



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 01:55 PM UTC

PDB ID : 3NRO / pdb_00003nro
Title : Crystal Structure of putative transcriptional factor Lmo1026 from *Listeria monocytogenes* (FRAGMENT 52-321), Northeast Structural Genomics Consortium Target LmR194
Authors : Kuzin, A.; Su, M.; Seetharaman, J.; Mao, M.; Xiao, R.; Ciccocanti, C.; Lee, D.; Everett, J.K.; Nair, R.; Acton, T.B.; Rost, B.; Montelione, G.T.; Hunt, J.F.; Tong, L.; Northeast Structural Genomics Consortium (NESG)
Deposited on : 2010-06-30
Resolution : 2.90 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)

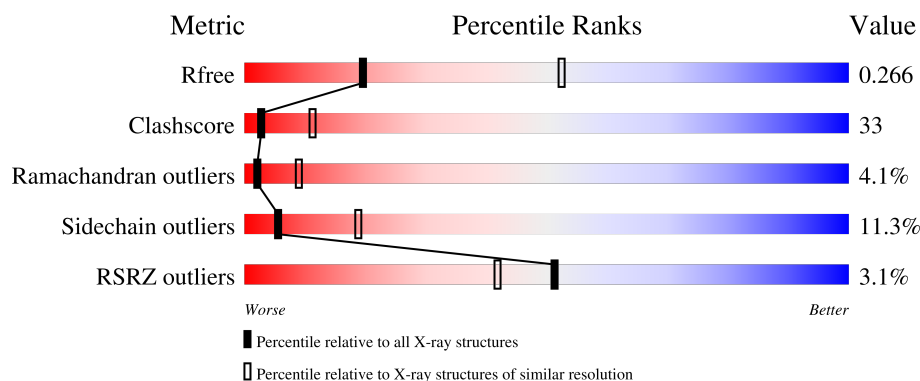
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (2.90-2.90)

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	279	4% 42% 34% 7% 16%
1	B	279	% 39% 39% 6% 15%

2 Entry composition [i](#)

There is only 1 type of molecule in this entry. The entry contains 3647 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 Lmo1026 protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	235	Total	C	N	O	Se	0	0	0
			1814	1139	303	363	9			
1	B	237	Total	C	N	O	Se	0	0	0
			1833	1152	304	368	9			

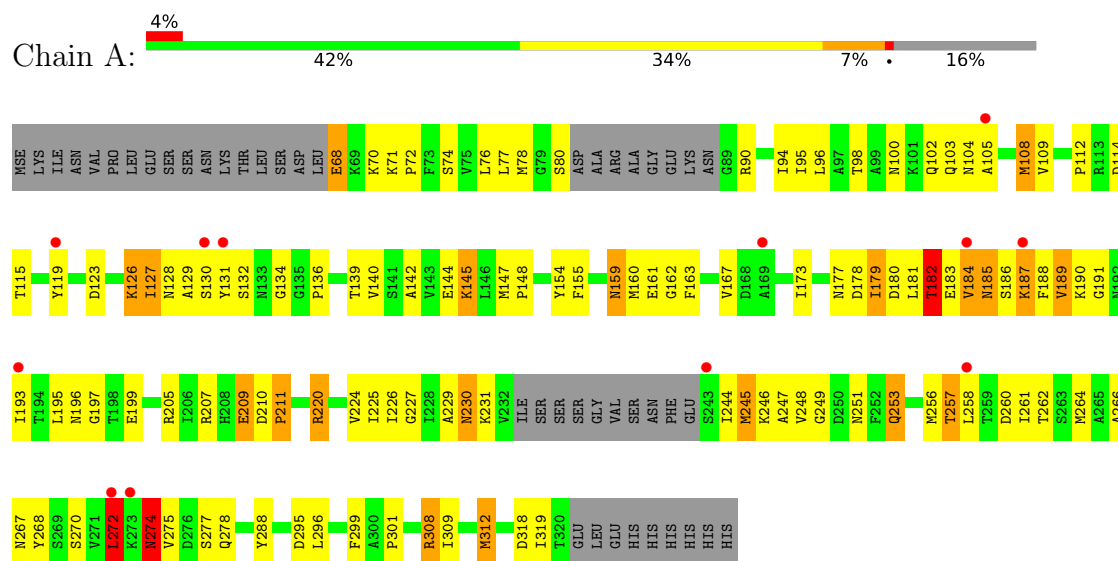
There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	51	MSE	-	expression tag	UNP Q8Y889
A	322	LEU	-	expression tag	UNP Q8Y889
A	323	GLU	-	expression tag	UNP Q8Y889
A	324	HIS	-	expression tag	UNP Q8Y889
A	325	HIS	-	expression tag	UNP Q8Y889
A	326	HIS	-	expression tag	UNP Q8Y889
A	327	HIS	-	expression tag	UNP Q8Y889
A	328	HIS	-	expression tag	UNP Q8Y889
A	329	HIS	-	expression tag	UNP Q8Y889
B	51	MSE	-	expression tag	UNP Q8Y889
B	322	LEU	-	expression tag	UNP Q8Y889
B	323	GLU	-	expression tag	UNP Q8Y889
B	324	HIS	-	expression tag	UNP Q8Y889
B	325	HIS	-	expression tag	UNP Q8Y889
B	326	HIS	-	expression tag	UNP Q8Y889
B	327	HIS	-	expression tag	UNP Q8Y889
B	328	HIS	-	expression tag	UNP Q8Y889
B	329	HIS	-	expression tag	UNP Q8Y889

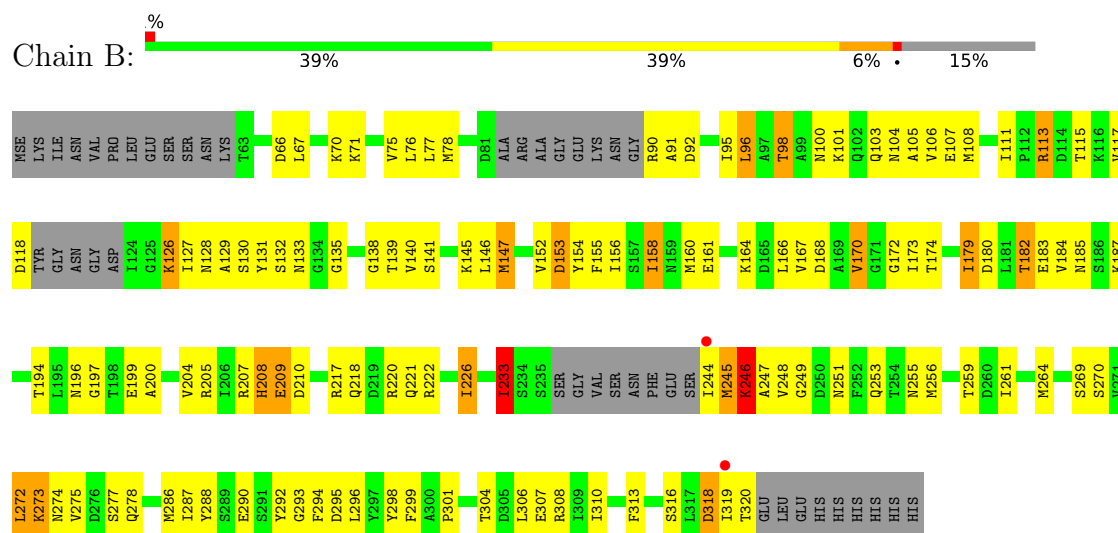
3 Residue-property plots

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: Lmo1026 protein



• Molecule 1: Lmo1026 protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	44.24Å 90.30Å 118.04Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.66 – 2.90 29.66 – 2.90	Depositor EDS
% Data completeness (in resolution range)	99.5 (29.66-2.90) 99.4 (29.66-2.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.28 (at 2.90Å)	Xtriage
Refinement program	PHENIX 1.6.1 _357	Depositor
R, R_{free}	0.234 , 0.264 0.233 , 0.266	Depositor DCC
R_{free} test set	526 reflections (4.80%)	wwPDB-VP
Wilson B-factor (Å ²)	66.3	Xtriage
Anisotropy	0.496	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 74.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	3647	wwPDB-VP
Average B, all atoms (Å ²)	70.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.37% 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](#)

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	0.55	0/1829	0.96	2/2449 (0.1%)
1	B	0.60	0/1846	1.00	6/2472 (0.2%)
All	All	0.58	0/3675	0.98	8/4921 (0.2%)

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	247	ALA	N-CA-C	-13.28	97.56	112.57
1	A	247	ALA	N-CA-C	-11.94	97.19	112.23
1	B	255	ASN	N-CA-C	-7.02	104.03	112.59
1	B	147	MSE	CA-C-N	6.53	126.76	119.32
1	B	147	MSE	C-N-CA	6.53	126.76	119.32
1	B	269	SER	N-CA-C	-6.09	103.08	110.88
1	B	246	LYS	N-CA-C	5.70	122.93	110.80
1	A	245	MSE	N-CA-C	5.67	117.72	108.76

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	1814	0	1804	128	0
1	B	1833	0	1833	115	0
All	All	3647	0	3637	242	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 33.

All (242) 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:187:LYS:H	1:A:187:LYS:HD3	1.13	1.10
1:B:108:MSE:HE1	1:B:222:ARG:HG3	1.37	1.03
1:A:244:ILE:HD12	1:A:244:ILE:O	1.67	0.95
1:A:159:ASN:H	1:A:159:ASN:HD22	1.19	0.90
1:B:115:THR:HA	1:B:299:PHE:HB3	1.55	0.88
1:A:159:ASN:HD21	1:A:251:ASN:HB3	1.38	0.86
1:A:187:LYS:HD3	1:A:187:LYS:N	1.90	0.86
1:A:188:PHE:HA	1:A:193:ILE:HD13	1.61	0.82
1:A:167:VAL:HG13	1:A:173:ILE:HG12	1.62	0.82
1:A:159:ASN:HD21	1:A:162:GLY:H	1.28	0.81
1:B:273:LYS:HE2	1:B:273:LYS:HA	1.61	0.80
1:A:90:ARG:HD2	1:A:128:ASN:HD22	1.46	0.79
1:A:159:ASN:ND2	1:A:162:GLY:H	1.80	0.79
1:A:185:ASN:H	1:A:185:ASN:ND2	1.82	0.77
1:A:185:ASN:HD22	1:A:186:SER:H	1.32	0.77
1:B:160:MSE:HG3	1:B:205:ARG:HH21	1.50	0.76
1:B:113:ARG:H	1:B:113:ARG:HE	1.35	0.75
1:A:100:ASN:HD21	1:A:102:GLN:HB2	1.51	0.74
1:B:286:MSE:HG3	1:B:295:ASP:HB3	1.68	0.74
1:B:75:VAL:HG22	1:B:154:TYR:HB2	1.70	0.74
1:B:319:ILE:HG12	1:B:320:THR:H	1.52	0.73
1:A:179:ILE:HD13	1:A:180:ASP:N	2.04	0.73
1:A:105:ALA:HA	1:A:274:ASN:HB3	1.70	0.72
1:A:71:LYS:HE3	1:A:72:PRO:O	1.89	0.72
1:B:96:LEU:HD12	1:B:313:PHE:CE2	2.25	0.72
1:A:72:PRO:HG3	1:A:100:ASN:ND2	2.06	0.71
1:A:185:ASN:H	1:A:185:ASN:HD22	1.37	0.70
1:A:90:ARG:HD2	1:A:128:ASN:ND2	2.06	0.70
1:A:188:PHE:HA	1:A:193:ILE:CD1	2.20	0.70
1:B:98:THR:HG21	1:B:316:SER:HB3	1.73	0.70
1:B:208:HIS:O	1:B:210:ASP:N	2.25	0.69
1:A:126:LYS:O	1:A:130:SER:HB2	1.92	0.69
1:B:113:ARG:O	1:B:128:ASN:HB3	1.93	0.69
1:A:248:VAL:HG22	1:A:249:GLY:H	1.55	0.69
1:B:179:ILE:HD12	1:B:180:ASP:H	1.57	0.68
1:A:187:LYS:H	1:A:187:LYS:CD	2.00	0.68
1:A:128:ASN:CG	1:A:129:ALA:N	2.52	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:167:VAL:HB	1:A:197:GLY:HA2	1.76	0.68
1:A:264:MSE:HA	1:A:268:TYR:HD2	1.58	0.68
1:A:159:ASN:ND2	1:A:251:ASN:HB3	2.07	0.68
1:B:244:ILE:HG13	1:B:245:MSE:HG3	1.77	0.67
1:A:182:THR:C	1:A:184:VAL:H	2.01	0.66
1:B:272:LEU:HD13	1:B:273:LYS:H	1.60	0.66
1:A:246:LYS:HD2	1:A:246:LYS:N	2.11	0.66
1:A:257:THR:O	1:A:261:ILE:HG13	1.95	0.66
1:B:115:THR:HG23	1:B:301:PRO:HG3	1.77	0.65
1:A:72:PRO:HG3	1:A:100:ASN:HD22	1.61	0.65
1:A:132:SER:HB2	1:A:139:THR:OG1	1.97	0.64
1:A:180:ASP:OD2	1:A:182:THR:HG22	1.97	0.64
1:A:159:ASN:HD22	1:A:159:ASN:N	1.84	0.64
1:B:179:ILE:HD12	1:B:180:ASP:N	2.12	0.64
1:A:76:LEU:HD23	1:A:140:VAL:HG22	1.79	0.64
1:A:159:ASN:H	1:A:159:ASN:ND2	1.94	0.64
1:B:130:SER:OG	1:B:139:THR:HA	1.98	0.64
1:B:104:ASN:O	1:B:274:ASN:HB2	1.98	0.63
1:B:182:THR:C	1:B:184:VAL:H	2.04	0.63
1:A:145:LYS:NZ	1:A:145:LYS:HB3	2.12	0.63
1:A:109:VAL:HG11	1:A:309:ILE:HD13	1.81	0.63
1:A:112:PRO:HG2	1:A:299:PHE:CE2	2.34	0.63
1:B:98:THR:HG21	1:B:316:SER:CB	2.29	0.63
1:B:272:LEU:CD1	1:B:273:LYS:H	2.12	0.62
1:B:278:GLN:NE2	1:B:308:ARG:HH22	1.97	0.62
1:A:185:ASN:ND2	1:A:185:ASN:N	2.42	0.62
1:A:128:ASN:CG	1:A:129:ALA:H	2.08	0.62
1:A:179:ILE:HG23	1:A:181:LEU:HG	1.82	0.62
1:A:76:LEU:HG	1:A:78:MSE:HE2	1.82	0.61
1:A:108:MSE:O	1:A:277:SER:HA	2.01	0.61
1:B:167:VAL:O	1:B:170:VAL:HG12	2.01	0.60
1:A:108:MSE:HG3	1:A:277:SER:HB2	1.82	0.60
1:A:248:VAL:HG22	1:A:249:GLY:N	2.16	0.60
1:B:167:VAL:HG13	1:B:173:ILE:HG12	1.83	0.60
1:B:287:ILE:HB	1:B:298:TYR:CD2	2.36	0.60
1:B:272:LEU:O	1:B:274:ASN:N	2.29	0.60
1:B:319:ILE:HG12	1:B:320:THR:N	2.17	0.60
1:B:306:LEU:O	1:B:310:ILE:HG13	2.02	0.59
1:B:184:VAL:CG2	1:B:209:GLU:HG2	2.32	0.59
1:A:148:PRO:HB3	1:B:290:GLU:CD	2.28	0.59
1:B:155:PHE:C	1:B:155:PHE:CD1	2.81	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:156:ILE:HG22	1:B:256:MSE:HE2	1.84	0.59
1:A:76:LEU:CD2	1:A:78:MSE:HE2	2.32	0.59
1:A:159:ASN:HD21	1:A:162:GLY:N	1.99	0.59
1:B:209:GLU:H	1:B:209:GLU:CD	2.10	0.59
1:A:163:PHE:HZ	1:A:224:VAL:HG11	1.68	0.58
1:B:160:MSE:O	1:B:164:LYS:HE2	2.03	0.58
1:B:76:LEU:HD13	1:B:96:LEU:HD23	1.86	0.58
1:B:91:ALA:O	1:B:128:ASN:HB2	2.04	0.58
1:B:172:GLY:O	1:B:173:ILE:HD13	2.04	0.58
1:A:186:SER:HB3	1:A:188:PHE:CE2	2.39	0.57
1:A:187:LYS:O	1:A:187:LYS:HG2	2.04	0.57
1:A:173:ILE:HD12	1:A:227:GLY:C	2.30	0.57
1:A:226:ILE:HG22	1:A:230:ASN:ND2	2.20	0.56
1:A:74:SER:HA	1:A:98:THR:HA	1.86	0.56
1:A:182:THR:O	1:A:184:VAL:N	2.39	0.56
1:A:94:ILE:C	1:A:95:ILE:HD12	2.31	0.56
1:A:189:VAL:HG22	1:A:190:LYS:N	2.21	0.56
1:B:278:GLN:HE22	1:B:308:ARG:HH22	1.53	0.56
1:B:108:MSE:HE3	1:B:277:SER:HB2	1.87	0.56
1:A:260:ASP:OD2	1:A:260:ASP:C	2.49	0.55
1:B:196:ASN:C	1:B:196:ASN:OD1	2.50	0.55
1:A:173:ILE:HD12	1:A:227:GLY:O	2.08	0.54
1:A:119:TYR:HD2	1:A:145:LYS:HE3	1.72	0.54
1:B:156:ILE:CG2	1:B:256:MSE:HE2	2.37	0.54
1:A:185:ASN:HD22	1:A:186:SER:N	2.02	0.54
1:B:158:ILE:HD13	1:B:158:ILE:H	1.73	0.53
1:A:251:ASN:N	1:A:251:ASN:HD22	2.07	0.53
1:A:190:LYS:HG2	1:A:191:GLY:N	2.24	0.53
1:A:193:ILE:HG23	1:A:193:ILE:O	2.08	0.52
1:B:92:ASP:O	1:B:113:ARG:HB3	2.10	0.52
1:B:248:VAL:O	1:B:251:ASN:HB2	2.09	0.52
1:B:292:TYR:HB3	1:B:294:PHE:CE2	2.45	0.52
1:A:244:ILE:O	1:A:244:ILE:CD1	2.51	0.52
1:B:105:ALA:O	1:B:270:SER:HB3	2.10	0.52
1:A:77:LEU:HD13	1:A:225:ILE:HD12	1.92	0.52
1:A:145:LYS:HB3	1:A:145:LYS:HZ2	1.75	0.52
1:B:133:ASN:C	1:B:135:GLY:H	2.19	0.51
1:A:154:TYR:CD1	1:A:256:MSE:HG2	2.45	0.51
1:B:77:LEU:HD22	1:B:156:ILE:HD11	1.92	0.51
1:A:182:THR:C	1:A:184:VAL:N	2.64	0.51
1:B:77:LEU:HD23	1:B:156:ILE:HG12	1.92	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:207:ARG:HG3	1:B:207:ARG:HH11	1.77	0.50
1:B:307:GLU:O	1:B:308:ARG:C	2.55	0.50
1:A:74:SER:CB	1:A:98:THR:HG22	2.42	0.50
1:A:173:ILE:HG23	1:A:227:GLY:HA3	1.94	0.50
1:A:196:ASN:ND2	1:A:199:GLU:HG3	2.27	0.50
1:B:107:GLU:HG3	1:B:316:SER:OG	2.12	0.49
1:A:180:ASP:CG	1:A:182:THR:HG22	2.37	0.49
1:A:76:LEU:CG	1:A:78:MSE:HE2	2.42	0.49
1:A:115:THR:HA	1:A:299:PHE:HB3	1.94	0.49
1:A:278:GLN:NE2	1:A:312:MSE:HE1	2.28	0.49
1:A:159:ASN:ND2	1:A:162:GLY:N	2.54	0.48
1:A:257:THR:H	1:A:260:ASP:CG	2.21	0.48
1:B:288:TYR:HE1	1:B:293:GLY:C	2.21	0.48
1:A:119:TYR:O	1:A:123:ASP:HB3	2.12	0.48
1:A:195:LEU:HD22	1:A:199:GLU:HB3	1.95	0.48
1:A:288:TYR:HA	1:A:295:ASP:HA	1.95	0.48
1:A:68:GLU:C	1:A:70:LYS:H	2.21	0.48
1:A:177:ASN:N	1:A:191:GLY:O	2.42	0.48
1:A:127:ILE:O	1:A:131:TYR:HB2	2.11	0.48
1:A:270:SER:C	1:A:272:LEU:H	2.20	0.48
1:B:160:MSE:HG3	1:B:205:ARG:NH2	2.25	0.48
1:B:185:ASN:HD22	1:B:187:LYS:HB3	1.78	0.48
1:A:229:ALA:C	1:A:231:LYS:H	2.22	0.48
1:A:74:SER:HB2	1:A:98:THR:HG22	1.96	0.48
1:A:96:LEU:O	1:A:108:MSE:HA	2.13	0.48
1:A:173:ILE:CG2	1:A:227:GLY:HA3	2.44	0.48
1:B:70:LYS:HD2	1:B:101:LYS:HD2	1.95	0.48
1:B:272:LEU:C	1:B:274:ASN:H	2.19	0.48
1:B:96:LEU:HD12	1:B:313:PHE:CZ	2.49	0.47
1:B:100:ASN:OD1	1:B:103:GLN:HG3	2.14	0.47
1:B:158:ILE:HD13	1:B:158:ILE:O	2.14	0.47
1:B:246:LYS:C	1:B:248:VAL:H	2.21	0.47
1:B:319:ILE:CG1	1:B:320:THR:H	2.24	0.47
1:B:106:VAL:HG23	1:B:270:SER:O	2.14	0.47
1:B:90:ARG:HB3	1:B:131:TYR:HB3	1.95	0.47
1:A:100:ASN:ND2	1:A:102:GLN:H	2.12	0.47
1:A:131:TYR:CD1	1:A:142:ALA:HB2	2.50	0.47
1:A:257:THR:O	1:A:260:ASP:OD2	2.33	0.47
1:A:270:SER:C	1:A:272:LEU:N	2.72	0.47
1:A:179:ILE:HD13	1:A:179:ILE:C	2.40	0.47
1:B:204:VAL:HG12	1:B:204:VAL:O	2.13	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:155:PHE:CD1	1:B:155:PHE:O	2.67	0.46
1:A:115:THR:HG23	1:A:301:PRO:HG3	1.97	0.46
1:A:189:VAL:HG12	1:A:193:ILE:HD12	1.98	0.46
1:B:182:THR:C	1:B:184:VAL:N	2.70	0.46
1:B:185:ASN:C	1:B:187:LYS:H	2.22	0.46
1:B:249:GLY:C	1:B:251:ASN:H	2.22	0.46
1:A:147:MSE:HE1	1:A:309:ILE:HG21	1.96	0.46
1:B:141:SER:O	1:B:145:LYS:HG3	2.16	0.46
1:B:158:ILE:HD13	1:B:158:ILE:N	2.31	0.46
1:A:253:GLN:HE21	1:A:253:GLN:HB2	1.55	0.46
1:A:76:LEU:HD23	1:A:78:MSE:HE2	1.98	0.46
1:A:159:ASN:ND2	1:A:159:ASN:N	2.57	0.46
1:B:261:ILE:O	1:B:264:MSE:HB2	2.16	0.46
1:A:264:MSE:HA	1:A:268:TYR:CD2	2.45	0.45
1:B:182:THR:HG22	1:B:183:GLU:H	1.81	0.45
1:B:208:HIS:O	1:B:209:GLU:C	2.59	0.45
1:B:155:PHE:O	1:B:155:PHE:CG	2.69	0.45
1:B:161:GLU:HB3	1:B:251:ASN:HD21	1.81	0.45
1:B:226:ILE:HD11	1:B:275:VAL:HG21	1.98	0.45
1:B:118:ASP:C	1:B:145:LYS:HZ1	2.24	0.45
1:A:167:VAL:HG13	1:A:173:ILE:CG1	2.41	0.45
1:B:168:ASP:OD1	1:B:196:ASN:HB2	2.17	0.45
1:A:80:SER:C	1:A:160:MSE:HE2	2.43	0.45
1:A:308:ARG:C	1:A:308:ARG:HD3	2.42	0.45
1:B:111:ILE:HD11	1:B:147:MSE:HE1	1.98	0.45
1:A:126:LYS:H	1:A:126:LYS:HG2	1.34	0.44
1:A:163:PHE:O	1:A:167:VAL:HG23	2.17	0.44
1:B:66:ASP:HB3	1:B:71:LYS:HD3	1.99	0.44
1:B:139:THR:O	1:B:140:VAL:C	2.59	0.44
1:A:100:ASN:ND2	1:A:100:ASN:C	2.74	0.44
1:A:196:ASN:H	1:A:199:GLU:HB2	1.81	0.44
1:A:264:MSE:C	1:A:266:ALA:H	2.26	0.44
1:A:140:VAL:O	1:A:144:GLU:HG3	2.17	0.44
1:B:249:GLY:C	1:B:251:ASN:N	2.75	0.44
1:A:220:ARG:O	1:A:224:VAL:HG23	2.17	0.44
1:B:173:ILE:O	1:B:194:THR:HA	2.18	0.44
1:B:173:ILE:HG22	1:B:174:THR:N	2.33	0.44
1:B:275:VAL:HG13	1:B:275:VAL:O	2.17	0.44
1:B:117:VAL:HB	1:B:146:LEU:HD13	1.99	0.44
1:A:181:LEU:O	1:A:182:THR:O	2.36	0.43
1:B:146:LEU:HD23	1:B:147:MSE:HE2	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:182:THR:O	1:B:184:VAL:N	2.51	0.43
1:B:108:MSE:HE1	1:B:222:ARG:CG	2.27	0.43
1:A:209:GLU:OE1	1:A:209:GLU:N	2.49	0.43
1:B:153:ASP:OD1	1:B:153:ASP:N	2.50	0.43
1:A:210:ASP:CB	1:A:220:ARG:HH21	2.32	0.43
1:B:126:LYS:HG3	1:B:298:TYR:CE1	2.53	0.42
1:A:109:VAL:HA	1:A:278:GLN:O	2.18	0.42
1:A:114:ASP:O	1:A:115:THR:C	2.61	0.42
1:B:170:VAL:O	1:B:170:VAL:HG13	2.16	0.42
1:A:181:LEU:O	1:A:188:PHE:HB2	2.20	0.42
1:B:76:LEU:HG	1:B:78:MSE:HE2	2.01	0.42
1:A:109:VAL:HG22	1:A:278:GLN:HG3	2.02	0.42
1:A:196:ASN:HD21	1:A:199:GLU:HG3	1.83	0.42
1:B:130:SER:C	1:B:132:SER:H	2.27	0.42
1:B:233:ILE:HD12	1:B:233:ILE:O	2.18	0.42
1:B:275:VAL:O	1:B:275:VAL:HG22	2.19	0.42
1:B:100:ASN:ND2	1:B:316:SER:O	2.39	0.42
1:B:200:ALA:O	1:B:204:VAL:HG23	2.20	0.42
1:B:287:ILE:HB	1:B:298:TYR:HD2	1.82	0.42
1:B:158:ILE:N	1:B:158:ILE:CD1	2.83	0.41
1:A:178:ASP:OD2	1:A:178:ASP:N	2.54	0.41
1:B:95:ILE:HD11	1:B:221:GLN:OE1	2.20	0.41
1:B:209:GLU:CD	1:B:209:GLU:N	2.77	0.41
1:A:210:ASP:HA	1:A:211:PRO:HD2	1.87	0.41
1:A:78:MSE:SE	1:A:136:PRO:HB3	2.71	0.41
1:B:111:ILE:CD1	1:B:147:MSE:HE1	2.51	0.41
1:B:217:ARG:HA	1:B:220:ARG:CZ	2.50	0.41
1:A:103:GLN:O	1:A:104:ASN:C	2.63	0.41
1:B:152:VAL:O	1:B:152:VAL:HG12	2.21	0.41
1:B:154:TYR:CD1	1:B:256:MSE:HG2	2.55	0.41
1:B:184:VAL:HG23	1:B:209:GLU:HG2	2.03	0.41
1:A:155:PHE:CD1	1:A:155:PHE:C	2.99	0.40
1:B:127:ILE:HG23	1:B:128:ASN:N	2.35	0.40
1:B:196:ASN:OD1	1:B:199:GLU:HG3	2.22	0.40
1:A:163:PHE:CZ	1:A:224:VAL:HG11	2.51	0.40
1:B:133:ASN:HB3	1:B:138:GLY:HA3	2.03	0.40
1:B:299:PHE:O	1:B:301:PRO:HD3	2.22	0.40
1:A:100:ASN:ND2	1:A:102:GLN:HB2	2.30	0.40
1:A:226:ILE:HD11	1:A:275:VAL:HG11	2.02	0.40
1:B:126:LYS:O	1:B:129:ALA:HB3	2.21	0.40
1:B:184:VAL:HG21	1:B:209:GLU:HG2	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	229/279 (82%)	181 (79%)	38 (17%)	10 (4%)	2	8
1	B	229/279 (82%)	187 (82%)	33 (14%)	9 (4%)	2	10
All	All	458/558 (82%)	368 (80%)	71 (16%)	19 (4%)	2	9

All (19) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	182	THR
1	A	183	GLU
1	A	272	LEU
1	B	208	HIS
1	B	209	GLU
1	B	246	LYS
1	B	153	ASP
1	B	245	MSE
1	B	273	LYS
1	A	134	GLY
1	A	274	ASN
1	A	318	ASP
1	A	319	ILE
1	A	230	ASN
1	B	197	GLY
1	B	233	ILE
1	B	318	ASP
1	A	189	VAL
1	A	211	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	201/231 (87%)	174 (87%)	27 (13%)	4	12
1	B	206/231 (89%)	187 (91%)	19 (9%)	8	27
All	All	407/462 (88%)	361 (89%)	46 (11%)	5	19

All (46) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	68	GLU
1	A	108	MSE
1	A	126	LYS
1	A	127	ILE
1	A	145	LYS
1	A	159	ASN
1	A	161	GLU
1	A	179	ILE
1	A	182	THR
1	A	184	VAL
1	A	185	ASN
1	A	187	LYS
1	A	205	ARG
1	A	207	ARG
1	A	209	GLU
1	A	220	ARG
1	A	245	MSE
1	A	253	GLN
1	A	257	THR
1	A	258	LEU
1	A	262	THR
1	A	267	ASN
1	A	272	LEU
1	A	274	ASN
1	A	296	LEU
1	A	308	ARG
1	A	312	MSE

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Mol	Chain	Res	Type
1	B	67	LEU
1	B	96	LEU
1	B	98	THR
1	B	113	ARG
1	B	126	LYS
1	B	158	ILE
1	B	166	LEU
1	B	170	VAL
1	B	179	ILE
1	B	182	THR
1	B	218	GLN
1	B	226	ILE
1	B	233	ILE
1	B	253	GLN
1	B	259	THR
1	B	272	LEU
1	B	296	LEU
1	B	304	THR
1	B	318	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	100	ASN
1	A	104	ASN
1	A	121	ASN
1	A	159	ASN
1	A	185	ASN
1	A	253	GLN
1	A	278	GLN
1	B	185	ASN
1	B	267	ASN
1	B	278	GLN

5.3.3 RNA ⓘ

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	226/279 (81%)	0.22	12 (5%) 32 25	40, 71, 117, 151	0
1	B	228/279 (81%)	-0.00	2 (0%) 81 75	27, 60, 98, 128	0
All	All	454/558 (81%)	0.11	14 (3%) 51 42	27, 66, 115, 151	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	130	SER	6.0
1	A	131	TYR	4.0
1	B	244	ILE	3.8
1	A	193	ILE	3.2
1	B	319	ILE	3.2
1	A	184	VAL	2.8
1	A	105	ALA	2.8
1	A	258	LEU	2.7
1	A	243	SER	2.4
1	A	119	TYR	2.2
1	A	272	LEU	2.2
1	A	169	ALA	2.2
1	A	273	LYS	2.1
1	A	187	LYS	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.