



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 07:14 PM UTC

PDB ID : 7U9K / pdb_00007u9k
Title : Staphylococcus aureus D-alanine-D-alanine ligase in complex with ATP, D-ala-D-ala, Mg²⁺ and K⁺
Authors : Pederick, J.L.; Bruning, J.B.
Deposited on : 2022-03-10
Resolution : 2.00 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

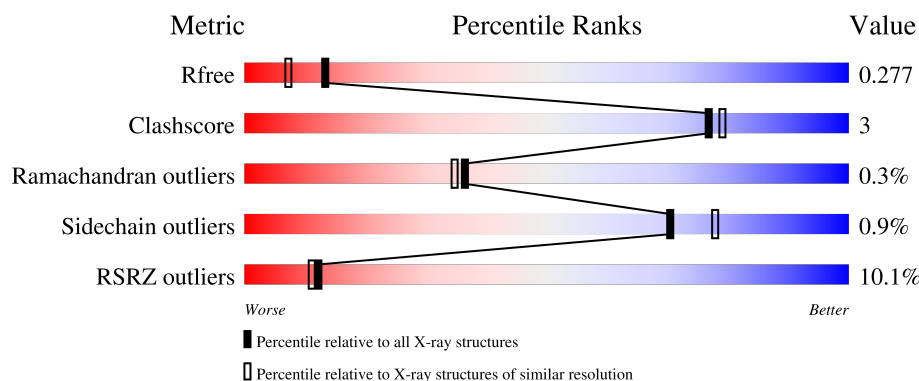
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	364	 5% 85% 7% 9%
1	B	364	 13% 82% 9% 9%
2	F	2	 50% 50%

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 5487 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 D-alanine–D-alanine ligase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	332	Total	C	N	O	S	0	0	0
			2568	1631	426	502	9			
1	B	330	Total	C	N	O	S	0	0	0
			2562	1638	417	498	9			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	357	ARG	-	expression tag	UNP Q2FWH3
A	358	SER	-	expression tag	UNP Q2FWH3
A	359	HIS	-	expression tag	UNP Q2FWH3
A	360	HIS	-	expression tag	UNP Q2FWH3
A	361	HIS	-	expression tag	UNP Q2FWH3
A	362	HIS	-	expression tag	UNP Q2FWH3
A	363	HIS	-	expression tag	UNP Q2FWH3
A	364	HIS	-	expression tag	UNP Q2FWH3
B	357	ARG	-	expression tag	UNP Q2FWH3
B	358	SER	-	expression tag	UNP Q2FWH3
B	359	HIS	-	expression tag	UNP Q2FWH3
B	360	HIS	-	expression tag	UNP Q2FWH3
B	361	HIS	-	expression tag	UNP Q2FWH3
B	362	HIS	-	expression tag	UNP Q2FWH3
B	363	HIS	-	expression tag	UNP Q2FWH3
B	364	HIS	-	expression tag	UNP Q2FWH3

- Molecule 2 is a protein (with D amino acids) called DAL-DAL.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	F	2	Total	C	N	O	0	0	0
			11	6	2	3			

- Molecule 3 is POTASSIUM ION (CCD ID: K) (formula: K) (labeled as "Ligand of Interest")

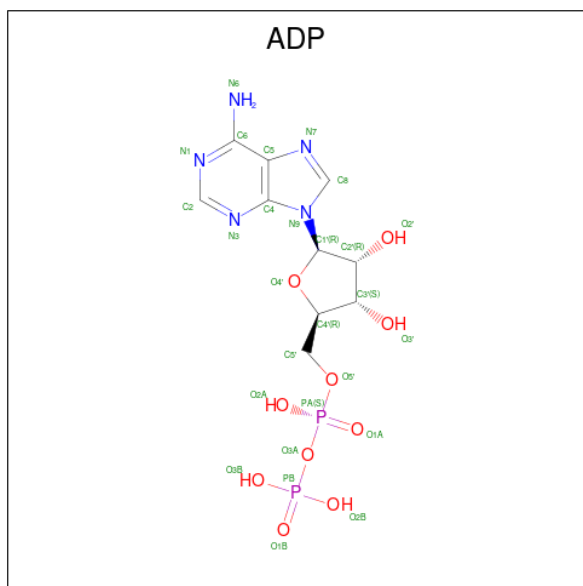
by depositor).

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

- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

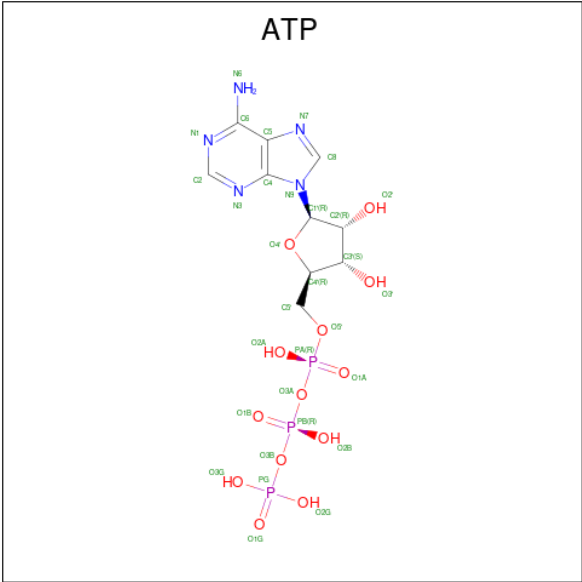
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Mg	0	0
			1	1		
4	B	1	Total	Mg	0	0
			1	1		

- Molecule 5 is ADENOSINE-5'-DIPHOSPHATE (CCD ID: ADP) (formula: C₁₀H₁₅N₅O₁₀P₂).



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

- Molecule 6 is ADENOSINE-5'-TRIPHOSPHATE (CCD ID: ATP) (formula: C₁₀H₁₆N₅O₁₃P₃) (labeled as "Ligand of Interest" by depositor).

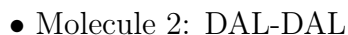


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	B	1	Total	C	N	O	P	0	0
			31	10	5	13	3		

- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	169	Total	O	0	0
			169	169		
7	B	112	Total	O	0	0
			112	112		
7	F	1	Total	O	0	0
			1	1		

- Molecule 1: D-alanine–D-alanine ligase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	52.45Å 70.06Å 91.60Å 90.00° 102.58° 90.00°	Depositor
Resolution (Å)	37.68 – 2.00 37.68 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.8 (37.68-2.00) 98.4 (37.68-2.00)	Depositor EDS
R_{merge}	0.22	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.67 (at 2.00Å)	Xtriage
Refinement program	PHENIX 1.18.2-3874	Depositor
R, R_{free}	0.239 , 0.277 0.240 , 0.277	Depositor DCC
R_{free} test set	2160 reflections (4.92%)	wwPDB-VP
Wilson B-factor (Å ²)	27.1	Xtriage
Anisotropy	0.392	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 44.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	5487	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.99% 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: K, MG, DAL, ATP, ADP

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	0.08	0/2616	0.26	0/3550
1	B	0.09	0/2613	0.27	0/3550
All	All	0.08	0/5229	0.26	0/7100

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2568	0	2471	13	0
1	B	2562	0	2446	17	0
2	F	11	0	11	0	0
3	A	1	0	0	0	0
3	B	3	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	27	0	12	1	0
6	B	31	0	12	2	0
7	A	169	0	0	1	0
7	B	112	0	0	0	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	F	1	0	0	0	0
All	All	5487	0	4952	29	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 3.

All (29) 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:23:THR:HG23	1:B:313:PHE:HB2	1.72	0.71
1:B:222:GLU:HB2	1:B:318:MET:HE2	1.80	0.64
1:B:148:ILE:HD11	1:B:169:LYS:HD3	1.88	0.56
1:B:8:ILE:HD13	1:B:92:PHE:HB3	1.89	0.55
1:A:269:VAL:HG13	1:A:302:ILE:HD11	1.88	0.54
1:B:251:LYS:NZ	6:B:405:ATP:O1G	2.41	0.51
1:A:207:ASP:OD1	1:A:208:ARG:N	2.45	0.49
1:A:217:ASN:ND2	7:A:508:HOH:O	2.39	0.49
1:A:96:HIS:CE1	1:A:310:MET:HG3	2.48	0.48
1:A:130:LYS:NZ	5:A:403:ADP:O2B	2.44	0.47
1:A:8:ILE:HD13	1:A:92:PHE:HB3	1.98	0.46
1:B:313:PHE:CD2	1:B:320:PRO:HG3	2.52	0.45
1:B:299:ASP:OD2	1:B:301:GLN:NE2	2.40	0.45
1:A:122:LEU:HD21	1:B:132:VAL:HG21	1.99	0.45
1:A:178:PRO:HG2	1:A:181:LEU:HD12	1.98	0.44
1:B:140:ARG:NE	1:B:279:GLU:OE1	2.43	0.44
1:B:207:ASP:OD1	1:B:208:ARG:N	2.50	0.44
1:A:259:LEU:HD22	1:A:318:MET:HE2	2.01	0.43
1:B:220:GLU:OE1	6:B:405:ATP:O2'	2.35	0.42
1:A:330:TYR:HB3	1:A:331:PRO:HD3	2.02	0.41
1:B:345:HIS:O	1:B:349:GLN:HG2	2.20	0.41
1:A:319:TYR:HB3	1:A:320:PRO:HD3	2.02	0.41
1:B:7:CYS:HB2	1:B:88:TYR:CE1	2.56	0.41
1:A:61:SER:O	1:A:64:GLU:HG2	2.21	0.41
1:B:318:MET:HG2	1:B:322:LEU:HG	2.03	0.41
1:B:318:MET:HE3	1:B:322:LEU:HD21	2.02	0.41
1:B:330:TYR:HB3	1:B:331:PRO:HD3	2.03	0.41
1:A:204:PHE:HD1	1:A:204:PHE:HA	1.81	0.41
1:B:143:PRO:O	1:B:304:ILE:HG22	2.22	0.40

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	326/364 (90%)	318 (98%)	7 (2%)	1 (0%)	36	35
1	B	322/364 (88%)	313 (97%)	8 (2%)	1 (0%)	36	35
All	All	648/728 (89%)	631 (97%)	15 (2%)	2 (0%)	36	35

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	54	ASN
1	A	100	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	273/319 (86%)	272 (100%)	1 (0%)	84	89
1	B	270/319 (85%)	266 (98%)	4 (2%)	57	64
All	All	543/638 (85%)	538 (99%)	5 (1%)	70	78

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	290	VAL
1	B	45	ASN
1	B	62	THR
1	B	153	SER

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	300	ASN

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	26	ASN
1	A	167	ASN
1	A	171	ASN
1	B	5	ASN
1	B	66	HIS
1	B	214	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

2 non-standard protein/DNA/RNA residues 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	DAL	F	1	2	3,4,5	0.67	0	2,4,6	0.76	0
2	DAL	F	2	2	5,5,5	1.18	1 (20%)	6,6,6	1.40	2 (33%)

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	DAL	F	1	2	-	0/1/2/4	-
2	DAL	F	2	2	-	0/4/4/4	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	2	DAL	OXT-C	-2.20	1.23	1.30

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	2	DAL	OXT-C-O	-2.66	118.05	124.08
2	F	2	DAL	OXT-C-CA	2.02	120.75	113.79

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 6 are monoatomic - leaving 2 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
5	ADP	A	403	4	28,29,29	1.43	4 (14%)	43,45,45	1.85	8 (18%)
6	ATP	B	405	4,3	32,33,33	1.47	7 (21%)	48,52,52	1.76	9 (18%)

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
5	ADP	A	403	4	-	1/16/32/32	0/3/3/3
6	ATP	B	405	4,3	-	4/22/38/38	0/3/3/3

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	B	405	ATP	C5-C4	4.79	1.47	1.39
5	A	403	ADP	C5-C4	4.73	1.47	1.39
5	A	403	ADP	C5-C6	2.74	1.48	1.41
6	B	405	ATP	C5-C6	2.71	1.48	1.41
5	A	403	ADP	C5-N7	-2.35	1.34	1.39
6	B	405	ATP	PB-O3A	2.35	1.62	1.59
6	B	405	ATP	C8-N7	2.34	1.36	1.31
5	A	403	ADP	C8-N7	2.34	1.36	1.31
6	B	405	ATP	C5-N7	-2.31	1.34	1.39
6	B	405	ATP	PB-O3B	2.22	1.61	1.59
6	B	405	ATP	PA-O3A	2.05	1.61	1.59

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	B	405	ATP	C5-C4-N3	-6.02	118.43	126.72
5	A	403	ADP	C5-C4-N3	-5.93	118.56	126.72
6	B	405	ATP	N3-C4-N9	4.75	135.24	127.17
5	A	403	ADP	N3-C4-N9	4.70	135.17	127.17
5	A	403	ADP	C2-N3-C4	3.71	120.89	111.83
6	B	405	ATP	C2-N3-C4	3.70	120.87	111.83
6	B	405	ATP	C4-C5-N7	-3.52	106.56	110.58
5	A	403	ADP	C4-C5-N7	-3.45	106.64	110.58
5	A	403	ADP	N3-C2-N1	-3.18	123.77	128.58
6	B	405	ATP	N3-C2-N1	-3.11	123.88	128.58
6	B	405	ATP	C5-N7-C8	2.60	107.54	103.45
5	A	403	ADP	C4-N9-C8	2.60	108.47	105.74
5	A	403	ADP	C5-N7-C8	2.57	107.48	103.45
6	B	405	ATP	C4-N9-C8	2.51	108.37	105.74
5	A	403	ADP	C3'-C2'-C1'	2.28	105.77	101.46
6	B	405	ATP	C2'-C1'-N9	-2.17	107.92	113.30
6	B	405	ATP	C3'-C2'-C1'	2.04	105.32	101.46

There are no chirality outliers.

All (5) torsion outliers are listed below:

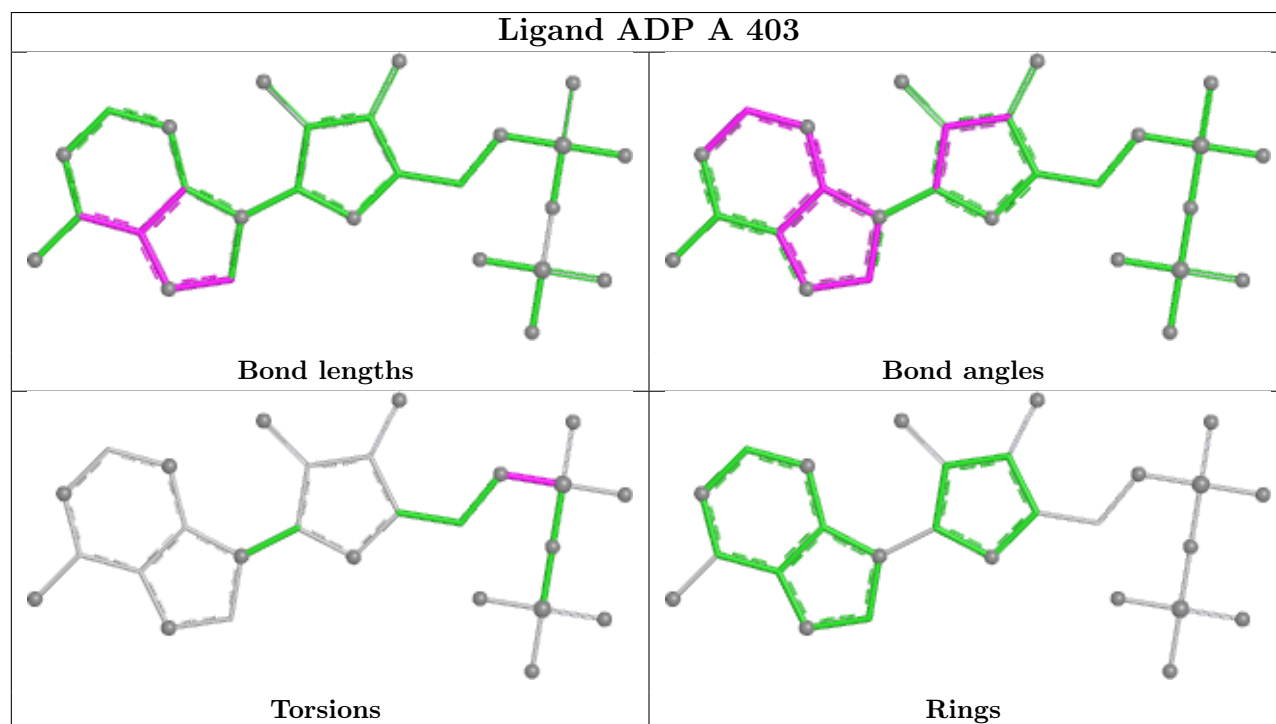
Mol	Chain	Res	Type	Atoms
6	B	405	ATP	C5'-O5'-PA-O3A
5	A	403	ADP	C5'-O5'-PA-O1A
6	B	405	ATP	C5'-O5'-PA-O1A
6	B	405	ATP	PB-O3A-PA-O2A
6	B	405	ATP	PB-O3A-PA-O1A

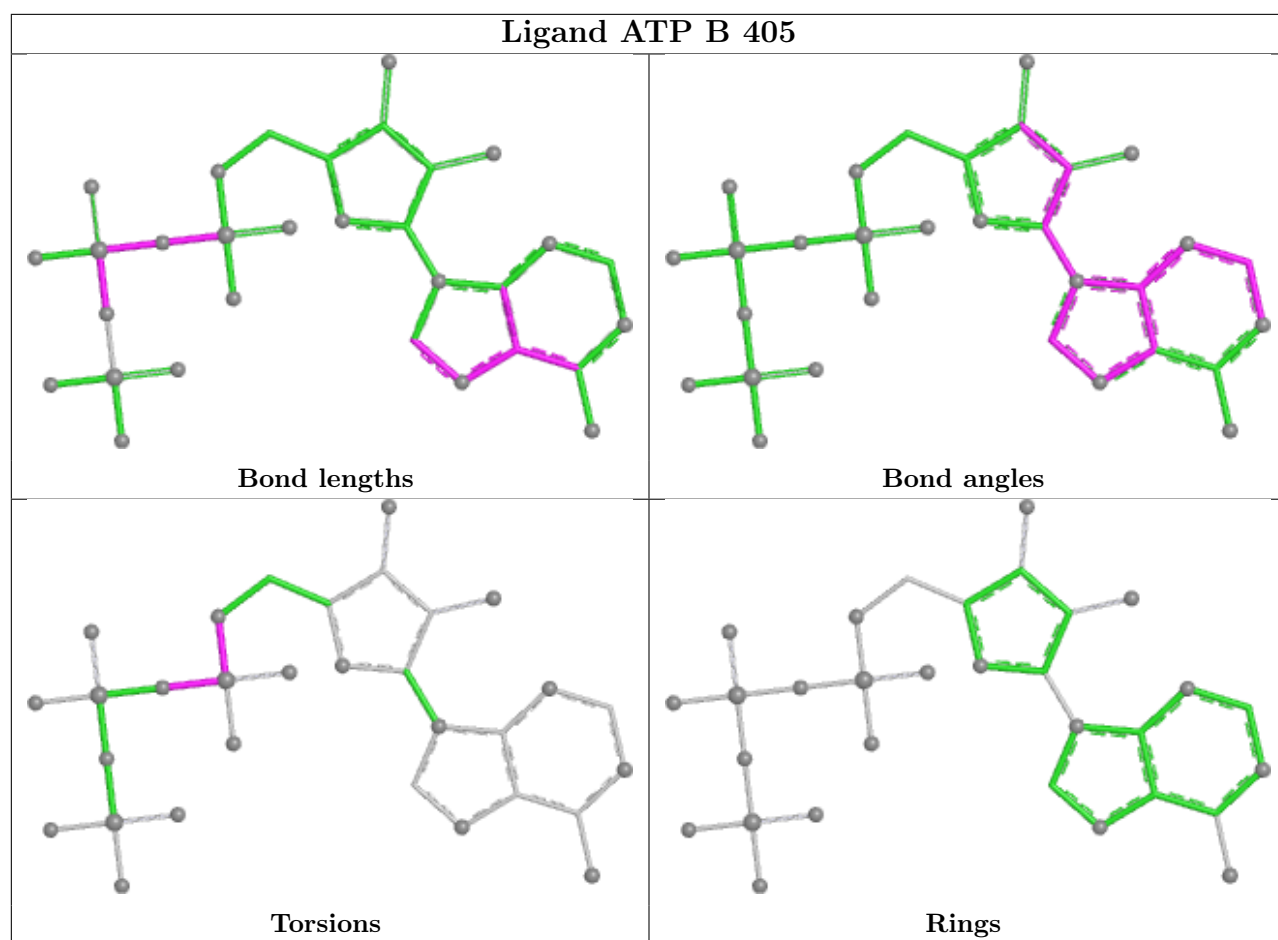
There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	403	ADP	1	0
6	B	405	ATP	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.





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	332/364 (91%)	0.56	20 (6%) 27 26	17, 28, 48, 56	0
1	B	330/364 (90%)	1.00	47 (14%) 6 5	19, 33, 52, 62	0
2	F	0/2	-	-	-	-
All	All	662/730 (90%)	0.78	67 (10%) 12 11	17, 30, 50, 62	0

All (67) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	90	ALA	4.3
1	B	92	PHE	4.0
1	A	206	PHE	4.0
1	A	180	ASN	3.7
1	B	65	LEU	3.6
1	B	76	SER	3.6
1	B	271	LEU	3.5
1	B	56	THR	3.4
1	A	257	VAL	3.4
1	A	184	SER	3.2
1	B	59	ILE	3.1
1	B	328	LEU	2.9
1	A	85	GLY	2.9
1	A	194	ALA	2.9
1	B	86	GLN	2.9
1	B	255	GLY	2.9
1	A	240	VAL	2.8
1	B	8	ILE	2.8
1	A	73	LEU	2.8
1	B	62	THR	2.8
1	B	55	ILE	2.7
1	B	148	ILE	2.7
1	A	15	ALA	2.7

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	60	LYS	2.7
1	B	230	TYR	2.7
1	B	257	VAL	2.7
1	B	254	ASP	2.7
1	A	84	SER	2.6
1	B	158	TYR	2.6
1	B	334	ILE	2.6
1	B	316	PHE	2.6
1	A	83	SER	2.6
1	B	6	ILE	2.6
1	B	27	VAL	2.6
1	A	182	GLY	2.6
1	B	36	TYR	2.5
1	B	37	HIS	2.5
1	B	327	GLY	2.5
1	A	347	ASP	2.4
1	B	330	TYR	2.4
1	A	181	LEU	2.4
1	A	97	GLY	2.4
1	A	98	PRO	2.4
1	B	57	ALA	2.3
1	B	237	GLY	2.3
1	B	263	ALA	2.3
1	B	79	LEU	2.3
1	B	117	VAL	2.3
1	B	115	PRO	2.3
1	B	335	THR	2.2
1	B	322	LEU	2.2
1	A	183	SER	2.2
1	B	61	SER	2.2
1	B	258	GLN	2.2
1	B	40	ILE	2.1
1	B	91	VAL	2.1
1	B	95	LEU	2.1
1	B	153	SER	2.1
1	B	63	ASP	2.1
1	B	31	ILE	2.1
1	B	235	TRP	2.1
1	B	74	GLU	2.0
1	B	53	ASN	2.0
1	B	88	TYR	2.0
1	A	47	GLY	2.0

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	158	TYR	2.0
1	A	230	TYR	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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	DAL	F	1	5/6	0.89	0.09	23,24,30,32	0
2	DAL	F	2	6/6	0.91	0.09	28,31,32,32	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

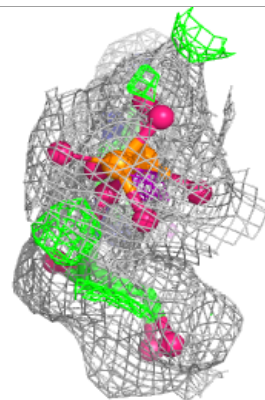
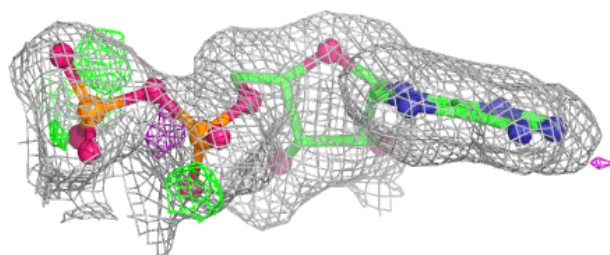
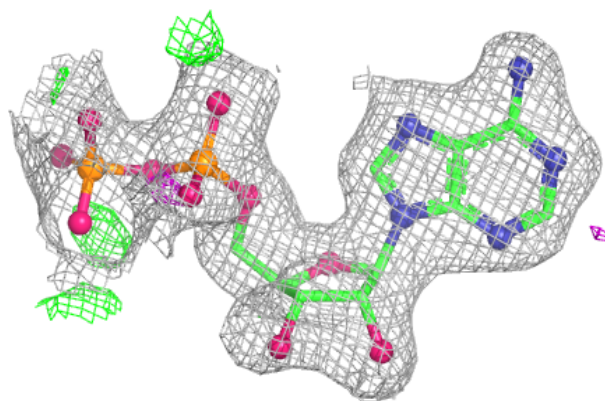
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	ADP	A	403	27/27	0.88	0.10	17,28,45,46	0
3	K	B	401	1/1	0.91	0.08	29,29,29,29	1
4	MG	A	402	1/1	0.92	0.05	28,28,28,28	1
3	K	B	402	1/1	0.92	0.06	23,23,23,23	0
6	ATP	B	405	31/31	0.93	0.10	17,25,30,36	0
3	K	B	403	1/1	0.94	0.06	25,25,25,25	1
3	K	A	401	1/1	0.98	0.04	25,25,25,25	1
4	MG	B	404	1/1	0.98	0.07	19,19,19,19	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.

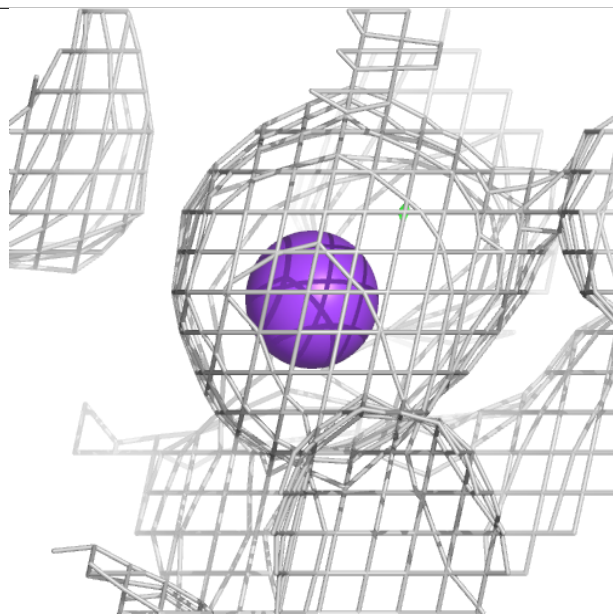
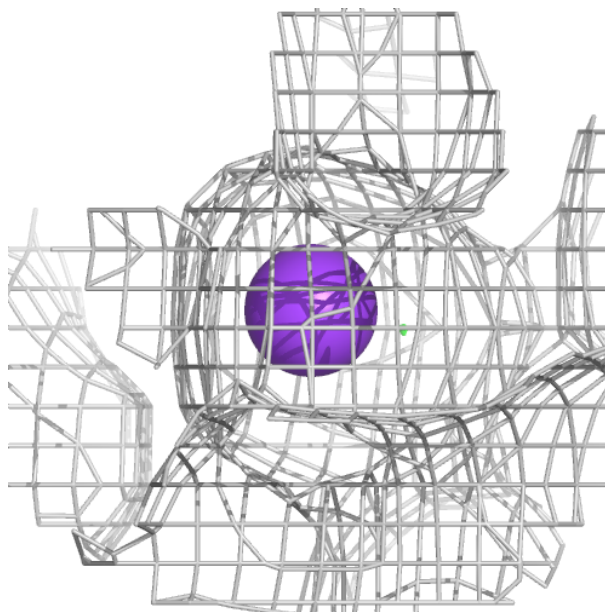
Electron density around ADP A 403:

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



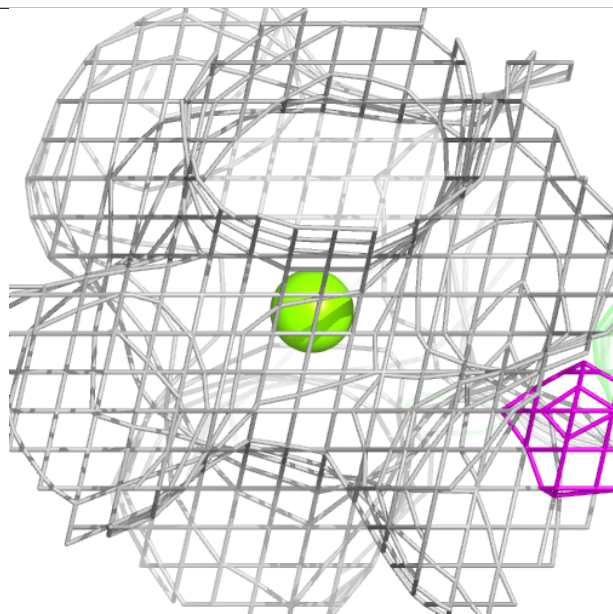
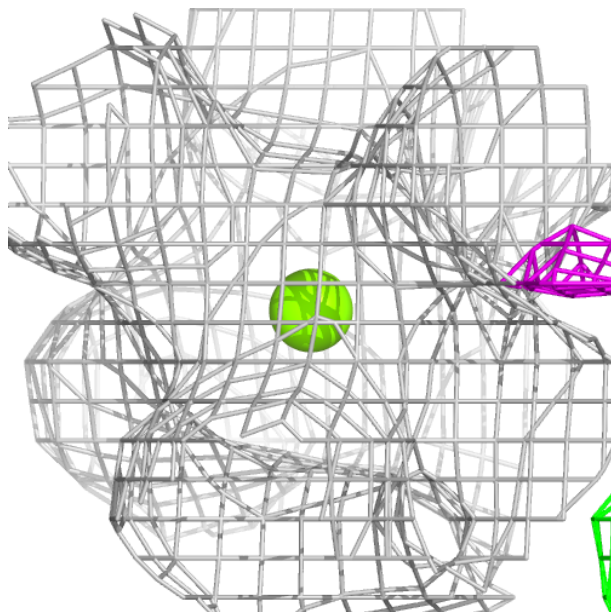
Electron density around K B 401:

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



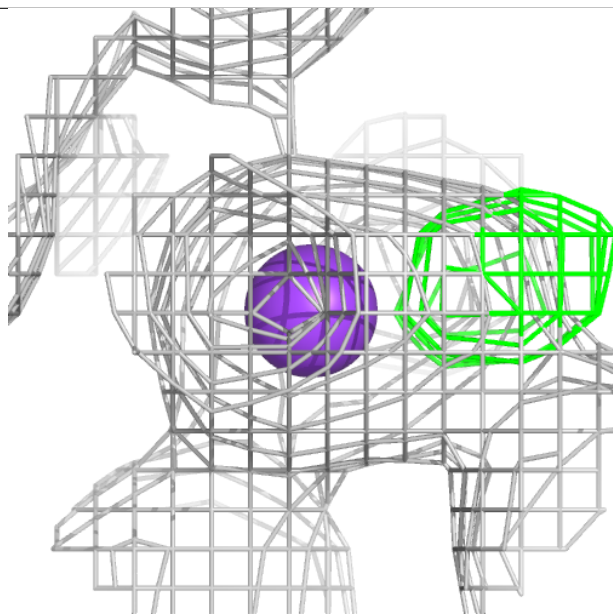
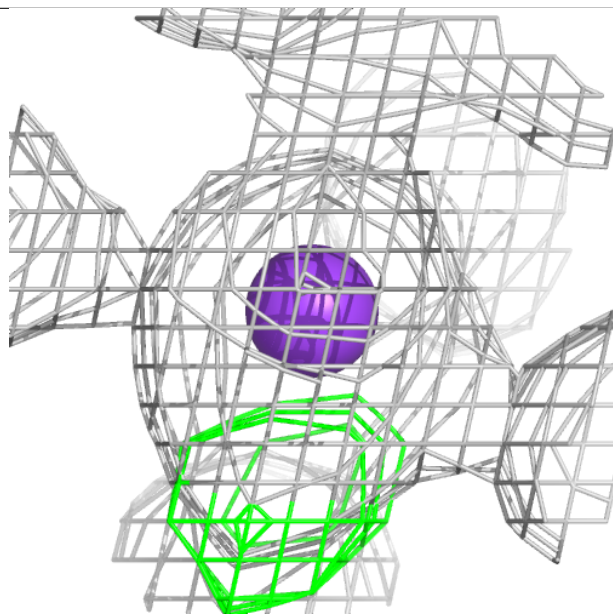
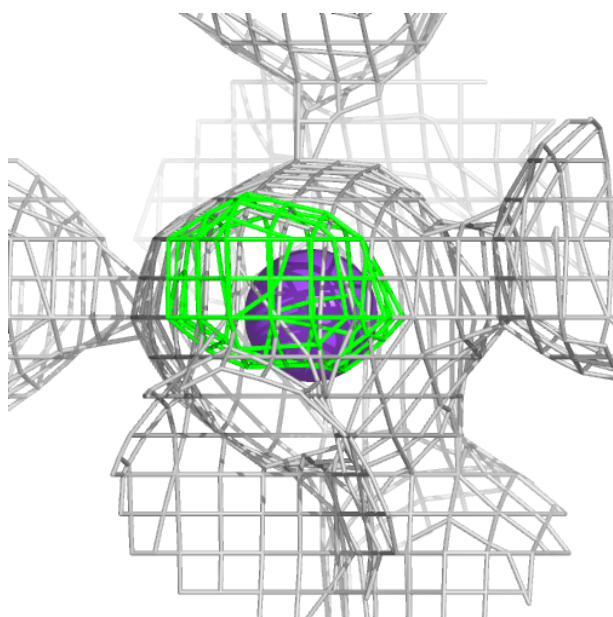
Electron density around MG A 402:

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



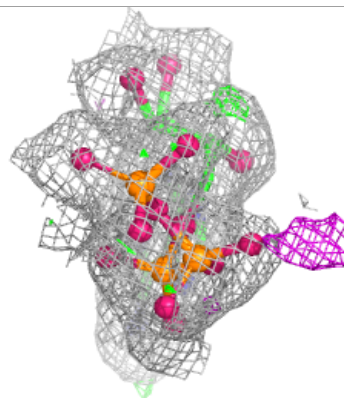
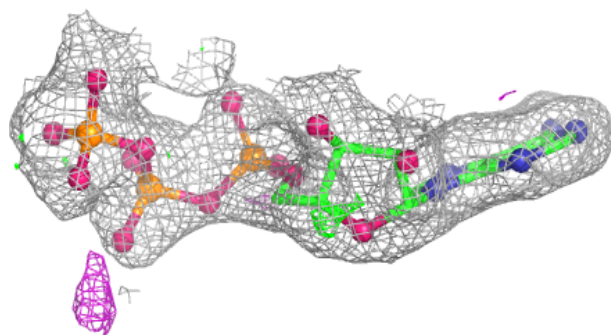
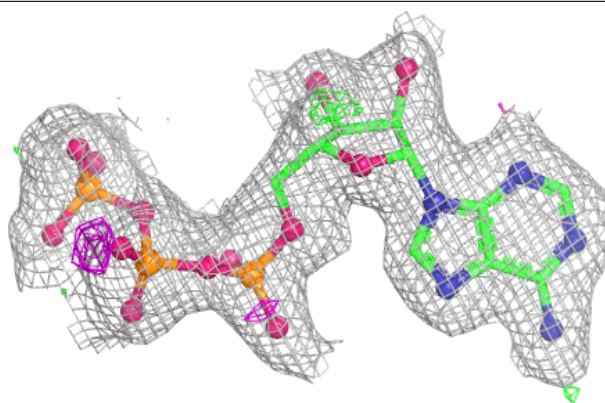
Electron density around K B 402:

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



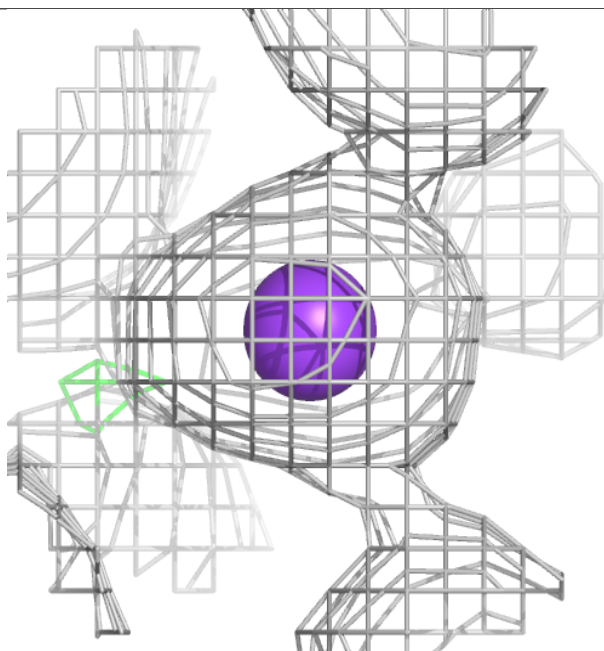
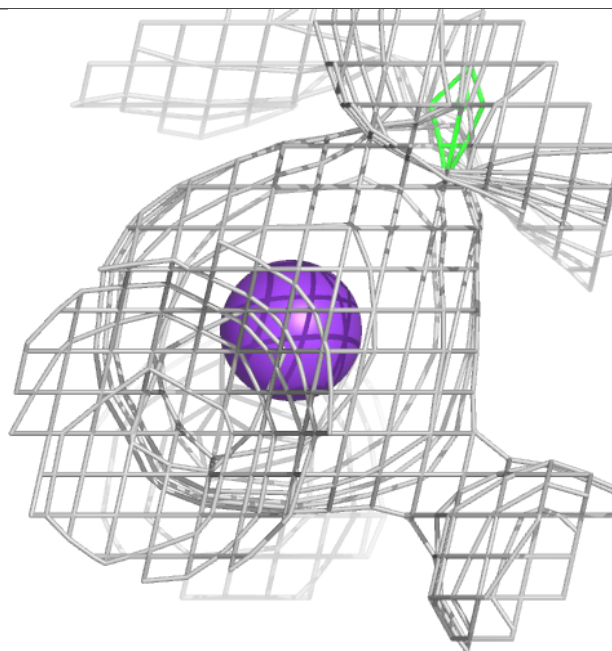
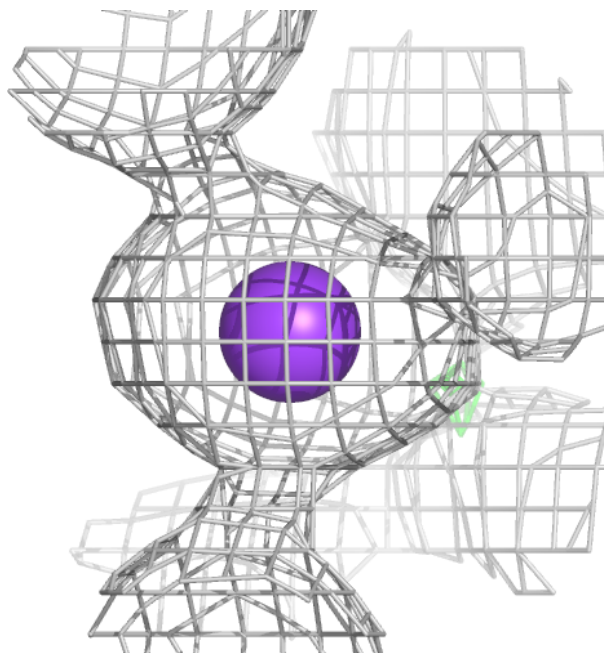
Electron density around ATP B 405:

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



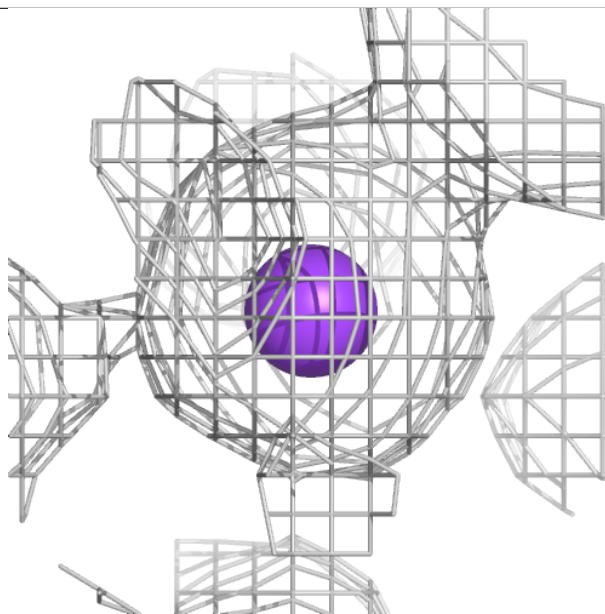
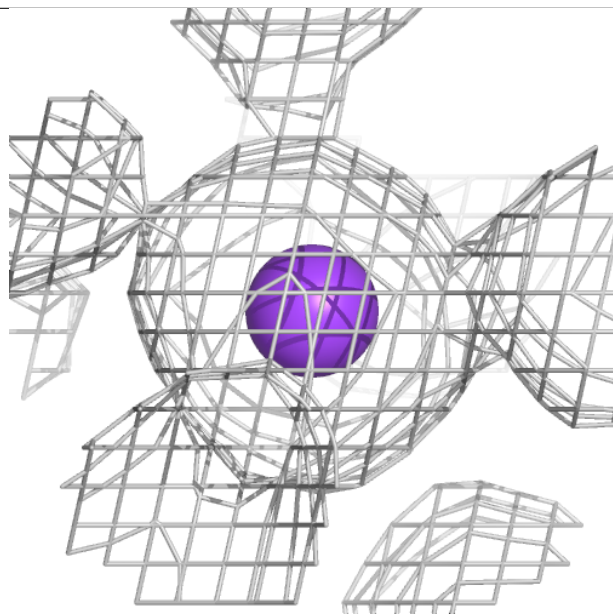
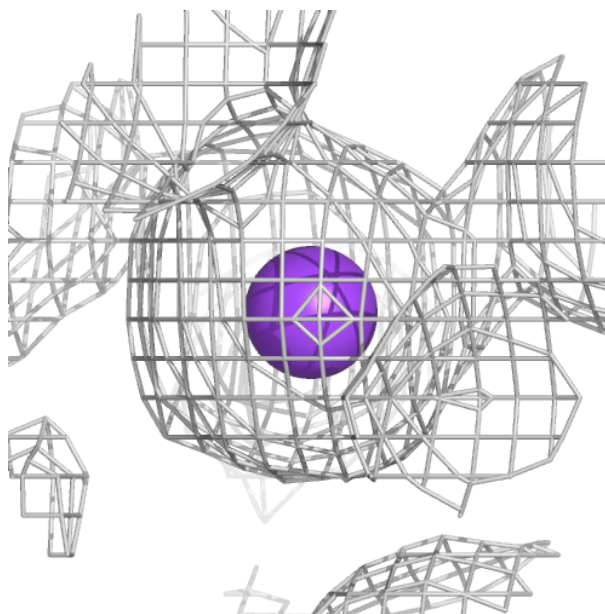
Electron density around K B 403:

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



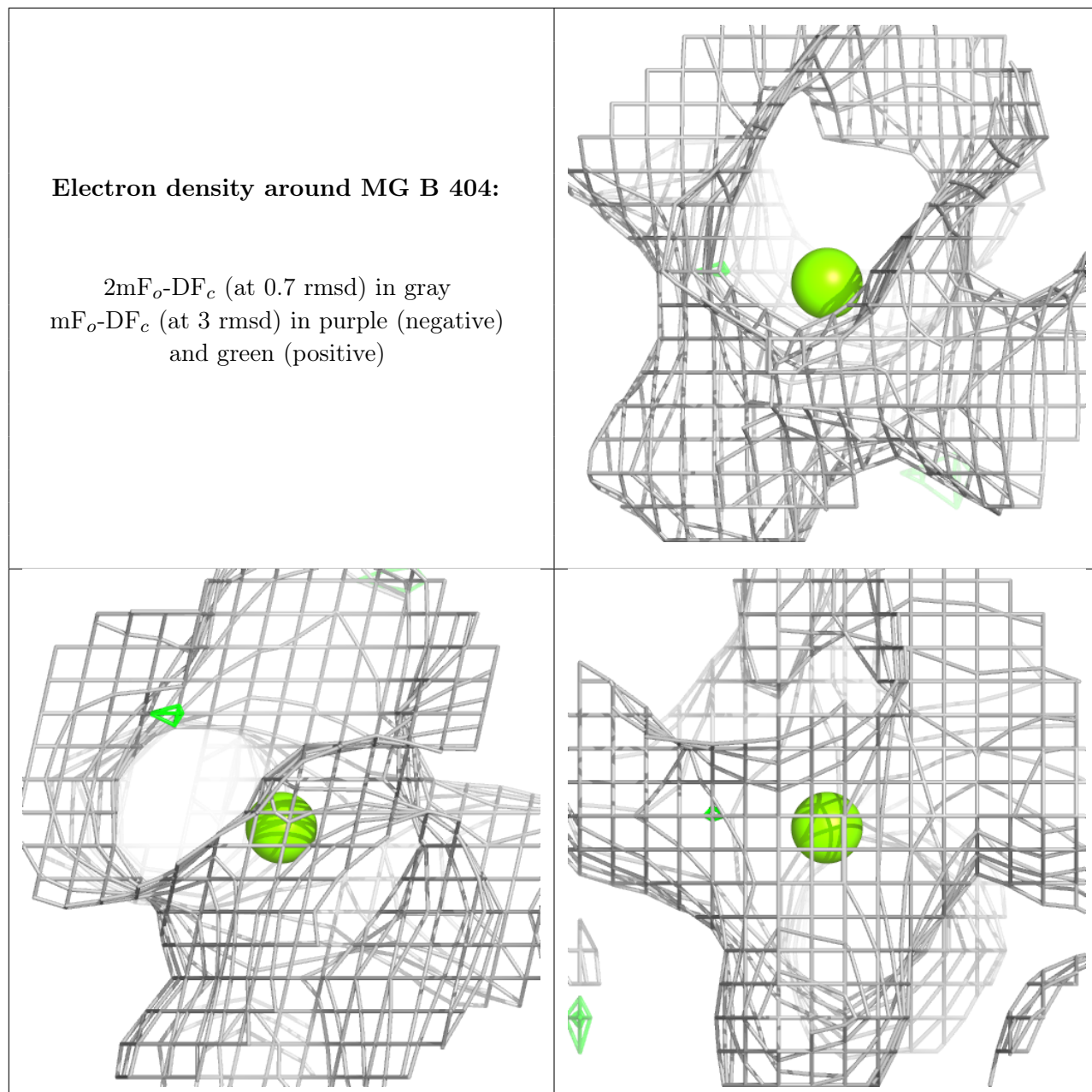
Electron density around K A 401:

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



Electron density around MG B 404:

$2mF_o-DF_c$ (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.