Full wwPDB X-ray Structure Validation Report (i)

Apr 28, 2024 – 04:13 pm BST

PDB ID	:	1E4Y
Title	:	Mutant P9L of adenylate kinase from E. coli, modified in the Gly-loop
Authors	:	Mueller, C.W.; Schulz, G.E.
Deposited on	:	2000-07-12
Resolution	:	1.85 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 (i) 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 (i)) were used in the production of this report:

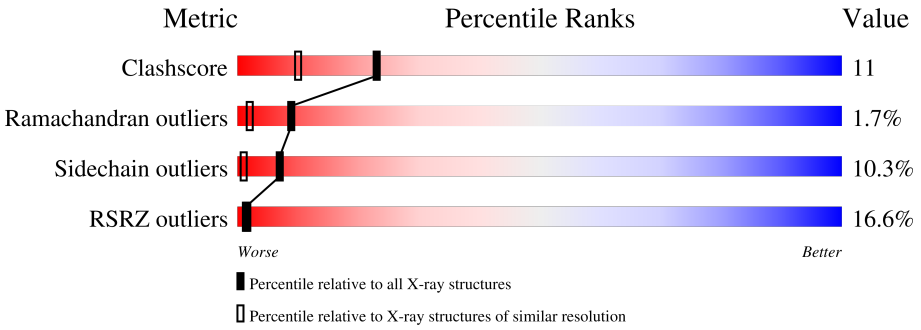
MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36.2
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

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 1.85 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))
Clashscore	141614	2625 (1.86-1.86)
Ramachandran outliers	138981	2592 (1.86-1.86)
Sidechain outliers	138945	2592 (1.86-1.86)
RSRZ outliers	127900	2436 (1.86-1.86)

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 <=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	214	16% 63% 30% 6% •
1	B	214	17% 61% 34% • •

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 3428 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.

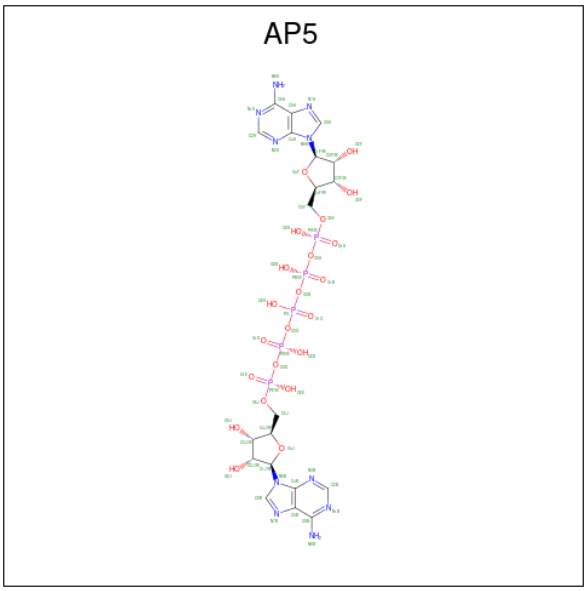
- Molecule 1 is a protein called Adenylate kinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	214	1657	1041	289	320	7	0	0	0
1	B	214	1657	1041	289	320	7	0	0	0

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	9	LEU	PRO	engineered mutation	UNP P69441
A	10	VAL	GLY	conflict	UNP P69441
B	9	LEU	PRO	engineered mutation	UNP P69441
B	10	VAL	GLY	conflict	UNP P69441

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



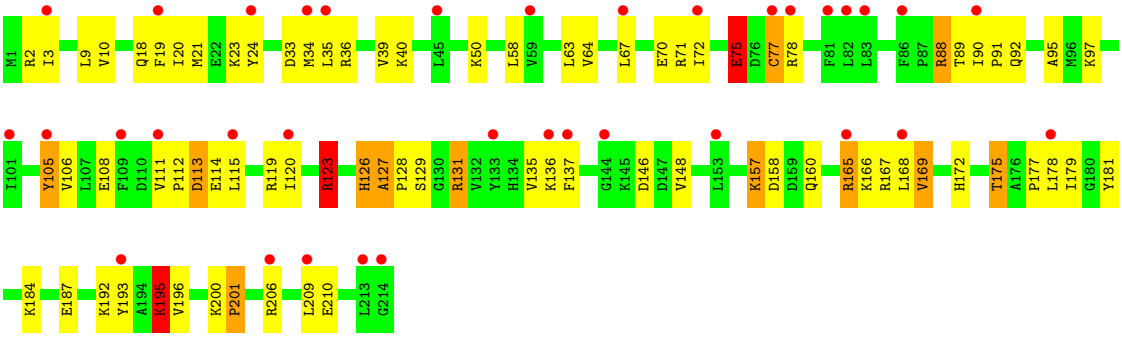
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
2	A	1	57	20	10	22	5	0	0
			Total	C	N	O	P		
2	B	1	57	20	10	22	5	0	0

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.

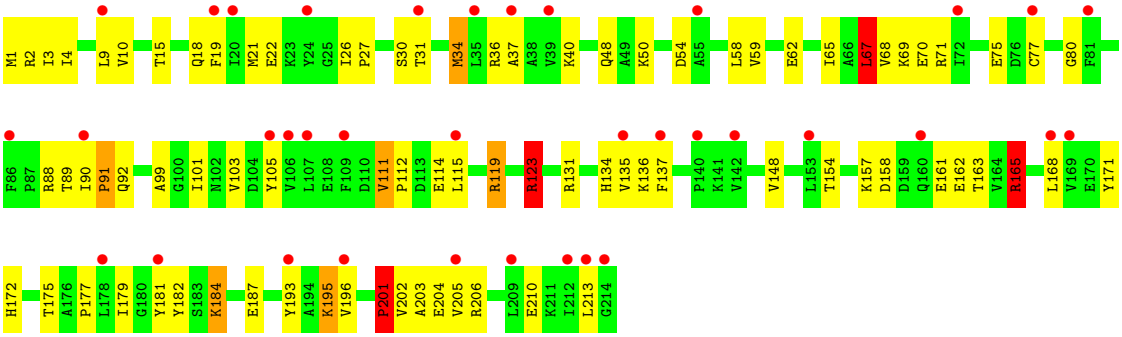
- Molecule 1: Adenylate kinase

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	60.30Å 77.80Å 57.70Å 90.00° 94.30° 90.00°	Depositor
Resolution (Å)	10.00 – 1.85 77.80 – 3.42	Depositor EDS
% Data completeness (in resolution range)	90.0 (10.00-1.85) 99.7 (77.80-3.42)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	0.10	Depositor
$< I/\sigma(I) >$	-	Xtriage
Refinement program	X-PLOR 1.5	Depositor
R, R_{free}	0.178 , (Not available) 0.160 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å^2)	35.0	Xtriage
Anisotropy	0.073	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.32 , 83.8	EDS
L-test for twinning[1]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Estimated twinning fraction	0.032 for l,-k,h	Xtriage
F_o,F_c correlation	0.93	EDS
Total number of atoms	3428	wwPDB-VP
Average B, all atoms (Å^2)	13.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.53% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]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 (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: 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	1.00	1/1680 (0.1%)	1.74	27/2261 (1.2%)
1	B	1.01	0/1680	1.70	25/2261 (1.1%)
All	All	1.01	1/3360 (0.0%)	1.72	52/4522 (1.1%)

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	B	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	195	LYS	CA-CB	-5.13	1.42	1.53

All (52) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	71	ARG	NE-CZ-NH2	-11.33	114.63	120.30
1	A	34	MET	CG-SD-CE	-11.32	82.09	100.20
1	A	2	ARG	NE-CZ-NH2	-10.43	115.09	120.30
1	A	71	ARG	NE-CZ-NH2	-9.07	115.77	120.30
1	B	50	LYS	CA-CB-CG	-8.74	94.16	113.40
1	A	181	TYR	CB-CG-CD2	-8.18	116.09	121.00
1	A	123	ARG	NE-CZ-NH1	7.81	124.20	120.30
1	A	165	ARG	NE-CZ-NH1	7.62	124.11	120.30
1	B	67	LEU	CA-CB-CG	7.41	132.35	115.30
1	B	184	LYS	CA-CB-CG	7.37	129.62	113.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	111	VAL	CA-CB-CG2	-7.35	99.87	110.90
1	B	34	MET	CG-SD-CE	-7.30	88.53	100.20
1	A	50	LYS	CA-CB-CG	-6.87	98.28	113.40
1	B	165	ARG	NE-CZ-NH1	6.86	123.73	120.30
1	B	119	ARG	NE-CZ-NH1	6.70	123.65	120.30
1	A	2	ARG	NE-CZ-NH1	6.67	123.64	120.30
1	B	181	TYR	CB-CG-CD2	-6.44	117.13	121.00
1	B	68	VAL	CG1-CB-CG2	-6.38	100.70	110.90
1	B	15	THR	OG1-CB-CG2	-6.37	95.36	110.00
1	A	33	ASP	CB-CG-OD1	6.35	124.02	118.30
1	A	165	ARG	NE-CZ-NH2	-6.30	117.15	120.30
1	A	67	LEU	CA-CB-CG	6.20	129.57	115.30
1	B	48	GLN	CB-CA-C	-6.10	98.20	110.40
1	B	123	ARG	NE-CZ-NH1	6.05	123.33	120.30
1	B	111	VAL	CA-CB-CG1	6.00	119.90	110.90
1	B	22	GLU	CB-CA-C	-5.99	98.42	110.40
1	A	169	VAL	CA-CB-CG2	-5.98	101.93	110.90
1	B	54	ASP	CB-CG-OD2	-5.93	112.97	118.30
1	A	63	LEU	CB-CA-C	-5.91	98.98	110.20
1	A	167	ARG	NE-CZ-NH2	-5.90	117.35	120.30
1	B	184	LYS	CB-CA-C	-5.82	98.76	110.40
1	A	50	LYS	CB-CG-CD	5.73	126.49	111.60
1	B	134	HIS	CA-CB-CG	-5.69	103.93	113.60
1	B	203	ALA	N-CA-CB	-5.67	102.16	110.10
1	A	119	ARG	NE-CZ-NH1	5.67	123.13	120.30
1	A	114	GLU	CA-CB-CG	5.61	125.75	113.40
1	A	131	ARG	NE-CZ-NH2	5.53	123.06	120.30
1	A	169	VAL	CA-CB-CG1	5.52	119.19	110.90
1	B	182	TYR	CB-CG-CD1	-5.51	117.69	121.00
1	A	64	VAL	CG1-CB-CG2	-5.49	102.11	110.90
1	A	77	CYS	CA-CB-SG	-5.49	104.12	114.00
1	B	36	ARG	NE-CZ-NH2	5.48	123.04	120.30
1	A	119	ARG	NE-CZ-NH2	-5.45	117.58	120.30
1	B	77	CYS	CA-CB-SG	-5.43	104.22	114.00
1	B	105	TYR	CB-CG-CD1	-5.35	117.79	121.00
1	B	114	GLU	CA-CB-CG	5.34	125.15	113.40
1	A	123	ARG	NE-CZ-NH2	-5.32	117.64	120.30
1	A	106	VAL	N-CA-C	-5.31	96.67	111.00
1	A	126	HIS	N-CA-C	-5.23	96.87	111.00
1	B	195	LYS	N-CA-CB	-5.21	101.22	110.60
1	A	127	ALA	CA-C-N	5.20	131.65	117.10
1	A	105	TYR	CB-CG-CD1	-5.13	117.92	121.00

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	171	TYR	Sidechain

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	1657	0	1687	41	0
1	B	1657	0	1687	38	0
2	A	57	0	22	8	0
2	B	57	0	22	6	0
All	All	3428	0	3418	78	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 11.

All (78) 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:92:GLN:HE22	2:B:215:AP5:H62B	1.28	0.81
1:A:19:PHE:HB2	1:A:206:ARG:HH21	1.49	0.78
1:A:137:PHE:HB3	2:A:215:AP5:H2A	1.75	0.68
1:B:19:PHE:HZ	1:B:210:GLU:HG2	1.59	0.68
1:B:27:PRO:HD2	1:B:80:GLY:O	1.97	0.64
1:A:135:VAL:HG23	1:A:136:LYS:HG3	1.80	0.64
1:A:92:GLN:HE22	2:A:215:AP5:H62B	1.45	0.64
1:A:131:ARG:NH1	1:A:148:VAL:HB	2.13	0.64
1:B:168:LEU:HD22	1:B:172:HIS:CE1	2.34	0.63
1:B:137:PHE:HB3	2:B:215:AP5:H2A	1.81	0.62
1:B:161:GLU:O	1:B:165:ARG:HG3	2.01	0.61
1:B:131:ARG:NH1	1:B:148:VAL:HB	2.16	0.60
1:A:126:HIS:HD1	1:A:129:SER:HG	1.50	0.59
1:A:36:ARG:O	1:A:40:LYS:HG2	2.03	0.59
1:A:123:ARG:HG3	2:A:215:AP5:H3F	1.85	0.59
1:B:9:LEU:HA	2:B:215:AP5:O1G	2.03	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:58:LEU:HD21	1:B:88:ARG:HD3	1.85	0.58
1:A:9:LEU:HA	2:A:215:AP5:O2G	2.02	0.58
1:B:123:ARG:HG3	2:B:215:AP5:H3F	1.86	0.58
1:B:18:GLN:O	1:B:21:MET:HB3	2.05	0.56
1:A:137:PHE:HB3	2:A:215:AP5:C2A	2.36	0.55
1:A:20:ILE:HD11	1:A:209:LEU:HD13	1.89	0.55
1:A:3:ILE:HG12	1:A:105:TYR:HB2	1.90	0.54
1:B:59:VAL:O	2:B:215:AP5:H2B	2.08	0.54
1:A:18:GLN:O	1:A:21:MET:HB3	2.08	0.54
1:A:157:LYS:O	1:A:160:GLN:HG2	2.08	0.53
1:B:19:PHE:CZ	1:B:210:GLU:HG2	2.42	0.53
1:B:30:SER:O	1:B:34:MET:HG3	2.08	0.53
1:B:158:ASP:HA	1:B:163:THR:HG21	1.88	0.53
1:B:37:ALA:HA	1:B:40:LYS:HG2	1.89	0.53
1:B:135:VAL:HG23	1:B:136:LYS:HG3	1.91	0.53
1:A:112:PRO:HG2	1:A:115:LEU:HD13	1.91	0.53
1:A:23:LYS:HD3	1:A:24:TYR:CZ	2.44	0.52
1:A:108:GLU:HB3	1:A:195:LYS:HG2	1.90	0.52
1:A:168:LEU:HD22	1:A:172:HIS:CE1	2.45	0.51
1:B:34:MET:SD	1:B:67:LEU:CD2	2.99	0.51
1:B:131:ARG:HH12	1:B:148:VAL:HB	1.74	0.51
1:A:88:ARG:HA	1:A:175:THR:HG23	1.92	0.51
1:B:19:PHE:HB2	1:B:206:ARG:HH21	1.76	0.50
1:A:75:GLU:O	1:A:78:ARG:HB2	2.11	0.50
1:B:3:ILE:HD12	1:B:26:ILE:HD11	1.94	0.50
1:B:2:ARG:HB3	1:B:103:VAL:HG12	1.95	0.49
1:A:19:PHE:HZ	1:A:210:GLU:CG	2.25	0.49
1:B:112:PRO:HG2	1:B:115:LEU:HD13	1.94	0.48
1:B:123:ARG:HG3	2:B:215:AP5:C3F	2.43	0.48
1:A:19:PHE:HZ	1:A:210:GLU:HG2	1.79	0.48
1:A:123:ARG:HG3	2:A:215:AP5:C3F	2.42	0.47
1:A:35:LEU:O	1:A:39:VAL:HG23	2.14	0.47
1:B:34:MET:HB3	1:B:67:LEU:HD21	1.98	0.46
1:B:90:ILE:HB	1:B:91:PRO:HD3	1.97	0.46
1:A:126:HIS:CE1	1:A:128:PRO:HG2	2.51	0.45
1:B:88:ARG:HA	1:B:175:THR:HG23	1.98	0.45
1:A:19:PHE:HB2	1:A:206:ARG:NH2	2.26	0.45
1:B:4:ILE:HG13	1:B:103:VAL:HG21	1.97	0.45
1:A:88:ARG:H	1:A:92:GLN:NE2	2.14	0.45
1:A:179:ILE:HG23	1:A:193:TYR:OH	2.18	0.44
1:A:92:GLN:NE2	2:A:215:AP5:H62B	2.14	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:179:ILE:HG23	1:B:193:TYR:OH	2.18	0.43
1:A:184:LYS:O	1:A:187:GLU:HB2	2.18	0.43
1:A:131:ARG:HH12	1:A:148:VAL:HB	1.82	0.43
1:B:201:PRO:O	1:B:205:VAL:HG23	2.17	0.43
1:A:90:ILE:HD11	1:A:177:PRO:HB2	2.00	0.43
1:A:131:ARG:NH1	1:A:146:ASP:OD2	2.52	0.43
1:A:200:LYS:O	2:A:215:AP5:N6A	2.51	0.42
1:B:202:VAL:HG12	1:B:206:ARG:NH1	2.34	0.42
1:A:90:ILE:HD13	1:A:178:LEU:HD23	2.02	0.42
1:A:95:ALA:HA	1:B:40:LYS:HA	2.01	0.41
1:A:72:ILE:HG21	1:A:72:ILE:HD13	1.85	0.41
1:A:72:ILE:O	1:A:77:CYS:HB2	2.21	0.41
1:B:184:LYS:O	1:B:187:GLU:HB2	2.21	0.41
1:B:1:MET:O	1:B:80:GLY:HA3	2.21	0.41
1:A:58:LEU:HD21	1:A:88:ARG:HD3	2.01	0.41
1:B:99:ALA:HB3	1:B:101:ILE:HD12	2.03	0.41
1:B:119:ARG:HD3	1:B:119:ARG:HA	1.81	0.40
1:A:120:ILE:HG23	1:A:123:ARG:HH21	1.85	0.40
1:B:201:PRO:HB2	1:B:204:GLU:HG3	2.03	0.40
1:A:166:LYS:O	1:A:169:VAL:HG12	2.22	0.40
1:B:65:ILE:HG21	1:B:65:ILE:HD13	1.80	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	212/214 (99%)	198 (93%)	9 (4%)	5 (2%)	6 1
1	B	212/214 (99%)	202 (95%)	8 (4%)	2 (1%)	17 6
All	All	424/428 (99%)	400 (94%)	17 (4%)	7 (2%)	9 2

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	10	VAL
1	B	10	VAL
1	A	113	ASP
1	A	201	PRO
1	A	75	GLU
1	A	127	ALA
1	B	201	PRO

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	175/176 (99%)	158 (90%)	17 (10%)	8 1
1	B	175/176 (99%)	156 (89%)	19 (11%)	6 1
All	All	350/352 (99%)	314 (90%)	36 (10%)	7 1

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	70	GLU
1	A	75	GLU
1	A	88	ARG
1	A	89	THR
1	A	91	PRO
1	A	97	LYS
1	A	111	VAL
1	A	113	ASP
1	A	123	ARG
1	A	157	LYS
1	A	158	ASP
1	A	165	ARG
1	A	175	THR
1	A	192	LYS
1	A	195	LYS
1	A	196	VAL

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Mol	Chain	Res	Type
1	A	201	PRO
1	B	31	THR
1	B	62	GLU
1	B	67	LEU
1	B	69	LYS
1	B	70	GLU
1	B	75	GLU
1	B	89	THR
1	B	91	PRO
1	B	111	VAL
1	B	123	ARG
1	B	154	THR
1	B	157	LYS
1	B	162	GLU
1	B	165	ARG
1	B	177	PRO
1	B	195	LYS
1	B	196	VAL
1	B	201	PRO
1	B	213	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	74	GLN
1	A	92	GLN
1	A	160	GLN
1	A	172	HIS
1	B	92	GLN
1	B	172	HIS

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)

2 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	$\#\|Z\| > 2$
2	AP5	A	215	-	48,62,62	1.02	2 (4%)	51,98,98	1.45	6 (11%)
2	AP5	B	215	-	48,62,62	1.14	3 (6%)	51,98,98	1.52	7 (13%)

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	A	215	-	-	2/36/76/76	0/6/6/6
2	AP5	B	215	-	-	2/36/76/76	0/6/6/6

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	215	AP5	C8A-N7A	-2.47	1.30	1.34
2	B	215	AP5	C8B-N7B	-2.40	1.30	1.34
2	B	215	AP5	C2J-C1J	-2.26	1.50	1.53
2	A	215	AP5	C2A-N3A	2.25	1.35	1.32
2	A	215	AP5	C8B-N7B	-2.24	1.30	1.34

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	215	AP5	C3J-C2J-C1J	-5.60	92.55	100.98

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	215	AP5	C3J-C2J-C1J	-4.42	94.32	100.98
2	A	215	AP5	O4J-C1J-C2J	4.32	113.23	106.93
2	B	215	AP5	O4J-C1J-C2J	3.45	111.96	106.93
2	B	215	AP5	N3A-C2A-N1A	2.83	133.11	128.68
2	B	215	AP5	O5J-PE-O1E	-2.39	99.73	109.07
2	B	215	AP5	O5J-C5J-C4J	2.36	117.13	108.99
2	A	215	AP5	C4A-C5A-N7A	-2.31	106.99	109.40
2	A	215	AP5	PB-O3A-PA	-2.24	125.15	132.83
2	B	215	AP5	C5B-C6B-N1B	-2.23	115.30	120.35
2	A	215	AP5	N3A-C2A-N1A	2.15	132.05	128.68
2	A	215	AP5	O3F-C3F-C4F	-2.09	105.00	111.05
2	B	215	AP5	C1F-N9A-C4A	2.05	130.24	126.64

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	215	AP5	PD-O3G-PG-O1G
2	A	215	AP5	C5F-O5F-PA-O3A
2	B	215	AP5	PB-O3B-PG-O2G
2	A	215	AP5	C5F-O5F-PA-O1A

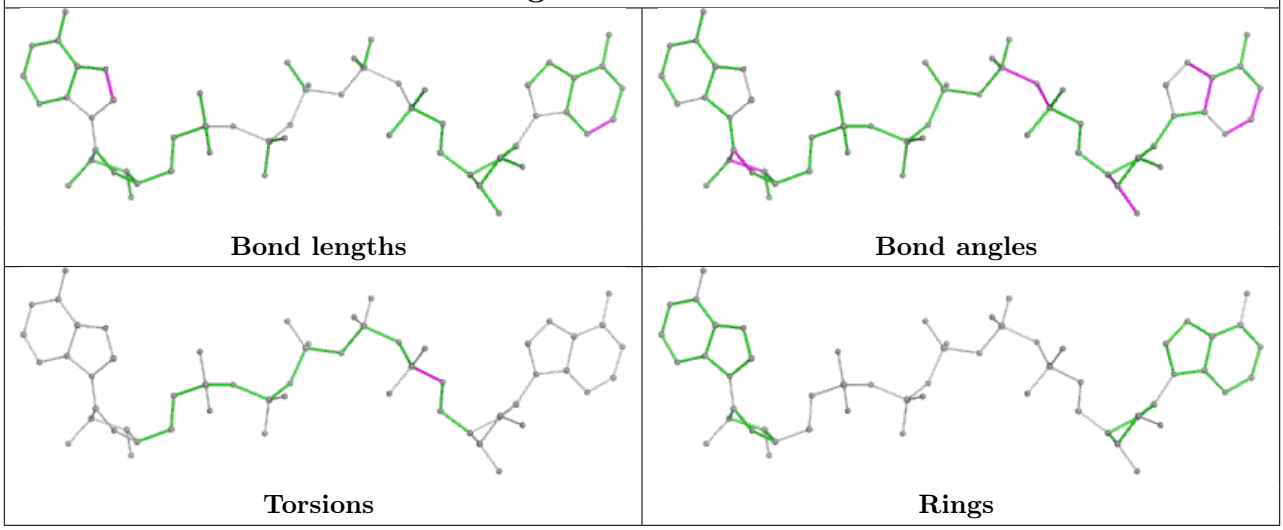
There are no ring outliers.

2 monomers are involved in 14 short contacts:

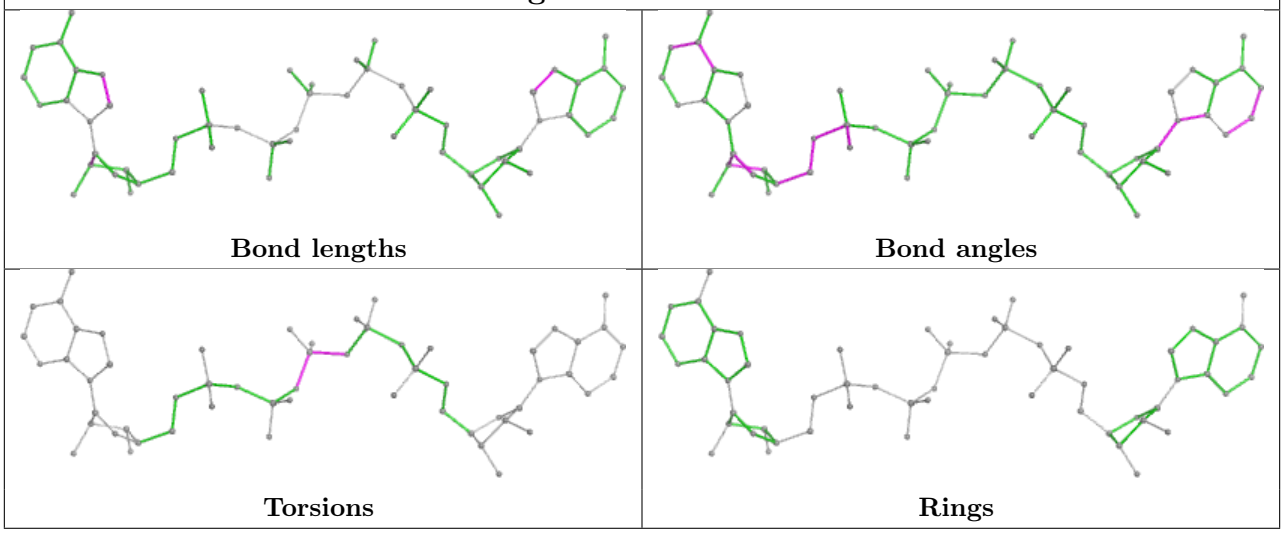
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	215	AP5	8	0
2	B	215	AP5	6	0

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 A 215



Ligand AP5 B 215



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 (i)

6.1 Protein, DNA and RNA chains (i)

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	214/214 (100%)	1.19	35 (16%)	1	2	2, 13, 24, 35	0
1	B	214/214 (100%)	1.22	36 (16%)	1	1	2, 13, 24, 38	0
All	All	428/428 (100%)	1.20	71 (16%)	1	1	2, 13, 24, 38	0

All (71) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	214	GLY	4.6
1	B	209	LEU	4.2
1	A	19	PHE	4.2
1	A	83	LEU	4.0
1	B	39	VAL	3.8
1	A	120	ILE	3.8
1	A	213	LEU	3.7
1	B	19	PHE	3.5
1	B	24	TYR	3.4
1	A	193	TYR	3.4
1	B	35	LEU	3.4
1	A	35	LEU	3.3
1	B	105	TYR	3.3
1	A	67	LEU	3.3
1	A	72	ILE	3.2
1	A	111	VAL	3.2
1	B	213	LEU	3.2
1	B	205	VAL	3.2
1	B	115	LEU	3.1
1	A	214	GLY	3.1
1	A	45	LEU	3.1
1	B	137	PHE	3.0
1	B	9	LEU	3.0
1	A	3	ILE	2.9

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Mol	Chain	Res	Type	RSRZ
1	B	20	ILE	2.9
1	B	168	LEU	2.9
1	A	144	GLY	2.9
1	B	153	LEU	2.8
1	A	77	CYS	2.8
1	A	109	PHE	2.8
1	B	193	TYR	2.8
1	B	178	LEU	2.7
1	A	209	LEU	2.7
1	A	137	PHE	2.7
1	A	105	TYR	2.7
1	A	78	ARG	2.6
1	A	165	ARG	2.6
1	A	59	VAL	2.6
1	B	109	PHE	2.6
1	B	169	VAL	2.6
1	B	72	ILE	2.6
1	A	133	TYR	2.5
1	B	181	TYR	2.5
1	A	86	PHE	2.5
1	A	101	ILE	2.5
1	A	90	ILE	2.4
1	B	86	PHE	2.4
1	A	24	TYR	2.4
1	B	31	THR	2.4
1	B	142	VAL	2.4
1	B	140	PRO	2.4
1	B	77	CYS	2.3
1	B	135	VAL	2.3
1	B	37	ALA	2.3
1	A	153	LEU	2.3
1	B	90	ILE	2.2
1	B	160	GLN	2.2
1	A	178	LEU	2.2
1	B	196	VAL	2.2
1	A	82	LEU	2.1
1	A	168	LEU	2.1
1	A	136	LYS	2.1
1	B	107	LEU	2.1
1	A	81	PHE	2.1
1	A	115	LEU	2.1
1	A	206	ARG	2.0

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Mol	Chain	Res	Type	RSRZ
1	B	106	VAL	2.0
1	A	34	MET	2.0
1	B	55	ALA	2.0
1	B	81	PHE	2.0
1	B	212	ILE	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)

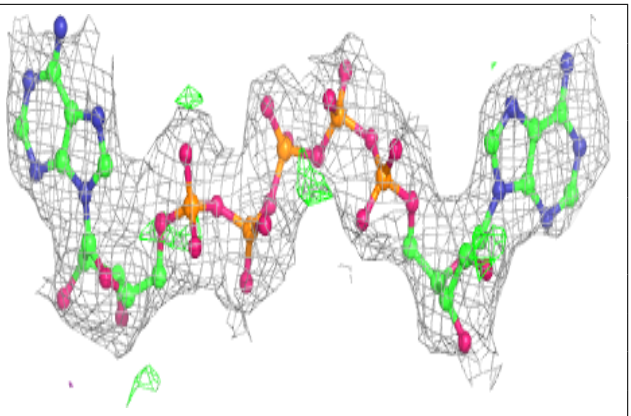
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

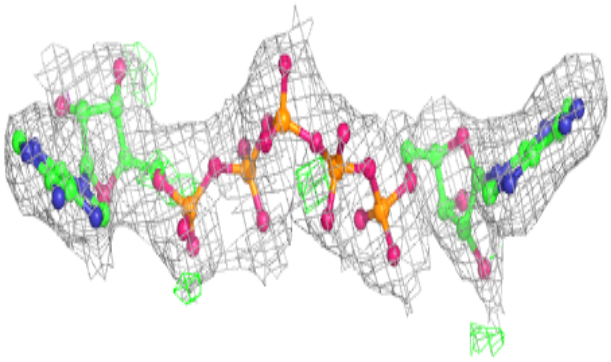
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	AP5	A	215	57/57	0.97	0.15	3,7,15,15	0
2	AP5	B	215	57/57	0.98	0.14	6,8,9,9	0

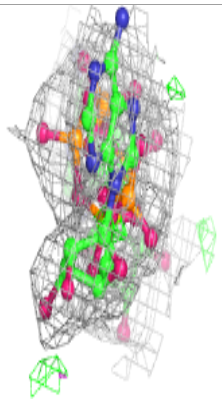
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around AP5 A 215:

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

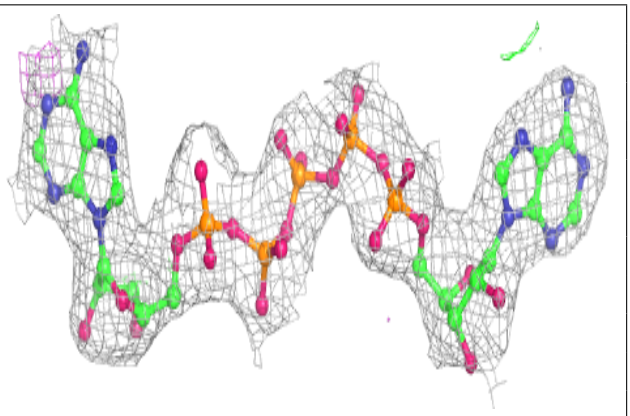


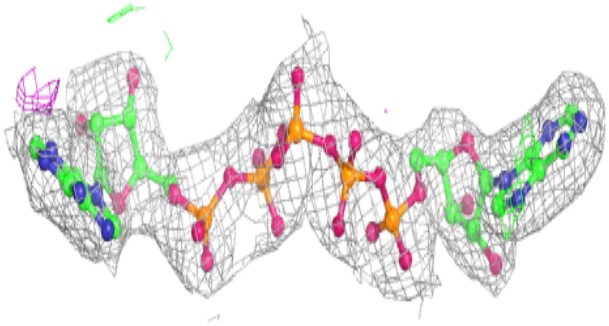


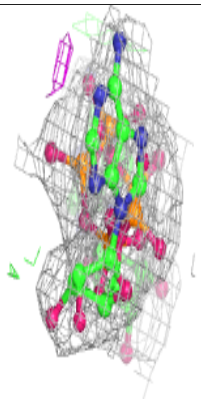


Electron density around AP5 B 215:

$2mF_o$-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.