



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 06:18 PM UTC

PDB ID : 4UKD / pdb_00004ukd
Title : UMP/CMP KINASE FROM SLIME MOLD COMPLEXED WITH ADP,
UDP, BERYLLIUM FLUORIDE
Authors : Schlichting, I.; Reinstein, J.
Deposited on : 1997-05-20
Resolution : 2.00 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
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

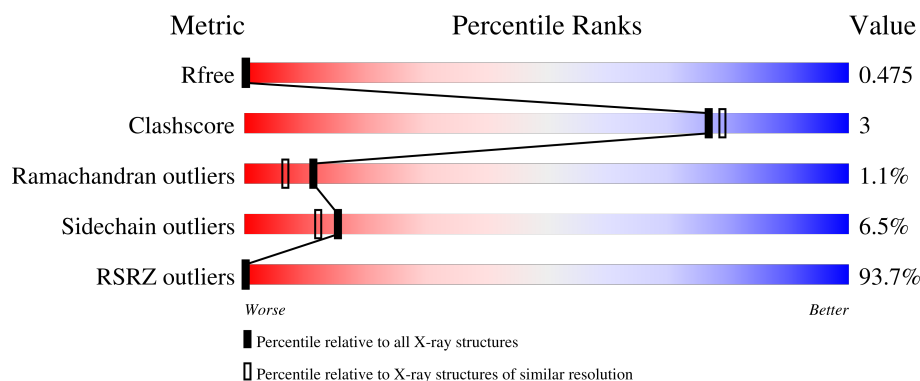
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	194	92% 80%15%...

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 1679 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.

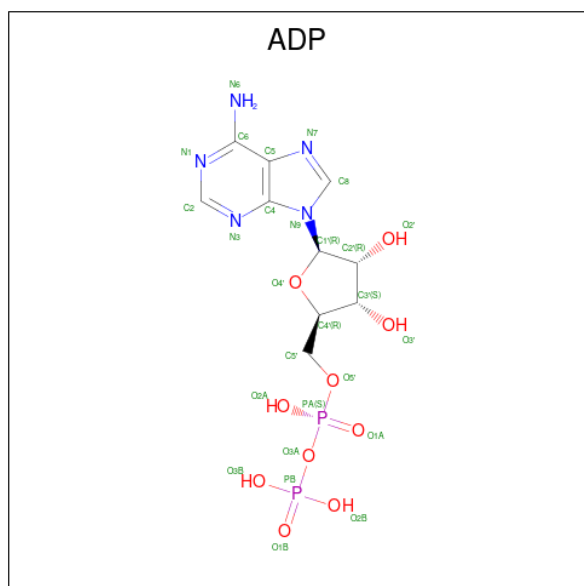
- Molecule 1 is a protein called URIDYLMONOPHOSPHATE/CYTIDYLMONOPHOSPHATE KINASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	191	Total	C	N	O	S	0	0	0
			1516	952	264	293	7			

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

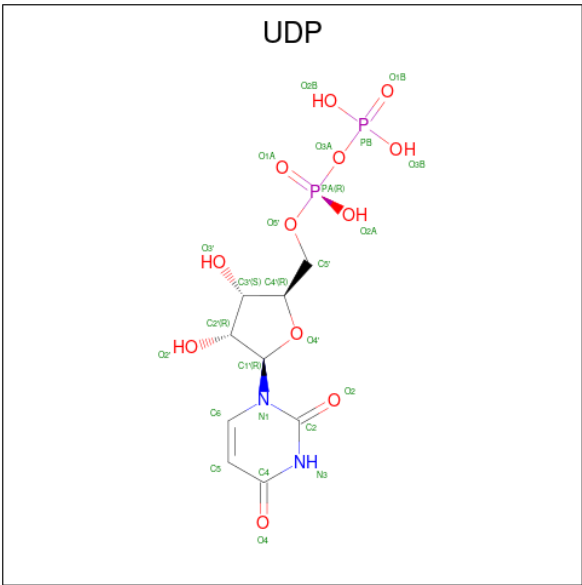
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Mg	0	0
			1	1		

- Molecule 3 is ADENOSINE-5'-DIPHOSPHATE (CCD ID: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$).



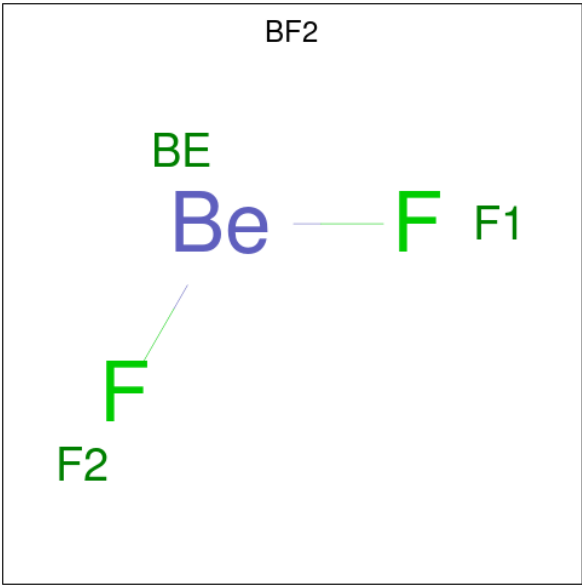
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			27	10	5	10	2		

- Molecule 4 is URIDINE-5'-DIPHOSPHATE (CCD ID: UDP) (formula: $C_9H_{14}N_2O_{12}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			25	9	2	12	2		

- Molecule 5 is BERYLLIUM DIFLUORIDE (CCD ID: BF2) (formula: BeF₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	Be	F	0	0
			3	1	2		

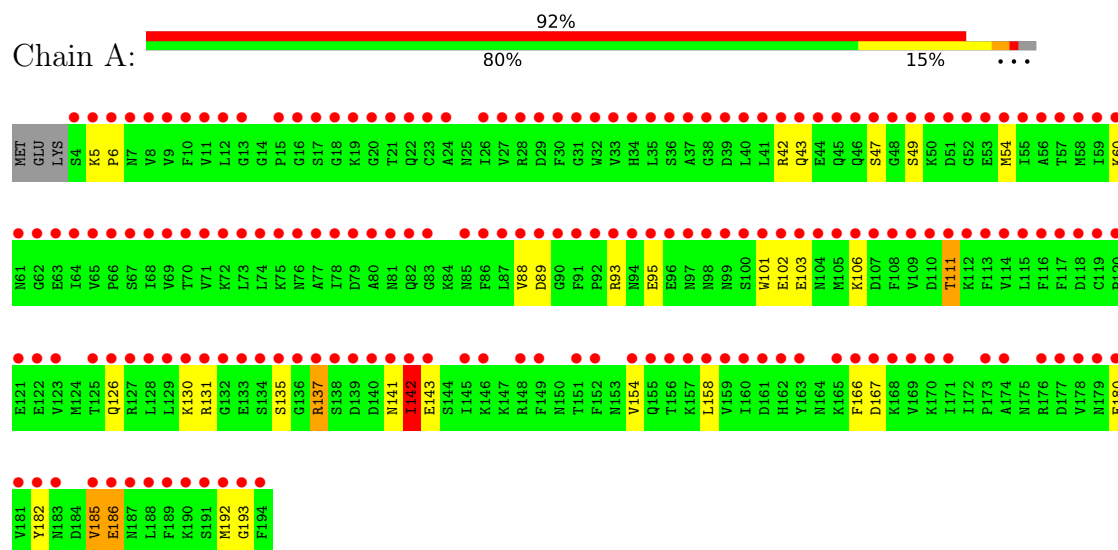
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	107	Total 107	O 107	0	0

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: URIDYLMONOPHOSPHATE/CYTIDYLMONOPHOSPHATE KINASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	78.10Å 78.10Å 101.90Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.00 – 2.00 39.00 – 2.00	Depositor EDS
% Data completeness (in resolution range)	95.0 (39.00-2.00) 86.5 (39.00-2.00)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.40 (at 1.98Å)	Xtriage
Refinement program	X-PLOR 3.851	Depositor
R, R_{free}	0.194 , 0.212 0.467 , 0.475	Depositor DCC
R_{free} test set	913 reflections (4.76%)	wwPDB-VP
Wilson B-factor (Å ²)	24.2	Xtriage
Anisotropy	0.080	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 30.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.67	EDS
Total number of atoms	1679	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.23% 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: UDP, ADP, MG, BF2

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	0.88	1/1542 (0.1%)	0.97	8/2073 (0.4%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	141	ASN	C-N	23.27	1.60	1.33

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	141	ASN	CA-C-N	-8.28	107.06	120.64
1	A	141	ASN	C-N-CA	-8.28	107.06	120.64
1	A	142	ILE	O-C-N	-6.72	112.84	122.12
1	A	137	ARG	NE-CZ-NH2	5.65	124.29	119.20
1	A	143	GLU	N-CA-C	5.62	117.08	111.07
1	A	185	VAL	N-CA-C	-5.45	101.78	109.63
1	A	42	ARG	NE-CZ-NH2	5.17	123.85	119.20
1	A	166	PHE	CA-CB-CG	5.13	118.93	113.80

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	142	ILE	Mainchain

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	1516	0	1493	9	6
2	A	1	0	0	0	0
3	A	27	0	12	0	0
4	A	25	0	11	0	0
5	A	3	0	0	0	0
6	A	107	0	0	1	3
All	All	1679	0	1516	9	6

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 3.

All (9) 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:182:TYR:O	1:A:185:VAL:O	2.25	0.55
1:A:101:TRP:HE1	1:A:111:THR:HG21	1.76	0.50
1:A:131:ARG:HG2	1:A:135:SER:HB3	1.95	0.48
1:A:185:VAL:O	1:A:186:GLU:CB	2.62	0.48
1:A:102:GLU:O	1:A:106:LYS:HB2	2.16	0.46
1:A:182:TYR:O	1:A:182:TYR:CG	2.71	0.44
1:A:5:LYS:HA	1:A:6:PRO:HD3	1.90	0.42
1:A:93:ARG:HG3	6:A:199:HOH:O	2.20	0.42
1:A:54:MET:HE2	1:A:54:MET:HB3	1.86	0.42

All (6) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:47:SER:OG	1:A:95:GLU:OE1[3_655]	1.82	0.38
1:A:193:GLY:O	6:A:532:HOH:O[6_564]	1.89	0.31

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:154:VAL:CG1	6:A:207:HOH:O[4_464]	1.95	0.25
1:A:47:SER:CB	1:A:95:GLU:OE1[3_655]	1.97	0.23
1:A:126:GLN:NE2	1:A:192:MET:O[6_464]	2.02	0.18
1:A:43:GLN:NE2	6:A:577:HOH:O[3_655]	2.10	0.10

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	189/194 (97%)	182 (96%)	5 (3%)	2 (1%)	11 7

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	49	SER
1	A	186	GLU

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	170/173 (98%)	159 (94%)	11 (6%)	15 12

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	60	LYS
1	A	88	VAL
1	A	89	ASP
1	A	103	GLU
1	A	111	THR
1	A	130	LYS
1	A	137	ARG
1	A	142	ILE
1	A	158	LEU
1	A	167	ASP
1	A	180	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	22	GLN
1	A	43	GLN
1	A	46	GLN
1	A	97	ASN
1	A	98	ASN
1	A	99	ASN
1	A	155	GLN
1	A	164	ASN
1	A	175	ASN

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 1 is monoatomic - leaving 3 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	$\# Z > 2$
4	UDP	A	196	5	25,26,26	2.42	10 (40%)	38,40,40	2.73	8 (21%)
5	BF2	A	501	3,2,4	0,2,2	-	-	-	-	-
3	ADP	A	195	2,5	28,29,29	1.40	5 (17%)	43,45,45	0.84	1 (2%)

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
4	UDP	A	196	5	-	2/16/32/32	0/2/2/2
3	ADP	A	195	2,5	-	1/16/32/32	0/3/3/3

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	196	UDP	O2-C2	6.88	1.35	1.23
4	A	196	UDP	O4-C4	5.24	1.34	1.24
3	A	195	ADP	PA-O3A	-3.78	1.55	1.59
4	A	196	UDP	C5-C4	-3.65	1.35	1.43
4	A	196	UDP	C6-N1	3.45	1.46	1.38
4	A	196	UDP	C2-N3	-2.96	1.32	1.38
4	A	196	UDP	PB-O2B	-2.95	1.43	1.54
3	A	195	ADP	PA-O2A	-2.83	1.42	1.55
3	A	195	ADP	PB-O2B	-2.72	1.44	1.54
4	A	196	UDP	O4'-C4'	-2.32	1.39	1.45
4	A	196	UDP	C4-N3	-2.31	1.34	1.38
4	A	196	UDP	PA-O2A	-2.30	1.44	1.55
4	A	196	UDP	PB-O1B	-2.25	1.43	1.50
3	A	195	ADP	O2'-C2'	-2.16	1.37	1.43
3	A	195	ADP	PB-O1B	-2.12	1.43	1.50

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	196	UDP	N3-C2-N1	9.43	127.16	114.89
4	A	196	UDP	C6-N1-C2	-8.61	110.52	121.00
4	A	196	UDP	C4-N3-C2	-5.27	120.07	126.61
4	A	196	UDP	C6-C5-C4	4.50	125.28	119.53
4	A	196	UDP	O2-C2-N1	-4.33	117.16	122.80
4	A	196	UDP	O2-C2-N3	-3.60	114.86	121.49
4	A	196	UDP	C1'-N1-C6	2.48	126.08	120.78
4	A	196	UDP	O4'-C4'-C3'	-2.41	100.37	105.15
3	A	195	ADP	O4'-C1'-C2'	-2.05	102.22	106.62

There are no chirality outliers.

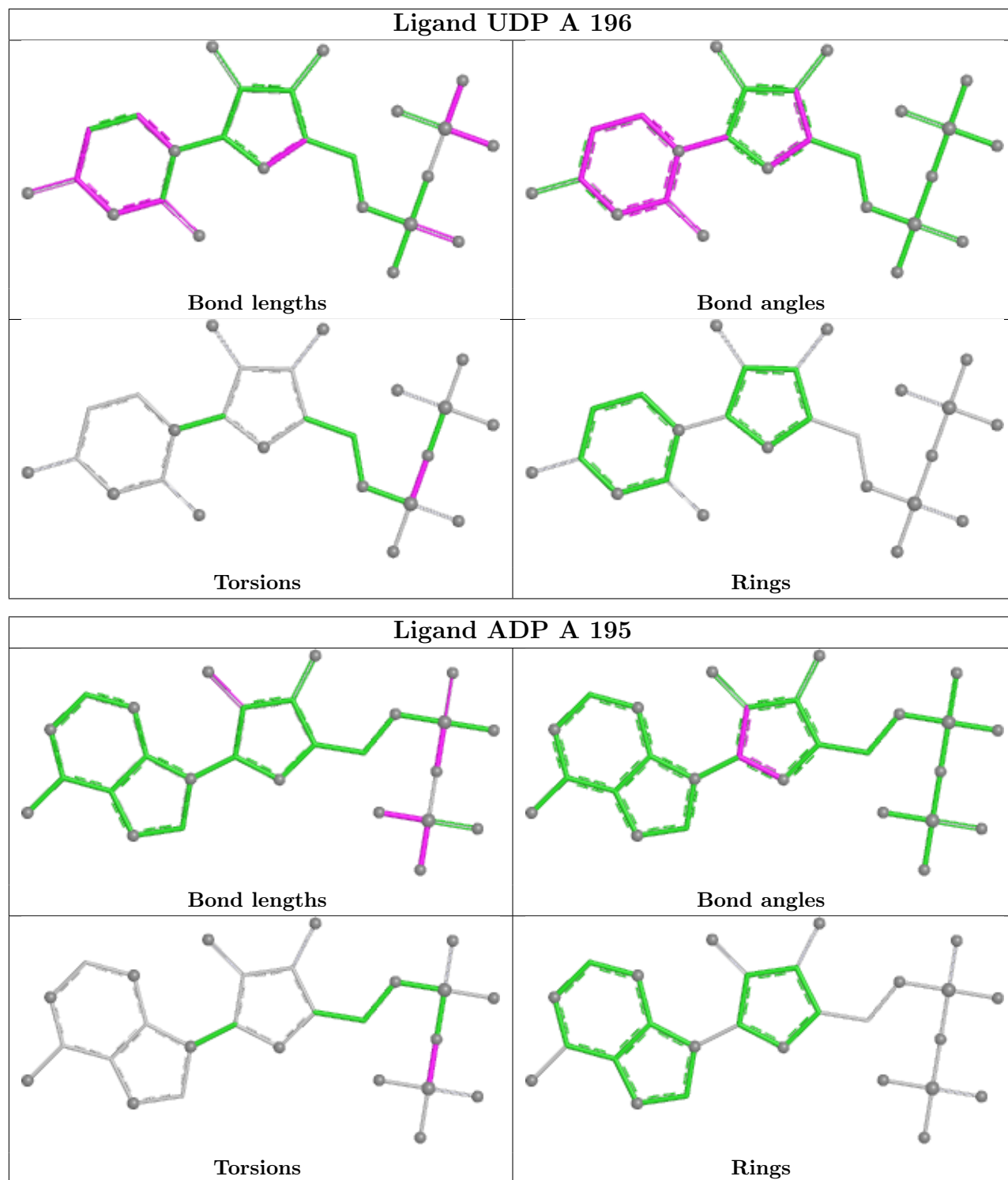
All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	196	UDP	PB-O3A-PA-O1A
4	A	196	UDP	PB-O3A-PA-O2A
3	A	195	ADP	PA-O3A-PB-O1B

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.



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	A	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	141:ASN	C	142:ILE	N	1.60

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	191/194 (98%)	4.06	179 (93%) 0 0	15, 26, 48, 54	0

All (179) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	48	GLY	8.9
1	A	128	LEU	8.5
1	A	37	ALA	8.4
1	A	132	GLY	7.8
1	A	47	SER	7.7
1	A	136	GLY	7.4
1	A	165	LYS	7.3
1	A	103	GLU	7.0
1	A	174	ALA	6.9
1	A	41	LEU	6.4
1	A	154	VAL	6.4
1	A	73	LEU	6.4
1	A	158	LEU	6.4
1	A	113	PHE	6.3
1	A	27	VAL	6.3
1	A	20	GLY	6.2
1	A	50	LYS	6.1
1	A	13	GLY	6.0
1	A	12	LEU	6.0
1	A	194	PHE	5.9
1	A	117	PHE	5.9
1	A	49	SER	5.8
1	A	108	PHE	5.8
1	A	166	PHE	5.8
1	A	15	PRO	5.8
1	A	6	PRO	5.6
1	A	52	GLY	5.6

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Mol	Chain	Res	Type	RSRZ
1	A	65	VAL	5.6
1	A	80	ALA	5.6
1	A	91	PHE	5.5
1	A	145	ILE	5.5
1	A	116	PHE	5.4
1	A	125	THR	5.4
1	A	87	LEU	5.4
1	A	71	VAL	5.4
1	A	95	GLU	5.4
1	A	54	MET	5.3
1	A	40	LEU	5.3
1	A	104	ASN	5.3
1	A	149	PHE	5.2
1	A	64	ILE	5.1
1	A	33	VAL	5.1
1	A	46	GLN	5.1
1	A	51	ASP	5.1
1	A	111	THR	5.0
1	A	5	LYS	5.0
1	A	173	PRO	5.0
1	A	186	GLU	5.0
1	A	92	PRO	4.9
1	A	35	LEU	4.9
1	A	159	VAL	4.9
1	A	57	THR	4.9
1	A	105	MET	4.8
1	A	11	VAL	4.8
1	A	109	VAL	4.8
1	A	123	VAL	4.8
1	A	185	VAL	4.8
1	A	74	LEU	4.7
1	A	101	TRP	4.7
1	A	178	VAL	4.6
1	A	10	PHE	4.6
1	A	110	ASP	4.5
1	A	129	LEU	4.5
1	A	55	ILE	4.5
1	A	97	ASN	4.5
1	A	32	TRP	4.5
1	A	31	GLY	4.5
1	A	90	GLY	4.4
1	A	189	PHE	4.4

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Mol	Chain	Res	Type	RSRZ
1	A	61	ASN	4.4
1	A	106	LYS	4.4
1	A	139	ASP	4.4
1	A	177	ASP	4.4
1	A	26	ILE	4.4
1	A	24	ALA	4.3
1	A	121	GLU	4.3
1	A	148	ARG	4.2
1	A	69	VAL	4.2
1	A	134	SER	4.2
1	A	4	SER	4.2
1	A	75	LYS	4.2
1	A	191	SER	4.1
1	A	76	ASN	4.1
1	A	23	CYS	4.0
1	A	60	LYS	4.0
1	A	79	ASP	4.0
1	A	78	ILE	4.0
1	A	99	ASN	4.0
1	A	107	ASP	3.9
1	A	151	THR	3.9
1	A	94	ASN	3.9
1	A	59	ILE	3.9
1	A	193	GLY	3.8
1	A	138	SER	3.8
1	A	88	VAL	3.8
1	A	83	GLY	3.8
1	A	21	THR	3.8
1	A	188	LEU	3.8
1	A	146	LYS	3.8
1	A	81	ASN	3.7
1	A	135	SER	3.7
1	A	119	CYS	3.7
1	A	133	GLU	3.7
1	A	56	ALA	3.7
1	A	58	MET	3.7
1	A	63	GLU	3.7
1	A	93	ARG	3.7
1	A	168	LYS	3.7
1	A	155	GLN	3.7
1	A	160	ILE	3.6
1	A	70	THR	3.6

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Mol	Chain	Res	Type	RSRZ
1	A	9	VAL	3.6
1	A	72	LYS	3.6
1	A	38	GLY	3.5
1	A	8	VAL	3.5
1	A	34	HIS	3.5
1	A	156	THR	3.5
1	A	86	PHE	3.5
1	A	171	ILE	3.4
1	A	16	GLY	3.4
1	A	118	ASP	3.4
1	A	169	VAL	3.4
1	A	89	ASP	3.4
1	A	176	ARG	3.4
1	A	137	ARG	3.4
1	A	30	PHE	3.4
1	A	152	PHE	3.3
1	A	44	GLU	3.3
1	A	53	GLU	3.3
1	A	122	GLU	3.3
1	A	43	GLN	3.3
1	A	126	GLN	3.3
1	A	68	ILE	3.2
1	A	85	ASN	3.2
1	A	140	ASP	3.2
1	A	39	ASP	3.2
1	A	45	GLN	3.2
1	A	19	LYS	3.2
1	A	42	ARG	3.1
1	A	157	LYS	3.1
1	A	143	GLU	3.1
1	A	163	TYR	3.1
1	A	77	ALA	3.1
1	A	62	GLY	3.0
1	A	180	GLU	3.0
1	A	114	VAL	3.0
1	A	67	SER	3.0
1	A	142	ILE	3.0
1	A	183	ASN	3.0
1	A	130	LYS	3.0
1	A	28	ARG	3.0
1	A	66	PRO	3.0
1	A	182	TYR	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	179	ASN	2.9
1	A	29	ASP	2.9
1	A	17	SER	2.9
1	A	112	LYS	2.9
1	A	181	VAL	2.8
1	A	82	GLN	2.8
1	A	162	HIS	2.8
1	A	167	ASP	2.8
1	A	141	ASN	2.8
1	A	190	LYS	2.7
1	A	161	ASP	2.7
1	A	100	SER	2.7
1	A	120	PRO	2.6
1	A	102	GLU	2.6
1	A	7	ASN	2.6
1	A	170	LYS	2.5
1	A	98	ASN	2.5
1	A	187	ASN	2.5
1	A	192	MET	2.4
1	A	18	GLY	2.4
1	A	127	ARG	2.2
1	A	22	GLN	2.2
1	A	36	SER	2.2
1	A	115	LEU	2.1
1	A	131	ARG	2.1
1	A	96	GLU	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](#)

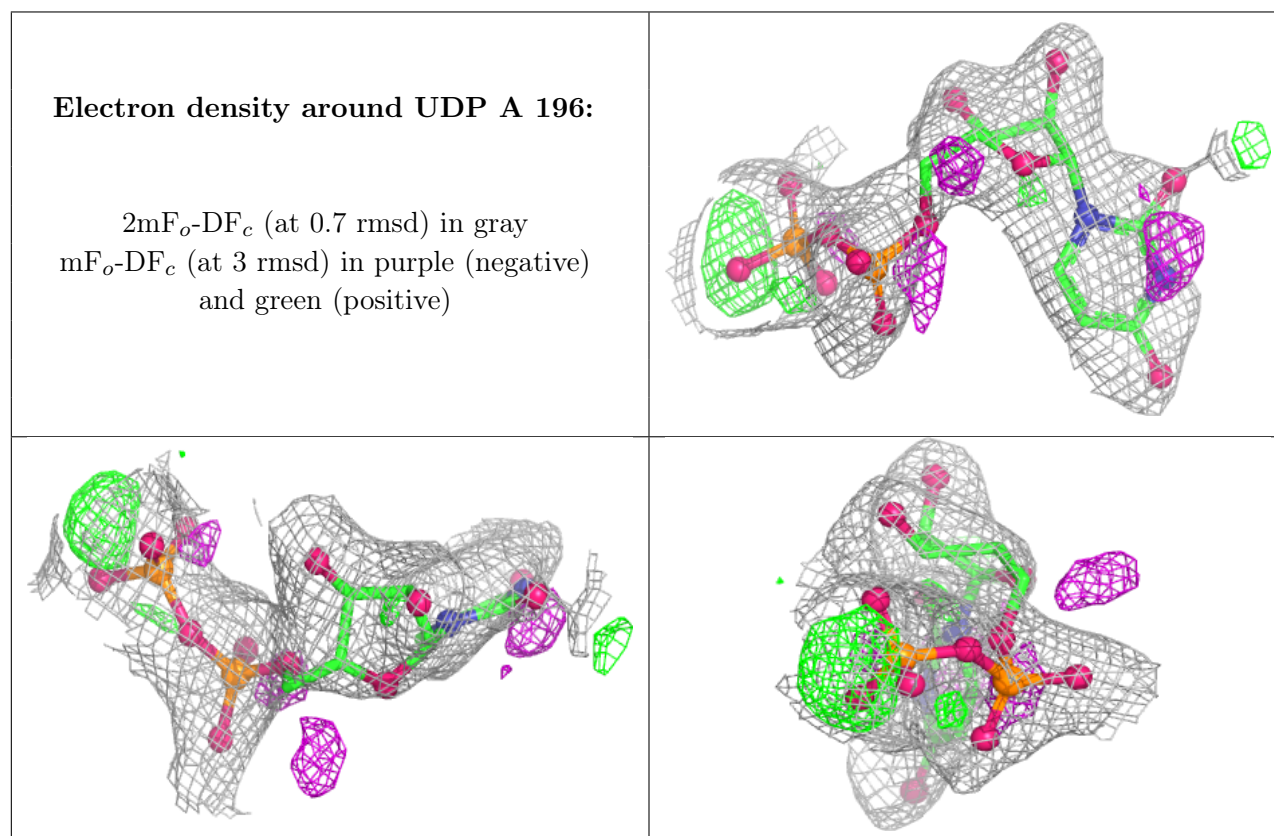
There are no oligosaccharides in this entry.

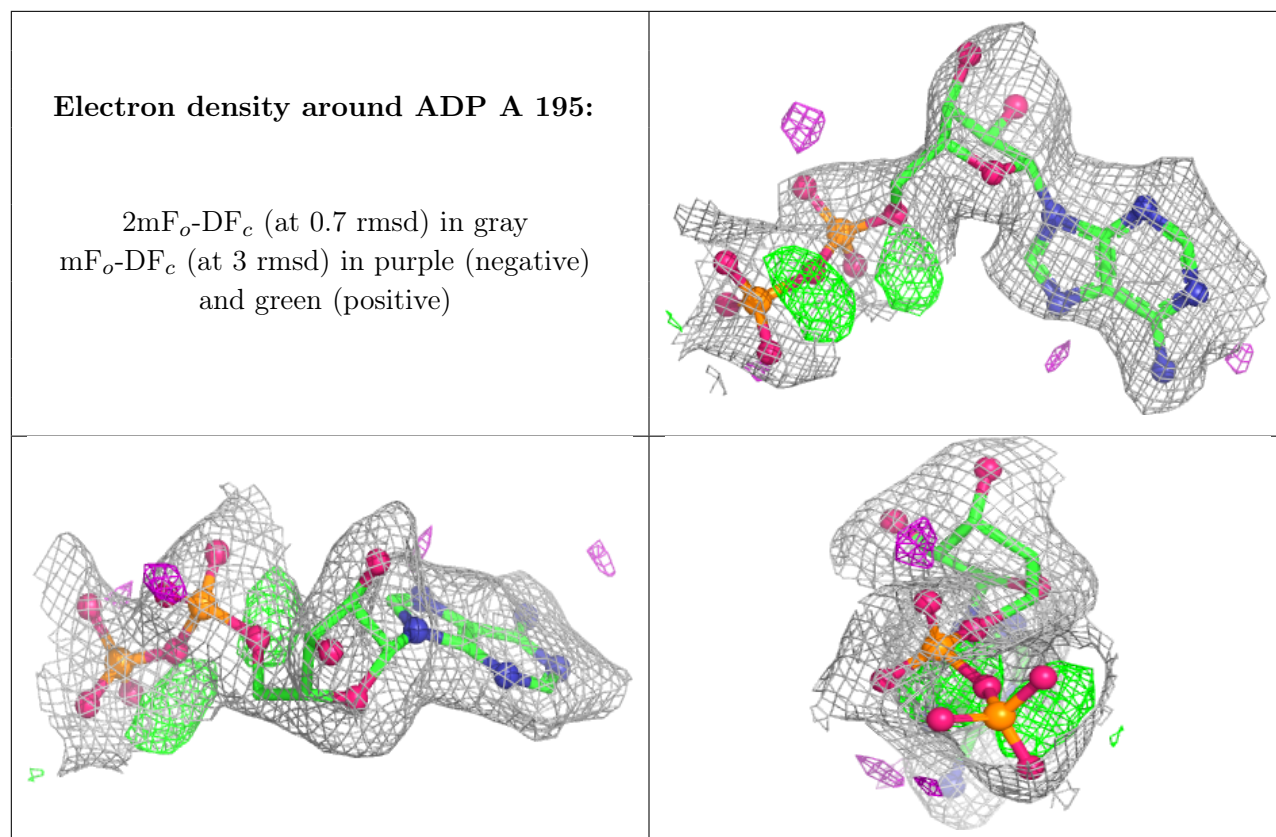
6.4 Ligands [i](#)

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($\text{\AA}^2$)	Q<0.9
2	MG	A	500	1/1	0.74	0.10	18,18,18,18	0
4	UDP	A	196	25/25	0.75	0.20	18,20,21,22	0
3	ADP	A	195	27/27	0.80	0.16	17,19,21,24	0
5	BF2	A	501	3/3	0.84	0.17	15,15,19,19	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.