Full wwPDB X-ray Structure Validation Report (i)

Nov 14, 2023 – 11:32 PM JST

PDB ID	:	6IGF
Title	:	Crystal structure of Human Papillomavirus type 52 pentamer
Authors	:	Li, Z.H.; Song, S.; He, M.Z.; Gu, Y.; Li, S.W.
Deposited on	:	2018-09-25
Resolution	:	2.75 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 (i) symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

The following versions of software and data (see references (1)) were used in the production of this report:

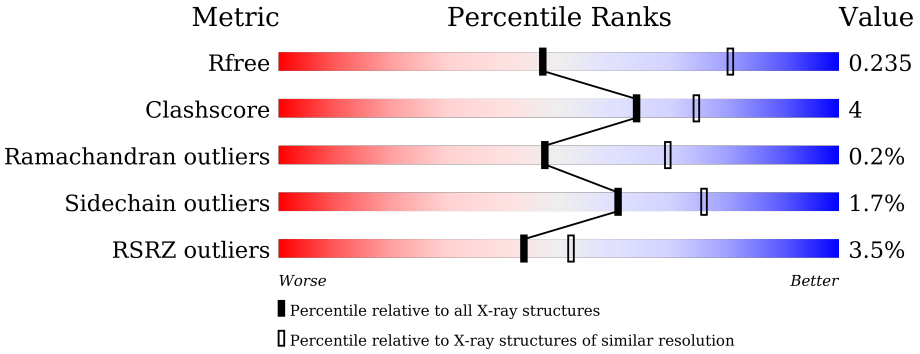
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation 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 2.75 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	1235 (2.78-2.74)
Clashscore	141614	1277 (2.78-2.74)
Ramachandran outliers	138981	1257 (2.78-2.74)
Sidechain outliers	138945	1257 (2.78-2.74)
RSRZ outliers	127900	1207 (2.78-2.74)

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	529	4% 69% 9% • 22%
1	B	529	3% 72% 7% 20%
1	C	529	3% 73% 8% 19%
1	D	529	2% 68% 10% • 22%
1	E	529	2% 73% 8% 19%
1	F	529	3% 69% 9% • 21%

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Mol	Chain	Length	Quality of chain
1	G	529	2% 69% 9% • 21%
1	H	529	3% 68% 9% • 21%
1	I	529	2% 70% 8% • 21%
1	J	529	2% 70% 9% 21%

2 Entry composition (i)

There is only 1 type of molecule in this entry. The entry contains 33255 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 Major capsid protein L1.

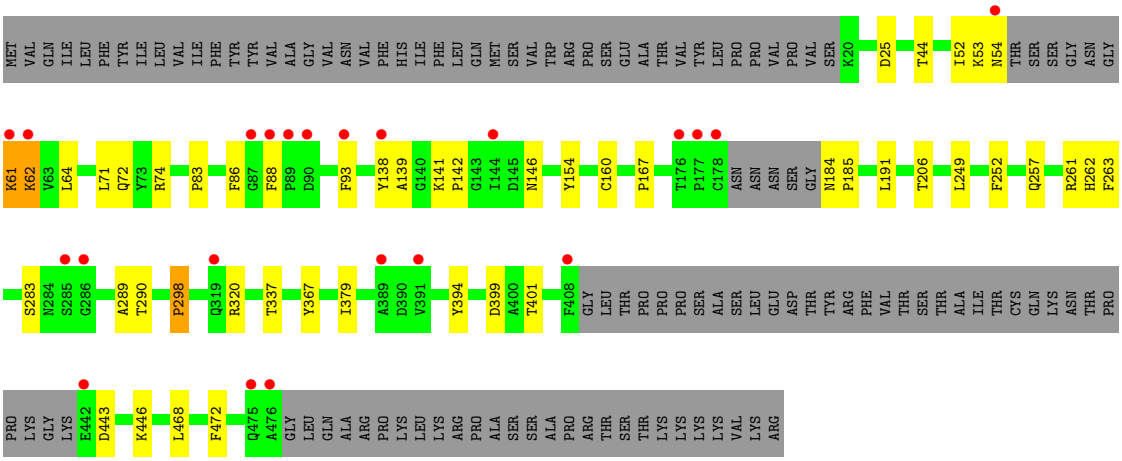
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	413	3293	2105	544	621	23	0	1	0
1	B	421	3344	2139	552	631	22	0	1	0
1	C	427	3377	2149	562	644	22	0	0	0
1	D	414	3291	2103	545	621	22	0	0	0
1	E	429	3400	2168	567	642	23	0	1	0
1	F	417	3316	2119	548	626	23	0	1	0
1	G	418	3322	2120	550	630	22	0	1	0
1	H	416	3305	2111	547	625	22	0	0	0
1	I	416	3300	2108	547	623	22	0	0	0
1	J	417	3307	2113	548	625	21	0	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.

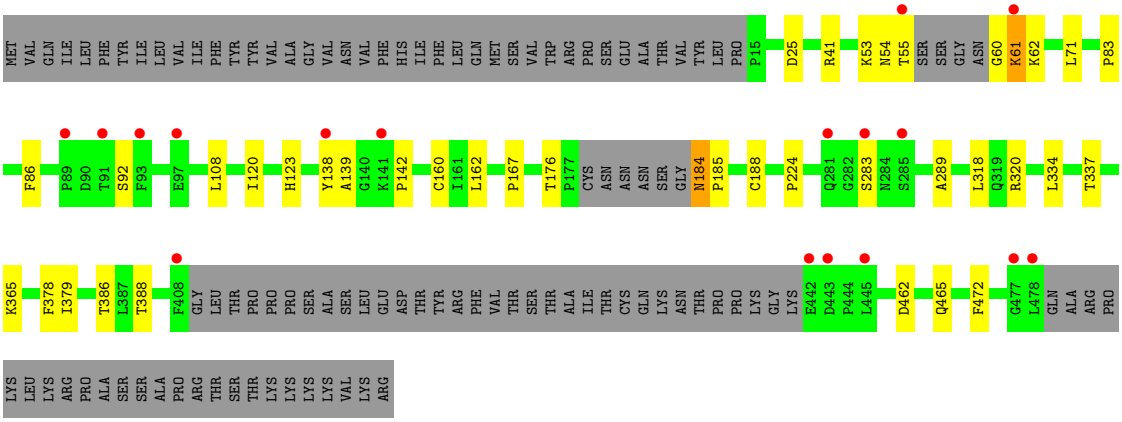
• Molecule 1: Major capsid protein L1

Chain A:



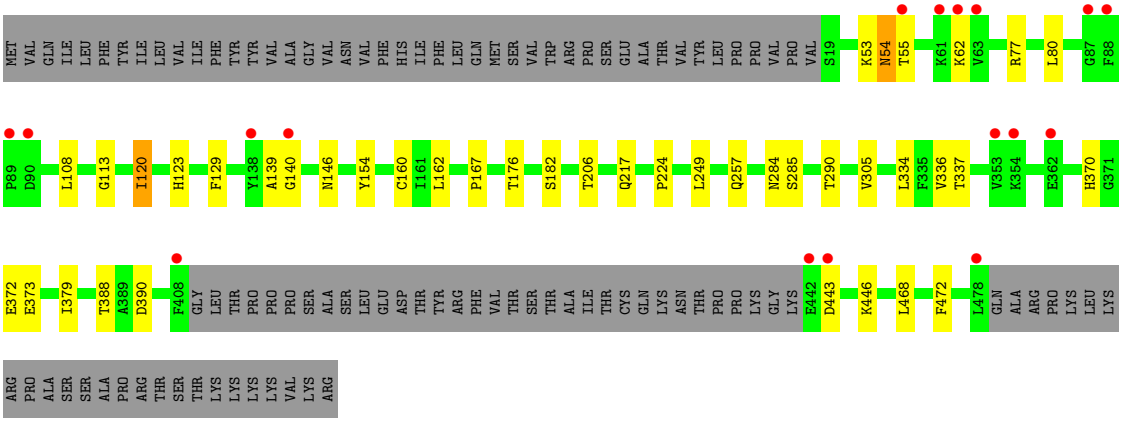
• Molecule 1: Major capsid protein L1

Chain B:



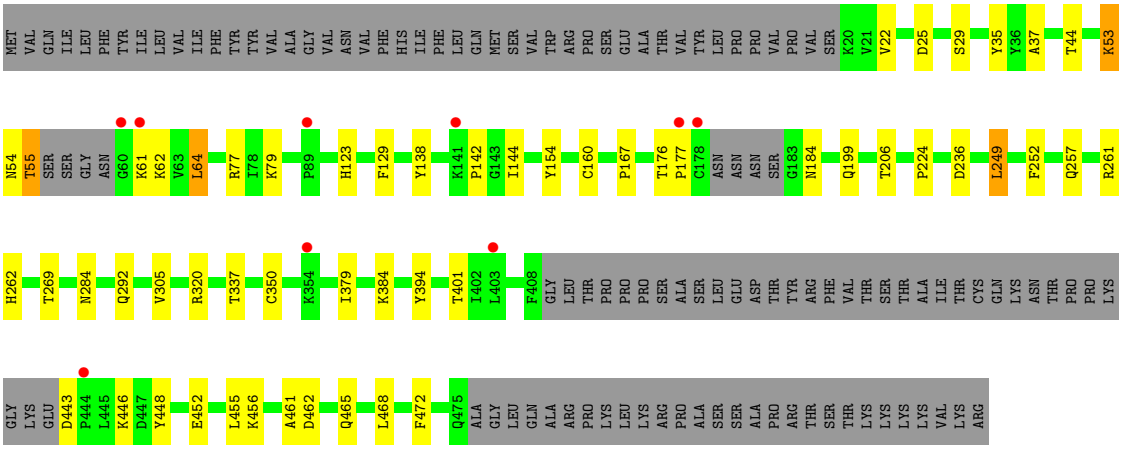
• Molecule 1: Major capsid protein L1

Chain C:



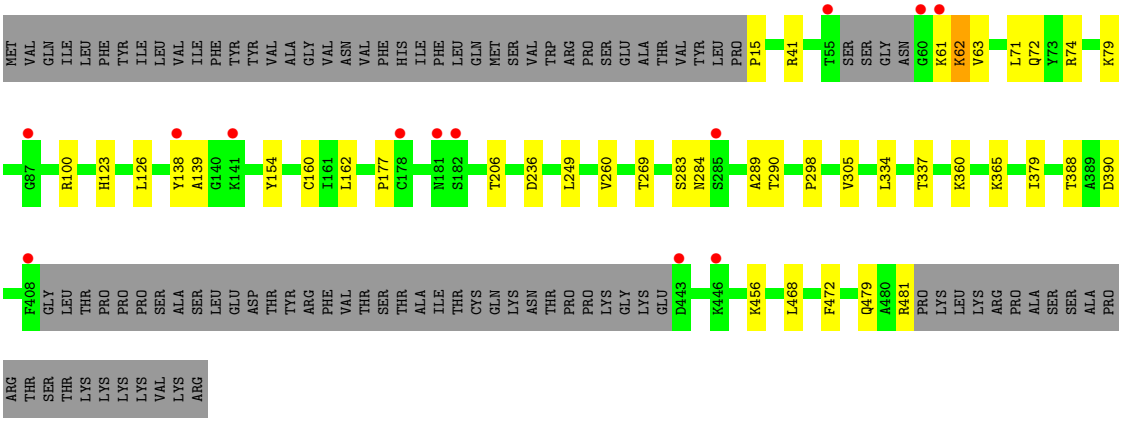
• Molecule 1: Major capsid protein L1

Chain D:



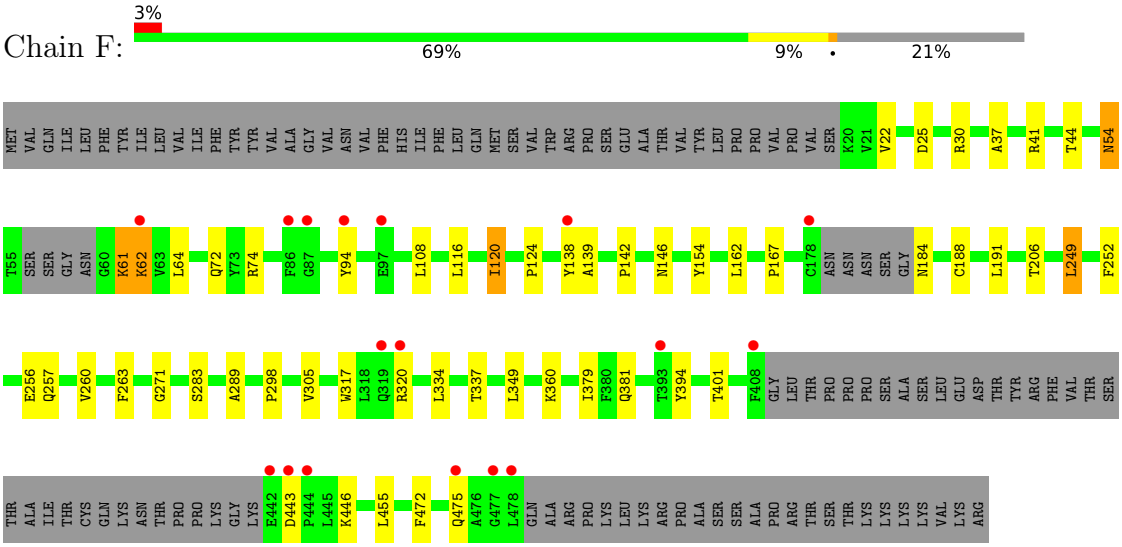
• Molecule 1: Major capsid protein L1

Chain E:



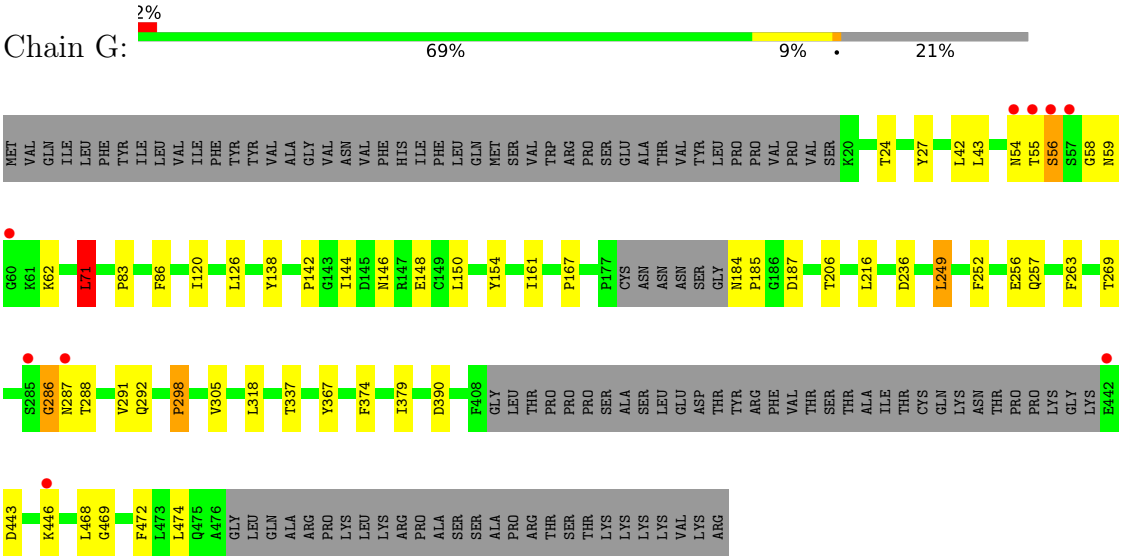
• Molecule 1: Major capsid protein L1

Chain F:



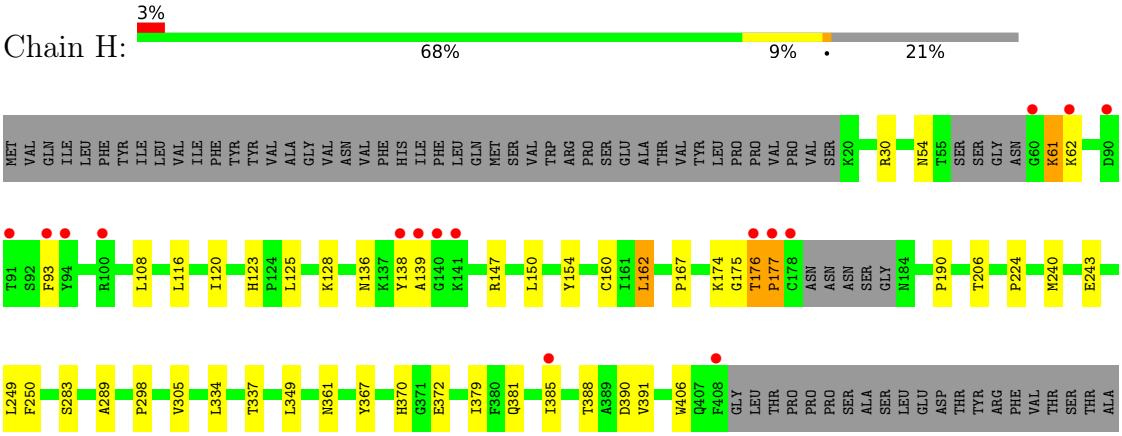
• Molecule 1: Major capsid protein L1

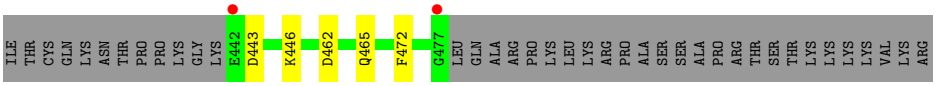
Chain G:



• Molecule 1: Major capsid protein L1

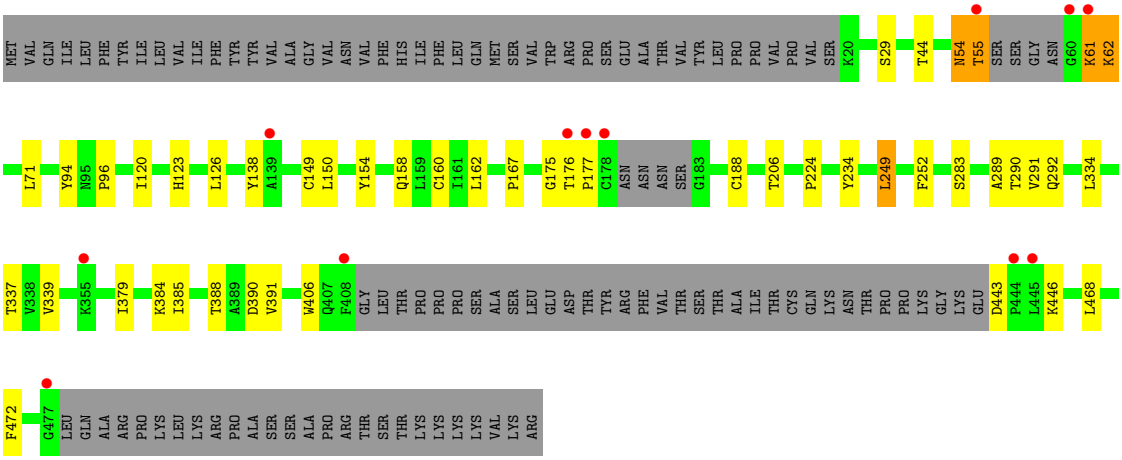
Chain H:





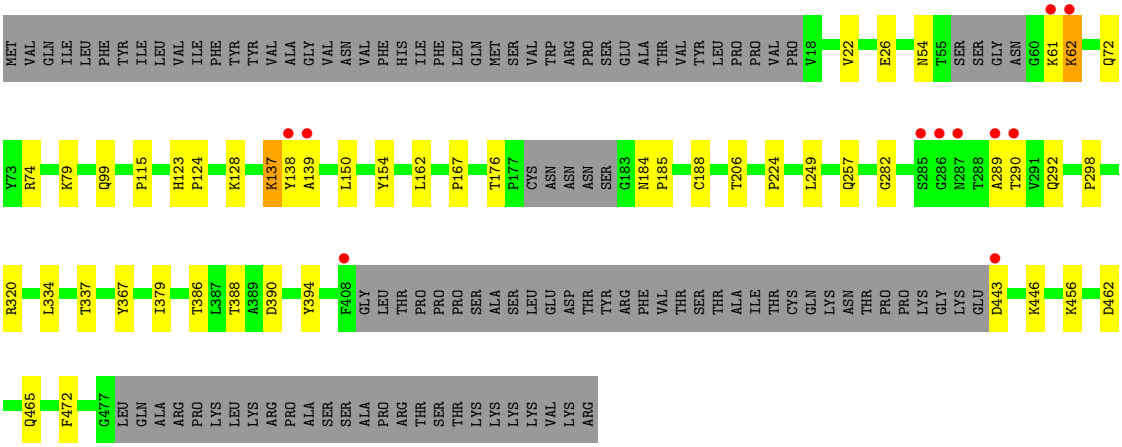
• Molecule 1: Major capsid protein L1

Chain I:



• Molecule 1: Major capsid protein L1

Chain J:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	306.77Å 105.07Å 196.90Å 90.00° 125.80° 90.00°	Depositor
Resolution (Å)	42.95 – 2.75 42.95 – 2.75	Depositor EDS
% Data completeness (in resolution range)	99.8 (42.95-2.75) 99.8 (42.95-2.75)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.20	Depositor
$< I/\sigma(I) >$ [1]	2.35 (at 2.77Å)	Xtriage
Refinement program	PHENIX (1.11.1_2575: ???)	Depositor
R, R_{free}	0.187 , 0.235 0.187 , 0.235	Depositor DCC
R_{free} test set	6697 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å^2)	43.8	Xtriage
Anisotropy	0.865	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.30 , 37.7	EDS
L-test for twinning[2]	$< \|L\| > = 0.51$, $< L^2 > = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	33255	wwPDB-VP
Average B, all atoms (Å^2)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.89% 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	A	0.36	0/3381	0.58	1/4578 (0.0%)
1	B	0.37	0/3434	0.58	2/4652 (0.0%)
1	C	0.34	0/3464	0.55	1/4693 (0.0%)
1	D	0.37	0/3376	0.60	0/4571
1	E	0.39	0/3491	0.58	0/4730
1	F	0.37	0/3404	0.56	0/4609
1	G	0.40	0/3411	0.62	4/4620 (0.1%)
1	H	0.36	0/3390	0.58	0/4590
1	I	0.38	0/3385	0.58	1/4583 (0.0%)
1	J	0.41	0/3392	0.60	1/4593 (0.0%)
All	All	0.38	0/34128	0.58	10/46219 (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	C	0	1

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	G	71	LEU	CA-CB-CG	-7.45	98.16	115.30
1	G	318	LEU	CA-CB-CG	6.70	130.71	115.30
1	B	71	LEU	CA-CB-CG	-6.12	101.23	115.30
1	J	137	LYS	CD-CE-NZ	5.93	125.33	111.70
1	G	390	ASP	CB-CG-OD1	-5.73	113.14	118.30
1	I	175	GLY	C-N-CA	5.68	135.89	121.70
1	G	298	PRO	C-N-CA	-5.41	108.18	121.70
1	C	140	GLY	C-N-CA	-5.34	108.34	121.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	298	PRO	C-N-CA	-5.11	108.92	121.70
1	B	318	LEU	CA-CB-CG	5.08	126.99	115.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	139	ALA	Peptide

5.2 Too-close contacts (i)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3293	0	3191	31	0
1	B	3344	0	3248	26	1
1	C	3377	0	3262	23	0
1	D	3291	0	3188	35	1
1	E	3400	0	3300	24	0
1	F	3316	0	3215	38	1
1	G	3322	0	3216	38	0
1	H	3305	0	3199	42	0
1	I	3300	0	3196	33	0
1	J	3307	0	3205	31	1
All	All	33255	0	32220	273	2

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 (273) 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:54:ASN:HB2	1:A:64:LEU:HD12	1.54	0.89
1:J:79:LYS:HE2	1:J:456:LYS:HE3	1.56	0.87
1:B:120:ILE:HD12	1:E:298:PRO:HB3	1.54	0.87
1:A:25:ASP:OD2	1:A:320:ARG:NH1	2.11	0.84

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:443:ASP:HB3	1:G:446:LYS:HG3	1.61	0.82
1:F:142:PRO:O	1:F:146:ASN:ND2	2.18	0.77
1:F:72:GLN:OE1	1:F:74:ARG:NH2	2.18	0.76
1:G:286:GLY:O	1:G:288:THR:N	2.19	0.75
1:J:388:THR:OG1	1:J:390:ASP:OD1	2.05	0.74
1:D:25:ASP:OD2	1:D:320:ARG:NH1	2.22	0.73
1:E:379:ILE:HD13	1:E:472:PHE:HB2	1.71	0.72
1:I:54:ASN:HD22	1:I:55:THR:H	1.35	0.72
1:I:388:THR:OG1	1:I:390:ASP:OD1	2.08	0.72
1:B:54:ASN:OD1	1:B:55:THR:N	2.23	0.71
1:I:162:LEU:HD11	1:I:334:LEU:HD12	1.71	0.71
1:E:79:LYS:HE2	1:E:456:LYS:HE3	1.73	0.70
1:C:370:HIS:NE2	1:C:372:GLU:OE2	2.25	0.69
1:B:162:LEU:HD11	1:B:334:LEU:HD12	1.74	0.69
1:H:162:LEU:HD21	1:H:334:LEU:HD12	1.74	0.67
1:C:162:LEU:HD11	1:C:334:LEU:HD12	1.77	0.67
1:C:443:ASP:HB3	1:C:446:LYS:HG3	1.77	0.66
1:H:443:ASP:HB3	1:H:446:LYS:HG3	1.78	0.65
1:F:298:PRO:HB3	1:H:120:ILE:HD12	1.78	0.65
1:B:379:ILE:HD13	1:B:472:PHE:HB2	1.79	0.65
1:F:120:ILE:HD12	1:G:298:PRO:HB3	1.79	0.64
1:G:379:ILE:HD13	1:G:472:PHE:HB2	1.80	0.64
1:D:79:LYS:HE2	1:D:456:LYS:HE3	1.79	0.63
1:C:388:THR:OG1	1:C:390:ASP:OD1	2.19	0.61
1:H:283:SER:O	1:H:289:ALA:HB2	2.01	0.60
1:D:236:ASP:OD2	1:E:41:ARG:NH2	2.35	0.60
1:F:260:VAL:HG11	1:H:120:ILE:HD11	1.83	0.60
1:I:443:ASP:HB3	1:I:446:LYS:HE3	1.82	0.60
1:B:365:LYS:HG2	1:E:269:THR:OG1	2.01	0.60
1:A:443:ASP:HB3	1:A:446:LYS:HG3	1.85	0.59
1:B:120:ILE:HD11	1:E:260:VAL:HG11	1.84	0.58
1:C:379:ILE:HD13	1:C:472:PHE:HB2	1.85	0.58
1:D:54:ASN:CA	1:D:64:LEU:HD11	2.32	0.58
1:J:22:VAL:HG22	1:J:394:TYR:CE2	2.39	0.58
1:J:184:ASN:HB2	1:J:185:PRO:HD2	1.85	0.57
1:H:370:HIS:NE2	1:H:372:GLU:OE2	2.37	0.57
1:G:146:ASN:O	1:J:137:LYS:HE2	2.04	0.57
1:G:146:ASN:HB3	1:J:137:LYS:NZ	2.20	0.57
1:D:176:THR:HG22	1:D:177:PRO:HD2	1.87	0.56
1:F:167:PRO:HG3	1:F:337:THR:OG1	2.05	0.56
1:F:298:PRO:HB3	1:H:120:ILE:CD1	2.35	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:379:ILE:HD13	1:J:472:PHE:HB2	1.88	0.56
1:I:379:ILE:HD13	1:I:472:PHE:HB2	1.89	0.55
1:E:283:SER:O	1:E:289:ALA:HB2	2.06	0.55
1:A:72:GLN:OE1	1:A:74:ARG:NH2	2.40	0.55
1:A:53:LYS:HD3	1:A:61:LYS:HG2	1.89	0.55
1:A:184:ASN:HB2	1:A:185:PRO:HD3	1.89	0.55
1:J:167:PRO:HG3	1:J:337:THR:OG1	2.06	0.55
1:D:54:ASN:OD1	1:D:55:THR:N	2.41	0.54
1:C:108:LEU:HD22	1:C:162:LEU:HD22	1.90	0.54
1:I:126:LEU:HD23	1:I:150:LEU:HD12	1.90	0.54
1:G:71:LEU:HD22	1:G:206:THR:HG22	1.90	0.54
1:B:462:ASP:OD2	1:B:465:GLN:HG2	2.09	0.53
1:D:167:PRO:HG3	1:D:337:THR:OG1	2.08	0.53
1:H:390:ASP:OD1	1:H:391:VAL:N	2.41	0.53
1:J:22:VAL:HB	1:J:26:GLU:OE1	2.07	0.53
1:C:160:CYS:HA	1:C:337:THR:O	2.09	0.53
1:F:22:VAL:HG22	1:F:394:TYR:CE2	2.44	0.53
1:G:167:PRO:HG3	1:G:337:THR:OG1	2.08	0.53
1:F:25:ASP:OD2	1:F:320:ARG:NH1	2.39	0.52
1:D:269:THR:OG1	1:E:365:LYS:HG2	2.09	0.52
1:A:298:PRO:HB3	1:C:120:ILE:HD12	1.92	0.52
1:F:108:LEU:HD22	1:F:162:LEU:HD22	1.92	0.52
1:F:188:CYS:HB2	1:H:367:TYR:CD2	2.45	0.52
1:F:120:ILE:HD11	1:G:263:PHE:CD2	2.45	0.52
1:H:298:PRO:HB3	1:I:120:ILE:CD1	2.40	0.52
1:H:379:ILE:HD13	1:H:472:PHE:HB2	1.91	0.52
1:D:138:TYR:CE2	1:D:292:GLN:HG3	2.45	0.51
1:G:146:ASN:HB3	1:J:137:LYS:HZ1	1.74	0.51
1:J:154:TYR:CD1	1:J:206:THR:HB	2.46	0.51
1:A:52:ILE:HG22	1:A:54:ASN:ND2	2.26	0.51
1:D:144:ILE:HD11	1:E:360:LYS:HG2	1.92	0.51
1:D:443:ASP:HB3	1:D:446:LYS:HE3	1.93	0.51
1:F:54:ASN:HB2	1:F:62:LYS:HB3	1.93	0.51
1:A:167:PRO:HG3	1:A:337:THR:OG1	2.11	0.50
1:H:54:ASN:HD21	1:H:61:LYS:HB2	1.76	0.50
1:D:54:ASN:CB	1:D:64:LEU:HD11	2.40	0.50
1:F:154:TYR:CD1	1:F:206:THR:HB	2.46	0.50
1:H:388:THR:OG1	1:H:390:ASP:OD1	2.29	0.50
1:C:154:TYR:CD1	1:C:206:THR:HB	2.47	0.50
1:H:136:ASN:ND2	1:I:149:CYS:HB3	2.26	0.50
1:H:162:LEU:HD13	1:H:250:PHE:HD2	1.77	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:283:SER:O	1:I:289:ALA:HB2	2.11	0.50
1:C:284:ASN:O	1:C:285:SER:OG	2.25	0.50
1:I:167:PRO:HG3	1:I:337:THR:OG1	2.12	0.50
1:F:379:ILE:HD13	1:F:472:PHE:HB2	1.95	0.49
1:A:249:LEU:HD21	1:A:252:PHE:HB3	1.94	0.49
1:C:123:HIS:HB2	1:C:224:PRO:HA	1.93	0.49
1:F:41:ARG:NH2	1:G:236:ASP:OD2	2.46	0.49
1:F:443:ASP:HB3	1:F:446:LYS:HE3	1.94	0.49
1:D:462:ASP:OD2	1:D:465:GLN:HG2	2.12	0.49
1:F:317:TRP:NE1	1:F:475:GLN:OE1	2.46	0.49
1:A:257:GLN:HA	1:C:305:VAL:O	2.13	0.49
1:G:469:GLY:HA2	1:G:472:PHE:HB3	1.94	0.48
1:C:77:ARG:NH2	1:C:443:ASP:OD2	2.43	0.48
1:H:125:LEU:HD13	1:H:147:ARG:NH2	2.28	0.48
1:I:154:TYR:CD1	1:I:206:THR:HB	2.47	0.48
1:J:22:VAL:HG22	1:J:394:TYR:CZ	2.48	0.48
1:I:54:ASN:HD22	1:I:55:THR:N	2.09	0.48
1:I:62:LYS:HE2	1:I:62:LYS:HA	1.96	0.48
1:G:54:ASN:HD21	1:G:62:LYS:HD2	1.79	0.48
1:H:154:TYR:CD1	1:H:206:THR:HB	2.48	0.48
1:H:462:ASP:OD2	1:H:465:GLN:HG2	2.13	0.48
1:I:94:TYR:CZ	1:I:96:PRO:HG3	2.48	0.48
1:C:113:GLY:HA3	1:C:373:GLU:CD	2.34	0.48
1:F:443:ASP:HB3	1:F:446:LYS:HG3	1.95	0.48
1:A:160[A]:CYS:HA	1:A:337:THR:O	2.13	0.48
1:E:379:ILE:HD13	1:E:472:PHE:CB	2.43	0.48
1:A:138:TYR:HA	1:A:139:ALA:HA	1.61	0.48
1:J:282:GLY:HA3	1:J:289:ALA:HB3	1.94	0.48
1:A:142:PRO:O	1:A:146:ASN:ND2	2.40	0.47
1:F:116:LEU:HD22	1:G:256:GLU:HB2	1.96	0.47
1:D:22:VAL:HG22	1:D:394:TYR:CE2	2.49	0.47
1:J:162:LEU:HD11	1:J:334:LEU:HD21	1.95	0.47
1:B:54:ASN:HB3	1:B:62:LYS:HB3	1.96	0.47
1:G:184:ASN:HB2	1:G:187:ASP:OD1	2.15	0.47
1:I:291:VAL:HG12	1:J:124:PRO:HB2	1.96	0.47
1:D:54:ASN:HA	1:D:64:LEU:HD21	1.97	0.47
1:F:256:GLU:HB2	1:H:116:LEU:HD22	1.97	0.47
1:G:154:TYR:CD1	1:G:206:THR:HB	2.49	0.47
1:B:167:PRO:HG3	1:B:337:THR:OG1	2.14	0.47
1:E:123:HIS:HB3	1:E:126:LEU:HB2	1.96	0.47
1:E:162:LEU:HD11	1:E:334:LEU:HD21	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:191:LEU:HD11	1:H:349:LEU:HD11	1.95	0.47
1:J:462:ASP:OD2	1:J:465:GLN:HG2	2.15	0.47
1:D:160:CYS:HA	1:D:337:THR:O	2.14	0.47
1:I:390:ASP:OD1	1:I:391:VAL:N	2.48	0.47
1:A:71:LEU:HG	1:A:154:TYR:HD2	1.80	0.46
1:D:29:SER:HB2	1:D:384:LYS:HG3	1.95	0.46
1:D:54:ASN:HA	1:D:64:LEU:HD11	1.97	0.46
1:H:138:TYR:HA	1:H:139:ALA:HA	1.70	0.46
1:D:53:LYS:HA	1:D:62:LYS:O	2.14	0.46
1:D:54:ASN:HB2	1:D:64:LEU:HD11	1.98	0.46
1:F:61:LYS:HD2	1:F:61:LYS:O	2.15	0.46
1:E:72:GLN:OE1	1:E:74:ARG:NH2	2.49	0.46
1:J:54:ASN:HD22	1:J:62:LYS:HG2	1.81	0.46
1:A:263:PHE:CD2	1:C:120:ILE:HD11	2.51	0.46
1:G:148:GLU:OE2	1:J:137:LYS:HE3	2.16	0.46
1:C:167:PRO:HG3	1:C:337:THR:OG1	2.15	0.46
1:H:61:LYS:HB2	1:H:62:LYS:H	1.57	0.46
1:A:160[B]:CYS:HA	1:A:337:THR:O	2.16	0.46
1:B:138:TYR:HA	1:B:139:ALA:HA	1.55	0.46
1:J:72:GLN:OE1	1:J:74:ARG:NH2	2.48	0.46
1:A:141:LYS:HB3	1:A:142:PRO:HD2	1.98	0.45
1:B:83:PRO:HA	1:B:86:PHE:HB2	1.98	0.45
1:B:160[A]:CYS:HA	1:B:337:THR:O	2.17	0.45
1:B:283:SER:O	1:B:289:ALA:HB2	2.16	0.45
1:F:305:VAL:O	1:G:257:GLN:HA	2.16	0.45
1:H:160:CYS:HA	1:H:337:THR:O	2.16	0.45
1:D:123:HIS:HB2	1:D:224:PRO:HA	1.98	0.45
1:G:55:THR:HA	1:G:56:SER:HA	1.65	0.45
1:H:108:LEU:HD22	1:H:162:LEU:HG	1.98	0.45
1:C:162:LEU:HD13	1:C:336:VAL:HG22	1.99	0.45
1:F:138:TYR:HA	1:F:139:ALA:HA	1.60	0.45
1:F:263:PHE:CD2	1:H:120:ILE:HD13	2.51	0.45
1:F:283:SER:O	1:F:289:ALA:HB2	2.15	0.45
1:B:25:ASP:OD2	1:B:320:ARG:NH1	2.33	0.45
1:E:71:LEU:HG	1:E:154:TYR:HD2	1.80	0.45
1:G:138:TYR:CE2	1:G:292:GLN:HG3	2.51	0.45
1:B:53:LYS:NZ	1:B:61:LYS:HE2	2.32	0.45
1:H:162:LEU:HD13	1:H:250:PHE:CD2	2.52	0.45
1:I:123:HIS:HB2	1:I:224:PRO:HA	1.98	0.45
1:I:160:CYS:HA	1:I:337:THR:O	2.16	0.45
1:G:120:ILE:CD1	1:J:298:PRO:HB3	2.47	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:136:ASN:HD22	1:I:149:CYS:HB3	1.81	0.45
1:I:249:LEU:HD21	1:I:252:PHE:HB3	1.99	0.45
1:C:129:PHE:HE1	1:C:146:ASN:HD22	1.65	0.45
1:D:379:ILE:HD12	1:D:468:LEU:HG	1.99	0.45
1:D:37:ALA:HB1	1:D:455:LEU:HD13	1.98	0.44
1:A:83:PRO:HA	1:A:86:PHE:HB2	1.98	0.44
1:A:379:ILE:HD12	1:A:468:LEU:HG	1.99	0.44
1:H:385:ILE:HD12	1:H:406:TRP:CH2	2.53	0.44
1:J:138:TYR:HA	1:J:139:ALA:HA	1.63	0.44
1:E:138:TYR:HA	1:E:139:ALA:HA	1.60	0.44
1:H:123:HIS:HB2	1:H:224:PRO:HA	1.98	0.44
1:H:240:MET:O	1:H:243:GLU:HG2	2.18	0.44
1:J:99:GLN:HG2	1:J:386:THR:HA	1.99	0.44
1:A:261:ARG:HG3	1:A:262:HIS:CD2	2.52	0.44
1:F:257:GLN:HA	1:H:305:VAL:O	2.17	0.44
1:H:385:ILE:HD12	1:H:406:TRP:HH2	1.83	0.44
1:I:29:SER:HB2	1:I:384:LYS:HG3	1.99	0.44
1:J:443:ASP:HA	1:J:446:LYS:HE3	1.99	0.44
1:G:54:ASN:HB3	1:G:58:GLY:HA3	1.98	0.44
1:I:234:TYR:CD1	1:J:115:PRO:HB3	2.52	0.44
1:I:62:LYS:HA	1:I:62:LYS:CE	2.47	0.44
1:J:184:ASN:HB2	1:J:185:PRO:CD	2.48	0.44
1:E:62:LYS:HG3	1:E:63:VAL:N	2.31	0.44
1:G:42:LEU:HB2	1:G:374:PHE:HB2	2.00	0.44
1:G:83:PRO:HA	1:G:86:PHE:HB2	1.99	0.44
1:H:176:THR:N	1:H:177:PRO:CD	2.81	0.44
1:G:161:ILE:HB	1:G:337:THR:HB	2.00	0.43
1:J:123:HIS:HB2	1:J:224:PRO:HA	2.00	0.43
1:B:41:ARG:NH2	1:E:236:ASP:OD2	2.52	0.43
1:G:291:VAL:HG12	1:G:292:GLN:O	2.18	0.43
1:B:54:ASN:ND2	1:B:62:LYS:H	2.16	0.43
1:H:167:PRO:HG3	1:H:337:THR:OG1	2.18	0.43
1:H:298:PRO:HB3	1:I:120:ILE:HD12	2.01	0.43
1:B:386:THR:HG22	1:B:388:THR:HG23	2.00	0.43
1:E:388:THR:OG1	1:E:390:ASP:OD1	2.26	0.43
1:G:184:ASN:N	1:G:185:PRO:HD3	2.34	0.43
1:G:126:LEU:HD23	1:G:150:LEU:HD12	2.01	0.43
1:G:379:ILE:CD1	1:G:468:LEU:HG	2.49	0.43
1:F:30:ARG:HB3	1:F:381:GLN:NE2	2.33	0.43
1:G:43:LEU:HD23	1:G:43:LEU:HA	1.84	0.43
1:B:184:ASN:HA	1:B:185:PRO:HD3	1.91	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:217:GLN:HA	1:D:350:CYS:HB3	2.01	0.43
1:D:35:TYR:CZ	1:D:461:ALA:HB2	2.54	0.43
1:C:54:ASN:OD1	1:C:62:LYS:HD3	2.18	0.42
1:D:249:LEU:HD21	1:D:252:PHE:HB3	2.01	0.42
1:D:379:ILE:HD13	1:D:472:PHE:HB2	2.00	0.42
1:F:124:PRO:HB2	1:G:291:VAL:HG13	2.01	0.42
1:F:188:CYS:HB2	1:H:367:TYR:CE2	2.55	0.42
1:H:174:LYS:HG2	1:H:175:GLY:O	2.19	0.42
1:A:64:LEU:HD23	1:A:64:LEU:HA	1.85	0.42
1:B:60:GLY:O	1:B:61:LYS:HD2	2.19	0.42
1:D:129:PHE:CZ	1:D:142:PRO:HA	2.55	0.42
1:A:191:LEU:HD23	1:A:191:LEU:HA	1.91	0.42
1:D:199:GLN:NE2	1:D:448:TYR:HD1	2.18	0.42
1:G:305:VAL:O	1:J:257:GLN:HA	2.19	0.42
1:I:379:ILE:HD11	1:I:468:LEU:O	2.19	0.42
1:A:399:ASP:OD1	1:A:401:THR:HG22	2.19	0.42
1:D:154:TYR:CD1	1:D:206:THR:HB	2.55	0.42
1:D:261:ARG:HG3	1:D:262:HIS:CD2	2.54	0.42
1:F:146:ASN:HD22	1:H:361:ASN:HD21	1.68	0.42
1:H:30:ARG:HB3	1:H:381:GLN:NE2	2.35	0.42
1:I:61:LYS:HB3	1:I:61:LYS:HE2	1.85	0.42
1:B:120:ILE:CD1	1:E:298:PRO:HB3	2.36	0.42
1:E:154:TYR:CD1	1:E:206:THR:HB	2.55	0.42
1:F:37:ALA:HB1	1:F:455:LEU:HD13	2.02	0.42
1:F:249:LEU:HD21	1:F:252:PHE:HB3	2.02	0.42
1:B:379:ILE:CD1	1:B:472:PHE:HB2	2.49	0.41
1:C:257:GLN:HA	1:D:305:VAL:O	2.19	0.41
1:E:100:ARG:HD3	1:E:100:ARG:HA	1.93	0.41
1:E:379:ILE:HD12	1:E:468:LEU:HG	2.02	0.41
1:H:128:LYS:HD2	1:H:150:LEU:HD11	2.02	0.41
1:B:123:HIS:HB2	1:B:224:PRO:HA	2.01	0.41
1:C:379:ILE:HD11	1:C:468:LEU:O	2.20	0.41
1:G:24:THR:HA	1:G:27:TYR:CE1	2.55	0.41
1:G:379:ILE:HD12	1:G:468:LEU:HG	2.01	0.41
1:I:138:TYR:CE2	1:I:292:GLN:HG3	2.54	0.41
1:D:64:LEU:CD1	1:D:64:LEU:N	2.84	0.41
1:I:71:LEU:HG	1:I:154:TYR:HD2	1.85	0.41
1:A:71:LEU:HD13	1:A:71:LEU:HA	1.93	0.41
1:J:138:TYR:CZ	1:J:292:GLN:HG3	2.56	0.41
1:A:86:PHE:O	1:A:88:PHE:N	2.53	0.41
1:A:283:SER:O	1:A:289:ALA:HB2	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:271:GLY:HA3	1:H:367:TYR:CE2	2.56	0.41
1:G:249:LEU:HD21	1:G:252:PHE:HB3	2.03	0.41
1:G:367:TYR:CD2	1:J:188:CYS:HB2	2.56	0.41
1:A:62:LYS:HB2	1:A:62:LYS:HE3	1.57	0.41
1:A:367:TYR:CD2	1:B:188:CYS:HB2	2.56	0.41
1:F:349:LEU:HD22	1:G:216:LEU:HB3	2.02	0.41
1:I:158:GLN:HA	1:I:339:VAL:O	2.20	0.41
1:I:188:CYS:HB2	1:J:367:TYR:CD2	2.56	0.41
1:A:154:TYR:CD1	1:A:206:THR:HB	2.56	0.41
1:A:379:ILE:HD13	1:A:472:PHE:HB2	2.02	0.41
1:F:360:LYS:HG2	1:G:144:ILE:HD12	2.02	0.40
1:I:123:HIS:HB3	1:I:126:LEU:HB2	2.01	0.40
1:B:108:LEU:HD13	1:B:378:PHE:CE2	2.55	0.40
1:F:162:LEU:HD11	1:F:334:LEU:HD21	2.02	0.40
1:H:177:PRO:HD2	1:H:190:PRO:CG	2.52	0.40
1:I:385:ILE:HD12	1:I:406:TRP:HH2	1.85	0.40
1:B:53:LYS:HZ3	1:B:61:LYS:HG3	1.86	0.40
1:E:160[B]:CYS:HA	1:E:337:THR:O	2.21	0.40
1:J:128:LYS:HD2	1:J:150:LEU:HD11	2.04	0.40
1:C:80:LEU:HD11	1:C:334:LEU:HD22	2.04	0.40
1:D:77:ARG:HD3	1:D:452:GLU:OE1	2.20	0.40
1:D:257:GLN:HA	1:E:305:VAL:O	2.22	0.40
1:I:176:THR:HG22	1:I:177:PRO:HD2	2.04	0.40

All (2) 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)
1:D:284:ASN:ND2	1:F:94:TYR:O[3_555]	1.95	0.25
1:B:92:SER:OG	1:J:320:ARG:NH2[2_556]	2.16	0.04

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	406/529 (77%)	391 (96%)	15 (4%)	0	100	100
1	B	414/529 (78%)	401 (97%)	12 (3%)	1 (0%)	47	69
1	C	423/529 (80%)	404 (96%)	19 (4%)	0	100	100
1	D	406/529 (77%)	392 (97%)	14 (3%)	0	100	100
1	E	424/529 (80%)	405 (96%)	18 (4%)	1 (0%)	47	69
1	F	410/529 (78%)	395 (96%)	15 (4%)	0	100	100
1	G	413/529 (78%)	393 (95%)	17 (4%)	3 (1%)	22	39
1	H	408/529 (77%)	391 (96%)	15 (4%)	2 (0%)	29	47
1	I	408/529 (77%)	392 (96%)	16 (4%)	0	100	100
1	J	409/529 (77%)	394 (96%)	15 (4%)	0	100	100
All	All	4121/5290 (78%)	3958 (96%)	156 (4%)	7 (0%)	47	69

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	G	287	ASN
1	H	177	PRO
1	B	142	PRO
1	E	177	PRO
1	H	176	THR
1	G	142	PRO
1	G	286	GLY

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	368/469 (78%)	362 (98%)	6 (2%)	62	77
1	B	374/469 (80%)	371 (99%)	3 (1%)	81	88
1	C	377/469 (80%)	369 (98%)	8 (2%)	53	71

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	D	367/469 (78%)	359 (98%)	8 (2%)	52	70
1	E	380/469 (81%)	372 (98%)	8 (2%)	53	71
1	F	370/469 (79%)	361 (98%)	9 (2%)	49	68
1	G	371/469 (79%)	365 (98%)	6 (2%)	62	77
1	H	368/469 (78%)	364 (99%)	4 (1%)	73	84
1	I	367/469 (78%)	360 (98%)	7 (2%)	57	73
1	J	368/469 (78%)	363 (99%)	5 (1%)	67	79
All	All	3710/4690 (79%)	3646 (98%)	64 (2%)	60	76

All (64) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	44	THR
1	A	61	LYS
1	A	62	LYS
1	A	93	PHE
1	A	290	THR
1	A	394	TYR
1	B	61	LYS
1	B	176	THR
1	B	184	ASN
1	C	53	LYS
1	C	54	ASN
1	C	55	THR
1	C	120	ILE
1	C	176	THR
1	C	182	SER
1	C	249	LEU
1	C	290	THR
1	D	44	THR
1	D	53	LYS
1	D	55	THR
1	D	61	LYS
1	D	64	LEU
1	D	184	ASN
1	D	249	LEU
1	D	401	THR
1	E	15	PRO
1	E	61	LYS
1	E	62	LYS

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Mol	Chain	Res	Type
1	E	249	LEU
1	E	284	ASN
1	E	290	THR
1	E	479	GLN
1	E	481	ARG
1	F	44	THR
1	F	54	ASN
1	F	61	LYS
1	F	62	LYS
1	F	64	LEU
1	F	120	ILE
1	F	184	ASN
1	F	249	LEU
1	F	401	THR
1	G	56	SER
1	G	59	ASN
1	G	71	LEU
1	G	249	LEU
1	G	269	THR
1	G	474	LEU
1	H	61	LYS
1	H	93	PHE
1	H	162	LEU
1	H	249	LEU
1	I	44	THR
1	I	54	ASN
1	I	55	THR
1	I	61	LYS
1	I	62	LYS
1	I	249	LEU
1	I	290	THR
1	J	61	LYS
1	J	62	LYS
1	J	176	THR
1	J	249	LEU
1	J	290	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	54	ASN
1	A	146	ASN

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Mol	Chain	Res	Type
1	B	184	ASN
1	D	184	ASN
1	D	239	GLN
1	F	54	ASN
1	H	156	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	A	413/529 (78%)	0.05	22 (5%)	26	31	35, 55, 113, 200	0
1	B	421/529 (79%)	-0.13	17 (4%)	38	45	27, 47, 104, 163	0
1	C	427/529 (80%)	0.06	17 (3%)	38	45	33, 53, 112, 170	0
1	D	414/529 (78%)	-0.20	9 (2%)	62	70	27, 47, 96, 195	0
1	E	429/529 (81%)	-0.16	13 (3%)	50	59	26, 39, 98, 154	0
1	F	417/529 (78%)	-0.00	17 (4%)	37	44	28, 51, 111, 166	0
1	G	418/529 (79%)	-0.33	9 (2%)	62	70	21, 38, 97, 154	0
1	H	416/529 (78%)	-0.11	18 (4%)	35	42	29, 53, 105, 151	0
1	I	416/529 (78%)	-0.22	12 (2%)	51	61	24, 43, 93, 169	0
1	J	417/529 (78%)	-0.26	11 (2%)	56	65	20, 36, 92, 150	0
All	All	4188/5290 (79%)	-0.13	145 (3%)	44	52	20, 47, 103, 200	0

All (145) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	I	408	PHE	8.8
1	H	477	GLY	7.3
1	A	178	CYS	7.1
1	I	178	CYS	7.1
1	F	478	LEU	6.8
1	G	442	GLU	6.5
1	B	141	LYS	6.4
1	G	55	THR	6.4
1	B	478	LEU	6.3
1	A	88	PHE	6.0
1	A	177	PRO	5.9
1	A	90	ASP	5.7
1	A	89	PRO	5.2

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Mol	Chain	Res	Type	RSRZ
1	G	56	SER	5.2
1	C	478	LEU	5.2
1	D	61	LYS	5.1
1	H	442	GLU	5.1
1	A	61	LYS	5.0
1	C	442	GLU	4.9
1	F	138	TYR	4.9
1	A	93	PHE	4.9
1	I	176	THR	4.6
1	J	138	TYR	4.5
1	E	446	LYS	4.5
1	J	408	PHE	4.5
1	D	60	GLY	4.5
1	B	442	GLU	4.3
1	E	55	THR	4.2
1	E	408	PHE	4.2
1	A	442	GLU	4.2
1	B	285	SER	4.1
1	I	177	PRO	4.0
1	E	182	SER	4.0
1	E	60	GLY	4.0
1	A	475	GLN	3.9
1	H	60	GLY	3.9
1	C	61	LYS	3.8
1	F	178	CYS	3.7
1	B	89	PRO	3.6
1	D	89	PRO	3.6
1	E	61	LYS	3.5
1	F	442	GLU	3.5
1	I	60	GLY	3.4
1	F	87	GLY	3.4
1	H	140	GLY	3.4
1	C	89	PRO	3.4
1	J	287	ASN	3.3
1	H	93	PHE	3.3
1	C	140	GLY	3.3
1	B	443	ASP	3.2
1	D	177	PRO	3.2
1	F	86	PHE	3.1
1	H	138	TYR	3.1
1	B	61	LYS	3.1
1	I	55	THR	3.1

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Mol	Chain	Res	Type	RSRZ
1	F	477	GLY	3.1
1	H	177	PRO	3.1
1	A	62	LYS	3.1
1	I	477	GLY	3.0
1	H	408	PHE	3.0
1	D	178	CYS	3.0
1	F	408	PHE	3.0
1	H	178	CYS	3.0
1	C	87	GLY	2.9
1	G	60	GLY	2.9
1	B	408	PHE	2.9
1	B	138	TYR	2.9
1	B	477	GLY	2.8
1	D	444	PRO	2.8
1	I	444	PRO	2.8
1	A	476	ALA	2.8
1	J	443	ASP	2.8
1	G	285	SER	2.8
1	E	138	TYR	2.8
1	E	178	CYS	2.8
1	J	61	LYS	2.7
1	A	87	GLY	2.7
1	B	93	PHE	2.7
1	C	354	LYS	2.7
1	A	176	THR	2.6
1	G	57	SER	2.6
1	I	61	LYS	2.6
1	C	62	LYS	2.6
1	F	475	GLN	2.5
1	H	141	LYS	2.5
1	A	54	ASN	2.5
1	G	287	ASN	2.5
1	J	285	SER	2.5
1	E	181	ASN	2.5
1	H	139	ALA	2.5
1	E	141	LYS	2.5
1	G	446	LYS	2.5
1	C	63	VAL	2.4
1	J	286	GLY	2.4
1	I	445	LEU	2.4
1	C	138	TYR	2.3
1	G	54	ASN	2.3

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Mol	Chain	Res	Type	RSRZ
1	F	393	THR	2.3
1	B	445	LEU	2.3
1	F	444	PRO	2.3
1	B	283	SER	2.3
1	E	285	SER	2.3
1	H	385	ILE	2.3
1	H	100	ARG	2.3
1	D	141	LYS	2.3
1	F	443	ASP	2.3
1	A	285	SER	2.3
1	B	55	THR	2.3
1	J	290	THR	2.2
1	A	286	GLY	2.2
1	F	62	LYS	2.2
1	A	138	TYR	2.2
1	I	139	ALA	2.2
1	H	176	THR	2.2
1	F	319	GLN	2.2
1	J	139	ALA	2.2
1	A	144	ILE	2.2
1	B	281	GLN	2.2
1	H	91	THR	2.2
1	F	320	ARG	2.2
1	H	94	TYR	2.2
1	C	55	THR	2.2
1	C	90	ASP	2.1
1	C	408	PHE	2.1
1	D	354	LYS	2.1
1	I	355	LYS	2.1
1	C	362	GLU	2.1
1	F	97	GLU	2.1
1	F	94	TYR	2.1
1	J	62	LYS	2.1
1	A	389	ALA	2.1
1	J	289	ALA	2.1
1	H	90	ASP	2.1
1	B	97	GLU	2.1
1	C	443	ASP	2.1
1	H	62	LYS	2.1
1	E	87	GLY	2.1
1	C	353	VAL	2.0
1	B	91	THR	2.0

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Mol	Chain	Res	Type	RSRZ
1	E	443	ASP	2.0
1	A	408	PHE	2.0
1	A	319	GLN	2.0
1	D	403	LEU	2.0
1	C	88	PHE	2.0
1	A	391	VAL	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.