



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 10:02 AM UTC

PDB ID : 1E4Y / pdb_00001e4y
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 ⓘ 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

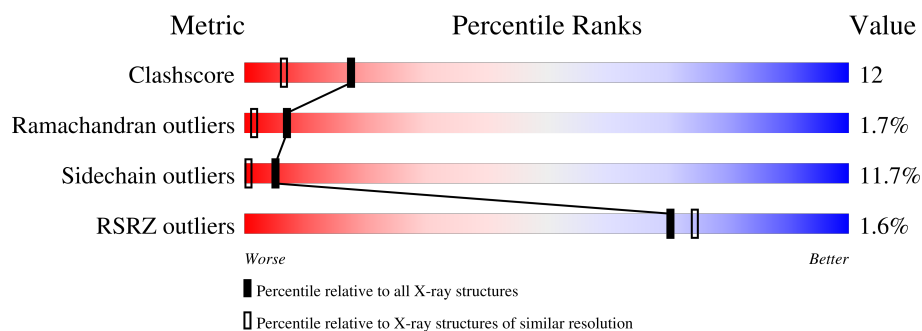
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 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	190562	3579 (1.86-1.86)
Ramachandran outliers	187476	3553 (1.86-1.86)
Sidechain outliers	187428	3553 (1.86-1.86)
RSRZ outliers	180081	3429 (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 $\leq 5\%$. 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	214	
1	B	214	

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.

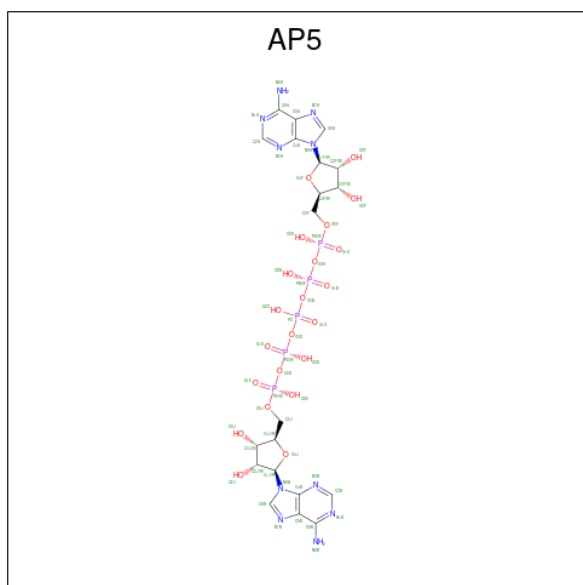
- Molecule 1 is a protein called Adenylate kinase.

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

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 (CCD ID: 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		

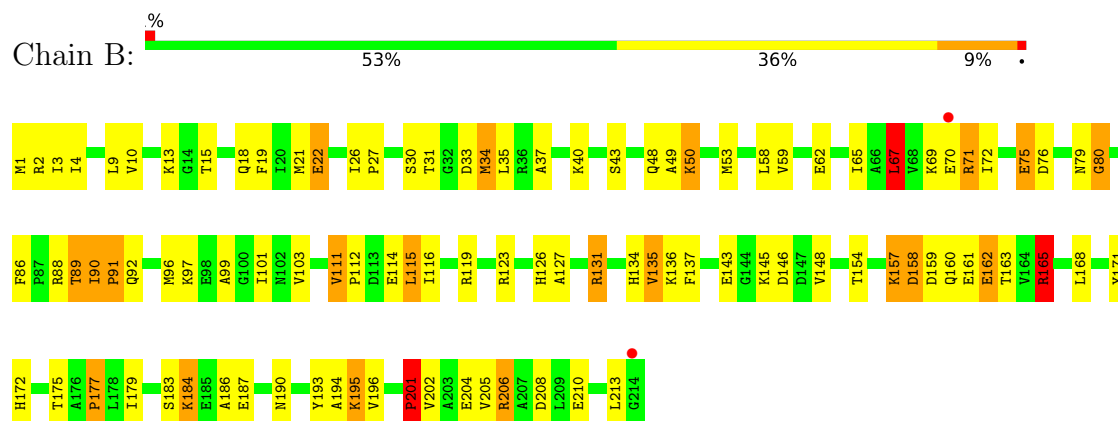
3 Residue-property plots [i](#)

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



• Molecule 1: Adenylate kinase



4 Data and refinement statistics

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 10.00 – 3.42	Depositor EDS
% Data completeness (in resolution range)	90.0 (10.00-1.85) 95.6 (10.00-3.42)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$	-	Xtriage
Refinement program	X-PLOR 1.5	Depositor
R, R_{free}	0.178 , (Not available) 0.166 , (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)	0.25 , 58.3	EDS
L-test for twinning ¹	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	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.*

¹Theoretical 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: 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	1.29	6/1680 (0.4%)	2.15	72/2261 (3.2%)
1	B	1.30	6/1680 (0.4%)	2.11	62/2261 (2.7%)
All	All	1.29	12/3360 (0.4%)	2.13	134/4522 (3.0%)

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 (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	134	HIS	CD2-NE2	-7.99	1.29	1.37
1	A	52	ILE	CA-CB	-7.98	1.44	1.54
1	A	134	HIS	CD2-NE2	-7.10	1.30	1.37
1	B	126	HIS	CD2-NE2	-7.03	1.30	1.37
1	A	126	HIS	CD2-NE2	-6.64	1.30	1.37
1	B	177	PRO	CA-CB	-6.25	1.45	1.53
1	A	108	GLU	CA-CB	-6.08	1.45	1.53
1	B	126	HIS	CG-ND1	-5.94	1.31	1.38
1	A	175	THR	CA-CB	5.16	1.61	1.53
1	B	206	ARG	NE-CZ	5.04	1.38	1.33
1	B	131	ARG	CA-CB	-5.02	1.46	1.53
1	A	117	VAL	CA-CB	5.02	1.61	1.54

All (134) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	79	ASN	OD1-CG-ND2	-11.59	111.01	122.60
1	B	159	ASP	CA-CB-CG	10.35	122.95	112.60
1	B	50	LYS	CA-CB-CG	-9.97	94.16	114.10
1	B	134	HIS	CA-CB-CG	-9.87	103.93	113.80
1	A	49	ALA	N-CA-C	8.75	120.82	111.28
1	A	63	LEU	N-CA-C	8.57	122.23	111.69
1	A	34	MET	CG-SD-CE	-8.55	82.09	100.90
1	A	118	ASP	CA-CB-CG	8.51	121.11	112.60
1	B	90	ILE	CB-CA-C	-8.37	105.48	114.35
1	A	160	GLN	OE1-CD-NE2	-8.27	114.33	122.60
1	B	134	HIS	CB-CG-CD2	-8.24	120.49	131.20
1	A	102	ASN	OD1-CG-ND2	-8.11	114.49	122.60
1	B	48	GLN	CB-CA-C	-8.08	98.20	110.88
1	A	106	VAL	N-CA-C	-7.95	96.67	108.12
1	A	50	LYS	CA-CB-CG	-7.91	98.28	114.10
1	A	84	ASP	CA-CB-CG	7.87	120.47	112.60
1	B	115	LEU	N-CA-C	7.81	119.79	111.28
1	B	184	LYS	CA-CB-CG	7.76	129.62	114.10
1	A	75	GLU	CB-CG-CD	7.62	125.56	112.60
1	A	104	ASP	N-CA-C	7.38	120.77	111.69
1	B	22	GLU	CB-CA-C	-7.37	98.42	110.79
1	B	80	GLY	N-CA-C	-7.22	104.78	112.08
1	B	162	GLU	CB-CG-CD	7.17	124.80	112.60
1	A	127	ALA	CA-C-O	-7.14	110.38	120.16
1	B	184	LYS	CB-CA-C	-7.01	98.76	110.68
1	B	15	THR	OG1-CB-CG2	-6.97	95.36	109.30
1	A	50	LYS	N-CA-C	6.97	118.88	111.28
1	B	53	MET	O-C-N	6.92	129.20	122.07
1	B	208	ASP	CA-CB-CG	6.89	119.49	112.60
1	B	190	ASN	N-CA-C	6.87	121.41	112.89
1	A	203	ALA	N-CA-C	6.87	118.42	111.07
1	B	79	ASN	N-CA-C	6.84	120.93	112.59
1	A	76	ASP	CA-CB-CG	6.81	119.41	112.60
1	B	75	GLU	CB-CG-CD	6.80	124.16	112.60
1	A	116	ILE	N-CA-C	6.77	117.52	110.62
1	A	111	VAL	N-CA-CB	-6.72	101.80	111.21
1	A	208	ASP	CA-CB-CG	6.63	119.23	112.60
1	A	162	GLU	CB-CG-CD	6.61	123.83	112.60
1	A	50	LYS	CB-CG-CD	6.60	126.49	111.30
1	B	67	LEU	N-CA-C	-6.57	104.20	111.36
1	A	190	ASN	N-CA-C	6.54	120.99	112.89
1	B	186	ALA	N-CA-C	-6.53	104.09	111.07
1	A	63	LEU	CB-CA-C	-6.41	98.98	110.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	48	GLN	CB-CA-C	-6.38	101.14	110.96
1	B	134	HIS	CB-CG-ND1	6.34	132.21	122.70
1	A	126	HIS	N-CA-C	-6.32	96.87	107.99
1	A	132	VAL	N-CA-CB	-6.31	103.42	110.99
1	B	157	LYS	N-CA-C	-6.28	104.67	112.90
1	B	43	SER	CA-CB-OG	6.26	123.63	111.10
1	B	183	SER	CA-C-O	-6.26	113.86	120.55
1	A	142	VAL	N-CA-C	-6.21	99.41	108.11
1	B	76	ASP	CA-CB-CG	6.20	118.80	112.60
1	B	111	VAL	CA-CB-CG2	-6.19	99.87	110.40
1	A	186	ALA	N-CA-C	-6.17	104.61	111.71
1	A	134	HIS	CB-CG-CD2	-6.17	123.18	131.20
1	A	74	GLN	OE1-CD-NE2	-6.16	116.44	122.60
1	B	90	ILE	N-CA-CB	6.15	114.37	110.50
1	A	13	LYS	CB-CA-C	-6.13	101.52	110.96
1	B	35	LEU	N-CA-C	-6.07	104.58	111.14
1	A	79	ASN	CB-CG-ND2	6.02	125.42	116.40
1	A	135	VAL	N-CA-C	6.01	116.75	110.62
1	B	50	LYS	N-CA-C	5.99	118.58	111.33
1	A	174	MET	O-C-N	5.96	129.18	122.32
1	A	126	HIS	O-C-N	5.94	130.23	122.68
1	B	79	ASN	CB-CG-ND2	5.90	125.25	116.40
1	B	86	PHE	CA-CB-CG	-5.86	107.94	113.80
1	A	164	VAL	N-CA-C	-5.85	104.22	110.36
1	B	143	GLU	N-CA-C	5.85	118.68	109.96
1	B	145	LYS	CA-C-O	5.84	126.90	120.71
1	A	114	GLU	CA-CB-CG	5.83	125.75	114.10
1	A	159	ASP	CA-CB-CG	5.79	118.39	112.60
1	B	22	GLU	N-CA-CB	5.79	118.70	110.13
1	A	132	VAL	N-CA-C	5.71	116.52	108.36
1	A	90	ILE	CB-CA-C	-5.68	108.32	114.35
1	A	125	VAL	O-C-N	-5.66	117.08	123.20
1	A	160	GLN	CG-CD-NE2	5.64	124.85	116.40
1	B	34	MET	CG-SD-CE	-5.62	88.53	100.90
1	B	111	VAL	CA-CB-CG1	5.59	119.90	110.40
1	A	161	GLU	N-CA-C	5.58	120.58	111.37
1	B	195	LYS	N-CA-CB	-5.56	101.22	110.23
1	B	131	ARG	N-CA-C	-5.55	100.24	109.46
1	B	89	THR	CA-C-N	5.55	123.74	120.24
1	B	89	THR	C-N-CA	5.55	123.74	120.24
1	B	114	GLU	CA-CB-CG	5.52	125.15	114.10
1	A	131	ARG	CA-CB-CG	-5.52	103.07	114.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	97	LYS	CG-CD-CE	-5.50	98.64	111.30
1	B	49	ALA	CA-C-O	-5.47	115.08	120.82
1	A	102	ASN	CB-CG-ND2	5.44	124.56	116.40
1	A	26	ILE	CA-C-N	5.44	125.70	119.83
1	A	26	ILE	C-N-CA	5.44	125.70	119.83
1	A	128	PRO	N-CA-C	5.43	123.66	112.47
1	B	184	LYS	CG-CD-CE	5.43	123.80	111.30
1	B	127	ALA	CA-C-O	-5.41	112.74	120.16
1	A	86	PHE	CA-C-O	-5.40	114.67	119.86
1	B	158	ASP	CA-CB-CG	5.40	118.00	112.60
1	B	48	GLN	N-CA-CB	5.40	117.84	110.01
1	B	146	ASP	N-CA-CB	-5.40	102.39	110.17
1	A	22	GLU	CB-CA-C	-5.39	102.38	110.90
1	B	135	VAL	CB-CA-C	-5.39	102.44	111.29
1	A	184	LYS	CB-CA-C	-5.39	101.51	110.68
1	B	72	ILE	N-CA-C	-5.36	104.51	110.62
1	B	127	ALA	CA-C-N	5.31	125.38	119.32
1	B	127	ALA	C-N-CA	5.31	125.38	119.32
1	B	13	LYS	CB-CA-C	-5.29	102.57	110.88
1	A	172	HIS	CB-CG-CD2	-5.29	124.32	131.20
1	A	163	THR	CA-C-O	-5.27	114.84	120.42
1	B	184	LYS	N-CA-CB	5.25	117.93	110.06
1	A	159	ASP	N-CA-C	-5.24	107.42	112.97
1	A	148	VAL	N-CA-C	5.22	116.63	111.67
1	A	152	GLU	CB-CG-CD	5.21	121.45	112.60
1	B	194	ALA	CA-C-O	5.20	126.94	120.54
1	A	169	VAL	CA-C-N	5.19	127.55	120.54
1	A	169	VAL	C-N-CA	5.19	127.55	120.54
1	B	13	LYS	CB-CG-CD	-5.19	99.36	111.30
1	A	86	PHE	O-C-N	-5.17	116.42	121.38
1	A	169	VAL	CA-CB-CG1	5.17	119.19	110.40
1	A	154	THR	N-CA-C	5.14	117.09	109.69
1	A	101	ILE	CA-C-O	5.14	126.70	120.74
1	B	50	LYS	CB-CG-CD	5.14	123.12	111.30
1	A	66	ALA	CA-C-N	5.13	127.42	120.65
1	A	66	ALA	C-N-CA	5.13	127.42	120.65
1	A	179	ILE	CB-CG1-CD1	-5.12	103.05	113.80
1	B	165	ARG	CB-CG-CD	5.11	123.06	111.30
1	B	116	ILE	CA-C-O	-5.11	115.95	121.27
1	A	134	HIS	N-CA-C	-5.09	100.15	108.76
1	A	207	ALA	CA-C-O	-5.07	115.47	120.70
1	B	71	ARG	NE-CZ-NH2	-5.07	114.63	119.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	67	LEU	N-CA-C	-5.07	105.44	110.97
1	A	134	HIS	CA-CB-CG	-5.04	108.75	113.80
1	A	126	HIS	CB-CG-CD2	-5.04	124.65	131.20
1	A	50	LYS	CG-CD-CE	-5.03	99.73	111.30
1	B	33	ASP	CB-CA-C	-5.03	102.44	110.79
1	B	160	GLN	CG-CD-NE2	5.01	123.91	116.40
1	A	13	LYS	N-CA-CB	5.00	117.22	109.91

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 ⓘ

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	45	0
1	B	1657	0	1687	41	0
2	A	57	0	22	8	0
2	B	57	0	22	6	0
All	All	3428	0	3418	85	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 12.

All (85) 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:A:4:ILE:HD11	1:A:96:MET:HE1	1.77	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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: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
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:A:126:HIS:HD1	1:A:129:SER:HG	1.52	0.58
1:B:123:ARG:HG3	2:B:215:AP5:H3F	1.86	0.58
1:B:34:MET:HE1	1:B:71:ARG:HG2	1.84	0.57
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:1:MET:HE3	1:B:3:ILE:HD11	1.88	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:53:MET:HB3	1:A:53:MET:HE2	1.77	0.44
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
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:36:ARG:HG3	1:A:53:MET:HE1	2.02	0.42
1:A:36:ARG:CG	1:A:53:MET:HE1	2.50	0.42
1:B:4:ILE:HD11	1:B:96:MET:HE1	2.00	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%)	4	1
1	B	212/214 (99%)	202 (95%)	8 (4%)	2 (1%)	14	4
All	All	424/428 (99%)	400 (94%)	17 (4%)	7 (2%)	7	1

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%)	154 (88%)	21 (12%)	5	0
1	B	175/176 (99%)	155 (89%)	20 (11%)	5	1
All	All	350/352 (99%)	309 (88%)	41 (12%)	5	1

All (41) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	23	LYS
1	A	31	THR
1	A	62	GLU
1	A	70	GLU
1	A	75	GLU
1	A	89	THR
1	A	91	PRO
1	A	97	LYS
1	A	111	VAL
1	A	123	ARG
1	A	157	LYS
1	A	158	ASP
1	A	162	GLU
1	A	165	ARG
1	A	175	THR
1	A	184	LYS
1	A	192	LYS
1	A	195	LYS
1	A	196	VAL
1	A	201	PRO
1	A	213	LEU
1	B	22	GLU
1	B	31	THR
1	B	50	LYS
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	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 oligosaccharides 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	B	215	-	62,62,62	1.17	3 (4%)	89,98,98	2.06	22 (24%)
2	AP5	A	215	-	62,62,62	1.12	3 (4%)	89,98,98	1.98	21 (23%)

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	215	-	-	5/44/76/76	0/6/6/6
2	AP5	A	215	-	-	2/44/76/76	0/6/6/6

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	215	AP5	PD-O3D	3.65	1.63	1.59
2	A	215	AP5	PD-O3G	3.44	1.63	1.59
2	A	215	AP5	C5A-C4A	3.23	1.44	1.39
2	B	215	AP5	PE-O3D	2.90	1.62	1.59
2	A	215	AP5	PD-O3D	2.35	1.62	1.59
2	B	215	AP5	PD-O3G	2.35	1.62	1.59

All (43) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	215	AP5	O4F-C1F-N9A	8.52	124.45	108.09
2	A	215	AP5	O4F-C1F-N9A	7.85	123.17	108.09
2	B	215	AP5	O4J-C1J-N9B	6.01	119.63	108.09
2	A	215	AP5	O4J-C1J-N9B	5.98	119.58	108.09
2	B	215	AP5	C2J-C1J-N9B	4.74	125.07	113.30
2	B	215	AP5	C3J-C2J-C1J	-4.71	92.55	101.46
2	A	215	AP5	O3A-PB-O1B	-4.38	97.54	110.70
2	A	215	AP5	C2J-C1J-N9B	4.23	123.82	113.30
2	A	215	AP5	C2F-C1F-N9A	4.19	123.72	113.30
2	B	215	AP5	O2B-PB-O3B	4.01	118.11	107.27
2	A	215	AP5	C3J-C2J-C1J	-3.78	94.32	101.46
2	B	215	AP5	O3G-PD-O1D	-3.74	99.45	110.70
2	A	215	AP5	C5A-C4A-N3A	-3.66	121.67	126.72
2	B	215	AP5	O2G-PG-O3B	3.61	117.03	107.27
2	A	215	AP5	O3G-PD-O1D	-3.60	99.87	110.70
2	B	215	AP5	O3A-PB-O1B	-3.37	100.57	110.70
2	A	215	AP5	C4A-C5A-N7A	-3.14	106.99	110.58
2	A	215	AP5	C4J-O4J-C1J	-3.11	102.61	109.47
2	B	215	AP5	C5A-C4A-N3A	-3.10	122.45	126.72
2	A	215	AP5	O4J-C1J-C2J	3.09	113.23	106.62
2	B	215	AP5	C4J-O4J-C1J	-3.09	102.65	109.47
2	B	215	AP5	O3B-PG-O1G	-3.07	101.46	110.70
2	B	215	AP5	N3A-C2A-N1A	2.99	133.11	128.58
2	A	215	AP5	O2B-PB-O3B	2.71	114.61	107.27
2	B	215	AP5	C6A-C5A-C4A	2.67	120.83	117.18
2	B	215	AP5	C2F-C1F-N9A	2.67	119.94	113.30
2	B	215	AP5	O2A-PA-O3A	2.61	114.34	107.27

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	215	AP5	O3D-PD-O1D	-2.60	102.87	110.70
2	A	215	AP5	O2D-PD-O3D	2.52	114.09	107.27
2	B	215	AP5	O2D-PD-O3D	2.50	114.04	107.27
2	B	215	AP5	O4J-C1J-C2J	2.50	111.96	106.62
2	A	215	AP5	N3A-C4A-N9A	2.42	131.29	127.17
2	B	215	AP5	N3A-C4A-N9A	2.41	131.26	127.17
2	B	215	AP5	O5J-C5J-C4J	2.39	117.13	108.99
2	B	215	AP5	O5J-PE-O1E	-2.32	99.73	108.94
2	A	215	AP5	N3A-C2A-N1A	2.29	132.05	128.58
2	A	215	AP5	O2D-PD-O3G	2.29	113.45	107.27
2	B	215	AP5	C4A-C5A-N7A	-2.18	108.09	110.58
2	A	215	AP5	O3F-C3F-C4F	-2.12	105.00	111.08
2	A	215	AP5	O2A-PA-O3A	2.11	112.99	107.27
2	B	215	AP5	N3B-C4B-N9B	2.10	130.74	127.17
2	A	215	AP5	O2G-PG-O3G	-2.09	101.62	107.27
2	A	215	AP5	C5A-N7A-C8A	2.08	106.72	103.45

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	215	AP5	PD-O3G-PG-O1G
2	B	215	AP5	C2J-C1J-N9B-C4B
2	B	215	AP5	PB-O3B-PG-O2G
2	B	215	AP5	O4F-C1F-N9A-C8A
2	A	215	AP5	O4F-C1F-N9A-C8A
2	A	215	AP5	PD-O3D-PE-O2E
2	B	215	AP5	PD-O3D-PE-O2E

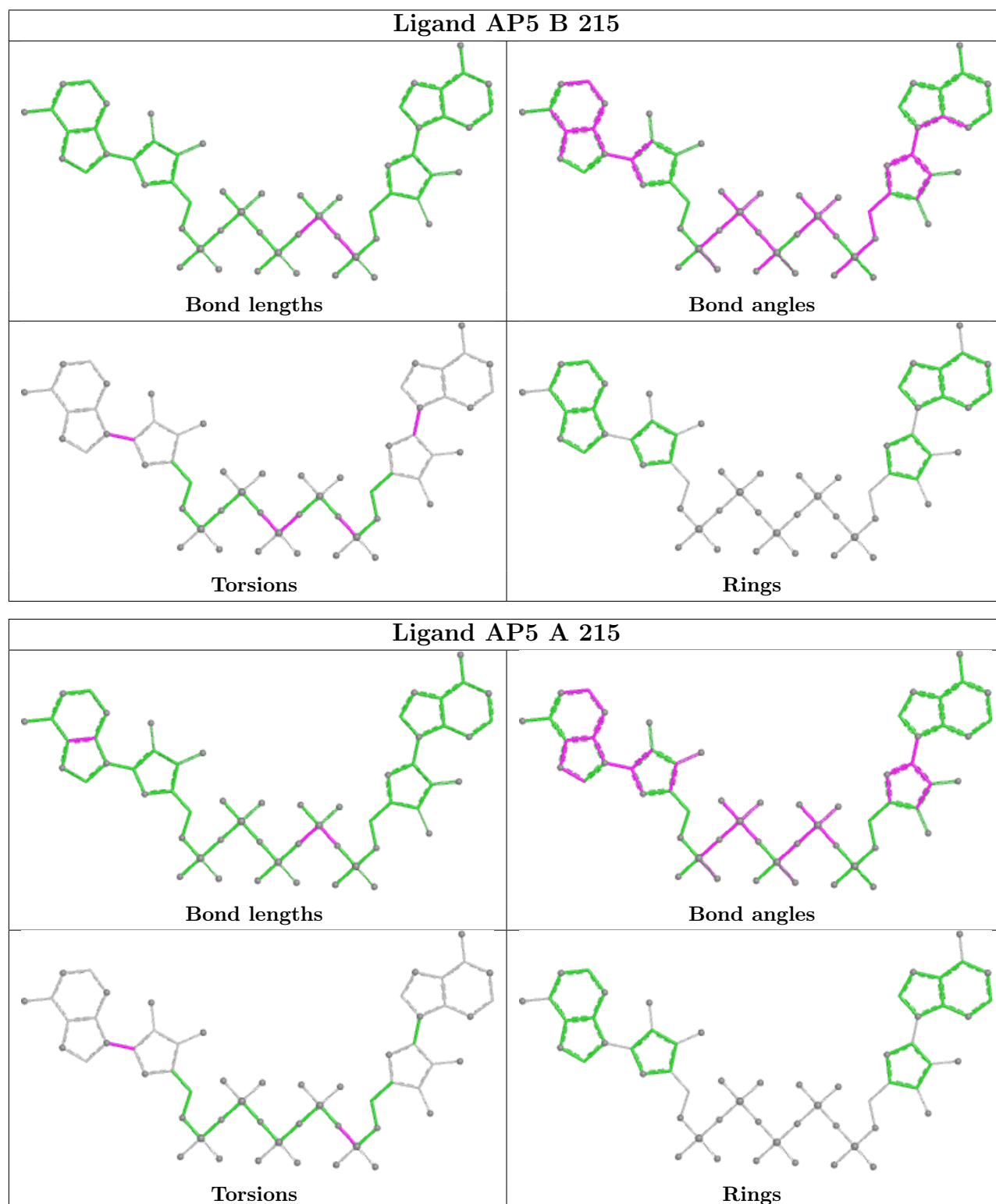
There are no ring outliers.

2 monomers are involved in 14 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	215	AP5	6	0
2	A	215	AP5	8	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 [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%)	-0.00	5 (2%) 61 64	2, 13, 24, 35	0
1	B	214/214 (100%)	0.01	2 (0%) 81 84	2, 13, 24, 38	0
All	All	428/428 (100%)	0.01	7 (1%) 70 74	2, 13, 24, 38	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	19	PHE	3.4
1	A	78	ARG	3.0
1	B	214	GLY	2.5
1	B	70	GLU	2.3
1	A	214	GLY	2.2
1	A	169	VAL	2.2
1	A	162	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](#)

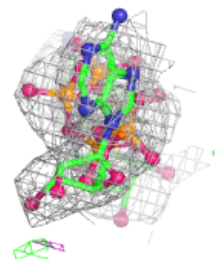
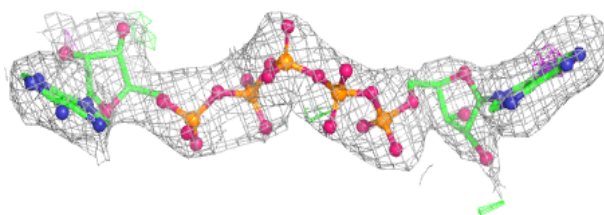
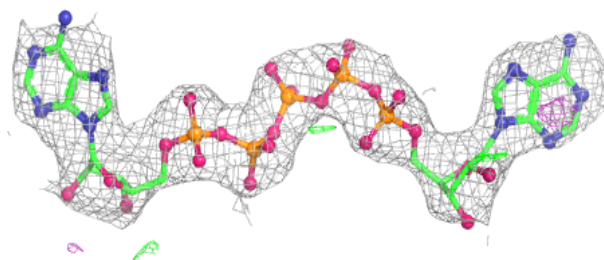
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q < 0.9’ lists 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	AP5	A	215	57/57	0.98	0.06	3,7,15,15	0
2	AP5	B	215	57/57	0.98	0.06	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.

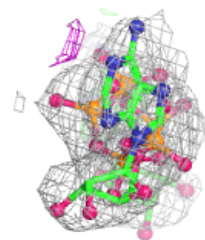
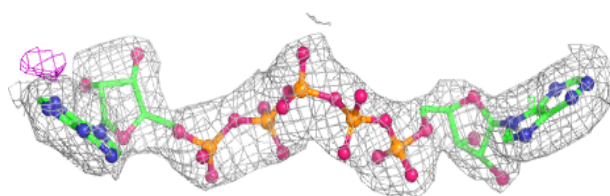
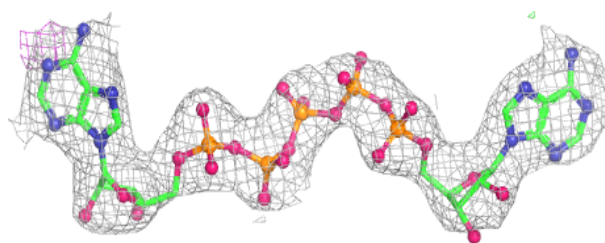
Electron density around AP5 A 215:

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.