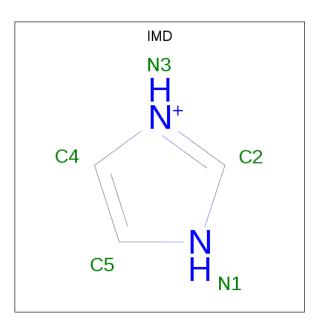
Chain	Residue	Modelled	Actual	Comment	Reference
A	71	SER		EXPRESSION TAG	v i
А	72	ASN		EXPRESSION TAG	•
A	73	ALA	-	EXPRESSION TAG	UNP Q6LXX7
В	71	SER		EXPRESSION TAG	•
В	72	ASN	-	EXPRESSION TAG	UNP Q6LXX7
В	73	ALA	-	EXPRESSION TAG	UNP Q6LXX7

• Molecule 2 is NICKEL (II) ION (three-letter code: NI) (formula: Ni).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	А	1	Total Ni 1 1	0	0

• Molecule 3 is IMIDAZOLE (three-letter code: IMD) (formula: $C_3H_5N_2$).





Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	А	1	$\begin{array}{ccc} \text{Total} \text{C} \text{N} \\ 5 3 2 \end{array}$	0	0
3	А	1	$\begin{array}{ccc} \text{Total} \text{C} \text{N} \\ 5 3 2 \end{array}$	0	0
3	В	1	$\begin{array}{ccc} \text{Total} \text{C} \text{N} \\ 5 3 2 \end{array}$	0	0
3	В	1	$\begin{array}{ccc} \text{Total} \text{C} \text{N} \\ 5 3 2 \end{array}$	0	0

• Molecule 4 is water.

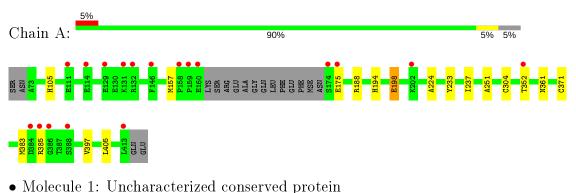
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	А	147	Total O 147 147	0	0
4	В	173	Total O 173 173	0	0

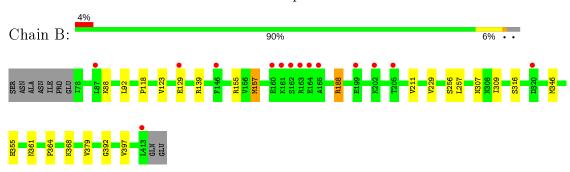


3 Residue-property plots (i)

These plots are drawn for all protein, RNA and DNA 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: Uncharacterized conserved protein







4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants	119.59Å 119.59Å 161.74Å	Depositor
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	96.23 - 2.00	Depositor
Resolution (A)	33.27 - 2.00	EDS
% Data completeness	99.8 (96.23-2.00)	Depositor
(in resolution range)	99.9(33.27-2.00)	EDS
R _{merge}	0.06	Depositor
R _{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	$7.36 (at 2.00 \text{\AA})$	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
D D	0.191 , 0.217	Depositor
R, R_{free}	0.193 , 0.220	DCC
R_{free} test set	3994 reflections $(5.02%)$	wwPDB-VP
Wilson B-factor $(Å^2)$	26.3	Xtriage
Anisotropy	0.016	Xtriage
Bulk solvent $k_{sol}(e/Å^3), B_{sol}(Å^2)$	0.40 , 52.0	EDS
L-test for twinning ²	$ \langle L \rangle = 0.51, \langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5386	wwPDB-VP
Average B, all atoms $(Å^2)$	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 3.31% of the height of the origin peak. No significant pseudotranslation is detected.

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.



¹Intensities estimated from amplitudes.

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: NI, IMD

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	Bo	nd lengths	Bond angles		
	Unam	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	А	0.71	1/2514~(0.0%)	0.69	0/3385	
1	В	0.72	0/2586	0.69	0/3477	
All	All	0.72	1/5100~(0.0%)	0.69	0/6862	

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	А	0	1

All (1) bond length outliers are listed below:

Mol	Chain	\mathbf{Res}	Type	Atoms	Z	Observed(A)	Ideal(Å)
1	А	198	GLU	CD-OE1	5.20	1.31	1.25

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	А	251	ALA	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



Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	А	2487	0	2584	18	0
1	В	2558	0	2653	19	0
2	А	1	0	0	0	0
3	А	10	0	10	3	0
3	В	10	0	10	4	0
4	А	147	0	0	2	0
4	В	173	0	0	1	0
All	All	5386	0	5257	37	0

the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

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 4.

All (37) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic	Clash
Atom-1	Atom-2	${ m distance}~({ m \AA})$	overlap (Å)
1:B:397:VAL:H	3:B:502:IMD:H2	1.05	1.18
1:B:188:ARG:HG2	1:B:188:ARG:HH11	1.11	1.09
1:B:397:VAL:N	3:B:502:IMD:H2	1.88	0.87
1:A:397:VAL:H	3:A:501:IMD:H2	1.39	0.87
1:B:188:ARG:HG2	1:B:188:ARG:NH1	1.91	0.78
1:A:361:ASN:ND2	1:A:385:ARG:HD3	2.07	0.69
1:A:371:CYS:SG	1:A:383:MSE:CE	2.81	0.68
1:A:371:CYS:SG	1:A:383:MSE:HE3	2.38	0.64
1:B:211:VAL:CG1	1:B:355:HIS:HB3	2.30	0.62
1:A:383:MSE:HE2	1:A:383:MSE:HA	1.82	0.60
1:A:397:VAL:H	3:A:501:IMD:C2	2.12	0.60
1:A:194:HIS:O	1:A:198:GLU:HG2	2.01	0.60
1:A:105:HIS:ND1	1:A:157:MSE:HE3	2.16	0.60
1:B:92:LEU:HD13	1:B:123:VAL:HG11	1.84	0.59
1:B:307:ASN:HB2	1:B:309:ILE:HD12	1.89	0.54
1:A:385:ARG:CZ	1:A:385:ARG:HB3	2.38	0.54
1:B:361:ASN:OD1	4:B:668:HOH:O	2.19	0.54
1:B:188:ARG:CG	1:B:188:ARG:HH11	2.00	0.52
1:A:188:ARG:HG3	4:A:586:HOH:O	2.11	0.51
1:B:307:ASN:HB2	1:B:309:ILE:CD1	2.41	0.50
1:B:397:VAL:H	3:B:502:IMD:C2	1.98	0.49
1:B:346:MSE:HE3	1:B:368:LYS:HG2	1.95	0.48
1:A:371:CYS:SG	1:A:383:MSE:HE1	2.54	0.46
1:A:237:ILE:HD12	1:A:304:CYS:SG	2.55	0.46
1:A:397:VAL:N	3:A:501:IMD:H2	2.20	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:155:ARG:HG2	1:B:157:MSE:SE	2.68	0.44
1:B:307:ASN:CB	1:B:309:ILE:HD12	2.48	0.43
1:B:118:PRO:HG2	1:B:364:PRO:HB3	1.99	0.43
1:A:157:MSE:HE2	1:A:157:MSE:HB2	1.84	0.42
1:A:224:ALA:HA	1:A:405:LEU:HD22	2.01	0.42
1:A:361:ASN:HD22	1:A:385:ARG:HD3	1.80	0.42
1:A:105:HIS:ND1	1:A:157:MSE:CE	2.82	0.42
1:B:256:SER:O	1:B:257:LEU:HB2	2.21	0.41
1:B:397:VAL:HG12	3:B:502:IMD:C2	2.51	0.41
1:A:233:TYR:OH	4:A:613:HOH:O	2.19	0.40
1:B:229:VAL:HG13	1:B:309:ILE:HD13	2.03	0.40
1:B:379:VAL:HG11	1:B:392:GLY:HA3	2.03	0.40

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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	Perce	\mathbf{ntiles}
1	А	324/345~(94%)	318~(98%)	6(2%)	0	100	100
1	В	334/345~(97%)	328~(98%)	6 (2%)	0	100	100
All	All	658/690 (95%)	646 (98%)	12 (2%)	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



Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	А	275/277~(99%)	273~(99%)	2(1%)	84 88
1	В	282/277~(102%)	276~(98%)	6 (2%)	53 57
All	All	557/554~(100%)	549~(99%)	8 (1%)	67 72

analysed, and the total number of residues.

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	А	175	GLU
1	А	352	THR
1	В	88	LYS
1	В	129	GLU
1	В	139	ARG
1	В	157	MSE
1	В	188	ARG
1	В	316	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	А	361	ASN
1	А	375	GLN
1	В	109	ASN
1	В	152	ASN

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)

Of 5 ligands modelled in this entry, 1 is 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	B	ond leng	gths	B	Bond ang	gles
	туре	Cham	nes		Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	IMD	А	502	-	3,5,5	0.45	0	4,5,5	0.58	0
3	IMD	В	502	-	3, 5, 5	0.13	0	4,5,5	0.78	0
3	IMD	В	503	-	3, 5, 5	0.32	0	4,5,5	0.57	0
3	IMD	А	501	-	3, 5, 5	0.72	0	4,5,5	0.41	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
3	IMD	А	502	-	-	-	0/1/1/1
3	IMD	В	502	-	-	-	0/1/1/1
3	IMD	В	503	-	-	-	0/1/1/1
3	IMD	А	501	-	-	-	0/1/1/1

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.

2 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	В	502	IMD	4	0
3	А	501	IMD	3	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	А	316/345~(91%)	0.04	18 (5%) 23 23	14, 25, 44, 67	0
1	В	323/345~(93%)	0.02	14 (4%) 35 34	14, 24, 43, 54	0
All	All	639/690~(92%)	0.03	32 (5%) 28 28	14, 24, 44, 67	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ	
1	А	160	GLU	8.8	
1	В	161	LYS	5.2	
1	А	159	PRO	4.8	
1	В	165	ALA	4.8	
1	В	162	SER	4.7	
1	В	413	LEU	4.5	
1	А	413	LEU	4.2	
1	А	174	SER	4.2	
1	А	129	GLU	3.6	
1	А	175	GLU	3.4	
1	А	386	GLY	3.4	
1	А	388	SER	3.2	
1	В	160	GLU	3.2	
1	1 A		ARG	3.0	
1	1 B		ARG	2.9	
1	А	158	PRO	2.7	
1	В	129	GLU	2.6	
1	А	131	LYS	2.6	
1	А	202	LYS	2.5	
1	В	87	LEU	2.5	
1	А	352	THR	2.5	
1	А	384	ASP	2.4	
1	А	111	GLU	2.4	
1	1 B		GLU	2.4	

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Mol	Chain	Res	Type	RSRZ
1	А	132	ARG	2.3
1	В	146	PHE	2.2
1	В	199	GLU	2.1
1	В	205	THR	2.1
1	В	202	LYS	2.1
1	А	114	GLU	2.1
1	А	146	PHE	2.0
1	В	320	ASP	2.0

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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)

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, 95^{th} 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	$\mathbf{B} ext{-factors}(\mathbf{A}^2)$	$Q{<}0.9$
3	IMD	А	501	5/5	0.55	0.40	$25,\!33,\!34,\!37$	0
3	IMD	В	502	5/5	0.95	0.11	$20,\!21,\!23,\!26$	0
2	NI	А	500	1/1	0.96	0.20	$75,\!75,\!75,\!75$	0
3	IMD	В	503	5/5	0.96	0.09	$27,\!28,\!28,\!31$	0
3	IMD	А	502	5/5	0.96	0.15	$29,\!29,\!31,\!31$	0

6.5 Other polymers (i)

There are no such residues in this entry.

