



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 11:26 PM UTC

PDB ID : 4NRX / pdb_00004nrx
Title : Crystal Structure of HIV-1 Neutralizing Antibody m66 in complex with gp41 MPER peptide
Authors : Ofek, G.; Yang, Y.; Kwong, P.D.
Deposited on : 2013-11-27
Resolution : 2.21 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
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

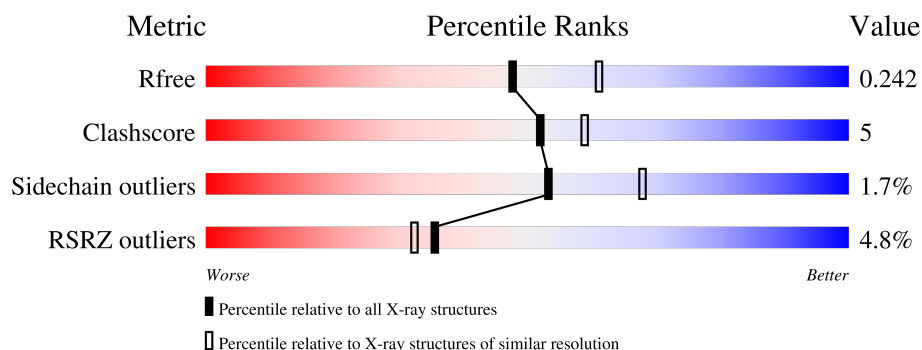
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.21 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	7682 (2.24-2.20)
Clashscore	190562	8402 (2.24-2.20)
Sidechain outliers	187428	8304 (2.24-2.20)
RSRZ outliers	180081	7683 (2.24-2.20)

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	C	19	5% 68% 32%
1	P	19	11% 68% 5% 26%
2	A	234	2% 90% 7% ..
2	D	234	7% 89% 8% .
2	H	234	2% 89% 10% .
3	B	213	2% 92% 5% .

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Mol	Chain	Length	Quality of chain
3	E	213	14% 85% 12% ..
3	L	213	% 94% 6%

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 10772 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 Envelope glycoprotein gp41.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
1	P	14	Total	C	N	O	0	0	0
			112	72	17	23			
1	C	13	Total	C	N	O	0	0	0
			103	66	16	21			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
P	654	LYS	GLU	engineered mutation	UNP Q9QQN5
C	654	LYS	GLU	engineered mutation	UNP Q9QQN5

- Molecule 2 is a protein called m66 Heavy Chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	231	Total	C	N	O	S	0	0	0
			1752	1117	286	342	7			
2	A	229	Total	C	N	O	S	0	0	0
			1742	1112	284	339	7			
2	D	228	Total	C	N	O	S	0	0	0
			1716	1089	280	339	8			

- Molecule 3 is a protein called m66 Light Chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	L	213	Total	C	N	O	S	0	0	0
			1626	1015	273	334	4			
3	B	207	Total	C	N	O	S	0	0	0
			1582	990	265	323	4			
3	E	211	Total	C	N	O	S	0	0	0
			1609	1006	271	328	4			

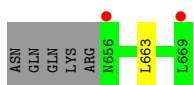
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	P	9	Total 9	O 9	0	0
4	C	6	Total 6	O 6	0	0
4	H	102	Total 102	O 102	0	0
4	A	110	Total 110	O 110	0	0
4	D	36	Total 36	O 36	0	0
4	L	90	Total 90	O 90	0	0
4	B	121	Total 121	O 121	0	0
4	E	56	Total 56	O 56	0	0

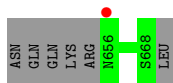
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

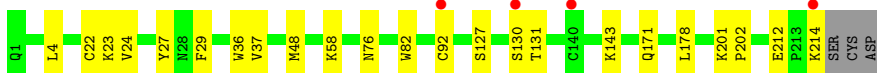
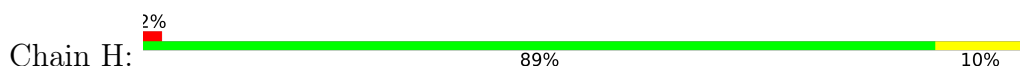
- Molecule 1: Envelope glycoprotein gp41



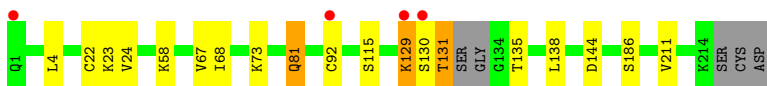
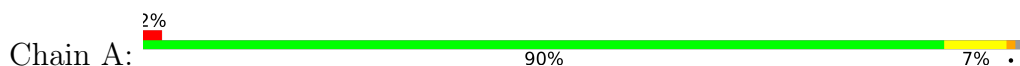
- Molecule 1: Envelope glycoprotein gp41



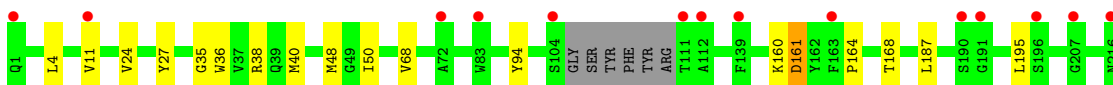
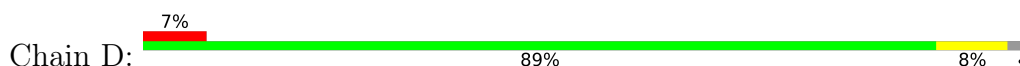
- Molecule 2: m66 Heavy Chain

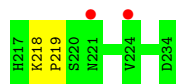


- Molecule 2: m66 Heavy Chain



- Molecule 2: m66 Heavy Chain

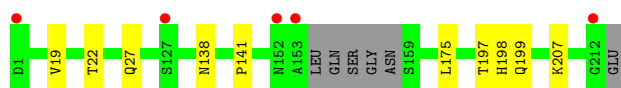




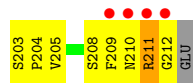
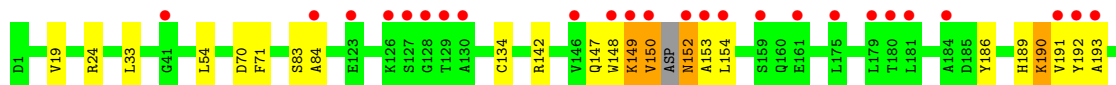
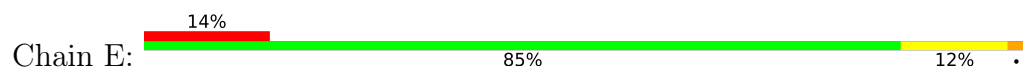
● Molecule 3: m66 Light Chain



● Molecule 3: m66 Light Chain



● Molecule 3: m66 Light Chain



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	197.26Å 93.50Å 89.83Å 90.00° 111.80° 90.00°	Depositor
Resolution (Å)	35.02 – 2.21 35.02 – 2.21	Depositor EDS
% Data completeness (in resolution range)	94.2 (35.02-2.21) 90.7 (35.02-2.21)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.05 (at 2.01Å)	Xtriage
Refinement program	PHENIX 1.8.2_1309	Depositor
R, R_{free}	0.188 , 0.215 (Not available) , 0.242	Depositor DCC
R_{free} test set	4693 reflections (4.72%)	wwPDB-VP
Wilson B-factor (Å ²)	27.3	Xtriage
Anisotropy	0.216	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 42.9	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	0.95	EDS
Total number of atoms	10772	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.04% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	C	0.64	0/104	0.59	0/140
1	P	0.67	0/113	0.65	0/151
2	A	0.58	1/1789 (0.1%)	0.76	2/2436 (0.1%)
2	D	0.48	0/1760	0.73	1/2397 (0.0%)
2	H	0.63	1/1800 (0.1%)	0.74	2/2452 (0.1%)
3	B	0.61	0/1616	0.75	1/2196 (0.0%)
3	E	0.55	1/1643 (0.1%)	0.77	2/2232 (0.1%)
3	L	0.56	0/1661	0.74	0/2258
All	All	0.57	3/10486 (0.0%)	0.75	8/14262 (0.1%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	H	82	TRP	C-N	13.72	1.65	1.34
2	A	135	THR	C-N	5.98	1.41	1.33
3	E	189	HIS	CG-ND1	-5.17	1.32	1.38

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	82	TRP	CA-C-N	-6.94	101.94	117.20
2	D	161	ASP	N-CA-C	6.55	120.12	111.28
2	H	82	TRP	O-C-N	6.04	132.37	122.70
2	A	144	ASP	N-CA-C	6.00	119.92	111.52
3	E	189	HIS	N-CA-C	5.96	119.42	109.95
3	B	138	ASN	N-CA-C	5.76	119.59	111.52
2	A	81	GLN	N-CA-C	5.37	117.55	109.23
3	E	153	ALA	N-CA-C	5.12	117.09	109.25

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	103	0	93	0	0
1	P	112	0	104	1	0
2	A	1742	0	1703	19	0
2	D	1716	0	1675	12	0
2	H	1752	0	1712	22	0
3	B	1582	0	1540	13	0
3	E	1609	0	1569	40	0
3	L	1626	0	1580	11	0
4	A	110	0	0	0	0
4	B	121	0	0	6	0
4	C	6	0	0	0	0
4	D	36	0	0	0	0
4	E	56	0	0	0	0
4	H	102	0	0	4	0
4	L	90	0	0	0	0
4	P	9	0	0	0	0
All	All	10772	0	9976	103	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 5.

All (103) 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)
2:A:131:THR:HG21	3:B:207:LYS:NZ	1.41	1.34
3:L:154:LEU:CD2	3:E:154:LEU:HD23	1.75	1.17
3:L:154:LEU:HD22	3:E:154:LEU:HD23	1.24	1.09
2:A:130:SER:O	2:A:131:THR:HB	1.55	1.07
2:D:38:ARG:NH2	2:D:40:MET:SD	2.37	0.97
2:A:131:THR:HG21	3:B:207:LYS:HZ2	1.02	0.97
2:A:131:THR:HG21	3:B:207:LYS:HZ1	1.29	0.93
3:E:190:LYS:CG	3:E:191:VAL:HG23	2.01	0.89
3:E:190:LYS:HG2	3:E:191:VAL:HG23	1.53	0.88
2:H:212:GLU:OE2	2:A:23:LYS:HE3	1.74	0.86
2:A:131:THR:CG2	3:B:207:LYS:HZ2	1.89	0.85

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:131:THR:CG2	3:B:207:LYS:NZ	2.34	0.84
2:H:36:TRP:HB3	2:H:48:MET:HE3	1.60	0.82
3:L:154:LEU:HD23	3:E:154:LEU:HD23	1.62	0.80
2:H:130:SER:HB3	4:H:393:HOH:O	1.82	0.79
3:E:150:VAL:HG12	3:E:150:VAL:O	1.82	0.77
3:E:190:LYS:HG2	3:E:191:VAL:N	1.98	0.77
3:E:209:PHE:O	3:E:209:PHE:CD2	2.38	0.77
2:A:129:LYS:O	2:A:130:SER:HB2	1.84	0.74
3:E:211:ARG:O	3:E:212:GLY:C	2.30	0.73
3:E:83:SER:O	3:E:84:ALA:HB2	1.88	0.72
2:A:130:SER:O	2:A:131:THR:CB	2.34	0.71
2:D:11:VAL:HG22	2:D:164:PRO:HG3	1.73	0.70
3:L:149:LYS:NZ	3:L:195:GLU:OE1	2.22	0.68
3:B:27:GLN:NE2	4:B:336:HOH:O	2.26	0.67
2:A:131:THR:CG2	3:B:207:LYS:HZ1	2.05	0.67
2:H:37:VAL:C	2:H:48:MET:HE2	2.19	0.67
3:E:134:CYS:HB2	3:E:148:TRP:CZ2	2.29	0.67
2:D:36:TRP:HB3	2:D:48:MET:HE3	1.79	0.65
3:E:152:ASN:N	3:E:152:ASN:OD1	2.30	0.64
3:E:190:LYS:CG	3:E:191:VAL:N	2.57	0.63
2:A:130:SER:OG	2:A:186:SER:HB2	1.97	0.63
3:E:192:TYR:HB2	3:E:209:PHE:CZ	2.34	0.63
3:L:24:ARG:NH1	3:L:70:ASP:OD2	2.29	0.63
3:E:150:VAL:O	3:E:150:VAL:CG1	2.46	0.63
2:H:23:LYS:NZ	4:H:367:HOH:O	2.26	0.63
3:L:156:SER:OG	3:E:147:GLN:NE2	2.32	0.62
3:E:134:CYS:CB	3:E:148:TRP:CZ2	2.84	0.61
2:H:127:SER:HB3	4:H:393:HOH:O	2.03	0.59
3:E:186:TYR:CD1	3:E:192:TYR:CZ	2.90	0.59
3:E:193:ALA:HB2	3:E:208:SER:HB3	1.85	0.59
3:E:134:CYS:HB2	3:E:148:TRP:CH2	2.37	0.58
3:E:83:SER:O	3:E:84:ALA:CB	2.53	0.57
2:D:24:VAL:HG11	2:D:27:TYR:CE1	2.40	0.57
2:H:4:LEU:HD22	2:H:24:VAL:HG22	1.88	0.56
3:E:209:PHE:CD2	3:E:209:PHE:C	2.83	0.56
3:L:24:ARG:HD3	3:L:70:ASP:OD2	2.05	0.55
2:H:36:TRP:CB	2:H:48:MET:HE3	2.34	0.55
2:H:131:THR:HG22	4:H:391:HOH:O	2.06	0.55
2:A:22:CYS:CB	2:A:92:CYS:HG	2.20	0.54
3:B:197:THR:HG23	4:B:332:HOH:O	2.08	0.53
2:H:36:TRP:HB3	2:H:48:MET:CE	2.36	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:160:LYS:HG2	2:D:161:ASP:OD1	2.09	0.52
3:E:210:ASN:O	3:E:210:ASN:OD1	2.28	0.51
2:D:160:LYS:HG2	2:D:161:ASP:CG	2.36	0.51
3:E:186:TYR:CD1	3:E:192:TYR:OH	2.64	0.51
3:L:154:LEU:HD23	3:E:154:LEU:CD2	2.37	0.51
3:E:190:LYS:CD	3:E:191:VAL:HG23	2.41	0.51
2:H:36:TRP:C	2:H:48:MET:HE3	2.37	0.49
3:E:150:VAL:O	3:E:152:ASN:HB3	2.13	0.49
3:E:186:TYR:HD1	3:E:192:TYR:CZ	2.30	0.49
3:B:27:GLN:HG3	4:B:336:HOH:O	2.13	0.48
2:H:214:LYS:NZ	3:L:122:ASP:OD1	2.47	0.47
3:L:143:GLU:OE2	3:E:142:ARG:NH2	2.48	0.47
3:E:190:LYS:O	3:E:211:ARG:O	2.33	0.46
3:B:22:THR:HG23	4:B:334:HOH:O	2.16	0.46
3:E:203:SER:OG	3:E:204:PRO:HD2	2.15	0.46
3:E:149:LYS:HG2	3:E:193:ALA:HB3	1.97	0.46
2:H:143:LYS:NZ	2:H:171:GLN:OE1	2.49	0.45
2:D:48:MET:HE1	2:D:94:TYR:HD1	1.81	0.45
3:E:192:TYR:HB2	3:E:209:PHE:CE1	2.50	0.45
3:E:149:LYS:O	3:E:193:ALA:N	2.48	0.45
2:H:36:TRP:CZ3	2:H:92:CYS:HB2	2.52	0.45
2:H:201:LYS:N	2:H:202:PRO:CD	2.80	0.45
1:P:663:LEU:HD11	2:H:58:LYS:HD2	1.99	0.44
2:A:4:LEU:HD23	2:A:92:CYS:O	2.17	0.44
2:A:131:THR:HG22	4:B:301:HOH:O	2.16	0.44
2:H:22:CYS:SG	2:H:92:CYS:SG	3.16	0.43
2:D:195:LEU:HD12	2:D:195:LEU:C	2.43	0.43
2:H:178:LEU:C	2:H:178:LEU:HD12	2.44	0.43
2:D:218:LYS:N	2:D:219:PRO:CD	2.82	0.43
3:E:210:ASN:O	3:E:210:ASN:CG	2.61	0.43
2:A:129:LYS:H	2:A:129:LYS:HG3	1.60	0.43
2:D:168:THR:O	2:D:168:THR:HG23	2.18	0.43
3:E:33:LEU:HD22	3:E:71:PHE:CG	2.53	0.42
2:D:24:VAL:CG1	2:D:27:TYR:CE1	3.02	0.42
3:E:24:ARG:NH1	3:E:70:ASP:OD2	2.52	0.42
3:E:186:TYR:HD1	3:E:192:TYR:OH	2.01	0.42
3:L:54:LEU:HD11	3:L:58:VAL:CG1	2.50	0.42
2:H:24:VAL:CG1	2:H:27:TYR:CE1	3.03	0.42
2:H:22:CYS:HG	2:H:92:CYS:HG	1.63	0.42
2:A:68:ILE:HB	2:A:81:GLN:HG2	2.02	0.41
2:A:4:LEU:HG	2:A:24:VAL:HG22	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:138:LEU:HB2	2:A:211:VAL:HG11	2.02	0.41
3:B:199:GLN:OE1	4:B:388:HOH:O	2.21	0.41
2:H:29:PHE:CD2	2:H:76:ASN:HA	2.56	0.41
3:B:141:PRO:O	3:B:198:HIS:HE1	2.04	0.41
2:A:22:CYS:CB	2:A:92:CYS:SG	3.09	0.41
3:B:175:LEU:C	3:B:175:LEU:HD23	2.46	0.41
2:H:4:LEU:HB3	2:H:22:CYS:SG	2.62	0.40
2:D:35:GLY:HA3	2:D:50:ILE:HG22	2.02	0.40
3:E:186:TYR:CE1	3:E:192:TYR:CZ	3.09	0.40
3:E:192:TYR:CD1	3:E:192:TYR:N	2.90	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

There are no protein backbone outliers to report in this entry.

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	C	10/18 (56%)	10 (100%)	0	100	100
1	P	11/18 (61%)	11 (100%)	0	100	100
2	A	194/198 (98%)	188 (97%)	6 (3%)	35	46
2	D	193/198 (98%)	190 (98%)	3 (2%)	55	70
2	H	195/198 (98%)	195 (100%)	0	100	100
3	B	182/187 (97%)	181 (100%)	1 (0%)	81	89
3	E	185/187 (99%)	177 (96%)	8 (4%)	26	33
3	L	187/187 (100%)	185 (99%)	2 (1%)	65	78
All	All	1157/1191 (97%)	1137 (98%)	20 (2%)	53	68

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	A	58	LYS
2	A	67	VAL
2	A	73	LYS
2	A	115	SER
2	A	129	LYS
2	A	131	THR
2	D	4	LEU
2	D	68	VAL
2	D	187	LEU
3	L	19	VAL
3	L	136	LEU
3	B	19	VAL
3	E	19	VAL
3	E	54	LEU
3	E	149	LYS
3	E	150	VAL
3	E	152	ASN
3	E	190	LYS
3	E	205	VAL
3	E	211	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
2	H	39	GLN
2	A	39	GLN
2	D	39	GLN
2	D	77	ASN
3	L	27	GLN
3	L	38	GLN
3	L	79	GLN
3	B	38	GLN
3	B	79	GLN
3	B	198	HIS
3	E	38	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](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
2	H	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	H	82:TRP	C	82(A):SER	N	1.65

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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	C	13/19 (68%)	0.50	1 (7%) 19 17	27, 37, 66, 74	0
1	P	14/19 (73%)	0.27	2 (14%) 6 4	25, 36, 60, 66	0
2	A	229/234 (97%)	0.09	4 (1%) 69 67	25, 36, 59, 107	0
2	D	228/234 (97%)	0.69	16 (7%) 22 19	32, 54, 98, 133	0
2	H	231/234 (98%)	0.06	4 (1%) 69 67	25, 38, 62, 86	0
3	B	207/213 (97%)	0.02	5 (2%) 59 57	21, 33, 58, 85	0
3	E	211/213 (99%)	0.72	29 (13%) 6 5	26, 44, 106, 128	0
3	L	213/213 (100%)	0.28	3 (1%) 73 72	23, 39, 77, 97	0
All	All	1346/1379 (97%)	0.31	64 (4%) 35 32	21, 40, 86, 133	0

All (64) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	E	150	VAL	5.6
2	H	92	CYS	4.6
3	E	129	THR	3.9
2	H	130	SER	3.9
3	E	209	PHE	3.8
3	E	154	LEU	3.8
2	A	1	GLN	3.7
2	D	112	ALA	3.6
3	E	148	TRP	3.6
3	E	146	VAL	3.3
3	B	153	ALA	3.2
2	D	163	PHE	3.1
1	C	656	ASN	3.0
2	D	191	GLY	3.0
3	E	191	VAL	2.9
3	E	130	ALA	2.8

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Mol	Chain	Res	Type	RSRZ
3	E	159	SER	2.8
3	E	153	ALA	2.7
3	L	204	PRO	2.7
3	E	123	GLU	2.7
3	E	161	GLU	2.6
3	E	192	TYR	2.6
3	E	184	ALA	2.6
3	E	193	ALA	2.6
2	D	104	SER	2.6
2	A	129	LYS	2.5
2	D	111	THR	2.5
2	A	130	SER	2.5
3	E	181	LEU	2.5
3	E	179	LEU	2.5
2	H	140	CYS	2.4
2	D	83	TRP	2.4
3	E	128	GLY	2.4
2	D	11	VAL	2.4
3	E	41	GLY	2.4
3	L	154	LEU	2.4
3	B	1	ASP	2.4
2	D	139	PHE	2.4
3	L	181	LEU	2.3
3	E	212	GLY	2.3
3	E	180	THR	2.3
2	D	196	SER	2.3
2	D	190	SER	2.3
2	D	1	GLN	2.2
3	B	152	ASN	2.2
3	E	152	ASN	2.2
2	D	72	ALA	2.2
2	H	214	LYS	2.2
2	A	92	CYS	2.1
3	E	84	ALA	2.1
3	E	127	SER	2.1
1	P	656	ASN	2.1
2	D	221	ASN	2.1
3	E	211	ARG	2.1
3	E	149	LYS	2.1
3	E	175	LEU	2.1
2	D	207	GLY	2.1
3	B	127	SER	2.1

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Mol	Chain	Res	Type	RSRZ
3	E	126	LYS	2.1
2	D	216	ASN	2.1
3	E	210	ASN	2.1
2	D	224	VAL	2.1
3	B	212	GLY	2.0
1	P	669	LEU	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.