



Full wwPDB X-ray Structure Validation Report ⓘ

May 14, 2020 – 11:29 pm BST

PDB ID : 5HYJ
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 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 : 2.11
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) : 2.11

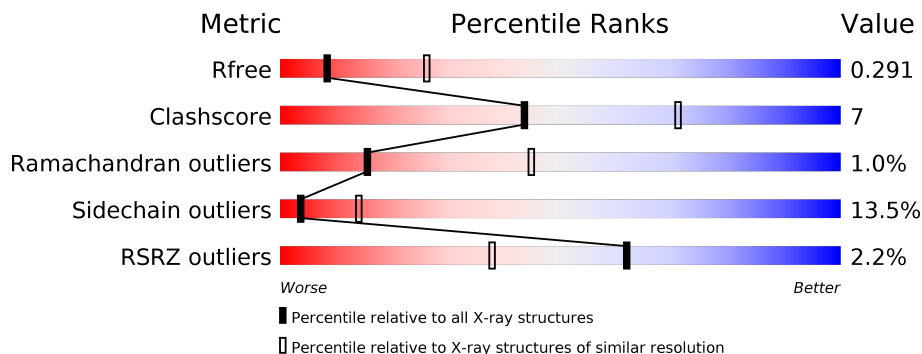
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.06 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}	130704	1754 (3.10-3.02)
Clashscore	141614	1864 (3.10-3.02)
Ramachandran outliers	138981	1794 (3.10-3.02)
Sidechain outliers	138945	1793 (3.10-3.02)
RSRZ outliers	127900	1713 (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 on 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	 75% 21% .
1	F	276	 77% 18% .
2	B	100	 76% 21% .
2	G	100	 76% 17% 7%
3	C	10	 60% 30% 10%
3	H	10	 70% 30%

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Mol	Chain	Length	Quality of chain
4	D	193	
4	I	193	
5	E	246	
5	J	246	

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
			Total	C	N	O	S			
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
			Total	C	N	O	S			
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
			Total	C	N	O			
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
			Total	C	N	O	S			
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			

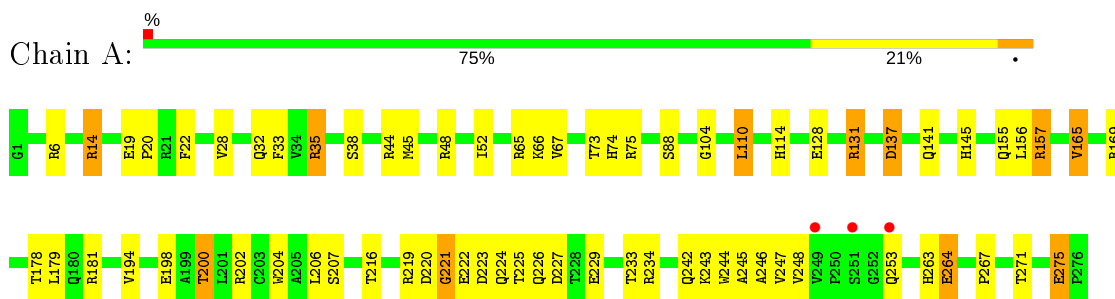
- Molecule 6 is water.

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		

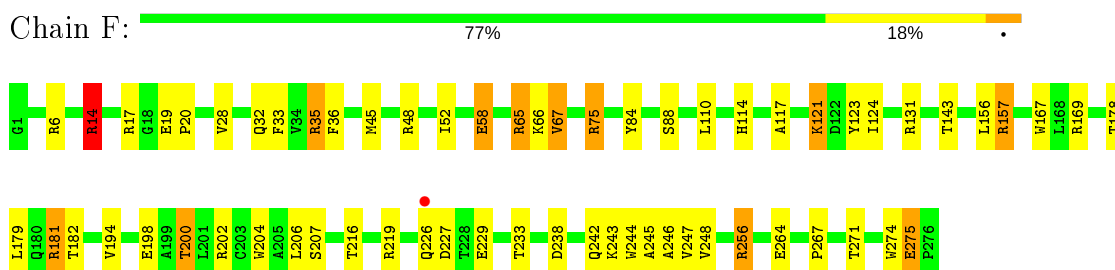
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

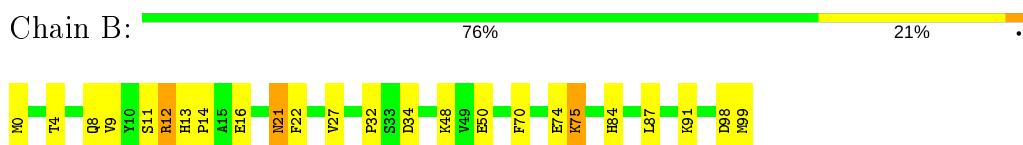
- Molecule 1: HLA class I histocompatibility antigen, A-2 alpha chain



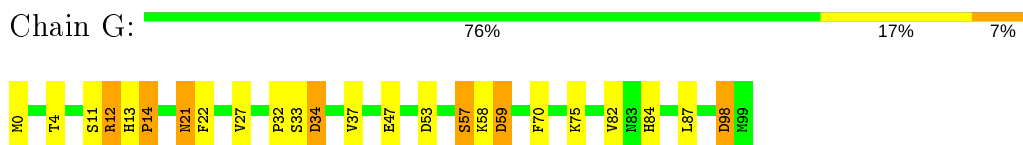
- Molecule 1: HLA class I histocompatibility antigen, A-2 alpha chain



- Molecule 2: Beta-2-microglobulin

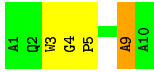


- Molecule 2: Beta-2-microglobulin



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

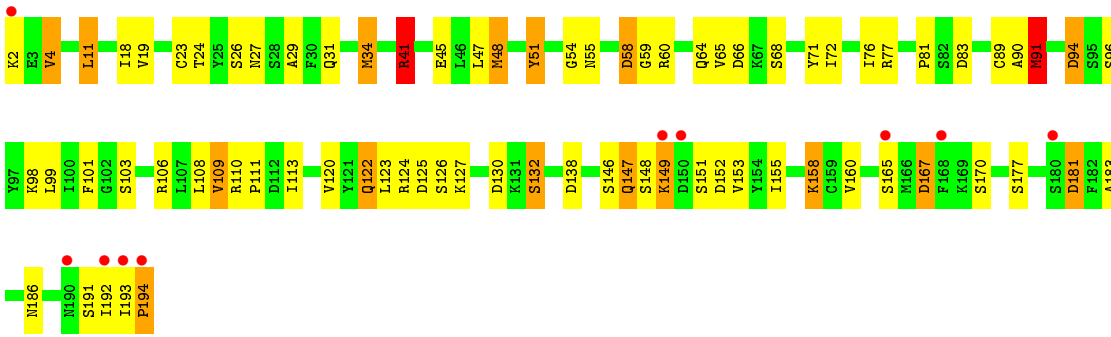




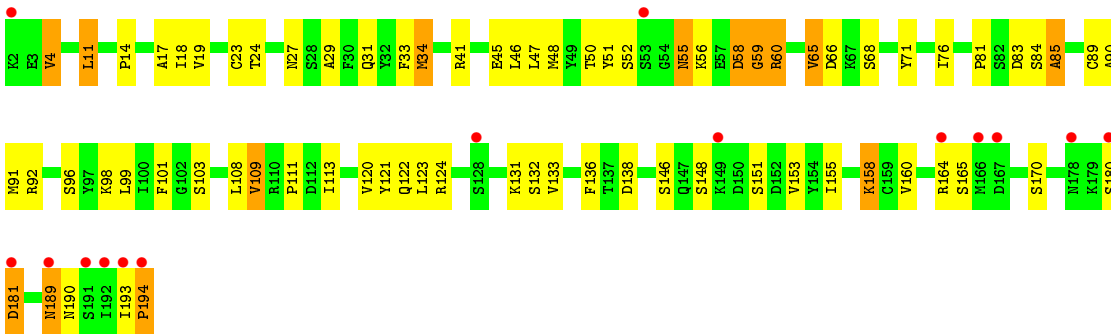
- Molecule 3: ALA-GLN-TRP-GLY-PRO-ASP-PRO-ALA-ALA-ALA



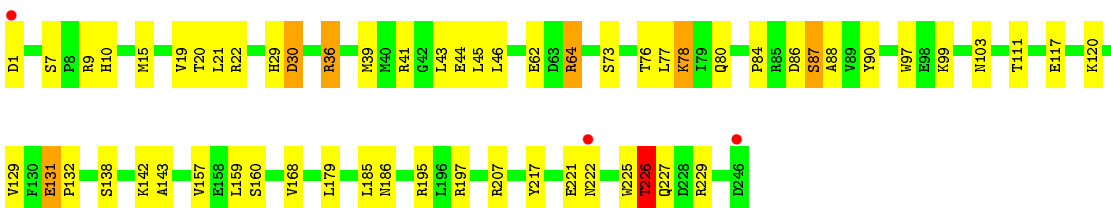
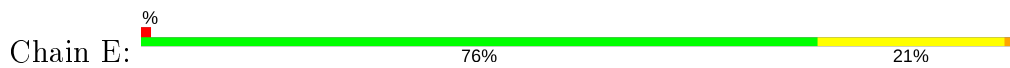
- Molecule 4: Human T-cell Receptor, Class I, Light alpha Chain




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

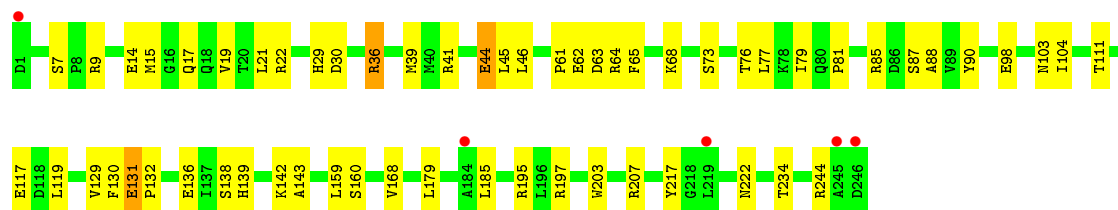


- Molecule 5: Human T-cell Receptor, Class I, Heavy beta Chain



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

Chain J:  2% 76% 23%



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)	Xtrriage
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	Xtrriage
Anisotropy	0.285	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 56.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	13392	wwPDB-VP
Average B, all atoms (Å ²)	73.0	wwPDB-VP

Xtrriage'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 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	0.89	0/2320	1.05	10/3149 (0.3%)
1	F	0.95	2/2320 (0.1%)	1.09	15/3149 (0.5%)
2	B	0.90	0/860	1.02	2/1162 (0.2%)
2	G	0.97	1/860 (0.1%)	1.12	5/1162 (0.4%)
3	C	1.14	0/73	0.96	0/100
3	H	1.15	0/73	1.13	0/100
4	D	0.91	1/1553 (0.1%)	1.08	10/2100 (0.5%)
4	I	0.85	1/1553 (0.1%)	1.01	4/2100 (0.2%)
5	E	0.84	1/2029 (0.0%)	1.00	4/2759 (0.1%)
5	J	0.75	1/2029 (0.0%)	0.97	6/2759 (0.2%)
All	All	0.88	7/13670 (0.1%)	1.04	56/18540 (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 (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	194	PRO	C-O	10.49	1.44	1.23
4	I	194	PRO	C-O	10.06	1.43	1.23
1	F	58	GLU	CB-CG	7.30	1.66	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	F	58	GLU	CG-CD	7.08	1.62	1.51
5	J	44	GLU	CD-OE1	5.14	1.31	1.25
2	G	57	SER	CB-OG	5.11	1.48	1.42
5	E	44	GLU	CD-OE2	5.03	1.31	1.25

All (56) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	14	ARG	NE-CZ-NH2	-10.03	115.28	120.30
1	A	44	ARG	NE-CZ-NH1	8.98	124.79	120.30
2	G	12	ARG	NE-CZ-NH2	-8.25	116.17	120.30
1	F	14	ARG	NE-CZ-NH1	8.18	124.39	120.30
1	A	48	ARG	NE-CZ-NH2	8.00	124.30	120.30
4	I	11	LEU	CA-CB-CG	7.89	133.46	115.30
2	G	34	ASP	CB-CG-OD1	7.66	125.19	118.30
4	D	11	LEU	CA-CB-CG	6.76	130.85	115.30
5	J	36	ARG	NE-CZ-NH1	6.64	123.62	120.30
1	F	6	ARG	NE-CZ-NH2	6.60	123.60	120.30
2	G	59	ASP	N-CA-C	-6.49	93.47	111.00
2	B	12	ARG	NE-CZ-NH2	-6.42	117.09	120.30
4	D	148	SER	N-CA-C	-6.32	93.93	111.00
4	D	130	ASP	CB-CA-C	6.30	123.00	110.40
1	A	35	ARG	NE-CZ-NH1	6.19	123.39	120.30
5	E	30	ASP	CB-CG-OD1	6.14	123.83	118.30
2	G	34	ASP	CB-CG-OD2	-6.12	112.79	118.30
1	F	256	ARG	NE-CZ-NH2	5.97	123.29	120.30
1	F	124	ILE	N-CA-C	5.89	126.91	111.00
5	E	229	ARG	N-CA-C	-5.76	95.43	111.00
1	A	44	ARG	NE-CZ-NH2	-5.76	117.42	120.30
1	A	157	ARG	NE-CZ-NH1	5.74	123.17	120.30
5	J	44	GLU	CA-CB-CG	5.73	126.01	113.40
5	J	36	ARG	NE-CZ-NH2	-5.72	117.44	120.30
1	A	145	HIS	CB-CA-C	5.67	121.74	110.40
1	A	131	ARG	NE-CZ-NH1	5.62	123.11	120.30
4	D	54	GLY	N-CA-C	-5.59	99.11	113.10
5	J	30	ASP	CB-CG-OD1	5.56	123.30	118.30
1	F	181	ARG	NE-CZ-NH2	5.52	123.06	120.30
4	D	41	ARG	NE-CZ-NH1	5.47	123.04	120.30
4	D	94	ASP	CB-CG-OD2	5.46	123.22	118.30
4	D	110	ARG	NE-CZ-NH1	5.46	123.03	120.30
1	A	110	LEU	CB-CG-CD2	5.46	120.28	111.00
1	F	123	TYR	N-CA-C	-5.45	96.28	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	34	ASP	CB-CG-OD1	-5.41	113.44	118.30
1	F	17	ARG	NE-CZ-NH2	5.40	123.00	120.30
5	E	41	ARG	NE-CZ-NH1	5.39	122.99	120.30
1	F	157	ARG	CG-CD-NE	5.34	123.02	111.80
4	D	91	MET	CB-CG-SD	5.34	128.42	112.40
1	F	65	ARG	NE-CZ-NH1	5.33	122.96	120.30
4	I	60	ARG	NE-CZ-NH1	5.31	122.95	120.30
1	F	58	GLU	OE1-CD-OE2	-5.24	117.01	123.30
4	I	58	ASP	N-CA-C	-5.23	96.88	111.00
1	A	6	ARG	NE-CZ-NH1	-5.23	117.69	120.30
1	F	35	ARG	NE-CZ-NH1	5.18	122.89	120.30
5	E	36	ARG	NE-CZ-NH1	5.17	122.89	120.30
1	F	65	ARG	NE-CZ-NH2	-5.17	117.72	120.30
4	D	106	ARG	NE-CZ-NH1	5.17	122.88	120.30
2	G	53	ASP	CB-CG-OD1	-5.17	113.65	118.30
5	J	244	ARG	NE-CZ-NH1	5.15	122.87	120.30
5	J	41	ARG	NE-CZ-NH1	5.11	122.85	120.30
4	D	167	ASP	CB-CG-OD1	5.06	122.86	118.30
4	I	181	ASP	N-CA-C	-5.03	97.42	111.00
1	A	48	ARG	NE-CZ-NH1	-5.01	117.79	120.30
1	F	84	TYR	CB-CG-CD1	5.01	124.01	121.00
1	F	157	ARG	NE-CZ-NH1	5.01	122.81	120.30

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 [\(i\)](#)

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	2254	0	2103	26	0
1	F	2254	0	2103	21	0
2	B	837	0	803	12	0
2	G	837	0	803	10	0
3	C	70	0	61	4	0
3	H	70	0	61	2	0
4	D	1520	0	1447	36	0
4	I	1520	0	1447	30	0
5	E	1974	0	1887	28	0
5	J	1974	0	1887	25	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	174	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 (174) 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	1.58	0.85
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
5:J:15:MET:HE2	5:J:117:GLU:HA	1.67	0.76
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:15:MET:HE2	5:E:117:GLU:HA	1.73	0.70
5:E:120:LYS:HD3	5:E:227:GLN:HE22	1.57	0.70
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: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
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:125:ASP:OD1	4:D:126:SER:N	2.36	0.58
1:F:19:GLU:CD	1:F:75:ARG:HD2	2.24	0.58
2:B:9:VAL:O	2:B:99:MET:CE	2.52	0.57
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
4:D:60:ARG:NH2	4:D:83:ASP:OD2	2.37	0.57
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
1:A:202:ARG:HG3	1:A:246:ALA:HB2	1.88	0.55
1:A:14:ARG:HD3	1:A:19:GLU:O	2.08	0.54
4:I:89:CYS:O	4:I:89:CYS:SG	2.66	0.54
2:G:58:LYS:HA	6:G:102:HOH:O	2.07	0.54
4:D:123:LEU:HD11	5:E:132:PRO:C	2.29	0.54
4:I:46:LEU:HD23	5:J:104:ILE:HG12	1.89	0.53
2:B:9:VAL:O	2:B:99:MET:HE2	2.08	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: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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
3:C:4:GLY:HA2	3:C:5:PRO:C	2.32	0.50
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:15:MET:HE3	5:E:117:GLU:HA	1.90	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
1:A:65:ARG:O	1:A:66:LYS:C	2.51	0.49
1:F:65:ARG:O	1:F:66:LYS:C	2.47	0.49
4:D:151:SER:HB3	4:D:153:VAL:HG23	1.94	0.49
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
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
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
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
4:D:72:ILE:HD11	4:D:91:MET:HE3	1.95	0.46
1:A:223:ASP:O	1:A:224:GLN:HG2	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
5:J:62:GLU:OE2	5:J:62:GLU:N	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
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:34:MET:HE2	5:E:103:ASN:HB2	1.98	0.46
4:D:111:PRO:HG3	4:D:160:VAL:HG11	1.98	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
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
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
4:D:66:ASP:OD1	4:D:66:ASP:C	2.56	0.44
5:E:142:LYS:HD3	5:E:197:ARG:HD3	2.00	0.44
5:J:159:LEU:HD23	5:J:160:SER:N	2.33	0.44
5:J:21:LEU:O	5:J:76:THR:HA	2.18	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:4:VAL:HG11	4:I:91:MET:HE3	2.00	0.44
1:F:28:VAL:HG11	1:F:179:LEU:HD13	2.00	0.43
2:G:98:ASP:OD1	2:G:98:ASP:N	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: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: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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:I:58:ASP:O	4:I:59:GLY:C	2.58	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
2:B:99:MET:HA	2:B:99:MET:HE3	2.02	0.42
4:I:84:SER:O	4:I:85:ALA:HB2	2.19	0.42
2:B:13:HIS:H	2:B:21:ASN:HD21	1.66	0.42
1:A:156:LEU:HG	3:C:3:TRP:CZ2	2.55	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
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
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: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
5:J:14:GLU:HB3	5:J:119:LEU:HG	2.01	0.41
4:D:51:TYR:CD2	4:D:51:TYR:O	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
4:I:66:ASP:OD1	4:I:66:ASP:C	2.59	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
2:B:12:ARG:HB2	2:B:12:ARG:HE	1.67	0.41
1:A:73:THR:HG23	3:C:9:ALA:HB2	2.03	0.41
1:F:274:TRP:NE1	6:F:301:HOH:O	2.36	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
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:N	4:D:181:ASP:OD1	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
4:D:90:ALA:C	4:D:91:MET:HG2	2.40	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:22:PHE:HB3	1:A:38:SER:HB3	2.04	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	274/276 (99%)	245 (89%)	25 (9%)	4 (2%)	10	35
1	F	274/276 (99%)	253 (92%)	19 (7%)	2 (1%)	22	52
2	B	98/100 (98%)	91 (93%)	6 (6%)	1 (1%)	15	45
2	G	98/100 (98%)	96 (98%)	1 (1%)	1 (1%)	15	45
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%)	9	33
4	I	191/193 (99%)	167 (87%)	19 (10%)	5 (3%)	5	22
5	E	244/246 (99%)	226 (93%)	17 (7%)	1 (0%)	34	64
5	J	244/246 (99%)	225 (92%)	19 (8%)	0	100	100
All	All	1630/1650 (99%)	1484 (91%)	129 (8%)	17 (1%)	15	45

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

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

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	232/232 (100%)	206 (89%)	26 (11%)	6	21
1	F	232/232 (100%)	207 (89%)	25 (11%)	6	22
2	B	95/95 (100%)	84 (88%)	11 (12%)	5	19
2	G	95/95 (100%)	84 (88%)	11 (12%)	5	19
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%)	137 (79%)	37 (21%)	1	4
4	I	174/174 (100%)	139 (80%)	35 (20%)	1	4
5	E	216/216 (100%)	191 (88%)	25 (12%)	5	19
5	J	216/216 (100%)	191 (88%)	25 (12%)	5	19
All	All	1444/1444 (100%)	1249 (86%)	195 (14%)	4	14

All (195) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	14	ARG
1	A	32	GLN

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Mol	Chain	Res	Type
1	A	35	ARG
1	A	45	MET
1	A	67	VAL
1	A	74	HIS
1	A	88	SER
1	A	128	GLU
1	A	131	ARG
1	A	141	GLN
1	A	155	GLN
1	A	157	ARG
1	A	165	VAL
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	219	ARG
1	A	222	GLU
1	A	225	THR
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	11	SER
2	B	16	GLU
2	B	21	ASN
2	B	27	VAL
2	B	48	LYS
2	B	50	GLU
2	B	70	PHE
2	B	75	LYS
2	B	91	LYS
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

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Mol	Chain	Res	Type
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
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	127	LYS
4	D	132	SER
4	D	138	ASP
4	D	147	GLN
4	D	152	ASP
4	D	165	SER
4	D	167	ASP
4	D	170	SER
4	D	177	SER
4	D	192	ILE
5	E	7	SER
5	E	19	VAL
5	E	22	ARG
5	E	29	HIS
5	E	30	ASP
5	E	36	ARG
5	E	39	MET
5	E	62	GLU
5	E	73	SER
5	E	78	LYS
5	E	80	GLN
5	E	87	SER

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Mol	Chain	Res	Type
5	E	97	TRP
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
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	45	MET
1	F	48	ARG
1	F	58	GLU
1	F	67	VAL
1	F	75	ARG
1	F	88	SER
1	F	110	LEU
1	F	121	LYS
1	F	131	ARG
1	F	157	ARG
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	219	ARG
1	F	226	GLN
1	F	256	ARG
1	F	271	THR
1	F	275	GLU
2	G	4	THR
2	G	11	SER
2	G	14	PRO
2	G	21	ASN

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Mol	Chain	Res	Type
2	G	27	VAL
2	G	33	SER
2	G	34	ASP
2	G	47	GLU
2	G	70	PHE
2	G	75	LYS
2	G	98	ASP
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	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	138	ASP
4	I	146	SER
4	I	151	SER
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	17	GLN
5	J	19	VAL
5	J	22	ARG
5	J	29	HIS
5	J	36	ARG
5	J	39	MET
5	J	44	GLU
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

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (33) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	54	GLN
1	A	72	GLN
1	A	174	ASN
1	A	180	GLN
1	A	262	GLN
2	B	21	ASN
2	B	24	ASN
2	B	84	HIS
2	B	89	GLN
4	D	64	GLN
4	D	144	ASN
4	D	171	ASN
4	D	186	ASN

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Mol	Chain	Res	Type
5	E	18	GLN
5	E	29	HIS
5	E	141	GLN
1	F	54	GLN
1	F	74	HIS
1	F	174	ASN
1	F	180	GLN
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	213	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 carbohydrates 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	276/276 (100%)	-0.18	3 (1%) 80 60	45, 65, 108, 138	0
1	F	276/276 (100%)	-0.18	1 (0%) 92 82	30, 57, 123, 169	0
2	B	100/100 (100%)	-0.33	0 100 100	47, 60, 88, 109	0
2	G	100/100 (100%)	-0.45	0 100 100	38, 55, 76, 91	0
3	C	10/10 (100%)	-0.40	0 100 100	47, 54, 56, 61	0
3	H	10/10 (100%)	-0.23	0 100 100	42, 50, 59, 60	0
4	D	193/193 (100%)	0.17	10 (5%) 27 11	46, 77, 130, 155	0
4	I	193/193 (100%)	0.30	15 (7%) 13 4	51, 86, 144, 173	0
5	E	246/246 (100%)	-0.18	3 (1%) 79 58	34, 63, 120, 143	0
5	J	246/246 (100%)	0.03	5 (2%) 65 41	49, 80, 117, 151	0
All	All	1650/1650 (100%)	-0.08	37 (2%) 62 38	30, 67, 126, 173	0

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
4	D	150	ASP	7.4
4	D	194	PRO	6.7
4	I	180	SER	5.0
4	I	194	PRO	4.9
5	J	1	ASP	4.7
4	I	53	SER	4.5
5	J	246	ASP	3.8
5	J	184	ALA	3.7
4	D	193	ILE	3.6
4	I	192	ILE	3.6
4	I	2	LYS	3.3
4	D	192	ILE	3.3
4	I	193	ILE	3.2

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Mol	Chain	Res	Type	RSRZ
4	D	165	SER	3.1
4	I	189	ASN	2.9
5	E	246	ASP	2.8
4	D	168	PHE	2.7
4	I	164	ARG	2.6
5	E	1	ASP	2.6
4	D	180	SER	2.5
5	J	219	LEU	2.5
1	A	253	GLN	2.5
4	I	178	ASN	2.5
4	I	181	ASP	2.4
4	I	149	LYS	2.4
1	A	251	SER	2.3
4	D	2	LYS	2.3
4	D	149	LYS	2.3
4	D	190	ASN	2.3
4	I	128	SER	2.3
5	J	245	ALA	2.2
5	E	222	ASN	2.2
4	I	166	MET	2.1
1	F	226	GLN	2.1
4	I	167	ASP	2.1
1	A	249	VAL	2.0
4	I	191	SER	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 carbohydrates 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.