



Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 05:27 PM GMT

PDB ID : 3MLW
Title : Crystal structure of anti-HIV-1 V3 Fab 1006-15D in complex with an MN V3 peptide
Authors : Kong, X.-P.
Deposited on : 2010-04-18
Resolution : 2.70 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>

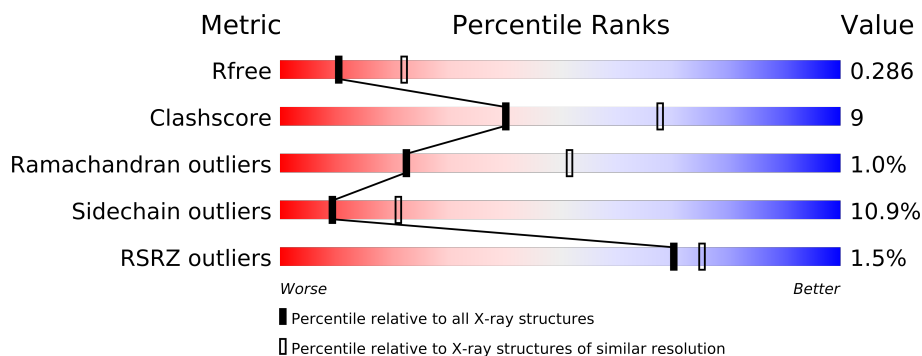
The following versions of software and data (see [references](#)) were used in the production of this report:

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 2.70 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	1557 (2.70-2.70)
Clashscore	79885	1939 (2.70-2.70)
Ramachandran outliers	78287	1905 (2.70-2.70)
Sidechain outliers	78261	1905 (2.70-2.70)
RSRZ outliers	66119	1559 (2.70-2.70)

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	215	
1	M	215	
2	H	228	
2	I	228	
3	P	23	
3	Q	23	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 7170 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 1006-15D Fab light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	215	Total	C	N	O	S	0	0	0
			1621	1010	275	332	4			
1	M	215	Total	C	N	O	S	0	0	0
			1621	1010	275	332	4			

- Molecule 2 is a protein called Human monoclonal anti-HIV-1 gp120 V3 antibody 1006-15D Fab heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	228	Total	C	N	O	S	0	0	0
			1726	1088	289	340	9			
2	I	228	Total	C	N	O	S	0	0	0
			1726	1088	289	340	9			

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

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	P	17	Total	C	N	O	0	0	0
			143	92	31	20			
3	Q	17	Total	C	N	O	0	0	0
			142	90	31	21			

- Molecule 4 is PHOSPHATE ION (three-letter code: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	L	1	Total	O	P	0	0
			5	4	1		
4	H	1	Total	O	P		
			5	4	1		
4	H	1	Total	O	P		
			5	4	1		
4	M	1	Total	O	P	0	0
			5	4	1		
4	I	1	Total	O	P		
			5	4	1		
4	I	1	Total	O	P		
			5	4	1		

- Molecule 5 is water.

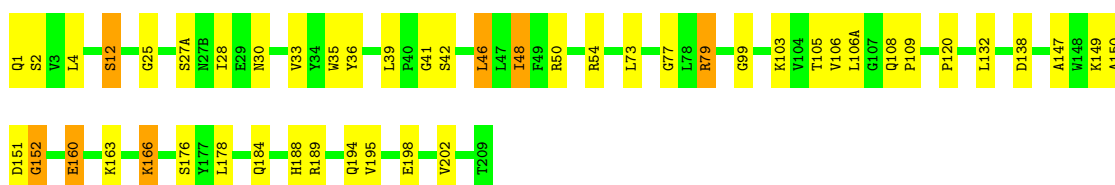
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	L	36	Total	O	0	0
			36	36		
5	H	53	Total	O		
			53	53		
5	P	2	Total	O		
			2	2		
5	M	24	Total	O	0	0
			24	24		
5	I	41	Total	O		
			41	41		
5	Q	5	Total	O		
			5	5		

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.

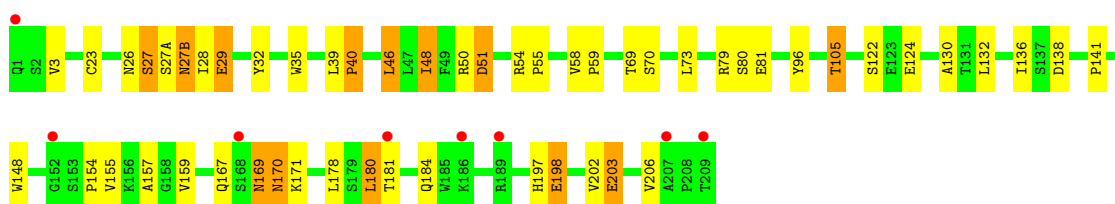
- Molecule 1: Human monoclonal anti-HIV-1 gp120 V3 antibody 1006-15D Fab light chain

Chain L:



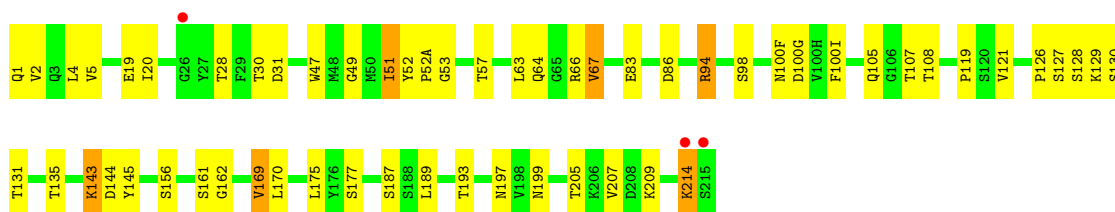
- Molecule 1: Human monoclonal anti-HIV-1 gp120 V3 antibody 1006-15D Fab light chain

Chain M:



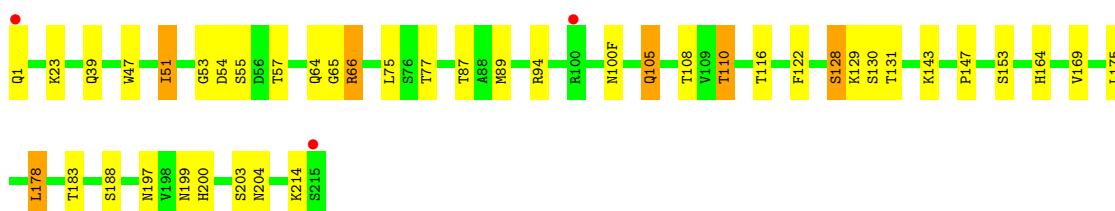
- Molecule 2: Human monoclonal anti-HIV-1 gp120 V3 antibody 1006-15D Fab heavy chain

Chain H:



- Molecule 2: Human monoclonal anti-HIV-1 gp120 V3 antibody 1006-15D Fab heavy chain

Chain I:



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

Chain P: 

TYR	ASN	K303	R304	K305	R306	R315	T320	K321	ASN	ILE	ILE	GLY
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- Molecule 3: HIV-1 gp120 third variable region (V3) crown

Chain Q: 

TYR	N302	K303	R304	K305	R306	T320	LYS	ASN	ILE	ILE	GLY
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4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	98.92Å 82.31Å 149.83Å 90.00° 110.35° 90.00°	Depositor
Resolution (Å)	42.27 – 2.70 42.27 – 2.70	Depositor EDS
% Data completeness (in resolution range)	93.3 (42.27-2.70) 93.4 (42.27-2.70)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.13 (at 2.69Å)	Xtriage
Refinement program	REFMAC 5.5.0091	Depositor
R, R_{free}	0.214 , 0.291 0.212 , 0.286	Depositor DCC
R_{free} test set	1473 reflections (5.34%)	DCC
Wilson B-factor (Å ²)	53.1	Xtriage
Anisotropy	0.041	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 21.1	EDS
Estimated twinning fraction	0.010 for h,-k,-h-l	Xtriage
L-test for twinning	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 29059 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	7170	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.62% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: PO4

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	0.47	0/1662	0.65	1/2269 (0.0%)
1	M	0.45	0/1662	0.59	0/2269
2	H	0.48	0/1771	0.64	0/2411
2	I	0.49	0/1771	0.62	0/2411
3	P	0.47	0/146	0.70	0/192
3	Q	0.47	0/145	0.68	0/192
All	All	0.47	0/7157	0.63	1/9744 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	L	46	LEU	CA-CB-CG	6.24	129.66	115.30

There are no chirality outliers.

There are no planarity outliers.

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	1621	0	1559	32	0
1	M	1621	0	1559	31	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	H	1726	0	1676	33	0
2	I	1726	0	1676	27	0
3	P	143	0	156	12	0
3	Q	142	0	149	5	0
4	H	10	0	0	1	0
4	I	10	0	0	0	0
4	L	5	0	0	1	0
4	M	5	0	0	0	0
5	H	53	0	0	2	0
5	I	41	0	0	4	0
5	L	36	0	0	0	0
5	M	24	0	0	0	0
5	P	2	0	0	0	0
5	Q	5	0	0	0	0
All	All	7170	0	6775	129	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 9.

All (129) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:L:166:LYS:H	1:L:166:LYS:HD2	1.30	0.95
2:H:51:ILE:HG12	2:H:57:THR:HG22	1.50	0.93
2:I:105:GLN:HE21	2:I:105:GLN:H	1.08	0.93
2:I:105:GLN:NE2	2:I:105:GLN:H	1.77	0.82
3:Q:303:LYS:HG2	3:Q:304:ARG:H	1.43	0.81
1:M:27(B):ASN:HD22	1:M:28:ILE:H	1.26	0.81
2:H:100(F):ASN:HA	5:H:268:HOH:O	1.80	0.80
2:I:51:ILE:HG23	2:I:57:THR:HG22	1.65	0.77
1:L:50:ARG:NH1	2:H:100(G):ASP:OD1	2.18	0.76
2:H:156:SER:H	2:H:197:ASN:HD21	1.31	0.75
1:L:166:LYS:HD2	1:L:166:LYS:N	2.06	0.69
1:L:30:ASN:HB2	3:P:315:ARG:HH21	1.57	0.69
3:Q:303:LYS:CG	3:Q:304:ARG:H	2.06	0.68
2:I:105:GLN:HE21	2:I:105:GLN:N	1.88	0.68
1:L:108:GLN:HB2	1:L:109:PRO:HD2	1.76	0.68
2:H:127:SER:H	2:H:130:SER:HB3	1.59	0.66
1:L:30:ASN:HB2	3:P:315:ARG:NH2	2.10	0.66
2:H:2:VAL:HG11	2:H:94:ARG:NH2	2.12	0.64
3:P:303:LYS:HD2	3:P:304:ARG:HG2	1.80	0.64
2:H:20:ILE:HG21	2:H:107:THR:HG21	1.81	0.62
1:L:132:LEU:HD12	1:L:178:LEU:HD23	1.81	0.61

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:M:48:ILE:HD11	1:M:73:LEU:HD13	1.83	0.61
2:H:30:THR:OG1	2:H:53:GLY:HA2	2.01	0.60
3:P:303:LYS:N	3:P:321:LYS:HG3	2.16	0.60
1:L:108:GLN:HB2	1:L:109:PRO:CD	2.32	0.59
3:Q:302:ASN:HD22	3:Q:302:ASN:N	2.01	0.59
3:Q:303:LYS:HG2	3:Q:304:ARG:N	2.14	0.58
2:I:51:ILE:CG2	2:I:57:THR:HG22	2.32	0.58
1:L:4:LEU:HB2	1:L:99:GLY:HA2	1.86	0.57
1:L:12:SER:HB2	1:L:198:GLU:OE1	2.04	0.57
2:I:23:LYS:HG3	2:I:77:THR:OG1	2.05	0.57
2:H:143:LYS:HE3	2:H:144:ASP:OD1	2.04	0.57
1:L:25:GLY:N	1:L:28:ILE:HD11	2.20	0.56
3:P:320:THR:O	3:P:320:THR:HG23	2.06	0.56
1:M:132:LEU:HD12	1:M:178:LEU:HD23	1.89	0.55
1:L:41:GLY:H	4:L:210:PO4:P	2.29	0.54
1:L:28:ILE:HG22	1:L:33:VAL:HG22	1.89	0.54
1:L:48:ILE:HD11	1:L:73:LEU:HD13	1.88	0.54
1:M:202:VAL:HG22	1:M:203:GLU:H	1.72	0.54
1:M:35:TRP:HB2	1:M:48:ILE:CG1	2.38	0.54
2:I:200:HIS:HD2	2:I:203:SER:HB3	1.73	0.53
1:L:79:ARG:O	1:L:106:VAL:HG21	2.09	0.53
1:M:27:SER:HA	1:M:29:GLU:HB2	1.90	0.53
1:M:202:VAL:HG22	1:M:203:GLU:N	2.23	0.53
1:L:151:ASP:OD2	1:L:189:ARG:HB2	2.09	0.52
2:I:64:GLN:HB3	2:I:66:ARG:HG2	1.91	0.52
2:I:188:SER:HB3	5:I:220:HOH:O	2.09	0.52
2:H:30:THR:HG1	2:H:53:GLY:HA2	1.74	0.52
1:M:169:ASN:O	1:M:171:LYS:N	2.43	0.52
1:M:169:ASN:H	1:M:169:ASN:HD22	1.56	0.52
1:L:28:ILE:HG22	1:L:33:VAL:CG2	2.40	0.52
1:L:30:ASN:CB	3:P:315:ARG:HH21	2.21	0.51
3:P:304:ARG:C	3:P:306:ARG:H	2.14	0.51
1:L:35:TRP:HB2	1:L:48:ILE:HG12	1.92	0.51
1:L:77:GLY:O	1:L:79:ARG:HG2	2.11	0.51
2:I:200:HIS:HD2	2:I:203:SER:CB	2.24	0.50
1:M:96:TYR:HB3	2:I:47:TRP:CG	2.47	0.50
1:M:26:ASN:HB3	1:M:27(A):SER:H	1.77	0.50
1:M:181:THR:H	1:M:184:GLN:HG3	1.78	0.49
2:H:30:THR:HA	2:H:52(A):PRO:HB2	1.94	0.49
2:I:147:PRO:O	2:I:200:HIS:HE1	1.96	0.49
2:I:51:ILE:HG23	2:I:57:THR:CG2	2.39	0.48
1:M:155:VAL:HG22	1:M:157:ALA:H	1.78	0.48

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:M:105:THR:HG21	1:M:141:PRO:HB3	1.95	0.48
2:H:209:LYS:HE3	5:H:254:HOH:O	2.12	0.48
2:I:87:THR:HG23	2:I:110:THR:HA	1.94	0.48
2:H:52:TYR:CZ	3:P:305:LYS:HB2	2.48	0.48
2:H:119:PRO:HB3	2:H:145:TYR:HB3	1.95	0.47
2:I:64:GLN:C	2:I:66:ARG:H	2.17	0.47
1:M:181:THR:OG1	1:M:184:GLN:HG2	2.14	0.47
2:H:4:LEU:HD11	2:H:94:ARG:HB3	1.96	0.47
4:H:216:PO4:O3	3:P:304:ARG:NH1	2.47	0.47
1:M:138:ASP:HA	1:M:171:LYS:HB3	1.96	0.47
2:I:39:GLN:HA	5:I:258:HOH:O	2.13	0.47
2:I:100(F):ASN:ND2	5:I:225:HOH:O	2.47	0.47
2:I:51:ILE:CG2	2:I:57:THR:CG2	2.93	0.47
1:M:28:ILE:HD11	1:M:70:SER:HA	1.97	0.47
1:L:184:GLN:O	1:L:188:HIS:HD2	1.97	0.47
3:P:320:THR:HA	3:P:321:LYS:HA	1.67	0.46
2:I:128:SER:HA	2:I:131:THR:HG22	1.97	0.46
1:L:25:GLY:H	1:L:28:ILE:HD11	1.79	0.46
1:M:124:GLU:HG3	2:I:122:PHE:CE2	2.50	0.46
2:H:156:SER:N	2:H:197:ASN:HD21	2.07	0.46
1:L:166:LYS:CD	1:L:166:LYS:N	2.75	0.46
2:I:64:GLN:O	2:I:66:ARG:N	2.49	0.45
3:Q:303:LYS:CG	3:Q:304:ARG:N	2.78	0.45
2:H:66:ARG:NH2	2:H:86:ASP:OD2	2.48	0.45
1:L:120:PRO:HD3	1:L:132:LEU:HD23	1.98	0.45
1:M:39:LEU:O	1:M:40:PRO:C	2.55	0.45
1:M:46:LEU:HD13	1:M:55:PRO:HG3	1.99	0.44
1:L:147:ALA:O	1:L:194:GLN:HB3	2.17	0.44
1:L:35:TRP:HB2	1:L:48:ILE:CG1	2.47	0.44
1:M:148:TRP:CD1	1:M:159:VAL:HG12	2.52	0.44
2:H:63:LEU:HD22	2:H:67:VAL:HG13	1.99	0.44
2:I:178:LEU:C	2:I:178:LEU:HD12	2.38	0.44
2:I:55:SER:HB2	5:I:228:HOH:O	2.17	0.44
1:M:197:HIS:O	1:M:198:GLU:C	2.56	0.44
2:H:2:VAL:HG11	2:H:94:ARG:HH21	1.83	0.44
1:M:35:TRP:HB2	1:M:48:ILE:HG13	1.99	0.43
3:P:320:THR:CG2	3:P:320:THR:O	2.66	0.43
1:L:36:TYR:OH	2:H:100(I):PHE:HB2	2.18	0.43
1:L:166:LYS:H	1:L:166:LYS:CD	2.14	0.43
2:H:47:TRP:CZ2	2:H:49:GLY:HA2	2.53	0.43
1:M:58:VAL:HA	1:M:59:PRO:HD3	1.87	0.42
1:M:35:TRP:CD1	1:M:48:ILE:HG13	2.55	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:L:150:ALA:O	1:L:152:GLY:N	2.48	0.42
2:H:143:LYS:HB2	2:H:177:SER:OG	2.20	0.42
2:H:121:VAL:HG13	2:H:207:VAL:HG11	2.02	0.42
1:L:160:GLU:HB2	2:H:169:VAL:HG22	2.01	0.42
2:H:28:THR:HB	2:H:31:ASP:OD1	2.20	0.42
2:I:153:SER:OG	2:I:197:ASN:HB2	2.19	0.42
1:M:130:ALA:HB3	1:M:180:LEU:O	2.20	0.42
1:M:23:CYS:HB2	1:M:35:TRP:CH2	2.54	0.41
2:H:63:LEU:HD22	2:H:67:VAL:CG1	2.50	0.41
2:I:200:HIS:CD2	2:I:203:SER:CB	3.03	0.41
2:H:127:SER:N	2:H:130:SER:HB3	2.32	0.41
2:I:54:ASP:OD1	2:I:54:ASP:N	2.52	0.41
1:L:160:GLU:HB2	2:H:169:VAL:CG2	2.50	0.41
1:M:32:TYR:CG	1:M:50:ARG:HG2	2.55	0.41
3:P:303:LYS:HD2	3:P:304:ARG:H	1.86	0.41
1:M:154:PRO:O	1:M:155:VAL:HB	2.21	0.41
2:H:66:ARG:NH2	2:H:83:GLU:HG2	2.36	0.41
1:M:54:ARG:HD2	1:M:58:VAL:HG12	2.02	0.41
2:H:126:PRO:O	2:H:214:LYS:HE3	2.20	0.41
2:H:161:SER:HA	2:H:162:GLY:HA2	1.81	0.41
1:L:120:PRO:HD3	1:L:132:LEU:CD2	2.51	0.40
2:H:64:GLN:NE2	2:H:66:ARG:HD3	2.36	0.40
2:H:127:SER:O	2:H:131:THR:HG22	2.21	0.40
1:M:167:GLN:HE21	2:I:164:HIS:CE1	2.39	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	213/215 (99%)	194 (91%)	17 (8%)	2 (1%)	25	55
1	M	213/215 (99%)	190 (89%)	19 (9%)	4 (2%)	12	29
2	H	226/228 (99%)	215 (95%)	11 (5%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	I	226/228 (99%)	213 (94%)	11 (5%)	2 (1%)	25	55
3	P	15/23 (65%)	14 (93%)	1 (7%)	0	100	100
3	Q	15/23 (65%)	14 (93%)	0	1 (7%)	2	2
All	All	908/932 (97%)	840 (92%)	59 (6%)	9 (1%)	22	51

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	M	40	PRO
1	L	2	SER
1	M	170	ASN
2	I	65	GLY
1	M	198	GLU
3	Q	303	LYS
1	M	51	ASP
1	L	152	GLY
2	I	53	GLY

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%)	163 (89%)	20 (11%)	9	21
1	M	183/183 (100%)	164 (90%)	19 (10%)	10	23
2	H	194/194 (100%)	172 (89%)	22 (11%)	9	19
2	I	194/194 (100%)	173 (89%)	21 (11%)	9	21
3	P	14/19 (74%)	14 (100%)	0	100	100
3	Q	14/19 (74%)	11 (79%)	3 (21%)	1	4
All	All	782/792 (99%)	697 (89%)	85 (11%)	9	21

All (85) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	L	1	GLN

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Mol	Chain	Res	Type
1	L	12	SER
1	L	27(A)	SER
1	L	39	LEU
1	L	42	SER
1	L	46	LEU
1	L	48	ILE
1	L	54	ARG
1	L	79	ARG
1	L	103	LYS
1	L	105	THR
1	L	106(A)	LEU
1	L	138	ASP
1	L	149	LYS
1	L	160	GLU
1	L	163	LYS
1	L	166	LYS
1	L	176	SER
1	L	195	VAL
1	L	202	VAL
2	H	1	GLN
2	H	5	VAL
2	H	19	GLU
2	H	51	ILE
2	H	67	VAL
2	H	94	ARG
2	H	98	SER
2	H	105	GLN
2	H	108	THR
2	H	128	SER
2	H	129	LYS
2	H	135	THR
2	H	143	LYS
2	H	169	VAL
2	H	170	LEU
2	H	175	LEU
2	H	187	SER
2	H	189	LEU
2	H	193	THR
2	H	199	ASN
2	H	205	THR
2	H	214	LYS
1	M	3	VAL

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Mol	Chain	Res	Type
1	M	27	SER
1	M	27(B)	ASN
1	M	29	GLU
1	M	46	LEU
1	M	48	ILE
1	M	51	ASP
1	M	69	THR
1	M	79	ARG
1	M	80	SER
1	M	81	GLU
1	M	105	THR
1	M	122	SER
1	M	136	ILE
1	M	169	ASN
1	M	170	ASN
1	M	180	LEU
1	M	203	GLU
1	M	206	VAL
2	I	1	GLN
2	I	51	ILE
2	I	66	ARG
2	I	75	LEU
2	I	89	MET
2	I	94	ARG
2	I	105	GLN
2	I	108	THR
2	I	110	THR
2	I	116	THR
2	I	128	SER
2	I	129	LYS
2	I	130	SER
2	I	143	LYS
2	I	169	VAL
2	I	175	LEU
2	I	178	LEU
2	I	183	THR
2	I	199	ASN
2	I	204	ASN
2	I	214	LYS
3	Q	302	ASN
3	Q	303	LYS
3	Q	306	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	L	108	GLN
1	L	126	GLN
1	L	188	HIS
2	H	64	GLN
2	H	171	GLN
2	H	197	ASN
3	P	308	HIS
1	M	26	ASN
1	M	27(B)	ASN
1	M	167	GLN
1	M	169	ASN
1	M	170	ASN
1	M	194	GLN
2	I	1	GLN
2	I	32	HIS
2	I	81	GLN
2	I	96	HIS
2	I	101	HIS
2	I	105	GLN
2	I	200	HIS
3	Q	308	HIS

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 ⓘ

6 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	PO4	H	216	-	4,4,4	0.30	0	6,6,6	0.31	0
4	PO4	H	217	-	4,4,4	0.23	0	6,6,6	0.31	0
4	PO4	I	216	-	4,4,4	0.23	0	6,6,6	0.32	0
4	PO4	I	217	-	4,4,4	0.22	0	6,6,6	0.31	0
4	PO4	L	210	-	4,4,4	0.27	0	6,6,6	0.31	0
4	PO4	M	210	-	4,4,4	0.27	0	6,6,6	0.31	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	PO4	H	216	-	-	0/0/0/0	0/0/0/0
4	PO4	H	217	-	-	0/0/0/0	0/0/0/0
4	PO4	I	216	-	-	0/0/0/0	0/0/0/0
4	PO4	I	217	-	-	0/0/0/0	0/0/0/0
4	PO4	L	210	-	-	0/0/0/0	0/0/0/0
4	PO4	M	210	-	-	0/0/0/0	0/0/0/0

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.

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	215/215 (100%)	-0.13	0 100 100	30, 44, 54, 61	0
1	M	215/215 (100%)	0.32	8 (3%) 39 44	36, 54, 75, 78	0
2	H	228/228 (100%)	-0.16	3 (1%) 74 79	30, 42, 55, 69	0
2	I	228/228 (100%)	0.06	3 (1%) 74 79	32, 51, 64, 76	0
3	P	17/23 (73%)	-0.02	0 100 100	44, 48, 57, 58	0
3	Q	17/23 (73%)	0.15	0 100 100	46, 48, 55, 57	0
All	All	920/932 (98%)	0.02	14 (1%) 70 75	30, 47, 68, 78	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	H	215	SER	4.5
2	I	100	ARG	4.0
2	I	1	GLN	3.9
1	M	168	SER	3.9
1	M	152	GLY	3.6
1	M	209	THR	3.2
1	M	186	LYS	3.1
1	M	189	ARG	2.7
2	H	214	LYS	2.6
1	M	181	THR	2.4
1	M	1	GLN	2.3
1	M	207	ALA	2.1
2	I	215	SER	2.0
2	H	26	GLY	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

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	RSR	LLDF	B-factors(Å ²)	Q<0.9
4	PO4	I	216	5/5	0.15	-	75,76,76,76	0
4	PO4	H	217	5/5	0.17	-	84,85,85,85	0
4	PO4	I	217	5/5	0.09	-	79,80,80,80	0
4	PO4	L	210	5/5	0.17	-	112,112,112,112	0
4	PO4	M	210	5/5	0.21	-	94,95,95,95	0
4	PO4	H	216	5/5	0.15	-	93,93,93,93	0

6.5 Other polymers

There are no such residues in this entry.