



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 23, 2026 – 04:13 AM UTC

PDB ID : 4CLT / pdb_00004clt
Title : Crystal structure of human soluble Adenylyl Cyclase with adenosine-3', 5'-cyclic-monophosphate and pyrophosphate
Authors : Kleinboelting, S.; Weyand, M.; Steegborn, C.
Deposited on : 2014-01-15
Resolution : 1.95 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 : **FAILED**
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

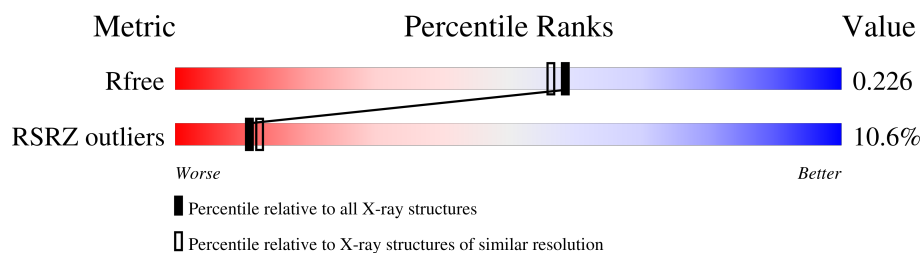
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.95 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	3494 (1.96-1.96)
RSRZ outliers	180081	3495 (1.96-1.96)

MolProbity failed to run properly - the sequence quality summary graphics cannot be shown.

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 3838 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 ADENYLATE CYCLASE TYPE 10.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	462	Total	C	N	O	S	0	3	0
			3673	2380	596	663	34			

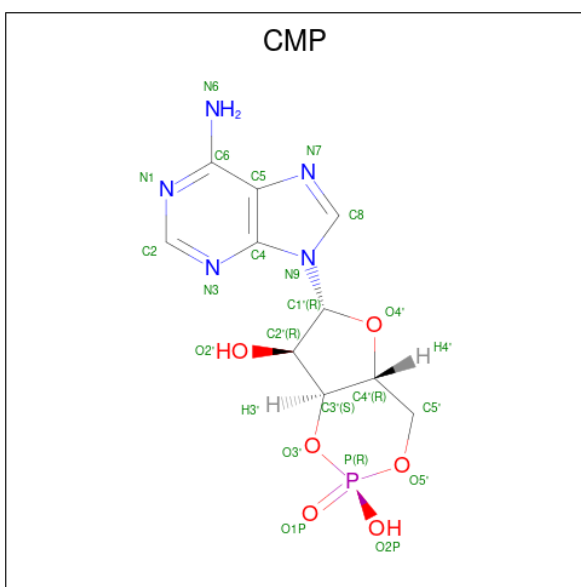
There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	470	HIS	-	expression tag	UNP Q96PN6
A	471	HIS	-	expression tag	UNP Q96PN6
A	472	HIS	-	expression tag	UNP Q96PN6
A	473	HIS	-	expression tag	UNP Q96PN6
A	474	HIS	-	expression tag	UNP Q96PN6
A	475	HIS	-	expression tag	UNP Q96PN6

- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

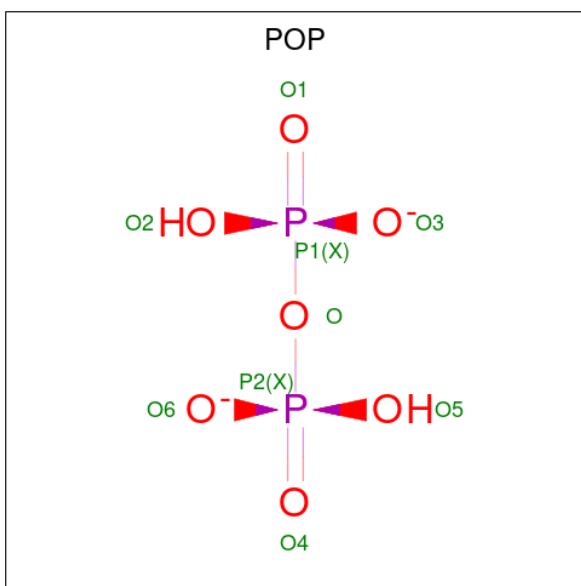
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Mg	0	0
			1	1		

- Molecule 3 is ADENOSINE-3',5'-CYCLIC-MONOPHOSPHATE (CCD ID: CMP) (formula: C₁₀H₁₂N₅O₆P).



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

- Molecule 4 is PYROPHOSPHATE 2- (CCD ID: POP) (formula: $\text{H}_2\text{O}_7\text{P}_2$).

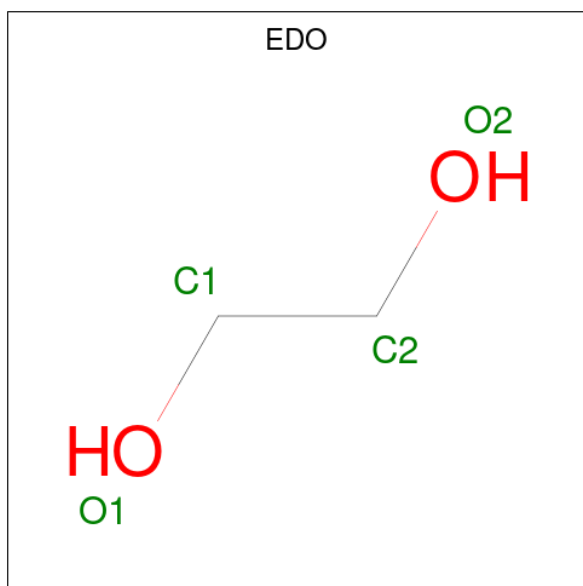


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	P	0	0
			9	7	2		

- Molecule 5 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total Cl 1 1	0	0

- Molecule 6 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	1	Total C O 4 2 2	0	0

- Molecule 7 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	106	Total O 106 106	0	0

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3 Data and refinement statistics

Property	Value	Source
Space group	P 63	Depositor
Cell constants a, b, c, α , β , γ	100.97Å 100.97Å 97.21Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	87.44 – 1.95 87.44 – 1.95	Depositor EDS
% Data completeness (in resolution range)	99.9 (87.44-1.95) 99.9 (87.44-1.95)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.97 (at 1.95Å)	Xtriage
Refinement program	REFMAC 5.8.0049	Depositor
R, R_{free}	0.173 , 0.220 0.183 , 0.226	Depositor DCC
R_{free} test set	2065 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	23.7	Xtriage
Anisotropy	0.230	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 36.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.048 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	3838	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.54% 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.

4 Model quality [i](#)

4.1 Standard geometry [i](#)

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4.2 Too-close contacts [i](#)

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4.3 Torsion angles [i](#)

4.3.1 Protein backbone [i](#)

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4.3.2 Protein sidechains [i](#)

MolProbity failed to run properly - this section is therefore empty.

4.3.3 RNA [i](#)

MolProbity failed to run properly - this section is therefore empty.

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

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	$\# Z > 2$
1	CME	A	253	1	8,9,10	0.78	0	6,9,11	1.85	2 (33%)

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
1	CME	A	253	1	-	0/5/8/10	-

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	253	CME	OH-CZ-CE	-3.71	96.33	110.82
1	A	253	CME	CE-SD-SG	-2.45	92.68	103.46

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

4.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

4.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 2 are monoatomic - leaving 4 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
6	EDO	A	1476	-	3,3,3	0.34	0	2,2,2	0.08	0
3	CMP	A	1472	-	25,25,25	1.96	6 (24%)	37,39,39	2.23	17 (45%)
4	POP	A	1474	2	6,8,8	0.94	0	12,13,13	1.24	2 (16%)
3	CMP	A	1473	-	25,25,25	1.94	7 (28%)	37,39,39	2.37	15 (40%)

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	EDO	A	1476	-	-	1/1/1/1	-
3	CMP	A	1472	-	-	1/4/31/31	0/4/4/4
4	POP	A	1474	2	-	0/6/6/6	-
3	CMP	A	1473	-	-	0/4/31/31	0/4/4/4

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1472	CMP	P-O5'	5.78	1.64	1.57
3	A	1472	CMP	C5-C4	5.19	1.48	1.39
3	A	1473	CMP	C5-C4	4.77	1.47	1.39
3	A	1473	CMP	P-O5'	3.67	1.61	1.57
3	A	1473	CMP	C5-C6	3.18	1.49	1.41
3	A	1473	CMP	P-O3'	3.03	1.62	1.57
3	A	1473	CMP	C8-N7	2.96	1.37	1.31
3	A	1472	CMP	C4-N9	-2.77	1.31	1.37
3	A	1473	CMP	C4-N9	-2.68	1.32	1.37
3	A	1473	CMP	O5'-C5'	-2.56	1.42	1.46
3	A	1472	CMP	C8-N7	2.39	1.36	1.31
3	A	1472	CMP	C5-N7	-2.14	1.35	1.39
3	A	1472	CMP	C5-C6	2.11	1.46	1.41

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1473	CMP	O3'-C3'-C4'	-5.54	106.53	110.71
3	A	1473	CMP	O2P-P-O1P	5.35	125.00	108.56
3	A	1472	CMP	O5'-P-O3'	4.70	111.99	105.70
3	A	1473	CMP	C5-C4-N3	-4.59	120.39	126.72
3	A	1473	CMP	C4-C5-N7	-4.43	105.52	110.58
3	A	1472	CMP	C5-C4-N3	-4.31	120.78	126.72
3	A	1472	CMP	O2P-P-O3'	-4.05	97.61	107.04
3	A	1473	CMP	C6-C5-N7	3.60	139.04	132.09
3	A	1472	CMP	N3-C2-N1	-3.55	123.21	128.58
3	A	1472	CMP	N3-C4-N9	3.43	133.01	127.17
3	A	1472	CMP	C2-N3-C4	3.10	119.41	111.83
3	A	1472	CMP	C4-C5-N7	-3.10	107.04	110.58
3	A	1472	CMP	C4-N9-C8	3.08	108.97	105.74

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1473	CMP	O2P-P-O5'	-2.99	99.86	107.16
3	A	1473	CMP	C2-N3-C4	2.93	118.99	111.83
3	A	1473	CMP	O4'-C1'-N9	-2.89	102.54	108.09
3	A	1473	CMP	C4'-O4'-C1'	-2.82	103.24	109.47
3	A	1472	CMP	O2P-P-O1P	2.71	116.89	108.56
3	A	1472	CMP	C5'-C4'-C3'	2.69	117.62	112.56
3	A	1473	CMP	N3-C4-N9	2.69	131.73	127.17
3	A	1473	CMP	C5-N7-C8	2.68	107.67	103.45
3	A	1472	CMP	C2-N1-C6	2.63	123.05	118.73
3	A	1472	CMP	C6-C5-N7	2.58	137.06	132.09
3	A	1473	CMP	C4-N9-C8	2.57	108.44	105.74
3	A	1472	CMP	O3'-C3'-C2'	-2.50	113.15	115.61
3	A	1473	CMP	N9-C8-N7	-2.44	110.48	113.94
3	A	1473	CMP	O3'-C3'-C2'	2.39	117.94	115.61
3	A	1472	CMP	O5'-P-O1P	-2.22	105.33	110.44
3	A	1472	CMP	C4'-O4'-C1'	-2.19	104.63	109.47
3	A	1472	CMP	C5-N7-C8	2.17	106.87	103.45
4	A	1474	POP	O3-P1-O1	2.14	119.18	110.83
3	A	1472	CMP	N9-C8-N7	-2.10	110.95	113.94
4	A	1474	POP	O-P1-O1	-2.10	100.01	111.04
3	A	1473	CMP	N3-C2-N1	-2.09	125.41	128.58

There are no chirality outliers.

All (2) torsion outliers are listed below:

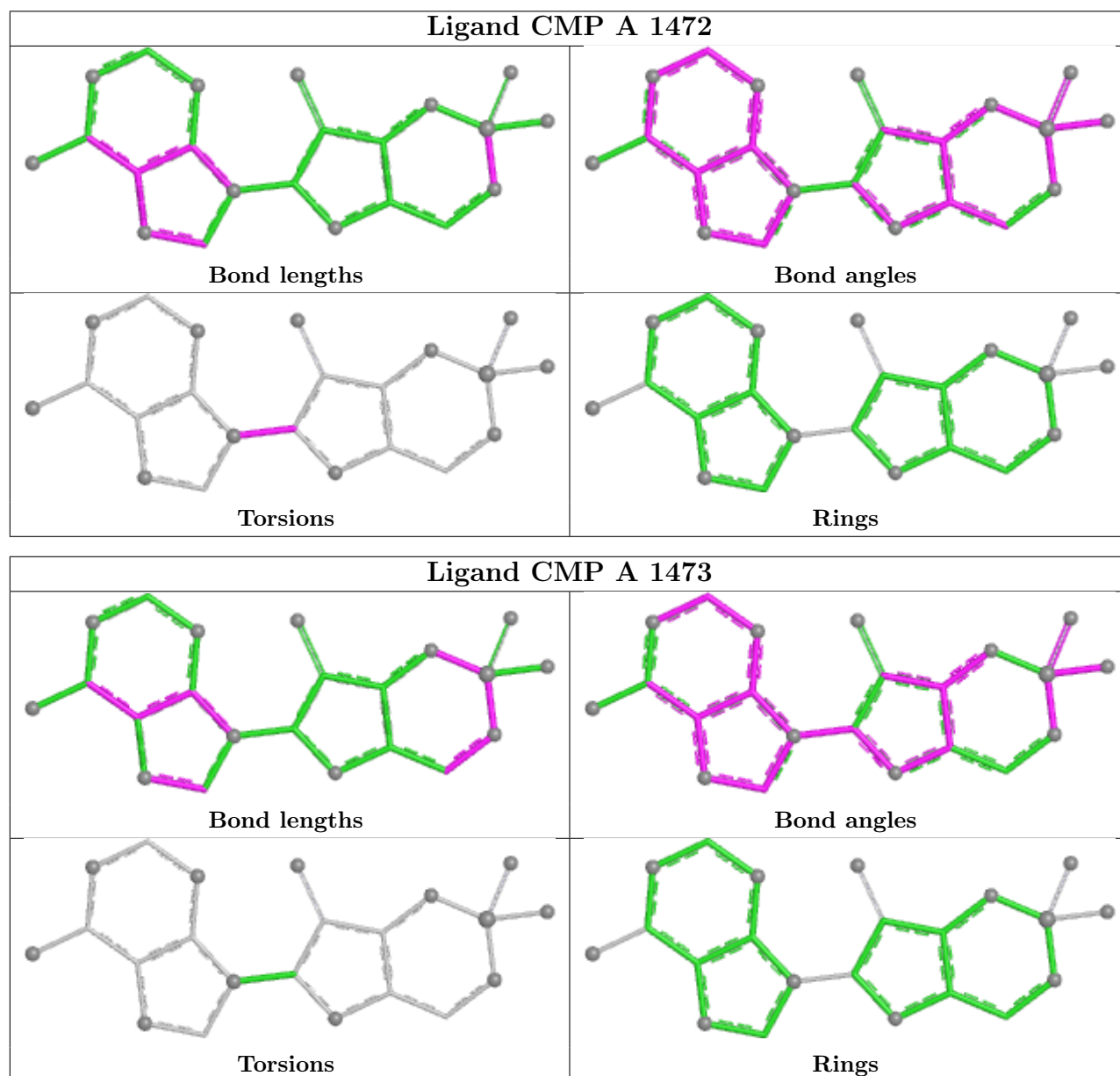
Mol	Chain	Res	Type	Atoms
6	A	1476	EDO	O1-C1-C2-O2
3	A	1472	CMP	C2'-C1'-N9-C8

There are no ring outliers.

No monomer is involved in short contacts.

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.



4.7 Other polymers [i](#)

There are no such residues in this entry.

4.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

5 Fit of model and data

5.1 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	461/475 (97%)	0.29	49 (10%)	11 13	10, 27, 66, 96	3 (0%)

All (49) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	356	PRO	7.0
1	A	138	GLY	6.4
1	A	127	HIS	6.3
1	A	469	VAL	6.2
1	A	355	VAL	5.7
1	A	351	PRO	5.7
1	A	468	LYS	5.6
1	A	139	LEU	5.4
1	A	130	PHE	5.4
1	A	8	PHE	5.2
1	A	466	THR	4.6
1	A	350	PHE	4.1
1	A	352	GLY	4.1
1	A	285	TYR	4.0
1	A	454	ALA	4.0
1	A	449	VAL	3.5
1	A	51	PHE	3.5
1	A	353	GLU	3.4
1	A	185	ASN	3.4
1	A	184	MET	3.3
1	A	451	LYS	3.3
1	A	283	GLN	3.2
1	A	349	GLY	3.1
1	A	467	GLU	2.9
1	A	141	ILE	2.7
1	A	354	LYS	2.7
1	A	456	SER	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	447	LYS	2.7
1	A	377	HIS	2.7
1	A	465	ARG	2.6
1	A	223	ASN	2.6
1	A	53	ALA	2.6
1	A	140	ASP	2.6
1	A	339	ASP	2.5
1	A	340	LYS	2.5
1	A	452	GLY	2.4
1	A	5	LYS	2.4
1	A	338	PHE	2.3
1	A	50	GLY	2.3
1	A	49	SER	2.3
1	A	302	GLU	2.2
1	A	453	VAL	2.2
1	A	128	GLY	2.2
1	A	48	ILE	2.1
1	A	7	GLU	2.1
1	A	9	GLN	2.1
1	A	142	ARG	2.1
1	A	457	GLY	2.1
1	A	28	HIS	2.0

5.2 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($\text{\AA}^2$)	Q<0.9
1	CME	A	253	10/11	0.97	0.09	16,21,36,37	0

5.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.4 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,

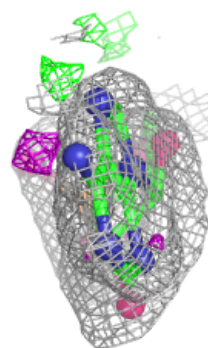
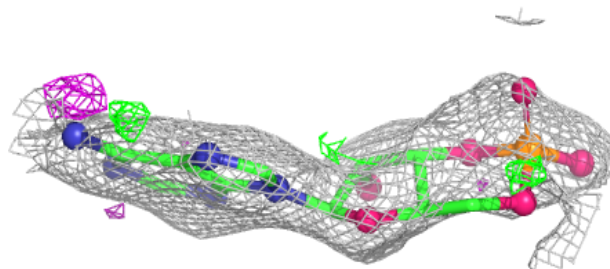
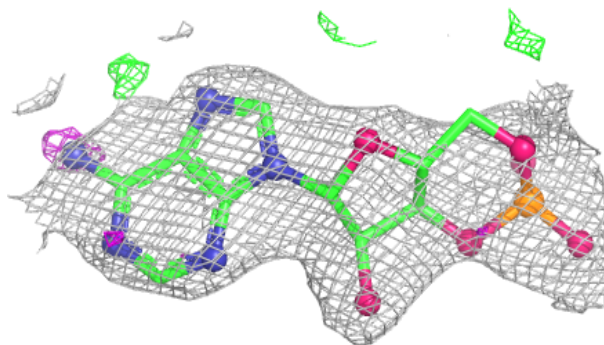
median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	CMP	A	1473	22/22	0.81	0.15	43,61,91,96	0
6	EDO	A	1476	4/4	0.86	0.15	39,42,50,57	0
4	POP	A	1474	9/9	0.87	0.09	48,56,62,64	0
3	CMP	A	1472	22/22	0.94	0.08	34,40,46,49	0
2	MG	A	1471	1/1	0.94	0.07	40,40,40,40	0
5	CL	A	1475	1/1	0.99	0.03	23,23,23,23	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.

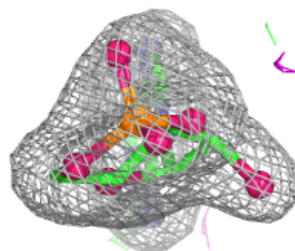
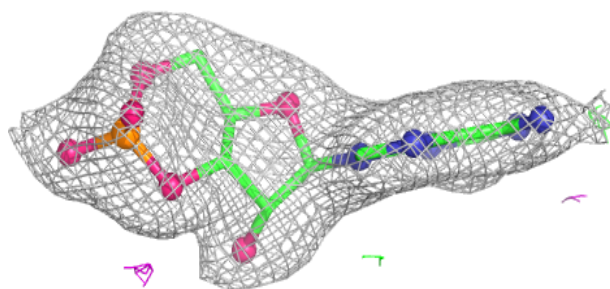
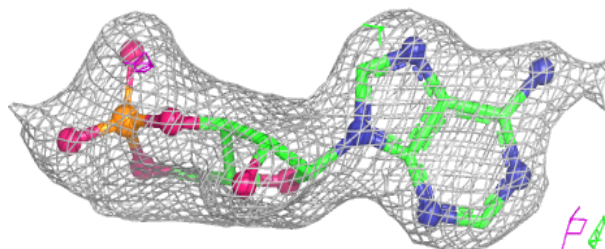
Electron density around CMP A 1473:

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



Electron density around CMP A 1472:

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



5.5 Other polymers [i](#)

There are no such residues in this entry.