Full wwPDB X-ray Structure Validation Report (i)

Jun 19, 2024 – 05:00 AM EDT

PDB ID	:	4G6F
Title	:	Crystal Structure of 10E8 Fab in Complex with an HIV-1 gp41 Peptide
Authors	:	Ofek, G.; Huang, J.; Connors, M.; Kwong, P.D.
Deposited on	:	2012-07-19
Resolution	:	2.10 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:

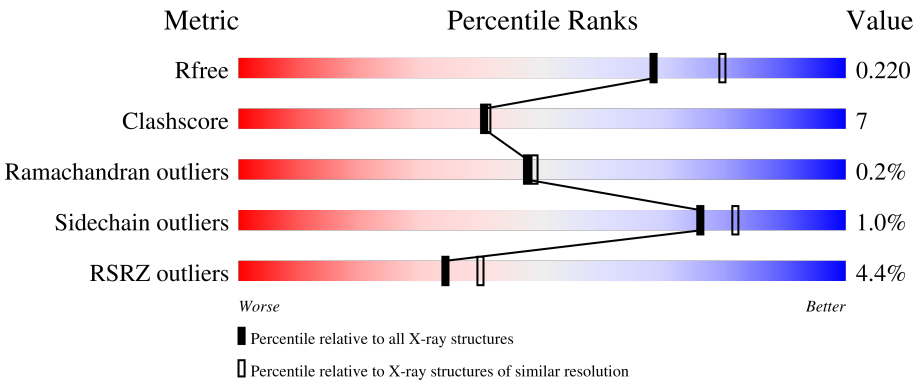
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.20.1
EDS	:	2.37.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation 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 2.10 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	5197 (2.10-2.10)
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (2.10-2.10)

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	B	236	% 84% 14% •
1	H	236	6% 83% 14% •
2	D	215	5% 86% 11% • •
2	L	215	% 86% 13% •
3	F	34	24% 71% 6% • 21%

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Mol	Chain	Length	Quality of chain
3	P	34	18% 85% 15%

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 7728 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 10E8 Heavy Chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	H	232	1763	1122	294	340	7	0	0	0
1	B	235	1783	1132	297	346	8	0	0	0

- Molecule 2 is a protein called 10E8 Light Chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	L	212	1590	988	276	322	4	0	0	0
2	D	211	1581	983	275	319	4	0	0	0

- Molecule 3 is a protein called gp41 MPER Peptide.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
3	P	29	268	180	42	46	0	0	0
3	F	27	259	175	44	40	0	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	H	109	109	109	0	0
4	B	161	161	161	0	0
4	L	104	104	104	0	0
4	D	98	98	98	0	0

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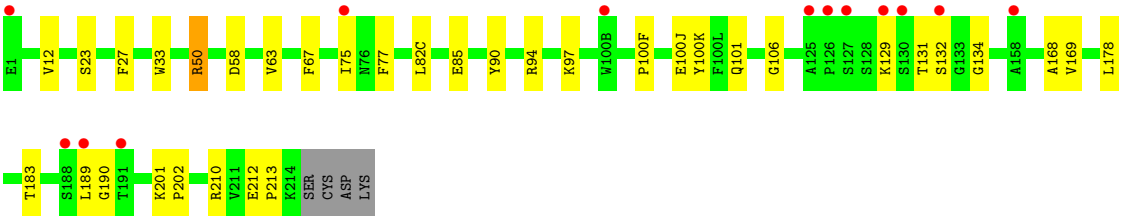
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	P	7	Total O 7 7	0	0
4	F	5	Total O 5 5	0	0

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.

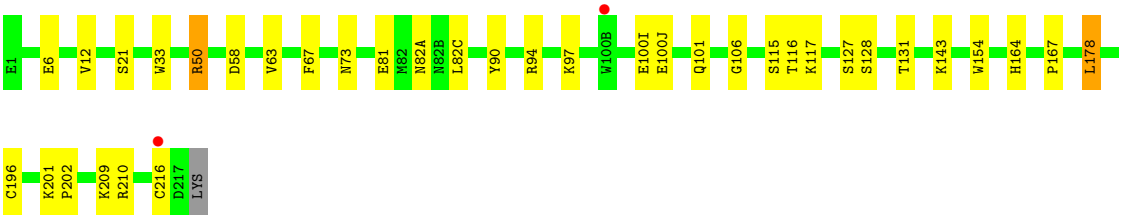
• Molecule 1: 10E8 Heavy Chain

Chain H:



• Molecule 1: 10E8 Heavy Chain

Chain B:

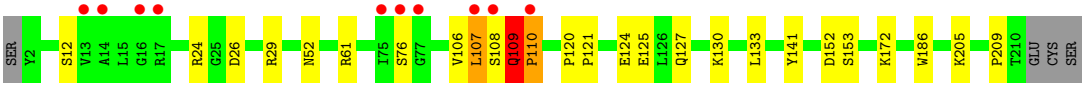


• Molecule 2: 10E8 Light Chain

Chain L:

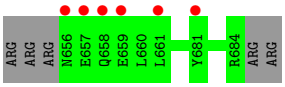
• Molecule 2: 10E8 Light Chain

Chain D:



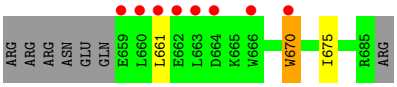
• Molecule 3: gp41 MPER Peptide

Chain P:



• Molecule 3: gp41 MPER Peptide

Chain F:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	68.02Å 71.43Å 129.24Å 90.00° 104.50° 90.00°	Depositor
Resolution (Å)	42.59 – 2.10 42.59 – 2.10	Depositor EDS
% Data completeness (in resolution range)	88.4 (42.59-2.10) 84.4 (42.59-2.10)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.42 (at 2.10Å)	Xtriage
Refinement program	PHENIX 1.8_1069	Depositor
R, R_{free}	0.180 , 0.218 0.186 , 0.220	Depositor DCC
R_{free} test set	2012 reflections (3.25%)	wwPDB-VP
Wilson B-factor (Å^2)	36.3	Xtriage
Anisotropy	0.569	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.35 , 57.5	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.021 for h,-k,-h-l	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	7728	wwPDB-VP
Average B, all atoms (Å^2)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.87% 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	B	0.43	0/1834	0.58	0/2499
1	H	0.36	0/1814	0.52	0/2472
2	D	0.40	1/1616 (0.1%)	0.55	1/2198 (0.0%)
2	L	0.37	0/1625	0.52	0/2210
3	F	0.31	0/270	0.41	0/369
3	P	0.28	0/279	0.36	0/383
All	All	0.38	1/7438 (0.0%)	0.53	1/10131 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	110	PRO	N-CD	5.36	1.55	1.47

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	D	109	GLN	C-N-CD	5.57	140.09	128.40

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	B	1783	0	1714	28	0
1	H	1763	0	1700	29	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	D	1581	0	1542	26	0
2	L	1590	0	1548	23	0
3	F	259	0	245	5	0
3	P	268	0	241	0	0
4	B	161	0	0	9	0
4	D	98	0	0	9	0
4	F	5	0	0	0	0
4	H	109	0	0	5	0
4	L	104	0	0	2	0
4	P	7	0	0	0	0
All	All	7728	0	6990	102	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 (102) 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)
4:H:387:HOH:O	2:L:166:SER:HB3	1.64	0.94
1:B:178:LEU:HD12	1:B:178:LEU:C	1.95	0.86
1:H:50:ARG:NH2	1:H:58:ASP:OD2	2.11	0.82
1:B:209:LYS:NZ	2:D:124:GLU:OE2	2.11	0.80
1:H:210:ARG:HD3	1:H:212:GLU:OE2	1.87	0.74
2:D:24:ARG:NH2	4:D:330:HOH:O	2.25	0.69
1:B:117:LYS:NZ	4:B:355:HOH:O	2.26	0.68
2:D:121:PRO:HD3	2:D:133:LEU:CD2	2.23	0.68
2:L:210:THR:O	2:L:211:GLU:HB2	1.93	0.68
2:D:172:LYS:NZ	4:D:382:HOH:O	2.31	0.64
1:B:167:PRO:HG2	4:D:310:HOH:O	1.97	0.64
1:H:189:LEU:HD12	1:H:190:GLY:N	2.13	0.64
1:B:94:ARG:NH2	4:B:322:HOH:O	2.30	0.64
3:F:661:LEU:HD12	3:F:661:LEU:H	1.63	0.63
1:H:169:VAL:HG21	2:L:161:GLU:HB3	1.80	0.63
2:D:26:ASP:O	2:D:29:ARG:HG2	1.99	0.63
1:B:33:TRP:CE2	1:B:97:LYS:HD3	2.34	0.62
2:D:127:GLN:NE2	4:D:372:HOH:O	2.34	0.61
1:H:169:VAL:CG2	2:L:161:GLU:HB3	2.30	0.61
1:B:178:LEU:C	1:B:178:LEU:CD1	2.67	0.60
2:D:121:PRO:HD3	2:D:133:LEU:HD21	1.83	0.60
1:H:50:ARG:HD3	1:H:100(J):GLU:OE1	2.02	0.59
2:L:182:THR:CG2	2:L:185:GLN:H	2.16	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:178:LEU:HD12	1:B:178:LEU:O	2.03	0.59
1:H:129:LYS:NZ	2:L:208:ALA:H	2.02	0.58
1:H:23:SER:HB3	1:H:77:PHE:CZ	2.39	0.58
2:D:205:LYS:NZ	4:D:388:HOH:O	2.35	0.57
2:L:172:LYS:NZ	4:L:397:HOH:O	2.36	0.57
1:H:33:TRP:CE2	1:H:97:LYS:HD3	2.40	0.57
1:B:128:SER:N	4:B:428:HOH:O	2.37	0.57
1:B:82(A):ASN:ND2	4:B:386:HOH:O	2.37	0.56
2:D:121:PRO:HD3	2:D:133:LEU:HD23	1.88	0.56
2:L:38:LYS:HD3	2:L:44:PRO:HG3	1.88	0.55
2:D:109:GLN:OE1	2:D:110:PRO:HD2	2.06	0.55
1:H:183:THR:HG21	4:H:368:HOH:O	2.06	0.55
2:L:51:LYS:NZ	4:L:377:HOH:O	2.37	0.55
2:L:186:TRP:CZ2	2:L:209:PRO:HA	2.42	0.54
1:H:131:THR:HG21	1:B:115:SER:HB3	1.90	0.54
1:B:216:CYS:SG	1:B:216:CYS:O	2.66	0.53
2:L:120:PRO:HA	2:L:133:LEU:HD23	1.91	0.53
2:D:106:VAL:HG12	2:D:106:VAL:O	2.09	0.53
2:D:29:ARG:NH2	4:D:363:HOH:O	2.39	0.52
1:B:50:ARG:NH2	1:B:58:ASP:OD2	2.34	0.51
2:L:189:HIS:O	2:L:209:PRO:HG2	2.09	0.51
1:B:143:LYS:HG3	4:B:426:HOH:O	2.11	0.51
1:H:132:SER:HB3	4:H:388:HOH:O	2.11	0.50
1:H:27:PHE:CD2	1:H:94:ARG:HD2	2.46	0.50
2:D:152:ASP:O	2:D:153:SER:HB2	2.11	0.50
1:H:63:VAL:HG22	1:H:67:PHE:CE2	2.48	0.49
2:D:61:ARG:HB2	2:D:76:SER:O	2.12	0.49
1:H:168:ALA:HA	1:H:178:LEU:HB3	1.95	0.49
1:B:50:ARG:HD3	1:B:100(J):GLU:OE1	2.13	0.49
1:H:129:LYS:HZ3	2:L:208:ALA:H	1.60	0.48
1:H:75:ILE:O	1:H:75:ILE:HG13	2.13	0.48
2:D:107:LEU:O	2:D:108:SER:OG	2.27	0.48
1:H:131:THR:HG23	1:B:116:THR:O	2.13	0.48
2:D:172:LYS:CE	4:D:382:HOH:O	2.62	0.47
2:L:182:THR:HG22	2:L:185:GLN:H	1.80	0.47
1:B:73:ASN:O	4:B:301:HOH:O	2.21	0.47
2:D:108:SER:O	2:D:141:TYR:CE1	2.67	0.47
3:F:661:LEU:HD12	3:F:661:LEU:N	2.29	0.46
1:B:63:VAL:HG13	1:B:67:PHE:HB2	1.96	0.46
3:F:661:LEU:H	3:F:661:LEU:CD1	2.27	0.46
2:D:29:ARG:NE	4:D:363:HOH:O	2.23	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:110:PRO:HD2	2:D:127:GLN:OE1	2.16	0.46
1:B:90:TYR:O	1:B:106:GLY:HA2	2.17	0.45
1:H:100(K):TYR:CE2	1:H:101:GLN:HG2	2.52	0.45
2:D:172:LYS:HE2	4:D:382:HOH:O	2.17	0.45
2:D:120:PRO:HA	2:D:133:LEU:CD2	2.48	0.44
1:H:85:GLU:H	1:H:85:GLU:CD	2.21	0.44
1:B:81:GLU:HG3	4:B:321:HOH:O	2.18	0.44
2:D:186:TRP:CZ2	2:D:209:PRO:HA	2.53	0.44
1:H:90:TYR:O	1:H:106:GLY:HA2	2.18	0.44
1:H:33:TRP:CE3	1:H:50:ARG:HD2	2.53	0.43
2:L:83:ASP:OD1	2:L:103:LYS:NZ	2.50	0.43
2:L:181:LEU:HD22	2:L:185:GLN:HB3	1.99	0.43
2:L:116:VAL:O	2:L:205:LYS:HE3	2.19	0.43
1:B:201:LYS:N	1:B:202:PRO:CD	2.81	0.43
3:F:670:TRP:CD2	3:F:675:ILE:HD11	2.53	0.43
3:F:670:TRP:CG	3:F:675:ILE:HD11	2.53	0.43
2:D:108:SER:O	2:D:109:GLN:HG2	2.19	0.43
1:B:6:GLU:HA	1:B:21:SER:O	2.20	0.42
2:L:26:ASP:O	2:L:29:ARG:HG3	2.20	0.42
1:H:12:VAL:HG11	1:H:82(C):LEU:HD13	2.02	0.42
1:B:127:SER:O	1:B:131:THR:HG23	2.19	0.42
2:L:182:THR:HG23	2:L:185:GLN:H	1.84	0.42
1:H:63:VAL:HG13	1:H:67:PHE:HB2	2.02	0.42
1:H:134:GLY:C	4:H:398:HOH:O	2.57	0.42
1:B:12:VAL:HG11	1:B:82(C):LEU:HD13	2.02	0.42
1:B:101:GLN:H	1:B:101:GLN:HG2	1.63	0.42
1:B:210:ARG:HD3	4:B:347:HOH:O	2.19	0.42
1:H:201:LYS:N	1:H:202:PRO:CD	2.84	0.41
1:H:100(F):PRO:O	2:L:95:GLY:HA2	2.19	0.41
2:L:36:TYR:HA	2:L:45:ILE:O	2.20	0.41
1:B:63:VAL:HG22	1:B:67:PHE:CE2	2.55	0.41
1:B:154:TRP:CH2	1:B:196:CYS:HB3	2.56	0.41
4:B:366:HOH:O	2:D:205:LYS:HE3	2.21	0.41
2:L:208:ALA:HA	2:L:209:PRO:HD3	1.91	0.41
1:H:189:LEU:HD22	1:H:213:PRO:HG2	2.02	0.41
1:H:132:SER:HB2	4:H:386:HOH:O	2.20	0.41
2:D:125:GLU:HG2	2:D:130:LYS:O	2.21	0.40
2:D:12:SER:HB2	2:D:107:LEU:HD22	2.03	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	B	233/236 (99%)	230 (99%)	3 (1%)	0	100	100
1	H	230/236 (98%)	221 (96%)	9 (4%)	0	100	100
2	D	209/215 (97%)	205 (98%)	2 (1%)	2 (1%)	15	11
2	L	210/215 (98%)	205 (98%)	5 (2%)	0	100	100
3	F	25/34 (74%)	25 (100%)	0	0	100	100
3	P	27/34 (79%)	27 (100%)	0	0	100	100
All	All	934/970 (96%)	913 (98%)	19 (2%)	2 (0%)	47	49

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	D	52	ASN
2	D	109	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	B	197/198 (100%)	193 (98%)	4 (2%)	55	60
1	H	194/198 (98%)	193 (100%)	1 (0%)	88	92
2	D	175/180 (97%)	173 (99%)	2 (1%)	73	79
2	L	176/180 (98%)	176 (100%)	0	100	100
3	F	26/33 (79%)	25 (96%)	1 (4%)	33	34

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
3	P	27/33 (82%)	27 (100%)	0	100	100
All	All	795/822 (97%)	787 (99%)	8 (1%)	76	82

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	H	50	ARG
1	B	50	ARG
1	B	100(I)	GLU
1	B	164	HIS
1	B	178	LEU
2	D	107	LEU
2	D	109	GLN
3	F	670	TRP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
2	D	79	GLN

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	B	235/236 (99%)	-0.26	2 (0%)	84	86	26, 37, 60, 103	0
1	H	232/236 (98%)	0.25	13 (5%)	24	29	33, 51, 95, 127	0
2	D	211/215 (98%)	0.12	10 (4%)	31	37	26, 49, 79, 113	0
2	L	212/215 (98%)	-0.10	3 (1%)	75	78	30, 47, 78, 106	0
3	F	27/34 (79%)	1.65	8 (29%)	0	0	41, 67, 121, 129	0
3	P	29/34 (85%)	0.98	6 (20%)	1	1	51, 67, 131, 136	0
All	All	946/970 (97%)	0.08	42 (4%)	34	40	26, 47, 86, 136	0

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	108	SER	9.2
3	F	661	LEU	7.5
3	F	660	LEU	6.0
1	B	216	CYS	5.8
1	H	129	LYS	5.8
3	P	657	GLU	5.6
3	P	656	ASN	5.4
3	F	662	GLU	5.3
3	F	659	GLU	5.3
3	P	658	GLN	5.1
1	H	188	SER	4.7
3	F	663	LEU	4.5
3	F	666	TRP	4.5
2	D	107	LEU	4.2
3	P	659	GLU	4.0
1	H	126	PRO	3.8
1	H	130	SER	3.4
2	D	76	SER	3.2
1	H	127	SER	3.0

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Mol	Chain	Res	Type	RSRZ
1	H	132	SER	3.0
2	D	17	ARG	3.0
1	H	1	GLU	3.0
2	L	211	GLU	3.0
2	D	16	GLY	2.9
1	H	158	ALA	2.9
2	D	13	VAL	2.7
1	H	191	THR	2.7
2	D	75	ILE	2.6
2	D	77	GLY	2.6
2	D	110	PRO	2.6
3	F	664	ASP	2.5
2	L	76	SER	2.5
1	H	100(B)	TRP	2.5
2	D	14	ALA	2.5
1	B	100(B)	TRP	2.4
1	H	75	ILE	2.3
3	P	661	LEU	2.2
1	H	125	ALA	2.2
3	P	681	TYR	2.0
3	F	670	TRP	2.0
1	H	189	LEU	2.0
2	L	157	LYS	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.