



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 18, 2026 – 05:54 PM UTC

PDB ID : 4QRP / pdb_00004qrp
Title : Crystal Structure of HLA B*0801 in complex with HSKKKCDEL and DD31 TCR
Authors : Gras, S.; Berry, R.; Lucet, I.S.; Rossjohn, J.
Deposited on : 2014-07-02
Resolution : 2.90 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

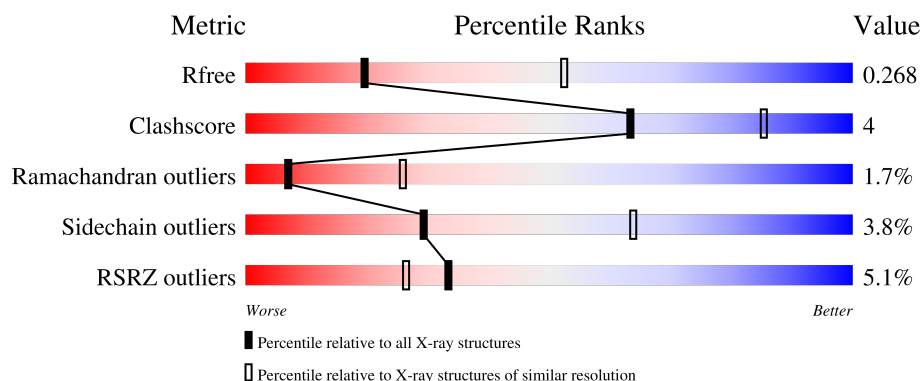
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (2.90-2.90)

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	8% 85% 14% .
1	F	276	8% 88% 11% .
2	B	100	16% 80% 14% . ..
2	G	100	% 92% 8%
3	C	9	100%

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Mol	Chain	Length	Quality of chain
3	H	9	89% 11%
4	D	206	5% 83% 15% .
4	J	206	7% 85% 13% .
4	K	206	% 86% 11% ..
5	E	245	% 89% 9% ..
5	I	245	3% 84% 12% ..
5	L	245	4% 83% 15% ..

2 Entry composition [i](#)

There are 8 unique types of molecules in this entry. The entry contains 16924 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, B-8 alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	276	Total	C	N	O	S	0	0	0
			2251	1395	411	438	7			
1	F	276	Total	C	N	O	S	0	0	0
			2251	1395	411	438	7			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	99	Total	C	N	O	S	0	0	0
			829	528	140	158	3			
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	-	expression tag	UNP P61769
G	0	MET	-	expression tag	UNP P61769

- Molecule 3 is a protein called NS3-4A protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	9	Total	C	N	O	S	0	0	0
			75	45	14	15	1			
3	H	9	Total	C	N	O	S	0	0	0
			75	45	14	15	1			

- Molecule 4 is a protein called DD31 TCR alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	203	Total	C	N	O	S	0	0	0
			1573	987	254	322	10			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	J	203	Total	C	N	O	S	0	0	0
			1573	987	254	322	10			
4	K	202	Total	C	N	O	S	0	2	0
			1578	990	257	322	9			

- Molecule 5 is a protein called DD31 TCR beta chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	E	243	Total	C	N	O	S	0	0	0
			1937	1223	344	365	5			
5	I	243	Total	C	N	O	S	0	0	0
			1937	1223	344	365	5			
5	L	243	Total	C	N	O	S	0	0	0
			1937	1223	344	365	5			

- Molecule 6 is IODIDE ION (CCD ID: IOD) (formula: I).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	D	1	Total	I	0	0
			1	1		
6	E	2	Total	I	0	0
			2	2		
6	G	1	Total	I	0	0
			1	1		
6	I	2	Total	I	0	0
			2	2		
6	J	1	Total	I	0	0
			1	1		
6	K	2	Total	I	0	0
			2	2		
6	L	3	Total	I	0	0
			3	3		

- Molecule 7 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	E	1	Total	Na	0	0
			1	1		
7	G	1	Total	Na	0	0
			1	1		
7	L	1	Total	Na	0	0
			1	1		

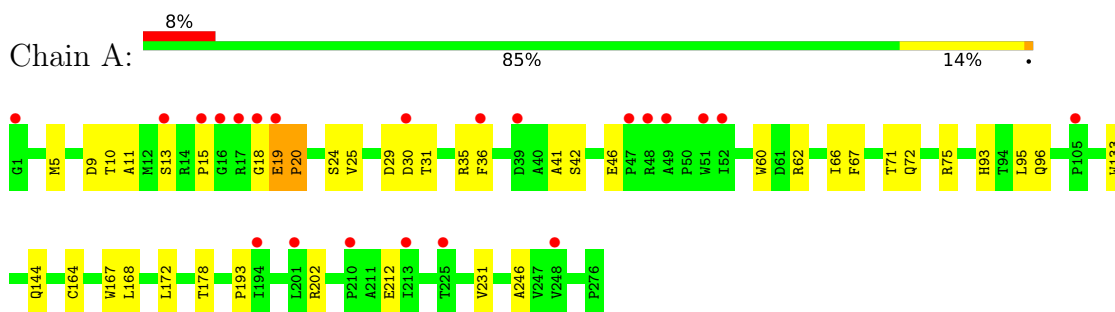
- Molecule 8 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	D	5	Total O 5 5	0	0
8	E	9	Total O 9 9	0	0
8	F	10	Total O 10 10	0	0
8	G	6	Total O 6 6	0	0
8	H	1	Total O 1 1	0	0
8	I	3	Total O 3 3	0	0
8	J	3	Total O 3 3	0	0
8	K	14	Total O 14 14	0	0
8	L	5	Total O 5 5	0	0

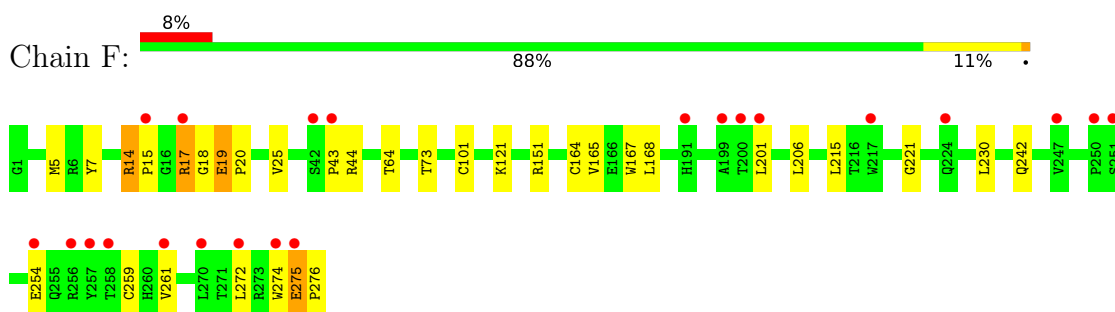
3 Residue-property plots [i](#)

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.

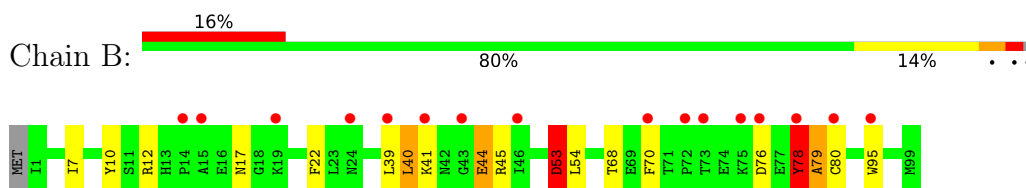
- Molecule 1: HLA class I histocompatibility antigen, B-8 alpha chain



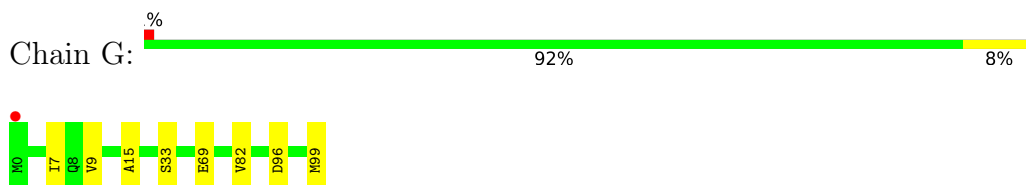
- Molecule 1: HLA class I histocompatibility antigen, B-8 alpha chain



- Molecule 2: Beta-2-microglobulin



- Molecule 2: Beta-2-microglobulin



- Molecule 3: NS3-4A protein

Chain C:  100%

There are no outlier residues recorded for this chain.

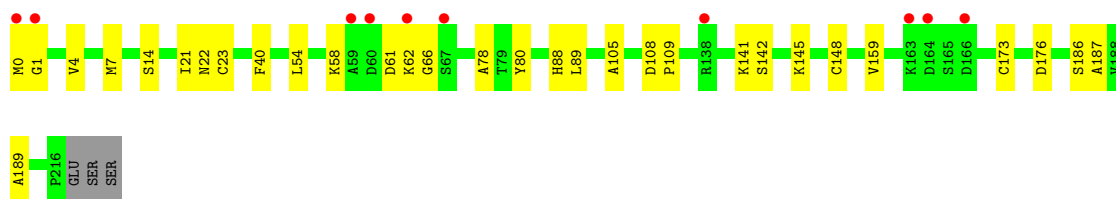
- Molecule 3: NS3-4A protein

Chain H:  89% 11%



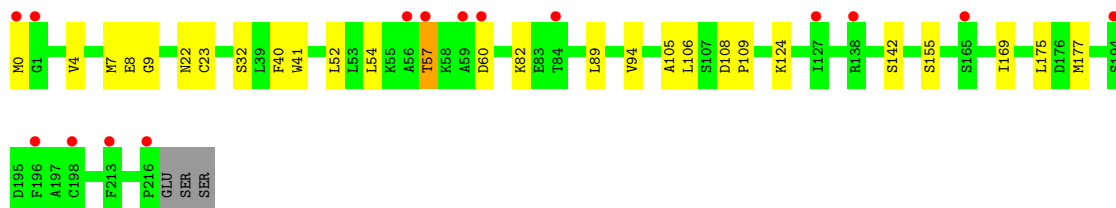
- Molecule 4: DD31 TCR alpha chain

Chain D:  5% 83% 15%



- Molecule 4: DD31 TCR alpha chain

Chain J:  7% 85% 13%



- Molecule 4: DD31 TCR alpha chain

Chain K:  0% 86% 11%



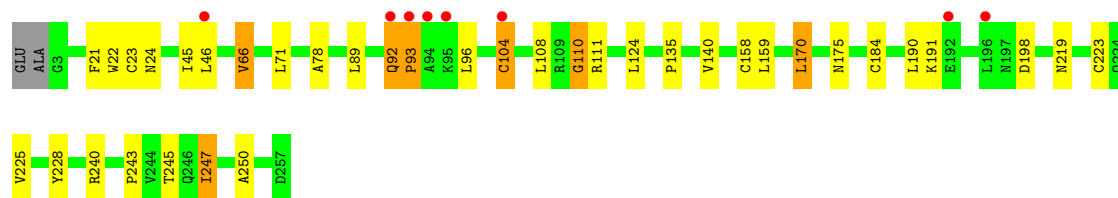
- Molecule 5: DD31 TCR beta chain

Chain E:  0% 89% 9%

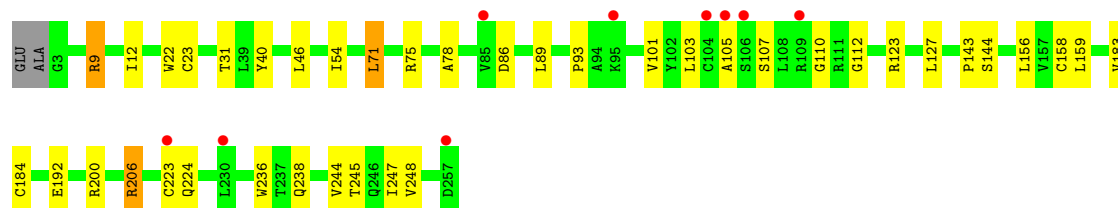
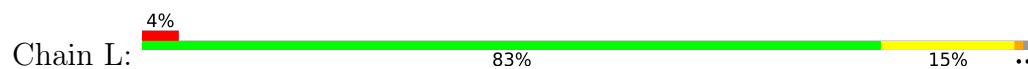


- Molecule 5: DD31 TCR beta chain

Chain I:  3% 84% 12%



• Molecule 5: DD31 TCR beta chain



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	74.16Å 252.19Å 79.45Å 90.00° 101.97° 90.00°	Depositor
Resolution (Å)	19.85 – 2.90 19.85 – 2.90	Depositor EDS
% Data completeness (in resolution range)	99.9 (19.85-2.90) 99.6 (19.85-2.90)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.43 (at 2.88Å)	Xtriage
Refinement program	BUSTER-TNT BUSTER 2.10.0, BUSTER 2.10.0	Depositor
R, R_{free}	0.206 , 0.248 0.229 , 0.268	Depositor DCC
R_{free} test set	3181 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	49.0	Xtriage
Anisotropy	0.297	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.26 , 35.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	16924	wwPDB-VP
Average B, all atoms (Å ²)	61.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.54% 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

Bond lengths and bond angles in the following residue types are not validated in this section: NA, IOD

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	0.67	0/2313	1.22	2/3146 (0.1%)
1	F	0.67	0/2313	1.24	4/3146 (0.1%)
2	B	0.75	0/852	1.36	10/1152 (0.9%)
2	G	0.64	0/860	1.14	1/1162 (0.1%)
3	C	0.55	0/75	1.03	0/95
3	H	0.59	0/75	1.07	0/95
4	D	0.67	0/1608	1.15	1/2182 (0.0%)
4	J	0.71	0/1608	1.15	0/2182
4	K	0.69	0/1613	1.16	3/2189 (0.1%)
5	E	0.64	0/1988	1.15	6/2702 (0.2%)
5	I	0.64	0/1988	1.13	7/2702 (0.3%)
5	L	0.65	0/1988	1.13	2/2702 (0.1%)
All	All	0.67	0/17281	1.18	36/23455 (0.2%)

There are no bond length outliers.

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	78	TYR	CA-C-N	10.94	141.40	121.70
2	B	78	TYR	C-N-CA	10.94	141.40	121.70
5	I	110	GLY	CA-C-N	6.75	133.85	121.70
5	I	110	GLY	C-N-CA	6.75	133.85	121.70
5	E	110	GLY	CA-C-N	6.45	133.30	121.70
5	E	110	GLY	C-N-CA	6.45	133.30	121.70
1	A	20	PRO	N-CA-C	6.35	125.55	112.47
2	B	70	PHE	CA-CB-CG	5.94	119.74	113.80
2	B	76	ASP	CA-C-N	5.83	132.20	121.70
2	B	76	ASP	C-N-CA	5.83	132.20	121.70
5	E	211	ALA	CA-C-N	5.64	128.10	120.38
5	E	211	ALA	C-N-CA	5.64	128.10	120.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	54	LEU	N-CA-C	5.57	118.43	110.24
5	I	66	VAL	CB-CA-C	5.56	116.36	110.91
5	I	111	ARG	CA-C-N	5.55	126.25	120.03
5	I	111	ARG	C-N-CA	5.55	126.25	120.03
2	B	44	GLU	CA-C-N	5.43	131.92	121.54
2	B	44	GLU	C-N-CA	5.43	131.92	121.54
5	I	96	LEU	CA-C-N	5.39	127.51	120.28
5	I	96	LEU	C-N-CA	5.39	127.51	120.28
1	A	9	ASP	CA-CB-CG	5.35	117.95	112.60
2	G	33	SER	N-CA-C	5.34	117.85	111.71
4	D	176	ASP	CA-CB-CG	5.33	117.93	112.60
4	K	191	ASN	CA-C-N	5.33	128.67	121.05
4	K	191	ASN	C-N-CA	5.33	128.67	121.05
5	E	238	GLN	CA-C-N	5.30	127.65	120.65
5	E	238	GLN	C-N-CA	5.30	127.65	120.65
4	K	153	ASP	CA-CB-CG	5.29	117.89	112.60
1	F	151	ARG	CA-C-N	5.21	127.12	120.56
1	F	151	ARG	C-N-CA	5.21	127.12	120.56
1	F	64	THR	CA-C-N	5.17	127.17	120.44
1	F	64	THR	C-N-CA	5.17	127.17	120.44
5	L	238	GLN	CA-C-N	5.05	127.04	120.28
5	L	238	GLN	C-N-CA	5.05	127.04	120.28
2	B	53	ASP	CA-C-N	5.03	128.25	121.05
2	B	53	ASP	C-N-CA	5.03	128.25	121.05

There are no chirality outliers.

There are no planarity outliers.

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	2251	0	2097	25	0
1	F	2251	0	2097	30	0
2	B	829	0	796	8	0
2	G	837	0	805	2	0
3	C	75	0	79	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	H	75	0	79	1	0
4	D	1573	0	1498	11	0
4	J	1573	0	1498	9	0
4	K	1578	0	1498	12	0
5	E	1937	0	1875	11	0
5	I	1937	0	1875	14	0
5	L	1937	0	1875	17	0
6	D	1	0	0	0	0
6	E	2	0	0	0	0
6	G	1	0	0	0	0
6	I	2	0	0	0	0
6	J	1	0	0	0	0
6	K	2	0	0	1	0
6	L	3	0	0	0	0
7	E	1	0	0	0	0
7	G	1	0	0	0	0
7	L	1	0	0	0	0
8	D	5	0	0	0	0
8	E	9	0	0	0	0
8	F	10	0	0	0	0
8	G	6	0	0	0	0
8	H	1	0	0	0	0
8	I	3	0	0	0	0
8	J	3	0	0	0	0
8	K	14	0	0	0	0
8	L	5	0	0	0	0
All	All	16924	0	16072	128	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 4.

All (128) 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:F:254:GLU:CB	1:F:275:GLU:OE1	1.79	1.30
1:F:201:LEU:HD11	1:F:275:GLU:OE2	1.13	1.30
1:A:19:GLU:OE1	1:A:75:ARG:NH1	1.74	1.19
2:B:78:TYR:HB2	2:B:79:ALA:HB2	1.13	1.12
1:F:254:GLU:HB2	1:F:275:GLU:CD	1.76	1.11
1:F:254:GLU:HB2	1:F:275:GLU:OE1	1.43	1.11
1:F:201:LEU:CD1	1:F:275:GLU:OE2	2.03	1.05

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:201:LEU:HD21	1:F:275:GLU:HG2	1.40	1.00
1:A:19:GLU:CG	1:A:75:ARG:HG2	1.91	0.99
1:F:254:GLU:HB3	1:F:275:GLU:OE1	1.63	0.97
1:A:19:GLU:HG2	1:A:75:ARG:HG2	1.46	0.97
5:E:158:CYS:HG	5:E:223:CYS:HG	1.06	0.97
1:A:19:GLU:CD	1:A:75:ARG:HG2	1.91	0.94
2:B:78:TYR:HB2	2:B:79:ALA:CB	1.98	0.92
1:F:201:LEU:HD21	1:F:275:GLU:CG	2.08	0.82
5:E:23:CYS:HG	5:E:104:CYS:HG	1.21	0.80
4:D:0:MET:HG2	4:D:1:GLY:H	1.53	0.73
1:F:254:GLU:CD	1:F:275:GLU:OE1	2.34	0.70
5:I:22:TRP:CH2	5:I:24:ASN:HB2	2.26	0.70
4:K:174:LEU:HB3	5:L:184:CYS:HB2	1.73	0.70
4:J:7:MET:HE3	4:J:22:ASN:H	1.57	0.69
5:L:9:ARG:HE	5:L:9:ARG:H	1.41	0.68
1:F:254:GLU:CG	1:F:275:GLU:OE1	2.42	0.67
1:F:201:LEU:CD2	1:F:275:GLU:HG2	2.21	0.65
4:K:22:CYS:HG	4:K:103:CYS:HG	0.69	0.65
1:A:19:GLU:CD	1:A:75:ARG:CG	2.71	0.63
1:F:201:LEU:HD21	1:F:275:GLU:CD	2.24	0.62
4:K:32:LEU:HD13	4:K:86:PHE:HB2	1.81	0.61
5:L:40:TYR:HB2	5:L:105:ALA:HB3	1.82	0.61
4:D:7:MET:HE3	4:D:22:ASN:H	1.65	0.61
5:E:92:GLN:HB3	5:E:93:PRO:HD2	1.83	0.61
1:A:46:GLU:HB3	1:A:60:TRP:HE1	1.67	0.59
4:J:40:PHE:HB2	4:J:105:ALA:HB3	1.84	0.59
1:A:13:SER:HB2	1:A:93:HIS:H	1.67	0.59
1:A:18:GLY:O	1:A:19:GLU:HG3	2.04	0.58
4:D:108:ASP:HB2	4:D:109:PRO:HD2	1.85	0.58
5:E:92:GLN:O	5:E:93:PRO:C	2.48	0.57
5:L:143:PRO:HD3	5:L:156:LEU:HG	1.86	0.57
5:I:170:LEU:HG	5:I:225:VAL:HG22	1.87	0.56
2:B:39:LEU:HA	2:B:80:CYS:HA	1.88	0.56
1:F:215:LEU:HD22	1:F:261:VAL:HG12	1.88	0.55
4:K:172:CYS:HB3	5:L:184:CYS:SG	2.46	0.55
1:A:5:MET:HB2	1:A:168:LEU:HD13	1.87	0.55
1:F:254:GLU:CB	1:F:275:GLU:CD	2.52	0.54
5:I:175:ASN:HD21	5:I:219:ASN:HD22	1.54	0.54
5:L:158:CYS:HG	5:L:223:CYS:HG	1.53	0.54
2:G:7:ILE:HG12	2:G:82:VAL:HG21	1.90	0.54
1:A:19:GLU:OE2	1:A:75:ARG:HB3	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:201:LEU:HD21	1:F:275:GLU:OE2	2.08	0.53
5:L:54:ILE:HB	5:L:71:LEU:HD13	1.90	0.53
1:A:19:GLU:CD	1:A:75:ARG:HH11	2.09	0.53
5:I:92:GLN:HG3	5:I:93:PRO:HD2	1.90	0.53
4:J:108:ASP:HB2	4:J:109:PRO:HD2	1.91	0.53
4:D:21:ILE:HD12	4:D:89:LEU:HD23	1.90	0.53
5:E:247:ILE:H	1:F:17:ARG:HH21	1.57	0.53
4:K:192:LYS:HE3	4:K:194:ASP:HB2	1.90	0.52
4:D:40:PHE:HB2	4:D:105:ALA:HB3	1.91	0.52
1:A:11:ALA:HB3	1:A:95:LEU:HB3	1.92	0.52
1:A:35:ARG:CZ	2:B:53:ASP:HB2	2.40	0.51
1:A:167:TRP:HZ2	4:D:0:MET:HB2	1.75	0.51
5:E:247:ILE:HD12	1:F:17:ARG:HG3	1.92	0.51
4:D:23:CYS:H	4:D:88:HIS:HD2	1.59	0.51
1:A:24:SER:HB3	1:A:36:PHE:HB2	1.93	0.50
2:G:96:ASP:HB3	2:G:99:MET:HB2	1.92	0.50
1:A:72:GLN:HA	1:A:75:ARG:HD2	1.94	0.49
4:K:153:ASP:HB2	6:K:301:IOD:I	2.82	0.49
2:B:39:LEU:HD11	2:B:68:THR:HG22	1.95	0.49
5:L:158:CYS:CB	5:L:223:CYS:HG	2.26	0.49
1:A:62:ARG:O	1:A:66:ILE:HG12	2.13	0.49
4:K:141:SER:HB2	4:K:142:SER:HA	1.95	0.49
5:I:92:GLN:O	5:I:93:PRO:C	2.54	0.49
5:I:228:TYR:HA	5:I:245:THR:HG23	1.95	0.48
1:A:18:GLY:C	1:A:19:GLU:HG3	2.38	0.48
1:F:259:CYS:HB3	1:F:272:LEU:HB2	1.95	0.48
1:A:10:THR:HG22	1:A:96:GLN:HG2	1.96	0.47
1:A:41:ALA:HA	1:A:42:SER:HA	1.68	0.47
1:F:5:MET:HB2	1:F:168:LEU:HD13	1.96	0.47
4:D:148:CYS:HB2	4:D:189:ALA:HB3	1.97	0.47
5:L:22:TRP:HE1	5:L:86:ASP:HB3	1.80	0.47
4:D:58:LYS:O	4:D:61:ASP:HB2	2.15	0.47
5:E:78:ALA:HB2	5:E:89:LEU:HD12	1.97	0.46
5:L:184:CYS:HB3	5:L:206:ARG:HB2	1.97	0.46
5:L:192:GLU:HG3	5:L:200:ARG:HB2	1.96	0.46
1:A:19:GLU:HG2	1:A:75:ARG:CG	2.31	0.46
1:A:133:TRP:HB2	1:A:144:GLN:HG3	1.97	0.46
5:E:21:PHE:HZ	5:E:124:LEU:HD22	1.82	0.45
2:B:40:LEU:HD11	2:B:44:GLU:C	2.41	0.45
5:E:247:ILE:HD13	1:F:15:PRO:HB2	1.98	0.45
5:L:75:ARG:HA	5:L:93:PRO:HD2	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:L:78:ALA:HB2	5:L:89:LEU:HD12	1.98	0.45
2:B:10:TYR:HA	2:B:95:TRP:CZ3	2.52	0.45
5:E:193:GLN:HG3	5:E:196:LEU:HD13	2.00	0.44
5:I:158:CYS:SG	5:I:223:CYS:SG	3.03	0.44
1:F:206:LEU:HD23	1:F:242:GLN:HG2	2.01	0.43
5:L:101:VAL:HG22	5:L:123:ARG:HD2	2.00	0.43
5:I:135:PRO:HD3	5:I:243:PRO:HB3	1.99	0.43
4:J:8:GLU:HG2	4:J:9:GLY:N	2.33	0.43
5:L:31:THR:HG21	5:L:110:GLY:HA3	2.01	0.43
4:D:159:VAL:HG21	4:D:187:ALA:HB2	2.01	0.43
4:K:85:SER:HB2	4:K:87:HIS:CE1	2.54	0.43
4:D:54:LEU:HD11	4:D:78:ALA:HB3	1.99	0.43
5:L:12:ILE:HG23	5:L:127:LEU:HD13	2.00	0.43
5:I:140:VAL:HG23	5:I:250:ALA:HB3	2.00	0.43
4:K:163:ASP:HB3	4:K:166:VAL:HG22	2.00	0.43
5:I:21:PHE:HZ	5:I:124:LEU:HD22	1.84	0.42
1:A:202:ARG:HG3	1:A:246:ALA:HB2	2.02	0.42
1:F:19:GLU:HA	1:F:20:PRO:HD3	1.94	0.42
4:K:195:PHE:HA	4:K:199:ASN:HD21	1.84	0.42
5:I:78:ALA:HB2	5:I:89:LEU:HD12	2.01	0.42
1:F:73:THR:HG21	3:H:6:CYS:HA	2.01	0.42
4:J:124:LYS:HD2	4:J:155:SER:HB3	2.01	0.42
1:A:67:PHE:O	1:A:71:THR:HG23	2.20	0.41
5:L:224:GLN:HG3	5:L:247:ILE:HG23	2.03	0.41
5:I:184:CYS:SG	4:J:175:LEU:HB3	2.60	0.41
1:F:14:ARG:HG3	1:F:18:GLY:H	1.85	0.41
1:F:167:TRP:HZ2	4:J:0:MET:HB2	1.86	0.41
2:B:12:ARG:HB2	2:B:22:PHE:HB2	2.02	0.41
1:A:15:PRO:HB2	5:I:247:ILE:HD11	2.01	0.41
1:F:101:CYS:SG	1:F:164:CYS:SG	3.13	0.41
4:J:32:SER:HA	4:J:57:THR:HA	2.02	0.41
5:E:247:ILE:H	1:F:17:ARG:HE	1.68	0.41
1:F:5:MET:HE3	1:F:7:TYR:HE1	1.86	0.41
5:I:23:CYS:CB	5:I:104:CYS:HG	2.33	0.40
4:K:192:LYS:HE2	4:K:195:PHE:HB2	2.04	0.40
1:F:201:LEU:CD2	1:F:275:GLU:OE2	2.69	0.40
4:J:41:TRP:CE2	4:J:89:LEU:HB2	2.56	0.40
4:K:1:ASP:O	4:K:113:ARG:HD2	2.21	0.40
1:F:275:GLU:H	1:F:276:PRO:CD	2.33	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%)	249 (91%)	20 (7%)	5 (2%)	6	25
1	F	274/276 (99%)	249 (91%)	18 (7%)	7 (3%)	4	17
2	B	97/100 (97%)	83 (86%)	10 (10%)	4 (4%)	2	9
2	G	98/100 (98%)	94 (96%)	3 (3%)	1 (1%)	12	39
3	C	7/9 (78%)	7 (100%)	0	0	100	100
3	H	7/9 (78%)	7 (100%)	0	0	100	100
4	D	201/206 (98%)	189 (94%)	9 (4%)	3 (2%)	8	28
4	J	201/206 (98%)	189 (94%)	10 (5%)	2 (1%)	12	39
4	K	202/206 (98%)	181 (90%)	18 (9%)	3 (2%)	8	28
5	E	241/245 (98%)	222 (92%)	16 (7%)	3 (1%)	10	34
5	I	241/245 (98%)	217 (90%)	19 (8%)	5 (2%)	5	21
5	L	241/245 (98%)	224 (93%)	14 (6%)	3 (1%)	10	34
All	All	2084/2123 (98%)	1911 (92%)	137 (7%)	36 (2%)	7	26

All (36) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	20	PRO
2	B	45	ARG
5	E	46	LEU
5	E	93	PRO
1	F	17	ARG
1	F	275	GLU
5	I	93	PRO
4	K	7	GLU
4	K	141	SER
1	A	19	GLU
2	B	79	ALA
5	E	110	GLY

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Mol	Chain	Res	Type
1	F	44	ARG
5	I	46	LEU
5	I	110	GLY
5	I	240	ARG
4	K	62	GLY
1	A	29	ASP
4	D	142	SER
2	B	17	ASN
2	B	78	TYR
4	D	62	LYS
1	F	19	GLU
1	F	221	GLY
1	F	274	TRP
2	G	15	ALA
4	J	142	SER
5	L	236	TRP
1	A	212	GLU
5	L	112	GLY
5	L	245	THR
4	J	57	THR
4	D	66	GLY
1	A	193	PRO
5	I	45	ILE
1	F	43	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	235/235 (100%)	228 (97%)	7 (3%)	36	70
1	F	235/235 (100%)	230 (98%)	5 (2%)	47	77
2	B	94/95 (99%)	90 (96%)	4 (4%)	26	60
2	G	95/95 (100%)	93 (98%)	2 (2%)	47	77
3	C	9/9 (100%)	9 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
3	H	9/9 (100%)	9 (100%)	0	100	100
4	D	179/182 (98%)	172 (96%)	7 (4%)	28	63
4	J	179/182 (98%)	169 (94%)	10 (6%)	19	50
4	K	178/182 (98%)	174 (98%)	4 (2%)	45	77
5	E	210/211 (100%)	202 (96%)	8 (4%)	29	64
5	I	210/211 (100%)	199 (95%)	11 (5%)	21	52
5	L	210/211 (100%)	198 (94%)	12 (6%)	18	49
All	All	1843/1857 (99%)	1773 (96%)	70 (4%)	29	64

All (70) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	25	VAL
1	A	30	ASP
1	A	31	THR
1	A	164	CYS
1	A	172	LEU
1	A	178	THR
1	A	231	VAL
2	B	7	ILE
2	B	40	LEU
2	B	41	LYS
2	B	53	ASP
4	D	4	VAL
4	D	14	SER
4	D	80	TYR
4	D	141	LYS
4	D	145	LYS
4	D	173	CYS
4	D	186	SER
5	E	13	ILE
5	E	130	LEU
5	E	159	LEU
5	E	161	THR
5	E	168	VAL
5	E	183	VAL
5	E	196	LEU
5	E	206	ARG
1	F	14	ARG
1	F	25	VAL

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Mol	Chain	Res	Type
1	F	121	LYS
1	F	165	VAL
1	F	230	LEU
2	G	9	VAL
2	G	69	GLU
5	I	66	VAL
5	I	71	LEU
5	I	92	GLN
5	I	104	CYS
5	I	108	LEU
5	I	159	LEU
5	I	170	LEU
5	I	190	LEU
5	I	191	LYS
5	I	198	ASP
5	I	247	ILE
4	J	4	VAL
4	J	23	CYS
4	J	52	LEU
4	J	54	LEU
4	J	60	ASP
4	J	82	LYS
4	J	94	VAL
4	J	106	LEU
4	J	169	ILE
4	J	177	MET
4	K	49	LEU
4	K	53	LEU
4	K	79	TYR
4	K	172	CYS
5	L	9	ARG
5	L	23	CYS
5	L	46	LEU
5	L	71	LEU
5	L	103	LEU
5	L	107	SER
5	L	144	SER
5	L	159	LEU
5	L	183	VAL
5	L	206	ARG
5	L	244	VAL
5	L	248	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (38) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	54	GLN
1	A	174	ASN
1	A	242	GLN
2	B	8	GLN
2	B	84	HIS
4	D	44	GLN
4	D	88	HIS
4	D	158	ASN
4	D	200	ASN
5	E	44	GLN
5	E	55	GLN
5	E	57	GLN
5	E	117	HIS
5	E	220	HIS
5	E	226	GLN
1	F	174	ASN
2	G	13	HIS
5	I	44	GLN
5	I	57	GLN
5	I	117	HIS
5	I	219	ASN
5	I	226	GLN
5	I	238	GLN
4	J	22	ASN
4	J	44	GLN
4	J	68	ASN
4	J	161	GLN
4	J	200	ASN
4	K	50	GLN
4	K	199	ASN
4	K	202	ASN
5	L	17	GLN
5	L	48	GLN
5	L	92	GLN
5	L	132	ASN
5	L	215	GLN
5	L	226	GLN
5	L	238	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 15 ligands modelled in this entry, 15 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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%)	0.88	22 (7%)	18 15	39, 72, 101, 124	0
1	F	276/276 (100%)	0.38	22 (7%)	18 15	22, 48, 105, 124	0
2	B	99/100 (99%)	1.25	16 (16%)	4 4	58, 85, 115, 119	0
2	G	100/100 (100%)	-0.04	1 (1%)	79 73	26, 44, 66, 73	0
3	C	9/9 (100%)	0.48	0	100 100	53, 58, 62, 62	0
3	H	9/9 (100%)	-0.12	0	100 100	36, 39, 49, 53	0
4	D	203/206 (98%)	0.42	10 (4%)	35 27	31, 54, 83, 96	0
4	J	203/206 (98%)	0.74	15 (7%)	20 17	36, 61, 93, 109	0
4	K	202/206 (98%)	0.07	3 (1%)	72 64	29, 47, 68, 81	2 (0%)
5	E	243/245 (99%)	0.14	2 (0%)	82 77	22, 48, 82, 93	1 (0%)
5	I	243/245 (99%)	0.56	8 (3%)	49 40	29, 63, 98, 117	0
5	L	243/245 (99%)	0.53	9 (3%)	45 37	28, 62, 92, 104	0
All	All	2106/2123 (99%)	0.48	108 (5%)	33 26	22, 58, 98, 124	3 (0%)

All (108) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	19	GLU	6.6
4	K	143	ASP	5.9
1	F	275	GLU	5.1
1	A	17	ARG	4.9
4	D	59	ALA	4.7
4	J	1	GLY	4.3
1	F	17	ARG	3.8
4	J	59	ALA	3.8
2	B	73	THR	3.7
4	J	138	ARG	3.6
5	I	92	GLN	3.5

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Mol	Chain	Res	Type	RSRZ
4	D	0	MET	3.5
1	A	47	PRO	3.4
1	F	261	VAL	3.4
1	F	201	LEU	3.3
1	F	254	GLU	3.2
4	D	1	GLY	3.2
2	B	76	ASP	3.2
1	A	225	THR	3.2
4	J	57	THR	3.2
1	A	13	SER	3.1
5	I	93	PRO	3.1
4	J	0	MET	3.1
5	L	109	ARG	3.1
2	B	14	PRO	3.1
2	B	19	LYS	3.1
2	G	0	MET	3.0
4	J	194	SER	3.0
2	B	15	ALA	3.0
4	J	165	SER	3.0
5	L	257	ASP	3.0
4	D	62	LYS	3.0
4	J	56	ALA	3.0
5	E	46	LEU	2.9
1	A	1	GLY	2.9
1	A	16	GLY	2.9
1	A	201	LEU	2.8
1	F	270	LEU	2.8
4	D	67	SER	2.8
2	B	78	TYR	2.7
5	L	106	SER	2.7
1	F	43	PRO	2.7
1	F	274	TRP	2.7
1	F	15	PRO	2.7
5	I	94	ALA	2.6
1	F	257	TYR	2.6
2	B	80	CYS	2.6
2	B	72	PRO	2.6
1	A	194	ILE	2.6
1	A	18	GLY	2.6
2	B	41	LYS	2.6
4	J	196	PHE	2.6
5	L	95	LYS	2.6

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Mol	Chain	Res	Type	RSRZ
5	L	230	LEU	2.5
4	J	60	ASP	2.5
1	F	250	PRO	2.5
5	L	104	CYS	2.5
1	A	52	ILE	2.5
4	D	138	ARG	2.5
1	A	39	ASP	2.5
2	B	39	LEU	2.5
1	A	210	PRO	2.5
2	B	46	ILE	2.5
5	E	9	ARG	2.5
4	J	127	ILE	2.4
1	F	251	SER	2.4
2	B	70	PHE	2.4
4	D	60	ASP	2.4
1	A	48	ARG	2.4
2	B	95	TRP	2.4
4	J	84	THR	2.4
1	F	191	HIS	2.4
1	F	272	LEU	2.4
1	A	15	PRO	2.4
4	D	164	ASP	2.4
4	K	141	SER	2.3
5	I	46	LEU	2.3
1	A	49	ALA	2.3
1	A	213	ILE	2.3
2	B	75	LYS	2.3
1	A	105	PRO	2.3
5	L	223	CYS	2.3
4	D	163	LYS	2.3
1	F	42	SER	2.3
1	F	200	THR	2.3
1	F	258	THR	2.3
4	J	198	CYS	2.2
2	B	43	GLY	2.2
4	K	27	THR	2.2
1	A	30	ASP	2.2
1	F	217	TRP	2.2
5	I	196	LEU	2.2
1	F	256	ARG	2.1
2	B	24	ASN	2.1
5	I	95	LYS	2.1

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Mol	Chain	Res	Type	RSRZ
1	F	199	ALA	2.1
1	A	248	VAL	2.1
4	D	166	ASP	2.1
4	J	213	PHE	2.1
1	F	224	GLN	2.1
4	J	216	PRO	2.1
5	I	192	GLU	2.1
5	I	104	CYS	2.0
1	A	51	TRP	2.0
5	L	85	VAL	2.0
1	A	36	PHE	2.0
5	L	105	ALA	2.0
1	F	247	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](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
6	IOD	K	302	1/1	0.85	0.10	80,80,80,80	1
7	NA	G	102	1/1	0.86	0.11	43,43,43,43	0
6	IOD	I	302	1/1	0.91	0.09	106,106,106,106	0
7	NA	L	303	1/1	0.92	0.28	55,55,55,55	0
6	IOD	G	101	1/1	0.93	0.07	90,90,90,90	1
7	NA	E	303	1/1	0.94	0.07	39,39,39,39	0
6	IOD	L	301	1/1	0.95	0.07	89,89,89,89	1
6	IOD	L	302	1/1	0.98	0.03	60,60,60,60	0
6	IOD	L	304	1/1	0.98	0.04	60,60,60,60	1
6	IOD	J	301	1/1	0.98	0.06	69,69,69,69	1

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
6	IOD	D	301	1/1	0.98	0.07	81,81,81,81	0
6	IOD	E	301	1/1	0.98	0.03	57,57,57,57	0
6	IOD	I	301	1/1	0.99	0.03	76,76,76,76	1
6	IOD	K	301	1/1	0.99	0.04	51,51,51,51	1
6	IOD	E	302	1/1	0.99	0.03	67,67,67,67	1

6.5 Other polymers [i](#)

There are no such residues in this entry.