Full wwPDB X-ray Structure Validation Report ⓘ

May 19, 2025 – 01:11 PM EDT

PDB ID	:	9MJD / pdb_00009mjd
Title	:	Crystal structure of the VRC01-class antibody 3G08
Authors	:	Agrawal, S.; Wilson, I.A.
Deposited on	:	2024-12-14
Resolution	:	3.50 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:

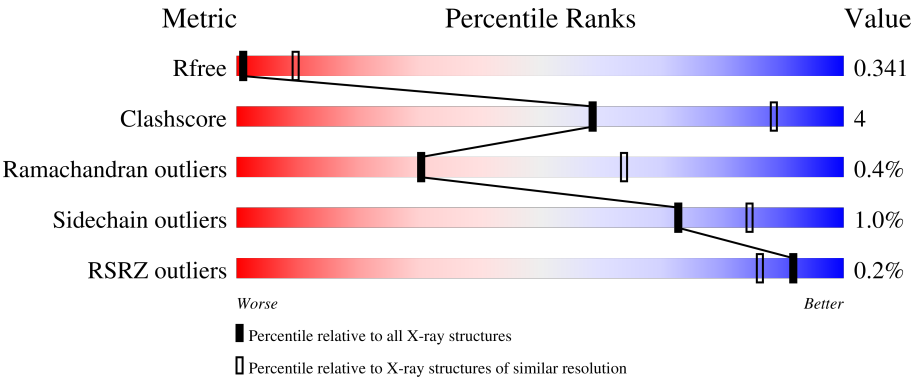
MolProbity	:	4-5-2 with Phenix2.0rc1
Xtriage (Phenix)	:	2.0rc1
EDS	:	3.0
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.006 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.43.1

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 3.50 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}	164625	1094 (3.56-3.44)
Clashscore	180529	1045 (3.54-3.46)
Ramachandran outliers	177936	1032 (3.54-3.46)
Sidechain outliers	177891	1033 (3.54-3.46)
RSRZ outliers	164620	1093 (3.56-3.44)

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	E	223	79% 14% • 6%
1	I	223	86% 9% 6%
1	M	223	86% 9% 6%
1	P	223	% 79% 13% • 6%
2	F	210	86% 7% 7%

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Mol	Chain	Length	Quality of chain
2	J	210	
2	N	210	
2	R	210	

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 12328 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 3G08 Fab heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	P	210	Total	C	N	O	S	0	0	0
			1621	1031	274	308	8			
1	E	210	Total	C	N	O	S	0	0	0
			1621	1031	274	308	8			
1	I	210	Total	C	N	O	S	0	0	0
			1621	1031	274	308	8			
1	M	210	Total	C	N	O	S	0	0	0
			1621	1031	274	308	8			

- Molecule 2 is a protein called 3G08 Fab light chain.

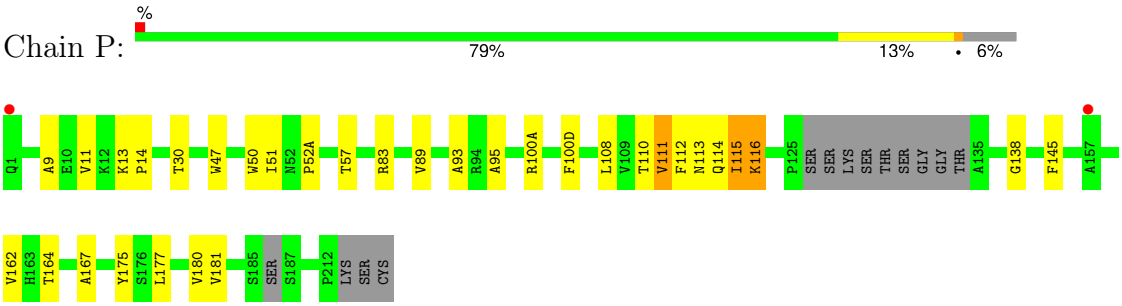
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	R	198	Total	C	N	O	S	0	0	0
			1471	922	244	301	4			
2	F	196	Total	C	N	O	S	0	0	0
			1451	909	241	297	4			
2	J	198	Total	C	N	O	S	0	0	0
			1471	922	244	301	4			
2	N	196	Total	C	N	O	S	0	0	0
			1451	909	241	297	4			

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.

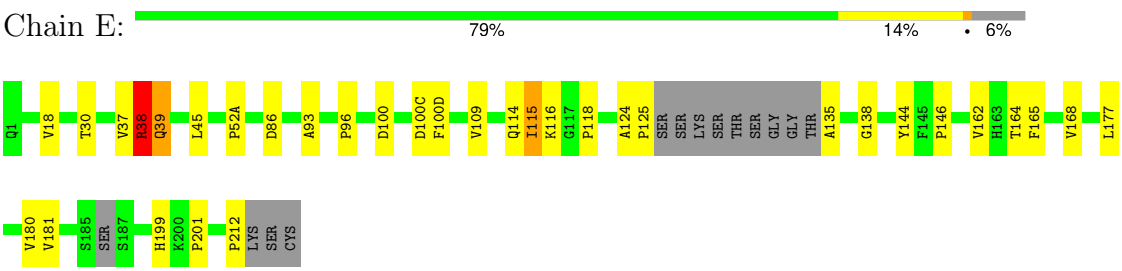
• Molecule 1: 3G08 Fab heavy chain

Chain P:



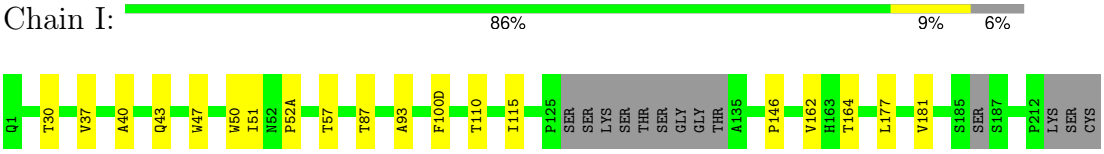
• Molecule 1: 3G08 Fab heavy chain

Chain E:



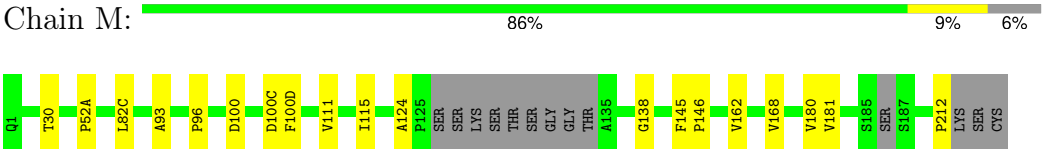
• Molecule 1: 3G08 Fab heavy chain

Chain I:



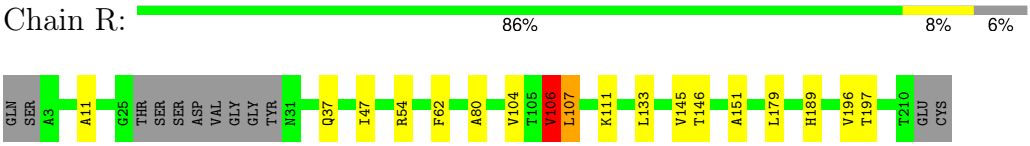
• Molecule 1: 3G08 Fab heavy chain

Chain M:



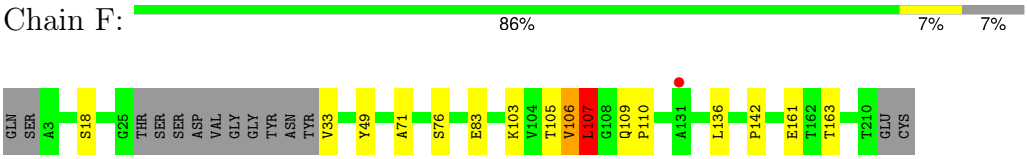
• Molecule 2: 3G08 Fab light chain

Chain R:



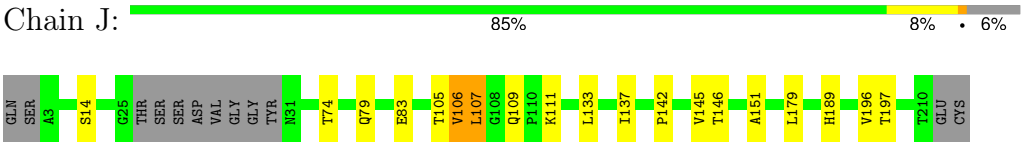
• Molecule 2: 3G08 Fab light chain

Chain F:



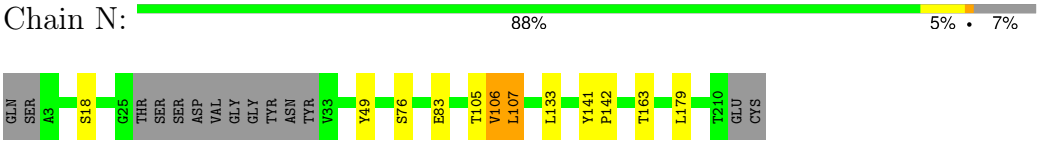
• Molecule 2: 3G08 Fab light chain

Chain J:



• Molecule 2: 3G08 Fab light chain

Chain N:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 32	Depositor
Cell constants a, b, c, α, β, γ	107.93Å 107.93Å 223.56Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	39.59 – 3.50 39.59 – 3.50	Depositor EDS
% Data completeness (in resolution range)	94.6 (39.59-3.50) 82.2 (39.59-3.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	0.49 (at 3.48Å)	Xtriage
Refinement program	PHENIX 1.21.2_5419	Depositor
R, R_{free}	0.306 , 0.343 0.309 , 0.341	Depositor DCC
R_{free} test set	28484 reflections (5.60%)	wwPDB-VP
Wilson B-factor (Å^2)	81.4	Xtriage
Anisotropy	0.711	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.32 , 195.6	EDS
L-test for twinning[2]	$< \|L\| > = 0.44$, $< L^2 > = 0.26$	Xtriage
Estimated twinning fraction	0.066 for -h,-k,l 0.340 for h,-h-k,-l 0.069 for -k,-h,-l	Xtriage
F_o,F_c correlation	0.90	EDS
Total number of atoms	12328	wwPDB-VP
Average B, all atoms (Å^2)	139.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.47% 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

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	E	0.25	0/1666	0.34	1/2272 (0.0%)
1	I	0.07	0/1666	0.21	0/2272
1	M	0.07	0/1666	0.21	0/2272
1	P	0.20	0/1666	0.40	0/2272
2	F	0.20	0/1487	0.31	0/2031
2	J	0.18	0/1508	0.29	0/2060
2	N	0.17	0/1487	0.28	0/2031
2	R	0.21	0/1508	0.34	0/2060
All	All	0.18	0/12654	0.30	1/17270 (0.0%)

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
1	E	0	1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	E	39	GLN	CB-CA-C	-5.09	102.63	110.78

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	E	38	ARG	Sidechain

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	E	1621	0	1558	21	0
1	I	1621	0	1558	10	0
1	M	1621	0	1558	11	0
1	P	1621	0	1558	28	0
2	F	1451	0	1406	12	0
2	J	1471	0	1421	13	0
2	N	1451	0	1406	9	0
2	R	1471	0	1421	9	0
All	All	12328	0	11886	104	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 4.

All (104) 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:P:113:ASN:CG	1:P:114:GLN:HE21	1.65	1.03
1:P:113:ASN:OD1	1:P:114:GLN:NE2	1.92	1.00
1:P:113:ASN:ND2	1:P:114:GLN:HE21	1.70	0.90
1:P:30:THR:HA	1:P:52(A):PRO:HB2	1.72	0.71
1:I:30:THR:HA	1:I:52(A):PRO:HB2	1.73	0.71
1:P:14:PRO:HD3	1:P:112:PHE:O	1.92	0.69
1:P:14:PRO:HG2	1:P:113:ASN:CG	2.18	0.68
2:R:80:ALA:HA	2:R:106:VAL:HG21	1.76	0.68
2:J:83:GLU:HG3	2:J:105:THR:HA	1.81	0.62
1:P:93:ALA:HB1	1:P:100(D):PHE:HB3	1.81	0.62
1:I:93:ALA:HB1	1:I:100(D):PHE:HB3	1.81	0.62
2:F:18:SER:HB3	2:F:76:SER:HA	1.83	0.59
2:J:105:THR:HB	2:J:142:PRO:HB3	1.84	0.59
2:F:105:THR:O	2:F:106:VAL:C	2.46	0.58
2:J:107:LEU:HD11	2:J:111:LYS:HE3	1.85	0.58
1:M:138:GLY:HA3	1:M:180:VAL:HG12	1.85	0.58
1:P:113:ASN:CG	1:P:114:GLN:H	2.10	0.58
1:E:30:THR:HA	1:E:52(A):PRO:HB2	1.86	0.58
1:I:87:THR:HG23	1:I:110:THR:HA	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:P:113:ASN:HD21	1:P:114:GLN:HE21	1.49	0.57
2:R:146:THR:HB	2:R:197:THR:HB	1.87	0.57
1:P:110:THR:O	1:P:111:VAL:C	2.48	0.57
2:F:83:GLU:HG3	2:F:105:THR:HA	1.88	0.56
1:I:115:ILE:HG12	1:I:146:PRO:HD3	1.88	0.56
1:P:113:ASN:CG	1:P:114:GLN:NE2	2.44	0.56
2:J:146:THR:HB	2:J:197:THR:HB	1.86	0.56
1:P:138:GLY:HA3	1:P:180:VAL:HG12	1.86	0.55
1:E:138:GLY:HA3	1:E:180:VAL:HG12	1.89	0.55
2:R:37:GLN:HB2	2:R:47:ILE:HD13	1.89	0.55
1:E:168:VAL:HB	2:F:163:THR:HG22	1.90	0.54
1:E:38:ARG:NH2	1:E:86:ASP:OD1	2.41	0.52
2:F:103:LYS:HE3	2:F:105:THR:HG22	1.91	0.51
1:P:162:VAL:HG22	1:P:181:VAL:HG12	1.91	0.51
1:M:30:THR:HA	1:M:52(A):PRO:HB2	1.91	0.51
2:F:106:VAL:O	2:F:107:LEU:C	2.53	0.51
1:P:83:ARG:O	1:P:111:VAL:HG11	2.10	0.51
1:P:83:ARG:C	1:P:111:VAL:HG11	2.35	0.51
1:E:162:VAL:HG22	1:E:181:VAL:HG12	1.91	0.50
1:P:115:ILE:HG23	1:P:145:PHE:O	2.13	0.49
2:J:83:GLU:OE2	2:J:105:THR:HG23	2.13	0.49
1:M:115:ILE:HD12	1:M:146:PRO:HD3	1.94	0.49
1:M:162:VAL:HG22	1:M:181:VAL:HG12	1.94	0.48
1:M:96:PRO:HG2	1:M:100:ASP:HB3	1.94	0.48
1:I:47:TRP:HZ2	1:I:50:TRP:HD1	1.61	0.48
2:J:133:LEU:HD12	2:J:179:LEU:HD23	1.95	0.48
2:J:14:SER:HA	2:J:107:LEU:HB2	1.95	0.48
1:P:13:LYS:HA	1:P:112:PHE:O	2.14	0.48
1:E:93:ALA:HB1	1:E:100(D):PHE:HB3	1.95	0.48
2:N:105:THR:O	2:N:106:VAL:C	2.57	0.48
1:P:95:ALA:HB1	1:P:100(A):ARG:HA	1.98	0.46
1:P:47:TRP:HZ2	1:P:50:TRP:HD1	1.62	0.46
1:E:199:HIS:CD2	1:E:201:PRO:HD2	2.50	0.46
1:I:162:VAL:HG22	1:I:181:VAL:HG12	1.97	0.46
1:M:124:ALA:HB1	1:M:212:PRO:HA	1.97	0.45
1:M:168:VAL:HB	2:N:163:THR:HG22	1.97	0.45
1:I:37:VAL:HG22	1:I:47:TRP:HA	1.98	0.45
1:M:115:ILE:HA	1:M:145:PHE:HD2	1.82	0.45
1:P:115:ILE:HG22	1:P:116:LYS:N	2.32	0.45
2:R:133:LEU:HD12	2:R:179:LEU:HD23	1.99	0.45
1:E:39:GLN:HB2	1:E:45:LEU:HD23	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:38:ARG:O	1:E:38:ARG:HG3	2.16	0.44
1:I:40:ALA:HB3	1:I:43:GLN:HB2	1.98	0.44
1:M:93:ALA:HB1	1:M:100(D):PHE:HB3	1.98	0.44
2:F:105:THR:OG1	2:F:142:PRO:HB3	2.17	0.44
2:N:133:LEU:HD12	2:N:179:LEU:HD23	1.98	0.44
1:M:100(C):ASP:OD2	2:N:49:TYR:HB2	2.18	0.44
1:P:113:ASN:CG	1:P:114:GLN:N	2.75	0.44
2:R:106:VAL:O	2:R:107:LEU:C	2.61	0.44
1:M:82(C):LEU:HB3	1:M:111:VAL:HG21	2.00	0.44
2:J:145:VAL:HB	2:J:196:VAL:HG23	2.00	0.43
2:N:83:GLU:HG3	2:N:105:THR:HA	2.01	0.43
2:R:11:ALA:O	2:R:104:VAL:HA	2.18	0.43
1:E:18:VAL:HG11	1:E:109:VAL:HG11	2.00	0.43
1:I:51:ILE:HB	1:I:57:THR:HG22	2.00	0.43
1:E:168:VAL:HG21	2:F:161:GLU:HB3	2.01	0.43
2:J:79:GLN:O	2:J:106:VAL:HG21	2.18	0.42
2:J:137:ILE:HD13	2:J:196:VAL:HG21	2.01	0.42
2:R:145:VAL:HB	2:R:196:VAL:HG23	2.01	0.42
2:R:54:ARG:HD3	2:R:62:PHE:O	2.19	0.42
1:P:11:VAL:H	1:E:114:GLN:HE22	1.67	0.42
1:P:164:THR:HG23	1:P:177:LEU:HD21	2.01	0.42
1:E:164:THR:HG23	1:E:177:LEU:HD21	2.01	0.42
2:N:18:SER:HB3	2:N:76:SER:HA	2.02	0.42
2:N:107:LEU:HD12	2:N:107:LEU:HA	1.79	0.41
1:P:89:VAL:HG22	1:P:108:LEU:HD13	2.02	0.41
1:P:167:ALA:HB1	1:P:175:TYR:HB3	2.01	0.41
1:E:118:PRO:HB3	1:E:144:TYR:HB3	2.02	0.41
2:J:105:THR:O	2:J:106:VAL:C	2.62	0.41
1:E:124:ALA:HB1	1:E:212:PRO:HA	2.02	0.41
1:P:9:ALA:HB3	1:E:115:ILE:HB	2.02	0.41
1:E:96:PRO:HG2	1:E:100:ASP:HB3	2.03	0.41
1:E:100(C):ASP:OD2	2:F:49:TYR:HB2	2.20	0.41
2:N:105:THR:OG1	2:N:142:PRO:HB3	2.20	0.41
1:P:112:PHE:CE1	1:P:145:PHE:HE2	2.38	0.41
1:E:115:ILE:HD13	1:E:146:PRO:HD3	2.02	0.41
1:I:164:THR:HG23	1:I:177:LEU:HD21	2.02	0.41
1:E:165:PHE:CZ	2:F:136:LEU:HB3	2.56	0.40
2:F:110:PRO:HG3	2:J:74:THR:HB	2.02	0.40
2:J:151:ALA:HB1	2:J:189:HIS:CD2	2.56	0.40
1:P:51:ILE:HB	1:P:57:THR:HG22	2.03	0.40
1:E:125:PRO:HB2	1:E:135:ALA:HB1	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:33:VAL:HG21	2:F:71:ALA:HB2	2.03	0.40
2:R:151:ALA:HB1	2:R:189:HIS:CD2	2.56	0.40
2:N:107:LEU:HD12	2:N:141:TYR:HE2	1.87	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	E	204/223 (92%)	199 (98%)	5 (2%)	0	100	100
1	I	204/223 (92%)	199 (98%)	5 (2%)	0	100	100
1	M	204/223 (92%)	200 (98%)	4 (2%)	0	100	100
1	P	204/223 (92%)	194 (95%)	9 (4%)	1 (0%)	25	59
2	F	192/210 (91%)	185 (96%)	5 (3%)	2 (1%)	13	46
2	J	194/210 (92%)	182 (94%)	10 (5%)	2 (1%)	13	46
2	N	192/210 (91%)	183 (95%)	8 (4%)	1 (0%)	25	59
2	R	194/210 (92%)	184 (95%)	9 (5%)	1 (0%)	25	59
All	All	1588/1732 (92%)	1526 (96%)	55 (4%)	7 (0%)	30	64

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	F	106	VAL
1	P	111	VAL
2	F	107	LEU
2	N	106	VAL
2	J	107	LEU
2	J	106	VAL
2	R	106	VAL

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	E	177/189 (94%)	173 (98%)	4 (2%)	45 69
1	I	177/189 (94%)	177 (100%)	0	100 100
1	M	177/189 (94%)	177 (100%)	0	100 100
1	P	177/189 (94%)	175 (99%)	2 (1%)	70 83
2	F	164/177 (93%)	162 (99%)	2 (1%)	67 82
2	J	166/177 (94%)	165 (99%)	1 (1%)	84 91
2	N	164/177 (93%)	163 (99%)	1 (1%)	84 91
2	R	166/177 (94%)	163 (98%)	3 (2%)	54 74
All	All	1368/1464 (93%)	1355 (99%)	13 (1%)	73 84

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	P	115	ILE
1	P	116	LYS
2	R	106	VAL
2	R	107	LEU
2	R	111	LYS
1	E	37	VAL
1	E	38	ARG
1	E	115	ILE
1	E	116	LYS
2	F	107	LEU
2	F	109	GLN
2	J	109	GLN
2	N	107	LEU

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

Mol	Chain	Res	Type
1	P	113	ASN

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Mol	Chain	Res	Type
1	P	114	GLN
2	R	31	ASN
2	R	109	GLN
2	R	198	HIS
1	E	114	GLN
1	I	58	ASN
2	J	109	GLN
1	M	58	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 oligosaccharides 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, 95^{th} 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	E	210/223 (94%)	-0.83	0 100 100	30, 138, 178, 198	0
1	I	210/223 (94%)	-0.88	0 100 100	48, 134, 209, 238	0
1	M	210/223 (94%)	-0.83	0 100 100	73, 142, 180, 200	0
1	P	210/223 (94%)	-0.79	2 (0%) 79 61	44, 134, 207, 234	0
2	F	196/210 (93%)	-0.72	1 (0%) 87 75	65, 135, 204, 242	0
2	J	198/210 (94%)	-0.90	0 100 100	67, 127, 203, 229	0
2	N	196/210 (93%)	-0.63	0 100 100	87, 145, 201, 233	0
2	R	198/210 (94%)	-0.89	0 100 100	48, 123, 197, 220	0
All	All	1628/1732 (93%)	-0.81	3 (0%) 92 86	30, 135, 199, 242	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	P	157	ALA	3.1
1	P	1	GLN	2.6
2	F	131	ALA	2.4

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 monosaccharides 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.