



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 02:41 AM UTC

PDB ID : 5HYJ / pdb_00005hyj
Title : 1E6 TCR in Complex with HLA-A02 carrying AQWGPDPAAA
Authors : Rizkallah, P.J.; Bulek, A.M.; Cole, D.K.; Sewell, A.K.
Deposited on : 2016-02-01
Resolution : 3.06 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

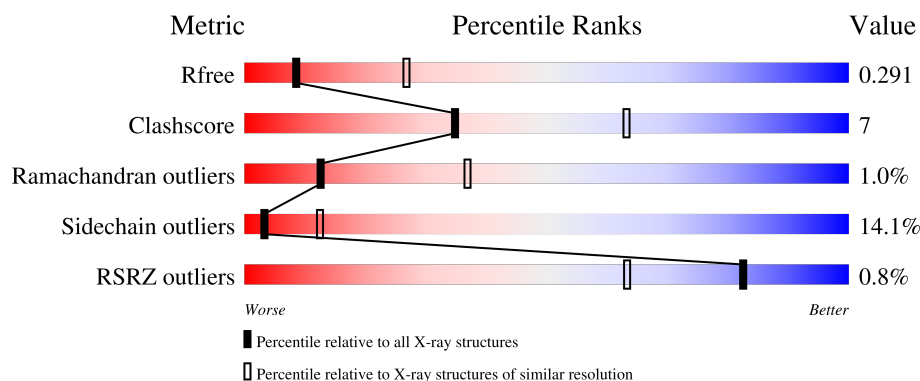
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.06 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}	180053	2469 (3.10-3.02)
Clashscore	190562	2569 (3.10-3.02)
Ramachandran outliers	187476	2424 (3.10-3.02)
Sidechain outliers	187428	2423 (3.10-3.02)
RSRZ outliers	180081	2469 (3.10-3.02)

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 $\leq 5\%$. 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	276	74% 22% • •
1	F	276	71% 25% •
2	B	100	69% 26% 5%
2	G	100	71% 23% 6%
3	C	10	60% 30% 10%

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Mol	Chain	Length	Quality of chain
3	H	10	60%40%
4	D	193	% 59%30%10%
4	I	193	3% 58%34%8%
5	E	246	76%20%
5	J	246	76%22%

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 13392 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 HLA class I histocompatibility antigen, A-2 alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	276	Total	C	N	O	S	0	0	0
			2254	1408	410	427	9			
1	F	276	Total	C	N	O	S	0	0	0
			2254	1408	410	427	9			

- Molecule 2 is a protein called Beta-2-microglobulin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	100	Total	C	N	O	S	0	0	0
			837	533	141	159	4			
2	G	100	Total	C	N	O	S	0	0	0
			837	533	141	159	4			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	0	MET	-	initiating methionine	UNP P61769
G	0	MET	-	initiating methionine	UNP P61769

- Molecule 3 is a protein called ALA-GLN-TRP-GLY-PRO-ASP-PRO-ALA-ALA-ALA.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	10	Total	C	N	O	0	0	0
			70	44	12	14			
3	H	10	Total	C	N	O	0	0	0
			70	44	12	14			

- Molecule 4 is a protein called Human T-cell Receptor, Class I, Light alpha Chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	193	Total	C	N	O	S	0	0	0
			1520	950	253	307	10			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	I	193	Total	C	N	O	S	0	0	0
			1520	950	253	307	10			

- Molecule 5 is a protein called Human T-cell Receptor, Class I, Heavy beta Chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	E	246	Total	C	N	O	S	0	0	0
			1974	1249	341	374	10			
5	J	246	Total	C	N	O	S	0	0	0
			1974	1249	341	374	10			

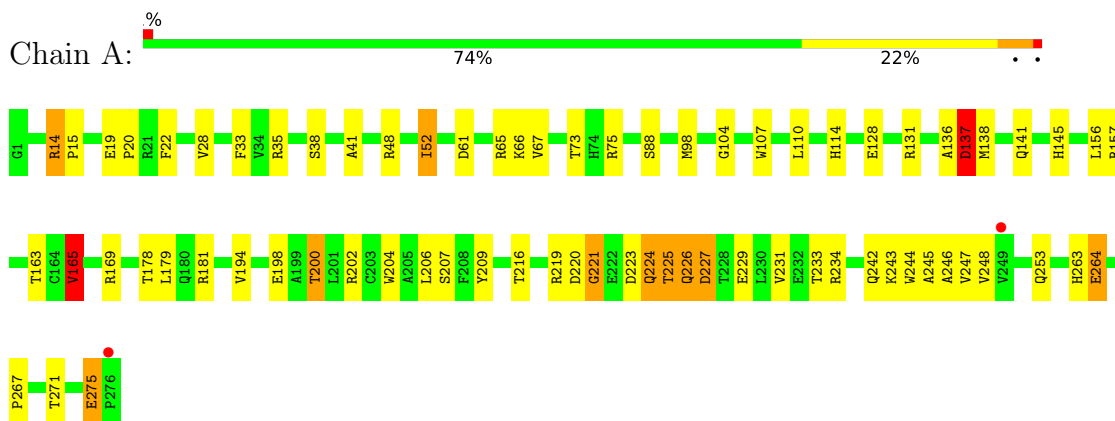
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	12	Total	O	0	0
			12	12		
6	B	2	Total	O	0	0
			2	2		
6	D	14	Total	O	0	0
			14	14		
6	E	13	Total	O	0	0
			13	13		
6	F	20	Total	O	0	0
			20	20		
6	G	2	Total	O	0	0
			2	2		
6	I	11	Total	O	0	0
			11	11		
6	J	8	Total	O	0	0
			8	8		

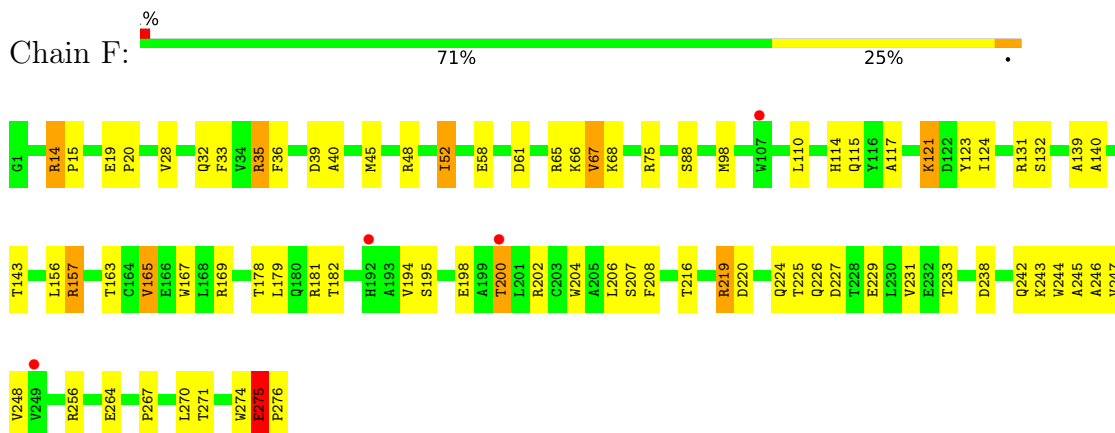
3 Residue-property plots

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.

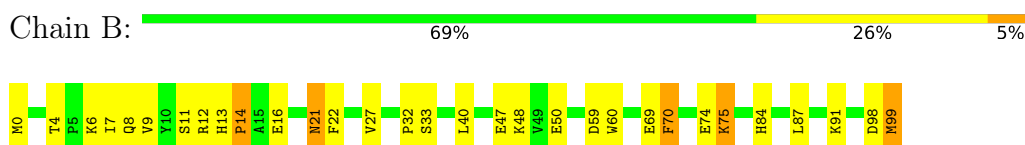
- Molecule 1: HLA class I histocompatibility antigen, A-2 alpha chain



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- Molecule 2: Beta-2-microglobulin



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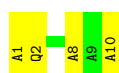




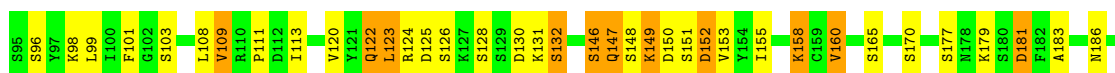
- Molecule 3: ALA-GLN-TRP-GLY-PRO-ASP-PRO-ALA-ALA-ALA



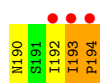
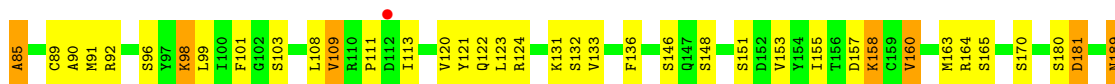
- Molecule 3: ALA-GLN-TRP-GLY-PRO-ASP-PRO-ALA-ALA-ALA



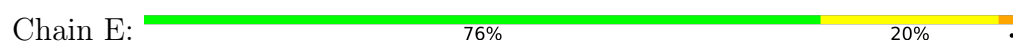
- Molecule 4: Human T-cell Receptor, Class I, Light alpha Chain

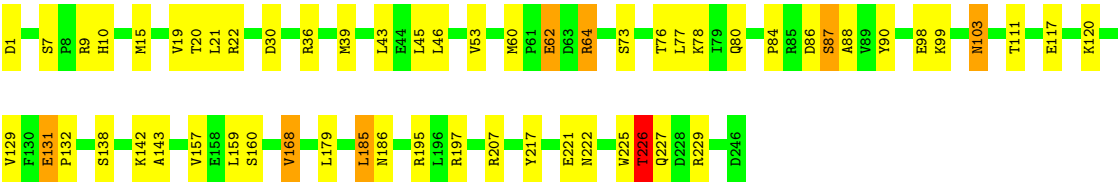


- Molecule 4: Human T-cell Receptor, Class I, Light alpha Chain

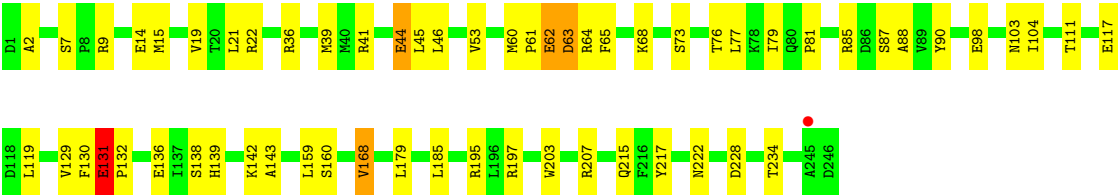


- Molecule 5: Human T-cell Receptor, Class I, Heavy beta Chain





• Molecule 5: Human T-cell Receptor, Class I, Heavy beta Chain



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	41.44Å 97.31Å 121.37Å 97.29° 97.66° 92.73°	Depositor
Resolution (Å)	38.63 – 3.06 38.63 – 3.06	Depositor EDS
% Data completeness (in resolution range)	97.2 (38.63-3.06) 97.2 (38.63-3.06)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.36 (at 3.06Å)	Xtriage
Refinement program	REFMAC 5.8.0135	Depositor
R, R_{free}	0.212 , 0.293 0.213 , 0.291	Depositor DCC
R_{free} test set	1726 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	56.0	Xtriage
Anisotropy	0.285	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 63.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	13392	wwPDB-VP
Average B, all atoms (Å ²)	73.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.76% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ 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	1.19	6/2320 (0.3%)	1.24	7/3149 (0.2%)
1	F	1.23	5/2320 (0.2%)	1.30	12/3149 (0.4%)
2	B	1.15	2/860 (0.2%)	1.23	4/1162 (0.3%)
2	G	1.19	0/860	1.32	6/1162 (0.5%)
3	C	1.41	0/73	1.37	0/100
3	H	1.47	0/73	1.48	0/100
4	D	1.19	7/1553 (0.5%)	1.26	11/2100 (0.5%)
4	I	1.14	5/1553 (0.3%)	1.19	4/2100 (0.2%)
5	E	1.10	1/2029 (0.0%)	1.18	4/2759 (0.1%)
5	J	0.97	0/2029	1.16	9/2759 (0.3%)
All	All	1.15	26/13670 (0.2%)	1.23	57/18540 (0.3%)

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	A	0	2
2	G	0	1
3	C	0	1
3	H	0	1
4	D	0	2
4	I	0	2
5	E	0	2
All	All	0	11

All (26) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	194	PRO	C-O	10.35	1.44	1.23
4	I	194	PRO	C-O	9.91	1.43	1.23
4	D	128	SER	N-CA	6.34	1.53	1.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	E	185	LEU	CA-C	6.20	1.60	1.52
4	I	193	ILE	N-CA	6.16	1.54	1.46
4	I	193	ILE	CA-C	6.06	1.59	1.52
4	D	193	ILE	CA-C	6.01	1.59	1.52
1	F	68	LYS	CA-C	-5.97	1.45	1.52
1	F	225	THR	CA-C	5.95	1.60	1.52
1	A	224	GLN	CA-C	5.90	1.60	1.52
1	A	209	TYR	CA-C	5.83	1.57	1.52
2	B	59	ASP	N-CA	5.66	1.53	1.46
1	F	40	ALA	N-CA	-5.56	1.38	1.45
2	B	47	GLU	CA-CB	5.52	1.62	1.53
4	D	193	ILE	CA-CB	5.51	1.61	1.54
1	F	35	ARG	CA-C	-5.44	1.46	1.52
4	D	192	ILE	CA-CB	5.41	1.61	1.53
1	A	227	ASP	CA-CB	5.32	1.57	1.53
1	A	98	MET	CA-C	-5.30	1.46	1.52
4	D	150	ASP	CA-C	5.19	1.60	1.52
1	F	98	MET	C-O	-5.19	1.18	1.23
4	I	57	GLU	CA-C	5.17	1.58	1.52
1	A	226	GLN	CA-C	5.09	1.59	1.52
1	A	107	TRP	N-CA	5.08	1.52	1.46
4	I	192	ILE	CA-C	5.07	1.58	1.53
4	D	130	ASP	CA-CB	5.06	1.60	1.53

All (57) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	G	59	ASP	N-CA-C	-13.22	93.47	110.53
1	F	124	ILE	N-CA-C	12.17	126.91	109.51
5	E	229	ARG	N-CA-C	-11.70	95.43	110.53
5	J	62	GLU	N-CA-C	9.90	121.76	110.97
1	F	224	GLN	N-CA-C	8.76	122.47	109.24
4	D	54	GLY	N-CA-C	-8.03	99.11	111.00
1	F	123	TYR	N-CA-C	-7.95	96.28	109.80
4	D	148	SER	N-CA-C	-7.92	93.93	110.80
4	I	181	ASP	N-CA-C	-7.81	97.42	109.52
1	F	61	ASP	N-CA-C	-7.78	102.39	111.03
4	D	130	ASP	CB-CA-C	7.75	123.00	110.29
2	G	47	GLU	N-CA-C	7.63	121.07	111.69
4	D	123	LEU	N-CA-C	7.32	120.40	108.26
2	B	70	PHE	N-CA-C	7.07	120.10	110.35
1	F	165	VAL	N-CA-C	6.95	117.65	110.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	61	ASP	N-CA-C	-6.74	103.55	111.03
1	A	41	ALA	N-CA-C	6.68	119.12	111.11
1	A	145	HIS	CB-CA-C	6.59	121.74	110.79
5	J	131	GLU	CA-C-N	6.32	126.34	119.89
5	J	131	GLU	C-N-CA	6.32	126.34	119.89
1	A	165	VAL	N-CA-C	6.17	116.84	110.36
4	I	58	ASP	N-CA-C	-6.17	96.88	107.61
4	D	152	ASP	CA-C-N	6.10	130.75	123.19
4	D	152	ASP	C-N-CA	6.10	130.75	123.19
2	G	59	ASP	CA-C-O	-6.08	114.94	121.94
1	A	138	MET	N-CA-C	6.07	118.42	111.02
5	J	63	ASP	N-CA-C	-6.06	101.81	111.02
5	J	44	GLU	CA-CB-CG	5.96	126.01	114.10
2	B	69	GLU	N-CA-C	5.86	118.76	108.56
4	D	18	ILE	N-CA-C	-5.85	99.74	108.46
1	F	39	ASP	N-CA-C	5.78	120.19	113.20
5	E	168	VAL	CB-CA-C	-5.69	102.96	110.98
1	F	219	ARG	CB-CG-CD	5.68	124.35	111.30
1	A	48	ARG	NE-CZ-NH2	5.67	124.30	119.20
5	J	2	ALA	N-CA-C	-5.65	102.12	110.48
2	B	7	ILE	CB-CA-C	-5.50	102.99	110.42
4	I	11	LEU	CB-CA-C	5.47	118.78	109.75
5	J	215	GLN	N-CA-C	-5.44	101.11	109.76
1	A	163	THR	N-CA-C	5.42	116.87	111.07
2	G	32	PRO	CA-C-N	5.40	128.27	120.38
2	G	32	PRO	C-N-CA	5.40	128.27	120.38
1	F	163	THR	N-CA-C	5.38	116.83	111.07
5	J	228	ASP	N-CA-CB	5.32	116.31	110.35
5	J	168	VAL	CB-CA-C	-5.29	103.96	111.21
5	E	62	GLU	N-CA-C	5.28	116.72	111.07
2	B	60	TRP	N-CA-C	5.26	122.01	110.80
4	D	146	SER	N-CA-C	-5.26	100.11	109.06
4	D	91	MET	CB-CG-SD	5.24	128.42	112.70
1	F	132	SER	N-CA-C	5.19	117.29	109.41
4	D	58	ASP	N-CA-C	-5.13	98.69	107.61
1	F	208	PHE	N-CA-C	5.09	118.11	109.76
2	G	9	VAL	CB-CA-C	-5.09	103.55	110.42
4	D	131	LYS	N-CA-C	-5.08	103.69	110.55
4	I	157	ASP	N-CA-C	-5.08	102.86	110.23
1	F	195	SER	N-CA-C	5.04	114.90	108.24
1	F	157	ARG	CG-CD-NE	5.01	123.02	112.00
5	E	103	ASN	N-CA-C	5.00	121.45	110.80

There are no chirality outliers.

All (11) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	137	ASP	Peptide
1	A	226	GLN	Peptide
3	C	9	ALA	Peptide
4	D	149	LYS	Peptide
4	D	158	LYS	Peptide
5	E	1	ASP	Peptide
5	E	226	THR	Peptide
2	G	0	MET	Peptide
3	H	8	ALA	Peptide
4	I	158	LYS	Peptide
4	I	180	SER	Peptide

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	2254	0	2103	30	0
1	F	2254	0	2103	26	0
2	B	837	0	803	11	0
2	G	837	0	803	11	0
3	C	70	0	61	4	0
3	H	70	0	61	3	0
4	D	1520	0	1447	38	0
4	I	1520	0	1447	33	0
5	E	1974	0	1887	30	0
5	J	1974	0	1887	29	0
6	A	12	0	0	0	0
6	B	2	0	0	0	0
6	D	14	0	0	1	0
6	E	13	0	0	1	0
6	F	20	0	0	1	0
6	G	2	0	0	1	0
6	I	11	0	0	0	0
6	J	8	0	0	0	0
All	All	13392	0	12602	193	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 7.

All (193) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:41:ARG:NH2	5:E:157:VAL:O	2.06	0.87
4:I:34:MET:HE2	5:J:103:ASN:HB2	1.60	0.84
2:B:21:ASN:HD22	2:B:22:PHE:H	1.27	0.82
4:D:183:ALA:HB3	4:D:186:ASN:HD21	1.46	0.80
5:E:15:MET:CE	5:E:117:GLU:HA	2.15	0.77
2:G:21:ASN:HD22	2:G:22:PHE:H	1.32	0.77
2:G:57:SER:O	2:G:59:ASP:O	2.01	0.76
4:D:123:LEU:HD11	5:E:131:GLU:O	1.87	0.74
5:J:15:MET:CE	5:J:117:GLU:HA	2.17	0.73
4:I:98:LYS:HB2	5:J:45:LEU:HD22	1.71	0.73
4:D:48:MET:SD	4:D:58:ASP:HB2	2.30	0.71
1:F:143:THR:HG23	3:H:10:ALA:HA	1.72	0.71
4:D:147:GLN:HB3	4:D:155:ILE:HD13	1.73	0.70
5:E:88:ALA:HB3	5:E:90:TYR:CE1	2.27	0.70
5:E:120:LYS:HD3	5:E:227:GLN:HE22	1.57	0.70
5:J:15:MET:HE2	5:J:117:GLU:HA	1.75	0.68
2:B:9:VAL:O	2:B:99:MET:HE1	1.94	0.68
5:E:15:MET:HE2	5:E:117:GLU:HA	1.75	0.68
5:J:88:ALA:HB3	5:J:90:TYR:CE1	2.34	0.63
4:D:111:PRO:CG	4:D:160:VAL:HG11	2.28	0.63
1:F:19:GLU:CD	1:F:75:ARG:HD2	2.24	0.62
1:F:194:VAL:HG21	1:F:248:VAL:HG11	1.81	0.62
4:I:111:PRO:CG	4:I:160:VAL:HG11	2.29	0.62
5:E:84:PRO:O	5:E:87:SER:OG	2.18	0.61
5:J:61:PRO:HD2	5:J:65:PHE:CD2	2.35	0.61
1:A:194:VAL:HG21	1:A:248:VAL:HG11	1.82	0.60
5:E:132:PRO:HG3	5:E:143:ALA:HB1	1.84	0.60
1:F:45:MET:HE2	3:H:2:GLN:HE22	1.66	0.60
5:J:132:PRO:HG3	5:J:143:ALA:HB1	1.83	0.60
4:I:55:ASN:N	4:I:55:ASN:OD1	2.36	0.59
4:D:123:LEU:HD11	5:E:132:PRO:C	2.29	0.58
4:D:125:ASP:OD1	4:D:126:SER:N	2.36	0.58
5:J:65:PHE:CE1	5:J:79:ILE:HG23	2.38	0.57
1:A:227:ASP:O	1:A:247:VAL:HG23	2.04	0.57
2:B:9:VAL:O	2:B:99:MET:CE	2.52	0.57
4:D:60:ARG:NH2	4:D:83:ASP:OD2	2.37	0.57
4:D:34:MET:HE2	5:E:103:ASN:HB2	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:202:ARG:HG3	1:F:246:ALA:HB2	1.88	0.56
4:D:72:ILE:HD11	4:D:91:MET:CE	2.35	0.56
1:F:227:ASP:O	1:F:247:VAL:HG23	2.06	0.56
2:G:21:ASN:HD22	2:G:22:PHE:N	2.03	0.56
2:B:21:ASN:HD22	2:B:22:PHE:N	1.98	0.56
4:D:72:ILE:HD11	4:D:91:MET:HE3	1.86	0.55
1:A:202:ARG:HG3	1:A:246:ALA:HB2	1.88	0.55
3:C:4:GLY:HA2	3:C:5:PRO:C	2.32	0.54
1:A:14:ARG:HD3	1:A:19:GLU:O	2.08	0.54
1:A:65:ARG:O	1:A:66:LYS:C	2.51	0.54
4:I:89:CYS:SG	4:I:89:CYS:O	2.66	0.54
2:G:58:LYS:HA	6:G:102:HOH:O	2.07	0.54
1:F:65:ARG:O	1:F:66:LYS:C	2.47	0.54
4:I:46:LEU:HD23	5:J:104:ILE:HG12	1.89	0.53
4:D:90:ALA:HB2	4:D:101:PHE:CE1	2.44	0.53
4:D:111:PRO:HG2	4:D:160:VAL:HG11	1.90	0.53
4:D:64:GLN:HE22	4:I:18:ILE:HG13	1.73	0.53
5:E:15:MET:HE3	5:E:117:GLU:HA	1.87	0.52
5:E:20:THR:HG23	5:E:78:LYS:HE3	1.90	0.52
5:J:132:PRO:CB	5:J:143:ALA:HB1	2.39	0.52
5:E:132:PRO:CB	5:E:143:ALA:HB1	2.40	0.52
5:J:64:ARG:HB3	5:J:65:PHE:HD1	1.74	0.52
4:D:89:CYS:O	4:D:89:CYS:SG	2.68	0.52
5:J:64:ARG:HB3	5:J:65:PHE:CD1	2.44	0.52
1:A:220:ASP:O	1:A:221:GLY:O	2.28	0.52
1:F:233:THR:OG1	1:F:243:LYS:HD2	2.10	0.52
1:F:182:THR:HG22	1:F:264:GLU:OE1	2.10	0.51
4:D:81:PRO:HA	4:D:109:VAL:HG22	1.92	0.51
4:I:121:TYR:CE2	5:J:136:GLU:HG3	2.45	0.51
4:I:81:PRO:HA	4:I:109:VAL:HG22	1.92	0.51
4:I:60:ARG:NH2	4:I:83:ASP:OD2	2.42	0.51
4:I:111:PRO:HG2	4:I:160:VAL:HG11	1.92	0.51
1:A:263:HIS:ND1	1:A:264:GLU:O	2.44	0.51
1:F:274:TRP:NE1	6:F:301:HOH:O	2.36	0.51
4:I:193:ILE:HB	4:I:194:PRO:CD	2.42	0.50
5:J:132:PRO:CG	5:J:143:ALA:HB1	2.41	0.50
1:A:233:THR:OG1	1:A:243:LYS:HD2	2.12	0.50
5:E:132:PRO:CG	5:E:143:ALA:HB1	2.41	0.50
4:D:123:LEU:CD1	5:E:131:GLU:O	2.59	0.49
5:J:60:MET:HE1	5:J:77:LEU:HD13	1.93	0.49
4:D:151:SER:HB3	4:D:153:VAL:HG23	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:E:99:LYS:NZ	6:E:301:HOH:O	2.46	0.49
4:D:146:SER:O	4:D:191:SER:OG	2.26	0.49
4:I:24:THR:HG22	4:I:71:TYR:HD1	1.76	0.49
4:I:189:ASN:HD22	4:I:190:ASN:N	2.10	0.48
4:I:90:ALA:HB2	4:I:101:PHE:CE1	2.48	0.48
1:F:36:PHE:CD2	1:F:67:VAL:HG11	2.48	0.48
4:D:66:ASP:OD1	4:D:66:ASP:C	2.56	0.48
5:J:46:LEU:HB3	5:J:60:MET:HE2	1.96	0.48
1:F:194:VAL:HG23	1:F:200:THR:OG1	2.13	0.48
2:B:98:ASP:O	2:B:99:MET:HB2	2.14	0.47
4:D:77:ARG:NH2	6:D:201:HOH:O	2.47	0.47
2:B:74:GLU:HB3	2:B:75:LYS:HD3	1.95	0.47
1:F:33:PHE:CD2	1:F:52:ILE:HD13	2.49	0.47
4:I:58:ASP:O	4:I:59:GLY:C	2.58	0.47
1:F:114:HIS:CD2	1:F:156:LEU:HD11	2.50	0.47
1:A:194:VAL:HG23	1:A:200:THR:OG1	2.14	0.47
1:F:229:GLU:O	1:F:245:ALA:HA	2.14	0.47
5:J:15:MET:HE3	5:J:117:GLU:HA	1.94	0.47
1:A:33:PHE:CD2	1:A:52:ILE:HD13	2.50	0.46
4:D:24:THR:HG22	4:D:71:TYR:HD1	1.79	0.46
1:A:223:ASP:O	1:A:224:GLN:HG2	2.15	0.46
4:I:133:VAL:HG11	5:J:130:PHE:CG	2.51	0.46
5:E:120:LYS:CD	5:E:227:GLN:HE22	2.27	0.46
4:D:90:ALA:C	4:D:91:MET:HG2	2.40	0.46
5:J:62:GLU:N	5:J:62:GLU:OE2	2.49	0.46
5:E:21:LEU:O	5:E:76:THR:HA	2.16	0.46
5:E:225:TRP:O	5:E:226:THR:HG23	2.16	0.46
4:I:66:ASP:C	4:I:66:ASP:OD1	2.59	0.46
1:A:28:VAL:HG11	1:A:179:LEU:HD13	1.98	0.46
1:A:229:GLU:O	1:A:245:ALA:HA	2.15	0.46
4:D:101:PHE:CE2	5:E:43:LEU:HD13	2.50	0.46
4:D:111:PRO:HG3	4:D:160:VAL:HG11	1.98	0.45
5:E:46:LEU:HB3	5:E:60:MET:HE2	1.97	0.45
5:E:60:MET:HE1	5:E:77:LEU:HD13	1.96	0.45
2:G:13:HIS:H	2:G:21:ASN:HD21	1.65	0.45
4:I:27:ASN:HD21	4:I:29:ALA:HB3	1.81	0.45
4:I:34:MET:CE	5:J:103:ASN:HB2	2.40	0.45
1:A:206:LEU:HD23	1:A:242:GLN:HG2	1.99	0.45
1:A:104:GLY:CA	1:A:110:LEU:HD12	2.47	0.45
4:D:193:ILE:HB	4:D:194:PRO:CD	2.46	0.45
5:E:159:LEU:HD23	5:E:160:SER:N	2.32	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:219:ARG:O	1:A:220:ASP:C	2.58	0.44
5:J:21:LEU:CD2	5:J:111:THR:HG21	2.48	0.44
2:G:84:HIS:HB3	2:G:87:LEU:HD12	1.99	0.44
5:E:9:ARG:NE	5:E:9:ARG:HA	2.33	0.44
1:A:165:VAL:CG1	1:A:169:ARG:NH1	2.81	0.44
1:F:238:ASP:HB3	2:G:12:ARG:HD3	2.00	0.44
5:E:142:LYS:HD3	5:E:197:ARG:HD3	2.00	0.44
5:J:21:LEU:O	5:J:76:THR:HA	2.18	0.44
5:J:159:LEU:HD23	5:J:160:SER:N	2.33	0.44
5:J:64:ARG:HG2	5:J:81:PRO:O	2.17	0.44
5:E:21:LEU:CD2	5:E:111:THR:HG21	2.48	0.44
4:I:25:TYR:HB3	4:I:91:MET:HE1	1.99	0.43
1:A:14:ARG:O	1:A:15:PRO:C	2.61	0.43
1:F:28:VAL:HG11	1:F:179:LEU:HD13	2.00	0.43
2:G:98:ASP:N	2:G:98:ASP:OD1	2.51	0.43
2:G:32:PRO:O	2:G:84:HIS:HE1	2.02	0.43
1:A:19:GLU:OE1	1:A:75:ARG:HD2	2.18	0.43
4:D:18:ILE:HA	4:D:76:ILE:O	2.18	0.43
1:A:263:HIS:CE1	1:A:264:GLU:O	2.72	0.43
4:I:153:VAL:HG12	4:I:155:ILE:HD11	2.01	0.43
4:I:4:VAL:HG23	4:I:23:CYS:HB3	2.00	0.43
5:E:64:ARG:NH1	5:E:86:ASP:OD2	2.52	0.43
1:F:139:ALA:O	1:F:140:ALA:C	2.61	0.43
1:A:204:TRP:HA	1:A:244:TRP:HB3	2.01	0.43
4:D:153:VAL:HG12	4:D:155:ILE:HD11	2.01	0.43
4:I:25:TYR:CB	4:I:91:MET:HE1	2.49	0.43
4:D:122:GLN:OE1	4:D:132:SER:OG	2.25	0.42
4:I:33:PHE:CG	4:I:65:VAL:HG21	2.55	0.42
4:I:50:THR:O	4:I:50:THR:HG23	2.19	0.42
4:I:111:PRO:HG3	4:I:160:VAL:HG11	1.99	0.42
1:A:224:GLN:C	1:A:225:THR:HG22	2.45	0.42
2:G:99:MET:HE3	2:G:99:MET:HB3	1.87	0.42
4:I:84:SER:O	4:I:85:ALA:HB2	2.19	0.42
1:A:136:ALA:O	1:A:137:ASP:C	2.62	0.42
1:A:156:LEU:HG	3:C:3:TRP:CZ2	2.55	0.42
2:B:13:HIS:H	2:B:21:ASN:HD21	1.66	0.42
4:I:136:PHE:CZ	4:I:193:ILE:HG22	2.54	0.42
1:F:14:ARG:HD2	1:F:20:PRO:HA	2.02	0.42
2:G:37:VAL:HG22	2:G:82:VAL:HG22	2.02	0.42
1:F:117:ALA:HB1	1:F:121:LYS:O	2.20	0.42
4:D:98:LYS:HB2	5:E:45:LEU:HD22	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:I:18:ILE:HA	4:I:76:ILE:O	2.19	0.41
4:D:27:ASN:HD21	4:D:29:ALA:HB3	1.84	0.41
1:F:219:ARG:O	1:F:220:ASP:C	2.61	0.41
5:J:46:LEU:CD1	5:J:77:LEU:HD21	2.50	0.41
1:A:114:HIS:CD2	1:A:156:LEU:HD11	2.55	0.41
4:D:25:TYR:CG	4:D:91:MET:HE1	2.56	0.41
4:D:147:GLN:OE1	4:D:149:LYS:HB2	2.21	0.41
1:F:167:TRP:CG	3:H:1:ALA:HB3	2.56	0.41
4:I:78:ASP:O	4:I:79:SER:C	2.63	0.41
5:J:14:GLU:HB3	5:J:119:LEU:HG	2.01	0.41
4:D:51:TYR:O	4:D:51:TYR:CD2	2.74	0.41
4:I:14:PRO:HD2	4:I:17:ALA:HB2	2.01	0.41
5:E:10:HIS:HB2	5:E:217:TYR:CG	2.56	0.41
1:F:206:LEU:HD23	1:F:242:GLN:HG2	2.02	0.41
5:J:131:GLU:HB3	5:J:203:TRP:CH2	2.56	0.41
1:A:234:ARG:HD3	2:B:8:GLN:OE1	2.21	0.41
5:E:46:LEU:CD1	5:E:77:LEU:HD21	2.51	0.41
1:F:14:ARG:O	1:F:15:PRO:C	2.64	0.41
1:A:73:THR:HG23	3:C:9:ALA:HB2	2.03	0.41
2:B:12:ARG:HE	2:B:12:ARG:HB2	1.67	0.41
1:A:66:LYS:HD3	3:C:4:GLY:HA3	2.03	0.41
4:D:4:VAL:HG23	4:D:23:CYS:HB3	2.02	0.41
4:I:4:VAL:HG11	4:I:91:MET:CE	2.51	0.41
2:B:32:PRO:O	2:B:84:HIS:HE1	2.03	0.40
5:J:142:LYS:HD3	5:J:197:ARG:HD3	2.02	0.40
4:D:78:ASP:O	4:D:79:SER:C	2.64	0.40
5:J:217:TYR:HA	5:J:234:THR:HG23	2.03	0.40
2:B:84:HIS:HB3	2:B:87:LEU:HD12	2.03	0.40
4:D:181:ASP:OD1	4:D:181:ASP:N	2.51	0.40
1:F:204:TRP:HA	1:F:244:TRP:HB3	2.03	0.40
1:A:14:ARG:CD	1:A:20:PRO:HA	2.51	0.40
1:A:22:PHE:HB3	1:A:38:SER:HB3	2.04	0.40
1:F:275:GLU:O	1:F:276:PRO:C	2.65	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	274/276 (99%)	245 (89%)	25 (9%)	4 (2%)	8	28
1	F	274/276 (99%)	253 (92%)	19 (7%)	2 (1%)	18	45
2	B	98/100 (98%)	91 (93%)	6 (6%)	1 (1%)	12	37
2	G	98/100 (98%)	96 (98%)	1 (1%)	1 (1%)	12	37
3	C	8/10 (80%)	8 (100%)	0	0	100	100
3	H	8/10 (80%)	7 (88%)	1 (12%)	0	100	100
4	D	191/193 (99%)	166 (87%)	22 (12%)	3 (2%)	7	26
4	I	191/193 (99%)	167 (87%)	19 (10%)	5 (3%)	4	17
5	E	244/246 (99%)	226 (93%)	17 (7%)	1 (0%)	30	57
5	J	244/246 (99%)	225 (92%)	19 (8%)	0	100	100
All	All	1630/1650 (99%)	1484 (91%)	129 (8%)	17 (1%)	12	37

All (17) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	221	GLY
1	A	137	ASP
4	D	181	ASP
4	I	158	LYS
4	D	158	LYS
1	F	275	GLU
4	I	85	ALA
1	A	275	GLU
1	F	267	PRO
4	I	52	SER
4	I	148	SER
5	E	64	ARG
2	B	14	PRO
4	D	59	GLY

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Mol	Chain	Res	Type
1	A	267	PRO
4	I	59	GLY
2	G	14	PRO

5.3.2 Protein sidechains ⓘ

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	232/232 (100%)	210 (90%)	22 (10%)	8	27
1	F	232/232 (100%)	205 (88%)	27 (12%)	5	19
2	B	95/95 (100%)	79 (83%)	16 (17%)	2	8
2	G	95/95 (100%)	80 (84%)	15 (16%)	2	10
3	C	5/5 (100%)	5 (100%)	0	100	100
3	H	5/5 (100%)	5 (100%)	0	100	100
4	D	174/174 (100%)	138 (79%)	36 (21%)	1	4
4	I	174/174 (100%)	136 (78%)	38 (22%)	1	4
5	E	216/216 (100%)	192 (89%)	24 (11%)	6	21
5	J	216/216 (100%)	191 (88%)	25 (12%)	5	19
All	All	1444/1444 (100%)	1241 (86%)	203 (14%)	3	12

All (203) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	14	ARG
1	A	35	ARG
1	A	52	ILE
1	A	67	VAL
1	A	88	SER
1	A	128	GLU
1	A	131	ARG
1	A	141	GLN
1	A	157	ARG
1	A	165	VAL

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Mol	Chain	Res	Type
1	A	178	THR
1	A	181	ARG
1	A	198	GLU
1	A	200	THR
1	A	207	SER
1	A	216	THR
1	A	225	THR
1	A	231	VAL
1	A	253	GLN
1	A	264	GLU
1	A	271	THR
1	A	275	GLU
2	B	0	MET
2	B	4	THR
2	B	6	LYS
2	B	11	SER
2	B	14	PRO
2	B	16	GLU
2	B	21	ASN
2	B	27	VAL
2	B	33	SER
2	B	40	LEU
2	B	48	LYS
2	B	50	GLU
2	B	70	PHE
2	B	75	LYS
2	B	91	LYS
2	B	99	MET
4	D	2	LYS
4	D	4	VAL
4	D	11	LEU
4	D	19	VAL
4	D	26	SER
4	D	31	GLN
4	D	34	MET
4	D	41	ARG
4	D	45	GLU
4	D	47	LEU
4	D	48	MET
4	D	51	TYR
4	D	55	ASN
4	D	58	ASP

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Mol	Chain	Res	Type
4	D	65	VAL
4	D	68	SER
4	D	91	MET
4	D	94	ASP
4	D	96	SER
4	D	99	LEU
4	D	103	SER
4	D	108	LEU
4	D	109	VAL
4	D	113	ILE
4	D	120	VAL
4	D	122	GLN
4	D	124	ARG
4	D	132	SER
4	D	147	GLN
4	D	152	ASP
4	D	160	VAL
4	D	165	SER
4	D	170	SER
4	D	177	SER
4	D	179	LYS
4	D	192	ILE
5	E	7	SER
5	E	19	VAL
5	E	22	ARG
5	E	30	ASP
5	E	36	ARG
5	E	39	MET
5	E	53	VAL
5	E	62	GLU
5	E	73	SER
5	E	80	GLN
5	E	87	SER
5	E	98	GLU
5	E	129	VAL
5	E	131	GLU
5	E	138	SER
5	E	168	VAL
5	E	179	LEU
5	E	185	LEU
5	E	186	ASN
5	E	195	ARG

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Mol	Chain	Res	Type
5	E	207	ARG
5	E	221	GLU
5	E	222	ASN
5	E	226	THR
1	F	14	ARG
1	F	32	GLN
1	F	35	ARG
1	F	48	ARG
1	F	52	ILE
1	F	58	GLU
1	F	67	VAL
1	F	88	SER
1	F	110	LEU
1	F	115	GLN
1	F	121	LYS
1	F	131	ARG
1	F	157	ARG
1	F	165	VAL
1	F	169	ARG
1	F	178	THR
1	F	181	ARG
1	F	198	GLU
1	F	200	THR
1	F	207	SER
1	F	216	THR
1	F	226	GLN
1	F	231	VAL
1	F	256	ARG
1	F	270	LEU
1	F	271	THR
1	F	275	GLU
2	G	4	THR
2	G	6	LYS
2	G	11	SER
2	G	14	PRO
2	G	16	GLU
2	G	21	ASN
2	G	27	VAL
2	G	33	SER
2	G	34	ASP
2	G	40	LEU
2	G	47	GLU

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Mol	Chain	Res	Type
2	G	50	GLU
2	G	70	PHE
2	G	75	LYS
2	G	98	ASP
4	I	2	LYS
4	I	4	VAL
4	I	11	LEU
4	I	19	VAL
4	I	31	GLN
4	I	34	MET
4	I	41	ARG
4	I	45	GLU
4	I	47	LEU
4	I	48	MET
4	I	51	TYR
4	I	55	ASN
4	I	56	LYS
4	I	65	VAL
4	I	68	SER
4	I	92	ARG
4	I	96	SER
4	I	98	LYS
4	I	99	LEU
4	I	103	SER
4	I	108	LEU
4	I	109	VAL
4	I	113	ILE
4	I	120	VAL
4	I	122	GLN
4	I	123	LEU
4	I	124	ARG
4	I	131	LYS
4	I	132	SER
4	I	146	SER
4	I	151	SER
4	I	160	VAL
4	I	163	MET
4	I	164	ARG
4	I	165	SER
4	I	170	SER
4	I	181	ASP
4	I	189	ASN

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Mol	Chain	Res	Type
5	J	7	SER
5	J	9	ARG
5	J	19	VAL
5	J	22	ARG
5	J	36	ARG
5	J	39	MET
5	J	41	ARG
5	J	44	GLU
5	J	53	VAL
5	J	63	ASP
5	J	68	LYS
5	J	73	SER
5	J	85	ARG
5	J	87	SER
5	J	98	GLU
5	J	129	VAL
5	J	131	GLU
5	J	138	SER
5	J	139	HIS
5	J	168	VAL
5	J	179	LEU
5	J	185	LEU
5	J	195	ARG
5	J	207	ARG
5	J	222	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (49) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	54	GLN
1	A	86	ASN
1	A	141	GLN
1	A	174	ASN
1	A	180	GLN
1	A	262	GLN
2	B	13	HIS
2	B	21	ASN
2	B	24	ASN
2	B	84	HIS
2	B	89	GLN
4	D	27	ASN
4	D	64	GLN

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Mol	Chain	Res	Type
4	D	144	ASN
4	D	171	ASN
4	D	186	ASN
4	D	190	ASN
5	E	18	GLN
5	E	29	HIS
5	E	51	ASN
5	E	121	ASN
5	E	141	GLN
5	E	177	GLN
5	E	222	ASN
5	E	235	GLN
1	F	32	GLN
1	F	54	GLN
1	F	87	GLN
1	F	141	GLN
1	F	174	ASN
1	F	180	GLN
2	G	13	HIS
2	G	21	ASN
2	G	24	ASN
2	G	84	HIS
2	G	89	GLN
3	H	2	GLN
4	I	27	ASN
4	I	64	GLN
4	I	144	ASN
4	I	171	ASN
4	I	178	ASN
4	I	189	ASN
5	J	29	HIS
5	J	51	ASN
5	J	177	GLN
5	J	213	GLN
5	J	222	ASN
5	J	235	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 oligosaccharides 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, 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	276/276 (100%)	-0.16	2 (0%) 84 66	45, 65, 108, 138	0
1	F	276/276 (100%)	-0.19	4 (1%) 73 51	30, 57, 123, 169	0
2	B	100/100 (100%)	-0.23	0 100 100	47, 60, 88, 109	0
2	G	100/100 (100%)	-0.37	0 100 100	38, 55, 76, 90	0
3	C	10/10 (100%)	-0.38	0 100 100	47, 54, 56, 61	0
3	H	10/10 (100%)	-0.10	0 100 100	42, 50, 59, 60	0
4	D	193/193 (100%)	0.05	2 (1%) 79 59	46, 77, 130, 155	0
4	I	193/193 (100%)	0.20	5 (2%) 57 34	51, 86, 144, 173	0
5	E	246/246 (100%)	-0.16	0 100 100	34, 63, 120, 143	0
5	J	246/246 (100%)	0.01	1 (0%) 88 75	49, 80, 117, 151	0
All	All	1650/1650 (100%)	-0.09	14 (0%) 82 63	30, 67, 126, 173	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
4	D	194	PRO	3.1
5	J	245	ALA	2.8
4	I	194	PRO	2.5
4	I	53	SER	2.4
1	A	276	PRO	2.4
4	I	192	ILE	2.2
4	D	2	LYS	2.2
1	F	192	HIS	2.2
1	A	249	VAL	2.1
4	I	112	ASP	2.1
1	F	200	THR	2.1
1	F	107	TRP	2.1
4	I	193	ILE	2.0

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Mol	Chain	Res	Type	RSRZ
1	F	249	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 oligosaccharides 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.