



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 11, 2016 – 05:39 AM EDT

PDB ID : 5DSS
Title : MP-4 contributes to snake venom neutralization by Mucuna pruriens seeds through stimulation of cross-reactive antibodies
Authors : Kumar, A.; Nair, D.T.; Salunke, D.M.
Deposited on : 2015-09-17
Resolution : 2.80 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 : unknown
Xtriage (Phenix) : 1.9-1692
EDS : rb-20027107
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) : rb-20027107

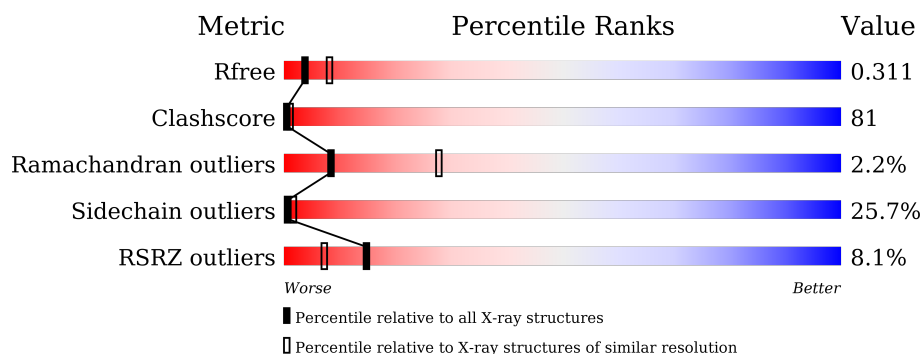
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	2393 (2.80-2.80)
Clashscore	102246	2827 (2.80-2.80)
Ramachandran outliers	100387	2782 (2.80-2.80)
Sidechain outliers	100360	2784 (2.80-2.80)
RSRZ outliers	91569	2404 (2.80-2.80)

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	B	185	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 1533 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.

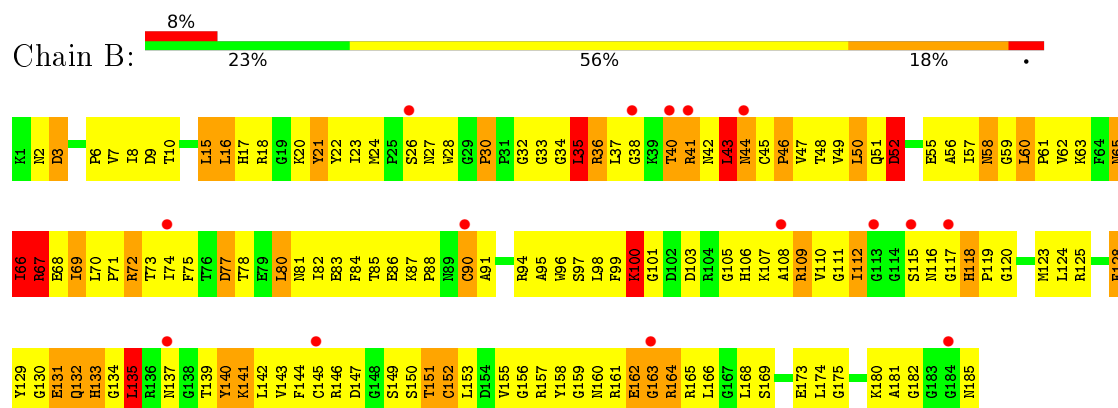
- Molecule 1 is a protein called MP-4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	185	Total	C	N	O	S	0	0	0
			1411	873	266	266	6			

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	122	Total	O	0	0
			122	122		

- Molecule 1: MP-4



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	51.00Å 69.18Å 45.32Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	37.91 – 2.80 37.91 – 2.80	Depositor EDS
% Data completeness (in resolution range)	94.7 (37.91-2.80) 94.8 (37.91-2.80)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	13.43 (at 2.81Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.267 , 0.284 0.267 , 0.311	Depositor DCC
R_{free} test set	179 reflections (4.64%)	DCC
Wilson B-factor (Å ²)	28.7	Xtriage
Anisotropy	0.796	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 133.9	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.42$, $\langle L^2 \rangle = 0.25$	Xtriage
Outliers	0 of 4048 reflections	Xtriage
F_o, F_c correlation	0.81	EDS
Total number of atoms	1533	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 16.05% 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	B	0.84	2/1440 (0.1%)	1.34	20/1936 (1.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	2

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	185	ASN	C-OXT	14.49	1.50	1.23
1	B	96	TRP	CD2-CE2	5.26	1.47	1.41

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	151	THR	N-CA-C	-11.78	79.18	111.00
1	B	43	LEU	N-CA-C	9.43	136.47	111.00
1	B	66	ILE	N-CA-C	-9.03	86.62	111.00
1	B	117	GLY	N-CA-C	-8.54	91.75	113.10
1	B	133	HIS	N-CA-C	-7.71	90.18	111.00
1	B	115	SER	N-CA-C	-7.66	90.32	111.00
1	B	67	ARG	N-CA-C	6.98	129.85	111.00
1	B	100	LYS	N-CA-C	6.53	128.64	111.00
1	B	35	LEU	CA-CB-CG	6.41	130.03	115.30
1	B	15	LEU	CA-CB-CG	6.37	129.95	115.30
1	B	43	LEU	CA-CB-CG	6.28	129.74	115.30
1	B	52	ASP	CB-CG-OD1	6.10	123.79	118.30
1	B	107	LYS	N-CA-C	-5.94	94.97	111.00
1	B	141	LYS	N-CA-C	5.76	126.56	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	135	LEU	N-CA-C	5.74	126.49	111.00
1	B	88	PRO	N-CA-C	-5.52	97.76	112.10
1	B	60	LEU	CA-CB-CG	5.50	127.94	115.30
1	B	163	GLY	N-CA-C	-5.36	99.71	113.10
1	B	69	ILE	CB-CA-C	-5.28	101.05	111.60
1	B	100	LYS	CB-CA-C	-5.06	100.28	110.40

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	100	LYS	Peptide
1	B	44	ASN	Peptide

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	B	1411	0	1362	225	0
2	B	122	0	0	29	0
All	All	1533	0	1362	225	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 81.

All (225) 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:62:VAL:HG12	1:B:82:ILE:HD11	1.21	1.14
1:B:118:HIS:HB3	1:B:119:PRO:CD	1.80	1.12
1:B:66:ILE:HB	1:B:69:ILE:HD12	1.29	1.12
1:B:8:ILE:HG22	1:B:10:THR:H	1.08	1.11
1:B:118:HIS:CG	1:B:119:PRO:HD3	1.87	1.09
1:B:118:HIS:HB3	1:B:119:PRO:HD3	1.37	1.03
1:B:118:HIS:CB	1:B:119:PRO:HD3	1.87	1.03

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:10:THR:HG22	1:B:10:THR:O	1.58	1.01
1:B:118:HIS:CB	1:B:119:PRO:CD	2.41	0.99
1:B:144:PHE:O	1:B:152:CYS:HB3	1.62	0.97
1:B:87:LYS:HB3	1:B:94:ARG:HG2	1.42	0.97
1:B:8:ILE:HG22	1:B:10:THR:N	1.78	0.97
1:B:18:ARG:HG2	1:B:63:LYS:HE3	1.46	0.95
1:B:146:ARG:HB2	1:B:150:SER:CB	1.96	0.95
1:B:35:LEU:O	1:B:59:GLY:HA3	1.66	0.94
1:B:146:ARG:HB2	1:B:150:SER:HB3	1.47	0.94
1:B:144:PHE:O	1:B:152:CYS:CB	2.17	0.92
1:B:156:GLY:HA3	1:B:169:SER:OG	1.69	0.92
1:B:41:ARG:HH11	1:B:41:ARG:HG3	1.35	0.91
1:B:146:ARG:HD3	1:B:150:SER:HB2	1.49	0.91
1:B:180:LYS:HE3	1:B:182:GLY:O	1.69	0.91
1:B:63:LYS:HD3	2:B:207:HOH:O	1.69	0.90
1:B:132:GLN:HG2	1:B:134:GLY:N	1.88	0.89
1:B:33:GLY:HA3	1:B:52:ASP:O	1.74	0.88
1:B:62:VAL:CG1	1:B:82:ILE:HD11	2.02	0.88
1:B:62:VAL:HG12	1:B:82:ILE:CD1	2.04	0.87
1:B:145:CYS:HA	1:B:152:CYS:HB3	1.56	0.87
1:B:87:LYS:HB3	1:B:94:ARG:CG	2.05	0.86
1:B:99:PHE:CE1	1:B:111:GLY:HA3	2.11	0.85
1:B:69:ILE:HG22	1:B:70:LEU:H	1.39	0.85
1:B:157:ARG:HD2	1:B:164:ARG:HD3	1.59	0.84
1:B:26:SER:HB2	1:B:175:GLY:O	1.77	0.84
1:B:105:GLY:HA3	1:B:109:ARG:NE	1.92	0.83
1:B:41:ARG:HG3	1:B:41:ARG:NH1	1.89	0.83
1:B:47:VAL:HG12	1:B:111:GLY:HA2	1.59	0.82
1:B:75:PHE:O	1:B:78:THR:HG22	1.81	0.81
1:B:77:ASP:O	2:B:201:HOH:O	1.98	0.81
1:B:15:LEU:HG	1:B:21:TYR:CE2	2.16	0.80
1:B:151:THR:O	1:B:151:THR:CG2	2.30	0.79
1:B:57:ILE:HG22	1:B:59:GLY:H	1.47	0.79
1:B:156:GLY:HA2	1:B:173:GLU:HB3	1.65	0.78
1:B:37:LEU:HD23	1:B:60:LEU:HD22	1.65	0.78
1:B:61:PRO:HB2	1:B:85:THR:OG1	1.84	0.78
1:B:118:HIS:HB3	1:B:119:PRO:HD2	1.66	0.78
1:B:10:THR:CG2	1:B:10:THR:O	2.32	0.78
1:B:146:ARG:CD	1:B:150:SER:HB2	2.14	0.77
1:B:41:ARG:HH11	1:B:41:ARG:CG	1.96	0.77
1:B:124:LEU:HD13	1:B:146:ARG:HD2	1.67	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:20:LYS:O	1:B:180:LYS:HD2	1.85	0.77
1:B:63:LYS:HD2	2:B:272:HOH:O	1.85	0.76
1:B:8:ILE:CG2	1:B:10:THR:H	1.93	0.76
1:B:132:GLN:HG2	1:B:134:GLY:H	1.49	0.76
1:B:7:VAL:HB	1:B:74:ILE:O	1.84	0.75
1:B:50:LEU:HD11	1:B:163:GLY:HA3	1.69	0.74
1:B:69:ILE:HG22	1:B:70:LEU:N	2.02	0.74
1:B:105:GLY:HA3	1:B:109:ARG:HE	1.51	0.74
1:B:119:PRO:HB3	2:B:247:HOH:O	1.89	0.73
1:B:118:HIS:ND1	1:B:119:PRO:HD3	2.03	0.72
1:B:146:ARG:H	1:B:150:SER:HB3	1.54	0.72
1:B:151:THR:O	1:B:151:THR:HG23	1.87	0.72
1:B:118:HIS:CG	1:B:119:PRO:CD	2.67	0.72
1:B:46:PRO:HB3	2:B:233:HOH:O	1.87	0.72
1:B:67:ARG:HB2	1:B:81:ASN:HB2	1.72	0.71
1:B:26:SER:O	2:B:203:HOH:O	2.07	0.71
1:B:99:PHE:O	1:B:109:ARG:NH2	2.24	0.71
1:B:30:PRO:HD3	1:B:157:ARG:HH22	1.56	0.71
1:B:37:LEU:HD23	1:B:60:LEU:CD2	2.21	0.70
1:B:20:LYS:C	1:B:21:TYR:HD1	1.94	0.70
1:B:99:PHE:HE1	1:B:111:GLY:HA3	1.55	0.70
1:B:66:ILE:HB	1:B:69:ILE:CD1	2.15	0.70
1:B:159:GLY:HA3	2:B:218:HOH:O	1.91	0.69
1:B:65:ASN:HB2	1:B:67:ARG:HA	1.73	0.69
1:B:146:ARG:CB	1:B:150:SER:HB3	2.22	0.68
1:B:132:GLN:HB2	1:B:140:TYR:HB2	1.75	0.68
1:B:87:LYS:CB	1:B:94:ARG:HG2	2.21	0.68
1:B:100:LYS:HG2	1:B:101:GLY:N	2.08	0.68
1:B:66:ILE:CB	1:B:69:ILE:HD12	2.18	0.67
1:B:16:LEU:HG	1:B:72:ARG:O	1.95	0.67
1:B:87:LYS:N	2:B:206:HOH:O	2.23	0.66
1:B:145:CYS:CA	1:B:152:CYS:HB3	2.26	0.66
1:B:125:ARG:NH2	2:B:209:HOH:O	2.24	0.66
1:B:36:ARG:HD3	1:B:36:ARG:N	2.11	0.66
1:B:100:LYS:CE	2:B:256:HOH:O	2.45	0.65
1:B:125:ARG:HB3	1:B:147:ASP:CG	2.17	0.65
1:B:90:CYS:O	1:B:90:CYS:SG	2.55	0.65
1:B:180:LYS:HG3	1:B:182:GLY:H	1.63	0.64
1:B:65:ASN:HB3	2:B:260:HOH:O	1.96	0.64
1:B:57:ILE:N	1:B:57:ILE:HD12	2.13	0.63
1:B:131:GLU:N	1:B:141:LYS:O	2.30	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:40:THR:HG23	1:B:162:GLU:OE1	1.99	0.63
1:B:20:LYS:C	1:B:21:TYR:CD1	2.72	0.63
1:B:132:GLN:HA	1:B:140:TYR:HB3	1.79	0.63
1:B:144:PHE:O	1:B:152:CYS:HB2	1.97	0.62
1:B:30:PRO:HD3	1:B:157:ARG:NH2	2.13	0.62
1:B:47:VAL:CG1	1:B:111:GLY:HA2	2.30	0.62
1:B:95:ALA:HB2	2:B:232:HOH:O	1.98	0.62
1:B:173:GLU:O	1:B:174:LEU:HB2	2.00	0.61
1:B:51:GLN:HB2	1:B:164:ARG:HB3	1.82	0.61
1:B:110:VAL:HG13	1:B:110:VAL:O	2.00	0.60
1:B:67:ARG:NH2	1:B:68:GLU:OE2	2.34	0.60
1:B:141:LYS:HE2	2:B:252:HOH:O	2.02	0.59
1:B:67:ARG:HD3	1:B:81:ASN:HB3	1.84	0.59
1:B:49:VAL:HB	1:B:110:VAL:HG11	1.83	0.59
1:B:16:LEU:HD23	1:B:17:HIS:H	1.68	0.59
1:B:32:GLY:O	1:B:55:GLU:HG3	2.02	0.59
1:B:50:LEU:CD1	1:B:163:GLY:HA3	2.33	0.58
1:B:34:GLY:O	1:B:52:ASP:HB2	2.03	0.58
1:B:18:ARG:CG	1:B:63:LYS:HE3	2.28	0.58
1:B:51:GLN:HE21	1:B:166:LEU:HD21	1.70	0.57
1:B:40:THR:OG1	1:B:162:GLU:HB2	2.05	0.57
1:B:153:LEU:N	1:B:153:LEU:HD12	2.19	0.57
1:B:15:LEU:HD23	1:B:74:ILE:HD12	1.87	0.57
1:B:72:ARG:NH1	2:B:213:HOH:O	2.37	0.57
1:B:49:VAL:H	1:B:110:VAL:HG13	1.70	0.56
1:B:55:GLU:N	1:B:55:GLU:OE1	2.33	0.56
1:B:43:LEU:O	1:B:44:ASN:C	2.42	0.56
1:B:100:LYS:HE2	2:B:256:HOH:O	2.03	0.56
1:B:128:PHE:HD1	1:B:143:VAL:O	1.87	0.56
1:B:67:ARG:CZ	1:B:67:ARG:HB3	2.35	0.56
1:B:17:HIS:CE1	1:B:74:ILE:HG12	2.41	0.56
1:B:180:LYS:HD2	1:B:181:ALA:H	1.72	0.55
1:B:7:VAL:O	1:B:15:LEU:HB3	2.07	0.55
1:B:36:ARG:CD	1:B:36:ARG:N	2.70	0.54
1:B:49:VAL:HB	1:B:110:VAL:CG1	2.37	0.54
1:B:26:SER:C	1:B:27:ASN:HD22	2.10	0.54
1:B:85:THR:O	1:B:85:THR:HG22	2.07	0.54
1:B:67:ARG:HD2	1:B:83:GLU:CD	2.27	0.54
1:B:27:ASN:N	1:B:27:ASN:HD22	2.05	0.54
1:B:30:PRO:CD	1:B:30:PRO:O	2.56	0.54
1:B:99:PHE:C	1:B:109:ARG:NH2	2.61	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:50:LEU:HD13	1:B:164:ARG:H	1.73	0.53
1:B:67:ARG:HG3	1:B:81:ASN:O	2.08	0.53
1:B:146:ARG:NH2	1:B:149:SER:HB3	2.23	0.53
1:B:23:ILE:HD12	1:B:35:LEU:HD21	1.91	0.53
1:B:146:ARG:HB2	1:B:150:SER:HB2	1.84	0.53
1:B:99:PHE:HA	1:B:123:MET:HA	1.90	0.53
1:B:45:CYS:O	1:B:47:VAL:HG23	2.08	0.52
1:B:47:VAL:HG13	1:B:112:ILE:O	2.09	0.52
1:B:37:LEU:CD2	1:B:60:LEU:HD22	2.37	0.52
1:B:69:ILE:CG2	1:B:70:LEU:H	2.17	0.52
1:B:47:VAL:CG1	1:B:111:GLY:CA	2.88	0.52
1:B:58:ASN:OD1	1:B:58:ASN:N	2.30	0.52
1:B:129:TYR:N	1:B:143:VAL:O	2.43	0.52
1:B:36:ARG:HD3	1:B:36:ARG:H	1.75	0.51
1:B:141:LYS:CE	2:B:252:HOH:O	2.58	0.51
1:B:6:PRO:HB3	1:B:73:THR:HG21	1.92	0.51
1:B:7:VAL:O	1:B:15:LEU:CB	2.59	0.51
1:B:160:ASN:O	1:B:160:ASN:OD1	2.29	0.50
1:B:60:LEU:HB3	1:B:84:PHE:CD1	2.47	0.50
1:B:155:VAL:HG22	1:B:168:LEU:CD2	2.41	0.50
1:B:30:PRO:HD2	1:B:30:PRO:O	2.10	0.50
1:B:86:GLU:HA	2:B:206:HOH:O	2.11	0.50
1:B:105:GLY:HA3	1:B:109:ARG:CZ	2.40	0.50
1:B:18:ARG:HG2	1:B:63:LYS:CE	2.31	0.50
1:B:20:LYS:HB3	1:B:61:PRO:HB3	1.92	0.49
1:B:21:TYR:CD1	1:B:21:TYR:N	2.80	0.49
1:B:16:LEU:HD23	1:B:17:HIS:N	2.27	0.49
1:B:134:GLY:C	1:B:135:LEU:HG	2.25	0.49
1:B:132:GLN:HG2	1:B:134:GLY:CA	2.43	0.48
1:B:23:ILE:HD11	1:B:62:VAL:HG11	1.94	0.48
1:B:38:GLY:HA2	2:B:222:HOH:O	2.12	0.48
1:B:99:PHE:CD1	1:B:99:PHE:N	2.81	0.48
1:B:146:ARG:N	1:B:150:SER:HB3	2.27	0.48
1:B:21:TYR:N	1:B:21:TYR:HD1	2.12	0.48
1:B:42:ASN:HB2	2:B:255:HOH:O	2.12	0.48
1:B:72:ARG:NH2	2:B:214:HOH:O	2.38	0.47
1:B:109:ARG:NH1	1:B:120:GLY:HA3	2.29	0.47
1:B:110:VAL:CG1	1:B:110:VAL:O	2.62	0.47
1:B:129:TYR:CD2	1:B:145:CYS:SG	3.07	0.47
1:B:156:GLY:CA	1:B:173:GLU:HB3	2.40	0.47
1:B:145:CYS:HB3	2:B:208:HOH:O	2.13	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:47:VAL:HG12	1:B:111:GLY:CA	2.37	0.47
1:B:146:ARG:CB	1:B:150:SER:CB	2.82	0.46
1:B:17:HIS:NE2	1:B:74:ILE:HG12	2.31	0.46
1:B:66:ILE:HA	1:B:80:LEU:HA	1.97	0.46
1:B:87:LYS:HB3	1:B:94:ARG:HG3	1.93	0.46
1:B:109:ARG:H	1:B:109:ARG:HG3	1.53	0.46
1:B:41:ARG:HB2	2:B:259:HOH:O	2.15	0.46
1:B:56:ALA:HB3	1:B:57:ILE:HD12	1.97	0.46
1:B:21:TYR:O	1:B:61:PRO:HA	2.16	0.46
1:B:20:LYS:H	1:B:180:LYS:NZ	2.14	0.46
1:B:161:ARG:NH1	2:B:220:HOH:O	2.48	0.46
1:B:116:ASN:ND2	1:B:119:PRO:O	2.49	0.45
1:B:99:PHE:O	1:B:108:ALA:HA	2.16	0.45
1:B:10:THR:OG1	1:B:139:THR:HG23	2.16	0.45
1:B:22:TYR:CD1	1:B:58:ASN:HB2	2.52	0.45
1:B:48:THR:CG2	1:B:165:ARG:HG2	2.46	0.45
1:B:173:GLU:O	1:B:174:LEU:CB	2.65	0.45
1:B:130:GLY:HA2	1:B:142:LEU:HA	1.99	0.45
1:B:132:GLN:CG	1:B:134:GLY:HA3	2.46	0.45
1:B:66:ILE:CG2	1:B:69:ILE:HD11	2.47	0.45
1:B:157:ARG:HG3	1:B:174:LEU:CD1	2.47	0.44
1:B:27:ASN:ND2	1:B:27:ASN:N	2.64	0.44
1:B:151:THR:O	1:B:151:THR:HG22	2.15	0.44
1:B:155:VAL:HA	1:B:168:LEU:HA	1.99	0.44
1:B:43:LEU:O	1:B:44:ASN:O	2.36	0.44
1:B:97:SER:HA	2:B:254:HOH:O	2.17	0.44
1:B:18:ARG:O	1:B:18:ARG:HG2	2.18	0.44
1:B:55:GLU:N	1:B:55:GLU:CD	2.71	0.44
1:B:21:TYR:CE1	1:B:180:LYS:HD3	2.52	0.44
1:B:3:ASP:OD1	1:B:3:ASP:N	2.51	0.44
1:B:10:THR:HG23	2:B:285:HOH:O	2.18	0.44
1:B:123:MET:HG3	1:B:123:MET:O	2.18	0.44
1:B:129:TYR:HD2	1:B:145:CYS:SG	2.41	0.44
1:B:47:VAL:HG13	1:B:112:ILE:N	2.33	0.43
1:B:100:LYS:HG2	1:B:101:GLY:CA	2.48	0.43
1:B:63:LYS:HB2	2:B:217:HOH:O	2.19	0.43
1:B:20:LYS:HB2	1:B:20:LYS:HE2	1.60	0.43
1:B:36:ARG:NH2	1:B:50:LEU:HB3	2.34	0.43
1:B:85:THR:O	1:B:85:THR:CG2	2.67	0.43
1:B:99:PHE:HD1	1:B:99:PHE:H	1.67	0.42
1:B:65:ASN:ND2	2:B:207:HOH:O	2.23	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:132:GLN:HG3	1:B:134:GLY:HA3	2.00	0.42
1:B:146:ARG:HD3	1:B:150:SER:CB	2.36	0.42
1:B:41:ARG:NH1	1:B:41:ARG:CG	2.60	0.42
1:B:132:GLN:HA	1:B:140:TYR:CB	2.49	0.42
1:B:82:ILE:HG22	1:B:128:PHE:HD2	1.85	0.41
1:B:98:LEU:CD2	1:B:108:ALA:O	2.69	0.41
1:B:2:ASN:N	1:B:3:ASP:HA	2.35	0.41
1:B:99:PHE:C	1:B:109:ARG:HH22	2.22	0.41
1:B:100:LYS:HE3	2:B:256:HOH:O	2.15	0.41
1:B:86:GLU:HG2	2:B:206:HOH:O	2.20	0.41
1:B:91:ALA:HB3	1:B:112:ILE:HD11	2.02	0.41
1:B:144:PHE:C	1:B:152:CYS:HB3	2.37	0.41
1:B:118:HIS:CG	1:B:119:PRO:N	2.90	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	B	183/185 (99%)	159 (87%)	20 (11%)	4 (2%)	8	28

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	30	PRO
1	B	71	PRO
1	B	118	HIS
1	B	131	GLU

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	B	144/146 (99%)	107 (74%)	37 (26%)	0 2

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	3	ASP
1	B	9	ASP
1	B	16	LEU
1	B	21	TYR
1	B	24	MET
1	B	28	TRP
1	B	35	LEU
1	B	36	ARG
1	B	40	THR
1	B	41	ARG
1	B	43	LEU
1	B	46	PRO
1	B	50	LEU
1	B	52	ASP
1	B	58	ASN
1	B	65	ASN
1	B	66	ILE
1	B	67	ARG
1	B	72	ARG
1	B	77	ASP
1	B	80	LEU
1	B	90	CYS
1	B	100	LYS
1	B	103	ASP
1	B	106	HIS
1	B	109	ARG
1	B	112	ILE
1	B	128	PHE
1	B	132	GLN
1	B	133	HIS

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Mol	Chain	Res	Type
1	B	135	LEU
1	B	137	ASN
1	B	140	TYR
1	B	152	CYS
1	B	158	TYR
1	B	162	GLU
1	B	164	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	27	ASN
1	B	132	GLN
1	B	160	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 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 [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	B	185/185 (100%)	0.51	15 (8%) 15 7	7, 23, 37, 44	0

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	117	GLY	7.9
1	B	41	ARG	3.8
1	B	163	GLY	3.0
1	B	115	SER	2.8
1	B	44	ASN	2.6
1	B	74	ILE	2.6
1	B	108	ALA	2.5
1	B	40	THR	2.5
1	B	137	ASN	2.4
1	B	90	CYS	2.2
1	B	184	GLY	2.2
1	B	145	CYS	2.2
1	B	113	GLY	2.0
1	B	38	GLY	2.0
1	B	26	SER	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

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.