



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 03:01 AM UTC

PDB ID : 3EWF / pdb_00003ewf
Title : Crystal Structure Analysis of human HDAC8 H143A variant complexed with substrate.
Authors : Dowling, D.P.; Gantt, S.L.; Gattis, S.G.; Fierke, C.A.; Christianson, D.W.
Deposited on : 2008-10-14
Resolution : 2.50 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

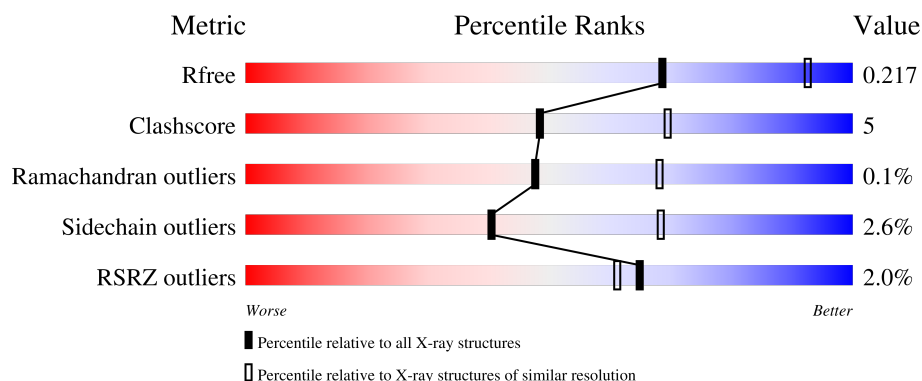
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	388	 2% 80% 12% • 6%
1	B	388	 2% 79% 13% • 6%
1	C	388	 3% 78% 14% • 6%
1	D	388	 % 79% 13% • 6%
2	I	5	 80% 20%

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Mol	Chain	Length	Quality of chain
2	J	5	 80% 20%
2	K	5	 40% 60%
2	L	5	 80% 20%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	MCM	I	6	-	X	-	-
5	MCM	J	6	-	X	-	-

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 11893 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 Histone deacetylase 8.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	364	Total	C	N	O	S	0	0	0
			2834	1816	470	529	19			
1	B	364	Total	C	N	O	S	6	0	0
			2834	1816	470	529	19			
1	C	364	Total	C	N	O	S	5	0	0
			2834	1816	470	529	19			
1	D	364	Total	C	N	O	S	1	0	0
			2834	1816	470	529	19			

There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	143	ALA	HIS	engineered mutation	UNP Q9BY41
A	378	ILE	-	expression tag	UNP Q9BY41
A	379	GLU	-	expression tag	UNP Q9BY41
A	380	GLY	-	expression tag	UNP Q9BY41
A	381	ARG	-	expression tag	UNP Q9BY41
A	382	SER	-	expression tag	UNP Q9BY41
A	383	HIS	-	expression tag	UNP Q9BY41
A	384	HIS	-	expression tag	UNP Q9BY41
A	385	HIS	-	expression tag	UNP Q9BY41
A	386	HIS	-	expression tag	UNP Q9BY41
A	387	HIS	-	expression tag	UNP Q9BY41
A	388	HIS	-	expression tag	UNP Q9BY41
B	143	ALA	HIS	engineered mutation	UNP Q9BY41
B	378	ILE	-	expression tag	UNP Q9BY41
B	379	GLU	-	expression tag	UNP Q9BY41
B	380	GLY	-	expression tag	UNP Q9BY41
B	381	ARG	-	expression tag	UNP Q9BY41
B	382	SER	-	expression tag	UNP Q9BY41
B	383	HIS	-	expression tag	UNP Q9BY41
B	384	HIS	-	expression tag	UNP Q9BY41
B	385	HIS	-	expression tag	UNP Q9BY41

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Chain	Residue	Modelled	Actual	Comment	Reference
B	386	HIS	-	expression tag	UNP Q9BY41
B	387	HIS	-	expression tag	UNP Q9BY41
B	388	HIS	-	expression tag	UNP Q9BY41
C	143	ALA	HIS	engineered mutation	UNP Q9BY41
C	378	ILE	-	expression tag	UNP Q9BY41
C	379	GLU	-	expression tag	UNP Q9BY41
C	380	GLY	-	expression tag	UNP Q9BY41
C	381	ARG	-	expression tag	UNP Q9BY41
C	382	SER	-	expression tag	UNP Q9BY41
C	383	HIS	-	expression tag	UNP Q9BY41
C	384	HIS	-	expression tag	UNP Q9BY41
C	385	HIS	-	expression tag	UNP Q9BY41
C	386	HIS	-	expression tag	UNP Q9BY41
C	387	HIS	-	expression tag	UNP Q9BY41
C	388	HIS	-	expression tag	UNP Q9BY41
D	143	ALA	HIS	engineered mutation	UNP Q9BY41
D	378	ILE	-	expression tag	UNP Q9BY41
D	379	GLU	-	expression tag	UNP Q9BY41
D	380	GLY	-	expression tag	UNP Q9BY41
D	381	ARG	-	expression tag	UNP Q9BY41
D	382	SER	-	expression tag	UNP Q9BY41
D	383	HIS	-	expression tag	UNP Q9BY41
D	384	HIS	-	expression tag	UNP Q9BY41
D	385	HIS	-	expression tag	UNP Q9BY41
D	386	HIS	-	expression tag	UNP Q9BY41
D	387	HIS	-	expression tag	UNP Q9BY41
D	388	HIS	-	expression tag	UNP Q9BY41

- Molecule 2 is a protein called PEPTIDIC SUBSTRATE.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	I	5	Total	C	N	O	0	0	0
			48	30	11	7			
2	J	5	Total	C	N	O	0	0	0
			48	30	11	7			
2	K	5	Total	C	N	O	0	0	0
			48	30	11	7			
2	L	5	Total	C	N	O	0	0	0
			48	30	11	7			

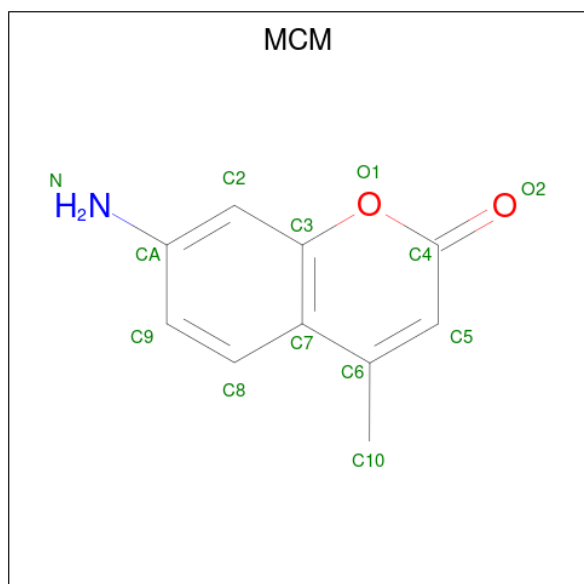
- Molecule 3 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	2	Total 2	Zn 2	0	0
3	B	1	Total 1	Zn 1	0	0
3	C	2	Total 2	Zn 2	0	0
3	D	1	Total 1	Zn 1	0	0

- Molecule 4 is POTASSIUM ION (CCD ID: K) (formula: K).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	2	Total 2	K 2	0	0
4	B	2	Total 2	K 2	0	0
4	C	2	Total 2	K 2	0	0
4	D	2	Total 2	K 2	0	0

- Molecule 5 is 7-AMINO-4-METHYL-CHROMEN-2-ONE (CCD ID: MCM) (formula: C₁₀H₉NO₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	I	1	Total 13	C 10	N 1	O 2	0	0

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	J	1	Total	C	N	O	0	0
			13	10	1	2		
5	K	1	Total	C	N	O	0	0
			13	10	1	2		
5	L	1	Total	C	N	O	0	0
			13	10	1	2		

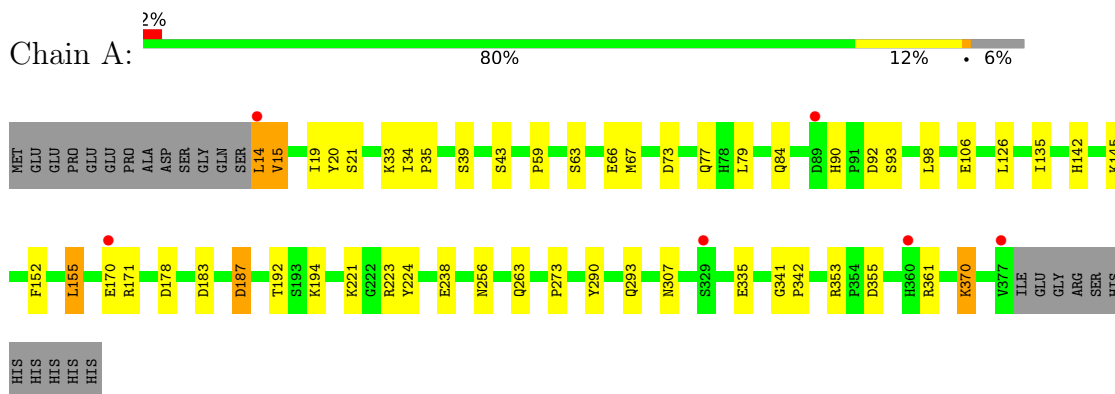
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	85	Total	O	0	0
			85	85		
6	B	82	Total	O	0	0
			82	82		
6	C	65	Total	O	0	0
			65	65		
6	D	58	Total	O	0	0
			58	58		
6	I	4	Total	O	0	0
			4	4		
6	J	3	Total	O	0	0
			3	3		
6	K	1	Total	O	0	0
			1	1		
6	L	1	Total	O	0	0
			1	1		

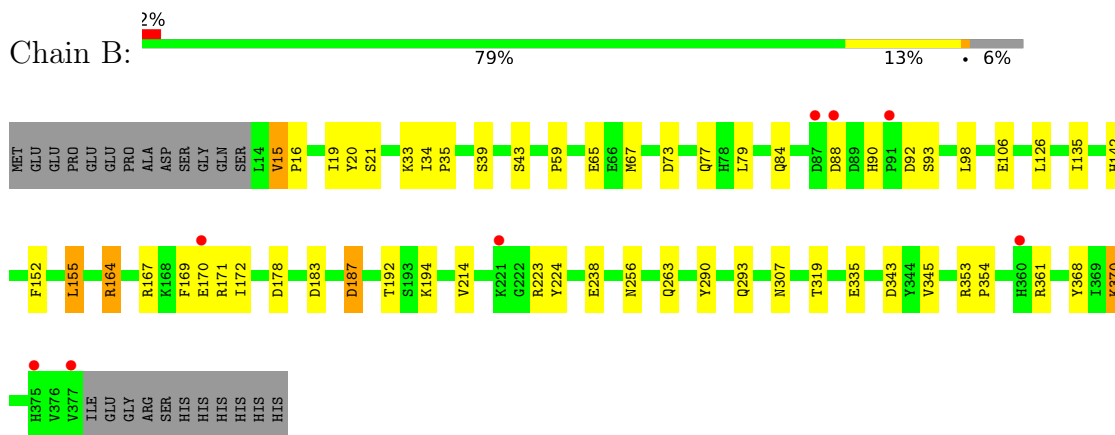
3 Residue-property plots [i](#)

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.

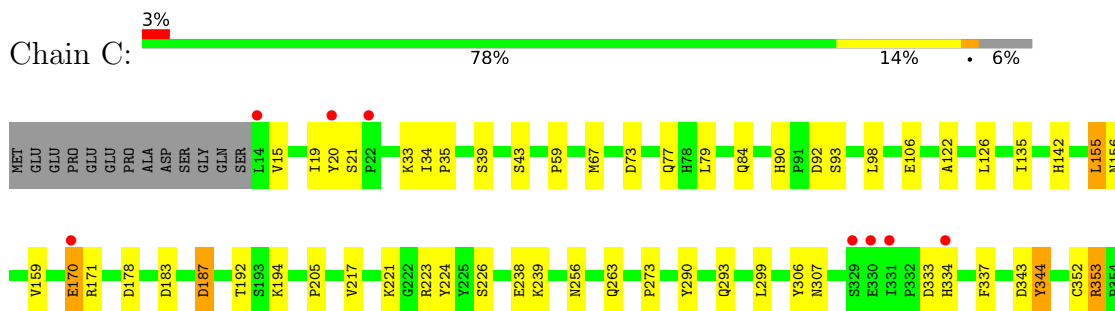
• Molecule 1: Histone deacetylase 8

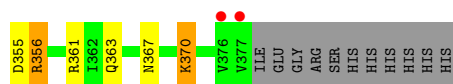


• Molecule 1: Histone deacetylase 8

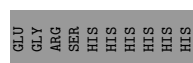
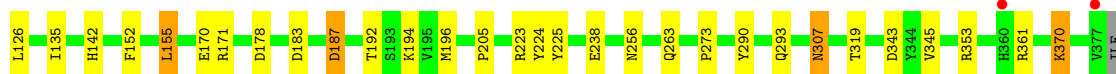
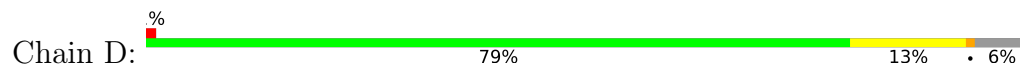


• Molecule 1: Histone deacetylase 8

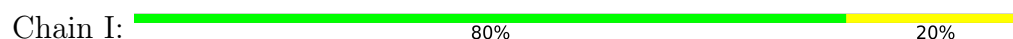




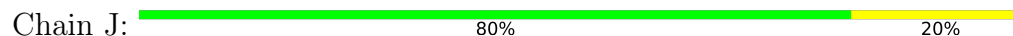
- Molecule 1: Histone deacetylase 8



- Molecule 2: PEPTIDIC SUBSTRATE



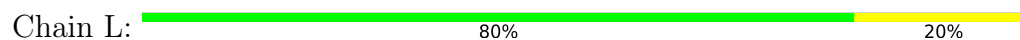
- Molecule 2: PEPTIDIC SUBSTRATE



- Molecule 2: PEPTIDIC SUBSTRATE



- Molecule 2: PEPTIDIC SUBSTRATE



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	82.90Å 91.84Å 196.58Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.15 – 2.50 49.15 – 2.50	Depositor EDS
% Data completeness (in resolution range)	93.6 (49.15-2.50) 99.4 (49.15-2.50)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.69 (at 2.48Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.198 , 0.229 0.206 , 0.217	Depositor DCC
R_{free} test set	2680 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å ²)	29.0	Xtriage
Anisotropy	0.595	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 42.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	11893	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.95% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ALY, MCM, ACE, ZN, K

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	0.44	0/2904	0.94	10/3941 (0.3%)
1	B	0.45	0/2904	0.95	15/3941 (0.4%)
1	C	0.49	3/2904 (0.1%)	0.94	12/3941 (0.3%)
1	D	0.43	0/2904	0.92	11/3941 (0.3%)
2	I	0.33	0/22	0.69	0/28
2	J	0.20	0/22	0.57	0/28
2	K	0.32	0/22	0.64	0/28
2	L	0.23	0/22	0.46	0/28
All	All	0.45	3/11704 (0.0%)	0.94	48/15876 (0.3%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	363	GLN	CD-OE1	-7.67	1.08	1.23
1	C	363	GLN	CG-CD	-6.95	1.34	1.52
1	C	367	ASN	CB-CG	-5.93	1.37	1.52

All (48) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	164	ARG	NE-CZ-NH2	-7.25	112.68	119.20
1	D	135	ILE	N-CA-C	6.87	117.78	108.17
1	B	15	VAL	N-CA-C	6.66	114.36	108.63
1	C	135	ILE	N-CA-C	6.64	117.41	108.11
1	B	135	ILE	N-CA-C	6.54	117.33	108.17
1	B	353	ARG	CA-C-N	6.53	126.48	119.76
1	B	353	ARG	C-N-CA	6.53	126.48	119.76
1	A	135	ILE	N-CA-C	6.47	117.17	108.11
1	C	353	ARG	CA-C-N	6.42	126.39	119.78
1	C	353	ARG	C-N-CA	6.42	126.39	119.78

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	353	ARG	CA-C-N	6.34	126.31	119.78
1	A	353	ARG	C-N-CA	6.34	126.31	119.78
1	C	15	VAL	CA-C-N	6.22	126.17	119.76
1	C	15	VAL	C-N-CA	6.22	126.17	119.76
1	D	307	ASN	N-CA-C	-6.21	95.39	107.62
1	B	307	ASN	N-CA-C	-6.12	95.56	107.62
1	A	142	HIS	N-CA-C	6.08	120.31	113.01
1	C	142	HIS	N-CA-C	5.92	120.11	113.01
1	A	335	GLU	N-CA-C	5.91	118.48	111.33
1	D	142	HIS	N-CA-C	5.87	120.06	113.01
1	C	307	ASN	N-CA-C	-5.84	96.11	107.62
1	B	142	HIS	N-CA-C	5.74	119.89	113.01
1	D	353	ARG	CA-C-N	5.73	125.69	119.78
1	D	353	ARG	C-N-CA	5.73	125.69	119.78
1	D	15	VAL	CA-C-N	5.70	125.71	119.90
1	D	15	VAL	C-N-CA	5.70	125.71	119.90
1	B	164	ARG	NE-CZ-NH1	5.64	127.14	121.50
1	A	307	ASN	N-CA-C	-5.58	96.13	107.41
1	B	354	PRO	N-CA-C	5.51	119.73	111.14
1	D	361	ARG	NE-CZ-NH2	5.48	124.13	119.20
1	B	164	ARG	CD-NE-CZ	5.46	132.04	124.40
1	B	319	THR	N-CA-C	-5.43	105.45	111.36
1	B	335	GLU	N-CA-C	5.41	117.87	111.33
1	C	238	GLU	N-CA-C	5.37	116.82	110.97
1	D	238	GLU	N-CA-C	5.36	116.82	110.97
1	A	15	VAL	CA-C-N	5.35	125.36	119.90
1	A	15	VAL	C-N-CA	5.35	125.36	119.90
1	D	361	ARG	NE-CZ-NH1	-5.29	116.21	121.50
1	A	355	ASP	N-CA-C	-5.21	99.30	108.20
1	B	238	GLU	N-CA-C	5.20	116.64	111.07
1	C	356	ARG	NE-CZ-NH2	-5.20	114.52	119.20
1	B	361	ARG	NE-CZ-NH1	-5.16	116.34	121.50
1	A	238	GLU	N-CA-C	5.16	116.59	110.97
1	D	319	THR	N-CA-C	-5.14	105.76	111.36
1	C	355	ASP	N-CA-C	-5.05	99.56	108.20
1	C	170	GLU	N-CA-C	5.03	116.57	111.14
1	C	306	TYR	CA-CB-CG	-5.01	104.87	113.90
1	B	361	ARG	NE-CZ-NH2	5.01	123.71	119.20

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2834	0	2786	32	0
1	B	2834	0	2785	28	0
1	C	2834	0	2786	35	0
1	D	2834	0	2784	34	0
2	I	48	0	51	1	0
2	J	48	0	51	1	0
2	K	48	0	51	4	0
2	L	48	0	50	1	0
3	A	2	0	0	0	0
3	B	1	0	0	0	0
3	C	2	0	0	0	0
3	D	1	0	0	0	0
4	A	2	0	0	0	0
4	B	2	0	0	0	0
4	C	2	0	0	0	0
4	D	2	0	0	0	0
5	I	13	0	7	1	0
5	J	13	0	7	2	0
5	K	13	0	7	2	0
5	L	13	0	7	2	0
6	A	85	0	0	2	0
6	B	82	0	0	1	0
6	C	65	0	0	1	0
6	D	58	0	0	2	0
6	I	4	0	0	0	0
6	J	3	0	0	0	0
6	K	1	0	0	1	0
6	L	1	0	0	0	0
All	All	11893	0	11372	125	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 5.

All (125) 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:B:192:THR:HG22	1:B:194:LYS:H	1.45	0.80
1:D:192:THR:HG22	1:D:194:LYS:H	1.47	0.80
1:C:334:HIS:H	1:C:337:PHE:HB2	1.45	0.79
1:A:33:LYS:HG3	5:J:6:MCM:H5	1.66	0.78
1:A:192:THR:HG22	1:A:194:LYS:H	1.54	0.73
1:C:192:THR:HG22	1:C:194:LYS:H	1.54	0.73
1:C:333:ASP:HA	1:C:337:PHE:CD2	2.27	0.70
1:B:73:ASP:O	1:B:77:GLN:HG2	1.92	0.69
1:B:33:LYS:HG3	5:I:6:MCM:H5	1.75	0.68
1:C:170:GLU:OE2	1:C:170:GLU:HA	1.93	0.67
1:A:73:ASP:O	1:A:77:GLN:HG2	1.95	0.66
1:C:33:LYS:HG3	5:L:6:MCM:H5	1.79	0.63
1:A:221:LYS:HB3	2:K:1:ACE:H3	1.81	0.63
1:D:73:ASP:O	1:D:77:GLN:HG2	1.99	0.63
1:C:73:ASP:O	1:C:77:GLN:HG2	2.00	0.61
1:D:90:HIS:HD2	1:D:92:ASP:H	1.49	0.60
1:C:343:ASP:O	1:C:344:TYR:HB2	2.01	0.60
1:C:90:HIS:HD2	1:C:92:ASP:H	1.49	0.59
6:C:885:HOH:O	2:K:2:ARG:HG2	2.03	0.58
1:B:90:HIS:HD2	1:B:92:ASP:H	1.51	0.58
1:D:90:HIS:CD2	1:D:92:ASP:H	2.21	0.58
1:D:92:ASP:HA	6:D:1023:HOH:O	2.03	0.58
1:A:90:HIS:HD2	1:A:92:ASP:H	1.51	0.58
1:A:84:GLN:NE2	1:A:84:GLN:HA	2.19	0.57
1:C:67:MET:SD	1:C:155:LEU:HD21	2.44	0.57
1:A:90:HIS:CD2	1:A:92:ASP:H	2.23	0.57
1:B:67:MET:SD	1:B:155:LEU:HD21	2.45	0.57
1:D:290:TYR:O	1:D:293:GLN:HG3	2.04	0.57
1:C:205:PRO:HA	6:K:994:HOH:O	2.04	0.57
1:B:90:HIS:CD2	1:B:92:ASP:H	2.23	0.56
1:D:67:MET:SD	1:D:155:LEU:HD21	2.45	0.56
1:D:273:PRO:HB2	5:K:6:MCM:H9	1.87	0.56
1:C:90:HIS:CD2	1:C:92:ASP:H	2.23	0.56
1:D:171:ARG:NH1	1:D:256:ASN:O	2.39	0.56
1:A:290:TYR:O	1:A:293:GLN:HG3	2.07	0.55
1:C:290:TYR:O	1:C:293:GLN:HG3	2.06	0.55
1:D:170:GLU:OE2	1:D:170:GLU:HA	2.07	0.55
1:A:170:GLU:OE2	1:A:170:GLU:HA	2.06	0.55
1:A:341:GLY:C	2:J:1:ACE:H3	2.32	0.55
1:B:290:TYR:O	1:B:293:GLN:HG3	2.06	0.55
1:C:178:ASP:HB2	1:C:263:GLN:OE1	2.07	0.54
1:A:93:SER:HB3	1:A:98:LEU:HB2	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:93:SER:HB3	1:D:98:LEU:HB2	1.90	0.53
1:C:352:CYS:O	1:D:205:PRO:HG2	2.08	0.53
1:D:33:LYS:HG3	5:K:6:MCM:H5	1.91	0.53
1:D:223:ARG:O	1:D:224:TYR:HB2	2.09	0.53
1:D:84:GLN:NE2	1:D:84:GLN:HA	2.24	0.52
1:A:370:LYS:NZ	1:A:370:LYS:HB2	2.25	0.52
1:A:342:PRO:HD2	6:A:912:HOH:O	2.09	0.52
1:B:170:GLU:OE2	1:B:170:GLU:HA	2.08	0.52
1:B:84:GLN:NE2	1:B:84:GLN:HA	2.24	0.52
1:B:343:ASP:OD1	1:C:192:THR:HG23	2.10	0.52
1:A:171:ARG:NH1	1:A:256:ASN:O	2.42	0.51
1:B:178:ASP:HB2	1:B:263:GLN:OE1	2.10	0.51
1:B:214:VAL:HG22	1:B:368:TYR:HD2	1.76	0.51
1:C:93:SER:HB3	1:C:98:LEU:HB2	1.93	0.51
1:B:93:SER:HB3	1:B:98:LEU:HB2	1.93	0.51
1:A:223:ARG:O	1:A:224:TYR:HB2	2.09	0.51
1:A:67:MET:SD	1:A:155:LEU:HD21	2.51	0.50
1:C:84:GLN:NE2	1:C:84:GLN:HA	2.26	0.50
1:A:14:LEU:N	1:A:14:LEU:HD13	2.27	0.50
1:B:223:ARG:O	1:B:224:TYR:HB2	2.11	0.50
1:C:171:ARG:NH1	1:C:256:ASN:O	2.45	0.49
1:A:34:ILE:N	1:A:35:PRO:HD3	2.27	0.49
1:C:170:GLU:OE2	1:C:170:GLU:CA	2.58	0.49
1:D:29:ASP:OD2	1:D:38:ALA:HB3	2.13	0.49
1:D:152:PHE:CE1	2:L:5:ALY:HD2	2.47	0.49
1:B:370:LYS:HB2	1:B:370:LYS:NZ	2.27	0.49
1:C:223:ARG:O	1:C:224:TYR:HB2	2.12	0.48
1:D:370:LYS:HB2	1:D:370:LYS:NZ	2.28	0.48
1:C:221:LYS:HB3	2:I:1:ACE:H3	1.94	0.48
1:A:192:THR:HG23	1:D:343:ASP:OD1	2.14	0.47
1:D:34:ILE:N	1:D:35:PRO:HD3	2.29	0.47
1:B:34:ILE:N	1:B:35:PRO:HD3	2.30	0.47
1:A:178:ASP:HB2	1:A:263:GLN:OE1	2.15	0.47
1:C:183:ASP:O	1:C:187:ASP:HB2	2.14	0.47
1:D:178:ASP:HB2	1:D:263:GLN:OE1	2.14	0.47
1:D:39:SER:O	1:D:43:SER:HB2	2.15	0.47
1:A:183:ASP:O	1:A:187:ASP:HB2	2.15	0.47
1:C:34:ILE:N	1:C:35:PRO:HD3	2.29	0.46
1:C:273:PRO:HB2	5:L:6:MCM:H9	1.97	0.46
1:C:20:TYR:CG	1:C:21:SER:N	2.83	0.46
1:A:39:SER:O	1:A:43:SER:HB2	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:370:LYS:HB2	1:C:370:LYS:NZ	2.30	0.46
1:D:307:ASN:HA	2:K:4:ALY:HD3	1.97	0.46
1:B:183:ASP:O	1:B:187:ASP:HB2	2.16	0.46
1:B:345:VAL:HG11	1:C:194:LYS:HE2	1.97	0.46
1:B:171:ARG:NH1	1:B:256:ASN:O	2.48	0.46
1:D:20:TYR:CG	1:D:21:SER:N	2.85	0.45
1:C:19:ILE:CG2	1:C:59:PRO:HB3	2.47	0.45
1:A:194:LYS:HE2	1:D:345:VAL:HG11	1.99	0.45
1:B:20:TYR:CG	1:B:21:SER:N	2.85	0.45
1:D:183:ASP:O	1:D:187:ASP:HB2	2.18	0.44
1:B:15:VAL:HA	1:B:16:PRO:HD3	1.90	0.44
1:C:39:SER:O	1:C:43:SER:HB2	2.17	0.44
6:D:917:HOH:O	2:K:4:ALY:HH33	2.18	0.43
1:D:192:THR:HG22	1:D:194:LYS:N	2.26	0.43
1:C:334:HIS:O	1:C:337:PHE:HB2	2.17	0.43
1:D:19:ILE:CG2	1:D:59:PRO:HB3	2.49	0.43
1:B:39:SER:O	1:B:43:SER:HB2	2.18	0.43
1:D:63:SER:OG	1:D:66:GLU:HG3	2.19	0.43
1:A:20:TYR:CG	1:A:21:SER:N	2.86	0.43
1:C:353:ARG:HD3	1:D:205:PRO:O	2.19	0.42
1:B:192:THR:HG22	1:B:194:LYS:N	2.24	0.42
1:A:19:ILE:CG2	1:A:59:PRO:HB3	2.49	0.42
1:C:217:VAL:C	1:C:226:SER:HB2	2.45	0.42
1:A:145:LYS:NZ	6:A:924:HOH:O	2.53	0.42
1:C:337:PHE:O	1:C:344:TYR:HE1	2.03	0.42
1:A:84:GLN:HA	1:A:84:GLN:HE21	1.84	0.42
1:B:88:ASP:HA	6:B:834:HOH:O	2.19	0.42
1:B:19:ILE:CG2	1:B:59:PRO:HB3	2.49	0.42
1:A:273:PRO:HB2	5:J:6:MCM:H9	2.02	0.42
1:B:164:ARG:HD2	1:B:167:ARG:HD2	2.02	0.42
1:D:60:LYS:O	1:D:117:GLY:HA3	2.20	0.42
1:C:156:ASN:OD1	1:C:159:VAL:HG23	2.20	0.41
1:B:169:PHE:HB2	1:B:172:ILE:HD11	2.03	0.41
1:D:152:PHE:N	1:D:152:PHE:CD2	2.89	0.41
1:A:63:SER:OG	1:A:66:GLU:HG3	2.21	0.41
1:D:29:ASP:OD1	1:D:39:SER:OG	2.33	0.41
1:A:171:ARG:NH1	1:A:256:ASN:HB3	2.36	0.41
1:B:152:PHE:CD2	1:B:152:PHE:N	2.88	0.41
1:C:122:ALA:HB1	1:C:299:LEU:HD13	2.03	0.41
1:A:152:PHE:CD2	1:A:152:PHE:N	2.88	0.40
1:A:84:GLN:HE21	1:A:84:GLN:CA	2.33	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:196:MET:HE2	1:D:225:TYR:HA	2.03	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	362/388 (93%)	350 (97%)	12 (3%)	0	100	100
1	B	362/388 (93%)	351 (97%)	11 (3%)	0	100	100
1	C	362/388 (93%)	348 (96%)	13 (4%)	1 (0%)	36	55
1	D	362/388 (93%)	349 (96%)	13 (4%)	0	100	100
2	I	2/5 (40%)	2 (100%)	0	0	100	100
2	J	2/5 (40%)	2 (100%)	0	0	100	100
2	K	2/5 (40%)	2 (100%)	0	0	100	100
2	L	2/5 (40%)	2 (100%)	0	0	100	100
All	All	1456/1572 (93%)	1406 (97%)	49 (3%)	1 (0%)	48	68

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	344	TYR

5.3.2 Protein sidechains [i](#)

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	304/325 (94%)	295 (97%)	9 (3%)	36	64
1	B	304/325 (94%)	297 (98%)	7 (2%)	44	72
1	C	304/325 (94%)	295 (97%)	9 (3%)	36	64
1	D	304/325 (94%)	297 (98%)	7 (2%)	44	72
2	I	2/2 (100%)	2 (100%)	0	100	100
2	J	2/2 (100%)	2 (100%)	0	100	100
2	K	2/2 (100%)	2 (100%)	0	100	100
2	L	2/2 (100%)	2 (100%)	0	100	100
All	All	1224/1308 (94%)	1192 (97%)	32 (3%)	40	68

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	14	LEU
1	A	15	VAL
1	A	79	LEU
1	A	106	GLU
1	A	126	LEU
1	A	155	LEU
1	A	187	ASP
1	A	361	ARG
1	A	370	LYS
1	B	65	GLU
1	B	79	LEU
1	B	106	GLU
1	B	126	LEU
1	B	155	LEU
1	B	187	ASP
1	B	370	LYS
1	C	79	LEU
1	C	106	GLU
1	C	126	LEU
1	C	155	LEU
1	C	187	ASP
1	C	239	LYS
1	C	356	ARG
1	C	361	ARG
1	C	370	LYS

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Mol	Chain	Res	Type
1	D	15	VAL
1	D	79	LEU
1	D	106	GLU
1	D	126	LEU
1	D	155	LEU
1	D	187	ASP
1	D	370	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	84	GLN
1	A	90	HIS
1	A	242	GLN
1	A	256	ASN
1	B	80	GLN
1	B	84	GLN
1	B	90	HIS
1	B	242	GLN
1	B	256	ASN
1	C	80	GLN
1	C	84	GLN
1	C	90	HIS
1	C	242	GLN
1	C	256	ASN
1	D	84	GLN
1	D	90	HIS
1	D	256	ASN
2	L	3	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

8 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	5	3,5,2	10,11,12	0.43	0	7,12,14	0.62	0
2	ALY	J	4	2	10,11,12	0.43	0	7,12,14	0.55	0
2	ALY	J	5	3,5,2	10,11,12	0.43	0	7,12,14	0.54	0
2	ALY	K	4	2	10,11,12	0.47	0	7,12,14	0.81	0
2	ALY	L	4	2	10,11,12	0.42	0	7,12,14	0.78	0
2	ALY	L	5	3,5,2	10,11,12	0.46	0	7,12,14	0.57	0
2	ALY	I	4	2	10,11,12	0.39	0	7,12,14	0.70	0
2	ALY	K	5	3,5,2	10,11,12	0.45	0	7,12,14	0.67	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	5	3,5,2	-	0/9/10/12	-
2	ALY	J	4	2	-	2/9/10/12	-
2	ALY	J	5	3,5,2	-	0/9/10/12	-
2	ALY	K	4	2	-	2/9/10/12	-
2	ALY	L	4	2	-	0/9/10/12	-
2	ALY	L	5	3,5,2	-	0/9/10/12	-
2	ALY	I	4	2	-	1/9/10/12	-
2	ALY	K	5	3,5,2	-	0/9/10/12	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	K	4	ALY	O-C-CA-CB
2	J	4	ALY	CH3-CH-NZ-CE
2	J	4	ALY	OH-CH-NZ-CE
2	I	4	ALY	CE-CD-CG-CB
2	K	4	ALY	C-CA-CB-CG

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	K	4	ALY	2	0
2	L	5	ALY	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 18 ligands modelled in this entry, 14 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	MCM	I	6	2	14,14,14	1.98	6 (42%)	20,20,20	2.35	10 (50%)
5	MCM	J	6	2	14,14,14	2.00	5 (35%)	20,20,20	2.25	10 (50%)
5	MCM	K	6	2	14,14,14	1.86	4 (28%)	20,20,20	2.33	9 (45%)
5	MCM	L	6	2	14,14,14	2.02	5 (35%)	20,20,20	2.31	9 (45%)

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	MCM	I	6	2	-	-	0/2/2/2
5	MCM	J	6	2	-	-	0/2/2/2
5	MCM	K	6	2	-	-	0/2/2/2
5	MCM	L	6	2	-	-	0/2/2/2

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	K	6	MCM	C7-C6	3.74	1.50	1.45
5	L	6	MCM	C7-C6	3.38	1.50	1.45
5	L	6	MCM	C5-C4	3.24	1.51	1.44
5	J	6	MCM	C5-C4	3.22	1.51	1.44
5	I	6	MCM	C7-C6	3.20	1.50	1.45
5	J	6	MCM	C7-C6	3.00	1.49	1.45
5	K	6	MCM	C5-C4	2.93	1.50	1.44
5	I	6	MCM	C5-C4	2.80	1.50	1.44
5	J	6	MCM	C7-C3	2.73	1.45	1.40
5	I	6	MCM	C8-C7	2.68	1.44	1.39
5	L	6	MCM	C8-C7	2.68	1.44	1.39
5	L	6	MCM	C7-C3	2.67	1.45	1.40
5	J	6	MCM	C8-C7	2.64	1.43	1.39
5	I	6	MCM	C7-C3	2.63	1.45	1.40
5	J	6	MCM	C5-C6	2.55	1.38	1.35
5	L	6	MCM	C5-C6	2.55	1.38	1.35
5	K	6	MCM	C7-C3	2.38	1.45	1.40
5	I	6	MCM	C5-C6	2.34	1.38	1.35
5	K	6	MCM	C8-C7	2.16	1.43	1.39
5	I	6	MCM	C2-C3	2.04	1.42	1.38

All (38) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	K	6	MCM	O1-C4-C5	4.35	122.45	117.16
5	I	6	MCM	O1-C4-C5	4.11	122.16	117.16
5	L	6	MCM	O1-C4-C5	4.04	122.07	117.16
5	I	6	MCM	O2-C4-C5	-3.97	117.94	125.92
5	J	6	MCM	O1-C4-C5	3.88	121.88	117.16
5	K	6	MCM	O2-C4-C5	-3.84	118.20	125.92
5	L	6	MCM	O2-C4-C5	-3.84	118.20	125.92
5	J	6	MCM	O2-C4-C5	-3.80	118.29	125.92
5	K	6	MCM	C4-C5-C6	-3.80	119.84	123.06
5	I	6	MCM	C2-CA-N	3.65	126.72	120.56
5	L	6	MCM	C4-C5-C6	-3.59	120.02	123.06
5	I	6	MCM	C4-C5-C6	-3.54	120.06	123.06
5	K	6	MCM	C2-CA-N	3.54	126.53	120.56
5	L	6	MCM	C2-CA-N	3.40	126.29	120.56
5	J	6	MCM	C4-C5-C6	-3.31	120.26	123.06
5	J	6	MCM	C2-CA-N	3.31	126.14	120.56
5	L	6	MCM	CA-C2-C3	3.18	123.54	119.81
5	I	6	MCM	C9-CA-N	-3.05	115.29	120.90
5	K	6	MCM	CA-C2-C3	2.98	123.31	119.81

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	J	6	MCM	CA-C2-C3	2.97	123.29	119.81
5	I	6	MCM	CA-C2-C3	2.83	123.13	119.81
5	I	6	MCM	C3-O1-C4	-2.75	118.25	121.67
5	L	6	MCM	C3-O1-C4	-2.73	118.28	121.67
5	I	6	MCM	O1-C4-O2	2.69	119.89	116.40
5	K	6	MCM	C3-O1-C4	-2.67	118.35	121.67
5	J	6	MCM	C3-O1-C4	-2.65	118.36	121.67
5	J	6	MCM	O1-C4-O2	2.64	119.83	116.40
5	K	6	MCM	C9-CA-N	-2.63	116.05	120.90
5	K	6	MCM	C10-C6-C7	2.63	122.91	119.94
5	J	6	MCM	C9-CA-N	-2.61	116.10	120.90
5	L	6	MCM	C9-CA-N	-2.60	116.10	120.90
5	L	6	MCM	O1-C4-O2	2.55	119.72	116.40
5	K	6	MCM	O1-C4-O2	2.26	119.34	116.40
5	L	6	MCM	C10-C6-C7	2.24	122.47	119.94
5	I	6	MCM	O1-C3-C2	2.19	118.94	115.83
5	I	6	MCM	C10-C6-C7	2.19	122.41	119.94
5	J	6	MCM	C10-C6-C7	2.16	122.39	119.94
5	J	6	MCM	O1-C3-C7	2.00	123.39	121.57

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	I	6	MCM	1	0
5	J	6	MCM	2	0
5	K	6	MCM	2	0
5	L	6	MCM	2	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	364/388 (93%)	0.01	6 (1%) 70 67	14, 27, 45, 64	5 (1%)
1	B	364/388 (93%)	0.05	8 (2%) 62 58	14, 27, 45, 64	3 (0%)
1	C	364/388 (93%)	0.12	10 (2%) 56 51	14, 29, 46, 64	5 (1%)
1	D	364/388 (93%)	0.00	5 (1%) 73 70	16, 29, 46, 63	5 (1%)
2	I	2/5 (40%)	0.31	0 100 100	29, 29, 29, 44	0
2	J	2/5 (40%)	0.77	0 100 100	37, 37, 37, 50	0
2	K	2/5 (40%)	0.06	0 100 100	31, 31, 31, 39	0
2	L	2/5 (40%)	1.03	0 100 100	39, 39, 39, 56	0
All	All	1464/1572 (93%)	0.05	29 (1%) 65 61	14, 28, 46, 64	18 (1%)

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	377	VAL	4.5
1	C	377	VAL	4.2
1	D	377	VAL	3.7
1	A	14	LEU	3.3
1	D	14	LEU	3.3
1	C	334	HIS	3.0
1	C	170	GLU	2.9
1	C	22	PRO	2.7
1	B	91	PRO	2.7
1	B	377	VAL	2.7
1	B	88	ASP	2.6
1	B	170	GLU	2.6
1	D	94	ILE	2.5
1	B	360	HIS	2.5
1	B	87	ASP	2.4
1	C	331	ILE	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	375	HIS	2.4
1	A	360	HIS	2.3
1	A	329	SER	2.3
1	A	170	GLU	2.3
1	D	360	HIS	2.2
1	C	14	LEU	2.2
1	D	89	ASP	2.2
1	C	376	VAL	2.2
1	C	330	GLU	2.2
1	A	89	ASP	2.1
1	B	221	LYS	2.1
1	C	20	TYR	2.1
1	C	329	SER	2.1

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

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	K	4	12/13	0.87	0.14	20,33,49,50	0
2	ALY	I	4	12/13	0.88	0.15	20,35,54,55	0
2	ALY	L	4	12/13	0.92	0.10	23,31,42,43	0
2	ALY	K	5	12/13	0.93	0.09	9,15,19,20	0
2	ALY	J	4	12/13	0.93	0.08	21,28,33,33	0
2	ALY	L	5	12/13	0.94	0.10	20,23,29,30	0
2	ALY	I	5	12/13	0.95	0.09	13,19,24,25	0
2	ALY	J	5	12/13	0.96	0.07	16,19,21,21	0

6.3 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

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	MCM	K	6	13/13	0.92	0.09	18,23,26,30	0
5	MCM	I	6	13/13	0.95	0.07	17,21,21,24	0
5	MCM	L	6	13/13	0.95	0.07	23,25,28,28	0
5	MCM	J	6	13/13	0.96	0.07	16,18,19,21	0
4	K	D	411	1/1	0.97	0.03	28,28,28,28	0
4	K	D	410	1/1	0.98	0.03	23,23,23,23	0
3	ZN	A	413	1/1	0.98	0.03	49,49,49,49	0
4	K	A	402	1/1	0.98	0.02	20,20,20,20	0
4	K	B	404	1/1	0.98	0.03	21,21,21,21	0
4	K	B	405	1/1	0.98	0.05	24,24,24,24	0
4	K	C	408	1/1	0.98	0.04	23,23,23,23	0
3	ZN	C	412	1/1	0.99	0.03	46,46,46,46	0
4	K	A	401	1/1	0.99	0.02	16,16,16,16	0
3	ZN	A	400	1/1	1.00	0.01	14,14,14,14	0
3	ZN	D	409	1/1	1.00	0.02	17,17,17,17	0
4	K	C	407	1/1	1.00	0.01	20,20,20,20	0
3	ZN	B	403	1/1	1.00	0.01	14,14,14,14	0
3	ZN	C	406	1/1	1.00	0.01	16,16,16,16	0

6.5 Other polymers ⓘ

There are no such residues in this entry.