



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 08:23 PM UTC

PDB ID : 5M3H / pdb_00005m3h
Title : Bat influenza A/H17N10 polymerase bound to four heptad repeats of serine 5 phosphorylated Pol II CTD
Authors : Lukarska, M.; Pflug, A.; Cusack, S.
Deposited on : 2016-10-14
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

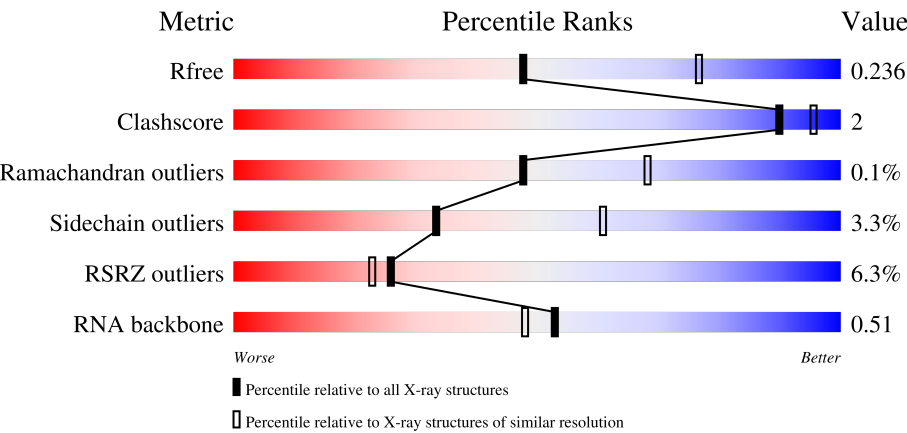
1 Overall quality at a glance i

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)
RNA backbone	3983	1003 (2.78-2.22)

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	728	11% 89% 6% • •
2	B	776	4% 90% 7% •
3	C	797	4% 86% 6% 8%
4	R	18	61% 11% 28%

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Mol	Chain	Length	Quality of chain
5	V	16	81% 12% 6%
6	X	28	4% 18% 79%
6	Y	28	4% 32% 64%

2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 18529 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 Polymerase acidic protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	699	Total	C	N	O	S	0	0	0
			5692	3617	959	1079	37			

There are 15 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	GLY	-	expression tag	UNP H6QM92
A	-2	SER	-	expression tag	UNP H6QM92
A	-1	GLY	-	expression tag	UNP H6QM92
A	0	SER	-	expression tag	UNP H6QM92
A	714	GLY	-	expression tag	UNP H6QM92
A	715	SER	-	expression tag	UNP H6QM92
A	716	GLY	-	expression tag	UNP H6QM92
A	717	SER	-	expression tag	UNP H6QM92
A	718	GLY	-	expression tag	UNP H6QM92
A	719	GLU	-	expression tag	UNP H6QM92
A	720	ASN	-	expression tag	UNP H6QM92
A	721	LEU	-	expression tag	UNP H6QM92
A	722	TYR	-	expression tag	UNP H6QM92
A	723	PHE	-	expression tag	UNP H6QM92
A	724	GLN	-	expression tag	UNP H6QM92

- Molecule 2 is a protein called RNA-directed RNA polymerase catalytic subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	750	Total	C	N	O	S	0	0	0
			5986	3766	1062	1118	40			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-8	GLY	-	expression tag	UNP H6QM91

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-7	SER	-	expression tag	UNP H6QM91
B	-6	GLY	-	expression tag	UNP H6QM91
B	-5	SER	-	expression tag	UNP H6QM91
B	-4	GLY	-	expression tag	UNP H6QM91
B	-3	SER	-	expression tag	UNP H6QM91
B	-2	GLY	-	expression tag	UNP H6QM91
B	-1	SER	-	expression tag	UNP H6QM91
B	0	GLY	-	expression tag	UNP H6QM91
B	757	GLY	-	expression tag	UNP H6QM91
B	758	SER	-	expression tag	UNP H6QM91
B	759	GLY	-	expression tag	UNP H6QM91
B	760	SER	-	expression tag	UNP H6QM91
B	761	GLY	-	expression tag	UNP H6QM91
B	762	GLU	-	expression tag	UNP H6QM91
B	763	ASN	-	expression tag	UNP H6QM91
B	764	LEU	-	expression tag	UNP H6QM91
B	765	TYR	-	expression tag	UNP H6QM91
B	766	PHE	-	expression tag	UNP H6QM91
B	767	GLN	-	expression tag	UNP H6QM91

- Molecule 3 is a protein called Polymerase basic protein 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	730	Total	C	N	O	S	0	0	0
			5796	3657	1025	1082	32			

There are 37 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	-8	GLY	-	expression tag	UNP H6QM90
C	-7	SER	-	expression tag	UNP H6QM90
C	-6	GLY	-	expression tag	UNP H6QM90
C	-5	SER	-	expression tag	UNP H6QM90
C	-4	GLY	-	expression tag	UNP H6QM90
C	-3	SER	-	expression tag	UNP H6QM90
C	-2	GLY	-	expression tag	UNP H6QM90
C	-1	SER	-	expression tag	UNP H6QM90
C	0	GLY	-	expression tag	UNP H6QM90
C	761	GLY	-	expression tag	UNP H6QM90
C	762	TRP	-	expression tag	UNP H6QM90
C	763	SER	-	expression tag	UNP H6QM90
C	764	HIS	-	expression tag	UNP H6QM90

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Chain	Residue	Modelled	Actual	Comment	Reference
C	765	PRO	-	expression tag	UNP H6QM90
C	766	GLN	-	expression tag	UNP H6QM90
C	767	PHE	-	expression tag	UNP H6QM90
C	768	GLU	-	expression tag	UNP H6QM90
C	769	LYS	-	expression tag	UNP H6QM90
C	770	GLY	-	expression tag	UNP H6QM90
C	771	TRP	-	expression tag	UNP H6QM90
C	772	SER	-	expression tag	UNP H6QM90
C	773	HIS	-	expression tag	UNP H6QM90
C	774	PRO	-	expression tag	UNP H6QM90
C	775	GLN	-	expression tag	UNP H6QM90
C	776	PHE	-	expression tag	UNP H6QM90
C	777	GLU	-	expression tag	UNP H6QM90
C	778	LYS	-	expression tag	UNP H6QM90
C	779	GLY	-	expression tag	UNP H6QM90
C	780	SER	-	expression tag	UNP H6QM90
C	781	GLY	-	expression tag	UNP H6QM90
C	782	SER	-	expression tag	UNP H6QM90
C	783	GLU	-	expression tag	UNP H6QM90
C	784	ASN	-	expression tag	UNP H6QM90
C	785	LEU	-	expression tag	UNP H6QM90
C	786	TYR	-	expression tag	UNP H6QM90
C	787	PHE	-	expression tag	UNP H6QM90
C	788	GLN	-	expression tag	UNP H6QM90

- Molecule 4 is a RNA chain called RNA (5'-R(*UP*AP*UP*AP*CP*CP*UP*CP*UP*GP*CP*UP*U)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	R	13	Total	C	N	O	P	0	0	0
			247	111	37	87	12			

- Molecule 5 is a RNA chain called RNA 5'-pAGUAGUAACAAGAGGG.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	V	16	Total	C	N	O	P	0	0	0
			353	157	72	108	16			

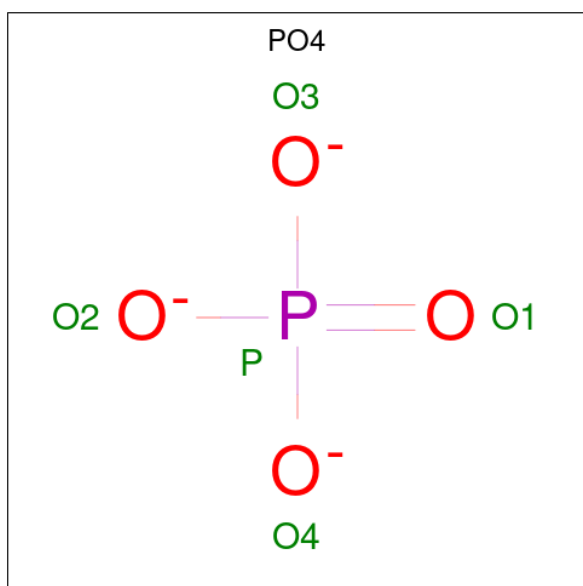
- Molecule 6 is a protein called TYR-SER-PRO-THR-SEP-PRO.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	X	6	Total 49	C 29	N 6	O 13	P 1	0	0	0
6	Y	10	Total 80	C 49	N 10	O 20	P 1	0	0	0

- Molecule 7 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	1	Total	Zn	0	0
			1	1		

- Molecule 8 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	A	1	Total	O	P	0	0
			5	4	1		
8	A	1	Total	O	P	0	0
			5	4	1		
8	A	1	Total	O	P	0	0
			5	4	1		
8	B	1	Total	O	P	0	0
			5	4	1		
8	B	1	Total	O	P	0	0
			5	4	1		
8	B	1	Total	O	P	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	B	1	Total	O	P	0	0
			5	4	1		
8	B	1	Total	O	P	0	0
			5	4	1		
8	B	1	Total	O	P	0	0
			5	4	1		
8	C	1	Total	O	P	0	0
			5	4	1		
8	C	1	Total	O	P	0	0
			5	4	1		
8	C	1	Total	O	P	0	0
			5	4	1		
8	C	1	Total	O	P	0	0
			5	4	1		

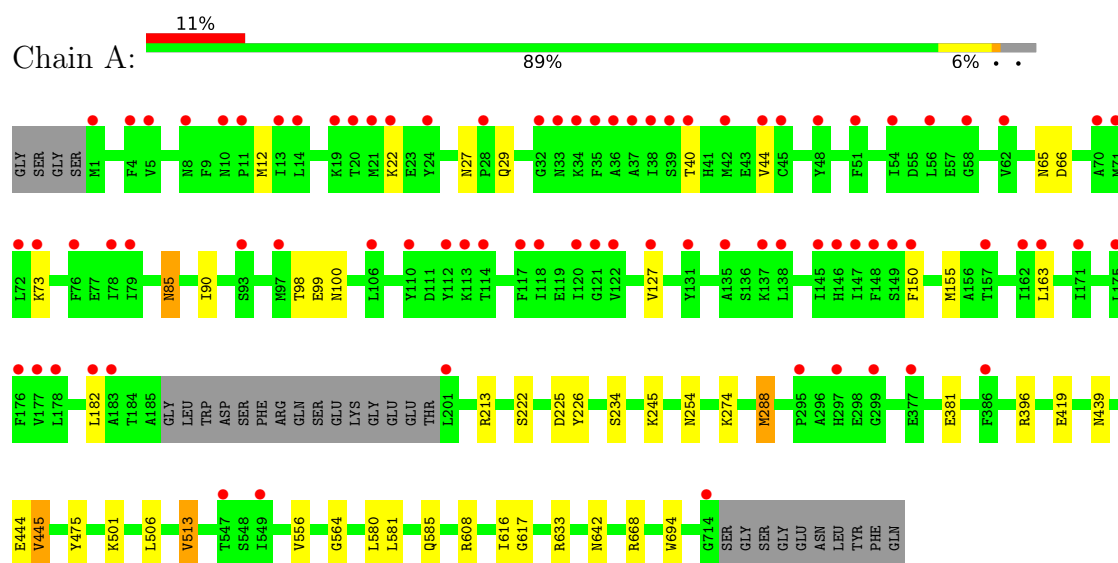
- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	78	Total	O	0	0
			78	78		
9	B	105	Total	O	0	0
			105	105		
9	C	56	Total	O	0	0
			56	56		
9	R	3	Total	O	0	0
			3	3		
9	V	7	Total	O	0	0
			7	7		
9	X	1	Total	O	0	0
			1	1		

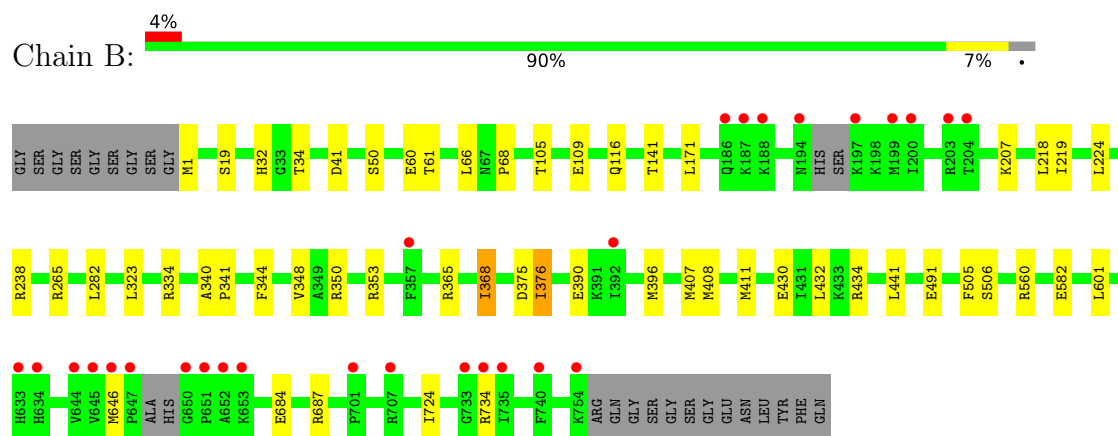
3 Residue-property plots [i](#)

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.

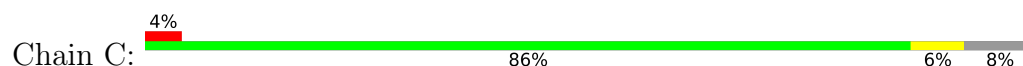
- Molecule 1: Polymerase acidic protein

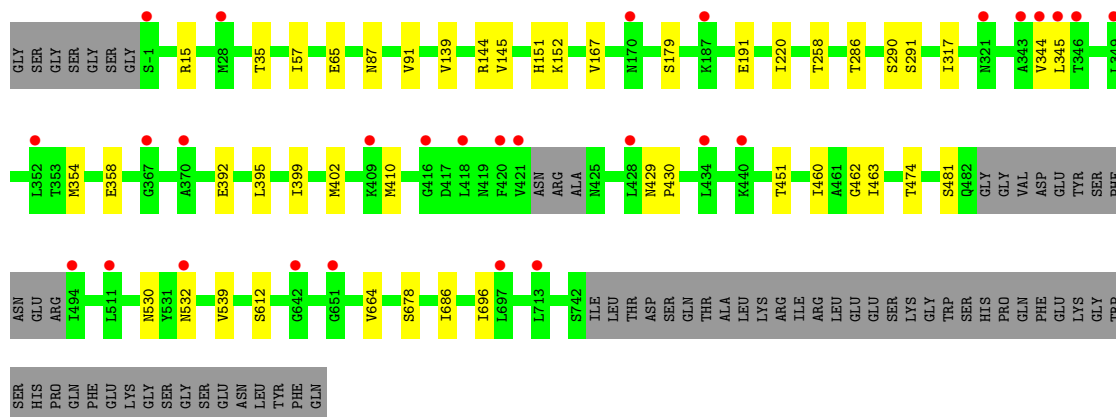


- Molecule 2: RNA-directed RNA polymerase catalytic subunit



- Molecule 3: Polymerase basic protein 2

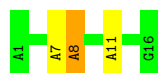
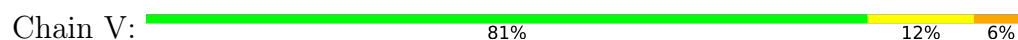




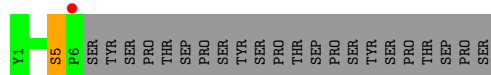
- Molecule 4: RNA (5'-R(*UP*AP*UP*AP*CP*CP*UP*CP*UP*GP*CP*UP*U)-3')



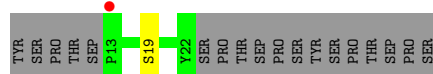
- Molecule 5: RNA 5'-pAGUAGUAACAAGAGGG



- Molecule 6: TYR-SER-PRO-THR-SEP-PRO



- Molecule 6: TYR-SER-PRO-THR-SEP-PRO



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	269.49Å 147.52Å 88.48Å 90.00° 97.22° 90.00°	Depositor
Resolution (Å)	49.53 – 2.50 49.53 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.6 (49.53-2.50) 83.1 (49.53-2.50)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.58 (at 2.51Å)	Xtriage
Refinement program	PHENIX (1.10_2155: ???)	Depositor
R, R_{free}	0.210 , 0.245 (Not available) , 0.236	Depositor DCC
R_{free} test set	5757 reflections (4.87%)	wwPDB-VP
Wilson B-factor (Å ²)	55.2	Xtriage
Anisotropy	0.219	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 50.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	18529	wwPDB-VP
Average B, all atoms (Å ²)	77.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.91% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: PO4, SEP, ZN

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.12	0/5812	0.30	0/7828
2	B	0.13	0/6100	0.40	3/8229 (0.0%)
3	C	0.11	0/5894	0.31	0/7955
4	R	0.20	0/273	0.42	0/421
5	V	0.11	0/397	0.30	0/617
6	X	0.28	0/40	0.20	0/53
6	Y	0.06	0/73	0.21	0/98
All	All	0.12	0/18589	0.34	3/25201 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	376	ILE	N-CA-C	15.89	142.38	109.34
2	B	375	ASP	CB-CA-C	9.44	129.20	110.42
2	B	376	ILE	N-CA-CB	-9.21	96.04	111.23

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	5692	0	5613	19	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	5986	0	6006	27	0
3	C	5796	0	5923	21	0
4	R	247	0	128	4	0
5	V	353	0	175	1	0
6	X	49	0	40	1	0
6	Y	80	0	64	0	0
7	A	1	0	0	0	0
8	A	15	0	0	0	0
8	B	35	0	0	0	0
8	C	25	0	0	1	0
9	A	78	0	0	0	0
9	B	105	0	0	0	0
9	C	56	0	0	1	0
9	R	3	0	0	0	0
9	V	7	0	0	0	0
9	X	1	0	0	0	0
All	All	18529	0	17949	65	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 2.

All (65) 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)
2:B:61:THR:HG22	2:B:218:LEU:HD23	1.36	1.05
4:R:4:A:H8	4:R:4:A:H5''	1.36	0.88
2:B:61:THR:CG2	2:B:218:LEU:HD23	2.06	0.84
4:R:4:A:H5''	4:R:4:A:C8	2.19	0.77
3:C:286:THR:HB	3:C:532:ASN:O	1.86	0.74
3:C:286:THR:O	3:C:532:ASN:ND2	2.19	0.73
2:B:61:THR:CG2	2:B:218:LEU:CD2	2.70	0.69
3:C:144:ARG:NH2	8:C:801:PO4:O1	2.28	0.67
2:B:61:THR:HG22	2:B:218:LEU:CD2	2.18	0.65
4:R:4:A:H2'	4:R:5:C:O4'	2.02	0.59
3:C:530:ASN:ND2	3:C:532:ASN:OD1	2.40	0.55
2:B:368:ILE:O	2:B:368:ILE:HG13	2.06	0.55
1:A:213:ARG:NH2	2:B:60:GLU:OE2	2.40	0.53
1:A:444:GLU:OE2	1:A:633:ARG:NH1	2.41	0.52
1:A:585:GLN:HG2	2:B:505:PHE:CD2	2.44	0.52
3:C:530:ASN:CG	3:C:532:ASN:OD1	2.53	0.51
2:B:34:THR:HG23	5:V:8:A:OP1	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:506:LEU:HD11	1:A:513:VAL:HG22	1.93	0.50
1:A:85:ASN:OD1	1:A:85:ASN:N	2.44	0.50
2:B:408:MET:HB3	2:B:411:MET:HE3	1.94	0.50
2:B:34:THR:HG22	2:B:353:ARG:HB2	1.94	0.49
2:B:344:PHE:CG	2:B:407:MET:HE1	2.47	0.49
3:C:290:SER:O	3:C:291:SER:OG	2.24	0.49
2:B:109:GLU:OE2	2:B:265:ARG:NH2	2.47	0.48
3:C:530:ASN:HD22	3:C:530:ASN:C	2.20	0.48
2:B:684:GLU:OE1	2:B:687:ARG:NH1	2.47	0.48
3:C:395:LEU:HD11	3:C:463:ILE:CD1	2.44	0.47
2:B:224:LEU:HD13	2:B:411:MET:HE2	1.96	0.47
2:B:19:SER:OG	2:B:506:SER:OG	2.28	0.47
3:C:57:ILE:O	3:C:91:VAL:N	2.48	0.46
1:A:98:THR:O	1:A:100:ASN:N	2.49	0.45
1:A:288:MET:HE1	1:A:556:VAL:HG13	1.98	0.45
2:B:61:THR:HG23	2:B:218:LEU:CD2	2.44	0.45
3:C:144:ARG:NH1	9:C:903:HOH:O	2.48	0.45
2:B:141:THR:HG21	3:C:35:THR:O	2.17	0.45
4:R:4:A:C8	4:R:4:A:C5'	2.96	0.45
2:B:323:LEU:O	2:B:334:ARG:NH1	2.51	0.44
1:A:633:ARG:NH2	6:X:5:SEP:O2P	2.49	0.44
3:C:429:ASN:HB2	3:C:430:PRO:HD2	2.00	0.44
3:C:145:VAL:HG21	3:C:220:ILE:HD11	1.99	0.43
1:A:225:ASP:OD1	1:A:226:TYR:N	2.52	0.43
1:A:513:VAL:O	1:A:564:GLY:HA2	2.19	0.43
3:C:151:HIS:ND1	3:C:151:HIS:O	2.52	0.43
1:A:642:ASN:OD1	1:A:694:TRP:NE1	2.46	0.42
1:A:445:VAL:CG1	1:A:580:LEU:HD22	2.49	0.42
1:A:475:TYR:HA	1:A:501:LYS:O	2.19	0.42
3:C:399:ILE:HD13	3:C:402:MET:CE	2.50	0.42
3:C:462:GLY:HA3	3:C:474:THR:HB	2.02	0.42
3:C:686:ILE:HA	3:C:696:ILE:O	2.20	0.42
2:B:282:LEU:HG	2:B:441:LEU:HD22	2.01	0.42
1:A:65:ASN:O	1:A:66:ASP:CB	2.67	0.41
1:A:90:ILE:HA	2:B:724:ILE:HD13	2.01	0.41
1:A:617:GLY:HA3	2:B:1:MET:HE3	2.01	0.41
2:B:34:THR:HG22	2:B:353:ARG:CB	2.51	0.41
3:C:429:ASN:HB2	3:C:430:PRO:CD	2.50	0.41
2:B:50:SER:HB3	2:B:68:PRO:HB3	2.03	0.41
1:A:274:LYS:O	1:A:396:ARG:HD2	2.20	0.41
1:A:234:SER:OG	2:B:334:ARG:NH2	2.54	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:61:THR:HB	2:B:219:ILE:HD11	2.03	0.41
3:C:344:VAL:HG12	3:C:345:LEU:N	2.36	0.41
1:A:40:THR:HG22	1:A:44:VAL:HG23	2.03	0.41
3:C:460:ILE:HD13	3:C:481:SER:HB2	2.03	0.40
3:C:286:THR:CB	3:C:532:ASN:O	2.64	0.40
2:B:340:ALA:HB3	2:B:341:PRO:HD3	2.03	0.40
2:B:432:LEU:HB3	2:B:434:ARG:HG2	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	
2	B	1/776 (0%)	0	1 (100%)	0	100	100
3	C	682/797 (86%)	661 (97%)	20 (3%)	1 (0%)	48	68
6	X	3/28 (11%)	3 (100%)	0	0	100	100
6	Y	7/28 (25%)	7 (100%)	0	0	100	100
All	All	693/1629 (42%)	671 (97%)	21 (3%)	1 (0%)	48	68

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	C	152	LYS

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	625/648 (96%)	600 (96%)	25 (4%)	28	54
2	B	659/676 (98%)	637 (97%)	22 (3%)	33	61
3	C	647/703 (92%)	629 (97%)	18 (3%)	38	66
6	X	5/24 (21%)	5 (100%)	0	100	100
6	Y	9/24 (38%)	9 (100%)	0	100	100
All	All	1945/2075 (94%)	1880 (97%)	65 (3%)	33	61

All (65) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	MET
1	A	22	LYS
1	A	27	ASN
1	A	29	GLN
1	A	73	LYS
1	A	85	ASN
1	A	99	GLU
1	A	127	VAL
1	A	150	PHE
1	A	155	MET
1	A	163	LEU
1	A	182	LEU
1	A	222	SER
1	A	245	LYS
1	A	254	ASN
1	A	288	MET
1	A	381	GLU
1	A	419	GLU
1	A	439	ASN
1	A	445	VAL
1	A	513	VAL
1	A	581	LEU
1	A	608	ARG
1	A	616	ILE
1	A	668	ARG
2	B	32	HIS
2	B	41	ASP
2	B	66	LEU

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Mol	Chain	Res	Type
2	B	105	THR
2	B	116	GLN
2	B	171	LEU
2	B	207	LYS
2	B	238	ARG
2	B	348	VAL
2	B	350	ARG
2	B	365	ARG
2	B	368	ILE
2	B	376	ILE
2	B	390	GLU
2	B	396	MET
2	B	430	GLU
2	B	491	GLU
2	B	560	ARG
2	B	582	GLU
2	B	601	LEU
2	B	646	MET
2	B	734	ARG
3	C	15	ARG
3	C	65	GLU
3	C	87	ASN
3	C	139	VAL
3	C	167	VAL
3	C	179	SER
3	C	191	GLU
3	C	258	THR
3	C	317	ILE
3	C	354	MET
3	C	358	GLU
3	C	392	GLU
3	C	410	MET
3	C	451	THR
3	C	539	VAL
3	C	612	SER
3	C	664	VAL
3	C	678	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	294	ASN

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Mol	Chain	Res	Type
1	A	416	ASN
2	B	116	GLN
2	B	137	GLN
2	B	184	HIS
2	B	186	GLN
2	B	562	HIS
2	B	679	GLN
3	C	160	GLN
3	C	383	GLN
3	C	530	ASN
3	C	540	ASN
3	C	548	ASN
3	C	563	GLN
3	C	601	GLN
3	C	624	HIS

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
4	R	11/18 (61%)	2 (18%)	1 (9%)
5	V	15/16 (93%)	3 (20%)	0
All	All	26/34 (76%)	5 (19%)	1 (3%)

All (5) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
4	R	4	A
4	R	5	C
5	V	7	A
5	V	8	A
5	V	11	A

All (1) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
4	R	4	A

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

2 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
6	SEP	X	5	6	8,9,10	1.61	1 (12%)	7,12,14	1.16	1 (14%)
6	SEP	Y	19	6	8,9,10	1.63	1 (12%)	7,12,14	1.41	2 (28%)

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
6	SEP	X	5	6	-	3/6/8/10	-
6	SEP	Y	19	6	-	3/6/8/10	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	Y	19	SEP	P-O1P	3.59	1.61	1.50
6	X	5	SEP	P-O1P	3.53	1.61	1.50

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	Y	19	SEP	OG-CB-CA	2.77	110.84	108.14
6	X	5	SEP	OG-CB-CA	2.49	110.57	108.14
6	Y	19	SEP	O2P-P-OG	2.18	112.34	106.67

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	X	5	SEP	C-CA-CB-OG
6	Y	19	SEP	CB-OG-P-O1P
6	Y	19	SEP	CB-OG-P-O2P
6	Y	19	SEP	CB-OG-P-O3P

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Mol	Chain	Res	Type	Atoms
6	X	5	SEP	N-CA-CB-OG
6	X	5	SEP	CA-CB-OG-P

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	X	5	SEP	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 16 ligands modelled in this entry, 1 is monoatomic - leaving 15 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	$\# Z > 2$
8	PO4	B	802	-	4,4,4	0.95	0	6,6,6	0.48	0
8	PO4	B	807	-	4,4,4	0.96	0	6,6,6	0.46	0
8	PO4	C	804	-	4,4,4	0.95	0	6,6,6	0.46	0
8	PO4	B	801	-	4,4,4	0.95	0	6,6,6	0.47	0
8	PO4	B	806	-	4,4,4	0.97	0	6,6,6	0.48	0
8	PO4	C	801	-	4,4,4	0.94	0	6,6,6	0.50	0
8	PO4	C	803	-	4,4,4	0.94	0	6,6,6	0.48	0
8	PO4	C	802	-	4,4,4	0.96	0	6,6,6	0.44	0
8	PO4	A	803	-	4,4,4	0.95	0	6,6,6	0.48	0
8	PO4	B	804	-	4,4,4	0.98	0	6,6,6	0.44	0
8	PO4	A	804	-	4,4,4	0.95	0	6,6,6	0.46	0
8	PO4	B	805	-	4,4,4	0.94	0	6,6,6	0.50	0
8	PO4	B	803	-	4,4,4	0.95	0	6,6,6	0.47	0
8	PO4	A	802	-	4,4,4	0.94	0	6,6,6	0.47	0
8	PO4	C	805	-	4,4,4	0.98	0	6,6,6	0.46	0

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	C	801	PO4	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	699/728 (96%)	0.57	81 (11%) 9 8	37, 63, 208, 256	0
2	B	750/776 (96%)	0.21	28 (3%) 45 40	36, 61, 114, 153	0
3	C	730/797 (91%)	0.41	28 (3%) 44 39	40, 70, 118, 157	0
4	R	13/18 (72%)	-0.20	0 100 100	52, 61, 88, 90	0
5	V	16/16 (100%)	-0.34	0 100 100	51, 54, 80, 141	0
6	X	5/28 (17%)	1.07	1 (20%) 3 2	66, 68, 77, 91	0
6	Y	9/28 (32%)	0.95	1 (11%) 10 8	58, 66, 107, 119	0
All	All	2222/2391 (92%)	0.39	139 (6%) 26 23	36, 65, 175, 256	0

All (139) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	121	GLY	6.6
1	A	40	THR	6.3
1	A	120	ILE	5.5
1	A	44	VAL	4.6
1	A	117	PHE	4.4
1	A	182	LEU	4.3
3	C	-1	SER	4.2
1	A	175	LEU	4.1
2	B	647	PRO	4.1
2	B	652	ALA	4.0
1	A	148	PHE	4.0
1	A	48	TYR	4.0
1	A	122	VAL	4.0
6	Y	13	PRO	3.8
1	A	73	LYS	3.8
3	C	494	ILE	3.8
1	A	5	VAL	3.7

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Mol	Chain	Res	Type	RSRZ
1	A	146	HIS	3.6
1	A	36	ALA	3.5
1	A	20	THR	3.5
1	A	145	ILE	3.5
3	C	345	LEU	3.5
3	C	170	ASN	3.4
1	A	299	GLY	3.4
1	A	38	ILE	3.4
1	A	72	LEU	3.3
1	A	39	SER	3.2
1	A	201	LEU	3.2
1	A	56	LEU	3.2
2	B	357	PHE	3.1
2	B	650	GLY	3.1
1	A	178	LEU	3.1
2	B	200	ILE	3.1
1	A	37	ALA	3.1
2	B	187	LYS	3.1
1	A	21	MET	3.1
1	A	297	HIS	3.1
1	A	32	GLY	3.0
1	A	70	ALA	3.0
1	A	42	MET	3.0
1	A	149	SER	3.0
1	A	549	ILE	3.0
3	C	420	PHE	3.0
2	B	186	GLN	3.0
1	A	71	MET	2.9
1	A	28	PRO	2.9
2	B	735	ILE	2.9
1	A	8	ASN	2.9
1	A	76	PHE	2.8
1	A	162	ILE	2.8
3	C	418	LEU	2.8
1	A	118	ILE	2.8
1	A	171	ILE	2.8
1	A	14	LEU	2.8
1	A	135	ALA	2.8
3	C	349	LEU	2.8
3	C	428	LEU	2.8
3	C	321	ASN	2.8
2	B	203	ARG	2.8

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Mol	Chain	Res	Type	RSRZ
3	C	440	LYS	2.8
1	A	97	MET	2.7
1	A	54	ILE	2.7
1	A	78	ILE	2.7
1	A	183	ALA	2.7
2	B	653	LYS	2.7
1	A	113	LYS	2.7
3	C	352	LEU	2.7
1	A	714	GLY	2.6
2	B	392	ILE	2.6
1	A	1	MET	2.6
3	C	434	LEU	2.6
1	A	137	LYS	2.6
2	B	754	LYS	2.6
2	B	707	ARG	2.6
2	B	651	PRO	2.5
2	B	645	VAL	2.5
3	C	344	VAL	2.5
3	C	532	ASN	2.5
1	A	22	LYS	2.5
1	A	51	PHE	2.5
1	A	19	LYS	2.5
2	B	194	ASN	2.5
3	C	346	THR	2.5
2	B	646	MET	2.5
6	X	6	PRO	2.5
3	C	367	GLY	2.5
1	A	177	VAL	2.5
1	A	295	PRO	2.4
2	B	197	LYS	2.4
1	A	79	ILE	2.4
1	A	112	TYR	2.4
2	B	204	THR	2.4
1	A	35	PHE	2.4
1	A	138	LEU	2.4
2	B	633	HIS	2.4
1	A	45	CYS	2.4
1	A	131	TYR	2.4
3	C	416	GLY	2.3
1	A	127	VAL	2.3
1	A	147	ILE	2.3
2	B	734	ARG	2.3

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Mol	Chain	Res	Type	RSRZ
3	C	28	MET	2.3
3	C	409	LYS	2.3
1	A	4	PHE	2.3
1	A	62	VAL	2.3
1	A	24	TYR	2.3
2	B	199	MET	2.2
1	A	547	THR	2.2
1	A	106	LEU	2.2
1	A	110	TYR	2.2
1	A	150	PHE	2.2
3	C	421	VAL	2.2
2	B	701	PRO	2.2
1	A	176	PHE	2.2
1	A	157	THR	2.1
1	A	386	PHE	2.1
1	A	58	GLY	2.1
1	A	114	THR	2.1
3	C	642	GLY	2.1
3	C	713	LEU	2.1
1	A	13	ILE	2.1
2	B	634	HIS	2.1
2	B	644	VAL	2.1
1	A	10	ASN	2.1
3	C	651	GLY	2.1
2	B	188	LYS	2.1
3	C	187	LYS	2.1
1	A	11	PRO	2.1
2	B	733	GLY	2.1
3	C	343	ALA	2.0
3	C	370	ALA	2.0
1	A	33	ASN	2.0
3	C	511	LEU	2.0
1	A	377	GLU	2.0
1	A	34	LYS	2.0
2	B	740	PHE	2.0
1	A	163	LEU	2.0
3	C	697	LEU	2.0
1	A	93	SER	2.0

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

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
6	SEP	X	5	10/11	0.91	0.08	65,74,89,92	0
6	SEP	Y	19	10/11	0.95	0.07	52,57,63,63	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
8	PO4	A	802	5/5	0.75	0.13	93,93,110,110	0
8	PO4	A	804	5/5	0.80	0.12	94,101,119,125	0
8	PO4	B	806	5/5	0.81	0.11	95,105,116,120	0
8	PO4	C	801	5/5	0.82	0.14	72,83,90,100	0
8	PO4	C	804	5/5	0.84	0.28	75,83,94,98	0
8	PO4	A	803	5/5	0.85	0.22	67,74,91,92	0
8	PO4	C	805	5/5	0.87	0.13	88,93,95,98	0
8	PO4	B	804	5/5	0.88	0.21	47,59,67,89	0
8	PO4	B	807	5/5	0.89	0.20	78,81,84,105	0
8	PO4	B	801	5/5	0.90	0.09	64,67,93,94	0
8	PO4	B	803	5/5	0.90	0.11	65,68,80,83	0
8	PO4	C	803	5/5	0.92	0.14	62,67,81,86	0
8	PO4	C	802	5/5	0.93	0.11	75,77,93,93	0
8	PO4	B	802	5/5	0.93	0.18	68,69,85,92	0
8	PO4	B	805	5/5	0.96	0.07	63,66,67,72	0
7	ZN	A	801	1/1	0.97	0.06	74,74,74,74	0

6.5 Other polymers [i](#)

There are no such residues in this entry.