Full wwPDB X-ray Structure Validation Report ⓘ

Nov 21, 2023 – 09:47 PM JST

PDB ID	:	7X7O
Title	:	SARS-CoV-2 spike RBD in complex with neutralizing antibody UT28K
Authors	:	Ozawa, T.; Tani, H.; Anraku, Y.; Kita, S.; Igarashi, E.; Saga, Y.; Inasaki, N.; Kawasuji, H.; Yamada, H.; Sasaki, S.; Somekawa, M.; Sasaki, J.; Hayakawa, Y.; Yamamoto, Y.; Morinaga, Y.; Kurosawa, N.; Isobe, M.; Fukuhara, H.; Maenaka, K.; Hashiguchi, T.; Kishi, H.; Kitajima, I.; Saito, S.; Niimi, H.
Deposited on	:	2022-03-10
Resolution	:	3.75 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:

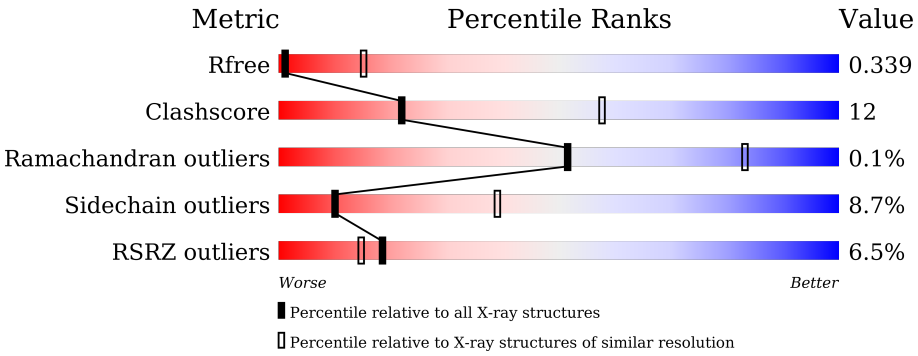
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

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 3.75 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	1039 (3.94-3.58)
Clashscore	141614	1051 (3.92-3.60)
Ramachandran outliers	138981	1015 (3.92-3.60)
Sidechain outliers	138945	1011 (3.92-3.60)
RSRZ outliers	127900	1050 (3.96-3.56)

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	215	7% 57% 20% • 22%
1	B	215	6% 48% 29% • 20%
1	C	215	5% 57% 14% • 27%
1	D	215	7% 53% 20% • 24%
2	E	257	% 30% 16% • 53%
2	G	257	3% 30% 16% • 53%

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Mol	Chain	Length	Quality of chain
2	H	257	3% 31% 13% 53%
2	J	257	2% 30% 16% 53%
3	F	235	4% 36% 9% 54%
3	I	235	3% 36% 10% 54%
3	K	235	2% 33% 12% 54%
3	L	235	3% 35% 10% 54%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 12279 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 Spike protein S1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	167	Total	C	N	O	S	0	0	0
			1334	854	225	249	6			
1	B	172	Total	C	N	O	S	0	0	0
			1371	880	231	253	7			
1	C	157	Total	C	N	O	S	0	0	0
			1261	809	213	233	6			
1	D	163	Total	C	N	O	S	0	0	0
			1309	840	220	243	6			

- Molecule 2 is a protein called UT28K Fab, heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	122	Total	C	N	O	S	0	0	0
			935	587	159	181	8			
2	G	122	Total	C	N	O	S	0	0	0
			935	587	159	181	8			
2	H	122	Total	C	N	O	S	0	0	0
			935	587	159	181	8			
2	J	122	Total	C	N	O	S	0	0	0
			935	587	159	181	8			

- Molecule 3 is a protein called UT28K Fab, light chain.

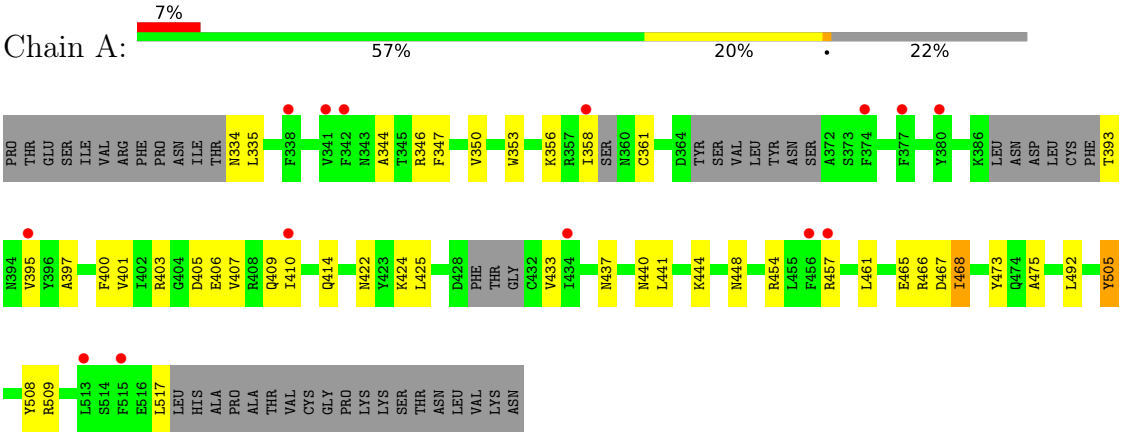
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	F	107	Total	C	N	O	S	0	0	0
			816	513	138	163	2			
3	I	107	Total	C	N	O	S	0	0	0
			816	513	138	163	2			
3	K	107	Total	C	N	O	S	0	0	0
			816	513	138	163	2			
3	L	107	Total	C	N	O	S	0	0	0
			816	513	138	163	2			

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.

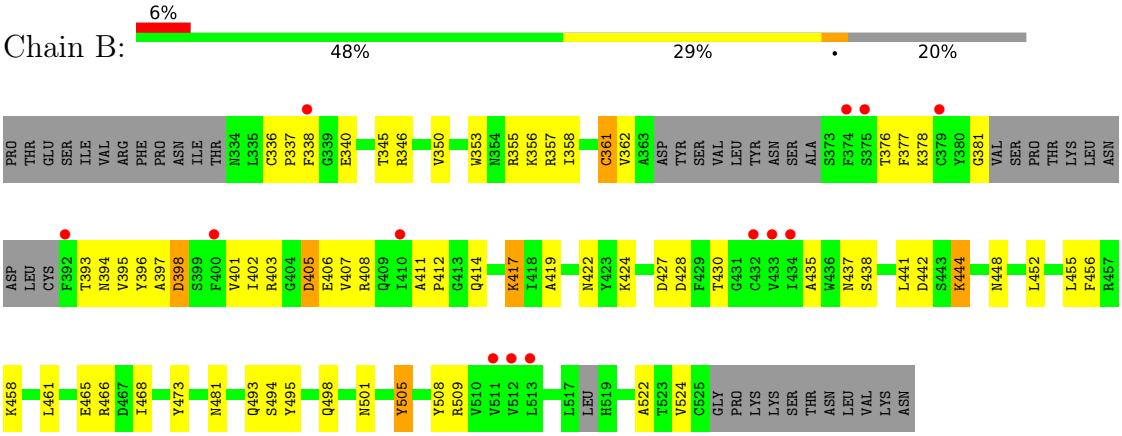
• Molecule 1: Spike protein S1

Chain A:



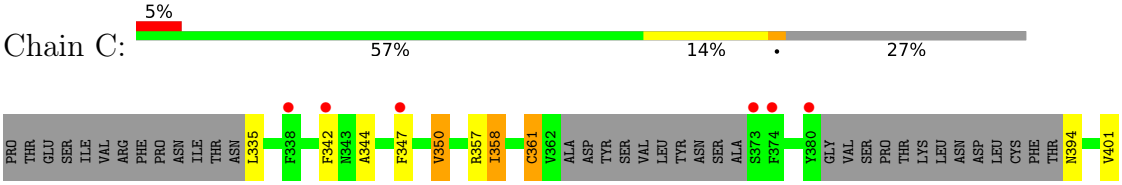
• Molecule 1: Spike protein S1

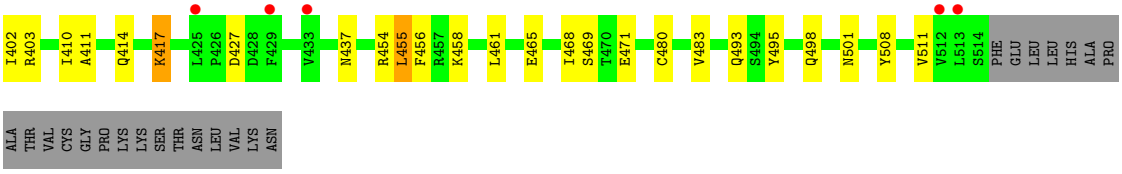
Chain B:



• Molecule 1: Spike protein S1

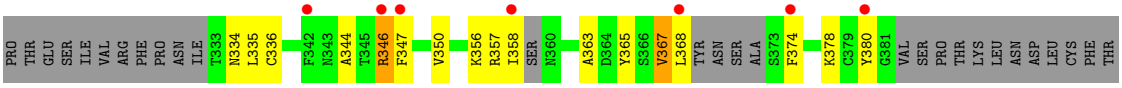
Chain C:

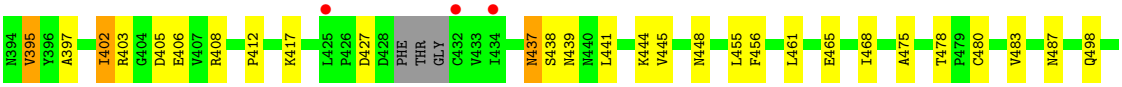


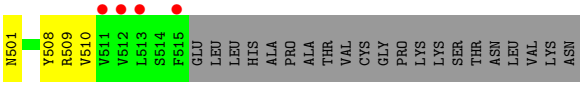


• Molecule 1: Spike protein S1

Chain D:

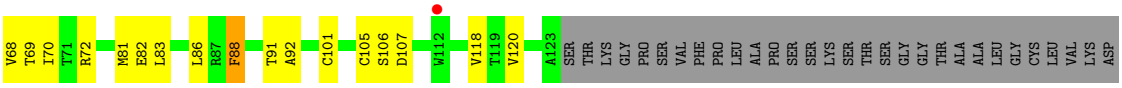






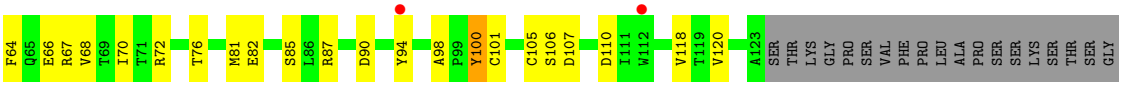
• Molecule 2: UT28K Fab, heavy chain

Chain E:



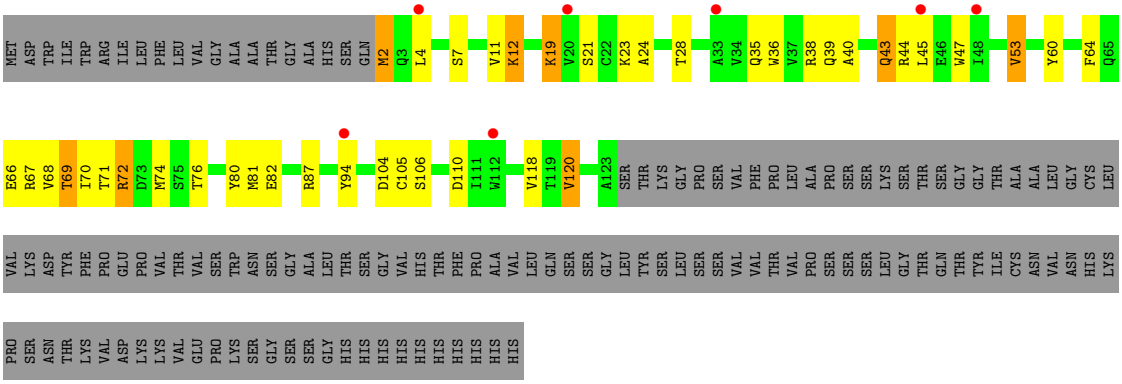
• Molecule 2: UT28K Fab, heavy chain

Chain G:



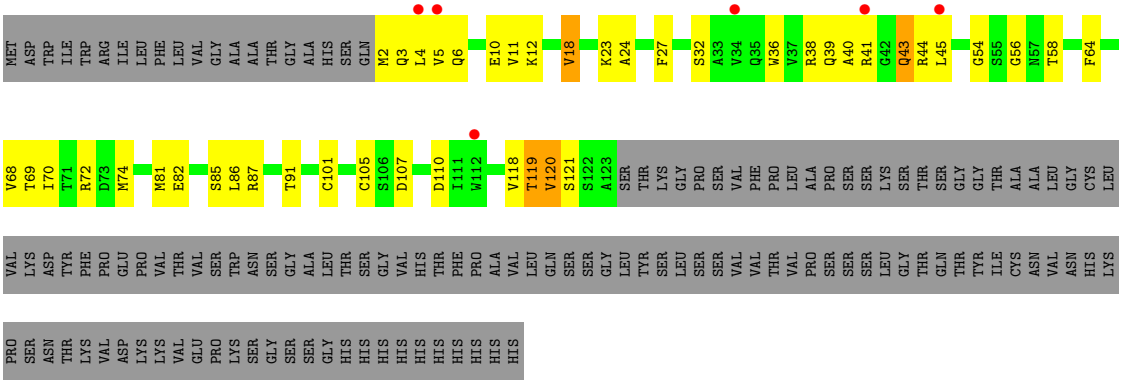
• Molecule 2: UT28K Fab, heavy chain

Chain H:



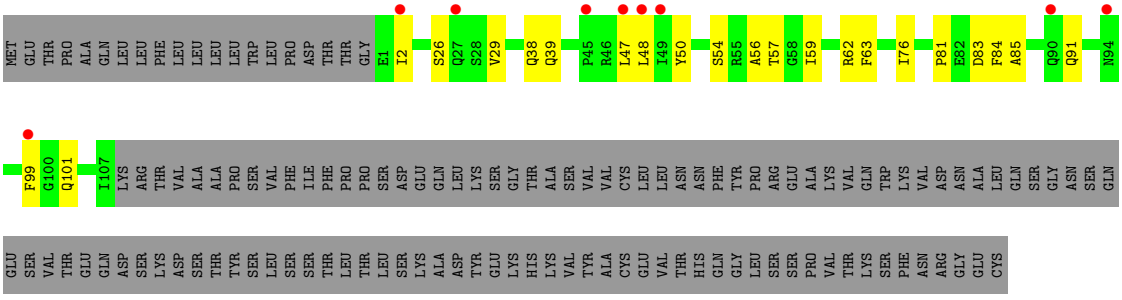
• Molecule 2: UT28K Fab, heavy chain

Chain J:



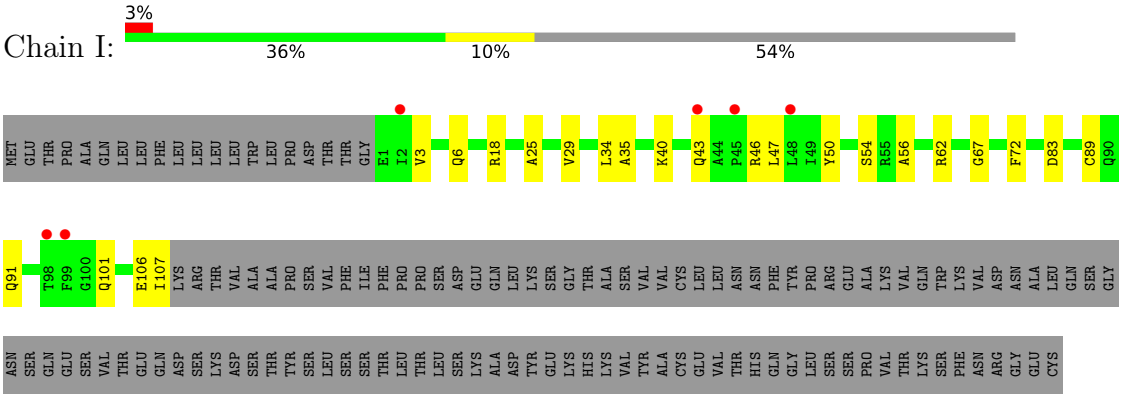
• Molecule 3: UT28K Fab, light chain

Chain F:



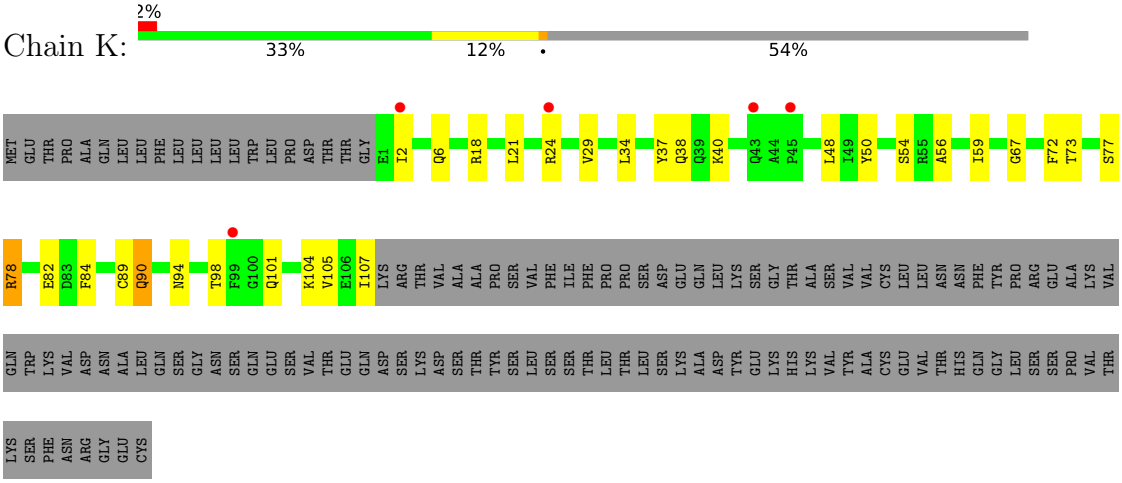
• Molecule 3: UT28K Fab, light chain

Chain I:



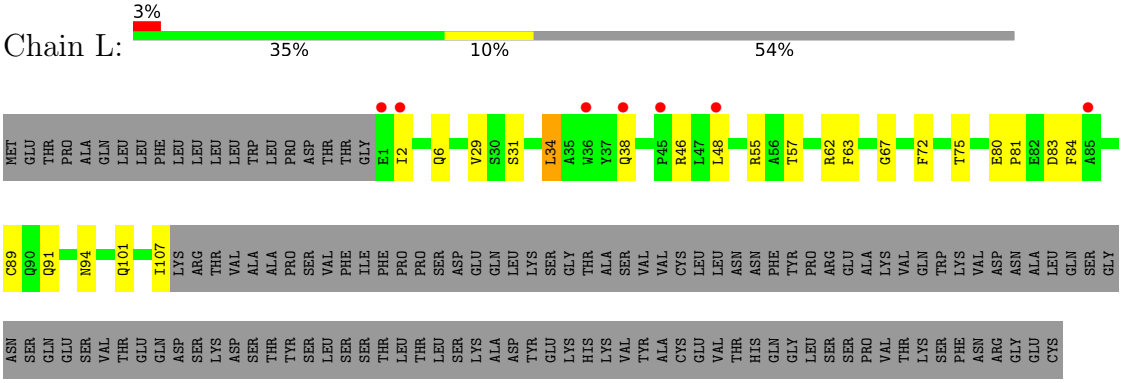
• Molecule 3: UT28K Fab, light chain

Chain K:



• Molecule 3: UT28K Fab, light chain

Chain L:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	I 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	138.25Å 138.41Å 221.11Å 90.00° 90.05° 90.00°	Depositor
Resolution (Å)	48.14 – 3.75 48.91 – 3.75	Depositor EDS
% Data completeness (in resolution range)	99.9 (48.14-3.75) 99.7 (48.91-3.75)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.60 (at 3.77Å)	Xtriage
Refinement program	REFMAC 5.8.0267, PHENIX 1.18_3845	Depositor
R, R_{free}	0.303 , 0.336 0.315 , 0.339	Depositor DCC
R_{free} test set	2193 reflections (5.11%)	wwPDB-VP
Wilson B-factor (Å^2)	111.2	Xtriage
Anisotropy	0.307	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.32 , 94.8	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.30$	Xtriage
Estimated twinning fraction	0.398 for k,h,-l 0.367 for -k,-h,-l 0.380 for h,-k,-l	Xtriage
F_o,F_c correlation	0.84	EDS
Total number of atoms	12279	wwPDB-VP
Average B, all atoms (Å^2)	133.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.79% 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	A	0.33	0/1367	0.48	0/1851
1	B	0.35	0/1408	0.51	0/1909
1	C	0.34	0/1295	0.48	0/1755
1	D	0.36	0/1342	0.50	0/1817
2	E	0.34	0/955	0.53	0/1293
2	G	0.33	0/955	0.54	0/1293
2	H	0.39	0/955	0.55	0/1293
2	J	0.36	0/955	0.55	0/1293
3	F	0.28	0/836	0.48	0/1137
3	I	0.29	0/836	0.48	0/1137
3	K	0.35	0/836	0.52	0/1137
3	L	0.31	0/836	0.50	0/1137
All	All	0.34	0/12576	0.51	0/17052

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	1334	0	1261	39	0
1	B	1371	0	1290	46	0
1	C	1261	0	1190	16	0
1	D	1309	0	1232	26	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	E	935	0	899	30	0
2	G	935	0	899	40	0
2	H	935	0	899	24	0
2	J	935	0	899	28	0
3	F	816	0	786	14	0
3	I	816	0	786	14	0
3	K	816	0	786	13	0
3	L	816	0	786	14	0
All	All	12279	0	11713	282	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 12.

All (282) 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:A:475:ALA:CB	2:G:31:ILE:HD11	1.71	1.20
1:A:475:ALA:HB1	2:G:31:ILE:HD11	1.10	1.09
1:A:473:TYR:HE2	2:G:31:ILE:HD12	1.20	1.05
2:G:31:ILE:HG13	2:G:31:ILE:O	1.61	0.97
2:G:31:ILE:HG12	2:G:101:CYS:SG	2.08	0.93
1:A:473:TYR:CE2	2:G:31:ILE:HD12	2.04	0.92
1:A:473:TYR:HE2	2:G:31:ILE:CD1	1.87	0.88
2:G:31:ILE:HG12	2:G:105:CYS:SG	2.13	0.87
1:A:475:ALA:CB	2:G:31:ILE:CD1	2.54	0.84
1:B:381:GLY:HA3	1:B:430:THR:HA	1.61	0.81
1:A:410:ILE:HD12	1:A:433:VAL:HG11	1.65	0.77
1:B:412:PRO:HB3	1:B:427:ASP:HA	1.70	0.73
2:G:101:CYS:HA	2:G:105:CYS:HA	1.70	0.72
1:A:475:ALA:HB1	2:G:31:ILE:CD1	2.05	0.72
1:D:358:ILE:HD12	1:D:395:VAL:HG22	1.71	0.72
2:G:31:ILE:O	2:G:31:ILE:CG1	2.36	0.72
2:G:31:ILE:CG1	2:G:105:CYS:SG	2.77	0.72
2:E:2:MET:N	2:E:25:SER:O	2.23	0.72
1:A:403:ARG:HB3	1:A:406:GLU:HG3	1.73	0.71
1:D:358:ILE:HB	1:D:395:VAL:HG13	1.74	0.70
3:K:18:ARG:HE	3:K:77:SER:HA	1.54	0.70
1:A:407:VAL:O	1:A:410:ILE:HG12	1.92	0.70
1:D:358:ILE:HD11	1:D:397:ALA:HB2	1.73	0.69
1:A:334:ASN:ND2	1:A:335:LEU:O	2.25	0.68
2:G:10:GLU:HB2	2:G:118:VAL:HG12	1.73	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:334:ASN:ND2	1:D:335:LEU:O	2.26	0.68
1:B:455:LEU:HD23	1:B:493:GLN:HB2	1.76	0.67
3:I:40:LYS:HB2	3:I:43:GLN:HB2	1.75	0.67
1:A:467:ASP:HA	1:B:481:ASN:HB3	1.77	0.67
1:B:455:LEU:HD12	1:B:456:PHE:CG	2.30	0.67
1:A:475:ALA:HB2	2:G:31:ILE:HD11	1.74	0.67
2:H:39:GLN:HB2	2:H:45:LEU:HD23	1.76	0.67
1:A:358:ILE:HD11	1:A:397:ALA:HB2	1.76	0.66
1:A:475:ALA:HB2	2:G:31:ILE:CD1	2.25	0.65
1:B:350:VAL:HG12	1:B:422:ASN:HB3	1.78	0.64
1:D:498:GLN:HE21	1:D:501:ASN:HD21	1.42	0.64
3:I:67:GLY:HA3	3:I:72:PHE:HA	1.80	0.64
2:J:39:GLN:HB2	2:J:45:LEU:HD23	1.80	0.63
2:E:27:PHE:HA	3:L:57:THR:HG22	1.80	0.63
1:B:442:ASP:OD2	1:B:509:ARG:NE	2.32	0.63
1:A:410:ILE:HG13	1:A:410:ILE:O	1.99	0.62
2:E:101:CYS:HA	2:E:105:CYS:HA	1.80	0.62
3:K:67:GLY:HA3	3:K:72:PHE:HA	1.81	0.62
2:H:40:ALA:N	2:H:43:GLN:OE1	2.30	0.61
3:K:37:TYR:HE1	3:K:90:GLN:HG2	1.65	0.61
3:L:55:ARG:NH1	3:L:63:PHE:O	2.32	0.61
2:G:19:LYS:HG3	2:G:82:GLU:HB2	1.82	0.60
1:D:402:ILE:HD11	1:D:510:VAL:HG21	1.85	0.59
1:A:350:VAL:HA	1:A:400:PHE:HB2	1.84	0.59
1:C:402:ILE:HD13	1:C:410:ILE:HD11	1.84	0.59
1:D:461:LEU:HD22	1:D:465:GLU:HB3	1.85	0.59
2:J:91:THR:HG22	2:J:120:VAL:H	1.68	0.59
1:B:461:LEU:HD22	1:B:465:GLU:HB3	1.83	0.59
2:E:44:ARG:HG3	3:F:101:GLN:HA	1.84	0.59
2:H:38:ARG:HE	2:H:64:PHE:HZ	1.50	0.58
1:B:395:VAL:HG23	1:B:524:VAL:HG11	1.85	0.58
1:A:358:ILE:HB	1:A:395:VAL:HG13	1.86	0.58
2:H:23:LYS:NZ	2:H:76:THR:O	2.23	0.58
2:H:19:LYS:HD2	2:H:80:TYR:HB3	1.84	0.58
2:J:10:GLU:HB2	2:J:118:VAL:HG22	1.86	0.58
1:C:357:ARG:NH2	1:C:394:ASN:OD1	2.38	0.57
1:D:455:LEU:HD12	1:D:456:PHE:CG	2.39	0.57
3:I:29:VAL:HG11	3:I:91:GLN:HB2	1.86	0.57
2:E:19:LYS:HA	2:E:82:GLU:HG2	1.86	0.57
1:A:409:GLN:HA	1:A:414:GLN:HG2	1.85	0.56
1:B:350:VAL:CG1	1:B:422:ASN:HB3	2.36	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:J:40:ALA:N	2:J:43:GLN:OE1	2.36	0.56
3:I:35:ALA:HB1	3:I:47:LEU:CD1	2.36	0.55
2:E:105:CYS:SG	2:E:106:SER:N	2.78	0.55
1:B:403:ARG:HB2	1:B:406:GLU:HG3	1.89	0.55
3:K:77:SER:OG	3:K:78:ARG:CZ	2.55	0.55
2:E:41:ARG:NH1	2:E:92:ALA:HA	2.22	0.55
2:J:44:ARG:HG2	2:J:45:LEU:N	2.22	0.55
1:B:438:SER:HB2	1:B:441:LEU:HB2	1.87	0.54
1:C:411:ALA:HB3	1:C:414:GLN:HG3	1.90	0.54
1:A:473:TYR:CE2	2:G:31:ILE:CD1	2.78	0.54
2:E:2:MET:HB3	2:H:2:MET:O	2.07	0.54
2:G:38:ARG:HG3	2:G:46:GLU:HB2	1.90	0.54
1:A:444:LYS:H	1:A:448:ASN:HB2	1.73	0.53
2:J:2:MET:N	2:J:3:GLN:OE1	2.40	0.53
1:B:455:LEU:HD11	1:B:456:PHE:CD2	2.43	0.53
1:C:461:LEU:HD22	1:C:465:GLU:HB3	1.91	0.53
2:J:18:VAL:HG13	2:J:86:LEU:HD11	1.90	0.53
2:E:44:ARG:NH1	3:F:99:PHE:O	2.41	0.53
1:B:455:LEU:CD1	1:B:456:PHE:CG	2.92	0.53
3:L:62:ARG:NE	3:L:83:ASP:OD2	2.41	0.52
1:A:356:LYS:HD3	1:A:358:ILE:HG12	1.91	0.52
1:B:358:ILE:HD11	1:B:397:ALA:HB2	1.90	0.52
1:B:455:LEU:HD12	1:B:455:LEU:C	2.29	0.52
1:B:345:THR:HG23	1:B:346:ARG:N	2.24	0.52
2:E:41:ARG:HH12	2:E:92:ALA:HA	1.74	0.52
1:B:411:ALA:HB3	1:B:414:GLN:CG	2.40	0.52
2:J:36:TRP:CE2	2:J:81:MET:HB2	2.43	0.52
1:A:410:ILE:CD1	1:A:433:VAL:HG11	2.35	0.52
1:A:440:ASN:OD1	1:A:441:LEU:HG	2.10	0.52
2:G:60:TYR:HE1	2:G:70:ILE:HG13	1.74	0.52
1:B:356:LYS:HD3	1:B:358:ILE:HG12	1.93	0.51
2:E:28:THR:HG23	3:L:57:THR:HG23	1.92	0.51
1:D:367:VAL:HG13	1:D:368:LEU:HD12	1.92	0.51
1:A:401:VAL:HG22	1:A:509:ARG:HD3	1.93	0.51
1:D:438:SER:HB2	1:D:441:LEU:HB2	1.93	0.51
2:H:60:TYR:HE2	2:H:70:ILE:HG13	1.76	0.51
1:B:393:THR:HA	1:B:522:ALA:HA	1.91	0.51
2:J:91:THR:HG22	2:J:119:THR:HA	1.93	0.51
3:L:38:GLN:HB2	3:L:48:LEU:HD11	1.93	0.51
2:G:36:TRP:CE2	2:G:81:MET:HB2	2.45	0.51
1:B:437:ASN:HB2	1:B:508:TYR:CZ	2.45	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:19:LYS:HG2	2:E:82:GLU:OE2	2.10	0.50
2:G:12:LYS:NZ	2:G:18:VAL:HA	2.27	0.50
1:B:338:PHE:HE1	1:B:358:ILE:HD13	1.76	0.50
3:K:6:GLN:H	3:K:101:GLN:HE22	1.59	0.50
1:D:444:LYS:H	1:D:448:ASN:HB2	1.76	0.50
2:H:35:GLN:HB2	2:H:47:TRP:HE1	1.76	0.50
2:J:85:SER:HB2	2:J:87:ARG:HH21	1.75	0.50
3:K:84:PHE:CD1	3:K:105:VAL:HG12	2.47	0.50
1:A:410:ILE:HD12	1:A:433:VAL:CG1	2.40	0.49
1:A:424:LYS:NZ	1:A:425:LEU:O	2.45	0.49
2:E:64:PHE:O	2:E:68:VAL:HG22	2.12	0.49
1:D:405:ASP:O	1:D:408:ARG:HG2	2.12	0.49
2:H:4:LEU:HD23	2:H:24:ALA:HA	1.94	0.49
3:K:38:GLN:HB2	3:K:48:LEU:HD11	1.94	0.49
1:D:403:ARG:HB3	1:D:406:GLU:HG3	1.92	0.49
2:J:32:SER:OG	2:J:101:CYS:HB2	2.12	0.49
2:G:12:LYS:O	2:G:120:VAL:HA	2.13	0.49
1:B:353:TRP:CD1	1:B:466:ARG:HD3	2.48	0.49
1:D:344:ALA:HB3	1:D:347:PHE:HE1	1.78	0.49
3:I:35:ALA:HB1	3:I:47:LEU:HD11	1.95	0.48
2:J:12:LYS:HG3	2:J:18:VAL:CG1	2.43	0.48
1:B:501:ASN:HB3	1:B:505:TYR:HB3	1.94	0.48
1:A:461:LEU:HG	1:A:465:GLU:HB3	1.95	0.48
2:E:36:TRP:CE2	2:E:81:MET:HB2	2.47	0.48
1:B:376:THR:CG2	1:B:435:ALA:HB3	2.43	0.48
1:C:344:ALA:HB3	1:C:347:PHE:HE1	1.77	0.48
3:L:2:ILE:HD11	3:L:94:ASN:HB2	1.95	0.48
1:A:393:THR:N	1:A:517:LEU:HA	2.29	0.48
1:D:456:PHE:CZ	2:J:54:GLY:HA3	2.48	0.48
2:G:38:ARG:NH2	2:G:94:TYR:OH	2.47	0.48
3:I:3:VAL:O	3:I:25:ALA:HA	2.14	0.48
3:L:67:GLY:HA3	3:L:72:PHE:HA	1.95	0.48
2:G:100:TYR:HB3	2:G:110:ASP:OD2	2.14	0.48
1:B:444:LYS:HB3	1:B:448:ASN:HB2	1.94	0.47
2:E:2:MET:N	2:E:26:GLY:HA3	2.29	0.47
3:F:63:PHE:CD1	3:F:76:ILE:HD12	2.49	0.47
2:G:64:PHE:HB3	2:G:68:VAL:HG13	1.96	0.47
1:C:417:LYS:HD2	1:C:455:LEU:HA	1.97	0.47
2:H:43:GLN:HG2	2:H:44:ARG:N	2.30	0.47
2:G:110:ASP:O	3:I:46:ARG:NH2	2.47	0.47
1:C:342:PHE:CE1	1:C:511:VAL:HG11	2.49	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:F:48:LEU:HA	3:F:59:ILE:HG12	1.96	0.47
2:H:110:ASP:O	3:L:46:ARG:NH1	2.47	0.47
3:I:6:GLN:H	3:I:101:GLN:HE22	1.60	0.47
1:B:411:ALA:HB3	1:B:414:GLN:HG3	1.97	0.47
1:B:455:LEU:CD1	1:B:456:PHE:CD2	2.97	0.47
3:F:38:GLN:HB2	3:F:48:LEU:HD11	1.97	0.47
2:G:63:LYS:HD3	2:G:64:PHE:CE2	2.50	0.47
2:E:18:VAL:HG11	2:E:118:VAL:HG11	1.97	0.47
1:B:403:ARG:HG3	1:B:495:TYR:CE1	2.49	0.46
2:J:43:GLN:HG2	2:J:44:ARG:N	2.29	0.46
3:L:6:GLN:H	3:L:101:GLN:HE22	1.63	0.46
1:A:457:ARG:HD2	1:A:457:ARG:HA	1.51	0.46
1:C:456:PHE:CZ	2:E:54:GLY:HA3	2.51	0.46
1:B:336:CYS:N	1:B:361:CYS:SG	2.88	0.46
1:D:438:SER:HB3	1:D:509:ARG:HG3	1.96	0.46
2:E:39:GLN:NE2	2:E:43:GLN:O	2.49	0.46
1:D:444:LYS:NZ	1:D:448:ASN:HA	2.31	0.46
2:H:43:GLN:HG2	2:H:44:ARG:H	1.81	0.46
2:E:60:TYR:HE2	2:E:70:ILE:HG13	1.80	0.46
1:B:417:LYS:HG2	1:B:455:LEU:HA	1.97	0.46
1:B:419:ALA:O	1:B:424:LYS:HD3	2.16	0.46
2:G:68:VAL:HA	2:G:82:GLU:O	2.15	0.46
2:J:12:LYS:O	2:J:120:VAL:HA	2.16	0.46
1:C:437:ASN:HB2	1:C:508:TYR:CZ	2.51	0.46
3:I:47:LEU:HD21	3:I:50:TYR:HB3	1.97	0.46
2:J:4:LEU:HD23	2:J:24:ALA:HA	1.98	0.45
1:B:411:ALA:O	1:B:414:GLN:HG2	2.15	0.45
3:K:6:GLN:H	3:K:101:GLN:NE2	2.13	0.45
1:A:437:ASN:HB2	1:A:508:TYR:CZ	2.51	0.45
3:I:6:GLN:H	3:I:101:GLN:NE2	2.13	0.45
1:A:353:TRP:CE2	1:A:466:ARG:HD2	2.52	0.45
1:D:363:ALA:HB1	1:D:365:TYR:CE1	2.51	0.45
2:G:13:LYS:O	2:G:16:THR:HG22	2.16	0.45
1:A:468:ILE:H	1:A:468:ILE:HG12	1.66	0.45
1:B:455:LEU:HD12	1:B:456:PHE:N	2.31	0.45
2:E:4:LEU:HD23	2:E:24:ALA:HA	1.99	0.45
2:H:69:THR:HG23	2:H:82:GLU:HG3	1.99	0.45
2:H:38:ARG:HD3	2:H:94:TYR:CZ	2.52	0.44
1:A:454:ARG:HA	1:A:492:LEU:HD23	1.99	0.44
1:B:357:ARG:NH2	1:B:394:ASN:ND2	2.65	0.44
1:B:405:ASP:O	1:B:408:ARG:HG2	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:J:18:VAL:O	2:J:82:GLU:HA	2.16	0.44
1:B:337:PRO:HB2	1:B:340:GLU:HG3	1.99	0.44
3:F:29:VAL:HG11	3:F:91:GLN:HB2	1.99	0.44
1:A:347:PHE:CD2	1:A:509:ARG:HD2	2.52	0.44
1:A:344:ALA:HB3	1:A:347:PHE:HE1	1.81	0.44
2:E:35:GLN:HB2	2:E:47:TRP:HE1	1.82	0.44
1:B:378:LYS:HD2	1:B:378:LYS:HA	1.70	0.44
2:E:88:PHE:O	2:E:91:THR:HG22	2.18	0.44
2:J:64:PHE:O	2:J:68:VAL:HG22	2.18	0.44
2:J:68:VAL:HA	2:J:82:GLU:O	2.17	0.44
2:H:7:SER:HB3	2:H:21:SER:H	1.83	0.43
2:H:36:TRP:CE2	2:H:81:MET:HB2	2.52	0.43
2:J:101:CYS:HA	2:J:105:CYS:HA	2.00	0.43
3:L:84:PHE:CE2	3:L:107:ILE:HD11	2.53	0.43
2:G:28:THR:OG1	2:G:101:CYS:O	2.35	0.43
3:K:59:ILE:HD13	3:K:59:ILE:HA	1.85	0.43
1:C:350:VAL:HG22	1:C:401:VAL:O	2.18	0.43
2:E:18:VAL:HG12	2:E:86:LEU:HD21	1.99	0.43
3:K:2:ILE:HD13	3:K:29:VAL:HG12	1.99	0.43
1:B:498:GLN:H	1:B:501:ASN:ND2	2.16	0.43
3:F:56:ALA:HB3	3:F:59:ILE:HD13	2.00	0.43
3:I:50:TYR:O	3:I:54:SER:OG	2.29	0.43
1:B:452:LEU:HD23	1:B:494:SER:HA	2.01	0.43
1:D:344:ALA:HB3	1:D:347:PHE:CE1	2.54	0.43
1:D:417:LYS:HD3	1:D:417:LYS:HA	1.79	0.43
2:E:12:LYS:O	2:E:120:VAL:HA	2.19	0.43
2:G:100:TYR:O	2:G:106:SER:HB3	2.18	0.43
2:J:18:VAL:HG22	2:J:86:LEU:HD21	2.00	0.43
1:B:357:ARG:NH1	1:B:396:TYR:OH	2.52	0.43
1:C:344:ALA:HB3	1:C:347:PHE:CE1	2.53	0.43
2:H:53:VAL:O	2:H:72:ARG:HD3	2.18	0.43
2:J:43:GLN:HG2	2:J:44:ARG:H	1.84	0.43
3:L:29:VAL:HG11	3:L:91:GLN:HB2	2.00	0.42
3:L:81:PRO:HA	3:L:84:PHE:CE2	2.54	0.42
1:A:475:ALA:HB1	2:G:105:CYS:SG	2.59	0.42
3:F:63:PHE:HD1	3:F:76:ILE:HD12	1.85	0.42
2:G:67:ARG:NH2	2:G:90:ASP:OD2	2.52	0.42
2:G:85:SER:HB2	2:G:87:ARG:HH21	1.85	0.42
2:H:66:GLU:CD	2:H:66:GLU:H	2.23	0.42
2:H:71:THR:HG23	2:H:80:TYR:HB2	2.01	0.42
2:E:39:GLN:HB2	2:E:45:LEU:HD23	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:505:TYR:HD1	1:A:505:TYR:O	2.02	0.42
2:E:67:ARG:HE	2:E:83:LEU:HD11	1.83	0.42
3:I:62:ARG:NE	3:I:83:ASP:OD2	2.53	0.42
2:E:19:LYS:CB	2:E:82:GLU:HG2	2.50	0.42
2:G:37:VAL:HG11	2:G:45:LEU:HD23	2.01	0.42
1:B:402:ILE:HD11	1:B:407:VAL:HG12	2.01	0.42
3:F:47:LEU:HD21	3:F:50:TYR:HB3	2.02	0.42
3:I:106:GLU:HG2	3:I:107:ILE:H	1.84	0.42
3:L:84:PHE:CD2	3:L:107:ILE:HD11	2.54	0.42
1:C:357:ARG:NH1	1:C:358:ILE:O	2.53	0.42
2:H:38:ARG:HB3	2:H:94:TYR:CD2	2.55	0.41
2:J:56:GLY:HA3	2:J:72:ARG:HH21	1.85	0.41
3:F:2:ILE:HD13	3:F:29:VAL:HG22	2.01	0.41
1:C:498:GLN:H	1:C:501:ASN:ND2	2.18	0.41
2:J:27:PHE:HE2	2:J:32:SER:HG	1.68	0.41
3:L:34:LEU:HD13	3:L:72:PHE:CE1	2.55	0.41
1:D:475:ALA:O	1:D:487:ASN:HB3	2.20	0.41
3:K:40:LYS:HE2	3:K:82:GLU:O	2.21	0.41
1:C:403:ARG:HB2	1:C:495:TYR:CE1	2.56	0.41
2:E:68:VAL:HA	2:E:82:GLU:O	2.21	0.41
1:D:437:ASN:HB3	1:D:508:TYR:CZ	2.56	0.41
2:E:18:VAL:O	2:E:82:GLU:HA	2.21	0.41
1:B:456:PHE:HB3	1:B:473:TYR:CD2	2.56	0.41
1:C:335:LEU:HD22	1:C:361:CYS:HB2	2.01	0.41
1:D:498:GLN:HE21	1:D:501:ASN:ND2	2.16	0.41
2:J:81:MET:HE3	2:J:81:MET:HB3	1.99	0.41
1:B:361:CYS:SG	1:B:362:VAL:N	2.94	0.41
2:H:12:LYS:O	2:H:120:VAL:HA	2.21	0.41
2:H:67:ARG:HG2	2:H:87:ARG:HH22	1.86	0.41
2:J:6:GLN:HE21	2:J:6:GLN:HB3	1.68	0.41
3:K:50:TYR:CE2	3:K:56:ALA:HA	2.55	0.41
1:B:398:ASP:N	1:B:398:ASP:OD1	2.52	0.41
2:E:19:LYS:CA	2:E:82:GLU:HG2	2.51	0.41
2:G:64:PHE:O	2:G:68:VAL:HG22	2.21	0.41
1:D:412:PRO:HB3	1:D:427:ASP:HA	2.02	0.40
2:G:23:LYS:NZ	2:G:76:THR:O	2.51	0.40
3:I:50:TYR:CE2	3:I:56:ALA:HA	2.57	0.40
1:D:368:LEU:HA	1:D:374:PHE:CE2	2.56	0.40
3:F:39:GLN:O	3:F:85:ALA:HB1	2.21	0.40
3:F:57:THR:HG23	2:H:28:THR:HG23	2.02	0.40
3:F:62:ARG:NE	3:F:83:ASP:OD2	2.54	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:J:74:MET:H	2:J:74:MET:HG2	1.73	0.40
3:K:104:LYS:NZ	3:K:105:VAL:O	2.54	0.40
2:G:4:LEU:HD11	2:G:98:ALA:HB3	2.03	0.40
1:C:454:ARG:HH22	1:C:469:SER:N	2.19	0.40
2:H:68:VAL:HA	2:H:82:GLU:O	2.21	0.40
2:J:70:ILE:HG13	2:J:81:MET:HG2	2.04	0.40
1:A:350:VAL:HG22	1:A:422:ASN:HB3	2.03	0.40
1:B:350:VAL:CG2	1:B:401:VAL:O	2.69	0.40
1:D:346:ARG:HE	1:D:346:ARG:HB3	1.76	0.40
3:F:81:PRO:O	3:F:84:PHE:HD2	2.04	0.40
2:G:66:GLU:H	2:G:66:GLU:CD	2.24	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	A	157/215 (73%)	147 (94%)	10 (6%)	0	100	100
1	B	164/215 (76%)	154 (94%)	10 (6%)	0	100	100
1	C	151/215 (70%)	141 (93%)	10 (7%)	0	100	100
1	D	153/215 (71%)	143 (94%)	10 (6%)	0	100	100
2	E	120/257 (47%)	115 (96%)	5 (4%)	0	100	100
2	G	120/257 (47%)	109 (91%)	10 (8%)	1 (1%)	19	56
2	H	120/257 (47%)	114 (95%)	6 (5%)	0	100	100
2	J	120/257 (47%)	113 (94%)	7 (6%)	0	100	100
3	F	105/235 (45%)	99 (94%)	6 (6%)	0	100	100
3	I	105/235 (45%)	99 (94%)	6 (6%)	0	100	100
3	K	105/235 (45%)	98 (93%)	7 (7%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
3	L	105/235 (45%)	99 (94%)	6 (6%)	0	100 100
All	All	1525/2828 (54%)	1431 (94%)	93 (6%)	1 (0%)	51 83

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	G	100	TYR

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	144/188 (77%)	139 (96%)	5 (4%)	36 63
1	B	147/188 (78%)	136 (92%)	11 (8%)	13 44
1	C	136/188 (72%)	124 (91%)	12 (9%)	10 38
1	D	141/188 (75%)	124 (88%)	17 (12%)	5 25
2	E	101/216 (47%)	89 (88%)	12 (12%)	5 26
2	G	101/216 (47%)	92 (91%)	9 (9%)	9 38
2	H	101/216 (47%)	87 (86%)	14 (14%)	3 21
2	J	101/216 (47%)	87 (86%)	14 (14%)	3 21
3	F	89/204 (44%)	87 (98%)	2 (2%)	52 73
3	I	89/204 (44%)	86 (97%)	3 (3%)	37 64
3	K	89/204 (44%)	78 (88%)	11 (12%)	4 24
3	L	89/204 (44%)	84 (94%)	5 (6%)	21 53
All	All	1328/2432 (55%)	1213 (91%)	115 (9%)	10 39

All (115) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	346	ARG
1	A	361	CYS

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Mol	Chain	Res	Type
1	A	405	ASP
1	A	468	ILE
1	A	505	TYR
1	B	355	ARG
1	B	361	CYS
1	B	377	PHE
1	B	398	ASP
1	B	405	ASP
1	B	417	LYS
1	B	428	ASP
1	B	444	LYS
1	B	458	LYS
1	B	468	ILE
1	B	505	TYR
1	C	350	VAL
1	C	358	ILE
1	C	361	CYS
1	C	417	LYS
1	C	427	ASP
1	C	455	LEU
1	C	458	LYS
1	C	468	ILE
1	C	471	GLU
1	C	480	CYS
1	C	483	VAL
1	C	493	GLN
1	D	336	CYS
1	D	346	ARG
1	D	350	VAL
1	D	356	LYS
1	D	357	ARG
1	D	367	VAL
1	D	378	LYS
1	D	380	TYR
1	D	395	VAL
1	D	402	ILE
1	D	437	ASN
1	D	439	ASN
1	D	445	VAL
1	D	468	ILE
1	D	478	THR
1	D	480	CYS

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Mol	Chain	Res	Type
1	D	483	VAL
2	E	3	GLN
2	E	11	VAL
2	E	13	LYS
2	E	16	THR
2	E	19	LYS
2	E	41	ARG
2	E	46	GLU
2	E	63	LYS
2	E	69	THR
2	E	72	ARG
2	E	88	PHE
2	E	107	ASP
3	F	26	SER
3	F	54	SER
2	G	2	MET
2	G	5	VAL
2	G	20	VAL
2	G	31	ILE
2	G	38	ARG
2	G	41	ARG
2	G	44	ARG
2	G	72	ARG
2	G	107	ASP
2	H	2	MET
2	H	11	VAL
2	H	12	LYS
2	H	19	LYS
2	H	43	GLN
2	H	53	VAL
2	H	69	THR
2	H	72	ARG
2	H	74	MET
2	H	104	ASP
2	H	105	CYS
2	H	106	SER
2	H	118	VAL
2	H	120	VAL
3	I	18	ARG
3	I	34	LEU
3	I	89	CYS
2	J	5	VAL

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Mol	Chain	Res	Type
2	J	11	VAL
2	J	18	VAL
2	J	23	LYS
2	J	38	ARG
2	J	41	ARG
2	J	43	GLN
2	J	58	THR
2	J	69	THR
2	J	107	ASP
2	J	110	ASP
2	J	119	THR
2	J	120	VAL
2	J	121	SER
3	K	21	LEU
3	K	24	ARG
3	K	34	LEU
3	K	54	SER
3	K	73	THR
3	K	78	ARG
3	K	89	CYS
3	K	90	GLN
3	K	94	ASN
3	K	98	THR
3	K	107	ILE
3	L	31	SER
3	L	34	LEU
3	L	75	THR
3	L	80	GLU
3	L	89	CYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	501	ASN
1	C	493	GLN
1	D	498	GLN
2	E	39	GLN

5.3.3 RNA

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

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	167/215 (77%)	0.48	14 (8%)	11	10	85, 136, 214, 248	0
1	B	172/215 (80%)	0.47	13 (7%)	13	11	83, 140, 207, 267	0
1	C	157/215 (73%)	0.46	11 (7%)	16	13	81, 131, 200, 287	0
1	D	163/215 (75%)	0.54	14 (8%)	10	9	64, 137, 202, 283	0
2	E	122/257 (47%)	0.36	3 (2%)	57	51	72, 120, 172, 250	0
2	G	122/257 (47%)	0.55	7 (5%)	23	20	76, 122, 182, 202	0
2	H	122/257 (47%)	0.46	7 (5%)	23	20	62, 117, 172, 198	0
2	J	122/257 (47%)	0.45	6 (4%)	29	26	75, 120, 179, 209	0
3	F	107/235 (45%)	0.58	9 (8%)	11	10	80, 127, 178, 200	0
3	I	107/235 (45%)	0.49	6 (5%)	24	21	88, 125, 171, 198	0
3	K	107/235 (45%)	0.43	5 (4%)	31	28	77, 134, 176, 234	0
3	L	107/235 (45%)	0.60	7 (6%)	18	15	87, 128, 179, 203	0
All	All	1575/2828 (55%)	0.49	102 (6%)	18	15	62, 130, 193, 287	0

All (102) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	373	SER	5.5
2	G	48	ILE	4.4
1	A	341	VAL	4.4
1	A	342	PHE	4.3
1	D	432	CYS	4.0
1	B	513	LEU	4.0
1	C	347	PHE	4.0
1	D	346	ARG	4.0
1	B	432	CYS	3.9
1	A	457	ARG	3.9
1	A	358	ILE	3.9

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Mol	Chain	Res	Type	RSRZ
1	A	338	PHE	3.8
1	B	374	PHE	3.7
1	D	513	LEU	3.7
1	A	410	ILE	3.7
1	D	368	LEU	3.6
3	I	2	ILE	3.5
2	G	45	LEU	3.4
3	K	45	PRO	3.4
1	A	374	PHE	3.4
2	J	45	LEU	3.3
2	G	47	TRP	3.2
3	L	45	PRO	3.2
1	B	512	VAL	3.2
3	K	99	PHE	3.2
3	F	49	ILE	3.2
1	B	375	SER	3.2
1	B	379	CYS	3.1
1	C	374	PHE	3.1
3	F	47	LEU	3.1
1	A	515	PHE	3.1
2	J	112	TRP	3.1
3	I	45	PRO	3.1
3	F	27	GLN	3.0
2	G	112	TRP	3.0
1	C	338	PHE	3.0
1	D	347	PHE	3.0
1	D	425	LEU	2.9
1	D	342	PHE	2.9
1	B	410	ILE	2.9
1	D	512	VAL	2.9
1	B	433	VAL	2.8
1	D	515	PHE	2.8
2	H	45	LEU	2.8
1	B	392	PHE	2.8
3	I	98	THR	2.8
3	K	24	ARG	2.7
2	H	94	TYR	2.7
2	E	45	LEU	2.7
3	F	2	ILE	2.7
1	A	377	PHE	2.6
1	C	429	PHE	2.6
3	L	48	LEU	2.6

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Mol	Chain	Res	Type	RSRZ
1	C	342	PHE	2.5
1	B	511	VAL	2.5
1	A	513	LEU	2.5
1	C	380	TYR	2.5
2	G	34	VAL	2.5
3	I	48	LEU	2.5
3	F	94	ASN	2.5
3	F	99	PHE	2.5
2	H	48	ILE	2.5
2	E	112	TRP	2.5
3	F	48	LEU	2.4
1	A	380	TYR	2.4
3	L	2	ILE	2.4
1	D	380	TYR	2.4
1	D	434	ILE	2.4
1	D	511	VAL	2.4
2	J	4	LEU	2.4
2	G	94	TYR	2.3
2	J	5	VAL	2.3
3	I	43	GLN	2.3
1	B	434	ILE	2.3
1	C	512	VAL	2.3
3	L	38	GLN	2.3
1	B	338	PHE	2.3
1	D	374	PHE	2.3
3	L	1	GLU	2.3
1	C	425	LEU	2.3
1	B	400	PHE	2.3
1	A	434	ILE	2.3
2	E	41	ARG	2.3
2	G	49	GLY	2.3
2	J	34	VAL	2.3
1	C	513	LEU	2.2
1	D	358	ILE	2.2
2	H	4	LEU	2.2
1	A	456	PHE	2.2
1	C	433	VAL	2.2
2	H	20	VAL	2.1
2	J	41	ARG	2.1
3	L	36	TRP	2.1
3	F	45	PRO	2.1
1	A	395	VAL	2.1

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Mol	Chain	Res	Type	RSRZ
2	H	33	ALA	2.1
3	K	2	ILE	2.1
3	F	90	GLN	2.0
3	I	99	PHE	2.0
2	H	112	TRP	2.0
3	K	43	GLN	2.0
3	L	85	ALA	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.