



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:07 PM GMT

PDB ID : 4NRX  
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](mailto: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.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
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) : trunk26865

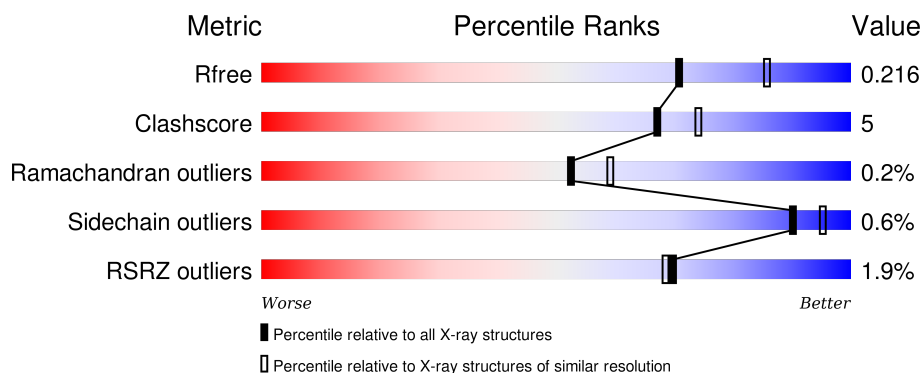
# 1 Overall quality at a glance ⓘ

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))
$R_{free}$	91344	4405 (2.24-2.20)
Clashscore	102246	5146 (2.24-2.20)
Ramachandran outliers	100387	5065 (2.24-2.20)
Sidechain outliers	100360	5066 (2.24-2.20)
RSRZ outliers	91569	4414 (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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	C	19	<div> <div>68%</div> <div>32%</div> </div>
1	P	19	<div> <div>68%</div> <div>5%</div> <div>26%</div> </div>
2	A	234	<div> <div>2%</div> <div>91%</div> <div>6%</div> <div>••</div> </div>
2	D	234	<div> <div>3%</div> <div>92%</div> <div>6%</div> <div>•</div> </div>
2	H	234	<div> <div>90%</div> <div>9%</div> <div>•</div> </div>

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Mol	Chain	Length	Quality of chain
3	B	213	<div><div></div><div>94%</div><div></div><div></div><div></div></div>
3	E	213	<div><div>5%</div><div></div><div>87%</div><div></div><div>10%</div><div></div><div></div></div>
3	L	213	<div><div>%</div><div></div><div>95%</div><div></div><div>5%</div><div></div><div></div></div>

## 2 Entry composition

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	O	0	0
			9	9		
4	C	6	Total	O	0	0
			6	6		
4	H	102	Total	O	0	0
			102	102		
4	A	110	Total	O	0	0
			110	110		
4	D	36	Total	O	0	0
			36	36		
4	L	90	Total	O	0	0
			90	90		
4	B	121	Total	O	0	0
			121	121		
4	E	56	Total	O	0	0
			56	56		

### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

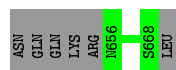
- Molecule 1: Envelope glycoprotein gp41

Chain P: 



- Molecule 1: Envelope glycoprotein gp41

Chain C: 



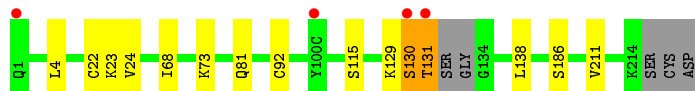
- Molecule 2: m66 Heavy Chain

Chain H: 




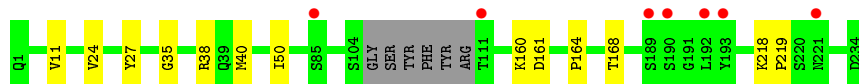
- Molecule 2: m66 Heavy Chain

Chain A: 



- Molecule 2: m66 Heavy Chain

Chain D: 



- Molecule 3: m66 Light Chain

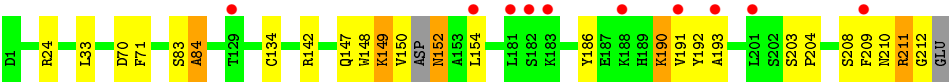
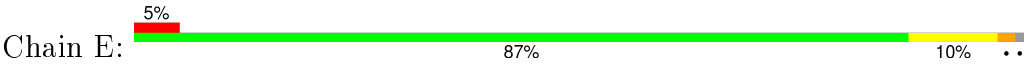
Chain L: 



● 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, $\alpha$ , $\beta$ , $\gamma$	197.26Å 93.50Å 89.83Å 90.00° 111.80° 90.00°	Depositor
Resolution (Å)	35.02 – 2.21 35.02 – 2.02	Depositor EDS
% Data completeness (in resolution range)	94.2 (35.02-2.21) 85.6 (35.02-2.02)	Depositor EDS
$R_{merge}$	0.17	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.87 (at 2.01Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE: 1.8.2_1309)	Depositor
R, $R_{free}$	0.188 , 0.215 0.187 , 0.216	Depositor DCC
$R_{free}$ test set	3471 reflections (5.04%)	DCC
Wilson B-factor (Å <sup>2</sup> )	27.2	Xtriage
Anisotropy	0.225	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 48.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	4 of 93755 reflections (0.004%)	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	10772	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.03% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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	C	0.46	0/104	0.49	0/140
1	P	0.48	0/113	0.53	0/151
2	A	0.48	0/1789	0.59	0/2436
2	D	0.40	0/1760	0.56	0/2397
2	H	0.56	1/1800 (0.1%)	0.61	2/2452 (0.1%)
3	B	0.49	0/1616	0.61	0/2196
3	E	0.44	0/1643	0.58	0/2232
3	L	0.44	0/1661	0.58	0/2258
All	All	0.47	1/10486 (0.0%)	0.59	2/14262 (0.0%)

All (1) 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

All (2) 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	H	82	TRP	O-C-N	6.04	132.37	122.70

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	20	0
2	D	1716	0	1675	9	0
2	H	1752	0	1712	18	0
3	B	1582	0	1540	12	0
3	E	1609	0	1569	39	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	3	0
4	L	90	0	0	0	0
4	P	9	0	0	0	0
All	All	10772	0	9976	95	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 (95) 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
2:A:131:THR:CG2	3:B:207:LYS:NZ	2.34	0.84
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:O	3:E:150:VAL:HG12	1.82	0.77
3:E:190:LYS:HG2	3:E:191:VAL:N	1.98	0.77
3:E:209:PHE:CD2	3:E:209:PHE:O	2.38	0.77
2:A:129:LYS:O	2:A:130:SER:HB2	1.84	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:E:83:SER:O	3:E:84:ALA:HB2	1.88	0.73
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:E:211:ARG:O	3:E:212:GLY:C	2.30	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
3:E:134:CYS:HB2	3:E:148:TRP:CZ2	2.29	0.67
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	0.64
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:36:TRP:HB3	2:H:48:MET:HE1	1.80	0.62
3:L:156:SER:OG	3:E:147:GLN:NE2	2.32	0.62
3:L:149:LYS:NZ	3:L:195:GLU:OE1	2.22	0.61
3:E:134:CYS:CB	3:E:148:TRP:CZ2	2.84	0.61
2:A:22:CYS:HG	2:A:92:CYS:HG	1.43	0.60
2:H:36:TRP:C	2:H:48:MET:HE3	2.23	0.60
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
2:H:36:TRP:HB3	2:H:48:MET:CE	2.36	0.56
3:L:24:ARG:HD3	3:L:70:ASP:OD2	2.05	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:D:160:LYS:HG2	2:D:161:ASP:OD1	2.09	0.52
3:E:209:PHE:CD2	3:E:209:PHE:C	2.83	0.51
3:E:210:ASN:OD1	3:E:210:ASN:O	2.28	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
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:37:VAL:N	2:H:48:MET:HE3	2.29	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
2:D:160:LYS:HG2	2:D:161:ASP:CG	2.36	0.46
3:E:149:LYS:HG2	3:E:193:ALA:HB3	1.97	0.46
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:143:LYS:NZ	2:H:171:GLN:OE1	2.49	0.45
2:H:201:LYS:N	2:H:202:PRO:CD	2.80	0.45
2:A:131:THR:HG22	4:B:301:HOH:O	2.16	0.44
2:A:4:LEU:HD23	2:A:92:CYS:O	2.17	0.44
1:P:663:LEU:HD11	2:H:58:LYS:HD2	1.99	0.44
2:H:22:CYS:SG	2:H:92:CYS:SG	3.16	0.43
2:D:168:THR:O	2:D:168:THR:HG23	2.18	0.43
2:D:218:LYS:N	2:D:219:PRO:CD	2.82	0.43
2:A:129:LYS:H	2:A:129:LYS:HG3	1.60	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:186:TYR:HD1	3:E:192:TYR:OH	2.01	0.42
3:E:24:ARG:NH1	3:E:70:ASP:OD2	2.52	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: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
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
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:192:TYR:CD1	3:E:192:TYR:N	2.90	0.40
3:E:186:TYR:CE1	3:E:192:TYR:CZ	3.09	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	C	11/19 (58%)	11 (100%)	0	0	100	100
1	P	12/19 (63%)	12 (100%)	0	0	100	100
2	A	225/234 (96%)	222 (99%)	2 (1%)	1 (0%)	39	41
2	D	224/234 (96%)	219 (98%)	5 (2%)	0	100	100
2	H	229/234 (98%)	228 (100%)	1 (0%)	0	100	100
3	B	203/213 (95%)	200 (98%)	3 (2%)	0	100	100
3	E	207/213 (97%)	202 (98%)	4 (2%)	1 (0%)	34	34
3	L	211/213 (99%)	204 (97%)	7 (3%)	0	100	100
All	All	1322/1379 (96%)	1298 (98%)	22 (2%)	2 (0%)	52	59

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	E	84	ALA
2	A	130	SER

### 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	C	10/18 (56%)	10 (100%)	0	100	100
1	P	11/18 (61%)	11 (100%)	0	100	100
2	A	194/198 (98%)	191 (98%)	3 (2%)	72	84

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	D	193/198 (98%)	193 (100%)	0	100	100
2	H	195/198 (98%)	195 (100%)	0	100	100
3	B	182/187 (97%)	182 (100%)	0	100	100
3	E	185/187 (99%)	181 (98%)	4 (2%)	60	72
3	L	187/187 (100%)	187 (100%)	0	100	100
All	All	1157/1191 (97%)	1150 (99%)	7 (1%)	90	95

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	A	73	LYS
2	A	115	SER
2	A	131	THR
3	E	149	LYS
3	E	152	ASN
3	E	190	LYS
3	E	211	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
3	B	198	HIS

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

## 5.5 Carbohydrates ⓘ

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 ⓘ

### 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	C	13/19 (68%)	-0.36	0 100 100	27, 37, 66, 74	0
1	P	14/19 (73%)	-0.41	0 100 100	25, 36, 60, 66	0
2	A	229/234 (97%)	-0.35	4 (1%) 73 72	25, 36, 59, 107	0
2	D	228/234 (97%)	0.09	7 (3%) 52 51	32, 54, 98, 133	0
2	H	231/234 (98%)	-0.39	0 100 100	25, 38, 62, 86	0
3	B	207/213 (97%)	-0.37	1 (0%) 91 91	21, 33, 58, 85	0
3	E	211/213 (99%)	0.06	10 (4%) 35 34	26, 44, 107, 128	0
3	L	213/213 (100%)	-0.17	3 (1%) 78 77	23, 39, 77, 97	0
All	All	1346/1379 (97%)	-0.19	25 (1%) 70 68	21, 40, 86, 133	0

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	111	THR	5.5
2	A	100(C)	TYR	4.4
3	E	188	LYS	3.5
2	D	221	ASN	3.4
3	E	129	THR	3.2
3	L	126	LYS	3.2
3	B	1	ASP	3.0
3	E	181	LEU	3.0
3	E	191	VAL	3.0
2	D	189	SER	2.9
2	D	190	SER	2.6
2	D	85	SER	2.6
3	E	183	LYS	2.6
3	E	182	SER	2.5
2	D	192	LEU	2.5
3	E	209	PHE	2.4

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Mol	Chain	Res	Type	RSRZ
3	E	154	LEU	2.3
3	E	201	LEU	2.3
2	A	1	GLN	2.2
3	L	213	GLU	2.2
3	E	193	ALA	2.2
3	L	154	LEU	2.1
2	A	130	SER	2.1
2	D	193	TYR	2.0
2	A	131	THR	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 [i](#)

There are no ligands in this entry.

## 6.5 Other polymers [i](#)

There are no such residues in this entry.