



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 06:09 AM UTC

PDB ID : 6RXM / pdb_00006rxm
Title : Crystal structure of CobB Ac2 (A76G, I131C, V162G) in complex with H4K16-Acetyl peptide
Authors : Spinck, M.; Gasper, R.; Neumann, H.
Deposited on : 2019-06-08
Resolution : 1.92 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

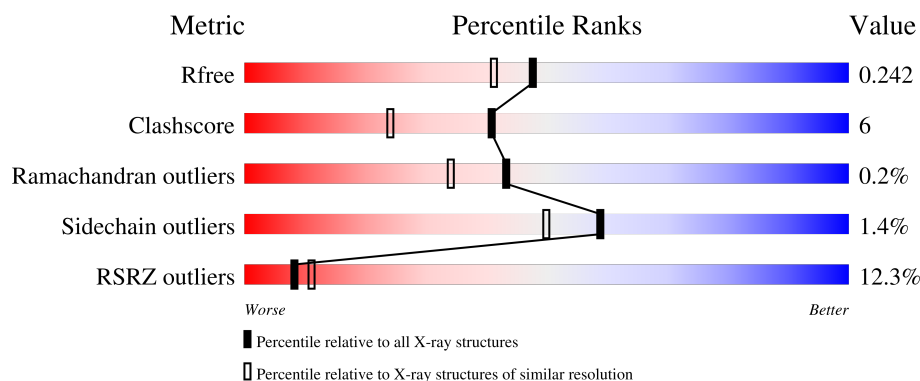
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.92 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	1188 (1.92-1.92)
Clashscore	190562	1209 (1.92-1.92)
Ramachandran outliers	187476	1195 (1.92-1.92)
Sidechain outliers	187428	1195 (1.92-1.92)
RSRZ outliers	180081	1188 (1.92-1.92)

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	254	8% 80% 11% 7%
1	B	254	15% 79% 13% 8%
1	C	254	6% 84% 8% 7%
1	D	254	11% 76% 13% 11%
1	E	254	6% 85% 6% 8%

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
1	F	254	22% 77% 13% 8%
2	G	11	9% 27% 36% 36%
2	H	11	9% 27% 27% 9% 36%
2	I	11	9% 36% 27% 36%
2	J	11	55% 18% 27%
2	K	11	9% 55% 18% 9% 18%
2	L	11	9% 64% 36%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 12179 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 NAD-dependent protein deacylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	235	Total	C	N	O	S	0	2	0
			1838	1163	325	339	11			
1	B	234	Total	C	N	O	S	0	1	0
			1824	1152	323	338	11			
1	C	235	Total	C	N	O	S	0	2	0
			1837	1161	325	341	10			
1	D	227	Total	C	N	O	S	0	0	0
			1763	1115	310	328	10			
1	E	233	Total	C	N	O	S	0	0	0
			1813	1144	322	337	10			
1	F	233	Total	C	N	O	S	0	0	0
			1813	1144	322	337	10			

There are 102 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-14	MET	-	initiating methionine	UNP P75960
A	-13	GLY	-	expression tag	UNP P75960
A	-12	SER	-	expression tag	UNP P75960
A	-11	SER	-	expression tag	UNP P75960
A	-10	HIS	-	expression tag	UNP P75960
A	-9	HIS	-	expression tag	UNP P75960
A	-8	HIS	-	expression tag	UNP P75960
A	-7	HIS	-	expression tag	UNP P75960
A	-6	HIS	-	expression tag	UNP P75960
A	-5	HIS	-	expression tag	UNP P75960
A	-4	SER	-	expression tag	UNP P75960
A	-3	GLN	-	expression tag	UNP P75960
A	-2	ASP	-	expression tag	UNP P75960
A	-1	PRO	-	expression tag	UNP P75960
A	76	GLY	ALA	engineered mutation	UNP P75960
A	131	CYS	ILE	engineered mutation	UNP P75960
A	161	ALA	VAL	engineered mutation	UNP P75960

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
B	-14	MET	-	initiating methionine	UNP P75960
B	-13	GLY	-	expression tag	UNP P75960
B	-12	SER	-	expression tag	UNP P75960
B	-11	SER	-	expression tag	UNP P75960
B	-10	HIS	-	expression tag	UNP P75960
B	-9	HIS	-	expression tag	UNP P75960
B	-8	HIS	-	expression tag	UNP P75960
B	-7	HIS	-	expression tag	UNP P75960
B	-6	HIS	-	expression tag	UNP P75960
B	-5	HIS	-	expression tag	UNP P75960
B	-4	SER	-	expression tag	UNP P75960
B	-3	GLN	-	expression tag	UNP P75960
B	-2	ASP	-	expression tag	UNP P75960
B	-1	PRO	-	expression tag	UNP P75960
B	76	GLY	ALA	engineered mutation	UNP P75960
B	131	CYS	ILE	engineered mutation	UNP P75960
B	161	ALA	VAL	engineered mutation	UNP P75960
C	-14	MET	-	initiating methionine	UNP P75960
C	-13	GLY	-	expression tag	UNP P75960
C	-12	SER	-	expression tag	UNP P75960
C	-11	SER	-	expression tag	UNP P75960
C	-10	HIS	-	expression tag	UNP P75960
C	-9	HIS	-	expression tag	UNP P75960
C	-8	HIS	-	expression tag	UNP P75960
C	-7	HIS	-	expression tag	UNP P75960
C	-6	HIS	-	expression tag	UNP P75960
C	-5	HIS	-	expression tag	UNP P75960
C	-4	SER	-	expression tag	UNP P75960
C	-3	GLN	-	expression tag	UNP P75960
C	-2	ASP	-	expression tag	UNP P75960
C	-1	PRO	-	expression tag	UNP P75960
C	76	GLY	ALA	engineered mutation	UNP P75960
C	131	CYS	ILE	engineered mutation	UNP P75960
C	161	ALA	VAL	engineered mutation	UNP P75960
D	-14	MET	-	initiating methionine	UNP P75960
D	-13	GLY	-	expression tag	UNP P75960
D	-12	SER	-	expression tag	UNP P75960
D	-11	SER	-	expression tag	UNP P75960
D	-10	HIS	-	expression tag	UNP P75960
D	-9	HIS	-	expression tag	UNP P75960
D	-8	HIS	-	expression tag	UNP P75960
D	-7	HIS	-	expression tag	UNP P75960

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
D	-6	HIS	-	expression tag	UNP P75960
D	-5	HIS	-	expression tag	UNP P75960
D	-4	SER	-	expression tag	UNP P75960
D	-3	GLN	-	expression tag	UNP P75960
D	-2	ASP	-	expression tag	UNP P75960
D	-1	PRO	-	expression tag	UNP P75960
D	76	GLY	ALA	engineered mutation	UNP P75960
D	131	CYS	ILE	engineered mutation	UNP P75960
D	161	ALA	VAL	engineered mutation	UNP P75960
E	-14	MET	-	initiating methionine	UNP P75960
E	-13	GLY	-	expression tag	UNP P75960
E	-12	SER	-	expression tag	UNP P75960
E	-11	SER	-	expression tag	UNP P75960
E	-10	HIS	-	expression tag	UNP P75960
E	-9	HIS	-	expression tag	UNP P75960
E	-8	HIS	-	expression tag	UNP P75960
E	-7	HIS	-	expression tag	UNP P75960
E	-6	HIS	-	expression tag	UNP P75960
E	-5	HIS	-	expression tag	UNP P75960
E	-4	SER	-	expression tag	UNP P75960
E	-3	GLN	-	expression tag	UNP P75960
E	-2	ASP	-	expression tag	UNP P75960
E	-1	PRO	-	expression tag	UNP P75960
E	76	GLY	ALA	engineered mutation	UNP P75960
E	131	CYS	ILE	engineered mutation	UNP P75960
E	161	ALA	VAL	engineered mutation	UNP P75960
F	-14	MET	-	initiating methionine	UNP P75960
F	-13	GLY	-	expression tag	UNP P75960
F	-12	SER	-	expression tag	UNP P75960
F	-11	SER	-	expression tag	UNP P75960
F	-10	HIS	-	expression tag	UNP P75960
F	-9	HIS	-	expression tag	UNP P75960
F	-8	HIS	-	expression tag	UNP P75960
F	-7	HIS	-	expression tag	UNP P75960
F	-6	HIS	-	expression tag	UNP P75960
F	-5	HIS	-	expression tag	UNP P75960
F	-4	SER	-	expression tag	UNP P75960
F	-3	GLN	-	expression tag	UNP P75960
F	-2	ASP	-	expression tag	UNP P75960
F	-1	PRO	-	expression tag	UNP P75960
F	76	GLY	ALA	engineered mutation	UNP P75960
F	131	CYS	ILE	engineered mutation	UNP P75960

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
F	161	ALA	VAL	engineered mutation	UNP P75960

- Molecule 2 is a protein called Histone H4.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	G	7	Total	C	N	O	0	0	0
			57	33	16	8			
2	H	7	Total	C	N	O	0	1	0
			65	38	19	8			
2	I	7	Total	C	N	O	0	0	0
			57	33	16	8			
2	J	8	Total	C	N	O	0	0	0
			66	39	18	9			
2	K	9	Total	C	N	O	0	0	0
			74	45	19	10			
2	L	11	Total	C	N	O	0	0	0
			87	54	21	12			

- Molecule 3 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Zn	0	0
			1	1		
3	B	1	Total	Zn	0	0
			1	1		
3	C	1	Total	Zn	0	0
			1	1		
3	D	1	Total	Zn	0	0
			1	1		
3	E	1	Total	Zn	0	0
			1	1		
3	F	1	Total	Zn	0	0
			1	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	161	Total	O	0	0
			161	161		
4	B	89	Total	O	0	0
			89	89		

Continued on next page...

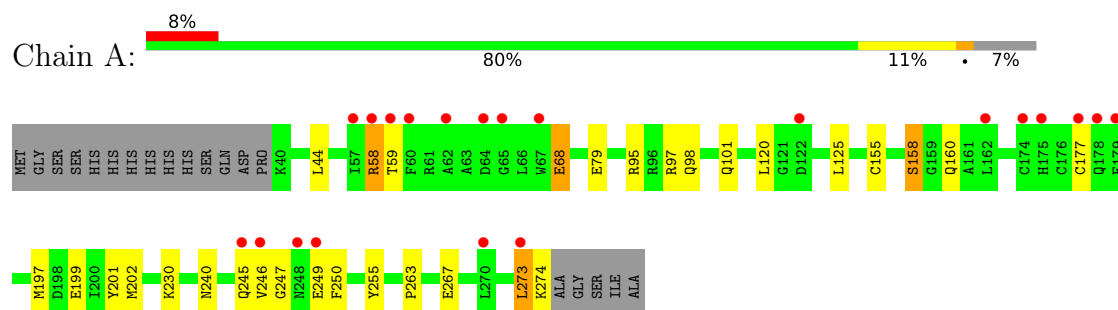
Continued from previous page...

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	C	185	Total 185	O 185	0	0
4	D	85	Total 85	O 85	0	0
4	E	211	Total 211	O 211	0	0
4	F	95	Total 95	O 95	0	0
4	G	7	Total 7	O 7	0	0
4	H	7	Total 7	O 7	0	0
4	I	10	Total 10	O 10	0	0
4	J	7	Total 7	O 7	0	0
4	K	10	Total 10	O 10	0	0
4	L	12	Total 12	O 12	0	0

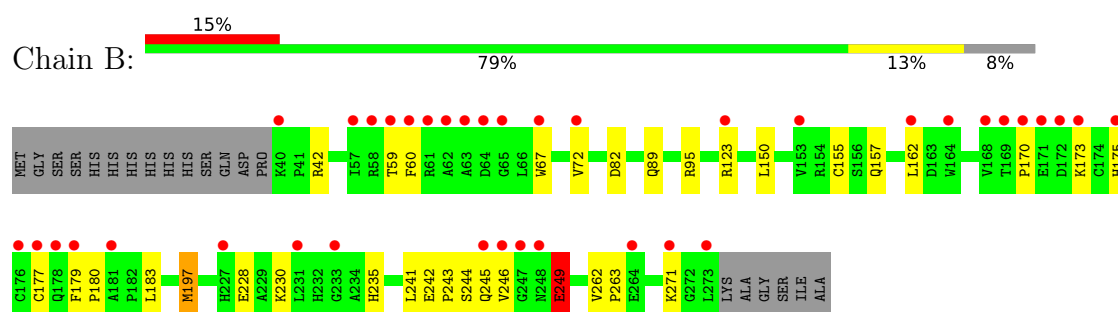
3 Residue-property plots [i](#)

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.

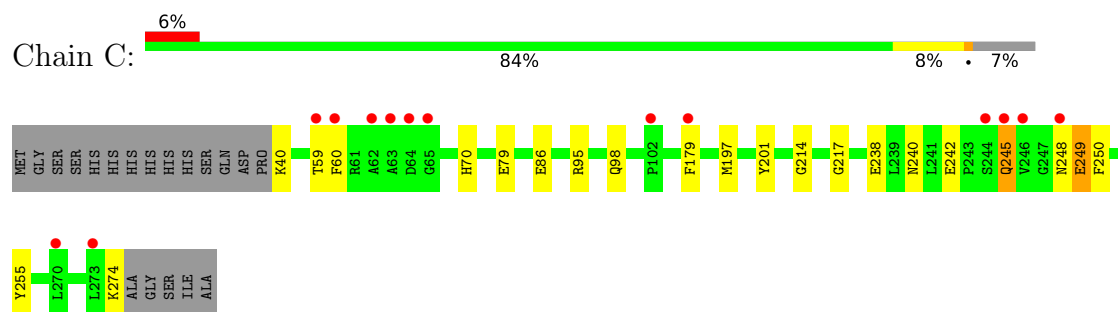
- Molecule 1: NAD-dependent protein deacetylase



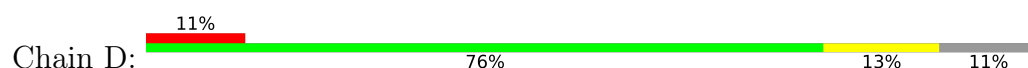
- Molecule 1: NAD-dependent protein deacetylase

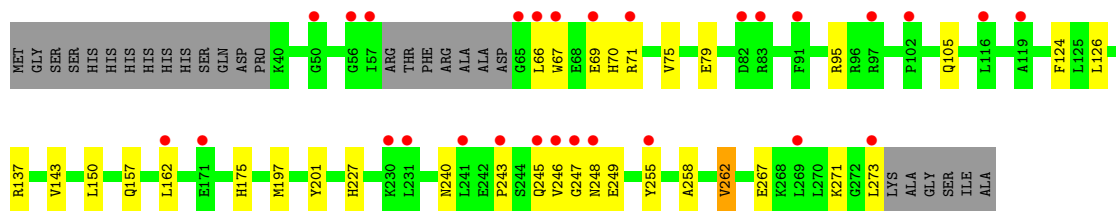


- Molecule 1: NAD-dependent protein deacetylase

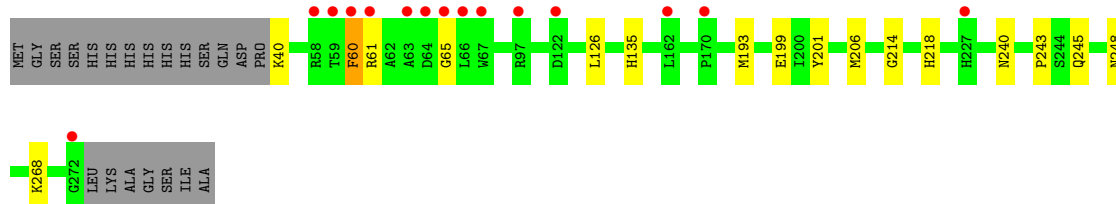
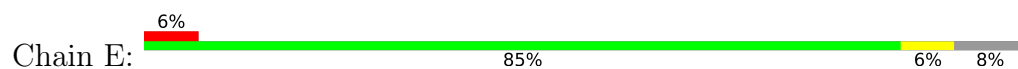


- Molecule 1: NAD-dependent protein deacetylase

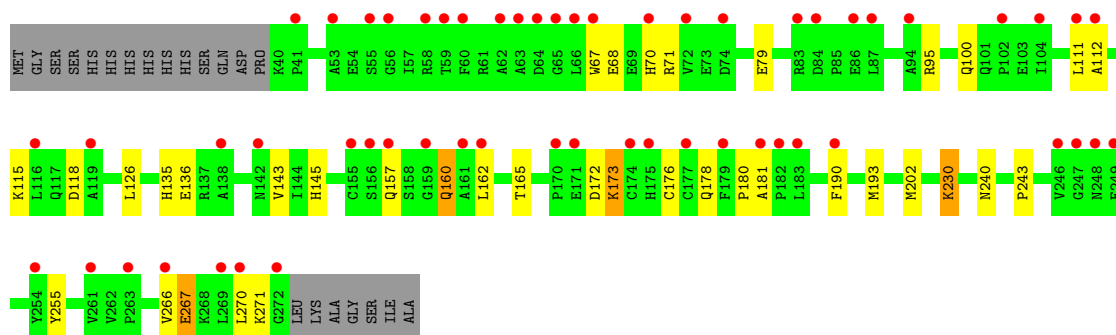
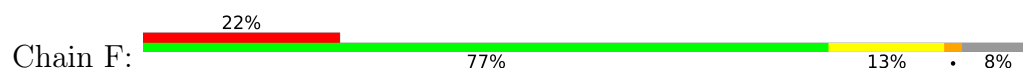




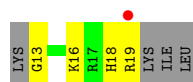
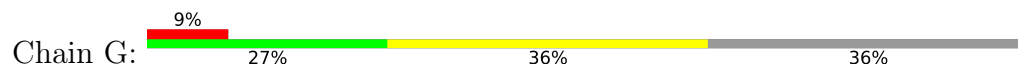
• Molecule 1: NAD-dependent protein deacylase



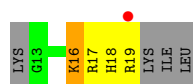
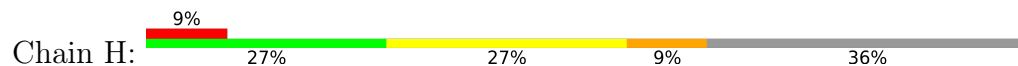
• Molecule 1: NAD-dependent protein deacylase



• Molecule 2: Histone H4

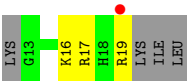


• Molecule 2: Histone H4

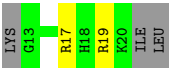


• Molecule 2: Histone H4

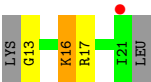




● Molecule 2: Histone H4



● Molecule 2: Histone H4



● Molecule 2: Histone H4



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	92.30Å 95.38Å 168.83Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.47 – 1.92 48.47 – 1.92	Depositor EDS
% Data completeness (in resolution range)	100.0 (48.47-1.92) 100.0 (48.47-1.92)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.91 (at 1.92Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.204 , 0.242 0.207 , 0.242	Depositor DCC
R_{free} test set	1998 reflections (1.75%)	wwPDB-VP
Wilson B-factor (Å ²)	32.3	Xtriage
Anisotropy	0.191	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 36.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.055 for k,h,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	12179	wwPDB-VP
Average B, all atoms (Å ²)	41.0	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 38.63 % 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 3.5796e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: ALY, ZN

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	0.61	0/1889	0.93	3/2564 (0.1%)
1	B	0.59	0/1872	0.93	4/2542 (0.2%)
1	C	0.63	0/1888	0.91	3/2563 (0.1%)
1	D	0.52	0/1806	0.88	0/2452
1	E	0.65	0/1858	0.89	0/2523
1	F	0.58	0/1858	0.87	0/2523
2	G	0.44	0/44	0.79	0/54
2	H	0.35	0/54	0.71	0/65
2	I	0.51	0/44	0.85	0/54
2	J	0.40	0/53	0.82	0/65
2	K	0.40	0/61	0.56	0/76
2	L	0.52	0/74	0.62	0/94
All	All	0.60	0/11501	0.90	10/15575 (0.1%)

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
2	H	0	1
2	K	0	1
All	All	0	2

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	179	PHE	CA-C-N	8.22	128.71	119.83
1	B	179	PHE	C-N-CA	8.22	128.71	119.83

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	247	GLY	N-CA-C	7.99	121.22	110.43
1	A	58	ARG	N-CA-C	-7.21	99.78	110.23
1	A	177	CYS	N-CA-C	6.97	120.86	109.85
1	C	179	PHE	CA-C-N	6.78	127.18	119.93
1	C	179	PHE	C-N-CA	6.78	127.18	119.93
1	B	249	GLU	N-CA-C	-6.59	105.24	113.28
1	C	249	GLU	N-CA-C	-5.54	106.37	113.23
1	B	241	LEU	N-CA-C	-5.09	105.19	111.40

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	H	16	ALY	Mainchain
2	K	16	ALY	Mainchain

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	1838	0	1792	20	0
1	B	1824	0	1769	30	0
1	C	1837	0	1788	17	0
1	D	1763	0	1706	34	0
1	E	1813	0	1752	17	0
1	F	1813	0	1752	29	0
2	G	57	0	57	2	0
2	H	65	0	70	3	0
2	I	57	0	57	2	0
2	J	66	0	70	2	0
2	K	74	0	81	3	0
2	L	87	0	94	5	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
3	E	1	0	0	0	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	F	1	0	0	0	0
4	A	161	0	0	3	2
4	B	89	0	0	4	1
4	C	185	0	0	3	1
4	D	85	0	0	2	0
4	E	211	0	0	3	0
4	F	95	0	0	1	0
4	G	7	0	0	1	0
4	H	7	0	0	1	0
4	I	10	0	0	0	0
4	J	7	0	0	0	0
4	K	10	0	0	1	0
4	L	12	0	0	0	0
All	All	12179	0	10988	144	2

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 6.

All (144) 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:79:GLU:OE2	2:K:17:ARG:NH2	1.98	0.96
1:C:248:ASN:HD21	1:E:61:ARG:HH12	1.11	0.93
1:B:170:PRO:O	1:B:173:LYS:NZ	2.03	0.89
1:D:69:GLU:OE2	1:F:95:ARG:NH2	2.11	0.83
1:D:67:TRP:CD1	1:D:70:HIS:HD1	1.99	0.81
1:B:157:GLN:NE2	1:B:177:CYS:SG	2.55	0.80
1:D:79:GLU:OE2	2:I:17:ARG:NH1	2.14	0.79
1:F:230:LYS:NZ	4:F:401:HOH:O	2.08	0.78
1:D:67:TRP:HE1	1:D:69:GLU:HB2	1.50	0.77
1:A:230:LYS:HE3	1:A:249:GLU:HA	1.66	0.76
1:B:246:VAL:HG21	1:D:245:GLN:HG3	1.68	0.75
1:E:201:TYR:HH	2:K:13:GLY:N	1.84	0.75
1:D:67:TRP:NE1	1:D:69:GLU:HB2	2.01	0.74
1:A:101:GLN:NE2	4:A:402:HOH:O	2.22	0.73
2:G:13:GLY:N	4:G:101:HOH:O	2.23	0.72
1:A:79:GLU:OE1	2:H:17:ARG:NH2	2.16	0.72
1:B:230:LYS:NZ	1:B:249:GLU:HA	2.03	0.72
1:D:69:GLU:HG2	1:F:95:ARG:HH22	1.54	0.70
1:B:155:CYS:SG	1:B:157:GLN:HG3	2.32	0.69
1:B:60:PHE:CD2	1:B:72:VAL:HG12	2.28	0.69

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:248:ASN:ND2	1:E:61:ARG:HH12	1.87	0.69
1:A:245:GLN:HG3	1:A:249:GLU:HG2	1.75	0.68
1:A:97:ARG:NH1	4:A:404:HOH:O	2.27	0.67
1:B:60:PHE:HD2	1:B:72:VAL:HG12	1.60	0.66
1:A:249:GLU:N	1:A:249:GLU:OE1	2.28	0.66
1:C:79:GLU:OE1	2:J:17:ARG:NH2	2.24	0.65
1:B:82:ASP:OD1	4:B:401:HOH:O	2.13	0.65
1:B:230:LYS:HZ3	1:B:249:GLU:HA	1.61	0.64
2:K:13:GLY:N	4:K:101:HOH:O	2.31	0.62
1:B:60:PHE:CZ	1:B:67:TRP:CE3	2.91	0.59
1:B:95:ARG:HB3	1:B:150:LEU:HD21	1.86	0.58
1:D:227:HIS:HD2	1:D:249:GLU:OE1	1.86	0.58
1:B:60:PHE:CD2	1:B:67:TRP:HB2	2.38	0.57
1:C:249:GLU:OE1	1:C:249:GLU:N	2.32	0.57
1:D:246:VAL:C	1:D:248:ASN:H	2.13	0.56
1:F:190:PHE:HD1	2:L:17:ARG:HA	1.71	0.56
1:F:126:LEU:HD23	1:F:143:VAL:HG22	1.88	0.55
1:B:197:MET:HE2	1:B:197:MET:HA	1.87	0.55
1:C:217:GLY:O	2:I:19:ARG:NH2	2.40	0.55
1:A:58:ARG:HD2	1:A:95:ARG:HE	1.72	0.55
1:D:267:GLU:O	1:D:271:LYS:HD2	2.06	0.54
1:D:267:GLU:OE1	1:D:271:LYS:NZ	2.36	0.54
1:A:58:ARG:HE	1:A:98:GLN:HG3	1.71	0.54
1:A:199:GLU:HA	1:A:202:MET:HE2	1.89	0.53
1:F:67:TRP:HZ2	1:F:190:PHE:CE2	2.27	0.53
1:A:245:GLN:HG3	1:A:249:GLU:CG	2.38	0.53
1:C:197:MET:HE3	1:C:201:TYR:CE1	2.44	0.53
1:E:268:LYS:HE2	4:E:423:HOH:O	2.08	0.53
1:E:40:LYS:N	4:E:411:HOH:O	2.41	0.53
1:C:70:HIS:HD2	4:C:556:HOH:O	1.91	0.53
1:C:86:GLU:HG2	4:C:428:HOH:O	2.09	0.53
1:D:258:ALA:O	1:D:262:VAL:HG13	2.09	0.52
1:F:172:ASP:O	1:F:173:LYS:HD3	2.09	0.52
1:D:67:TRP:HD1	1:D:70:HIS:HD1	1.51	0.52
1:C:248:ASN:HD21	1:E:61:ARG:NH1	1.93	0.52
1:D:67:TRP:CD1	1:D:69:GLU:H	2.28	0.51
1:D:95:ARG:HB3	1:D:150:LEU:HD21	1.93	0.50
1:B:230:LYS:HZ1	1:B:249:GLU:HA	1.76	0.50
1:D:240:ASN:O	1:D:255:TYR:HA	2.12	0.50
1:D:71:ARG:HG3	1:F:71:ARG:HG2	1.94	0.50
1:B:157:GLN:HE22	1:B:177:CYS:HB2	1.78	0.49

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:95:ARG:HH21	1:C:98:GLN:NE2	2.11	0.48
1:F:67:TRP:HZ2	1:F:190:PHE:CD2	2.30	0.48
1:D:246:VAL:O	1:D:248:ASN:N	2.46	0.48
1:A:120:LEU:HD21	1:A:273:LEU:HD11	1.95	0.48
1:E:193:MET:HG2	1:F:193:MET:HE3	1.95	0.48
1:A:249:GLU:H	1:A:249:GLU:CD	2.21	0.48
1:B:245:GLN:NE2	1:B:249:GLU:HG3	2.29	0.48
1:F:160:GLN:OE1	1:F:176:CYS:SG	2.73	0.47
2:H:18:HIS:ND1	2:H:19[A]:ARG:HG3	2.29	0.47
1:C:255:TYR:CG	1:E:243:PRO:HD3	2.49	0.47
1:E:65:GLY:HA2	2:L:22:LEU:CD1	2.45	0.47
1:E:60:PHE:CD1	1:E:60:PHE:N	2.77	0.47
2:J:19:ARG:O	2:J:19:ARG:HG2	2.15	0.47
1:D:105:GLN:OE1	1:D:137:ARG:NH2	2.49	0.46
1:D:66:LEU:CD1	1:D:75:VAL:HG23	2.45	0.46
1:B:95:ARG:HA	1:B:95:ARG:HD3	1.83	0.46
1:A:263:PRO:O	1:A:267:GLU:HG3	2.17	0.45
1:B:245:GLN:HE22	1:B:249:GLU:HG3	1.81	0.45
1:D:67:TRP:CG	1:D:95:ARG:HH21	2.34	0.45
1:E:214:GLY:N	1:E:240:ASN:OD1	2.47	0.45
1:B:235:HIS:ND1	4:B:408:HOH:O	2.36	0.45
1:C:245:GLN:H	1:C:245:GLN:CD	2.19	0.45
1:F:100:GLN:HE22	1:F:165:THR:HB	1.81	0.45
1:E:218:HIS:NE2	1:E:245:GLN:HG3	2.32	0.45
1:A:160:GLN:OE1	4:A:401:HOH:O	2.20	0.45
1:D:67:TRP:CD1	1:D:70:HIS:ND1	2.77	0.45
1:B:228:GLU:OE2	4:B:402:HOH:O	2.21	0.45
1:B:59:THR:OG1	1:B:60:PHE:N	2.50	0.44
1:F:240:ASN:O	1:F:255:TYR:HA	2.17	0.44
1:E:126:LEU:HD21	1:E:135:HIS:CG	2.53	0.44
1:F:202:MET:HE3	1:F:202:MET:HB2	1.77	0.44
1:D:197:MET:HE3	1:D:201:TYR:CE1	2.53	0.44
1:F:115:LYS:O	1:F:118:ASP:HB2	2.17	0.44
1:F:126:LEU:HD21	1:F:135:HIS:CG	2.53	0.44
1:F:266:VAL:O	1:F:270:LEU:HG	2.18	0.43
1:D:126:LEU:HD23	1:D:143:VAL:HG22	2.01	0.43
1:E:206:MET:HE3	1:E:206:MET:HB3	1.84	0.43
1:F:157:GLN:NE2	1:F:181:ALA:HB2	2.34	0.43
2:G:18:HIS:O	2:G:19:ARG:HB2	2.18	0.43
2:L:21:ILE:HG22	2:L:22:LEU:HD23	2.00	0.43
1:F:267:GLU:O	1:F:271:LYS:HG3	2.19	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:89:GLN:NE2	1:B:183:LEU:O	2.49	0.43
1:B:173:LYS:HB3	1:B:180:PRO:HB3	2.01	0.42
1:D:227:HIS:CD2	1:D:249:GLU:OE1	2.71	0.42
1:F:162:LEU:HD23	1:F:162:LEU:HA	1.85	0.42
1:E:65:GLY:HA2	2:L:22:LEU:HD12	2.01	0.42
1:E:199:GLU:OE1	4:E:401:HOH:O	2.22	0.42
1:A:44[A]:LEU:HD13	1:A:125:LEU:HB3	2.02	0.42
1:B:42:ARG:HG2	1:B:123:ARG:HD3	2.01	0.42
1:A:155:CYS:SG	1:A:158:SER:HB3	2.60	0.42
1:C:242:GLU:HG2	1:E:248:ASN:HD21	1.84	0.42
1:D:175:HIS:HA	4:D:440:HOH:O	2.18	0.42
1:F:112:ALA:O	1:F:115:LYS:HB3	2.20	0.42
2:H:19[A]:ARG:NH2	4:H:101:HOH:O	2.40	0.42
1:D:249:GLU:HG2	1:D:249:GLU:O	2.20	0.42
1:F:79:GLU:CD	1:F:79:GLU:H	2.27	0.42
1:A:197:MET:HE3	1:A:201:TYR:CE1	2.54	0.41
1:D:67:TRP:HB2	1:D:95:ARG:HH21	1.85	0.41
1:C:98:GLN:HG3	4:C:454:HOH:O	2.19	0.41
1:F:136:GLU:OE2	1:F:145:HIS:NE2	2.43	0.41
1:F:178:GLN:O	1:F:180:PRO:HD3	2.20	0.41
1:C:238:GLU:HB2	1:C:250:PHE:CG	2.56	0.41
1:A:59:THR:HG22	1:A:68:GLU:HG3	2.02	0.41
1:D:243:PRO:HA	1:D:255:TYR:CE1	2.56	0.41
1:F:68:GLU:OE2	1:F:70:HIS:HD2	2.03	0.41
1:A:240:ASN:O	1:A:255:TYR:HA	2.21	0.41
1:D:69:GLU:CG	1:F:95:ARG:HH22	2.30	0.41
1:D:124:PHE:CZ	1:D:126:LEU:HB2	2.55	0.41
1:D:157:GLN:HE21	1:D:157:GLN:HB2	1.63	0.41
1:D:162:LEU:HD22	1:D:175:HIS:CE1	2.56	0.41
2:L:21:ILE:HD13	2:L:21:ILE:HA	1.87	0.41
1:B:244:SER:O	4:B:403:HOH:O	2.22	0.41
1:C:214:GLY:N	1:C:240:ASN:OD1	2.53	0.41
1:A:245:GLN:CD	1:A:250:PHE:HD2	2.29	0.40
1:B:157:GLN:HE22	1:B:177:CYS:CB	2.34	0.40
1:D:273:LEU:O	4:D:401:HOH:O	2.22	0.40
1:B:242:GLU:HG3	1:B:243:PRO:HD2	2.02	0.40
1:B:262:VAL:HB	1:B:263:PRO:HD3	2.02	0.40
1:C:59:THR:HG23	1:C:60:PHE:O	2.20	0.40
1:B:95:ARG:CB	1:B:150:LEU:HD21	2.49	0.40
1:F:67:TRP:CZ2	1:F:190:PHE:CE2	3.07	0.40
1:F:243:PRO:HA	1:F:255:TYR:CE1	2.57	0.40

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:162:LEU:HD21	1:B:175:HIS:CE1	2.57	0.40

All (2) 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)
4:A:405:HOH:O	4:B:413:HOH:O[3_544]	2.10	0.10
4:A:431:HOH:O	4:C:424:HOH:O[3_644]	2.12	0.08

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	235/254 (92%)	231 (98%)	3 (1%)	1 (0%)	30	19
1	B	233/254 (92%)	231 (99%)	2 (1%)	0	100	100
1	C	235/254 (92%)	231 (98%)	3 (1%)	1 (0%)	30	19
1	D	223/254 (88%)	221 (99%)	1 (0%)	1 (0%)	30	19
1	E	231/254 (91%)	229 (99%)	2 (1%)	0	100	100
1	F	231/254 (91%)	228 (99%)	3 (1%)	0	100	100
2	G	4/11 (36%)	4 (100%)	0	0	100	100
2	H	4/11 (36%)	4 (100%)	0	0	100	100
2	I	4/11 (36%)	4 (100%)	0	0	100	100
2	J	5/11 (46%)	5 (100%)	0	0	100	100
2	K	6/11 (54%)	6 (100%)	0	0	100	100
2	L	8/11 (73%)	8 (100%)	0	0	100	100
All	All	1419/1590 (89%)	1402 (99%)	14 (1%)	3 (0%)	43	34

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	245	GLN
1	A	68	GLU
1	D	247	GLY

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	194/207 (94%)	190 (98%)	4 (2%)	47	33
1	B	192/207 (93%)	189 (98%)	3 (2%)	55	44
1	C	194/207 (94%)	192 (99%)	2 (1%)	68	64
1	D	186/207 (90%)	185 (100%)	1 (0%)	81	80
1	E	190/207 (92%)	189 (100%)	1 (0%)	81	80
1	F	190/207 (92%)	185 (97%)	5 (3%)	40	24
2	G	3/7 (43%)	3 (100%)	0	100	100
2	H	4/7 (57%)	4 (100%)	0	100	100
2	I	3/7 (43%)	3 (100%)	0	100	100
2	J	4/7 (57%)	4 (100%)	0	100	100
2	K	5/7 (71%)	5 (100%)	0	100	100
2	L	6/7 (86%)	6 (100%)	0	100	100
All	All	1171/1284 (91%)	1155 (99%)	16 (1%)	59	50

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	158	SER
1	A	246	VAL
1	A	273	LEU
1	A	274	LYS
1	B	197	MET
1	B	249	GLU
1	B	271	LYS
1	C	40	LYS

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	C	274	LYS
1	D	262	VAL
1	E	60	PHE
1	F	111	LEU
1	F	160	GLN
1	F	173	LYS
1	F	230	LYS
1	F	267	GLU

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

Mol	Chain	Res	Type
1	A	105	GLN
1	A	129	GLN
1	A	160	GLN
1	A	248	ASN
1	B	70	HIS
1	B	157	GLN
1	B	160	GLN
1	B	175	HIS
1	B	245	GLN
1	C	129	GLN
1	C	175	HIS
1	C	248	ASN
1	D	101	GLN
1	D	129	GLN
1	D	157	GLN
1	D	160	GLN
1	D	175	HIS
1	D	227	HIS
1	E	175	HIS
1	E	248	ASN
1	F	70	HIS
1	F	100	GLN
1	F	157	GLN
1	F	218	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

6 non-standard protein/DNA/RNA residues are modelled in this entry.

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
2	ALY	I	16	2	10,11,12	0.83	0	7,12,14	0.93	1 (14%)
2	ALY	G	16	2	10,11,12	1.58	1 (10%)	7,12,14	1.94	2 (28%)
2	ALY	K	16	2	10,11,12	1.49	1 (10%)	7,12,14	1.38	1 (14%)
2	ALY	H	16	2	10,11,12	1.54	1 (10%)	7,12,14	1.13	1 (14%)
2	ALY	L	16	2	10,11,12	1.58	1 (10%)	7,12,14	0.99	1 (14%)
2	ALY	J	16	2	10,11,12	0.83	0	7,12,14	0.86	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
2	ALY	I	16	2	-	0/9/10/12	-
2	ALY	G	16	2	-	0/9/10/12	-
2	ALY	K	16	2	-	0/9/10/12	-
2	ALY	H	16	2	-	0/9/10/12	-
2	ALY	L	16	2	-	0/9/10/12	-
2	ALY	J	16	2	-	0/9/10/12	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	G	16	ALY	O-C	4.43	1.36	1.20
2	L	16	ALY	O-C	4.30	1.36	1.20
2	H	16	ALY	O-C	4.08	1.35	1.20
2	K	16	ALY	O-C	4.07	1.35	1.20

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
2	G	16	ALY	CE-NZ-CH	-3.86	116.86	122.56
2	G	16	ALY	CD-CG-CB	-2.50	104.18	113.62
2	H	16	ALY	CD-CG-CB	-2.49	104.25	113.62
2	K	16	ALY	CD-CG-CB	-2.38	104.64	113.62
2	I	16	ALY	CD-CG-CB	-2.36	104.74	113.62
2	L	16	ALY	CD-CG-CB	-2.30	104.94	113.62

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 6 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	235/254 (92%)	0.62	21 (8%) 15 19	17, 34, 72, 84	2 (0%)
1	B	234/254 (92%)	1.04	38 (16%) 4 6	19, 40, 80, 97	1 (0%)
1	C	235/254 (92%)	0.46	14 (5%) 27 32	19, 32, 65, 92	2 (0%)
1	D	227/254 (89%)	0.97	28 (12%) 8 11	24, 43, 71, 93	0
1	E	233/254 (91%)	0.29	15 (6%) 25 29	20, 29, 51, 88	0
1	F	233/254 (91%)	1.25	56 (24%) 2 2	22, 44, 73, 99	0
2	G	6/11 (54%)	0.83	1 (16%) 4 6	27, 33, 39, 58	0
2	H	6/11 (54%)	0.83	1 (16%) 4 6	26, 31, 43, 47	1 (16%)
2	I	6/11 (54%)	0.78	1 (16%) 4 6	23, 29, 42, 59	0
2	J	7/11 (63%)	0.72	0 100 100	26, 34, 55, 59	0
2	K	8/11 (72%)	0.81	1 (12%) 8 10	26, 34, 54, 62	0
2	L	10/11 (90%)	0.59	1 (10%) 12 16	26, 37, 54, 55	0
All	All	1440/1590 (90%)	0.77	177 (12%) 8 11	17, 37, 71, 99	6 (0%)

All (177) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	60	PHE	6.9
1	B	59	THR	5.9
1	E	272	GLY	5.3
1	A	273	LEU	5.1
1	D	246	VAL	5.1
1	E	59	THR	5.0
1	F	272	GLY	5.0
1	C	65	GLY	4.8
1	E	64	ASP	4.8
1	B	60	PHE	4.7
1	B	273	LEU	4.7

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	246	VAL	4.6
1	F	62	ALA	4.5
1	B	179	PHE	4.5
1	A	246	VAL	4.5
1	D	65	GLY	4.5
1	C	246	VAL	4.4
1	C	60	PHE	4.3
1	C	63	ALA	4.2
1	D	67	TRP	4.2
1	A	177	CYS	4.1
1	B	62	ALA	4.1
1	B	176	CYS	4.1
1	A	179	PHE	4.1
1	A	59	THR	4.0
1	F	175	HIS	3.9
1	B	67	TRP	3.9
1	F	111	LEU	3.9
1	F	60	PHE	3.9
1	F	162	LEU	3.7
1	D	247	GLY	3.7
1	A	270	LEU	3.7
1	A	248	ASN	3.7
1	F	55	SER	3.7
1	A	245	GLN	3.7
1	C	62	ALA	3.6
1	F	112	ALA	3.6
1	D	57	ILE	3.5
1	F	183	LEU	3.5
1	C	64	ASP	3.5
1	B	248	ASN	3.5
1	B	40	LYS	3.5
1	F	171	GLU	3.4
1	C	59	THR	3.4
1	D	269	LEU	3.4
1	D	273	LEU	3.4
1	E	65	GLY	3.3
1	D	66	LEU	3.3
1	F	190	PHE	3.3
1	B	233	GLY	3.3
1	B	178	GLN	3.2
1	B	63	ALA	3.1
1	F	246	VAL	3.1

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	58	ARG	3.1
1	B	177	CYS	3.1
1	B	61	ARG	3.1
1	F	66	LEU	3.1
1	A	60	PHE	3.1
1	F	179	PHE	3.1
1	F	65	GLY	3.0
1	D	231	LEU	3.0
1	B	172	ASP	3.0
1	F	269	LEU	2.9
1	B	247	GLY	2.9
1	F	56	GLY	2.9
2	K	21	ILE	2.9
1	B	58	ARG	2.9
1	C	270	LEU	2.9
2	I	19	ARG	2.9
1	F	87	LEU	2.9
1	A	175	HIS	2.8
1	B	123	ARG	2.8
1	D	56	GLY	2.8
1	B	169	THR	2.8
1	C	245	GLN	2.8
1	F	247	GLY	2.8
1	F	63	ALA	2.8
1	E	67	TRP	2.8
1	B	245	GLN	2.7
1	B	173	LYS	2.7
2	G	19	ARG	2.7
1	F	116	LEU	2.7
1	F	155	CYS	2.6
1	C	273	LEU	2.6
1	E	66	LEU	2.6
1	F	266	VAL	2.6
1	F	181	ALA	2.6
1	D	248	ASN	2.6
1	C	248	ASN	2.6
1	F	263	PRO	2.6
1	F	104	ILE	2.6
1	F	261	VAL	2.6
1	F	59	THR	2.6
1	B	175	HIS	2.5
1	D	243	PRO	2.5

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	E	58	ARG	2.5
1	A	162	LEU	2.5
1	F	270	LEU	2.5
1	D	230	LYS	2.5
1	F	70	HIS	2.5
1	F	83	ARG	2.5
1	B	72	VAL	2.5
1	D	245	GLN	2.5
1	D	171	GLU	2.5
1	F	41	PRO	2.5
1	F	159	GLY	2.5
2	L	22	LEU	2.5
1	A	57	ILE	2.5
1	E	63	ALA	2.5
1	B	64	ASP	2.5
1	D	116	LEU	2.4
1	F	64	ASP	2.4
1	A	65	GLY	2.4
1	B	227	HIS	2.4
1	B	181	ALA	2.4
1	F	72	VAL	2.4
1	F	74	ASP	2.3
1	B	168	VAL	2.3
1	C	244	SER	2.3
1	D	82	ASP	2.3
1	F	58	ARG	2.3
1	F	157	GLN	2.3
1	C	179	PHE	2.3
1	B	171	GLU	2.3
1	A	178	GLN	2.3
1	F	67	TRP	2.3
1	F	53	ALA	2.3
1	F	119	ALA	2.3
1	E	61	ARG	2.3
1	B	264	GLU	2.3
1	D	162	LEU	2.3
1	B	164	TRP	2.2
1	F	161	ALA	2.2
1	F	156	SER	2.2
1	A	64	ASP	2.2
1	B	57	ILE	2.2
1	D	50	GLY	2.2

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	F	170	PRO	2.2
1	F	182	PRO	2.2
1	A	249	GLU	2.2
1	D	241	LEU	2.2
1	A	67	TRP	2.2
1	E	97	ARG	2.2
1	B	231	LEU	2.2
1	F	142	ASN	2.1
1	A	122	ASP	2.1
1	B	153	VAL	2.1
1	A	174	CYS	2.1
1	B	271	LYS	2.1
1	D	119	ALA	2.1
1	F	94	ALA	2.1
1	D	97	ARG	2.1
2	H	19[A]	ARG	2.1
1	C	102	PRO	2.1
1	E	227	HIS	2.1
1	F	102	PRO	2.1
1	F	174	CYS	2.1
1	F	248	ASN	2.1
1	E	122	ASP	2.1
1	F	254	TYR	2.1
1	D	83	ARG	2.1
1	F	249	GLU	2.1
1	B	170	PRO	2.1
1	E	170	PRO	2.1
1	B	162	LEU	2.1
1	F	138	ALA	2.1
1	F	84	ASP	2.1
1	D	69	GLU	2.0
1	D	91	PHE	2.0
1	F	86	GLU	2.0
1	D	255	TYR	2.0
1	E	162	LEU	2.0
1	F	177	CYS	2.0
1	A	62	ALA	2.0
1	B	65	GLY	2.0
1	D	102	PRO	2.0
1	D	71	ARG	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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
2	ALY	L	16	12/13	0.91	0.09	22,27,36,45	0
2	ALY	H	16	12/13	0.95	0.07	21,24,33,39	0
2	ALY	I	16	12/13	0.95	0.07	18,23,35,40	0
2	ALY	J	16	12/13	0.95	0.08	22,27,38,39	0
2	ALY	K	16	12/13	0.95	0.07	23,25,35,39	0
2	ALY	G	16	12/13	0.95	0.07	16,23,31,33	0

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
3	ZN	B	301	1/1	0.95	0.06	59,59,59,59	0
3	ZN	A	300	1/1	0.97	0.07	51,51,51,51	0
3	ZN	F	301	1/1	0.97	0.06	54,54,54,54	0
3	ZN	D	301	1/1	0.99	0.03	40,40,40,40	0
3	ZN	E	301	1/1	0.99	0.03	28,28,28,28	0
3	ZN	C	300	1/1	0.99	0.04	41,41,41,41	0

6.5 Other polymers [i](#)

There are no such residues in this entry.