



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 14, 2026 – 04:21 PM UTC

PDB ID : 8Q3E / pdb_00008q3e
Title : High Resolution Structure of Nucleosome Core with Bound Foamy Virus GAG Peptide
Authors : De Falco, L.; Batchelor, L.K.; Dyson, P.J.; Davey, C.A.
Deposited on : 2023-08-04
Resolution : 2.17 Å(reported)

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

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

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

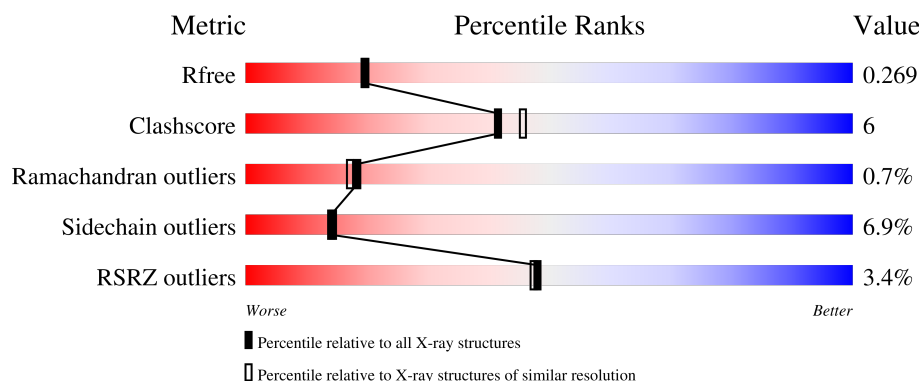
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.17 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.



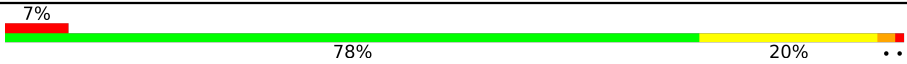
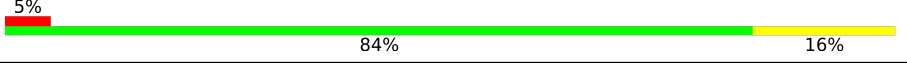
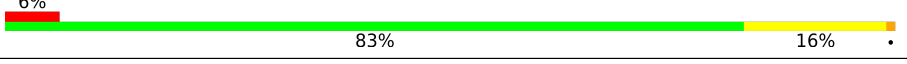
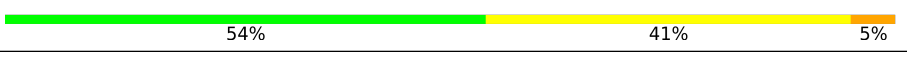
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	8975 (2.20-2.16)
Clashscore	190562	9786 (2.20-2.16)
Ramachandran outliers	187476	9664 (2.20-2.16)
Sidechain outliers	187428	9664 (2.20-2.16)
RSRZ outliers	180081	8979 (2.20-2.16)

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	MMM	18	28% 72% 22% 6%
2	AAA	98	5% 86% 13% .
2	EEE	98	3% 83% 16% .
3	BBB	87	3% 83% 9% 6% .
3	FFF	87	86% 11% .

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Mol	Chain	Length	Quality of chain
4	CCC	107	
4	GGG	107	
5	DDD	95	
5	HHH	95	
6	III	145	
7	JJJ	145	

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 12197 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 GLY-GLY-TYR-ASN-LEU-ARG-PRO-ARG-THR-TYR-GLN-PRO-GLN-ARG-TYR-GLY-GLY-GLY.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
1	MMM	18	Total	C	N	O	0	0	1
			141	87	30	24			

- Molecule 2 is a protein called Histone H3.1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	AAA	98	Total	C	N	O	S	0	0	0
			807	508	156	139	4			
2	EEE	98	Total	C	N	O	S	0	0	0
			807	508	156	139	4			

- Molecule 3 is a protein called Histone H4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	BBB	82	Total	C	N	O	S	0	0	0
			653	412	127	113	1			
3	FFF	87	Total	C	N	O	S	0	0	0
			703	442	142	118	1			

- Molecule 4 is a protein called Histone H2A type 1-B/E.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	CCC	107	Total	C	N	O	0	0	0
			828	523	162	143			
4	GGG	107	Total	C	N	O	0	0	0
			828	523	162	143			

- Molecule 5 is a protein called Histone H2B type 1-K.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	DDD	95	Total	C	N	O	S	0	0	0
			745	467	136	140	2			
5	HHH	95	Total	C	N	O	S	0	0	0
			745	467	136	140	2			

- Molecule 6 is a DNA chain called DNA (145-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	III	145	Total	C	N	O	P	0	0	0
			2970	1421	538	867	144			

- Molecule 7 is a DNA chain called DNA (145-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
7	JJJ	145	Total	C	N	O	P	0	0	0
			2969	1421	535	869	144			

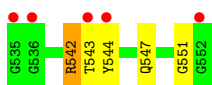
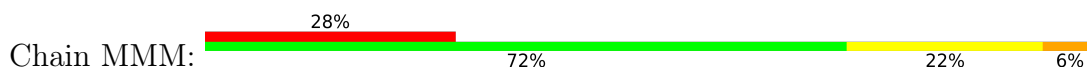
- Molecule 8 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	EEE	1	Total	Mg	0	0
			1	1		

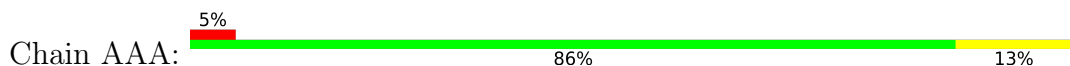
3 Residue-property plots [i](#)

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.

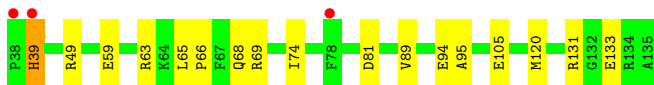
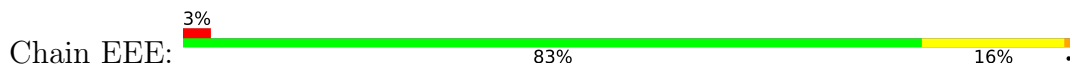
- Molecule 1: GLY-GLY-TYR-ASN-LEU-ARG-PRO-ARG-THR-TYR-GLN-PRO-GLN-ARG-TYR-GLY-GLY-GLY



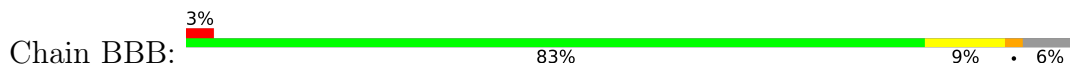
- Molecule 2: Histone H3.1



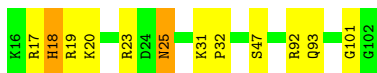
- Molecule 2: Histone H3.1



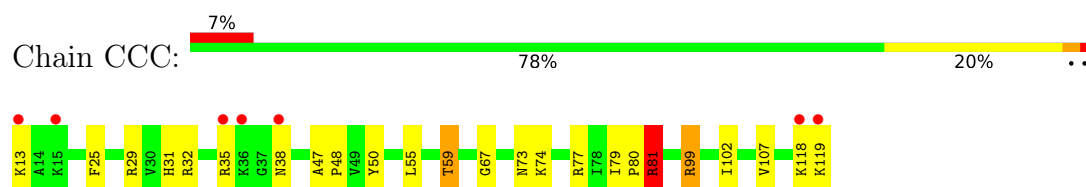
- Molecule 3: Histone H4



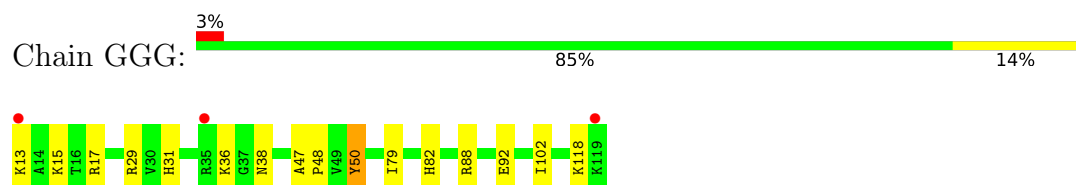
- Molecule 3: Histone H4



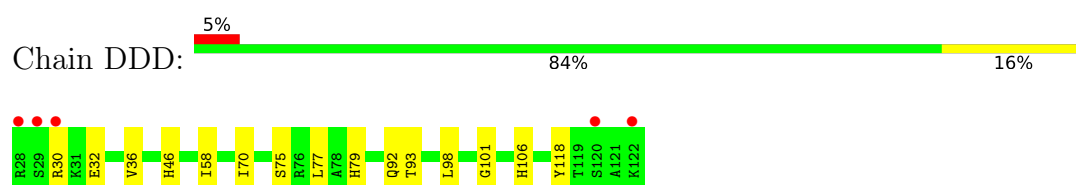
- Molecule 4: Histone H2A type 1-B/E



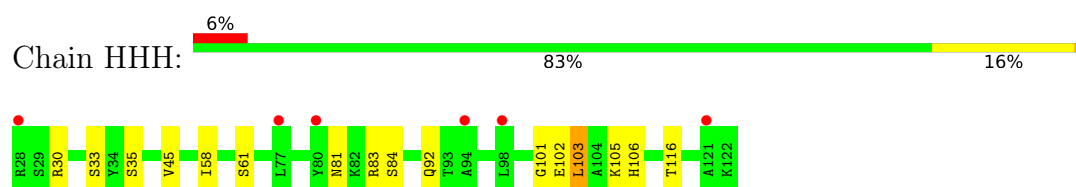
- Molecule 4: Histone H2A type 1-B/E



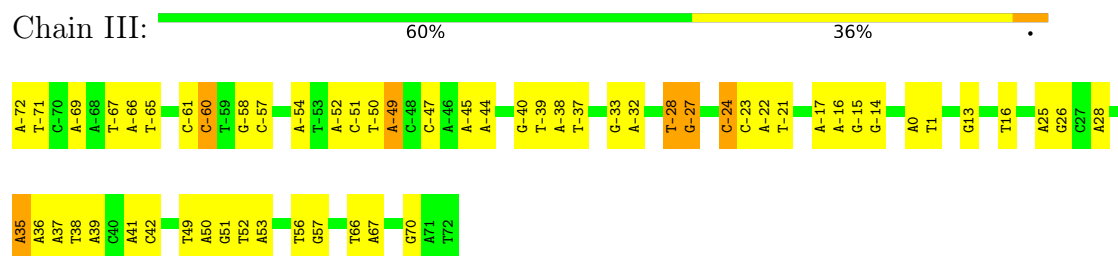
- Molecule 5: Histone H2B type 1-K



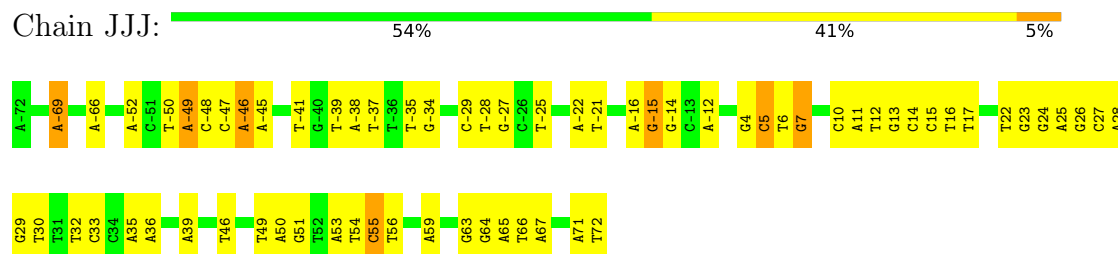
- Molecule 5: Histone H2B type 1-K



- Molecule 6: DNA (145-MER)



- Molecule 7: DNA (145-MER)



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	107.17Å 109.64Å 183.56Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.85 – 2.17 48.85 – 2.17	Depositor EDS
% Data completeness (in resolution range)	98.1 (48.85-2.17) 98.1 (48.85-2.17)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.06 (at 2.18Å)	Xtriage
Refinement program	REFMAC 5.8.0258	Depositor
R, R_{free}	0.234 , 0.270 0.239 , 0.269	Depositor DCC
R_{free} test set	2303 reflections (2.02%)	wwPDB-VP
Wilson B-factor (Å ²)	55.8	Xtriage
Anisotropy	0.361	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 40.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.019 for k,h,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	12197	wwPDB-VP
Average B, all atoms (Å ²)	88.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.85% 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: MG

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	MMM	1.06	0/145	1.20	0/195
2	AAA	1.06	0/819	1.46	1/1097 (0.1%)
2	EEE	1.19	0/819	1.52	4/1097 (0.4%)
3	BBB	1.10	0/660	1.53	2/883 (0.2%)
3	FFF	1.13	0/711	1.53	3/948 (0.3%)
4	CCC	1.12	0/838	1.53	2/1129 (0.2%)
4	GGG	1.09	1/838 (0.1%)	1.47	0/1129
5	DDD	1.18	1/756 (0.1%)	1.57	4/1014 (0.4%)
5	HHH	1.15	0/756	1.54	0/1014
6	III	0.44	0/3332	0.98	16/5141 (0.3%)
7	JJJ	0.45	0/3330	0.93	13/5138 (0.3%)
All	All	0.85	2/13004 (0.0%)	1.24	45/18785 (0.2%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	DDD	70	ILE	C-O	6.25	1.31	1.24
4	GGG	50	TYR	C-O	5.23	1.30	1.24

All (45) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	JJJ	-45	DA	C2'-C3'-O3'	-9.73	96.90	111.50
6	III	-67	DT	C2'-C3'-O3'	-8.81	98.28	111.50
6	III	16	DT	C4'-C3'-O3'	-8.81	96.78	110.00
6	III	-49	DA	N9-C1'-C2'	8.24	125.87	113.50
6	III	-40	DG	N9-C1'-C2'	8.15	125.72	113.50
7	JJJ	-45	DA	C4'-C3'-O3'	7.65	121.47	110.00
6	III	-49	DA	C4-N9-C1'	-7.56	115.72	127.05
6	III	-49	DA	C8-N9-C1'	7.51	138.31	127.05

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	III	-28	DT	C2'-C3'-O3'	-6.81	101.28	111.50
6	III	66	DT	C2'-C3'-O3'	6.61	121.42	111.50
3	FFF	92	ARG	CB-CA-C	-6.37	99.86	110.68
6	III	-27	DG	C2'-C3'-O3'	-6.34	101.99	111.50
7	JJJ	5	DC	C4'-C3'-O3'	6.31	119.46	110.00
6	III	-69	DA	C2'-C3'-O3'	-6.28	102.08	111.50
5	DDD	75	SER	CA-C-N	6.21	128.85	120.65
5	DDD	75	SER	C-N-CA	6.21	128.85	120.65
3	BBB	71	THR	CA-CB-OG1	-5.94	100.69	109.60
7	JJJ	7	DG	P-O5'-C5'	-5.93	111.11	120.00
6	III	-60	DC	C2'-C3'-O3'	5.88	120.32	111.50
4	CCC	73	ASN	N-CA-C	-5.84	106.07	112.72
7	JJJ	55	DC	C2'-C3'-O3'	5.79	120.19	111.50
7	JJJ	4	DG	C2'-C3'-O3'	-5.70	102.95	111.50
6	III	-49	DA	C1'-O4'-C4'	-5.68	101.18	109.70
3	BBB	101	GLY	CA-C-O	-5.60	118.10	122.52
7	JJJ	30	DT	C2'-C3'-O3'	-5.59	103.11	111.50
6	III	-65	DT	C2'-C3'-O3'	5.57	119.85	111.50
6	III	28	DA	C4'-C3'-O3'	-5.56	101.66	110.00
7	JJJ	-15	DG	C2'-C3'-O3'	5.44	119.66	111.50
7	JJJ	-25	DT	C2'-C3'-O3'	-5.41	103.39	111.50
3	FFF	25	ASN	CA-CB-CG	-5.40	107.20	112.60
7	JJJ	-46	DA	C2'-C3'-O3'	5.39	119.58	111.50
7	JJJ	-69	DA	C2'-C3'-O3'	-5.37	103.44	111.50
2	EEE	105	GLU	N-CA-C	-5.36	105.34	111.07
6	III	-24	DC	C2'-C3'-O3'	-5.36	103.47	111.50
2	EEE	94	GLU	N-CA-C	-5.34	105.53	111.36
2	EEE	95	ALA	CA-C-O	-5.32	114.78	120.42
7	JJJ	-49	DA	C4'-C3'-O3'	5.31	117.96	110.00
2	AAA	129	ARG	CB-CA-C	5.26	119.21	110.90
7	JJJ	7	DG	C5'-C4'-C3'	-5.17	107.14	114.90
2	EEE	74	ILE	N-CA-C	-5.16	105.68	110.53
3	FFF	101	GLY	CA-C-O	-5.12	118.47	122.52
5	DDD	118	TYR	CA-C-N	5.06	127.32	120.44
5	DDD	118	TYR	C-N-CA	5.06	127.32	120.44
6	III	35	DA	C4'-C3'-O3'	5.06	117.59	110.00
4	CCC	81	ARG	CB-CG-CD	5.05	122.92	111.30

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	MMM	141	0	131	3	0
2	AAA	807	0	844	11	0
2	EEE	807	0	844	7	0
3	BBB	653	0	696	10	0
3	FFF	703	0	755	3	0
4	CCC	828	0	892	17	0
4	GGG	828	0	892	10	0
5	DDD	745	0	769	8	0
5	HHH	745	0	769	7	0
6	III	2970	0	1640	46	0
7	JJJ	2969	0	1641	60	0
8	EEE	1	0	0	0	0
All	All	12197	0	9873	142	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 6.

All (142) 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)
7:JJJ:-22:DA:H2''	7:JJJ:-21:DT:O5'	1.78	0.81
7:JJJ:-49:DA:H4'	7:JJJ:-48:DC:OP1	1.80	0.81
2:AAA:96:CYS:SG	3:BBB:58:LEU:CD1	2.70	0.80
6:III:-50:DT:C4	6:III:-49:DA:N6	2.57	0.73
4:GGG:102:ILE:HG23	5:HHH:58:ILE:HD13	1.72	0.72
7:JJJ:66:DT:H2''	7:JJJ:67:DA:OP2	1.93	0.69
2:AAA:96:CYS:SG	3:BBB:58:LEU:HD13	2.37	0.64
6:III:53:DA:C2	7:JJJ:-52:DA:C2	2.86	0.64
2:EEE:49:ARG:HH11	2:EEE:49:ARG:HG3	1.63	0.64
4:CCC:47:ALA:HB3	4:CCC:48:PRO:HD3	1.80	0.62
6:III:-58:DG:N2	7:JJJ:59:DA:C2	2.67	0.62
2:AAA:96:CYS:SG	3:BBB:58:LEU:HD11	2.40	0.61
6:III:-50:DT:C4	6:III:-49:DA:C6	2.88	0.60
2:AAA:73:GLU:OE1	3:BBB:25:ASN:HB2	2.01	0.60
7:JJJ:-15:DG:H2''	7:JJJ:-14:DG:OP2	2.02	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:JJJ:-28:DT:H3'	7:JJJ:-27:DG:H5''	1.84	0.59
6:III:49:DT:H2''	6:III:50:DA:C8	2.37	0.59
6:III:-52:DA:C2	7:JJJ:53:DA:N1	2.71	0.59
6:III:-39:DT:H2''	6:III:-38:DA:H8	1.69	0.58
2:EEE:69:ARG:HD2	3:FFF:25:ASN:OD1	2.04	0.58
7:JJJ:50:DA:H2''	7:JJJ:51:DG:OP2	2.03	0.58
6:III:-22:DA:C2	7:JJJ:23:DG:N2	2.72	0.57
4:GGG:29:ARG:NH1	5:HHH:33:SER:O	2.37	0.57
6:III:-52:DA:C2	7:JJJ:53:DA:C2	2.93	0.57
7:JJJ:24:DG:H4'	7:JJJ:25:DA:OP1	2.04	0.57
7:JJJ:13:DG:H4'	7:JJJ:14:DC:OP1	2.05	0.57
4:CCC:99:ARG:HH11	4:CCC:99:ARG:HG3	1.70	0.56
6:III:-50:DT:O4	6:III:-49:DA:N6	2.37	0.56
4:GGG:13:LYS:HD3	7:JJJ:-41:DT:OP1	2.06	0.55
4:CCC:25:PHE:CZ	4:CCC:59:THR:HG21	2.43	0.54
5:DDD:77:LEU:HD21	5:DDD:93:THR:HB	1.89	0.54
7:JJJ:54:DT:H2''	7:JJJ:55:DC:OP2	2.07	0.54
6:III:13:DG:C2	7:JJJ:-12:DA:C2	2.95	0.54
7:JJJ:-28:DT:H2'	7:JJJ:-27:DG:C8	2.42	0.54
2:AAA:76:GLN:HE21	2:AAA:80:THR:HG22	1.73	0.54
7:JJJ:5:DC:H4'	7:JJJ:6:DT:OP1	2.07	0.54
7:JJJ:-28:DT:C3'	7:JJJ:-27:DG:H5''	2.39	0.53
4:CCC:32:ARG:HH22	5:DDD:32:GLU:CD	2.18	0.52
5:DDD:79:HIS:NE2	4:GGG:38:ASN:OD1	2.43	0.52
6:III:-28:DT:H4'	6:III:-27:DG:OP1	2.10	0.51
6:III:-24:DC:H1'	6:III:-23:DC:C6	2.45	0.51
6:III:-24:DC:H1'	6:III:-23:DC:C5	2.46	0.50
5:HHH:33:SER:HG	5:HHH:35:SER:HG	1.49	0.50
7:JJJ:35:DA:H2''	7:JJJ:36:DA:OP2	2.12	0.50
4:GGG:17:ARG:HH12	4:GGG:31:HIS:HD2	1.60	0.50
6:III:-61:DC:H2''	6:III:-60:DC:O5'	2.12	0.49
6:III:67:DA:C2	7:JJJ:-66:DA:C2	3.00	0.49
7:JJJ:-49:DA:C4'	7:JJJ:-48:DC:OP1	2.55	0.49
6:III:-22:DA:H2''	6:III:-21:DT:OP2	2.12	0.49
7:JJJ:65:DA:H2''	7:JJJ:66:DT:OP2	2.12	0.49
4:CCC:81:ARG:NH2	4:CCC:107:VAL:O	2.36	0.49
6:III:-39:DT:H2''	6:III:-38:DA:C8	2.48	0.49
7:JJJ:55:DC:C6	7:JJJ:56:DT:H72	2.48	0.48
7:JJJ:14:DC:H4'	7:JJJ:15:DC:OP1	2.13	0.48
4:CCC:47:ALA:HB3	4:CCC:48:PRO:CD	2.42	0.48
4:CCC:55:LEU:O	4:CCC:59:THR:CG2	2.61	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:GGG:17:ARG:HH12	4:GGG:31:HIS:CD2	2.32	0.48
7:JJJ:50:DA:C6	7:JJJ:51:DG:C6	3.02	0.48
2:AAA:40:ARG:HG3	7:JJJ:10:DC:H5''	1.96	0.48
7:JJJ:11:DA:H2''	7:JJJ:12:DT:O5'	2.14	0.47
4:CCC:79:ILE:HB	4:CCC:80:PRO:CD	2.44	0.47
7:JJJ:53:DA:C2	7:JJJ:54:DT:C2	3.01	0.47
1:MMM:544:TYR:CD1	1:MMM:544:TYR:N	2.82	0.47
4:GGG:79:ILE:HG12	4:GGG:82:HIS:CE1	2.50	0.47
6:III:13:DG:N2	7:JJJ:-12:DA:C2	2.83	0.47
7:JJJ:15:DC:H2''	7:JJJ:16:DT:OP2	2.13	0.47
4:CCC:77:ARG:HD3	6:III:-54:DA:H5''	1.95	0.47
6:III:39:DA:C2	7:JJJ:-38:DA:C2	3.01	0.47
3:BBB:46:ILE:O	7:JJJ:7:DG:H3'	2.15	0.47
3:BBB:22:LEU:HD22	3:BBB:22:LEU:C	2.39	0.47
7:JJJ:22:DT:H2''	7:JJJ:23:DG:C8	2.50	0.46
7:JJJ:-34:DG:H2'	7:JJJ:-34:DG:OP2	2.16	0.46
6:III:-51:DC:H5''	6:III:-51:DC:H6	1.81	0.46
4:CCC:50:TYR:OH	5:DDD:92:GLN:NE2	2.35	0.46
7:JJJ:12:DT:H2''	7:JJJ:13:DG:C8	2.50	0.46
6:III:-15:DG:H4'	6:III:-14:DG:OP1	2.14	0.46
6:III:70:DG:N2	7:JJJ:-69:DA:C2	2.83	0.46
4:CCC:67:GLY:HA3	5:DDD:46:HIS:CD2	2.50	0.46
6:III:25:DA:C2	6:III:26:DG:C2	3.04	0.46
6:III:41:DA:C4	6:III:42:DC:C5	3.03	0.46
4:CCC:55:LEU:O	4:CCC:59:THR:HG23	2.15	0.46
7:JJJ:-47:DC:O2	7:JJJ:-46:DA:C2	2.69	0.46
1:MMM:542:ARG:HD2	1:MMM:544:TYR:HE1	1.80	0.45
6:III:38:DT:H1'	6:III:39:DA:C8	2.52	0.45
7:JJJ:-35:DT:H2''	7:JJJ:-34:DG:C8	2.51	0.45
2:AAA:63:ARG:CZ	7:JJJ:17:DT:H4'	2.47	0.45
4:GGG:50:TYR:OH	5:HHH:92:GLN:NE2	2.35	0.45
6:III:-45:DA:H1'	6:III:-44:DA:H5'	1.98	0.44
4:CCC:32:ARG:O	4:CCC:35:ARG:HB3	2.17	0.44
2:AAA:59:GLU:OE2	3:BBB:40:ARG:NH2	2.50	0.44
4:CCC:80:PRO:HB3	5:DDD:58:ILE:CD1	2.47	0.44
6:III:-47:DC:N4	7:JJJ:46:DT:O4	2.51	0.44
4:CCC:79:ILE:HB	4:CCC:80:PRO:HD2	2.00	0.44
3:FFF:31:LYS:N	3:FFF:32:PRO:HD2	2.32	0.44
7:JJJ:-16:DA:H4'	7:JJJ:-15:DG:OP1	2.17	0.44
4:CCC:102:ILE:HG23	5:DDD:58:ILE:HD13	1.99	0.44
2:EEE:39:HIS:O	2:EEE:39:HIS:ND1	2.51	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:EEE:81:ASP:HA	3:FFF:19:ARG:HH22	1.83	0.44
5:HHH:81:ASN:O	5:HHH:83:ARG:HG2	2.17	0.44
7:JJJ:63:DG:H2''	7:JJJ:64:DG:OP2	2.18	0.44
4:GGG:92:GLU:HB3	5:HHH:103:LEU:HD22	1.99	0.43
6:III:0:DA:H2''	6:III:1:DT:OP2	2.18	0.43
6:III:51:DG:H2''	6:III:52:DT:O5'	2.19	0.43
7:JJJ:32:DT:H2''	7:JJJ:33:DC:C5	2.53	0.43
2:AAA:63:ARG:NH1	7:JJJ:17:DT:H4'	2.33	0.43
6:III:35:DA:H4'	6:III:36:DA:OP1	2.16	0.43
7:JJJ:-38:DA:H2''	7:JJJ:-37:DT:OP2	2.19	0.43
6:III:-66:DA:C2	7:JJJ:67:DA:C2	3.07	0.43
6:III:-38:DA:C2	6:III:-37:DT:C2	3.06	0.43
7:JJJ:53:DA:H4'	7:JJJ:54:DT:OP1	2.18	0.43
1:MMM:542:ARG:HD2	1:MMM:544:TYR:CE1	2.53	0.43
6:III:-50:DT:H5'	6:III:-49:DA:OP2	2.19	0.43
7:JJJ:-50:DT:C4	7:JJJ:-49:DA:N6	2.87	0.43
6:III:56:DT:H2''	6:III:57:DG:N7	2.34	0.43
7:JJJ:-22:DA:H2''	7:JJJ:-21:DT:C5'	2.48	0.43
6:III:37:DA:C5	6:III:38:DT:C4	3.07	0.42
6:III:-24:DC:C2	6:III:-23:DC:C4	3.07	0.42
2:EEE:131:ARG:HD3	2:EEE:133:GLU:OE2	2.19	0.42
6:III:-58:DG:N3	6:III:-57:DC:O2	2.53	0.42
6:III:-17:DA:C2	6:III:-16:DA:C5	3.08	0.42
6:III:-33:DG:N2	6:III:-32:DA:C2	2.88	0.42
7:JJJ:27:DC:C4	7:JJJ:28:DA:C6	3.07	0.42
2:AAA:83:ARG:O	3:BBB:80:THR:HA	2.20	0.41
2:AAA:63:ARG:HB2	2:AAA:66:PRO:HG2	2.02	0.41
7:JJJ:-15:DG:C2'	7:JJJ:-14:DG:OP2	2.68	0.41
7:JJJ:49:DT:H2''	7:JJJ:50:DA:C8	2.55	0.41
6:III:-47:DC:O2	6:III:-47:DC:O4'	2.37	0.41
7:JJJ:71:DA:C5	7:JJJ:72:DT:C4	3.08	0.41
7:JJJ:-39:DT:H2''	7:JJJ:-38:DA:H8	1.85	0.41
6:III:-38:DA:C2	7:JJJ:39:DA:C2	3.08	0.41
7:JJJ:28:DA:H1'	7:JJJ:29:DG:C8	2.55	0.41
6:III:-72:DA:C2	6:III:-71:DT:C2	3.09	0.41
3:BBB:47:SER:HA	7:JJJ:7:DG:H5'	2.03	0.41
4:CCC:32:ARG:NH2	5:DDD:32:GLU:OE1	2.48	0.41
2:EEE:68:GLN:HG3	2:EEE:89:VAL:HG11	2.03	0.41
6:III:35:DA:C4'	6:III:36:DA:OP1	2.68	0.41
7:JJJ:-29:DC:H2'	7:JJJ:-28:DT:C6	2.56	0.41
7:JJJ:25:DA:C5	7:JJJ:26:DG:C6	3.09	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:EEE:65:LEU:HB3	2:EEE:66:PRO:HD3	2.02	0.40
4:GGG:47:ALA:N	4:GGG:48:PRO:HD2	2.37	0.40
3:BBB:98:TYR:CD1	5:HHH:61:SER:HB3	2.56	0.40
6:III:56:DT:O3'	6:III:57:DG:C8	2.75	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	MMM	16/18 (89%)	13 (81%)	2 (12%)	1 (6%)	1	0
2	AAA	96/98 (98%)	94 (98%)	2 (2%)	0	100	100
2	EEE	96/98 (98%)	96 (100%)	0	0	100	100
3	BBB	80/87 (92%)	79 (99%)	1 (1%)	0	100	100
3	FFF	85/87 (98%)	83 (98%)	1 (1%)	1 (1%)	10	7
4	CCC	105/107 (98%)	99 (94%)	5 (5%)	1 (1%)	12	10
4	GGG	105/107 (98%)	100 (95%)	5 (5%)	0	100	100
5	DDD	93/95 (98%)	86 (92%)	6 (6%)	1 (1%)	11	9
5	HHH	93/95 (98%)	90 (97%)	2 (2%)	1 (1%)	11	9
All	All	769/792 (97%)	740 (96%)	24 (3%)	5 (1%)	18	17

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	FFF	18	HIS
5	DDD	101	GLY
4	CCC	118	LYS
5	HHH	101	GLY
1	MMM	551	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	MMM	13/13 (100%)	10 (77%)	3 (23%)	1	0
2	AAA	85/85 (100%)	80 (94%)	5 (6%)	18	19
2	EEE	85/85 (100%)	81 (95%)	4 (5%)	23	28
3	BBB	67/72 (93%)	65 (97%)	2 (3%)	36	46
3	FFF	72/72 (100%)	66 (92%)	6 (8%)	10	10
4	CCC	85/85 (100%)	76 (89%)	9 (11%)	6	5
4	GGG	85/85 (100%)	81 (95%)	4 (5%)	23	28
5	DDD	81/81 (100%)	77 (95%)	4 (5%)	22	26
5	HHH	81/81 (100%)	73 (90%)	8 (10%)	7	6
All	All	654/659 (99%)	609 (93%)	45 (7%)	14	14

All (45) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	MMM	542	ARG
1	MMM	543	THR
1	MMM	547	GLN
2	AAA	48	LEU
2	AAA	59	GLU
2	AAA	86	SER
2	AAA	122	LYS
2	AAA	134	ARG
3	BBB	22	LEU
3	BBB	47	SER
4	CCC	13	LYS
4	CCC	29	ARG
4	CCC	31	HIS
4	CCC	38	ASN
4	CCC	59	THR
4	CCC	74	LYS
4	CCC	81	ARG
4	CCC	99	ARG

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Mol	Chain	Res	Type
4	CCC	119	LYS
5	DDD	30	ARG
5	DDD	36	VAL
5	DDD	98	LEU
5	DDD	106	HIS
2	EEE	39	HIS
2	EEE	59	GLU
2	EEE	63	ARG
2	EEE	120	MET
3	FFF	17	ARG
3	FFF	18	HIS
3	FFF	20	LYS
3	FFF	23	ARG
3	FFF	47	SER
3	FFF	93	GLN
4	GGG	15	LYS
4	GGG	36	LYS
4	GGG	88	ARG
4	GGG	118	LYS
5	HHH	30	ARG
5	HHH	45	VAL
5	HHH	84	SER
5	HHH	102	GLU
5	HHH	103	LEU
5	HHH	105	LYS
5	HHH	106	HIS
5	HHH	116	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 1 ligands modelled in this entry, 1 is 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	MMM	18/18 (100%)	1.36	5 (27%) 1 1	57, 83, 107, 113	0
2	AAA	98/98 (100%)	0.42	5 (5%) 33 33	42, 57, 93, 139	0
2	EEE	98/98 (100%)	0.08	3 (3%) 51 51	35, 46, 71, 116	0
3	BBB	82/87 (94%)	0.32	3 (3%) 45 44	42, 56, 80, 125	0
3	FFF	87/87 (100%)	0.11	0 100 100	36, 45, 76, 146	0
4	CCC	107/107 (100%)	0.25	7 (6%) 25 24	36, 53, 94, 150	0
4	GGG	107/107 (100%)	0.41	3 (2%) 55 54	43, 59, 104, 137	0
5	DDD	95/95 (100%)	0.41	5 (5%) 32 31	41, 57, 95, 149	0
5	HHH	95/95 (100%)	0.67	6 (6%) 26 25	43, 60, 107, 146	0
6	III	145/145 (100%)	0.16	0 100 100	62, 114, 159, 215	0
7	JJJ	145/145 (100%)	0.17	0 100 100	64, 113, 161, 206	0
All	All	1077/1082 (99%)	0.31	37 (3%) 48 47	35, 62, 146, 215	0

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	AAA	38	PRO	5.8
2	EEE	38	PRO	4.8
3	BBB	21	VAL	4.7
1	MMM	535	GLY	4.7
3	BBB	22	LEU	4.6
2	AAA	39	HIS	4.5
1	MMM	552	GLY	4.5
2	AAA	135	ALA	4.1
2	EEE	39	HIS	3.6
1	MMM	544	TYR	3.5
4	CCC	36	LYS	3.0
2	AAA	84	PHE	2.8

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Mol	Chain	Res	Type	RSRZ
5	HHH	77	LEU	2.8
4	GGG	35	ARG	2.8
4	CCC	119	LYS	2.8
2	AAA	134	ARG	2.7
5	DDD	28	ARG	2.6
5	DDD	122	LYS	2.6
5	DDD	30	ARG	2.6
1	MMM	536	GLY	2.5
4	CCC	38	ASN	2.5
4	CCC	15	LYS	2.5
4	CCC	118	LYS	2.5
3	BBB	80	THR	2.4
2	EEE	78	PHE	2.3
5	HHH	121	ALA	2.2
5	HHH	98	LEU	2.2
5	HHH	80	TYR	2.2
5	HHH	28	ARG	2.2
5	DDD	29	SER	2.2
4	CCC	35	ARG	2.2
4	CCC	13	LYS	2.2
4	GGG	119	LYS	2.2
5	DDD	120	SER	2.2
4	GGG	13	LYS	2.2
5	HHH	94	ALA	2.1
1	MMM	543	THR	2.1

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

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

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
8	MG	EEE	201	1/1	0.87	0.22	50,50,50,50	0

6.5 Other polymers [i](#)

There are no such residues in this entry.