



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 10:00 AM UTC

PDB ID : 7BU9 / pdb_00007bu9
Title : Crystal Structure of Spindlin1-H3(K4me3-K9me2) complex
Authors : Zhao, F.; Li, H.
Deposited on : 2020-04-05
Resolution : 3.50 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

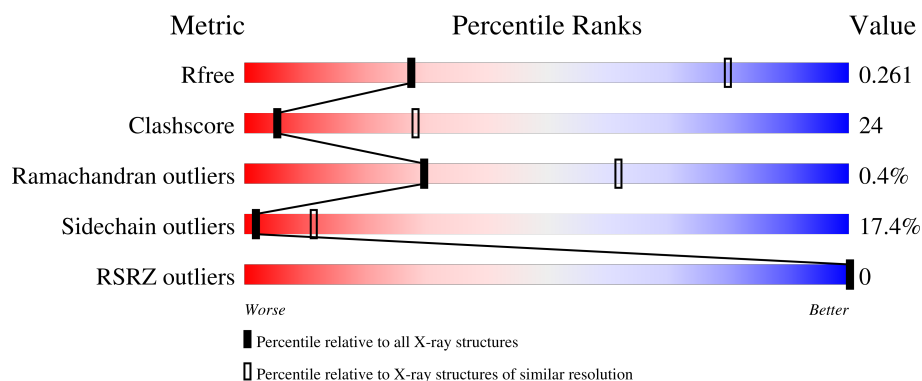
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.50 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	1085 (3.54-3.46)
Clashscore	190562	1140 (3.54-3.46)
Ramachandran outliers	187476	1113 (3.54-3.46)
Sidechain outliers	187428	1114 (3.54-3.46)
RSRZ outliers	180081	1084 (3.54-3.46)

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	220	
1	C	220	
1	E	220	
1	G	220	
2	B	15	

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Mol	Chain	Length	Quality of chain
2	D	15	
2	F	15	
2	H	15	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	M3L	D	4	-	-	X	-
2	MLY	D	9	-	-	X	-
2	M3L	F	4	-	-	X	-
2	MLY	F	9	-	-	X	-
2	M3L	H	4	-	-	X	-

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6911 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 Spindlin-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	204	Total	C	N	O	S	0	0	0
			1646	1051	273	313	9			
1	C	205	Total	C	N	O	S	0	0	0
			1655	1056	274	316	9			
1	E	206	Total	C	N	O	S	0	0	0
			1662	1060	276	317	9			
1	G	200	Total	C	N	O	S	0	0	0
			1618	1035	269	305	9			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	43	GLY	-	expression tag	UNP Q9Y657
A	44	SER	-	expression tag	UNP Q9Y657
C	43	GLY	-	expression tag	UNP Q9Y657
C	44	SER	-	expression tag	UNP Q9Y657
E	43	GLY	-	expression tag	UNP Q9Y657
E	44	SER	-	expression tag	UNP Q9Y657
G	43	GLY	-	expression tag	UNP Q9Y657
G	44	SER	-	expression tag	UNP Q9Y657

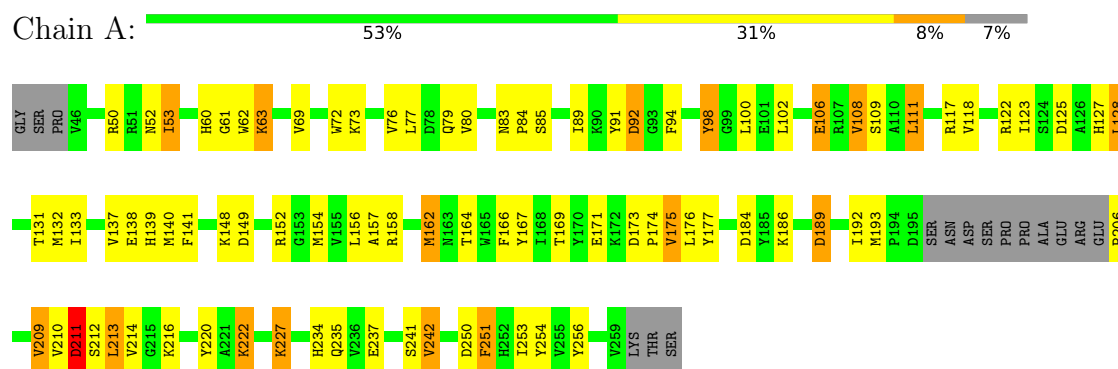
- Molecule 2 is a protein called H3(K4me3-K9me2) peptide.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	B	9	Total	C	N	O	0	0	0
			78	48	18	12			
2	D	10	Total	C	N	O	0	0	0
			84	51	19	14			
2	F	10	Total	C	N	O	0	0	0
			84	51	19	14			
2	H	10	Total	C	N	O	0	0	0
			84	51	19	14			

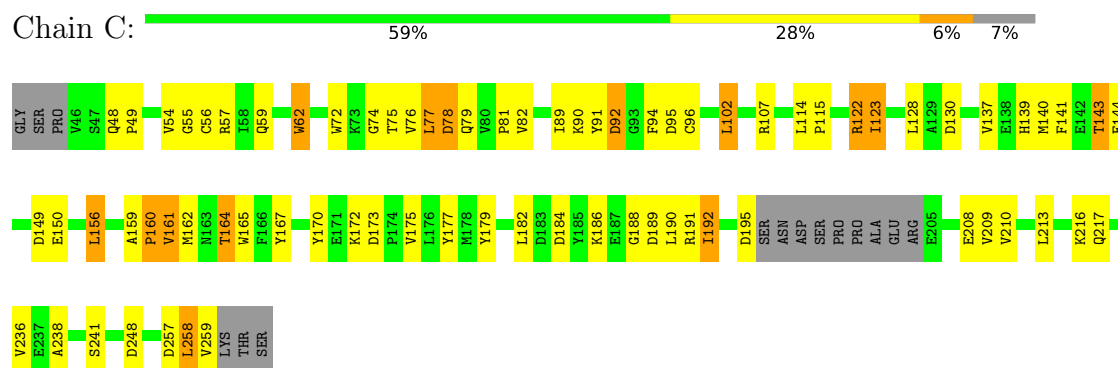
3 Residue-property plots [i](#)

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.

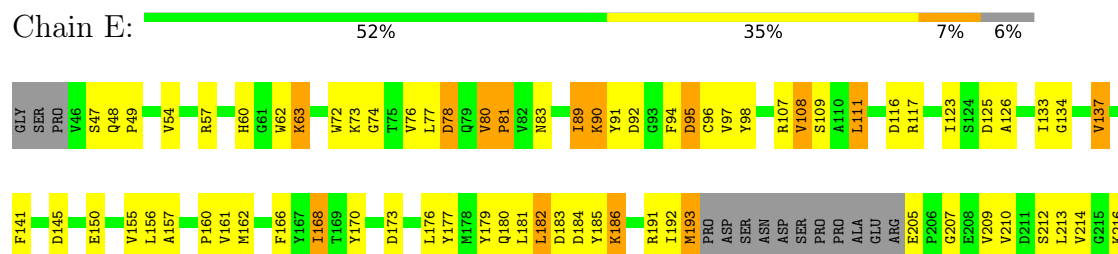
• Molecule 1: Spindlin-1



• Molecule 1: Spindlin-1



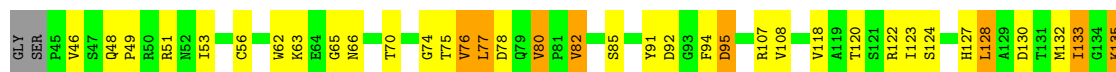
• Molecule 1: Spindlin-1





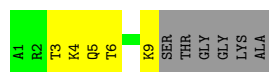
• Molecule 1: Spindlin-1

Chain G: 55% 28% 9% 9%



• Molecule 2: H3(K4me3-K9me2) peptide

Chain B: 27% 33% 40%



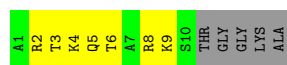
• Molecule 2: H3(K4me3-K9me2) peptide

Chain D: 27% 33% 7% 33%



• Molecule 2: H3(K4me3-K9me2) peptide

Chain F: 20% 47% 33%



• Molecule 2: H3(K4me3-K9me2) peptide

Chain H: 53% 13% 33%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	40.47Å 143.78Å 129.99Å 90.00° 95.86° 90.00°	Depositor
Resolution (Å)	48.07 – 3.50 48.07 – 3.50	Depositor EDS
% Data completeness (in resolution range)	98.7 (48.07-3.50) 84.6 (48.07-3.50)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.41 (at 3.48Å)	Xtriage
Refinement program	PHENIX 1.14_3247	Depositor
R, R_{free}	0.246 , 0.264 0.251 , 0.261	Depositor DCC
R_{free} test set	876 reflections (4.68%)	wwPDB-VP
Wilson B-factor (Å ²)	71.6	Xtriage
Anisotropy	1.095	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 194.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.36$, $\langle L^2 \rangle = 0.19$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	6911	wwPDB-VP
Average B, all atoms (Å ²)	80.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.38% 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: MLY, M3L

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	0.47	0/1685	0.80	3/2283 (0.1%)
1	C	0.42	0/1694	0.69	2/2296 (0.1%)
1	E	0.37	0/1700	0.64	0/2302
1	G	0.50	0/1656	0.82	1/2243 (0.0%)
2	B	0.15	0/54	0.46	0/71
2	D	0.78	0/60	1.10	0/79
2	F	0.62	0/60	1.03	0/79
2	H	0.73	0/60	1.14	0/79
All	All	0.45	0/6969	0.75	6/9432 (0.1%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	122	ARG	CA-C-N	6.04	128.60	120.50
1	A	122	ARG	C-N-CA	6.04	128.60	120.50
1	C	160	PRO	N-CA-C	5.95	124.72	112.47
1	C	143	THR	CB-CA-C	5.53	122.99	110.67
1	A	211	ASP	CB-CA-C	-5.41	109.87	117.23
1	G	75	THR	CB-CA-C	-5.08	102.43	110.81

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	1646	0	1601	81	0
1	C	1655	0	1606	91	0
1	E	1662	0	1620	93	0
1	G	1618	0	1581	83	0
2	B	78	0	95	11	0
2	D	84	0	99	27	0
2	F	84	0	100	29	0
2	H	84	0	100	9	0
All	All	6911	0	6802	328	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 (328) 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:C:141:PHE:CE2	2:D:4:M3L:HM32	1.71	1.23
1:E:72:TRP:CH2	1:E:94:PHE:HE2	1.61	1.18
1:E:207:GLY:HA2	1:G:158:ARG:HG3	1.39	1.03
1:E:72:TRP:HH2	1:E:94:PHE:CE2	1.77	1.00
1:E:72:TRP:CH2	1:E:94:PHE:CE2	2.50	0.99
1:G:213:LEU:CD1	1:G:243:TYR:CE2	2.44	0.98
1:C:77:LEU:HB3	1:C:161:VAL:CG2	1.94	0.98
1:C:123:ILE:O	1:G:124:SER:HA	1.62	0.98
1:G:213:LEU:HD12	1:G:243:TYR:CE2	2.00	0.97
1:A:220:TYR:O	1:A:227:LYS:HB3	1.68	0.94
1:C:141:PHE:CD2	2:D:4:M3L:HB3	2.03	0.92
1:C:141:PHE:CZ	2:D:4:M3L:HM32	2.04	0.92
1:G:213:LEU:HD12	1:G:213:LEU:H	1.34	0.92
2:H:4:M3L:HG3	2:H:4:M3L:HM33	1.52	0.91
1:A:213:LEU:HB3	1:C:81:PRO:HB2	1.52	0.90
1:C:141:PHE:HE2	2:D:4:M3L:HM32	1.14	0.90
1:C:141:PHE:CE2	2:D:4:M3L:HD3	2.06	0.89
1:G:135:LYS:HD3	1:G:192:ILE:HD11	1.54	0.89
1:E:62:TRP:CD1	2:F:9:MLY:CH2	2.57	0.88
1:E:72:TRP:HH2	1:E:94:PHE:HE2	0.89	0.88
1:E:78:ASP:OD2	1:G:210:VAL:HG11	1.78	0.84
1:G:56:CYS:O	1:G:76:VAL:HG12	1.78	0.83
1:C:62:TRP:CZ2	2:D:9:MLY:CH1	2.61	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:91:TYR:HE2	2:F:9:MLY:CH1	1.92	0.83
1:G:48:GLN:HG3	1:G:49:PRO:CD	2.09	0.82
1:C:55:GLY:HA2	1:C:161:VAL:HG11	1.62	0.82
1:C:95:ASP:OD2	1:C:179:TYR:HE1	1.63	0.81
1:C:122:ARG:HB2	1:G:124:SER:HB2	1.62	0.81
1:G:156:LEU:CD1	1:G:167:TYR:HB3	2.10	0.81
1:G:159:ALA:HB3	1:G:162:MET:O	1.81	0.80
1:E:91:TYR:HE2	2:F:9:MLY:HH12	1.47	0.80
1:C:62:TRP:CZ2	2:D:9:MLY:HH13	2.18	0.79
1:G:169:THR:HG21	1:G:255:VAL:HG11	1.65	0.79
1:E:168:ILE:HD12	1:E:181:LEU:HD11	1.64	0.79
1:A:212:SER:OG	1:A:214:VAL:HG23	1.84	0.78
1:G:48:GLN:HG3	1:G:49:PRO:HD2	1.66	0.77
1:C:95:ASP:OD2	1:C:179:TYR:CE1	2.37	0.77
1:C:141:PHE:HE2	2:D:4:M3L:HD3	1.49	0.76
1:E:62:TRP:CD1	2:F:9:MLY:HH23	2.21	0.76
1:A:213:LEU:HD12	1:A:213:LEU:H	1.51	0.75
1:G:213:LEU:HD12	1:G:213:LEU:N	2.02	0.75
1:C:173:ASP:OD1	2:D:8:ARG:NH1	2.20	0.74
1:E:78:ASP:CG	1:G:210:VAL:HG11	2.12	0.74
1:A:61:GLY:N	1:A:109:SER:OG	2.21	0.74
1:A:213:LEU:C	1:A:213:LEU:HD13	2.13	0.73
1:C:76:VAL:O	1:C:161:VAL:HG21	1.87	0.73
1:G:213:LEU:CD1	1:G:243:TYR:CZ	2.72	0.72
1:C:144:GLU:OE1	1:C:144:GLU:N	2.21	0.72
1:C:77:LEU:HB3	1:C:161:VAL:HG23	1.73	0.71
1:G:144:GLU:OE1	1:G:144:GLU:N	2.18	0.70
1:G:156:LEU:HD21	1:G:169:THR:HG22	1.72	0.70
1:C:141:PHE:CD2	2:D:4:M3L:HD3	2.26	0.70
1:E:81:PRO:HG2	1:G:212:SER:O	1.92	0.70
1:G:213:LEU:HD12	1:G:243:TYR:HE2	1.57	0.69
1:G:156:LEU:HD12	1:G:167:TYR:HB3	1.74	0.69
1:A:213:LEU:O	1:A:213:LEU:HD22	1.93	0.69
1:C:141:PHE:HE2	2:D:4:M3L:CM3	1.99	0.69
1:E:184:ASP:OD1	2:F:2:ARG:NH1	2.26	0.69
2:H:4:M3L:HM33	2:H:4:M3L:CG	2.22	0.69
1:G:132:MET:HE2	1:G:166:PHE:CE2	2.28	0.68
1:G:123:ILE:HG12	1:G:182:LEU:HD13	1.75	0.68
1:G:141:PHE:CZ	2:H:4:M3L:CM3	2.76	0.68
1:G:94:PHE:CE2	2:H:9:MLY:HA	2.28	0.68
1:E:137:VAL:HG11	1:E:155:VAL:HG23	1.75	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:4:M3L:O	2:D:4:M3L:HG2	1.95	0.67
1:A:60:HIS:ND1	1:A:108:VAL:HG22	2.09	0.67
1:G:63:LYS:HE3	1:G:66:ASN:H	1.59	0.67
1:G:213:LEU:HD11	1:G:243:TYR:CZ	2.30	0.67
1:E:62:TRP:CE2	2:F:9:MLY:HH21	2.30	0.66
1:A:60:HIS:ND1	1:A:108:VAL:CG2	2.60	0.65
1:A:213:LEU:HB3	1:C:81:PRO:CB	2.25	0.65
1:E:245:ILE:HD11	1:E:258:LEU:HD11	1.79	0.65
1:G:51:ARG:HG2	1:G:51:ARG:HH21	1.61	0.65
1:C:77:LEU:HB3	1:C:161:VAL:HG22	1.75	0.65
1:A:72:TRP:CH2	1:A:94:PHE:CZ	2.85	0.65
1:A:186:LYS:HG2	1:E:186:LYS:O	1.97	0.65
1:G:133:ILE:CD1	1:G:157:ALA:HA	2.27	0.65
1:C:48:GLN:HB3	1:C:49:PRO:HD2	1.79	0.65
1:E:91:TYR:CE2	2:F:9:MLY:CH1	2.80	0.64
1:E:166:PHE:N	1:E:179:TYR:O	2.29	0.64
1:G:220:TYR:HB2	1:G:228:ARG:HB2	1.78	0.64
1:A:156:LEU:HD21	1:A:169:THR:HG22	1.80	0.64
1:A:234:HIS:NE2	1:C:208:GLU:OE2	2.22	0.64
1:C:122:ARG:CB	1:G:124:SER:HB2	2.28	0.63
1:C:139:HIS:ND1	1:C:170:TYR:OH	2.29	0.63
1:E:78:ASP:OD2	1:G:210:VAL:CG1	2.47	0.62
1:C:77:LEU:CB	1:C:161:VAL:CG2	2.76	0.62
1:C:91:TYR:CE2	2:D:9:MLY:HH22	2.34	0.62
2:D:2:ARG:CG	2:D:2:ARG:HH11	2.12	0.62
1:G:132:MET:CE	1:G:166:PHE:CE2	2.83	0.62
1:E:95:ASP:N	1:E:95:ASP:OD1	2.33	0.62
1:C:123:ILE:HA	1:C:164:THR:HG23	1.80	0.61
1:E:80:VAL:HG12	1:E:83:ASN:H	1.66	0.61
1:E:207:GLY:HA2	1:G:158:ARG:CG	2.25	0.60
1:A:128:LEU:HG	1:E:185:TYR:OH	2.01	0.60
1:C:137:VAL:HG12	1:C:192:ILE:HA	1.82	0.60
1:E:62:TRP:NE1	2:F:9:MLY:HH23	2.15	0.60
1:G:141:PHE:CZ	2:H:4:M3L:HM33	2.36	0.60
1:E:62:TRP:CG	2:F:9:MLY:CH2	2.84	0.60
1:C:77:LEU:CB	1:C:161:VAL:HG22	2.31	0.59
1:A:61:GLY:C	1:A:109:SER:HG	2.09	0.59
1:E:57:ARG:NH2	1:E:92:ASP:OD2	2.35	0.59
1:E:259:VAL:HG13	1:E:259:VAL:O	2.01	0.59
1:E:72:TRP:CZ3	1:E:94:PHE:CE2	2.90	0.59
1:A:157:ALA:HB3	1:C:208:GLU:HG3	1.85	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:2:ARG:HH11	2:D:2:ARG:HG2	1.66	0.59
1:G:210:VAL:HG22	1:G:210:VAL:O	2.02	0.59
1:E:212:SER:HB2	1:E:214:VAL:HG23	1.84	0.58
1:E:62:TRP:CD2	2:F:9:MLY:HH21	2.39	0.58
1:A:213:LEU:CB	1:C:81:PRO:HB2	2.29	0.58
1:G:48:GLN:CG	1:G:49:PRO:HD2	2.33	0.58
1:C:94:PHE:CE2	2:D:9:MLY:HA	2.39	0.57
1:A:137:VAL:HG12	1:A:192:ILE:HG13	1.85	0.57
1:C:189:ASP:OD2	2:D:1:ALA:N	2.32	0.57
1:E:62:TRP:NE1	2:F:9:MLY:CH2	2.66	0.57
1:C:241:SER:OG	1:C:257:ASP:OD1	2.23	0.57
1:A:213:LEU:H	1:A:213:LEU:CD1	2.16	0.57
1:A:98:TYR:CD1	1:A:98:TYR:N	2.72	0.56
1:G:135:LYS:CD	1:G:192:ILE:HD11	2.32	0.56
1:G:156:LEU:HD21	1:G:169:THR:CG2	2.35	0.56
1:G:46:VAL:HG12	1:G:46:VAL:O	2.04	0.56
1:A:212:SER:HB2	1:A:235:GLN:OE1	2.05	0.56
1:E:77:LEU:HD22	1:E:160:PRO:HD2	1.87	0.56
1:A:89:ILE:HD11	1:A:102:LEU:HD21	1.88	0.56
1:A:210:VAL:HG21	1:C:79:GLN:O	2.06	0.56
1:C:96:CYS:HA	1:C:177:TYR:HD1	1.71	0.56
1:C:77:LEU:HD11	1:C:90:LYS:HB3	1.87	0.55
1:A:133:ILE:HG23	1:A:157:ALA:HA	1.89	0.55
1:C:77:LEU:HA	1:C:161:VAL:HG22	1.87	0.55
1:A:250:ASP:OD2	1:A:254:TYR:OH	2.22	0.55
1:E:78:ASP:CG	1:G:210:VAL:CG1	2.79	0.55
1:G:141:PHE:CE2	2:H:4:M3L:HG3	2.41	0.55
1:G:48:GLN:HG3	1:G:49:PRO:HD3	1.87	0.55
1:A:62:TRP:CD1	2:B:9:MLY:HH11	2.42	0.54
1:A:72:TRP:CH2	1:A:94:PHE:HZ	2.24	0.54
2:H:4:M3L:CG	2:H:4:M3L:CM3	2.86	0.54
1:E:177:TYR:CE2	2:F:4:M3L:HM11	2.43	0.54
1:C:159:ALA:HA	1:C:167:TYR:CE2	2.43	0.54
1:A:133:ILE:HD12	1:A:157:ALA:HA	1.89	0.54
1:E:60:HIS:HB3	1:E:111:LEU:HA	1.90	0.54
1:A:139:HIS:NE2	1:A:184:ASP:OD2	2.30	0.53
1:G:192:ILE:HD13	1:G:193:MET:N	2.22	0.53
1:A:72:TRP:CH2	1:A:94:PHE:CE2	2.95	0.53
1:A:89:ILE:HD12	1:A:100:LEU:HD12	1.90	0.53
1:E:210:VAL:HG11	1:G:80:VAL:HG13	1.88	0.53
1:G:136:ALA:HA	1:G:154:MET:HA	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:177:TYR:HE2	2:F:4:M3L:HM11	1.73	0.53
1:G:80:VAL:HG12	1:G:82:VAL:HG23	1.91	0.53
1:C:91:TYR:HE2	2:D:9:MLY:HH22	1.73	0.53
1:E:91:TYR:CE2	2:F:9:MLY:HH13	2.44	0.53
1:G:94:PHE:HE2	2:H:9:MLY:HA	1.74	0.53
1:G:95:ASP:HB2	1:G:178:MET:O	2.09	0.53
1:A:140:MET:HB3	1:A:189:ASP:HB3	1.91	0.53
1:G:248:ASP:N	1:G:248:ASP:OD1	2.42	0.53
1:C:91:TYR:CE2	2:D:9:MLY:CH2	2.91	0.52
1:A:77:LEU:HD21	1:A:162:MET:HE3	1.91	0.52
1:E:91:TYR:CE2	2:F:9:MLY:HH12	2.36	0.52
2:F:3:THR:HG22	2:F:5:GLN:H	1.74	0.52
1:C:141:PHE:CE2	2:D:4:M3L:HB3	2.43	0.52
1:A:158:ARG:HA	1:A:166:PHE:HD1	1.74	0.52
1:E:96:CYS:HA	1:E:177:TYR:HD1	1.73	0.52
1:G:48:GLN:CG	1:G:49:PRO:CD	2.84	0.52
1:A:94:PHE:CE2	2:B:9:MLY:HB2	2.44	0.52
1:E:233:ILE:HG12	1:E:244:PHE:HB2	1.92	0.52
1:C:77:LEU:HA	1:C:161:VAL:CG2	2.39	0.52
1:A:173:ASP:OD2	2:B:4:M3L:HM31	2.10	0.52
1:C:62:TRP:CE2	2:D:9:MLY:CH1	2.93	0.51
1:G:133:ILE:HD12	1:G:157:ALA:HA	1.92	0.51
1:E:78:ASP:OD1	1:G:210:VAL:HG13	2.10	0.51
1:G:213:LEU:HD13	1:G:243:TYR:CE2	2.42	0.51
1:E:123:ILE:HG21	1:E:126:ALA:HB2	1.91	0.51
1:A:73:LYS:O	1:A:92:ASP:OD1	2.28	0.51
1:A:117:ARG:CZ	1:A:117:ARG:HB3	2.39	0.51
1:C:57:ARG:HH12	1:C:92:ASP:CG	2.19	0.51
1:C:94:PHE:CZ	2:D:9:MLY:HA	2.46	0.51
1:G:156:LEU:HD12	1:G:167:TYR:CB	2.41	0.51
1:E:182:LEU:C	1:E:182:LEU:HD22	2.35	0.51
1:A:98:TYR:CE2	2:B:9:MLY:HE2	2.46	0.51
1:C:156:LEU:HD12	1:C:167:TYR:HB3	1.93	0.51
1:E:156:LEU:HD11	1:E:255:VAL:HG21	1.92	0.51
1:G:77:LEU:HD11	1:G:178:MET:SD	2.51	0.51
1:G:51:ARG:HG2	1:G:51:ARG:NH2	2.26	0.50
1:C:74:GLY:HA3	1:C:91:TYR:HD1	1.76	0.50
1:A:79:GLN:NE2	1:A:84:PRO:O	2.43	0.50
1:A:94:PHE:HA	2:B:6:THR:O	2.12	0.50
1:C:95:ASP:CG	1:C:179:TYR:CE1	2.89	0.50
2:D:2:ARG:CG	2:D:2:ARG:NH1	2.73	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:248:ASP:N	1:E:248:ASP:OD1	2.44	0.50
1:A:141:PHE:HE2	2:B:4:M3L:HM23	1.77	0.49
1:A:80:VAL:HG12	1:C:210:VAL:HG11	1.94	0.49
1:C:89:ILE:HD11	1:C:102:LEU:HD11	1.93	0.49
1:E:62:TRP:CD1	2:F:9:MLY:HH22	2.46	0.49
1:G:137:VAL:HG12	1:G:192:ILE:HA	1.94	0.49
1:C:62:TRP:CE2	2:D:9:MLY:HH12	2.48	0.49
1:A:61:GLY:CA	1:A:109:SER:OG	2.60	0.49
1:E:170:TYR:CD2	2:F:4:M3L:HM21	2.47	0.49
1:G:211:ASP:OD1	1:G:211:ASP:N	2.46	0.49
1:E:182:LEU:C	1:E:182:LEU:CD2	2.85	0.49
1:G:167:TYR:CD1	1:G:176:LEU:HD11	2.48	0.49
1:A:213:LEU:CD1	1:A:213:LEU:N	2.76	0.49
1:C:55:GLY:HA2	1:C:161:VAL:CG1	2.39	0.48
1:C:77:LEU:CA	1:C:161:VAL:HG22	2.43	0.48
1:E:173:ASP:OD2	2:F:4:M3L:CM1	2.61	0.48
1:A:72:TRP:HH2	1:A:94:PHE:HZ	1.59	0.48
1:G:63:LYS:HG2	1:G:65:GLY:H	1.79	0.48
1:A:189:ASP:OD1	1:A:189:ASP:N	2.46	0.48
1:E:150:GLU:OE1	1:E:191:ARG:NH2	2.35	0.48
1:G:56:CYS:O	1:G:76:VAL:CG1	2.56	0.48
1:A:222:LYS:HA	1:A:222:LYS:HD3	1.62	0.48
1:G:74:GLY:HA3	1:G:91:TYR:HA	1.96	0.48
1:A:251:PHE:CD1	1:A:251:PHE:N	2.82	0.48
1:E:78:ASP:OD1	1:G:210:VAL:CG1	2.62	0.48
1:A:209:VAL:HG13	1:C:160:PRO:HB3	1.94	0.48
1:E:98:TYR:CZ	2:F:9:MLY:HG2	2.49	0.48
1:E:173:ASP:OD2	2:F:4:M3L:HM11	2.14	0.48
1:E:182:LEU:HD22	1:E:182:LEU:O	2.14	0.48
1:E:262:SER:O	1:E:262:SER:OG	2.31	0.48
1:E:168:ILE:CD1	1:E:181:LEU:HD11	2.40	0.47
1:C:90:LYS:HE2	1:C:94:PHE:O	2.15	0.47
1:C:140:MET:SD	1:C:150:GLU:OE1	2.72	0.47
1:A:127:HIS:O	1:A:131:THR:HG23	2.14	0.47
1:E:177:TYR:CE2	2:F:4:M3L:HM31	2.49	0.47
1:G:156:LEU:CD2	1:G:169:THR:HG22	2.44	0.47
1:E:62:TRP:CE2	2:F:9:MLY:CH2	2.96	0.47
1:A:158:ARG:HA	1:A:166:PHE:CD1	2.50	0.47
1:A:175:VAL:HG23	1:A:177:TYR:CE2	2.50	0.47
1:C:114:LEU:HD23	1:C:115:PRO:HD2	1.96	0.47
1:E:123:ILE:CG2	1:E:126:ALA:HB2	2.44	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:60:HIS:O	1:E:109:SER:O	2.33	0.47
1:G:167:TYR:HD1	1:G:176:LEU:HD11	1.80	0.47
1:C:62:TRP:HZ3	1:C:107:ARG:HD2	1.79	0.46
1:G:245:ILE:HD12	1:G:256:TYR:HD2	1.79	0.46
1:C:90:LYS:HE3	1:C:165:TRP:CZ3	2.50	0.46
1:C:160:PRO:HD3	1:C:167:TYR:CE2	2.49	0.46
1:G:62:TRP:CH2	1:G:107:ARG:HD2	2.50	0.46
1:E:73:LYS:O	1:E:92:ASP:OD1	2.33	0.46
1:E:90:LYS:HE3	1:E:94:PHE:O	2.15	0.46
1:A:213:LEU:HD23	1:A:216:LYS:HD3	1.98	0.46
1:C:48:GLN:HB3	1:C:49:PRO:CD	2.46	0.46
1:C:92:ASP:N	1:C:92:ASP:OD1	2.48	0.46
1:E:62:TRP:CG	2:F:9:MLY:HH22	2.50	0.46
1:E:48:GLN:HG3	1:E:49:PRO:HD2	1.97	0.46
1:E:251:PHE:O	1:E:251:PHE:CD2	2.69	0.46
1:A:212:SER:CB	1:A:235:GLN:OE1	2.65	0.45
1:C:72:TRP:CH2	2:D:9:MLY:HD3	2.51	0.45
1:C:141:PHE:CZ	2:D:4:M3L:CM3	2.91	0.45
1:E:96:CYS:HB3	1:E:98:TYR:CE1	2.51	0.45
1:E:185:TYR:CG	1:E:185:TYR:O	2.69	0.45
1:E:185:TYR:O	1:E:185:TYR:CD1	2.70	0.45
1:A:154:MET:HG2	1:A:156:LEU:HD23	1.99	0.45
1:A:213:LEU:HD12	1:A:213:LEU:N	2.24	0.45
1:A:152:ARG:HD3	1:A:193:MET:HE1	1.97	0.45
1:C:62:TRP:O	1:C:62:TRP:CD1	2.69	0.45
1:C:139:HIS:NE2	1:C:184:ASP:OD2	2.48	0.45
1:E:213:LEU:HB3	1:E:216:LYS:HD3	1.97	0.45
1:A:210:VAL:CG2	1:C:81:PRO:HD3	2.47	0.45
1:E:63:LYS:HG3	1:E:107:ARG:O	2.17	0.45
1:E:141:PHE:CE2	2:F:4:M3L:HM32	2.52	0.45
1:A:210:VAL:HG21	1:C:81:PRO:HD3	1.98	0.45
1:C:90:LYS:HE3	1:C:165:TRP:HZ3	1.82	0.44
1:G:184:ASP:HB3	1:G:189:ASP:HB2	1.99	0.44
1:C:236:VAL:HG12	1:C:238:ALA:H	1.83	0.44
1:E:141:PHE:HE2	2:F:4:M3L:HM32	1.82	0.44
1:A:53:ILE:HD13	1:A:53:ILE:H	1.82	0.44
2:B:3:THR:HG22	2:B:5:GLN:H	1.82	0.44
1:C:75:THR:HB	1:C:162:MET:CE	2.47	0.44
1:A:125:ASP:OD2	1:E:193:MET:N	2.50	0.44
1:A:211:ASP:O	1:C:81:PRO:HG2	2.16	0.44
1:C:57:ARG:NH1	1:C:75:THR:OG1	2.49	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:74:GLY:HA2	1:E:92:ASP:OD1	2.17	0.44
1:C:188:GLY:O	1:C:191:ARG:NH1	2.50	0.44
1:G:228:ARG:NH1	1:G:256:TYR:OH	2.50	0.44
1:A:141:PHE:CE2	2:B:4:M3L:HM23	2.53	0.44
2:F:9:MLY:HH12	2:F:9:MLY:HD2	1.64	0.44
1:E:76:VAL:HA	1:E:89:ILE:HG23	1.98	0.43
1:C:149:ASP:OD1	1:C:150:GLU:N	2.52	0.43
1:A:91:TYR:CE2	2:B:9:MLY:HE3	2.53	0.43
1:A:152:ARG:HH12	1:A:171:GLU:HG2	1.83	0.43
1:C:123:ILE:O	1:G:124:SER:O	2.35	0.43
1:E:176:LEU:HD22	1:E:253:ILE:HG21	2.01	0.43
1:E:60:HIS:ND1	1:E:108:VAL:CG1	2.81	0.43
1:A:111:LEU:HD13	1:A:111:LEU:HA	1.92	0.43
1:A:128:LEU:HD13	1:A:132:MET:HE2	2.01	0.43
1:C:123:ILE:O	1:G:124:SER:CA	2.50	0.43
1:E:63:LYS:O	1:E:107:ARG:HB3	2.19	0.43
1:A:213:LEU:HD23	1:A:216:LYS:CD	2.49	0.42
1:E:168:ILE:HG22	1:E:177:TYR:HB2	2.01	0.42
1:E:210:VAL:N	1:G:78:ASP:OD2	2.43	0.42
1:A:52:ASN:O	1:A:52:ASN:ND2	2.50	0.42
1:C:173:ASP:CG	2:D:8:ARG:HH11	2.26	0.42
1:E:185:TYR:CD1	1:E:185:TYR:C	2.97	0.42
1:G:170:TYR:CD2	2:H:4:M3L:HM21	2.54	0.42
1:A:106:GLU:N	1:A:106:GLU:OE2	2.52	0.42
1:C:213:LEU:HD22	1:C:258:LEU:HD12	2.00	0.42
1:A:242:VAL:HA	1:A:256:TYR:O	2.19	0.42
1:A:167:TYR:CD1	1:A:176:LEU:HD11	2.55	0.42
1:E:176:LEU:HD12	1:E:176:LEU:HA	1.74	0.42
1:C:248:ASP:OD1	1:C:248:ASP:N	2.53	0.42
1:G:219:GLU:OE1	1:G:227:LYS:HD3	2.19	0.42
1:G:227:LYS:HE2	1:G:227:LYS:HB3	1.83	0.42
1:C:77:LEU:CA	1:C:161:VAL:CG2	2.97	0.42
1:E:60:HIS:ND1	1:E:108:VAL:HG13	2.34	0.42
1:E:98:TYR:OH	2:F:9:MLY:HG2	2.20	0.42
1:A:213:LEU:HD13	1:A:213:LEU:O	2.18	0.41
1:C:213:LEU:HD23	1:C:216:LYS:HD3	2.01	0.41
1:G:95:ASP:N	1:G:95:ASP:OD1	2.50	0.41
1:G:184:ASP:OD1	1:G:184:ASP:N	2.52	0.41
1:A:253:ILE:HD13	1:A:253:ILE:HA	1.82	0.41
1:C:78:ASP:OD2	1:C:160:PRO:HG2	2.20	0.41
1:E:168:ILE:HD13	1:E:170:TYR:OH	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:237:GLU:HB3	1:C:208:GLU:OE1	2.20	0.41
1:C:62:TRP:HB2	1:C:107:ARG:O	2.21	0.41
1:E:180:GLN:C	1:E:182:LEU:N	2.79	0.41
1:C:122:ARG:HA	1:C:122:ARG:HD3	1.82	0.41
1:E:133:ILE:HG13	1:E:157:ALA:HA	2.02	0.41
1:E:244:PHE:CD1	1:E:255:VAL:HG22	2.55	0.41
1:G:246:LYS:HE2	1:G:247:PHE:O	2.21	0.41
1:A:212:SER:HB2	1:A:235:GLN:HB2	2.03	0.41
1:C:186:LYS:HA	1:C:186:LYS:HD2	1.89	0.40
1:G:249:ASP:OD1	1:G:249:ASP:N	2.53	0.40
2:B:9:MLY:HD3	2:B:9:MLY:HH13	1.85	0.40
1:A:106:GLU:OE2	1:A:106:GLU:CA	2.70	0.40
1:G:128:LEU:HD13	1:G:128:LEU:HA	1.78	0.40
1:G:186:LYS:HE3	1:G:186:LYS:HB2	1.78	0.40
1:E:95:ASP:OD1	2:F:6:THR:O	2.40	0.40
1:A:63:LYS:CG	1:A:69:VAL:HG22	2.51	0.40
1:A:91:TYR:HE2	2:B:9:MLY:HE3	1.87	0.40
1:E:134:GLY:HA3	1:E:238:ALA:HB1	2.03	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	200/220 (91%)	178 (89%)	21 (10%)	1 (0%)	24	57
1	C	201/220 (91%)	182 (90%)	19 (10%)	0	100	100
1	E	202/220 (92%)	181 (90%)	20 (10%)	1 (0%)	24	57
1	G	196/220 (89%)	183 (93%)	12 (6%)	1 (0%)	24	57
2	B	6/15 (40%)	6 (100%)	0	0	100	100
2	D	6/15 (40%)	6 (100%)	0	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	F	6/15 (40%)	6 (100%)	0	0	100	100
2	H	6/15 (40%)	5 (83%)	1 (17%)	0	100	100
All	All	823/940 (88%)	747 (91%)	73 (9%)	3 (0%)	30	62

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	174	PRO
1	E	81	PRO
1	G	160	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	180/194 (93%)	150 (83%)	30 (17%)	2	13
1	C	181/194 (93%)	154 (85%)	27 (15%)	3	17
1	E	182/194 (94%)	147 (81%)	35 (19%)	1	8
1	G	177/194 (91%)	143 (81%)	34 (19%)	1	8
2	B	5/8 (62%)	5 (100%)	0	100	100
2	D	6/8 (75%)	4 (67%)	2 (33%)	0	2
2	F	6/8 (75%)	5 (83%)	1 (17%)	2	13
2	H	6/8 (75%)	6 (100%)	0	100	100
All	All	743/808 (92%)	614 (83%)	129 (17%)	2	11

All (129) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	50	ARG
1	A	53	ILE
1	A	63	LYS
1	A	76	VAL

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Mol	Chain	Res	Type
1	A	83	ASN
1	A	85	SER
1	A	92	ASP
1	A	98	TYR
1	A	106	GLU
1	A	108	VAL
1	A	111	LEU
1	A	118	VAL
1	A	123	ILE
1	A	128	LEU
1	A	138	GLU
1	A	148	LYS
1	A	149	ASP
1	A	162	MET
1	A	164	THR
1	A	175	VAL
1	A	189	ASP
1	A	206	PRO
1	A	209	VAL
1	A	211	ASP
1	A	213	LEU
1	A	222	LYS
1	A	227	LYS
1	A	241	SER
1	A	242	VAL
1	A	251	PHE
1	C	54	VAL
1	C	56	CYS
1	C	59	GLN
1	C	62	TRP
1	C	77	LEU
1	C	78	ASP
1	C	82	VAL
1	C	92	ASP
1	C	102	LEU
1	C	122	ARG
1	C	123	ILE
1	C	128	LEU
1	C	130	ASP
1	C	143	THR
1	C	156	LEU
1	C	161	VAL

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Mol	Chain	Res	Type
1	C	164	THR
1	C	172	LYS
1	C	175	VAL
1	C	182	LEU
1	C	190	LEU
1	C	192	ILE
1	C	195	ASP
1	C	209	VAL
1	C	217	GLN
1	C	258	LEU
1	C	259	VAL
2	D	2	ARG
2	D	10	SER
1	E	47	SER
1	E	54	VAL
1	E	63	LYS
1	E	78	ASP
1	E	80	VAL
1	E	89	ILE
1	E	90	LYS
1	E	95	ASP
1	E	97	VAL
1	E	108	VAL
1	E	111	LEU
1	E	116	ASP
1	E	117	ARG
1	E	125	ASP
1	E	137	VAL
1	E	145	ASP
1	E	161	VAL
1	E	162	MET
1	E	168	ILE
1	E	182	LEU
1	E	183	ASP
1	E	186	LYS
1	E	192	ILE
1	E	193	MET
1	E	205	GLU
1	E	209	VAL
1	E	222	LYS
1	E	223	GLU
1	E	226	SER

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Mol	Chain	Res	Type
1	E	229	THR
1	E	232	VAL
1	E	235	GLN
1	E	242	VAL
1	E	252	HIS
1	E	259	VAL
2	F	8	ARG
1	G	53	ILE
1	G	70	THR
1	G	76	VAL
1	G	77	LEU
1	G	80	VAL
1	G	82	VAL
1	G	85	SER
1	G	92	ASP
1	G	95	ASP
1	G	108	VAL
1	G	118	VAL
1	G	120	THR
1	G	122	ARG
1	G	127	HIS
1	G	128	LEU
1	G	130	ASP
1	G	133	ILE
1	G	135	LYS
1	G	156	LEU
1	G	158	ARG
1	G	162	MET
1	G	164	THR
1	G	169	THR
1	G	192	ILE
1	G	209	VAL
1	G	210	VAL
1	G	212	SER
1	G	213	LEU
1	G	222	LYS
1	G	227	LYS
1	G	237	GLU
1	G	248	ASP
1	G	249	ASP
1	G	259	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such

sidechains are listed below:

Mol	Chain	Res	Type
1	A	180	GLN
2	B	5	GLN
1	C	52	ASN
1	C	234	HIS
1	E	163	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

8 non-standard protein/DNA/RNA residues are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	M3L	D	4	2	10,11,12	0.57	0	9,14,16	0.14	0
2	M3L	F	4	2	10,11,12	0.54	0	9,14,16	0.11	0
2	M3L	B	4	2	10,11,12	0.55	0	9,14,16	0.56	0
2	MLY	H	9	2	9,10,11	0.44	0	6,11,13	0.22	0
2	MLY	B	9	2	9,10,11	0.47	0	6,11,13	0.97	0
2	M3L	H	4	2	10,11,12	0.47	0	9,14,16	0.10	0
2	MLY	F	9	2	9,10,11	0.43	0	6,11,13	0.14	0
2	MLY	D	9	2	9,10,11	0.46	0	6,11,13	0.20	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	M3L	D	4	2	-	6/9/10/12	-
2	M3L	F	4	2	-	5/9/10/12	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	M3L	B	4	2	-	0/9/10/12	-
2	MLY	H	9	2	-	4/8/9/11	-
2	MLY	B	9	2	-	1/8/9/11	-
2	M3L	H	4	2	-	4/9/10/12	-
2	MLY	F	9	2	-	4/8/9/11	-
2	MLY	D	9	2	-	4/8/9/11	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (28) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	D	4	M3L	C-CA-CB-CG
2	H	4	M3L	N-CA-CB-CG
2	H	4	M3L	C-CA-CB-CG
2	H	9	MLY	C-CA-CB-CG
2	F	9	MLY	CE-CD-CG-CB
2	H	4	M3L	CG-CD-CE-NZ
2	H	9	MLY	CE-CD-CG-CB
2	D	9	MLY	CD-CE-NZ-CH1
2	D	9	MLY	CD-CE-NZ-CH2
2	F	9	MLY	CG-CD-CE-NZ
2	D	9	MLY	CG-CD-CE-NZ
2	F	9	MLY	CD-CE-NZ-CH1
2	H	9	MLY	CD-CE-NZ-CH2
2	H	4	M3L	CA-CB-CG-CD
2	D	4	M3L	CE-CD-CG-CB
2	D	9	MLY	CE-CD-CG-CB
2	F	4	M3L	CE-CD-CG-CB
2	F	4	M3L	CA-CB-CG-CD
2	F	4	M3L	CD-CE-NZ-CM3
2	F	4	M3L	CD-CE-NZ-CM2
2	D	4	M3L	CD-CE-NZ-CM2
2	F	4	M3L	CD-CE-NZ-CM1
2	D	4	M3L	CA-CB-CG-CD
2	D	4	M3L	CD-CE-NZ-CM1
2	D	4	M3L	CD-CE-NZ-CM3
2	B	9	MLY	N-CA-CB-CG
2	F	9	MLY	N-CA-CB-CG

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Mol	Chain	Res	Type	Atoms
2	H	9	MLY	N-CA-CB-CG

There are no ring outliers.

8 monomers are involved in 65 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	4	M3L	11	0
2	F	4	M3L	8	0
2	B	4	M3L	3	0
2	H	9	MLY	2	0
2	B	9	MLY	6	0
2	H	4	M3L	7	0
2	F	9	MLY	18	0
2	D	9	MLY	10	0

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	204/220 (92%)	-0.99	0 100 100	30, 79, 128, 176	0
1	C	205/220 (93%)	-0.97	0 100 100	30, 78, 131, 181	0
1	E	206/220 (93%)	-1.02	0 100 100	37, 74, 113, 181	0
1	G	200/220 (90%)	-1.04	0 100 100	30, 77, 132, 172	0
2	B	7/15 (46%)	-1.01	0 100 100	68, 84, 94, 101	0
2	D	8/15 (53%)	-1.03	0 100 100	79, 88, 108, 110	0
2	F	8/15 (53%)	-1.18	0 100 100	68, 75, 83, 96	0
2	H	8/15 (53%)	-1.09	0 100 100	62, 82, 90, 94	0
All	All	846/940 (90%)	-1.01	0 100 100	30, 78, 129, 181	0

There are no RSRZ outliers to report.

6.2 Non-standard residues in protein, DNA, RNA chains [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(Å ²)	Q<0.9
2	MLY	H	9	11/12	0.98	0.09	60,69,91,101	0
2	MLY	B	9	11/12	0.99	0.07	74,85,100,103	0
2	M3L	D	4	12/13	0.99	0.06	55,73,97,103	0
2	MLY	D	9	11/12	0.99	0.06	96,105,114,117	0
2	M3L	F	4	12/13	0.99	0.08	42,83,90,95	0
2	MLY	F	9	11/12	0.99	0.05	62,73,91,92	0
2	M3L	H	4	12/13	0.99	0.05	39,67,86,91	0
2	M3L	B	4	12/13	0.99	0.05	42,66,78,81	0

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.