



wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 1, 2016 – 03:17 PM GMT

PDB ID : 4BXX
Title : Structures of RNA polymerase II complexes with Bye1, a chromatin- binding PHF3 DIDO homologue
Authors : Kinkelin, K.; Wozniak, G.G.; Rothbart, S.B.; Lidschreiber, M.; Strahl, B.D.; Cramer, P.
Deposited on : 2013-07-16
Resolution : 3.28 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
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) : trunk26865

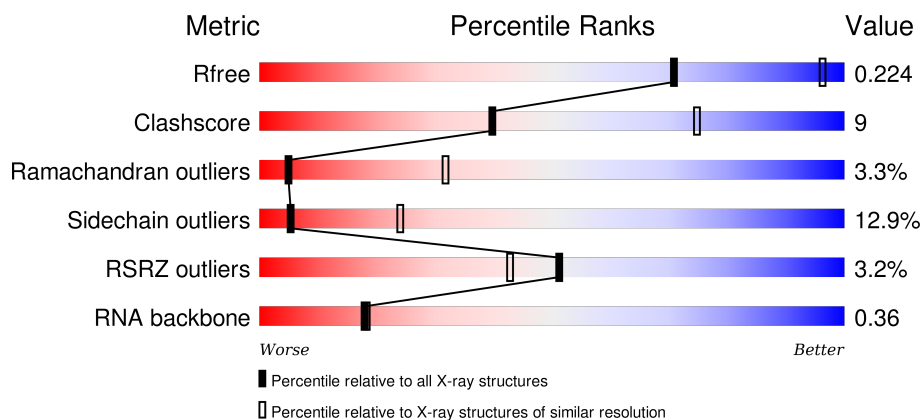
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.28 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}	91344	1756 (3.36-3.20)
Clashscore	102246	1941 (3.36-3.20)
Ramachandran outliers	100387	1905 (3.36-3.20)
Sidechain outliers	100360	1903 (3.36-3.20)
RSRZ outliers	91569	1764 (3.36-3.20)
RNA backbone	2183	1022 (3.80-2.76)

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	1733	<div> <div>1%</div> <div>57% 21% 18%</div> </div>
2	B	1224	<div> <div>2%</div> <div>63% 24% 10%</div> </div>
3	C	318	<div> <div>59% 20% 16%</div> </div>
4	D	221	<div> <div>59% 18% 20%</div> </div>

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Mol	Chain	Length	Quality of chain
5	E	215	
6	F	155	
7	G	171	
8	H	146	
9	I	122	
10	J	70	
11	K	120	
12	L	70	
13	N	14	
14	P	11	
15	T	27	
16	X	146	

2 Entry composition

There are 18 unique types of molecules in this entry. The entry contains 32756 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 DNA-DIRECTED RNA POLYMERASE II SUBUNIT RPB1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	1429	Total	C	N	O	S	0	0	0
			11249	7087	1964	2136	62			

- Molecule 2 is a protein called DNA-DIRECTED RNA POLYMERASE II SUBUNIT RPB2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	1097	Total	C	N	O	S	0	0	0
			8720	5526	1523	1617	54			

- Molecule 3 is a protein called DNA-DIRECTED RNA POLYMERASE II SUBUNIT RPB3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	266	Total	C	N	O	S	0	0	0
			2095	1317	348	417	13			

- Molecule 4 is a protein called DNA-DIRECTED RNA POLYMERASE II SUBUNIT RPB4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	177	Total	C	N	O	S	0	0	0
			1356	840	241	273	2			

- Molecule 5 is a protein called DNA-DIRECTED RNA POLYMERASES I, II, AND III SUBUNIT RPABC 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	E	214	Total	C	N	O	S	0	0	0
			1752	1111	309	321	11			

- Molecule 6 is a protein called DNA-DIRECTED RNA POLYMERASES I, II, AND III SUBUNIT RPABC 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	F	84	Total	C	N	O	S	0	0	0
			679	434	115	127	3			

- Molecule 7 is a protein called DNA-DIRECTED RNA POLYMERASE II SUBUNIT RPB7.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
7	G	171	Total	C	N	O	S	0	0	0
			1340	861	222	249	8			

- Molecule 8 is a protein called DNA-DIRECTED RNA POLYMERASES I, II, AND III SUBUNIT RPABC 3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
8	H	133	Total	C	N	O	S	0	0	0
			1068	673	180	211	4			

- Molecule 9 is a protein called DNA-DIRECTED RNA POLYMERASE II SUBUNIT RPB9.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
9	I	119	Total	C	N	O	S	0	0	0
			971	596	179	186	10			

- Molecule 10 is a protein called DNA-DIRECTED RNA POLYMERASES I, II, AND III SUBUNIT RPABC 5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
10	J	65	Total	C	N	O	S	0	0	0
			532	339	93	94	6			

- Molecule 11 is a protein called DNA-DIRECTED RNA POLYMERASE II SUBUNIT RPB11.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
11	K	115	Total	C	N	O	S	0	0	1
			920	590	157	171	2			

- Molecule 12 is a protein called DNA-DIRECTED RNA POLYMERASES I, II, AND III SUBUNIT RPABC 4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
12	L	46	Total	C	N	O	S	0	0	0
			363	224	72	63	4			

- Molecule 13 is a DNA chain called 5'-D(*GP*AP*GP*GP*TP*AP*AP*GP*CP*TP*AP*

GP*CP*TP)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
13	N	10	Total	C	N	O	P	0	0	0
			209	99	42	58	10			

- Molecule 14 is a RNA chain called 5'-D(*CP*CP*CP*CP*CP*CP*CP*CP*CP*CP)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
14	P	11	Total	C	N	O	P	0	0	0
			220	99	33	77	11			

- Molecule 15 is a DNA chain called 5'-D(*AP*GP*CP*TP*AP*GP*CP*TP*TP*AP*CP*CP*TP*GP *GP*TP*GP* BRUP*TP*GP*CP*TP*CP*TP*AP*AP*DC)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
15	T	14	Total	Br	C	N	O	P	0	0
			287	1	136	49	87	14		

- Molecule 16 is a protein called TRANSCRIPTION FACTOR BYE1.

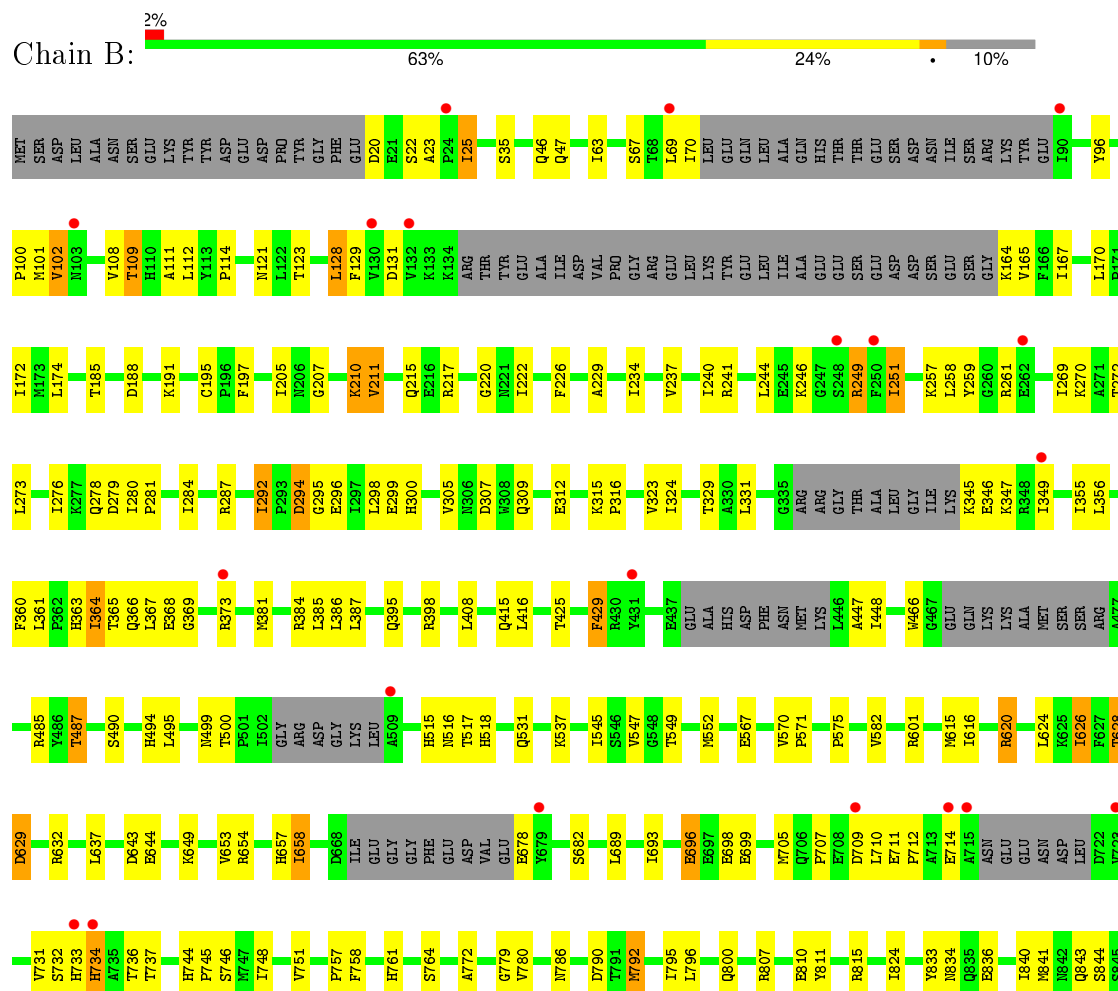
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
16	X	120	Total	C	N	O	S	0	0	0
			986	634	164	185	3			

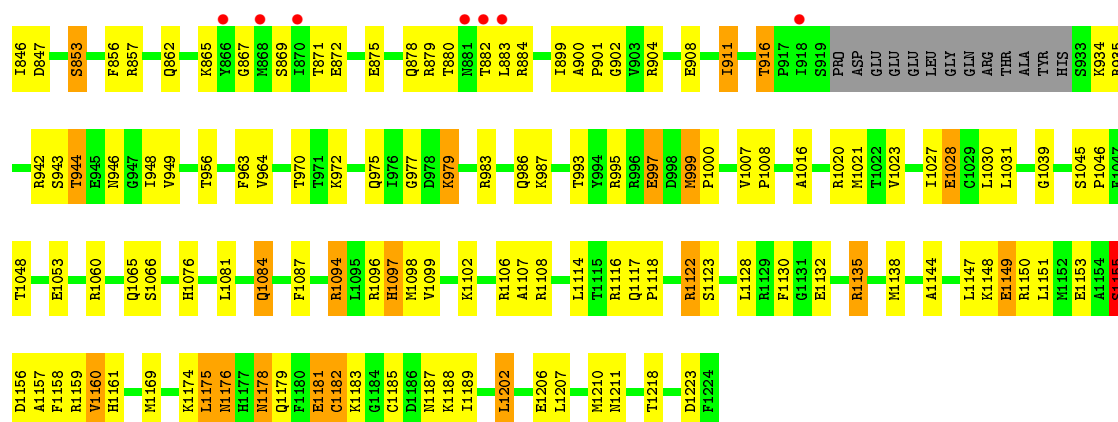
- Molecule 17 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
17	J	1	Total	Zn	0	0
			1	1		
17	B	1	Total	Zn	0	0
			1	1		
17	I	2	Total	Zn	0	0
			2	2		
17	C	1	Total	Zn	0	0
			1	1		
17	A	2	Total	Zn	0	0
			2	2		
17	L	1	Total	Zn	0	0
			1	1		

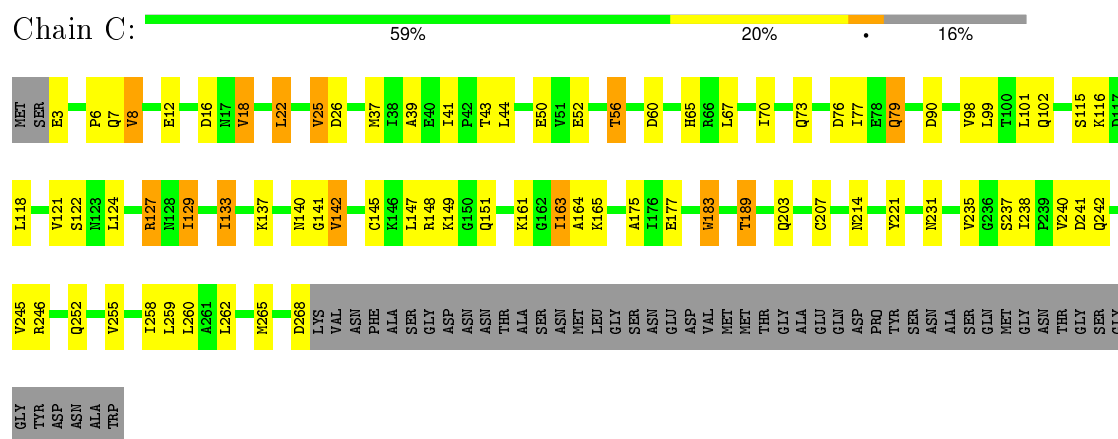
- Molecule 18 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
18	A	1	Total	Mg	0	0
			1	1		

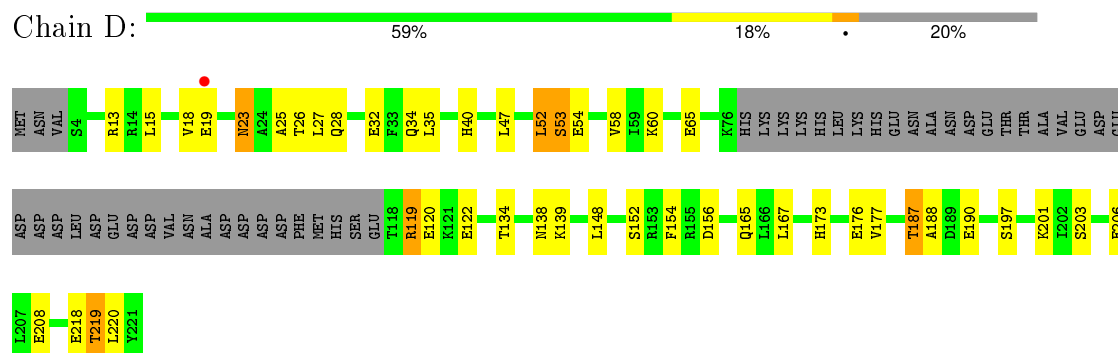




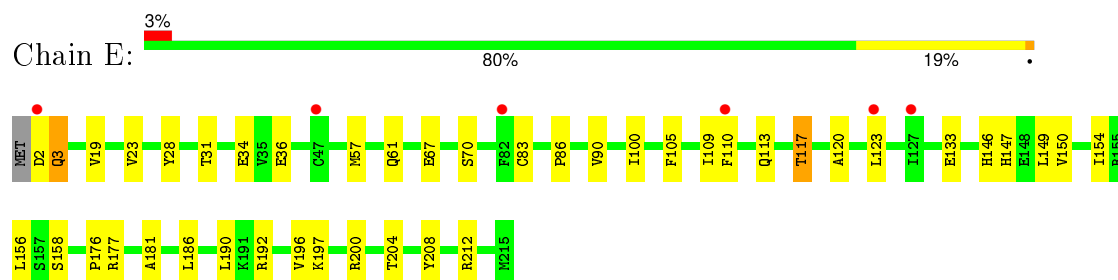
• Molecule 3: DNA-DIRECTED RNA POLYMERASE II SUBUNIT RPB3



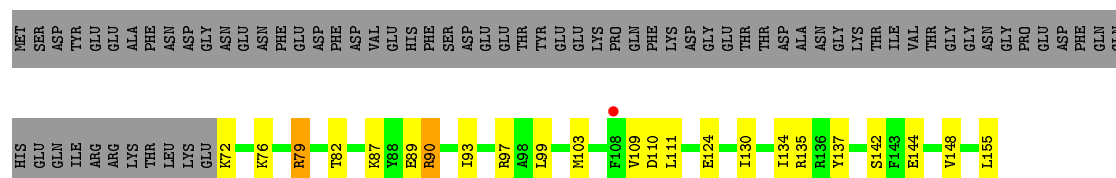
• Molecule 4: DNA-DIRECTED RNA POLYMERASE II SUBUNIT RPB4



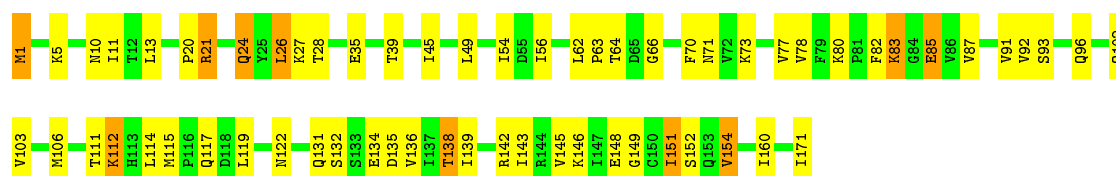
• Molecule 5: DNA-DIRECTED RNA POLYMERASES I, II, AND III SUBUNIT RPABC 1



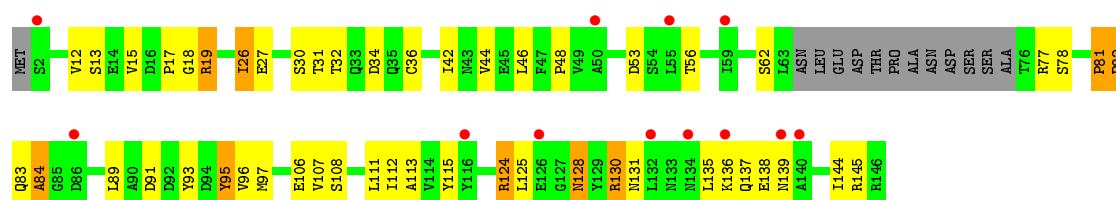
• Molecule 6: DNA-DIRECTED RNA POLYMERASES I, II, AND III SUBUNIT RPABC 2



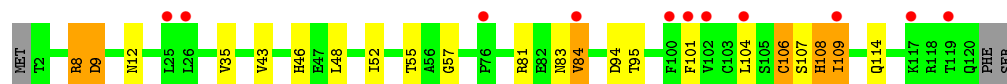
• Molecule 7: DNA-DIRECTED RNA POLYMERASE II SUBUNIT RPB7



• Molecule 8: DNA-DIRECTED RNA POLYMERASES I, II, AND III SUBUNIT RPABC 3



• Molecule 9: DNA-DIRECTED RNA POLYMERASE II SUBUNIT RPB9

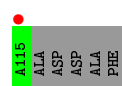


• Molecule 10: DNA-DIRECTED RNA POLYMERASES I, II, AND III SUBUNIT RPABC 5

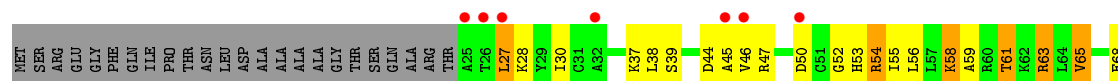
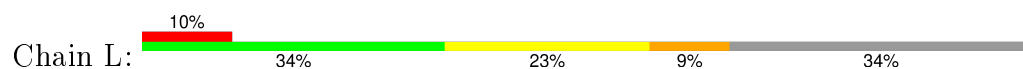


• Molecule 11: DNA-DIRECTED RNA POLYMERASE II SUBUNIT RPB11

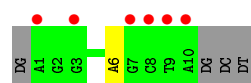
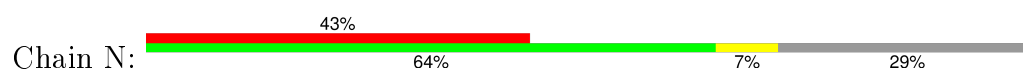




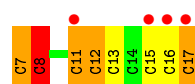
- Molecule 12: DNA-DIRECTED RNA POLYMERASES I, II, AND III SUBUNIT RPABC 4



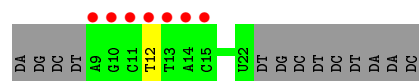
- Molecule 13: 5'-D(*GP*AP*GP*GP*TP*AP*AP*GP*CP*TP*AP*GP*CP*TP)-3'



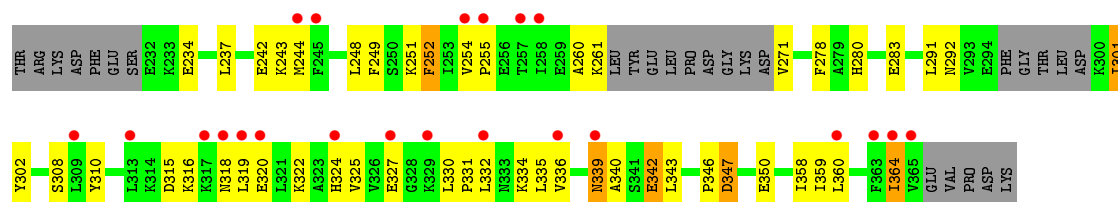
- Molecule 14: 5'-D(*CP*CP*CP*CP*CP*CP*CP*CP*CP*CP)-3'



- Molecule 15: 5'-D(*AP*GP*CP*TP*AP*GP*CP*TP*TP*AP*CP*CP*TP*GP *GP*TP*GP*BRUP*TP*GP*CP*TP*CP*TP*AP*AP*DC)-3'



- Molecule 16: TRANSCRIPTION FACTOR BYE1



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	222.92Å 392.67Å 281.04Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.08 – 3.28 49.08 – 3.28	Depositor EDS
% Data completeness (in resolution range)	100.0 (49.08-3.28) 100.0 (49.08-3.28)	Depositor EDS
R_{merge}	0.01	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.00 (at 3.25Å)	Xtriage
Refinement program	BUSTER 2.11.5	Depositor
R, R_{free}	0.180 , 0.208 0.200 , 0.224	Depositor DCC
R_{free} test set	3698 reflections (2.02%)	DCC
Wilson B-factor (Å ²)	98.3	Xtriage
Anisotropy	0.446	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 100.9	EDS
Estimated twinning fraction	0.012 for 1/2*h-1/2*k,-3/2*h-1/2*k,-l 0.018 for 1/2*h+1/2*k,3/2*h-1/2*k,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	1 of 187168 reflections (0.001%)	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	32756	wwPDB-VP
Average B, all atoms (Å ²)	120.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.80% of the height of the origin peak. No significant pseudotranslation is detected.*

¹ Intensities estimated from amplitudes.

² Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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, ZN, BRU

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	0.54	0/11450	0.82	10/15483 (0.1%)
2	B	0.51	0/8889	0.77	3/11987 (0.0%)
3	C	0.50	0/2133	0.79	1/2891 (0.0%)
4	D	0.53	0/1365	0.84	2/1837 (0.1%)
5	E	0.45	0/1788	0.67	0/2406
6	F	0.60	0/691	0.78	0/933
7	G	0.53	0/1368	0.82	1/1844 (0.1%)
8	H	0.49	0/1086	0.76	0/1470
9	I	0.45	0/989	0.72	0/1331
10	J	0.54	0/541	0.82	0/727
11	K	0.48	0/938	0.70	0/1267
12	L	0.56	0/365	0.90	0/485
13	N	0.99	0/235	0.89	0/361
14	P	1.59	5/241 (2.1%)	1.04	0/370
15	T	1.27	0/298	0.99	0/458
16	X	0.50	0/999	0.75	0/1336
All	All	0.55	5/33376 (0.0%)	0.79	17/45186 (0.0%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
14	P	17	C	C1'-N1	5.93	1.57	1.48
14	P	15	C	C1'-N1	5.59	1.57	1.48
14	P	11	C	C3'-O3'	5.20	1.49	1.42
14	P	8	C	C1'-N1	5.16	1.56	1.48
14	P	7	C	C1'-N1	5.16	1.56	1.48

The worst 5 of 17 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	3	GLY	C-N-CA	8.18	142.14	121.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	25	ALA	C-N-CA	7.72	141.00	121.70
1	A	56	PRO	C-N-CA	6.95	139.07	121.70
1	A	399	HIS	N-CA-CB	6.93	123.07	110.60
1	A	54	ASN	C-N-CA	6.55	138.09	121.70

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	11249	0	11312	256	0
2	B	8720	0	8744	151	0
3	C	2095	0	2051	40	0
4	D	1356	0	1319	20	0
5	E	1752	0	1776	21	0
6	F	679	0	701	19	0
7	G	1340	0	1357	40	0
8	H	1068	0	1040	31	0
9	I	971	0	927	13	0
10	J	532	0	542	13	0
11	K	920	0	929	20	0
12	L	363	0	386	10	0
13	N	209	0	113	1	0
14	P	220	0	122	2	0
15	T	287	0	157	1	0
16	X	986	0	1009	18	0
17	A	2	0	0	0	0
17	B	1	0	0	0	0
17	C	1	0	0	0	0
17	I	2	0	0	0	0
17	J	1	0	0	0	0
17	L	1	0	0	0	0
18	A	1	0	0	0	0
All	All	32756	0	32485	571	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 9.

The worst 5 of 571 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:A:867:ILE:CD1	1:A:867:ILE:CG1	1.81	1.55
1:A:53:LEU:HD23	1:A:54:ASN:H	1.06	1.10
1:A:567:LYS:HB3	8:H:96:VAL:H	1.21	1.03
1:A:855:THR:HG21	1:A:857:ARG:HE	1.22	1.02
6:F:93:ILE:HD11	6:F:134:ILE:HD11	1.39	1.01

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	1419/1733 (82%)	1259 (89%)	111 (8%)	49 (4%)	4	30
2	B	1075/1224 (88%)	962 (90%)	78 (7%)	35 (3%)	5	32
3	C	264/318 (83%)	242 (92%)	15 (6%)	7 (3%)	6	37
4	D	173/221 (78%)	151 (87%)	17 (10%)	5 (3%)	6	35
5	E	212/215 (99%)	201 (95%)	9 (4%)	2 (1%)	21	63
6	F	82/155 (53%)	75 (92%)	7 (8%)	0	100	100
7	G	169/171 (99%)	153 (90%)	15 (9%)	1 (1%)	30	71
8	H	129/146 (88%)	100 (78%)	16 (12%)	13 (10%)	1	4
9	I	117/122 (96%)	99 (85%)	16 (14%)	2 (2%)	11	50
10	J	63/70 (90%)	54 (86%)	5 (8%)	4 (6%)	2	14
11	K	113/120 (94%)	108 (96%)	5 (4%)	0	100	100
12	L	44/70 (63%)	27 (61%)	11 (25%)	6 (14%)	0	2

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
16	X	114/146 (78%)	99 (87%)	9 (8%)	6 (5%)	2	18
All	All	3974/4711 (84%)	3530 (89%)	314 (8%)	130 (3%)	5	32

5 of 130 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	4	GLN
1	A	44	THR
1	A	54	ASN
1	A	57	ARG
1	A	70	CYS

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	1251/1520 (82%)	1084 (87%)	167 (13%)	5	22
2	B	952/1061 (90%)	835 (88%)	117 (12%)	6	26
3	C	234/274 (85%)	204 (87%)	30 (13%)	5	24
4	D	140/200 (70%)	123 (88%)	17 (12%)	6	27
5	E	196/197 (100%)	183 (93%)	13 (7%)	21	60
6	F	74/137 (54%)	68 (92%)	6 (8%)	15	49
7	G	152/152 (100%)	130 (86%)	22 (14%)	4	19
8	H	117/128 (91%)	99 (85%)	18 (15%)	3	16
9	I	113/116 (97%)	102 (90%)	11 (10%)	10	38
10	J	60/65 (92%)	50 (83%)	10 (17%)	3	13
11	K	99/102 (97%)	88 (89%)	11 (11%)	8	32
12	L	40/57 (70%)	28 (70%)	12 (30%)	0	1
16	X	110/136 (81%)	88 (80%)	22 (20%)	1	7
All	All	3538/4145 (85%)	3082 (87%)	456 (13%)	5	24

5 of 456 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	537	LYS
2	B	1053	GLU
11	K	113	THR
2	B	601	ARG
2	B	795	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 79 such sidechains are listed below:

Mol	Chain	Res	Type
2	B	300	HIS
2	B	744	HIS
11	K	65	HIS
2	B	325	GLN
2	B	516	ASN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
14	P	10/11 (90%)	5 (50%)	2 (20%)

All (5) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
14	P	8	C
14	P	11	C
14	P	12	C
14	P	16	C
14	P	17	C

All (2) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
14	P	11	C
14	P	16	C

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

1 non-standard protein/DNA/RNA residue is 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	$\# Z > 2$
15	BRU	T	22	15	13,21,22	2.67	1 (7%)	16,30,33	2.28	3 (18%)

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
15	BRU	T	22	15	-	0/3/21/22	0/2/2/2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
15	T	22	BRU	C4-C5	9.41	1.50	1.38

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
15	T	22	BRU	C5-C4-N3	-4.42	119.28	124.00
15	T	22	BRU	O4'-C1'-N1	2.60	112.21	107.72
15	T	22	BRU	C4-N3-C2	7.25	121.52	115.25

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry

Of 9 ligands modelled in this entry, 9 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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	1429/1733 (82%)	0.02	24 (1%)	73	65	59, 101, 164, 229	0
2	B	1097/1224 (89%)	0.15	27 (2%)	61	52	59, 114, 175, 208	0
3	C	266/318 (83%)	-0.13	0	100	100	77, 101, 141, 173	0
4	D	177/221 (80%)	-0.05	1 (0%)	90	87	80, 111, 165, 179	0
5	E	214/215 (99%)	0.09	6 (2%)	56	48	76, 135, 181, 211	0
6	F	84/155 (54%)	-0.22	1 (1%)	81	74	63, 83, 110, 134	0
7	G	171/171 (100%)	0.06	0	100	100	73, 100, 139, 166	0
8	H	133/146 (91%)	0.63	12 (9%)	12	8	112, 148, 183, 209	0
9	I	119/122 (97%)	0.38	11 (9%)	11	8	116, 148, 186, 203	0
10	J	65/70 (92%)	-0.10	0	100	100	82, 101, 138, 146	0
11	K	115/120 (95%)	0.03	2 (1%)	73	65	70, 103, 141, 155	0
12	L	46/70 (65%)	0.93	7 (15%)	3	2	86, 180, 196, 201	0
13	N	10/14 (71%)	2.32	6 (60%)	0	0	239, 248, 270, 275	0
14	P	11/11 (100%)	2.30	4 (36%)	0	0	204, 214, 222, 224	0
15	T	13/27 (48%)	2.01	7 (53%)	0	0	165, 186, 260, 263	0
16	X	120/146 (82%)	0.98	22 (18%)	2	1	172, 203, 219, 229	0
All	All	4070/4763 (85%)	0.13	130 (3%)	51	43	59, 110, 188, 275	0

The worst 5 of 130 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
11	K	115	ALA	7.9
12	L	25	ALA	7.9
14	P	16	C	6.8
12	L	26	THR	6.1
1	A	69	THR	5.8

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. 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
15	BRU	T	22	20/21	0.64	0.32	-	191,208,212,215	0

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
17	ZN	B	2225	1/1	0.99	0.22	0.89	83,83,83,83	0
17	ZN	I	1121	1/1	0.99	0.15	-0.06	124,124,124,124	0
17	ZN	J	1066	1/1	0.99	0.23	-0.54	90,90,90,90	0
17	ZN	A	2457	1/1	1.00	0.17	-0.83	79,79,79,79	0
17	ZN	C	1269	1/1	1.00	0.12	-0.88	81,81,81,81	0
17	ZN	A	2456	1/1	0.99	0.06	-1.78	141,141,141,141	0
18	MG	A	2458	1/1	0.92	0.10	-1.82	77,77,77,77	0
17	ZN	I	1122	1/1	0.97	0.03	-1.95	196,196,196,196	0
17	ZN	L	1071	1/1	0.98	0.06	-2.50	207,207,207,207	0

6.5 Other polymers [i](#)

There are no such residues in this entry.