



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:20 AM GMT

PDB ID : 2GP4
Title : Structure of [FeS]cluster-free Apo Form of 6-Phosphogluconate Dehydratase from *Shewanella oneidensis*
Authors : Schormann, N.; Symersky, J.; Southeast Collaboratory for Structural Genomics (SECSG)
Deposited on : 2006-04-16
Resolution : 2.49 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
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
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) : trunk26865

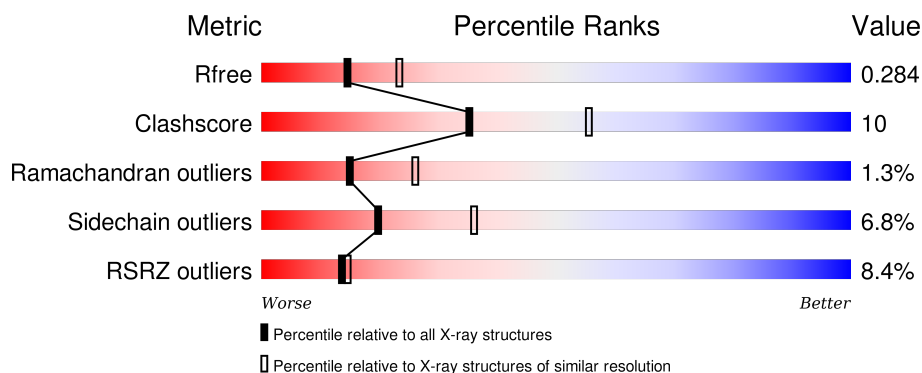
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.49 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}	91344	3553 (2.50-2.50)
Clashscore	102246	4242 (2.50-2.50)
Ramachandran outliers	100387	4156 (2.50-2.50)
Sidechain outliers	100360	4158 (2.50-2.50)
RSRZ outliers	91569	3562 (2.50-2.50)

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	628	<div> <div>7%</div> <div>64%</div> <div>15%</div> <div>•</div> <div>17%</div> </div>
1	B	628	<div> <div>7%</div> <div>65%</div> <div>12%</div> <div>•</div> <div>21%</div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8116 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 6-phosphogluconate dehydratase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	523	Total	C	N	O	S	Se	0	0	0
			3892	2445	685	743	5	14			
1	B	495	Total	C	N	O	S	Se	0	0	0
			3671	2307	649	697	4	14			

There are 74 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MSE	-	EXPRESSION TAG	GB 24348501
A	-18	GLY	-	EXPRESSION TAG	GB 24348501
A	-17	SER	-	EXPRESSION TAG	GB 24348501
A	-16	SER	-	EXPRESSION TAG	GB 24348501
A	-15	HIS	-	EXPRESSION TAG	GB 24348501
A	-14	HIS	-	EXPRESSION TAG	GB 24348501
A	-13	HIS	-	EXPRESSION TAG	GB 24348501
A	-12	HIS	-	EXPRESSION TAG	GB 24348501
A	-11	HIS	-	EXPRESSION TAG	GB 24348501
A	-10	HIS	-	EXPRESSION TAG	GB 24348501
A	-9	SER	-	EXPRESSION TAG	GB 24348501
A	-8	SER	-	EXPRESSION TAG	GB 24348501
A	-7	GLY	-	EXPRESSION TAG	GB 24348501
A	-6	LEU	-	EXPRESSION TAG	GB 24348501
A	-5	VAL	-	EXPRESSION TAG	GB 24348501
A	-4	PRO	-	EXPRESSION TAG	GB 24348501
A	-3	ARG	-	EXPRESSION TAG	GB 24348501
A	-2	GLY	-	EXPRESSION TAG	GB 24348501
A	-1	SER	-	EXPRESSION TAG	GB 24348501
A	0	HIS	-	EXPRESSION TAG	GB 24348501
A	1	MSE	MET	MODIFIED RESIDUE	GB 24348501
A	77	MSE	MET	MODIFIED RESIDUE	GB 24348501
A	111	MSE	MET	MODIFIED RESIDUE	GB 24348501
A	122	MSE	MET	MODIFIED RESIDUE	GB 24348501
A	134	MSE	MET	MODIFIED RESIDUE	GB 24348501

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Chain	Residue	Modelled	Actual	Comment	Reference
A	144	MSE	MET	MODIFIED RESIDUE	GB 24348501
A	173	MSE	MET	MODIFIED RESIDUE	GB 24348501
A	181	MSE	MET	MODIFIED RESIDUE	GB 24348501
A	233	MSE	MET	MODIFIED RESIDUE	GB 24348501
A	237	MSE	MET	MODIFIED RESIDUE	GB 24348501
A	260	MSE	MET	MODIFIED RESIDUE	GB 24348501
A	305	MSE	MET	MODIFIED RESIDUE	GB 24348501
A	353	MSE	MET	MODIFIED RESIDUE	GB 24348501
A	484	MSE	MET	MODIFIED RESIDUE	GB 24348501
A	507	MSE	MET	MODIFIED RESIDUE	GB 24348501
A	512	MSE	MET	MODIFIED RESIDUE	GB 24348501
A	579	MSE	MET	MODIFIED RESIDUE	GB 24348501
B	-19	MSE	-	EXPRESSION TAG	GB 24348501
B	-18	GLY	-	EXPRESSION TAG	GB 24348501
B	-17	SER	-	EXPRESSION TAG	GB 24348501
B	-16	SER	-	EXPRESSION TAG	GB 24348501
B	-15	HIS	-	EXPRESSION TAG	GB 24348501
B	-14	HIS	-	EXPRESSION TAG	GB 24348501
B	-13	HIS	-	EXPRESSION TAG	GB 24348501
B	-12	HIS	-	EXPRESSION TAG	GB 24348501
B	-11	HIS	-	EXPRESSION TAG	GB 24348501
B	-10	HIS	-	EXPRESSION TAG	GB 24348501
B	-9	SER	-	EXPRESSION TAG	GB 24348501
B	-8	SER	-	EXPRESSION TAG	GB 24348501
B	-7	GLY	-	EXPRESSION TAG	GB 24348501
B	-6	LEU	-	EXPRESSION TAG	GB 24348501
B	-5	VAL	-	EXPRESSION TAG	GB 24348501
B	-4	PRO	-	EXPRESSION TAG	GB 24348501
B	-3	ARG	-	EXPRESSION TAG	GB 24348501
B	-2	GLY	-	EXPRESSION TAG	GB 24348501
B	-1	SER	-	EXPRESSION TAG	GB 24348501
B	0	HIS	-	EXPRESSION TAG	GB 24348501
B	1	MSE	MET	MODIFIED RESIDUE	GB 24348501
B	77	MSE	MET	MODIFIED RESIDUE	GB 24348501
B	111	MSE	MET	MODIFIED RESIDUE	GB 24348501
B	122	MSE	MET	MODIFIED RESIDUE	GB 24348501
B	134	MSE	MET	MODIFIED RESIDUE	GB 24348501
B	144	MSE	MET	MODIFIED RESIDUE	GB 24348501
B	173	MSE	MET	MODIFIED RESIDUE	GB 24348501
B	181	MSE	MET	MODIFIED RESIDUE	GB 24348501
B	233	MSE	MET	MODIFIED RESIDUE	GB 24348501
B	237	MSE	MET	MODIFIED RESIDUE	GB 24348501

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Chain	Residue	Modelled	Actual	Comment	Reference
B	260	MSE	MET	MODIFIED RESIDUE	GB 24348501
B	305	MSE	MET	MODIFIED RESIDUE	GB 24348501
B	353	MSE	MET	MODIFIED RESIDUE	GB 24348501
B	484	MSE	MET	MODIFIED RESIDUE	GB 24348501
B	507	MSE	MET	MODIFIED RESIDUE	GB 24348501
B	512	MSE	MET	MODIFIED RESIDUE	GB 24348501
B	579	MSE	MET	MODIFIED RESIDUE	GB 24348501

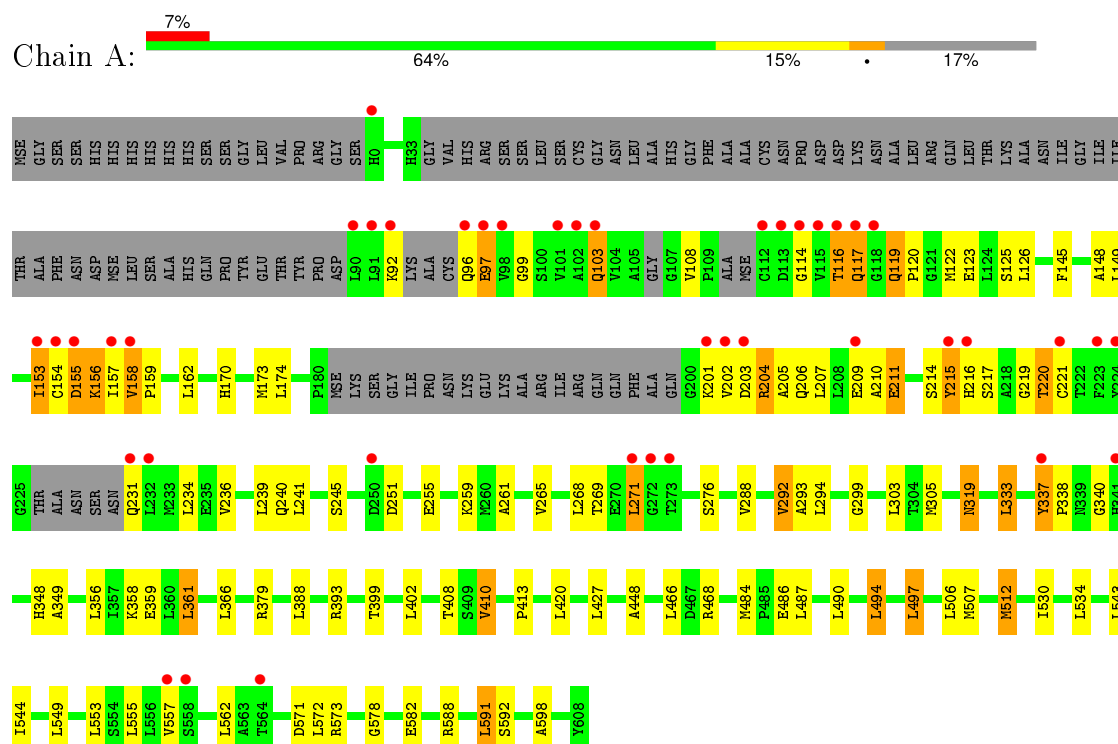
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	292	Total O 292 292	0	0
2	B	261	Total O 261 261	0	0

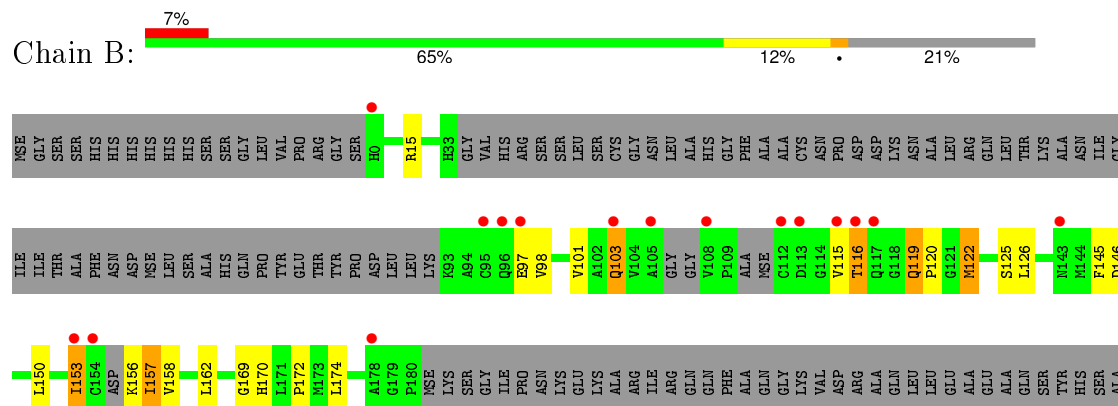
3 Residue-property plots

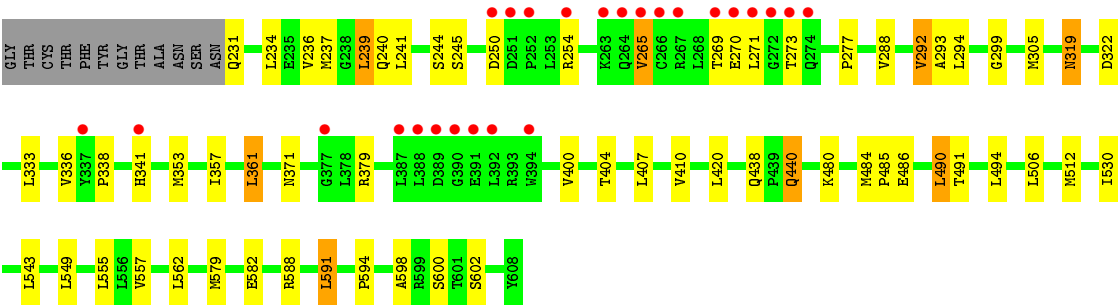
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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: 6-phosphogluconate dehydratase



- Molecule 1: 6-phosphogluconate dehydratase





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	61.36Å 118.65Å 160.02Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.68 – 2.49 47.65 – 2.49	Depositor EDS
% Data completeness (in resolution range)	100.0 (48.68-2.49) 99.5 (47.65-2.49)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.39 (at 2.48Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.230 , 0.291 0.226 , 0.284	Depositor DCC
R_{free} test set	2086 reflections (5.30%)	DCC
Wilson B-factor (Å ²)	31.8	Xtriage
Anisotropy	0.077	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 53.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 41472 reflections	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	8116	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.77% 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, 0.375 respectively for untwinned datasets, and 0.333, 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	0/3934	0.55	0/5304
1	B	0.36	0/3710	0.56	0/5003
All	All	0.37	0/7644	0.55	0/10307

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	3892	0	3949	85	0
1	B	3671	0	3735	65	0
2	A	292	0	0	7	0
2	B	261	0	0	2	0
All	All	8116	0	7684	150	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 10.

All (150) 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:B:122:MSE:HE2	1:B:486:GLU:H	1.17	1.08

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:305:MSE:HE2	1:B:598:ALA:HB2	1.34	1.05
1:A:119:GLN:HB2	1:A:120:PRO:HD3	1.40	1.04
1:B:305:MSE:CE	1:B:598:ALA:HB2	1.89	1.02
1:B:305:MSE:HE2	1:B:598:ALA:CB	1.97	0.94
1:A:122:MSE:HE2	1:A:486:GLU:H	1.31	0.93
1:A:158:VAL:HG11	1:A:234:LEU:HD21	1.48	0.93
1:B:440:GLN:H	1:B:440:GLN:HE21	1.19	0.87
1:B:158:VAL:HG11	1:B:234:LEU:HD21	1.57	0.86
1:A:507:MSE:HE1	1:A:544:ILE:HG21	1.58	0.86
1:B:122:MSE:HE2	1:B:486:GLU:N	1.94	0.81
1:A:507:MSE:CE	1:A:544:ILE:HG21	2.12	0.80
1:A:158:VAL:CG1	1:A:234:LEU:HD21	2.13	0.79
1:A:507:MSE:HE1	1:A:544:ILE:CG2	2.12	0.78
1:B:119:GLN:HB2	1:B:120:PRO:HD3	1.68	0.73
1:B:353:MSE:O	1:B:357:ILE:HG12	1.88	0.73
1:A:209:GLU:OE1	1:A:251:ASP:HB3	1.89	0.72
1:A:210:ALA:O	1:A:211:GLU:HB3	1.89	0.72
1:A:379:ARG:HD2	2:A:712:HOH:O	1.90	0.72
1:A:305:MSE:SE	1:A:484:MSE:HE1	2.40	0.72
1:A:269:THR:HG22	1:A:271:LEU:H	1.56	0.70
1:B:588:ARG:HA	1:B:591:LEU:HD22	1.73	0.70
1:A:214:SER:HB2	1:A:216:HIS:ND1	2.06	0.70
1:B:555:LEU:HG	1:B:557:VAL:HG22	1.73	0.70
1:A:96:GLN:O	1:A:97:GLU:HB2	1.91	0.69
1:A:206:GLN:O	1:A:209:GLU:HG2	1.91	0.69
1:B:122:MSE:CE	1:B:486:GLU:H	2.02	0.68
1:A:571:ASP:OD1	1:A:573:ARG:HD3	1.94	0.68
1:A:119:GLN:HB2	1:A:120:PRO:CD	2.20	0.67
1:B:122:MSE:HE3	1:B:122:MSE:O	1.95	0.67
1:A:103:GLN:HG3	1:A:145:PHE:CE1	2.30	0.66
1:B:440:GLN:H	1:B:440:GLN:NE2	1.94	0.65
1:B:353:MSE:HE3	1:B:357:ILE:HD11	1.78	0.64
1:A:156:LYS:HA	2:A:879:HOH:O	1.97	0.64
1:A:211:GLU:HB2	1:A:388:LEU:HD22	1.79	0.64
1:A:174:LEU:HD13	1:A:265:VAL:HG22	1.79	0.63
1:B:305:MSE:HE3	1:B:594:PRO:HA	1.82	0.62
1:A:255:GLU:O	1:A:259:LYS:HG2	2.01	0.61
1:A:216:HIS:O	1:A:220:THR:HG22	2.00	0.61
1:A:123:GLU:HG3	2:A:781:HOH:O	2.01	0.60
1:A:204:ARG:HG3	1:A:221:CYS:SG	2.41	0.60
1:B:357:ILE:HG22	1:B:361:LEU:HD22	1.82	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:15:ARG:NH2	1:B:322:ASP:OD2	2.34	0.60
1:A:103:GLN:HG2	1:A:108:VAL:O	2.02	0.59
1:B:174:LEU:HD13	1:B:265:VAL:HG13	1.83	0.59
1:A:96:GLN:N	2:A:658:HOH:O	2.35	0.59
1:A:149:LEU:HD21	1:A:265:VAL:HG21	1.85	0.59
1:A:333:LEU:HD21	1:A:410:VAL:HG12	1.83	0.59
1:B:440:GLN:HE21	1:B:440:GLN:N	1.98	0.58
1:A:236:VAL:HG11	1:A:293:ALA:HB2	1.83	0.58
1:A:534:LEU:HD11	1:A:562:LEU:HD13	1.86	0.58
1:A:388:LEU:HD12	1:A:393:ARG:HD3	1.86	0.58
1:A:588:ARG:HA	1:A:591:LEU:HD22	1.87	0.57
1:A:261:ALA:O	1:A:265:VAL:HG23	2.03	0.57
1:B:236:VAL:HG11	1:B:293:ALA:HB2	1.86	0.56
1:A:157:ILE:O	1:A:158:VAL:HB	2.05	0.56
1:A:361:LEU:HD13	1:A:366:LEU:HB2	1.87	0.56
1:B:116:THR:CG2	1:B:125:SER:HA	2.36	0.55
1:B:305:MSE:SE	1:B:484:MSE:HE1	2.55	0.55
1:A:211:GLU:HB2	1:A:388:LEU:CD2	2.37	0.55
1:B:319:ASN:C	1:B:319:ASN:HD22	2.10	0.55
1:A:116:THR:HG23	1:A:125:SER:HA	1.87	0.55
1:B:244:SER:O	1:B:254:ARG:NH2	2.40	0.55
1:B:490:LEU:CB	1:B:512:MSE:HE1	2.36	0.54
1:A:153:ILE:HD13	1:A:157:ILE:HD12	1.89	0.54
1:A:507:MSE:HE2	1:A:553:LEU:HD11	1.89	0.53
1:B:122:MSE:HE2	1:B:485:PRO:HA	1.89	0.53
1:A:103:GLN:HB2	1:A:145:PHE:CD1	2.43	0.53
1:B:494:LEU:HD22	1:B:506:LEU:HD22	1.91	0.52
1:B:490:LEU:HB2	1:B:512:MSE:HE1	1.91	0.52
1:A:337:TYR:O	1:A:338:PRO:C	2.46	0.52
1:A:507:MSE:HE1	1:A:544:ILE:HG23	1.93	0.51
1:B:103:GLN:HB3	1:B:145:PHE:CD2	2.46	0.51
1:A:216:HIS:CD2	1:A:217:SER:HG	2.28	0.51
1:B:116:THR:HG23	1:B:125:SER:HA	1.92	0.51
1:A:231:GLN:HA	1:A:234:LEU:HD23	1.93	0.51
1:A:204:ARG:HA	1:A:207:LEU:HD12	1.92	0.51
1:A:240:GLN:HE22	1:A:245:SER:HB3	1.76	0.50
1:B:294:LEU:HD22	1:B:299:GLY:HA3	1.93	0.50
1:B:15:ARG:HD3	2:B:847:HOH:O	2.12	0.50
1:B:120:PRO:HB2	1:B:579:MSE:HG3	1.94	0.50
1:A:97:GLU:CG	1:A:114:GLY:HA3	2.42	0.50
1:B:237:MSE:HE3	1:B:239:LEU:CD2	2.42	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:215:TYR:CG	1:A:402:LEU:HD11	2.47	0.49
1:A:337:TYR:O	1:A:340:GLY:O	2.30	0.49
1:B:404:THR:HA	1:B:407:LEU:O	2.13	0.49
1:A:319:ASN:C	1:A:319:ASN:HD22	2.15	0.49
1:B:122:MSE:CE	1:B:485:PRO:HA	2.43	0.49
1:B:491:THR:HG23	1:B:512:MSE:HE2	1.95	0.48
1:B:357:ILE:HG22	1:B:361:LEU:CD2	2.42	0.48
1:B:231:GLN:HE21	1:B:245:SER:HB2	1.78	0.48
1:A:494:LEU:HD13	1:A:506:LEU:HD21	1.96	0.48
1:A:158:VAL:HB	1:A:159:PRO:HD3	1.96	0.47
1:B:169:GLY:O	1:B:277:PRO:HB2	2.14	0.47
1:A:149:LEU:CD2	1:A:265:VAL:HG21	2.45	0.47
1:A:211:GLU:HG2	1:A:211:GLU:O	2.15	0.47
1:A:148:ALA:HB3	1:A:173:MSE:HG2	1.96	0.47
1:A:122:MSE:HE2	1:A:486:GLU:N	2.14	0.47
1:A:97:GLU:HG3	1:A:114:GLY:HA3	1.97	0.47
1:A:530:ILE:HG21	1:A:572:LEU:HD11	1.96	0.47
1:B:288:VAL:O	1:B:292:VAL:HG12	2.14	0.47
1:A:348:HIS:HD2	2:A:687:HOH:O	1.98	0.46
1:A:408:THR:HG21	1:A:413:PRO:HA	1.97	0.45
1:B:237:MSE:HE3	1:B:239:LEU:HD22	1.99	0.45
1:B:484:MSE:HE2	1:B:484:MSE:HB3	1.86	0.45
1:A:359:GLU:HB3	1:A:410:VAL:HG13	1.99	0.45
1:B:240:GLN:NE2	1:B:245:SER:HB3	2.32	0.45
1:A:170:HIS:CD2	1:A:170:HIS:H	2.35	0.45
1:A:582:GLU:CD	1:A:582:GLU:H	2.18	0.44
1:B:305:MSE:HE2	1:B:598:ALA:HB3	1.90	0.44
1:B:240:GLN:HE22	1:B:245:SER:HB3	1.83	0.44
1:B:269:THR:HG22	1:B:270:GLU:N	2.33	0.44
1:B:582:GLU:H	1:B:582:GLU:CD	2.21	0.44
1:B:170:HIS:H	1:B:170:HIS:CD2	2.34	0.44
1:A:408:THR:CG2	1:A:413:PRO:HA	2.48	0.44
1:A:358:LYS:HE3	1:A:399:THR:HG22	1.99	0.44
1:B:480:LYS:HB3	1:B:530:ILE:HD11	2.00	0.44
1:A:487:LEU:O	1:A:512:MSE:HG3	2.18	0.43
1:A:448:ALA:HB2	1:A:544:ILE:HD12	2.00	0.43
1:B:153:ILE:HG13	1:B:153:ILE:H	1.58	0.43
1:B:156:LYS:HA	2:B:739:HOH:O	2.17	0.43
1:B:591:LEU:HD13	1:B:600:SER:HB3	2.01	0.43
1:B:101:VAL:HG13	1:B:150:LEU:HD23	2.00	0.43
1:A:215:TYR:CD2	1:A:402:LEU:HD11	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:153:ILE:HD12	1:B:157:ILE:HG13	2.01	0.43
1:A:201:LYS:O	1:A:204:ARG:HD3	2.18	0.42
1:A:268:LEU:HG	1:A:276:SER:HB2	2.01	0.42
1:A:294:LEU:HD22	1:A:299:GLY:HA3	2.00	0.42
1:A:427:LEU:HD13	1:A:507:MSE:SE	2.70	0.42
1:A:288:VAL:O	1:A:292:VAL:HG13	2.20	0.42
1:A:149:LEU:HD21	1:A:265:VAL:CG2	2.47	0.42
1:A:204:ARG:HG2	1:A:205:ALA:N	2.35	0.42
1:B:103:GLN:CB	1:B:145:PHE:CD2	3.02	0.42
1:A:466:LEU:HB3	1:A:497:LEU:HD21	2.02	0.42
1:A:99:GLY:N	2:A:852:HOH:O	2.51	0.41
1:A:555:LEU:HG	1:A:557:VAL:HG22	2.02	0.41
1:B:336:VAL:HG13	1:B:338:PRO:HD2	2.02	0.41
1:A:219:GLY:HA2	1:A:349:ALA:HB1	2.03	0.41
1:B:438:GLN:HB3	1:B:440:GLN:NE2	2.36	0.41
1:A:592:SER:O	1:A:598:ALA:HA	2.20	0.41
1:A:578:GLY:N	1:A:582:GLU:OE2	2.50	0.41
1:A:117:GLN:HB2	2:A:883:HOH:O	2.20	0.41
1:B:116:THR:HG21	1:B:125:SER:HA	2.02	0.40
1:A:154:CYS:SG	1:A:155:ASP:N	2.94	0.40
1:B:146:ASP:O	1:B:172:PRO:HD2	2.22	0.40
1:B:557:VAL:HG21	1:B:562:LEU:HD21	2.04	0.40
1:B:115:VAL:HG12	1:B:157:ILE:HD11	2.03	0.40
1:A:204:ARG:HG2	1:A:205:ALA:H	1.86	0.40
1:B:371:ASN:HD22	1:B:371:ASN:HA	1.75	0.40
1:B:379:ARG:HA	1:B:379:ARG:HD3	1.90	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	509/628 (81%)	486 (96%)	16 (3%)	7 (1%)	14	24
1	B	483/628 (77%)	461 (95%)	16 (3%)	6 (1%)	16	29
All	All	992/1256 (79%)	947 (96%)	32 (3%)	13 (1%)	15	26

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	97	GLU
1	A	158	VAL
1	B	98	VAL
1	B	153	ILE
1	A	153	ILE
1	A	155	ASP
1	B	97	GLU
1	B	273	THR
1	A	119	GLN
1	A	337	TYR
1	B	119	GLN
1	B	271	LEU
1	A	202	VAL

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	416/482 (86%)	385 (92%)	31 (8%)	17	31
1	B	392/482 (81%)	368 (94%)	24 (6%)	23	42
All	All	808/964 (84%)	753 (93%)	55 (7%)	20	36

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

Mol	Chain	Res	Type
1	A	92	LYS
1	A	103	GLN
1	A	116	THR

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Mol	Chain	Res	Type
1	A	117	GLN
1	A	126	LEU
1	A	156	LYS
1	A	162	LEU
1	A	203	ASP
1	A	204	ARG
1	A	211	GLU
1	A	215	TYR
1	A	220	THR
1	A	239	LEU
1	A	241	LEU
1	A	271	LEU
1	A	292	VAL
1	A	303	LEU
1	A	319	ASN
1	A	333	LEU
1	A	356	LEU
1	A	361	LEU
1	A	410	VAL
1	A	420	LEU
1	A	468	ARG
1	A	490	LEU
1	A	494	LEU
1	A	497	LEU
1	A	512	MSE
1	A	543	LEU
1	A	549	LEU
1	A	591	LEU
1	B	103	GLN
1	B	116	THR
1	B	122	MSE
1	B	126	LEU
1	B	157	ILE
1	B	162	LEU
1	B	239	LEU
1	B	241	LEU
1	B	250	ASP
1	B	265	VAL
1	B	292	VAL
1	B	319	ASN
1	B	333	LEU
1	B	341	HIS

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Mol	Chain	Res	Type
1	B	361	LEU
1	B	400	VAL
1	B	410	VAL
1	B	420	LEU
1	B	440	GLN
1	B	490	LEU
1	B	543	LEU
1	B	549	LEU
1	B	591	LEU
1	B	602	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	28	ASN
1	A	142	HIS
1	A	170	HIS
1	A	231	GLN
1	A	240	GLN
1	A	319	ASN
1	A	438	GLN
1	B	6	GLN
1	B	96	GLN
1	B	142	HIS
1	B	170	HIS
1	B	231	GLN
1	B	240	GLN
1	B	248	ASN
1	B	319	ASN
1	B	371	ASN
1	B	440	GLN

5.3.3 RNA ⓘ

There are no RNA molecules 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 [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	509/628 (81%)	0.24	42 (8%)	14 15	16, 26, 49, 70	0
1	B	481/628 (76%)	0.23	41 (8%)	13 14	16, 25, 48, 62	0
All	All	990/1256 (78%)	0.24	83 (8%)	14 14	16, 25, 49, 70	0

All (83) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	389	ASP	6.8
1	A	215	TYR	6.7
1	B	273	THR	6.6
1	B	271	LEU	6.4
1	A	91	LEU	6.3
1	B	116	THR	5.8
1	B	391	GLU	5.1
1	B	95	CYS	4.9
1	B	388	LEU	4.8
1	B	272	GLY	4.5
1	A	155	ASP	4.4
1	A	271	LEU	4.3
1	A	103	GLN	4.3
1	A	154	CYS	4.1
1	B	265	VAL	4.0
1	A	114	GLY	4.0
1	B	112	CYS	4.0
1	A	116	THR	3.9
1	B	96	GLN	3.9
1	B	103	GLN	3.9
1	A	273	THR	3.8
1	B	117	GLN	3.8
1	A	0	HIS	3.8
1	A	153	ILE	3.4

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Mol	Chain	Res	Type	RSRZ
1	A	98	VAL	3.4
1	A	202	VAL	3.4
1	A	337	TYR	3.3
1	A	118	GLY	3.3
1	A	272	GLY	3.3
1	B	387	LEU	3.3
1	B	115	VAL	3.2
1	B	154	CYS	3.1
1	A	557	VAL	3.1
1	A	341	HIS	3.1
1	A	97	GLU	3.1
1	B	337	TYR	3.1
1	B	153	ILE	3.1
1	B	274	GLN	3.1
1	A	113	ASP	3.0
1	A	112	CYS	3.0
1	B	392	LEU	3.0
1	A	224	TYR	3.0
1	A	102	ALA	2.9
1	A	90	LEU	2.9
1	B	0	HIS	2.9
1	A	117	GLN	2.9
1	A	558	SER	2.8
1	A	115	VAL	2.8
1	A	216	HIS	2.8
1	B	267	ARG	2.8
1	B	113	ASP	2.8
1	A	203	ASP	2.7
1	A	223	PHE	2.6
1	A	96	GLN	2.6
1	B	251	ASP	2.6
1	B	105	ALA	2.5
1	B	269	THR	2.5
1	B	250	ASP	2.5
1	A	221	CYS	2.4
1	B	252	PRO	2.4
1	B	264	GLN	2.3
1	B	254	ARG	2.3
1	A	101	VAL	2.2
1	A	250	ASP	2.2
1	A	232	LEU	2.2
1	A	231	GLN	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	97	GLU	2.2
1	A	92	LYS	2.2
1	B	377	GLY	2.2
1	B	266	CYS	2.2
1	A	564	THR	2.1
1	B	263	LYS	2.1
1	A	157	ILE	2.1
1	B	270	GLU	2.1
1	A	201	LYS	2.1
1	B	390	GLY	2.1
1	B	178	ALA	2.1
1	B	143	ASN	2.1
1	B	394	TRP	2.1
1	B	341	HIS	2.1
1	A	209	GLU	2.1
1	A	158	VAL	2.0
1	B	108	VAL	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.