



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 10:17 AM UTC

PDB ID : 3L8M / pdb_00003l8m
Title : Crystal Structure of a probable thiamine pyrophosphokinase from *Staphylococcus saprophyticus* subsp. *saprophyticus*. Northeast Structural Genomics Consortium target id SyR86
Authors : Seetharaman, J.; Lew, S.; Wang, D.; Janjua, H.; Cunningham, K.; Owens, L.; Xiao, R.; Liu, J.; Baran, M.C.; Acton, T.B.; Montelione, G.T.; Hunt, J.F.; Tong, L.; Northeast Structural Genomics Consortium (NESG)
Deposited on : 2009-12-31
Resolution : 2.40 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)

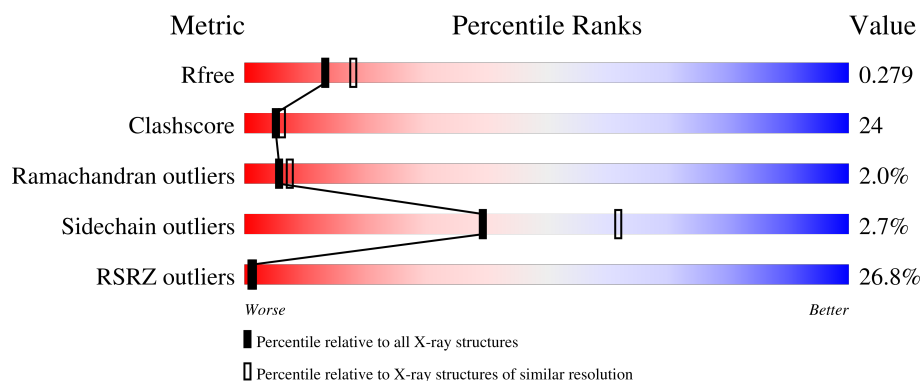
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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	212	23% 53% 35% 9%
1	B	212	23% 50% 32% 17%

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 2904 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 probable thiamine pyrophosphokinase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	193	Total	C	N	O	S	Se	0	0	0
			1500	948	259	287	2	4			
1	B	177	Total	C	N	O	S	Se	0	0	0
			1305	826	227	246	2	4			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	212	GLU	-	expression tag	UNP Q49X04
B	212	GLU	-	expression tag	UNP Q49X04

- Molecule 2 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Na	0	0
			2	2		
2	B	1	Total	Na	0	0
			1	1		

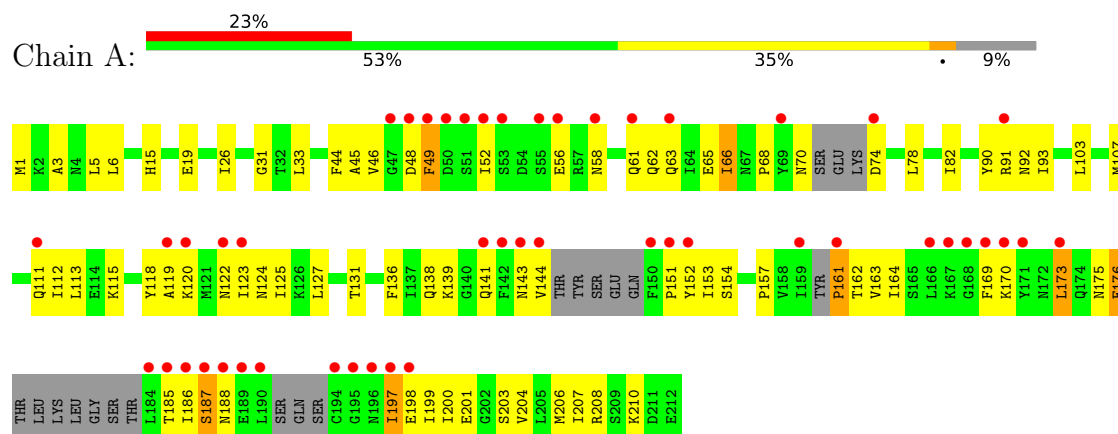
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	58	Total	O	0	0
			58	58		
3	B	38	Total	O	0	0
			38	38		

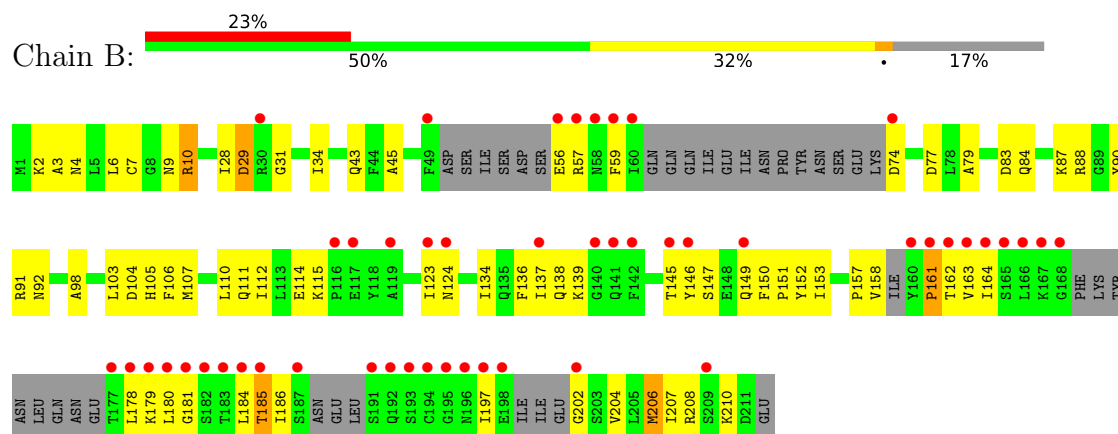
3 Residue-property plots [i](#)

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: probable thiamine pyrophosphokinase



- Molecule 1: probable thiamine pyrophosphokinase



4 Data and refinement statistics

Property	Value	Source
Space group	P 63	Depositor
Cell constants a, b, c, α , β , γ	101.87Å 101.87Å 88.02Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	33.35 – 2.40 33.35 – 2.40	Depositor EDS
% Data completeness (in resolution range)	92.6 (33.35-2.40) 99.0 (33.35-2.40)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.45 (at 2.39Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.238 , 0.271 0.257 , 0.279	Depositor DCC
R_{free} test set	1831 reflections (4.58%)	wwPDB-VP
Wilson B-factor (Å ²)	35.9	Xtriage
Anisotropy	0.127	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 67.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.037 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	2904	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.95% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: NA

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	0.50	2/1518 (0.1%)	1.01	5/2039 (0.2%)
1	B	0.48	1/1317 (0.1%)	0.87	1/1773 (0.1%)
All	All	0.49	3/2835 (0.1%)	0.95	6/3812 (0.2%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	206	MSE	SE-CE	-6.64	1.75	1.95
1	B	206	MSE	SE-CE	-6.17	1.76	1.95
1	A	206	MSE	CG-SE	-5.54	1.78	1.95

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	48	ASP	N-CA-C	-16.69	81.92	109.96
1	A	49	PHE	N-CA-C	-10.15	93.50	109.23
1	A	161	PRO	CA-N-CD	-9.29	99.00	112.00
1	A	48	ASP	CB-CA-C	-7.85	96.57	109.53
1	B	43	GLN	N-CA-C	-5.97	104.78	111.28
1	A	143	ASN	CB-CA-C	-5.11	109.02	116.53

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	A	1500	0	1455	76	0
1	B	1305	0	1257	64	0
2	A	2	0	0	0	0
2	B	1	0	0	0	0
3	A	58	0	0	1	0
3	B	38	0	0	2	0
All	All	2904	0	2712	135	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 24.

All (135) 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:164:ILE:HG12	1:A:173:LEU:H	1.13	1.09
1:B:134:ILE:HG12	1:B:207:ILE:HD12	1.37	1.02
1:A:185:THR:HB	1:A:188:ASN:HD21	1.28	0.94
1:B:10:ARG:HH11	1:B:10:ARG:HB3	1.34	0.92
1:B:7:CYS:HB3	1:B:28:ILE:HB	1.57	0.84
1:B:178:LEU:HG	1:B:180:LEU:HD22	1.59	0.84
1:A:175:ASN:HD22	1:A:176:GLU:H	1.29	0.79
1:A:62:GLN:HG3	1:A:63:GLN:HG3	1.64	0.79
1:A:161:PRO:HG2	1:A:176:GLU:CB	2.12	0.79
1:B:9:ASN:HB3	1:B:34:ILE:HD13	1.63	0.79
1:B:3:ALA:HB2	1:B:90:TYR:CD1	2.18	0.78
1:A:164:ILE:HG12	1:A:173:LEU:N	1.96	0.78
1:A:164:ILE:CG1	1:A:173:LEU:H	1.97	0.76
1:B:206:MSE:HE2	1:B:208:ARG:HD3	1.69	0.74
1:A:45:ALA:O	1:A:66:ILE:HD13	1.89	0.73
1:B:2:LYS:HE2	1:B:4:ASN:HD21	1.51	0.73
1:A:185:THR:HB	1:A:188:ASN:ND2	2.05	0.69
1:A:161:PRO:HG2	1:A:176:GLU:HB2	1.73	0.69
1:B:57:ARG:HA	3:B:237:HOH:O	1.93	0.69
1:A:91:ARG:NH1	1:A:122:ASN:HD22	1.90	0.69
1:A:175:ASN:ND2	1:A:176:GLU:H	1.92	0.68
1:B:28:ILE:O	1:B:29:ASP:HB2	1.93	0.68
1:A:151:PRO:HD2	1:A:210:LYS:HA	1.77	0.67
1:A:163:VAL:HG22	1:A:199:ILE:HD13	1.78	0.65
1:A:161:PRO:HG2	1:A:176:GLU:HB3	1.77	0.65
1:A:175:ASN:HD22	1:A:176:GLU:N	1.94	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:9:ASN:HB3	1:B:34:ILE:CD1	2.27	0.65
1:A:3:ALA:HB2	1:A:90:TYR:CD1	2.32	0.65
1:A:61:GLN:O	1:A:61:GLN:HG3	1.96	0.65
1:B:138:GLN:HG2	1:B:139:LYS:H	1.63	0.63
1:A:91:ARG:O	1:A:123:ILE:HD12	1.98	0.62
1:B:83:ASP:OD2	1:B:87:LYS:HE2	2.00	0.62
1:B:153:ILE:O	1:B:186:ILE:HG22	2.00	0.61
1:B:111:GLN:HG3	3:B:235:HOH:O	2.00	0.61
1:A:66:ILE:HG12	1:A:68:PRO:HD3	1.81	0.61
1:B:180:LEU:H	1:B:180:LEU:HD23	1.66	0.61
1:A:175:ASN:ND2	1:A:176:GLU:N	2.49	0.60
1:B:98:ALA:HA	1:B:105:HIS:HE1	1.67	0.60
1:B:106:PHE:CD2	1:B:107:MSE:HE3	2.36	0.60
1:A:66:ILE:HD13	1:A:66:ILE:H	1.65	0.60
3:A:228:HOH:O	1:B:157:PRO:HG2	2.00	0.59
1:A:197:ILE:C	1:A:197:ILE:HD13	2.27	0.59
1:A:186:ILE:O	1:A:187:SER:HB2	2.02	0.59
1:A:111:GLN:CD	1:B:178:LEU:O	2.47	0.58
1:A:161:PRO:HA	1:A:201:GLU:O	2.05	0.57
1:B:6:LEU:O	1:B:31:GLY:HA3	2.03	0.57
1:B:9:ASN:HD22	1:B:34:ILE:HD11	1.68	0.57
1:B:10:ARG:HH11	1:B:10:ARG:CB	2.15	0.56
1:A:131:THR:HG23	1:A:210:LYS:HG2	1.87	0.56
1:B:161:PRO:HA	1:B:202:GLY:N	2.21	0.56
1:A:207:ILE:HD11	1:B:103:LEU:HB2	1.87	0.55
1:B:123:ILE:HG12	1:B:124:ASN:N	2.22	0.55
1:B:10:ARG:HB3	1:B:10:ARG:NH1	2.14	0.55
1:A:141:GLN:HG3	1:A:198:GLU:HB2	1.89	0.55
1:B:153:ILE:N	1:B:153:ILE:HD12	2.23	0.54
1:B:147:SER:OG	1:B:150:PHE:HB2	2.07	0.54
1:A:5:LEU:HG	1:A:26:ILE:HB	1.90	0.54
1:A:61:GLN:HA	1:A:65:GLU:HB3	1.90	0.53
1:A:52:ILE:HG22	1:A:56:GLU:OE2	2.08	0.53
1:B:157:PRO:CB	1:B:161:PRO:HG2	2.39	0.53
1:A:93:ILE:HB	1:A:125:ILE:HD13	1.90	0.53
1:B:181:GLY:HA2	1:B:184:LEU:HD12	1.90	0.53
1:A:207:ILE:HD13	1:B:104:ASP:N	2.25	0.52
1:A:153:ILE:HD12	1:A:153:ILE:N	2.25	0.51
1:A:119:ALA:HB1	1:A:138:GLN:OE1	2.10	0.51
1:A:136:PHE:CE1	1:A:203:SER:HB2	2.46	0.51
1:B:84:GLN:O	1:B:88:ARG:HG3	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:157:PRO:HA	1:B:204:VAL:HG12	1.93	0.51
1:B:152:TYR:C	1:B:153:ILE:HD12	2.36	0.51
1:A:103:LEU:O	1:A:107:MSE:HG2	2.11	0.50
1:A:186:ILE:O	1:A:186:ILE:HG22	2.11	0.50
1:B:163:VAL:C	1:B:164:ILE:HD12	2.37	0.50
1:A:33:LEU:HD21	1:A:49:PHE:HB2	1.94	0.50
1:A:58:ASN:O	1:A:62:GLN:HG2	2.13	0.49
1:A:91:ARG:HD2	1:A:122:ASN:O	2.12	0.49
1:A:131:THR:O	1:A:131:THR:HG22	2.12	0.49
1:A:26:ILE:HD12	1:A:44:PHE:CE2	2.49	0.48
1:A:46:VAL:HG22	1:A:66:ILE:HD11	1.95	0.48
1:A:33:LEU:HD22	1:A:33:LEU:N	2.29	0.48
1:A:70:ASN:O	1:A:74:ASP:HB2	2.13	0.48
1:A:161:PRO:O	1:A:161:PRO:HD2	2.13	0.48
1:B:164:ILE:HD12	1:B:164:ILE:N	2.28	0.47
1:A:92:ASN:C	1:A:93:ILE:HD13	2.39	0.47
1:A:152:TYR:C	1:A:153:ILE:HD12	2.39	0.47
1:A:162:THR:HG21	1:A:173:LEU:HD13	1.97	0.47
1:A:15:HIS:O	1:A:19:GLU:HG3	2.15	0.47
1:B:163:VAL:HG12	1:B:164:ILE:N	2.29	0.47
1:A:1:MSE:HG2	1:A:90:TYR:CD1	2.50	0.46
1:B:157:PRO:HB3	1:B:161:PRO:HG2	1.97	0.46
1:B:106:PHE:HD2	1:B:107:MSE:HE3	1.79	0.46
1:B:114:GLU:O	1:B:115:LYS:C	2.59	0.45
1:B:151:PRO:HD2	1:B:210:LYS:HA	1.98	0.45
1:A:131:THR:HA	1:A:210:LYS:CD	2.46	0.45
1:B:106:PHE:CD2	1:B:107:MSE:CE	2.98	0.45
1:A:138:GLN:HG2	1:A:139:LYS:H	1.82	0.45
1:B:184:LEU:O	1:B:185:THR:C	2.57	0.45
1:A:118:TYR:HB3	1:A:123:ILE:HG21	1.98	0.45
1:A:118:TYR:HB3	1:A:123:ILE:CG2	2.48	0.44
1:A:197:ILE:HD13	1:A:198:GLU:N	2.32	0.44
1:B:84:GLN:NE2	1:B:84:GLN:HA	2.32	0.44
1:B:180:LEU:H	1:B:180:LEU:CD2	2.29	0.44
1:B:206:MSE:C	1:B:207:ILE:HD13	2.43	0.44
1:B:138:GLN:CG	1:B:139:LYS:H	2.30	0.44
1:A:92:ASN:O	1:A:93:ILE:HD13	2.17	0.44
1:A:200:ILE:HD12	1:A:200:ILE:N	2.33	0.44
1:A:33:LEU:HD22	1:A:33:LEU:H	1.83	0.43
1:B:107:MSE:HE2	1:B:110:LEU:HD12	2.00	0.43
1:A:111:GLN:OE1	1:B:179:LYS:HA	2.17	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:138:GLN:HG2	1:B:139:LYS:N	2.30	0.43
1:A:144:VAL:O	1:A:144:VAL:HG12	2.19	0.43
1:B:91:ARG:HG3	1:B:92:ASN:N	2.33	0.43
1:B:79:ALA:HB2	1:B:112:ILE:HD11	2.01	0.42
1:A:136:PHE:CZ	1:A:203:SER:HB2	2.55	0.42
1:A:123:ILE:HG13	1:A:124:ASN:N	2.33	0.42
1:B:137:ILE:HD12	1:B:137:ILE:N	2.35	0.42
1:A:93:ILE:HG12	1:A:123:ILE:HD11	2.02	0.42
1:B:45:ALA:O	1:B:56:GLU:HA	2.20	0.42
1:A:45:ALA:O	1:A:66:ILE:CD1	2.65	0.42
1:A:154:SER:HB2	1:A:207:ILE:HG23	2.02	0.42
1:B:134:ILE:HG23	1:B:207:ILE:CD1	2.50	0.41
1:A:120:LYS:H	1:A:120:LYS:HG2	1.51	0.41
1:A:207:ILE:CD1	1:B:103:LEU:HB2	2.49	0.41
1:B:137:ILE:HG22	1:B:197:ILE:HD12	2.02	0.41
1:A:61:GLN:NE2	1:A:65:GLU:HG2	2.36	0.41
1:A:78:LEU:O	1:A:82:ILE:HG12	2.20	0.41
1:B:136:PHE:CE1	1:B:158:VAL:HG21	2.56	0.41
1:B:29:ASP:OD1	1:B:77:ASP:OD2	2.39	0.41
1:B:145:THR:HG22	1:B:146:TYR:N	2.36	0.41
1:B:137:ILE:HG22	1:B:197:ILE:CD1	2.51	0.40
1:A:112:ILE:HA	1:A:115:LYS:HG3	2.03	0.40
1:A:6:LEU:O	1:A:31:GLY:HA3	2.22	0.40
1:A:113:LEU:HD11	1:A:127:LEU:HG	2.02	0.40
1:B:74:ASP:OD2	1:B:74:ASP:N	2.53	0.40
1:B:107:MSE:O	1:B:111:GLN:HG2	2.22	0.40
1:A:157:PRO:HA	1:A:204:VAL:HG12	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	181/212 (85%)	162 (90%)	16 (9%)	3 (2%)	7	10
1	B	163/212 (77%)	152 (93%)	7 (4%)	4 (2%)	4	5
All	All	344/424 (81%)	314 (91%)	23 (7%)	7 (2%)	6	7

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	187	SER
1	B	59	PHE
1	A	170	LYS
1	B	29	ASP
1	B	185	THR
1	B	161	PRO
1	A	169	PHE

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	160/185 (86%)	155 (97%)	5 (3%)	35	57
1	B	133/185 (72%)	130 (98%)	3 (2%)	44	66
All	All	293/370 (79%)	285 (97%)	8 (3%)	39	62

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

Mol	Chain	Res	Type
1	A	66	ILE
1	A	173	LEU
1	A	176	GLU
1	A	197	ILE
1	A	208	ARG
1	B	10	ARG
1	B	149	GLN
1	B	162	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	20	HIS
1	A	61	GLN
1	A	62	GLN
1	A	67	ASN
1	A	105	HIS
1	A	122	ASN
1	A	143	ASN
1	A	172	ASN
1	A	175	ASN
1	A	188	ASN
1	B	4	ASN
1	B	9	ASN
1	B	11	ASN
1	B	24	HIS
1	B	92	ASN
1	B	149	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](#)

Of 3 ligands modelled in this entry, 3 are monoatomic - leaving 0 for Mogul analysis.

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.

No monomer is involved in short contacts.

5.7 Other polymers

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	189/212 (89%)	1.02	48 (25%) 1 1	17, 40, 77, 83	0
1	B	173/212 (81%)	1.20	49 (28%) 1 1	21, 43, 81, 87	0
All	All	362/424 (85%)	1.11	97 (26%) 1 1	17, 42, 79, 87	0

All (97) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	186	ILE	7.0
1	B	163	VAL	6.6
1	A	184	LEU	6.4
1	A	144	VAL	5.6
1	B	56	GLU	5.5
1	B	49	PHE	5.5
1	A	52	ILE	5.4
1	B	194	CYS	5.4
1	B	60	ILE	5.2
1	B	202	GLY	5.1
1	B	178	LEU	4.7
1	A	168	GLY	4.6
1	A	195	GLY	4.6
1	A	169	PHE	4.6
1	B	160	TYR	4.6
1	B	58	ASN	4.4
1	A	48	ASP	4.4
1	A	185	THR	4.4
1	B	57	ARG	4.4
1	A	171	TYR	4.2
1	A	170	LYS	4.2
1	B	181	GLY	4.1
1	A	161	PRO	4.0
1	B	164	ILE	4.0

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Mol	Chain	Res	Type	RSRZ
1	A	50	ASP	3.9
1	B	187	SER	3.7
1	B	198	GLU	3.7
1	B	162	THR	3.6
1	A	190	LEU	3.6
1	B	193	SER	3.6
1	B	59	PHE	3.6
1	B	166	LEU	3.6
1	B	177	THR	3.6
1	A	69	TYR	3.5
1	B	140	GLY	3.5
1	A	196	ASN	3.5
1	B	197	ILE	3.5
1	B	165	SER	3.4
1	B	168	GLY	3.4
1	A	194	CYS	3.3
1	B	180	LEU	3.3
1	A	61	GLN	3.3
1	B	184	LEU	3.3
1	B	185	THR	3.3
1	B	146	TYR	3.2
1	A	188	ASN	3.0
1	B	167	LYS	3.0
1	A	47	GLY	2.9
1	A	58	ASN	2.9
1	B	196	ASN	2.9
1	B	145	THR	2.8
1	A	151	PRO	2.8
1	B	141	GLN	2.8
1	B	161	PRO	2.7
1	B	142	PHE	2.7
1	A	111	GLN	2.7
1	B	123	ILE	2.7
1	A	63	GLN	2.7
1	A	143	ASN	2.6
1	A	122	ASN	2.6
1	B	195	GLY	2.6
1	A	119	ALA	2.5
1	A	189	GLU	2.5
1	A	49	PHE	2.5
1	B	182	SER	2.5
1	B	183	THR	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	123	ILE	2.5
1	A	166	LEU	2.4
1	B	192	GLN	2.4
1	A	141	GLN	2.4
1	A	142	PHE	2.4
1	B	116	PRO	2.4
1	A	55	SER	2.4
1	A	187	SER	2.4
1	B	30	ARG	2.4
1	B	124	ASN	2.4
1	B	179	LYS	2.4
1	A	197	ILE	2.3
1	B	149	GLN	2.3
1	A	167	LYS	2.3
1	A	74	ASP	2.3
1	A	53	SER	2.3
1	A	159	ILE	2.3
1	B	117	GLU	2.2
1	B	209	SER	2.2
1	B	191	SER	2.2
1	A	198	GLU	2.2
1	A	150	PHE	2.2
1	A	173	LEU	2.2
1	A	152	TYR	2.1
1	B	137	ILE	2.1
1	A	51	SER	2.1
1	B	119	ALA	2.1
1	A	56	GLU	2.1
1	A	120	LYS	2.1
1	B	74	ASP	2.1
1	A	91	ARG	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($\text{\AA}^2$)	Q<0.9
2	NA	A	8003	1/1	0.96	0.07	14,14,14,14	0
2	NA	B	8002	1/1	0.98	0.03	16,16,16,16	0
2	NA	A	8001	1/1	0.99	0.05	16,16,16,16	0

6.5 Other polymers [i](#)

There are no such residues in this entry.