



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 09:36 pm GMT

PDB ID : 4XUJ
Title : Nucleosome core particle containing adducts from treatment with a thiomorpholine-substituted [(eta-6-p-cymene)Ru(3-hydroxy-2-pyridone)Cl] compound
Authors : Adhireksan, Z.; Davey, C.A.
Deposited on : 2015-01-26
Resolution : 3.18 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

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : recalc28949

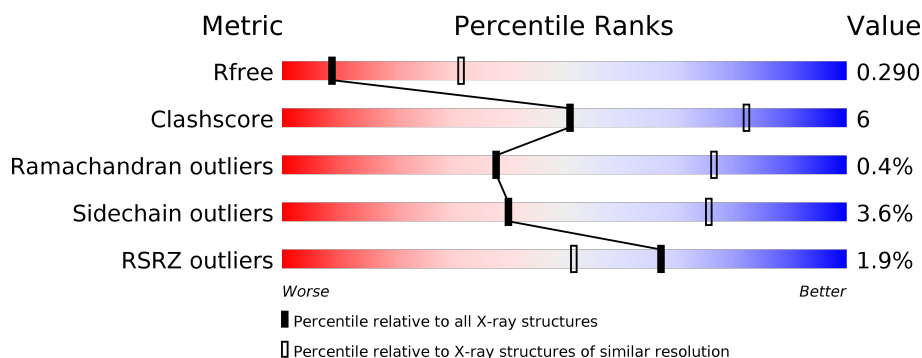
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.18 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}	100719	1123 (3.20-3.16)
Clashscore	112137	1255 (3.20-3.16)
Ramachandran outliers	110173	1233 (3.20-3.16)
Sidechain outliers	110143	1232 (3.20-3.16)
RSRZ outliers	101464	1128 (3.20-3.16)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	135	<div> <div>0% (Poor fit)</div> <div> <div>59% (Green)</div> <div>13% (Yellow)</div> <div>28% (Grey)</div> </div> </div>
1	E	135	<div> <div>59% (Green)</div> <div>13% (Yellow)</div> <div>28% (Grey)</div> </div>
2	B	102	<div> <div>2% (Poor fit)</div> <div> <div>63% (Green)</div> <div>16% (Yellow)</div> <div>20% (Grey)</div> </div> </div>
2	F	102	<div> <div>3% (Poor fit)</div> <div> <div>72% (Green)</div> <div>13% (Yellow)</div> <div>15% (Grey)</div> </div> </div>
3	C	128	<div> <div>2% (Poor fit)</div> <div> <div>66% (Green)</div> <div>16% (Yellow)</div> <div>17% (Grey)</div> </div> </div>
3	G	128	<div> <div>4% (Poor fit)</div> <div> <div>70% (Green)</div> <div>13% (Yellow)</div> <div>17% (Grey)</div> </div> </div>

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Mol	Chain	Length	Quality of chain
4	D	125	<div><div><div></div><div></div><div></div></div><div><div>2%</div><div>64%</div><div>12%</div><div>24%</div></div></div>
4	H	125	<div><div><div></div><div></div><div></div></div><div><div>2%</div><div>61%</div><div>14%</div><div>24%</div></div></div>
5	I	145	<div><div><div></div><div></div><div></div></div><div><div></div><div>53%</div><div>39%</div><div>8%</div></div></div>
6	J	145	<div><div><div></div><div></div><div></div></div><div><div></div><div>59%</div><div>32%</div><div>8%</div></div></div>

2 Entry composition [i](#)

There are 10 unique types of molecules in this entry. The entry contains 12075 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 H3.2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	97	Total	C	N	O	S	0	0	0
			802	506	155	138	3			
1	E	97	Total	C	N	O	S	0	0	0
			802	506	155	138	3			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	102	ALA	GLY	variant	UNP P84233
E	102	ALA	GLY	variant	UNP P84233

- Molecule 2 is a protein called Histone H4.

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

- Molecule 3 is a protein called Histone H2A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	106	Total	C	N	O		0	0	0
			818	516	160	142				
3	G	106	Total	C	N	O		0	0	0
			818	516	160	142				

- Molecule 4 is a protein called Histone H2B 1.1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	95	Total	C	N	O	S	0	0	0
			745	469	134	140	2			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	H	95	Total	C	N	O	S	0	0	0
			745	469	134	140	2			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	29	THR	SER	variant	UNP P02281
H	29	THR	SER	variant	UNP P02281

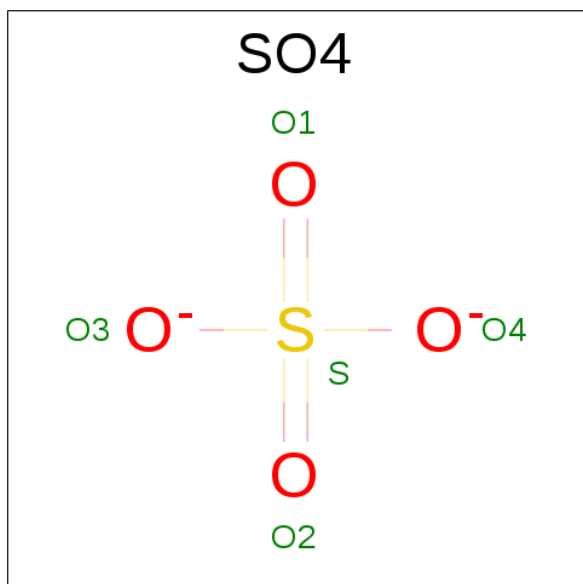
- Molecule 5 is a DNA chain called DNA (145-MER).

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

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

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

- Molecule 7 is SULFATE ION (three-letter code: SO₄) (formula: O₄S).



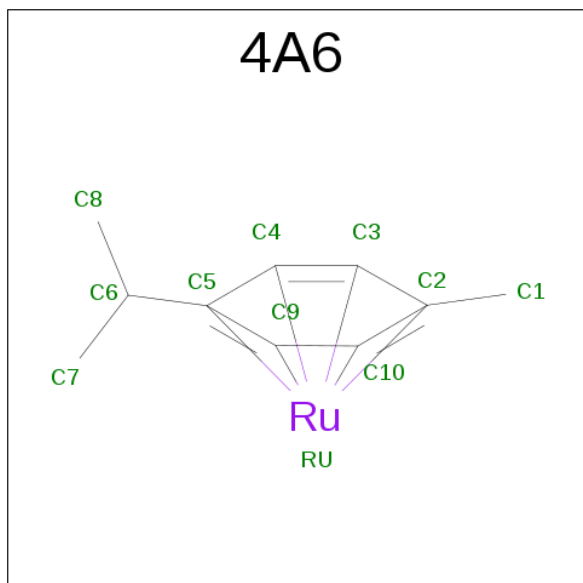
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	C	1	Total	O S	0	0
			5 4 1			

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	G	1	Total	O	S	0	0
			5	4	1		
7	H	1	Total	O	S	0	0
			5	4	1		

- Molecule 8 is [(1,2,3,4,5,6-eta)-1-methyl-4-(propan-2-yl)benzene]ruthenium (three-letter code: 4A6) (formula: C₁₀H₁₄Ru).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	D	1	Total	C	Ru	0	0
			11	10	1		
8	H	1	Total	C	Ru	0	0
			11	10	1		
8	H	1	Total	C	Ru	0	0
			11	10	1		

- Molecule 9 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	E	1	Total	Mg	0	0
			1	1		

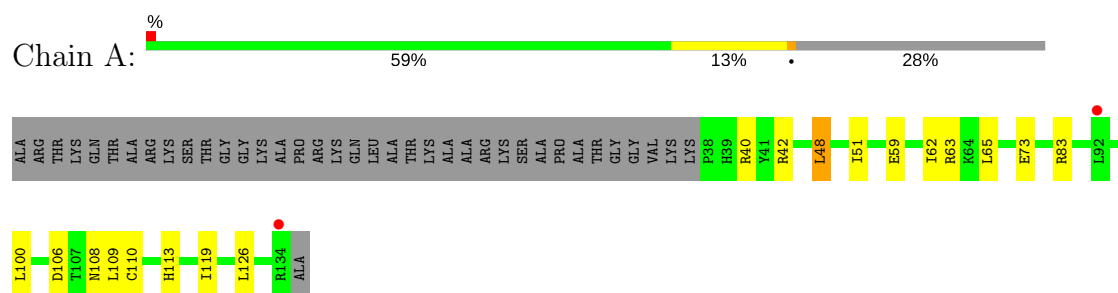
- Molecule 10 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	H	1	Total	O	0	0
			1	1		

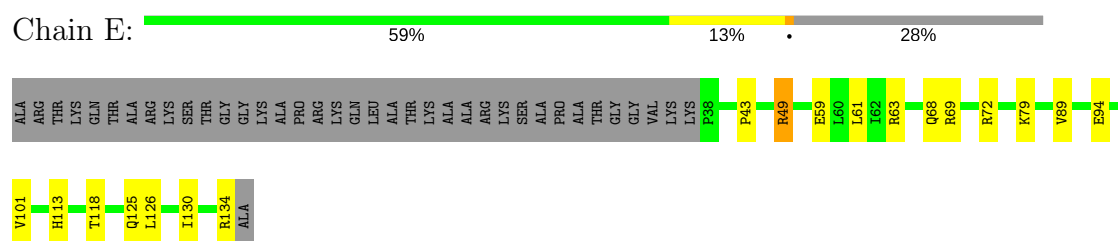
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

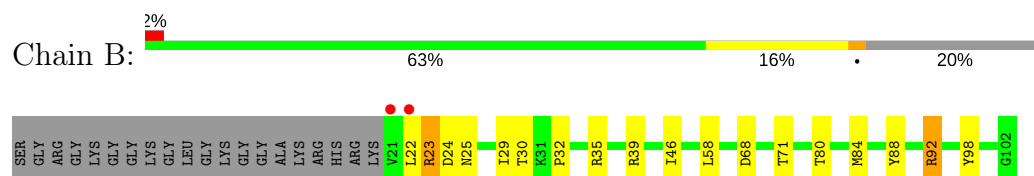
- Molecule 1: Histone H3.2



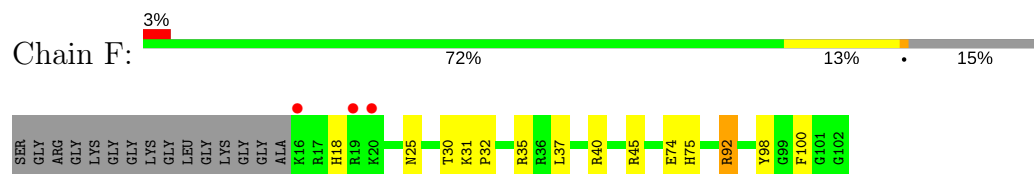
- Molecule 1: Histone H3.2



- Molecule 2: Histone H4

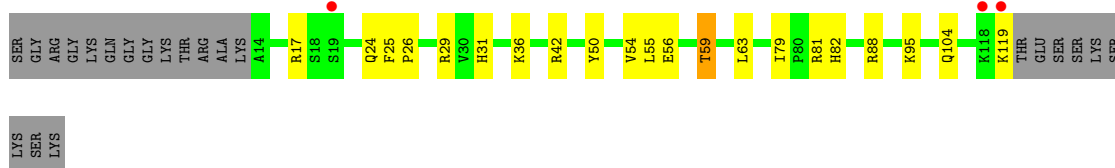


- Molecule 2: Histone H4



- Molecule 3: Histone H2A





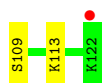
• Molecule 3: Histone H2A



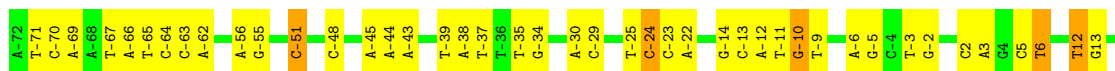
• Molecule 4: Histone H2B 1.1



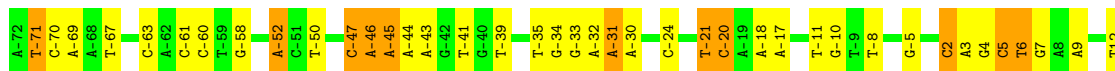
• Molecule 4: Histone H2B 1.1



• Molecule 5: DNA (145-MER)



• Molecule 6: 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	106.43Å 109.75Å 181.28Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	93.89 – 3.18 93.89 – 3.18	Depositor EDS
% Data completeness (in resolution range)	100.0 (93.89-3.18) 98.1 (93.89-3.18)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.53 (at 3.19Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.263 , 0.299 0.252 , 0.290	Depositor DCC
R_{free} test set	1783 reflections (5.27%)	DCC
Wilson B-factor (Å ²)	83.9	Xtriage
Anisotropy	0.282	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.22 , 23.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	0.045 for k,h,-l	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	12075	wwPDB-VP
Average B, all atoms (Å ²)	105.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.89% 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: 4A6, MG, SO4

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	0.34	0/814	0.51	0/1092
1	E	0.38	0/814	0.56	0/1092
2	B	0.33	0/660	0.56	0/883
2	F	0.38	0/711	0.56	0/948
3	C	0.35	0/828	0.54	0/1117
3	G	0.32	0/828	0.50	0/1117
4	D	0.36	0/756	0.52	0/1015
4	H	0.36	0/756	0.50	0/1015
5	I	0.67	0/3332	1.37	35/5141 (0.7%)
6	J	0.67	0/3330	1.39	35/5138 (0.7%)
All	All	0.54	0/12829	1.09	70/18558 (0.4%)

There are no bond length outliers.

All (70) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	J	-67	DT	O4'-C1'-N1	8.38	113.86	108.00
6	J	-45	DA	O4'-C1'-N9	8.13	113.69	108.00
5	I	-5	DG	O4'-C1'-N9	7.93	113.55	108.00
5	I	-67	DT	O4'-C1'-N1	7.77	113.44	108.00
6	J	64	DG	O4'-C1'-N9	7.71	113.40	108.00
6	J	22	DT	O4'-C1'-N1	7.70	113.39	108.00
6	J	36	DA	O4'-C1'-N9	7.63	113.34	108.00
5	I	-30	DA	O4'-C1'-N9	7.35	113.15	108.00
5	I	32	DT	O4'-C1'-N1	7.32	113.13	108.00
6	J	62	DT	O4'-C4'-C3'	-7.03	101.69	104.50
5	I	30	DT	O4'-C1'-N1	6.97	112.88	108.00
5	I	53	DA	P-O3'-C3'	6.97	128.06	119.70
5	I	64	DG	O4'-C1'-N9	6.96	112.87	108.00
5	I	37	DA	O4'-C1'-N9	6.95	112.86	108.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	I	-14	DG	O4'-C1'-N9	-6.94	103.14	108.00
5	I	45	DT	P-O3'-C3'	6.93	128.02	119.70
5	I	-71	DT	P-O3'-C3'	6.89	127.97	119.70
5	I	-14	DG	O4'-C4'-C3'	-6.89	101.74	104.50
5	I	30	DT	C1'-O4'-C4'	-6.74	103.36	110.10
6	J	6	DT	P-O3'-C3'	6.64	127.67	119.70
6	J	-31	DA	P-O3'-C3'	6.62	127.64	119.70
6	J	-58	DG	P-O3'-C3'	6.45	127.44	119.70
5	I	42	DC	P-O3'-C3'	6.44	127.43	119.70
5	I	-51	DC	P-O3'-C3'	6.43	127.41	119.70
6	J	43	DT	O4'-C1'-N1	6.43	112.50	108.00
6	J	-39	DT	P-O3'-C3'	6.40	127.38	119.70
5	I	-24	DC	O4'-C1'-N1	6.39	112.47	108.00
6	J	-47	DC	C1'-O4'-C4'	-6.38	103.72	110.10
6	J	65	DA	O4'-C1'-N9	6.38	112.46	108.00
6	J	-69	DA	O4'-C1'-N9	6.34	112.44	108.00
6	J	2	DC	O4'-C1'-N1	6.33	112.43	108.00
5	I	-29	DC	O4'-C1'-N1	6.29	112.40	108.00
5	I	-10	DG	P-O3'-C3'	6.25	127.19	119.70
5	I	32	DT	C1'-O4'-C4'	-6.23	103.87	110.10
5	I	71	DA	O4'-C1'-N9	6.22	112.35	108.00
5	I	40	DC	O4'-C1'-N1	6.19	112.33	108.00
5	I	45	DT	O4'-C1'-N1	6.19	112.33	108.00
6	J	-61	DC	O4'-C1'-N1	5.98	112.19	108.00
6	J	-71	DT	P-O3'-C3'	5.88	126.75	119.70
6	J	-20	DC	P-O3'-C3'	5.80	126.66	119.70
5	I	-48	DC	O4'-C1'-N1	5.78	112.05	108.00
6	J	-32	DA	P-O3'-C3'	5.76	126.62	119.70
6	J	5	DC	P-O3'-C3'	5.76	126.62	119.70
6	J	62	DT	O4'-C1'-N1	5.76	112.03	108.00
5	I	63	DG	P-O3'-C3'	5.75	126.59	119.70
5	I	43	DT	P-O3'-C3'	5.73	126.58	119.70
6	J	52	DT	O4'-C1'-N1	5.71	112.00	108.00
5	I	42	DC	O4'-C1'-N1	5.70	111.99	108.00
6	J	20	DG	O4'-C1'-N9	5.68	111.98	108.00
6	J	49	DT	P-O3'-C3'	5.67	126.51	119.70
5	I	-37	DT	P-O3'-C3'	5.58	126.39	119.70
6	J	33	DC	O4'-C1'-N1	5.56	111.89	108.00
6	J	18	DT	N3-C4-O4	5.52	123.21	119.90
6	J	-50	DT	P-O3'-C3'	5.48	126.28	119.70
5	I	-67	DT	C3'-C2'-C1'	-5.47	95.94	102.50
5	I	12	DT	O4'-C1'-N1	5.43	111.80	108.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	J	46	DT	O4'-C1'-N1	5.42	111.79	108.00
6	J	-46	DA	P-O3'-C3'	5.40	126.18	119.70
6	J	-63	DC	O4'-C1'-N1	5.38	111.77	108.00
5	I	6	DT	P-O3'-C3'	5.36	126.13	119.70
6	J	-8	DT	O4'-C1'-N1	5.33	111.73	108.00
6	J	22	DT	C1'-O4'-C4'	-5.31	104.79	110.10
6	J	-52	DA	P-O3'-C3'	5.25	126.00	119.70
5	I	-30	DA	C1'-O4'-C4'	-5.13	104.97	110.10
6	J	-41	DT	P-O3'-C3'	5.10	125.82	119.70
5	I	26	DG	O4'-C1'-N9	5.10	111.57	108.00
5	I	41	DA	O4'-C1'-N9	5.04	111.53	108.00
5	I	45	DT	O4'-C4'-C3'	-5.02	102.49	104.50
5	I	-45	DA	O4'-C1'-N9	5.02	111.51	108.00
6	J	-21	DT	O4'-C1'-N1	5.01	111.50	108.00

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	802	0	841	18	0
1	E	802	0	841	19	0
2	B	653	0	696	14	0
2	F	703	0	755	13	0
3	C	818	0	877	19	0
3	G	818	0	877	12	0
4	D	745	0	773	13	0
4	H	745	0	773	13	0
5	I	2970	0	1640	38	0
6	J	2969	0	1641	33	0
7	C	5	0	0	0	0
7	G	5	0	0	0	0
7	H	5	0	0	0	0
8	D	11	0	10	0	0
8	H	22	0	20	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
9	E	1	0	0	0	0
10	H	1	0	0	0	0
All	All	12075	0	9744	139	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 (139) 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:E:43:PRO:HG2	6:J:-5:DG:H5'	1.54	0.89
2:B:23:ARG:HH11	2:B:23:ARG:HG2	1.38	0.88
1:E:49:ARG:HG3	1:E:49:ARG:HH11	1.41	0.82
2:F:31:LYS:HE2	2:F:35:ARG:HH22	1.51	0.76
3:G:35:ARG:NH2	5:I:38:DT:OP2	2.17	0.74
4:D:77:LEU:HD21	4:D:93:THR:HB	1.68	0.74
1:E:69:ARG:HD2	2:F:25:ASN:OD1	1.93	0.68
4:H:28:LYS:HB3	6:J:29:DG:H5''	1.76	0.67
6:J:2:DC:H2''	6:J:3:DA:C8	2.31	0.65
5:I:-65:DT:H3	6:J:65:DA:H61	1.44	0.65
4:H:45:VAL:HG13	4:H:46:HIS:HD2	1.63	0.64
3:C:55:LEU:O	3:C:59:THR:HG22	1.99	0.61
4:D:117:LYS:HG3	4:D:117:LYS:O	2.00	0.61
1:E:79:LYS:HG3	2:F:74:GLU:OE2	2.00	0.61
3:G:35:ARG:HH22	5:I:38:DT:P	2.24	0.60
6:J:-18:DA:H2''	6:J:-17:DA:C8	2.37	0.60
2:B:23:ARG:HH11	2:B:23:ARG:CG	2.12	0.60
3:C:17:ARG:HH12	3:C:31:HIS:HD2	1.50	0.59
3:C:55:LEU:O	3:C:59:THR:CG2	2.51	0.59
2:B:98:TYR:OH	4:H:65:ASP:OD2	2.16	0.59
2:B:68:ASP:OD2	2:B:92:ARG:HD3	2.03	0.59
5:I:-11:DT:H2''	5:I:-10:DG:H5'	1.85	0.59
1:E:125:GLN:O	1:E:134:ARG:NH2	2.36	0.58
1:A:63:ARG:NH2	6:J:17:DT:H4'	2.18	0.58
5:I:56:DT:H2''	5:I:57:DG:N7	2.19	0.57
3:G:31:HIS:CD2	3:G:48:PRO:HG3	2.40	0.57
3:G:32:ARG:NH2	4:H:32:GLU:OE2	2.33	0.57
3:C:42:ARG:HB2	4:D:85:THR:HG23	1.87	0.57
5:I:-3:DT:H2''	5:I:-2:DG:C8	2.40	0.56
5:I:59:DA:H2''	5:I:60:DG:H5''	1.86	0.56
5:I:-63:DC:H2''	5:I:-62:DA:C8	2.40	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:51:ILE:HG21	3:G:111:ILE:HG12	1.88	0.56
5:I:42:DC:H2''	5:I:43:DT:OP2	2.05	0.56
5:I:5:DC:H2''	5:I:6:DT:H71	1.88	0.56
3:G:69:ALA:O	3:G:73:ASN:ND2	2.39	0.55
3:C:29:ARG:NH2	4:D:37:TYR:HE1	2.05	0.55
3:C:29:ARG:HH21	4:D:37:TYR:HE1	1.55	0.55
3:C:25:PHE:CZ	3:C:59:THR:HG21	2.42	0.55
5:I:-64:DC:H2''	5:I:-63:DC:H5''	1.88	0.55
3:C:63:LEU:HD22	4:D:42:LEU:HD13	1.89	0.54
5:I:-63:DC:C2'	5:I:-62:DA:C8	2.91	0.54
1:A:126:LEU:HD22	1:E:113:HIS:CG	2.43	0.54
6:J:5:DC:H4'	6:J:6:DT:OP1	2.06	0.54
3:C:55:LEU:HD22	4:D:63:VAL:HG13	1.90	0.54
5:I:-25:DT:H2'	5:I:-24:DC:C6	2.43	0.54
2:B:71:THR:HG22	4:D:93:THR:HG23	1.90	0.53
1:A:119:ILE:O	1:A:119:ILE:HG13	2.07	0.53
1:E:94:GLU:HG3	2:F:100:PHE:CE1	2.45	0.52
5:I:20:DG:H2''	5:I:21:DA:OP2	2.10	0.52
1:A:73:GLU:OE1	2:B:25:ASN:HB2	2.09	0.52
5:I:52:DT:H2''	5:I:53:DA:O4'	2.09	0.52
5:I:-56:DA:H2''	5:I:-55:DG:C8	2.46	0.51
5:I:12:DT:H2''	5:I:13:DG:C8	2.46	0.51
2:F:30:THR:HB	2:F:32:PRO:HD2	1.92	0.51
1:E:94:GLU:HG3	2:F:100:PHE:CZ	2.46	0.51
6:J:-71:DT:H6	6:J:-71:DT:OP2	1.94	0.51
1:E:118:THR:HA	2:F:45:ARG:HB3	1.93	0.51
3:C:17:ARG:HH12	3:C:31:HIS:CD2	2.28	0.51
3:C:104:GLN:HG2	1:E:94:GLU:OE2	2.11	0.51
5:I:-44:DA:H2'	5:I:-43:DA:C8	2.46	0.51
2:B:30:THR:HB	2:B:32:PRO:HD2	1.94	0.50
3:C:17:ARG:HH22	3:C:31:HIS:CD2	2.30	0.50
3:C:54:VAL:HG21	4:D:95:VAL:HG21	1.94	0.50
3:C:24:GLN:N	3:C:56:GLU:OE1	2.45	0.49
6:J:-71:DT:H4'	6:J:-70:DC:OP1	2.12	0.49
6:J:-31:DA:H2''	6:J:-30:DA:O5'	2.12	0.49
6:J:-45:DA:H2''	6:J:-44:DA:C8	2.47	0.49
6:J:-44:DA:H2'	6:J:-43:DA:C8	2.48	0.49
1:A:62:ILE:HD13	2:B:29:ILE:HD12	1.95	0.48
4:H:109:SER:O	4:H:113:LYS:HB2	2.13	0.48
5:I:-35:DT:H2''	5:I:-34:DG:N7	2.27	0.48
3:C:79:ILE:HG12	3:C:82:HIS:CE1	2.48	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:I:62:DT:H2''	5:I:63:DG:O5'	2.14	0.48
1:A:110:CYS:SG	1:E:130:ILE:HD11	2.55	0.47
1:E:49:ARG:HH11	1:E:49:ARG:CG	2.20	0.47
4:D:73:GLU:O	4:D:77:LEU:HB2	2.15	0.47
1:E:49:ARG:HG3	1:E:49:ARG:NH1	2.19	0.47
2:F:75:HIS:HD2	4:H:81:ASN:ND2	2.12	0.47
5:I:-23:DC:H2''	5:I:-22:DA:C8	2.50	0.47
1:A:106:ASP:HB3	1:E:130:ILE:HG12	1.97	0.47
1:A:63:ARG:CZ	6:J:17:DT:H4'	2.44	0.47
6:J:-35:DT:H2''	6:J:-34:DG:N7	2.30	0.46
1:A:65:LEU:HB3	6:J:17:DT:H3'	1.98	0.46
2:B:46:ILE:O	6:J:7:DG:H3'	2.15	0.46
1:A:40:ARG:NH2	6:J:9:DA:N3	2.64	0.46
3:G:85:LEU:O	3:G:89:ASN:HB2	2.16	0.46
5:I:-13:DC:H2''	5:I:-12:DA:N7	2.31	0.46
2:B:84:MET:HE3	2:B:88:TYR:CZ	2.50	0.46
4:H:84:SER:H	6:J:-33:DG:P	2.39	0.45
5:I:24:DG:N2	6:J:-24:DC:O2	2.46	0.45
1:A:106:ASP:HA	1:A:109:LEU:HD12	1.98	0.45
6:J:12:DT:H2''	6:J:13:DG:C5	2.51	0.45
1:A:48:LEU:HD21	3:G:117:PRO:HD3	1.97	0.45
6:J:3:DA:H2''	6:J:4:DG:C8	2.52	0.45
5:I:-55:DG:H1	6:J:55:DC:H42	1.65	0.44
1:E:68:GLN:HG3	1:E:89:VAL:HG11	1.99	0.44
6:J:-11:DT:H2''	6:J:-10:DG:C8	2.52	0.44
3:G:84:GLN:HG2	3:G:105:GLY:O	2.18	0.44
2:F:92:ARG:HH21	4:H:98:LEU:HD23	1.82	0.44
5:I:-39:DT:H2''	5:I:-38:DA:H8	1.82	0.44
6:J:-21:DT:H2''	6:J:-20:DC:C6	2.53	0.44
6:J:-44:DA:C2'	6:J:-43:DA:C8	3.00	0.44
6:J:-47:DC:H2''	6:J:-46:DA:N7	2.33	0.44
1:A:83:ARG:O	2:B:80:THR:HA	2.18	0.43
4:D:58:ILE:HG23	2:F:98:TYR:HB3	1.99	0.43
1:A:108:ASN:HD21	3:G:115:LEU:HD11	1.83	0.43
4:H:43:LYS:HA	4:H:43:LYS:HD3	1.73	0.43
5:I:19:DT:H2''	5:I:20:DG:C8	2.53	0.43
3:G:30:VAL:HG13	4:H:67:PHE:HE1	1.84	0.43
1:E:101:VAL:HG11	2:F:40:ARG:HG2	2.01	0.43
4:H:106:HIS:HE1	8:H:202:4A6:C10	2.29	0.42
1:A:113:HIS:CG	1:E:126:LEU:HD22	2.54	0.42
1:E:68:GLN:O	1:E:72:ARG:HG3	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:I:53:DA:C2	6:J:-52:DA:C2	3.08	0.42
5:I:5:DC:H2''	5:I:6:DT:C7	2.48	0.42
2:B:23:ARG:NH1	2:B:23:ARG:CG	2.79	0.42
1:A:100:LEU:HD11	2:B:58:LEU:HD13	2.02	0.42
3:C:26:PRO:HD3	4:D:37:TYR:CD2	2.55	0.42
5:I:-10:DG:H2''	5:I:-9:DT:OP2	2.20	0.42
2:B:35:ARG:O	2:B:39:ARG:HG2	2.20	0.41
6:J:26:DG:H2''	6:J:27:DC:C5	2.55	0.41
1:A:83:ARG:HD2	6:J:26:DG:H5'	2.02	0.41
3:C:88:ARG:HD3	3:C:88:ARG:HA	1.69	0.41
2:F:31:LYS:HB3	2:F:32:PRO:HD3	2.02	0.41
4:H:37:TYR:OH	5:I:47:DG:H5''	2.20	0.41
5:I:-6:DA:C2	6:J:7:DG:N2	2.89	0.41
5:I:64:DG:OP2	5:I:64:DG:H2'	2.20	0.41
3:G:27:VAL:HG11	3:G:49:VAL:HG22	2.02	0.41
5:I:2:DC:H2''	5:I:3:DA:C8	2.56	0.41
5:I:37:DA:H2''	5:I:38:DT:OP2	2.20	0.41
5:I:60:DG:N2	6:J:-60:DC:O2	2.49	0.41
3:C:17:ARG:NH1	3:C:31:HIS:HD2	2.17	0.40
3:C:95:LYS:HD3	4:D:100:PRO:HB3	2.03	0.40
4:H:99:LEU:HA	4:H:100:PRO:HD3	1.86	0.40
5:I:-51:DC:H6	5:I:-51:DC:H2'	1.79	0.40
6:J:12:DT:H2''	6:J:13:DG:N7	2.37	0.40
1:E:61:LEU:HD12	2:F:37:LEU:HD23	2.03	0.40
5:I:-70:DC:H2''	5:I:-69:DA:OP2	2.20	0.40
5:I:-66:DA:C2	6:J:67:DA:C2	3.10	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	95/135 (70%)	94 (99%)	1 (1%)	0	100	100
1	E	95/135 (70%)	93 (98%)	2 (2%)	0	100	100
2	B	80/102 (78%)	79 (99%)	1 (1%)	0	100	100
2	F	85/102 (83%)	82 (96%)	2 (2%)	1 (1%)	15	56
3	C	104/128 (81%)	100 (96%)	4 (4%)	0	100	100
3	G	104/128 (81%)	102 (98%)	2 (2%)	0	100	100
4	D	93/125 (74%)	92 (99%)	0	1 (1%)	17	58
4	H	93/125 (74%)	92 (99%)	0	1 (1%)	17	58
All	All	749/980 (76%)	734 (98%)	12 (2%)	3 (0%)	38	76

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
4	D	101	GLY
4	H	101	GLY
2	F	18	HIS

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	85/110 (77%)	82 (96%)	3 (4%)	41	76
1	E	85/110 (77%)	82 (96%)	3 (4%)	41	76
2	B	67/78 (86%)	63 (94%)	4 (6%)	22	60
2	F	72/78 (92%)	71 (99%)	1 (1%)	71	90
3	C	84/101 (83%)	79 (94%)	5 (6%)	22	60
3	G	84/101 (83%)	83 (99%)	1 (1%)	75	92
4	D	81/105 (77%)	78 (96%)	3 (4%)	39	75
4	H	81/105 (77%)	78 (96%)	3 (4%)	39	75
All	All	639/788 (81%)	616 (96%)	23 (4%)	40	75

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	42	ARG
1	A	48	LEU
1	A	59	GLU
2	B	22	LEU
2	B	23	ARG
2	B	24	ASP
2	B	92	ARG
3	C	36	LYS
3	C	50	TYR
3	C	59	THR
3	C	81	ARG
3	C	119	LYS
4	D	28	LYS
4	D	29	THR
4	D	79	HIS
1	E	49	ARG
1	E	59	GLU
1	E	63	ARG
2	F	92	ARG
3	G	88	ARG
4	H	29	THR
4	H	45	VAL
4	H	85	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	108	ASN
2	B	93	GLN
3	C	31	HIS
3	G	31	HIS
4	H	46	HIS
4	H	81	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 7 ligands modelled in this entry, 1 is monoatomic - leaving 6 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	$\# Z > 2$
7	SO4	C	1101	-	4,4,4	0.15	0	6,6,6	0.16	0
8	4A6	D	201	-	0,16,16	0.00	-	2,42,42	1.31	0
7	SO4	G	201	-	4,4,4	0.15	0	6,6,6	0.12	0
7	SO4	H	201	-	4,4,4	0.15	0	6,6,6	0.09	0
8	4A6	H	202	10,4	0,16,16	0.00	-	2,42,42	1.00	0
8	4A6	H	203	-	0,16,16	0.00	-	2,42,42	0.78	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
7	SO4	C	1101	-	-	0/0/0/0	0/0/0/0
8	4A6	D	201	-	-	0/0/90/90	0/0/6/6
7	SO4	G	201	-	-	0/0/0/0	0/0/0/0
7	SO4	H	201	-	-	0/0/0/0	0/0/0/0
8	4A6	H	202	10,4	-	0/0/90/90	0/0/6/6
8	4A6	H	203	-	-	0/0/90/90	0/0/6/6

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.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	H	202	4A6	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ > 2	OWAB(Å ²)	Q < 0.9
1	A	97/135 (71%)	-0.06	2 (2%) 64 48	62, 79, 97, 110	0
1	E	97/135 (71%)	0.03	0 100 100	56, 67, 89, 101	0
2	B	82/102 (80%)	0.08	2 (2%) 59 44	67, 74, 97, 111	0
2	F	87/102 (85%)	0.10	3 (3%) 46 30	55, 66, 94, 117	0
3	C	106/128 (82%)	0.07	3 (2%) 53 38	59, 73, 95, 99	0
3	G	106/128 (82%)	0.15	5 (4%) 32 19	67, 85, 102, 107	0
4	D	95/125 (76%)	0.09	2 (2%) 64 48	68, 78, 103, 109	0
4	H	95/125 (76%)	0.07	3 (3%) 48 31	74, 87, 107, 118	0
5	I	145/145 (100%)	-0.65	0 100 100	79, 135, 164, 173	0
6	J	145/145 (100%)	-0.49	0 100 100	80, 133, 163, 179	0
All	All	1055/1270 (83%)	-0.11	20 (1%) 67 52	55, 83, 155, 179	0

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	C	119	LYS	8.0
3	G	118	LYS	6.9
2	B	21	VAL	6.1
4	D	122	LYS	4.2
4	H	122	LYS	4.1
3	C	118	LYS	3.9
3	G	117	PRO	3.8
3	G	119	LYS	3.8
2	F	19	ARG	3.6
2	F	16	LYS	3.5
4	D	121	ALA	3.1
1	A	134	ARG	2.9
1	A	92	LEU	2.8

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Mol	Chain	Res	Type	RSRZ
3	G	83	LEU	2.7
4	H	30	ARG	2.7
4	H	86	ILE	2.6
3	C	19	SER	2.6
2	F	20	LYS	2.5
2	B	22	LEU	2.2
3	G	112	GLN	2.0

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
7	SO4	G	201	5/5	0.92	0.22	1.54	120,120,120,121	0
7	SO4	C	1101	5/5	0.94	0.31	1.01	115,115,116,116	0
9	MG	E	1001	1/1	0.94	0.26	0.70	66,66,66,66	0
8	4A6	D	201	11/11	0.95	0.15	-1.22	156,157,157,158	0
8	4A6	H	203	11/11	0.94	0.16	-1.63	178,179,179,179	0
8	4A6	H	202	11/11	0.97	0.15	-	126,126,127,127	0
7	SO4	H	201	5/5	0.84	0.18	-	129,129,130,130	0

6.5 Other polymers [i](#)

There are no such residues in this entry.