Full wwPDB X-ray Structure Validation Report ⓘ

May 14, 2020 – 02:13 pm BST

PDB ID	:	4QBH
Title	:	Crystal structure of a stable adenylate kinase variant AKlse5
Authors	:	Moon, S.; Bae, E.
Deposited on	:	2014-05-08
Resolution	:	1.67 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:

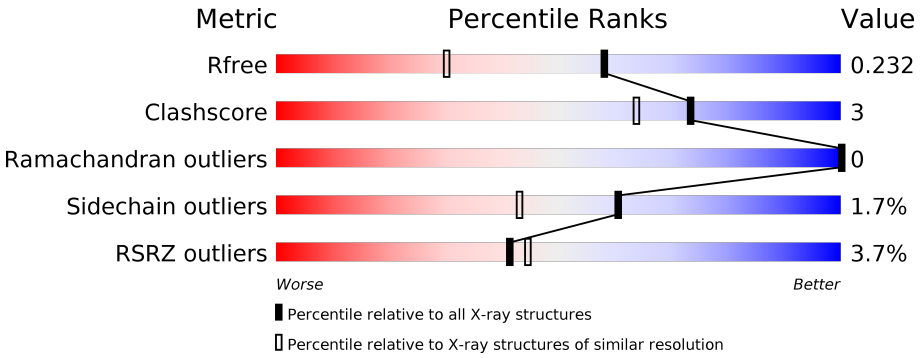
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.11
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.67 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	6780 (1.70-1.66)
Clashscore	141614	7310 (1.70-1.66)
Ramachandran outliers	138981	7173 (1.70-1.66)
Sidechain outliers	138945	7172 (1.70-1.66)
RSRZ outliers	127900	6661 (1.70-1.66)

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 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 <=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	217	4% 92% 7%
1	B	217	4% 84% 13%

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 3912 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 kinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	217	Total	C	N	O	S	0	0	0
			1707	1065	307	321	14			
1	B	215	Total	C	N	O	S	0	0	0
			1686	1053	301	318	14			

There are 42 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	19	ARG	LYS	engineered mutation	UNP P27142
A	22	GLU	ALA	engineered mutation	UNP P27142
A	73	GLY	SER	engineered mutation	UNP P27142
A	78	GLU	GLN	engineered mutation	UNP P27142
A	79	ARG	ASN	engineered mutation	UNP P27142
A	98	GLU	THR	engineered mutation	UNP P27142
A	99	ILE	MET	engineered mutation	UNP P27142
A	101	GLU	ALA	engineered mutation	UNP P27142
A	102	GLU	ASP	engineered mutation	UNP P27142
A	106	PRO	LYS	engineered mutation	UNP P27142
A	107	ILE	LEU	engineered mutation	UNP P27142
A	114	GLU	ASP	engineered mutation	UNP P27142
A	166	GLU	ALA	engineered mutation	UNP P27142
A	170	LYS	ASN	engineered mutation	UNP P27142
A	183	LEU	VAL	engineered mutation	UNP P27142
A	195	VAL	ILE	engineered mutation	UNP P27142
A	198	GLN	GLU	engineered mutation	UNP P27142
A	202	GLN	GLU	engineered mutation	UNP P27142
A	203	ASP	LYS	engineered mutation	UNP P27142
A	208	VAL	ILE	engineered mutation	UNP P27142
A	216	LYS	ALA	engineered mutation	UNP P27142
B	19	ARG	LYS	engineered mutation	UNP P27142
B	22	GLU	ALA	engineered mutation	UNP P27142
B	73	GLY	SER	engineered mutation	UNP P27142
B	78	GLU	GLN	engineered mutation	UNP P27142

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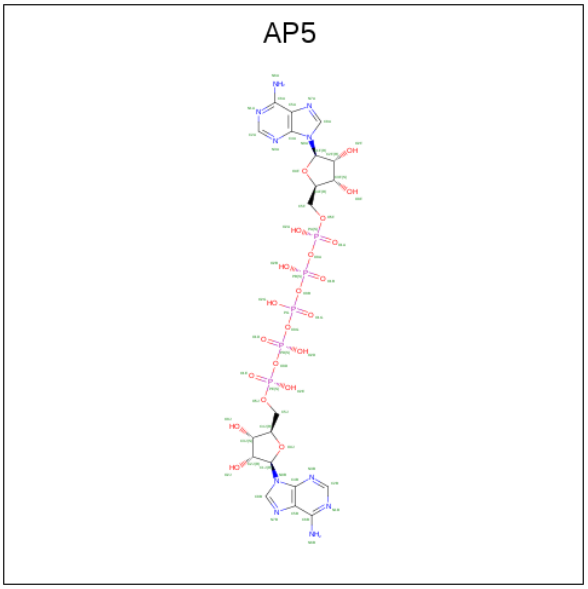
Chain	Residue	Modelled	Actual	Comment	Reference
B	79	ARG	ASN	engineered mutation	UNP P27142
B	98	GLU	THR	engineered mutation	UNP P27142
B	99	ILE	MET	engineered mutation	UNP P27142
B	101	GLU	ALA	engineered mutation	UNP P27142
B	102	GLU	ASP	engineered mutation	UNP P27142
B	106	PRO	LYS	engineered mutation	UNP P27142
B	107	ILE	LEU	engineered mutation	UNP P27142
B	114	GLU	ASP	engineered mutation	UNP P27142
B	166	GLU	ALA	engineered mutation	UNP P27142
B	170	LYS	ASN	engineered mutation	UNP P27142
B	183	LEU	VAL	engineered mutation	UNP P27142
B	195	VAL	ILE	engineered mutation	UNP P27142
B	198	GLN	GLU	engineered mutation	UNP P27142
B	202	GLN	GLU	engineered mutation	UNP P27142
B	203	ASP	LYS	engineered mutation	UNP P27142
B	208	VAL	ILE	engineered mutation	UNP P27142
B	216	LYS	ALA	engineered mutation	UNP P27142

- Molecule 2 is BIS(ADENOSINE)-5'-PENTAPHOSPHATE (three-letter code: AP5) (formula: $C_{20}H_{29}N_{10}O_{22}P_5$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			57	20	10	22	5		
2	B	1	Total	C	N	O	P	0	0
			57	20	10	22	5		

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	1	Total Zn 1 1	0	0
3	A	1	Total Zn 1 1	0	0

- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

- Molecule 5 is water.

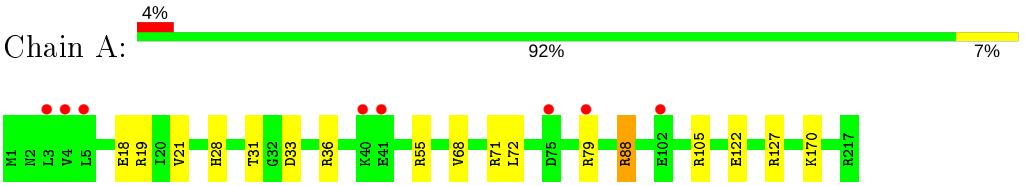
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	219	Total O 219 219	0	0
5	B	182	Total O 182 182	0	0

3 Residue-property plots (i)

These plots are drawn for all protein, RNA and DNA 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.

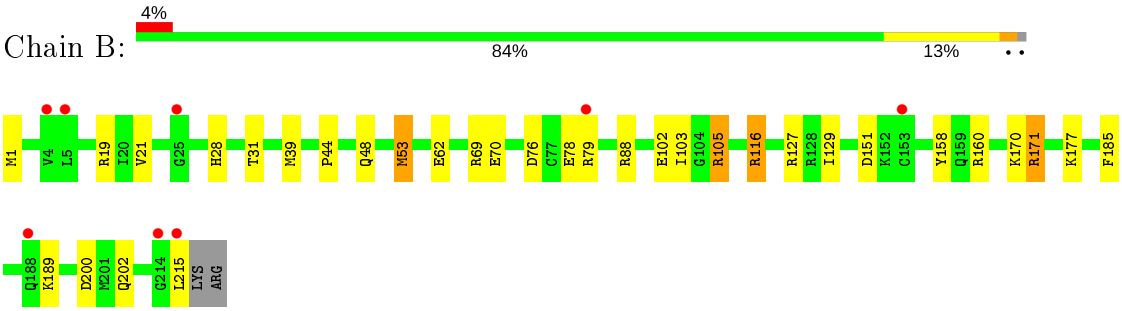
- Molecule 1: Adenylate kinase

Chain A:



- Molecule 1: Adenylate kinase

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	36.09Å 75.36Å 81.98Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 1.67 27.09 – 1.67	Depositor EDS
% Data completeness (in resolution range)	97.0 (50.00-1.67) 97.0 (27.09-1.67)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.29 (at 1.67Å)	Xtriage
Refinement program	REFMAC 5.8.0049	Depositor
R, R_{free}	0.176 , 0.223 0.187 , 0.232	Depositor DCC
R_{free} test set	2516 reflections (5.09%)	wwPDB-VP
Wilson B-factor (Å^2)	20.6	Xtriage
Anisotropy	0.158	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.39 , 41.5	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	0.038 for h,-k,-l	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	3912	wwPDB-VP
Average B, all atoms (Å^2)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.50% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 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: ZN, MG, AP5

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	0.96	1/1733 (0.1%)	1.06	7/2329 (0.3%)
1	B	0.98	1/1712 (0.1%)	1.17	13/2304 (0.6%)
All	All	0.97	2/3445 (0.1%)	1.12	20/4633 (0.4%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	62	GLU	CD-OE2	-7.92	1.17	1.25
1	A	18	GLU	CD-OE2	-5.07	1.20	1.25

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	171	ARG	NE-CZ-NH2	-13.20	113.70	120.30
1	B	88	ARG	NE-CZ-NH2	-12.59	114.00	120.30
1	B	53	MET	CG-SD-CE	-10.61	83.23	100.20
1	A	19	ARG	NE-CZ-NH1	8.53	124.56	120.30
1	B	69	ARG	NE-CZ-NH1	7.95	124.28	120.30
1	A	88	ARG	NE-CZ-NH1	7.20	123.90	120.30
1	B	19	ARG	NE-CZ-NH2	-7.18	116.71	120.30
1	A	19	ARG	NE-CZ-NH2	-7.00	116.80	120.30
1	B	151	ASP	CB-CG-OD1	6.66	124.29	118.30
1	B	171	ARG	NE-CZ-NH1	6.63	123.62	120.30
1	B	88	ARG	NE-CZ-NH1	6.61	123.60	120.30
1	A	88	ARG	NE-CZ-NH2	-6.43	117.08	120.30
1	A	71	ARG	NE-CZ-NH1	5.55	123.08	120.30
1	B	170	LYS	CD-CE-NZ	-5.24	99.64	111.70
1	A	105	ARG	NE-CZ-NH2	-5.20	117.70	120.30
1	A	79	ARG	NE-CZ-NH1	5.17	122.89	120.30
1	B	116	ARG	NE-CZ-NH2	-5.13	117.73	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	105	ARG	NE-CZ-NH1	5.07	122.83	120.30
1	B	76	ASP	CB-CG-OD1	5.06	122.86	118.30
1	B	160	ARG	NE-CZ-NH2	-5.01	117.80	120.30

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	1707	0	1707	6	0
1	B	1686	0	1681	17	0
2	A	57	0	24	2	0
2	B	57	0	24	2	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	219	0	0	0	0
5	B	182	0	0	2	0
All	All	3912	0	3436	23	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 3.

All (23) 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:B:39:MET:HE1	1:B:53:MET:HE3	1.26	1.09
1:B:39:MET:HE1	1:B:53:MET:CE	2.02	0.88
1:B:39:MET:CE	1:B:53:MET:HE3	2.05	0.84
1:A:33:ASP:OD1	1:A:36:ARG:NH1	2.15	0.80
1:A:127:ARG:HH12	2:A:301:AP5:PG	2.10	0.74
1:B:39:MET:CE	1:B:53:MET:CE	2.65	0.71
1:B:48:GLN:HB3	5:B:573:HOH:O	1.95	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:127:ARG:HH12	2:B:301:AP5:PG	2.23	0.62
1:A:88:ARG:NH2	2:A:301:AP5:O1D	2.34	0.60
1:B:103:ILE:HD12	1:B:105:ARG:HD2	1.85	0.57
1:B:200:ASP:OD2	1:B:202:GLN:HG2	2.05	0.55
1:A:21:VAL:HG21	1:A:28:HIS:HB2	1.93	0.50
1:B:129:ILE:HG13	1:B:158:TYR:CE1	2.48	0.49
1:B:177:LYS:O	1:B:177:LYS:HG2	2.14	0.48
1:B:1:MET:HG2	1:B:215:LEU:HD23	1.96	0.48
1:B:171:ARG:NH2	2:B:301:AP5:O1G	2.48	0.47
1:B:185:PHE:CZ	1:B:189:LYS:HE3	2.51	0.46
1:B:78:GLU:HG3	1:B:79:ARG:HD2	1.99	0.45
1:B:39:MET:HB2	1:B:39:MET:HE2	1.63	0.43
1:A:68:VAL:HG12	1:A:72:LEU:HD12	2.00	0.42
1:B:21:VAL:HG11	1:B:28:HIS:HB2	2.03	0.41
1:A:55:ARG:O	1:A:170:LYS:HG2	2.21	0.41
1:B:116:ARG:NE	5:B:468:HOH:O	2.54	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	215/217 (99%)	213 (99%)	2 (1%)	0	100	100
1	B	213/217 (98%)	210 (99%)	3 (1%)	0	100	100
All	All	428/434 (99%)	423 (99%)	5 (1%)	0	100	100

There are no Ramachandran outliers to report.

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	180/180 (100%)	178 (99%)	2 (1%)	73 61
1	B	178/180 (99%)	174 (98%)	4 (2%)	52 32
All	All	358/360 (99%)	352 (98%)	6 (2%)	60 43

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	31	THR
1	A	122	GLU
1	B	31	THR
1	B	44	PRO
1	B	70	GLU
1	B	102	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 carbohydrates in this entry.

5.6 Ligand geometry

Of 6 ligands modelled in this entry, 4 are monoatomic - leaving 2 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	$\#\|Z\| > 2$
2	AP5	B	301	4	48,62,62	1.13	4 (8%)	51,98,98	1.68	12 (23%)
2	AP5	A	301	4	48,62,62	1.23	6 (12%)	51,98,98	1.94	13 (25%)

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
2	AP5	B	301	4	-	3/36/76/76	0/6/6/6
2	AP5	A	301	4	-	4/36/76/76	0/6/6/6

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	301	AP5	C2B-N3B	3.44	1.37	1.32
2	B	301	AP5	O4F-C1F	3.30	1.45	1.41
2	A	301	AP5	C2B-N3B	3.29	1.37	1.32
2	A	301	AP5	O4J-C1J	3.18	1.45	1.41
2	B	301	AP5	C2J-C1J	-2.60	1.49	1.53
2	A	301	AP5	C2A-N3A	2.43	1.36	1.32
2	B	301	AP5	C2A-N3A	2.31	1.35	1.32
2	A	301	AP5	C2J-C1J	-2.27	1.50	1.53
2	A	301	AP5	O4F-C1F	2.27	1.44	1.41
2	A	301	AP5	C8A-N7A	2.06	1.38	1.34

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	301	AP5	N3B-C2B-N1B	-7.19	117.44	128.68
2	A	301	AP5	PE-O3D-PD	-5.12	115.27	132.83
2	B	301	AP5	N3B-C2B-N1B	-4.24	122.05	128.68
2	B	301	AP5	PD-O3G-PG	-4.04	118.96	132.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	301	AP5	N3A-C2A-N1A	-3.70	122.89	128.68
2	A	301	AP5	C5B-C6B-N6B	-3.56	114.94	120.35
2	A	301	AP5	O2G-PG-O1G	3.56	129.82	112.24
2	B	301	AP5	N6A-C6A-N1A	3.33	125.49	118.57
2	B	301	AP5	C2A-N1A-C6A	3.01	123.90	118.75
2	B	301	AP5	PE-O3D-PD	-2.95	122.70	132.83
2	A	301	AP5	C2A-N1A-C6A	2.89	123.69	118.75
2	A	301	AP5	C3F-C2F-C1F	2.79	105.19	100.98
2	B	301	AP5	C1J-N9B-C4B	-2.76	121.79	126.64
2	B	301	AP5	C3J-C2J-C1J	2.76	105.13	100.98
2	A	301	AP5	PD-O3G-PG	-2.70	123.56	132.83
2	B	301	AP5	C3F-C2F-C1F	2.66	104.99	100.98
2	B	301	AP5	N3A-C2A-N1A	-2.65	124.54	128.68
2	B	301	AP5	C5A-C6A-N1A	-2.50	114.69	120.35
2	A	301	AP5	N6B-C6B-N1B	2.49	123.74	118.57
2	A	301	AP5	C2B-N1B-C6B	2.47	122.98	118.75
2	B	301	AP5	C2B-N1B-C6B	2.27	122.64	118.75
2	B	301	AP5	C2J-C3J-C4J	-2.27	98.23	102.64
2	A	301	AP5	O2D-PD-O1D	2.07	122.49	112.24
2	A	301	AP5	C1J-N9B-C4B	-2.05	123.03	126.64
2	A	301	AP5	C1F-N9A-C4A	-2.01	123.11	126.64

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	301	AP5	PD-O3D-PE-O1E
2	B	301	AP5	PA-O3A-PB-O1B
2	A	301	AP5	PA-O3A-PB-O1B
2	A	301	AP5	PG-O3B-PB-O2B
2	A	301	AP5	PB-O3B-PG-O2G
2	B	301	AP5	PD-O3D-PE-O2E
2	A	301	AP5	PA-O3A-PB-O2B

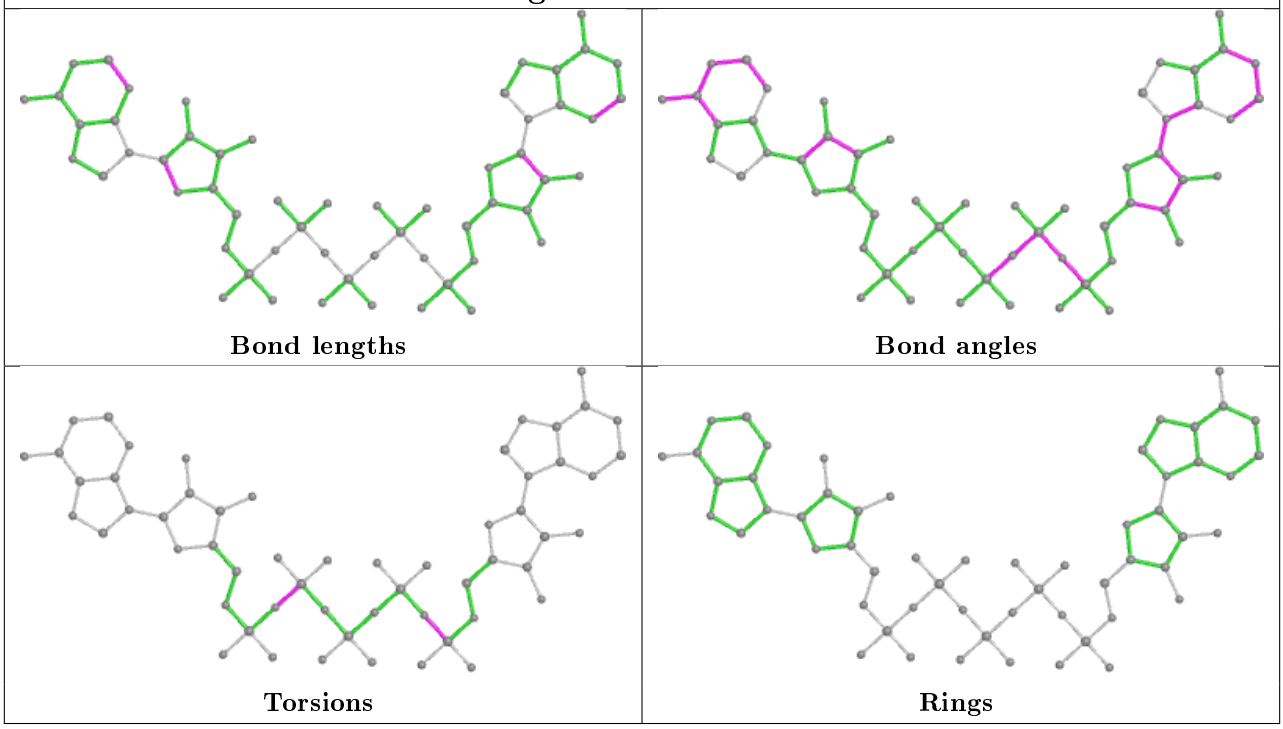
There are no ring outliers.

2 monomers are involved in 4 short contacts:

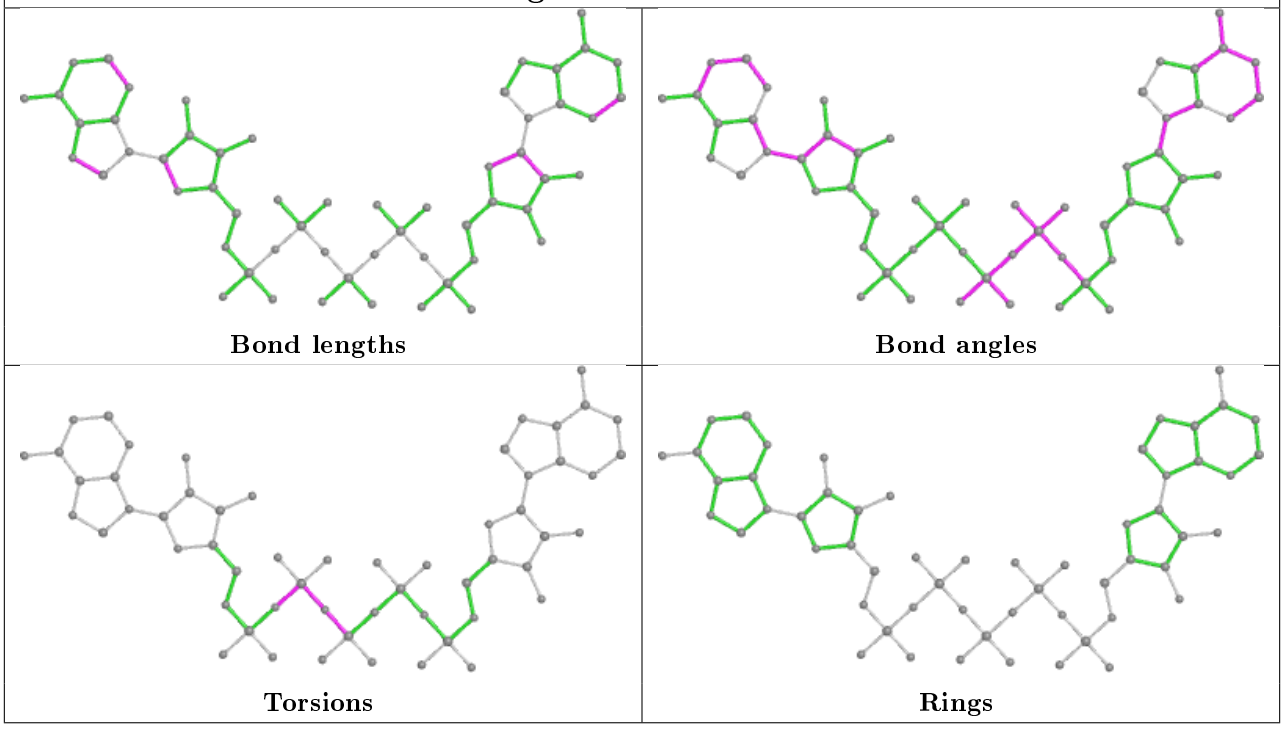
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	301	AP5	2	0
2	A	301	AP5	2	0

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 then 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.

Ligand AP5 B 301



Ligand AP5 A 301



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^{th} 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	217/217 (100%)	0.01	8 (3%) 41 44	14, 23, 37, 51	0
1	B	215/217 (99%)	0.20	8 (3%) 41 44	14, 26, 45, 56	0
All	All	432/434 (99%)	0.10	16 (3%) 41 44	14, 24, 41, 56	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	214	GLY	4.1
1	B	153	CYS	3.8
1	B	215	LEU	3.8
1	B	5	LEU	3.3
1	B	79	ARG	3.0
1	A	79	ARG	2.8
1	B	4	VAL	2.7
1	A	40	LYS	2.5
1	A	41	GLU	2.4
1	A	4	VAL	2.3
1	A	3	LEU	2.2
1	B	25	GLY	2.2
1	B	188	GLN	2.1
1	A	5	LEU	2.1
1	A	75	ASP	2.1
1	A	102	GLU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains

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

6.3 Carbohydrates (i)

There are no carbohydrates in this entry.

6.4 Ligands (i)

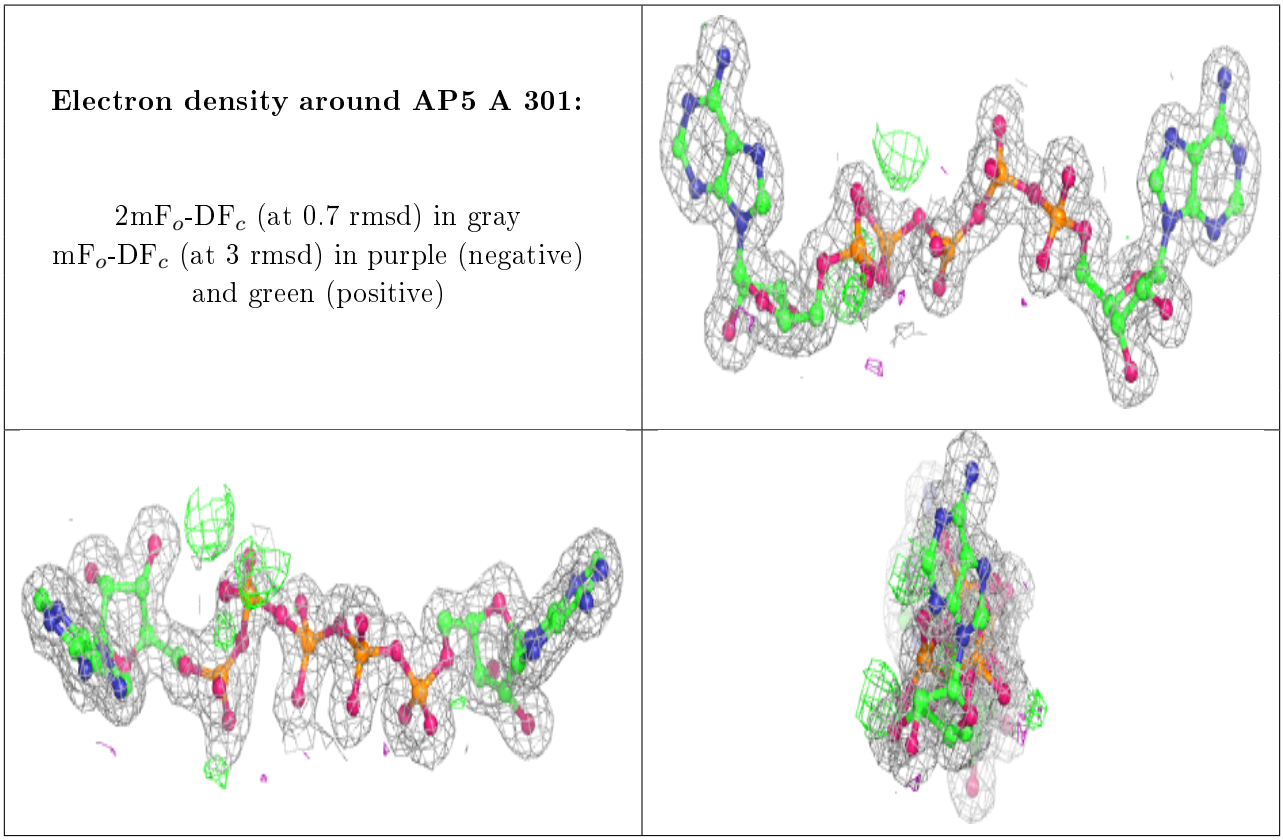
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' 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	ZN	B	302	1/1	0.83	0.10	43,43,43,43	0
4	MG	A	303	1/1	0.94	0.09	30,30,30,30	0
2	AP5	A	301	57/57	0.97	0.10	13,16,26,29	8
2	AP5	B	301	57/57	0.97	0.09	13,16,21,29	7
4	MG	B	303	1/1	0.98	0.06	35,35,35,35	0
3	ZN	A	302	1/1	0.98	0.06	26,26,26,26	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.

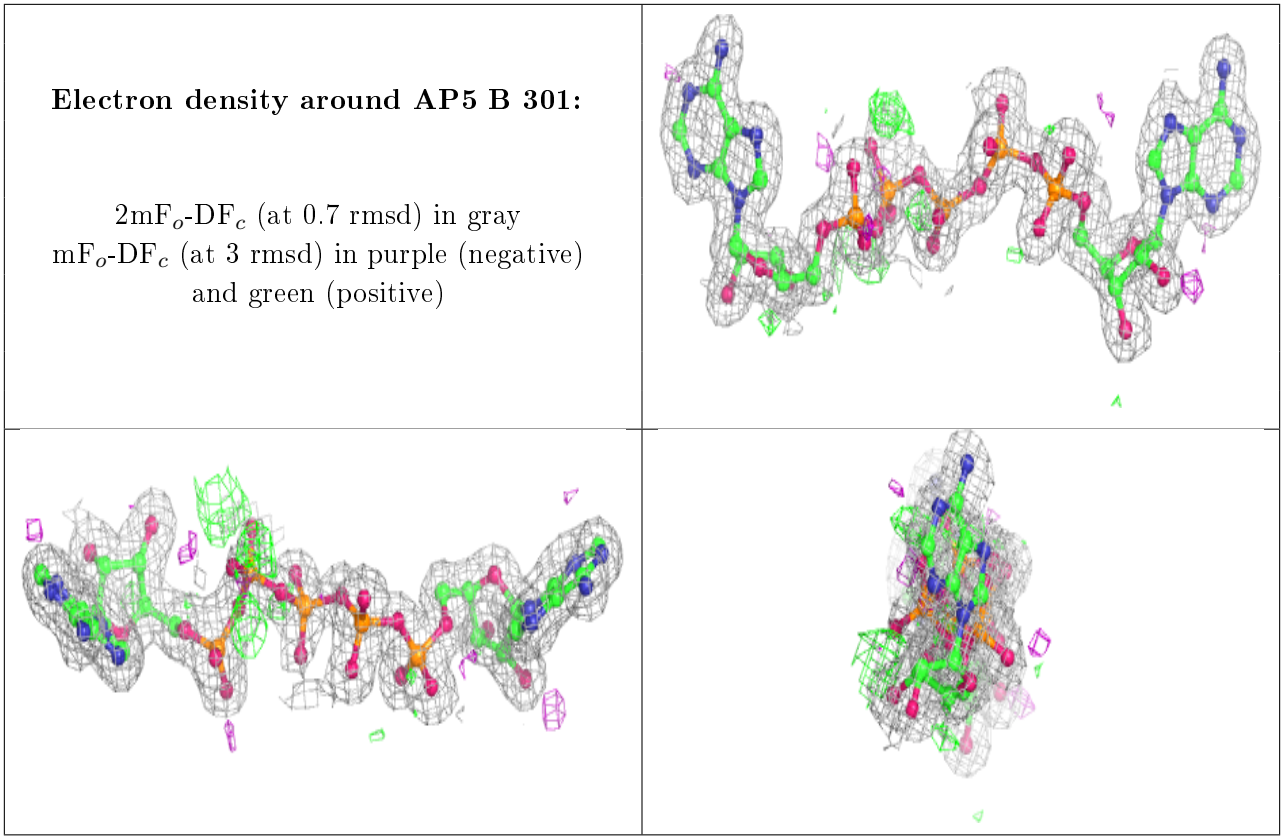
Electron density around AP5 A 301:

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



Electron density around AP5 B 301:

$2mF_o$-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.