



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 3, 2017 – 09:07 PM EDT

PDB ID : 4QRP
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 : unknown
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

<http://wwpdb.org/validation/2016/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)	:	1.9-1692
EDS	:	rb-20030345
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	rb-20030345

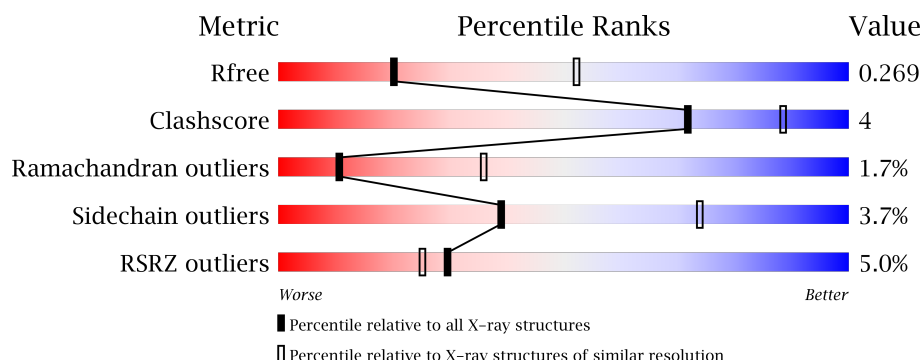
1 Overall quality at a glance

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.





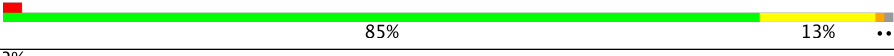

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	100719	1586 (2.90-2.90)
Clashscore	112137	1807 (2.90-2.90)
Ramachandran outliers	110173	1768 (2.90-2.90)
Sidechain outliers	110143	1770 (2.90-2.90)
RSRZ outliers	101464	1596 (2.90-2.90)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	<div> <div>7%</div> <div>85%</div> <div>14%</div> <div>.</div> </div>
1	F	276	<div> <div>9%</div> <div>90%</div> <div>9%</div> <div>.</div> </div>
2	B	100	<div> <div>13%</div> <div>83%</div> <div>12%</div> <div>...</div> </div>
2	G	100	<div> <div>%</div> <div>94%</div> <div>6%</div> </div>
3	C	9	<div> <div>100%</div> </div>

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Mol	Chain	Length	Quality of chain
3	H	9	 78%22%
4	D	206	 5%84%13%..
4	J	206	 7%88%10%.
4	K	206	 2%86%11%..
5	E	245	 2%90%8%..
5	I	245	 2%85%13%..
5	L	245	 2%82%16%..

2 Entry composition

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 (three-letter code: IOD) (formula: I).

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

- Molecule 7 is SODIUM ION (three-letter code: NA) (formula: Na).

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

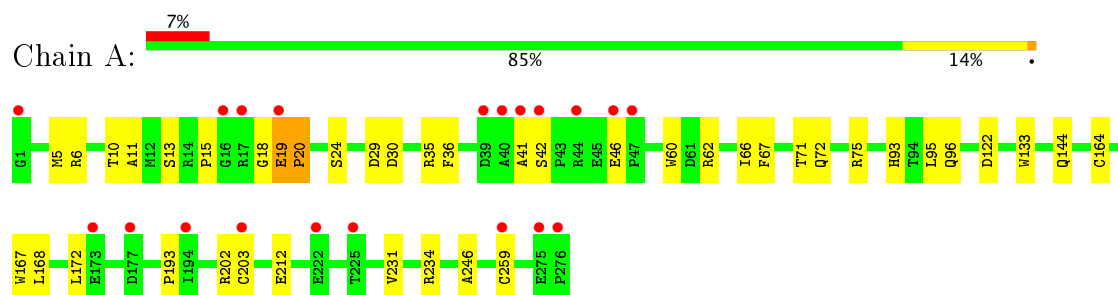
- Molecule 8 is water.

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

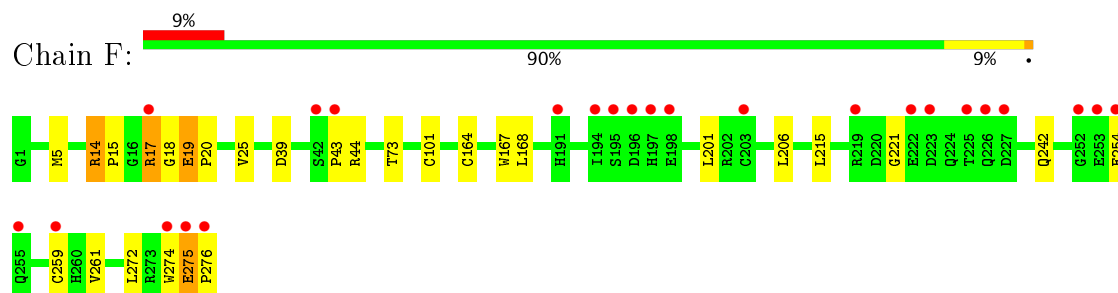
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA 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.

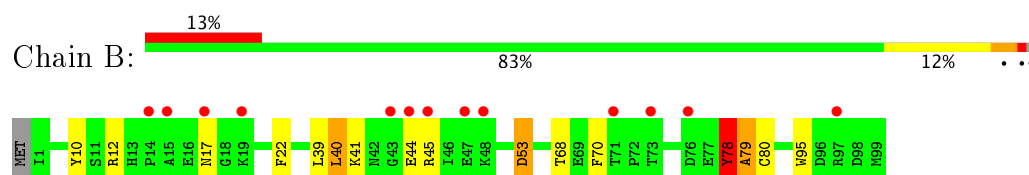
- Molecule 1: HLA class I histocompatibility antigen, B-8 alpha chain



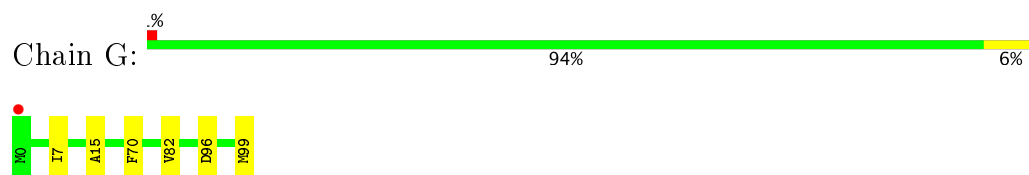
- Molecule 1: HLA class I histocompatibility antigen, B-8 alpha chain



- Molecule 2: Beta-2-microglobulin



- Molecule 2: Beta-2-microglobulin




- Molecule 3: NS3-4A protein




There are no outlier residues recorded for this chain.

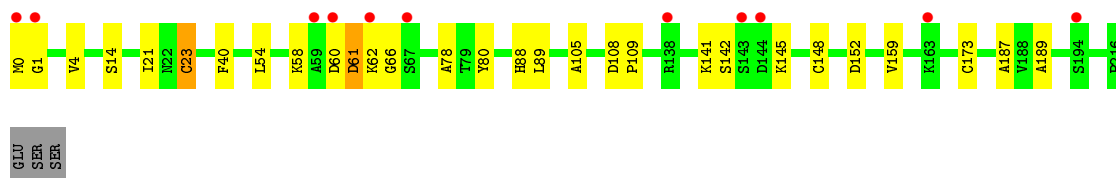
- Molecule 3: NS3-4A protein

Chain H:  78% 22%




- Molecule 4: DD31 TCR alpha chain

Chain D:  5% 84% 13% ..




- Molecule 4: DD31 TCR alpha chain

Chain J:  7% 88% 10% ..

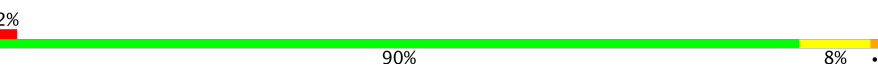


- Molecule 4: DD31 TCR alpha chain

Chain K:  2% 86% 11% ..




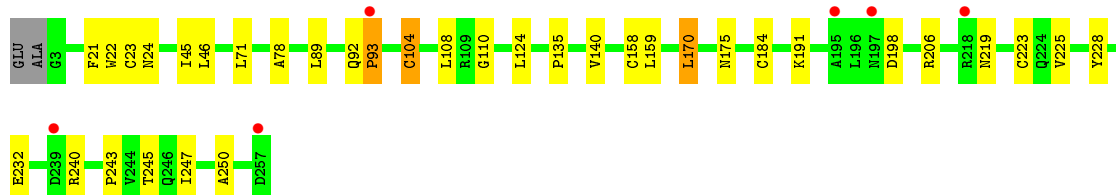
- Molecule 5: DD31 TCR beta chain

Chain E:  2% 90% 8% ..

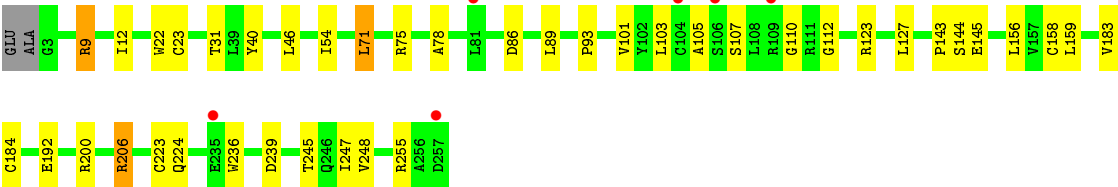
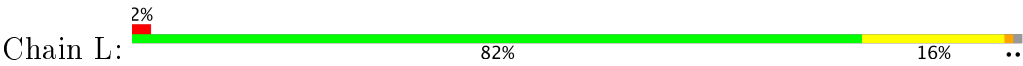


- Molecule 5: DD31 TCR beta chain

Chain I:  2% 85% 13% ..



● 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	74.16 Å 252.19 Å 79.45 Å 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) 99.9 (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 2.88 Å)	Xtriage
Refinement program	BUSTER-TNT BUSTER 2.10.0, BUSTER 2.10.0	Depositor
R, R_{free}	0.206 , 0.248 0.231 , 0.269	Depositor DCC
R_{free} test set	3179 reflections (5.07%)	DCC
Wilson B-factor (Å ²)	49.0	Xtriage
Anisotropy	0.297	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 39.1	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 [i](#)

5.1 Standard geometry [i](#)

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	0.37	0/2313	0.60	1/3146 (0.0%)
1	F	0.36	0/2313	0.61	0/3146
2	B	0.41	0/852	0.74	1/1152 (0.1%)
2	G	0.36	0/860	0.61	0/1162
3	C	0.32	0/75	0.52	0/95
3	H	0.34	0/75	0.56	0/95
4	D	0.37	0/1608	0.63	0/2182
4	J	0.38	0/1608	0.61	0/2182
4	K	0.37	0/1613	0.65	1/2189 (0.0%)
5	E	0.36	0/1988	0.62	0/2702
5	I	0.35	0/1988	0.61	0/2702
5	L	0.35	0/1988	0.58	0/2702
All	All	0.36	0/17281	0.62	3/23455 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	78	TYR	C-N-CA	7.88	141.40	121.70
4	K	143	ASP	CB-CG-OD2	5.20	122.98	118.30
1	A	20	PRO	N-CA-C	5.17	125.55	112.10

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	26	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
3	H	75	0	79	1	0
4	D	1573	0	1498	10	0
4	J	1573	0	1498	8	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	18	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:OE1	1.43	1.11
1:F:254:GLU:HB2	1:F:275:GLU:CD	1.76	1.06
1:F:201:LEU:CD1	1:F:275:GLU:OE2	2.03	1.05
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
2:B:78:TYR:HB2	2:B:79:ALA:CB	1.98	0.92
1:A:19:GLU:CD	1:A:75:ARG:HG2	1.91	0.89
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
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
5:L:9:ARG:H	5:L:9:ARG:HE	1.41	0.68
1:F:254:GLU:CG	1:F:275:GLU:OE1	2.42	0.67
1:F:254:GLU:CD	1:F:275:GLU:OE1	2.34	0.66
1:F:201:LEU:CD2	1:F:275:GLU:HG2	2.21	0.65
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
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:19:GLU:CD	1:A:75:ARG:CG	2.71	0.58
1:F:201:LEU:HD21	1:F:275:GLU:CD	2.24	0.58
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:L:143:PRO:HD3	5:L:156:LEU:HG	1.86	0.57
4:K:22:CYS:HG	4:K:103:CYS:HG	0.63	0.57
5:E:158:CYS:CB	5:E:223:CYS:HG	2.18	0.56
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
5:I:175:ASN:HD21	5:I:219:ASN:HD22	1.54	0.54
2:G:7:ILE:HG12	2:G:82:VAL:HG21	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:19:GLU:OE2	1:A:75:ARG:HB3	2.08	0.53
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
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:92:GLN:O	5:E:93:PRO:C	2.48	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:167:TRP:HZ2	4:D:0:MET:HB2	1.75	0.51
1:A:35:ARG:CZ	2:B:53:ASP:HB2	2.40	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
5:L:158:CYS:HG	5:L:223:CYS:HG	1.53	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
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
1:A:19:GLU:CD	1:A:75:ARG:HH11	2.09	0.48
5:I:228:TYR:HA	5:I:245:THR:HG23	1.95	0.48
5:L:158:CYS:SG	5:L:223:CYS:SG	3.09	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:158:CYS:CB	5:L:223:CYS:HG	2.28	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:192:GLU:HG3	5:L:200:ARG:HB2	1.96	0.46
5:L:184:CYS:HB3	5:L:206:ARG:HB2	1.97	0.46
1:A:19:GLU:HG2	1:A:75:ARG:CG	2.31	0.46
1:F:254:GLU:CB	1:F:275:GLU:CD	2.52	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:E:247:ILE:HD13	1:F:15:PRO:HB2	1.98	0.45
1:F:254:GLU:HB2	1:F:275:GLU:OE2	2.14	0.45
1:A:203:CYS:HG	1:A:259:CYS:HG	1.65	0.45
5:L:75:ARG:HA	5:L:93:PRO:HD2	1.99	0.45
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:I:92:GLN:O	5:I:93:PRO:C	2.54	0.44
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
1:A:18:GLY:C	1:A:19:GLU:HG3	2.38	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
2:B:40:LEU:HD11	2:B:44:GLU:C	2.41	0.41
5:E:247:ILE:H	1:F:17:ARG:HE	1.68	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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 [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%)	249 (91%)	20 (7%)	5 (2%)	10	34
1	F	274/276 (99%)	249 (91%)	18 (7%)	7 (3%)	6	24
2	B	97/100 (97%)	83 (86%)	10 (10%)	4 (4%)	3	13
2	G	98/100 (98%)	94 (96%)	3 (3%)	1 (1%)	18	51
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%)	12	39
4	J	201/206 (98%)	189 (94%)	10 (5%)	2 (1%)	18	51
4	K	202/206 (98%)	181 (90%)	18 (9%)	3 (2%)	12	39
5	E	241/245 (98%)	222 (92%)	16 (7%)	3 (1%)	15	46
5	I	241/245 (98%)	217 (90%)	19 (8%)	5 (2%)	8	30
5	L	241/245 (98%)	224 (93%)	14 (6%)	3 (1%)	15	46
All	All	2084/2123 (98%)	1911 (92%)	137 (7%)	36 (2%)	11	36

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
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%)	46	80
1	F	235/235 (100%)	232 (99%)	3 (1%)	73	93
2	B	94/95 (99%)	90 (96%)	4 (4%)	33	68
2	G	95/95 (100%)	94 (99%)	1 (1%)	78	94
3	C	9/9 (100%)	9 (100%)	0	100	100
3	H	9/9 (100%)	8 (89%)	1 (11%)	7	21
4	D	179/182 (98%)	169 (94%)	10 (6%)	25	57
4	J	179/182 (98%)	172 (96%)	7 (4%)	37	72
4	K	178/182 (98%)	173 (97%)	5 (3%)	49	82
5	E	210/211 (100%)	202 (96%)	8 (4%)	38	73
5	I	210/211 (100%)	201 (96%)	9 (4%)	33	68
5	L	210/211 (100%)	196 (93%)	14 (7%)	19	48
All	All	1843/1857 (99%)	1774 (96%)	69 (4%)	39	74

All (69) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	ARG
1	A	30	ASP
1	A	122	ASP
1	A	164	CYS
1	A	172	LEU
1	A	231	VAL
1	A	234	ARG
2	B	40	LEU
2	B	41	LYS
2	B	53	ASP
2	B	70	PHE
4	D	4	VAL
4	D	14	SER
4	D	23	CYS
4	D	60	ASP
4	D	61	ASP
4	D	80	TYR
4	D	141	LYS
4	D	145	LYS
4	D	152	ASP

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Mol	Chain	Res	Type
4	D	173	CYS
5	E	13	ILE
5	E	109	ARG
5	E	130	LEU
5	E	159	LEU
5	E	161	THR
5	E	183	VAL
5	E	196	LEU
5	E	206	ARG
1	F	14	ARG
1	F	25	VAL
1	F	39	ASP
2	G	70	PHE
3	H	8	GLU
5	I	71	LEU
5	I	104	CYS
5	I	108	LEU
5	I	159	LEU
5	I	170	LEU
5	I	191	LYS
5	I	198	ASP
5	I	206	ARG
5	I	232	GLU
4	J	4	VAL
4	J	23	CYS
4	J	60	ASP
4	J	82	LYS
4	J	94	VAL
4	J	106	LEU
4	J	169	ILE
4	K	1	ASP
4	K	49	LEU
4	K	79	TYR
4	K	151	ASP
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

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Mol	Chain	Res	Type
5	L	145	GLU
5	L	159	LEU
5	L	183	VAL
5	L	206	ARG
5	L	239	ASP
5	L	248	VAL
5	L	255	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	54	GLN
1	A	174	ASN
2	B	8	GLN
2	B	84	HIS
4	D	22	ASN
4	D	44	GLN
4	D	88	HIS
4	D	204	ASN
5	E	44	GLN
5	E	57	GLN
5	I	43	GLN
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	44	GLN
4	J	68	ASN
4	K	43	GLN
4	K	157	ASN
4	K	199	ASN
5	L	215	GLN
5	L	238	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 carbohydrates 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%)	0.45	20 (7%) 16 11	39, 72, 101, 124	0
1	F	276/276 (100%)	0.07	24 (8%) 11 8	22, 48, 105, 124	0
2	B	99/100 (99%)	0.79	13 (13%) 4 2	58, 85, 115, 119	0
2	G	100/100 (100%)	-0.27	1 (1%) 82 81	26, 44, 66, 73	0
3	C	9/9 (100%)	-0.41	0 100 100	53, 58, 62, 62	0
3	H	9/9 (100%)	-0.57	0 100 100	36, 39, 49, 53	0
4	D	203/206 (98%)	0.09	11 (5%) 26 22	31, 54, 83, 96	0
4	J	203/206 (98%)	0.39	14 (6%) 18 13	36, 61, 93, 109	0
4	K	202/206 (98%)	-0.11	4 (1%) 65 62	29, 47, 68, 81	0
5	E	243/245 (99%)	-0.12	6 (2%) 58 53	22, 48, 82, 93	1 (0%)
5	I	243/245 (99%)	0.19	6 (2%) 58 53	29, 63, 98, 117	0
5	L	243/245 (99%)	0.16	6 (2%) 58 53	28, 62, 92, 104	0
All	All	2106/2123 (99%)	0.15	105 (4%) 30 25	22, 58, 98, 124	1 (0%)

All (105) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
4	D	59	ALA	6.7
4	J	194	SER	6.1
4	J	59	ALA	5.8
2	B	76	ASP	5.4
4	J	1	GLY	5.2
4	K	143	ASP	5.2
1	A	17	ARG	4.9
2	B	73	THR	4.7
1	A	41	ALA	4.6
1	F	17	ARG	4.3
1	F	275	GLU	4.3

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Mol	Chain	Res	Type	RSRZ
5	L	257	ASP	4.3
1	A	19	GLU	4.3
1	A	1	GLY	4.2
4	D	62	LYS	4.0
5	I	257	ASP	4.0
5	E	237	THR	4.0
1	A	42	SER	3.9
5	E	239	ASP	3.8
5	L	81	LEU	3.7
1	A	44	ARG	3.7
1	A	225	THR	3.7
2	B	44	GLU	3.6
1	F	43	PRO	3.5
4	D	0	MET	3.5
1	A	47	PRO	3.5
1	F	274	TRP	3.5
4	J	60	ASP	3.5
4	J	138	ARG	3.4
4	K	141	SER	3.4
4	D	143	SER	3.2
5	L	235	GLU	3.2
2	B	14	PRO	3.2
1	F	227	ASP	3.2
1	A	39	ASP	3.1
4	D	1	GLY	3.1
2	B	19	LYS	3.0
4	J	193	LYS	3.0
1	F	223	ASP	3.0
1	F	254	GLU	3.0
2	B	15	ALA	3.0
2	B	47	GLU	3.0
4	D	163	LYS	3.0
1	A	275	GLU	3.0
4	D	144	ASP	2.9
5	I	197	ASN	2.9
2	B	97	ARG	2.8
1	A	203	CYS	2.8
1	A	16	GLY	2.8
4	J	142	SER	2.8
5	I	239	ASP	2.7
4	D	60	ASP	2.7
5	I	93	PRO	2.7

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Mol	Chain	Res	Type	RSRZ
5	L	104	CYS	2.6
4	J	216	PRO	2.6
2	B	45	ARG	2.6
4	D	67	SER	2.6
5	E	195	ALA	2.6
5	E	257	ASP	2.6
5	I	195	ALA	2.6
4	J	0	MET	2.5
5	I	218	ARG	2.5
1	F	198	GLU	2.5
1	F	42	SER	2.5
2	B	17	ASN	2.5
1	F	225	THR	2.5
4	D	194	SER	2.5
1	F	196	ASP	2.5
2	G	0	MET	2.5
1	F	195	SER	2.4
1	F	197	HIS	2.4
1	A	276	PRO	2.4
4	K	142	SER	2.4
1	F	194	ILE	2.4
5	L	109	ARG	2.4
1	F	259	CYS	2.3
5	L	106	SER	2.3
1	F	203	CYS	2.3
1	A	46	GLU	2.3
4	K	194	ASP	2.3
1	A	194	ILE	2.2
1	A	222	GLU	2.2
1	F	253	GLU	2.2
1	F	226	GLN	2.2
1	F	219	ARG	2.2
5	E	16	ARG	2.2
1	A	40	ALA	2.2
1	F	191	HIS	2.1
2	B	43	GLY	2.1
2	B	71	THR	2.1
4	J	57	THR	2.1
1	F	222	GLU	2.1
4	J	163	LYS	2.1
1	F	276	PRO	2.1
4	J	56	ALA	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	259	CYS	2.1
1	F	255	GLN	2.1
1	A	173	GLU	2.1
2	B	48	LYS	2.1
4	D	138	ARG	2.0
1	F	252	GLY	2.0
1	A	177	ASP	2.0
4	J	84	THR	2.0
5	E	233	ASN	2.0
4	J	195	ASP	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](#)

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
6	IOD	E	302	1/1	0.98	0.05	-1.36	67,67,67,67	1
6	IOD	L	304	1/1	0.98	0.06	-2.09	60,60,60,60	1
6	IOD	E	301	1/1	0.99	0.06	-2.96	57,57,57,57	0
6	IOD	I	301	1/1	0.99	0.02	-3.03	76,76,76,76	1
6	IOD	L	301	1/1	0.95	0.08	-4.03	89,89,89,89	1
6	IOD	L	302	1/1	0.99	0.05	-4.74	60,60,60,60	0
6	IOD	I	302	1/1	0.94	0.07	-5.43	106,106,106,106	0
6	IOD	J	301	1/1	0.98	0.04	-5.44	69,69,69,69	1
6	IOD	K	301	1/1	0.99	0.04	-5.72	51,51,51,51	1
6	IOD	D	301	1/1	0.98	0.04	-6.02	81,81,81,81	0
6	IOD	G	101	1/1	0.90	0.08	-	90,90,90,90	1

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
6	IOD	K	302	1/1	0.95	0.13	-	80,80,80,80	1
7	NA	L	303	1/1	0.97	0.49	-	55,55,55,55	0
7	NA	E	303	1/1	0.92	0.20	-	39,39,39,39	0
7	NA	G	102	1/1	0.97	0.12	-	43,43,43,43	0

6.5 Other polymers [i](#)

There are no such residues in this entry.