Full wwPDB X-ray Structure Validation Report ⓘ

Oct 9, 2023 – 06:57 AM EDT

PDB ID	:	8D6Z
Title	:	Crystal structure of SARS-CoV-2 spike stem fusion peptide in complex with neutralizing antibody COV91-27
Authors	:	Lee, C.C.D.; Lin, T.H.; Yuan, M.; Wilson, I.A.
Deposited on	:	2022-06-06
Resolution	:	2.30 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:

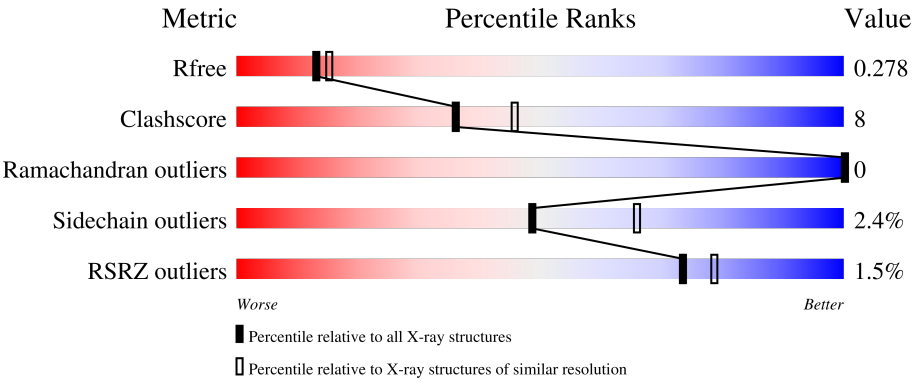
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.35.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35.1

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 2.30 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	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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	A	224	2% 83% 15%
1	C	224	4% 75% 21%
1	E	224	3% 82% 17%
1	G	224	3% 83% 16%
2	B	215	3% 79% 20%

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Mol	Chain	Length	Quality of chain
2	D	215	2% 84% 15%
2	F	215	3% 79% 21%
2	I	215	% 80% 20%
3	H	15	47% 27% 27%
3	J	15	67% 13% 20%
3	K	15	53% 20% 27%
3	L	15	7% 73% 7% 20%

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 13774 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 Neutralizing antibody COV91-27 heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	222	Total	C	N	O	S	0	0	0
			1662	1049	284	321	8			
1	C	221	Total	C	N	O	S	0	0	0
			1656	1046	283	320	7			
1	E	222	Total	C	N	O	S	0	0	0
			1662	1049	284	321	8			
1	G	222	Total	C	N	O	S	0	0	0
			1662	1049	284	321	8			

- Molecule 2 is a protein called Neutralizing antibody COV91-27 light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	214	Total	C	N	O	S	0	0	0
			1637	1020	285	328	4			
2	D	214	Total	C	N	O	S	0	0	0
			1637	1020	285	328	4			
2	F	214	Total	C	N	O	S	0	0	0
			1637	1020	285	328	4			
2	I	215	Total	C	N	O	S	0	0	0
			1640	1022	286	327	5			

- Molecule 3 is a protein called Spike protein S2 fusion peptide.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	H	11	Total	C	N	O	0	0	0
			96	63	15	18			
3	J	12	Total	C	N	O	0	0	0
			103	68	16	19			
3	K	11	Total	C	N	O	0	0	0
			96	63	15	18			
3	L	12	Total	C	N	O	0	0	0
			103	68	16	19			

- Molecule 4 is water.

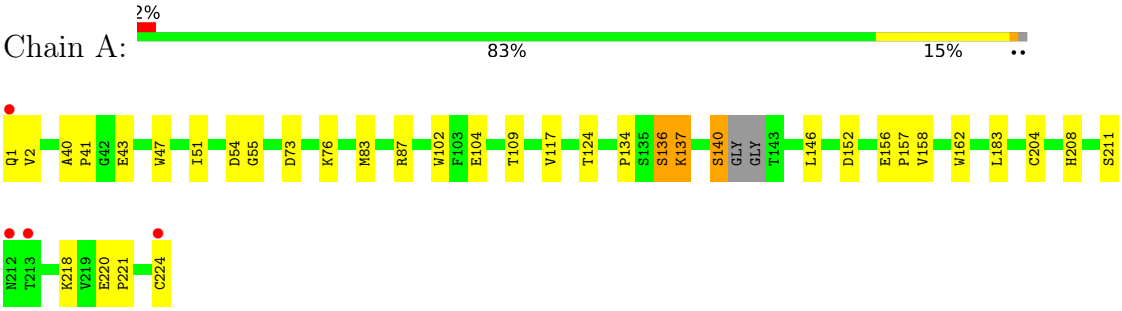
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	22	Total O 22 22	0	0
4	B	23	Total O 23 23	0	0
4	C	27	Total O 27 27	0	0
4	D	22	Total O 22 22	0	0
4	E	20	Total O 20 20	0	0
4	F	20	Total O 20 20	0	0
4	G	23	Total O 23 23	0	0
4	H	1	Total O 1 1	0	0
4	I	20	Total O 20 20	0	0
4	J	2	Total O 2 2	0	0
4	K	1	Total O 1 1	0	0
4	L	2	Total O 2 2	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.

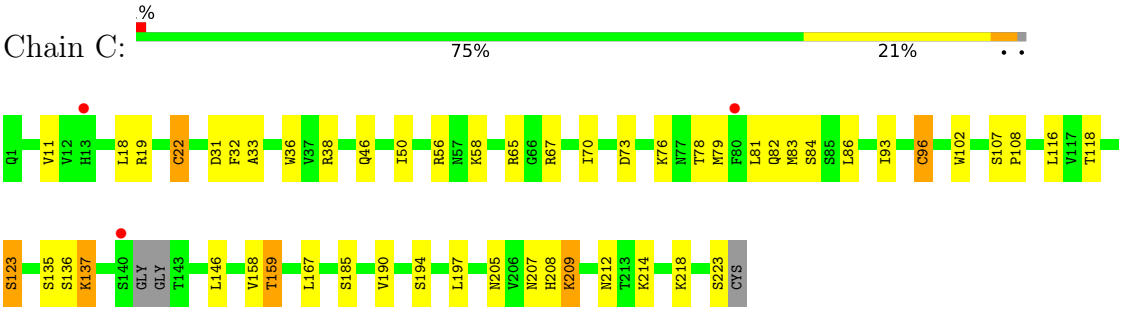
• Molecule 1: Neutralizing antibody COV91-27 heavy chain

Chain A:



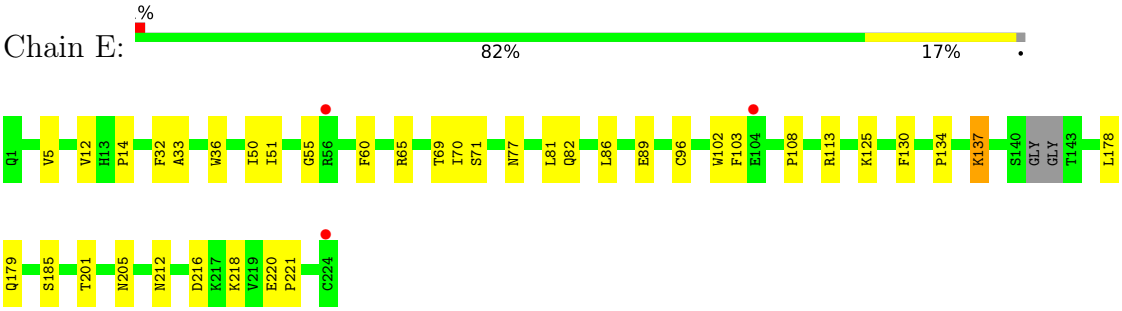
• Molecule 1: Neutralizing antibody COV91-27 heavy chain

Chain C:



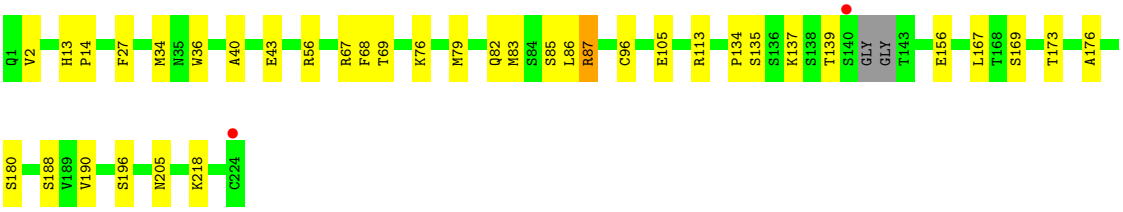
• Molecule 1: Neutralizing antibody COV91-27 heavy chain

Chain E:



• Molecule 1: Neutralizing antibody COV91-27 heavy chain

Chain G:

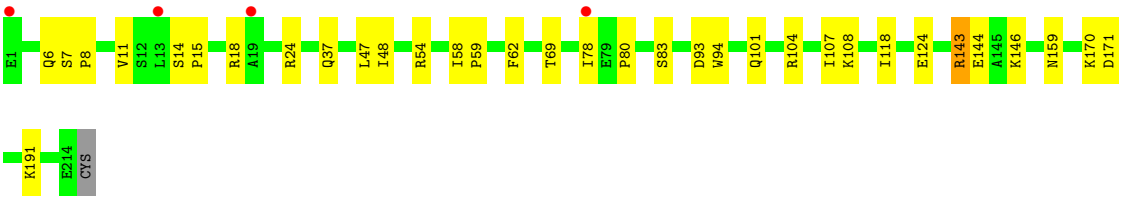


• Molecule 2: Neutralizing antibody COV91-27 light chain

Chain B:

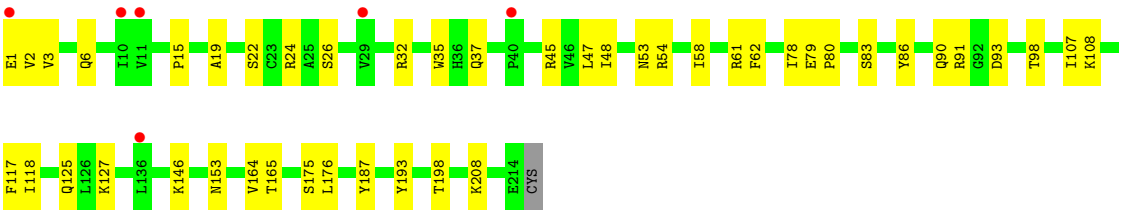
• Molecule 2: Neutralizing antibody COV91-27 light chain

Chain D:



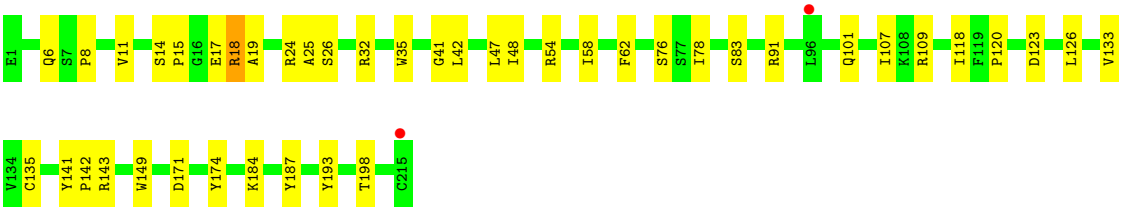
• Molecule 2: Neutralizing antibody COV91-27 light chain

Chain F:



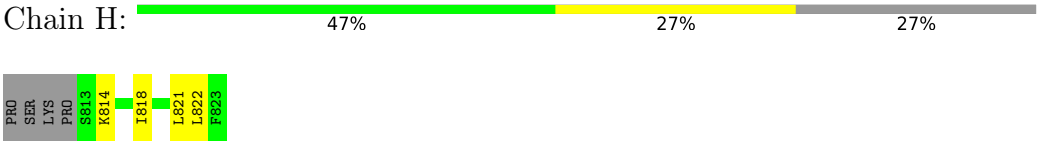
• Molecule 2: Neutralizing antibody COV91-27 light chain

Chain I:



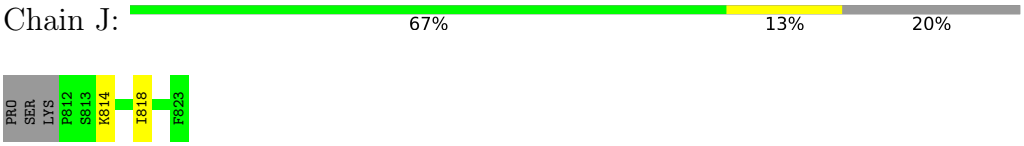
• Molecule 3: Spike protein S2 fusion peptide

Chain H:



• Molecule 3: Spike protein S2 fusion peptide

Chain J:

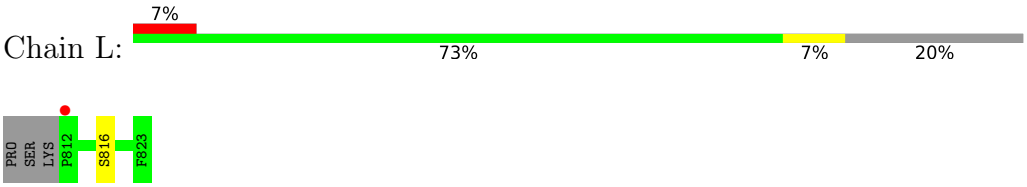


• Molecule 3: Spike protein S2 fusion peptide

Chain K:

• Molecule 3: Spike protein S2 fusion peptide

Chain L:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	196.17Å 60.31Å 186.51Å 90.00° 109.52° 90.00°	Depositor
Resolution (Å)	46.22 – 2.30 46.22 – 2.30	Depositor EDS
% Data completeness (in resolution range)	96.1 (46.22-2.30) 96.1 (46.22-2.30)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.41 (at 2.29Å)	Xtriage
Refinement program	PHENIX 1.19.2_4158	Depositor
R, R_{free}	0.235 , 0.279 0.235 , 0.278	Depositor DCC
R_{free} test set	4453 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å^2)	52.0	Xtriage
Anisotropy	0.334	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.31 , 45.7	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	13774	wwPDB-VP
Average B, all atoms (Å^2)	57.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 78.91 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 6.2579e-07. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

[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	A	0.34	0/1703	0.57	0/2317
1	C	0.33	0/1697	0.58	0/2309
1	E	0.31	0/1703	0.55	0/2317
1	G	0.34	0/1703	0.60	0/2317
2	B	0.31	0/1671	0.58	0/2271
2	D	0.31	0/1671	0.57	0/2271
2	F	0.32	0/1671	0.55	0/2271
2	I	0.31	0/1674	0.56	0/2275
3	H	0.32	0/97	0.50	0/126
3	J	0.35	0/105	0.48	0/137
3	K	0.29	0/97	0.48	0/126
3	L	0.29	0/105	0.45	0/137
All	All	0.32	0/13897	0.57	0/18874

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-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 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	A	1662	0	1624	26	0
1	C	1656	0	1621	39	0
1	E	1662	0	1624	29	0
1	G	1662	0	1624	22	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	1637	0	1603	33	0
2	D	1637	0	1603	26	1
2	F	1637	0	1603	30	1
2	I	1640	0	1606	25	0
3	H	96	0	96	4	0
3	J	103	0	104	1	0
3	K	96	0	96	2	0
3	L	103	0	104	1	0
4	A	22	0	0	2	0
4	B	23	0	0	3	0
4	C	27	0	0	7	0
4	D	22	0	0	4	0
4	E	20	0	0	2	0
4	F	20	0	0	3	0
4	G	23	0	0	2	0
4	H	1	0	0	0	0
4	I	20	0	0	1	0
4	J	2	0	0	0	0
4	K	1	0	0	0	0
4	L	2	0	0	0	0
All	All	13774	0	13308	214	1

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 (214) 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:C:96:CYS:SG	4:C:325:HOH:O	2.02	1.14
1:C:22:CYS:SG	4:C:325:HOH:O	2.03	1.14
2:D:144:GLU:O	4:D:301:HOH:O	1.83	0.96
1:C:185:SER:OG	4:C:301:HOH:O	1.92	0.86
2:I:198:THR:OG1	4:I:301:HOH:O	1.97	0.83
1:G:135:SER:H	1:G:137:LYS:HG2	1.45	0.82
2:F:32:ARG:NH1	4:F:301:HOH:O	1.95	0.81
2:F:47:LEU:HD12	2:F:58:ILE:HD12	1.63	0.80
2:I:14:SER:OG	2:I:17:GLU:OE2	2.01	0.78
2:I:47:LEU:HD12	2:I:58:ILE:HD12	1.66	0.78
2:F:3:VAL:HG12	2:F:26:SER:HB3	1.65	0.77
1:C:65:ARG:O	4:C:302:HOH:O	2.03	0.76
1:A:134:PRO:HG3	1:A:146:LEU:HB3	1.70	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:201:THR:HG23	1:E:218:LYS:HE2	1.70	0.73
2:D:7:SER:HB2	2:D:24:ARG:HH22	1.55	0.71
1:C:223:SER:OG	4:C:303:HOH:O	2.08	0.71
1:A:73:ASP:OD2	1:A:76:LYS:NZ	2.19	0.70
2:F:165:THR:HG22	2:F:175:SER:H	1.56	0.70
2:F:19:ALA:HB2	2:F:78:ILE:HD11	1.75	0.69
1:G:36:TRP:NE1	1:G:79:MET:SD	2.67	0.68
1:E:77:ASN:ND2	4:E:303:HOH:O	2.27	0.68
1:A:156:GLU:OE2	1:A:157:PRO:HA	1.95	0.67
1:C:31:ASP:OD1	4:C:304:HOH:O	2.12	0.67
1:C:33:ALA:HB1	1:C:50:ILE:HD11	1.76	0.67
1:A:40:ALA:HB3	1:A:43:GLU:HB2	1.77	0.67
2:D:104:ARG:NH1	4:D:303:HOH:O	2.17	0.67
1:C:159:THR:HG23	1:C:207:ASN:HB2	1.77	0.67
1:A:51:ILE:HD11	1:A:55:GLY:HA2	1.78	0.66
1:A:218:LYS:HE2	1:A:220:GLU:OE2	1.95	0.65
1:C:93:ILE:HG13	1:C:116:LEU:HD23	1.78	0.65
2:D:159:ASN:OD1	4:D:302:HOH:O	2.13	0.65
1:E:89:GLU:OE1	4:E:301:HOH:O	2.14	0.64
1:C:83:MET:HB3	1:C:86:LEU:HD21	1.80	0.63
2:D:8:PRO:HG2	2:D:11:VAL:HG22	1.79	0.63
2:B:164:VAL:HG22	2:B:176:LEU:HD13	1.80	0.62
2:I:19:ALA:HB2	2:I:78:ILE:HD11	1.81	0.62
1:A:134:PRO:HD2	1:A:221:PRO:HA	1.81	0.62
2:D:47:LEU:HB3	2:D:48:ILE:HD12	1.81	0.61
1:E:51:ILE:HD11	1:E:55:GLY:HA2	1.81	0.61
2:B:43:ALA:HB3	4:B:321:HOH:O	2.01	0.61
1:G:27:PHE:O	4:G:301:HOH:O	2.15	0.60
2:I:126:LEU:HB3	2:I:184:LYS:HD2	1.84	0.60
1:C:218:LYS:HB3	1:E:212:ASN:HB3	1.83	0.60
1:A:40:ALA:HB1	1:A:41:PRO:HD2	1.84	0.59
2:D:80:PRO:HA	2:D:107:ILE:HG13	1.85	0.59
2:F:32:ARG:HD3	4:F:301:HOH:O	2.02	0.58
2:B:49:TYR:OH	4:B:301:HOH:O	1.84	0.58
2:D:143:ARG:NH1	4:D:306:HOH:O	2.36	0.58
2:D:124:GLU:HG3	1:E:125:LYS:HA	1.86	0.58
1:C:214:LYS:HB2	1:E:216:ASP:HB2	1.86	0.58
1:A:140:SER:OG	4:A:301:HOH:O	2.16	0.57
1:E:51:ILE:HB	1:E:70:ILE:HG12	1.85	0.57
1:C:56:ARG:NH2	3:K:820:ASP:OD2	2.38	0.57
1:G:87:ARG:NH1	4:G:302:HOH:O	2.25	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:34:MET:HB3	1:G:79:MET:HE2	1.85	0.57
1:C:194:SER:HA	1:C:197:LEU:HD23	1.87	0.56
1:E:33:ALA:HB1	1:E:50:ILE:HD11	1.86	0.56
2:B:121:PRO:HB2	2:B:126:LEU:HD21	1.88	0.56
1:C:123:SER:HB2	2:F:127:LYS:HE3	1.87	0.56
1:G:156:GLU:OE2	1:G:176:ALA:HB3	2.06	0.56
1:G:56:ARG:NH1	3:L:816:SER:OG	2.40	0.55
1:E:218:LYS:HD3	1:E:220:GLU:HG3	1.88	0.55
1:E:137:LYS:HE2	2:F:118:ILE:HG22	1.89	0.55
1:E:12:VAL:HG11	1:E:86:LEU:HD13	1.89	0.55
1:G:83:MET:HB3	1:G:86:LEU:HD21	1.88	0.54
1:G:34:MET:HB3	1:G:79:MET:CE	2.37	0.54
2:B:150:LYS:HD2	2:B:155:LEU:HG	1.90	0.54
1:G:68:PHE:CZ	1:G:83:MET:HG2	2.44	0.53
2:I:54:ARG:NH1	2:I:58:ILE:O	2.40	0.53
2:I:135:CYS:HB2	2:I:149:TRP:CZ2	2.43	0.53
2:I:143:ARG:HD2	2:I:174:TYR:CE2	2.43	0.53
2:I:24:ARG:NH1	2:I:26:SER:HA	2.25	0.52
1:E:179:GLN:NE2	1:E:185:SER:HB2	2.25	0.52
2:F:80:PRO:HA	2:F:107:ILE:HG13	1.92	0.52
2:B:162:GLU:HB3	2:B:176:LEU:HD11	1.93	0.51
3:H:818:ILE:HD12	3:H:818:ILE:H	1.75	0.51
1:C:212:ASN:HB3	1:E:218:LYS:HB3	1.93	0.51
1:C:36:TRP:CE2	1:C:79:MET:HE1	2.46	0.51
1:C:73:ASP:OD2	1:C:76:LYS:HG3	2.11	0.51
2:B:168:ASP:OD2	2:B:170:LYS:N	2.39	0.51
2:B:176:LEU:HD12	2:B:177:SER:H	1.75	0.51
1:E:60:PHE:HB2	1:E:65:ARG:HG3	1.93	0.51
1:E:32:PHE:CZ	1:E:108:PRO:HD2	2.46	0.50
2:F:53:ASN:ND2	4:F:305:HOH:O	2.43	0.50
1:A:152:ASP:HB3	1:A:183:LEU:HD23	1.94	0.50
2:D:83:SER:HB2	2:D:107:ILE:HG12	1.93	0.50
1:C:36:TRP:NE1	1:C:81:LEU:HB2	2.26	0.50
1:A:137:LYS:O	4:A:302:HOH:O	2.18	0.50
2:I:8:PRO:HG2	2:I:11:VAL:HG22	1.94	0.50
1:C:58:LYS:HE3	1:C:70:ILE:HG22	1.95	0.49
1:E:137:LYS:HB3	2:F:117:PHE:HB3	1.93	0.49
2:B:15:PRO:HD3	2:B:108:LYS:O	2.12	0.49
1:C:36:TRP:NE1	1:C:79:MET:HE1	2.27	0.49
1:G:173:THR:HG23	1:G:188:SER:HB2	1.95	0.49
3:H:818:ILE:HA	3:H:821:LEU:HD12	1.93	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:134:PRO:HG2	1:E:221:PRO:HA	1.94	0.49
2:F:83:SER:HB2	2:F:107:ILE:HG12	1.95	0.48
1:C:22:CYS:O	1:C:78:THR:HA	2.13	0.48
2:F:15:PRO:HD3	2:F:108:LYS:O	2.13	0.48
1:A:136:SER:N	1:A:224:CYS:SG	2.86	0.48
1:A:137:LYS:HG2	2:B:118:ILE:HG22	1.96	0.48
2:B:83:SER:HB2	2:B:107:ILE:HG12	1.96	0.48
2:B:122:SER:O	2:B:126:LEU:HD23	2.14	0.48
2:B:176:LEU:HD12	2:B:177:SER:N	2.29	0.48
2:D:146:LYS:HB2	2:D:146:LYS:HE2	1.64	0.47
2:F:48:ILE:HD13	2:F:54:ARG:HA	1.96	0.47
2:F:90:GLN:NE2	2:F:98:THR:OG1	2.47	0.47
1:G:135:SER:N	1:G:137:LYS:HG2	2.23	0.47
2:B:148:GLN:NE2	2:B:149:TRP:O	2.48	0.47
1:G:40:ALA:HB3	1:G:43:GLU:HB2	1.95	0.47
2:B:187:TYR:O	2:B:193:TYR:OH	2.31	0.47
2:F:2:VAL:HB	2:F:90:GLN:NE2	2.30	0.47
1:G:167:LEU:HD21	1:G:190:VAL:HG21	1.96	0.47
2:I:187:TYR:HA	2:I:193:TYR:OH	2.15	0.47
3:J:814:LYS:HE3	3:J:818:ILE:HD11	1.97	0.47
2:B:126:LEU:HD11	2:B:187:TYR:CD2	2.50	0.47
1:C:107:SER:HB3	4:C:319:HOH:O	2.15	0.47
2:F:35:TRP:HB2	2:F:48:ILE:HB	1.97	0.47
1:G:67:ARG:HD2	1:G:85:SER:HB2	1.96	0.47
2:B:121:PRO:HD3	2:B:133:VAL:HG22	1.97	0.47
1:C:135:SER:O	1:C:136:SER:OG	2.25	0.47
1:A:104:GLU:CD	1:A:104:GLU:H	2.19	0.46
2:D:7:SER:CB	2:D:24:ARG:HH22	2.26	0.46
1:E:178:LEU:HD13	1:E:179:GLN:O	2.16	0.46
1:C:137:LYS:CG	2:D:118:ILE:H	2.28	0.46
1:A:54:ASP:OD1	1:A:54:ASP:N	2.37	0.46
3:H:814:LYS:O	3:H:818:ILE:HD12	2.16	0.46
1:C:137:LYS:HG2	2:D:118:ILE:O	2.16	0.46
1:C:158:VAL:HG12	1:C:208:HIS:HB2	1.97	0.46
2:I:18:ARG:HG3	2:I:76:SER:HA	1.98	0.46
2:I:24:ARG:HD3	2:I:25:ALA:N	2.31	0.46
2:I:41:GLY:O	2:I:42:LEU:HD13	2.16	0.45
2:B:78:ILE:HD12	2:B:105:LEU:HD21	1.98	0.45
2:B:155:LEU:HD12	2:B:155:LEU:H	1.82	0.45
1:C:137:LYS:HA	1:C:137:LYS:HD3	1.71	0.45
1:E:5:VAL:HG13	1:E:113:ARG:HD3	1.97	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:70:ILE:HG13	1:E:71:SER:N	2.31	0.45
2:F:37:GLN:OE1	2:F:45:ARG:NH2	2.47	0.45
1:G:135:SER:O	1:G:135:SER:OG	2.26	0.45
2:I:15:PRO:HA	2:I:78:ILE:HG22	1.98	0.45
2:D:6:GLN:H	2:D:101:GLN:HE22	1.64	0.45
2:I:32:ARG:HB3	2:I:91:ARG:HG3	1.99	0.45
2:B:47:LEU:HD21	2:B:62:PHE:CD2	2.52	0.45
1:A:83:MET:HE1	1:A:117:VAL:HG21	1.99	0.44
2:F:187:TYR:HA	2:F:193:TYR:OH	2.17	0.44
1:A:156:GLU:OE2	1:A:158:VAL:N	2.51	0.44
2:B:186:ASP:OD2	2:F:24:ARG:NH2	2.50	0.44
1:C:209:LYS:HA	1:C:209:LYS:HD2	1.61	0.44
2:D:78:ILE:HD13	2:D:78:ILE:HA	1.85	0.44
2:I:6:GLN:H	2:I:101:GLN:NE2	2.15	0.44
2:D:59:PRO:HG2	2:D:62:PHE:HE2	1.83	0.44
1:A:208:HIS:CE1	1:A:211:SER:HB3	2.53	0.43
2:D:170:LYS:HG3	2:D:171:ASP:N	2.32	0.43
1:E:130:PHE:CE2	2:F:125:GLN:HG3	2.52	0.43
2:I:83:SER:HB2	2:I:107:ILE:HG12	2.00	0.43
2:I:6:GLN:H	2:I:101:GLN:HE22	1.66	0.43
2:B:149:TRP:CE2	2:B:180:LEU:HB2	2.53	0.43
2:F:47:LEU:HD21	2:F:62:PHE:CD1	2.53	0.43
2:B:93:ASP:O	2:B:96:LEU:N	2.39	0.43
2:B:95:PRO:HG3	3:H:822:LEU:O	2.18	0.43
2:D:59:PRO:HG2	2:D:62:PHE:CE2	2.54	0.43
1:G:69:THR:HB	1:G:82:GLN:HB3	2.00	0.43
1:A:87:ARG:HD2	1:E:14:PRO:HG2	2.01	0.43
2:F:1:GLU:HG2	2:F:93:ASP:OD2	2.18	0.43
1:A:124:THR:HG22	1:A:211:SER:HB3	2.01	0.43
1:E:102:TRP:CZ2	1:E:103:PHE:HE1	2.37	0.42
2:I:54:ARG:HD3	2:I:62:PHE:O	2.18	0.42
2:I:109:ARG:NH1	2:I:171:ASP:O	2.51	0.42
1:C:32:PHE:CZ	1:C:108:PRO:HG2	2.54	0.42
2:F:61:ARG:CZ	2:F:79:GLU:HG3	2.49	0.42
3:K:818:ILE:HA	3:K:821:LEU:HD12	2.01	0.42
1:A:137:LYS:CG	2:B:118:ILE:H	2.32	0.42
1:C:11:VAL:HA	1:C:118:THR:O	2.19	0.42
2:D:24:ARG:HA	2:D:69:THR:O	2.19	0.42
1:E:36:TRP:NE1	1:E:81:LEU:HB2	2.34	0.42
2:I:118:ILE:HD11	2:I:133:VAL:HG12	2.01	0.42
1:E:32:PHE:CE2	1:E:108:PRO:HD2	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:102:TRP:HB3	2:B:94:TRP:CG	2.54	0.42
1:A:162:TRP:CH2	1:A:204:CYS:HB3	2.55	0.42
2:B:40:PRO:HB2	2:B:166:GLU:HG3	2.01	0.42
1:G:13:HIS:HD2	1:G:14:PRO:O	2.03	0.42
2:I:141:TYR:CG	2:I:142:PRO:HA	2.55	0.42
1:A:109:THR:HG21	2:B:36:HIS:HE1	1.85	0.42
2:F:47:LEU:HD21	2:F:62:PHE:CE1	2.55	0.42
2:B:141:TYR:CG	2:B:142:PRO:HA	2.55	0.41
1:C:19:ARG:HB2	1:C:82:GLN:NE2	2.34	0.41
2:D:54:ARG:NH1	2:D:58:ILE:O	2.53	0.41
2:F:6:GLN:HE21	2:F:6:GLN:HB3	1.65	0.41
2:F:32:ARG:HB3	2:F:91:ARG:HG3	2.02	0.41
2:F:164:VAL:HG22	2:F:176:LEU:HD12	2.02	0.41
2:D:54:ARG:HD3	2:D:62:PHE:O	2.20	0.41
1:E:137:LYS:HD3	1:E:137:LYS:HA	1.37	0.41
1:A:1:GLN:HG2	1:A:2:VAL:N	2.36	0.41
2:F:146:LYS:HB3	2:F:198:THR:HB	2.00	0.41
2:B:146:LYS:HB3	2:B:146:LYS:HE2	1.89	0.41
1:C:102:TRP:HB3	2:D:94:TRP:CG	2.55	0.41
2:B:118:ILE:HB	2:B:208:LYS:HG3	2.02	0.41
1:G:137:LYS:NZ	2:I:120:PRO:HD3	2.36	0.41
1:C:36:TRP:CG	1:C:81:LEU:HD22	2.56	0.41
1:C:218:LYS:HD3	1:E:212:ASN:OD1	2.21	0.41
1:C:38:ARG:HH21	1:C:46:GLN:HE22	1.67	0.41
1:C:167:LEU:HD21	1:C:190:VAL:HG21	2.03	0.41
2:D:37:GLN:HB2	2:D:47:LEU:HD22	2.03	0.41
1:E:69:THR:OG1	1:E:82:GLN:HB3	2.21	0.41
1:G:2:VAL:HG13	1:G:27:PHE:CD1	2.56	0.41
2:I:35:TRP:HB2	2:I:48:ILE:HD13	2.03	0.41
2:F:37:GLN:HG3	2:F:86:TYR:CE2	2.56	0.41
1:G:134:PRO:HA	1:G:137:LYS:HG3	2.02	0.41
2:B:143:ARG:HD3	4:B:313:HOH:O	2.21	0.40
1:G:105:GLU:OE1	1:G:105:GLU:N	2.46	0.40
1:A:47:TRP:CD1	2:B:97:ILE:HB	2.56	0.40
1:C:67:ARG:HB3	1:C:84:SER:O	2.21	0.40
1:C:137:LYS:HG2	2:D:118:ILE:H	1.86	0.40
2:D:15:PRO:HD3	2:D:108:LYS:O	2.22	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:191:LYS:NZ	2:F:153:ASN:OD1[1_545]	2.13	0.07

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	A	218/224 (97%)	213 (98%)	5 (2%)	0	100	100
1	C	217/224 (97%)	213 (98%)	4 (2%)	0	100	100
1	E	218/224 (97%)	215 (99%)	3 (1%)	0	100	100
1	G	218/224 (97%)	212 (97%)	6 (3%)	0	100	100
2	B	212/215 (99%)	205 (97%)	7 (3%)	0	100	100
2	D	212/215 (99%)	206 (97%)	6 (3%)	0	100	100
2	F	212/215 (99%)	205 (97%)	7 (3%)	0	100	100
2	I	213/215 (99%)	208 (98%)	5 (2%)	0	100	100
3	H	9/15 (60%)	8 (89%)	1 (11%)	0	100	100
3	J	10/15 (67%)	10 (100%)	0	0	100	100
3	K	9/15 (60%)	9 (100%)	0	0	100	100
3	L	10/15 (67%)	10 (100%)	0	0	100	100
All	All	1758/1816 (97%)	1714 (98%)	44 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	A	185/185 (100%)	182 (98%)	3 (2%)	62 78
1	C	184/185 (100%)	175 (95%)	9 (5%)	25 35
1	E	185/185 (100%)	182 (98%)	3 (2%)	62 78
1	G	185/185 (100%)	175 (95%)	10 (5%)	22 30
2	B	184/185 (100%)	181 (98%)	3 (2%)	62 78
2	D	184/185 (100%)	180 (98%)	4 (2%)	52 69
2	F	184/185 (100%)	182 (99%)	2 (1%)	73 86
2	I	184/185 (100%)	182 (99%)	2 (1%)	73 86
3	H	11/15 (73%)	11 (100%)	0	100 100
3	J	12/15 (80%)	12 (100%)	0	100 100
3	K	11/15 (73%)	11 (100%)	0	100 100
3	L	12/15 (80%)	12 (100%)	0	100 100
All	All	1521/1540 (99%)	1485 (98%)	36 (2%)	49 66

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	136	SER
1	A	137	LYS
1	A	140	SER
2	B	93	ASP
2	B	157	SER
2	B	204	SER
1	C	18	LEU
1	C	22	CYS
1	C	96	CYS
1	C	123	SER
1	C	137	LYS
1	C	146	LEU
1	C	159	THR
1	C	205	ASN
1	C	209	LYS
2	D	14	SER
2	D	18	ARG
2	D	93	ASP
2	D	143	ARG
1	E	96	CYS
1	E	137	LYS
1	E	205	ASN

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Mol	Chain	Res	Type
2	F	22	SER
2	F	208	LYS
1	G	76	LYS
1	G	87	ARG
1	G	96	CYS
1	G	113	ARG
1	G	139	THR
1	G	169	SER
1	G	180	SER
1	G	196	SER
1	G	205	ASN
1	G	218	LYS
2	I	18	ARG
2	I	123	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	148	GLN
2	B	153	ASN
2	D	27	GLN
1	G	3	GLN
1	G	13	HIS
1	G	207	ASN
2	I	38	GLN
2	I	148	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

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	A	222/224 (99%)	0.17	4 (1%) 68 74	38, 54, 70, 85	0
1	C	221/224 (98%)	0.22	3 (1%) 75 80	42, 56, 70, 81	0
1	E	222/224 (99%)	0.19	3 (1%) 75 80	40, 55, 68, 88	0
1	G	222/224 (99%)	0.21	2 (0%) 84 88	42, 55, 70, 95	0
2	B	214/215 (99%)	0.23	2 (0%) 84 88	40, 58, 73, 84	0
2	D	214/215 (99%)	0.23	4 (1%) 66 73	45, 57, 71, 77	0
2	F	214/215 (99%)	0.30	6 (2%) 53 60	42, 59, 71, 88	0
2	I	215/215 (100%)	0.22	2 (0%) 84 88	43, 57, 72, 84	0
3	H	11/15 (73%)	0.38	0 100 100	57, 66, 72, 81	0
3	J	12/15 (80%)	0.60	0 100 100	47, 63, 83, 93	0
3	K	11/15 (73%)	0.32	0 100 100	59, 62, 70, 74	0
3	L	12/15 (80%)	0.65	1 (8%) 11 15	51, 58, 78, 89	0
All	All	1790/1816 (98%)	0.23	27 (1%) 73 79	38, 56, 72, 95	0

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	224	CYS	8.6
2	B	1	GLU	3.7
1	G	140	SER	3.3
2	D	78	ILE	3.3
2	D	13	LEU	3.2
1	E	56	ARG	3.1
1	C	140	SER	2.7
1	C	13	HIS	2.7
3	L	812	PRO	2.7
2	D	1	GLU	2.6
1	A	213	THR	2.5

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Mol	Chain	Res	Type	RSRZ
2	F	1	GLU	2.5
2	F	10	ILE	2.5
1	E	224	CYS	2.5
1	E	104	GLU	2.4
2	D	19	ALA	2.3
1	C	80	PHE	2.2
2	I	215	CYS	2.2
1	A	1	GLN	2.2
2	I	96	LEU	2.1
2	B	148	GLN	2.1
1	A	224	CYS	2.1
2	F	40	PRO	2.1
1	A	212	ASN	2.1
2	F	11	VAL	2.1
2	F	29	VAL	2.0
2	F	136	LEU	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.