



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 8, 2026 – 12:50 PM UTC

PDB ID : 8K50 / pdb_00008k50
Title : Crystal structure of HLA-A*11:01 in complex with KRAS G12V peptide (VVVGARGVGK)
Authors : Xu, K.
Deposited on : 2023-07-20
Resolution : 2.80 Å(reported)

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

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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

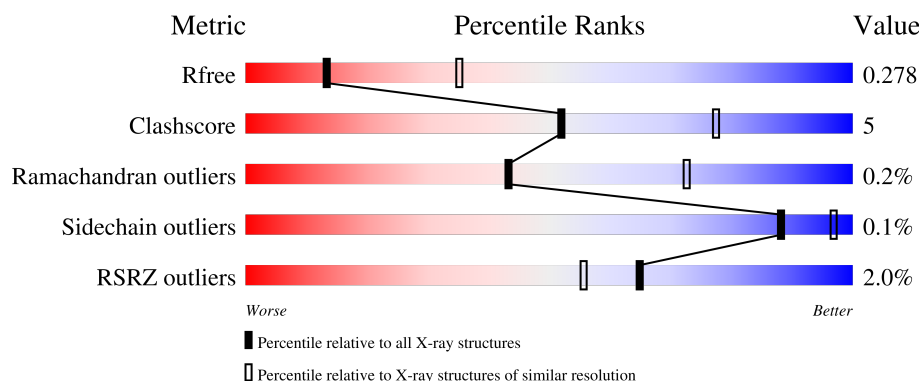
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	% 87% 12%
1	D	276	3% 83% 16%
1	G	276	2% 84% 16%
1	J	276	2% 88% 12%
2	B	122	% 76% 6% 18%

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Mol	Chain	Length	Quality of chain
2	E	122	72%10%18%
2	H	122	2% 68%14%18%
2	K	122	% 66%16%18%
3	C	10	80%20%
3	F	10	80%20%
3	I	10	10% 60%10%10%20%
3	L	10	20% 40%40%20%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 12530 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 alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	275	Total	C	N	O	S	0	0	0
			2240	1392	408	431	9			
1	D	275	Total	C	N	O	S	0	0	0
			2241	1393	408	431	9			
1	G	275	Total	C	N	O	S	0	0	0
			2244	1394	408	433	9			
1	J	275	Total	C	N	O	S	0	0	0
			2241	1393	408	431	9			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	24	MET	-	initiating methionine	UNP P04439
A	33	TYR	PHE	variant	UNP P04439
A	114	ASP	ALA	variant	UNP P04439
A	129	PRO	SER	variant	UNP P04439
D	24	MET	-	initiating methionine	UNP P04439
D	33	TYR	PHE	variant	UNP P04439
D	114	ASP	ALA	variant	UNP P04439
D	129	PRO	SER	variant	UNP P04439
G	24	MET	-	initiating methionine	UNP P04439
G	33	TYR	PHE	variant	UNP P04439
G	114	ASP	ALA	variant	UNP P04439
G	129	PRO	SER	variant	UNP P04439
J	24	MET	-	initiating methionine	UNP P04439
J	33	TYR	PHE	variant	UNP P04439
J	114	ASP	ALA	variant	UNP P04439
J	129	PRO	SER	variant	UNP P04439

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	100	Total	C	N	O	S	0	0	0
			832	530	140	158	4			
2	E	100	Total	C	N	O	S	0	0	0
			836	533	141	158	4			
2	H	100	Total	C	N	O	S	0	0	0
			836	533	141	158	4			
2	K	100	Total	C	N	O	S	0	0	0
			836	533	141	158	4			

There are 92 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	0	MET	-	initiating methionine	UNP P61769
B	100	GLY	-	expression tag	UNP P61769
B	101	SER	-	expression tag	UNP P61769
B	102	GLY	-	expression tag	UNP P61769
B	103	GLY	-	expression tag	UNP P61769
B	104	SER	-	expression tag	UNP P61769
B	105	GLY	-	expression tag	UNP P61769
B	106	ALA	-	expression tag	UNP P61769
B	107	GLY	-	expression tag	UNP P61769
B	108	LEU	-	expression tag	UNP P61769
B	109	ASN	-	expression tag	UNP P61769
B	110	ASP	-	expression tag	UNP P61769
B	111	ILE	-	expression tag	UNP P61769
B	112	PHE	-	expression tag	UNP P61769
B	113	GLU	-	expression tag	UNP P61769
B	114	ALA	-	expression tag	UNP P61769
B	115	GLN	-	expression tag	UNP P61769
B	116	LYS	-	expression tag	UNP P61769
B	117	ILE	-	expression tag	UNP P61769
B	118	GLU	-	expression tag	UNP P61769
B	119	TRP	-	expression tag	UNP P61769
B	120	HIS	-	expression tag	UNP P61769
B	121	GLU	-	expression tag	UNP P61769
E	0	MET	-	initiating methionine	UNP P61769
E	100	GLY	-	expression tag	UNP P61769
E	101	SER	-	expression tag	UNP P61769
E	102	GLY	-	expression tag	UNP P61769
E	103	GLY	-	expression tag	UNP P61769
E	104	SER	-	expression tag	UNP P61769
E	105	GLY	-	expression tag	UNP P61769
E	106	ALA	-	expression tag	UNP P61769

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Chain	Residue	Modelled	Actual	Comment	Reference
E	107	GLY	-	expression tag	UNP P61769
E	108	LEU	-	expression tag	UNP P61769
E	109	ASN	-	expression tag	UNP P61769
E	110	ASP	-	expression tag	UNP P61769
E	111	ILE	-	expression tag	UNP P61769
E	112	PHE	-	expression tag	UNP P61769
E	113	GLU	-	expression tag	UNP P61769
E	114	ALA	-	expression tag	UNP P61769
E	115	GLN	-	expression tag	UNP P61769
E	116	LYS	-	expression tag	UNP P61769
E	117	ILE	-	expression tag	UNP P61769
E	118	GLU	-	expression tag	UNP P61769
E	119	TRP	-	expression tag	UNP P61769
E	120	HIS	-	expression tag	UNP P61769
E	121	GLU	-	expression tag	UNP P61769
H	0	MET	-	initiating methionine	UNP P61769
H	100	GLY	-	expression tag	UNP P61769
H	101	SER	-	expression tag	UNP P61769
H	102	GLY	-	expression tag	UNP P61769
H	103	GLY	-	expression tag	UNP P61769
H	104	SER	-	expression tag	UNP P61769
H	105	GLY	-	expression tag	UNP P61769
H	106	ALA	-	expression tag	UNP P61769
H	107	GLY	-	expression tag	UNP P61769
H	108	LEU	-	expression tag	UNP P61769
H	109	ASN	-	expression tag	UNP P61769
H	110	ASP	-	expression tag	UNP P61769
H	111	ILE	-	expression tag	UNP P61769
H	112	PHE	-	expression tag	UNP P61769
H	113	GLU	-	expression tag	UNP P61769
H	114	ALA	-	expression tag	UNP P61769
H	115	GLN	-	expression tag	UNP P61769
H	116	LYS	-	expression tag	UNP P61769
H	117	ILE	-	expression tag	UNP P61769
H	118	GLU	-	expression tag	UNP P61769
H	119	TRP	-	expression tag	UNP P61769
H	120	HIS	-	expression tag	UNP P61769
H	121	GLU	-	expression tag	UNP P61769
K	0	MET	-	initiating methionine	UNP P61769
K	100	GLY	-	expression tag	UNP P61769
K	101	SER	-	expression tag	UNP P61769
K	102	GLY	-	expression tag	UNP P61769

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Chain	Residue	Modelled	Actual	Comment	Reference
K	103	GLY	-	expression tag	UNP P61769
K	104	SER	-	expression tag	UNP P61769
K	105	GLY	-	expression tag	UNP P61769
K	106	ALA	-	expression tag	UNP P61769
K	107	GLY	-	expression tag	UNP P61769
K	108	LEU	-	expression tag	UNP P61769
K	109	ASN	-	expression tag	UNP P61769
K	110	ASP	-	expression tag	UNP P61769
K	111	ILE	-	expression tag	UNP P61769
K	112	PHE	-	expression tag	UNP P61769
K	113	GLU	-	expression tag	UNP P61769
K	114	ALA	-	expression tag	UNP P61769
K	115	GLN	-	expression tag	UNP P61769
K	116	LYS	-	expression tag	UNP P61769
K	117	ILE	-	expression tag	UNP P61769
K	118	GLU	-	expression tag	UNP P61769
K	119	TRP	-	expression tag	UNP P61769
K	120	HIS	-	expression tag	UNP P61769
K	121	GLU	-	expression tag	UNP P61769

- Molecule 3 is a protein called KRAS G12V peptide (VVVGAVGVGK).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	10	Total	C	N	O	0	0	0
			62	40	11	11			
3	F	10	Total	C	N	O	0	0	0
			62	40	11	11			
3	I	8	Total	C	N	O	0	0	0
			50	32	9	9			
3	L	8	Total	C	N	O	0	0	0
			50	32	9	9			

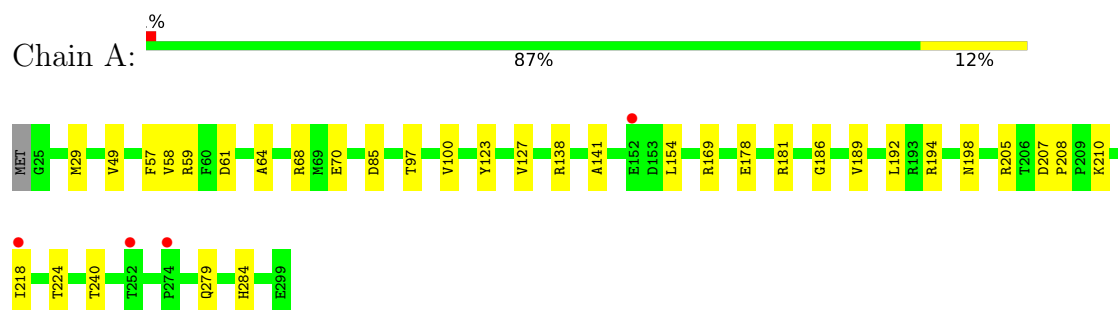
There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	6	VAL	GLY	engineered mutation	UNP P01116
F	6	VAL	GLY	engineered mutation	UNP P01116
I	6	VAL	GLY	engineered mutation	UNP P01116
L	6	VAL	GLY	engineered mutation	UNP P01116

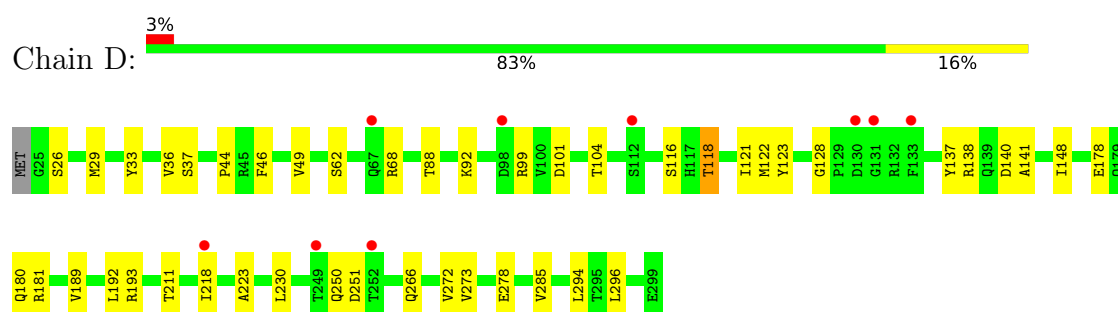
3 Residue-property plots [i](#)

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.

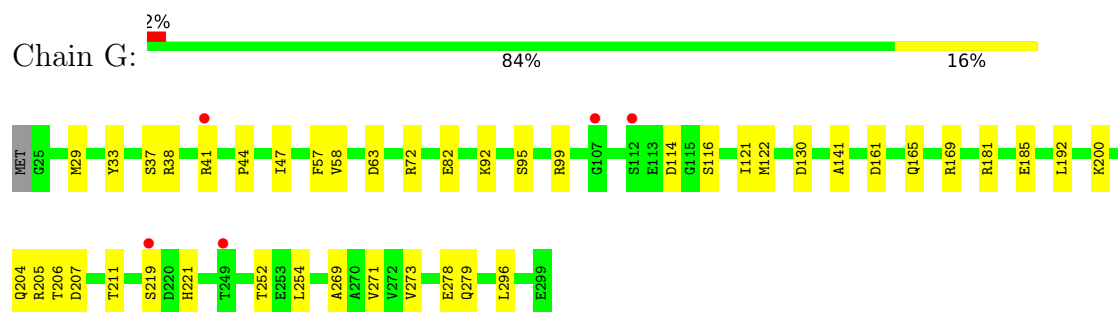
- Molecule 1: HLA class I histocompatibility antigen, A alpha chain



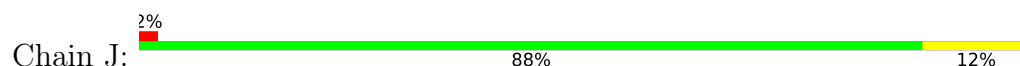
- Molecule 1: HLA class I histocompatibility antigen, A alpha chain



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



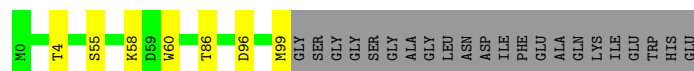
- Molecule 1: HLA class I histocompatibility antigen, A alpha chain





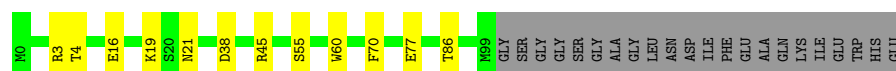
- Molecule 2: Beta-2-microglobulin

Chain B:



- Molecule 2: Beta-2-microglobulin

Chain E:



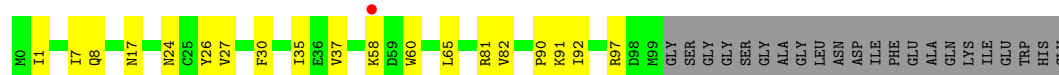
- Molecule 2: Beta-2-microglobulin

Chain H:



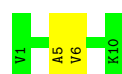
- Molecule 2: Beta-2-microglobulin

Chain K:



- Molecule 3: KRAS G12V peptide (VVVGAVGVGK)

Chain C:

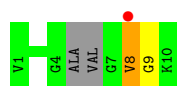


- Molecule 3: KRAS G12V peptide (VVVGAVGVGK)

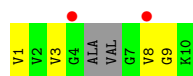
Chain F:



- Molecule 3: KRAS G12V peptide (VVVGAVGVGK)



- Molecule 3: KRAS G12V peptide (VVVGAVGVGK)



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	116.29Å 121.84Å 133.38Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	44.98 – 2.80 44.98 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.8 (44.98-2.80) 92.5 (44.98-2.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.79 (at 2.81Å)	Xtriage
Refinement program	PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.223 , 0.281 0.224 , 0.278	Depositor DCC
R_{free} test set	2000 reflections (4.21%)	wwPDB-VP
Wilson B-factor (Å ²)	43.7	Xtriage
Anisotropy	0.900	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 32.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.000 for k,h,-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	12530	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 29.42 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.5539e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² Theoretical 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	0.13	0/2301	0.33	0/3124
1	D	0.17	0/2302	0.34	0/3125
1	G	0.11	0/2305	0.32	0/3129
1	J	0.12	0/2302	0.32	0/3125
2	B	0.10	0/855	0.32	0/1158
2	E	0.11	0/859	0.31	0/1162
2	H	0.18	0/859	0.34	0/1162
2	K	0.15	0/859	0.36	0/1162
3	C	0.14	0/61	0.63	0/80
3	F	0.11	0/61	0.36	0/80
3	I	0.12	0/48	0.32	0/60
3	L	0.12	0/48	0.32	0/60
All	All	0.13	0/12860	0.33	0/17427

There are no bond length outliers.

There are no bond angle outliers.

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	2240	0	2084	19	1
1	D	2241	0	2086	30	0
1	G	2244	0	2088	28	0
1	J	2241	0	2086	21	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	832	0	792	5	0
2	E	836	0	803	8	0
2	H	836	0	803	12	0
2	K	836	0	803	13	0
3	C	62	0	74	0	0
3	F	62	0	74	3	0
3	I	50	0	59	1	0
3	L	50	0	59	4	0
All	All	12530	0	11811	123	1

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 5.

The worst 5 of 123 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:J:44:PRO:HD2	1:J:99:ARG:HD3	1.66	0.77
2:K:7:ILE:HD12	2:K:91:LYS:HD2	1.68	0.76
1:D:178:GLU:HG2	1:D:181:ARG:HH22	1.50	0.75
1:J:57:PHE:HA	1:J:72:ARG:O	1.89	0.71
1:D:29:MET:HB2	1:D:192:LEU:HD13	1.73	0.71

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:169:ARG:NH1	1:A:279:GLN:OE1[3_554]	2.01	0.19

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	273/276 (99%)	268 (98%)	5 (2%)	0	100	100
1	D	273/276 (99%)	269 (98%)	4 (2%)	0	100	100
1	G	273/276 (99%)	264 (97%)	9 (3%)	0	100	100
1	J	273/276 (99%)	265 (97%)	8 (3%)	0	100	100
2	B	98/122 (80%)	95 (97%)	3 (3%)	0	100	100
2	E	98/122 (80%)	94 (96%)	4 (4%)	0	100	100
2	H	98/122 (80%)	97 (99%)	1 (1%)	0	100	100
2	K	98/122 (80%)	97 (99%)	1 (1%)	0	100	100
3	C	8/10 (80%)	5 (62%)	1 (12%)	2 (25%)	0	0
3	F	8/10 (80%)	7 (88%)	1 (12%)	0	100	100
3	I	4/10 (40%)	3 (75%)	0	1 (25%)	0	0
3	L	4/10 (40%)	2 (50%)	2 (50%)	0	100	100
All	All	1508/1632 (92%)	1466 (97%)	39 (3%)	3 (0%)	43	72

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	C	5	ALA
3	I	8	VAL
3	C	6	VAL

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	231/233 (99%)	231 (100%)	0	100	100
1	D	231/233 (99%)	230 (100%)	1 (0%)	84	94
1	G	232/233 (100%)	232 (100%)	0	100	100
1	J	231/233 (99%)	231 (100%)	0	100	100
2	B	94/110 (86%)	94 (100%)	0	100	100
2	E	95/110 (86%)	95 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	H	95/110 (86%)	95 (100%)	0	100	100
2	K	95/110 (86%)	95 (100%)	0	100	100
3	C	6/6 (100%)	6 (100%)	0	100	100
3	F	6/6 (100%)	6 (100%)	0	100	100
3	I	5/6 (83%)	5 (100%)	0	100	100
3	L	5/6 (83%)	5 (100%)	0	100	100
All	All	1326/1396 (95%)	1325 (100%)	1 (0%)	88	97

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	118	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 14 such sidechains are listed below:

Mol	Chain	Res	Type
1	D	250	GLN
1	G	180	GLN
2	K	13	HIS
1	J	139	GLN
1	J	248	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](#)

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	275/276 (99%)	0.21	4 (1%) 72 63	33, 45, 73, 85	0
1	D	275/276 (99%)	0.58	9 (3%) 49 39	44, 59, 76, 93	0
1	G	275/276 (99%)	0.23	5 (1%) 67 58	31, 50, 65, 77	0
1	J	275/276 (99%)	0.52	6 (2%) 62 52	38, 57, 75, 84	0
2	B	100/122 (81%)	-0.09	0 100 100	34, 41, 53, 63	0
2	E	100/122 (81%)	0.52	0 100 100	45, 52, 65, 70	0
2	H	100/122 (81%)	0.21	3 (3%) 52 42	33, 44, 58, 61	0
2	K	100/122 (81%)	0.30	1 (1%) 79 72	39, 49, 62, 67	0
3	C	10/10 (100%)	0.38	0 100 100	43, 46, 54, 55	0
3	F	10/10 (100%)	0.96	0 100 100	59, 66, 72, 74	0
3	I	8/10 (80%)	0.66	1 (12%) 8 6	48, 50, 63, 64	0
3	L	8/10 (80%)	1.35	2 (25%) 2 1	56, 63, 69, 70	0
All	All	1536/1632 (94%)	0.36	31 (2%) 65 56	31, 51, 73, 93	0

The worst 5 of 31 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	H	99	MET	3.5
1	D	249	THR	3.3
1	D	252	THR	3.2
3	L	8	VAL	3.1
1	J	137	TYR	3.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.