



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 05:51 PM GMT

PDB ID : 4JO2
Title : Crystal structure of rabbit mAb R56 Fab in complex with V3 crown of HIV-1
Consensus A gp120
Authors : Pan, R.M.; Kong, X.P.
Deposited on : 2013-03-16
Resolution : 2.50 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

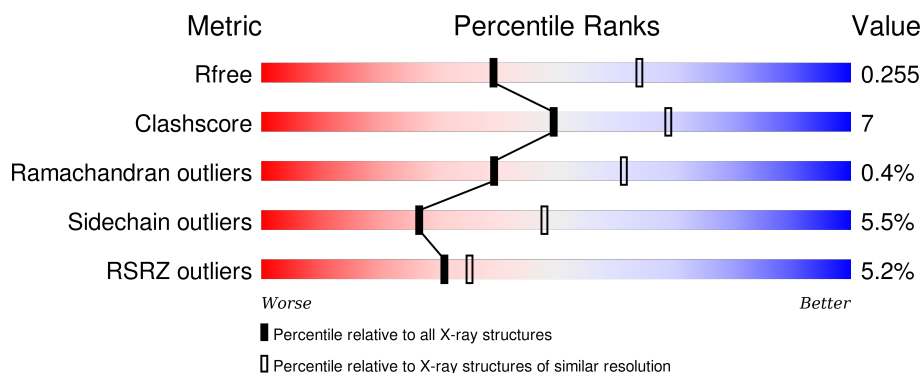
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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}	91344	3553 (2.50-2.50)
Clashscore	102246	4242 (2.50-2.50)
Ramachandran outliers	100387	4156 (2.50-2.50)
Sidechain outliers	100360	4158 (2.50-2.50)
RSRZ outliers	91569	3562 (2.50-2.50)

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. 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	L	216	<div> <div>6%</div> <div> <div></div> <div>85%</div> <div>13%</div> <div>•</div> </div> </div>
1	M	216	<div> <div>4%</div> <div> <div></div> <div>82%</div> <div>16%</div> <div>•</div> </div> </div>
2	H	210	<div> <div>5%</div> <div> <div></div> <div>81%</div> <div>15%</div> <div>• •</div> </div> </div>
2	I	210	<div> <div>5%</div> <div> <div></div> <div>80%</div> <div>14%</div> <div>• •</div> </div> </div>
3	P	23	<div> <div>4%</div> <div> <div>26%</div> <div>13%</div> <div>•</div> <div>57%</div> </div> </div>

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Mol	Chain	Length	Quality of chain
3	Q	23	<div><div><div></div><div></div><div></div></div><div>4%39%•57%</div></div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 6815 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 monoclonal anti-HIV-1 gp120 V3 antibody R56 light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	216	Total	C	N	O	S	0	0	0
			1596	988	265	334	9			
1	M	216	Total	C	N	O	S	0	0	0
			1596	988	265	334	9			

- Molecule 2 is a protein called monoclonal anti-HIV-1 gp120 V3 antibody R56 heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	204	Total	C	N	O	S	0	0	0
			1494	939	248	298	9			
2	I	204	Total	C	N	O	S	0	0	0
			1494	939	248	298	9			

- Molecule 3 is a protein called gp120.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	P	10	Total	C	N	O	0	0	0
			75	46	17	12			
3	Q	10	Total	C	N	O	0	0	0
			75	46	17	12			

- Molecule 4 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	H	1	Total	Ca	0	0
			1	1		
4	I	1	Total	Ca	0	0
			1	1		
4	L	2	Total	Ca	0	0
			2	2		
4	M	2	Total	Ca	0	0
			2	2		

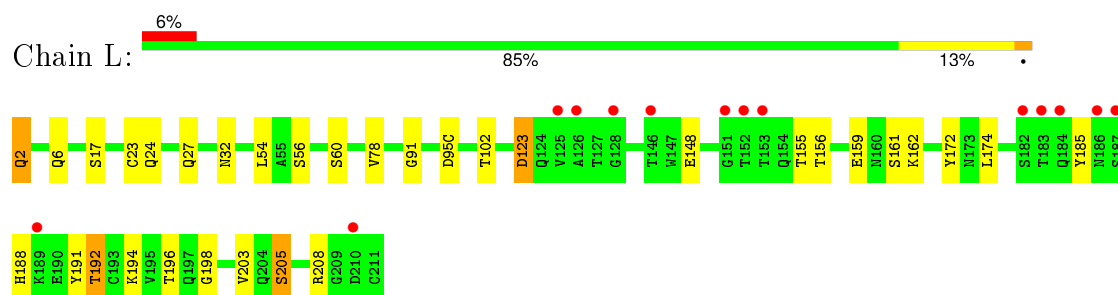
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	L	123	Total 123	O 123	0	0
5	H	105	Total 105	O 105	0	0
5	P	7	Total 7	O 7	0	0
5	M	136	Total 136	O 136	0	0
5	I	103	Total 103	O 103	0	0
5	Q	5	Total 5	O 5	0	0

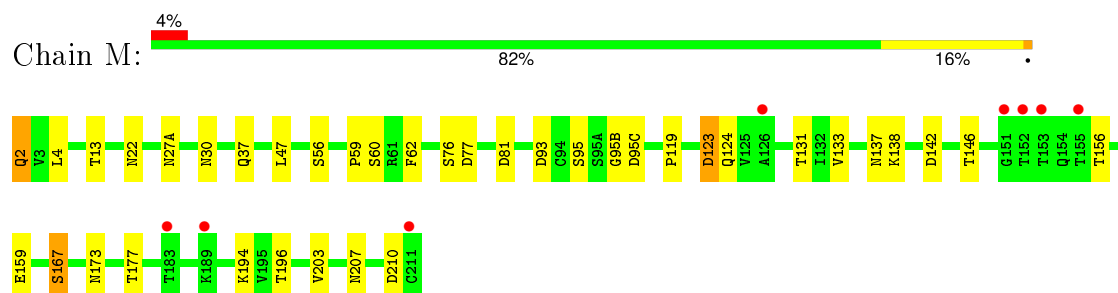
3 Residue-property plots [i](#)

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.

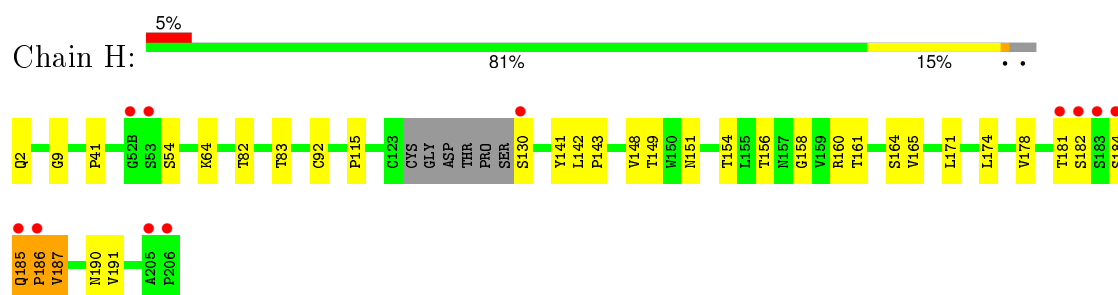
- Molecule 1: monoclonal anti-HIV-1 gp120 V3 antibody R56 light chain



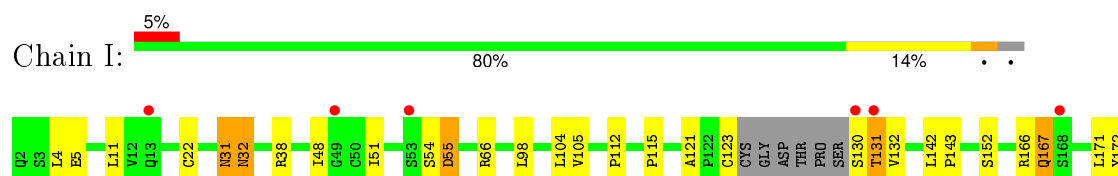
- Molecule 1: monoclonal anti-HIV-1 gp120 V3 antibody R56 light chain

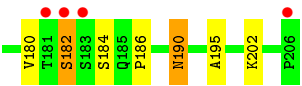


- Molecule 2: monoclonal anti-HIV-1 gp120 V3 antibody R56 heavy chain

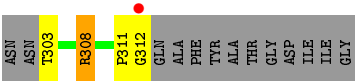


- Molecule 2: monoclonal anti-HIV-1 gp120 V3 antibody R56 heavy chain

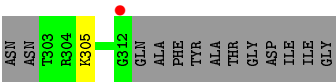




• Molecule 3: gp120



• Molecule 3: gp120



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	69.84Å 74.35Å 84.96Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	26.46 – 2.50 26.47 – 2.26	Depositor EDS
% Data completeness (in resolution range)	99.7 (26.46-2.50) 99.4 (26.47-2.26)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	0.16	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.24 (at 2.26Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7.3_928)	Depositor
R, R_{free}	0.190 , 0.265 0.179 , 0.255	Depositor DCC
R_{free} test set	1565 reflections (5.19%)	DCC
Wilson B-factor (Å ²)	25.5	Xtriage
Anisotropy	0.084	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 39.9	EDS
Estimated twinning fraction	0.098 for h,-k,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.42$, $\langle L^2 \rangle = 0.25$	Xtriage
Outliers	0 of 40911 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	6815	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.56% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CA

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	0.40	0/1628	0.57	0/2229
1	M	0.39	0/1628	0.58	0/2229
2	H	0.40	0/1528	0.59	0/2091
2	I	0.39	0/1528	0.59	0/2091
3	P	0.37	0/75	0.53	0/98
3	Q	0.38	0/75	0.56	0/98
All	All	0.39	0/6462	0.58	0/8836

There are no bond length outliers.

There are no bond angle outliers.

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	L	1596	0	1514	18	0
1	M	1596	0	1514	28	0
2	H	1494	0	1463	22	0
2	I	1494	0	1463	22	0
3	P	75	0	85	2	0
3	Q	75	0	85	0	0
4	H	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	I	1	0	0	0	0
4	L	2	0	0	0	0
4	M	2	0	0	0	0
5	H	105	0	0	8	0
5	I	103	0	0	5	0
5	L	123	0	0	2	0
5	M	136	0	0	8	0
5	P	7	0	0	1	0
5	Q	5	0	0	0	0
All	All	6815	0	6124	89	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 7.

All (89) 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:L:148:GLU:HB2	1:L:192:THR:HG23	1.55	0.88
1:L:198:GLY:O	5:L:437:HOH:O	1.96	0.83
1:M:119:PRO:O	5:M:446:HOH:O	1.98	0.81
2:H:64:LYS:NZ	5:H:450:HOH:O	2.14	0.79
2:H:2:GLN:N	5:H:428:HOH:O	2.19	0.75
1:M:138:LYS:H	1:M:173:ASN:ND2	1.85	0.74
1:L:188:HIS:O	1:L:208:ARG:NH2	2.19	0.74
1:L:56:SER:O	5:L:473:HOH:O	2.05	0.73
2:H:83:THR:OG1	5:H:415:HOH:O	2.07	0.72
2:I:152:SER:H	2:I:190:ASN:HD21	1.41	0.68
1:M:76:SER:HB3	5:M:436:HOH:O	1.95	0.67
1:M:124:GLN:OE1	1:M:131:THR:N	2.26	0.65
2:H:41:PRO:O	5:H:482:HOH:O	2.13	0.65
1:L:123:ASP:OD2	1:L:123:ASP:N	2.27	0.65
2:I:11:LEU:HD23	2:I:112:PRO:HG3	1.82	0.62
1:M:37:GLN:HB2	1:M:47:LEU:HD11	1.82	0.62
1:L:54:LEU:HD11	1:L:60:SER:HA	1.80	0.61
1:M:123:ASP:N	1:M:123:ASP:OD2	2.27	0.60
1:M:27(A):ASN:HD22	1:M:30:ASN:ND2	1.99	0.60
2:H:151:ASN:HD21	2:H:187:VAL:HG12	1.67	0.59
2:I:202:LYS:NZ	5:I:461:HOH:O	2.27	0.58
2:I:31:ASN:HB2	5:I:483:HOH:O	2.03	0.58
2:I:38:ARG:HB3	2:I:48:ILE:HD11	1.85	0.57
1:M:146:THR:HB	1:M:194:LYS:HB3	1.88	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:184:SER:HA	2:H:185:GLN:CB	2.36	0.55
2:H:9:GLY:O	5:H:427:HOH:O	2.18	0.55
2:H:148:VAL:HG22	2:H:191:VAL:HG22	1.89	0.54
1:M:2:GLN:NE2	1:M:95(C):ASP:HB2	2.22	0.54
1:M:2:GLN:NE2	5:M:487:HOH:O	2.28	0.54
2:H:151:ASN:HB2	2:H:154:THR:HB	1.90	0.54
3:P:308:ARG:HE	3:P:312:GLY:HA2	1.72	0.54
1:M:56:SER:OG	5:M:473:HOH:O	2.07	0.54
2:H:181:THR:OG1	2:H:182:SER:N	2.38	0.53
2:I:55:ASP:OD2	5:I:422:HOH:O	2.18	0.53
2:H:142:LEU:HD12	2:H:143:PRO:HA	1.90	0.52
2:I:142:LEU:HD12	2:I:143:PRO:HA	1.92	0.52
1:L:2:GLN:OE1	1:L:95(C):ASP:N	2.42	0.52
1:M:137:ASN:OD1	5:M:498:HOH:O	2.19	0.52
1:L:2:GLN:NE2	1:L:2:GLN:N	2.59	0.51
2:H:184:SER:HA	2:H:185:GLN:HB2	1.92	0.51
2:H:164:SER:HA	2:H:174:LEU:HB3	1.93	0.51
2:I:4:LEU:HD22	2:I:22:CYS:SG	2.51	0.51
1:M:119:PRO:HG3	2:I:123:CYS:HB2	1.93	0.50
2:H:54:SER:HB2	5:H:479:HOH:O	2.10	0.50
2:H:130:SER:O	5:H:484:HOH:O	2.20	0.50
1:M:77:ASP:N	5:M:436:HOH:O	2.44	0.50
3:P:311:PRO:O	5:P:405:HOH:O	2.20	0.49
1:L:6:GLN:NE2	1:L:102:THR:OG1	2.43	0.49
1:M:142:ASP:HB3	5:M:536:HOH:O	2.12	0.49
2:I:104:LEU:HD12	2:I:105:VAL:H	1.78	0.48
2:H:185:GLN:HE21	2:H:186:PRO:HD2	1.78	0.48
2:I:115:PRO:HD2	5:I:474:HOH:O	2.14	0.47
1:L:2:GLN:N	1:L:2:GLN:HE21	2.12	0.47
2:I:182:SER:OG	2:I:184:SER:N	2.45	0.47
1:L:192:THR:HB	1:L:205:SER:OG	2.15	0.46
2:H:115:PRO:HB3	2:H:141:TYR:HB3	1.97	0.46
1:L:162:LYS:HE3	1:L:172:TYR:CE2	2.50	0.46
1:M:93:ASP:OD1	1:M:95:SER:OG	2.26	0.45
2:I:142:LEU:HA	2:I:143:PRO:HA	1.74	0.45
2:I:180:VAL:HG13	2:I:182:SER:HB2	1.99	0.45
2:H:161:THR:N	5:H:434:HOH:O	2.17	0.45
1:L:17:SER:O	1:L:78:VAL:HG23	2.16	0.45
2:I:166:ARG:HB2	2:I:172:TYR:CE1	2.51	0.45
1:M:81:ASP:OD2	1:M:167:SER:HB2	2.17	0.44
2:I:143:PRO:HD2	2:I:195:ALA:CB	2.47	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:194:LYS:HG3	1:L:203:VAL:HG22	2.00	0.44
1:L:159:GLU:HG3	2:H:165:VAL:HG11	2.00	0.44
2:I:32:ASN:OD1	2:I:32:ASN:N	2.50	0.43
1:L:32:ASN:ND2	1:L:91:GLY:HA3	2.33	0.43
2:H:149:THR:OG1	2:H:190:ASN:HB2	2.18	0.43
1:M:2:GLN:N	1:M:2:GLN:CD	2.72	0.43
1:M:2:GLN:OE1	1:M:95(B):GLY:HA3	2.19	0.43
1:M:138:LYS:H	1:M:173:ASN:HD21	1.62	0.43
1:M:133:VAL:HG22	1:M:177:THR:HG22	2.00	0.43
1:M:124:GLN:NE2	5:M:428:HOH:O	2.51	0.43
1:L:2:GLN:HG3	1:L:27:GLN:CD	2.40	0.43
1:M:207:ASN:HB2	1:M:210:ASP:OD1	2.17	0.43
1:L:185:TYR:HA	1:L:191:TYR:OH	2.19	0.42
2:I:104:LEU:HD12	2:I:105:VAL:N	2.35	0.42
2:I:130:SER:HB2	2:I:131:THR:H	1.60	0.42
2:H:158:GLY:O	2:H:178:VAL:HA	2.20	0.42
1:M:137:ASN:HA	1:M:173:ASN:HD22	1.85	0.42
2:I:167:GLN:HG2	2:I:171:LEU:O	2.20	0.41
1:M:194:LYS:HG3	1:M:203:VAL:CG2	2.51	0.41
1:M:59:PRO:HG2	1:M:62:PHE:CE2	2.55	0.41
2:I:66:ARG:NE	5:I:423:HOH:O	2.38	0.41
2:H:142:LEU:HB2	2:H:171:LEU:HD22	2.02	0.41
1:M:4:LEU:HA	1:M:4:LEU:HD23	1.73	0.41
1:M:119:PRO:HD2	2:I:121:ALA:O	2.22	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	L	214/216 (99%)	202 (94%)	12 (6%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	M	214/216 (99%)	203 (95%)	11 (5%)	0	100	100
2	H	200/210 (95%)	188 (94%)	10 (5%)	2 (1%)	19	34
2	I	200/210 (95%)	189 (94%)	10 (5%)	1 (0%)	34	55
3	P	8/23 (35%)	7 (88%)	1 (12%)	0	100	100
3	Q	8/23 (35%)	7 (88%)	1 (12%)	0	100	100
All	All	844/898 (94%)	796 (94%)	45 (5%)	3 (0%)	39	61

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	H	186	PRO
2	H	185	GLN
2	I	186	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	L	183/183 (100%)	172 (94%)	11 (6%)	24	43
1	M	183/183 (100%)	174 (95%)	9 (5%)	31	55
2	H	170/175 (97%)	165 (97%)	5 (3%)	50	77
2	I	170/175 (97%)	158 (93%)	12 (7%)	18	34
3	P	8/17 (47%)	6 (75%)	2 (25%)	1	1
3	Q	8/17 (47%)	7 (88%)	1 (12%)	6	10
All	All	722/750 (96%)	682 (94%)	40 (6%)	27	48

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	L	2	GLN
1	L	23	CYS
1	L	24	GLN

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Mol	Chain	Res	Type
1	L	123	ASP
1	L	155	THR
1	L	156	THR
1	L	161	SER
1	L	174	LEU
1	L	192	THR
1	L	196	THR
1	L	205	SER
2	H	82	THR
2	H	92	CYS
2	H	156	THR
2	H	160	ARG
2	H	187	VAL
3	P	303	THR
3	P	308	ARG
1	M	2	GLN
1	M	13	THR
1	M	22	ASN
1	M	60	SER
1	M	123	ASP
1	M	156	THR
1	M	159	GLU
1	M	167	SER
1	M	196	THR
2	I	5	GLU
2	I	31	ASN
2	I	32	ASN
2	I	51	ILE
2	I	54	SER
2	I	55	ASP
2	I	98	LEU
2	I	131	THR
2	I	132	VAL
2	I	167	GLN
2	I	182	SER
2	I	190	ASN
3	Q	305	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	L	173	ASN

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Mol	Chain	Res	Type
2	H	32	ASN
2	H	185	GLN
1	M	30	ASN
1	M	137	ASN
1	M	173	ASN
2	I	190	ASN

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 ⓘ

Of 6 ligands modelled in this entry, 6 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	216/216 (100%)	0.07	14 (6%)	22	25	8, 19, 56, 72	0
1	M	216/216 (100%)	-0.08	8 (3%)	45	50	9, 19, 49, 59	0
2	H	204/210 (97%)	0.02	11 (5%)	29	33	9, 23, 43, 68	0
2	I	204/210 (97%)	0.13	10 (4%)	33	38	6, 22, 42, 61	0
3	P	10/23 (43%)	0.44	1 (10%)	9	10	16, 27, 41, 51	0
3	Q	10/23 (43%)	0.50	1 (10%)	9	10	15, 30, 47, 53	0
All	All	860/898 (95%)	0.05	45 (5%)	31	35	6, 21, 51, 72	0

All (45) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	I	182	SER	7.3
2	H	184	SER	5.3
2	H	130	SER	4.6
2	I	183	SER	4.2
2	I	131	THR	4.1
1	L	152	THR	3.7
1	M	151	GLY	3.7
1	M	153	THR	3.7
1	L	125	VAL	3.6
1	L	183	THR	3.4
2	H	205	ALA	3.4
2	I	130	SER	3.3
2	H	53	SER	3.2
1	L	186	ASN	3.2
2	I	181	THR	3.1
1	L	184	GLN	3.1
2	I	53	SER	3.1
1	L	187	SER	3.1
1	L	126	ALA	3.0

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Mol	Chain	Res	Type	RSRZ
1	M	183	THR	2.9
2	I	206	PRO	2.8
2	H	182	SER	2.8
3	Q	312	GLY	2.7
1	L	182	SER	2.7
1	L	189	LYS	2.7
1	L	151	GLY	2.7
2	I	168	SER	2.5
2	H	183	SER	2.5
2	H	206	PRO	2.4
1	M	155	THR	2.4
1	M	126	ALA	2.3
1	L	153	THR	2.3
3	P	312	GLY	2.3
1	L	146	THR	2.2
2	H	185	GLN	2.2
1	M	189	LYS	2.2
1	M	152	THR	2.2
1	L	128	GLY	2.1
2	H	186	PRO	2.1
2	H	52(B)	GLY	2.1
1	L	210	ASP	2.1
2	H	181	THR	2.1
2	I	13	GLN	2.0
1	M	211	CYS	2.0
2	I	49	GLY	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](#)

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(\AA^2)	Q<0.9
4	CA	I	301	1/1	0.97	0.11	-1.58	32,32,32,32	0
4	CA	H	301	1/1	0.97	0.10	-1.69	29,29,29,29	0
4	CA	L	301	1/1	0.99	0.08	-2.94	16,16,16,16	0
4	CA	M	301	1/1	0.99	0.09	-4.57	13,13,13,13	0
4	CA	L	302	1/1	0.92	0.07	-	37,37,37,37	0
4	CA	M	302	1/1	0.96	0.06	-	38,38,38,38	0

6.5 Other polymers [i](#)

There are no such residues in this entry.