

Full wwPDB X-ray Structure Validation Report (i)

May 26, 2020 – 02:33 pm BST

PDB ID	:	1UCR
Title	:	Three-dimensional crystal structure of dissimilatory sulfite reductase D $(DsrD)$
Authors	:	Mizuno, N.; Voordouw, G.; Miki, K.; Sarai, A.; Higuchi, Y.
Deposited on		
Resolution	:	1.20 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:

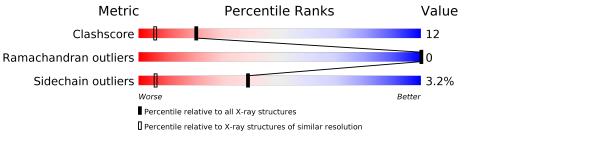
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (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 1.20 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	$egin{array}{c} { m Whole \ archive} \ (\#{ m Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries},{ m resolution\ range}({ m \AA}))$
Clashscore	141614	1286 (1.22-1.18)
Ramachandran outliers	138981	1240 (1.22-1.18)
Sidechain outliers	138945	1239 (1.22-1.18)

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%

Note EDS was not executed.

Mol	Chain	Length	Quality of chain		
1	А	78	79%	13%	•• 5%
1	В	78	76%	18%	••

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
2	SO4	А	205	-	-	Х	-
2	SO4	В	204	-	-	Х	-



 $\mathbf{2}$

Entry composition (i)

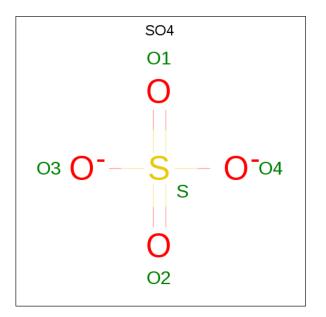
There are 3 unique types of molecules in this entry. The entry contains 1450 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Protein dsvD.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	1 A	74	Total	С	Ν	Ο	\mathbf{S}	0	1	0
			586	375	94	113	4	0		
1	В	B 75	Total	С	Ν	0	S	0	Б	0
	D	10	608	386	97	121	4	0	9	0

• Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	А	1	$\begin{array}{ccc} \text{Total} & \text{O} & \text{S} \\ 5 & 4 & 1 \end{array}$	0	0
2	А	1	$\begin{array}{ccc} \text{Total} & \text{O} & \text{S} \\ 5 & 4 & 1 \end{array}$	0	0
2	В	1	$\begin{array}{ccc} \text{Total} & \text{O} & \text{S} \\ 5 & 4 & 1 \end{array}$	0	0
2	В	1	$\begin{array}{ccc} \text{Total} & \text{O} & \text{S} \\ 5 & 4 & 1 \end{array}$	0	0

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	В	1	Total 5	0 4	S 1	0	0

• Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	А	143	Total O 143 143	0	0
3	В	88	Total O 88 88	0	0

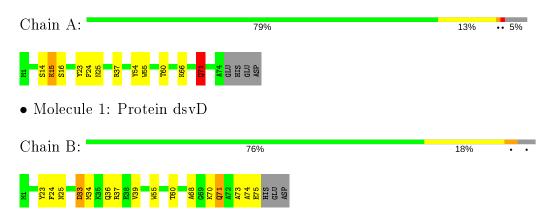


3 Residue-property plots (i)

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. Residues are colorcoded 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. 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.

Note EDS was not executed.

• Molecule 1: Protein dsvD





4 Data and refinement statistics (i)

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants	60.02Å 64.56 Å 45.22 Å	Depositor
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 - 1.20	Depositor
% Data completeness	89.7 (10.00-1.20)	Depositor
(in resolution range)	03.7 (10.00-1.20)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	SHELXL-97	Depositor
R, R_{free}	0.136 , 0.180	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	1450	wwPDB-VP
Average B, all atoms $(Å^2)$	21.0	wwPDB-VP



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: SO4

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		
MOI	Chain	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	А	0.71	0/601	1.40	7/803~(0.9%)	
1	В	0.75	0/641	1.38	7/855~(0.8%)	
All	All	0.73	0/1242	1.39	14/1658~(0.8%)	

There are no bond length outliers.

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
1	В	37	ARG	NE-CZ-NH1	-11.29	114.66	120.30
1	А	16	SER	CA-C-N	9.65	135.49	116.20
1	А	37	ARG	NE-CZ-NH2	9.36	124.98	120.30
1	А	16	SER	C-N-CA	-9.18	103.02	122.30
1	В	37	ARG	NH1-CZ-NH2	8.29	128.52	119.40
1	А	16	SER	O-C-N	-8.09	109.44	123.20
1	В	37	ARG	NE-CZ-NH2	-6.95	116.82	120.30
1	А	71	GLN	CA-CB-CG	6.55	127.80	113.40
1	А	71	GLN	CG-CD-NE2	6.30	131.83	116.70
1	В	33	ASP	CB-CG-OD2	6.16	123.84	118.30
1	В	36[A]	GLN	N-CA-CB	-6.13	99.56	110.60
1	В	36[B]	GLN	N-CA-CB	-6.13	99.56	110.60
1	В	55	TRP	CB-CG-CD1	5.75	134.47	127.00
1	А	71	GLN	CB-CG-CD	5.44	125.74	111.60

All (14) bond angle outliers are listed below:

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	А	586	0	585	15	0
1	В	608	0	606	21	0
2	А	10	0	0	6	0
2	В	15	0	0	8	0
3	А	143	0	0	3	0
3	В	88	0	0	2	0
All	All	1450	0	1191	30	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 12.

All (30) 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:71:GLN:HE21	1:B:71:GLN:HG2	1.47	0.79
1:A:71:GLN:HG3	3:A:464:HOH:O	1.91	0.70
1:A:66:LYS:HB3	3:B:454:HOH:O	1.92	0.69
1:B:25:ASN:HB2	2:B:204:SO4:O2	1.94	0.68
1:B:24:PHE:HB3	2:B:204:SO4:O1	1.93	0.68
1:A:71:GLN:NE2	1:B:71:GLN:HG2	2.09	0.67
1:B:23:TYR:HB3	2:B:204:SO4:O3	1.95	0.65
1:B:23:TYR:HB3	2:B:204:SO4:O2	1.98	0.64
1:A:14:SER:HA	1:B:70:LYS:HD2	1.83	0.61
1:A:25:ASN:HB2	2:A:205:SO4:O3	2.02	0.59
1:A:23:TYR:HB3	2:A:205:SO4:O4	2.03	0.58
1:B:23:TYR:HB3	2:B:204:SO4:S	2.47	0.55
1:B:71:GLN:O	1:B:71:GLN:HG3	2.00	0.55
1:B:34[B]:MET:CE	1:B:39:VAL:HG22	2.39	0.53
1:A:55:TRP:HE3	3:A:398:HOH:O	1.91	0.52
1:B:60:THR:OG1	2:B:204:SO4:O4	2.27	0.51
1:A:15:LYS:HE3	1:B:73:ALA:HB1	1.94	0.49
1:B:34[B]:MET:HE3	1:B:39:VAL:HG22	1.94	0.48
1:A:25:ASN:N	2:A:205:SO4:O3	2.43	0.48
1:B:74:ALA:O	1:B:75:GLU:HB2	2.13	0.47
1:B:25:ASN:N	2:B:204:SO4:O1	2.48	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)	
1:A:15:LYS:HG2	1:B:74:ALA:HB2	1.95	0.46	
1:A:24:PHE:N	2:A:205:SO4:O4	2.49	0.45	
1:B:24:PHE:N	2:B:204:SO4:O3	2.52	0.43	
2:A:205:SO4:O3	3:A:322:HOH:O	2.21	0.42	
1:A:71:GLN:OE1	1:B:68:ALA:HB1	2.19	0.42	
1:A:60:THR:HG21	2:A:205:SO4:O2	2.19	0.42	
1:A:15:LYS:HE3	1:B:73:ALA:CB	2.49	0.42	
1:B:34[B]:MET:HE3	1:B:34[B]:MET:HB3	1.98	0.41	
1:B:33:ASP:HB2	3:B:311:HOH:O	2.21	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	ntiles
1	А	73/78~(94%)	$71 \ (97\%)$	2(3%)	0	100	100
1	В	78/78~(100%)	77~(99%)	1 (1%)	0	100	100
All	All	151/156~(97%)	148 (98%)	3~(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 analysed, and the total number of residues.



Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	А	63/67~(94%)	60~(95%)	3~(5%)	25 2
1	В	69/67~(103%)	68~(99%)	1 (1%)	67 32
All	All	132/134~(98%)	128~(97%)	4(3%)	39 7

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	А	15	LYS
1	А	54	TYR
1	А	71	GLN
1	В	71	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	А	13	ASN
1	А	48	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)

5 ligands are modelled in this entry.

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



Mol	Type	Chain	Res	Link	B	ond leng	gths	E	ond ang	gles
	туре	Chain	nes		Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z >2
2	SO4	В	204	-	4,4,4	0.15	0	6,6,6	0.97	0
2	SO4	В	201	-	4,4,4	0.42	0	$6,\!6,\!6$	0.59	0
2	SO4	В	203	-	4,4,4	0.34	0	6,6,6	0.84	0
2	SO4	А	205	-	4,4,4	0.50	0	$6,\!6,\!6$	2.57	2 (33%)
2	SO4	А	202	-	4,4,4	0.33	0	6,6,6	0.27	0

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).

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
2	А	205	SO4	O4-S-O3	-4.74	88.82	109.06
2	А	205	SO4	O3-S-O1	3.82	129.25	109.31

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 14 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	В	204	SO4	8	0
2	А	205	SO4	6	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)

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates (i)

EDS was not executed - this section is therefore empty.

6.4 Ligands (i)

EDS was not executed - this section is therefore empty.

6.5 Other polymers (i)

EDS was not executed - this section is therefore empty.

