



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 04:00 PM UTC

PDB ID : 4JRX / pdb_00004jrx
Title : Crystal Structure of CA5 TCR-HLA B*3505-LPEP complex
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Deposited on : 2013-03-22
Resolution : 2.30 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

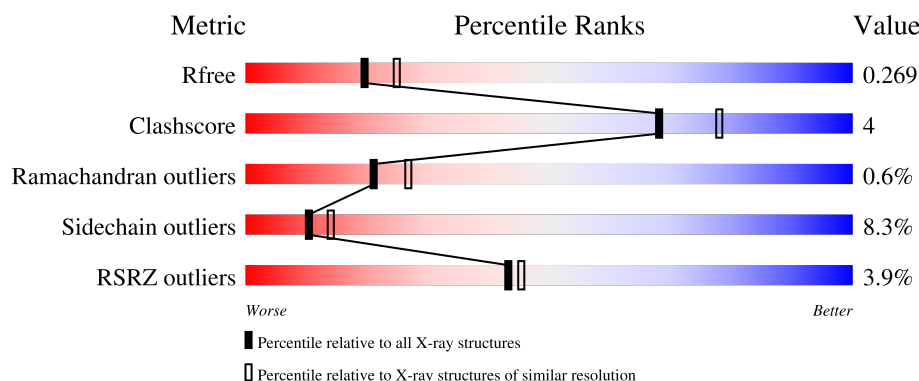
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	5% 85% 12% .
2	B	99	3% 85% 13% .
3	C	13	92% 8%
4	D	204	5% 82% 14% ..
5	E	238	2% 83% 15% .

2 Entry composition [i](#)

There are 8 unique types of molecules in this entry. The entry contains 6865 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 MHC class I antigen.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	276	Total	C	N	O	S	0	0	0
			2257	1405	414	431	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			

- Molecule 3 is a protein called Trans-activator protein BZLF1.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	13	Total	C	N	O	0	0	0
			101	66	15	20			

- Molecule 4 is a protein called CA5 TCR alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	204	Total	C	N	O	S	0	1	0
			1646	1041	266	332	7			

- Molecule 5 is a protein called CA5 TCR beta chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	E	238	Total	C	N	O	S	0	0	0
			1889	1185	329	366	9			

- Molecule 6 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	1	Total	Na	0	0
			1	1		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	D	1	Total 1	Na 1	0	0
6	E	1	Total 1	Na 1	0	0

- Molecule 7 is IODIDE ION (CCD ID: IOD) (formula: I).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	3	Total 3	I 3	0	0
7	C	1	Total 1	I 1	0	0
7	D	3	Total 3	I 3	0	0
7	E	1	Total 1	I 1	0	0

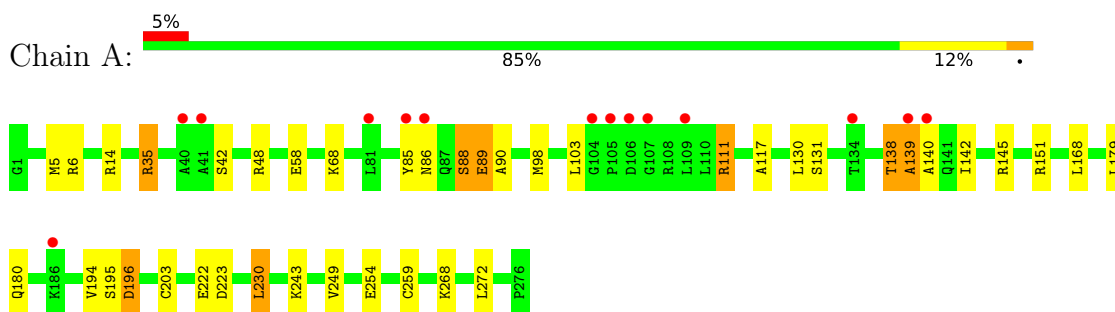
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	37	Total 37	O 37	0	0
8	B	18	Total 18	O 18	0	0
8	C	3	Total 3	O 3	0	0
8	D	36	Total 36	O 36	0	0
8	E	38	Total 38	O 38	0	0

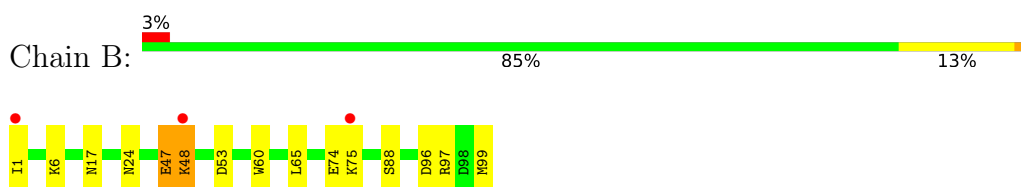
3 Residue-property plots [i](#)

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.

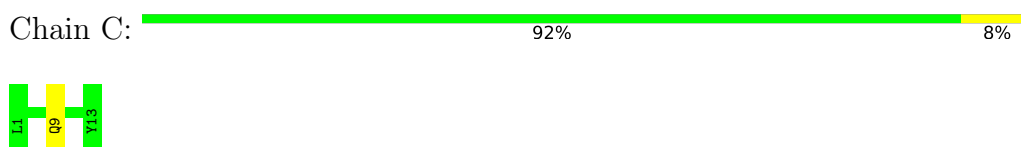
- Molecule 1: MHC class I antigen



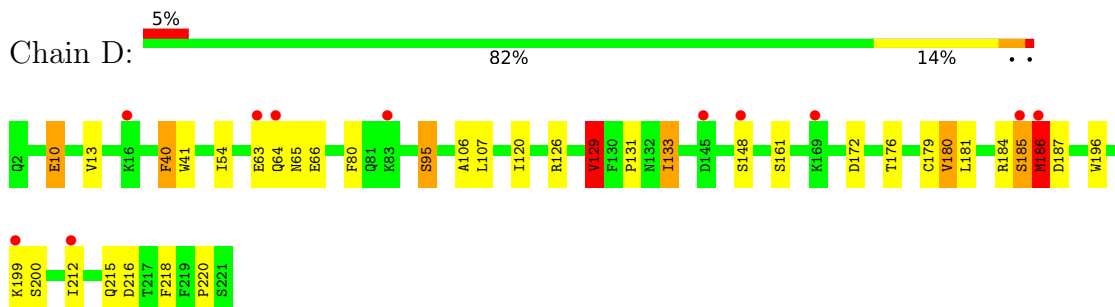
- Molecule 2: Beta-2-microglobulin



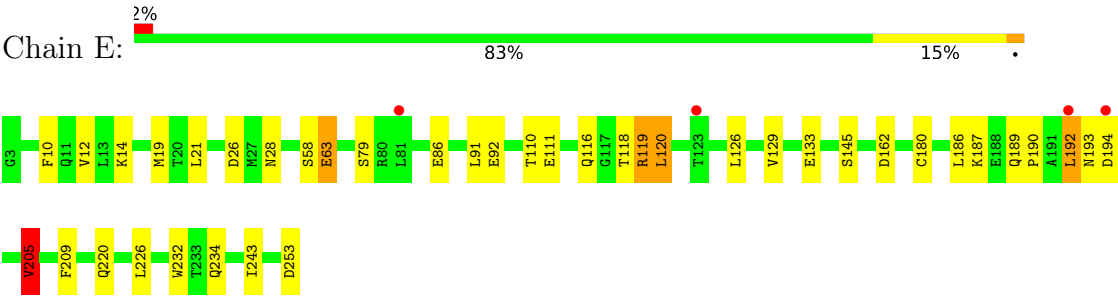
- Molecule 3: Trans-activator protein BZLF1



- Molecule 4: CA5 TCR alpha chain



- Molecule 5: CA5 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	55.10Å 78.41Å 105.34Å 90.00° 93.06° 90.00°	Depositor
Resolution (Å)	19.80 – 2.30 19.80 – 2.30	Depositor EDS
% Data completeness (in resolution range)	(Not available) (19.80-2.30) 99.8 (19.80-2.30)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.20 (at 2.30Å)	Xtriage
Refinement program	BUSTER-TNT, BUSTER	Depositor
R, R_{free}	0.205 , 0.257 0.219 , 0.269	Depositor DCC
R_{free} test set	2001 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	36.6	Xtriage
Anisotropy	0.567	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 35.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6865	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.77% 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: IOD, NA

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	0.75	0/2320	1.29	9/3153 (0.3%)
2	B	0.68	0/852	1.17	2/1152 (0.2%)
3	C	0.71	0/104	1.00	0/142
4	D	0.73	0/1684	1.22	13/2287 (0.6%)
5	E	0.71	0/1939	1.20	6/2638 (0.2%)
All	All	0.72	0/6899	1.23	30/9372 (0.3%)

There are no bond length outliers.

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	140	ALA	N-CA-C	-9.74	101.40	113.38
5	E	126	LEU	CA-C-N	7.34	131.09	120.38
5	E	126	LEU	C-N-CA	7.34	131.09	120.38
1	A	89	GLU	CA-C-N	6.71	129.50	120.65
1	A	89	GLU	C-N-CA	6.71	129.50	120.65
4	D	40	PHE	CA-CB-CG	6.20	120.00	113.80
5	E	86	GLU	N-CA-C	6.16	119.86	109.76
5	E	205	VAL	N-CA-CB	-5.99	104.14	112.52
5	E	116	GLN	N-CA-C	5.68	118.20	111.33
5	E	63	GLU	N-CA-C	5.64	118.16	111.33
1	A	138	THR	N-CA-C	-5.61	101.39	110.32
4	D	129	VAL	N-CA-CB	5.60	117.71	110.99
1	A	145	ARG	CA-C-N	5.47	127.61	120.28
1	A	145	ARG	C-N-CA	5.47	127.61	120.28
4	D	180	VAL	N-CA-C	5.42	115.91	107.99
1	A	142	ILE	N-CA-C	-5.36	105.49	110.53
4	D	186	MET	CA-C-N	5.32	129.96	122.46
4	D	186	MET	C-N-CA	5.32	129.96	122.46
2	B	74	GLU	CA-C-N	5.20	127.70	120.63

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	74	GLU	C-N-CA	5.20	127.70	120.63
4	D	199	LYS	CA-C-N	5.19	127.75	120.28
4	D	199	LYS	C-N-CA	5.19	127.75	120.28
4	D	66	GLU	N-CA-C	5.10	118.39	110.17
1	A	254	GLU	N-CA-C	5.10	117.50	111.33
4	D	66	GLU	CA-C-N	5.09	130.07	122.99
4	D	66	GLU	C-N-CA	5.09	130.07	122.99
4	D	95	SER	N-CA-C	5.05	117.79	110.42
1	A	85	TYR	N-CA-C	-5.03	100.59	108.63
4	D	185	SER	CA-C-N	5.01	131.12	121.54
4	D	185	SER	C-N-CA	5.01	131.12	121.54

There are no chirality outliers.

There are no planarity outliers.

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	2257	0	2121	12	0
2	B	829	0	796	7	0
3	C	101	0	102	1	0
4	D	1646	0	1555	13	0
5	E	1889	0	1789	15	0
6	A	1	0	0	0	0
6	D	1	0	0	0	0
6	E	1	0	0	0	0
7	A	3	0	0	1	0
7	C	1	0	0	0	0
7	D	3	0	0	1	0
7	E	1	0	0	0	0
8	A	37	0	0	0	0
8	B	18	0	0	0	0
8	C	3	0	0	0	0
8	D	36	0	0	0	0
8	E	38	0	0	0	0
All	All	6865	0	6363	43	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 (43) 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:161:SER:HB2	7:D:303:IOD:I	2.63	0.69
4:D:65:ASN:HD21	4:D:80:PHE:H	1.47	0.60
5:E:19:MET:SD	5:E:120:LEU:HD11	2.43	0.58
4:D:181:LEU:HB3	5:E:180:CYS:HB2	1.85	0.58
4:D:40:PHE:HB2	4:D:106:ALA:HB3	1.86	0.58
2:B:47:GLU:O	2:B:48:LYS:HG2	2.04	0.57
2:B:17:ASN:ND2	2:B:97:ARG:HH12	2.03	0.57
5:E:12:VAL:HG11	5:E:226:LEU:HD12	1.86	0.56
4:D:41:TRP:HB2	4:D:54:ILE:HG22	1.89	0.54
1:A:35:ARG:HG2	1:A:48:ARG:HD3	1.91	0.52
4:D:185:SER:O	4:D:186:MET:HB2	2.09	0.52
4:D:131:PRO:HG2	4:D:133:ILE:HD11	1.94	0.50
4:D:10:GLU:HG3	4:D:126:ARG:HB3	1.94	0.49
1:A:6:ARG:HD2	1:A:98:MET:SD	2.52	0.49
1:A:117:ALA:HB2	2:B:60:TRP:CE2	2.47	0.49
1:A:195:SER:OG	1:A:196:ASP:N	2.44	0.49
4:D:65:ASN:ND2	4:D:80:PHE:H	2.10	0.48
1:A:111:ARG:HE	1:A:111:ARG:HB2	1.53	0.47
1:A:138:THR:O	1:A:139:ALA:HB3	2.14	0.47
2:B:24:ASN:HB3	2:B:65:LEU:HD11	1.96	0.46
3:C:9:GLN:HG3	5:E:28:ASN:HD22	1.81	0.46
4:D:218:PHE:CE2	4:D:220:PRO:HG3	2.51	0.46
5:E:19:MET:HE3	5:E:19:MET:HB2	1.82	0.46
1:A:5:MET:HB2	1:A:168:LEU:HD13	1.98	0.45
5:E:192:LEU:HD23	5:E:194:ASP:H	1.82	0.44
1:A:35:ARG:NH1	2:B:53:ASP:OD1	2.49	0.44
1:A:88:SER:C	1:A:90:ALA:H	2.25	0.43
5:E:129:VAL:HG11	5:E:226:LEU:HD13	2.00	0.43
5:E:205:VAL:HG13	5:E:209:PHE:HB3	1.99	0.43
4:D:13:VAL:HG13	4:D:129:VAL:HB	2.00	0.43
5:E:220:GLN:HG3	5:E:243:ILE:HG23	2.00	0.43
5:E:110:THR:HG23	5:E:111:GLU:HG2	2.00	0.43
1:A:203:CYS:CB	1:A:259:CYS:SG	3.06	0.42
5:E:21:LEU:HD22	5:E:118:THR:HG21	2.01	0.42
4:D:176:THR:HG23	4:D:196:TRP:HZ3	1.85	0.41
1:A:230:LEU:HD23	1:A:243:LYS:HE3	2.02	0.41
2:B:17:ASN:HD21	2:B:97:ARG:HH12	1.66	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:E:10:PHE:CE1	5:E:119:ARG:HG3	2.56	0.41
1:A:68:LYS:HG2	7:A:304:IOD:I	2.91	0.40
5:E:187:LYS:HD2	5:E:190:PRO:HA	2.03	0.40
5:E:232:TRP:CE2	5:E:234:GLN:HB2	2.56	0.40
2:B:96:ASP:HB3	2:B:99:MET:HG3	2.03	0.40
4:D:176:THR:HG22	5:E:186:LEU:HD11	2.02	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%)	262 (96%)	10 (4%)	2 (1%)	18	23
2	B	97/99 (98%)	95 (98%)	2 (2%)	0	100	100
3	C	11/13 (85%)	10 (91%)	1 (9%)	0	100	100
4	D	203/204 (100%)	194 (96%)	7 (3%)	2 (1%)	12	15
5	E	236/238 (99%)	231 (98%)	4 (2%)	1 (0%)	30	38
All	All	821/830 (99%)	792 (96%)	24 (3%)	5 (1%)	21	27

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
4	D	186	MET
4	D	64	GLN
1	A	139	ALA
5	E	162	ASP
1	A	86	ASN

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	234/234 (100%)	213 (91%)	21 (9%)	9	12
2	B	94/94 (100%)	88 (94%)	6 (6%)	16	23
3	C	11/11 (100%)	11 (100%)	0	100	100
4	D	190/189 (100%)	171 (90%)	19 (10%)	7	9
5	E	206/206 (100%)	190 (92%)	16 (8%)	11	16
All	All	735/734 (100%)	673 (92%)	62 (8%)	10	14

All (62) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	14	ARG
1	A	35	ARG
1	A	42	SER
1	A	58	GLU
1	A	88	SER
1	A	89	GLU
1	A	103	LEU
1	A	111	ARG
1	A	130	LEU
1	A	131	SER
1	A	151	ARG
1	A	179	LEU
1	A	180	GLN
1	A	194	VAL
1	A	196	ASP
1	A	222	GLU
1	A	223	ASP
1	A	230	LEU
1	A	249	VAL
1	A	268	LYS
1	A	272	LEU
2	B	1	ILE
2	B	6	LYS

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Mol	Chain	Res	Type
2	B	47	GLU
2	B	48	LYS
2	B	75	LYS
2	B	88	SER
4	D	10	GLU
4	D	63	GLU
4	D	95	SER
4	D	107	LEU
4	D	120[A]	ILE
4	D	120[B]	ILE
4	D	129	VAL
4	D	133	ILE
4	D	148	SER
4	D	172	ASP
4	D	179	CYS
4	D	180	VAL
4	D	184	ARG
4	D	186	MET
4	D	187	ASP
4	D	200	SER
4	D	212	ILE
4	D	215	GLN
4	D	216	ASP
5	E	14	LYS
5	E	26	ASP
5	E	58	SER
5	E	63	GLU
5	E	79	SER
5	E	91	LEU
5	E	92	GLU
5	E	119	ARG
5	E	120	LEU
5	E	133	GLU
5	E	145	SER
5	E	189	GLN
5	E	192	LEU
5	E	193	ASN
5	E	205	VAL
5	E	253	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	80	ASN
1	A	144	GLN
1	A	180	GLN
1	A	224	GLN
2	B	17	ASN
2	B	24	ASN
3	C	7	GLN
4	D	2	GLN
4	D	65	ASN
4	D	142	GLN
5	E	17	GLN
5	E	22	GLN
5	E	28	ASN
5	E	128	ASN

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 11 ligands modelled in this entry, 11 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%)	0.37	14 (5%) 33 35	23, 38, 54, 73	43 (15%)
2	B	99/99 (100%)	0.07	3 (3%) 52 54	22, 34, 49, 59	14 (14%)
3	C	13/13 (100%)	-0.19	0 100 100	28, 32, 38, 40	0
4	D	204/204 (100%)	0.43	11 (5%) 31 33	16, 37, 57, 75	33 (16%)
5	E	238/238 (100%)	0.23	4 (1%) 69 70	20, 34, 50, 96	26 (10%)
All	All	830/830 (100%)	0.30	32 (3%) 43 45	16, 37, 54, 96	116 (13%)

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
5	E	81	LEU	4.4
1	A	107	GLY	4.0
2	B	1	ILE	3.6
4	D	63	GLU	3.3
4	D	148	SER	3.1
5	E	192	LEU	3.1
1	A	106	ASP	3.0
4	D	185	SER	2.8
4	D	64	GLN	2.7
1	A	104	GLY	2.7
1	A	41	ALA	2.6
1	A	140	ALA	2.6
5	E	123	THR	2.6
4	D	83	LYS	2.4
1	A	105	PRO	2.4
1	A	186	LYS	2.4
4	D	169	LYS	2.3
4	D	186	MET	2.3
4	D	16	LYS	2.3
1	A	134	THR	2.2

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Mol	Chain	Res	Type	RSRZ
2	B	48	LYS	2.2
1	A	40	ALA	2.2
1	A	81	LEU	2.2
1	A	86	ASN	2.2
1	A	109	LEU	2.1
1	A	139	ALA	2.1
4	D	145	ASP	2.1
4	D	199	LYS	2.1
1	A	85	TYR	2.1
2	B	75	LYS	2.1
4	D	212	ILE	2.1
5	E	194	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 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(Å ²)	Q<0.9
7	IOD	D	303	1/1	0.87	0.14	128,128,128,128	0
6	NA	A	301	1/1	0.89	0.17	49,49,49,49	0
7	IOD	E	302	1/1	0.95	0.24	109,109,109,109	1
6	NA	D	301	1/1	0.96	0.07	27,27,27,27	0
6	NA	E	301	1/1	0.97	0.16	30,30,30,30	0
7	IOD	C	101	1/1	0.98	0.04	38,38,38,38	1
7	IOD	D	302	1/1	0.98	0.07	67,67,67,67	1
7	IOD	A	302	1/1	0.98	0.03	37,37,37,37	0
7	IOD	D	304	1/1	0.98	0.04	64,64,64,64	1
7	IOD	A	304	1/1	0.98	0.05	56,56,56,56	0
7	IOD	A	303	1/1	0.99	0.03	44,44,44,44	0

6.5 Other polymers [i](#)

There are no such residues in this entry.