



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 06:24 AM UTC

PDB ID : 3DZB / pdb_00003dzb
Title : Crystal structure of Prephenate dehydrogenase from *Streptococcus thermophilus*
Authors : Zhang, Z.; Eswaramoorthy, S.; Burley, S.K.; Swaminathan, S.; New York SGX Research Center for Structural Genomics (NYSGXRC)
Deposited on : 2008-07-29
Resolution : 2.46 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

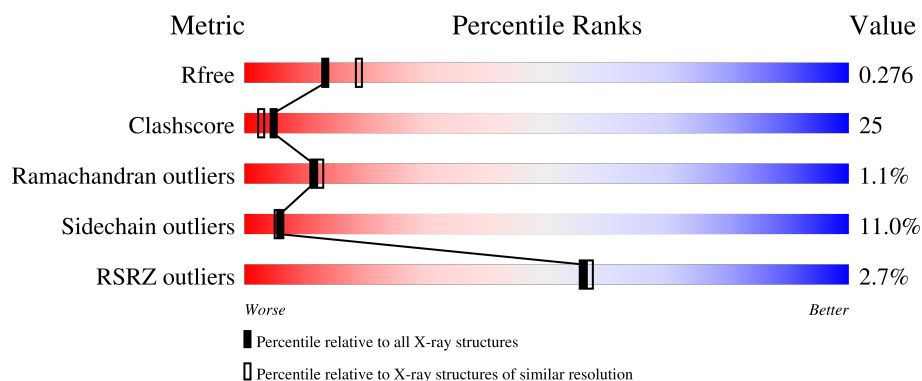
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.46 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	1190 (2.46-2.46)
Clashscore	190562	1229 (2.46-2.46)
Ramachandran outliers	187476	1218 (2.46-2.46)
Sidechain outliers	187428	1218 (2.46-2.46)
RSRZ outliers	180081	1190 (2.46-2.46)

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	317	
1	B	317	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4499 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 Prephenate dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	289	Total	C	N	O	Se	0	0	0
			2260	1432	386	434	8			
1	B	278	Total	C	N	O	Se	0	0	0
			2179	1380	373	418	8			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	MSE	-	expression tag	UNP Q5M554
A	0	SER	-	expression tag	UNP Q5M554
A	1	LEU	-	expression tag	UNP Q5M554
A	309	GLY	-	expression tag	UNP Q5M554
A	310	HIS	-	expression tag	UNP Q5M554
A	311	HIS	-	expression tag	UNP Q5M554
A	312	HIS	-	expression tag	UNP Q5M554
A	313	HIS	-	expression tag	UNP Q5M554
A	314	HIS	-	expression tag	UNP Q5M554
A	315	HIS	-	expression tag	UNP Q5M554
B	-1	MSE	-	expression tag	UNP Q5M554
B	0	SER	-	expression tag	UNP Q5M554
B	1	LEU	-	expression tag	UNP Q5M554
B	309	GLY	-	expression tag	UNP Q5M554
B	310	HIS	-	expression tag	UNP Q5M554
B	311	HIS	-	expression tag	UNP Q5M554
B	312	HIS	-	expression tag	UNP Q5M554
B	313	HIS	-	expression tag	UNP Q5M554
B	314	HIS	-	expression tag	UNP Q5M554
B	315	HIS	-	expression tag	UNP Q5M554

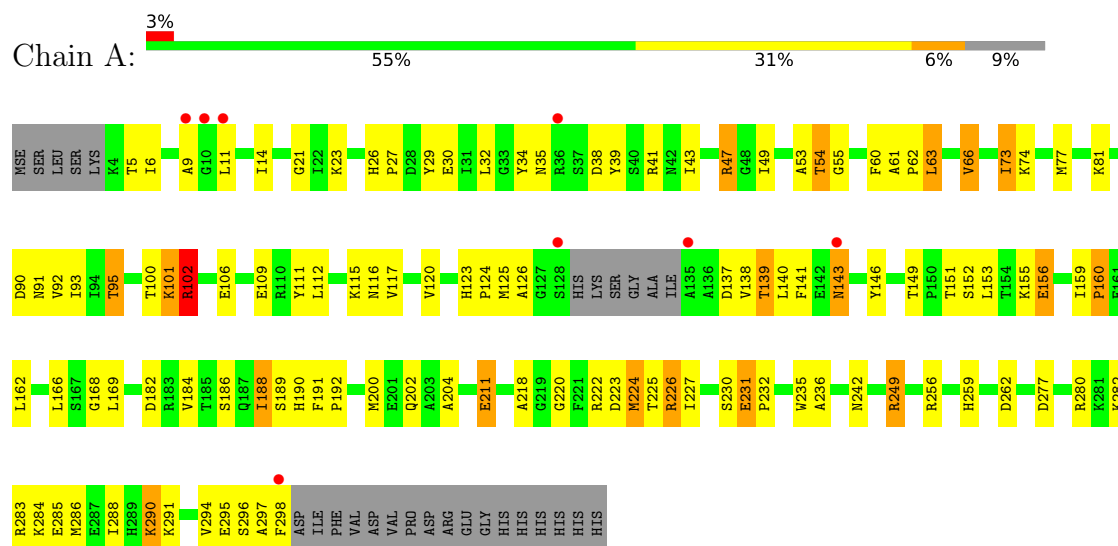
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	27	Total 27	O 27	0	0
2	B	33	Total 33	O 33	0	0

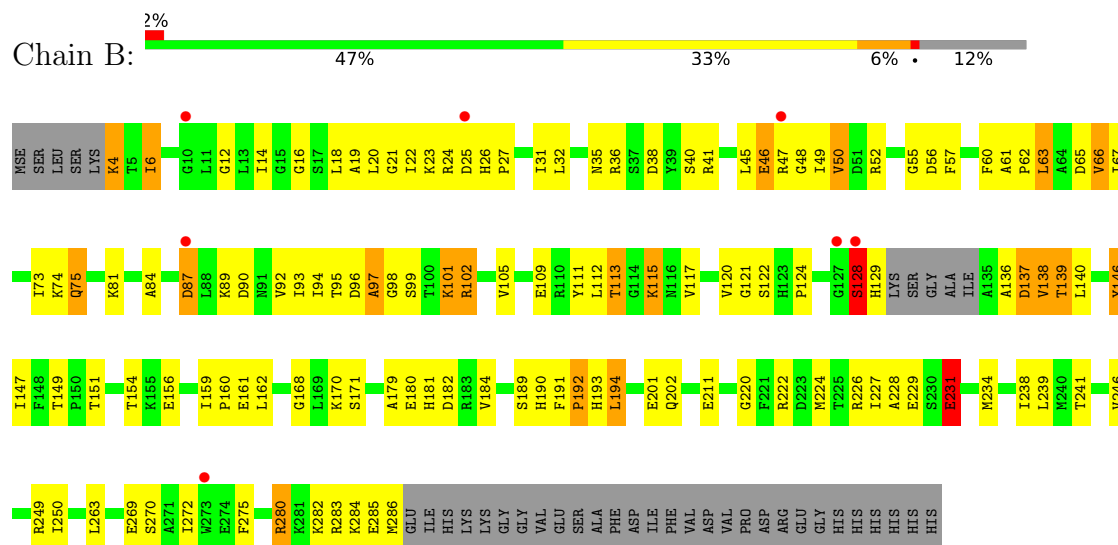
3 Residue-property plots [i](#)

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: Prephenate dehydrogenase



• Molecule 1: Prephenate dehydrogenase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	83.91Å 145.00Å 56.23Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.46 50.00 – 2.47	Depositor EDS
% Data completeness (in resolution range)	95.4 (50.00-2.46) 95.5 (50.00-2.47)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.73 (at 2.48Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.234 , 0.278 0.233 , 0.276	Depositor DCC
R_{free} test set	935 reflections (3.68%)	wwPDB-VP
Wilson B-factor (Å ²)	37.7	Xtriage
Anisotropy	0.307	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 30.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	4499	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.43% 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	0.52	0/2294	0.99	10/3082 (0.3%)
1	B	0.48	0/2212	1.00	11/2974 (0.4%)
All	All	0.50	0/4506	1.00	21/6056 (0.3%)

There are no bond length outliers.

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	115	LYS	N-CA-C	9.20	123.20	112.72
1	B	231	GLU	CA-C-N	7.45	126.72	118.97
1	B	231	GLU	C-N-CA	7.45	126.72	118.97
1	B	146	TYR	N-CA-C	-6.19	97.18	108.02
1	A	39	TYR	N-CA-C	6.16	117.69	110.97
1	B	128	SER	N-CA-C	6.14	120.32	112.34
1	B	12	GLY	N-CA-C	-6.03	103.98	112.60
1	B	147	ILE	N-CA-C	5.96	116.88	108.36
1	A	102	ARG	N-CA-C	5.95	118.27	111.02
1	A	112	LEU	N-CA-C	5.72	120.10	113.18
1	A	211	GLU	N-CA-C	5.69	118.25	111.71
1	A	100	THR	N-CA-C	-5.50	101.16	109.85
1	A	188	ILE	N-CA-C	5.41	117.37	112.29
1	A	73	ILE	N-CA-C	5.23	115.86	110.36
1	A	236	ALA	N-CA-C	-5.17	105.72	111.36
1	B	269	GLU	N-CA-C	5.17	116.92	111.28
1	B	14	ILE	N-CA-C	5.13	115.35	110.42
1	B	228	ALA	N-CA-C	-5.06	106.36	112.54
1	B	211	GLU	N-CA-C	5.03	118.88	112.34
1	A	231	GLU	CA-C-N	5.02	124.46	119.24
1	A	231	GLU	C-N-CA	5.02	124.46	119.24

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	2260	0	2241	131	0
1	B	2179	0	2158	108	0
2	A	27	0	0	3	0
2	B	33	0	0	2	0
All	All	4499	0	4399	224	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 25.

All (224) 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:226:ARG:HG2	1:A:226:ARG:NH1	1.65	1.04
1:A:151:THR:HG22	1:A:153:LEU:H	1.23	1.03
1:A:290:LYS:HE3	1:A:290:LYS:HA	1.40	1.02
1:A:226:ARG:HG2	1:A:226:ARG:HH11	0.85	0.99
1:A:226:ARG:HH11	1:A:226:ARG:CG	1.76	0.97
1:A:235:TRP:HE1	1:A:298:PHE:HE2	1.09	0.97
1:B:280:ARG:HH11	1:B:280:ARG:HB3	1.28	0.97
1:B:101:LYS:HG2	1:B:149:THR:HG21	1.47	0.96
1:A:11:LEU:HG	1:A:35:ASN:HB2	1.47	0.94
1:B:151:THR:O	1:B:154:THR:HG22	1.67	0.93
1:A:101:LYS:HG2	1:A:149:THR:HG21	1.51	0.91
1:A:226:ARG:HG3	1:A:227:ILE:N	1.85	0.91
1:A:95:THR:HG22	1:A:120:VAL:O	1.75	0.87
1:A:101:LYS:HE3	1:A:101:LYS:HA	1.58	0.86
1:A:200:MSE:HE2	1:A:225:THR:HB	1.58	0.85
1:A:11:LEU:CG	1:A:35:ASN:HB2	2.06	0.84
1:A:137:ASP:CG	1:A:139:THR:HG22	2.02	0.84
1:B:102:ARG:HB2	1:B:182:ASP:OD2	1.78	0.83
1:A:6:ILE:HD13	1:A:66:VAL:CG1	2.10	0.82
1:B:101:LYS:HE3	1:B:101:LYS:HA	1.62	0.82
1:A:5:THR:HG21	1:A:63:LEU:O	1.79	0.81
1:A:11:LEU:HD12	1:A:35:ASN:CB	2.13	0.79
1:B:280:ARG:HH11	1:B:280:ARG:CB	1.96	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:21:GLY:HA2	1:B:138:VAL:HG23	1.68	0.75
1:A:11:LEU:CD1	1:A:35:ASN:HB2	2.18	0.74
1:A:235:TRP:NE1	1:A:298:PHE:HE2	1.85	0.74
1:A:43:ILE:O	1:A:47:ARG:HB2	1.87	0.73
1:B:31:ILE:HG22	1:B:50:VAL:HG12	1.71	0.73
1:A:286:MSE:HE2	1:A:288:ILE:HD11	1.71	0.72
1:A:137:ASP:OD1	1:A:139:THR:HG22	1.89	0.71
1:B:6:ILE:HD13	1:B:66:VAL:CG1	2.20	0.70
1:B:101:LYS:HG2	1:B:149:THR:CG2	2.22	0.70
1:A:256:ARG:NH1	1:B:201:GLU:HG2	2.06	0.70
1:A:11:LEU:CD1	1:A:35:ASN:CB	2.69	0.69
1:A:186:SER:HA	1:A:190:HIS:HB2	1.74	0.69
1:A:11:LEU:HB2	1:A:35:ASN:ND2	2.07	0.69
1:A:102:ARG:HB2	1:A:182:ASP:OD2	1.91	0.69
1:A:6:ILE:HD13	1:A:66:VAL:HG11	1.74	0.68
1:B:194:LEU:HD12	1:B:194:LEU:O	1.94	0.67
1:B:222:ARG:HG2	1:B:222:ARG:HH21	1.59	0.67
1:A:11:LEU:HB2	1:A:35:ASN:HD22	1.59	0.67
1:A:101:LYS:HG2	1:A:149:THR:CG2	2.24	0.67
1:B:96:ASP:OD1	1:B:101:LYS:NZ	2.28	0.67
1:A:200:MSE:HE3	1:B:283:ARG:CD	2.24	0.66
1:A:61:ALA:HB3	1:A:62:PRO:HD3	1.78	0.66
1:B:280:ARG:HB3	1:B:280:ARG:NH1	2.07	0.66
1:A:226:ARG:NH1	1:A:226:ARG:CG	2.45	0.65
1:A:200:MSE:HE2	1:A:225:THR:CB	2.27	0.65
1:B:46:GLU:C	1:B:48:GLY:H	2.03	0.65
1:A:92:VAL:HG12	1:A:117:VAL:HG12	1.77	0.65
1:B:96:ASP:CG	1:B:97:ALA:H	2.06	0.63
1:A:26:HIS:HB3	1:A:29:TYR:CD1	2.33	0.63
1:A:90:ASP:OD1	1:A:116:ASN:ND2	2.32	0.63
1:A:11:LEU:HD12	1:A:35:ASN:HB3	1.81	0.63
1:B:61:ALA:HB3	1:B:62:PRO:HD3	1.80	0.63
1:A:290:LYS:HA	1:A:290:LYS:CE	2.25	0.62
1:A:218:ALA:O	1:A:222:ARG:HD2	1.99	0.62
1:A:60:PHE:O	1:A:63:LEU:HD22	1.99	0.62
1:A:200:MSE:HE3	1:B:283:ARG:HD2	1.81	0.61
1:B:220:GLY:O	1:B:224:MSE:HG2	2.01	0.61
1:B:6:ILE:HD13	1:B:66:VAL:HG11	1.83	0.60
1:B:23:LYS:NZ	1:B:49:ILE:O	2.24	0.59
1:A:126:ALA:HB3	1:A:294:VAL:HB	1.85	0.58
1:A:288:ILE:O	1:A:291:LYS:HB2	2.03	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:46:GLU:C	1:B:48:GLY:N	2.61	0.58
1:A:109:GLU:OE2	1:A:151:THR:HG23	2.03	0.58
1:B:193:HIS:CE1	1:B:227:ILE:HB	2.38	0.58
1:A:141:PHE:CD2	1:A:169:LEU:HD22	2.38	0.58
1:B:66:VAL:HA	1:B:93:ILE:O	2.03	0.58
1:A:224:MSE:HE3	1:B:224:MSE:O	2.04	0.58
1:B:84:ALA:O	1:B:115:LYS:HE3	2.03	0.58
1:B:231:GLU:O	1:B:231:GLU:HG3	2.03	0.58
1:A:249:ARG:HD3	2:A:324:HOH:O	2.02	0.57
1:B:65:ASP:O	1:B:92:VAL:HG23	2.04	0.57
1:A:125:MSE:HE3	1:A:295:GLU:HA	1.85	0.57
1:A:156:GLU:CD	1:A:156:GLU:H	2.12	0.57
1:A:182:ASP:HB3	1:A:242:ASN:OD1	2.05	0.56
1:A:137:ASP:C	1:A:139:THR:H	2.14	0.55
1:A:189:SER:C	1:A:192:PRO:HD2	2.31	0.55
1:A:102:ARG:HG3	1:A:102:ARG:HH11	1.72	0.55
1:A:91:ASN:H	1:A:116:ASN:ND2	2.05	0.55
1:A:137:ASP:OD2	1:A:139:THR:HB	2.07	0.55
1:B:35:ASN:ND2	1:B:40:SER:OG	2.40	0.55
1:B:46:GLU:O	1:B:48:GLY:N	2.40	0.55
1:A:277:ASP:OD1	1:A:280:ARG:NH1	2.39	0.55
1:B:113:THR:HG22	1:B:113:THR:O	2.06	0.55
1:A:91:ASN:HD21	1:A:155:LYS:NZ	2.05	0.54
1:B:4:LYS:NZ	1:B:4:LYS:HB2	2.22	0.54
1:A:6:ILE:CD1	1:A:66:VAL:HG11	2.37	0.54
1:A:14:ILE:HD12	1:A:14:ILE:N	2.21	0.54
1:B:96:ASP:CG	1:B:97:ALA:N	2.64	0.54
1:B:65:ASP:C	1:B:92:VAL:HG23	2.33	0.54
1:A:141:PHE:HD2	1:A:169:LEU:HD22	1.72	0.54
1:B:22:ILE:O	1:B:26:HIS:HB2	2.08	0.54
1:B:38:ASP:HA	1:B:41:ARG:NH1	2.23	0.53
1:B:234:MSE:O	1:B:238:ILE:HG13	2.09	0.53
1:B:93:ILE:HG21	1:B:162:LEU:HD21	1.90	0.53
1:B:75:GLN:CA	1:B:75:GLN:HE21	2.20	0.53
1:A:191:PHE:HB3	1:A:192:PRO:HD3	1.90	0.53
1:A:137:ASP:OD2	1:A:139:THR:CB	2.56	0.53
1:B:227:ILE:HG12	2:B:344:HOH:O	2.09	0.52
1:A:5:THR:CG2	1:A:63:LEU:O	2.54	0.52
1:B:16:GLY:HA3	1:B:136:ALA:HB2	1.91	0.52
1:B:112:LEU:HB3	1:B:117:VAL:HG21	1.90	0.52
1:A:137:ASP:OD1	1:A:139:THR:CG2	2.58	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:200:MSE:CE	1:B:283:ARG:HD2	2.39	0.52
1:B:4:LYS:HB2	1:B:4:LYS:HZ3	1.74	0.52
1:A:21:GLY:O	1:A:168:GLY:HA3	2.10	0.52
1:B:102:ARG:HG2	1:B:179:ALA:HA	1.92	0.52
1:A:32:LEU:CD1	1:A:63:LEU:HD21	2.40	0.52
1:A:41:ARG:HG2	1:A:53:ALA:HB1	1.92	0.51
1:B:60:PHE:O	1:B:63:LEU:HD22	2.09	0.51
1:B:105:VAL:O	1:B:109:GLU:HG3	2.10	0.51
1:B:191:PHE:HB3	1:B:192:PRO:HD3	1.91	0.51
1:A:54:THR:OG1	1:A:55:GLY:N	2.41	0.51
1:A:91:ASN:ND2	1:A:155:LYS:NZ	2.58	0.51
1:B:21:GLY:CA	1:B:138:VAL:HG23	2.38	0.51
1:B:246:VAL:O	1:B:250:ILE:HG13	2.11	0.51
1:A:226:ARG:CG	1:A:227:ILE:N	2.65	0.51
1:B:285:GLU:O	1:B:286:MSE:HB2	2.10	0.51
1:A:21:GLY:N	1:A:138:VAL:HG23	2.25	0.51
1:A:151:THR:CG2	1:A:152:SER:N	2.74	0.51
1:A:91:ASN:ND2	1:A:155:LYS:HZ1	2.08	0.51
1:A:224:MSE:CE	1:B:224:MSE:O	2.59	0.51
1:B:120:VAL:HA	1:B:151:THR:HG23	1.93	0.50
1:A:282:LYS:HA	1:A:285:GLU:HG2	1.94	0.50
2:A:327:HOH:O	1:B:74:LYS:HE2	2.11	0.50
1:B:6:ILE:HG23	1:B:66:VAL:HG13	1.93	0.50
1:A:9:ALA:O	1:A:34:TYR:O	2.30	0.50
1:A:159:ILE:HB	1:A:160:PRO:HD3	1.94	0.50
1:A:137:ASP:OD2	1:A:139:THR:HG22	2.12	0.50
1:B:75:GLN:HE21	1:B:75:GLN:C	2.19	0.50
1:A:200:MSE:HE3	1:B:283:ARG:HB2	1.93	0.50
1:A:41:ARG:HD3	1:A:54:THR:O	2.11	0.49
1:A:115:LYS:CG	1:A:116:ASN:N	2.76	0.49
1:B:20:LEU:HB3	1:B:138:VAL:HB	1.94	0.49
1:B:284:LYS:C	1:B:286:MSE:H	2.19	0.49
1:B:38:ASP:HA	1:B:41:ARG:HH12	1.78	0.49
1:B:63:LEU:HA	1:B:89:LYS:HE3	1.95	0.49
1:A:11:LEU:HD11	1:A:41:ARG:HG3	1.95	0.48
1:A:200:MSE:CE	1:B:283:ARG:CD	2.92	0.48
1:A:226:ARG:NH1	1:A:227:ILE:HG23	2.27	0.48
1:B:18:LEU:O	1:B:22:ILE:HG13	2.14	0.48
1:B:284:LYS:C	1:B:286:MSE:N	2.69	0.48
1:A:137:ASP:OD2	1:A:139:THR:N	2.41	0.48
1:B:75:GLN:CA	1:B:75:GLN:NE2	2.76	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:32:LEU:CD2	1:B:52:ARG:HB2	2.44	0.47
1:A:222:ARG:HA	1:B:283:ARG:HH12	1.79	0.47
1:B:121:GLY:HA3	1:B:149:THR:OG1	2.14	0.47
1:A:38:ASP:HA	1:A:41:ARG:CZ	2.44	0.47
1:A:137:ASP:C	1:A:139:THR:N	2.72	0.47
1:B:73:ILE:HG21	1:B:241:THR:HG21	1.96	0.47
1:A:11:LEU:CD1	1:A:35:ASN:HB3	2.41	0.47
1:A:32:LEU:HD12	1:A:63:LEU:HD21	1.97	0.46
1:B:24:ARG:NH1	1:B:170:LYS:HD2	2.30	0.46
1:B:137:ASP:OD2	1:B:139:THR:HB	2.15	0.46
1:B:189:SER:C	1:B:192:PRO:HD2	2.40	0.46
1:B:63:LEU:C	1:B:63:LEU:CD2	2.88	0.46
1:B:109:GLU:O	1:B:113:THR:HB	2.16	0.46
1:A:73:ILE:HG22	1:A:77:MSE:HE2	1.96	0.46
1:A:102:ARG:HG3	1:A:102:ARG:NH1	2.31	0.46
1:A:188:ILE:HD12	1:A:188:ILE:C	2.41	0.46
1:A:200:MSE:HE3	1:B:283:ARG:CB	2.45	0.45
1:A:220:GLY:O	1:A:224:MSE:HG3	2.16	0.45
1:B:194:LEU:HD12	1:B:194:LEU:C	2.41	0.45
1:B:87:ASP:CG	1:B:87:ASP:O	2.60	0.45
1:A:26:HIS:N	1:A:27:PRO:HD3	2.32	0.45
1:A:184:VAL:O	1:A:188:ILE:HG13	2.17	0.45
1:A:137:ASP:CG	1:A:139:THR:CG2	2.84	0.45
1:B:180:GLU:O	1:B:184:VAL:HG23	2.17	0.45
1:A:211:GLU:HB2	2:A:326:HOH:O	2.16	0.45
1:B:35:ASN:ND2	1:B:40:SER:HG	2.14	0.45
1:A:223:ASP:OD2	1:B:226:ARG:NH1	2.51	0.44
1:A:249:ARG:HH21	1:B:202:GLN:HG2	1.81	0.44
1:B:99:SER:O	1:B:181:HIS:HE1	2.01	0.44
1:A:115:LYS:CG	1:A:116:ASN:H	2.31	0.43
1:B:272:ILE:O	1:B:275:PHE:HB3	2.17	0.43
1:B:60:PHE:HA	1:B:63:LEU:HD11	2.01	0.43
1:B:124:PRO:HB3	1:B:146:TYR:CE1	2.53	0.43
1:B:67:ILE:HD12	1:B:94:ILE:HG12	2.00	0.43
1:A:21:GLY:HA2	1:A:138:VAL:HG23	2.00	0.43
1:A:102:ARG:NE	1:A:106:GLU:OE1	2.52	0.43
1:B:159:ILE:HB	1:B:160:PRO:HD3	2.00	0.43
1:A:66:VAL:HA	1:A:93:ILE:O	2.18	0.43
1:A:95:THR:CG2	1:A:120:VAL:O	2.58	0.43
1:A:81:LYS:HB2	1:A:111:TYR:CE1	2.54	0.42
1:B:81:LYS:HB2	1:B:111:TYR:CE1	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:137:ASP:C	1:A:137:ASP:OD2	2.60	0.42
1:B:6:ILE:HD13	1:B:66:VAL:HG13	1.99	0.42
1:B:138:VAL:O	1:B:138:VAL:HG22	2.19	0.42
1:A:146:TYR:CZ	1:A:166:LEU:HD13	2.54	0.42
1:B:191:PHE:N	1:B:192:PRO:CD	2.82	0.42
1:B:26:HIS:N	1:B:27:PRO:CD	2.83	0.42
1:A:21:GLY:CA	1:A:138:VAL:HG23	2.49	0.42
1:A:137:ASP:OD2	1:A:139:THR:CG2	2.68	0.42
1:B:23:LYS:O	1:B:24:ARG:C	2.63	0.42
1:A:143:ASN:O	1:A:143:ASN:CG	2.63	0.42
1:A:124:PRO:HB3	1:A:146:TYR:CD1	2.54	0.42
1:B:81:LYS:HG3	1:B:111:TYR:CD2	2.55	0.42
1:A:11:LEU:HG	1:A:34:TYR:O	2.20	0.41
1:A:32:LEU:HD12	1:A:63:LEU:CD2	2.50	0.41
1:A:38:ASP:HA	1:A:41:ARG:NH2	2.35	0.41
1:A:231:GLU:HA	1:A:232:PRO:HD3	1.93	0.41
1:B:190:HIS:CD2	1:B:190:HIS:N	2.87	0.41
1:A:14:ILE:HD11	1:A:297:ALA:O	2.20	0.41
1:A:191:PHE:N	1:A:192:PRO:CD	2.83	0.41
1:B:98:GLY:O	1:B:101:LYS:NZ	2.53	0.41
1:A:295:GLU:OE2	1:A:296:SER:OG	2.36	0.41
1:A:151:THR:HG22	1:A:152:SER:N	2.36	0.41
1:B:128:SER:O	1:B:129:HIS:C	2.64	0.41
1:A:202:GLN:CG	1:B:249:ARG:HH21	2.34	0.40
1:A:259:HIS:O	1:A:262:ASP:HB2	2.21	0.40
1:B:170:LYS:HE3	2:B:339:HOH:O	2.20	0.40
1:A:5:THR:HG21	1:A:63:LEU:HD23	2.03	0.40
1:A:91:ASN:N	1:A:116:ASN:ND2	2.69	0.40
1:A:115:LYS:HG2	1:A:116:ASN:N	2.36	0.40
1:A:123:HIS:CD2	1:A:123:HIS:C	2.99	0.40
1:A:204:ALA:HB1	1:B:282:LYS:HD2	2.02	0.40
1:B:16:GLY:O	1:B:19:ALA:HB3	2.22	0.40
1:B:41:ARG:HH21	1:B:55:GLY:HA3	1.87	0.40
1:B:56:ASP:O	1:B:57:PHE:C	2.63	0.40
1:A:283:ARG:HB3	1:B:229:GLU:CG	2.51	0.40
1:B:21:GLY:O	1:B:168:GLY:HA3	2.22	0.40
1:B:63:LEU:HD22	1:B:63:LEU:C	2.47	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	285/317 (90%)	260 (91%)	25 (9%)	0	100	100
1	B	274/317 (86%)	248 (90%)	20 (7%)	6 (2%)	5	3
All	All	559/634 (88%)	508 (91%)	45 (8%)	6 (1%)	11	12

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	47	ARG
1	B	97	ALA
1	B	156	GLU
1	B	102	ARG
1	B	128	SER
1	B	138	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	236/252 (94%)	213 (90%)	23 (10%)	8	8
1	B	228/252 (90%)	200 (88%)	28 (12%)	4	4
All	All	464/504 (92%)	413 (89%)	51 (11%)	6	5

All (51) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	23	LYS
1	A	30	GLU
1	A	47	ARG
1	A	49	ILE
1	A	54	THR
1	A	63	LEU
1	A	66	VAL
1	A	74	LYS
1	A	95	THR
1	A	101	LYS
1	A	102	ARG
1	A	139	THR
1	A	140	LEU
1	A	143	ASN
1	A	156	GLU
1	A	160	PRO
1	A	162	LEU
1	A	224	MSE
1	A	226	ARG
1	A	230	SER
1	A	249	ARG
1	A	284	LYS
1	A	290	LYS
1	B	4	LYS
1	B	6	ILE
1	B	25	ASP
1	B	36	ARG
1	B	45	LEU
1	B	46	GLU
1	B	50	VAL
1	B	63	LEU
1	B	66	VAL
1	B	75	GLN
1	B	87	ASP
1	B	90	ASP
1	B	95	THR
1	B	101	LYS
1	B	113	THR
1	B	122	SER
1	B	137	ASP
1	B	139	THR
1	B	140	LEU
1	B	161	GLU

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Mol	Chain	Res	Type
1	B	171	SER
1	B	192	PRO
1	B	194	LEU
1	B	231	GLU
1	B	239	LEU
1	B	263	LEU
1	B	270	SER
1	B	280	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	91	ASN
1	A	116	ASN
1	A	118	GLN
1	A	208	GLN
1	A	214	ASN
1	B	35	ASN
1	B	75	GLN
1	B	193	HIS
1	B	214	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 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 [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	281/317 (88%)	0.22	8 (2%) 55 56	18, 38, 55, 61	0
1	B	270/317 (85%)	0.20	7 (2%) 57 58	18, 38, 51, 55	0
All	All	551/634 (86%)	0.21	15 (2%) 56 57	18, 38, 54, 61	0

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	11	LEU	4.3
1	A	9	ALA	3.8
1	A	10	GLY	3.4
1	A	298	PHE	3.1
1	B	127	GLY	3.1
1	A	128	SER	2.7
1	B	128	SER	2.5
1	A	135	ALA	2.5
1	A	143	ASN	2.5
1	B	87	ASP	2.4
1	B	10	GLY	2.4
1	B	25	ASP	2.2
1	B	47	ARG	2.2
1	B	273	TRP	2.1
1	A	36	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.