



Full wwPDB X-ray Structure Validation Report

Feb 26, 2014 – 05:48 PM GMT

PDB ID : 3MLZ
Title : Crystal structure of anti-HIV-1 V3 Fab 3074 in complex with a VI191 V3 peptide
Authors : Kong, X.-P.
Deposited on : 2010-04-18
Resolution : 2.99 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

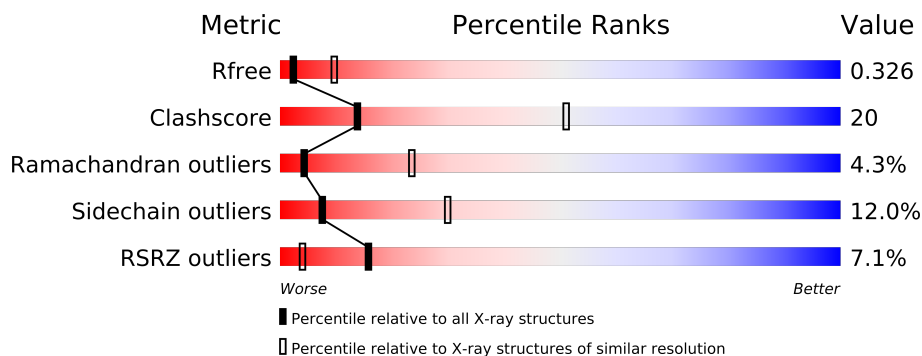
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.99 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}	66092	1216 (3.00-3.00)
Clashscore	79885	1594 (3.00-3.00)
Ramachandran outliers	78287	1537 (3.00-3.00)
Sidechain outliers	78261	1540 (3.00-3.00)
RSRZ outliers	66119	1217 (3.00-3.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	L	214	
2	H	230	
3	P	25	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 3663 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Human monoclonal anti-HIV-1 gp120 V3 antibody 3074 Fab light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	214	Total	C	N	O	S	0	0	0
			1591	991	271	323	6			

- Molecule 2 is a protein called Human monoclonal anti-HIV-1 gp120 V3 antibody 3074 Fab heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	230	Total	C	N	O	S	0	0	0
			1721	1086	283	345	7			

- Molecule 3 is a protein called HIV-1 gp120 third variable region (V3) crown.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	P	22	Total	C	N	O	0	0	0
			163	102	32	29			

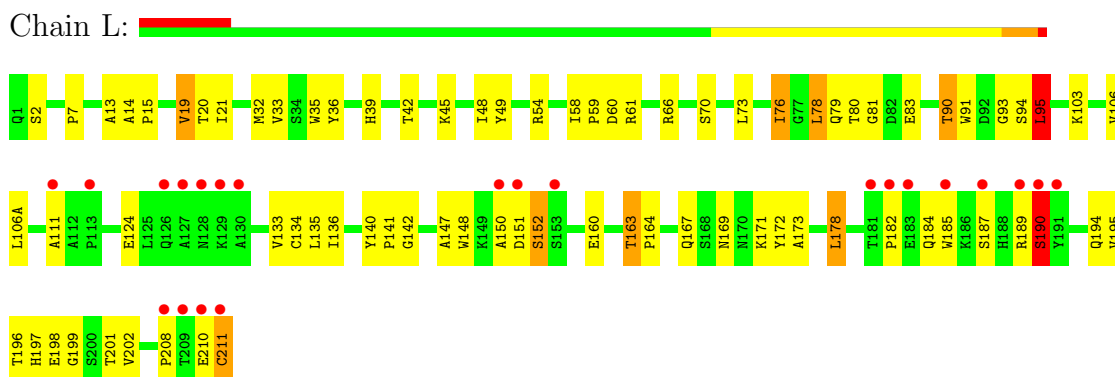
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	L	73	Total	O	0	0
			73	73		
4	H	106	Total	O	0	0
			106	106		
4	P	9	Total	O	0	0
			9	9		

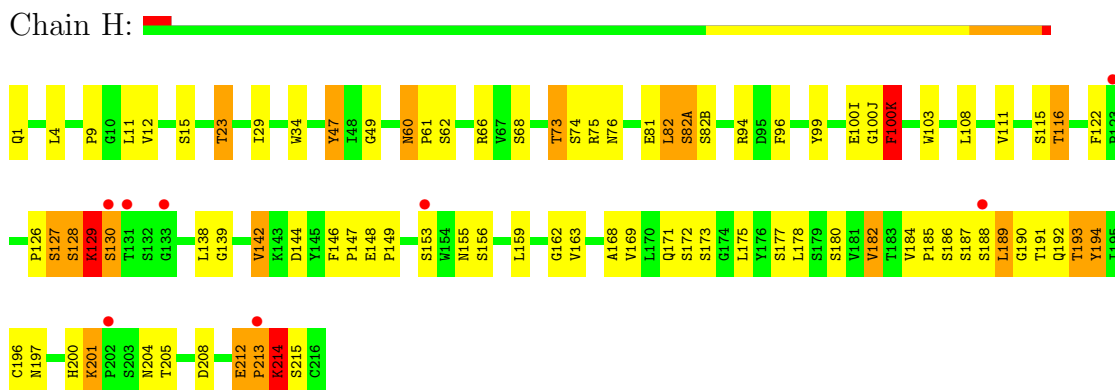
3 Residue-property plots

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.

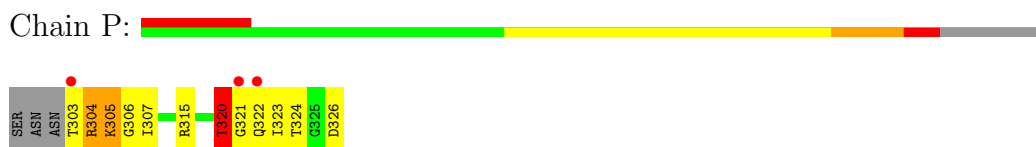
- Molecule 1: Human monoclonal anti-HIV-1 gp120 V3 antibody 3074 Fab light chain



- Molecule 2: Human monoclonal anti-HIV-1 gp120 V3 antibody 3074 Fab heavy chain



- Molecule 3: HIV-1 gp120 third variable region (V3) crown



4 Data and refinement statistics

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, α , β , γ	100.41 Å 100.41 Å 177.76 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.97 – 2.99 48.97 – 2.99	Depositor EDS
% Data completeness (in resolution range)	99.8 (48.97-2.99) 99.8 (48.97-2.99)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	7.67 (at 3.01 Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.191 , 0.330 0.190 , 0.326	Depositor DCC
R_{free} test set	538 reflections (5.01%)	DCC
Wilson B-factor (Å ²)	46.6	Xtriage
Anisotropy	0.105	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 24.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Outliers	0 of 11277 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	3663	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.78% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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	L	0.61	0/1629	0.76	0/2224
2	H	0.67	0/1766	0.82	0/2407
3	P	0.81	0/166	0.89	0/222
All	All	0.65	0/3561	0.80	0/4853

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	H	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	H	100(I)	GLU	Peptide

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	L	1591	0	1553	55	0
2	H	1721	0	1664	73	0
3	P	163	0	161	14	0
4	H	106	0	0	10	0
4	L	73	0	0	5	0
4	P	9	0	0	2	0
All	All	3663	0	3378	137	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 20.

All (137) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:H:82(A):SER:HB3	4:H:248:HOH:O	1.33	1.27
2:H:191:THR:HB	2:H:192:GLN:HA	1.06	1.05
3:P:304:ARG:HA	3:P:305:LYS:CB	1.86	1.03
1:L:61:ARG:HB2	1:L:76:ILE:HG22	1.41	0.99
2:H:193:THR:HA	2:H:194:TYR:HB2	1.46	0.97
2:H:191:THR:CB	2:H:192:GLN:HA	1.96	0.94
2:H:153:SER:HB3	2:H:197:ASN:HB2	1.48	0.94
2:H:116:THR:HA	2:H:146:PHE:O	1.66	0.94
3:P:304:ARG:HA	3:P:305:LYS:HB2	1.48	0.93
2:H:214:LYS:HG2	2:H:215:SER:H	1.35	0.92
2:H:191:THR:HB	2:H:192:GLN:CA	1.99	0.90
1:L:210:GLU:O	1:L:211:CYS:HB3	1.75	0.86
2:H:60:ASN:C	2:H:60:ASN:HD22	1.81	0.84
2:H:68:SER:HB3	2:H:81:GLU:HB2	1.59	0.83
1:L:189:ARG:O	1:L:190:SER:HB3	1.79	0.82
2:H:214:LYS:HG2	2:H:215:SER:N	1.97	0.78
2:H:148:GLU:HB2	2:H:149:PRO:HA	1.67	0.77
2:H:189:LEU:HB2	2:H:190:GLY:HA2	1.65	0.77
2:H:193:THR:HA	2:H:194:TYR:CB	2.15	0.75
2:H:66:ARG:HD2	2:H:82(A):SER:O	1.87	0.74
3:P:304:ARG:HA	3:P:305:LYS:HB3	1.68	0.73
1:L:182:PRO:HD3	4:L:281:HOH:O	1.90	0.71
2:H:60:ASN:HD22	2:H:61:PRO:N	1.88	0.70
1:L:185:TRP:O	1:L:208:PRO:HG3	1.90	0.70
1:L:133:VAL:HG12	1:L:135:LEU:HD21	1.73	0.69
2:H:189:LEU:CB	2:H:190:GLY:HA2	2.23	0.67
3:P:303:THR:HB	3:P:304:ARG:HH21	1.61	0.65
2:H:214:LYS:CG	2:H:215:SER:H	2.10	0.65
2:H:94:ARG:O	2:H:100(K):PHE:HA	1.95	0.65
2:H:116:THR:HG23	2:H:147:PRO:HD3	1.79	0.64

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:H:60:ASN:ND2	2:H:62:SER:H	1.94	0.64
2:H:82(A):SER:CB	4:H:248:HOH:O	2.11	0.63
2:H:29:ILE:HG12	2:H:76:ASN:OD1	2.01	0.61
1:L:111:ALA:HB2	1:L:171:LYS:HE2	1.82	0.60
3:P:315:ARG:NH2	4:P:3:HOH:O	2.33	0.60
2:H:163:VAL:HG22	2:H:182:VAL:HB	1.84	0.60
1:L:21:ILE:HD12	1:L:73:LEU:HD23	1.84	0.59
1:L:147:ALA:HB3	1:L:194:GLN:HB3	1.83	0.59
2:H:99:TYR:O	3:P:306:GLY:O	2.21	0.59
2:H:60:ASN:ND2	2:H:60:ASN:C	2.54	0.59
2:H:168:ALA:HA	2:H:178:LEU:HB3	1.85	0.59
1:L:32:MET:O	1:L:90:THR:HG23	2.04	0.58
1:L:94:SER:O	1:L:95:LEU:HB2	2.03	0.58
3:P:322:GLN:O	3:P:324:THR:N	2.33	0.57
2:H:142:VAL:HG13	2:H:178:LEU:HD12	1.87	0.56
2:H:126:PRO:HG2	2:H:213:PRO:HA	1.86	0.56
1:L:2:SER:HA	4:L:237:HOH:O	2.05	0.56
3:P:306:GLY:HA2	4:P:188:HOH:O	2.05	0.56
1:L:93:GLY:O	1:L:95:LEU:HD13	2.05	0.56
2:H:204:ASN:HA	4:H:320:HOH:O	2.05	0.56
2:H:193:THR:CA	2:H:194:TYR:HB2	2.31	0.55
1:L:133:VAL:CG1	1:L:135:LEU:HD21	2.37	0.55
1:L:160:GLU:HB3	2:H:169:VAL:HG21	1.88	0.55
1:L:60:ASP:O	1:L:61:ARG:CB	2.54	0.55
2:H:169:VAL:HG12	2:H:177:SER:O	2.06	0.55
1:L:36:TYR:OH	2:H:100(J):GLY:O	2.25	0.54
3:P:304:ARG:CA	3:P:305:LYS:CB	2.74	0.54
1:L:150:ALA:O	1:L:152:SER:N	2.35	0.54
1:L:148:TRP:CG	1:L:178:LEU:HD23	2.43	0.54
2:H:196:CYS:O	2:H:208:ASP:HA	2.09	0.53
2:H:126:PRO:O	2:H:128:SER:N	2.41	0.53
2:H:73:THR:N	4:H:247:HOH:O	2.42	0.53
1:L:7:PRO:HG2	4:L:270:HOH:O	2.09	0.52
2:H:9:PRO:HD2	4:H:249:HOH:O	2.09	0.52
1:L:111:ALA:CB	1:L:171:LYS:HE2	2.39	0.52
2:H:162:GLY:O	2:H:182:VAL:HA	2.09	0.52
2:H:156:SER:H	2:H:197:ASN:HD21	1.57	0.52
1:L:59:PRO:C	1:L:60:ASP:O	2.42	0.51
1:L:35:TRP:HB2	1:L:48:ILE:HB	1.93	0.50
2:H:189:LEU:CB	2:H:190:GLY:CA	2.89	0.50
2:H:23:THR:HB	4:H:297:HOH:O	2.11	0.50
1:L:83:GLU:HG3	1:L:106:VAL:H	1.77	0.50

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:L:169:ASN:OD1	1:L:171:LYS:HB2	2.12	0.49
1:L:45:LYS:HD3	4:L:214:HOH:O	2.12	0.49
2:H:144:ASP:HA	2:H:175:LEU:HB3	1.94	0.49
2:H:60:ASN:HD22	2:H:62:SER:H	1.58	0.49
1:L:148:TRP:CD2	1:L:178:LEU:HD23	2.48	0.49
2:H:159:LEU:HD21	2:H:182:VAL:HG21	1.95	0.49
2:H:201:LYS:HD3	2:H:201:LYS:O	2.13	0.48
1:L:210:GLU:OE2	2:H:129:LYS:HD2	2.13	0.48
3:P:303:THR:HB	3:P:304:ARG:NH2	2.28	0.47
2:H:47:TYR:CZ	2:H:49:GLY:HA2	2.49	0.47
2:H:189:LEU:HB2	2:H:190:GLY:CA	2.42	0.47
2:H:155:ASN:OD1	2:H:194:TYR:N	2.47	0.46
1:L:14:ALA:O	1:L:15:PRO:C	2.54	0.46
2:H:147:PRO:O	2:H:200:HIS:NE2	2.48	0.46
1:L:160:GLU:HA	1:L:160:GLU:OE2	2.15	0.46
2:H:96:PHE:CE1	2:H:100(J):GLY:HA2	2.51	0.46
2:H:171:GLN:C	2:H:173:SER:H	2.19	0.46
1:L:66:ARG:CD	4:L:266:HOH:O	2.63	0.46
1:L:134:CYS:C	1:L:135:LEU:HD23	2.35	0.46
1:L:142:GLY:HA3	1:L:172:TYR:CD1	2.51	0.46
2:H:188:SER:HB2	2:H:189:LEU:HA	1.98	0.45
2:H:185:PRO:C	2:H:187:SER:H	2.20	0.45
1:L:150:ALA:C	1:L:152:SER:N	2.70	0.45
4:H:224:HOH:O	3:P:320:THR:HG23	2.17	0.45
1:L:136:ILE:N	1:L:136:ILE:HD12	2.31	0.45
2:H:188:SER:H	2:H:189:LEU:HA	1.81	0.45
3:P:304:ARG:CA	3:P:305:LYS:HB2	2.33	0.45
1:L:60:ASP:O	1:L:61:ARG:HB3	2.15	0.45
2:H:189:LEU:HB3	2:H:191:THR:O	2.17	0.45
2:H:128:SER:HB2	2:H:214:LYS:HB3	1.99	0.44
1:L:39:HIS:HE1	1:L:81:GLY:O	2.01	0.44
2:H:12:VAL:O	2:H:111:VAL:HA	2.17	0.44
1:L:19:VAL:HG13	1:L:20:THR:N	2.33	0.43
1:L:210:GLU:O	1:L:211:CYS:CB	2.54	0.43
1:L:140:TYR:HA	1:L:141:PRO:C	2.39	0.43
1:L:195:VAL:O	1:L:201:THR:HA	2.18	0.43
2:H:188:SER:CB	2:H:189:LEU:HA	2.48	0.42
1:L:164:PRO:HA	1:L:173:ALA:O	2.19	0.42
1:L:48:ILE:HG22	1:L:49:TYR:N	2.35	0.42
2:H:82:LEU:HD23	2:H:82:LEU:HA	1.90	0.42
2:H:188:SER:N	2:H:189:LEU:HA	2.33	0.42
3:P:304:ARG:N	3:P:304:ARG:HD3	2.34	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:L:95:LEU:HD12	1:L:95:LEU:HA	1.83	0.42
1:L:163:THR:HG23	4:H:266:HOH:O	2.20	0.42
2:H:103:TRP:N	4:H:281:HOH:O	2.49	0.42
1:L:79:GLN:O	1:L:106:VAL:HG21	2.19	0.41
1:L:124:GLU:HG3	2:H:122:PHE:CE2	2.55	0.41
1:L:13:ALA:HB3	1:L:78:LEU:CD2	2.51	0.41
2:H:214:LYS:O	2:H:215:SER:CB	2.68	0.41
1:L:133:VAL:HG12	1:L:135:LEU:CD2	2.48	0.41
2:H:187:SER:HA	2:H:188:SER:HA	1.59	0.41
1:L:90:THR:HG22	1:L:91:TRP:H	1.85	0.41
2:H:193:THR:HA	2:H:194:TYR:CG	2.54	0.41
3:P:321:GLY:HA3	3:P:322:GLN:HA	1.56	0.41
1:L:54:ARG:HD2	1:L:58:ILE:O	2.21	0.41
2:H:212:GLU:HA	2:H:213:PRO:HD2	1.61	0.41
1:L:167:GLN:HB2	1:L:171:LYS:O	2.21	0.41
1:L:60:ASP:O	1:L:61:ARG:HG2	2.21	0.40
2:H:34:TRP:CH2	2:H:94:ARG:HD3	2.56	0.40
1:L:184:GLN:HA	1:L:187:SER:OG	2.21	0.40
2:H:139:GLY:HA3	2:H:180:SER:O	2.21	0.40
2:H:1:GLN:CA	4:H:241:HOH:O	2.65	0.40
1:L:197:HIS:C	1:L:199:GLY:H	2.24	0.40
2:H:108:LEU:HD13	2:H:108:LEU:C	2.42	0.40
2:H:96:PHE:CZ	2:H:100(J):GLY:HA2	2.57	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	L	212/214 (99%)	192 (91%)	14 (7%)	6 (3%)	8	37
2	H	228/230 (99%)	194 (85%)	24 (10%)	10 (4%)	4	22
3	P	20/25 (80%)	14 (70%)	2 (10%)	4 (20%)	0	0
All	All	460/469 (98%)	400 (87%)	40 (9%)	20 (4%)	4	23

All (20) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	L	95	LEU
1	L	152	SER
2	H	127	SER
2	H	129	LYS
2	H	194	TYR
2	H	213	PRO
2	H	214	LYS
3	P	305	LYS
3	P	323	ILE
1	L	190	SER
2	H	128	SER
2	H	172	SER
1	L	106(A)	LEU
2	H	130	SER
1	L	151	ASP
3	P	320	THR
2	H	100(K)	PHE
2	H	186	SER
3	P	307	ILE
1	L	198	GLU

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	L	180/180 (100%)	164 (91%)	16 (9%)	14	48
2	H	197/197 (100%)	169 (86%)	28 (14%)	5	22
3	P	15/18 (83%)	12 (80%)	3 (20%)	2	10
All	All	392/395 (99%)	345 (88%)	47 (12%)	7	30

All (47) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	L	19	VAL
1	L	33	VAL
1	L	42	THR

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Mol	Chain	Res	Type
1	L	70	SER
1	L	76	ILE
1	L	78	LEU
1	L	80	THR
1	L	90	THR
1	L	95	LEU
1	L	103	LYS
1	L	163	THR
1	L	178	LEU
1	L	190	SER
1	L	196	THR
1	L	202	VAL
1	L	211	CYS
2	H	4	LEU
2	H	11	LEU
2	H	15	SER
2	H	23	THR
2	H	47	TYR
2	H	60	ASN
2	H	73	THR
2	H	74	SER
2	H	75	ARG
2	H	82	LEU
2	H	82(A)	SER
2	H	82(B)	SER
2	H	100(K)	PHE
2	H	115	SER
2	H	116	THR
2	H	127	SER
2	H	129	LYS
2	H	130	SER
2	H	138	LEU
2	H	142	VAL
2	H	182	VAL
2	H	184	VAL
2	H	189	LEU
2	H	193	THR
2	H	201	LYS
2	H	205	THR
2	H	212	GLU
2	H	214	LYS
3	P	304	ARG

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Mol	Chain	Res	Type
3	P	320	THR
3	P	326	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	L	31	ASN
1	L	39	HIS
1	L	79	GLN
1	L	167	GLN
2	H	60	ASN
2	H	77	GLN
2	H	197	ASN
2	H	199	ASN
2	H	204	ASN

5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

There are no ligands in this entry.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

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, 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	L	214/214 (100%)	0.32	22 (10%) 7 2	16, 44, 77, 99	0
2	H	230/230 (100%)	0.18	8 (3%) 42 8	16, 38, 91, 96	0
3	P	22/25 (88%)	0.68	3 (13%) 4 1	22, 42, 70, 75	0
All	All	466/469 (99%)	0.27	33 (7%) 16 4	16, 43, 87, 99	0

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	L	211	CYS	6.5
1	L	129	LYS	4.8
2	H	130	SER	4.5
1	L	209	THR	4.4
3	P	321	GLY	4.2
1	L	210	GLU	4.1
1	L	150	ALA	3.8
1	L	126	GLN	3.6
1	L	153	SER	3.5
3	P	303	THR	3.2
2	H	188	SER	3.2
2	H	131	THR	2.9
1	L	190	SER	2.9
2	H	213	PRO	2.8
1	L	130	ALA	2.6
1	L	191	TYR	2.5
2	H	202	PRO	2.5
1	L	151	ASP	2.4
2	H	123	PRO	2.4
2	H	133	GLY	2.4
1	L	181	THR	2.4
1	L	128	ASN	2.3
1	L	185	TRP	2.3

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Mol	Chain	Res	Type	RSRZ
1	L	113	PRO	2.2
1	L	208	PRO	2.2
1	L	182	PRO	2.2
2	H	153	SER	2.2
3	P	322	GLN	2.2
1	L	127	ALA	2.2
1	L	111	ALA	2.1
1	L	189	ARG	2.1
1	L	187	SER	2.1
1	L	183	GLU	2.0

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

There are no ligands in this entry.

6.5 Other polymers ⓘ

There are no such residues in this entry.