



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 05:52 AM UTC

PDB ID : 2PW9 / pdb_00002pw9
Title : Crystal structure of a putative formate dehydrogenase accessory protein from *Desulfotalea psychrophila*
Authors : Bonanno, J.B.; Dickey, M.; Bain, K.T.; Logan, C.; Romero, R.; Smith, D.; Wasserman, S.; Sauder, J.M.; Burley, S.K.; Almo, S.C.; New York SGX Research Center for Structural Genomics (NYSGXRC)
Deposited on : 2007-05-10
Resolution : 2.10 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
Mogul	:	2022.3.0, CSD as543be (2022)
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

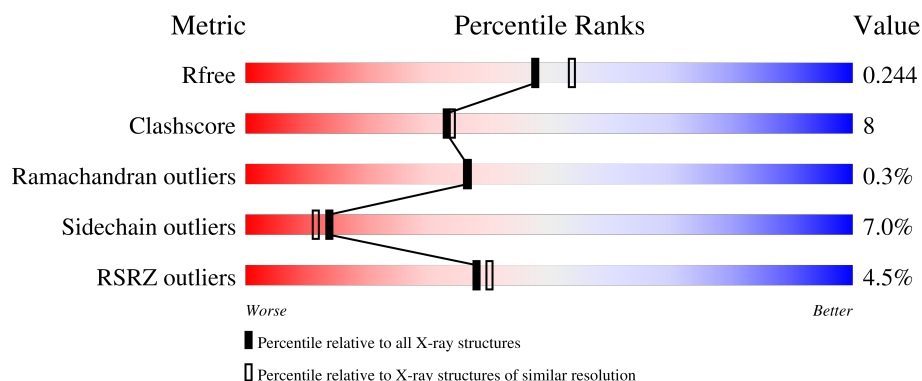
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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	268	 4% 71% 12% • 14%
1	B	268	 4% 71% 13% • 14%
1	C	268	 6% 63% 17% • 18%
1	D	268	 2% 69% 16% • 12%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7387 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 Putative formate dehydrogenase accessory protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	231	Total	C	N	O	S	0	1	0
			1780	1128	315	325	12			
1	B	230	Total	C	N	O	S	0	1	0
			1771	1123	314	322	12			
1	C	221	Total	C	N	O	S	0	0	0
			1693	1073	303	307	10			
1	D	237	Total	C	N	O	S	0	1	0
			1829	1157	322	338	12			

There are 44 discrepancies between the modelled and reference sequences:

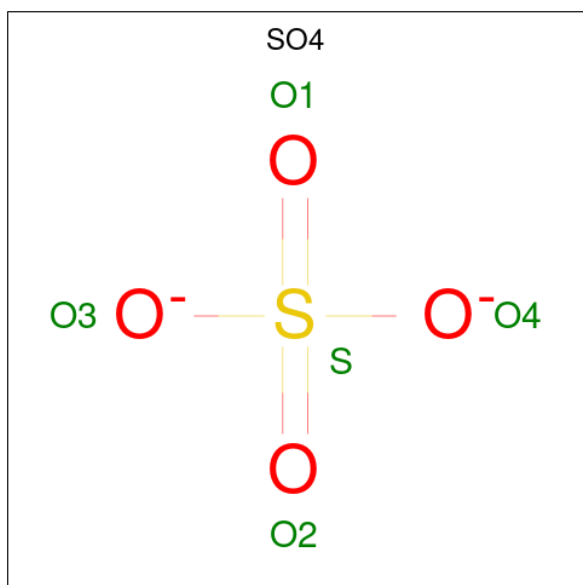
Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	MET	-	cloning artifact	UNP Q6AMB9
A	0	SER	-	cloning artifact	UNP Q6AMB9
A	1	LEU	-	cloning artifact	UNP Q6AMB9
A	259	GLU	-	cloning artifact	UNP Q6AMB9
A	260	GLY	-	cloning artifact	UNP Q6AMB9
A	261	HIS	-	cloning artifact	UNP Q6AMB9
A	262	HIS	-	cloning artifact	UNP Q6AMB9
A	263	HIS	-	cloning artifact	UNP Q6AMB9
A	264	HIS	-	cloning artifact	UNP Q6AMB9
A	265	HIS	-	cloning artifact	UNP Q6AMB9
A	266	HIS	-	cloning artifact	UNP Q6AMB9
B	-1	MET	-	cloning artifact	UNP Q6AMB9
B	0	SER	-	cloning artifact	UNP Q6AMB9
B	1	LEU	-	cloning artifact	UNP Q6AMB9
B	259	GLU	-	cloning artifact	UNP Q6AMB9
B	260	GLY	-	cloning artifact	UNP Q6AMB9
B	261	HIS	-	cloning artifact	UNP Q6AMB9
B	262	HIS	-	cloning artifact	UNP Q6AMB9
B	263	HIS	-	cloning artifact	UNP Q6AMB9
B	264	HIS	-	cloning artifact	UNP Q6AMB9
B	265	HIS	-	cloning artifact	UNP Q6AMB9

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Chain	Residue	Modelled	Actual	Comment	Reference
B	266	HIS	-	cloning artifact	UNP Q6AMB9
C	-1	MET	-	cloning artifact	UNP Q6AMB9
C	0	SER	-	cloning artifact	UNP Q6AMB9
C	1	LEU	-	cloning artifact	UNP Q6AMB9
C	259	GLU	-	cloning artifact	UNP Q6AMB9
C	260	GLY	-	cloning artifact	UNP Q6AMB9
C	261	HIS	-	cloning artifact	UNP Q6AMB9
C	262	HIS	-	cloning artifact	UNP Q6AMB9
C	263	HIS	-	cloning artifact	UNP Q6AMB9
C	264	HIS	-	cloning artifact	UNP Q6AMB9
C	265	HIS	-	cloning artifact	UNP Q6AMB9
C	266	HIS	-	cloning artifact	UNP Q6AMB9
D	-1	MET	-	cloning artifact	UNP Q6AMB9
D	0	SER	-	cloning artifact	UNP Q6AMB9
D	1	LEU	-	cloning artifact	UNP Q6AMB9
D	259	GLU	-	cloning artifact	UNP Q6AMB9
D	260	GLY	-	cloning artifact	UNP Q6AMB9
D	261	HIS	-	cloning artifact	UNP Q6AMB9
D	262	HIS	-	cloning artifact	UNP Q6AMB9
D	263	HIS	-	cloning artifact	UNP Q6AMB9
D	264	HIS	-	cloning artifact	UNP Q6AMB9
D	265	HIS	-	cloning artifact	UNP Q6AMB9
D	266	HIS	-	cloning artifact	UNP Q6AMB9

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		

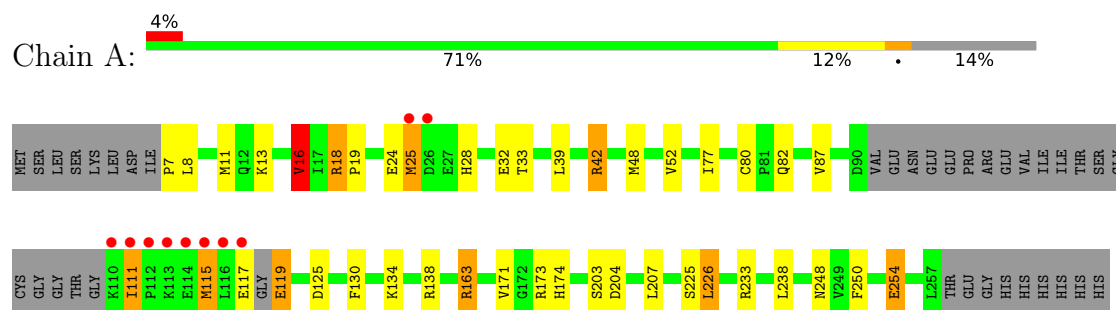
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	76	Total	O	0	0
			76	76		
3	B	88	Total	O	0	0
			88	88		
3	C	55	Total	O	0	0
			55	55		
3	D	85	Total	O	0	0
			85	85		

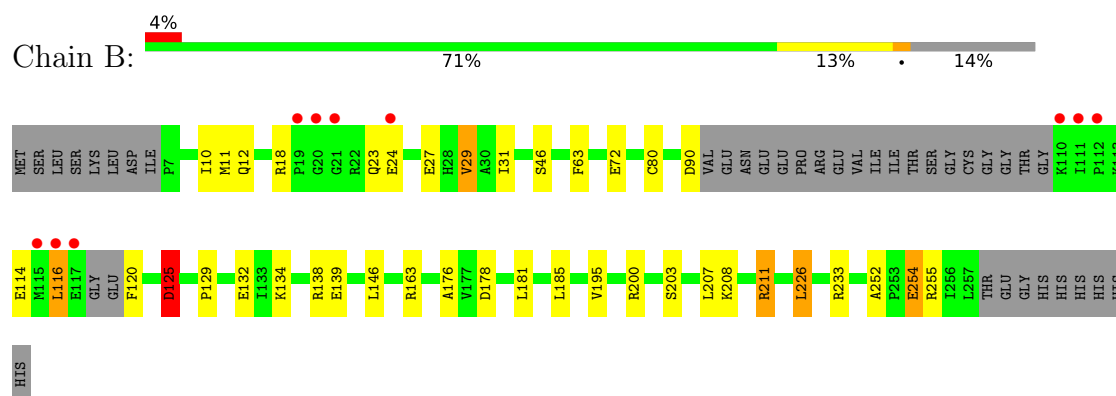
3 Residue-property plots [i](#)

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.

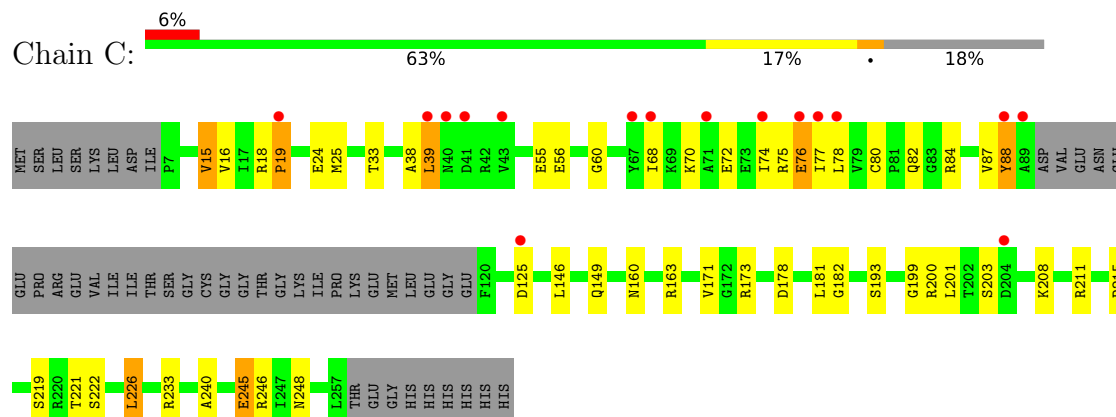
- Molecule 1: Putative formate dehydrogenase accessory protein



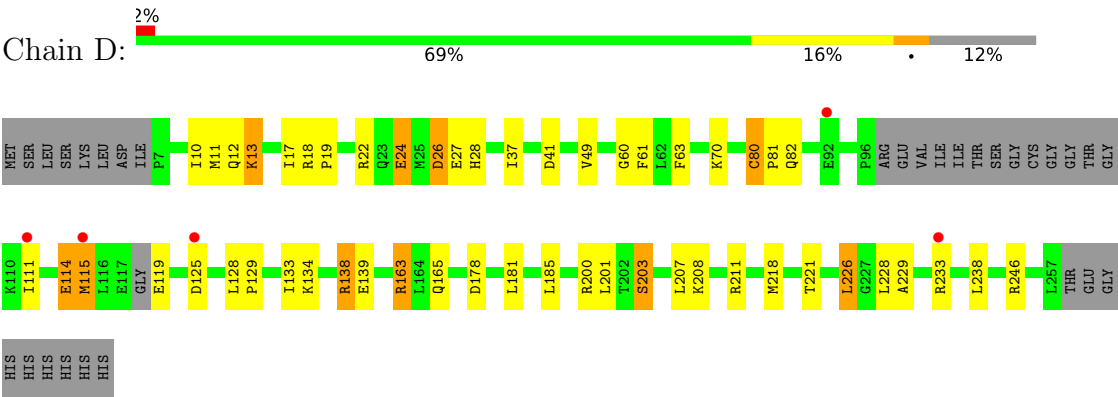
- Molecule 1: Putative formate dehydrogenase accessory protein



- Molecule 1: Putative formate dehydrogenase accessory protein



● Molecule 1: Putative formate dehydrogenase accessory protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	61.99Å 99.26Å 88.50Å 90.00° 109.70° 90.00°	Depositor
Resolution (Å)	20.00 – 2.10 20.00 – 2.10	Depositor EDS
% Data completeness (in resolution range)	98.2 (20.00-2.10) 98.1 (20.00-2.10)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	0.11	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.41 (at 2.10Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.196 , 0.259 (Not available) , 0.244	Depositor DCC
R_{free} test set	2941 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	24.6	Xtriage
Anisotropy	0.816	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 48.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.027 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	7387	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 15.66% 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

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	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.99	0/1816	1.05	1/2455 (0.0%)
1	B	1.03	3/1807 (0.2%)	1.11	3/2443 (0.1%)
1	C	0.94	0/1726	1.04	4/2336 (0.2%)
1	D	0.97	1/1866 (0.1%)	1.07	7/2524 (0.3%)
All	All	0.98	4/7215 (0.1%)	1.07	15/9758 (0.2%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	252	ALA	C-O	6.39	1.26	1.23
1	B	195	VAL	CA-CB	5.76	1.61	1.54
1	B	176	ALA	CA-CB	5.73	1.62	1.53
1	D	49	VAL	CA-CB	5.36	1.64	1.54

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	41	ASP	N-CA-C	6.89	121.44	113.38
1	B	125	ASP	N-CA-C	6.81	120.48	112.72
1	C	222	SER	N-CA-C	5.99	118.39	110.36
1	A	16	VAL	CB-CA-C	5.83	117.90	111.08
1	D	203	SER	N-CA-C	5.78	118.04	111.11
1	B	255	ARG	N-CA-C	-5.62	106.43	113.28
1	C	245	GLU	N-CA-C	5.55	118.00	110.06
1	B	139	GLU	N-CA-C	5.33	116.78	111.07
1	C	15	VAL	CB-CA-C	-5.28	103.30	110.42
1	D	61	PHE	N-CA-C	-5.08	105.63	111.07
1	C	76	GLU	N-CA-C	5.07	115.84	108.14
1	D	80[A]	CYS	CA-C-N	5.03	126.13	119.84
1	D	80[A]	CYS	C-N-CA	5.03	126.13	119.84

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	80[B]	CYS	CA-C-N	5.03	126.13	119.84
1	D	80[B]	CYS	C-N-CA	5.03	126.13	119.84

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	1780	0	1813	35	0
1	B	1771	0	1807	22	0
1	C	1693	0	1723	27	0
1	D	1829	0	1853	38	0
2	A	5	0	0	1	0
2	C	5	0	0	0	0
3	A	76	0	0	3	0
3	B	88	0	0	3	0
3	C	55	0	0	4	0
3	D	85	0	0	0	0
All	All	7387	0	7196	111	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 8.

All (111) 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:D:10:ILE:HD12	1:D:12:GLN:HE21	1.18	1.04
1:D:11:MET:HE1	1:D:28:HIS:CE1	2.01	0.96
1:A:32:GLU:HG2	1:A:48:MET:HE2	1.47	0.95
1:C:55:GLU:HG3	3:C:642:HOH:O	1.71	0.90
1:D:10:ILE:HD12	1:D:12:GLN:NE2	1.89	0.86
1:C:248:ASN:ND2	1:D:27:GLU:OE2	2.19	0.76
1:C:16:VAL:HG23	1:C:25:MET:HG3	1.68	0.75
1:A:32:GLU:HG2	1:A:48:MET:CE	2.21	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:80[B]:CYS:SG	1:D:82:GLN:NE2	2.63	0.71
1:A:225:SER:HB2	1:B:10:ILE:HD11	1.74	0.68
1:C:181:LEU:HD11	1:C:208:LYS:HB3	1.76	0.68
1:C:200:ARG:HD3	3:C:605:HOH:O	1.94	0.67
1:D:163:ARG:NH2	1:D:165:GLN:HG3	2.10	0.67
1:B:125:ASP:HA	3:B:271:HOH:O	1.96	0.66
1:D:119:GLU:O	1:D:211:ARG:NH1	2.21	0.66
1:D:181:LEU:HD11	1:D:208:LYS:HB3	1.79	0.65
1:D:134:LYS:O	1:D:138:ARG:HD3	1.97	0.65
1:D:229:ALA:O	1:D:233:ARG:HG2	1.98	0.64
3:A:657:HOH:O	1:B:200:ARG:HD3	1.98	0.63
1:C:149:GLN:NE2	3:C:639:HOH:O	2.32	0.62
1:D:10:ILE:CD1	1:D:12:GLN:HE21	2.06	0.62
1:C:171:VAL:O	1:D:200:ARG:NH2	2.27	0.61
1:B:120:PHE:HA	1:B:211:ARG:HH21	1.67	0.60
1:A:134:LYS:O	1:A:138:ARG:HD3	2.01	0.60
1:C:39:LEU:HA	1:C:87:VAL:O	2.01	0.60
1:A:171:VAL:O	1:B:200:ARG:NH2	2.26	0.59
1:C:70:LYS:HE3	1:C:72:GLU:HB2	1.85	0.59
1:A:39:LEU:O	1:A:42:ARG:HG2	2.04	0.57
3:C:645:HOH:O	1:D:200:ARG:HD3	2.03	0.57
1:D:114:GLU:HG2	1:D:115:MET:N	2.20	0.57
1:B:72:GLU:H	1:B:72:GLU:CD	2.13	0.56
1:D:18:ARG:HB3	1:D:19:PRO:HD2	1.88	0.55
1:A:80[B]:CYS:SG	1:A:82:GLN:HB2	2.46	0.55
1:D:13:LYS:HG2	1:D:24:GLU:OE2	2.07	0.55
1:D:203:SER:HB3	1:D:226:LEU:HD13	1.87	0.55
1:C:18:ARG:HH22	1:D:246:ARG:NH1	2.06	0.54
1:C:60:GLY:HA3	1:C:178:ASP:O	2.07	0.54
1:C:203:SER:HB3	1:C:226:LEU:HB3	1.88	0.54
1:D:18:ARG:HB3	1:D:19:PRO:CD	2.38	0.53
1:A:248:ASN:ND2	1:B:27:GLU:OE1	2.41	0.53
1:A:77:ILE:HG12	1:A:87:VAL:HG13	1.90	0.53
1:A:11:MET:HE1	1:A:28:HIS:NE2	2.24	0.53
1:B:178:ASP:OD1	3:B:340:HOH:O	2.19	0.51
1:B:181:LEU:HD11	1:B:208:LYS:HB3	1.93	0.51
1:D:17:ILE:HG12	1:D:22:ARG:HG2	1.93	0.51
1:A:163:ARG:HD2	3:A:662:HOH:O	2.10	0.51
1:B:72:GLU:CD	1:B:72:GLU:N	2.69	0.50
1:C:78:LEU:HD12	1:C:88:TYR:OH	2.11	0.50
1:A:48:MET:HA	1:A:48:MET:HE3	1.92	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:201:LEU:HG	1:C:219:SER:HB2	1.94	0.49
1:B:129:PRO:HB2	1:B:132:GLU:HG3	1.92	0.49
1:B:134:LYS:O	1:B:138:ARG:HD3	2.12	0.49
1:A:11:MET:HE1	1:A:28:HIS:CE1	2.49	0.48
1:A:238:LEU:C	1:A:238:LEU:HD23	2.39	0.48
1:B:116:LEU:HD12	1:B:207:LEU:HD22	1.95	0.48
1:A:111:ILE:HG21	1:A:207:LEU:HD23	1.94	0.48
1:A:174:HIS:NE2	1:A:204:ASP:HB3	2.28	0.48
1:A:7:PRO:HB3	1:D:80[B]:CYS:SG	2.55	0.47
1:A:250:PHE:CZ	1:B:27:GLU:HG3	2.49	0.47
1:B:11:MET:HA	1:B:27:GLU:O	2.15	0.47
1:C:77:ILE:HG12	1:C:87:VAL:HG22	1.96	0.47
1:A:11:MET:HE2	1:D:81:PRO:HG3	1.96	0.47
1:B:203:SER:HB3	1:B:226:LEU:HD13	1.97	0.47
1:D:115:MET:HB3	1:D:207:LEU:HD21	1.97	0.46
1:D:128:LEU:HD12	1:D:129:PRO:HD2	1.98	0.46
1:C:173:ARG:HD2	1:C:199:GLY:HA3	1.98	0.45
1:A:16:VAL:HG11	1:A:25:MET:HE2	1.99	0.45
1:C:18:ARG:HB3	1:C:19:PRO:HD2	1.98	0.45
1:A:117:GLU:O	1:A:119:GLU:N	2.49	0.45
1:C:80:CYS:HB2	1:C:84:ARG:HB3	1.99	0.45
1:A:203:SER:HB3	1:A:226:LEU:HD13	1.99	0.45
1:D:200:ARG:HG3	1:D:221:THR:CG2	2.47	0.45
1:D:128:LEU:HD21	1:D:133:ILE:HG12	1.98	0.45
1:C:38:ALA:O	1:C:87:VAL:N	2.43	0.45
1:D:11:MET:CE	1:D:28:HIS:CE1	2.89	0.45
1:B:134:LYS:NZ	3:B:309:HOH:O	2.50	0.44
1:A:32:GLU:CG	1:A:48:MET:HE2	2.34	0.44
1:A:39:LEU:O	1:A:42:ARG:CG	2.66	0.44
1:C:221:THR:O	1:C:240:ALA:HB1	2.17	0.44
1:A:8:LEU:HD21	1:A:52:VAL:HB	2.00	0.44
3:A:605:HOH:O	1:B:29:VAL:HA	2.17	0.44
1:C:80:CYS:HB3	1:C:82:GLN:OE1	2.18	0.44
1:D:63:PHE:HB2	1:D:185:LEU:HD11	1.99	0.43
1:A:130:PHE:CD2	1:A:130:PHE:C	2.96	0.43
1:B:134:LYS:O	1:B:138:ARG:CD	2.66	0.43
1:D:163:ARG:HH22	1:D:165:GLN:HG3	1.81	0.43
1:A:173:ARG:HD3	2:A:601:SO4:O4	2.18	0.43
1:D:218:MET:HE2	1:D:218:MET:HB3	1.89	0.42
1:A:18:ARG:HB2	1:A:19:PRO:CD	2.50	0.42
1:A:130:PHE:CD2	1:A:130:PHE:O	2.73	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:254:GLU:H	1:B:254:GLU:CD	2.26	0.42
1:D:37:ILE:HD12	1:D:37:ILE:N	2.34	0.42
1:A:254:GLU:H	1:A:254:GLU:HG3	1.46	0.42
1:A:8:LEU:CD2	1:A:52:VAL:HB	2.48	0.42
1:A:33:THR:CA	1:A:48:MET:HE1	2.50	0.42
1:C:246:ARG:HH11	1:C:246:ARG:HG2	1.84	0.42
1:D:60:GLY:HA3	1:D:178:ASP:O	2.19	0.42
1:B:63:PHE:HB2	1:B:185:LEU:HD11	2.02	0.41
1:C:193:SER:O	1:C:215:PRO:HD2	2.20	0.41
1:A:7:PRO:HA	1:D:80[B]:CYS:SG	2.60	0.41
1:C:68:ILE:HD13	1:C:74:ILE:HG12	2.01	0.41
1:D:13:LYS:HE3	1:D:26:ASP:OD2	2.20	0.41
1:C:56:GLU:O	1:C:182:GLY:HA3	2.20	0.41
1:A:8:LEU:H	1:D:82:GLN:NE2	2.18	0.41
1:A:115:MET:HB3	1:A:207:LEU:HD21	2.03	0.41
1:B:10:ILE:HD12	1:B:12:GLN:HE21	1.84	0.41
1:D:201:LEU:HD13	1:D:238:LEU:HD11	2.02	0.41
1:C:146:LEU:HD23	1:C:146:LEU:HA	1.90	0.41
1:C:160:ASN:OD1	1:C:163:ARG:NH2	2.52	0.41
1:D:11:MET:HE1	1:D:28:HIS:NE2	2.31	0.40
1:D:200:ARG:HG3	1:D:221:THR:HG21	2.04	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	226/268 (84%)	222 (98%)	4 (2%)	0	100	100
1	B	225/268 (84%)	220 (98%)	4 (2%)	1 (0%)	30	28
1	C	217/268 (81%)	207 (95%)	8 (4%)	2 (1%)	14	10
1	D	232/268 (87%)	228 (98%)	4 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	900/1072 (84%)	877 (97%)	20 (2%)	3 (0%)	36	36

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	75	ARG
1	B	23	GLN
1	C	19	PRO

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	194/224 (87%)	180 (93%)	14 (7%)	13	11
1	B	193/224 (86%)	176 (91%)	17 (9%)	9	7
1	C	183/224 (82%)	172 (94%)	11 (6%)	17	15
1	D	200/224 (89%)	187 (94%)	13 (6%)	15	13
All	All	770/896 (86%)	715 (93%)	55 (7%)	14	11

All (55) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	LYS
1	A	16	VAL
1	A	18	ARG
1	A	24	GLU
1	A	25	MET
1	A	42	ARG
1	A	111	ILE
1	A	115	MET
1	A	119	GLU
1	A	125	ASP
1	A	163	ARG
1	A	226	LEU

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Mol	Chain	Res	Type
1	A	233	ARG
1	A	254	GLU
1	B	18	ARG
1	B	24	GLU
1	B	29	VAL
1	B	31	ILE
1	B	46	SER
1	B	80[A]	CYS
1	B	80[B]	CYS
1	B	90	ASP
1	B	114	GLU
1	B	116	LEU
1	B	125	ASP
1	B	146	LEU
1	B	163	ARG
1	B	211	ARG
1	B	226	LEU
1	B	233	ARG
1	B	254	GLU
1	C	15	VAL
1	C	24	GLU
1	C	33	THR
1	C	39	LEU
1	C	76	GLU
1	C	88	TYR
1	C	125	ASP
1	C	211	ARG
1	C	226	LEU
1	C	233	ARG
1	C	245	GLU
1	D	13	LYS
1	D	24	GLU
1	D	26	ASP
1	D	70	LYS
1	D	111	ILE
1	D	114	GLU
1	D	115	MET
1	D	125	ASP
1	D	138	ARG
1	D	139	GLU
1	D	163	ARG
1	D	226	LEU

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Mol	Chain	Res	Type
1	D	228	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	65	GLN
1	B	174	HIS
1	C	23	GLN
1	C	174	HIS
1	D	12	GLN
1	D	82	GLN

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

2 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 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	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	SO4	A	601	-	4,4,4	0.38	0	6,6,6	0.31	0
2	SO4	C	602	-	4,4,4	0.35	0	6,6,6	0.54	0

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.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	601	SO4	1	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

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	231/268 (86%)	0.06	10 (4%) 40 41	18, 30, 50, 109	1 (0%)
1	B	230/268 (85%)	0.08	10 (4%) 40 41	19, 29, 67, 108	1 (0%)
1	C	221/268 (82%)	0.31	16 (7%) 21 23	22, 34, 65, 77	0
1	D	237/268 (88%)	-0.02	5 (2%) 63 66	18, 30, 56, 73	1 (0%)
All	All	919/1072 (85%)	0.11	41 (4%) 38 40	18, 30, 64, 109	3 (0%)

All (41) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	116	LEU	6.1
1	B	111	ILE	5.4
1	A	111	ILE	5.1
1	B	116	LEU	5.0
1	B	20	GLY	4.8
1	A	115	MET	4.5
1	B	115	MET	4.3
1	C	88	TYR	4.2
1	A	110	LYS	3.8
1	C	19	PRO	3.5
1	B	110	LYS	3.4
1	A	113	LYS	3.3
1	B	112	PRO	3.2
1	C	39	LEU	3.1
1	D	125	ASP	3.0
1	C	77	ILE	3.0
1	C	78	LEU	2.9
1	A	112	PRO	2.9
1	D	92	GLU	2.8
1	C	89	ALA	2.7
1	C	74	ILE	2.7

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Mol	Chain	Res	Type	RSRZ
1	B	117	GLU	2.7
1	B	21	GLY	2.5
1	D	111	ILE	2.4
1	C	43	VAL	2.4
1	C	71	ALA	2.4
1	C	67	TYR	2.3
1	C	40	ASN	2.3
1	C	125	ASP	2.3
1	B	24	GLU	2.3
1	B	19	PRO	2.2
1	A	26	ASP	2.2
1	A	114	GLU	2.2
1	C	41	ASP	2.2
1	D	233	ARG	2.1
1	A	117	GLU	2.1
1	A	25	MET	2.1
1	C	76	GLU	2.1
1	C	68	ILE	2.1
1	D	115	MET	2.0
1	C	204	ASP	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 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(Å ²)	Q<0.9
2	SO4	C	602	5/5	0.98	0.06	31,33,34,35	0
2	SO4	A	601	5/5	0.99	0.06	28,29,32,34	0

6.5 Other polymers [i](#)

There are no such residues in this entry.