



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 9, 2026 – 01:25 PM UTC

PDB ID : 3O8L / pdb_00003o8l
Title : Structure of phosphofructokinase from rabbit skeletal muscle
Authors : Banaszak, K.; Chang, S.H.; Rypniewski, W.
Deposited on : 2010-08-03
Resolution : 3.20 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.

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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

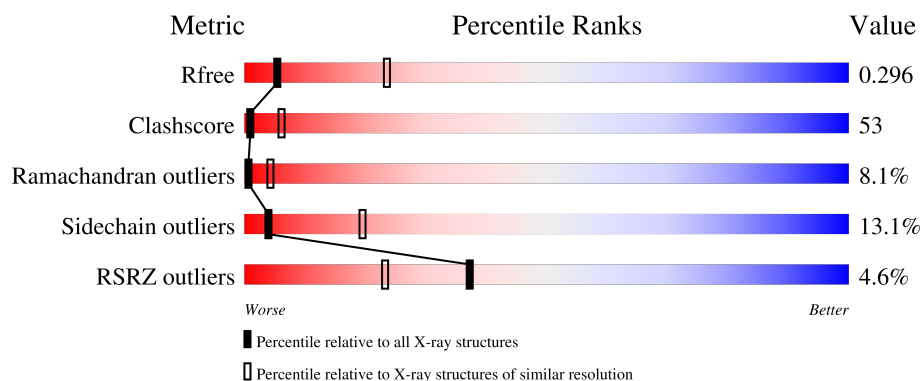
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1466 (3.20-3.20)
Clashscore	190562	1573 (3.20-3.20)
Ramachandran outliers	187476	1548 (3.20-3.20)
Sidechain outliers	187428	1547 (3.20-3.20)
RSRZ outliers	180081	1466 (3.20-3.20)

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	762	4% 31% 49% 16% ..
1	B	762	5% 32% 50% 15% ..

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	PO4	A	766	-	X	-	-
4	PO4	A	767	-	X	-	-
4	PO4	A	768	-	X	-	-
4	PO4	B	766	-	X	-	-
4	PO4	B	767	-	X	-	-
4	PO4	B	768	-	X	-	-

2 Entry composition [i](#)

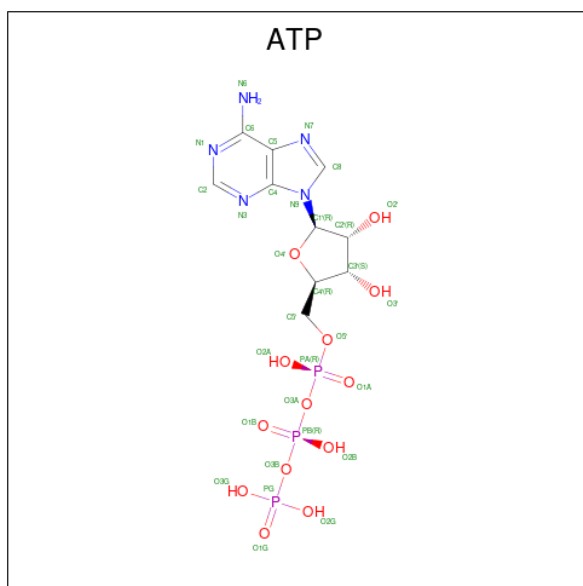
There are 4 unique types of molecules in this entry. The entry contains 11646 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 6-phosphofructokinase, muscle type.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	748	Total	C	N	O	S	406	0	0
			5719	3583	1031	1068	37			
1	B	748	Total	C	N	O	S	420	0	0
			5719	3583	1031	1068	37			

- Molecule 2 is ADENOSINE-5'-TRIPHOSPHATE (CCD ID: ATP) (formula: $C_{10}H_{16}N_5O_{13}P_3$).



- # ADP

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

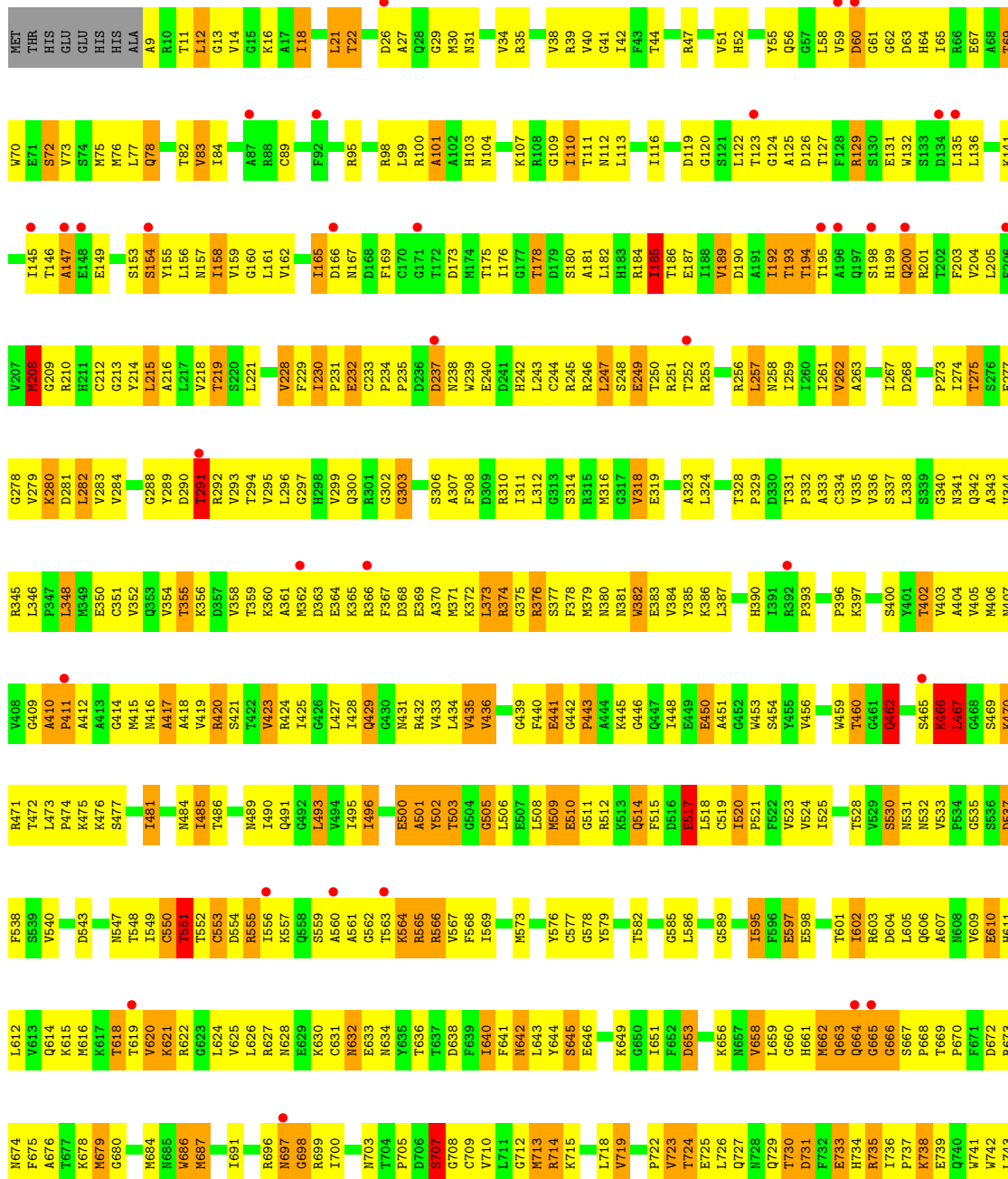
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- Diagram illustrating the Lewis structure of the phosphate ion (PO_4^{3-}). The central phosphorus atom (P) is bonded to four oxygen atoms (O). The bonds are as follows:
- Double bond to the right oxygen (O1).
 - Single bond to the top oxygen (O3).
 - Single bond to the bottom oxygen (O4).
 - Single bond to the left oxygen (O2).
- Each oxygen atom has a negative charge (O^-), and the phosphorus atom has a positive charge (P^+).

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	P	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	P	0	0
			5	4	1		
4	A	1	Total	O	P	0	0
			5	4	1		
4	B	1	Total	O	P	0	0
			5	4	1		
4	B	1	Total	O	P	0	0
			5	4	1		
4	B	1	Total	O	P	0	0
			5	4	1		



K744	L745	R746	P747	I748	L749	K750	I751	L752	A753	K754	Y755	E756	ILE	ASP	LEU	ASP	THR	SER
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4 Data and refinement statistics

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, α , β , γ	163.68Å 163.68Å 356.58Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	41.00 – 3.20 41.00 – 3.20	Depositor EDS
% Data completeness (in resolution range)	(Not available) (41.00-3.20) 99.9 (41.00-3.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.14	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.26 (at 3.19Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.238 , 0.302 0.235 , 0.296	Depositor DCC
R_{free} test set	2362 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	81.9	Xtriage
Anisotropy	0.290	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 117.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	11646	wwPDB-VP
Average B, all atoms (Å ²)	70.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.42% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: PO4, ADP, ATP

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	0.79	0/5819	1.24	67/7859 (0.9%)
1	B	0.79	4/5819 (0.1%)	1.21	59/7859 (0.8%)
All	All	0.79	4/11638 (0.0%)	1.22	126/15718 (0.8%)

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

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

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	436	VAL	CA-CB	6.44	1.62	1.54
1	B	417	ALA	CA-CB	-5.72	1.44	1.53
1	B	185	ILE	CA-CB	5.23	1.62	1.54
1	B	423	VAL	CA-CB	-5.01	1.48	1.54

The worst 5 of 126 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	533	VAL	CA-C-N	-13.21	106.26	120.66
1	B	533	VAL	C-N-CA	-13.21	106.26	120.66
1	A	501	ALA	N-CA-C	-11.41	97.75	112.93
1	A	441	GLU	N-CA-C	-10.58	100.40	113.20
1	A	299	VAL	N-CA-C	-10.57	103.17	113.53

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	385	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	5719	0	5742	524	0
1	B	5719	0	5742	601	0
2	A	62	0	24	5	0
2	B	62	0	24	1	0
3	A	27	0	12	2	0
3	B	27	0	12	0	0
4	A	15	0	0	1	0
4	B	15	0	0	1	0
All	All	11646	0	11556	1116	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 53.

The worst 5 of 1116 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:634:ASN:HD22	1:B:646:GLU:HG2	0.99	1.09
1:A:646:GLU:HG2	1:B:634:ASN:HD22	1.22	1.03
1:A:296:LEU:H	1:A:296:LEU:HD12	1.23	1.02
1:A:751:ILE:HD13	1:A:752:LEU:N	1.75	1.00
1:B:609:VAL:HG22	1:B:644:TYR:CE2	1.95	0.99

There are no symmetry-related clashes.

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	746/762 (98%)	564 (76%)	123 (16%)	59 (8%)	1	5
1	B	746/762 (98%)	565 (76%)	119 (16%)	62 (8%)	0	4
All	All	1492/1524 (98%)	1129 (76%)	242 (16%)	121 (8%)	1	4

5 of 121 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	43	PHE
1	A	51	VAL
1	A	54	GLY
1	A	60	ASP
1	A	153	SER

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	605/618 (98%)	525 (87%)	80 (13%)	4	19
1	B	605/618 (98%)	527 (87%)	78 (13%)	4	20
All	All	1210/1236 (98%)	1052 (87%)	158 (13%)	4	19

5 of 158 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	435	VAL

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Mol	Chain	Res	Type
1	B	687	MET
1	B	462	GLN
1	B	517	GLU
1	B	724	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 44 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	407	ASN
1	B	531	ASN
1	B	431	ASN
1	B	489	ASN
1	B	611	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](#)

12 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	ATP	B	763	-	32,33,33	1.16	3 (9%)	48,52,52	1.61	9 (18%)
4	PO4	B	767	-	4,4,4	2.01	3 (75%)	6,6,6	1.24	1 (16%)
4	PO4	A	768	-	4,4,4	1.99	3 (75%)	6,6,6	1.40	1 (16%)
4	PO4	A	767	-	4,4,4	1.99	3 (75%)	6,6,6	1.45	2 (33%)
2	ATP	A	763	-	32,33,33	1.21	3 (9%)	48,52,52	1.63	10 (20%)
2	ATP	A	765	-	32,33,33	1.19	3 (9%)	48,52,52	2.19	20 (41%)
3	ADP	B	764	-	28,29,29	1.15	3 (10%)	43,45,45	1.72	8 (18%)
4	PO4	B	768	-	4,4,4	2.00	3 (75%)	6,6,6	1.23	1 (16%)
4	PO4	B	766	-	4,4,4	2.04	3 (75%)	6,6,6	1.29	2 (33%)
3	ADP	A	764	-	28,29,29	1.16	3 (10%)	43,45,45	2.05	14 (32%)
2	ATP	B	765	-	32,33,33	1.15	3 (9%)	48,52,52	1.96	16 (33%)
4	PO4	A	766	-	4,4,4	2.01	3 (75%)	6,6,6	1.37	1 (16%)

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	ATP	B	763	-	-	4/22/38/38	0/3/3/3
2	ATP	A	763	-	-	4/22/38/38	0/3/3/3
2	ATP	A	765	-	-	5/22/38/38	0/3/3/3
3	ADP	B	764	-	-	2/16/32/32	0/3/3/3
3	ADP	A	764	-	-	2/16/32/32	0/3/3/3
2	ATP	B	765	-	-	4/22/38/38	0/3/3/3

The worst 5 of 36 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	765	ATP	C5-N7	-3.16	1.33	1.39
3	A	764	ADP	C5-N7	-3.15	1.33	1.39
2	A	765	ATP	C5-N7	-3.15	1.33	1.39
3	B	764	ADP	C5-N7	-3.13	1.33	1.39
2	A	763	ATP	C5-N7	-3.09	1.33	1.39

The worst 5 of 85 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	764	ADP	C5-C4-N3	-5.47	119.18	126.72
3	B	764	ADP	C5-C4-N3	-5.33	119.38	126.72
2	B	763	ATP	C5-C4-N3	-5.18	119.58	126.72
2	A	765	ATP	C5-C4-N3	-5.09	119.71	126.72
2	A	763	ATP	C5-C4-N3	-5.04	119.77	126.72

There are no chirality outliers.

5 of 21 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	763	ATP	PB-O3B-PG-O3G
2	A	765	ATP	C5'-O5'-PA-O1A
2	A	765	ATP	C5'-O5'-PA-O2A
2	A	765	ATP	C5'-O5'-PA-O3A
2	A	763	ATP	PA-O3A-PB-O1B

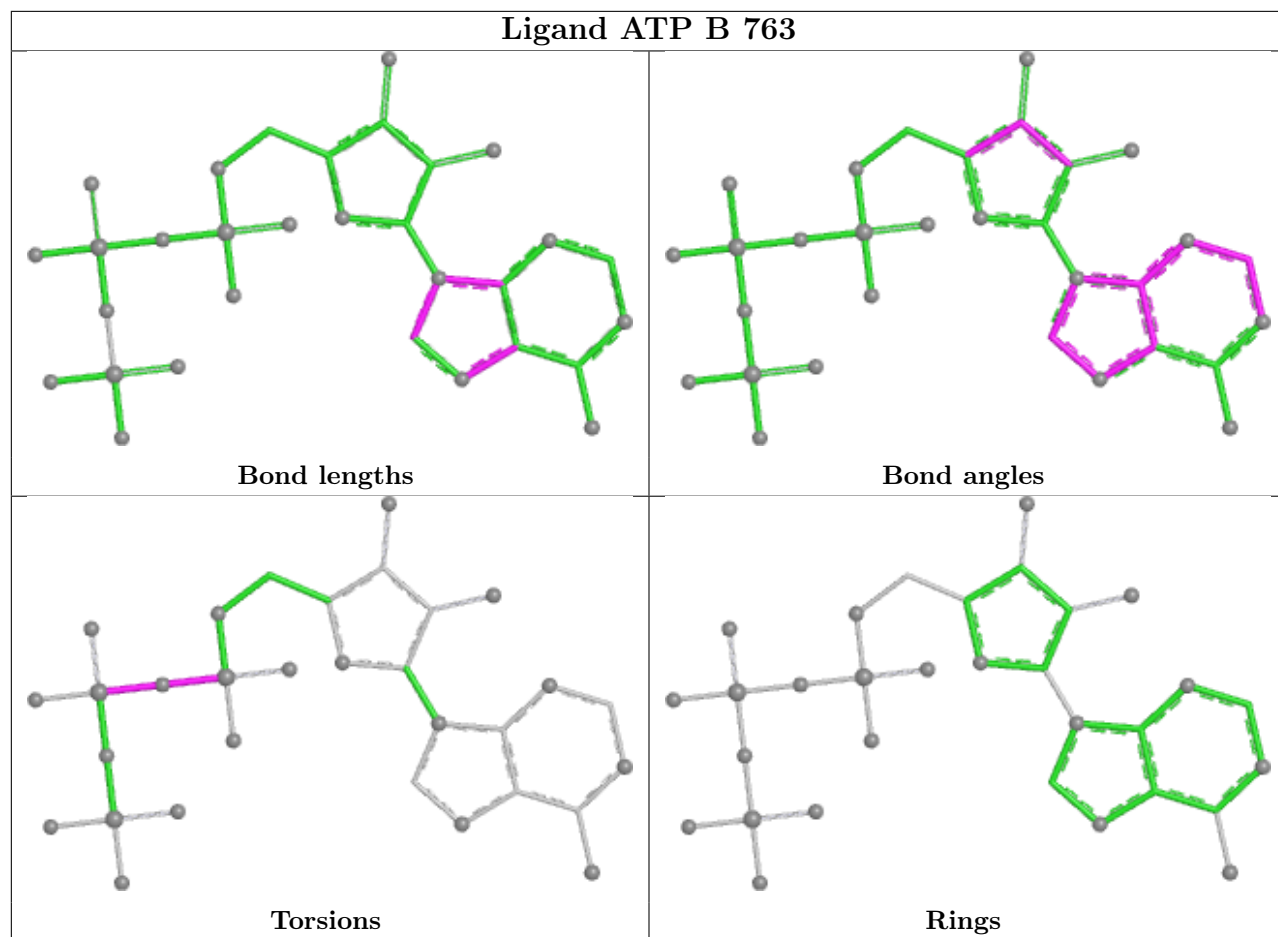
There are no ring outliers.

6 monomers are involved in 10 short contacts:

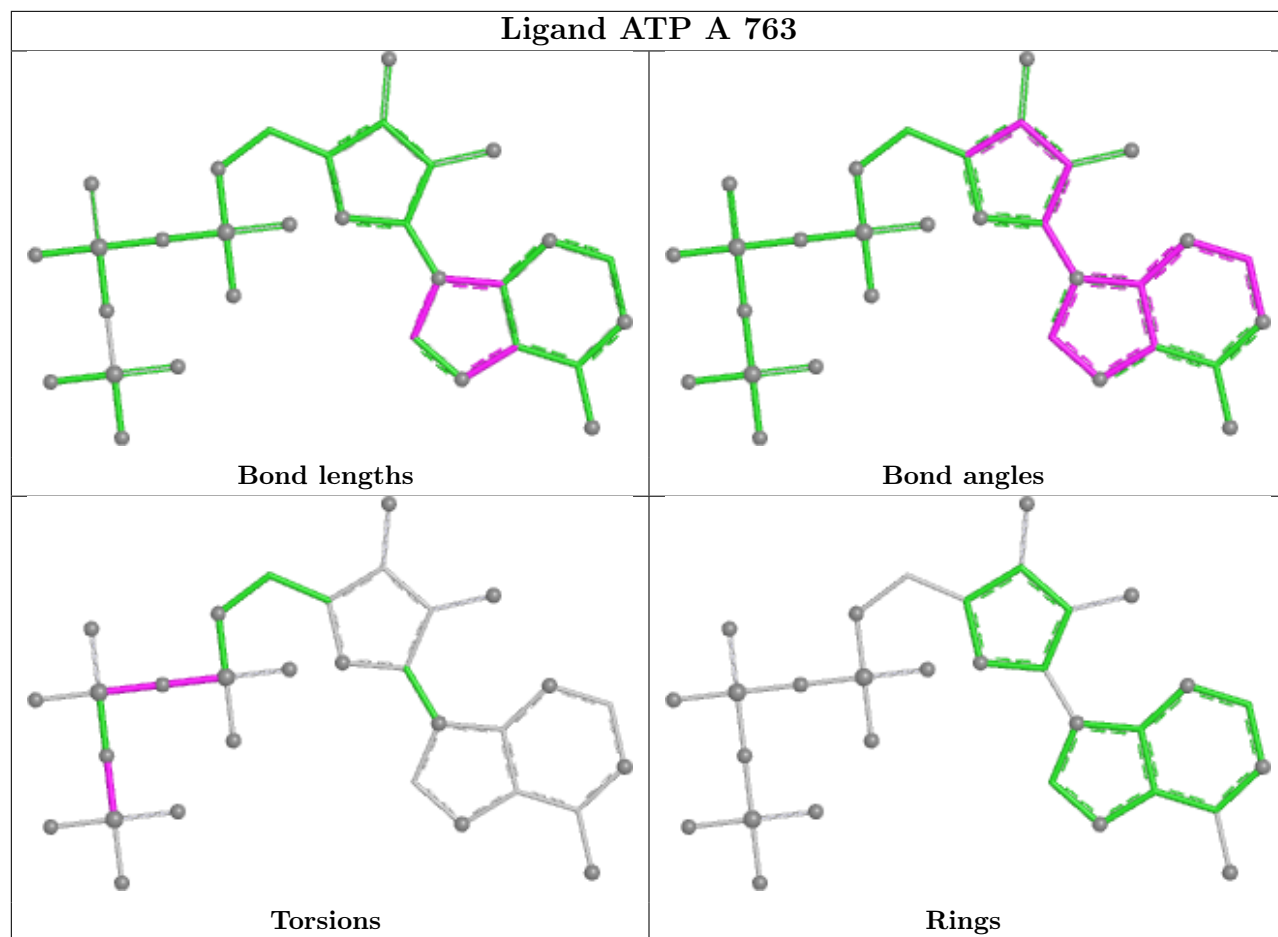
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	763	ATP	1	0
4	A	767	PO4	1	0
2	A	763	ATP	2	0
2	A	765	ATP	3	0
4	B	768	PO4	1	0
3	A	764	ADP	2	0

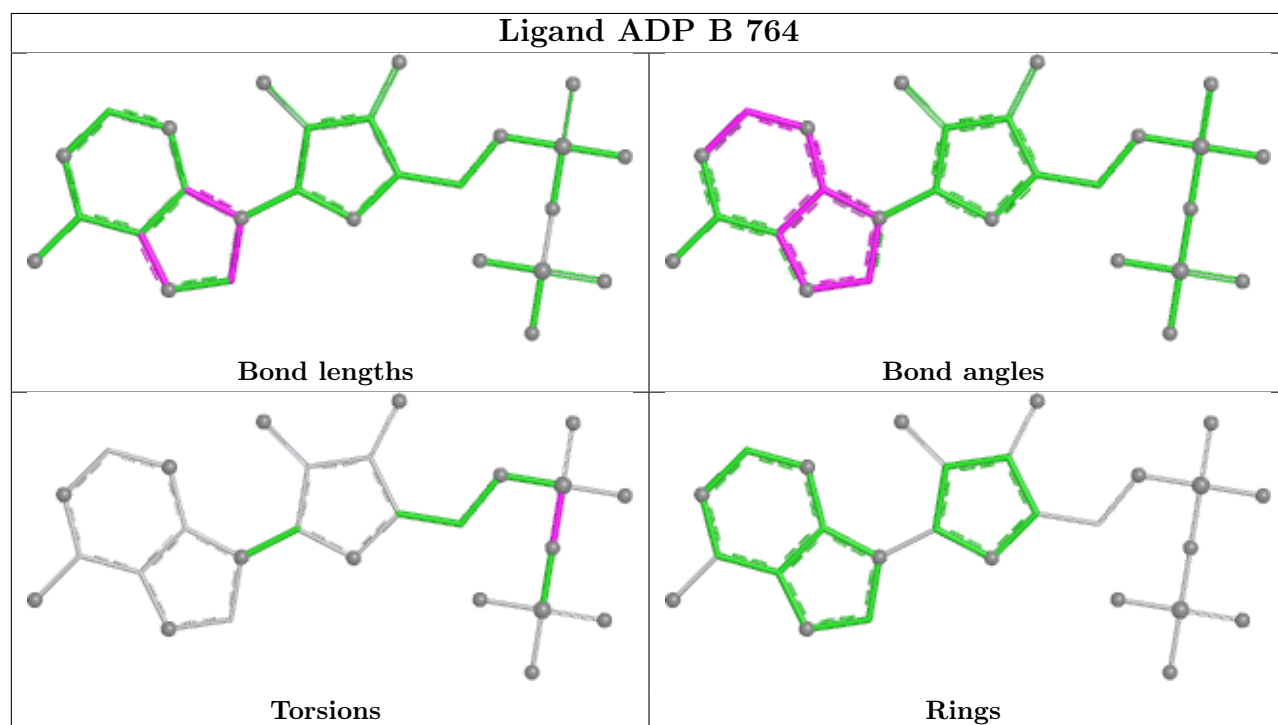
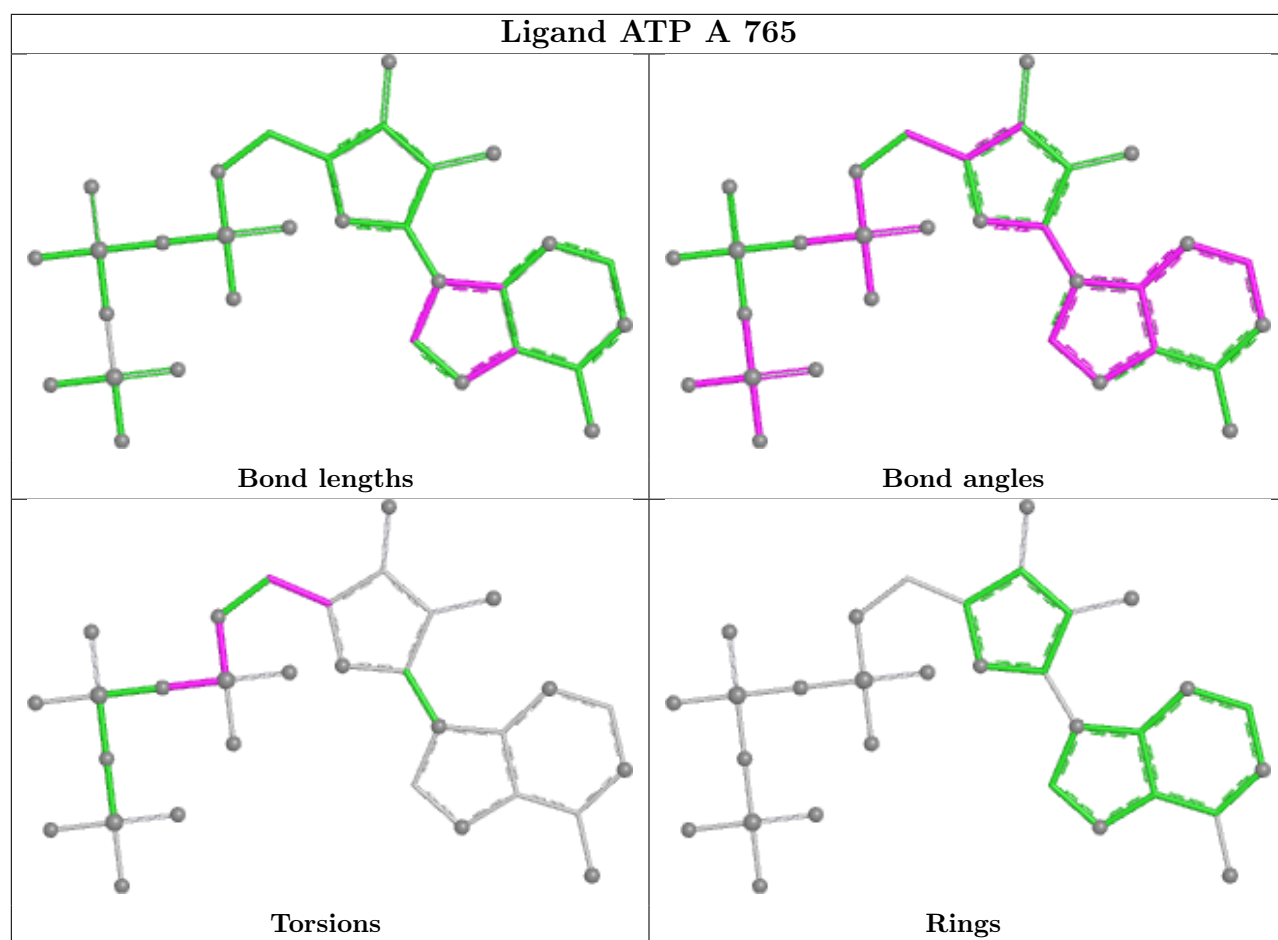
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less 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.

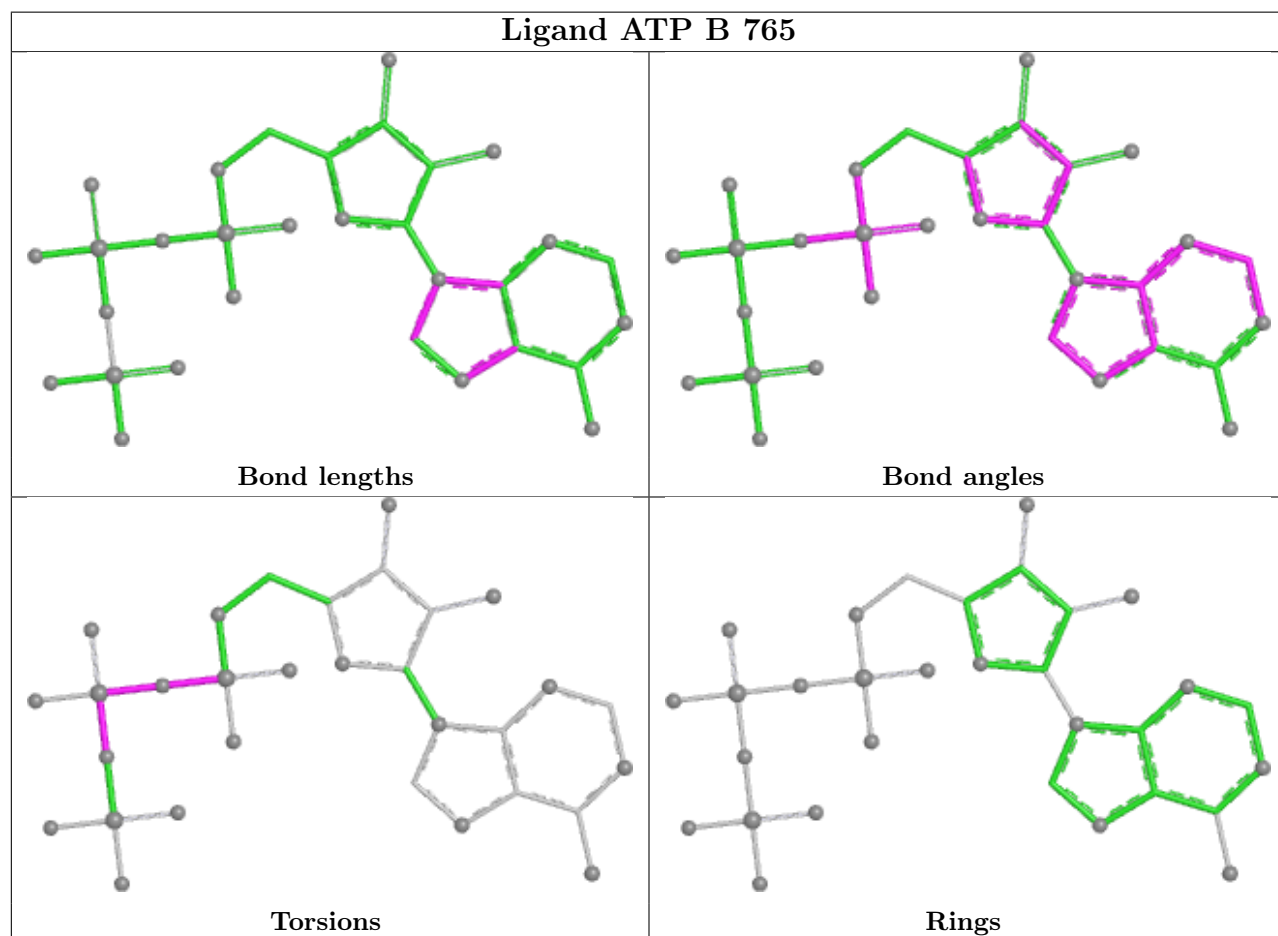
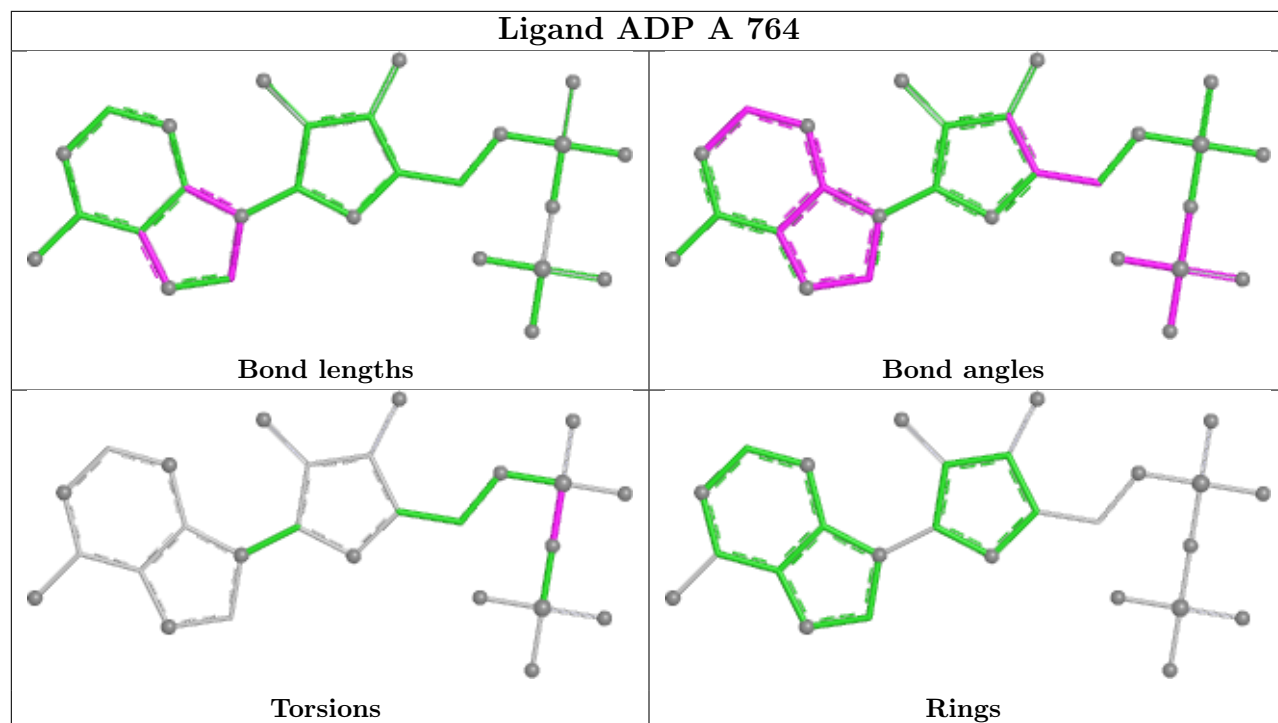
Ligand ATP B 763



Ligand ATP A 763







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	739/762 (96%)	0.08	33 (4%) 38 24	14, 52, 126, 184	76 (10%)
1	B	745/762 (97%)	0.12	35 (4%) 36 23	7, 54, 128, 202	86 (11%)
All	All	1484/1524 (97%)	0.10	68 (4%) 37 23	7, 53, 128, 202	162 (10%)

The worst 5 of 68 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	198	SER	8.1
1	A	148	GLU	4.8
1	B	563	THR	4.7
1	A	9	ALA	4.5
1	A	155	TYR	4.5

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](#)

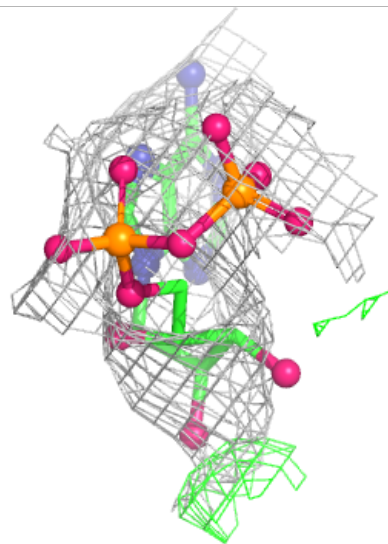
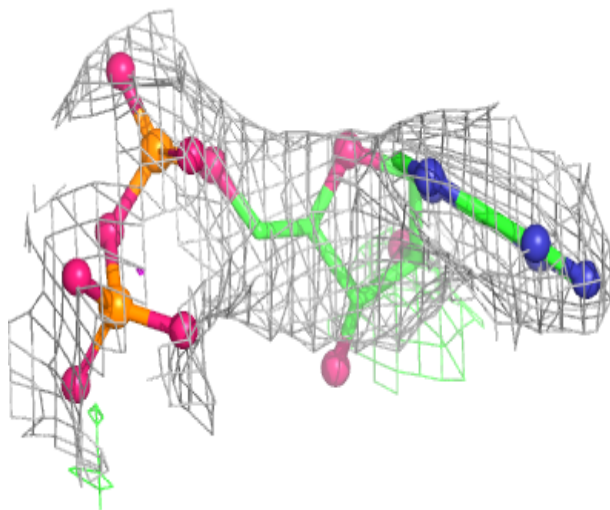
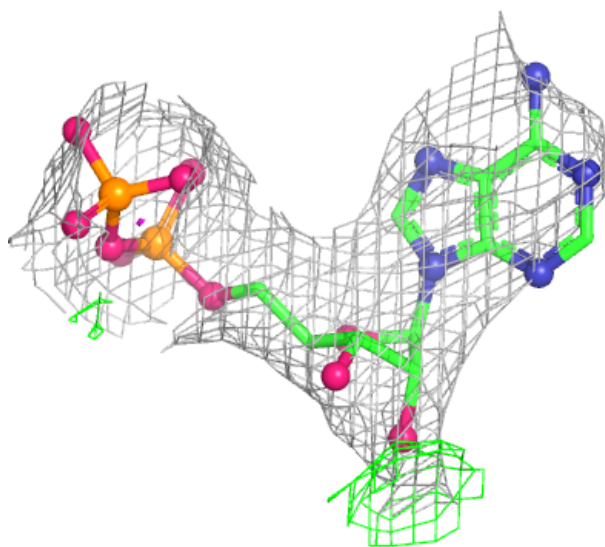
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q < 0.9’ lists 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
4	PO4	A	766	5/5	0.64	0.18	99,99,100,101	0
4	PO4	B	766	5/5	0.74	0.14	91,92,93,94	0
3	ADP	B	764	27/27	0.84	0.18	145,148,154,154	0
3	ADP	A	764	27/27	0.85	0.21	135,138,144,144	0
4	PO4	B	767	5/5	0.90	0.13	66,66,69,71	0
2	ATP	A	763	31/31	0.91	0.09	82,87,97,98	0
2	ATP	B	765	31/31	0.92	0.09	72,77,90,91	0
4	PO4	A	768	5/5	0.92	0.09	55,56,57,58	0
4	PO4	B	768	5/5	0.93	0.09	52,53,55,55	0
2	ATP	B	763	31/31	0.94	0.08	98,100,101,101	0
4	PO4	A	767	5/5	0.94	0.12	60,60,61,63	0
2	ATP	A	765	31/31	0.94	0.07	58,63,77,78	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.

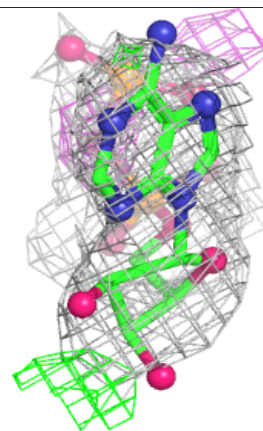
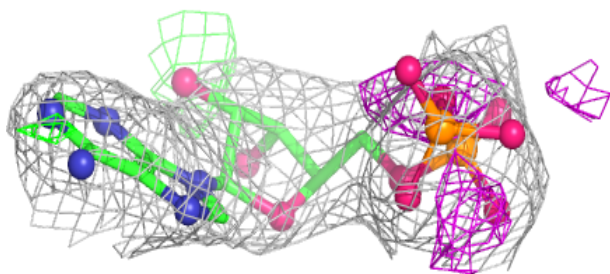
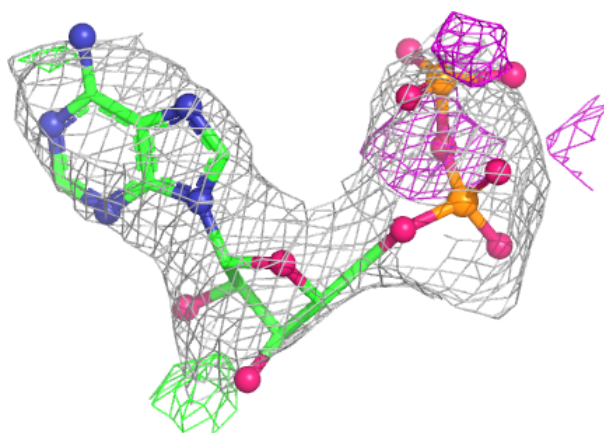
Electron density around ADP B 764:

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



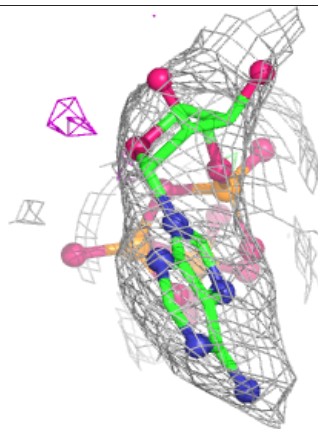
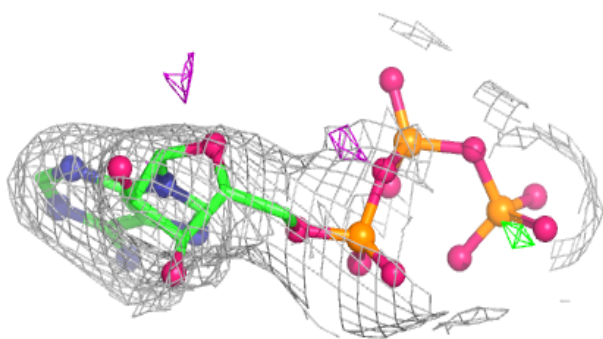
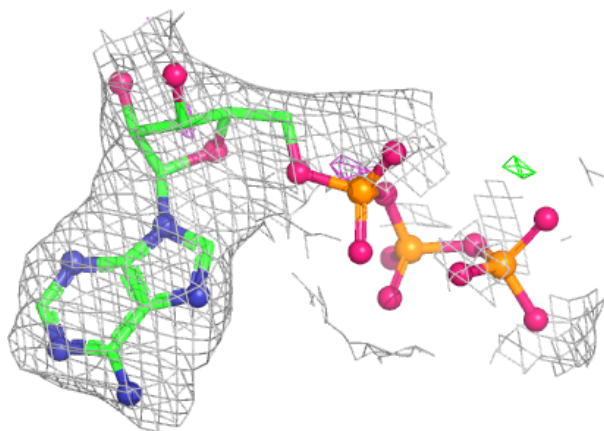
Electron density around ADP A 764:

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



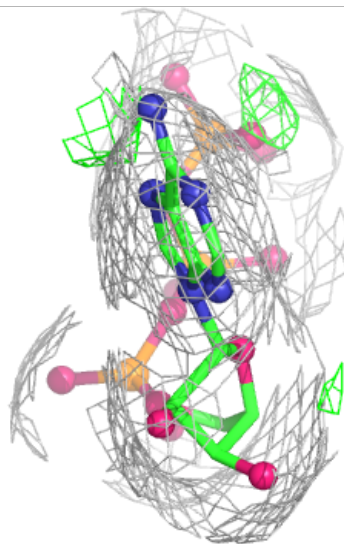
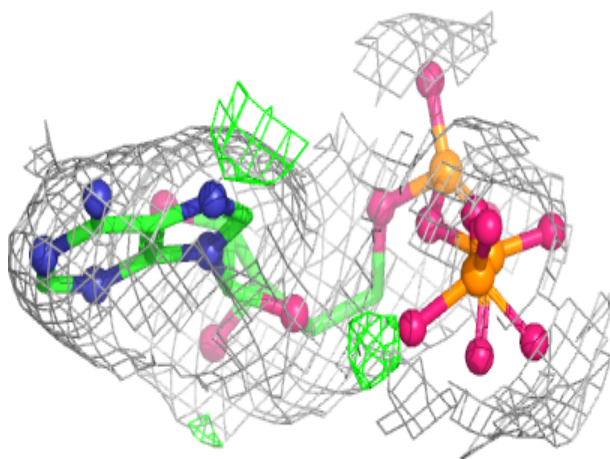
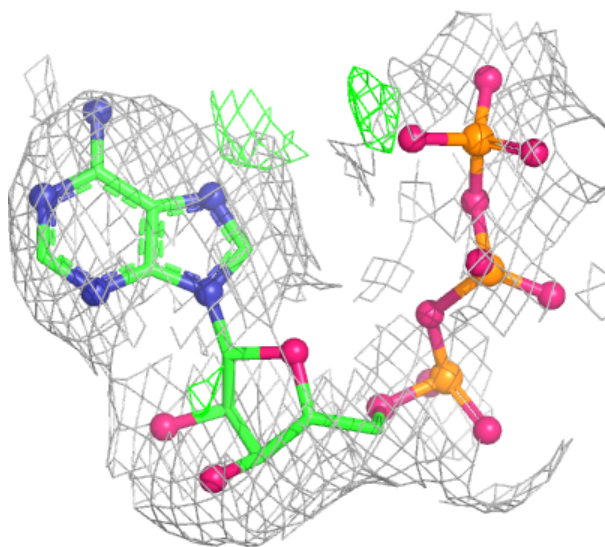
Electron density around ATP A 763:

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



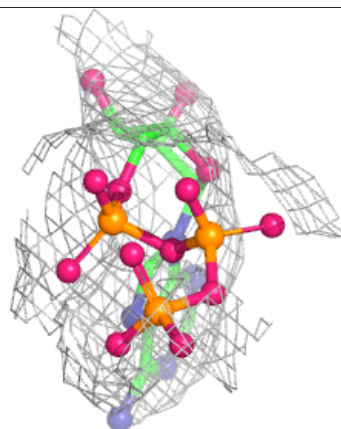
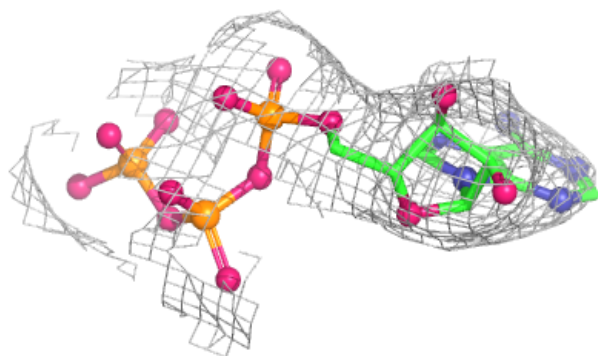
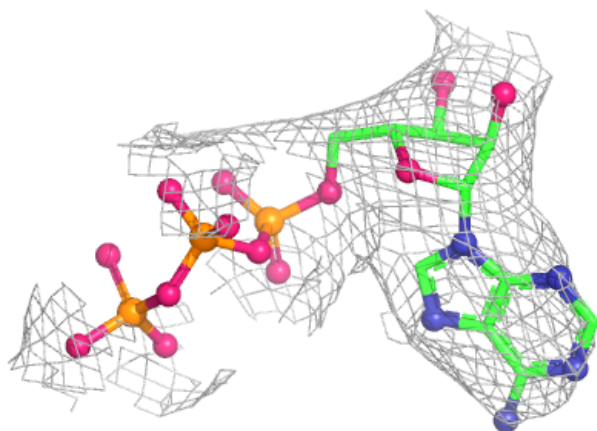
Electron density around ATP B 765:

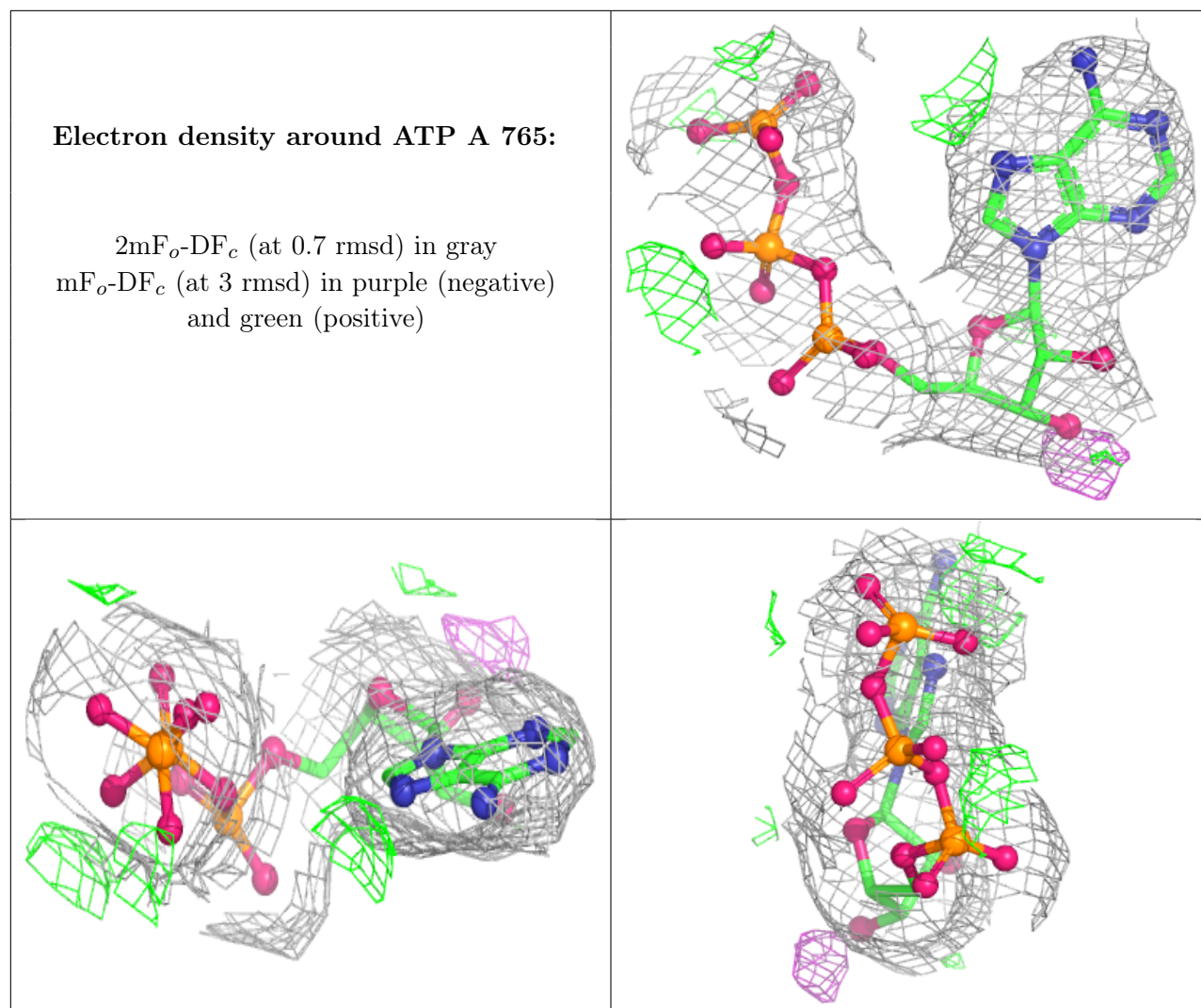
2mF_o-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around ATP B 763:

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





6.5 Other polymers ⓘ

There are no such residues in this entry.