



# Full wwPDB X-ray Structure Validation Report ⓘ

Nov 18, 2020 – 06:02 PM EST

PDB ID : 3N4M  
Title : E. coli RNA polymerase alpha subunit C-terminal domain in complex with CAP and DNA  
Authors : Lara-Gonzalez, S.; Birktoft, J.J.; Lawson, C.L.  
Deposited on : 2010-05-21  
Resolution : 2.99 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](mailto: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.

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

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.14.6  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.14.6

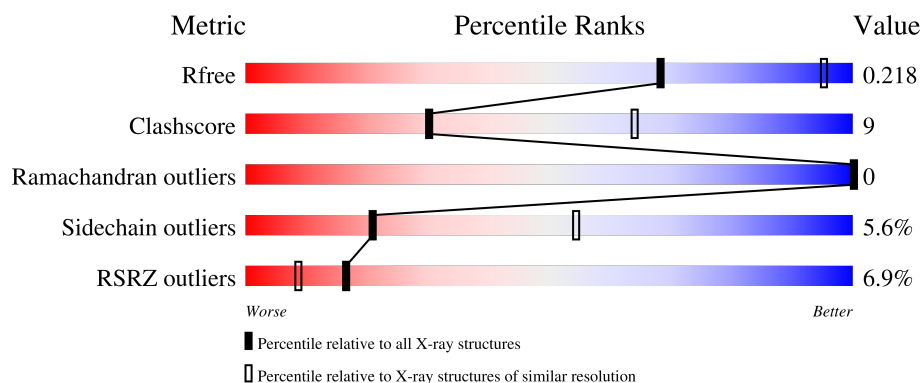
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*





The reported resolution of this entry is 2.99 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))
$R_{free}$	130704	2754 (3.00-2.96)
Clashscore	141614	3103 (3.00-2.96)
Ramachandran outliers	138981	2993 (3.00-2.96)
Sidechain outliers	138945	2996 (3.00-2.96)
RSRZ outliers	127900	2644 (3.00-2.96)

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  $\geq 3$ , 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	209	
2	B	84	
2	C	84	
3	D	20	
4	E	24	

## 2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 3737 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 Catabolite gene activator.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	203	Total	C	N	O	S	0	0	0
			1604	1014	282	299	9			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	74	Total	C	N	O	S	0	0	0
			583	371	101	109	2			
2	C	73	Total	C	N	O	S	0	0	0
			572	362	100	108	2			

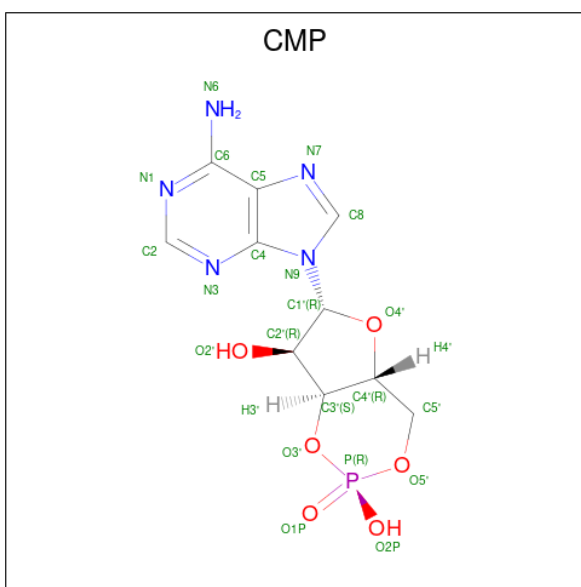
- Molecule 3 is a DNA chain called DNA (5'-D(\*CP\*TP\*TP\*TP\*TP\*TP\*TP\*CP\*CP\*TP\*AP\*AP\*AP\*AP\*TP\*GP\*TP\*GP\*AP\*T)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	20	Total	C	N	O	P	0	0	0
			403	197	64	123	19			

- Molecule 4 is a DNA chain called DNA (5'-D(\*CP\*TP\*AP\*GP\*AP\*TP\*CP\*AP\*CP\*AP\*TP\*TP\*TP\*TP\*AP\*GP\*GP\*AP\*AP\*AP\*AP\*AP\*AP\*G)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	E	24	Total	C	N	O	P	0	0	0
			493	237	96	137	23			

- Molecule 5 is ADENOSINE-3',5'-CYCLIC-MONOPHOSPHATE (three-letter code: CMP) (formula: C<sub>10</sub>H<sub>12</sub>N<sub>5</sub>O<sub>6</sub>P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	N	O	P	0	0
			22	10	5	6	1		
5	D	1	Total	C	N			0	0
			11	6	5				

- Molecule 6 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula:  $C_4H_{10}O_3$ ).



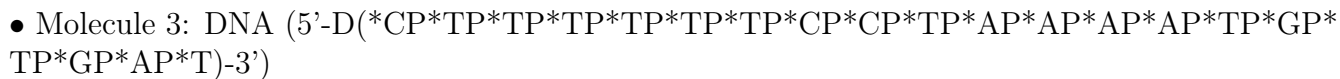
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			7	4	3		
6	A	1	Total	C	O	0	0
			7	4	3		

- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	11	Total 11	O 11	0	0
7	D	13	Total 13	O 13	0	0
7	E	11	Total 11	O 11	0	0



- Molecule 1: Catabolite gene activator



C21	T22	A23	G24	C27	A28	T33	G36	G44
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## 4 Data and refinement statistics

Property	Value	Source
Space group	P 62 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	175.73Å 175.73Å 160.10Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	38.51 – 2.99 38.51 – 2.99	Depositor EDS
% Data completeness (in resolution range)	96.5 (38.51-2.99) 96.5 (38.51-2.99)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	0.09	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.89 (at 3.01Å)	Xtriage
Refinement program	PHENIX	Depositor
R, $R_{free}$	0.198 , 0.224 0.198 , 0.218	Depositor DCC
$R_{free}$ test set	1484 reflections (5.10%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	87.8	Xtriage
Anisotropy	0.378	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.29 , 65.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	3737	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	110.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.58% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: PEG, CMP

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	0.29	0/1629	0.47	0/2193
2	B	0.21	0/591	0.41	0/800
2	C	0.21	0/579	0.41	0/784
3	D	0.53	0/449	1.35	11/691 (1.6%)
4	E	0.52	0/555	1.20	2/855 (0.2%)
All	All	0.35	0/3803	0.78	13/5323 (0.2%)

There are no bond length outliers.

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
3	D	18	DG	O4'-C4'-C3'	-6.68	101.83	104.50
3	D	18	DG	C4'-C3'-C2'	-6.01	97.69	103.10
4	E	36	DG	O4'-C1'-C2'	-5.54	101.47	105.90
3	D	2	DT	N3-C4-O4	5.49	123.19	119.90
3	D	7	DT	N3-C4-O4	5.31	123.09	119.90
3	D	4	DT	N3-C4-O4	5.21	123.03	119.90
4	E	33	DT	N3-C4-O4	5.17	123.00	119.90
3	D	2	DT	C5-C4-O4	-5.10	121.33	124.90
3	D	6	DT	N3-C4-O4	5.09	122.95	119.90
3	D	5	DT	O4'-C1'-N1	-5.07	104.45	108.00
3	D	20	DT	N3-C4-O4	5.06	122.94	119.90
3	D	6	DT	C5-C4-O4	-5.02	121.39	124.90
3	D	3	DT	C5-C4-O4	-5.00	121.40	124.90

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	1604	0	1637	29	0
2	B	583	0	611	21	0
2	C	572	0	602	13	0
3	D	403	0	232	3	0
4	E	493	0	272	7	0
5	A	22	0	11	1	0
5	D	11	0	4	1	0
6	A	14	0	20	2	0
7	A	11	0	0	1	0
7	D	13	0	0	0	0
7	E	11	0	0	0	0
All	All	3737	0	3389	66	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.

All (66) 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)
5:A:210:CMP:H2	5:A:210:CMP:C2	0.97	1.50
5:D:100:CMP:H2	5:D:100:CMP:C2	0.97	1.48
1:A:109:ASN:HD22	6:A:501:PEG:H22	1.46	0.80
2:B:285:THR:HG23	2:B:288:GLU:HB2	1.63	0.80
2:C:285:THR:HG23	2:C:288:GLU:HB2	1.63	0.78
1:A:160:PRO:HD3	2:B:285:THR:HG21	1.67	0.76
1:A:158:THR:HG23	2:B:286:GLU:HB2	1.73	0.69
4:E:23:DA:H2"	4:E:24:DG:H5"	1.76	0.68
2:B:253:LEU:HD13	2:B:318:LEU:HD21	1.78	0.65
2:C:253:LEU:HD13	2:C:318:LEU:HD21	1.78	0.65
1:A:76:PHE:HB2	7:A:218:HOH:O	1.99	0.63
1:A:208:THR:HG22	1:A:209:ARG:H	1.64	0.63
2:C:279:GLY:HA3	2:C:321:TRP:CZ2	2.34	0.62
2:B:279:GLY:HA3	2:B:321:TRP:CZ2	2.36	0.61
1:A:50:LEU:HD12	1:A:85:TRP:HB2	1.83	0.60
2:C:279:GLY:HA3	2:C:321:TRP:CH2	2.37	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:279:GLY:HA3	2:B:321:TRP:CH2	2.38	0.59
1:A:109:ASN:ND2	6:A:501:PEG:H22	2.18	0.57
4:E:23:DA:C2'	4:E:24:DG:H5''	2.34	0.57
1:A:209:ARG:HB3	2:B:317:ARG:HH21	1.70	0.56
1:A:169:ARG:NH2	1:A:200:GLY:O	2.38	0.56
4:E:21:DC:H2''	4:E:22:DT:H5'	1.87	0.56
4:E:23:DA:H2''	4:E:24:DG:C5'	2.35	0.56
1:A:158:THR:HG23	2:B:286:GLU:CB	2.36	0.55
3:D:17:DT:H2''	3:D:18:DG:H5'	1.88	0.54
1:A:75:LEU:HG	1:A:75:LEU:O	2.06	0.54
2:C:285:THR:HG23	2:C:288:GLU:CB	2.37	0.54
1:A:160:PRO:HD3	2:B:285:THR:CG2	2.38	0.54
2:B:285:THR:HG23	2:B:288:GLU:CB	2.37	0.54
1:A:143:ILE:HD11	1:A:183:VAL:HG22	1.92	0.52
1:A:137:LEU:HB2	1:A:142:ARG:HG2	1.93	0.50
1:A:159:HIS:HB3	1:A:162:GLY:O	2.12	0.50
1:A:160:PRO:CD	2:B:285:THR:HG21	2.40	0.49
4:E:27:DC:H2''	4:E:28:DA:H5'	1.93	0.49
1:A:159:HIS:CE1	1:A:160:PRO:HG2	2.48	0.49
2:C:286:GLU:HG3	2:C:300:LEU:HD11	1.96	0.48
1:A:159:HIS:CG	1:A:160:PRO:HD2	2.48	0.48
2:B:286:GLU:HG3	2:B:300:LEU:HD11	1.96	0.47
2:B:251:PRO:O	2:B:254:LEU:HG	2.15	0.47
2:C:251:PRO:O	2:C:254:LEU:HG	2.15	0.47
3:D:18:DG:H1'	3:D:19:DA:C8	2.50	0.47
1:A:138:ASP:HB2	4:E:24:DG:OP1	2.16	0.46
2:B:250:ASP:HB3	2:B:253:LEU:HG	1.97	0.46
4:E:27:DC:C2'	4:E:28:DA:H5'	2.46	0.46
1:A:42:ILE:CG1	1:A:70:ILE:HD11	2.46	0.45
2:C:250:ASP:HB3	2:C:253:LEU:HG	1.97	0.45
1:A:153:GLN:HB3	1:A:154:PRO:HD2	1.99	0.45
1:A:138:ASP:O	1:A:142:ARG:HG3	2.16	0.44
2:C:292:THR:HA	2:C:293:PRO:HD3	1.88	0.44
2:B:263:THR:HG23	2:B:302:GLU:OE2	2.16	0.44
1:A:158:THR:HG23	2:B:286:GLU:H	1.82	0.44
1:A:35:LYS:HD3	1:A:35:LYS:HA	1.84	0.44
2:C:263:THR:HG23	2:C:302:GLU:OE2	2.17	0.44
2:B:249:PHE:N	2:B:249:PHE:CD2	2.86	0.44
2:B:293:PRO:O	2:B:294:ASN:HB2	2.18	0.43
1:A:119:GLN:O	1:A:123:ARG:HG3	2.19	0.42
2:B:293:PRO:HB3	2:C:263:THR:HG21	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:293:PRO:O	2:C:294:ASN:HB2	2.18	0.42
1:A:155:ASP:OD1	1:A:155:ASP:N	2.48	0.42
3:D:4:DT:C6	3:D:5:DT:H72	2.55	0.41
1:A:195:LEU:O	1:A:206:TYR:HB2	2.20	0.41
2:B:250:ASP:HA	2:B:251:PRO:HD3	1.87	0.41
1:A:122:ARG:O	1:A:126:VAL:HG23	2.21	0.41
1:A:181:GLU:O	1:A:185:ARG:HG3	2.20	0.41
2:B:321:TRP:HA	2:B:322:PRO:C	2.42	0.41
2:C:321:TRP:HA	2:C:322:PRO:C	2.42	0.41

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	201/209 (96%)	193 (96%)	8 (4%)	0	100	100
2	B	72/84 (86%)	67 (93%)	5 (7%)	0	100	100
2	C	71/84 (84%)	66 (93%)	5 (7%)	0	100	100
All	All	344/377 (91%)	326 (95%)	18 (5%)	0	100	100

There are no Ramachandran outliers to report.

### 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	174/180 (97%)	168 (97%)	6 (3%)	37	70
2	B	66/74 (89%)	60 (91%)	6 (9%)	9	32
2	C	65/74 (88%)	60 (92%)	5 (8%)	13	40
All	All	305/328 (93%)	288 (94%)	17 (6%)	21	54

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	50	LEU
1	A	59	MET
1	A	70	ILE
1	A	122	ARG
1	A	158	THR
1	A	169	ARG
2	B	249	PHE
2	B	250	ASP
2	B	265	ARG
2	B	285	THR
2	B	317	ARG
2	B	318	LEU
2	C	250	ASP
2	C	265	ARG
2	C	285	THR
2	C	317	ARG
2	C	318	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	21	HIS
1	A	66	GLN
1	A	125	GLN
1	A	149	ASN
1	A	170	GLN
1	A	193	GLN
2	B	320	ASN
2	C	320	ASN

### 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 ⓘ

There are no monosaccharides in this entry.

## 5.6 Ligand geometry ⓘ

4 ligands 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
5	CMP	D	100	-	10,12,25	1.31	2 (20%)	8,17,39	1.67	2 (25%)
6	PEG	A	502	-	6,6,6	0.64	0	5,5,5	0.64	0
6	PEG	A	501	-	6,6,6	0.61	0	5,5,5	0.61	0
5	CMP	A	210	-	22,25,25	1.44	4 (18%)	24,39,39	1.41	5 (20%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	CMP	D	100	-	-	-	0/2/2/4
6	PEG	A	502	-	-	2/4/4/4	-
6	PEG	A	501	-	-	4/4/4/4	-
5	CMP	A	210	-	-	0/0/31/31	0/4/4/4

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	210	CMP	P-O5'	3.25	1.61	1.57
5	A	210	CMP	C5-C4	2.66	1.48	1.40
5	D	100	CMP	C5-C4	2.65	1.47	1.40
5	A	210	CMP	P-O3'	2.35	1.61	1.57
5	A	210	CMP	O5'-C5'	-2.32	1.42	1.46
5	D	100	CMP	C2-N3	2.02	1.35	1.32

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	D	100	CMP	N3-C2-N1	-3.42	123.33	128.68
5	A	210	CMP	N3-C2-N1	-3.10	123.84	128.68
5	A	210	CMP	C4-C5-N7	-2.34	106.96	109.40
5	A	210	CMP	O3'-C3'-C2'	2.32	117.88	115.61
5	D	100	CMP	C4-C5-N7	-2.27	107.04	109.40
5	A	210	CMP	O3'-C3'-C4'	-2.12	109.11	110.71
5	A	210	CMP	O5'-P-O3'	-2.01	102.91	105.68

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	501	PEG	O1-C1-C2-O2
6	A	502	PEG	O2-C3-C4-O4
6	A	501	PEG	O2-C3-C4-O4
6	A	501	PEG	C1-C2-O2-C3
6	A	501	PEG	C4-C3-O2-C2
6	A	502	PEG	C4-C3-O2-C2

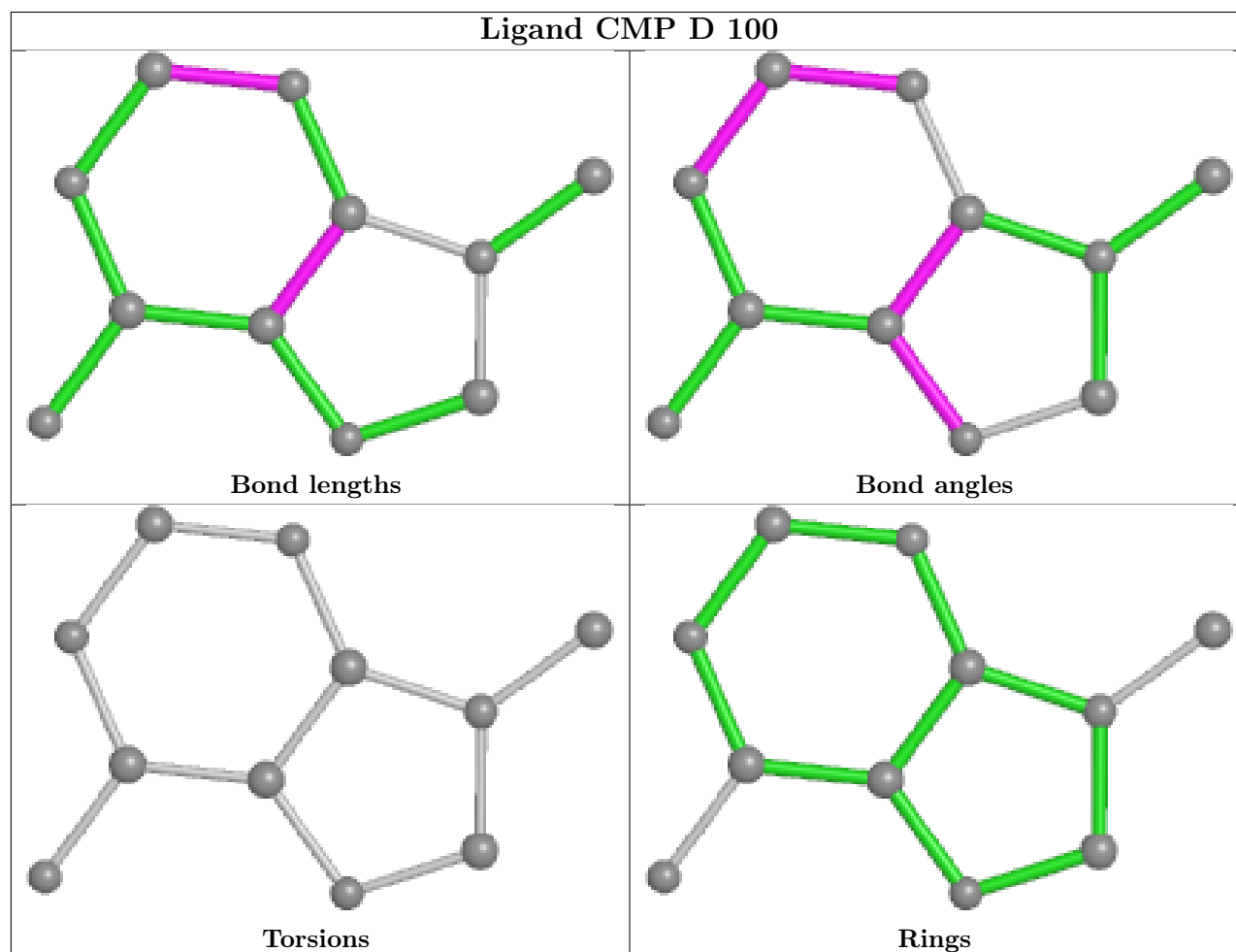
There are no ring outliers.

3 monomers are involved in 4 short contacts:

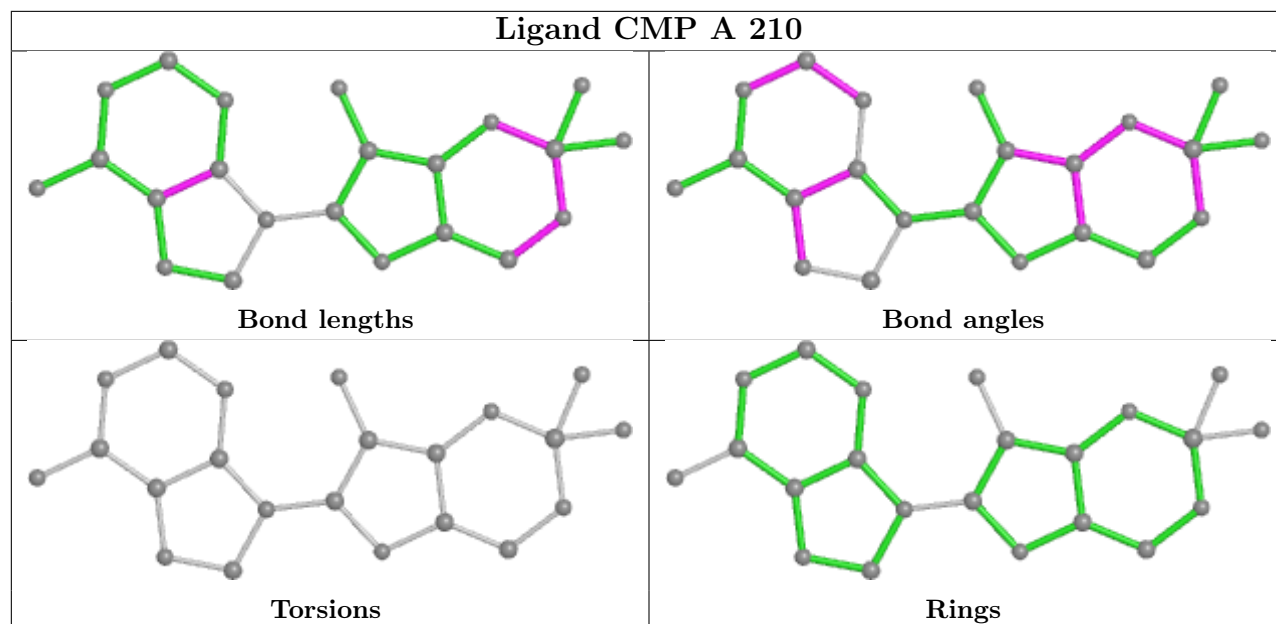
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	D	100	CMP	1	0
6	A	501	PEG	2	0
5	A	210	CMP	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will

also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.







## 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q < 0.9
1	A	203/209 (97%)	-0.06	0	100   100	55, 80, 123, 151	0
2	B	74/84 (88%)	0.25	4 (5%)	25   15	87, 121, 183, 198	0
2	C	73/84 (86%)	1.33	23 (31%)	0   0	96, 187, 269, 343	0
3	D	20/20 (100%)	-0.48	0	100   100	75, 92, 115, 131	0
4	E	24/24 (100%)	-0.64	0	100   100	75, 92, 112, 119	0
All	All	394/421 (93%)	0.20	27 (6%)	16   9	55, 99, 206, 343	0

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	C	321	TRP	4.9
2	C	317	ARG	4.6
2	C	259	ASP	4.5
2	C	318	LEU	4.2
2	C	320	ASN	3.9
2	C	316	MET	3.9
2	C	254	LEU	3.9
2	C	275	ILE	3.8
2	C	276	HIS	3.7
2	C	277	TYR	3.3
2	C	322	PRO	3.1
2	B	322	PRO	3.1
2	C	250	ASP	3.0
2	C	283	GLN	2.9
2	C	251	PRO	2.9
2	C	255	ARG	2.7
2	B	321	TRP	2.6
2	C	271	LYS	2.6
2	B	318	LEU	2.5
2	C	319	GLU	2.4

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Mol	Chain	Res	Type	RSRZ
2	C	269	CYS	2.4
2	C	253	LEU	2.4
2	C	279	GLY	2.3
2	C	278	ILE	2.2
2	B	320	ASN	2.1
2	C	252	ILE	2.0
2	C	310	ARG	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 monosaccharides in this entry.

## 6.4 Ligands [i](#)

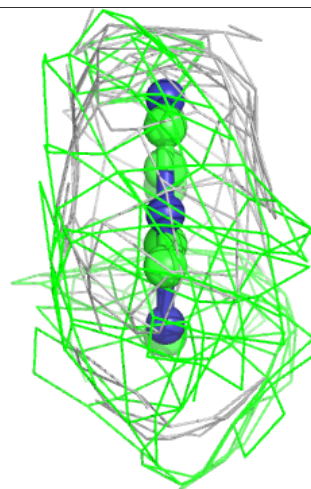
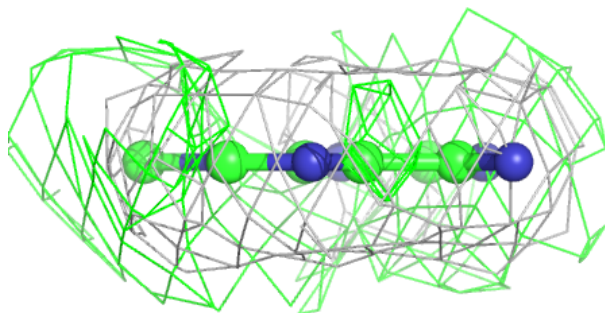
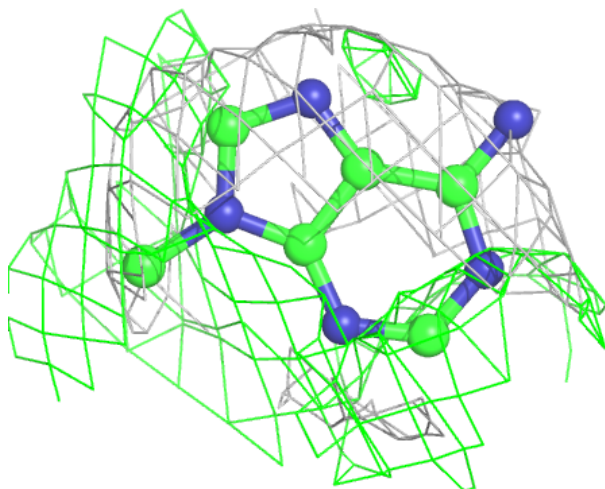
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

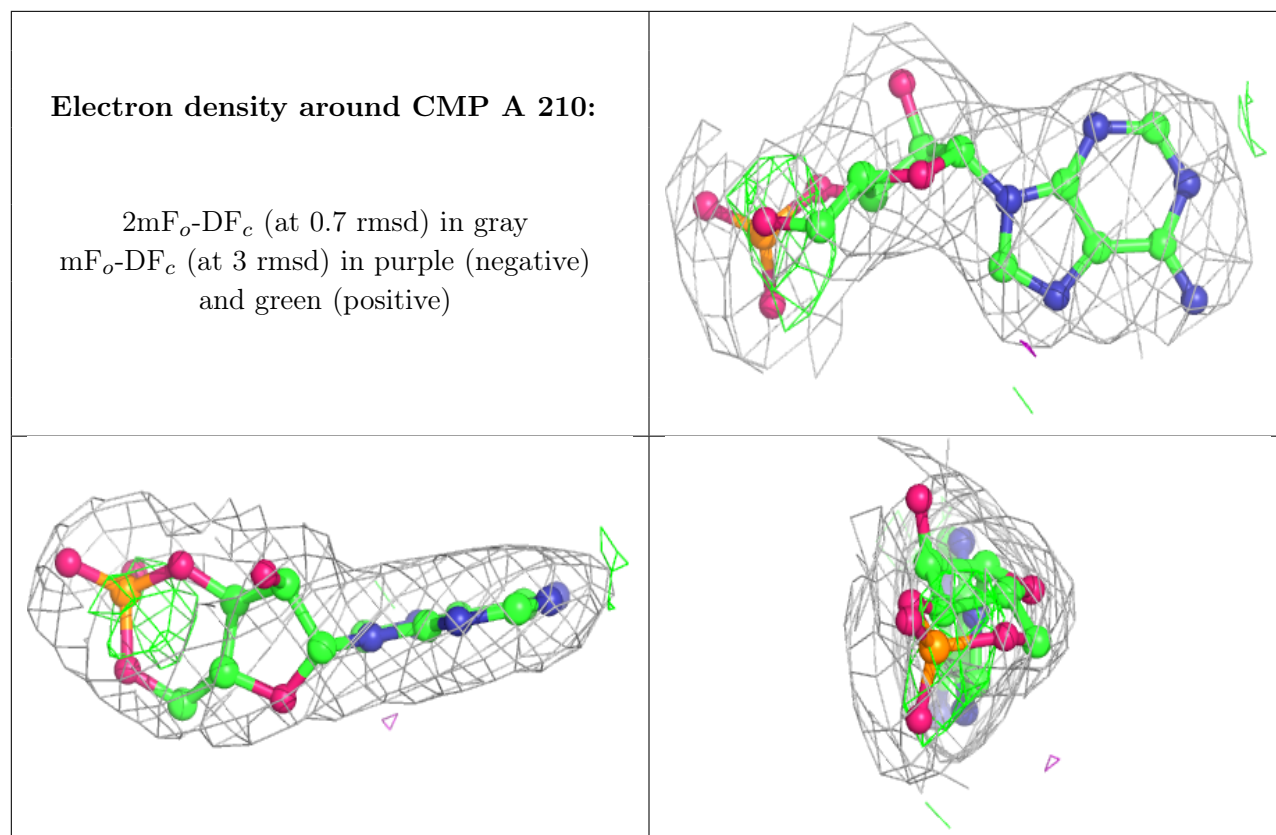
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
6	PEG	A	502	7/7	0.75	0.31	115,125,153,155	0
5	CMP	D	100	11/22	0.86	0.15	102,113,132,132	0
6	PEG	A	501	7/7	0.86	0.18	97,102,107,107	0
5	CMP	A	210	22/22	0.98	0.23	56,66,78,89	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

**Electron density around CMP D 100:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





## 6.5 Other polymers [i](#)

There are no such residues in this entry.