



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 07:54 AM UTC

PDB ID : 8JNK / pdb_00008jnk
Title : Crystal structure of human ALKBH3 bound to ssDNA through active site crosslink
Authors : Zhang, L.
Deposited on : 2023-06-06
Resolution : 2.69 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:

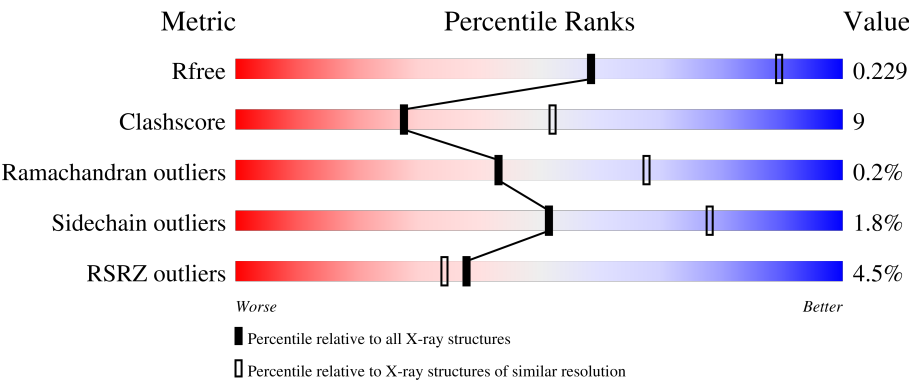
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.69 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	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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	232	4% 64% 20% • 12%
1	C	232	8% 64% 22% • 12%
1	E	232	2% 65% 22% • 12%
1	G	232	12% 62% 22% • 13%
2	B	6	17% 67% 33%

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Mol	Chain	Length	Quality of chain
2	D	6	
2	F	6	
2	H	6	
3	I	216	
3	J	216	
3	K	216	
3	L	216	
4	M	217	
4	N	217	
4	O	217	
4	P	217	

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 20379 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 Alpha-ketoglutarate-dependent dioxygenase alkB homolog 3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	203	Total	C	N	O	S	0	0	0
			1674	1061	299	308	6			
1	C	205	Total	C	N	O	S	1	0	0
			1694	1073	299	316	6			
1	E	203	Total	C	N	O	S	1	0	0
			1679	1065	297	311	6			
1	G	201	Total	C	N	O	S	2	0	0
			1654	1050	294	304	6			

There are 96 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	49	MET	-	initiating methionine	UNP Q96Q83
A	50	GLY	-	expression tag	UNP Q96Q83
A	51	SER	-	expression tag	UNP Q96Q83
A	52	SER	-	expression tag	UNP Q96Q83
A	53	HIS	-	expression tag	UNP Q96Q83
A	54	HIS	-	expression tag	UNP Q96Q83
A	55	HIS	-	expression tag	UNP Q96Q83
A	56	HIS	-	expression tag	UNP Q96Q83
A	57	HIS	-	expression tag	UNP Q96Q83
A	58	HIS	-	expression tag	UNP Q96Q83
A	59	SER	-	expression tag	UNP Q96Q83
A	60	SER	-	expression tag	UNP Q96Q83
A	61	GLY	-	expression tag	UNP Q96Q83
A	62	LEU	-	expression tag	UNP Q96Q83
A	63	VAL	-	expression tag	UNP Q96Q83
A	64	PRO	-	expression tag	UNP Q96Q83
A	65	ARG	-	expression tag	UNP Q96Q83
A	66	GLY	-	expression tag	UNP Q96Q83
A	67	SER	-	expression tag	UNP Q96Q83
A	68	HIS	-	expression tag	UNP Q96Q83
A	69	MET	-	expression tag	UNP Q96Q83

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Chain	Residue	Modelled	Actual	Comment	Reference
A	110	SER	CYS	engineered mutation	UNP Q96Q83
A	195	CYS	GLU	conflict	UNP Q96Q83
A	201	SER	CYS	engineered mutation	UNP Q96Q83
C	49	MET	-	initiating methionine	UNP Q96Q83
C	50	GLY	-	expression tag	UNP Q96Q83
C	51	SER	-	expression tag	UNP Q96Q83
C	52	SER	-	expression tag	UNP Q96Q83
C	53	HIS	-	expression tag	UNP Q96Q83
C	54	HIS	-	expression tag	UNP Q96Q83
C	55	HIS	-	expression tag	UNP Q96Q83
C	56	HIS	-	expression tag	UNP Q96Q83
C	57	HIS	-	expression tag	UNP Q96Q83
C	58	HIS	-	expression tag	UNP Q96Q83
C	59	SER	-	expression tag	UNP Q96Q83
C	60	SER	-	expression tag	UNP Q96Q83
C	61	GLY	-	expression tag	UNP Q96Q83
C	62	LEU	-	expression tag	UNP Q96Q83
C	63	VAL	-	expression tag	UNP Q96Q83
C	64	PRO	-	expression tag	UNP Q96Q83
C	65	ARG	-	expression tag	UNP Q96Q83
C	66	GLY	-	expression tag	UNP Q96Q83
C	67	SER	-	expression tag	UNP Q96Q83
C	68	HIS	-	expression tag	UNP Q96Q83
C	69	MET	-	expression tag	UNP Q96Q83
C	110	SER	CYS	engineered mutation	UNP Q96Q83
C	195	CYS	GLU	conflict	UNP Q96Q83
C	201	SER	CYS	engineered mutation	UNP Q96Q83
E	49	MET	-	initiating methionine	UNP Q96Q83
E	50	GLY	-	expression tag	UNP Q96Q83
E	51	SER	-	expression tag	UNP Q96Q83
E	52	SER	-	expression tag	UNP Q96Q83
E	53	HIS	-	expression tag	UNP Q96Q83
E	54	HIS	-	expression tag	UNP Q96Q83
E	55	HIS	-	expression tag	UNP Q96Q83
E	56	HIS	-	expression tag	UNP Q96Q83
E	57	HIS	-	expression tag	UNP Q96Q83
E	58	HIS	-	expression tag	UNP Q96Q83
E	59	SER	-	expression tag	UNP Q96Q83
E	60	SER	-	expression tag	UNP Q96Q83
E	61	GLY	-	expression tag	UNP Q96Q83
E	62	LEU	-	expression tag	UNP Q96Q83
E	63	VAL	-	expression tag	UNP Q96Q83

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Chain	Residue	Modelled	Actual	Comment	Reference
E	64	PRO	-	expression tag	UNP Q96Q83
E	65	ARG	-	expression tag	UNP Q96Q83
E	66	GLY	-	expression tag	UNP Q96Q83
E	67	SER	-	expression tag	UNP Q96Q83
E	68	HIS	-	expression tag	UNP Q96Q83
E	69	MET	-	expression tag	UNP Q96Q83
E	110	SER	CYS	engineered mutation	UNP Q96Q83
E	195	CYS	GLU	conflict	UNP Q96Q83
E	201	SER	CYS	engineered mutation	UNP Q96Q83
G	49	MET	-	initiating methionine	UNP Q96Q83
G	50	GLY	-	expression tag	UNP Q96Q83
G	51	SER	-	expression tag	UNP Q96Q83
G	52	SER	-	expression tag	UNP Q96Q83
G	53	HIS	-	expression tag	UNP Q96Q83
G	54	HIS	-	expression tag	UNP Q96Q83
G	55	HIS	-	expression tag	UNP Q96Q83
G	56	HIS	-	expression tag	UNP Q96Q83
G	57	HIS	-	expression tag	UNP Q96Q83
G	58	HIS	-	expression tag	UNP Q96Q83
G	59	SER	-	expression tag	UNP Q96Q83
G	60	SER	-	expression tag	UNP Q96Q83
G	61	GLY	-	expression tag	UNP Q96Q83
G	62	LEU	-	expression tag	UNP Q96Q83
G	63	VAL	-	expression tag	UNP Q96Q83
G	64	PRO	-	expression tag	UNP Q96Q83
G	65	ARG	-	expression tag	UNP Q96Q83
G	66	GLY	-	expression tag	UNP Q96Q83
G	67	SER	-	expression tag	UNP Q96Q83
G	68	HIS	-	expression tag	UNP Q96Q83
G	69	MET	-	expression tag	UNP Q96Q83
G	110	SER	CYS	engineered mutation	UNP Q96Q83
G	195	CYS	GLU	conflict	UNP Q96Q83
G	201	SER	CYS	engineered mutation	UNP Q96Q83

- Molecule 2 is a DNA chain called DNA.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	B	6	Total	C	N	O	P	S	0	0	0
			124	60	20	37	6	1			
2	D	5	Total	C	N	O	P	S	0	0	0
			104	50	18	30	5	1			
2	F	6	Total	C	N	O	P	S	0	0	0
			124	60	20	37	6	1			

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Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	H	5	Total	C	N	O	P	S	0	0	0
			104	50	18	30	5	1			

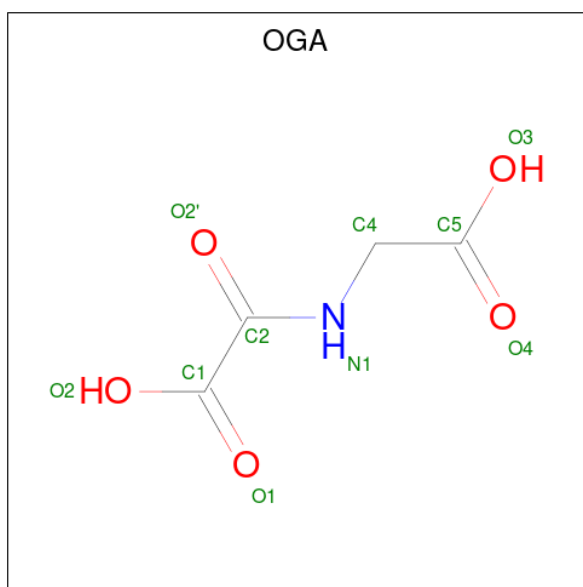
- Molecule 3 is a protein called Synthetic antibody heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	I	214	Total	C	N	O	S	0	0	0
			1601	1014	264	317	6			
3	J	212	Total	C	N	O	S	1	0	0
			1590	1008	262	314	6			
3	K	210	Total	C	N	O	S	0	0	0
			1572	997	259	310	6			
3	L	209	Total	C	N	O	S	0	0	0
			1566	994	258	308	6			

- Molecule 4 is a protein called Synthetic antibody light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	M	214	Total	C	N	O	S	1	0	0
			1643	1032	274	332	5			
4	N	214	Total	C	N	O	S	1	0	0
			1643	1032	274	332	5			
4	O	214	Total	C	N	O	S	2	0	0
			1643	1032	274	332	5			
4	P	214	Total	C	N	O	S	2	0	0
			1643	1032	274	332	5			

- Molecule 5 is N-OXALYLGLYCINE (CCD ID: OGA) (formula: C₄H₅NO₅).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	0	0
			10	4	1	5		
5	C	1	Total	C	N	O	0	0
			10	4	1	5		
5	E	1	Total	C	N	O	0	0
			10	4	1	5		
5	G	1	Total	C	N	O	0	0
			10	4	1	5		

- Molecule 6 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	1	Total	Mn	0	0
			1	1		
6	C	1	Total	Mn	0	0
			1	1		
6	E	1	Total	Mn	0	0
			1	1		
6	G	1	Total	Mn	0	0
			1	1		

- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	34	Total	O	0	0
			34	34		

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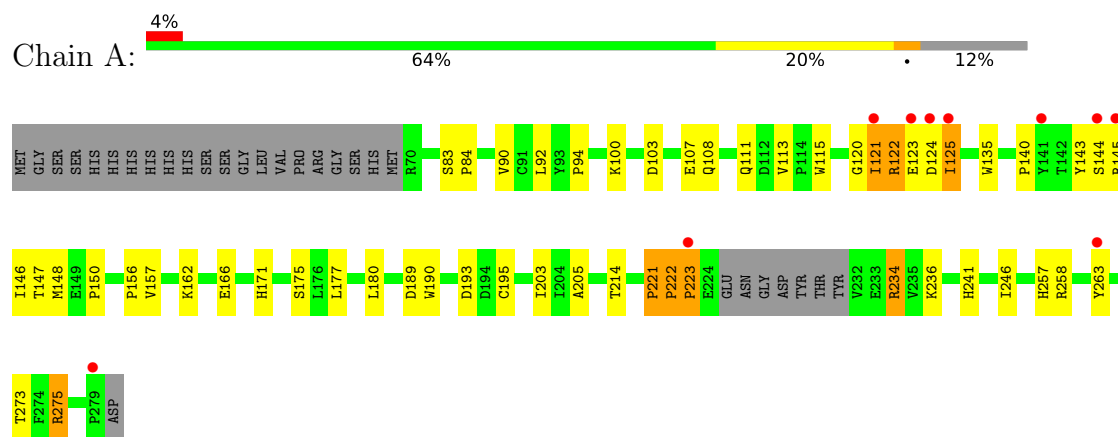
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	B	2	Total 2	O 2	0	0
7	C	4	Total 4	O 4	0	0
7	E	26	Total 26	O 26	0	0
7	F	2	Total 2	O 2	0	0
7	G	12	Total 12	O 12	0	0
7	H	1	Total 1	O 1	0	0
7	I	41	Total 41	O 41	0	0
7	J	30	Total 30	O 30	0	0
7	K	24	Total 24	O 24	0	0
7	L	7	Total 7	O 7	0	0
7	M	26	Total 26	O 26	0	0
7	N	20	Total 20	O 20	0	0
7	O	30	Total 30	O 30	0	0
7	P	18	Total 18	O 18	0	0

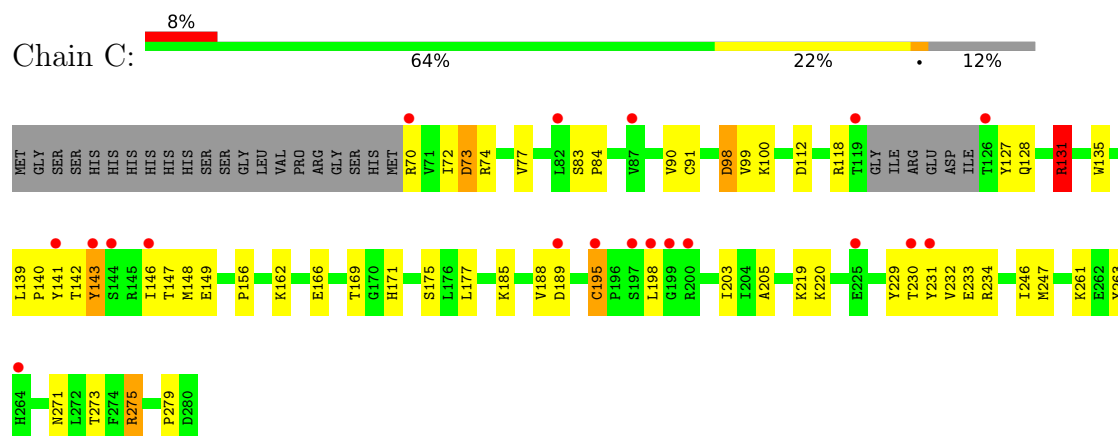
3 Residue-property plots [i](#)

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.

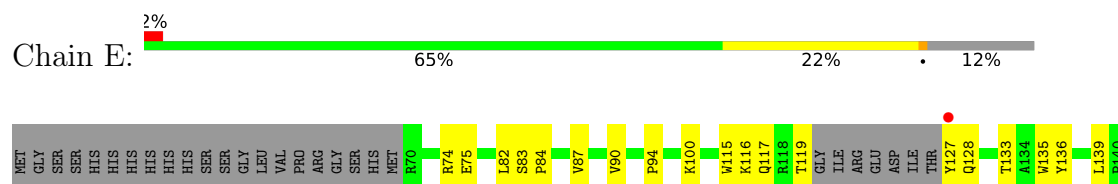
- Molecule 1: Alpha-ketoglutarate-dependent dioxygenase alkB homolog 3



- Molecule 1: Alpha-ketoglutarate-dependent dioxygenase alkB homolog 3

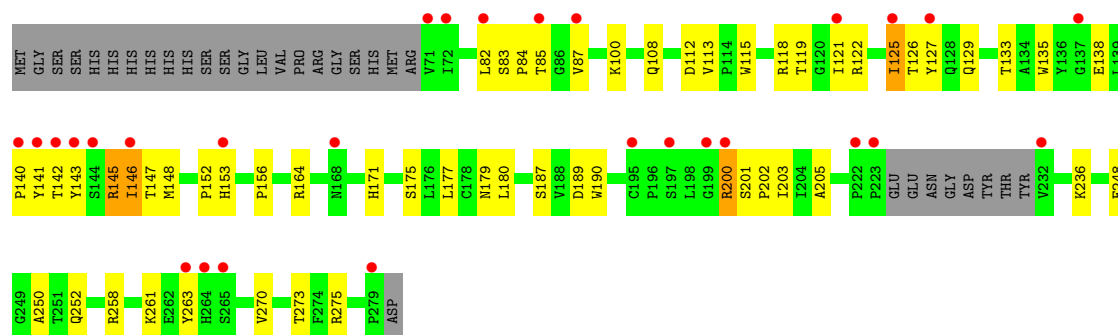


- Molecule 1: Alpha-ketoglutarate-dependent dioxygenase alkB homolog 3

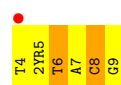




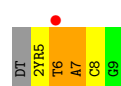
- Molecule 1: Alpha-ketoglutarate-dependent dioxygenase alkB homolog 3



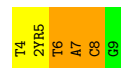
- Molecule 2: DNA



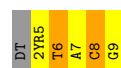
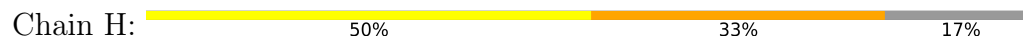
- Molecule 2: DNA



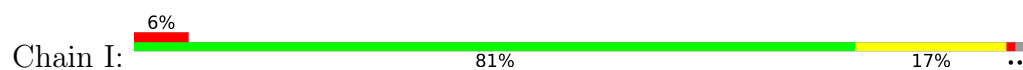
- Molecule 2: DNA

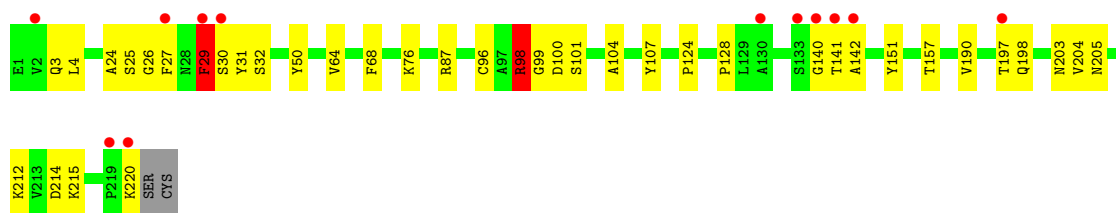


- Molecule 2: DNA

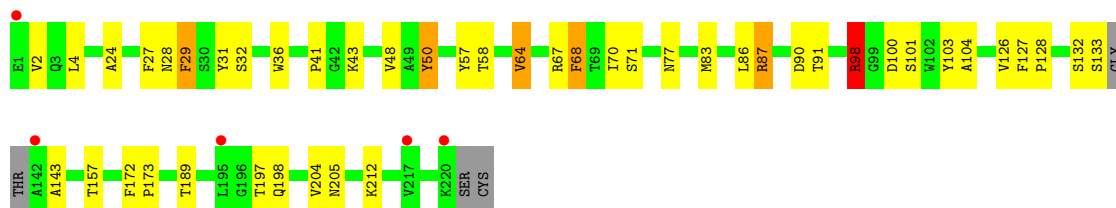
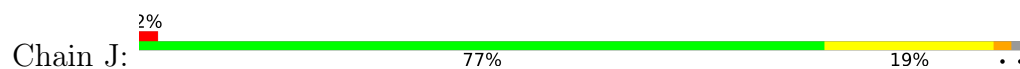


- Molecule 3: Synthetic antibody heavy chain

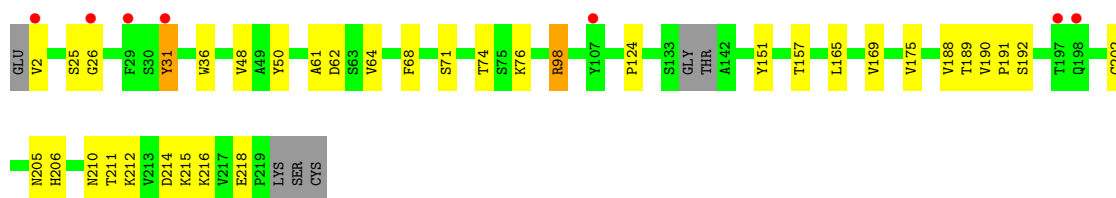
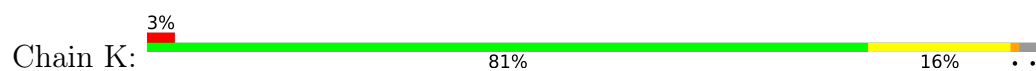




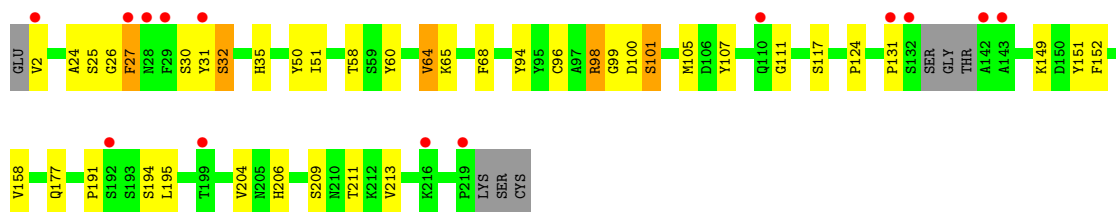
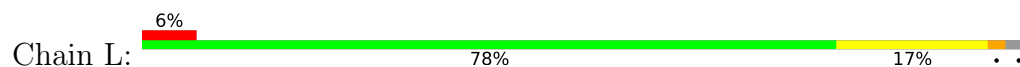
• Molecule 3: Synthetic antibody heavy chain



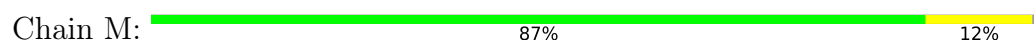
• Molecule 3: Synthetic antibody heavy chain



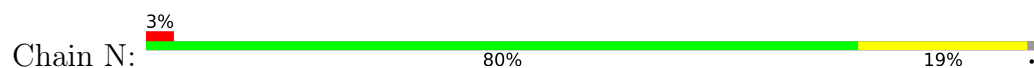
• Molecule 3: Synthetic antibody heavy chain

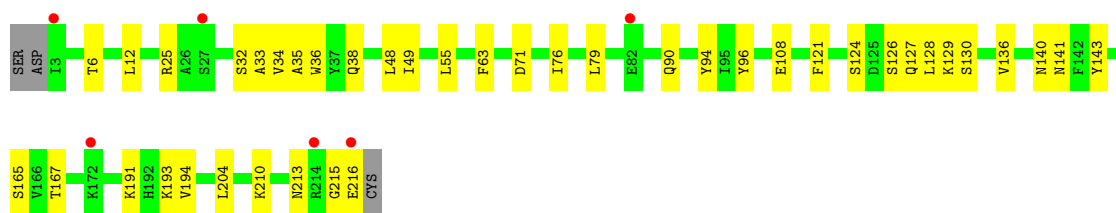


• Molecule 4: Synthetic antibody light chain

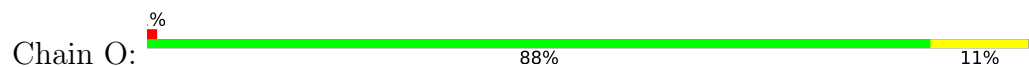


• Molecule 4: Synthetic antibody light chain

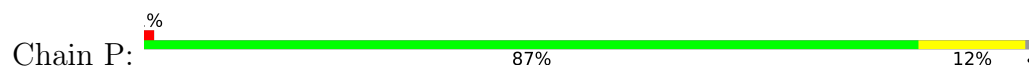




● Molecule 4: Synthetic antibody light chain



● Molecule 4: Synthetic antibody light chain



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	129.46Å 136.15Å 203.39Å 90.00° 107.11° 90.00°	Depositor
Resolution (Å)	46.90 – 2.69 46.90 – 2.69	Depositor EDS
% Data completeness (in resolution range)	99.1 (46.90-2.69) 99.2 (46.90-2.69)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.32 (at 2.69Å)	Xtriage
Refinement program	PHENIX (1.20.1_4487: ???)	Depositor
R, R_{free}	0.218 , 0.229 0.218 , 0.229	Depositor DCC
R_{free} test set	4689 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	52.2	Xtriage
Anisotropy	0.373	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 34.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.008 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	20379	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.33% 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: MN, OGA, 2YR

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	0.43	2/1722 (0.1%)	0.89	6/2345 (0.3%)
1	C	0.39	1/1744 (0.1%)	0.95	4/2377 (0.2%)
1	E	0.30	0/1729	0.87	6/2356 (0.3%)
1	G	0.46	0/1702	0.94	1/2319 (0.0%)
2	B	0.46	0/112	1.30	2/168 (1.2%)
2	D	0.45	0/91	1.21	3/138 (2.2%)
2	F	0.49	0/112	1.26	3/168 (1.8%)
2	H	0.47	0/91	1.27	3/138 (2.2%)
3	I	0.40	0/1643	0.70	4/2242 (0.2%)
3	J	0.76	1/1631 (0.1%)	1.00	9/2224 (0.4%)
3	K	0.57	0/1613	0.74	4/2201 (0.2%)
3	L	0.36	1/1607 (0.1%)	0.60	0/2193
4	M	0.22	1/1680 (0.1%)	0.44	0/2283
4	N	0.16	0/1680	0.45	0/2283
4	O	0.39	3/1680 (0.2%)	0.49	2/2283 (0.1%)
4	P	0.15	0/1680	0.46	0/2283
All	All	0.42	9/20517 (0.0%)	0.76	47/28001 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	C	0	2
1	G	0	2
3	I	0	1
3	J	0	2
All	All	0	8

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	O	133	ALA	C-O	-6.61	1.16	1.24
1	A	221	PRO	C-O	-5.94	1.17	1.24
4	M	129	LYS	CA-C	5.90	1.60	1.52
4	O	132	THR	C-O	-5.72	1.16	1.23
1	A	222	PRO	C-O	-5.72	1.18	1.24
4	O	127	GLN	C-O	-5.25	1.17	1.24
3	L	101	SER	C-O	-5.14	1.17	1.23
1	C	140	PRO	C-O	-5.11	1.17	1.23
3	J	32	SER	CA-CB	-5.08	1.45	1.53

All (47) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	117	GLN	CA-C-N	-9.32	106.80	122.64
1	E	117	GLN	C-N-CA	-9.32	106.80	122.64
3	J	100	ASP	N-CA-C	-8.98	102.32	113.28
3	I	141	THR	N-CA-C	-8.95	98.18	110.35
2	B	6	DT	P-O3'-C3'	-8.88	106.88	120.20
2	F	6	DT	P-O3'-C3'	-7.69	108.66	120.20
2	D	7	DA	P-O3'-C3'	-7.47	109.00	120.20
3	I	100	ASP	N-CA-C	-7.34	104.12	113.23
2	F	7	DA	P-O3'-C3'	-7.19	109.42	120.20
2	H	7	DA	P-O3'-C3'	-6.84	109.93	120.20
2	H	8	DC	P-O3'-C3'	-6.83	109.95	120.20
2	H	6	DT	P-O3'-C3'	-6.53	110.41	120.20
3	J	29	PHE	CA-CB-CG	6.49	120.29	113.80
3	J	128	PRO	N-CA-CB	-6.27	98.15	103.36
1	C	140	PRO	CB-CA-C	-6.20	102.83	110.95
3	K	74	THR	CA-C-N	6.10	128.45	120.28
3	K	74	THR	C-N-CA	6.10	128.45	120.28
1	C	73	ASP	N-CA-CB	-6.07	100.03	111.43
3	J	90	ASP	N-CA-C	-6.03	105.68	113.16
3	K	31	TYR	N-CA-C	-5.99	105.89	112.72
1	E	144	SER	CA-C-N	5.97	132.94	121.54
1	E	144	SER	C-N-CA	5.97	132.94	121.54
1	C	263	TYR	CA-CB-CG	5.92	124.56	113.90
3	J	29	PHE	CB-CA-C	-5.92	101.60	111.13
1	C	143	TYR	CA-CB-CG	5.86	124.45	113.90
2	D	8	DC	P-O3'-C3'	-5.82	111.47	120.20
3	I	29	PHE	CB-CA-C	-5.80	102.29	111.17
3	J	41	PRO	N-CA-CB	-5.74	97.22	103.25
1	A	148	MET	N-CA-C	-5.67	99.95	108.96

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	6	DT	P-O3'-C3'	-5.59	111.81	120.20
3	J	68	PHE	CA-CB-CG	5.58	119.39	113.80
1	E	145	ARG	CB-CG-CD	-5.53	98.58	111.30
4	O	129	LYS	CA-C-N	5.47	129.35	120.60
4	O	129	LYS	C-N-CA	5.47	129.35	120.60
1	A	147	THR	N-CA-C	5.43	117.26	108.41
2	F	8	DC	P-O3'-C3'	-5.42	112.08	120.20
3	J	29	PHE	N-CA-C	-5.40	105.13	112.26
3	I	140	GLY	CA-C-O	-5.34	115.32	121.47
1	A	150	PRO	N-CA-C	5.34	119.47	111.14
1	A	223	PRO	N-CA-C	-5.32	101.51	112.47
3	K	98	ARG	CG-CD-NE	-5.30	100.35	112.00
1	G	140	PRO	N-CA-C	5.27	120.40	111.03
3	J	91	THR	CB-CA-C	5.26	118.31	109.89
2	B	8	DC	P-O3'-C3'	-5.20	112.41	120.20
1	E	263	TYR	CA-CB-CG	5.02	122.94	113.90
1	A	223	PRO	CA-C-N	5.01	130.72	121.70
1	A	223	PRO	C-N-CA	5.01	130.72	121.70

There are no chirality outliers.

All (8) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	234	ARG	Sidechain
1	C	131	ARG	Sidechain
1	C	275	ARG	Sidechain
1	G	145	ARG	Sidechain
1	G	200	ARG	Sidechain
3	I	98	ARG	Sidechain
3	J	87	ARG	Sidechain
3	J	98	ARG	Sidechain

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	1674	0	1634	44	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	1694	0	1633	51	0
1	E	1679	0	1622	37	0
1	G	1654	0	1614	52	0
2	B	124	0	74	12	0
2	D	104	0	61	10	0
2	F	124	0	72	7	0
2	H	104	0	61	6	0
3	I	1601	0	1548	20	0
3	J	1590	0	1537	30	0
3	K	1572	0	1515	20	0
3	L	1566	0	1510	38	0
4	M	1643	0	1603	17	0
4	N	1643	0	1603	31	0
4	O	1643	0	1603	15	0
4	P	1643	0	1603	19	0
5	A	10	0	3	0	0
5	C	10	0	3	0	0
5	E	10	0	3	0	0
5	G	10	0	3	1	0
6	A	1	0	0	0	0
6	C	1	0	0	0	0
6	E	1	0	0	0	0
6	G	1	0	0	0	0
7	A	34	0	0	0	0
7	B	2	0	0	0	0
7	C	4	0	0	0	0
7	E	26	0	0	2	0
7	F	2	0	0	0	0
7	G	12	0	0	0	0
7	H	1	0	0	0	0
7	I	41	0	0	1	0
7	J	30	0	0	0	0
7	K	24	0	0	1	0
7	L	7	0	0	0	0
7	M	26	0	0	0	0
7	N	20	0	0	0	0
7	O	30	0	0	0	0
7	P	18	0	0	0	0
All	All	20379	0	19305	355	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 9.

All (355) 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:C:141:TYR:CE1	2:D:5:2YR:H16	1.65	1.31
1:A:195:CYS:SG	2:B:5:2YR:S	2.35	1.25
1:C:141:TYR:CZ	2:D:5:2YR:H16	1.80	1.17
3:L:2:VAL:HG22	3:L:26:GLY:HA3	1.46	0.98
1:E:195:CYS:N	2:F:5:2YR:S	2.33	0.95
1:A:125:ILE:HG21	2:B:6:DT:H71	1.52	0.91
3:J:27:PHE:HE2	3:J:98:ARG:HD3	1.37	0.88
3:J:64:VAL:HG13	3:J:68:PHE:HB2	1.53	0.88
1:C:141:TYR:CE1	2:D:5:2YR:C8	2.56	0.85
1:C:195:CYS:SG	1:C:198:LEU:CD2	2.67	0.82
1:G:121:ILE:O	1:G:121:ILE:HD12	1.82	0.79
3:L:51:ILE:HD12	3:L:58:THR:CG2	2.13	0.79
1:C:143:TYR:O	1:C:146:ILE:HG13	1.80	0.79
3:L:2:VAL:HA	3:L:25:SER:O	1.83	0.79
4:P:40:LYS:HB2	4:P:43:LYS:HD3	1.65	0.78
3:J:126:VAL:HG21	3:J:204:VAL:HG11	1.66	0.77
1:E:193:ASP:HA	2:F:5:2YR:H16	1.66	0.77
1:G:148:MET:HE1	1:G:177:LEU:HD13	1.66	0.76
1:A:125:ILE:HG21	2:B:6:DT:C7	2.17	0.75
4:N:215:GLY:O	4:N:216:GLU:HG3	1.86	0.75
3:L:98:ARG:HG2	3:L:98:ARG:HH11	1.51	0.75
4:P:38:GLN:HB2	4:P:48:LEU:HD11	1.70	0.74
3:L:51:ILE:CD1	3:L:58:THR:CG2	2.66	0.74
1:E:100:LYS:HD3	3:J:31:TYR:HB3	1.70	0.73
4:P:40:LYS:H	4:P:43:LYS:HZ2	1.34	0.73
3:I:142:ALA:HB3	3:I:190:VAL:O	1.89	0.73
1:G:189:ASP:HB3	2:H:5:2YR:H5	1.70	0.73
4:N:193:LYS:HD3	4:N:194:VAL:HG23	1.71	0.72
3:L:51:ILE:HD12	3:L:58:THR:HG22	1.71	0.72
1:C:141:TYR:HD2	1:C:148:MET:HE2	1.54	0.71
2:F:6:DT:H2'	2:F:7:DA:C8	2.24	0.71
3:J:83:MET:HB3	3:J:86:LEU:HD21	1.74	0.69
1:A:122:ARG:O	1:A:125:ILE:HG22	1.93	0.69
1:G:121:ILE:HG22	1:G:126:THR:HG22	1.75	0.69
2:B:7:DA:H4'	2:B:8:DC:C2	2.28	0.68
4:N:12:LEU:HD13	4:P:159:SER:HB2	1.75	0.68
1:A:189:ASP:HB3	2:B:5:2YR:H5	1.76	0.68
1:G:119:THR:O	1:G:146:ILE:HG12	1.94	0.68
3:J:157:THR:OG1	3:J:205:ASN:HB3	1.94	0.68
3:L:2:VAL:HG22	3:L:26:GLY:CA	2.23	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:121:ILE:HG13	1:G:145:ARG:HB2	1.77	0.67
3:L:51:ILE:CD1	3:L:58:THR:HG22	2.25	0.67
1:E:119:THR:O	1:E:146:ILE:HD11	1.94	0.67
1:G:121:ILE:CG1	1:G:145:ARG:HB2	2.26	0.66
1:C:141:TYR:OH	2:D:5:2YR:H18	1.94	0.66
1:G:177:LEU:HB3	1:G:273:THR:HG23	1.77	0.66
4:N:126:SER:HA	4:N:129:LYS:HD2	1.77	0.66
1:A:195:CYS:CB	2:B:5:2YR:S	2.84	0.66
3:L:2:VAL:CG2	3:L:26:GLY:HA3	2.25	0.65
1:G:118:ARG:NH1	1:G:133:THR:OG1	2.29	0.65
1:A:156:PRO:HG2	4:M:94:TYR:CD1	2.32	0.65
1:E:75:GLU:HB3	1:E:94:PRO:HD2	1.79	0.65
3:J:27:PHE:CE2	3:J:98:ARG:HD3	2.28	0.65
1:E:119:THR:HG22	1:E:128:GLN:HG2	1.79	0.65
3:I:3:GLN:HG2	3:I:4:LEU:H	1.61	0.64
1:E:119:THR:O	1:E:145:ARG:NH1	2.30	0.64
3:I:203:ASN:ND2	3:I:214:ASP:OD1	2.25	0.64
1:G:273:THR:HG21	1:G:275:ARG:HH21	1.63	0.63
1:C:135:TRP:HB2	1:C:148:MET:HE3	1.80	0.62
1:G:171:HIS:CD2	1:G:203:ILE:HD13	2.34	0.62
1:C:73:ASP:OD1	1:C:74:ARG:HG2	1.99	0.62
1:A:193:ASP:OD1	1:A:257:HIS:CE1	2.52	0.62
1:E:156:PRO:HG2	4:N:94:TYR:CD1	2.34	0.62
3:L:51:ILE:HD12	3:L:58:THR:HG23	1.81	0.62
4:M:38:GLN:HB2	4:M:48:LEU:HD11	1.81	0.62
3:L:98:ARG:HG3	3:L:99:GLY:N	2.14	0.61
4:M:25:ARG:NH2	4:O:183:THR:O	2.32	0.61
1:E:156:PRO:HG2	4:N:94:TYR:CG	2.36	0.61
3:L:51:ILE:CD1	3:L:58:THR:HG23	2.31	0.61
4:N:121:PHE:HB2	4:N:136:VAL:HG13	1.82	0.61
3:J:2:VAL:HG13	3:J:27:PHE:CD1	2.36	0.61
1:C:100:LYS:HD3	3:L:31:TYR:HB3	1.83	0.60
1:C:171:HIS:CD2	1:C:203:ILE:HG21	2.37	0.60
1:G:142:THR:HG22	1:G:147:THR:HG23	1.83	0.59
1:C:195:CYS:SG	1:C:198:LEU:HD23	2.41	0.59
1:G:142:THR:HA	1:G:146:ILE:O	2.02	0.59
3:K:169:VAL:HG22	3:K:188:VAL:HG22	1.83	0.59
1:E:193:ASP:OD2	1:E:275:ARG:NH2	2.35	0.59
1:C:156:PRO:HG2	4:P:94:TYR:CD1	2.37	0.59
3:L:124:PRO:HB3	3:L:151:TYR:HB3	1.84	0.59
1:C:142:THR:HG23	1:C:147:THR:OG1	2.03	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:263:TYR:CG	1:G:263:TYR:O	2.56	0.58
1:C:77:VAL:HG23	1:C:91:CYS:SG	2.43	0.58
3:J:64:VAL:HG13	3:J:68:PHE:CB	2.29	0.58
1:G:263:TYR:HB2	2:H:9:DG:O3'	2.02	0.58
3:K:61:ALA:O	3:K:64:VAL:HG12	2.04	0.58
1:G:156:PRO:HG2	4:O:94:TYR:CD1	2.38	0.58
3:L:60:TYR:HB2	3:L:65:LYS:HG3	1.86	0.58
3:K:175:VAL:HG21	4:O:163:GLN:HB3	1.86	0.58
1:A:263:TYR:HB2	2:B:9:DG:O3'	2.04	0.57
1:G:152:PRO:HG2	1:G:153:HIS:CD2	2.38	0.57
4:M:159:SER:HB2	4:O:12:LEU:HD13	1.87	0.57
2:H:8:DC:H2''	2:H:9:DG:O4'	2.05	0.57
3:L:27:PHE:HE2	3:L:107:TYR:CD2	2.23	0.57
3:I:87:ARG:NH1	7:I:301:HOH:O	2.36	0.57
1:C:70:ARG:NH1	1:C:72:ILE:HG12	2.21	0.56
3:I:128:PRO:HD3	3:I:215:LYS:HE2	1.87	0.56
1:E:115:TRP:O	1:E:116:LYS:HD2	2.06	0.56
1:C:141:TYR:CZ	2:D:5:2YR:C8	2.71	0.56
1:G:156:PRO:HG2	4:O:94:TYR:CG	2.40	0.56
4:M:171:SER:C	4:M:172:LYS:HD3	2.30	0.56
1:A:143:TYR:O	1:A:146:ILE:HG13	2.06	0.56
1:C:98:ASP:OD1	1:C:99:VAL:N	2.39	0.55
1:A:120:GLY:HA2	1:A:146:ILE:HG21	1.88	0.55
3:L:24:ALA:HB1	3:L:27:PHE:CE1	2.41	0.55
1:G:148:MET:HE1	1:G:177:LEU:CD1	2.34	0.55
1:G:200:ARG:O	1:G:202:PRO:HD3	2.07	0.55
1:A:122:ARG:O	1:A:125:ILE:N	2.28	0.55
1:E:171:HIS:ND1	7:E:1102:HOH:O	2.33	0.55
1:G:263:TYR:O	1:G:263:TYR:CD1	2.60	0.55
1:E:200:ARG:N	1:E:200:ARG:HD2	2.22	0.55
1:G:125:ILE:HD12	1:G:126:THR:H	1.72	0.54
1:C:118:ARG:NH2	1:C:149:GLU:OE1	2.40	0.54
1:G:141:TYR:HD2	1:G:148:MET:HE3	1.71	0.54
3:I:98:ARG:HG2	3:I:99:GLY:N	2.22	0.54
3:I:24:ALA:HB2	3:I:29:PHE:CZ	2.42	0.54
4:N:36:TRP:HB2	4:N:49:ILE:HB	1.88	0.54
1:C:143:TYR:HB3	2:D:5:2YR:H15	1.89	0.54
1:G:180:LEU:HD13	1:G:270:VAL:HG22	1.89	0.54
3:I:157:THR:OG1	3:I:205:ASN:HB2	2.07	0.54
1:C:70:ARG:NH1	1:C:72:ILE:CG1	2.71	0.53
3:J:103:TYR:HB2	4:N:90:GLN:NE2	2.23	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:179:ASN:ND2	5:G:1001:OGA:O1	2.38	0.53
1:G:119:THR:HA	1:G:127:TYR:O	2.09	0.53
1:C:83:SER:OG	1:C:84:PRO:HA	2.08	0.53
1:E:189:ASP:HB3	2:F:5:2YR:H5	1.90	0.53
3:J:101:SER:HG	3:J:104:ALA:H	1.54	0.53
3:I:197:THR:OG1	3:I:198:GLN:N	2.40	0.53
1:A:92:LEU:HG	1:A:94:PRO:HD3	1.91	0.53
4:P:35:ALA:HB3	4:P:90:GLN:HB3	1.91	0.53
1:E:200:ARG:HD2	1:E:200:ARG:H	1.74	0.52
3:L:64:VAL:HG13	3:L:68:PHE:CG	2.44	0.52
1:A:195:CYS:HB2	2:B:5:2YR:S	2.50	0.52
4:M:55:LEU:HD21	4:M:59:VAL:O	2.09	0.52
1:A:193:ASP:HA	2:B:5:2YR:H16	1.90	0.52
2:D:6:DT:H2''	2:D:7:DA:O5'	2.10	0.52
1:E:83:SER:HB2	1:E:84:PRO:HA	1.92	0.52
2:F:4:DT:H1'	2:F:5:2YR:H12	1.90	0.52
4:N:193:LYS:HZ2	4:N:213:ASN:HB3	1.73	0.52
3:I:124:PRO:HB3	3:I:151:TYR:HB3	1.91	0.52
3:J:64:VAL:HG13	3:J:68:PHE:CG	2.45	0.52
4:N:25:ARG:NH2	4:P:183:THR:O	2.41	0.52
1:A:122:ARG:O	1:A:123:GLU:C	2.53	0.51
1:G:261:LYS:HD2	2:H:8:DC:H4'	1.92	0.51
3:J:28:ASN:HA	3:J:77:ASN:HD21	1.74	0.51
4:P:145:ARG:HH11	4:P:145:ARG:HG2	1.75	0.51
1:A:156:PRO:HG2	4:M:94:TYR:CG	2.45	0.51
4:O:201:HIS:CD2	4:O:203:GLY:H	2.28	0.51
3:J:2:VAL:CG1	3:J:27:PHE:CD1	2.94	0.51
1:E:171:HIS:ND1	1:E:203:ILE:HD13	2.26	0.51
3:J:24:ALA:HB2	3:J:29:PHE:CZ	2.46	0.51
1:E:205:ALA:O	1:E:273:THR:HA	2.11	0.51
1:A:222:PRO:O	1:A:223:PRO:C	2.53	0.50
4:N:124:SER:O	4:N:128:LEU:HD12	2.12	0.50
3:K:216:LYS:HE2	3:K:218:GLU:OE1	2.11	0.50
1:C:220:LYS:HA	1:C:229:TYR:CE1	2.47	0.50
3:K:64:VAL:HG22	3:K:68:PHE:CG	2.47	0.50
3:L:32:SER:OG	3:L:98:ARG:NE	2.44	0.50
4:N:127:GLN:O	4:N:130:SER:OG	2.26	0.50
1:G:121:ILE:HD12	1:G:121:ILE:C	2.36	0.50
3:J:132:SER:OG	3:J:133:SER:N	2.39	0.50
4:M:13:SER:HB3	4:M:110:LYS:HG2	1.94	0.50
1:A:135:TRP:CH2	1:A:175:SER:HB3	2.48	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:118:ARG:O	1:G:129:GLN:HG3	2.12	0.49
4:N:38:GLN:HB2	4:N:48:LEU:HD11	1.93	0.49
1:C:205:ALA:O	1:C:273:THR:HA	2.13	0.49
3:K:36:TRP:O	3:K:48:VAL:HG22	2.12	0.49
1:A:122:ARG:HB2	1:A:122:ARG:HH11	1.78	0.49
1:E:221:PRO:HD3	1:E:229:TYR:CE1	2.47	0.49
1:C:118:ARG:HD3	1:C:146:ILE:HD13	1.94	0.49
3:I:27:PHE:CE2	3:I:98:ARG:HD2	2.47	0.49
3:K:206:HIS:HB3	3:K:211:THR:OG1	2.12	0.49
4:N:126:SER:O	4:N:129:LYS:HB2	2.12	0.49
1:A:90:VAL:HA	1:A:246:ILE:O	2.13	0.49
1:G:205:ALA:O	1:G:273:THR:HA	2.13	0.49
4:P:39:GLN:HA	4:P:43:LYS:NZ	2.27	0.49
1:C:148:MET:HE1	1:C:177:LEU:HD13	1.94	0.48
3:L:35:HIS:CG	3:L:105:MET:HE2	2.48	0.48
3:L:98:ARG:HB3	3:L:107:TYR:HB2	1.95	0.48
4:M:108:GLU:OE1	4:M:176:TYR:OH	2.29	0.48
3:J:143:ALA:HB2	3:J:189:THR:HG22	1.95	0.48
3:L:206:HIS:ND1	3:L:209:SER:HB2	2.28	0.48
1:C:156:PRO:HG2	4:P:94:TYR:CG	2.49	0.48
1:G:141:TYR:HD2	1:G:148:MET:CE	2.26	0.48
1:C:189:ASP:HB3	2:D:5:2YR:H5	1.94	0.48
1:A:83:SER:HB2	1:A:84:PRO:HA	1.95	0.48
1:G:85:THR:OG1	1:G:250:ALA:HB1	2.13	0.48
4:P:127:GLN:O	4:P:130:SER:OG	2.30	0.47
1:E:271:ASN:OD1	1:E:273:THR:HG23	2.14	0.47
4:N:191:LYS:HG3	4:N:191:LYS:O	2.15	0.47
1:C:230:THR:HG23	1:C:231:TYR:CD2	2.50	0.47
3:K:165:LEU:HD21	3:K:188:VAL:HG11	1.97	0.47
3:L:27:PHE:CE2	3:L:98:ARG:HD3	2.50	0.47
3:L:98:ARG:HG2	3:L:98:ARG:NH1	2.26	0.47
3:I:64:VAL:HG13	3:I:68:PHE:CG	2.51	0.46
3:J:28:ASN:HA	3:J:77:ASN:ND2	2.30	0.46
1:G:121:ILE:CG2	1:G:126:THR:HG22	2.45	0.46
4:N:193:LYS:NZ	4:N:213:ASN:HB3	2.31	0.46
1:C:271:ASN:OD1	1:C:273:THR:HG23	2.15	0.46
1:E:135:TRP:CH2	1:E:175:SER:HB3	2.51	0.46
1:C:127:TYR:CD1	1:C:127:TYR:N	2.84	0.46
1:E:74:ARG:HD3	1:E:74:ARG:HA	1.84	0.46
3:L:27:PHE:CD1	3:L:27:PHE:N	2.84	0.46
3:L:98:ARG:HH11	3:L:98:ARG:CG	2.22	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:191:PRO:HG2	3:L:194:SER:OG	2.15	0.46
1:C:233:GLU:N	1:C:233:GLU:OE1	2.48	0.46
4:M:13:SER:OG	4:M:108:GLU:OE2	2.28	0.46
1:A:125:ILE:HD12	1:A:125:ILE:HA	1.77	0.46
1:E:135:TRP:CZ2	1:E:175:SER:HB3	2.51	0.45
4:N:76:ILE:HG21	4:N:79:LEU:HD12	1.99	0.45
1:A:190:TRP:CE2	1:A:258:ARG:HD3	2.52	0.45
1:C:90:VAL:HG12	1:C:247:MET:HG2	1.98	0.45
1:G:118:ARG:HG3	1:G:118:ARG:HH11	1.81	0.45
1:A:177:LEU:HB3	1:A:275:ARG:HH11	1.82	0.45
1:E:135:TRP:CZ2	1:E:141:TYR:HB2	2.51	0.45
3:I:32:SER:HB2	3:I:99:GLY:O	2.16	0.45
3:J:4:LEU:HD21	3:J:27:PHE:HZ	1.80	0.45
1:C:141:TYR:OH	2:D:5:2YR:C9	2.63	0.45
3:L:158:VAL:HG22	3:L:204:VAL:HG22	1.98	0.45
1:C:135:TRP:CZ2	1:C:175:SER:HB3	2.51	0.45
4:O:14:ALA:O	4:O:109:ILE:HA	2.17	0.45
1:A:123:GLU:OE2	2:B:6:DT:H73	2.17	0.45
3:I:76:LYS:HB3	3:I:76:LYS:HE3	1.77	0.45
4:N:35:ALA:HB3	4:N:90:GLN:HB3	1.97	0.45
1:C:162:LYS:O	1:C:166:GLU:HG3	2.17	0.45
1:G:82:LEU:HA	1:G:87:VAL:HA	1.99	0.45
4:N:25:ARG:HD3	4:N:71:ASP:OD1	2.17	0.45
1:C:139:LEU:HD11	1:C:279:PRO:HD3	1.99	0.45
1:G:108:GLN:HG3	1:G:112:ASP:OD2	2.17	0.45
1:G:143:TYR:O	1:G:146:ILE:HB	2.18	0.45
3:J:173:PRO:O	4:N:165:SER:OG	2.35	0.44
1:A:121:ILE:CG2	1:A:145:ARG:HB3	2.47	0.44
3:J:43:LYS:HD2	4:N:6:THR:O	2.16	0.44
4:M:12:LEU:HD13	4:O:159:SER:HB2	1.98	0.44
1:G:203:ILE:HG12	1:G:248:GLU:HB2	1.98	0.44
3:J:197:THR:OG1	3:J:198:GLN:N	2.50	0.44
1:G:190:TRP:CE2	1:G:258:ARG:HD3	2.53	0.44
1:C:70:ARG:C	1:C:70:ARG:CD	2.91	0.44
1:E:176:LEU:HD11	1:E:272:LEU:HB3	1.98	0.44
1:E:177:LEU:HD23	1:E:275:ARG:HH12	1.83	0.44
3:L:27:PHE:CE2	3:L:107:TYR:CD2	3.04	0.44
4:O:144:PRO:O	4:O:201:HIS:HE1	2.01	0.44
1:A:108:GLN:OE1	1:A:157:VAL:HG21	2.17	0.44
1:A:171:HIS:ND1	1:A:203:ILE:HD13	2.32	0.44
1:G:135:TRP:CH2	1:G:175:SER:HB3	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:J:57:TYR:CE1	4:N:96:TYR:HA	2.53	0.44
3:K:215:LYS:NZ	7:K:303:HOH:O	2.49	0.44
4:P:82:GLU:H	4:P:82:GLU:CD	2.26	0.44
1:E:119:THR:C	1:E:145:ARG:HH12	2.26	0.44
1:E:139:LEU:HD11	1:E:279:PRO:HD3	1.99	0.44
1:A:100:LYS:HE2	3:I:31:TYR:O	2.18	0.44
3:J:64:VAL:CG1	3:J:68:PHE:HB2	2.37	0.44
1:A:103:ASP:OD1	1:A:241:HIS:NE2	2.51	0.44
1:A:205:ALA:O	1:A:273:THR:HA	2.18	0.43
1:E:214:THR:CG2	1:E:236:LYS:HG3	2.47	0.43
3:K:124:PRO:HB3	3:K:151:TYR:HB3	1.99	0.43
3:J:50:TYR:O	3:J:58:THR:HA	2.18	0.43
4:N:32:SER:O	4:N:34:VAL:N	2.50	0.43
4:O:43:LYS:HA	4:O:43:LYS:HD2	1.74	0.43
1:C:139:LEU:CD1	1:C:279:PRO:HD3	2.48	0.43
1:G:203:ILE:HG12	1:G:248:GLU:CB	2.49	0.43
4:M:149:VAL:HG22	4:M:199:VAL:HG22	2.00	0.43
1:A:122:ARG:C	1:A:123:GLU:HG2	2.44	0.43
1:A:144:SER:HA	2:B:4:DT:H2'	2.00	0.43
1:E:90:VAL:HA	1:E:246:ILE:O	2.18	0.43
3:K:62:ASP:C	3:K:64:VAL:H	2.27	0.43
4:P:145:ARG:HG2	4:P:145:ARG:NH1	2.33	0.43
4:P:147:ALA:HB2	4:P:201:HIS:HD2	1.84	0.43
1:A:263:TYR:CD1	1:A:263:TYR:C	2.97	0.43
1:C:169:THR:HG21	1:C:246:ILE:HD11	2.00	0.43
1:E:127:TYR:CD2	1:E:145:ARG:NH1	2.87	0.43
3:I:98:ARG:CD	3:I:107:TYR:HD2	2.31	0.43
4:P:66:SER:OG	4:P:67:ARG:N	2.52	0.43
1:G:122:ARG:NH1	2:H:6:DT:N3	2.66	0.43
3:J:127:PHE:CE2	4:N:127:GLN:HG3	2.54	0.43
4:P:115:ALA:HB2	4:P:203:GLY:O	2.19	0.43
2:F:7:DA:H2'	2:F:8:DC:C4	2.53	0.43
2:H:8:DC:H2'	2:H:9:DG:C8	2.54	0.43
3:K:157:THR:OG1	3:K:205:ASN:HB2	2.18	0.43
1:A:107:GLU:O	1:A:111:GLN:HG3	2.18	0.42
1:G:180:LEU:HD13	1:G:270:VAL:CG2	2.48	0.42
1:G:187:SER:OG	1:G:261:LYS:NZ	2.42	0.42
3:L:50:TYR:CD1	3:L:50:TYR:C	2.97	0.42
4:M:48:LEU:HA	4:M:59:VAL:HG21	2.00	0.42
1:A:162:LYS:O	1:A:166:GLU:HG3	2.19	0.42
1:G:141:TYR:CD2	1:G:148:MET:HE3	2.52	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:177:LEU:HB3	1:A:275:ARG:NH1	2.34	0.42
1:C:112:ASP:OD2	4:P:31:SER:OG	2.32	0.42
4:M:36:TRP:CZ3	4:M:89:CYS:HB2	2.55	0.42
1:A:171:HIS:CE1	1:A:203:ILE:HD13	2.54	0.42
4:O:38:GLN:HB2	4:O:48:LEU:HD11	2.02	0.42
1:C:131:ARG:HE	1:C:188:VAL:HB	1.85	0.42
4:P:139:LEU:HD22	4:P:178:LEU:HD22	2.01	0.42
1:C:229:TYR:O	1:C:232:VAL:HG22	2.19	0.42
1:E:136:TYR:HA	1:E:151:ASN:HB3	2.02	0.42
3:K:2:VAL:HA	3:K:26:GLY:HA3	2.02	0.42
3:L:131:PRO:HG3	3:L:195:LEU:HD22	2.02	0.42
3:L:211:THR:HG22	3:L:213:VAL:HG23	2.00	0.42
1:G:113:VAL:HB	1:G:115:TRP:CE2	2.55	0.42
4:N:140:ASN:ND2	4:N:141:ASN:OD1	2.53	0.42
1:C:70:ARG:NH1	1:C:72:ILE:HG13	2.35	0.42
1:E:82:LEU:HD22	1:E:87:VAL:HG22	2.01	0.42
3:K:212:LYS:HE2	3:K:212:LYS:HB2	1.77	0.42
1:A:115:TRP:CE2	1:A:180:LEU:HD12	2.55	0.42
3:I:50:TYR:CD1	3:I:50:TYR:C	2.98	0.42
4:M:117:SER:HB2	4:M:140:ASN:HB3	2.02	0.42
1:A:214:THR:HG21	1:A:236:LYS:HD2	2.01	0.41
3:K:190:VAL:HB	3:K:191:PRO:CD	2.50	0.41
4:M:39:GLN:O	4:M:85:ALA:HB1	2.20	0.41
4:N:55:LEU:HD21	4:N:63:PHE:HB2	2.02	0.41
1:C:149:GLU:OE1	1:C:149:GLU:N	2.41	0.41
1:E:142:THR:HG22	1:E:147:THR:OG1	2.19	0.41
1:G:100:LYS:HG2	3:K:31:TYR:HB3	2.02	0.41
3:K:50:TYR:C	3:K:50:TYR:CD1	2.98	0.41
2:F:7:DA:OP2	2:F:8:DC:N4	2.44	0.41
3:K:210:ASN:O	3:K:212:LYS:NZ	2.54	0.41
4:O:187:ALA:O	4:O:191:LYS:HG3	2.20	0.41
1:A:122:ARG:NH2	2:B:6:DT:C2	2.87	0.41
1:G:236:LYS:HD2	1:G:236:LYS:N	2.35	0.41
3:L:51:ILE:HD13	3:L:58:THR:CG2	2.49	0.41
4:P:166:VAL:HG12	4:P:167:THR:O	2.20	0.41
3:J:36:TRP:HD1	3:J:70:ILE:HD12	1.86	0.41
1:E:190:TRP:CE2	1:E:258:ARG:HD3	2.56	0.41
3:I:204:VAL:O	3:I:212:LYS:HA	2.21	0.41
3:L:94:TYR:O	3:L:111:GLY:HA2	2.21	0.41
3:I:101:SER:OG	3:I:104:ALA:N	2.42	0.41
3:J:67:ARG:C	3:J:68:PHE:HD1	2.29	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:J:172:PHE:CD1	4:N:167:THR:HG23	2.55	0.41
4:M:21:THR:OG1	4:M:75:THR:HG23	2.20	0.41
1:C:219:LYS:HZ2	1:C:219:LYS:HG2	1.69	0.41
3:K:189:THR:HG21	4:O:140:ASN:ND2	2.35	0.41
1:A:113:VAL:HB	1:A:115:TRP:CE2	2.55	0.41
1:C:234:ARG:HH11	1:C:234:ARG:HD3	1.73	0.41
3:K:202:CYS:O	3:K:214:ASP:HA	2.21	0.41
3:L:30:SER:O	3:L:31:TYR:C	2.63	0.41
3:L:149:LYS:NZ	3:L:177:GLN:HE22	2.18	0.41
1:A:121:ILE:HA	1:A:125:ILE:O	2.21	0.41
1:G:146:ILE:HD13	1:G:146:ILE:HA	1.83	0.41
1:G:164:ARG:NH1	4:O:96:TYR:CZ	2.88	0.41
4:N:204:LEU:HD23	4:N:204:LEU:HA	1.96	0.41
4:O:35:ALA:HA	4:O:49:ILE:O	2.21	0.41
1:C:185:LYS:O	1:C:261:LYS:HE3	2.21	0.40
3:L:27:PHE:H	3:L:27:PHE:HD1	1.69	0.40
3:L:117:SER:HB3	3:L:152:PHE:CZ	2.56	0.40
1:C:127:TYR:HB2	1:C:128:GLN:H	1.66	0.40
1:E:239:LEU:O	7:E:1101:HOH:O	2.22	0.40
1:E:133:THR:HA	1:E:178:CYS:O	2.21	0.40
3:I:220:LYS:HA	3:I:220:LYS:HD3	1.77	0.40
4:N:210:LYS:HA	4:N:210:LYS:HD3	1.92	0.40
1:C:141:TYR:OH	2:D:5:2YR:H16	2.16	0.40
1:G:83:SER:HB3	1:G:84:PRO:HA	2.04	0.40
3:J:205:ASN:OD1	3:J:212:LYS:HE2	2.21	0.40
4:N:108:GLU:OE1	4:N:143:TYR:HE2	2.04	0.40
1:C:195:CYS:SG	1:C:198:LEU:HD22	2.55	0.40
1:G:202:PRO:HG2	1:G:252:GLN:HG3	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	199/232 (86%)	196 (98%)	2 (1%)	1 (0%)	24	48
1	C	201/232 (87%)	194 (96%)	7 (4%)	0	100	100
1	E	199/232 (86%)	194 (98%)	5 (2%)	0	100	100
1	G	197/232 (85%)	195 (99%)	2 (1%)	0	100	100
3	I	212/216 (98%)	196 (92%)	15 (7%)	1 (0%)	24	48
3	J	208/216 (96%)	199 (96%)	9 (4%)	0	100	100
3	K	206/216 (95%)	193 (94%)	13 (6%)	0	100	100
3	L	205/216 (95%)	190 (93%)	14 (7%)	1 (0%)	24	48
4	M	212/217 (98%)	200 (94%)	12 (6%)	0	100	100
4	N	212/217 (98%)	200 (94%)	11 (5%)	1 (0%)	24	48
4	O	212/217 (98%)	206 (97%)	6 (3%)	0	100	100
4	P	212/217 (98%)	197 (93%)	15 (7%)	0	100	100
All	All	2475/2660 (93%)	2360 (95%)	111 (4%)	4 (0%)	43	68

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	I	26	GLY
4	N	33	ALA
1	A	140	PRO
3	L	64	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	187/212 (88%)	180 (96%)	7 (4%)	30	59
1	C	189/212 (89%)	185 (98%)	4 (2%)	47	75
1	E	187/212 (88%)	186 (100%)	1 (0%)	81	92
1	G	185/212 (87%)	181 (98%)	4 (2%)	45	74
3	I	177/179 (99%)	172 (97%)	5 (3%)	38	68

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
3	J	176/179 (98%)	170 (97%)	6 (3%)	32	62
3	K	174/179 (97%)	169 (97%)	5 (3%)	37	67
3	L	173/179 (97%)	167 (96%)	6 (4%)	32	61
4	M	189/192 (98%)	189 (100%)	0	100	100
4	N	189/192 (98%)	189 (100%)	0	100	100
4	O	189/192 (98%)	187 (99%)	2 (1%)	65	85
4	P	189/192 (98%)	189 (100%)	0	100	100
All	All	2204/2332 (94%)	2164 (98%)	40 (2%)	51	78

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	121	ILE
1	A	122	ARG
1	A	124	ASP
1	A	125	ILE
1	A	221	PRO
1	A	234	ARG
1	A	275	ARG
1	C	98	ASP
1	C	131	ARG
1	C	195	CYS
1	C	275	ARG
1	E	195	CYS
1	G	125	ILE
1	G	138	GLU
1	G	146	ILE
1	G	201	SER
3	I	25	SER
3	I	29	PHE
3	I	30	SER
3	I	96	CYS
3	I	98	ARG
3	J	48	VAL
3	J	50	TYR
3	J	64	VAL
3	J	71	SER
3	J	87	ARG
3	J	98	ARG
3	K	25	SER

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Mol	Chain	Res	Type
3	K	71	SER
3	K	76	LYS
3	K	98	ARG
3	K	192	SER
3	L	27	PHE
3	L	32	SER
3	L	96	CYS
3	L	98	ARG
3	L	100	ASP
3	L	101	SER
4	O	129	LYS
4	O	211	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	111	GLN
1	C	128	GLN
1	C	171	HIS
1	C	226	ASN
1	E	129	GLN
3	I	3	GLN
3	I	110	GLN
3	I	177	GLN
3	J	170	HIS
3	K	198	GLN
3	L	39	GLN
3	L	177	GLN
4	M	28	GLN
4	M	202	GLN
4	N	7	GLN
4	N	140	ASN
4	N	141	ASN
4	O	91	GLN
4	O	201	HIS
4	P	39	GLN
4	P	80	GLN
4	P	91	GLN
4	P	213	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 8 ligands modelled in this entry, 4 are monoatomic - leaving 4 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
5	OGA	C	1001	6	9,9,9	1.12	0	8,11,11	0.97	0
5	OGA	E	1001	6	9,9,9	1.11	0	8,11,11	1.04	0
5	OGA	A	1001	6	9,9,9	1.15	0	8,11,11	1.12	0
5	OGA	G	1001	6	9,9,9	1.13	0	8,11,11	1.16	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	OGA	C	1001	6	-	0/8/9/9	-
5	OGA	E	1001	6	-	2/8/9/9	-
5	OGA	A	1001	6	-	0/8/9/9	-
5	OGA	G	1001	6	-	0/8/9/9	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	E	1001	OGA	N1-C4-C5-O4
5	E	1001	OGA	N1-C4-C5-O3

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	G	1001	OGA	1	0

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	203/232 (87%)	0.25	10 (4%) 35 31	33, 45, 73, 92	1 (0%)
1	C	205/232 (88%)	0.78	19 (9%) 14 12	45, 63, 88, 108	0
1	E	203/232 (87%)	0.29	5 (2%) 58 55	35, 48, 71, 86	1 (0%)
1	G	201/232 (86%)	0.87	28 (13%) 6 5	41, 66, 94, 111	1 (0%)
2	B	5/6 (83%)	0.85	1 (20%) 3 2	66, 66, 73, 93	0
2	D	4/6 (66%)	1.16	1 (25%) 2 1	76, 77, 78, 82	0
2	F	5/6 (83%)	1.01	0 100 100	70, 71, 74, 93	0
2	H	4/6 (66%)	0.38	0 100 100	70, 75, 80, 80	0
3	I	214/216 (99%)	0.30	12 (5%) 30 26	30, 47, 84, 101	0
3	J	212/216 (98%)	0.19	5 (2%) 59 56	27, 47, 80, 93	1 (0%)
3	K	210/216 (97%)	0.34	7 (3%) 49 45	36, 52, 76, 86	0
3	L	209/216 (96%)	0.85	14 (6%) 24 21	38, 68, 91, 104	0
4	M	214/217 (98%)	0.18	0 100 100	28, 51, 66, 75	1 (0%)
4	N	214/217 (98%)	0.31	6 (2%) 55 51	31, 52, 70, 94	1 (0%)
4	O	214/217 (98%)	0.16	3 (1%) 73 72	35, 51, 64, 81	2 (0%)
4	P	214/217 (98%)	0.52	2 (0%) 81 80	41, 60, 76, 86	2 (0%)
All	All	2531/2684 (94%)	0.42	113 (4%) 38 34	27, 54, 84, 111	10 (0%)

All (113) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	I	141	THR	5.4
1	G	125	ILE	5.0
3	L	29	PHE	4.7
1	A	124	ASP	4.7
1	G	232	VAL	4.5

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Mol	Chain	Res	Type	RSRZ
1	G	146	ILE	4.4
3	L	142	ALA	4.2
1	G	143	TYR	4.0
1	G	264	HIS	4.0
1	A	144	SER	3.8
3	I	29	PHE	3.8
1	G	223	PRO	3.7
1	C	119	THR	3.6
4	N	3	ILE	3.6
4	P	3	ILE	3.4
3	L	192	SER	3.3
3	L	219	PRO	3.3
1	G	137	GLY	3.3
1	E	144	SER	3.2
1	C	146	ILE	3.2
3	I	140	GLY	3.2
3	K	107	TYR	3.2
3	I	2	VAL	3.1
3	K	2	VAL	3.1
1	C	144	SER	3.1
1	G	222	PRO	3.0
3	K	26	GLY	3.0
2	D	6	DT	3.0
4	N	214	ARG	3.0
1	G	141	TYR	2.9
1	G	82	LEU	2.9
3	K	29	PHE	2.9
1	G	144	SER	2.8
3	L	132	SER	2.8
1	C	264	HIS	2.8
3	I	220	LYS	2.7
1	G	279	PRO	2.7
4	P	4	GLN	2.7
1	C	82	LEU	2.7
3	L	28	ASN	2.7
1	C	141	TYR	2.6
1	C	126	THR	2.6
1	C	230	THR	2.6
1	E	227	GLY	2.6
3	I	142	ALA	2.6
3	J	142	ALA	2.6
1	C	199	GLY	2.6

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Mol	Chain	Res	Type	RSRZ
3	I	27	PHE	2.6
3	L	199	THR	2.6
1	E	127	TYR	2.6
1	C	195	CYS	2.6
1	G	265	SER	2.5
4	O	3	ILE	2.5
1	C	143	TYR	2.5
1	C	231	TYR	2.5
1	C	225	GLU	2.5
3	J	1	GLU	2.5
3	L	143	ALA	2.5
3	K	198	GLN	2.5
1	C	197	SER	2.5
1	G	140	PRO	2.4
1	A	123	GLU	2.4
1	G	195	CYS	2.4
3	L	2	VAL	2.4
1	G	121	ILE	2.3
4	N	172	LYS	2.3
3	I	133	SER	2.3
3	I	219	PRO	2.3
1	G	71	VAL	2.3
1	G	87	VAL	2.3
1	A	279	PRO	2.3
3	L	110	GLN	2.3
1	A	121	ILE	2.3
4	N	82	GLU	2.3
1	E	279	PRO	2.3
3	J	220	LYS	2.2
4	N	216	GLU	2.2
1	A	141	TYR	2.2
1	C	87	VAL	2.2
1	A	145	ARG	2.2
1	G	168	ASN	2.2
3	J	217	VAL	2.2
1	G	72	ILE	2.2
3	L	27	PHE	2.2
1	G	85	THR	2.2
1	G	142	THR	2.2
1	C	70	ARG	2.2
1	G	263	TYR	2.2
1	C	189	ASP	2.1

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Mol	Chain	Res	Type	RSRZ
4	N	27	SER	2.1
4	O	156	ALA	2.1
3	L	216	LYS	2.1
1	A	263	TYR	2.1
3	L	31	TYR	2.1
1	G	199	GLY	2.1
2	B	4	DT	2.1
3	I	30	SER	2.1
3	K	197	THR	2.1
1	A	223	PRO	2.1
3	L	131	PRO	2.1
1	G	127	TYR	2.1
1	E	145	ARG	2.1
1	C	198	LEU	2.1
1	A	125	ILE	2.1
1	G	200	ARG	2.0
3	I	130	ALA	2.0
3	I	197	THR	2.0
1	G	153	HIS	2.0
4	O	208	VAL	2.0
1	G	197	SER	2.0
3	J	195	LEU	2.0
1	C	200	ARG	2.0
3	K	31	TYR	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
5	OGA	G	1001	10/10	0.74	0.18	52,68,79,86	0
5	OGA	C	1001	10/10	0.79	0.18	54,66,79,83	0
5	OGA	A	1001	10/10	0.82	0.19	44,55,77,86	0
5	OGA	E	1001	10/10	0.83	0.18	44,55,58,69	0
6	MN	A	1002	1/1	0.99	0.03	46,46,46,46	0
6	MN	E	1002	1/1	0.99	0.03	44,44,44,44	0
6	MN	C	1002	1/1	1.00	0.02	61,61,61,61	0
6	MN	G	1002	1/1	1.00	0.02	68,68,68,68	0

6.5 Other polymers [i](#)

There are no such residues in this entry.