



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 22, 2024 – 05:00 AM EDT

PDB ID : 4I80
Title : Crystal structure of human menin in complex with a high-affinity macrocyclic peptidomimetics
Authors : Huang, J.; Lei, M.
Deposited on : 2012-12-01
Resolution : 3.10 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 : 4.02b-467
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 1.20.1
EDS : 3.0
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4 : 9.0.003 (Gargrove)
Density-Fitness : 1.0.11
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.39

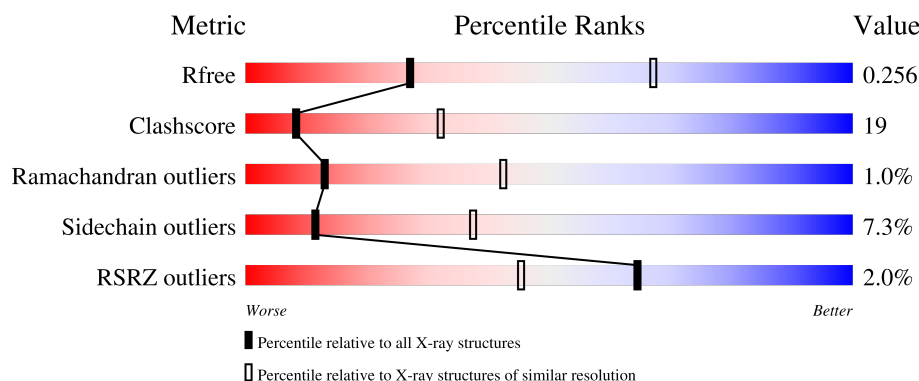
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.10 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}	164625	1351 (3.10-3.10)
Clashscore	180529	1454 (3.10-3.10)
Ramachandran outliers	177936	1391 (3.10-3.10)
Sidechain outliers	177891	1391 (3.10-3.10)
RSRZ outliers	164620	1351 (3.10-3.10)

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	550	
2	B	9	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3893 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 Menin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	484	Total	C	N	O	S	0	0	0
			3812	2442	653	703	14			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	SER	-	expression tag	UNP O00255

- Molecule 2 is a protein called macrocyclic peptidomimetic.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	B	9	Total	C	N	O	0	0	0
			81	56	16	9			

4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	141.38Å 141.38Å 92.88Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	44.71 – 3.10 44.71 – 3.10	Depositor EDS
% Data completeness (in resolution range)	92.2 (44.71-3.10) 92.5 (44.71-3.10)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	10.88 (at 3.12Å)	Xtriage
Refinement program	PHENIX 1.7_650	Depositor
R, R_{free}	0.215 , 0.245 0.229 , 0.256	Depositor DCC
R_{free} test set	1622 reflections (9.95%)	wwPDB-VP
Wilson B-factor (Å ²)	69.5	Xtriage
Anisotropy	1.193	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 94.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	3893	wwPDB-VP
Average B, all atoms (Å ²)	107.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.69% 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: ACE, 1E3

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	0.41	0/3900	0.58	0/5297
2	B	0.73	0/70	1.80	4/94 (4.3%)
All	All	0.42	0/3970	0.62	4/5391 (0.1%)

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
2	B	0	2

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
2	B	12	ARG	O-C-N	11.62	143.17	121.10
2	B	12	ARG	CA-C-N	-7.42	96.31	117.10
2	B	12	ARG	C-N-CD	7.26	143.65	128.40
2	B	12	ARG	C-N-CA	-5.48	99.00	122.00

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	7	TRP	Peptide
2	B	8	1E3	Mainchain

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	3812	0	3793	135	0
2	B	81	0	83	18	0
All	All	3893	0	3876	147	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 19.

All (147) 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:452:ARG:O	1:A:566:VAL:HG23	1.66	0.96
1:A:70:ASP:HB3	1:A:72:PRO:HD2	1.51	0.92
1:A:319:TYR:CD2	2:B:13:PRO:HG3	2.08	0.89
1:A:376:LEU:HB3	1:A:446:ARG:HG2	1.58	0.84
2:B:6:ARG:HB3	2:B:6:ARG:HH11	1.47	0.80
1:A:404:LEU:HD11	1:A:446:ARG:HG3	1.64	0.79
1:A:4:LYS:HD3	1:A:29:ARG:NH2	2.00	0.77
2:B:6:ARG:HB3	2:B:6:ARG:NH1	2.02	0.74
1:A:349:GLN:HB2	1:A:422:LYS:HB3	1.70	0.74
2:B:12:ARG:HG3	2:B:13:PRO:HD2	1.70	0.73
1:A:353:TYR:CE1	1:A:361:TYR:HB2	2.24	0.71
1:A:553:PHE:CE2	1:A:558:MET:HG3	2.27	0.70
1:A:345:ALA:HA	1:A:348:ILE:HG22	1.73	0.69
1:A:284:ALA:HB1	1:A:302:TYR:CE1	2.27	0.69
2:B:12:ARG:CG	2:B:13:PRO:HD2	2.23	0.68
1:A:334:VAL:O	1:A:338:LEU:HD12	1.93	0.68
1:A:259:LEU:O	1:A:259:LEU:HD12	1.94	0.67
2:B:12:ARG:HG2	2:B:13:PRO:N	2.10	0.67
1:A:23:PHE:O	1:A:27:LEU:HB2	1.95	0.66
1:A:319:TYR:CD2	2:B:13:PRO:CG	2.79	0.65
1:A:417:TYR:C	1:A:558:MET:HE3	2.18	0.64
1:A:130:SER:N	1:A:150:THR:HG22	2.13	0.63
1:A:185:VAL:HG12	1:A:193:THR:HG22	1.81	0.62
1:A:328:HIS:CD2	1:A:336:GLU:HB3	2.33	0.62
1:A:4:LYS:HD3	1:A:29:ARG:HH22	1.63	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:335:ARG:HG2	1:A:409:CYS:SG	2.39	0.62
1:A:136:ASP:O	2:B:7:TRP:HB3	1.98	0.62
1:A:137:ARG:HA	2:B:6:ARG:HH12	1.65	0.62
1:A:205:ASP:HB2	1:A:207:ARG:HG3	1.82	0.62
1:A:281:GLY:O	1:A:284:ALA:HB3	2.03	0.59
1:A:57:ASN:O	1:A:59:PRO:HD3	2.02	0.59
1:A:150:THR:HG22	1:A:150:THR:O	2.03	0.59
1:A:454:LYS:O	1:A:549:PRO:HD2	2.03	0.59
1:A:410:PHE:CE2	1:A:414:LEU:HD11	2.38	0.58
1:A:243:ILE:O	1:A:245:PRO:HD3	2.05	0.57
1:A:58:VAL:HG12	1:A:60:GLU:HG2	1.88	0.56
1:A:108:ARG:HB3	1:A:112:VAL:O	2.05	0.56
1:A:40:VAL:O	1:A:44:VAL:HG23	2.05	0.55
2:B:12:ARG:CG	2:B:13:PRO:CD	2.85	0.55
1:A:104:SER:HB2	1:A:105:LEU:HD23	1.88	0.55
1:A:183:TRP:HZ2	1:A:227:TYR:CD2	2.25	0.55
1:A:306:ILE:HD13	1:A:321:TYR:CE2	2.42	0.55
1:A:70:ASP:CB	1:A:72:PRO:HD2	2.32	0.54
1:A:78:PHE:CD2	1:A:79:PRO:HD2	2.42	0.54
1:A:363:GLU:O	1:A:367:VAL:HG23	2.08	0.54
1:A:353:TYR:HE1	1:A:361:TYR:HB2	1.71	0.53
1:A:31:GLU:OE1	1:A:232:ARG:HG3	2.09	0.53
1:A:177:LEU:HD21	1:A:238:PHE:CD2	2.45	0.52
1:A:346:THR:O	1:A:349:GLN:HG2	2.10	0.52
2:B:12:ARG:CG	2:B:13:PRO:N	2.72	0.52
1:A:318:ILE:HG23	1:A:344:THR:HG23	1.91	0.52
1:A:383:LEU:HD12	1:A:384:GLU:N	2.25	0.52
1:A:551:LEU:CD1	1:A:553:PHE:CE1	2.94	0.51
1:A:105:LEU:HD23	1:A:105:LEU:N	2.26	0.51
1:A:204:GLU:OE1	1:A:204:GLU:HA	2.10	0.51
2:B:6:ARG:HH11	2:B:6:ARG:CB	2.19	0.51
1:A:410:PHE:O	1:A:414:LEU:HG	2.11	0.51
1:A:137:ARG:HA	2:B:6:ARG:NH1	2.25	0.50
1:A:159:PHE:HZ	1:A:239:MET:CE	2.24	0.50
1:A:248:ASP:HB3	1:A:251:THR:HB	1.93	0.50
1:A:455:VAL:HG22	1:A:549:PRO:HG2	1.92	0.50
1:A:175:LEU:HG	1:A:177:LEU:HD13	1.94	0.50
1:A:149:GLY:O	1:A:150:THR:HB	2.12	0.50
1:A:159:PHE:CZ	1:A:239:MET:HE2	2.47	0.50
1:A:424:GLU:HG3	1:A:431:VAL:H	1.76	0.50
1:A:183:TRP:HB2	1:A:194:ALA:O	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:202:GLY:C	1:A:204:GLU:H	2.14	0.49
1:A:140:ILE:O	1:A:152:LEU:HA	2.13	0.49
2:B:12:ARG:HG2	2:B:13:PRO:CD	2.43	0.49
1:A:130:SER:H	1:A:150:THR:HG22	1.78	0.48
1:A:44:VAL:HG12	1:A:48:LEU:HD12	1.94	0.48
1:A:102:ASP:C	1:A:104:SER:H	2.17	0.48
1:A:128:SER:O	1:A:150:THR:HG23	2.13	0.48
1:A:173:VAL:O	1:A:232:ARG:NH2	2.46	0.48
1:A:221:LEU:HD23	1:A:221:LEU:HA	1.55	0.48
1:A:367:VAL:HG12	1:A:372:ILE:HD12	1.95	0.48
1:A:16:ILE:O	1:A:20:VAL:HG23	2.14	0.47
1:A:551:LEU:HD11	1:A:553:PHE:HE1	1.79	0.47
1:A:31:GLU:OE1	1:A:31:GLU:HA	2.15	0.47
1:A:410:PHE:CE1	1:A:452:ARG:HG2	2.49	0.47
1:A:265:TRP:O	1:A:266:LEU:C	2.50	0.47
1:A:551:LEU:HD11	1:A:553:PHE:CE1	2.50	0.47
1:A:316:GLU:HG3	1:A:351:TYR:OH	2.14	0.47
1:A:176:ALA:HB2	1:A:185:VAL:HG13	1.97	0.47
1:A:452:ARG:NH1	1:A:570:ILE:HD11	2.30	0.47
1:A:444:LEU:O	1:A:446:ARG:N	2.49	0.46
1:A:457:ILE:HG21	1:A:562:LYS:HB2	1.97	0.46
1:A:292:THR:HA	1:A:293:PRO:HD3	1.81	0.46
1:A:366:GLU:O	1:A:371:VAL:HG23	2.16	0.46
1:A:236:VAL:O	1:A:240:VAL:HG23	2.16	0.46
1:A:223:LEU:HD13	1:A:317:HIS:CD2	2.51	0.46
1:A:186:PHE:O	1:A:192:GLN:HB3	2.16	0.45
1:A:24:ALA:HA	1:A:266:LEU:HD21	1.98	0.45
1:A:175:LEU:O	1:A:230:CYS:HB2	2.17	0.45
1:A:58:VAL:HG12	1:A:60:GLU:CG	2.46	0.45
1:A:457:ILE:HG22	1:A:457:ILE:O	2.17	0.45
1:A:3:LEU:HD21	1:A:36:LEU:HD13	1.98	0.44
1:A:306:ILE:HD13	1:A:321:TYR:CD2	2.53	0.44
1:A:131:ARG:HD3	1:A:131:ARG:HA	1.80	0.44
1:A:383:LEU:HD12	1:A:384:GLU:H	1.83	0.44
1:A:71:PRO:N	1:A:72:PRO:CD	2.80	0.44
1:A:159:PHE:HZ	1:A:239:MET:HE1	1.82	0.44
1:A:133:TYR:CE1	1:A:135:LYS:HD2	2.53	0.44
1:A:553:PHE:CE2	1:A:558:MET:O	2.71	0.44
1:A:103:LEU:HD13	1:A:108:ARG:HH12	1.83	0.43
1:A:561:MET:O	1:A:565:LEU:HG	2.18	0.43
1:A:112:VAL:HG21	1:A:188:PRO:HG3	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:125:ILE:HD12	1:A:161:VAL:HA	2.01	0.43
1:A:155:SER:HG	2:B:7:TRP:HH2	1.64	0.43
1:A:410:PHE:CZ	1:A:452:ARG:HG2	2.53	0.43
1:A:319:TYR:N	1:A:320:PRO:CD	2.82	0.43
1:A:27:LEU:HD21	1:A:267:LEU:CD2	2.49	0.43
1:A:223:LEU:HD13	1:A:317:HIS:NE2	2.34	0.43
2:B:10:PRO:O	2:B:11:ALA:HB3	2.18	0.43
2:B:12:ARG:HG2	2:B:13:PRO:HD2	2.01	0.43
1:A:101:VAL:HG11	1:A:168:LEU:HD11	2.00	0.42
1:A:253:SER:HB3	1:A:256:LEU:HB2	2.00	0.42
1:A:348:ILE:HG13	1:A:423:TRP:CG	2.55	0.42
1:A:147:ILE:HA	1:A:147:ILE:HD13	1.76	0.42
1:A:239:MET:HE2	1:A:239:MET:HB2	1.61	0.42
1:A:71:PRO:HB2	1:A:72:PRO:HD3	2.01	0.42
1:A:231:ASP:O	1:A:232:ARG:C	2.58	0.42
1:A:284:ALA:HB1	1:A:302:TYR:CZ	2.53	0.42
1:A:267:LEU:HB3	1:A:273:LEU:HD13	2.02	0.42
1:A:358:GLU:O	1:A:362:LYS:HG3	2.20	0.42
1:A:448:GLU:HB3	1:A:451:VAL:HG23	2.01	0.41
1:A:553:PHE:CD2	1:A:558:MET:O	2.74	0.41
1:A:558:MET:HE2	1:A:558:MET:HB2	1.88	0.41
1:A:188:PRO:O	1:A:189:ASN:HB2	2.21	0.41
1:A:230:CYS:HB3	1:A:235:GLU:HG3	2.02	0.41
1:A:287:GLU:OE1	1:A:291:PRO:HA	2.20	0.41
1:A:561:MET:C	1:A:563:GLU:H	2.23	0.41
1:A:345:ALA:HB1	1:A:419:GLY:HA3	2.03	0.41
1:A:69:PRO:HG2	1:A:254:LEU:HD13	2.02	0.41
1:A:215:VAL:HG13	1:A:225:GLY:N	2.36	0.41
1:A:117:LEU:HD13	1:A:170:LEU:HD13	2.03	0.41
1:A:219:SER:HB3	1:A:354:CYS:SG	2.60	0.41
1:A:432:LEU:HG	1:A:579:LEU:HD22	2.02	0.41
1:A:71:PRO:N	1:A:72:PRO:HD2	2.36	0.41
1:A:375:LEU:HD23	1:A:375:LEU:HA	1.90	0.41
1:A:527:ARG:HE	1:A:527:ARG:CA	2.34	0.41
1:A:177:LEU:HA	1:A:177:LEU:HD12	1.69	0.40
1:A:54:ILE:HA	1:A:55:PRO:HD3	1.76	0.40
1:A:458:VAL:O	1:A:458:VAL:HG13	2.22	0.40
1:A:427:SER:HA	1:A:428:PRO:HD3	1.93	0.40
2:B:8:1E3:H15	2:B:13:PRO:O	2.21	0.40
1:A:62:THR:HG22	1:A:63:PHE:N	2.37	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	478/550 (87%)	434 (91%)	39 (8%)	5 (1%)	13	42
2	B	6/9 (67%)	6 (100%)	0	0	100	100
All	All	484/559 (87%)	440 (91%)	39 (8%)	5 (1%)	13	42

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	203	ASN
1	A	521	THR
1	A	452	ARG
1	A	572	SER
1	A	445	GLY

5.3.2 Protein sidechains [i](#)

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	406/461 (88%)	377 (93%)	29 (7%)	12	39
2	B	6/6 (100%)	5 (83%)	1 (17%)	2	7
All	All	412/467 (88%)	382 (93%)	30 (7%)	11	37

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	27	LEU

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Mol	Chain	Res	Type
1	A	66	SER
1	A	105	LEU
1	A	117	LEU
1	A	147	ILE
1	A	170	LEU
1	A	177	LEU
1	A	203	ASN
1	A	204	GLU
1	A	221	LEU
1	A	229	ARG
1	A	249	LEU
1	A	254	LEU
1	A	255	GLU
1	A	262	LYS
1	A	270	LEU
1	A	273	LEU
1	A	276	TYR
1	A	280	LEU
1	A	282	ASN
1	A	334	VAL
1	A	338	LEU
1	A	383	LEU
1	A	429	THR
1	A	438	THR
1	A	451	VAL
1	A	521	THR
1	A	522	VAL
1	A	580	THR
2	B	7	TRP

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

Mol	Chain	Res	Type
1	A	258	GLN
1	A	282	ASN
1	A	339	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

1 non-standard protein/DNA/RNA residue is 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	1E3	B	8	2	10,11,12	0.40	0	6,11,13	0.85	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	1E3	B	8	2	-	5/9/10/12	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	8	1E3	CAH-CAI-CAJ-CB
2	B	8	1E3	CAF-CAG-CAH-CAI
2	B	8	1E3	CAE-CAF-CAG-CAH
2	B	8	1E3	C-CA-CB-CAJ
2	B	8	1E3	N-CA-CB-CAJ

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	8	1E3	1	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	484/550 (88%)	0.03	9 (1%) 66 47	76, 103, 144, 183	0
2	B	7/9 (77%)	1.31	1 (14%) 7 4	91, 96, 113, 140	0
All	All	491/559 (87%)	0.05	10 (2%) 64 45	76, 103, 144, 183	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	6	ARG	4.0
1	A	69	PRO	4.0
1	A	576	LYS	3.3
1	A	572	SER	2.7
1	A	208	GLY	2.7
1	A	203	ASN	2.5
1	A	581	ALA	2.5
1	A	59	PRO	2.3
1	A	385	ALA	2.1
1	A	577	LEU	2.1

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	1E3	B	8	12/13	0.73	0.40	20,20,20,20	0

6.3 Carbohydrates [i](#)

There are no monosaccharides 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.