



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 08:41 PM UTC

PDB ID : 3GAF / pdb_00003gaf
Title : 2.2A Crystal Structure of 7-Alpha-Hydroxysteroid Dehydrogenase from Brucella Melitensis
Authors : Seattle Structural Genomics Center for Infectious Disease (SSGCID)
Deposited on : 2009-02-17
Resolution : 2.20 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 ⓘ 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
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

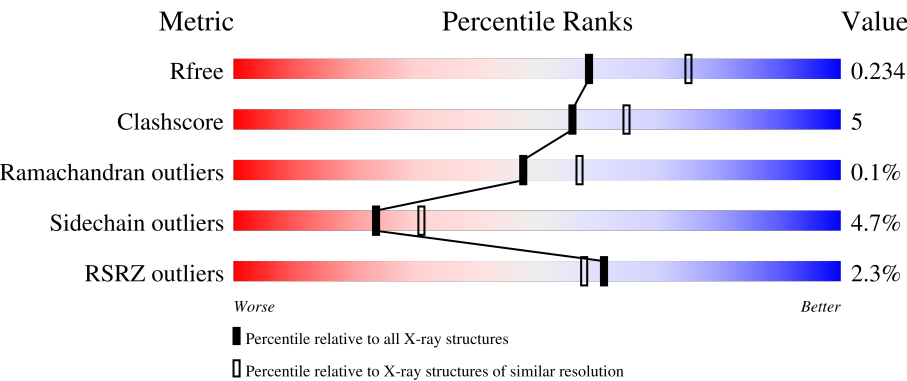
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R _{free}	180053	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	256	% 86%11%..
1	B	256	2% 80%14%.5%
1	C	256	2% 88%7%5%
1	D	256	2% 86%9%.
1	E	256	% 81%13%5%

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Mol	Chain	Length	Quality of chain
1	F	256	2% 83% 12%
1	G	256	2% 79% 15%
1	H	256	6% 81% 12% 6%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 14131 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 7-ALPHA-HYDROXYSTEROID DEHYDROGENASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	249	Total	C	N	O	S	0	0	0
			1763	1108	315	331	9			
1	B	243	Total	C	N	O	S	0	0	0
			1717	1078	306	324	9			
1	C	243	Total	C	N	O	S	0	0	0
			1706	1069	306	322	9			
1	D	245	Total	C	N	O	S	0	0	0
			1713	1076	308	320	9			
1	E	243	Total	C	N	O	S	0	0	0
			1716	1078	306	323	9			
1	F	246	Total	C	N	O	S	0	0	0
			1737	1090	310	328	9			
1	G	246	Total	C	N	O	S	0	0	0
			1733	1088	310	326	9			
1	H	241	Total	C	N	O	S	0	0	0
			1679	1054	300	316	9			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	SER	-	expression tag	UNP Q8YIN7
B	0	SER	-	expression tag	UNP Q8YIN7
C	0	SER	-	expression tag	UNP Q8YIN7
D	0	SER	-	expression tag	UNP Q8YIN7
E	0	SER	-	expression tag	UNP Q8YIN7
F	0	SER	-	expression tag	UNP Q8YIN7
G	0	SER	-	expression tag	UNP Q8YIN7
H	0	SER	-	expression tag	UNP Q8YIN7

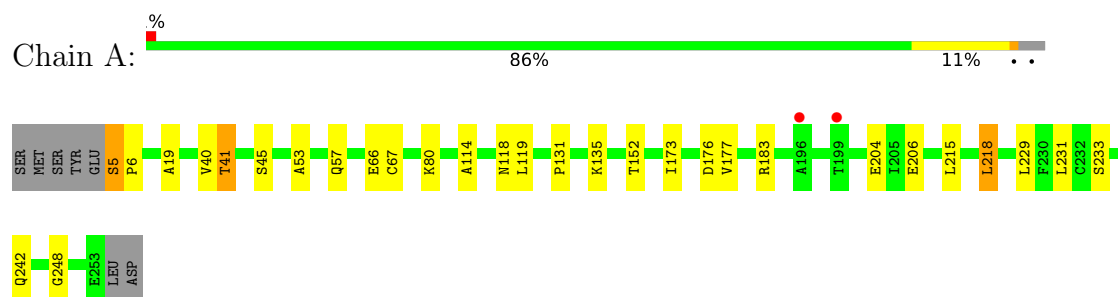
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	73	Total O 73 73	0	0
2	B	55	Total O 55 55	0	0
2	C	51	Total O 51 51	0	0
2	D	47	Total O 47 47	0	0
2	E	45	Total O 45 45	0	0
2	F	32	Total O 32 32	0	0
2	G	36	Total O 36 36	0	0
2	H	28	Total O 28 28	0	0

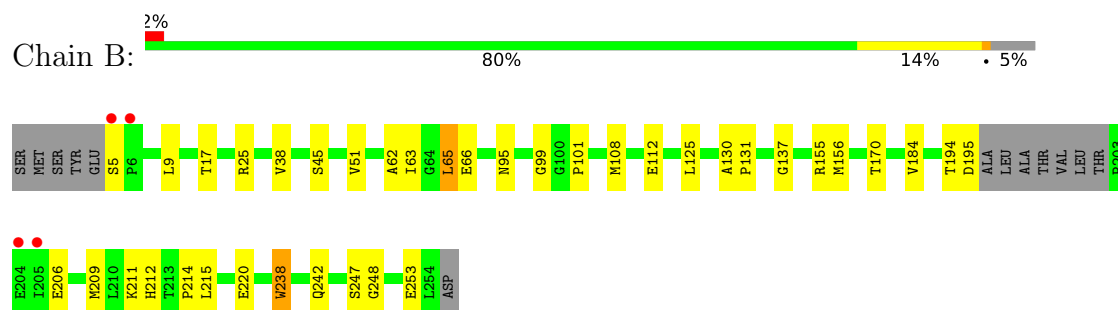
3 Residue-property plots [i](#)

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.

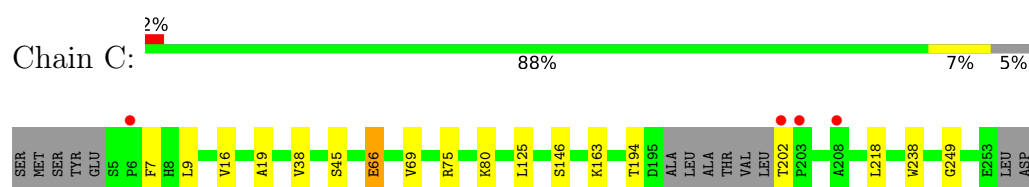
• Molecule 1: 7-ALPHA-HYDROXYSTEROID DEHYDROGENASE



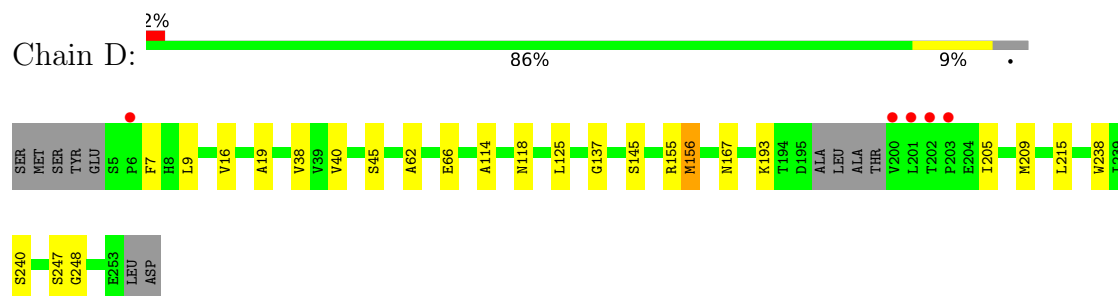
• Molecule 1: 7-ALPHA-HYDROXYSTEROID DEHYDROGENASE



• Molecule 1: 7-ALPHA-HYDROXYSTEROID DEHYDROGENASE

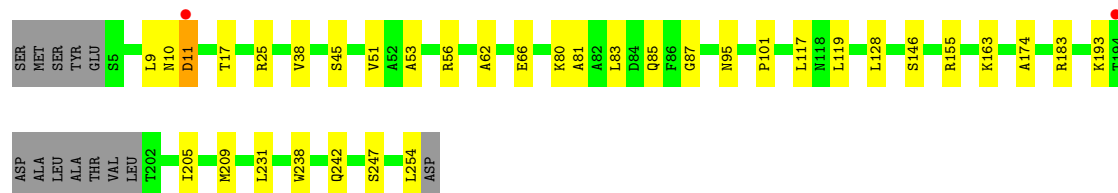


• Molecule 1: 7-ALPHA-HYDROXYSTEROID DEHYDROGENASE



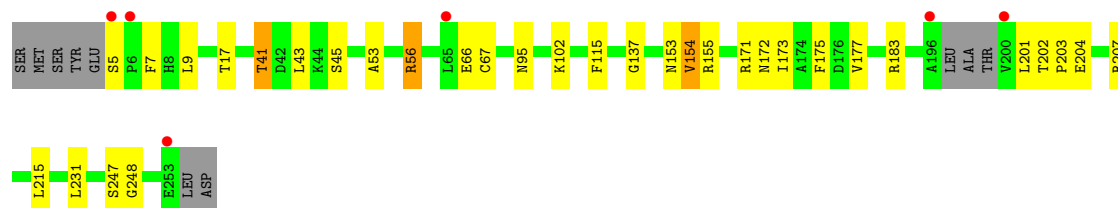
• Molecule 1: 7-ALPHA-HYDROXYSTEROID DEHYDROGENASE

Chain E: 



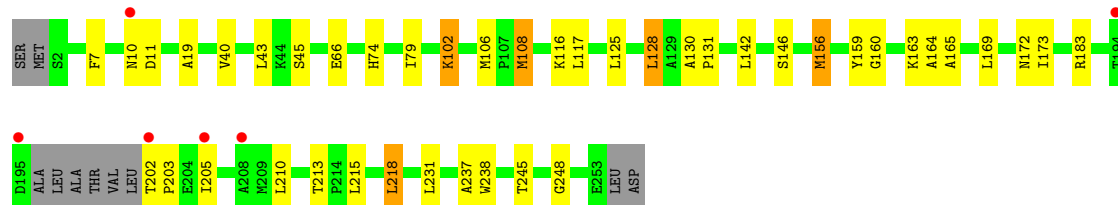
• Molecule 1: 7-ALPHA-HYDROXYSTEROID DEHYDROGENASE

Chain F: 



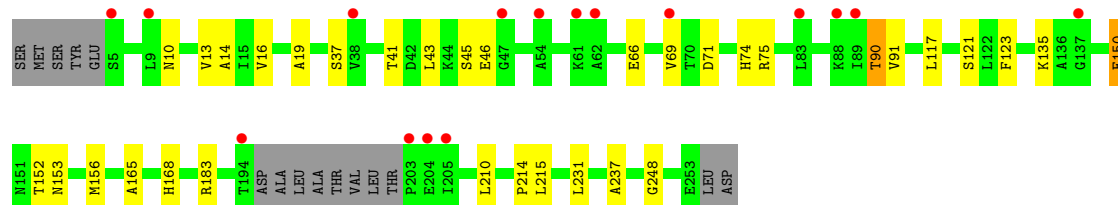
• Molecule 1: 7-ALPHA-HYDROXYSTEROID DEHYDROGENASE

Chain G: 



• Molecule 1: 7-ALPHA-HYDROXYSTEROID DEHYDROGENASE

Chain H: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	75.32Å 134.78Å 105.27Å 90.00° 109.52° 90.00°	Depositor
Resolution (Å)	48.51 – 2.20 48.51 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.1 (48.51-2.20) 99.0 (48.51-2.20)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.17 (at 2.20Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.196 , 0.244 (Not available) , 0.234	Depositor DCC
R_{free} test set	4955 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	39.5	Xtriage
Anisotropy	0.086	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 36.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.021 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	14131	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.62% 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	0.89	0/1792	1.01	5/2435 (0.2%)
1	B	0.82	0/1745	0.99	3/2368 (0.1%)
1	C	0.77	0/1734	0.96	1/2357 (0.0%)
1	D	0.75	0/1741	0.92	0/2365
1	E	0.75	0/1744	0.95	0/2368
1	F	0.76	1/1765 (0.1%)	0.93	1/2399 (0.0%)
1	G	0.73	0/1762	0.94	4/2395 (0.2%)
1	H	0.67	0/1707	0.95	0/2321
All	All	0.77	1/13990 (0.0%)	0.96	14/19008 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	F	154	VAL	CA-CB	5.06	1.60	1.54

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	5	SER	CA-C-N	-7.01	112.67	120.45
1	A	5	SER	C-N-CA	-7.01	112.67	120.45
1	G	213	THR	CA-C-N	6.20	126.38	119.32
1	G	213	THR	C-N-CA	6.20	126.38	119.32
1	B	99	GLY	N-CA-C	-5.62	104.30	112.17
1	F	172	ASN	N-CA-C	5.59	118.22	111.40
1	B	5	SER	CA-C-N	5.40	125.47	119.32
1	B	5	SER	C-N-CA	5.40	125.47	119.32
1	A	233	SER	CA-C-N	5.24	125.55	119.47
1	A	233	SER	C-N-CA	5.24	125.55	119.47
1	A	176	ASP	N-CA-C	5.08	116.51	111.07
1	G	202	THR	CA-C-N	5.03	126.13	119.84
1	G	202	THR	C-N-CA	5.03	126.13	119.84
1	C	38	VAL	N-CA-C	5.01	115.07	107.75

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	1763	0	1753	23	0
1	B	1717	0	1694	18	0
1	C	1706	0	1668	7	0
1	D	1713	0	1678	14	0
1	E	1716	0	1696	17	0
1	F	1737	0	1712	19	0
1	G	1733	0	1697	27	0
1	H	1679	0	1631	24	0
2	A	73	0	0	1	0
2	B	55	0	0	2	0
2	C	51	0	0	0	0
2	D	47	0	0	1	0
2	E	45	0	0	2	0
2	F	32	0	0	2	0
2	G	36	0	0	0	0
2	H	28	0	0	3	0
All	All	14131	0	13529	129	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 5.

All (129) 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:41:THR:HG23	1:A:67:CYS:HB3	1.55	0.89
1:F:41:THR:HG23	1:F:67:CYS:HB3	1.62	0.82
1:F:137:GLY:HA3	2:F:259:HOH:O	1.82	0.77
1:A:45:SER:HB3	1:A:66:GLU:HB2	1.69	0.72
1:H:183:ARG:NH2	1:H:231:LEU:O	2.17	0.72
1:F:45:SER:HB3	1:F:66:GLU:HB2	1.72	0.72
1:H:13:VAL:H	1:H:90:THR:HG22	1.58	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:146:SER:HB3	1:E:163:LYS:HG3	1.76	0.68
1:C:45:SER:HB3	1:C:66:GLU:HB2	1.76	0.68
1:D:156:MET:HG2	2:E:269:HOH:O	1.94	0.67
1:G:183:ARG:NH2	1:G:231:LEU:O	2.28	0.67
1:A:5:SER:CB	1:A:6:PRO:CD	2.73	0.66
1:G:108:MET:HA	1:G:108:MET:HE2	1.78	0.66
1:H:90:THR:HG23	1:H:91:VAL:HG23	1.76	0.66
1:B:137:GLY:HA3	2:B:256:HOH:O	1.97	0.65
1:G:172:ASN:ND2	1:H:153:ASN:H	1.96	0.63
1:E:10:ASN:O	1:E:11:ASP:HB2	1.99	0.62
1:A:5:SER:HB2	1:A:6:PRO:HD2	1.81	0.62
1:H:45:SER:HB3	1:H:66:GLU:HB2	1.82	0.62
1:H:156:MET:HG2	2:H:259:HOH:O	2.00	0.61
1:D:205:ILE:O	1:D:209:MET:HG3	2.01	0.60
1:A:53:ALA:O	1:A:57:GLN:HG2	2.01	0.60
1:B:156:MET:HE2	1:F:207:ARG:HE	1.66	0.59
1:F:183:ARG:NH2	1:F:231:LEU:O	2.36	0.59
1:G:45:SER:HB3	1:G:66:GLU:HB2	1.86	0.58
1:G:210:LEU:HD21	1:G:218:LEU:HD13	1.85	0.57
1:H:71:ASP:HB3	1:H:74:HIS:HB2	1.85	0.57
1:B:63:ILE:HD12	1:B:65:LEU:HD21	1.87	0.56
1:A:183:ARG:NH2	1:A:231:LEU:O	2.38	0.55
1:H:13:VAL:H	1:H:90:THR:CG2	2.20	0.55
1:G:172:ASN:HD22	1:H:153:ASN:H	1.54	0.55
1:A:131:PRO:O	1:A:135:LYS:HG2	2.06	0.55
1:A:5:SER:HB3	1:A:6:PRO:CD	2.38	0.54
1:B:156:MET:HE3	1:F:207:ARG:HH21	1.71	0.54
1:E:45:SER:HB3	1:E:66:GLU:HB2	1.89	0.54
1:D:45:SER:HB3	1:D:66:GLU:HB2	1.89	0.54
1:A:41:THR:HG22	2:A:327:HOH:O	2.08	0.54
1:B:130:ALA:HB3	1:B:131:PRO:HD3	1.89	0.54
1:D:137:GLY:HA2	2:D:273:HOH:O	2.08	0.52
1:H:183:ARG:HD2	1:H:237:ALA:O	2.09	0.52
1:B:45:SER:HB3	1:B:66:GLU:HB2	1.91	0.52
1:G:10:ASN:O	1:G:11:ASP:CB	2.57	0.51
1:G:102:LYS:HB3	1:G:106:MET:HE2	1.93	0.51
1:C:238:TRP:CZ3	1:D:247:SER:HA	2.46	0.51
1:G:146:SER:HB3	1:G:163:LYS:HG3	1.91	0.51
1:E:242:GLN:NE2	2:E:337:HOH:O	2.17	0.51
1:F:41:THR:CG2	1:F:67:CYS:HB3	2.38	0.51
1:B:242:GLN:OE1	2:B:284:HOH:O	2.20	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:215:LEU:HD12	1:A:248:GLY:HA2	1.92	0.50
1:D:16:VAL:HG12	1:D:19:ALA:HB2	1.93	0.50
1:G:183:ARG:HD2	1:G:237:ALA:O	2.12	0.50
1:A:41:THR:CG2	1:A:67:CYS:HB3	2.36	0.50
1:G:215:LEU:HD12	1:G:248:GLY:HA2	1.94	0.50
1:A:5:SER:CB	1:A:6:PRO:HD2	2.42	0.49
1:D:19:ALA:HB3	1:D:40:VAL:HG13	1.93	0.49
1:E:38:VAL:O	1:E:62:ALA:HA	2.12	0.49
1:B:17:THR:O	1:B:95:ASN:HB3	2.13	0.49
1:G:130:ALA:HB3	1:G:131:PRO:HD3	1.95	0.49
1:E:183:ARG:NH2	1:E:231:LEU:O	2.45	0.48
1:B:238:TRP:CZ3	1:E:247:SER:HA	2.48	0.48
1:H:156:MET:CG	2:H:259:HOH:O	2.58	0.48
1:E:83:LEU:O	1:E:87:GLY:HA2	2.14	0.47
1:A:5:SER:HB3	1:A:6:PRO:HD3	1.96	0.47
1:D:155:ARG:O	1:D:156:MET:HE3	2.15	0.47
1:A:152:THR:CG2	1:F:171:ARG:HB3	2.43	0.47
1:B:214:PRO:HB3	1:E:174:ALA:O	2.14	0.47
1:G:19:ALA:HB3	1:G:40:VAL:HG13	1.97	0.47
1:B:25:ARG:HG3	1:B:51:VAL:HG22	1.96	0.47
1:B:211:LYS:HE2	1:B:212:HIS:NE2	2.30	0.47
1:C:146:SER:HB3	1:C:163:LYS:HG3	1.95	0.47
1:B:215:LEU:HD12	1:B:248:GLY:HA2	1.97	0.47
1:C:7:PHE:CD1	1:D:7:PHE:CD1	3.03	0.47
1:G:10:ASN:O	1:G:11:ASP:HB3	2.15	0.46
1:A:5:SER:HB2	1:A:6:PRO:CD	2.41	0.46
1:D:38:VAL:O	1:D:62:ALA:HA	2.15	0.46
1:C:69:VAL:O	1:C:75:ARG:HD3	2.16	0.46
1:C:249:GLY:HA2	1:D:240:SER:O	2.16	0.46
1:E:17:THR:O	1:E:95:ASN:HB3	2.16	0.46
1:H:215:LEU:HD12	1:H:248:GLY:HA2	1.98	0.46
1:A:152:THR:HG21	1:F:171:ARG:HB3	1.98	0.45
1:E:25:ARG:HG3	1:E:51:VAL:CG2	2.46	0.45
1:B:38:VAL:O	1:B:62:ALA:HA	2.17	0.45
1:G:210:LEU:HD21	1:G:218:LEU:CD1	2.47	0.45
1:G:66:GLU:OE1	1:G:74:HIS:HE1	1.99	0.45
1:H:14:ALA:HA	1:H:91:VAL:O	2.17	0.45
1:A:119:LEU:HD21	1:F:115:PHE:CZ	2.53	0.44
1:E:25:ARG:HG3	1:E:51:VAL:HG22	1.99	0.44
1:G:108:MET:HA	1:G:108:MET:CE	2.44	0.44
1:A:206:GLU:HG3	1:A:218:LEU:HD22	1.98	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:108:MET:HE1	1:H:123:PHE:CE2	2.51	0.44
1:D:215:LEU:HD12	1:D:248:GLY:HA2	1.99	0.44
1:F:247:SER:OG	2:F:260:HOH:O	2.21	0.44
1:G:165:ALA:HB2	1:H:165:ALA:HB2	1.99	0.44
1:B:108:MET:O	1:B:112:GLU:HG3	2.18	0.43
1:F:154:VAL:HG22	1:F:155:ARG:HG3	1.98	0.43
1:H:69:VAL:O	1:H:75:ARG:NH1	2.46	0.43
1:D:145:SER:OG	1:D:167:ASN:ND2	2.52	0.43
1:A:242:GLN:NE2	1:G:245:THR:HB	2.34	0.43
1:A:173:ILE:O	1:A:177:VAL:HG22	2.17	0.43
1:B:101:PRO:HA	1:B:155:ARG:O	2.19	0.43
1:B:170:THR:HG23	1:B:184:VAL:HG12	1.99	0.43
1:E:101:PRO:HA	1:E:155:ARG:O	2.18	0.43
1:C:16:VAL:HG12	1:C:19:ALA:HB2	2.00	0.43
1:F:202:THR:HB	1:F:203:PRO:CD	2.50	0.42
1:F:215:LEU:HD12	1:F:248:GLY:HA2	1.99	0.42
1:G:156:MET:HG3	1:G:160:GLY:H	1.84	0.42
1:G:164:ALA:HB2	1:H:168:HIS:CD2	2.54	0.42
1:H:13:VAL:N	1:H:90:THR:HG22	2.30	0.42
1:G:79:ILE:HG13	1:G:128:LEU:HB3	2.02	0.42
1:E:53:ALA:HA	1:E:56:ARG:NH1	2.35	0.42
1:H:46:GLU:H	1:H:46:GLU:CD	2.27	0.42
1:H:150:GLU:OE2	2:H:275:HOH:O	2.22	0.42
1:F:175:PHE:CD2	1:H:214:PRO:HG3	2.55	0.41
1:G:169:LEU:O	1:G:173:ILE:HG12	2.20	0.41
1:F:53:ALA:HA	1:F:56:ARG:NH1	2.34	0.41
1:A:229:LEU:HD23	1:G:7:PHE:CE2	2.54	0.41
1:G:172:ASN:ND2	1:H:152:THR:HA	2.35	0.41
1:F:173:ILE:O	1:F:177:VAL:HG22	2.21	0.41
1:E:81:ALA:O	1:E:85:GLN:HB2	2.20	0.41
1:E:205:ILE:O	1:E:209:MET:HG3	2.21	0.41
1:A:114:ALA:O	1:A:118:ASN:HB2	2.21	0.41
1:B:247:SER:HA	1:E:238:TRP:CZ3	2.55	0.41
1:F:17:THR:O	1:F:95:ASN:HB3	2.20	0.41
1:D:114:ALA:O	1:D:118:ASN:HB2	2.21	0.40
1:H:16:VAL:HG12	1:H:19:ALA:HB2	2.03	0.40
1:F:5:SER:HB3	1:F:7:PHE:H	1.87	0.40
1:A:19:ALA:HB3	1:A:40:VAL:HG13	2.04	0.40
1:H:117:LEU:O	1:H:121:SER:OG	2.34	0.40
1:G:156:MET:HE2	1:G:159:TYR:HB3	2.03	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	247/256 (96%)	242 (98%)	5 (2%)	0	100	100
1	B	239/256 (93%)	233 (98%)	6 (2%)	0	100	100
1	C	239/256 (93%)	234 (98%)	5 (2%)	0	100	100
1	D	241/256 (94%)	233 (97%)	8 (3%)	0	100	100
1	E	239/256 (93%)	231 (97%)	8 (3%)	0	100	100
1	F	242/256 (94%)	231 (96%)	11 (4%)	0	100	100
1	G	242/256 (94%)	233 (96%)	8 (3%)	1 (0%)	30	34
1	H	237/256 (93%)	229 (97%)	8 (3%)	0	100	100
All	All	1926/2048 (94%)	1866 (97%)	59 (3%)	1 (0%)	48	57

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	G	203	PRO

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	170/184 (92%)	166 (98%)	4 (2%)	43	58
1	B	165/184 (90%)	155 (94%)	10 (6%)	17	20
1	C	162/184 (88%)	155 (96%)	7 (4%)	26	35
1	D	161/184 (88%)	156 (97%)	5 (3%)	35	48

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	E	165/184 (90%)	157 (95%)	8 (5%)	23	30
1	F	167/184 (91%)	159 (95%)	8 (5%)	23	30
1	G	165/184 (90%)	153 (93%)	12 (7%)	13	15
1	H	157/184 (85%)	149 (95%)	8 (5%)	21	27
All	All	1312/1472 (89%)	1250 (95%)	62 (5%)	23	31

All (62) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	41	THR
1	A	80	LYS
1	A	204	GLU
1	A	218	LEU
1	B	9	LEU
1	B	65	LEU
1	B	125	LEU
1	B	194	THR
1	B	195	ASP
1	B	206	GLU
1	B	209	MET
1	B	220	GLU
1	B	238	TRP
1	B	253	GLU
1	C	9	LEU
1	C	66	GLU
1	C	80	LYS
1	C	125	LEU
1	C	194	THR
1	C	202	THR
1	C	218	LEU
1	D	9	LEU
1	D	125	LEU
1	D	156	MET
1	D	193	LYS
1	D	238	TRP
1	E	9	LEU
1	E	11	ASP
1	E	80	LYS
1	E	117	LEU
1	E	119	LEU
1	E	128	LEU

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Mol	Chain	Res	Type
1	E	193	LYS
1	E	254	LEU
1	F	9	LEU
1	F	41	THR
1	F	43	LEU
1	F	56	ARG
1	F	102	LYS
1	F	153	ASN
1	F	201	LEU
1	F	204	GLU
1	G	43	LEU
1	G	102	LYS
1	G	108	MET
1	G	116	LYS
1	G	117	LEU
1	G	125	LEU
1	G	128	LEU
1	G	142	LEU
1	G	156	MET
1	G	205	ILE
1	G	218	LEU
1	G	238	TRP
1	H	10	ASN
1	H	37	SER
1	H	41	THR
1	H	43	LEU
1	H	90	THR
1	H	135	LYS
1	H	150	GLU
1	H	210	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	10	ASN
1	A	242	GLN
1	B	73	GLN
1	B	85	GLN
1	B	132	HIS
1	B	222	GLN
1	B	242	GLN
1	C	85	GLN

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Mol	Chain	Res	Type
1	C	132	HIS
1	C	222	GLN
1	D	10	ASN
1	D	167	ASN
1	E	8	HIS
1	E	10	ASN
1	E	94	ASN
1	E	143	ASN
1	F	94	ASN
1	F	222	GLN
1	G	10	ASN
1	G	94	ASN
1	G	172	ASN
1	H	8	HIS
1	H	73	GLN
1	H	212	HIS
1	H	222	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](#)

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 ⓘ

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	249/256 (97%)	-0.17	2 (0%) 82 80	22, 36, 54, 69	0
1	B	243/256 (94%)	-0.07	4 (1%) 70 67	23, 38, 55, 69	0
1	C	243/256 (94%)	0.06	4 (1%) 70 67	26, 42, 60, 89	0
1	D	245/256 (95%)	0.06	5 (2%) 65 62	26, 42, 64, 89	0
1	E	243/256 (94%)	0.10	2 (0%) 82 80	25, 41, 60, 65	0
1	F	246/256 (96%)	0.16	6 (2%) 59 56	26, 44, 66, 79	0
1	G	246/256 (96%)	0.28	6 (2%) 59 56	28, 47, 69, 89	0
1	H	241/256 (94%)	0.57	16 (6%) 24 21	33, 52, 75, 83	0
All	All	1956/2048 (95%)	0.12	45 (2%) 61 58	22, 42, 69, 89	0

All (45) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	200	VAL	4.7
1	H	203	PRO	4.4
1	F	200	VAL	4.3
1	C	208	ALA	3.8
1	C	203	PRO	3.3
1	D	201	LEU	3.2
1	F	196	ALA	3.1
1	A	196	ALA	3.1
1	F	6	PRO	3.1
1	G	205	ILE	2.7
1	D	202	THR	2.7
1	E	194	THR	2.7
1	G	195	ASP	2.7
1	H	38	VAL	2.7
1	B	205	ILE	2.7
1	G	194	THR	2.6

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Mol	Chain	Res	Type	RSRZ
1	G	202	THR	2.6
1	H	5	SER	2.6
1	A	199	THR	2.5
1	H	204	GLU	2.5
1	B	204	GLU	2.4
1	C	202	THR	2.4
1	H	137	GLY	2.3
1	D	203	PRO	2.2
1	E	11	ASP	2.2
1	G	208	ALA	2.2
1	C	6	PRO	2.2
1	H	62	ALA	2.2
1	F	5	SER	2.2
1	H	205	ILE	2.2
1	F	253	GLU	2.2
1	B	5	SER	2.2
1	H	194	THR	2.2
1	B	6	PRO	2.1
1	H	47	GLY	2.1
1	H	89	ILE	2.1
1	F	65	LEU	2.1
1	H	9	LEU	2.1
1	D	6	PRO	2.1
1	H	69	VAL	2.1
1	H	88	LYS	2.0
1	H	54	ALA	2.0
1	H	61	LYS	2.0
1	G	10	ASN	2.0
1	H	83	LEU	2.0

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

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.