



# Full wwPDB X-ray Structure Validation Report ⓘ

Oct 21, 2025 – 06:14 PM EDT

PDB ID : 9DSV / pdb\_00009dsv  
Title : Thermotoga maritima threonylcarbamoyl adenylate synthase (TsaC2) in complex with N-carboxy-L-threonine, magnesium and ATP  
Authors : Kutshuashvili, A.; Swairjo, M.A.  
Deposited on : 2024-09-29  
Resolution : 2.01 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](mailto: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>.

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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	:	1.1.7 (2018)
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
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)	:	2.46

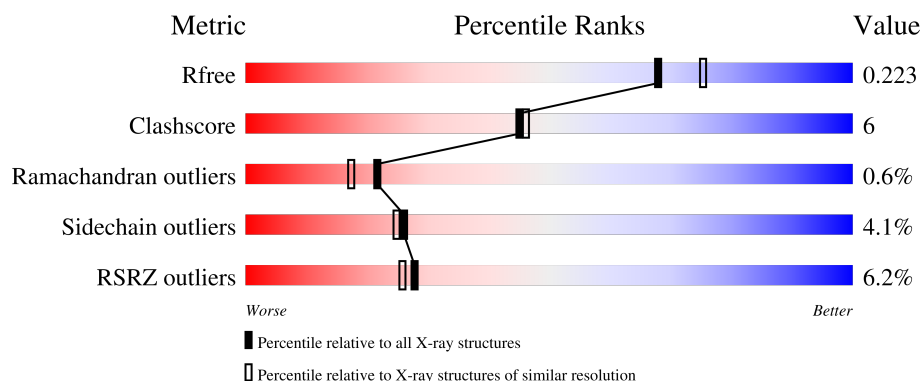
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.01 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))
$R_{free}$	164625	9409 (2.00-2.00)
Clashscore	180529	10737 (2.00-2.00)
Ramachandran outliers	177936	10628 (2.00-2.00)
Sidechain outliers	177891	10627 (2.00-2.00)
RSRZ outliers	164620	9409 (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  $\geq 3$ , 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	340	
1	B	340	

## 2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 5721 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.

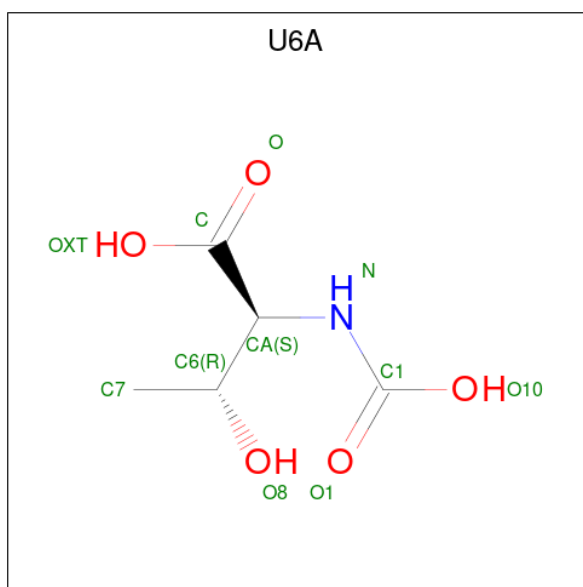
- Molecule 1 is a protein called Threonylcarbamoyl-AMP synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	323	Total	C	N	O	S	0	7	0
			2614	1706	434	465	9			
1	B	323	Total	C	N	O	S	0	11	0
			2667	1732	445	481	9			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-4	GLY	-	expression tag	UNP Q9WZV6
A	-3	SER	-	expression tag	UNP Q9WZV6
A	-2	HIS	-	expression tag	UNP Q9WZV6
A	-1	MET	-	expression tag	UNP Q9WZV6
A	0	ALA	-	expression tag	UNP Q9WZV6
A	1	SER	-	expression tag	UNP Q9WZV6
B	-4	GLY	-	expression tag	UNP Q9WZV6
B	-3	SER	-	expression tag	UNP Q9WZV6
B	-2	HIS	-	expression tag	UNP Q9WZV6
B	-1	MET	-	expression tag	UNP Q9WZV6
B	0	ALA	-	expression tag	UNP Q9WZV6
B	1	SER	-	expression tag	UNP Q9WZV6

- Molecule 2 is N-carboxy-L-threonine (CCD ID: U6A) (formula:  $C_5H_9NO_5$ ) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			11	5	1	5		
2	B	1	Total	C	N	O	11	1
			22	10	2	10		

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

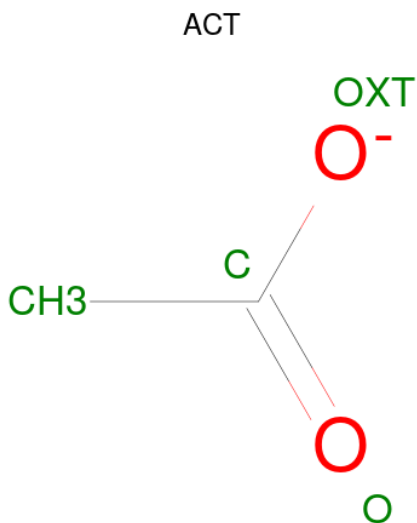
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Mg	0	0
			1	1		
3	B	1	Total	Mg	0	0
			1	1		

- Molecule 4 is ADENOSINE-5'-TRIPHOSPHATE (CCD ID: ATP) (formula: C<sub>10</sub>H<sub>16</sub>N<sub>5</sub>O<sub>13</sub>P<sub>3</sub>).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total 31	C 10	N 5	O 13	P 3	0	0
4	B	1	Total 62	C 20	N 10	O 26	P 6	31	1

- Molecule 5 is ACETATE ION (CCD ID: ACT) (formula:  $\text{C}_2\text{H}_3\text{O}_2$ ).



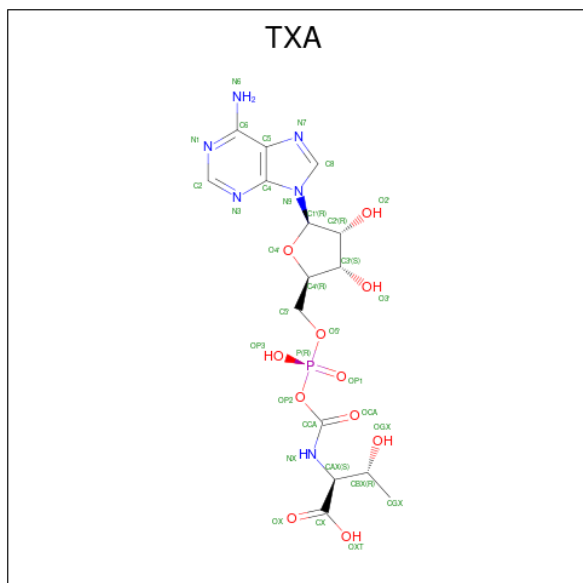
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total 4	C 2	O 2	0	0
5	A	1	Total 4	C 2	O 2	0	0

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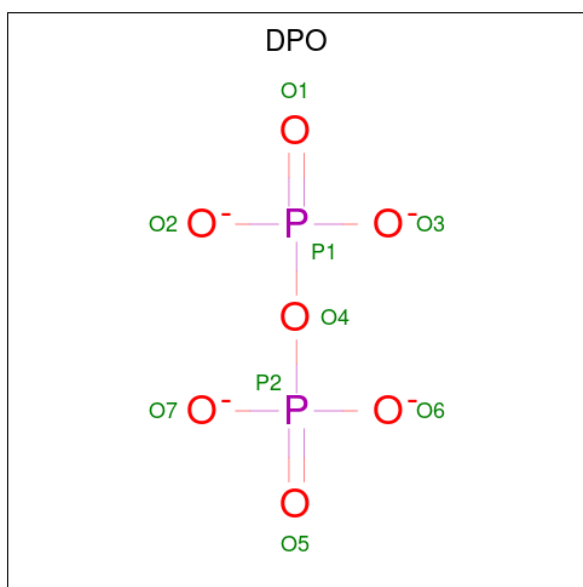
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	B	1	Total C O 4 2 2	0	0
5	B	1	Total C O 4 2 2	0	0

- Molecule 6 is threonylcarbamoyladenylate (CCD ID: TXA) (formula:  $\text{C}_{15}\text{H}_{21}\text{N}_6\text{O}_{11}\text{P}$ ) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	B	1	Total 66	C 30	N 12	O 22	P 2	33	1

- Molecule 7 is DIPHOSPHATE (CCD ID: DPO) (formula:  $\text{O}_7\text{P}_2$ ) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	B	1	Total	O	P	9	1
			18	14	4		

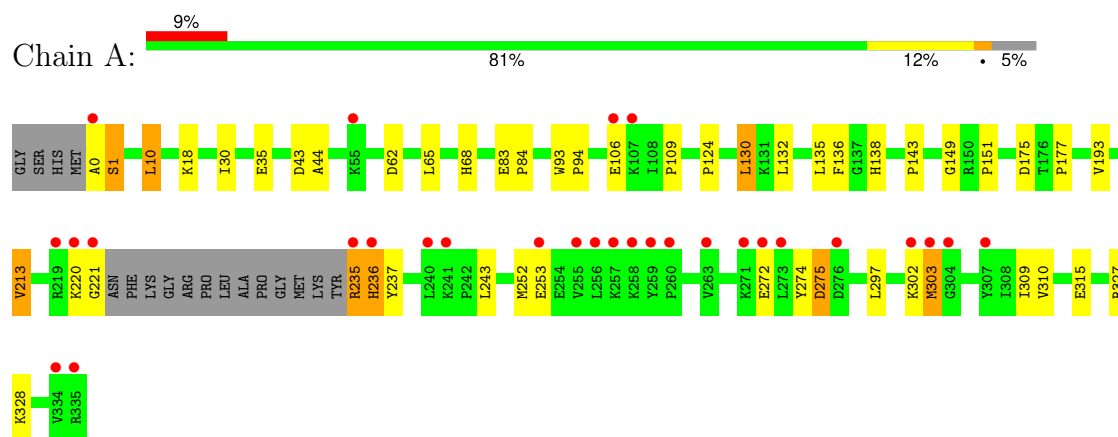
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	79	Total	O	0	0
			79	79		
8	B	133	Total	O	0	0
			133	133		

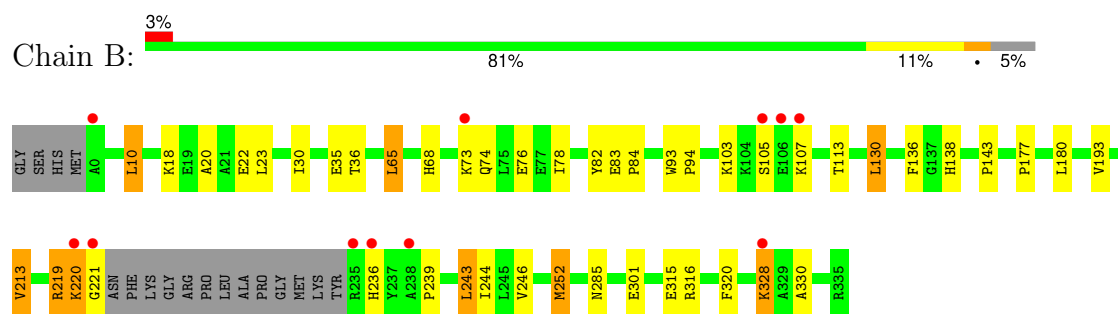
### 3 Residue-property plots [i](#)

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: Threonylcarbamoyl-AMP synthase



#### • Molecule 1: Threonylcarbamoyl-AMP synthase





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	153.60Å 153.60Å 86.55Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	44.38 – 2.01 44.38 – 2.01	Depositor EDS
% Data completeness (in resolution range)	100.0 (44.38-2.01) 100.0 (44.38-2.01)	Depositor EDS
$R_{merge}$	0.24	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.33 (at 2.01Å)	Xtriage
Refinement program	REFMAC 5.8.0430	Depositor
R, $R_{free}$	0.193 , 0.216 0.202 , 0.223	Depositor DCC
$R_{free}$ test set	3941 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	32.8	Xtriage
Anisotropy	0.352	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 40.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.029 for -h,-k,l	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	5721	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.26% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: MG, TXA, U6A, DPO, ACT, 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	0.49	0/2683	0.93	2/3632 (0.1%)
1	B	0.49	0/2727	0.92	0/3691
All	All	0.49	0/5410	0.92	2/7323 (0.0%)

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	2

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
1	A	62	ASP	CA-CB-CG	5.12	117.72	112.60
1	A	175	ASP	CA-CB-CG	5.02	117.62	112.60

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	219[B]	ARG	Mainchain
1	B	316	ARG	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	2614	0	2745	33	0
1	B	2667	0	2770	34	0
2	A	11	0	0	1	0
2	B	22	0	0	1	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	31	0	12	1	0
4	B	62	0	20	1	0
5	A	8	0	6	1	0
5	B	8	0	6	0	0
6	B	66	0	30	2	0
7	B	18	0	0	0	0
8	A	79	0	0	1	0
8	B	133	0	0	1	0
All	All	5721	0	5589	64	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 6.

All (64) 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:221:GLY:HA2	1:B:328:LYS:HE2	1.70	0.74
1:A:18:LYS:HG2	1:A:135[A]:LEU:HD21	1.75	0.67
1:B:76:GLU:HG2	1:B:82:TYR:OH	1.95	0.66
1:B:103:LYS:HE2	1:B:113:THR:HG21	1.80	0.64
1:B:136:PHE:CE2	1:B:138:HIS:HB2	2.38	0.59
1:B:243[B]:LEU:HD23	1:B:243[B]:LEU:C	2.29	0.58
1:B:243[A]:LEU:C	1:B:243[A]:LEU:CD1	2.78	0.57
1:A:10:LEU:CD1	1:B:10:LEU:CD1	2.82	0.57
1:B:220:LYS:O	1:B:221:GLY:C	2.49	0.56
1:A:221:GLY:C	1:A:328:LYS:HE3	2.30	0.55
1:A:193:VAL:HG13	1:A:213[A]:VAL:HG22	1.88	0.55
1:A:18:LYS:CG	1:A:135[A]:LEU:HD21	2.36	0.54
1:B:68:HIS:HE1	2:B:404[B]:U6A:O8	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:10:LEU:CD1	1:B:10:LEU:HD12	2.40	0.52
1:A:136:PHE:CE2	1:A:138:HIS:HB2	2.45	0.52
1:A:93:TRP:HA	1:A:94:PRO:C	2.35	0.52
1:B:143:PRO:HB3	4:B:405[B]:ATP:O1A	2.09	0.52
1:A:243:LEU:HD11	1:A:310:VAL:CG2	2.40	0.51
1:A:10:LEU:HD12	1:B:10:LEU:CD1	2.41	0.51
5:A:404:ACT:H2	8:A:572:HOH:O	2.10	0.51
1:A:30:ILE:HD11	1:A:132:LEU:HD21	1.93	0.50
1:A:252[A]:MET:HE1	1:A:309[A]:ILE:CG2	2.41	0.50
1:B:68:HIS:HE1	6:B:401[A]:TXA:OGX	1.94	0.50
1:B:65[A]:LEU:HG	1:B:143:PRO:HD3	1.93	0.49
1:A:315:GLU:CD	1:A:327:ARG:HH22	2.20	0.49
1:A:68:HIS:HE1	2:A:401:U6A:O8	1.94	0.49
1:B:246:VAL:HG21	1:B:252[A]:MET:CE	2.43	0.49
1:B:35:GLU:HA	1:B:177:PRO:HD2	1.95	0.48
1:A:35:GLU:HA	1:A:177:PRO:HD2	1.96	0.47
1:A:274:TYR:O	1:A:275:ASP:HB2	2.14	0.47
1:B:93:TRP:HA	1:B:94:PRO:C	2.39	0.47
1:A:83:GLU:N	1:A:84:PRO:HD2	2.30	0.47
1:A:235:ARG:O	1:A:236:HIS:HB2	2.15	0.46
1:A:43:ASP:OD1	1:A:138:HIS:HD2	2.00	0.45
1:B:20:ALA:HB1	1:B:30:ILE:HD13	1.98	0.45
1:B:246:VAL:HG11	1:B:252[A]:MET:HE3	1.97	0.45
1:A:315:GLU:CG	1:A:327:ARG:HH22	2.29	0.45
1:A:10:LEU:HD11	1:B:10:LEU:HD12	1.99	0.45
1:A:315:GLU:OE1	1:A:327:ARG:NH1	2.49	0.44
1:A:149:GLY:HA2	1:A:236:HIS:H	1.81	0.44
1:A:83:GLU:N	1:A:84:PRO:CD	2.81	0.44
1:B:83:GLU:N	1:B:84:PRO:CD	2.80	0.44
1:B:243[A]:LEU:C	1:B:243[A]:LEU:HD12	2.44	0.43
1:A:0:ALA:O	1:A:1:SER:HB3	2.18	0.43
1:B:130[A]:LEU:HD12	1:B:130[A]:LEU:HA	1.84	0.43
1:B:78:ILE:HA	1:B:105:SER:HB2	2.00	0.42
1:A:243:LEU:HD11	1:A:310:VAL:HG23	2.01	0.42
1:A:252[A]:MET:HE1	1:A:309[A]:ILE:HG21	2.01	0.42
1:A:44:ALA:O	1:A:109:PRO:HD2	2.19	0.42
1:A:65[A]:LEU:HG	1:A:143:PRO:HD3	2.01	0.42
1:B:74:GLN:NE2	8:B:510:HOH:O	2.53	0.42
1:B:243[A]:LEU:HD13	1:B:244:ILE:N	2.35	0.42
1:B:243[B]:LEU:HD22	1:B:330:ALA:HA	2.01	0.41
1:B:315:GLU:HG2	1:B:320:PHE:CD1	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:143:PRO:HB3	4:A:403:ATP:O1A	2.20	0.41
1:B:239:PRO:HA	1:B:301:GLU:OE1	2.21	0.41
1:B:18:LYS:O	1:B:22:GLU:HG2	2.20	0.41
1:B:105:SER:OG	1:B:107:LYS:HB2	2.20	0.41
1:B:193:VAL:HG22	1:B:213:VAL:HG13	2.03	0.41
1:A:124:PRO:O	1:A:130:LEU:HD13	2.20	0.41
1:A:151:PRO:HA	1:A:237:TYR:HA	2.02	0.41
1:B:36:THR:HB	1:B:180:LEU:O	2.21	0.40
1:B:143:PRO:HB3	6:B:401[A]:TXA:OP3	2.17	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	326/340 (96%)	311 (95%)	10 (3%)	5 (2%)	8	4
1	B	331/340 (97%)	328 (99%)	3 (1%)	0	100	100
All	All	657/680 (97%)	639 (97%)	13 (2%)	5 (1%)	22	12

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	236	HIS
1	A	1	SER
1	A	275	ASP
1	A	303[A]	MET
1	A	303[B]	MET

### 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	290/296 (98%)	278 (96%)	12 (4%)	26	25
1	B	295/296 (100%)	276 (94%)	19 (6%)	14	11
All	All	585/592 (99%)	554 (95%)	31 (5%)	26	16

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	10	LEU
1	A	106	GLU
1	A	130	LEU
1	A	213[A]	VAL
1	A	213[B]	VAL
1	A	220	LYS
1	A	235	ARG
1	A	253	GLU
1	A	272	GLU
1	A	297	LEU
1	A	303[A]	MET
1	A	303[B]	MET
1	B	10	LEU
1	B	23[A]	LEU
1	B	23[B]	LEU
1	B	65[A]	LEU
1	B	65[B]	LEU
1	B	73	LYS
1	B	130[A]	LEU
1	B	130[B]	LEU
1	B	213	VAL
1	B	219[A]	ARG
1	B	219[B]	ARG
1	B	220	LYS
1	B	236	HIS
1	B	243[A]	LEU
1	B	243[B]	LEU

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Mol	Chain	Res	Type
1	B	252[A]	MET
1	B	252[B]	MET
1	B	285	ASN
1	B	328	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	68	HIS
1	A	85	HIS
1	A	138	HIS
1	A	158	HIS
1	B	68	HIS
1	B	158	HIS
1	B	285	ASN

### 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](#)

Of 16 ligands modelled in this entry, 2 are monoatomic - leaving 14 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	DPO	B	403[B]	3,4	6,8,8	0.63	0	12,13,13	0.90	0
5	ACT	B	406	-	3,3,3	1.00	0	3,3,3	0.81	0
5	ACT	A	404	-	3,3,3	0.99	0	3,3,3	0.77	0
5	ACT	A	405	-	3,3,3	0.97	0	3,3,3	0.83	0
6	TXA	B	401[A]	3,2,4	30,35,35	0.75	0	36,52,52	0.94	2 (5%)
2	U6A	B	404[A]	6	10,10,10	3.24	3 (30%)	12,13,13	1.71	3 (25%)
2	U6A	A	401	-	10,10,10	2.97	2 (20%)	12,13,13	1.47	2 (16%)
5	ACT	B	407	-	3,3,3	1.05	0	3,3,3	0.79	0
4	ATP	A	403	3	28,33,33	0.77	0	34,52,52	0.87	1 (2%)
2	U6A	B	404[B]	6	10,10,10	2.84	3 (30%)	12,13,13	1.70	2 (16%)
4	ATP	B	405[A]	3,7,6	28,33,33	1.17	3 (10%)	34,52,52	0.80	1 (2%)
7	DPO	B	403[A]	3,4	6,8,8	0.79	0	12,13,13	0.80	0
6	TXA	B	401[B]	3,2,4	30,35,35	0.93	1 (3%)	36,52,52	1.23	2 (5%)
4	ATP	B	405[B]	3,7,6	28,33,33	0.85	0	34,52,52	0.82	1 (2%)

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	DPO	B	403[B]	3,4	-	3/6/6/6	-
6	TXA	B	401[A]	3,2,4	-	7/21/43/43	0/3/3/3
2	U6A	B	404[A]	6	-	1/12/12/12	-
2	U6A	A	401	-	-	0/12/12/12	-
4	ATP	A	403	3	-	8/18/38/38	0/3/3/3
2	U6A	B	404[B]	6	-	2/12/12/12	-
4	ATP	B	405[A]	3,7,6	-	8/18/38/38	0/3/3/3
7	DPO	B	403[A]	3,4	-	2/6/6/6	-
6	TXA	B	401[B]	3,2,4	-	5/21/43/43	0/3/3/3
4	ATP	B	405[B]	3,7,6	-	7/18/38/38	0/3/3/3

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	U6A	C1-N	8.67	1.50	1.35
2	B	404[A]	U6A	C1-N	8.67	1.50	1.35
2	B	404[B]	U6A	C1-N	7.38	1.48	1.35
4	B	405[A]	ATP	PA-O3A	3.76	1.63	1.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	404[A]	U6A	CA-C	3.45	1.58	1.52
2	B	404[B]	U6A	CA-C	3.11	1.57	1.52
2	B	404[A]	U6A	C7-C6	2.77	1.59	1.51
6	B	401[B]	TXA	P-OP2	2.73	1.65	1.60
4	B	405[A]	ATP	PB-O3A	2.37	1.62	1.59
2	A	401	U6A	O1-C1	2.26	1.25	1.21
2	B	404[B]	U6A	C7-C6	2.17	1.58	1.51
4	B	405[A]	ATP	PB-O3B	2.17	1.61	1.59

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	B	401[B]	TXA	OP2-CCA-NX	4.96	112.84	109.91
2	B	404[B]	U6A	O1-C1-N	-4.48	117.51	124.86
2	B	404[A]	U6A	O1-C1-N	-4.36	117.71	124.86
2	A	401	U6A	O1-C1-N	-3.08	119.81	124.86
6	B	401[A]	TXA	OP2-CCA-NX	2.48	111.38	109.91
4	A	403	ATP	C5-C6-N6	2.34	123.87	120.31
4	B	405[A]	ATP	C5-C6-N6	2.33	123.86	120.31
4	B	405[B]	ATP	C5-C6-N6	2.25	123.74	120.31
6	B	401[A]	TXA	C5-C6-N6	2.22	123.69	120.31
6	B	401[B]	TXA	C5-C6-N6	2.15	123.59	120.31
2	B	404[B]	U6A	C-CA-N	2.15	114.89	110.17
2	B	404[A]	U6A	C7-C6-CA	2.15	116.53	112.29
2	A	401	U6A	O-C-CA	-2.12	114.82	121.86
2	B	404[A]	U6A	C-CA-N	2.09	114.77	110.17

There are no chirality outliers.

All (43) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	403	ATP	PB-O3A-PA-O5'
4	A	403	ATP	C5'-O5'-PA-O1A
4	A	403	ATP	C5'-O5'-PA-O3A
4	B	405[A]	ATP	PB-O3B-PG-O3G
4	B	405[A]	ATP	C5'-O5'-PA-O1A
4	B	405[A]	ATP	C5'-O5'-PA-O3A
4	B	405[B]	ATP	PB-O3B-PG-O3G
4	B	405[B]	ATP	C5'-O5'-PA-O1A
4	B	405[B]	ATP	C5'-O5'-PA-O3A
6	B	401[A]	TXA	C5'-O5'-P-OP3
7	B	403[A]	DPO	P2-O4-P1-O3

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Mol	Chain	Res	Type	Atoms
7	B	403[B]	DPO	P2-O4-P1-O3
6	B	401[A]	TXA	O4'-C4'-C5'-O5'
6	B	401[A]	TXA	C3'-C4'-C5'-O5'
6	B	401[B]	TXA	C3'-C4'-C5'-O5'
6	B	401[B]	TXA	O4'-C4'-C5'-O5'
4	A	403	ATP	PB-O3B-PG-O1G
4	A	403	ATP	C5'-O5'-PA-O2A
4	B	405[A]	ATP	C5'-O5'-PA-O2A
4	B	405[B]	ATP	C5'-O5'-PA-O2A
6	B	401[A]	TXA	C5'-O5'-P-OP2
6	B	401[A]	TXA	C5'-O5'-P-OP1
6	B	401[B]	TXA	C5'-O5'-P-OP1
2	B	404[B]	U6A	C7-C6-CA-N
6	B	401[B]	TXA	NX-CAX-CBX-CGX
2	B	404[A]	U6A	O1-C1-N-CA
2	B	404[B]	U6A	O1-C1-N-CA
4	B	405[A]	ATP	PB-O3B-PG-O1G
4	B	405[B]	ATP	PB-O3B-PG-O1G
4	A	403	ATP	PB-O3B-PG-O2G
4	A	403	ATP	PB-O3B-PG-O3G
6	B	401[A]	TXA	NX-CAX-CBX-CGX
7	B	403[A]	DPO	P2-O4-P1-O2
7	B	403[B]	DPO	P2-O4-P1-O2
6	B	401[A]	TXA	CX-CAX-CBX-OGX
6	B	401[B]	TXA	CX-CAX-CBX-OGX
4	B	405[A]	ATP	PA-O3A-PB-O1B
4	B	405[A]	ATP	PA-O3A-PB-O2B
4	A	403	ATP	O4'-C4'-C5'-O5'
4	B	405[A]	ATP	O4'-C4'-C5'-O5'
7	B	403[B]	DPO	P2-O4-P1-O1
4	B	405[B]	ATP	PA-O3A-PB-O1B
4	B	405[B]	ATP	PA-O3A-PB-O2B

There are no ring outliers.

6 monomers are involved in 7 short contacts:

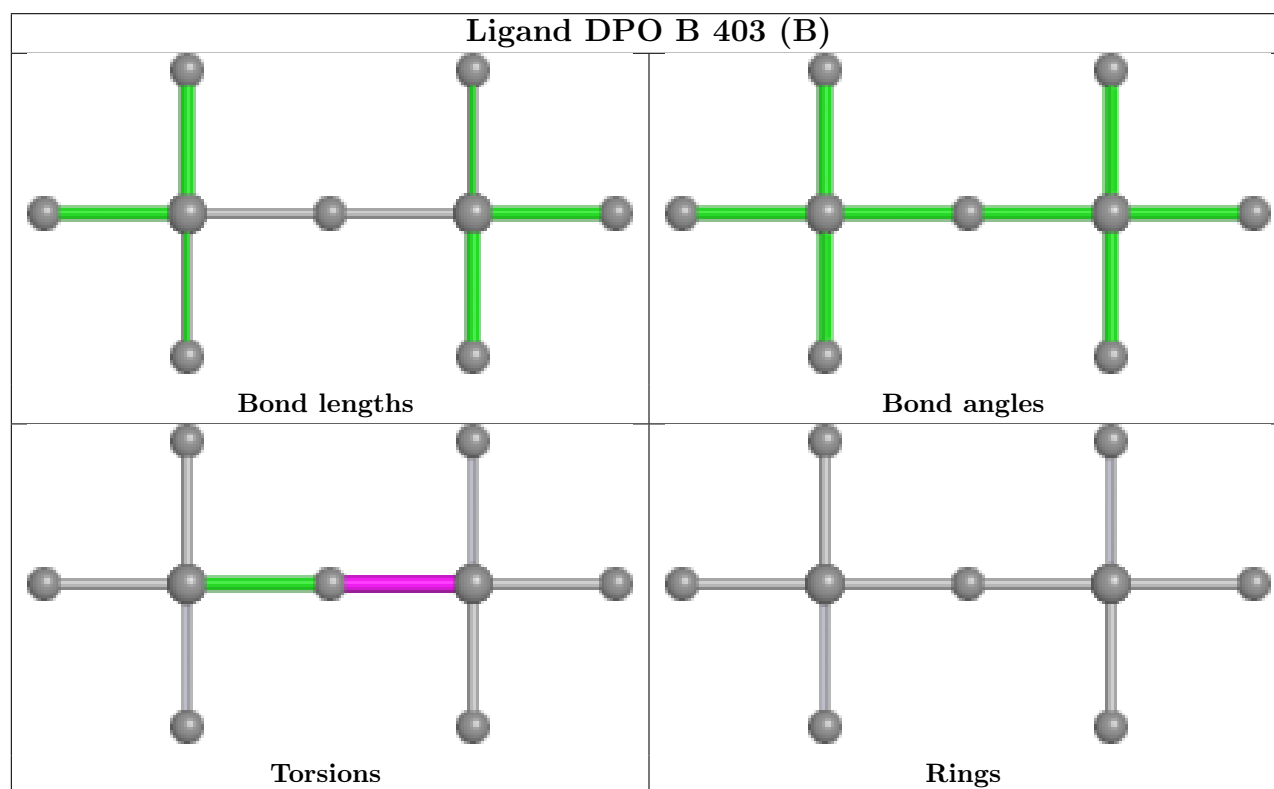
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	404	ACT	1	0
6	B	401[A]	TXA	2	0
2	A	401	U6A	1	0
4	A	403	ATP	1	0
2	B	404[B]	U6A	1	0

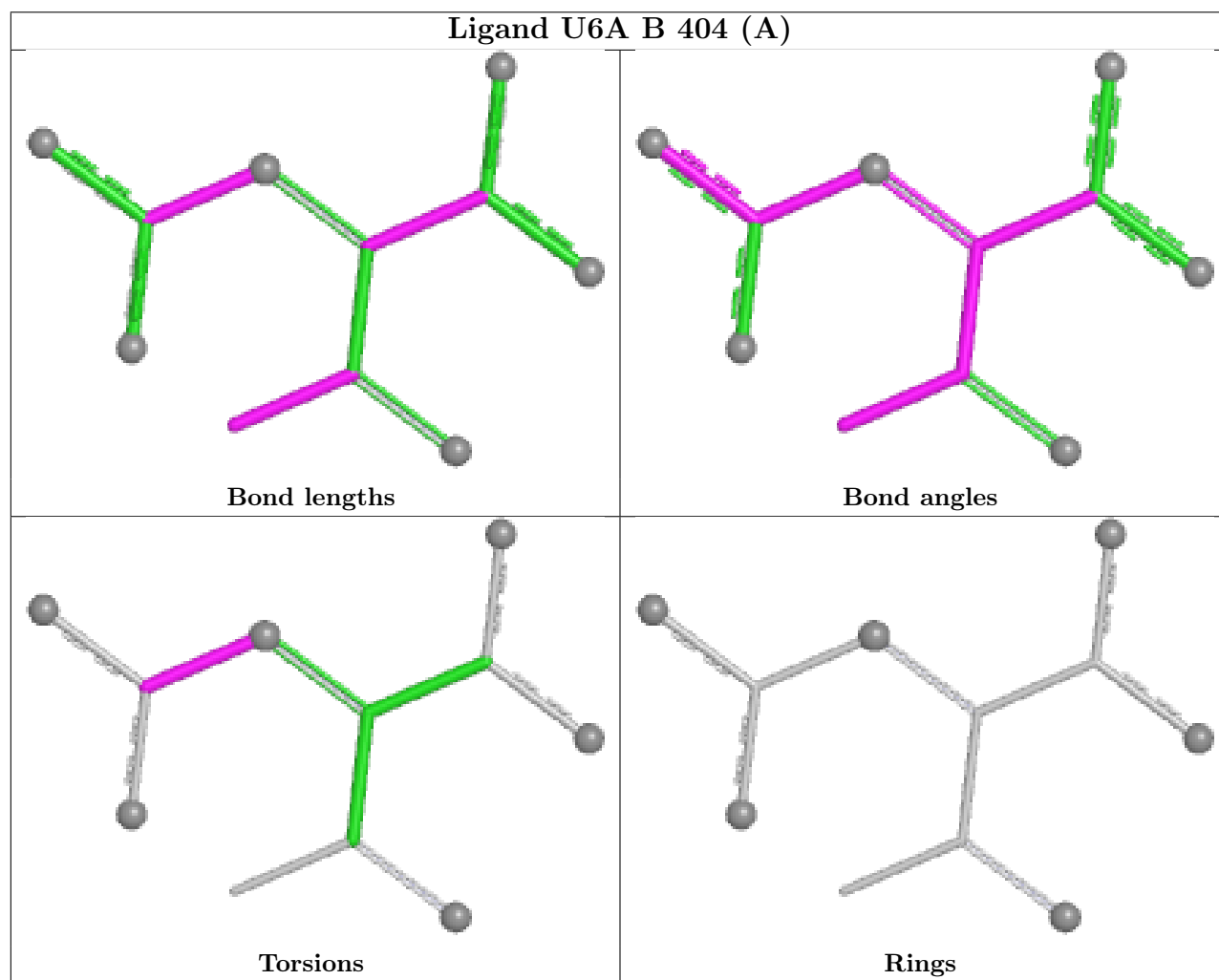
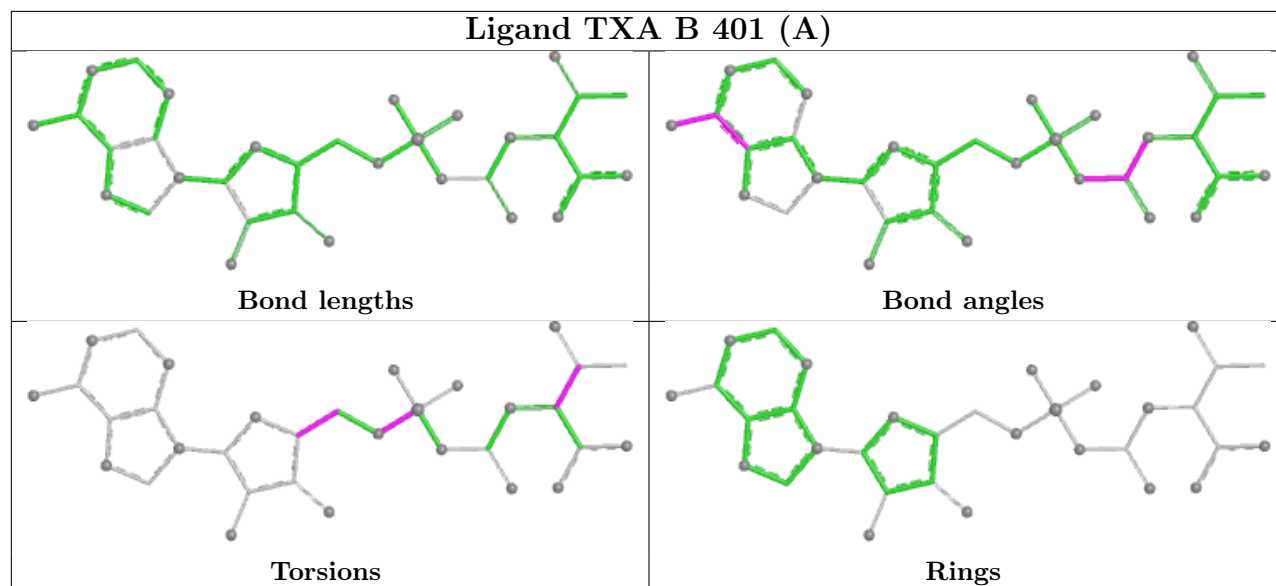
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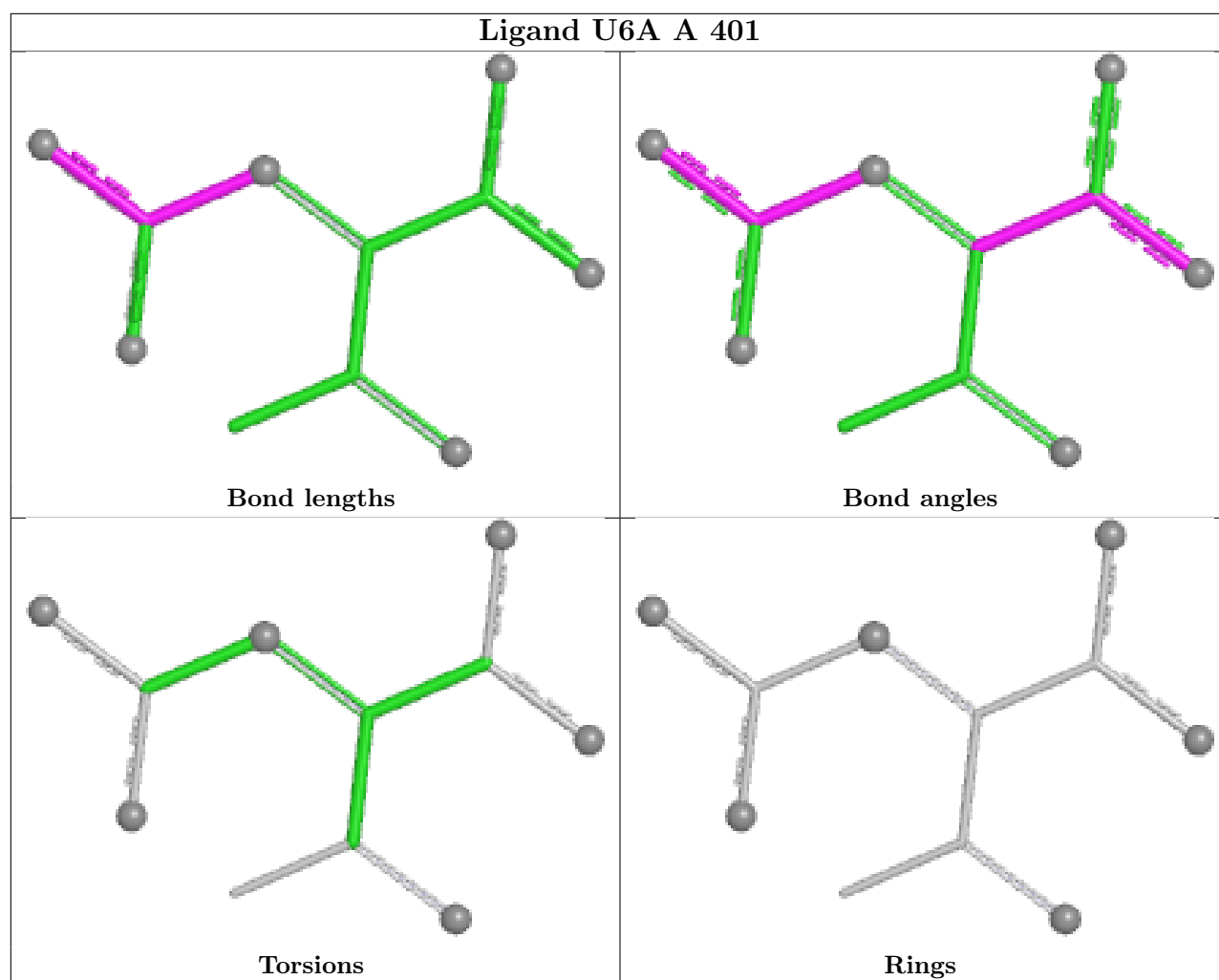
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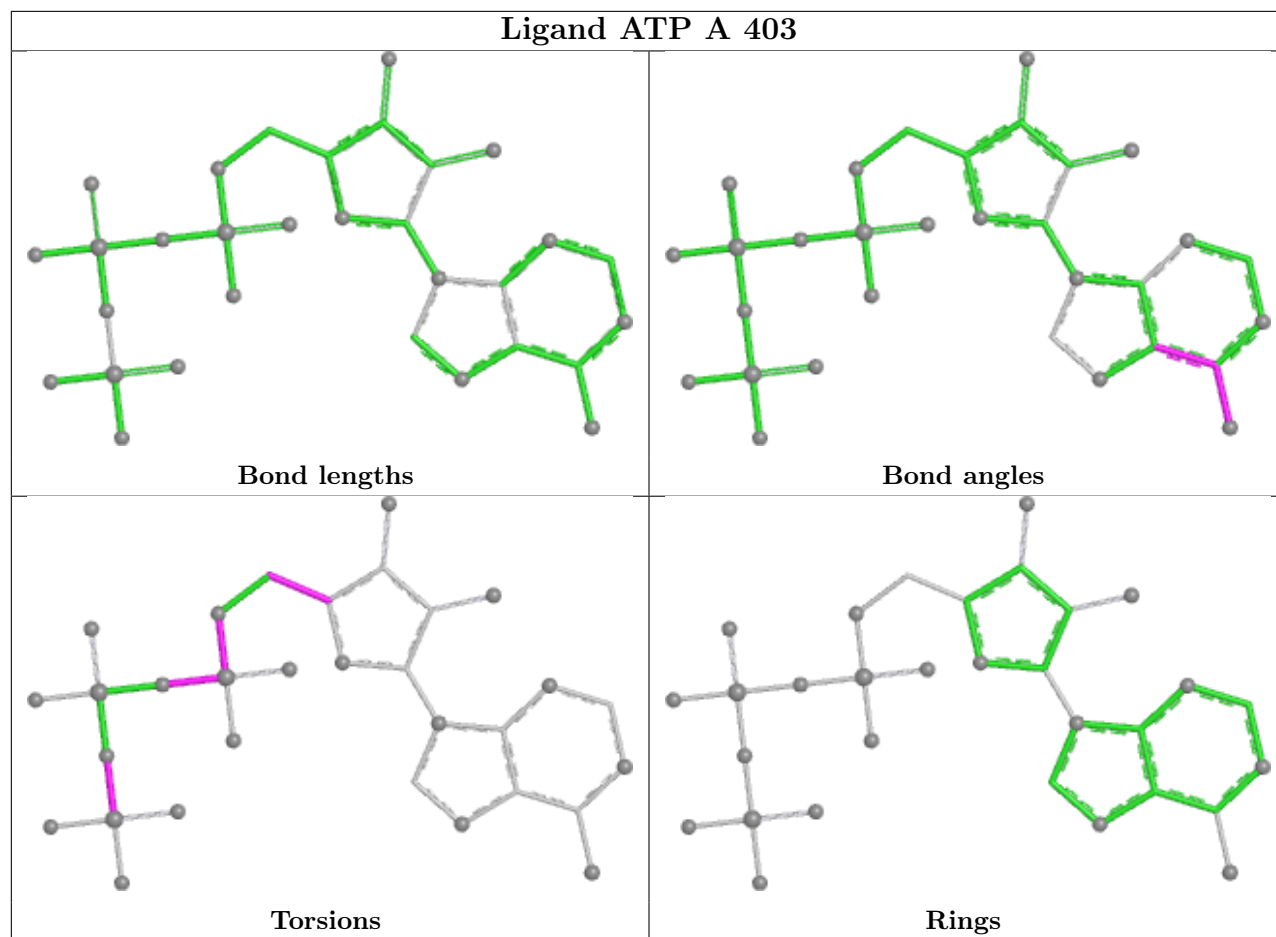
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	405[B]	ATP	1	0

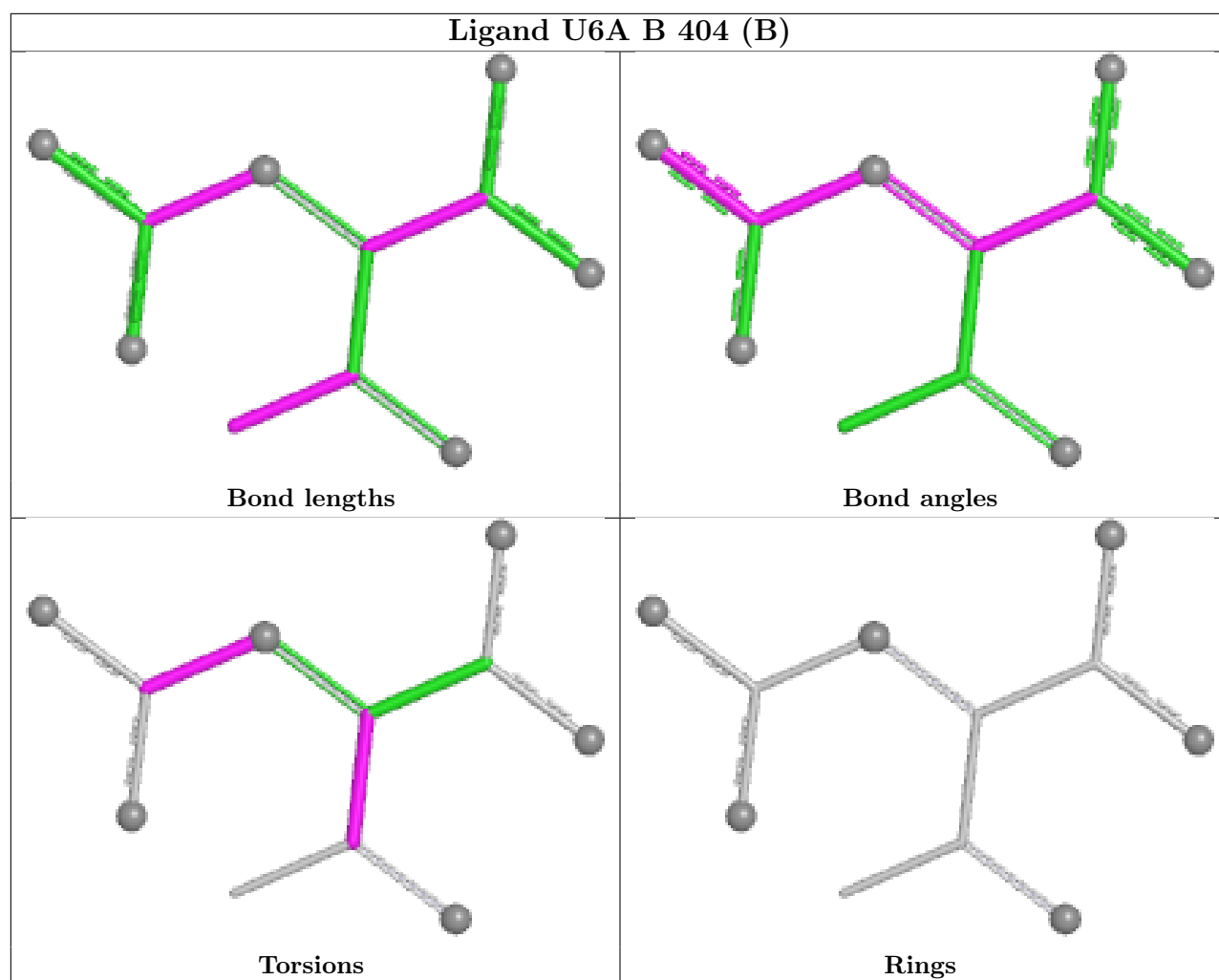
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

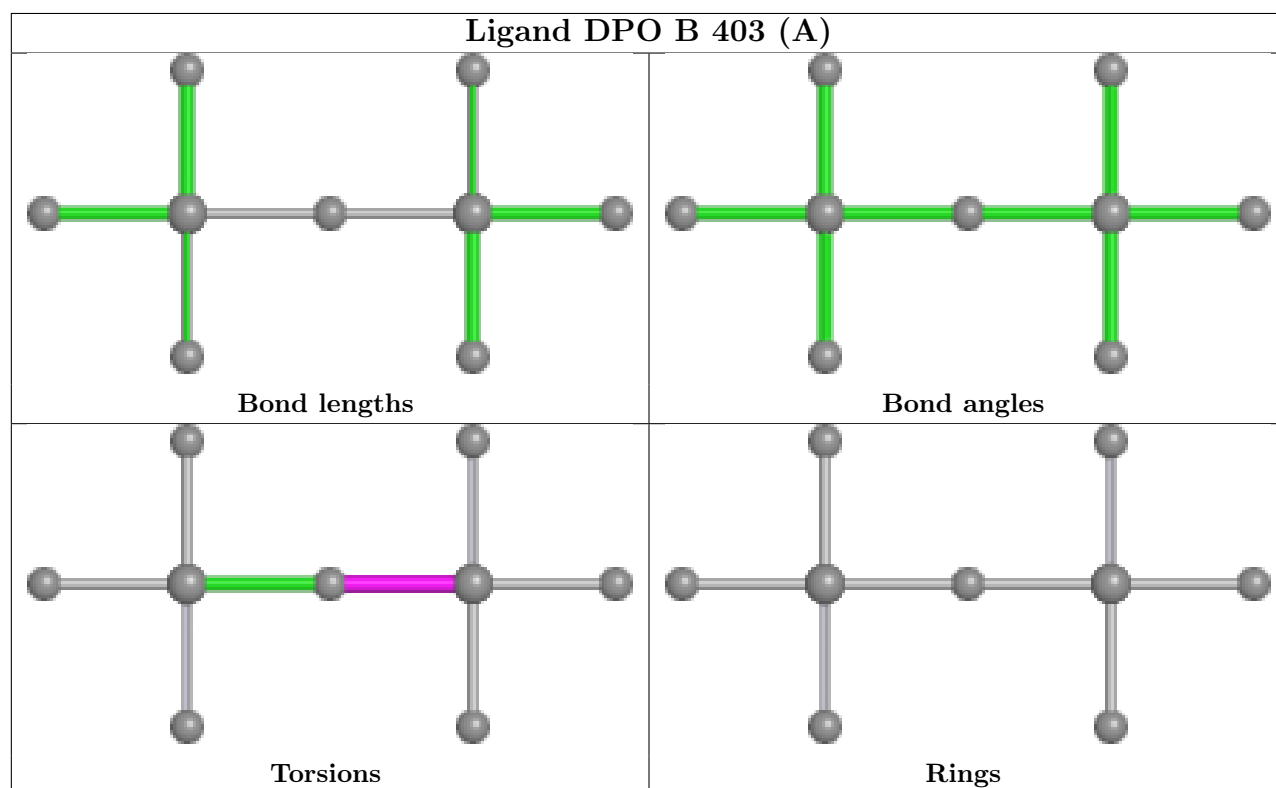
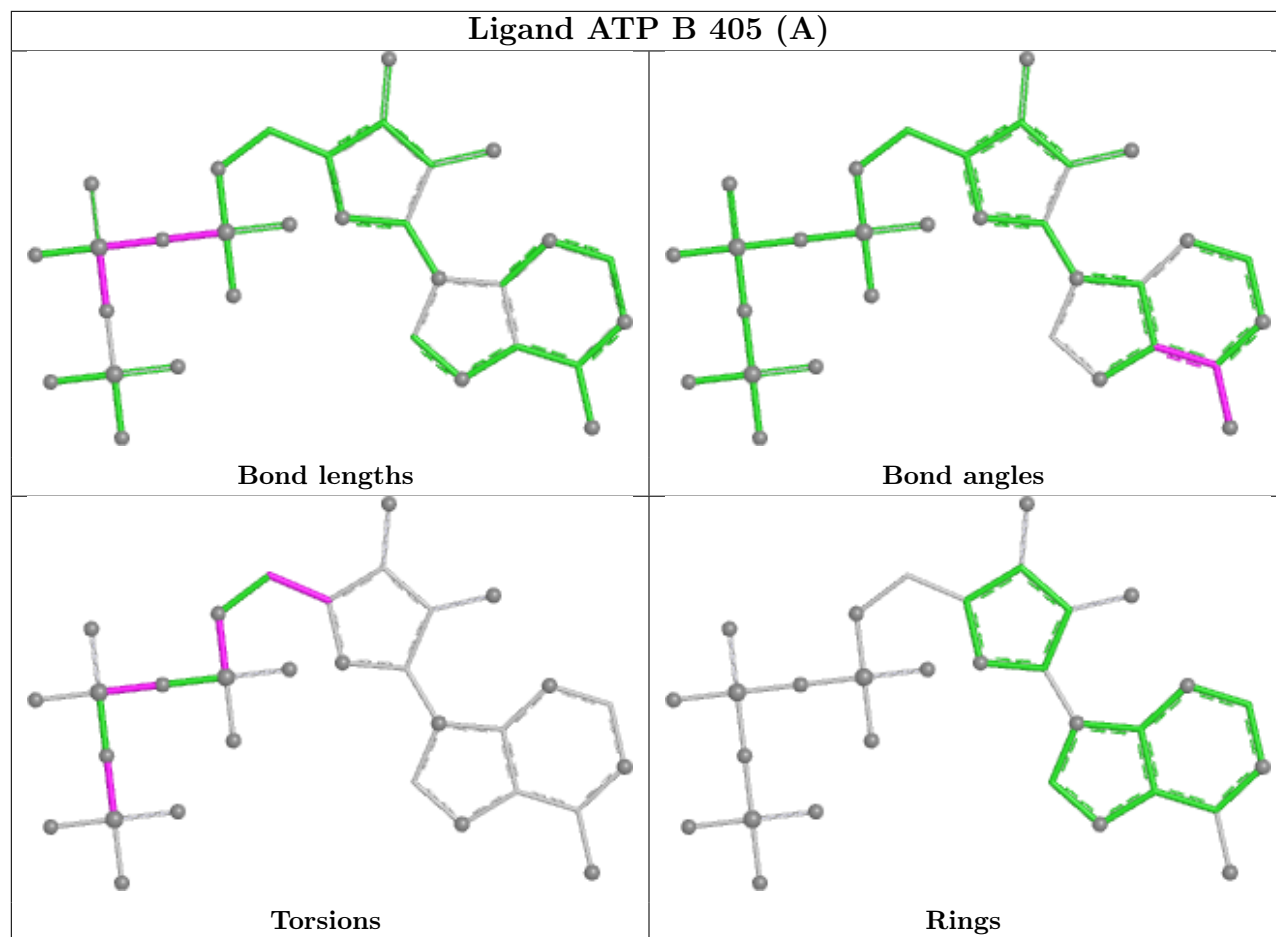




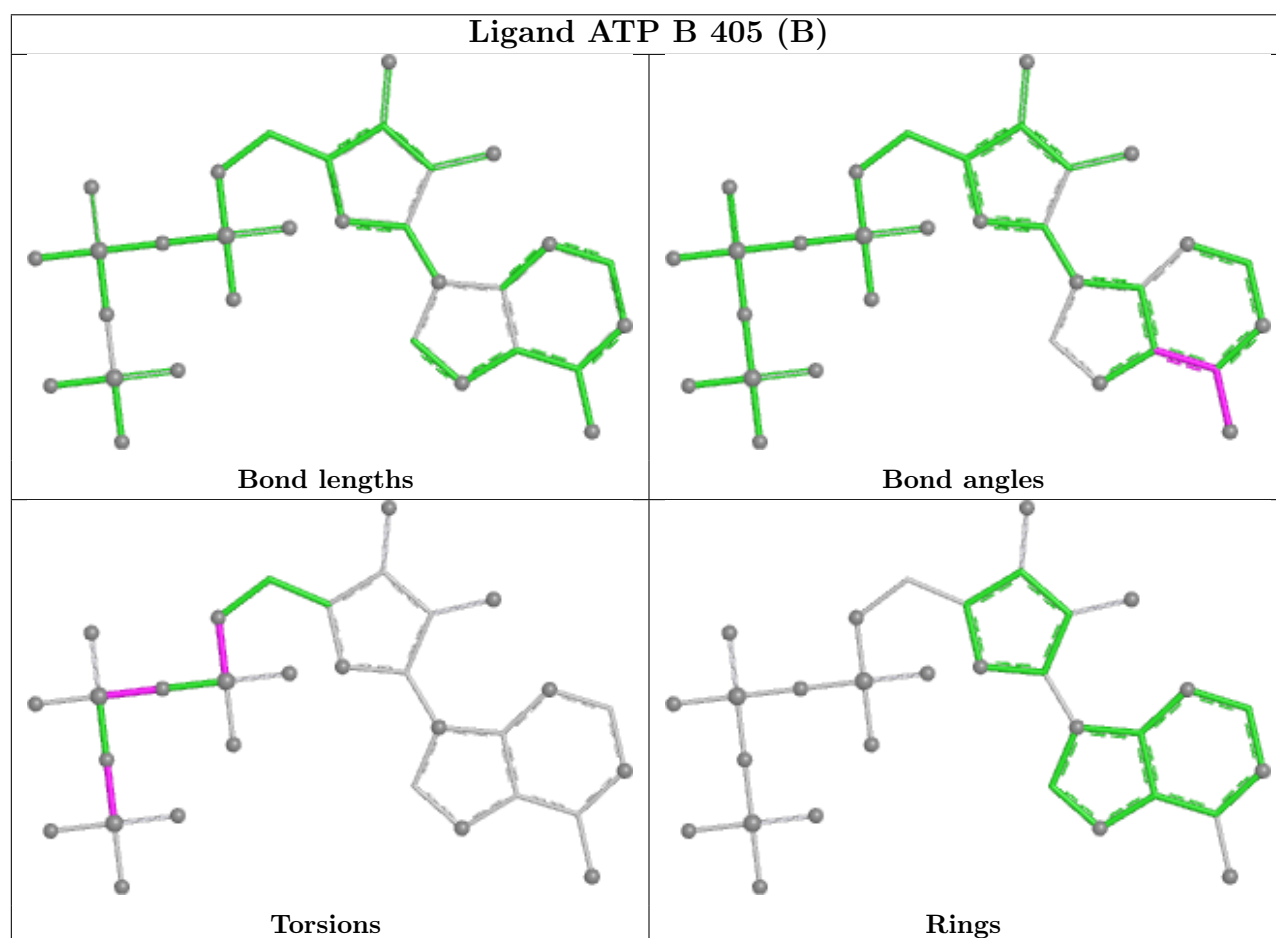
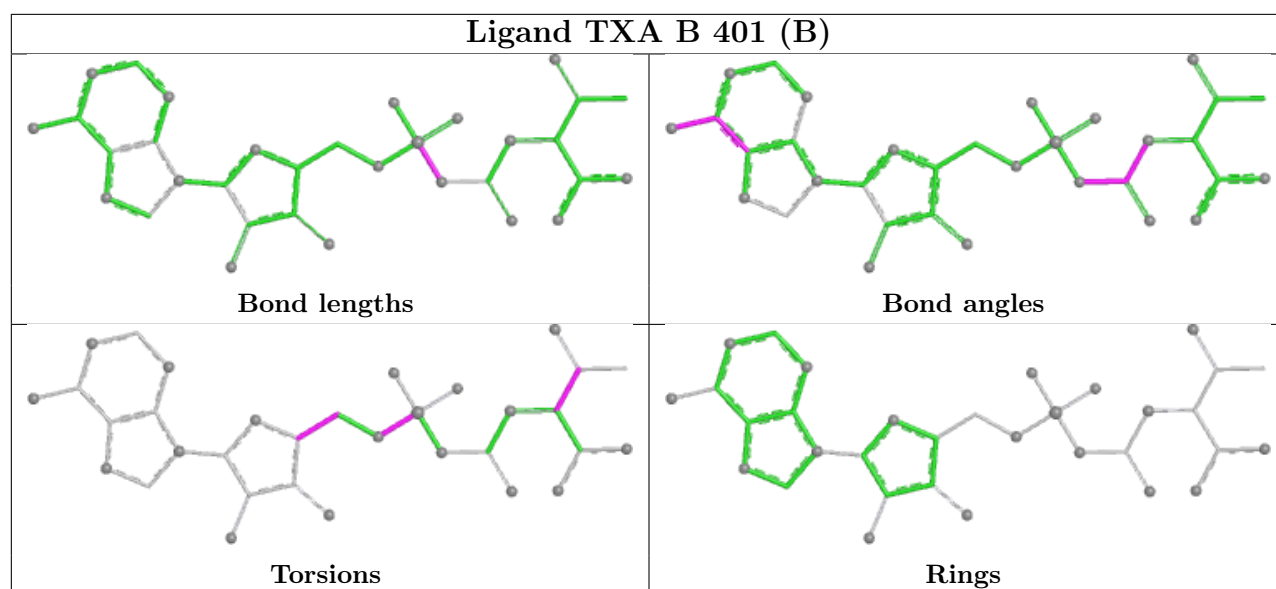












## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues ⓘ

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q < 0.9
1	A	323/340 (95%)	0.44	29 (8%)	17 15	19, 48, 89, 109	7 (2%)
1	B	323/340 (95%)	-0.10	11 (3%)	48 46	14, 36, 65, 119	11 (3%)
All	All	646/680 (95%)	0.17	40 (6%)	28 26	14, 41, 80, 119	18 (2%)

All (40) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	221	GLY	6.5
1	B	221	GLY	5.9
1	A	303[A]	MET	4.6
1	B	0	ALA	4.6
1	A	0	ALA	4.5
1	B	220	LYS	4.4
1	A	273	LEU	4.4
1	A	260	PRO	4.0
1	A	257	LYS	3.8
1	A	272	GLU	3.8
1	A	240	LEU	3.8
1	B	238	ALA	3.7
1	A	107	LYS	3.5
1	B	107	LYS	3.4
1	A	302	LYS	3.4
1	B	235	ARG	3.3
1	A	255	VAL	3.0
1	B	236	HIS	2.9
1	A	241	LYS	2.9
1	A	235	ARG	2.9
1	A	55	LYS	2.8
1	A	304	GLY	2.8
1	A	220	LYS	2.7
1	A	256	LEU	2.7

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Mol	Chain	Res	Type	RSRZ
1	B	106	GLU	2.6
1	A	335	ARG	2.6
1	A	236	HIS	2.5
1	A	258	LYS	2.4
1	A	219	ARG	2.3
1	A	307	TYR	2.3
1	B	105	SER	2.3
1	A	253	GLU	2.3
1	B	73	LYS	2.3
1	A	259	TYR	2.2
1	B	328	LYS	2.2
1	A	106	GLU	2.1
1	A	276	ASP	2.1
1	A	263	VAL	2.0
1	A	271	LYS	2.0
1	A	334	VAL	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, 95<sup>th</sup> 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.

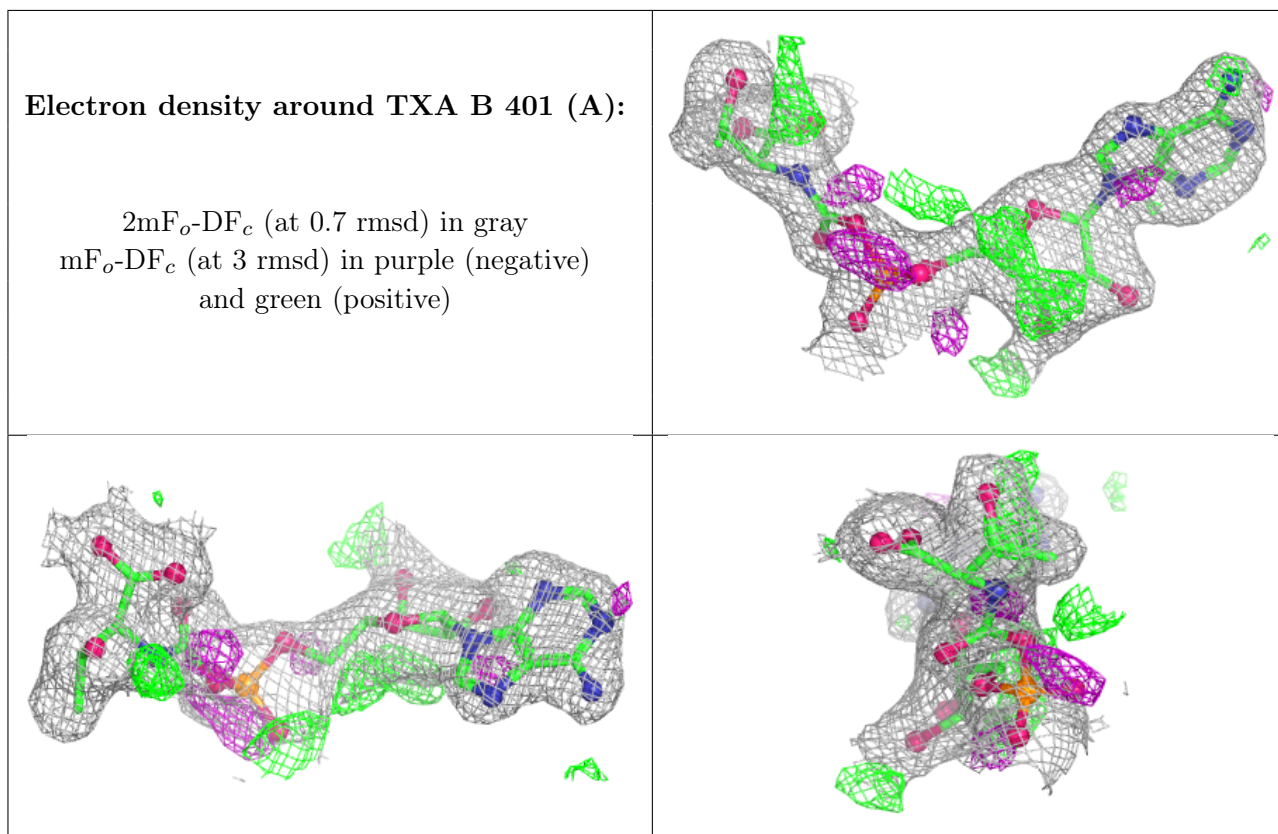
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
5	ACT	B	406	4/4	0.80	0.21	47,47,66,72	0
5	ACT	A	405	4/4	0.85	0.18	53,57,66,68	0
5	ACT	A	404	4/4	0.87	0.15	57,62,64,67	0
6	TXA	B	401[A]	33/33	0.90	0.13	22,47,74,105	33
6	TXA	B	401[B]	33/33	0.90	0.13	26,56,86,103	33
5	ACT	B	407	4/4	0.92	0.16	50,54,61,83	0
4	ATP	B	405[A]	31/31	-	-	32,49,66,97	31

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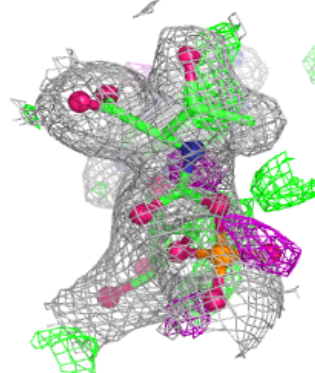
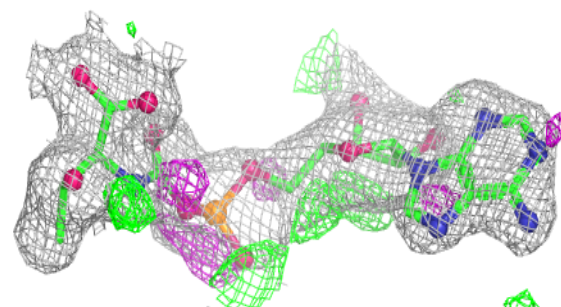
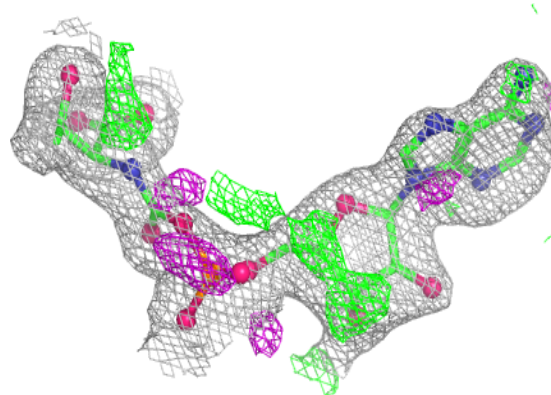
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
4	ATP	B	405[B]	31/31	-	-	30,50,70,84	31
4	ATP	A	403	31/31	0.95	0.11	37,51,81,96	31
2	U6A	A	401	11/11	0.96	0.08	26,32,50,72	0
7	DPO	B	403[A]	9/9	0.98	0.05	29,36,41,42	9
7	DPO	B	403[B]	9/9	0.98	0.05	34,39,49,53	9
2	U6A	B	404[B]	11/11	0.99	0.04	23,25,40,46	11
3	MG	A	402	1/1	0.99	0.05	40,40,40,40	0
3	MG	B	402	1/1	0.99	0.05	41,41,41,41	0
2	U6A	B	404[A]	11/11	0.99	0.04	28,31,46,50	11

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.

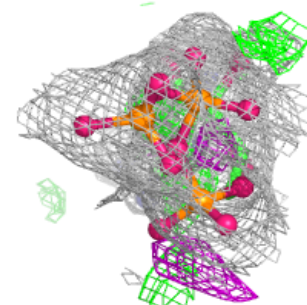
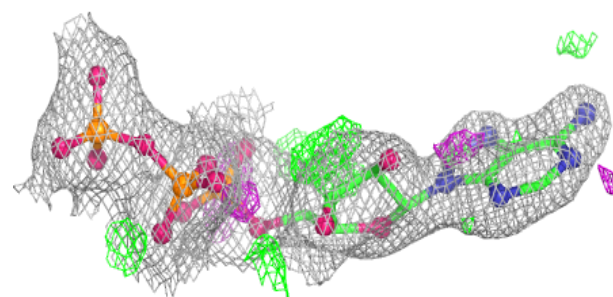
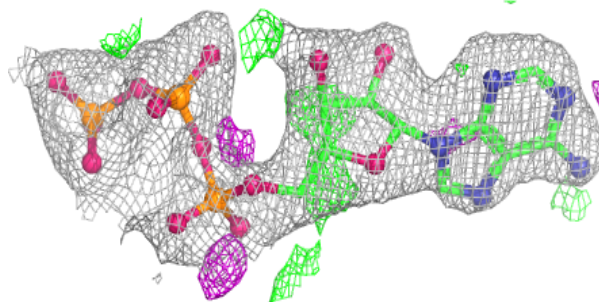


**Electron density around TXA B 401 (B):**

$2mF_o-DF_c$  (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 (A):**

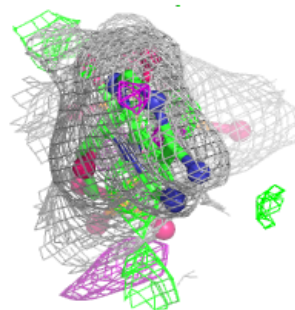
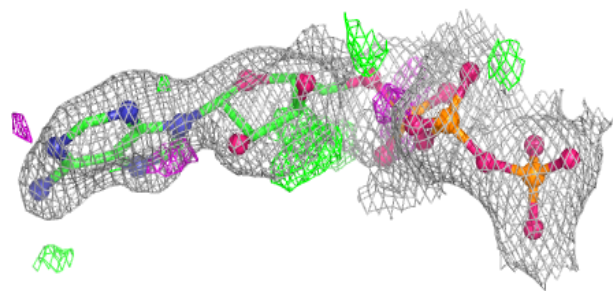
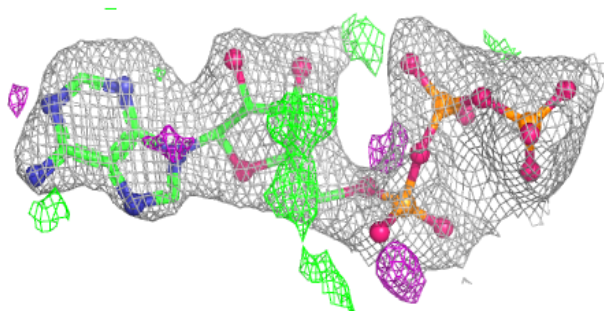
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



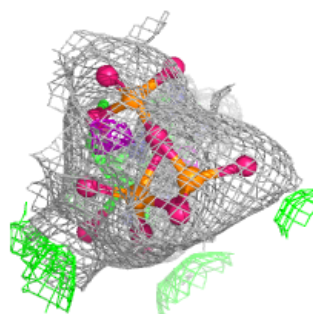
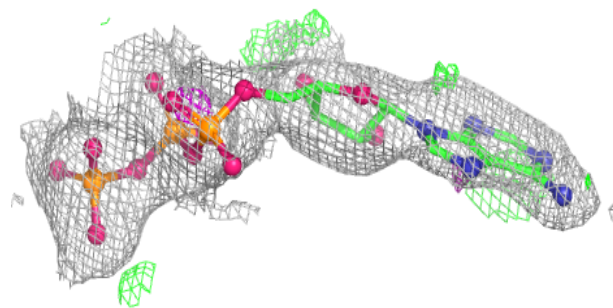
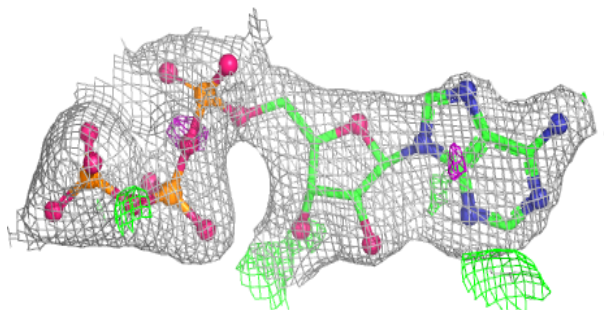


**Electron density around ATP B 405 (B):**

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

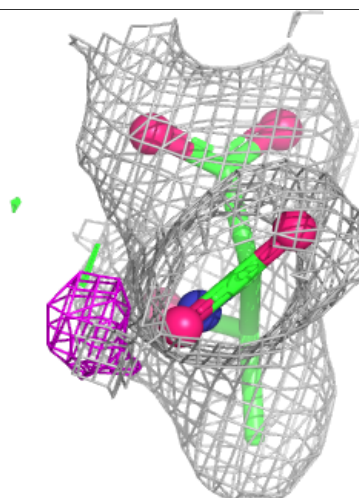
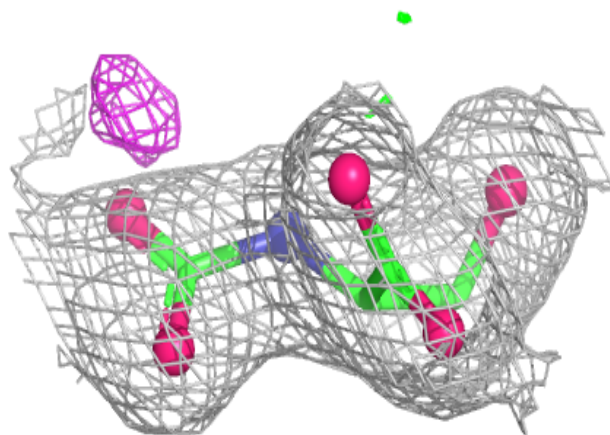
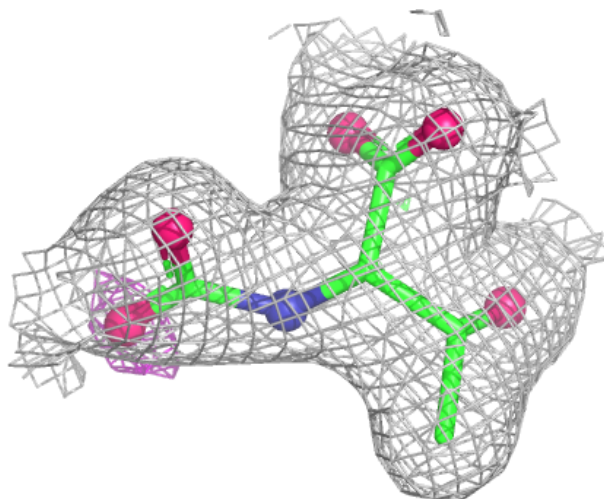
**Electron density around ATP A 403:**

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



**Electron density around U6A A 401:**

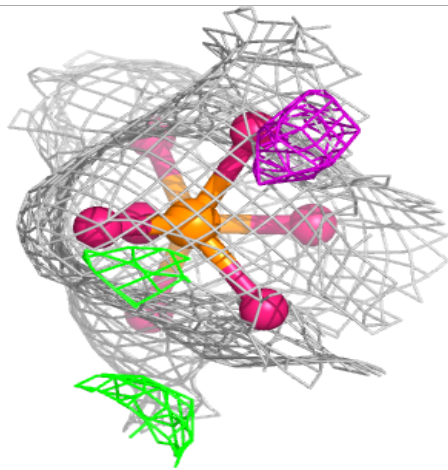
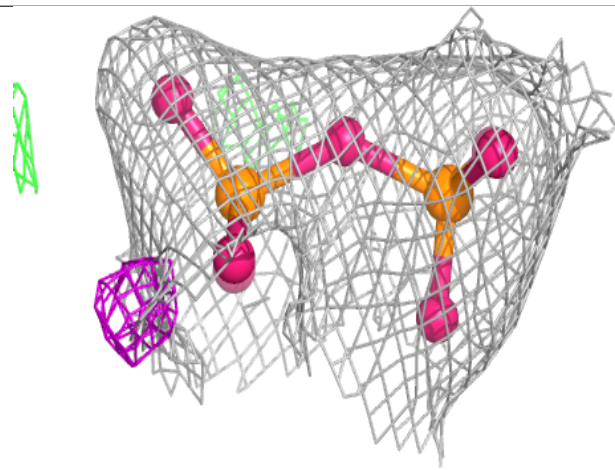
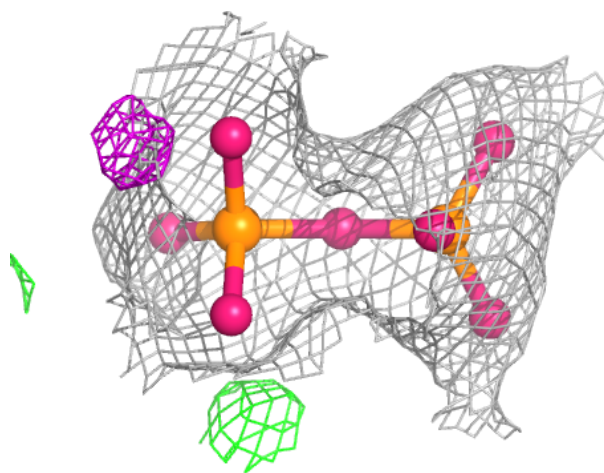
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





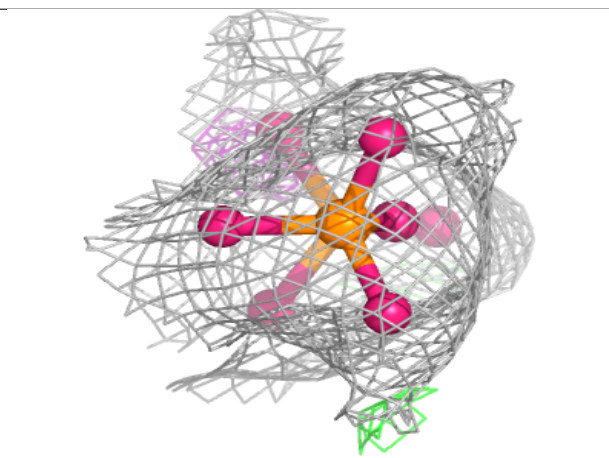
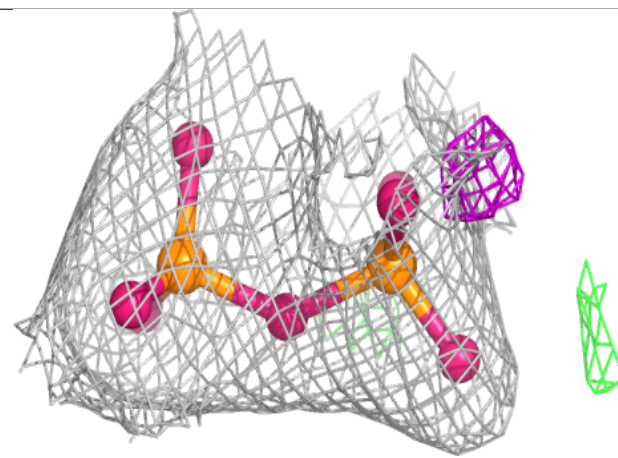
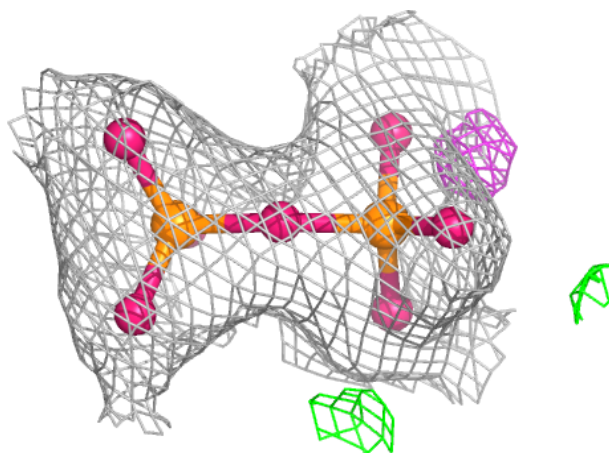
**Electron density around DPO B 403 (A):**

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



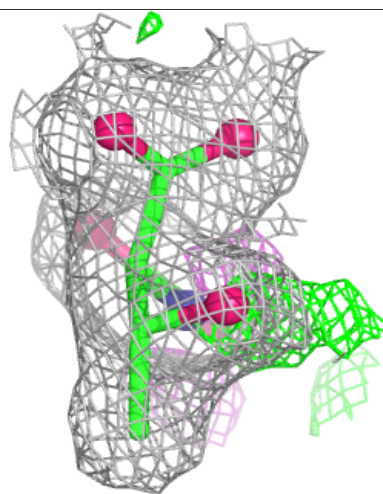
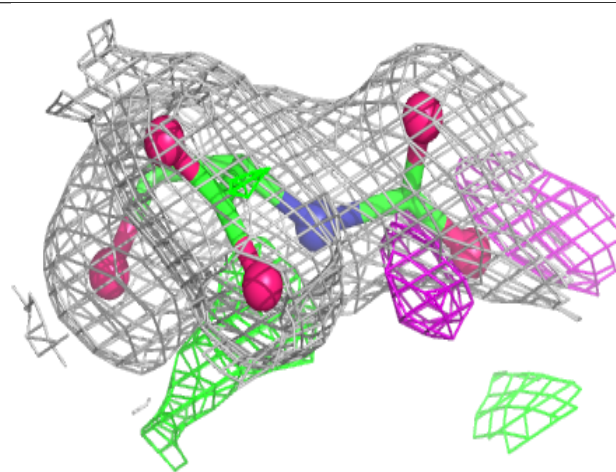
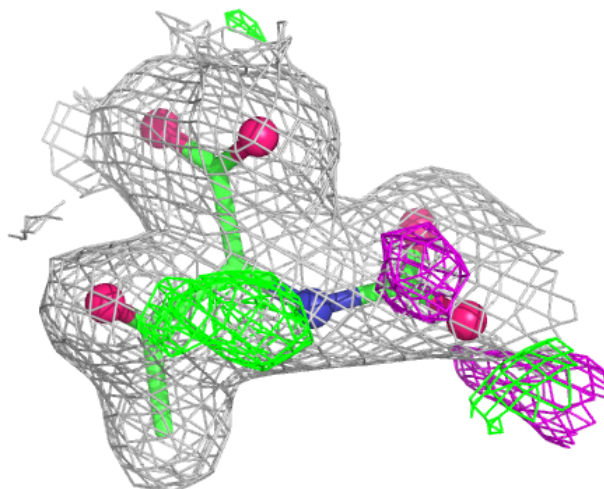
**Electron density around DPO B 403 (B):**

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



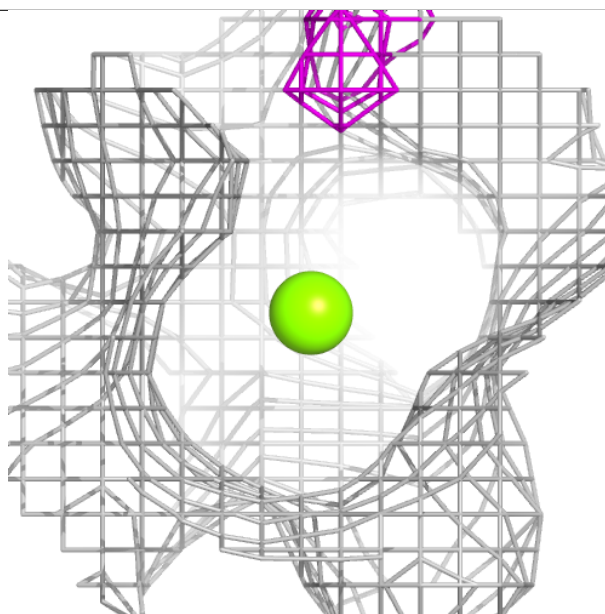
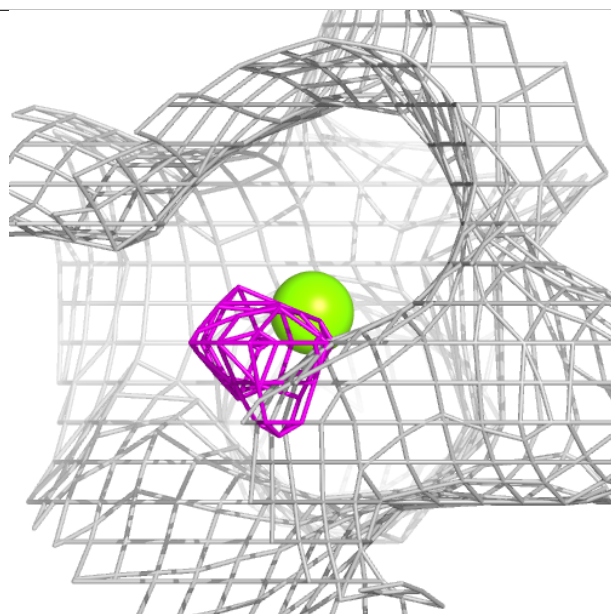
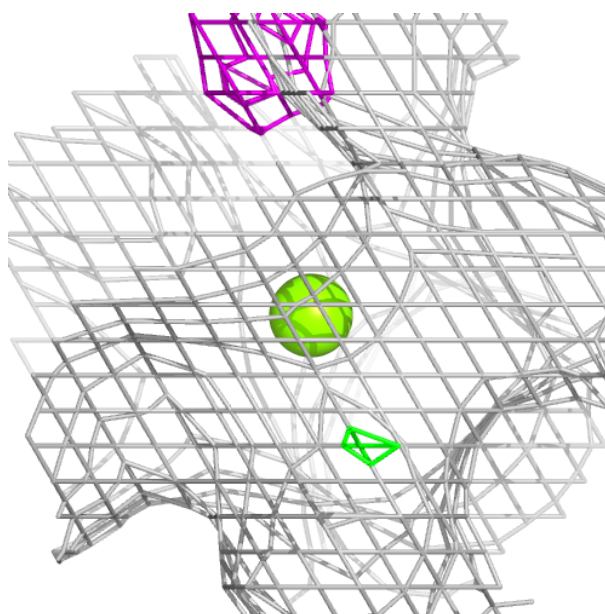
**Electron density around U6A B 404 (B):**

$2mF_o-DF_c$  (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:**

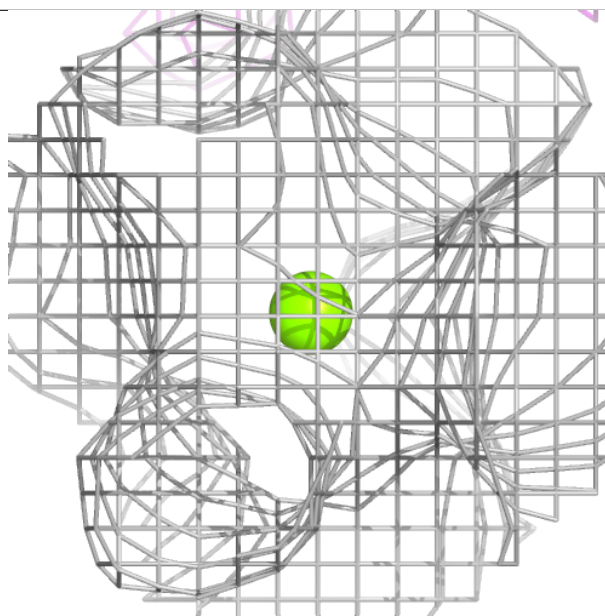
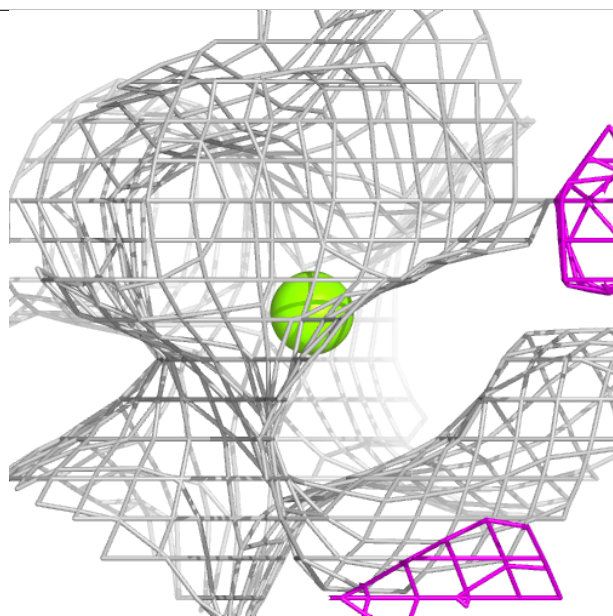
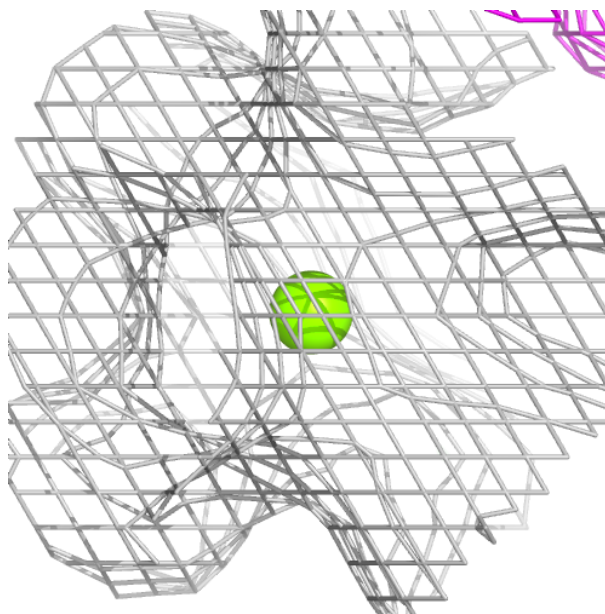
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





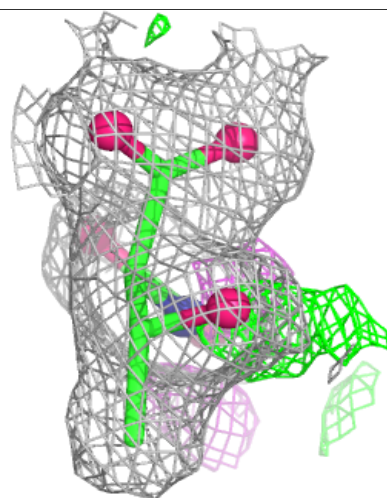
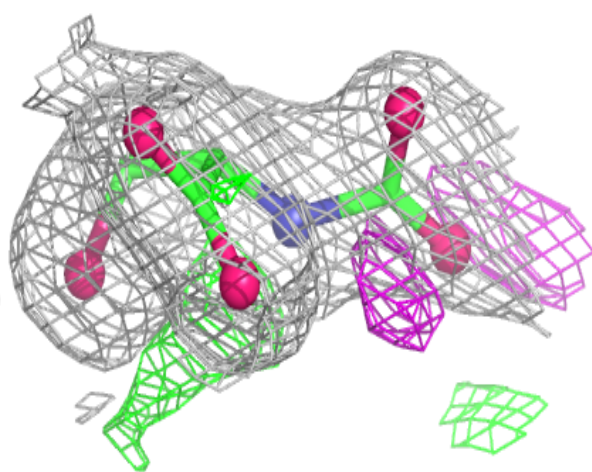
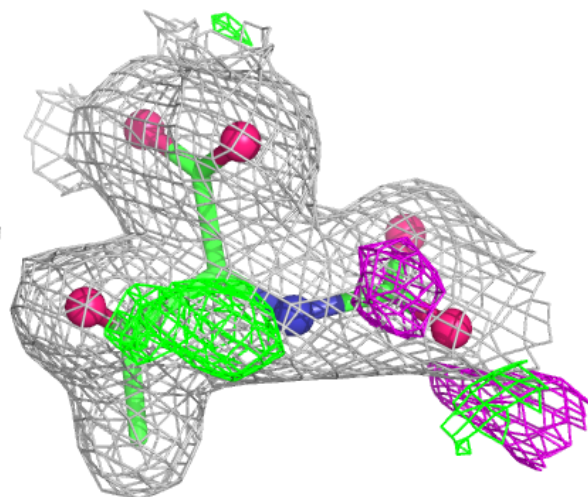
**Electron density around MG B 402:**

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



**Electron density around U6A B 404 (A):**

$2mF_o-DF_c$  (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.