Full wwPDB X-ray Structure Validation Report (i)

Jun 22, 2024 – 08:02 PM EDT

PDB ID	:	5EOC
Title	:	Crystal structure of Fab C2 in complex with a Cyclic variant of Hepatitis C Virus E2 epitope I
Authors	:	Berisio, R.; Ruggiero, A.; Sandomenico, A.; Patel, A.H.; Ruvo, M.; Vitagliano, L.
Deposited on	:	2015-11-10
Resolution	:	1.98 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 (i) 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 (1)) were used in the production of this report:

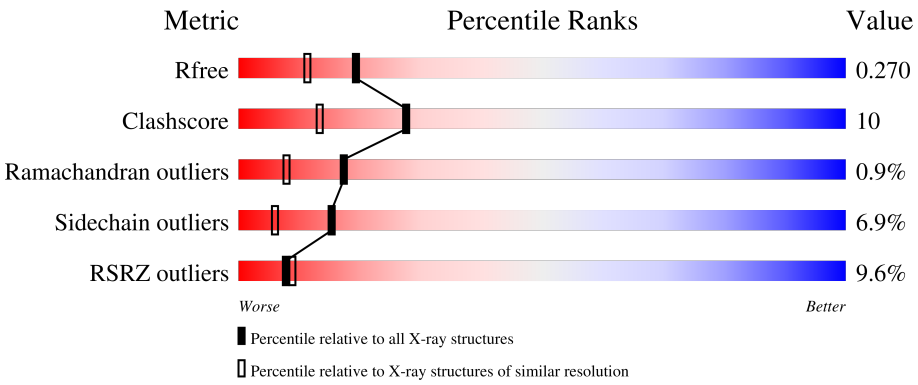
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.37.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.98 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}	130704	11647 (2.00-1.96)
Clashscore	141614	1014 (1.98-1.98)
Ramachandran outliers	138981	1006 (1.98-1.98)
Sidechain outliers	138945	1006 (1.98-1.98)
RSRZ outliers	127900	11410 (2.00-1.96)

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 <=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	H	217	5%, 76%, 20%, • •
2	L	216	8%, 81%, 15%, •
2	M	216	13%, 75%, 20%, •
3	J	217	6%, 79%, 18%, • •
4	P	14	43%, 50%, 36%, 7%, 7%

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Mol	Chain	Length	Quality of chain
4	Q	14	57% 71% 14% 7% 7%

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 7216 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 Fab fragment (Heavy chain).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	H	212	1597	1010	264	315	8	0	0	0

- Molecule 2 is a protein called Fab fragment (Light chain).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	L	216	1670	1038	291	335	6	0	1	0
2	M	215	1655	1031	285	333	6	0	0	0

- Molecule 3 is a protein called Fab fragment (Heavy chain).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	J	214	1614	1020	267	318	9	0	0	0

- Molecule 4 is a protein called ALA-CYS-GLN-LEU-ILE-ASN-THR-ASN-GLY-SER-TRP-HIS-ILE-CYS.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
4	P	13	102	63	19	18	2	0	0	0
4	Q	13	102	63	19	18	2	0	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
5	H	134	134	134	0	0

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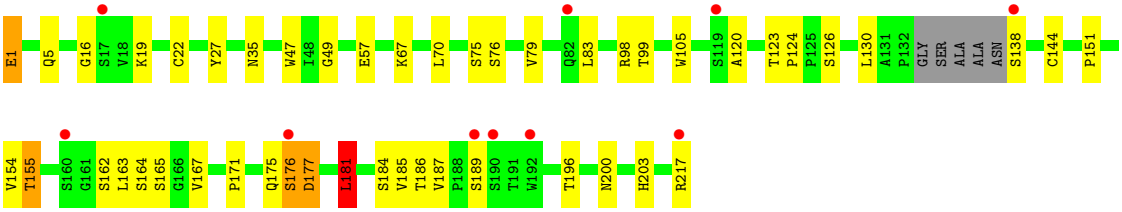
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	L	99	Total O 99 99	0	0
5	J	123	Total O 123 123	0	0
5	M	116	Total O 116 116	0	0
5	P	1	Total O 1 1	0	0
5	Q	3	Total O 3 3	0	0

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.

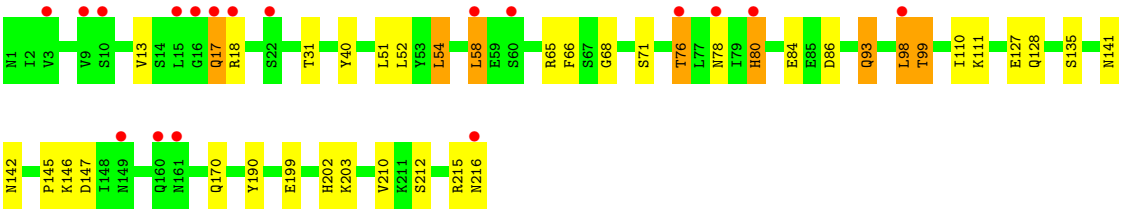
• Molecule 1: Fab fragment (Heavy chain)

Chain H:



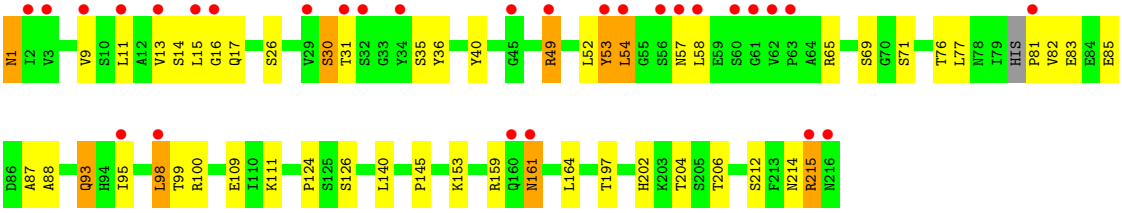
• Molecule 2: Fab fragment (Light chain)

Chain L:



• Molecule 2: Fab fragment (Light chain)

Chain M:

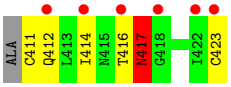


• Molecule 3: Fab fragment (Heavy chain)

Chain J:

- Molecule 4: ALA-CYS-GLN-LEU-ILE-ASN-THR-ASN-GLY-SER-TRP-HIS-ILE-CYS

Chain P:



- Molecule 4: ALA-CYS-GLN-LEU-ILE-ASN-THR-ASN-GLY-SER-TRP-HIS-ILE-CYS

Chain Q:

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α, β, γ	54.97Å 56.19Å 77.15Å 90.35° 90.18° 94.98°	Depositor
Resolution (Å)	15.00 – 1.98 10.00 – 1.98	Depositor EDS
% Data completeness (in resolution range)	92.3 (15.00-1.98) 92.9 (10.00-1.98)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.51 (at 1.98Å)	Xtriage
Refinement program	REFMAC 5.5.0110	Depositor
R, R_{free}	0.234 , 0.266 0.237 , 0.270	Depositor DCC
R_{free} test set	3000 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å^2)	26.7	Xtriage
Anisotropy	0.074	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.43 , 63.2	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.018 for -h,-k,l 0.013 for k,h,-l 0.032 for -k,-h,-l	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	7216	wwPDB-VP
Average B, all atoms (Å^2)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.11% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality (i)

5.1 Standard geometry (i)

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	H	0.70	2/1640 (0.1%)	0.75	1/2247 (0.0%)
2	L	0.63	0/1710	0.76	0/2320
2	M	0.67	0/1691	0.77	0/2293
3	J	0.66	0/1657	0.76	3/2267 (0.1%)
4	P	0.52	0/104	0.88	0/141
4	Q	0.48	0/104	0.93	1/141 (0.7%)
All	All	0.66	2/6906 (0.0%)	0.76	5/9409 (0.1%)

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	L	0	1
2	M	0	1
4	P	0	1
4	Q	0	1
All	All	0	4

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	H	22	CYS	CB-SG	-7.81	1.69	1.82
1	H	144	CYS	CB-SG	-6.12	1.71	1.82

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	J	181	LEU	CA-CB-CG	7.21	131.88	115.30
3	J	98	ARG	NE-CZ-NH1	6.17	123.39	120.30
1	H	181	LEU	CA-CB-CG	6.02	129.15	115.30
4	Q	417	ASN	CB-CA-C	5.30	120.99	110.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	J	35	MET	CB-CG-SD	-5.04	97.29	112.40

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	L	98	LEU	Peptide
2	M	16	GLY	Peptide
4	P	417	ASN	Peptide
4	Q	417	ASN	Peptide

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	H	1597	0	1549	30	0
2	L	1670	0	1608	27	0
2	M	1655	0	1597	42	0
3	J	1614	0	1572	39	0
4	P	102	0	92	4	0
4	Q	102	0	92	2	0
5	H	134	0	0	12	0
5	J	123	0	0	9	0
5	L	99	0	0	2	0
5	M	116	0	0	8	0
5	P	1	0	0	0	0
5	Q	3	0	0	0	0
All	All	7216	0	6510	132	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 10.

All (132) 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:H:1:GLU:N	1:H:1:GLU:OE1	1.69	1.25

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:J:98:ARG:NH1	3:J:99:THR:O	1.98	0.97
1:H:98:ARG:NH1	1:H:99:THR:O	1.99	0.95
1:H:35:ASN:HD22	1:H:47:TRP:HE1	1.21	0.89
2:M:95:ILE:HD12	2:M:95:ILE:O	1.72	0.88
1:H:196:THR:HG22	5:H:335:HOH:O	1.80	0.79
2:M:161:ASN:HD22	2:M:161:ASN:H	1.31	0.77
2:M:31:THR:HG21	4:Q:414:ILE:HG12	1.69	0.75
2:L:31:THR:CB	4:P:414:ILE:HG12	2.18	0.74
2:L:110:ILE:H	2:L:170:GLN:HE22	1.35	0.74
3:J:98:ARG:HG2	3:J:98:ARG:HH11	1.53	0.73
1:H:98:ARG:HG2	1:H:98:ARG:HH11	1.55	0.72
2:M:53:TYR:O	2:M:57:ASN:HB2	1.89	0.71
1:H:167:VAL:HG12	1:H:185:VAL:HG23	1.72	0.71
2:M:53:TYR:O	2:M:53:TYR:CG	2.43	0.71
3:J:203:HIS:HB3	3:J:208:THR:CG2	2.22	0.69
3:J:35:MET:HE2	3:J:106:LEU:HD21	1.76	0.68
3:J:93:VAL:HG23	5:J:356:HOH:O	1.95	0.66
3:J:206:SER:OG	3:J:208:THR:HG22	1.96	0.64
3:J:196:THR:HG22	5:J:382:HOH:O	1.99	0.63
3:J:203:HIS:HB3	3:J:208:THR:HG23	1.80	0.63
2:L:65:ARG:NH2	2:L:86:ASP:OD1	2.32	0.62
2:L:40:TYR:OH	2:L:93:GLN:NE2	2.32	0.62
3:J:35:MET:CE	3:J:106:LEU:HD21	2.30	0.62
2:L:215:ARG:O	2:L:216:ASN:HB2	2.01	0.61
2:M:161:ASN:HD22	2:M:161:ASN:N	1.98	0.61
4:P:416:THR:O	4:P:417:ASN:HB2	2.00	0.61
2:L:98:LEU:O	2:L:99:THR:OG1	2.17	0.60
3:J:173:VAL:HG11	2:M:164:LEU:HD13	1.84	0.59
3:J:59:ARG:HB3	2:M:98:LEU:HD11	1.84	0.59
2:L:141:ASN:HB3	2:L:142:ASN:HD22	1.69	0.58
3:J:59:ARG:CB	2:M:98:LEU:HD11	2.34	0.58
1:H:151:PRO:O	1:H:203:HIS:HE1	1.88	0.57
1:H:120:ALA:HB2	1:H:177:ASP:OD2	2.03	0.57
1:H:167:VAL:HG23	5:H:406:HOH:O	2.05	0.57
5:H:325:HOH:O	2:L:127:GLU:HG2	2.05	0.57
2:M:83:GLU:HB3	5:M:346:HOH:O	2.05	0.57
3:J:59:ARG:HB3	2:M:98:LEU:CD1	2.34	0.56
1:H:163:LEU:HD21	1:H:187:VAL:HG12	1.85	0.56
2:L:58:LEU:HD11	2:L:66:PHE:O	2.06	0.55
2:M:53:TYR:O	2:M:53:TYR:CD2	2.59	0.55
2:M:65:ARG:HD2	2:M:81:PRO:O	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:13:VAL:HG13	2:L:17:GLN:HB3	1.89	0.54
2:M:36:TYR:CD2	2:M:95:ILE:HD11	2.43	0.54
1:H:27:TYR:CZ	1:H:98:ARG:HD3	2.44	0.53
2:L:76:THR:CG2	2:L:78:ASN:ND2	2.71	0.53
3:J:47:TRP:CG	2:M:100:ARG:HB2	2.44	0.53
2:M:15:LEU:HD13	2:M:15:LEU:O	2.09	0.53
1:H:98:ARG:NH1	1:H:98:ARG:HG2	2.24	0.52
3:J:134:SER:HA	5:J:386:HOH:O	2.09	0.52
2:L:65:ARG:O	2:L:80:HIS:N	2.38	0.52
2:L:51:LEU:O	2:L:52:LEU:HD12	2.10	0.51
1:H:79:VAL:HG13	5:H:327:HOH:O	2.11	0.51
2:M:14:SER:HA	2:M:111:LYS:HB2	1.92	0.51
3:J:173:VAL:CG1	2:M:164:LEU:HD13	2.40	0.51
3:J:99:THR:HA	3:J:105:TRP:O	2.09	0.51
2:M:109:GLU:HG2	5:M:389:HOH:O	2.11	0.51
1:H:171:PRO:HD3	5:H:318:HOH:O	2.11	0.50
1:H:181:LEU:C	1:H:181:LEU:HD23	2.32	0.50
2:L:51:LEU:C	2:L:52:LEU:HD12	2.31	0.50
2:M:52:LEU:HD12	2:M:77:LEU:HD13	1.93	0.50
4:P:411:CYS:HA	4:P:423:CYS:HA	1.93	0.50
1:H:57:GLU:HG3	5:H:302:HOH:O	2.11	0.50
3:J:152:GLU:OE2	5:J:301:HOH:O	2.20	0.50
2:L:52:LEU:HD23	2:L:68:GLY:HA3	1.94	0.50
2:L:98:LEU:C	2:L:98:LEU:HD12	2.32	0.50
2:L:128:GLN:NE2	5:L:306:HOH:O	2.45	0.49
2:L:17:GLN:HG3	2:L:18:ARG:H	1.77	0.49
2:M:40:TYR:OH	2:M:93:GLN:NE2	2.46	0.49
2:L:199:GLU:HG2	2:L:210:VAL:HG22	1.94	0.49
1:H:19:LYS:NZ	5:H:308:HOH:O	2.46	0.49
3:J:72:LEU:HD12	3:J:72:LEU:O	2.13	0.49
2:M:202:HIS:HD2	2:M:204:THR:OG1	1.96	0.48
2:L:190:TYR:CE1	2:L:215:ARG:HD3	2.48	0.48
3:J:20:LEU:HD12	3:J:81:MET:HE2	1.96	0.48
2:M:214:ASN:O	2:M:215:ARG:HG2	2.13	0.48
2:L:128:GLN:NE2	2:L:135:SER:H	2.12	0.47
3:J:33:TRP:CE2	3:J:52:GLN:HB2	2.49	0.47
3:J:212:LYS:HE2	5:J:310:HOH:O	2.13	0.47
2:M:13:VAL:HG12	2:M:82:VAL:HG21	1.97	0.47
3:J:35:MET:CE	3:J:47:TRP:HE1	2.27	0.47
2:M:145:PRO:O	2:M:202:HIS:HE1	1.98	0.47
1:H:105:TRP:CZ3	2:L:54:LEU:HD22	2.50	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:J:72:LEU:HD12	3:J:72:LEU:C	2.35	0.47
3:J:176:SER:O	3:J:177:ASP:HB3	2.14	0.47
3:J:138:SER:HB3	5:J:418:HOH:O	2.14	0.46
3:J:151:PRO:O	3:J:203:HIS:HE1	1.98	0.46
1:H:163:LEU:HD21	1:H:187:VAL:CG1	2.46	0.46
2:M:95:ILE:O	2:M:95:ILE:CD1	2.56	0.46
1:H:175:GLN:O	1:H:176:SER:CB	2.65	0.45
1:H:186:THR:HG23	5:H:303:HOH:O	2.15	0.45
3:J:191:THR:O	3:J:195:GLU:HB2	2.16	0.45
1:H:67:LYS:HG3	5:H:421:HOH:O	2.16	0.45
2:L:145:PRO:O	2:L:202:HIS:HE1	2.00	0.45
2:L:98:LEU:HD23	4:P:417:ASN:ND2	2.32	0.44
2:M:17:GLN:CB	5:M:415:HOH:O	2.66	0.44
1:H:120:ALA:CB	1:H:177:ASP:OD2	2.66	0.44
3:J:47:TRP:CD2	2:M:100:ARG:HB2	2.52	0.44
3:J:51:ILE:HG13	3:J:58:THR:HG22	1.99	0.44
2:M:153:LYS:HB2	2:M:197:THR:HB	2.00	0.44
2:M:76:THR:HB	5:M:386:HOH:O	2.18	0.43
2:M:49:ARG:HE	2:M:49:ARG:HB2	1.67	0.43
1:H:164:SER:HB2	5:H:397:HOH:O	2.19	0.42
2:M:36:TYR:HA	2:M:54:LEU:HD23	2.01	0.42
2:M:206:THR:HG23	5:M:331:HOH:O	2.19	0.42
3:J:59:ARG:HB3	2:M:98:LEU:CD2	2.49	0.42
3:J:98:ARG:NH1	3:J:98:ARG:HG2	2.27	0.42
2:M:30:SER:HA	2:M:35:SER:HA	2.01	0.42
2:M:140:LEU:HD12	2:M:140:LEU:N	2.35	0.42
3:J:93:VAL:CG2	5:J:356:HOH:O	2.62	0.42
2:L:146:LYS:NZ	5:L:303:HOH:O	2.41	0.42
2:L:13:VAL:CG1	2:L:17:GLN:HB3	2.50	0.42
4:Q:417:ASN:C	4:Q:418:GLY:O	2.55	0.41
3:J:159:ASN:HD21	3:J:197:VAL:HA	1.85	0.41
1:H:155:THR:HG22	5:H:346:HOH:O	2.21	0.41
2:M:85:GLU:HG2	5:M:346:HOH:O	2.20	0.41
1:H:123:THR:HA	1:H:124:PRO:HD3	1.91	0.41
3:J:35:MET:SD	3:J:50:MET:HB3	2.60	0.41
2:M:52:LEU:HD21	2:M:58:LEU:HD13	2.02	0.41
3:J:159:ASN:ND2	5:J:313:HOH:O	2.54	0.41
3:J:164:SER:HB2	5:J:322:HOH:O	2.21	0.41
1:H:47:TRP:CH2	1:H:49:GLY:HA2	2.56	0.41
3:J:68:ALA:HA	3:J:82:GLN:O	2.20	0.41
2:M:87:ALA:O	2:M:88:ALA:HB2	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:M:124:PRO:HA	5:M:357:HOH:O	2.20	0.41
3:J:8:GLY:HA2	3:J:111:GLN:HE22	1.86	0.40
1:H:27:TYR:CZ	1:H:98:ARG:CD	3.05	0.40
1:H:165:SER:HB3	5:H:397:HOH:O	2.21	0.40
2:M:1:ASN:HB3	5:M:311:HOH:O	2.20	0.40
2:M:1:ASN:OD1	2:M:1:ASN:N	2.50	0.40
1:H:67:LYS:O	1:H:83:LEU:HA	2.22	0.40
2:L:40:TYR:HE2	2:L:93:GLN:HE21	1.61	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	H	208/217 (96%)	197 (95%)	9 (4%)	2 (1%)	15 6
2	L	215/216 (100%)	202 (94%)	11 (5%)	2 (1%)	17 8
2	M	211/216 (98%)	199 (94%)	10 (5%)	2 (1%)	17 8
3	J	210/217 (97%)	200 (95%)	10 (5%)	0	100 100
4	P	11/14 (79%)	9 (82%)	1 (9%)	1 (9%)	1 0
4	Q	11/14 (79%)	8 (73%)	2 (18%)	1 (9%)	1 0
All	All	866/894 (97%)	815 (94%)	43 (5%)	8 (1%)	17 8

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	H	176	SER
2	L	80	HIS
2	M	215	ARG
4	P	417	ASN

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Mol	Chain	Res	Type
1	H	16	GLY
2	M	53	TYR
4	Q	417	ASN
2	L	17	GLN

5.3.2 Protein sidechains (i)

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	H	182/188 (97%)	165 (91%)	17 (9%)	9 2
2	L	188/190 (99%)	177 (94%)	11 (6%)	19 8
2	M	187/190 (98%)	171 (91%)	16 (9%)	10 2
3	J	185/188 (98%)	177 (96%)	8 (4%)	29 17
4	P	12/12 (100%)	11 (92%)	1 (8%)	11 3
4	Q	12/12 (100%)	12 (100%)	0	100 100
All	All	766/780 (98%)	713 (93%)	53 (7%)	15 6

All (53) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	H	1	GLU
1	H	5	GLN
1	H	70	LEU
1	H	75	SER
1	H	76	SER
1	H	126	SER
1	H	130	LEU
1	H	138	SER
1	H	154	VAL
1	H	155	THR
1	H	162	SER
1	H	177	ASP
1	H	181	LEU
1	H	184	SER

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Mol	Chain	Res	Type
1	H	189	SER
1	H	200	ASN
1	H	217	ARG
2	L	54	LEU
2	L	58	LEU
2	L	71	SER
2	L	76	THR
2	L	84	GLU
2	L	93	GLN
2	L	99	THR
2	L	111	LYS
2	L	147	ASP
2	L	203	LYS
2	L	212	SER
3	J	62	GLN
3	J	63	LYS
3	J	72	LEU
3	J	77	SER
3	J	111	GLN
3	J	154	VAL
3	J	181	LEU
3	J	200	ASN
2	M	1	ASN
2	M	9	VAL
2	M	11	LEU
2	M	26	SER
2	M	30	SER
2	M	49	ARG
2	M	54	LEU
2	M	69	SER
2	M	71	SER
2	M	93	GLN
2	M	98	LEU
2	M	99	THR
2	M	126	SER
2	M	159	ARG
2	M	161	ASN
2	M	212	SER
4	P	412	GLN

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

Mol	Chain	Res	Type
1	H	35	ASN
1	H	52	GLN
1	H	159	ASN
1	H	203	HIS
2	L	57	ASN
2	L	93	GLN
2	L	94	HIS
2	L	128	GLN
2	L	142	ASN
2	L	170	GLN
2	L	193	HIS
2	L	202	HIS
3	J	52	GLN
3	J	111	GLN
3	J	159	ASN
3	J	203	HIS
2	M	93	GLN
2	M	149	ASN
2	M	161	ASN
2	M	170	GLN
2	M	202	HIS

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 monosaccharides 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 (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	H	212/217 (97%)	0.34	10 (4%)	31	33	19, 32, 46, 57	0
2	L	216/216 (100%)	0.65	18 (8%)	11	12	23, 36, 52, 62	0
2	M	215/216 (99%)	0.75	29 (13%)	3	3	18, 36, 55, 61	0
3	J	214/217 (98%)	0.46	14 (6%)	18	20	18, 32, 47, 53	0
4	P	13/14 (92%)	2.30	6 (46%)	0	0	57, 61, 66, 68	0
4	Q	13/14 (92%)	2.79	8 (61%)	0	0	61, 62, 69, 70	0
All	All	883/894 (98%)	0.61	85 (9%)	8	9	18, 34, 55, 70	0

All (85) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	M	9	VAL	7.2
2	L	9	VAL	6.3
2	L	216	ASN	5.9
4	P	414	ILE	5.3
2	L	80	HIS	5.1
4	Q	411	CYS	5.1
3	J	137	ASN	5.0
3	J	133	GLY	4.8
2	M	15	LEU	4.6
4	Q	412	GLN	4.5
2	M	16	GLY	4.2
2	M	53	TYR	4.1
3	J	134	SER	4.0
2	L	17	GLN	3.8
1	H	119	SER	3.7
4	Q	414	ILE	3.7
4	Q	422	ILE	3.7
2	L	15	LEU	3.7
2	M	32	SER	3.7

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Mol	Chain	Res	Type	RSRZ
4	P	422	ILE	3.7
2	L	98	LEU	3.6
2	L	18	ARG	3.6
2	M	60	SER	3.5
3	J	31	THR	3.5
4	Q	416	THR	3.5
2	M	54	LEU	3.5
2	M	34	TYR	3.4
3	J	119	SER	3.4
1	H	189	SER	3.2
4	P	416	THR	3.2
2	M	160	GLN	3.2
2	M	57	ASN	3.2
3	J	28	SER	3.1
2	M	215	ARG	3.1
3	J	32	TYR	3.1
4	P	418	GLY	3.1
1	H	190	SER	3.0
4	Q	417	ASN	3.0
4	Q	413	LEU	3.0
3	J	207	SER	2.9
4	P	412	GLN	2.9
2	M	62	VAL	2.9
2	L	58	LEU	2.9
1	H	176	SER	2.8
2	L	149	ASN	2.8
2	M	98	LEU	2.7
2	M	81	PRO	2.7
2	L	60	SER	2.7
3	J	24	ALA	2.7
2	M	161	ASN	2.6
2	L	160	GLN	2.6
2	M	31	THR	2.6
1	H	160	SER	2.5
2	L	78	ASN	2.5
3	J	100	GLY	2.4
2	M	95	ILE	2.4
2	M	58	LEU	2.4
1	H	217	ARG	2.4
3	J	177	ASP	2.4
2	L	76	THR	2.4
3	J	30	THR	2.4

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Mol	Chain	Res	Type	RSRZ
4	Q	420	TRP	2.3
3	J	5	GLN	2.3
2	M	63	PRO	2.3
1	H	17	SER	2.3
2	M	29	VAL	2.3
2	M	56	SER	2.3
2	L	161	ASN	2.3
2	M	45	GLY	2.3
2	L	3	VAL	2.2
4	P	423	CYS	2.2
2	M	11	LEU	2.2
2	M	49	ARG	2.2
2	L	22	SER	2.2
2	L	16	GLY	2.2
2	L	10	SER	2.2
2	M	61	GLY	2.2
1	H	138	SER	2.1
2	M	3	VAL	2.1
2	M	2	ILE	2.1
2	M	216	ASN	2.1
2	M	13	VAL	2.0
1	H	192	TRP	2.0
1	H	82	GLN	2.0
3	J	25	SER	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 monosaccharides 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.