



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 15, 2020 – 11:02 pm BST

PDB ID : 2RD5
Title : Structural basis for the regulation of N-acetylglutamate kinase by PII in *Arabidopsis thaliana*
Authors : Mizuno, Y.; Moorhead, G.B.G.; Ng, K.K.S.
Deposited on : 2007-09-21
Resolution : 2.51 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

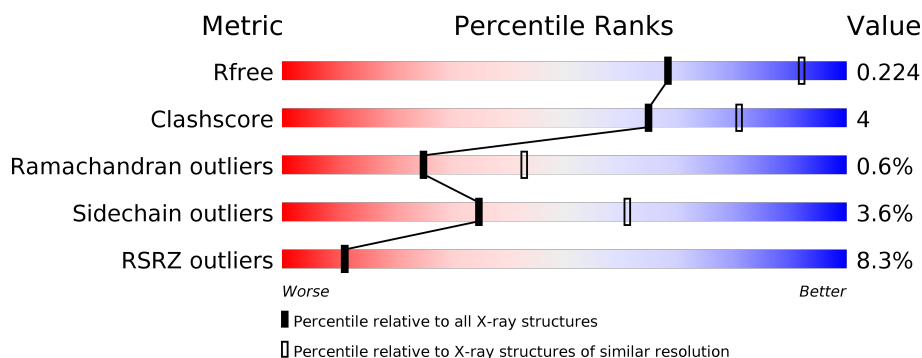
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.51 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	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)

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 $\leq 5\%$. 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	298	<div> <div>5%</div> <div> <div></div> <div>82%</div> <div>11%</div> <div>6%</div> </div> </div>
1	B	298	<div> <div>9%</div> <div> <div></div> <div>84%</div> <div>11%</div> <div>5%</div> </div> </div>
2	C	135	<div> <div>9%</div> <div> <div></div> <div>85%</div> <div>7%</div> <div>7%</div> </div> </div>
2	D	135	<div> <div>10%</div> <div> <div></div> <div>81%</div> <div>12%</div> <div>7%</div> </div> </div>

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 6328 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 Acetylglutamate kinase-like protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	281	Total	C	N	O	S	0	0	0
			2074	1309	360	395	10			
1	B	283	Total	C	N	O	S	0	0	0
			2087	1317	362	398	10			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	MET	-	INITIATING METHIONINE	UNP Q9SCL7
B	0	MET	-	INITIATING METHIONINE	UNP Q9SCL7

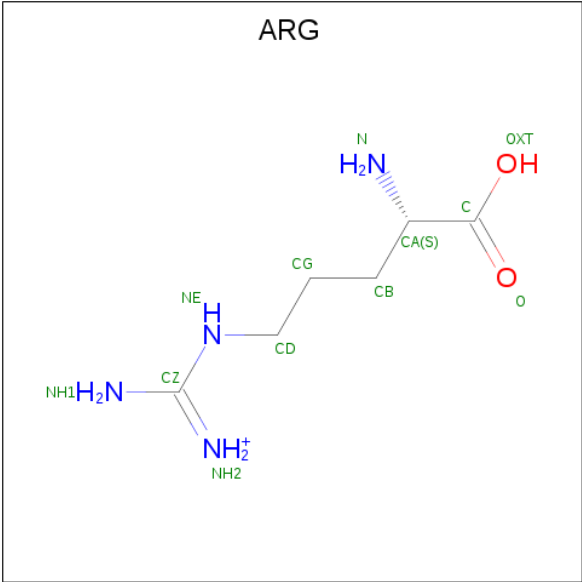
- Molecule 2 is a protein called PII protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	125	Total	C	N	O	S	0	0	0
			963	605	171	185	2			
2	D	126	Total	C	N	O	S	0	0	0
			971	610	172	186	3			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	0	MET	-	INITIATING METHIONINE	UNP Q9ZST4
D	0	MET	-	INITIATING METHIONINE	UNP Q9ZST4

- Molecule 3 is ARGinine (three-letter code: ARG) (formula: C₆H₁₅N₄O₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			12	6	4	2		
3	B	1	Total	C	N	O	0	0
			12	6	4	2		

- Molecule 4 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula: C₁₀H₁₅N₅O₁₀P₂).



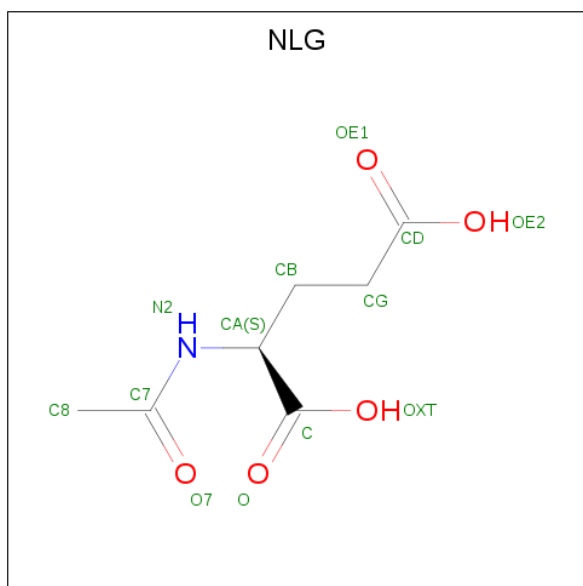
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			27	10	5	10	2		

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

- Molecule 5 is N-ACETYL-L-GLUTAMATE (three-letter code: NLG) (formula: $C_7H_{11}NO_5$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	0	0
			13	7	1	5		
5	B	1	Total	C	N	O	0	0
			13	7	1	5		

- Molecule 6 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	1	Total	Mg	0	0
			1	1		
6	D	1	Total	Mg	0	0
			1	1		
6	C	1	Total	Mg	0	0
			1	1		

- Molecule 7 is ADENOSINE-5'-TRIPHOSPHATE (three-letter code: ATP) (formula: $C_{10}H_{16}N_5O_{13}P_3$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
7	C	1	Total	C	N	O	P	0	0
			31	10	5	13	3		
7	D	1	Total	C	N	O	P	0	0
			31	10	5	13	3		

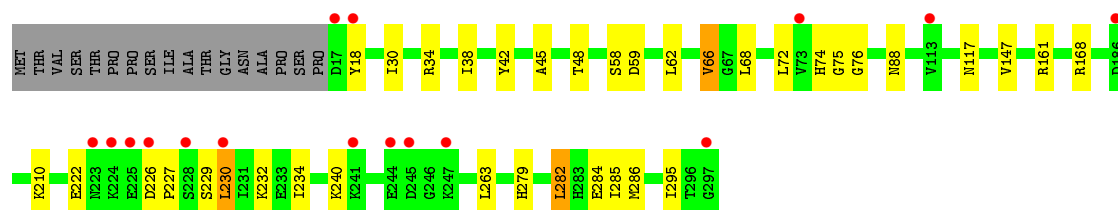
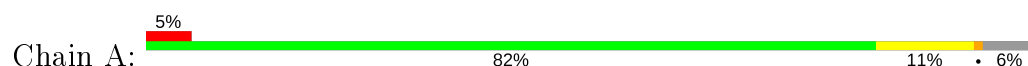
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	24	Total	O	0	0
			24	24		
8	B	25	Total	O	0	0
			25	25		
8	C	4	Total	O	0	0
			4	4		
8	D	11	Total	O	0	0
			11	11		

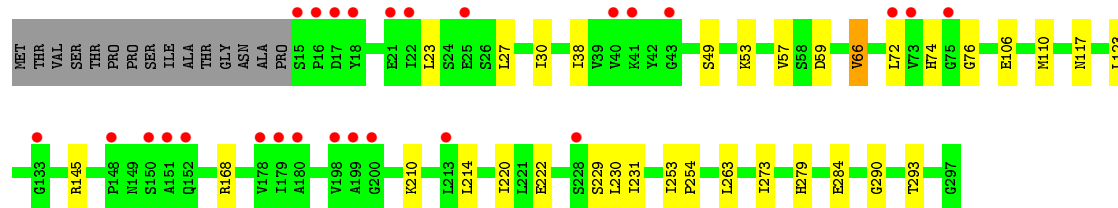
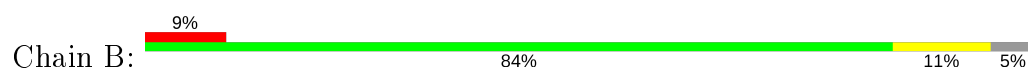
3 Residue-property plots [i](#)

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.

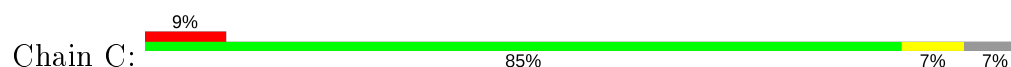
- Molecule 1: Acetylglutamate kinase-like protein



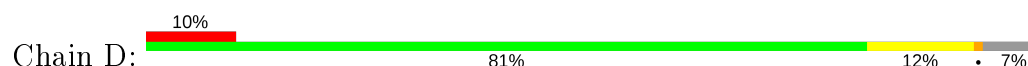
- Molecule 1: Acetylglutamate kinase-like protein



- Molecule 2: PII protein



- Molecule 2: PII protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 3	Depositor
Cell constants a, b, c, α , β , γ	171.13Å 171.13Å 171.13Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.00 – 2.51 32.93 – 2.51	Depositor EDS
% Data completeness (in resolution range)	99.9 (40.00-2.51) 99.9 (32.93-2.51)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	0.13	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.23 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.201 , 0.229 0.197 , 0.224	Depositor DCC
R_{free} test set	2907 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å ²)	49.6	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 54.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.031 for l,-k,h	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6328	wwPDB-VP
Average B, all atoms (Å ²)	67.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.31% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: MG, ATP, ADP, NLG

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	0.38	0/2099	0.55	1/2841 (0.0%)
1	B	0.38	0/2113	0.57	0/2861
2	C	0.41	0/975	0.55	0/1307
2	D	0.39	0/983	0.57	0/1317
All	All	0.39	0/6170	0.56	1/8326 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	230	LEU	CA-CB-CG	5.25	127.38	115.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2074	0	2168	20	0
1	B	2087	0	2180	14	0
2	C	963	0	985	6	1
2	D	971	0	994	10	0
3	A	12	0	12	2	0
3	B	12	0	12	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	27	0	12	0	0
4	B	27	0	12	0	0
5	A	13	0	8	0	0
5	B	13	0	8	0	0
6	B	1	0	0	0	0
6	C	1	0	0	0	0
6	D	1	0	0	0	0
7	C	31	0	12	4	1
7	D	31	0	12	5	0
8	A	24	0	0	0	0
8	B	25	0	0	0	0
8	C	4	0	0	1	0
8	D	11	0	0	2	0
All	All	6328	0	6415	52	1

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 4.

All (52) 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)
2:C:7:ILE:HG23	2:C:8:PRO:HD3	1.40	1.03
2:D:48:GLY:HA3	7:D:1000:ATP:O1A	1.66	0.94
1:A:74:HIS:HE1	1:A:117:ASN:HD22	1.16	0.92
2:C:48:GLY:HA3	7:C:1000:ATP:O1A	1.70	0.91
1:A:38:ILE:HD12	1:A:68:LEU:HD23	1.55	0.87
1:B:74:HIS:HE1	1:B:117:ASN:HD22	1.22	0.81
1:B:30:ILE:HG23	1:B:66:VAL:HG13	1.70	0.72
1:A:30:ILE:HG23	1:A:66:VAL:HG13	1.73	0.70
1:B:106:GLU:O	1:B:110:MET:HG2	1.93	0.67
1:B:222:GLU:HB2	1:B:231:ILE:HD11	1.77	0.66
2:D:99:GLY:HA2	7:D:1000:ATP:O2A	1.96	0.65
2:C:99:GLY:HA2	7:C:1000:ATP:O2A	1.98	0.63
1:A:285:ILE:HG22	1:A:286:MET:CE	2.30	0.61
1:A:74:HIS:CE1	1:A:117:ASN:HD22	2.08	0.59
2:C:7:ILE:CG2	2:C:8:PRO:HD3	2.24	0.57
2:D:77:VAL:HB	2:D:81:GLN:HG3	1.88	0.55
1:A:38:ILE:HG12	1:A:210:LYS:HB2	1.90	0.53
1:B:59:ASP:OD1	1:B:279:HIS:HD2	1.91	0.53
2:D:49:ALA:CB	2:D:64:GLU:HA	2.38	0.53
1:B:222:GLU:HB3	1:B:229:SER:HB2	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:285:ILE:HG22	1:A:286:MET:HE3	1.91	0.53
1:A:232:LYS:O	3:A:1000:ARG:NH2	2.44	0.51
2:D:102:LYS:HG3	7:D:1000:ATP:C8	2.46	0.51
1:A:285:ILE:HG22	1:A:286:MET:HE2	1.93	0.51
1:A:42:TYR:HB2	1:A:72:LEU:HD21	1.92	0.50
1:A:222:GLU:HB3	1:A:229:SER:HB2	1.93	0.49
1:B:253:ILE:HB	1:B:254:PRO:HD3	1.94	0.49
1:A:62:LEU:HD23	1:A:282:LEU:HD21	1.96	0.48
1:A:59:ASP:OD1	1:A:279:HIS:HD2	1.97	0.48
1:A:45:ALA:HA	1:A:48:THR:HG22	1.97	0.47
1:B:214:LEU:HD22	1:B:273:ILE:HD11	1.96	0.47
2:C:35:GLY:O	2:C:129:ASP:HB2	2.14	0.47
1:B:53:LYS:O	1:B:57:VAL:HG13	2.16	0.45
1:A:62:LEU:HD22	1:A:282:LEU:HD11	1.98	0.45
1:B:38:ILE:HD12	1:B:210:LYS:HB2	1.97	0.45
1:B:220:ILE:HD12	1:B:293:THR:HG21	1.99	0.45
1:A:240:LYS:HE3	8:D:1004:HOH:O	2.17	0.44
2:D:118:GLU:O	2:D:123:ALA:HB2	2.18	0.44
2:D:46:GLY:HA3	7:D:1000:ATP:O2'	2.18	0.44
2:C:102:LYS:HD2	7:C:1000:ATP:N7	2.33	0.43
7:C:1000:ATP:H5'2	8:C:1005:HOH:O	2.19	0.43
7:D:1000:ATP:H5'2	8:D:1012:HOH:O	2.17	0.43
1:B:74:HIS:CE1	1:B:117:ASN:HD22	2.15	0.43
1:A:74:HIS:HD2	1:A:75:GLY:O	2.01	0.43
1:B:59:ASP:OD2	1:B:279:HIS:HA	2.19	0.42
1:A:234:ILE:O	1:A:295:ILE:HA	2.20	0.42
2:D:49:ALA:HB2	2:D:64:GLU:HA	2.02	0.41
2:D:20:ARG:HA	2:D:21:PRO:HD3	1.94	0.41
1:A:284:GLU:O	3:A:1000:ARG:HB3	2.21	0.41
1:B:284:GLU:HA	1:B:290:GLY:HA2	2.03	0.40
2:D:41:VAL:HA	2:D:72:LYS:O	2.21	0.40
1:A:226:ASP:HA	1:A:227:PRO:HD3	1.77	0.40

All (1) 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)
2:C:115:ARG:NH1	7:C:1000:ATP:O1G[9_555]	2.19	0.01

5.3 Torsion angles

5.3.1 Protein backbone

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	279/298 (94%)	274 (98%)	4 (1%)	1 (0%)	34	54
1	B	281/298 (94%)	275 (98%)	5 (2%)	1 (0%)	34	54
2	C	123/135 (91%)	117 (95%)	5 (4%)	1 (1%)	19	35
2	D	124/135 (92%)	118 (95%)	4 (3%)	2 (2%)	9	17
All	All	807/866 (93%)	784 (97%)	18 (2%)	5 (1%)	25	43

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	C	51	GLY
1	A	76	GLY
1	B	76	GLY
2	D	64	GLU
2	D	128	GLY

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	225/239 (94%)	214 (95%)	11 (5%)	25	47
1	B	227/239 (95%)	217 (96%)	10 (4%)	28	52
2	C	104/114 (91%)	103 (99%)	1 (1%)	76	90
2	D	105/114 (92%)	103 (98%)	2 (2%)	57	80
All	All	661/706 (94%)	637 (96%)	24 (4%)	35	61

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	18	TYR
1	A	34	ARG
1	A	58	SER
1	A	66	VAL
1	A	88	ASN
1	A	147	VAL
1	A	161	ARG
1	A	168	ARG
1	A	230	LEU
1	A	263	LEU
1	A	282	LEU
1	B	23	LEU
1	B	27	LEU
1	B	49	SER
1	B	66	VAL
1	B	72	LEU
1	B	123	LEU
1	B	145	ARG
1	B	168	ARG
1	B	230	LEU
1	B	263	LEU
2	C	31	LEU
2	D	32	LEU
2	D	80	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	74	HIS
1	A	119	ASN
1	A	278	GLN
1	A	279	HIS
1	B	74	HIS
1	B	152	GLN
1	B	279	HIS
2	C	81	GLN
2	C	87	ASN
2	D	87	ASN

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 11 ligands modelled in this entry, 3 are monoatomic - leaving 8 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$
7	ATP	C	1000	6	26,33,33	1.12	2 (7%)	31,52,52	1.60	7 (22%)
5	NLG	B	2001	-	6,12,12	4.80	2 (33%)	7,15,15	1.09	0
4	ADP	B	2000	6	24,29,29	1.00	2 (8%)	29,45,45	1.31	4 (13%)
7	ATP	D	1000	6	26,33,33	1.01	2 (7%)	31,52,52	1.46	6 (19%)
5	NLG	A	2001	-	6,12,12	4.98	2 (33%)	7,15,15	1.72	1 (14%)
4	ADP	A	2000	-	24,29,29	1.01	1 (4%)	29,45,45	1.37	4 (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
7	ATP	C	1000	6	-	9/18/38/38	0/3/3/3
5	NLG	B	2001	-	-	1/7/13/13	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	ADP	B	2000	6	-	5/12/32/32	0/3/3/3
7	ATP	D	1000	6	-	6/18/38/38	0/3/3/3
5	NLG	A	2001	-	-	3/7/13/13	-
4	ADP	A	2000	-	-	4/12/32/32	0/3/3/3

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	2001	NLG	C8-C7	-8.81	1.32	1.50
5	B	2001	NLG	C8-C7	-8.48	1.32	1.50
5	A	2001	NLG	CA-N2	-8.42	1.35	1.46
5	B	2001	NLG	CA-N2	-8.14	1.35	1.46
7	C	1000	ATP	O4'-C1'	3.11	1.45	1.41
4	A	2000	ADP	C5-C4	2.70	1.48	1.40
7	C	1000	ATP	C5-C4	2.62	1.47	1.40
4	B	2000	ADP	C5-C4	2.57	1.47	1.40
7	D	1000	ATP	C5-C4	2.47	1.47	1.40
7	D	1000	ATP	O4'-C1'	2.27	1.44	1.41
4	B	2000	ADP	O4'-C1'	2.10	1.44	1.41

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	2001	NLG	CA-N2-C7	-4.25	117.76	122.44
4	B	2000	ADP	N3-C2-N1	-3.56	123.12	128.68
7	C	1000	ATP	PA-O3A-PB	-3.45	120.97	132.83
7	C	1000	ATP	O5'-C5'-C4'	3.42	120.76	108.99
4	A	2000	ADP	N3-C2-N1	-3.29	123.54	128.68
7	C	1000	ATP	N3-C2-N1	-3.22	123.64	128.68
7	D	1000	ATP	N3-C2-N1	-3.15	123.75	128.68
4	A	2000	ADP	C3'-C2'-C1'	3.12	105.67	100.98
7	D	1000	ATP	PA-O3A-PB	-2.98	122.60	132.83
4	A	2000	ADP	PA-O3A-PB	-2.64	123.77	132.83
7	C	1000	ATP	O4'-C4'-C5'	2.63	118.03	109.37
7	D	1000	ATP	PB-O3B-PG	-2.61	123.86	132.83
7	D	1000	ATP	O5'-C5'-C4'	2.55	117.77	108.99
4	A	2000	ADP	C4-C5-N7	-2.50	106.80	109.40
7	D	1000	ATP	C4-C5-N7	-2.43	106.87	109.40
4	B	2000	ADP	C3'-C2'-C1'	2.40	104.59	100.98
7	D	1000	ATP	O4'-C4'-C5'	2.40	117.26	109.37
4	B	2000	ADP	PA-O3A-PB	-2.33	124.84	132.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	C	1000	ATP	C4-C5-N7	-2.28	107.03	109.40
4	B	2000	ADP	C4-C5-N7	-2.26	107.04	109.40
7	C	1000	ATP	PB-O3B-PG	-2.08	125.70	132.83
7	C	1000	ATP	O2A-PA-O5'	2.01	117.08	107.75

There are no chirality outliers.

All (28) torsion outliers are listed below:

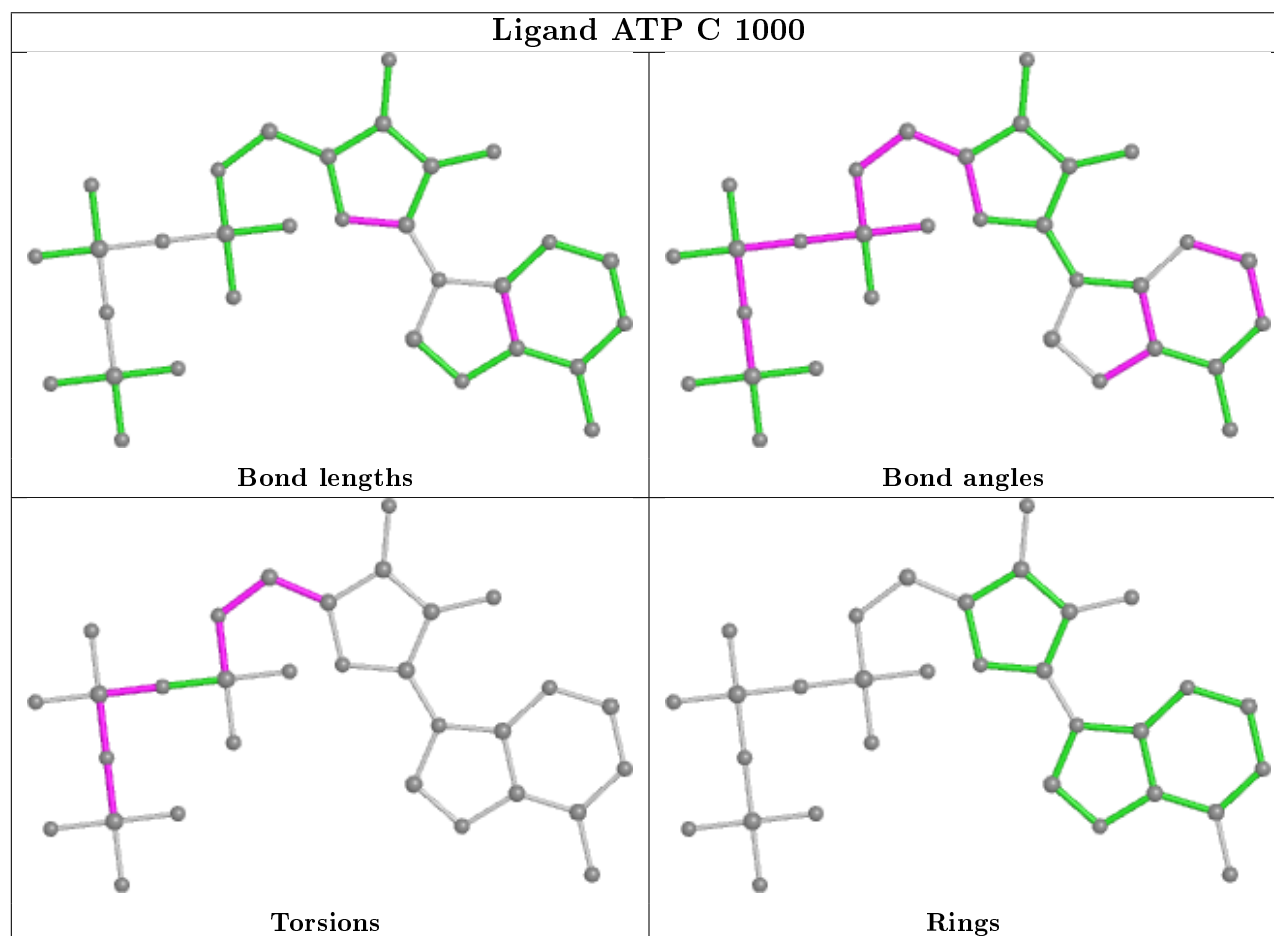
Mol	Chain	Res	Type	Atoms
7	C	1000	ATP	C5'-O5'-PA-O3A
4	A	2000	ADP	C5'-O5'-PA-O3A
7	C	1000	ATP	O4'-C4'-C5'-O5'
7	C	1000	ATP	C3'-C4'-C5'-O5'
7	D	1000	ATP	O4'-C4'-C5'-O5'
7	D	1000	ATP	C3'-C4'-C5'-O5'
5	A	2001	NLG	N2-CA-CB-CG
7	C	1000	ATP	PG-O3B-PB-O1B
4	B	2000	ADP	C4'-C5'-O5'-PA
7	D	1000	ATP	C4'-C5'-O5'-PA
5	A	2001	NLG	CA-CB-CG-CD
7	C	1000	ATP	C4'-C5'-O5'-PA
7	D	1000	ATP	PG-O3B-PB-O1B
4	A	2000	ADP	C4'-C5'-O5'-PA
4	A	2000	ADP	C5'-O5'-PA-O1A
7	C	1000	ATP	PB-O3B-PG-O1G
4	B	2000	ADP	PA-O3A-PB-O1B
7	C	1000	ATP	PG-O3B-PB-O2B
7	D	1000	ATP	PG-O3B-PB-O2B
5	B	2001	NLG	CA-CB-CG-CD
5	A	2001	NLG	C-CA-N2-C7
4	B	2000	ADP	PA-O3A-PB-O2B
4	B	2000	ADP	PA-O3A-PB-O3B
7	D	1000	ATP	C5'-O5'-PA-O3A
4	A	2000	ADP	O4'-C4'-C5'-O5'
7	C	1000	ATP	PA-O3A-PB-O2B
4	B	2000	ADP	PB-O3A-PA-O2A
7	C	1000	ATP	C5'-O5'-PA-O2A

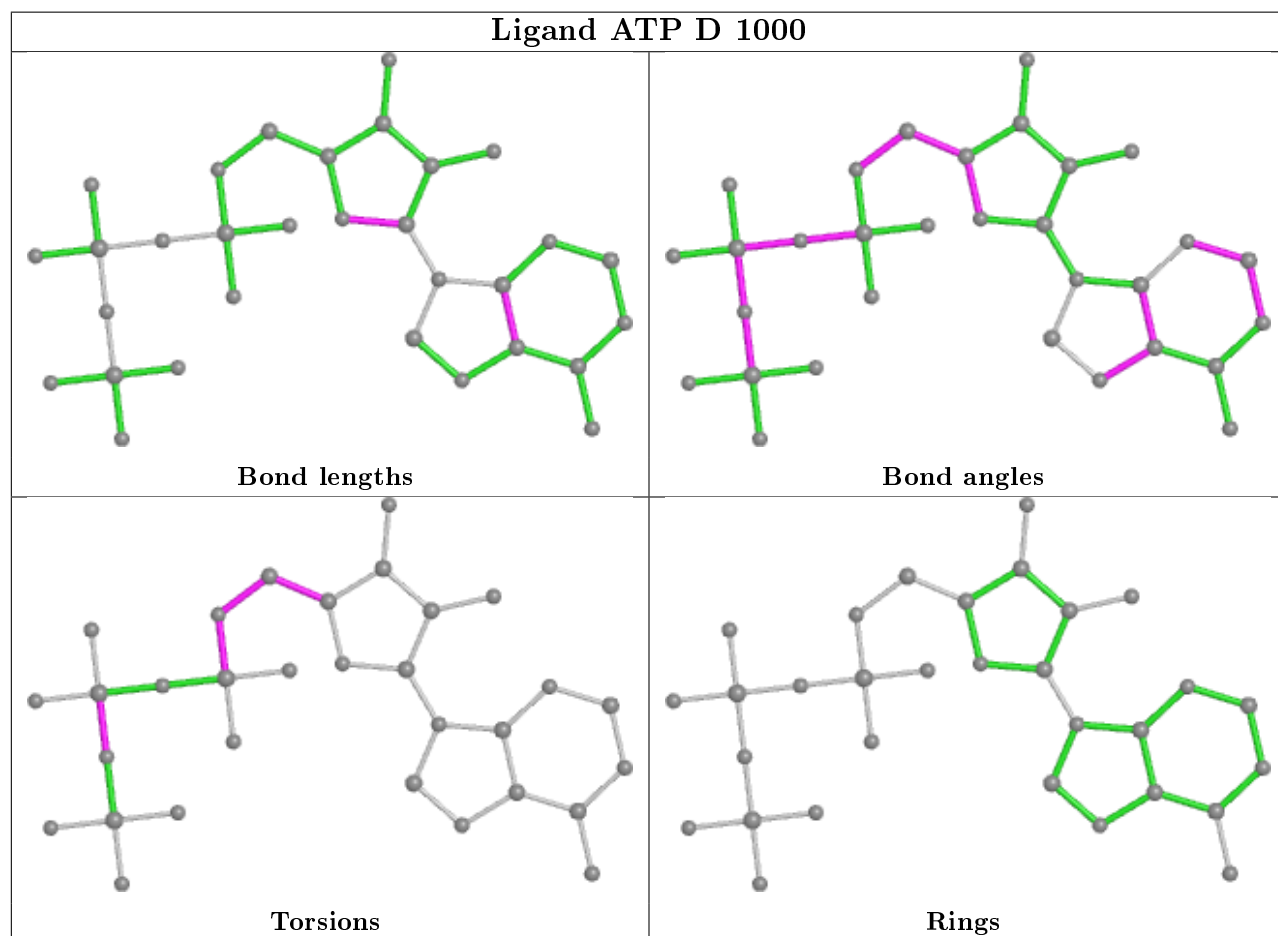
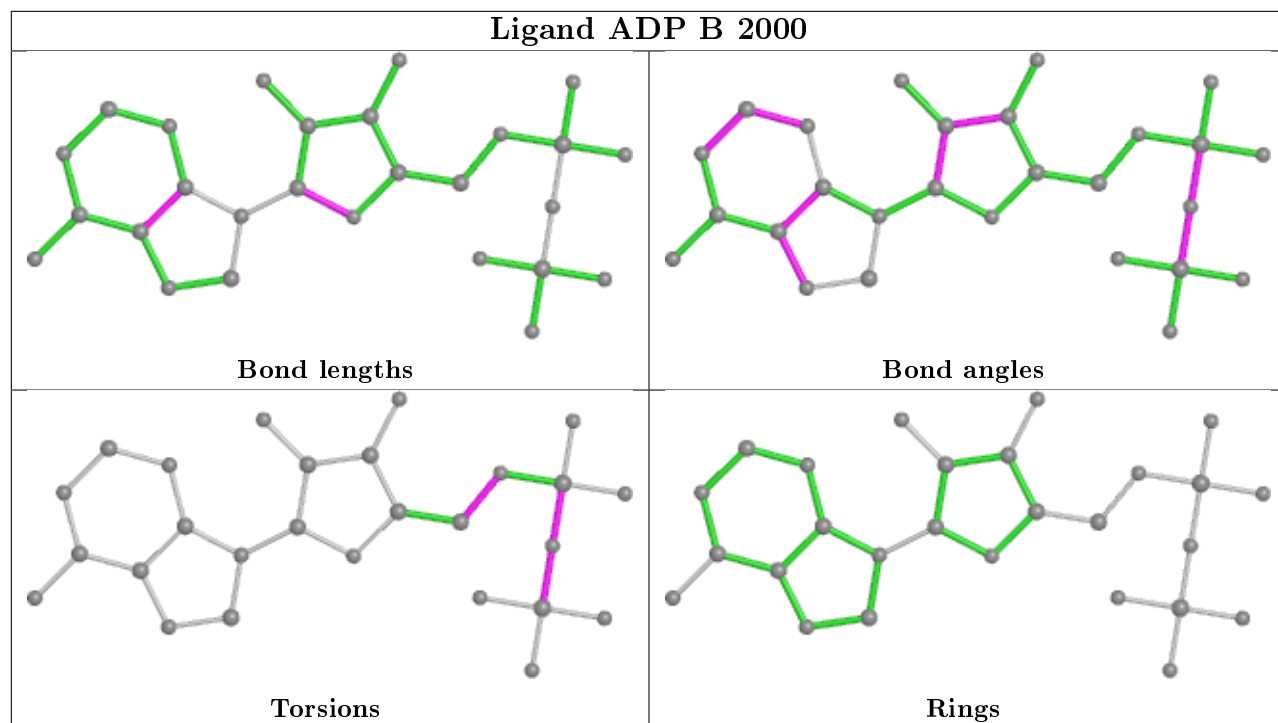
There are no ring outliers.

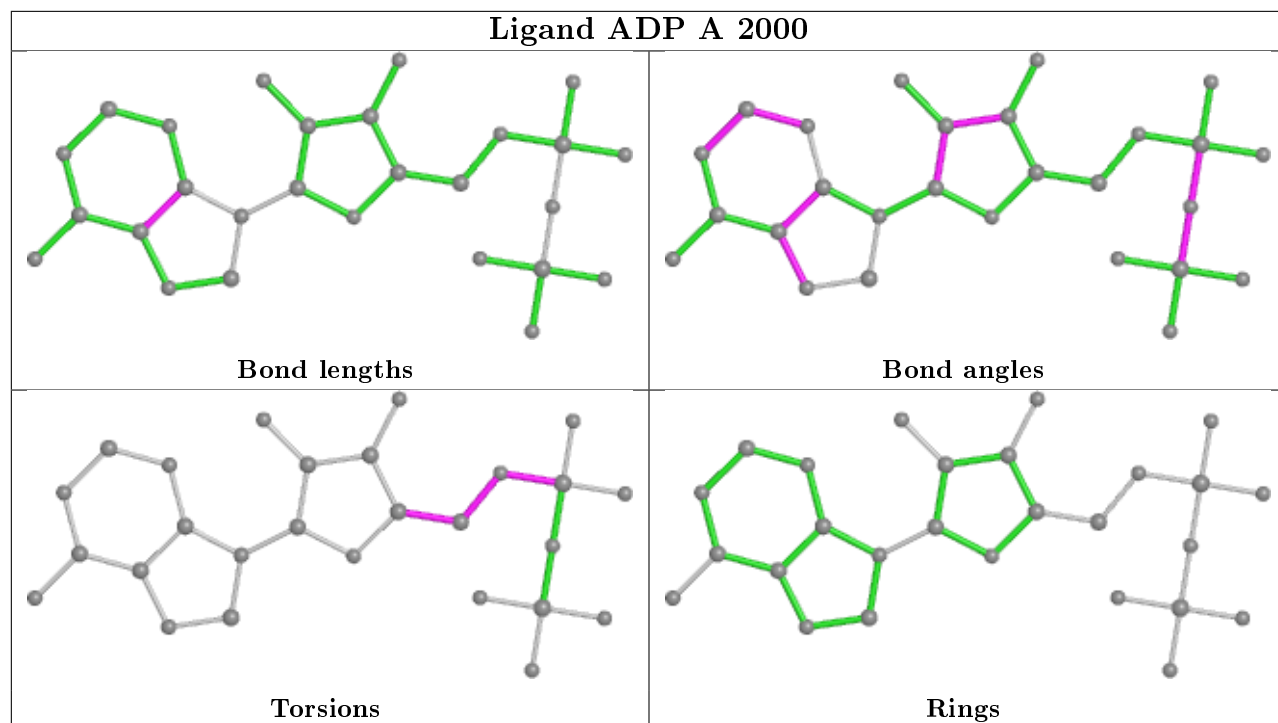
2 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	C	1000	ATP	4	1
7	D	1000	ATP	5	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.







5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	281/298 (94%)	0.31	16 (5%) 23 25	52, 64, 93, 110	0
1	B	283/298 (94%)	0.46	26 (9%) 9 9	54, 63, 84, 115	0
2	C	125/135 (92%)	0.54	12 (9%) 8 7	55, 62, 80, 104	0
2	D	126/135 (93%)	0.57	14 (11%) 5 5	56, 65, 84, 124	0
All	All	815/866 (94%)	0.44	68 (8%) 11 11	52, 64, 90, 124	0

All (68) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	C	128	GLY	8.6
2	C	7	ILE	8.1
1	B	18	TYR	7.9
2	D	130	MET	7.3
1	A	225	GLU	7.1
2	D	7	ILE	5.7
1	B	15	SER	5.4
2	D	128	GLY	5.3
2	C	129	ASP	5.3
1	B	151	ALA	5.1
2	C	127	THR	5.0
2	C	52	GLY	4.8
1	B	16	PRO	4.5
1	A	18	TYR	4.5
2	D	14	LYS	4.3
1	A	297	GLY	4.2
2	D	129	ASP	4.1
2	D	106	LEU	4.0
1	A	241	LYS	4.0
1	A	223	ASN	3.8
2	D	127	THR	3.7

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Mol	Chain	Res	Type	RSRZ
1	A	247	LYS	3.6
1	B	179	ILE	3.3
1	B	73	VAL	3.2
2	C	106	LEU	3.1
1	B	150	SER	3.0
1	B	178	VAL	3.0
1	B	22	ILE	3.0
1	B	199	ALA	2.9
2	D	5	ASP	2.8
1	B	75	GLY	2.8
1	A	73	VAL	2.8
1	A	228	SER	2.8
1	B	180	ALA	2.8
2	C	15	VAL	2.7
2	C	14	LYS	2.7
1	B	148	PRO	2.7
1	B	17	ASP	2.6
2	C	16	GLU	2.6
2	C	104	PHE	2.5
1	A	245	ASP	2.5
1	A	186	ASP	2.4
1	A	244	GLU	2.4
1	B	40	VAL	2.4
1	A	224	LYS	2.4
1	B	228	SER	2.4
1	A	113	VAL	2.3
1	B	41	LYS	2.3
1	B	72	LEU	2.3
2	C	51	GLY	2.3
2	D	122	LYS	2.3
1	B	198	VAL	2.2
1	A	226	ASP	2.2
1	A	230	LEU	2.2
2	D	15	VAL	2.2
2	D	76	VAL	2.2
2	C	125	LYS	2.2
1	A	17	ASP	2.2
2	D	17	ALA	2.2
1	B	200	GLY	2.2
1	B	25	GLU	2.1
2	D	105	VAL	2.1
1	B	133	GLY	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	21	GLU	2.0
1	B	152	GLN	2.0
1	B	213	LEU	2.0
2	D	72	LYS	2.0
1	B	43	GLY	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 carbohydrates in this entry.

6.4 Ligands [i](#)

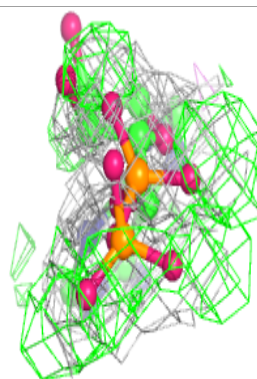
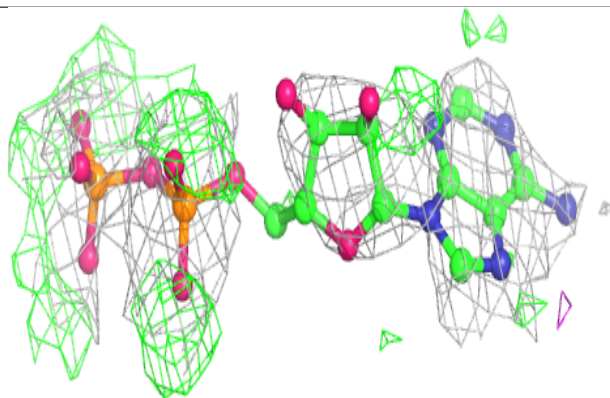
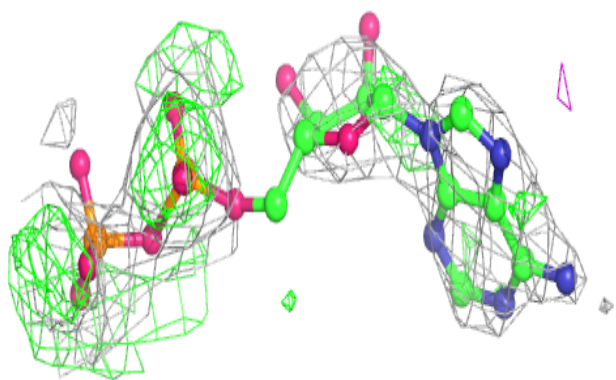
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
6	MG	C	1001	1/1	0.19	0.20	60,60,60,60	0
6	MG	D	1001	1/1	0.35	0.19	64,64,64,64	0
4	ADP	A	2000	27/27	0.53	0.38	71,75,95,98	27
6	MG	B	2002	1/1	0.73	0.33	65,65,65,65	0
3	ARG	B	1000	12/12	0.87	0.20	68,73,74,76	0
3	ARG	A	1000	12/12	0.90	0.18	73,77,80,85	0
5	NLG	A	2001	13/13	0.91	0.17	53,58,64,71	0
5	NLG	B	2001	13/13	0.93	0.13	41,51,55,55	0
7	ATP	C	1000	31/31	0.94	0.12	43,50,64,73	0
7	ATP	D	1000	31/31	0.95	0.12	55,60,66,78	0
4	ADP	B	2000	27/27	0.96	0.12	42,51,55,63	0

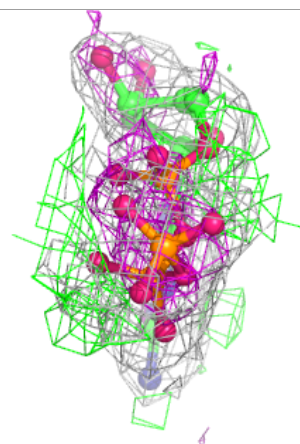
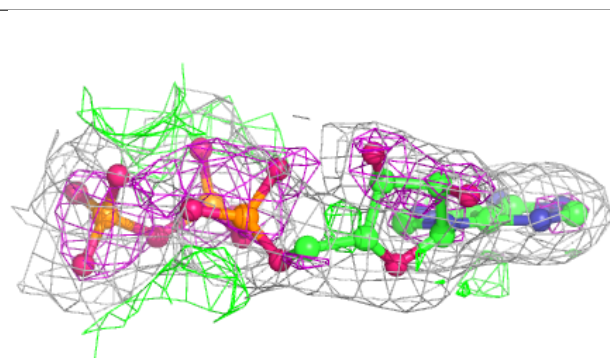
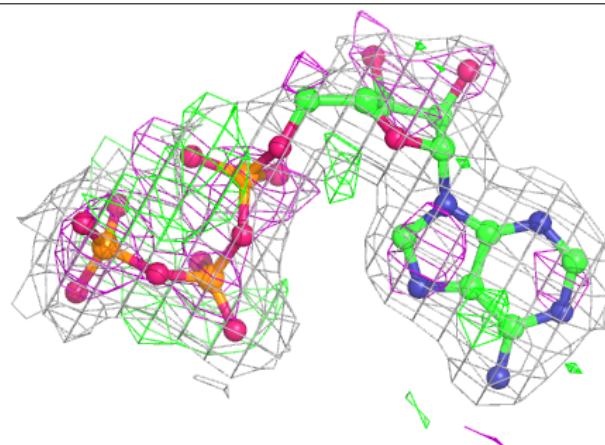
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around ADP A 2000:

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

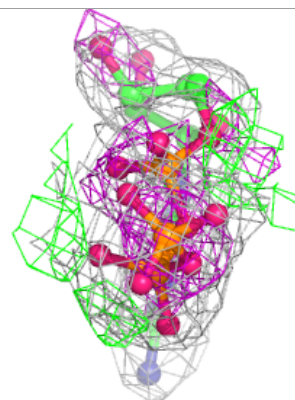
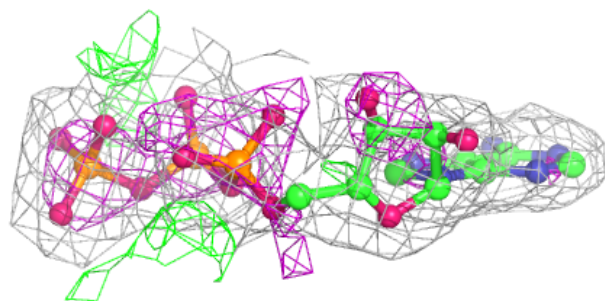
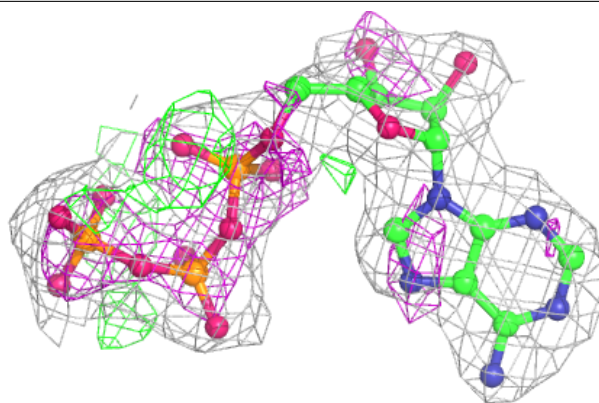
**Electron density around ATP C 1000:**

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

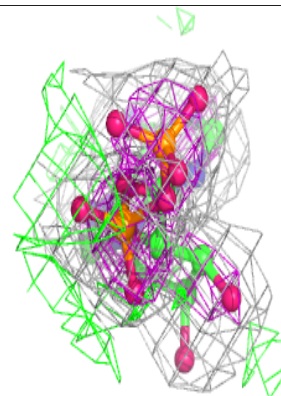
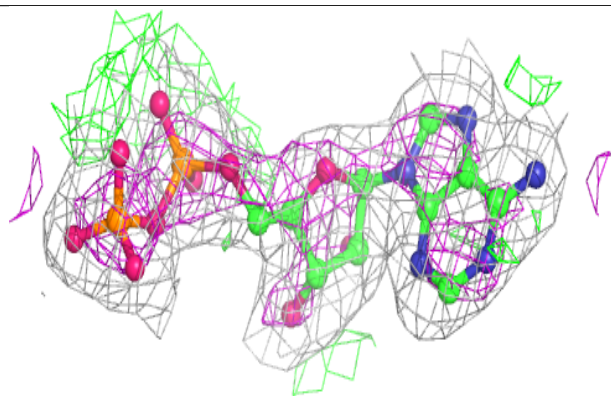
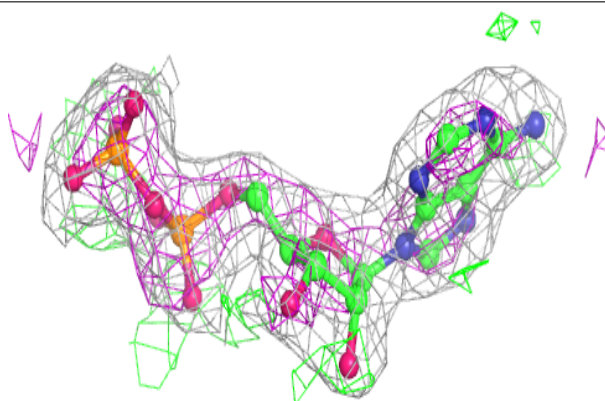


Electron density around ATP D 1000:

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

**Electron density around ADP B 2000:**

$2mF_o-DF_c$ (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.