



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 11, 2021 – 03:22 PM EDT

PDB ID : 2OUR
Title : crystal structure of PDE10A2 mutant D674A in complex with cAMP
Authors : Wang, H.C.; Liu, Y.D.; Hou, J.; Zheng, M.Y.; Robinson, H.
Deposited on : 2007-02-12
Resolution : 1.45 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 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 : 2.23.2
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) : 2.23.2

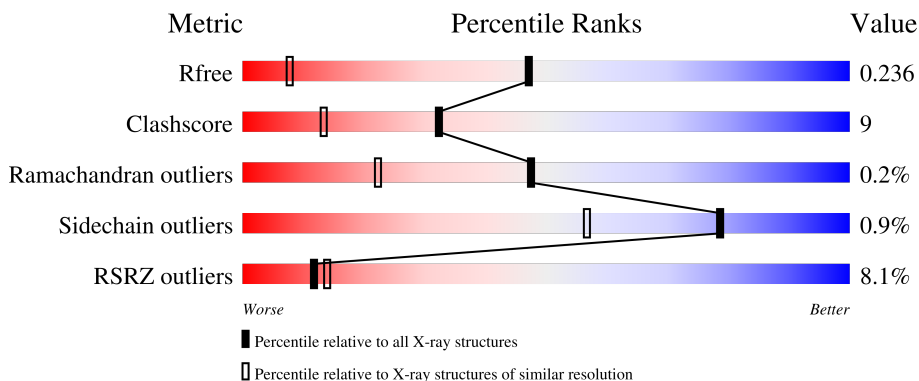
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.45 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	1156 (1.46-1.46)
Clashscore	141614	1202 (1.46-1.46)
Ramachandran outliers	138981	1178 (1.46-1.46)
Sidechain outliers	138945	1178 (1.46-1.46)
RSRZ outliers	127900	1139 (1.46-1.46)

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	331	 4% 80% 18% ••
1	B	331	 12% 78% 16% • 5%

2 Entry composition i

There are 4 unique types of molecules in this entry. The entry contains 5631 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 cAMP and cAMP-inhibited cGMP 3',5'-cyclic phosphodiesterase 10A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	324	2626	1677	447	477	25	0	0	0
1	B	315	2548	1624	434	465	25	0	0	0

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	446	SER	-	cloning artifact	UNP Q9Y233
A	447	HIS	-	cloning artifact	UNP Q9Y233
A	448	MET	-	cloning artifact	UNP Q9Y233
A	674	ALA	ASP	engineered mutation	UNP Q9Y233
B	446	SER	-	cloning artifact	UNP Q9Y233
B	447	HIS	-	cloning artifact	UNP Q9Y233
B	448	MET	-	cloning artifact	UNP Q9Y233
B	674	ALA	ASP	engineered mutation	UNP Q9Y233

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

- Molecule 3 is ADENOSINE-3',5'-CYCLIC-MONOPHOSPHATE (three-letter code: CMP) (formula: C₁₀H₁₂N₅O₆P).



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

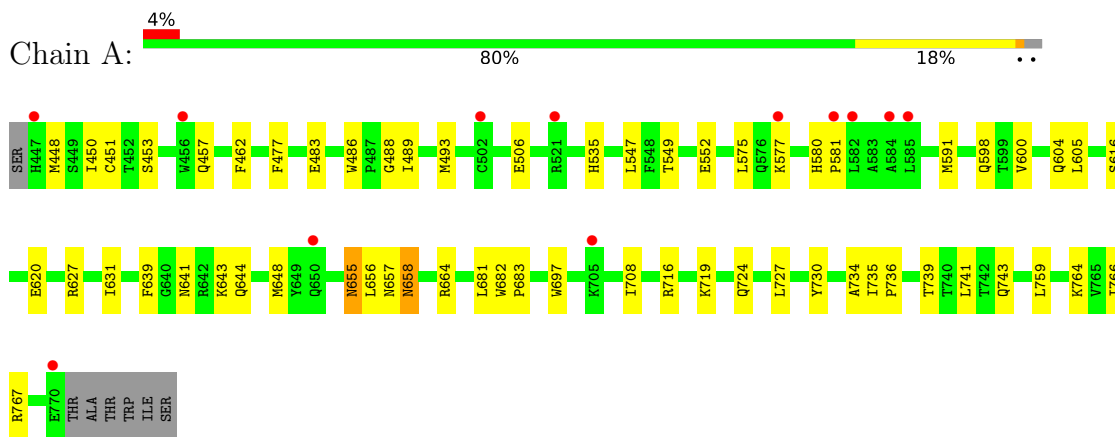
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	238	Total	O	0	0
			238	238		
4	B	195	Total	O	0	0
			195	195		

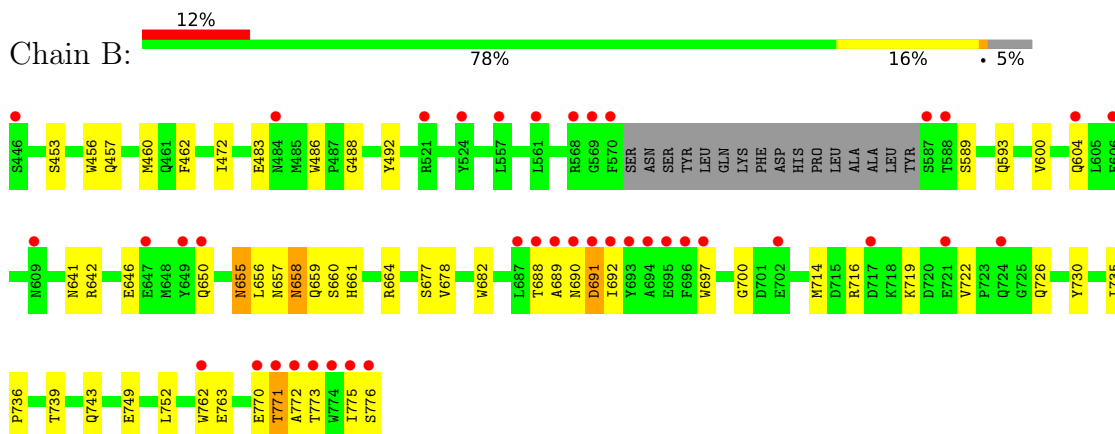
3 Residue-property plots [i](#)

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: cAMP and cAMP-inhibited cGMP 3',5'-cyclic phosphodiesterase 10A



- Molecule 1: cAMP and cAMP-inhibited cGMP 3',5'-cyclic phosphodiesterase 10A



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	49.30Å 82.28Å 154.00Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 1.45 29.22 – 1.45	Depositor EDS
% Data completeness (in resolution range)	(Not available) (30.00-1.45) 86.2 (29.22-1.45)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.96 (at 1.45Å)	Xtrriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.218 , 0.236 0.219 , 0.236	Depositor DCC
R_{free} test set	10158 reflections (10.02%)	wwPDB-VP
Wilson B-factor (Å ²)	14.4	Xtrriage
Anisotropy	0.691	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 36.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5631	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.31% 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: MG, 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	0/2691	0.53	0/3643
1	B	0.29	0/2609	0.53	0/3530
All	All	0.29	0/5300	0.53	0/7173

There are no bond length outliers.

There are no bond angle outliers.

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	2626	0	2589	43	0
1	B	2548	0	2513	49	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	22	0	11	1	0
4	A	238	0	0	1	0
4	B	195	0	0	1	0
All	All	5631	0	5113	91	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 (91) 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)
3:A:778:CMP:H2	3:A:778:CMP:C2	0.97	1.48
1:B:655:ASN:H	1:B:661:HIS:HD2	1.20	0.85
1:A:641:ASN:HD22	1:A:664:ARG:HH11	1.28	0.79
1:B:688:THR:HG22	1:B:691:ASP:OD2	1.84	0.76
1:B:655:ASN:H	1:B:661:HIS:CD2	2.05	0.74
1:B:453:SER:O	1:B:457:GLN:HG3	1.90	0.70
1:A:658:ASN:HD22	1:A:658:ASN:C	1.97	0.67
1:B:641:ASN:HD22	1:B:664:ARG:HH11	1.44	0.66
1:A:716:ARG:O	1:A:719:LYS:HG3	1.97	0.65
1:B:658:ASN:C	1:B:658:ASN:HD22	2.01	0.64
1:B:600:VAL:O	1:B:604:GLN:HG3	1.97	0.63
1:B:682:TRP:HE1	1:B:773:THR:HG22	1.64	0.63
1:B:655:ASN:ND2	1:B:657:ASN:H	1.98	0.61
1:B:656:LEU:N	1:B:656:LEU:HD22	2.15	0.61
1:A:739:THR:O	1:A:743:GLN:HG3	2.01	0.60
1:B:456:TRP:HZ3	1:B:460:MET:SD	2.26	0.58
1:A:450:ILE:HG13	1:A:605:LEU:HD13	1.85	0.58
1:B:770:GLU:HB3	1:B:773:THR:HB	1.84	0.58
1:A:453:SER:O	1:A:457:GLN:HG3	2.04	0.58
1:B:692:ILE:HG13	4:B:269:HOH:O	2.03	0.57
1:A:483:GLU:HA	1:A:486:TRP:CE2	2.40	0.56
1:A:506:GLU:CD	1:A:506:GLU:H	2.08	0.56
1:A:655:ASN:ND2	1:A:657:ASN:H	2.04	0.56
1:B:697:TRP:HB3	1:B:716:ARG:HE	1.71	0.56
1:B:483:GLU:HA	1:B:486:TRP:CE2	2.41	0.55
1:B:456:TRP:CZ3	1:B:460:MET:SD	3.00	0.55
1:B:472:ILE:HD13	1:B:492:TYR:CE2	2.42	0.55
1:B:658:ASN:HD22	1:B:659:GLN:N	2.06	0.54
1:A:697:TRP:CZ2	1:A:719:LYS:HG2	2.42	0.54
1:A:681:LEU:HD11	1:B:763:GLU:HG3	1.90	0.53
1:A:727:LEU:HD12	1:A:759:LEU:HD11	1.91	0.53
1:B:646:GLU:HG3	1:B:650:GLN:HE21	1.74	0.52
1:B:689:ALA:HB2	1:B:762:TRP:HH2	1.75	0.52
1:B:739:THR:O	1:B:743:GLN:HG3	2.09	0.51
1:B:682:TRP:NE1	1:B:773:THR:HG22	2.25	0.50
1:A:735:ILE:HB	1:A:736:PRO:HD3	1.94	0.50
1:B:483:GLU:HG3	1:B:486:TRP:CZ3	2.46	0.50
1:A:462:PHE:CD2	1:A:488:GLY:HA3	2.47	0.49
1:B:589:SER:O	1:B:593:GLN:HG3	2.13	0.49
1:B:678:VAL:HB	1:B:730:TYR:CZ	2.48	0.49

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:577:LYS:HE2	1:A:708:ILE:HG21	1.95	0.49
1:A:448:MET:O	1:A:450:ILE:HD12	2.13	0.49
1:B:656:LEU:HD22	1:B:656:LEU:H	1.78	0.48
1:A:547:LEU:HD11	1:A:656:LEU:HD13	1.94	0.48
1:A:655:ASN:HD21	1:A:657:ASN:HB2	1.79	0.48
1:A:450:ILE:HD12	1:A:450:ILE:N	2.29	0.48
1:A:639:PHE:O	1:A:643:LYS:HG3	2.14	0.48
1:B:472:ILE:HD13	1:B:492:TYR:CD2	2.49	0.47
1:A:730:TYR:HA	1:A:734:ALA:HB3	1.95	0.47
1:B:642:ARG:HD3	1:B:642:ARG:C	2.34	0.47
1:A:580:HIS:CG	1:A:581:PRO:HD2	2.49	0.47
1:B:697:TRP:CD2	1:B:716:ARG:HA	2.50	0.47
1:B:697:TRP:CE3	1:B:716:ARG:HG3	2.50	0.46
1:B:655:ASN:C	1:B:655:ASN:HD22	2.18	0.46
1:A:764:LYS:HG2	1:A:767:ARG:NH2	2.29	0.46
1:A:644:GLN:O	1:A:648:MET:HG3	2.16	0.46
1:B:719:LYS:HA	1:B:722:VAL:HG23	1.97	0.45
1:A:575:LEU:HD11	1:A:591:MET:SD	2.57	0.44
1:B:677:SER:HB2	1:B:688:THR:HG21	1.98	0.43
1:A:580:HIS:ND1	1:A:581:PRO:HD2	2.33	0.43
1:A:600:VAL:O	1:A:604:GLN:HG3	2.19	0.43
1:A:724:GLN:NE2	1:A:766:ILE:HD13	2.34	0.43
1:A:682:TRP:HB3	1:A:683:PRO:HD3	2.01	0.43
1:A:616:SER:O	1:A:620:GLU:HG3	2.19	0.43
1:A:741:LEU:HD23	4:A:278:HOH:O	2.19	0.42
1:A:627:ARG:O	1:A:631:ILE:HG12	2.19	0.42
1:B:775:ILE:O	1:B:776:SER:C	2.57	0.42
1:B:462:PHE:CD2	1:B:488:GLY:HA3	2.54	0.42
1:B:722:VAL:O	1:B:726:GLN:HG2	2.19	0.42
1:B:646:GLU:O	1:B:650:GLN:HG3	2.19	0.42
1:B:735:ILE:HB	1:B:736:PRO:HD3	2.02	0.42
1:B:472:ILE:HG21	1:B:492:TYR:CE2	2.54	0.42
1:A:489:ILE:O	1:A:493:MET:HG3	2.19	0.42
1:A:658:ASN:C	1:A:658:ASN:ND2	2.70	0.42
1:A:681:LEU:CD1	1:B:763:GLU:HG3	2.50	0.42
1:A:727:LEU:CD1	1:A:759:LEU:HD11	2.49	0.42
1:B:700:GLY:HA3	1:B:714:MET:O	2.20	0.41
1:B:771:THR:HG22	1:B:772:ALA:N	2.35	0.41
1:A:549:THR:OG1	1:A:552:GLU:HG3	2.19	0.41
1:A:697:TRP:CH2	1:A:719:LYS:HG2	2.55	0.41
1:A:580:HIS:CE1	1:A:581:PRO:HD2	2.55	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:658:ASN:ND2	1:B:660:SER:H	2.18	0.41
1:A:450:ILE:HG22	1:A:451:CYS:O	2.21	0.41
1:A:477:PHE:HB3	1:A:535:HIS:CE1	2.56	0.41
1:A:641:ASN:HD21	1:A:664:ARG:CB	2.34	0.41
1:B:749:GLU:OE1	1:B:752:LEU:HB3	2.21	0.41
1:B:775:ILE:O	1:B:775:ILE:HG13	2.21	0.41
1:B:688:THR:HG22	1:B:691:ASP:CG	2.39	0.40
1:B:690:ASN:HB3	1:B:697:TRP:CZ3	2.56	0.40
1:A:598:GLN:HA	1:A:598:GLN:NE2	2.36	0.40
1:B:453:SER:HA	1:B:456:TRP:NE1	2.36	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	322/331 (97%)	312 (97%)	10 (3%)	0	100	100
1	B	311/331 (94%)	301 (97%)	9 (3%)	1 (0%)	41	18
All	All	633/662 (96%)	613 (97%)	19 (3%)	1 (0%)	47	22

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	771	THR

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	292/298 (98%)	290 (99%)	2 (1%)	84	65
1	B	284/298 (95%)	281 (99%)	3 (1%)	73	48
All	All	576/596 (97%)	571 (99%)	5 (1%)	78	57

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	655	ASN
1	A	658	ASN
1	B	655	ASN
1	B	658	ASN
1	B	691	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	598	GLN
1	A	641	ASN
1	A	655	ASN
1	A	658	ASN
1	A	724	GLN
1	B	593	GLN
1	B	641	ASN
1	B	650	GLN
1	B	655	ASN
1	B	658	ASN
1	B	661	HIS
1	B	724	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 3 ligands modelled in this entry, 2 are monoatomic - leaving 1 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
3	CMP	A	778	-	22,25,25	1.12	2 (9%)	24,39,39	2.05	4 (16%)

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
3	CMP	A	778	-	-	0/0/31/31	0/4/4/4

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	778	CMP	P-O5'	2.80	1.60	1.57
3	A	778	CMP	O3'-C3'	-2.04	1.41	1.44

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	778	CMP	O3'-C3'-C4'	-7.52	105.03	110.71
3	A	778	CMP	O3'-C3'-C2'	3.94	119.47	115.61
3	A	778	CMP	O2P-P-O1P	3.55	119.84	108.73
3	A	778	CMP	C5-C6-N6	2.19	123.68	120.35

There are no chirality outliers.

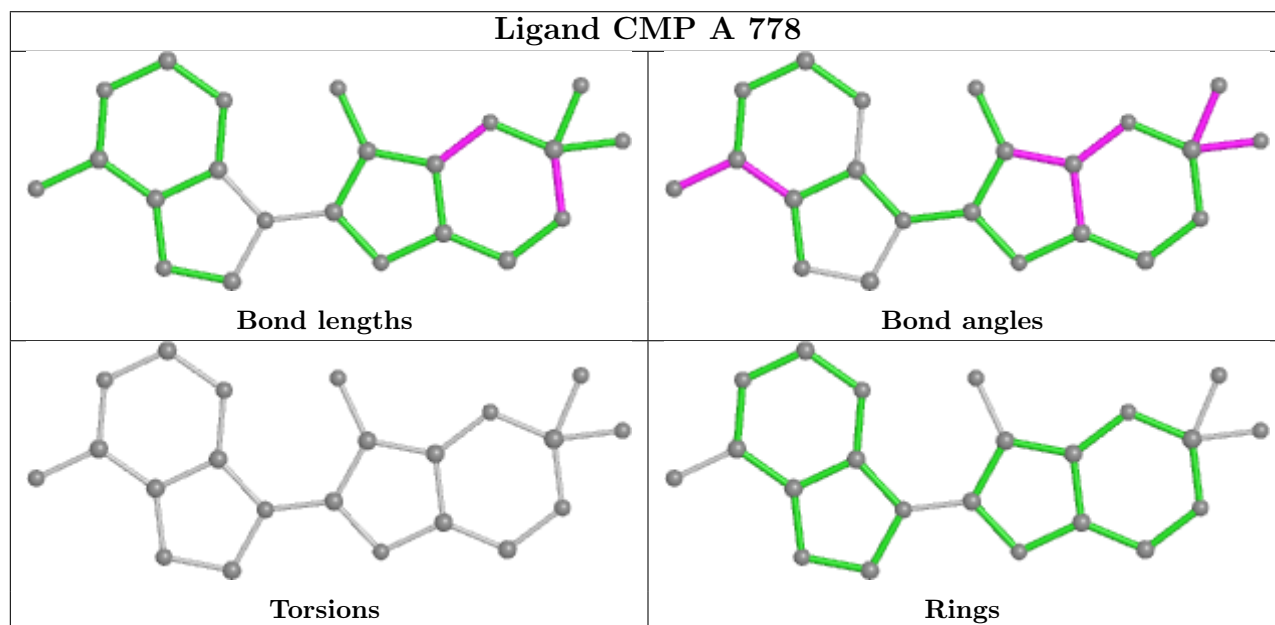
There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	778	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, 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	324/331 (97%)	0.42	12 (3%) 41 44	11, 17, 27, 40	0
1	B	315/331 (95%)	0.87	40 (12%) 3 4	12, 19, 35, 54	0
All	All	639/662 (96%)	0.64	52 (8%) 12 14	11, 18, 31, 54	0

All (52) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	771	THR	12.0
1	B	570	PHE	10.7
1	B	776	SER	9.6
1	B	772	ALA	7.9
1	B	773	THR	7.7
1	B	775	ILE	7.4
1	B	524	TYR	6.4
1	A	447	HIS	6.0
1	B	688	THR	5.8
1	B	770	GLU	5.7
1	B	446	SER	5.2
1	A	585	LEU	5.1
1	B	687	LEU	5.1
1	B	692	ILE	4.9
1	B	690	ASN	4.7
1	B	650	GLN	4.3
1	A	584	ALA	4.1
1	A	577	LYS	3.6
1	B	774	TRP	3.6
1	B	691	ASP	3.5
1	B	569	GLY	3.5
1	B	587	SER	3.5
1	A	650	GLN	3.4
1	B	695	GLU	3.3

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	581	PRO	3.2
1	A	770	GLU	3.2
1	B	521	ARG	3.1
1	B	696	PHE	3.1
1	A	521	ARG	3.0
1	B	588	THR	3.0
1	B	649	TYR	2.9
1	B	762	TRP	2.8
1	B	717	ASP	2.8
1	B	647	GLU	2.8
1	B	484	ASN	2.7
1	B	693	TYR	2.7
1	B	689	ALA	2.7
1	B	568	ARG	2.6
1	B	697	TRP	2.6
1	A	456	TRP	2.4
1	B	557	LEU	2.4
1	A	582	LEU	2.3
1	B	604	GLN	2.3
1	A	705	LYS	2.2
1	B	702	GLU	2.2
1	B	694	ALA	2.2
1	B	606	GLU	2.2
1	B	721	GLU	2.2
1	B	609	ASN	2.2
1	B	561	LEU	2.1
1	B	724	GLN	2.1
1	A	502	CYS	2.1

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.

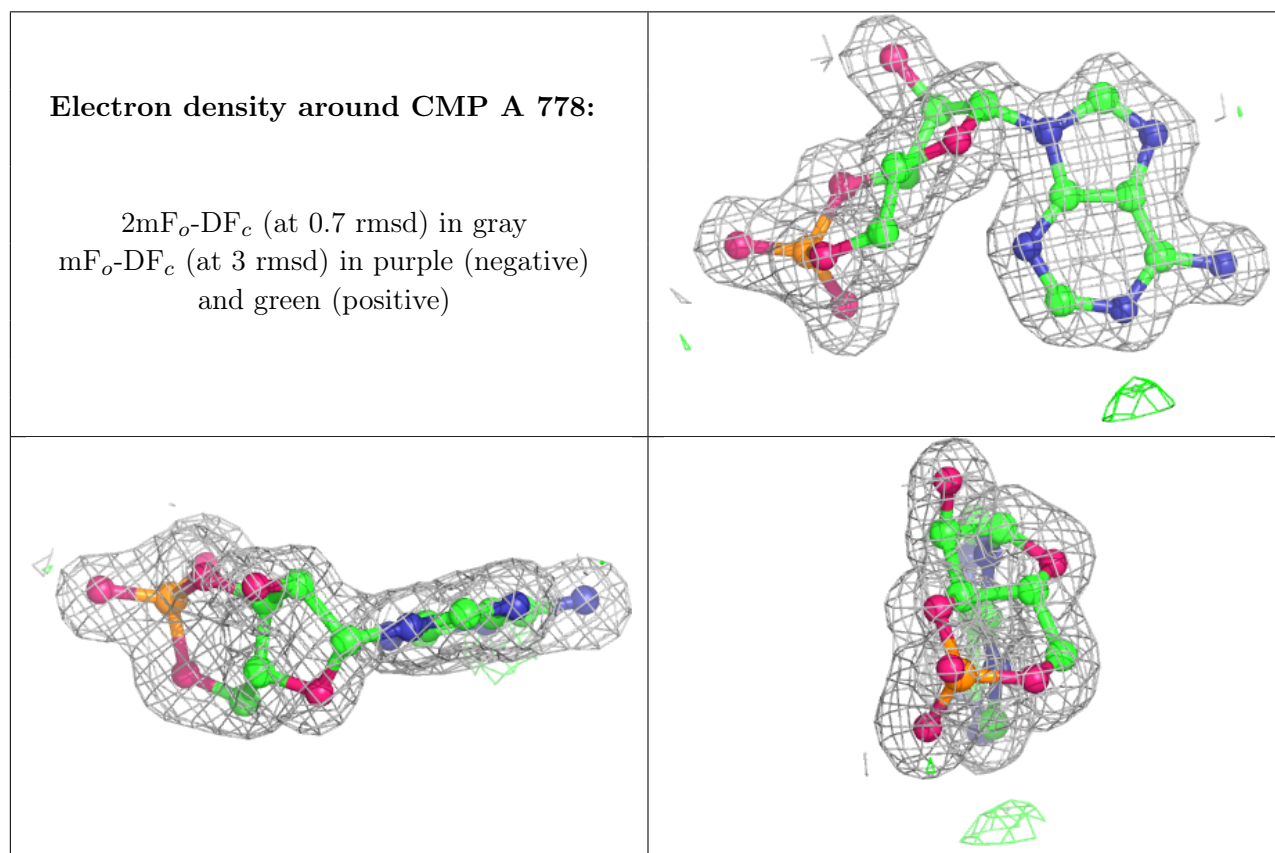
6.4 Ligands [i](#)

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, 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
3	CMP	A	778	22/22	0.95	0.08	18,19,19,19	0
2	MG	B	777	1/1	0.98	0.05	17,17,17,17	0
2	MG	A	777	1/1	0.99	0.12	13,13,13,13	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.



6.5 Other polymers [i](#)

There are no such residues in this entry.