



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 10:14 AM UTC

PDB ID : 4JO2 / pdb_00004jo2
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

<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

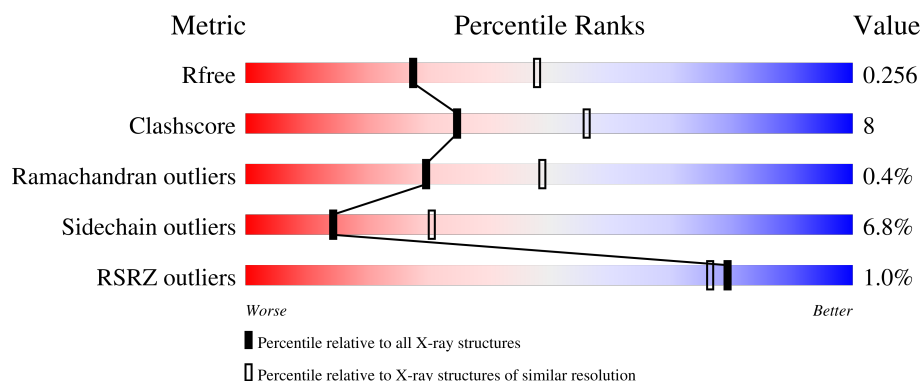
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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}	180053	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (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 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	L	216	 83% 15% .
1	M	216	 81% 17% .
2	H	210	 78% 17% . .
2	I	210	 78% 15% . .
3	P	23	 26% 13% . 57%

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Mol	Chain	Length	Quality of chain
3	Q	23	 A horizontal bar chart showing the quality of chain Q. The bar is divided into three segments: a green segment on the left labeled '39%', a yellow segment in the middle, and a grey segment on the right labeled '57%'. A small black dot is located at the boundary between the yellow and grey segments.

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 (CCD ID: CA) (formula: Ca).

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

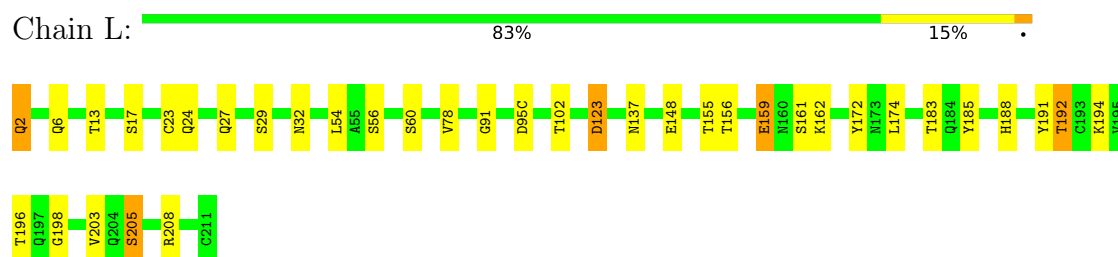
- Molecule 5 is water.

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

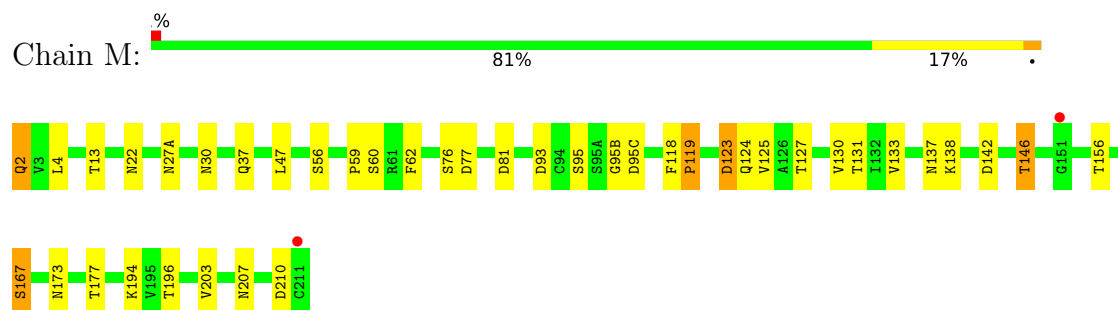
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, 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.

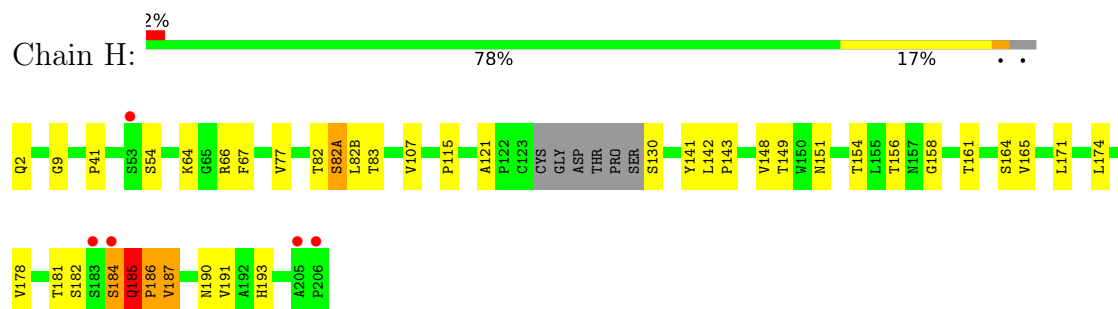
- Molecule 1: monoclonal anti-HIV-1 gp120 V3 antibody R56 light chain



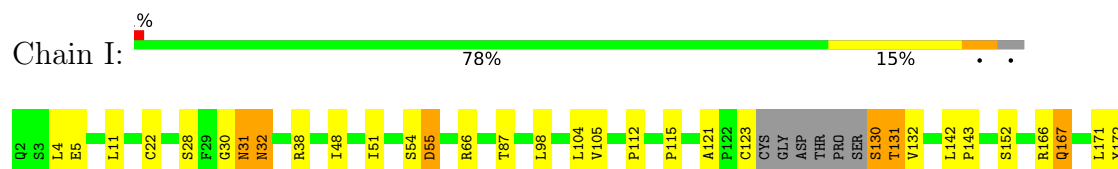
- Molecule 1: monoclonal anti-HIV-1 gp120 V3 antibody R56 light chain

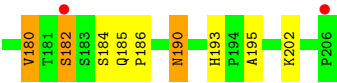


- Molecule 2: monoclonal anti-HIV-1 gp120 V3 antibody R56 heavy chain



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

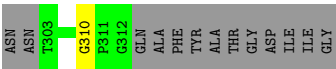




● Molecule 3: gp120



● 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 26.46 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.7 (26.46-2.50) 99.6 (26.46-2.50)	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 0.181 , 0.256	Depositor DCC
R_{free} test set	2127 reflections (5.17%)	wwPDB-VP
Wilson B-factor (Å ²)	25.5	Xtriage
Anisotropy	0.084	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 46.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.42$, $\langle L^2 \rangle = 0.25$	Xtriage
Estimated twinning fraction	0.098 for h,-k,-l	Xtriage
F_o, F_c correlation	0.92	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, 0.333 respectively for untwinned datasets, and 0.375, 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	0.52	0/1628	0.88	1/2229 (0.0%)
1	M	0.52	0/1628	0.92	3/2229 (0.1%)
2	H	0.53	0/1528	0.95	8/2091 (0.4%)
2	I	0.50	0/1528	0.90	4/2091 (0.2%)
3	P	0.55	0/75	0.94	0/98
3	Q	0.48	0/75	1.20	2/98 (2.0%)
All	All	0.52	0/6462	0.92	18/8836 (0.2%)

There are no bond length outliers.

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	I	185	GLN	CA-C-N	7.77	129.55	119.84
2	I	185	GLN	C-N-CA	7.77	129.55	119.84
3	Q	310	GLY	CA-C-N	6.75	126.79	119.90
3	Q	310	GLY	C-N-CA	6.75	126.79	119.90
2	H	121	ALA	CA-C-N	6.67	126.69	119.89
2	H	121	ALA	C-N-CA	6.67	126.69	119.89
2	H	184	SER	CA-C-N	6.55	133.49	121.70
2	H	184	SER	C-N-CA	6.55	133.49	121.70
1	M	118	PHE	CA-C-N	6.03	124.00	119.66
1	M	118	PHE	C-N-CA	6.03	124.00	119.66
2	H	185	GLN	CA-C-N	5.64	126.89	119.84
2	H	185	GLN	C-N-CA	5.64	126.89	119.84
2	H	193	HIS	CA-C-N	5.55	126.77	119.84
2	H	193	HIS	C-N-CA	5.55	126.77	119.84
1	L	137	ASN	N-CA-C	5.45	118.37	109.59
2	I	193	HIS	CA-C-N	5.30	126.46	119.84
2	I	193	HIS	C-N-CA	5.30	126.46	119.84
1	M	119	PRO	O-C-N	5.06	123.53	121.15

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	19	0
1	M	1596	0	1514	29	0
2	H	1494	0	1463	24	0
2	I	1494	0	1463	23	0
3	P	75	0	85	2	0
3	Q	75	0	85	0	0
4	H	1	0	0	0	0
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	9	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	94	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 8.

All (94) 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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:OD2	1:M:123:ASP:N	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
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: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
1:L:2:GLN:OE1	1:L:95(C):ASP:N	2.42	0.52
2:I:142:LEU:HD12	2:I:143:PRO:HA	1.92	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
2:H:54:SER:HB2	5:H:479:HOH:O	2.10	0.50
1:M:119:PRO:HG3	2:I:123:CYS:HB2	1.93	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
1:M:2:GLN:N	1:M:2:GLN:CD	2.72	0.48
2:H:181:THR:OG1	2:H:182:SER:N	2.38	0.48
1:L:2:GLN:HG3	1:L:27:GLN:CD	2.40	0.47
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:H:82(A):SER:OG	5:H:493:HOH:O	2.21	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
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
1:M:124:GLN:O	1:M:127:THR:OG1	2.35	0.43
2:H:149:THR:OG1	2:H:190:ASN:HB2	2.18	0.43
1:M:2:GLN:OE1	1:M:95(B):GLY:HA3	2.19	0.43
1:M:133:VAL:HG22	1:M:177:THR:HG22	2.00	0.43
1:M:138:LYS:H	1:M:173:ASN:HD21	1.62	0.43
1:M:124:GLN:NE2	5:M:428:HOH:O	2.51	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:28:SER:C	2:I:30:GLY:H	2.27	0.41
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:66:ARG:C	2:H:67:PHE:HD2	2.28	0.41
1:L:29:SER:O	1:L:29:SER:OG	2.39	0.41
2:H:142:LEU:HB2	2:H:171:LEU:HD22	2.02	0.41
1:M:4:LEU:HD23	1:M:4:LEU:HA	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
1	M	214/216 (99%)	203 (95%)	11 (5%)	0	100	100
2	H	200/210 (95%)	188 (94%)	10 (5%)	2 (1%)	12	24
2	I	200/210 (95%)	189 (94%)	10 (5%)	1 (0%)	24	43
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%)	30	49

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%)	169 (92%)	14 (8%)	12	25
1	M	183/183 (100%)	172 (94%)	11 (6%)	17	36
2	H	170/175 (97%)	163 (96%)	7 (4%)	27	53
2	I	170/175 (97%)	155 (91%)	15 (9%)	9	20
3	P	8/17 (47%)	6 (75%)	2 (25%)	0	1
3	Q	8/17 (47%)	8 (100%)	0	100	100
All	All	722/750 (96%)	673 (93%)	49 (7%)	14	31

All (49) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	L	2	GLN
1	L	13	THR
1	L	23	CYS
1	L	24	GLN
1	L	123	ASP
1	L	155	THR
1	L	156	THR
1	L	159	GLU
1	L	161	SER
1	L	174	LEU
1	L	183	THR
1	L	192	THR
1	L	196	THR
1	L	205	SER
2	H	77	VAL
2	H	82	THR
2	H	82(A)	SER
2	H	82(B)	LEU
2	H	107	VAL
2	H	156	THR
2	H	187	VAL
3	P	303	THR

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Mol	Chain	Res	Type
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	125	VAL
1	M	130	VAL
1	M	146	THR
1	M	156	THR
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	87	THR
2	I	98	LEU
2	I	130	SER
2	I	131	THR
2	I	132	VAL
2	I	167	GLN
2	I	180	VAL
2	I	182	SER
2	I	190	ASN

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

Mol	Chain	Res	Type
1	L	37	GLN
1	L	173	ASN
2	H	151	ASN
2	H	185	GLN
1	M	27	GLN
1	M	30	ASN
1	M	173	ASN
1	M	188	HIS
2	I	190	ASN

5.3.3 RNA [i](#)

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](#)

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 [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	L	216/216 (100%)	-0.25	0 100 100	8, 19, 56, 72	0
1	M	216/216 (100%)	-0.32	2 (0%) 81 78	9, 19, 49, 59	0
2	H	204/210 (97%)	-0.22	5 (2%) 58 54	8, 23, 43, 68	0
2	I	204/210 (97%)	-0.18	2 (0%) 79 76	6, 22, 42, 61	0
3	P	10/23 (43%)	0.13	0 100 100	16, 27, 40, 51	0
3	Q	10/23 (43%)	0.22	0 100 100	15, 30, 47, 53	0
All	All	860/898 (95%)	-0.23	9 (1%) 79 76	6, 21, 51, 72	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	I	182	SER	3.7
2	H	184	SER	3.3
2	H	206	PRO	2.8
2	I	206	PRO	2.7
2	H	183	SER	2.7
2	H	53	SER	2.3
1	M	211	CYS	2.3
1	M	151	GLY	2.2
2	H	205	ALA	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](#)

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
4	CA	L	302	1/1	0.96	0.05	37,37,37,37	0
4	CA	H	301	1/1	0.98	0.06	29,29,29,29	0
4	CA	M	302	1/1	0.98	0.06	38,38,38,38	0
4	CA	I	301	1/1	0.98	0.06	32,32,32,32	0
4	CA	L	301	1/1	0.99	0.06	16,16,16,16	0
4	CA	M	301	1/1	0.99	0.05	13,13,13,13	0

6.5 Other polymers [i](#)

There are no such residues in this entry.