



# wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 5, 2026 – 12:03 PM UTC

PDB ID : 7DAH / pdb\_00007dah  
Title : Adenosine triphosphate phosphoribosyltransferase from *Vibrio cholerae* in complex with ATP and PRPP  
Authors : Pal, R.K.; Gourinath, S.; Biswal, B.K.  
Deposited on : 2020-10-16  
Resolution : 2.92 Å (reported)

This is a wwPDB X-ray Structure Validation Summary 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>.

---

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)  
Xtrriage (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

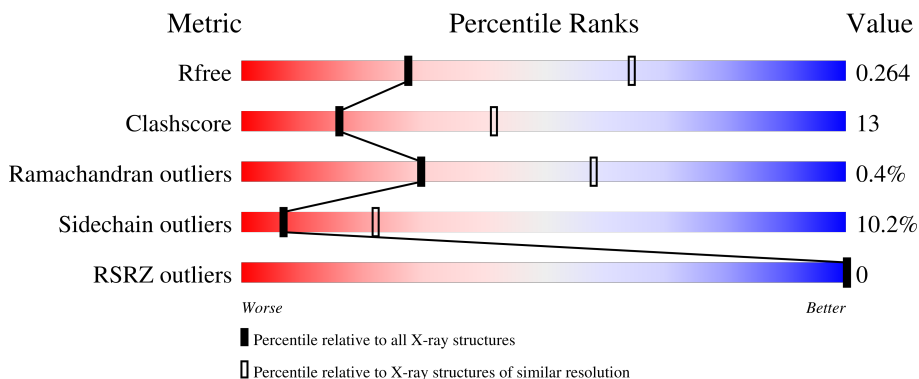
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.92 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	2995 (2.94-2.90)
Clashscore	190562	3213 (2.94-2.90)
Ramachandran outliers	187476	3128 (2.94-2.90)
Sidechain outliers	187428	3130 (2.94-2.90)
RSRZ outliers	180081	2995 (2.94-2.90)

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	310	 68% 21% • 7%
1	B	310	 66% 25% • 6%
1	C	310	 68% 21% • 7%
1	D	310	 70% 19% • 7%
1	E	310	 72% 20% • 5%

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Length	Quality of chain
1	F	310	 69% 19% 7%

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
2	PEG	D	401	-	-	X	-

## 2 Entry composition [i](#)

There are 9 unique types of molecules in this entry. The entry contains 13159 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 ATP phosphoribosyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	288	Total 2149	C 1362	N 372	O 399	S 16	0	0	0
1	B	292	Total 2131	C 1361	N 358	O 398	S 14	0	0	0
1	C	288	Total 2134	C 1350	N 370	O 398	S 16	0	0	0
1	D	287	Total 2142	C 1353	N 375	O 398	S 16	0	0	0
1	E	293	Total 2123	C 1350	N 365	O 394	S 14	0	0	0
1	F	287	Total 2105	C 1335	N 363	O 393	S 14	0	0	0

There are 36 discrepancies between the modelled and reference sequences:

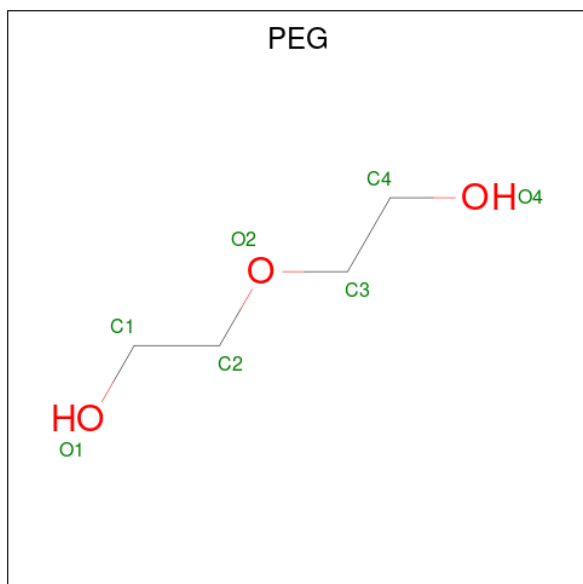
Chain	Residue	Modelled	Actual	Comment	Reference
A	-5	HIS	-	expression tag	UNP C3LU29
A	-4	HIS	-	expression tag	UNP C3LU29
A	-3	HIS	-	expression tag	UNP C3LU29
A	-2	HIS	-	expression tag	UNP C3LU29
A	-1	HIS	-	expression tag	UNP C3LU29
A	0	HIS	-	expression tag	UNP C3LU29
B	-5	HIS	-	expression tag	UNP C3LU29
B	-4	HIS	-	expression tag	UNP C3LU29
B	-3	HIS	-	expression tag	UNP C3LU29
B	-2	HIS	-	expression tag	UNP C3LU29
B	-1	HIS	-	expression tag	UNP C3LU29
B	0	HIS	-	expression tag	UNP C3LU29
C	-5	HIS	-	expression tag	UNP C3LU29
C	-4	HIS	-	expression tag	UNP C3LU29
C	-3	HIS	-	expression tag	UNP C3LU29
C	-2	HIS	-	expression tag	UNP C3LU29
C	-1	HIS	-	expression tag	UNP C3LU29

*Continued on next page...*

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
C	0	HIS	-	expression tag	UNP C3LU29
D	-5	HIS	-	expression tag	UNP C3LU29
D	-4	HIS	-	expression tag	UNP C3LU29
D	-3	HIS	-	expression tag	UNP C3LU29
D	-2	HIS	-	expression tag	UNP C3LU29
D	-1	HIS	-	expression tag	UNP C3LU29
D	0	HIS	-	expression tag	UNP C3LU29
E	-5	HIS	-	expression tag	UNP C3LU29
E	-4	HIS	-	expression tag	UNP C3LU29
E	-3	HIS	-	expression tag	UNP C3LU29
E	-2	HIS	-	expression tag	UNP C3LU29
E	-1	HIS	-	expression tag	UNP C3LU29
E	0	HIS	-	expression tag	UNP C3LU29
F	-5	HIS	-	expression tag	UNP C3LU29
F	-4	HIS	-	expression tag	UNP C3LU29
F	-3	HIS	-	expression tag	UNP C3LU29
F	-2	HIS	-	expression tag	UNP C3LU29
F	-1	HIS	-	expression tag	UNP C3LU29
F	0	HIS	-	expression tag	UNP C3LU29

- Molecule 2 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula:  $C_4H_{10}O_3$ ) (labeled as "Ligand of Interest" by depositor).



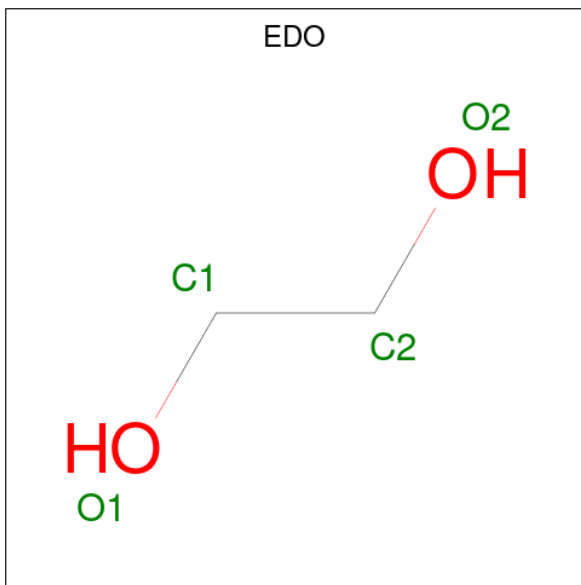
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	C O	0	0
			7	4 3		

Continued on next page...

Continued from previous page...

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	D	1	Total	C	O	0	0
			7	4	3		

- Molecule 3 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C<sub>2</sub>H<sub>6</sub>O<sub>2</sub>) (labeled as "Ligand of Interest" by depositor).



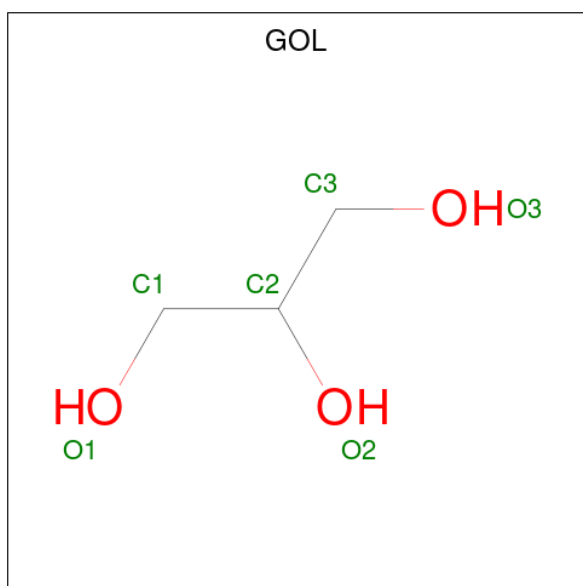
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		
3	C	1	Total	C	O	0	0
			4	2	2		
3	D	1	Total	C	O	0	0
			4	2	2		
3	D	1	Total	C	O	0	0
			4	2	2		
3	D	1	Total	C	O	0	0
			4	2	2		
3	E	1	Total	C	O	0	0
			4	2	2		
3	F	1	Total	C	O	0	0
			4	2	2		

- Molecule 4 is PHOSPHATE ION (CCD ID: PO4) (formula: O<sub>4</sub>P) (labeled as "Ligand of Interest" by depositor).



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

- Molecule 5 is GLYCEROL (CCD ID: GOL) (formula:  $C_3H_8O_3$ ) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total C O 6 3 3	0	0

*Continued on next page...*

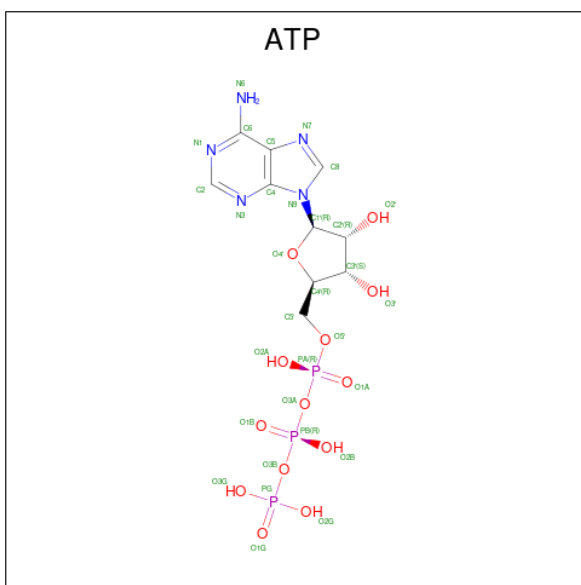
Continued from previous page...

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	D	1	Total	C	O	0	0
			6	3	3		

- Molecule 6 is MAGNESIUM ION (CCD ID: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

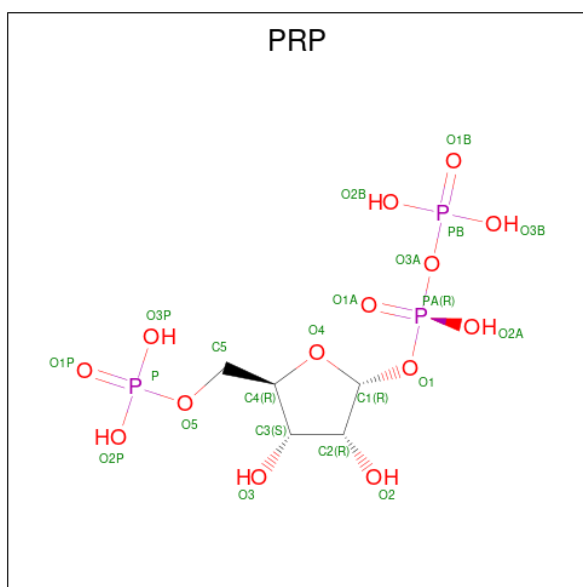
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	C	1	Total	Mg	0	0
			1	1		
6	F	1	Total	Mg	0	0
			1	1		

- Molecule 7 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>) (labeled as "Ligand of Interest" by depositor).



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

- Molecule 8 is 1-O-pyrophosphono-5-O-phosphono-alpha-D-ribofuranose (CCD ID: PRP) (formula: C<sub>5</sub>H<sub>13</sub>O<sub>14</sub>P<sub>3</sub>) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
8	C	1	Total	C	O	P	0	0
			22	5	14	3		
8	F	1	Total	C	O	P	0	0
			22	5	14	3		

- Molecule 9 is water.

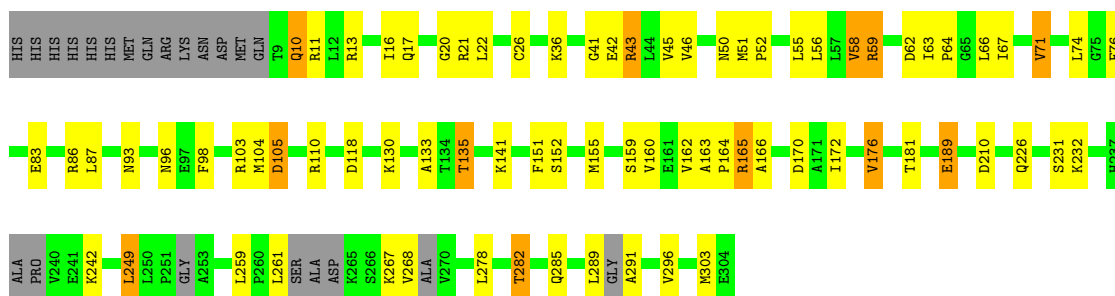
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	35	Total	O	0	0
			35	35		
9	B	27	Total	O	0	0
			27	27		
9	C	47	Total	O	0	0
			47	47		
9	D	42	Total	O	0	0
			42	42		
9	E	22	Total	O	0	0
			22	22		
9	F	26	Total	O	0	0
			26	26		

### 3 Residue-property plots

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.

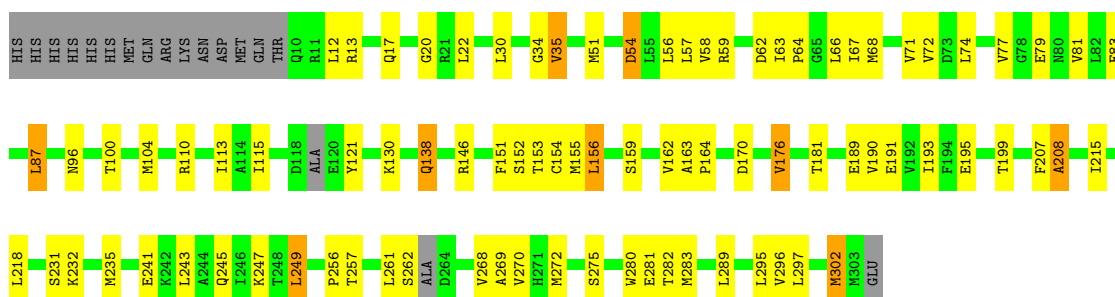
- Molecule 1: ATP phosphoribosyltransferase

Chain A: 



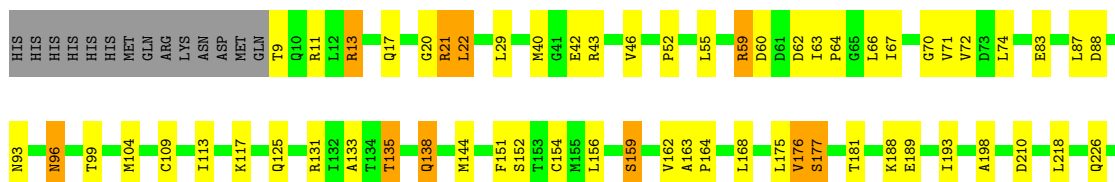
- Molecule 1: ATP phosphoribosyltransferase

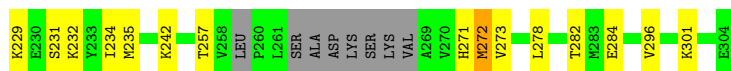
Chain B: 



- Molecule 1: ATP phosphoribosyltransferase

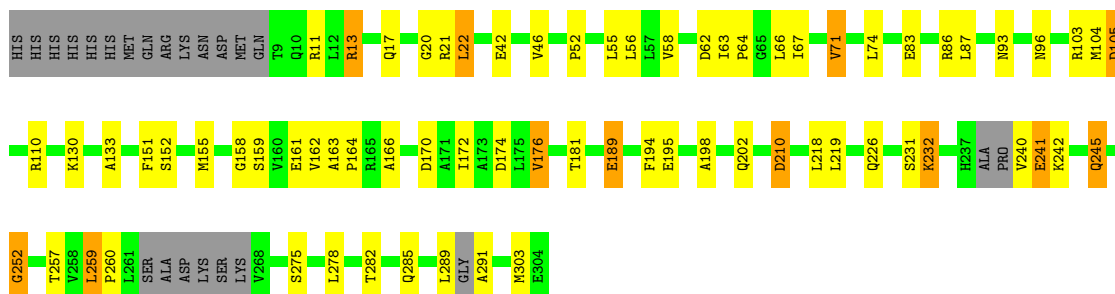
Chain C: 





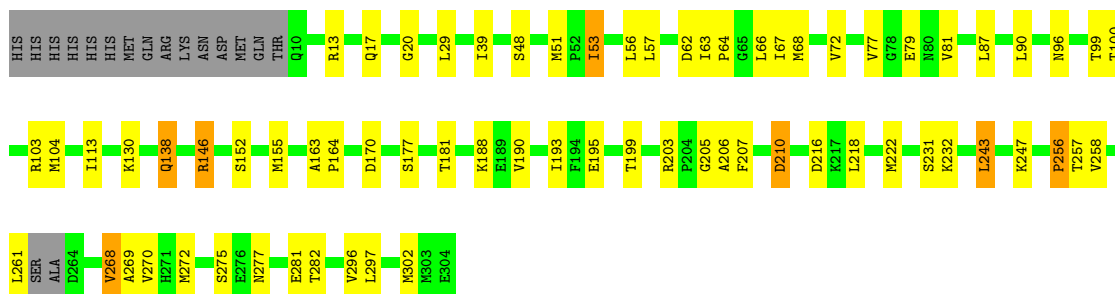
- Molecule 1: ATP phosphoribosyltransferase

Chain D: 70% 19% 7%



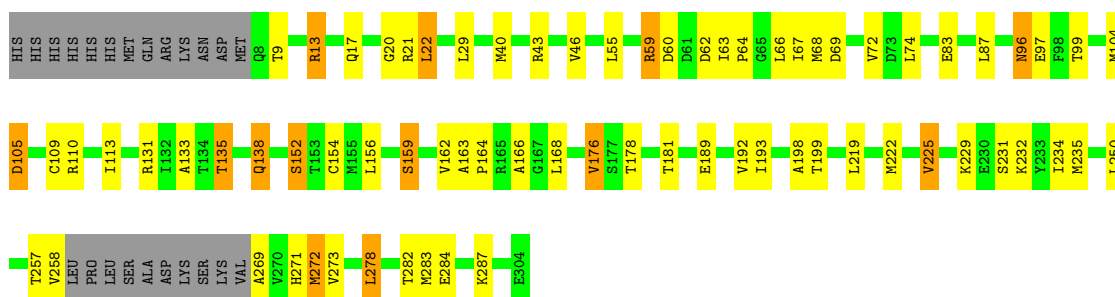
- Molecule 1: ATP phosphoribosyltransferase

Chain E: 72% 20% 5%



- Molecule 1: ATP phosphoribosyltransferase

Chain F: 69% 19% 7%



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 32	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	136.60Å 136.60Å 121.93Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	28.32 – 2.92 28.32 – 2.92	Depositor EDS
% Data completeness (in resolution range)	99.6 (28.32-2.92) 99.8 (28.32-2.92)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.03 (at 2.90Å)	Xtrriage
Refinement program	REFMAC 5.8.0257	Depositor
R, $R_{free}$	0.216 , 0.268 0.215 , 0.264	Depositor DCC
$R_{free}$ test set	2553 reflections (4.62%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	61.3	Xtrriage
Anisotropy	0.153	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.27 , 31.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.52$ , $\langle L^2 \rangle = 0.35$	Xtrriage
Estimated twinning fraction	0.009 for -h,-k,l 0.470 for h,-h-k,-l 0.010 for -k,-h,-l	Xtrriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	13159	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	72.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.78% 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 [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, EDO, ATP, PO4, PRP, PEG, MG

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 5$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z  >5	RMSZ	# Z  >5
1	A	1.14	1/2170 (0.0%)	1.55	6/2924 (0.2%)
1	B	1.16	0/2155	1.57	5/2916 (0.2%)
1	C	1.14	0/2157	1.54	2/2911 (0.1%)
1	D	1.17	2/2164 (0.1%)	1.54	9/2917 (0.3%)
1	E	1.18	3/2150 (0.1%)	1.57	6/2912 (0.2%)
1	F	1.15	0/2129	1.56	3/2876 (0.1%)
All	All	1.16	6/12925 (0.0%)	1.56	31/17456 (0.2%)

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	C	0	1
1	D	0	1
All	All	0	2

The worst 5 of 6 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	189	GLU	C-O	-6.65	1.16	1.23
1	A	189	GLU	C-O	-6.44	1.16	1.23
1	E	205	GLY	C-O	6.01	1.28	1.23
1	E	302	MET	C-O	5.55	1.29	1.23
1	D	158	GLY	C-O	5.42	1.30	1.24

The worst 5 of 31 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	105	ASP	CA-CB-CG	6.83	119.43	112.60
1	A	10	GLN	N-CA-C	-6.67	106.08	114.56
1	D	105	ASP	CA-CB-CG	5.99	118.59	112.60
1	E	190	VAL	CA-C-O	-5.79	114.95	119.46
1	A	105	ASP	CA-CB-CG	5.77	118.37	112.60

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	257	THR	Peptide
1	D	259	LEU	Peptide

## 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	2149	0	2130	57	0
1	B	2131	0	2087	77	0
1	C	2134	0	2101	67	0
1	D	2142	0	2124	48	0
1	E	2123	0	2056	49	0
1	F	2105	0	2057	62	0
2	A	7	0	10	0	0
2	D	7	0	10	4	0
3	A	8	0	12	0	0
3	C	4	0	6	0	0
3	D	12	0	18	1	0
3	E	4	0	6	0	0
3	F	4	0	6	0	0
4	A	5	0	0	0	0
4	D	5	0	0	0	0
5	A	6	0	8	0	0
5	D	6	0	8	1	0
6	C	1	0	0	0	0
6	F	1	0	0	0	0
7	C	31	0	12	2	0
7	F	31	0	12	2	0

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
8	C	22	0	8	2	0
8	F	22	0	8	2	0
9	A	35	0	0	2	0
9	B	27	0	0	4	0
9	C	47	0	0	6	0
9	D	42	0	0	1	0
9	E	22	0	0	3	0
9	F	26	0	0	0	0
All	All	13159	0	12679	336	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 13.

The worst 5 of 336 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:D:17:GLN:NE2	1:D:20:GLY:HA3	1.77	0.99
1:A:17:GLN:NE2	1:A:20:GLY:HA3	1.75	0.99
1:B:17:GLN:NE2	1:B:20:GLY:HA3	1.78	0.99
1:E:17:GLN:NE2	1:E:20:GLY:HA3	1.79	0.97
1:C:17:GLN:NE2	1:C:20:GLY:HA3	1.78	0.97

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	276/310 (89%)	260 (94%)	14 (5%)	2 (1%)	18	46
1	B	286/310 (92%)	271 (95%)	14 (5%)	1 (0%)	36	64
1	C	282/310 (91%)	269 (95%)	12 (4%)	1 (0%)	30	58

*Continued on next page...*

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	D	279/310 (90%)	262 (94%)	15 (5%)	2 (1%)	18	46
1	E	289/310 (93%)	275 (95%)	14 (5%)	0	100	100
1	F	283/310 (91%)	269 (95%)	14 (5%)	0	100	100
All	All	1695/1860 (91%)	1606 (95%)	83 (5%)	6 (0%)	30	58

5 of 6 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	242	LYS
1	B	208	ALA
1	C	242	LYS
1	D	260	PRO
1	D	259	LEU

### 5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	213/263 (81%)	193 (91%)	20 (9%)	8	25
1	B	201/263 (76%)	178 (89%)	23 (11%)	5	17
1	C	207/263 (79%)	184 (89%)	23 (11%)	6	19
1	D	212/263 (81%)	195 (92%)	17 (8%)	11	32
1	E	197/263 (75%)	177 (90%)	20 (10%)	7	22
1	F	201/263 (76%)	178 (89%)	23 (11%)	5	17
All	All	1231/1578 (78%)	1105 (90%)	126 (10%)	7	22

5 of 126 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	177	SER
1	F	135	THR
1	D	152	SER
1	F	99	THR

Continued on next page...

*Continued from previous page...*

Mol	Chain	Res	Type
1	F	219	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 37 such sidechains are listed below:

Mol	Chain	Res	Type
1	E	138	GLN
1	F	226	GLN
1	E	237	HIS
1	F	138	GLN
1	B	202	GLN

### 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 20 ligands modelled in this entry, 2 are monoatomic - leaving 18 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	$\# Z  > 2$
3	EDO	D	405	-	3,3,3	0.41	0	2,2,2	0.16	0
4	PO4	A	404	-	4,4,4	1.53	1 (25%)	6,6,6	0.38	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	PEG	A	401	-	6,6,6	0.19	0	5,5,5	0.22	0
3	EDO	A	403	-	3,3,3	0.27	0	2,2,2	0.14	0
4	PO4	D	406	-	4,4,4	0.76	0	6,6,6	0.61	0
3	EDO	A	402	-	3,3,3	0.25	0	2,2,2	0.20	0
8	PRP	C	403	-	20,22,22	0.98	1 (5%)	32,35,35	1.33	3 (9%)
3	EDO	F	402	-	3,3,3	0.37	0	2,2,2	0.34	0
3	EDO	C	404	-	3,3,3	0.11	0	2,2,2	0.10	0
3	EDO	E	401	-	3,3,3	0.19	0	2,2,2	0.19	0
7	ATP	F	404	6	32,33,33	0.78	1 (3%)	48,52,52	0.91	2 (4%)
2	PEG	D	401	-	6,6,6	0.23	0	5,5,5	0.18	0
8	PRP	F	403	-	20,22,22	1.44	1 (5%)	32,35,35	1.23	2 (6%)
3	EDO	D	404	-	3,3,3	0.22	0	2,2,2	0.18	0
7	ATP	C	402	6	32,33,33	0.96	3 (9%)	48,52,52	0.95	2 (4%)
3	EDO	D	402	-	3,3,3	0.14	0	2,2,2	0.06	0
5	GOL	D	403	-	5,5,5	0.17	0	5,5,5	0.35	0
5	GOL	A	405	-	5,5,5	0.17	0	5,5,5	0.34	0

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
3	EDO	D	405	-	-	1/1/1/1	-
2	PEG	A	401	-	-	0/4/4/4	-
3	EDO	A	403	-	-	1/1/1/1	-
8	PRP	C	403	-	-	4/16/33/33	0/1/1/1
3	EDO	A	402	-	-	1/1/1/1	-
3	EDO	F	402	-	-	0/1/1/1	-
3	EDO	C	404	-	-	1/1/1/1	-
3	EDO	E	401	-	-	1/1/1/1	-
7	ATP	F	404	6	-	5/22/38/38	0/3/3/3
2	PEG	D	401	-	-	3/4/4/4	-
8	PRP	F	403	-	-	3/16/33/33	0/1/1/1
3	EDO	D	404	-	-	0/1/1/1	-
7	ATP	C	402	6	-	8/22/38/38	0/3/3/3
3	EDO	D	402	-	-	0/1/1/1	-
5	GOL	D	403	-	-	3/4/4/4	-
5	GOL	A	405	-	-	0/4/4/4	-

The worst 5 of 7 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	F	403	PRP	PA-O3A	5.57	1.65	1.59
8	C	403	PRP	PA-O3A	3.13	1.62	1.59
4	A	404	PO4	P-O1	3.05	1.57	1.50
7	C	402	ATP	PA-O3A	2.92	1.62	1.59
7	C	402	ATP	C1'-N9	2.72	1.53	1.46

The worst 5 of 9 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	C	403	PRP	O5-P-O1P	-4.24	94.98	106.44
8	F	403	PRP	O2P-P-O5	-3.31	98.04	106.67
8	C	403	PRP	O1-C1-C2	3.00	111.52	106.65
7	F	404	ATP	C2'-C1'-N9	2.61	119.78	113.30
7	C	402	ATP	C2'-C1'-N9	2.56	119.66	113.30

There are no chirality outliers.

5 of 31 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	D	403	GOL	C1-C2-C3-O3
5	D	403	GOL	O2-C2-C3-O3
7	C	402	ATP	PB-O3B-PG-O3G
7	C	402	ATP	C5'-O5'-PA-O2A
7	C	402	ATP	O4'-C4'-C5'-O5'

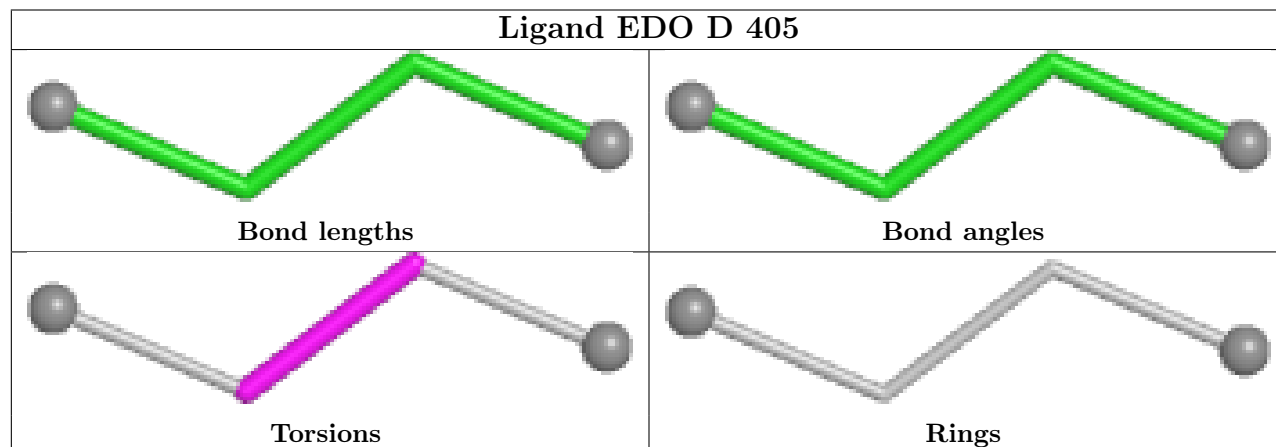
There are no ring outliers.

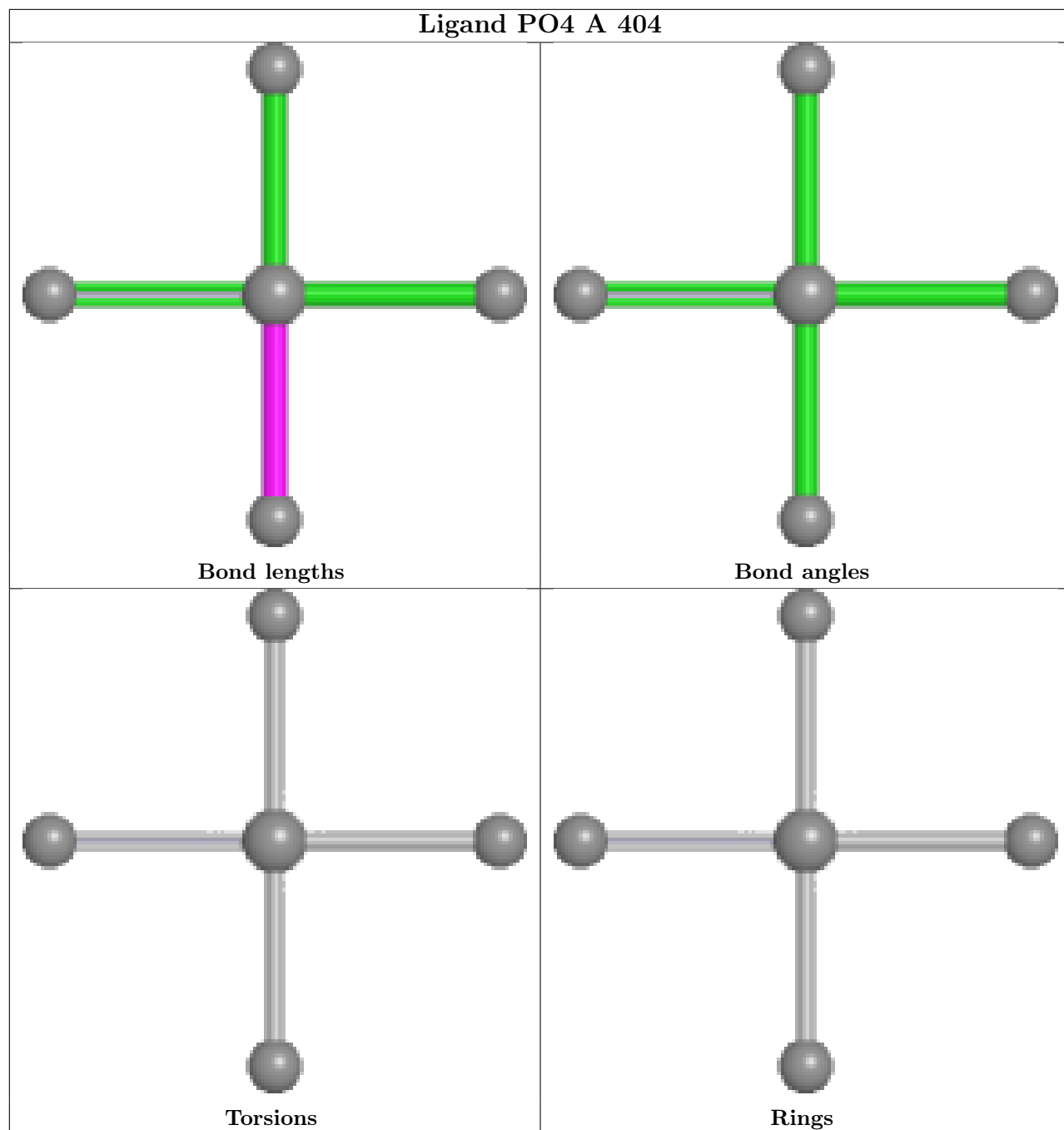
7 monomers are involved in 14 short contacts:

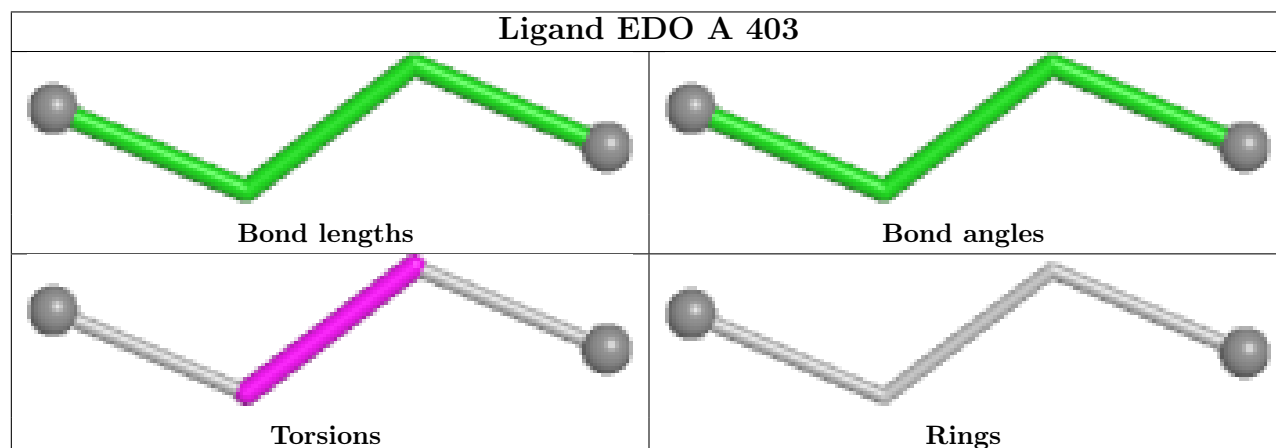
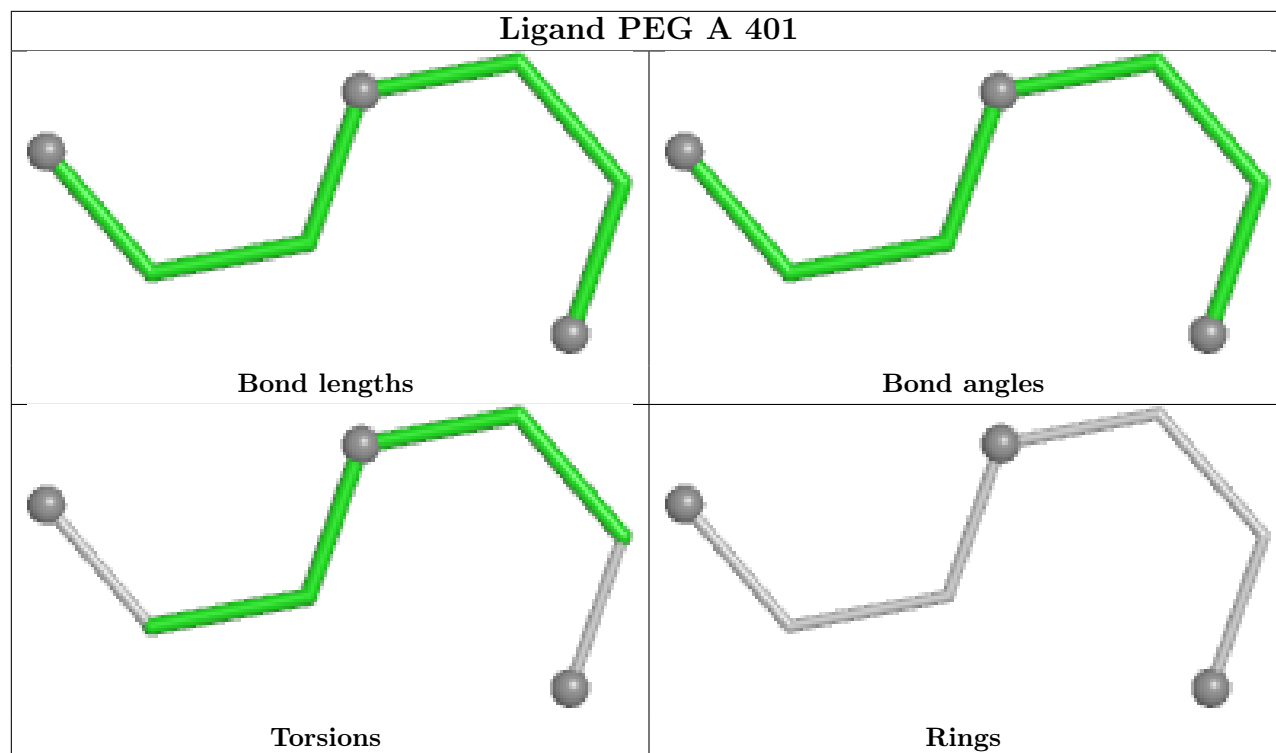
Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	C	403	PRP	2	0
7	F	404	ATP	2	0
2	D	401	PEG	4	0
8	F	403	PRP	2	0
7	C	402	ATP	2	0
3	D	402	EDO	1	0
5	D	403	GOL	1	0

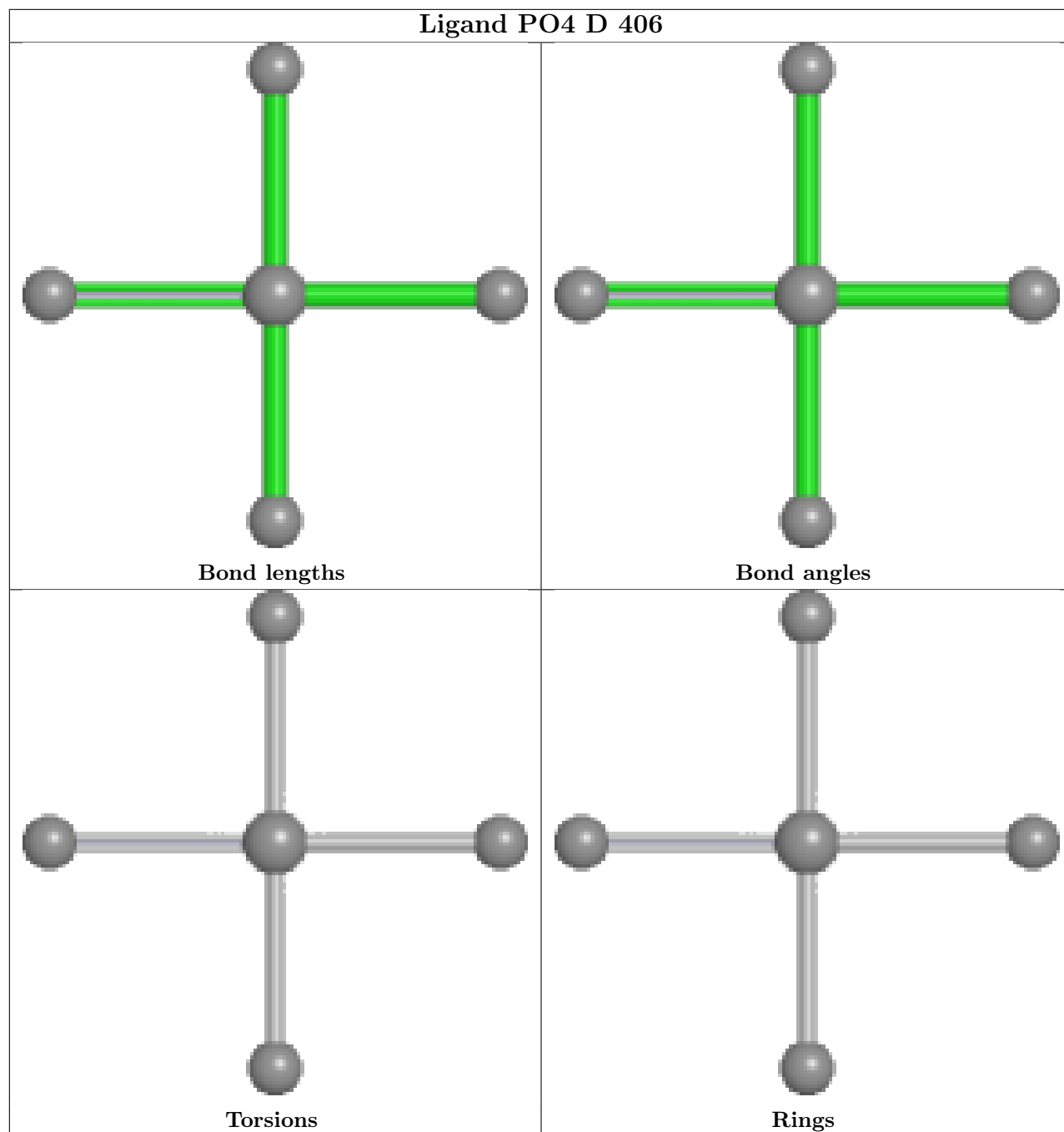
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is

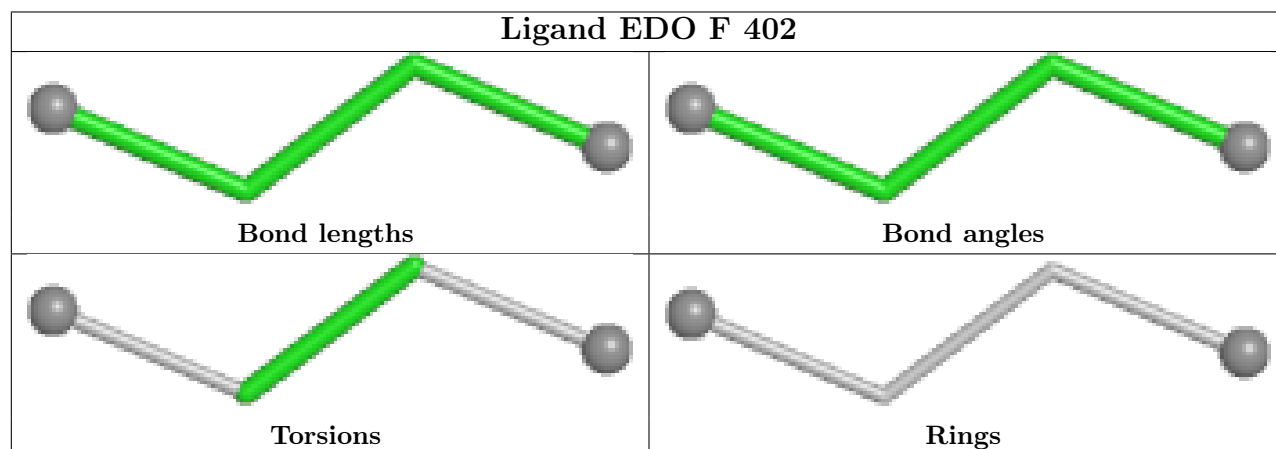
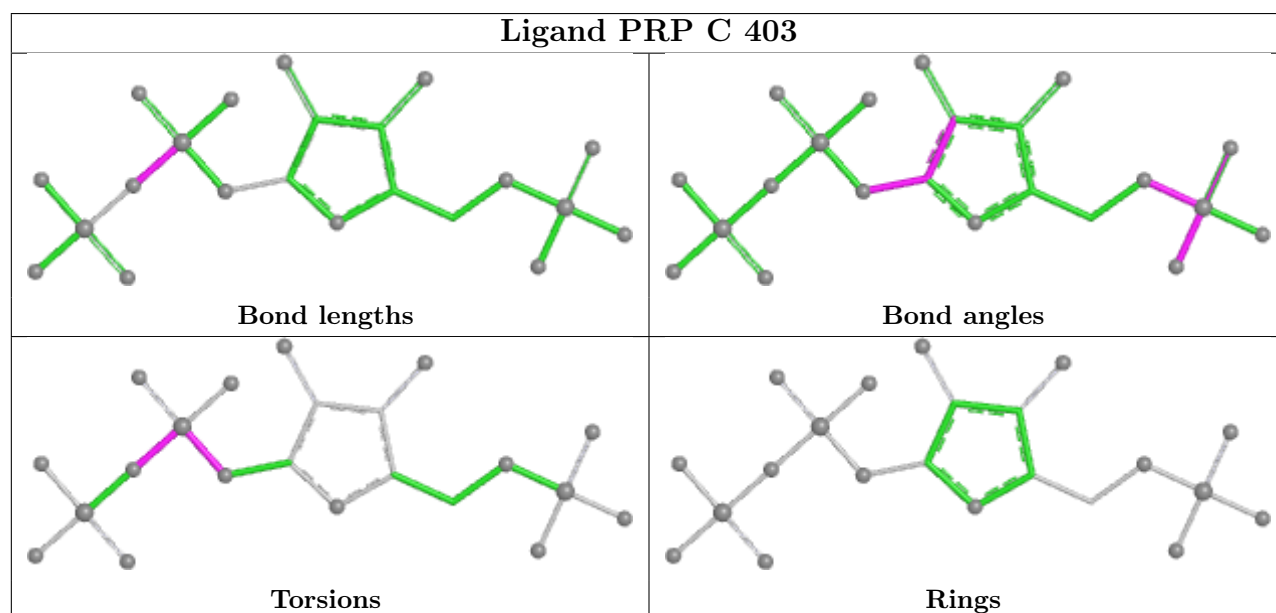
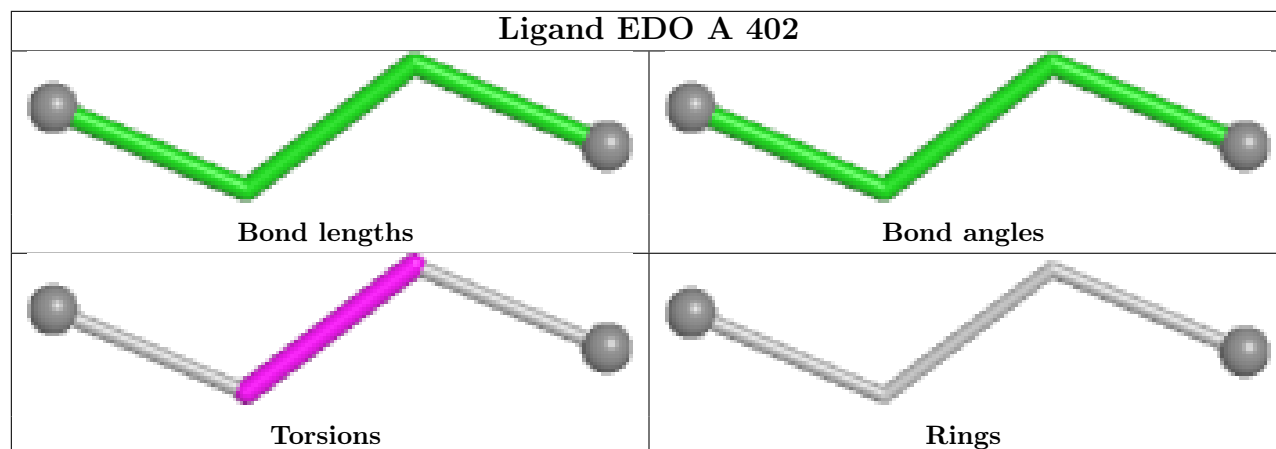
within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question 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.

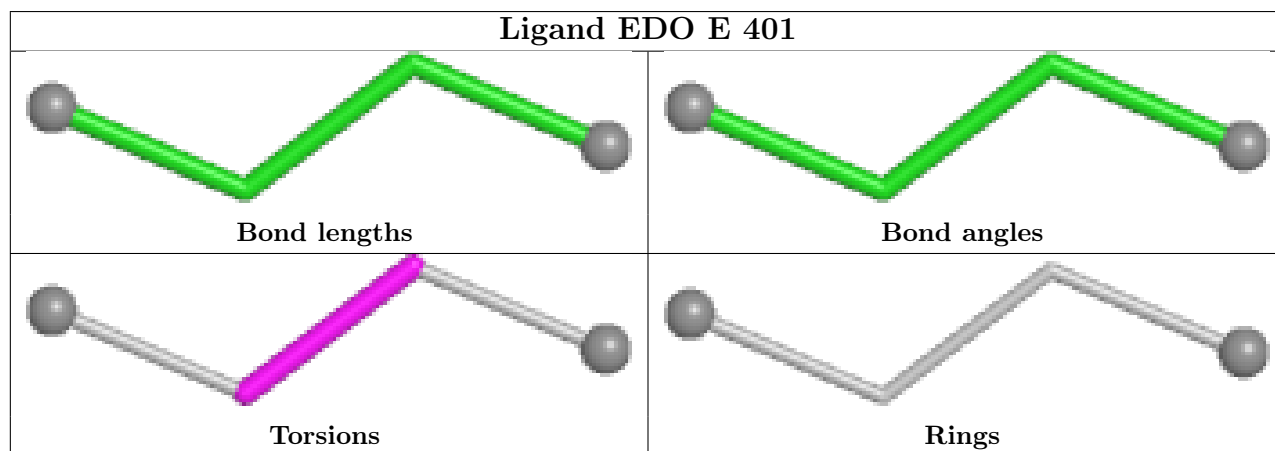
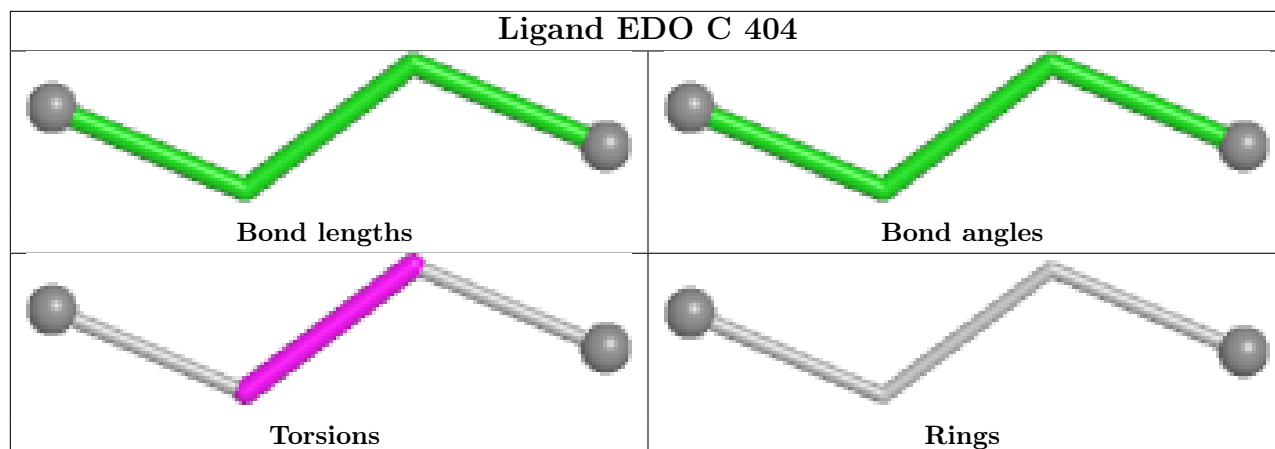


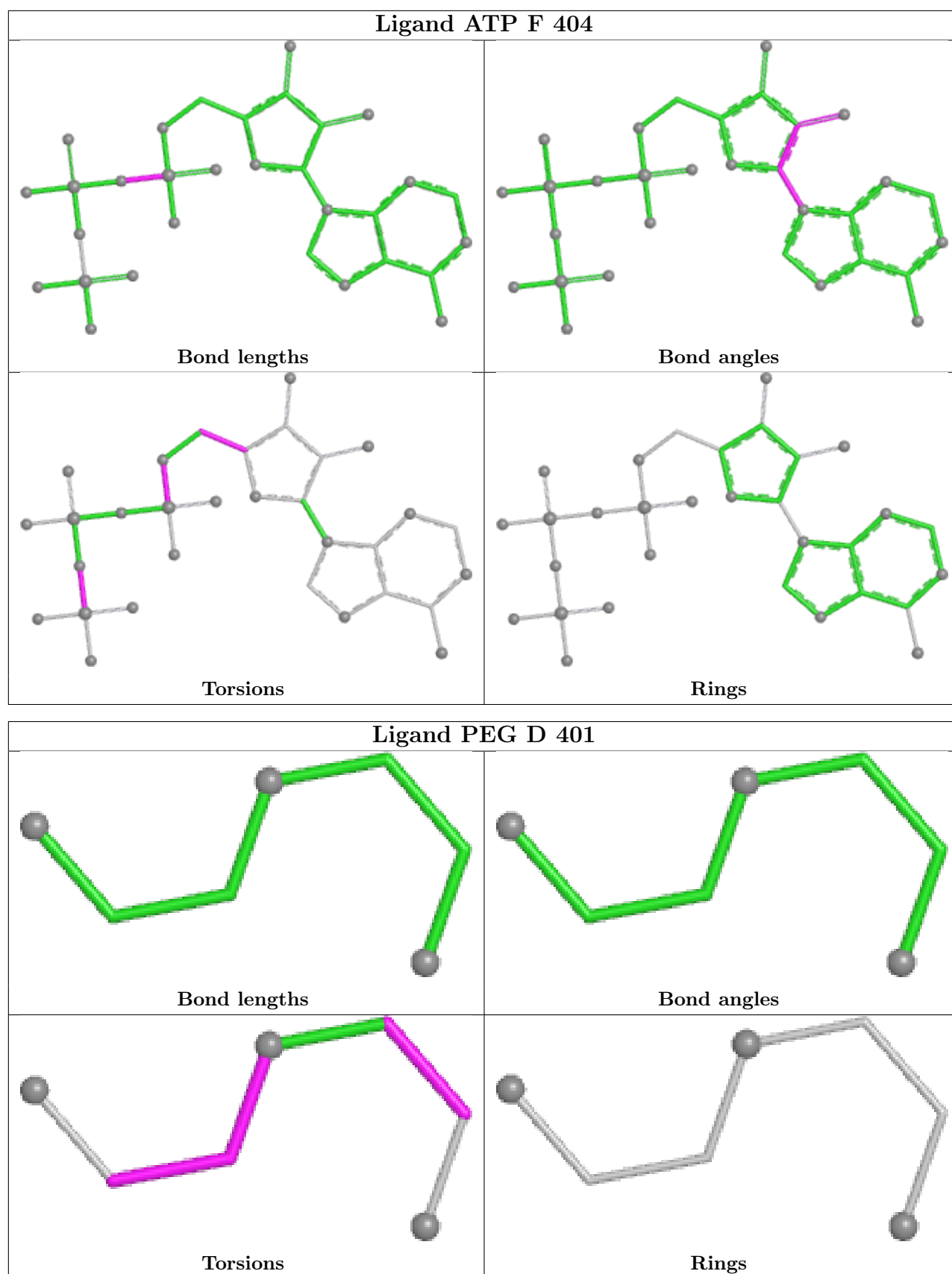


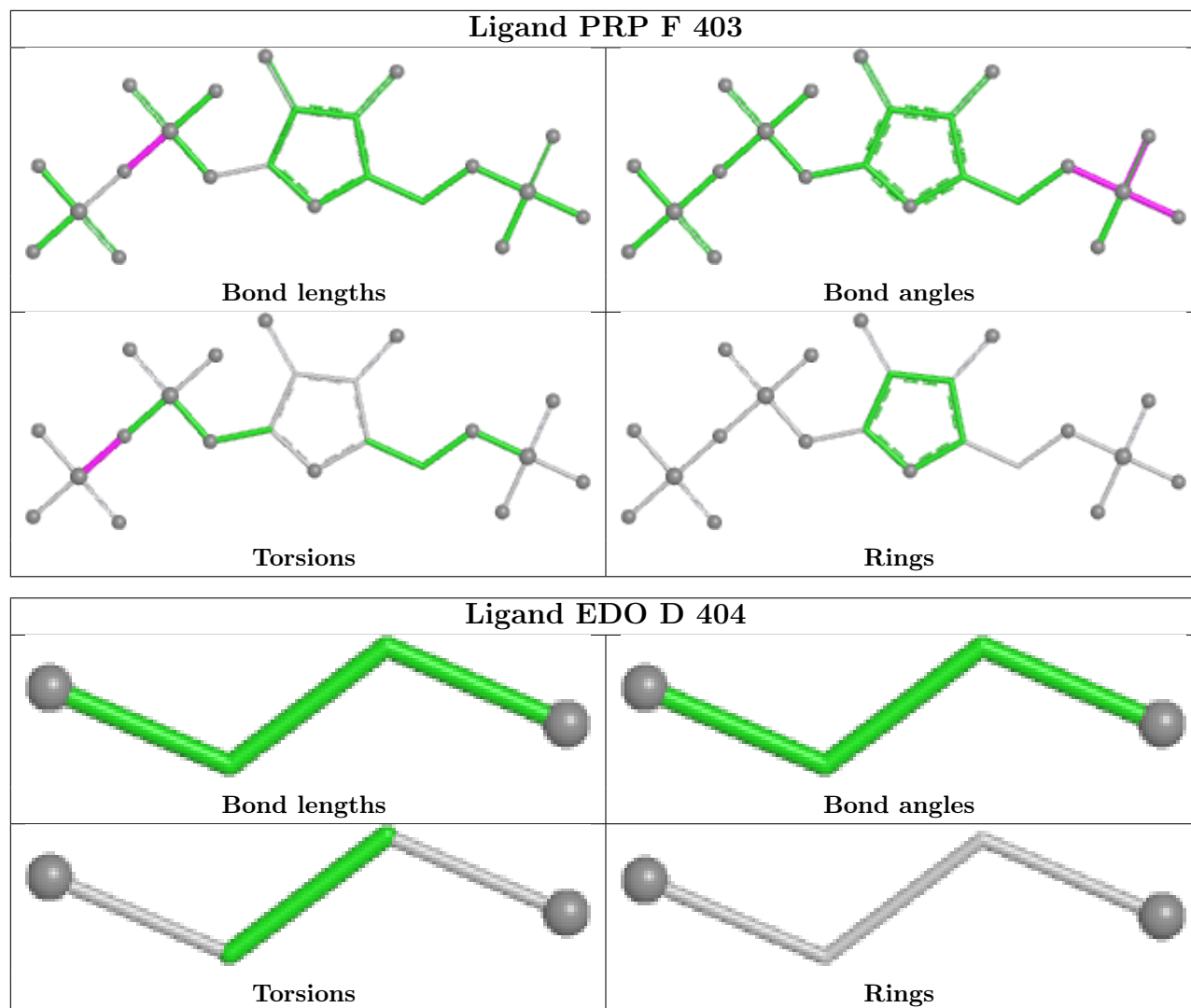


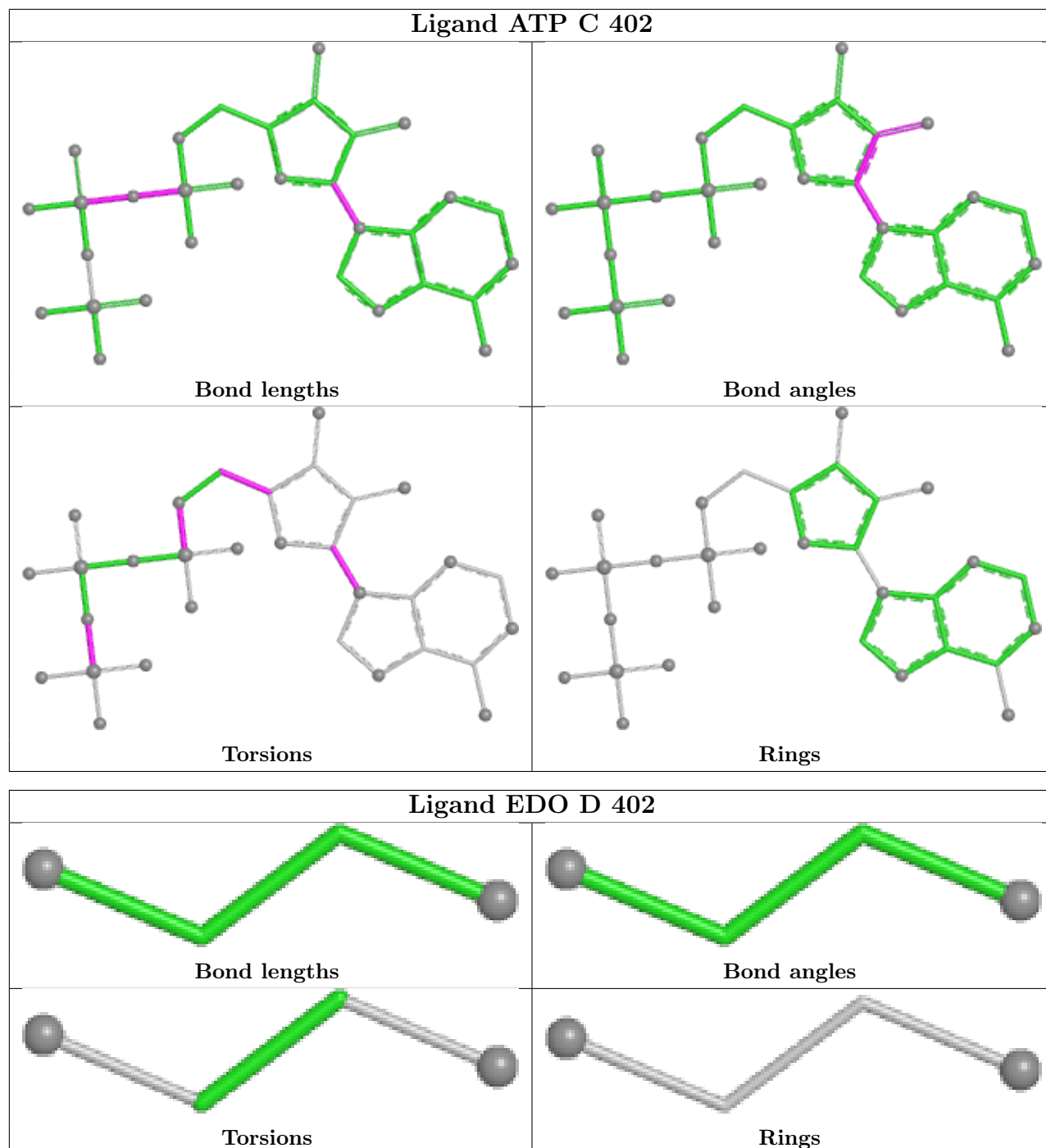


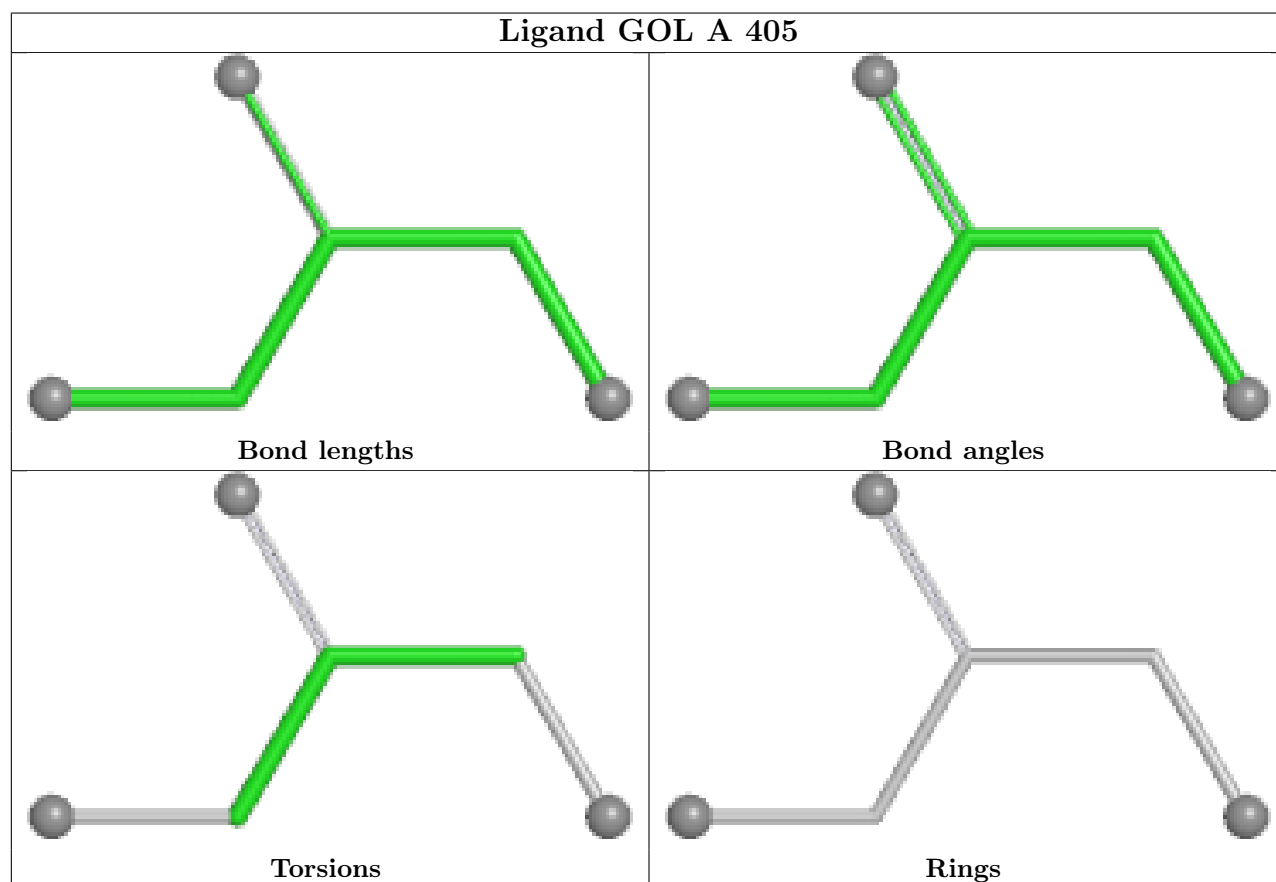
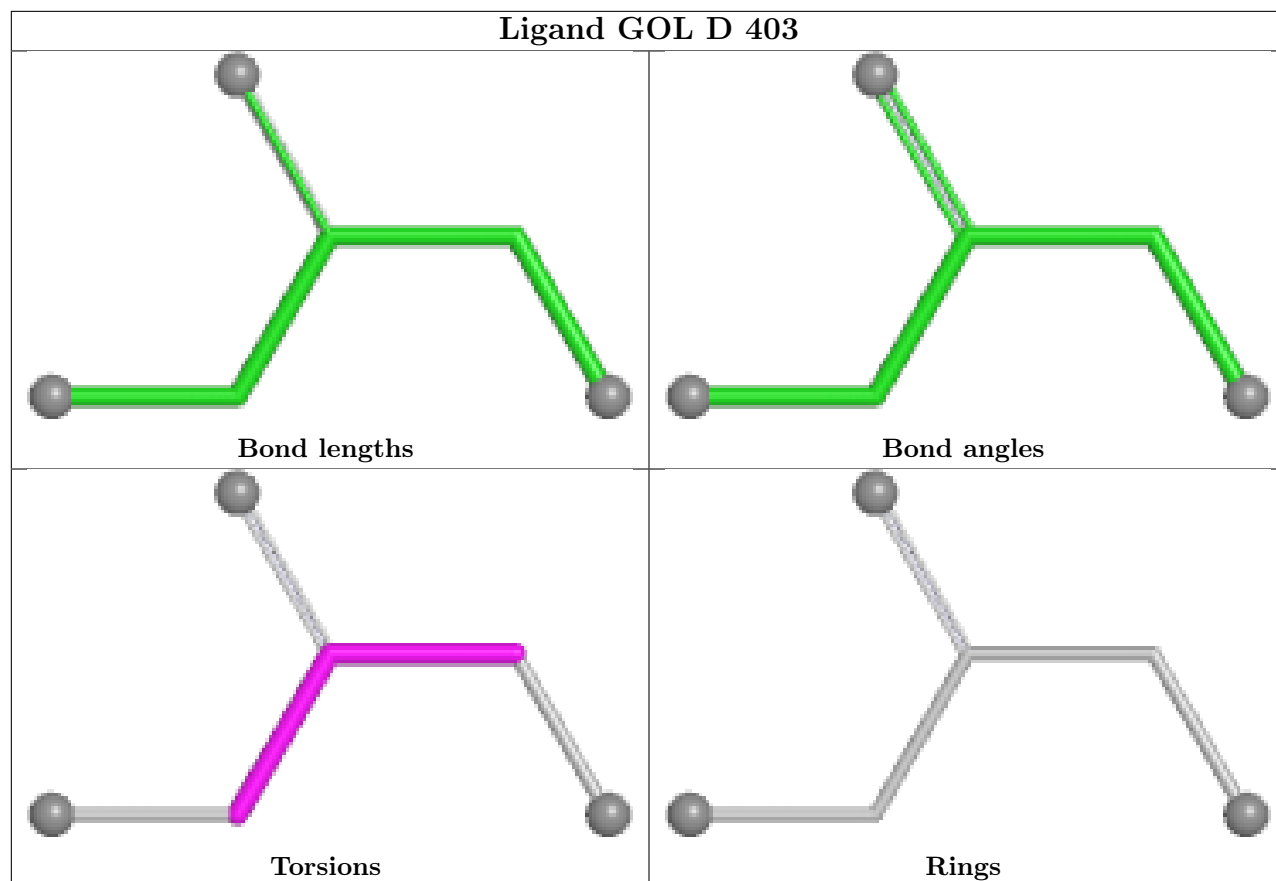












## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [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, 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	288/310 (92%)	-1.56	0 100 100	35, 62, 110, 136	1 (0%)
1	B	292/310 (94%)	-1.42	0 100 100	45, 78, 111, 128	1 (0%)
1	C	288/310 (92%)	-1.51	0 100 100	43, 67, 109, 137	0
1	D	287/310 (92%)	-1.56	0 100 100	38, 61, 108, 130	0
1	E	293/310 (94%)	-1.43	0 100 100	47, 78, 112, 125	1 (0%)
1	F	287/310 (92%)	-1.53	0 100 100	42, 68, 108, 138	1 (0%)
All	All	1735/1860 (93%)	-1.50	0 100 100	35, 69, 111, 138	4 (0%)

There are no RSRZ outliers to report.

### 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
2	PEG	A	401	7/7	0.97	0.10	84,87,89,90	0

*Continued on next page...*

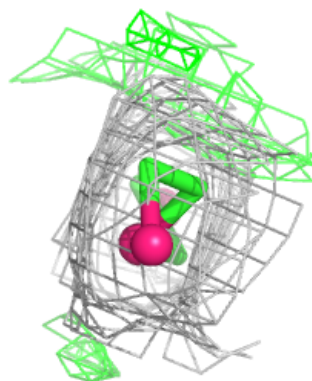
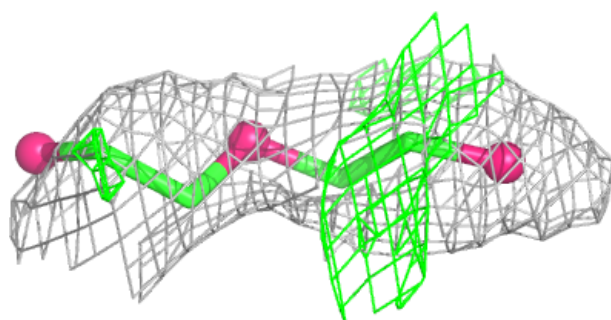
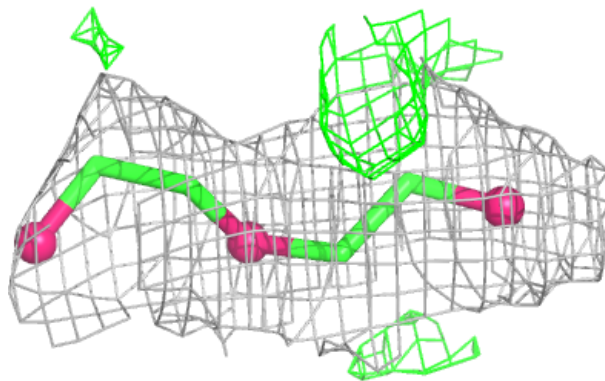
*Continued from previous page...*

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	PEG	D	401	7/7	0.97	0.08	80,89,102,102	0
3	EDO	C	404	4/4	0.98	0.07	91,91,92,92	0
3	EDO	F	402	4/4	0.98	0.06	77,82,85,85	0
3	EDO	A	402	4/4	0.99	0.07	74,75,77,77	0
3	EDO	D	402	4/4	0.99	0.07	91,96,96,97	0
3	EDO	D	405	4/4	0.99	0.05	71,76,81,87	0
3	EDO	E	401	4/4	0.99	0.05	89,89,89,90	0
3	EDO	A	403	4/4	0.99	0.06	74,75,76,77	0
4	PO4	A	404	5/5	0.99	0.04	97,104,113,114	0
4	PO4	D	406	5/5	0.99	0.04	106,116,123,123	0
5	GOL	A	405	6/6	0.99	0.04	69,72,77,80	0
5	GOL	D	403	6/6	0.99	0.04	67,68,70,71	0
7	ATP	C	402	31/31	0.99	0.04	81,111,120,124	1
7	ATP	F	404	31/31	0.99	0.04	78,97,116,118	12
8	PRP	C	403	22/22	0.99	0.04	104,159,200,202	0
8	PRP	F	403	22/22	0.99	0.03	106,168,191,198	0
6	MG	C	401	1/1	1.00	0.03	71,71,71,71	0
6	MG	F	401	1/1	1.00	0.02	76,76,76,76	0
3	EDO	D	404	4/4	1.00	0.04	69,75,78,79	0

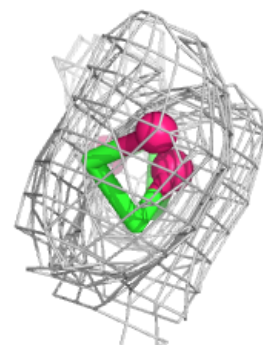
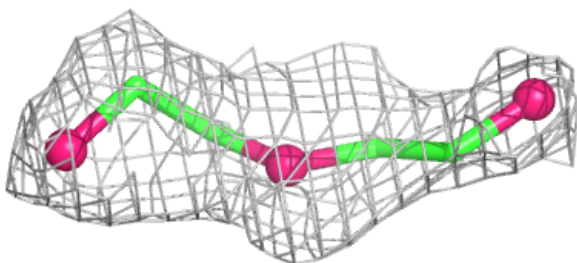
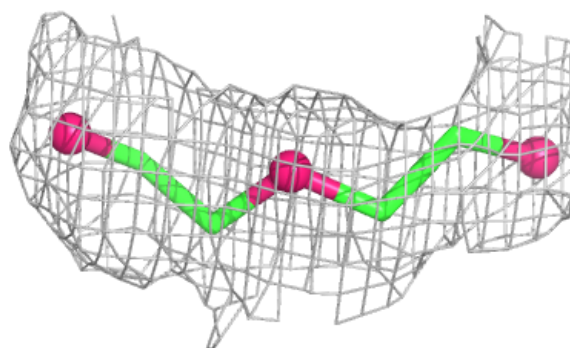
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

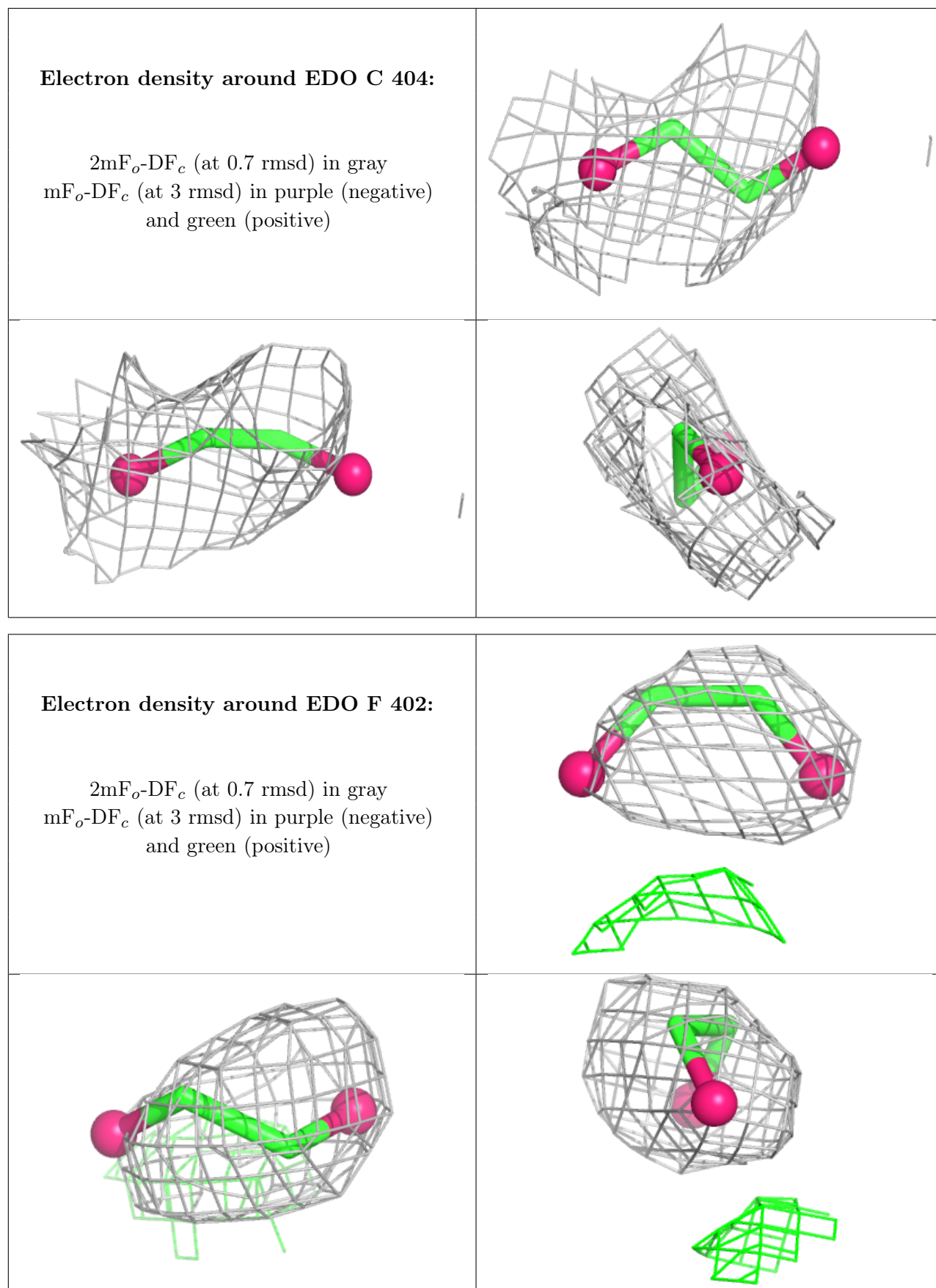
**Electron density around PEG A 401:**

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

**Electron density around PEG D 401:**

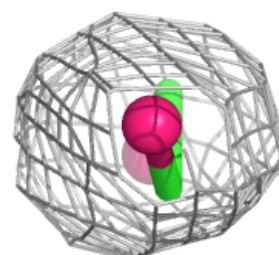
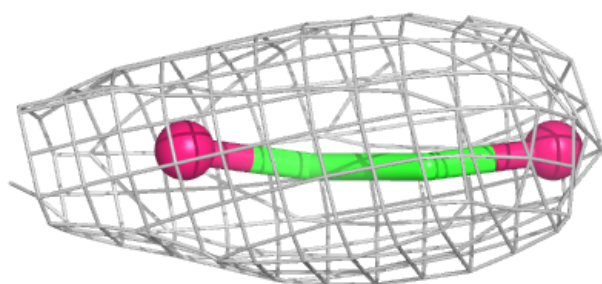
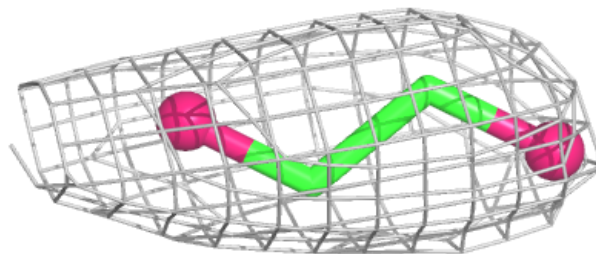
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



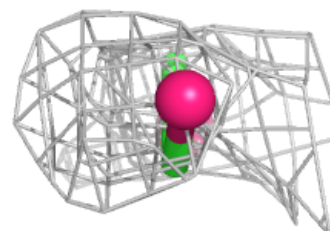
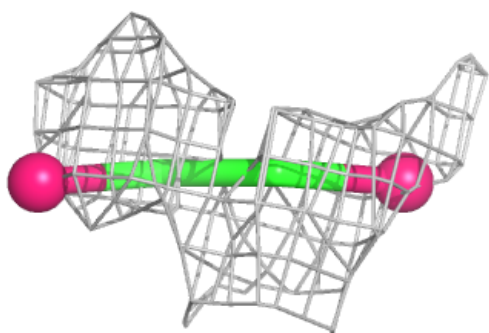
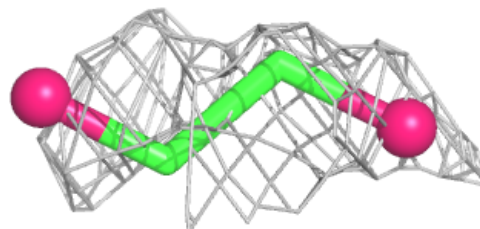


**Electron density around EDO A 402:**

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

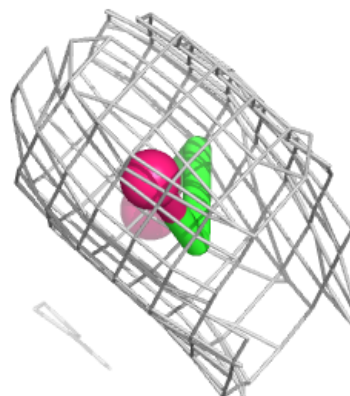
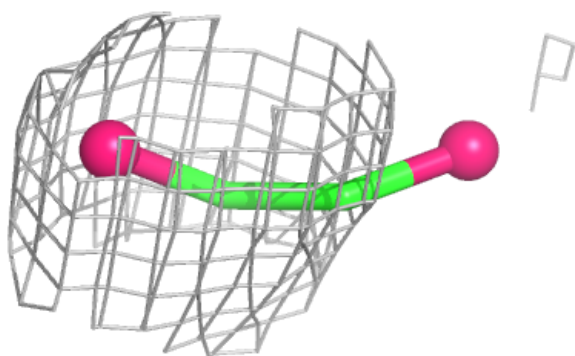
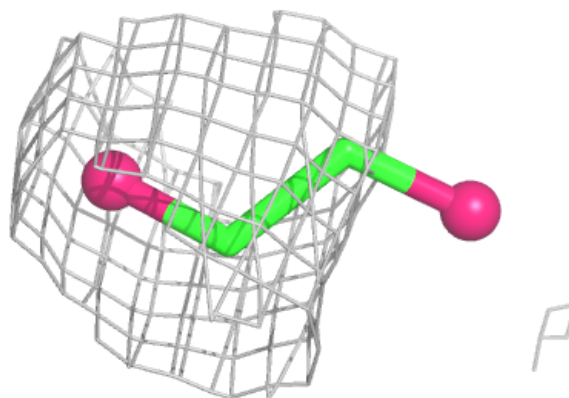
**Electron density around EDO D 402:**

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

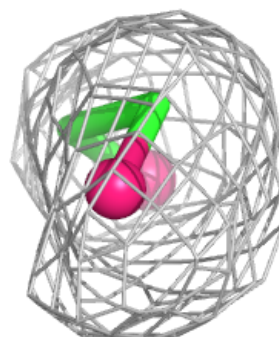
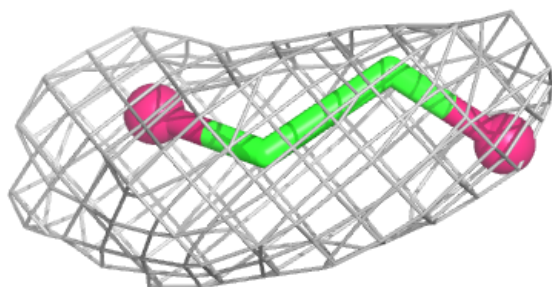
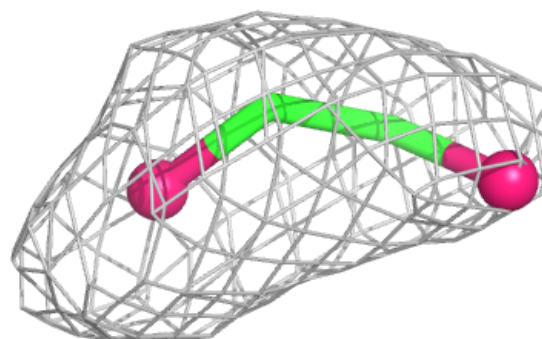


**Electron density around EDO D 405:**

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

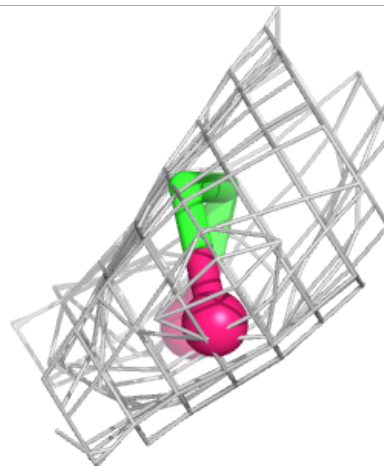
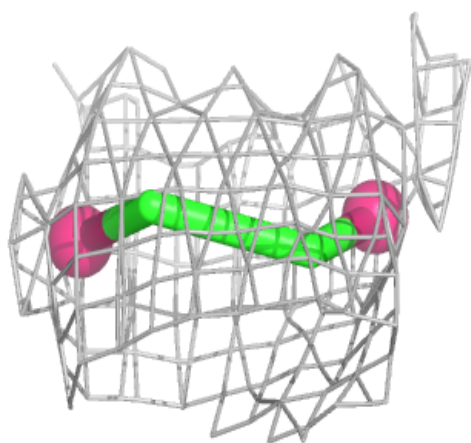
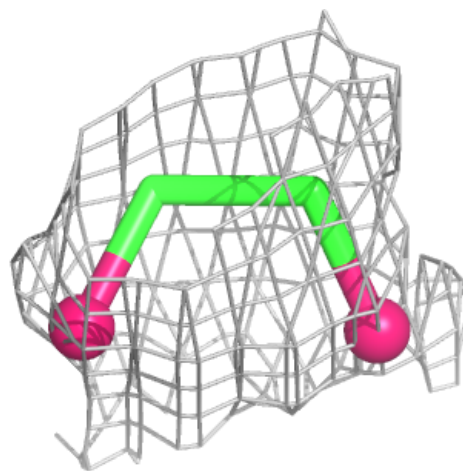
**Electron density around EDO E 401:**

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



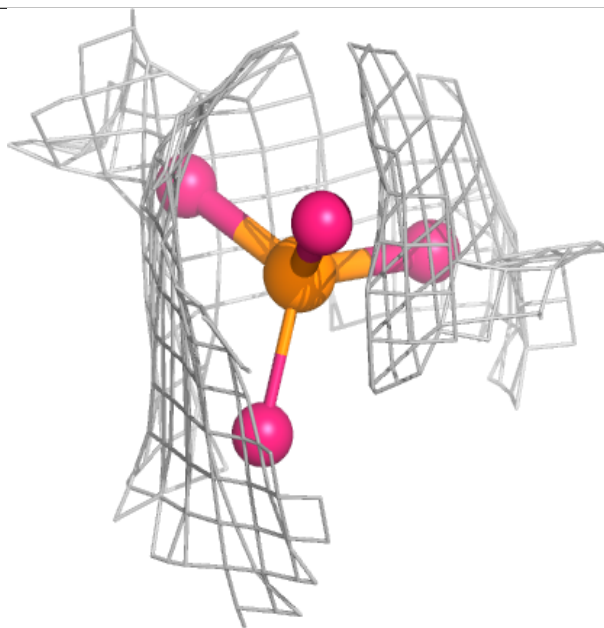
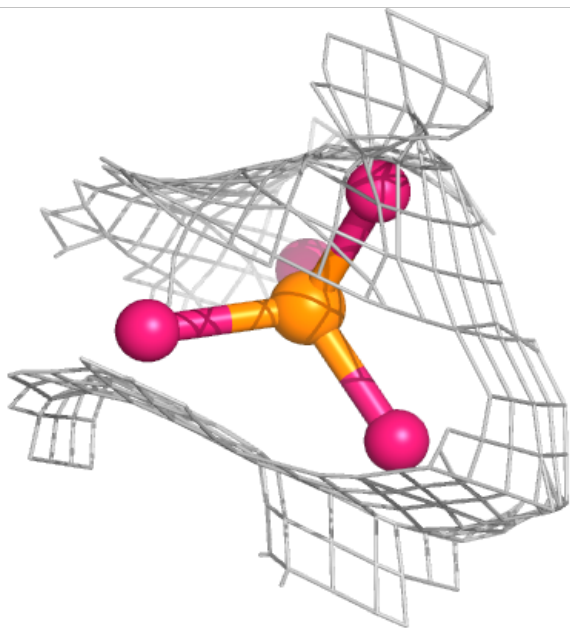
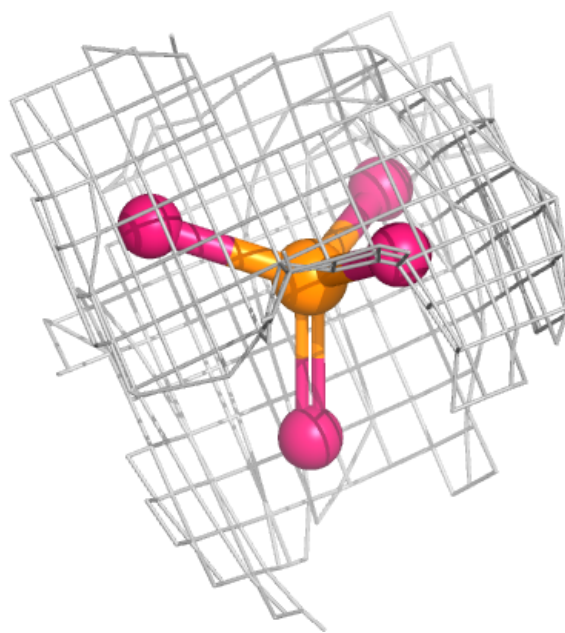
**Electron density around EDO A 403:**

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



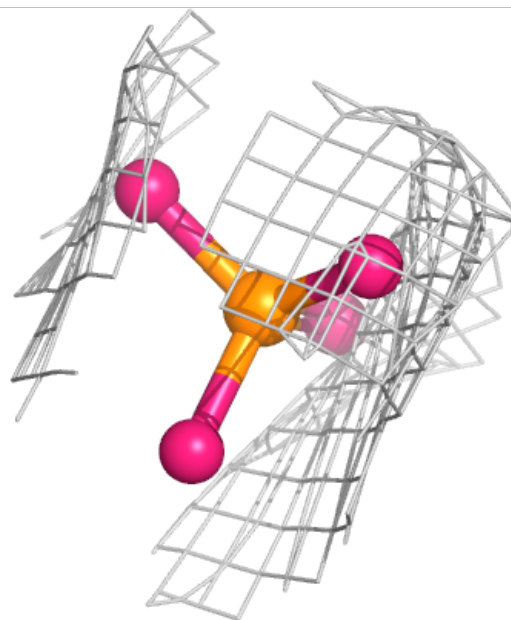
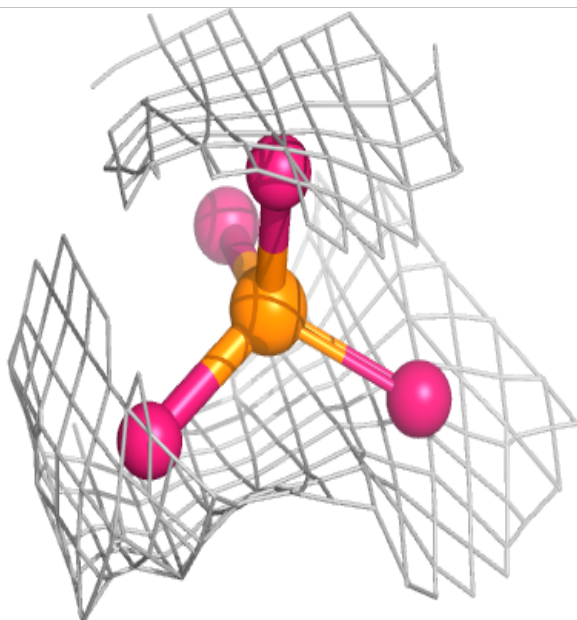
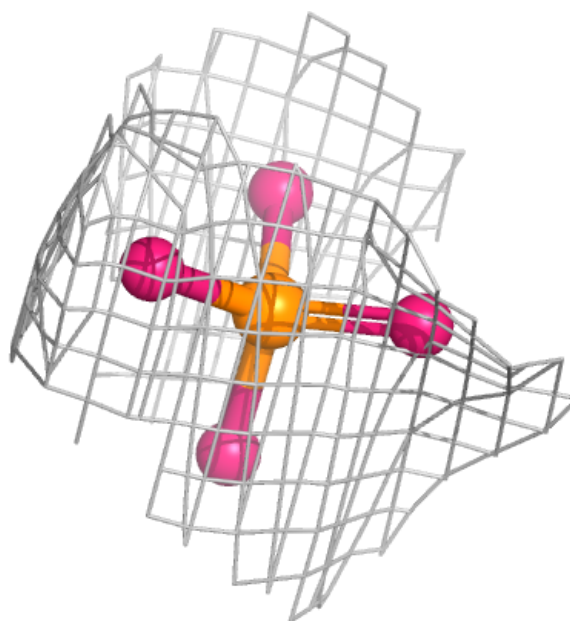
**Electron density around PO4 A 404:**

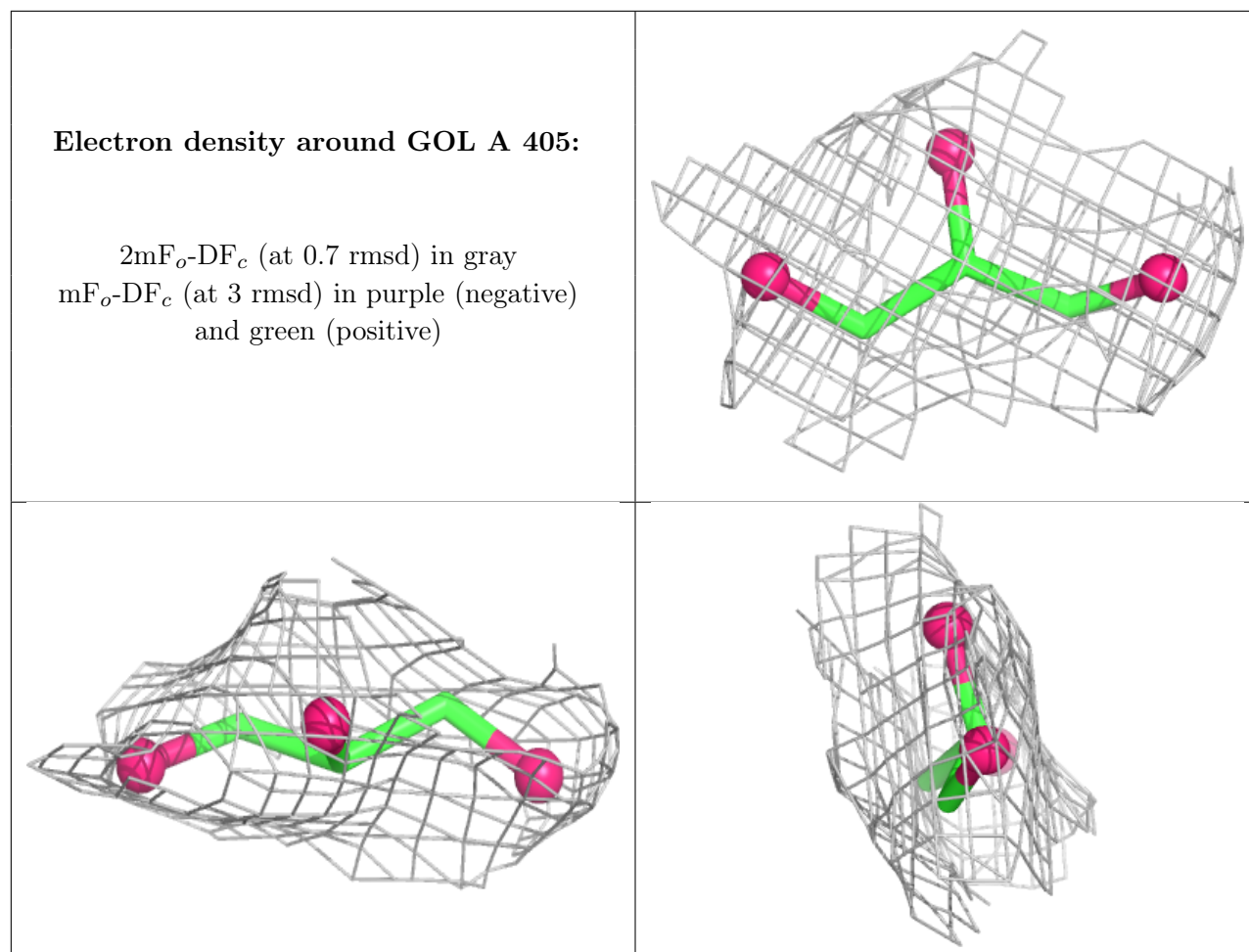
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



**Electron density around PO4 D 406:**

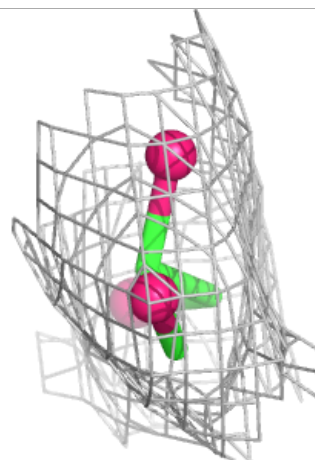
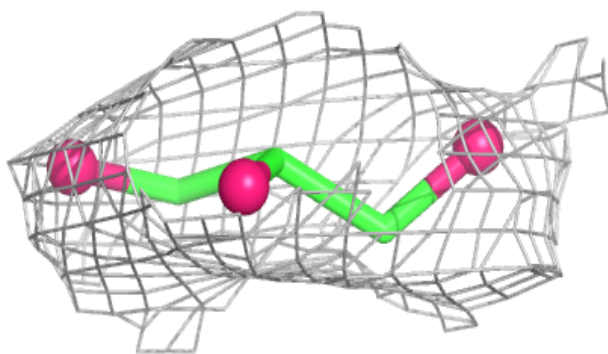
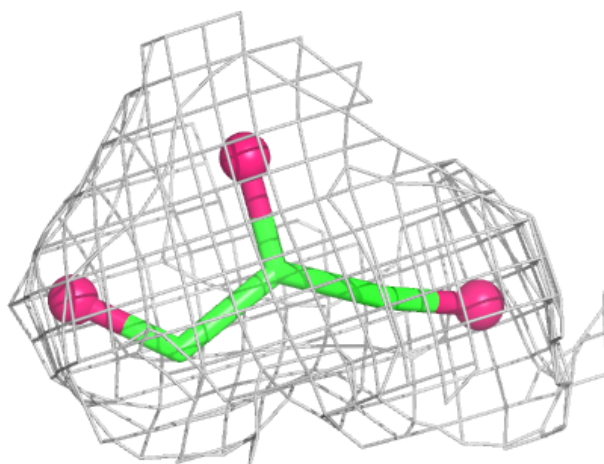
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





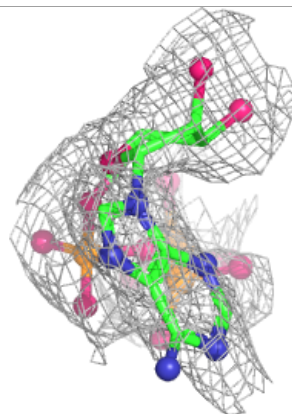
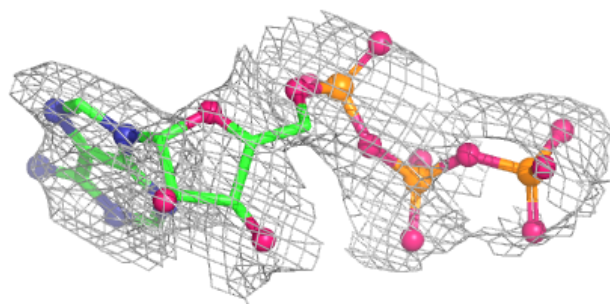
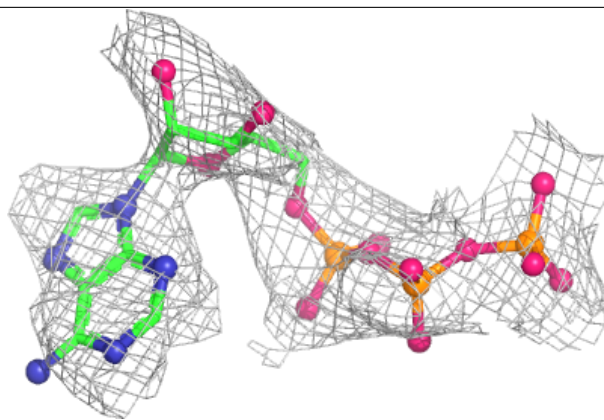
**Electron density around GOL D 403:**

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

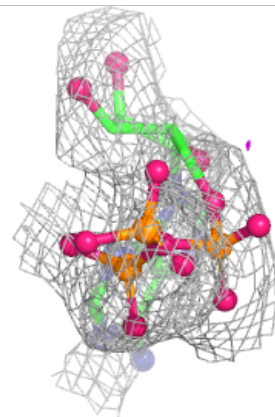
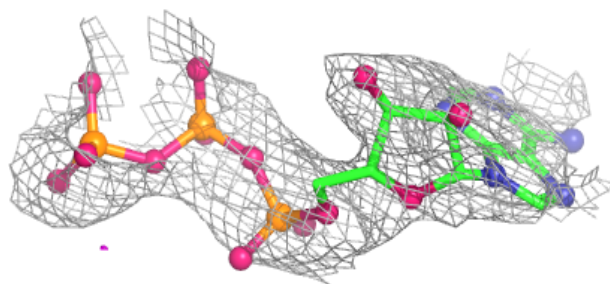
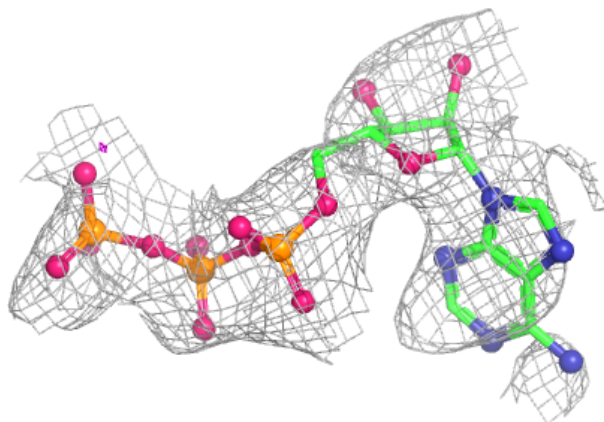


**Electron density around ATP C 402:**

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

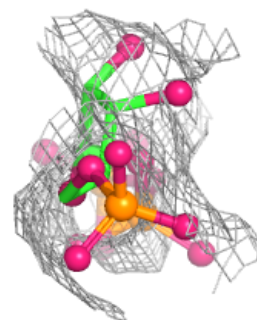
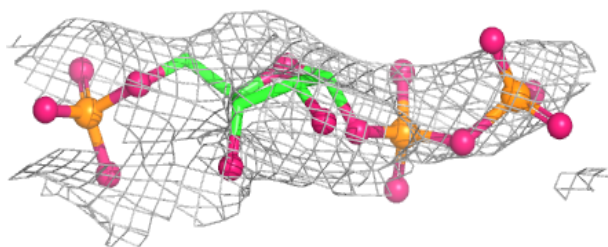
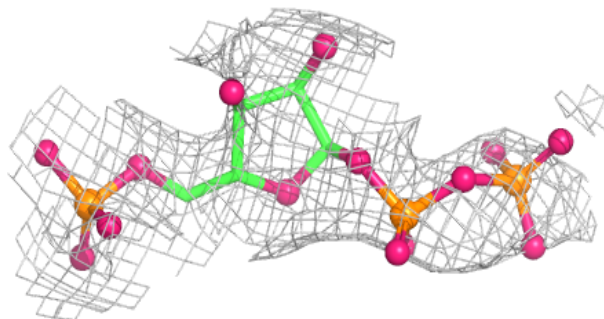
**Electron density around ATP F 404:**

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

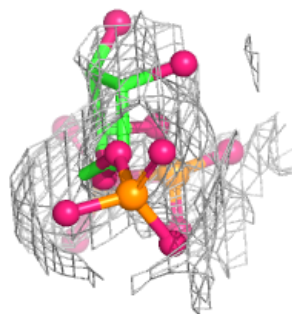
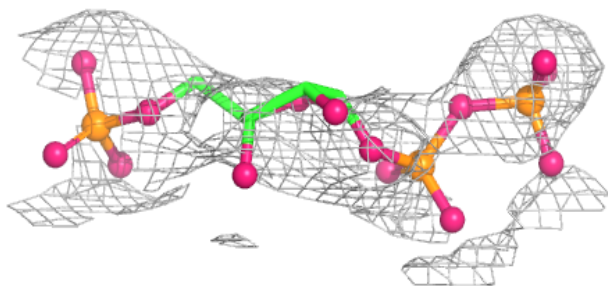
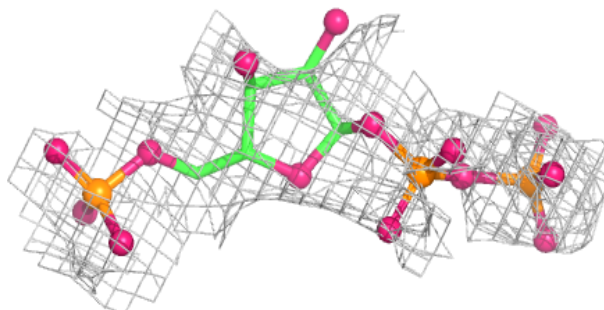


**Electron density around PRP C 403:**

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

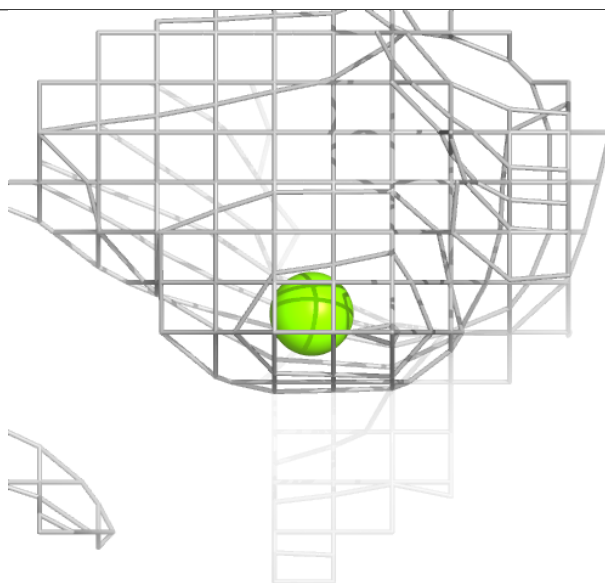
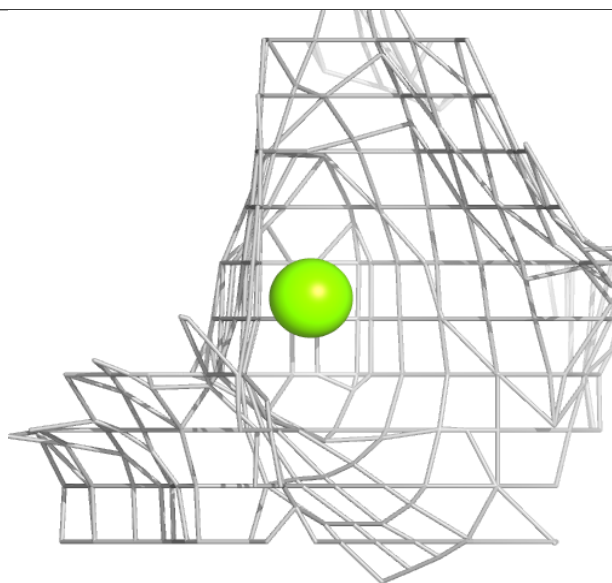
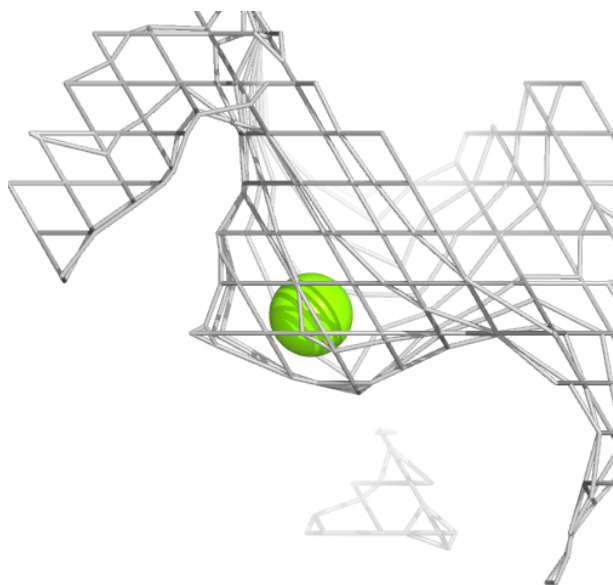
**Electron density around PRP F 403:**

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



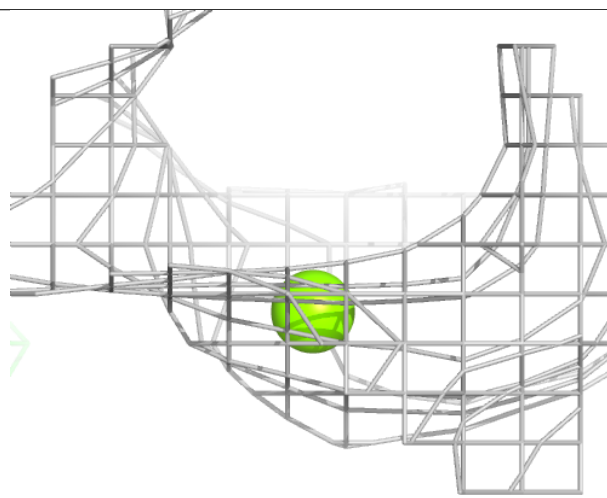
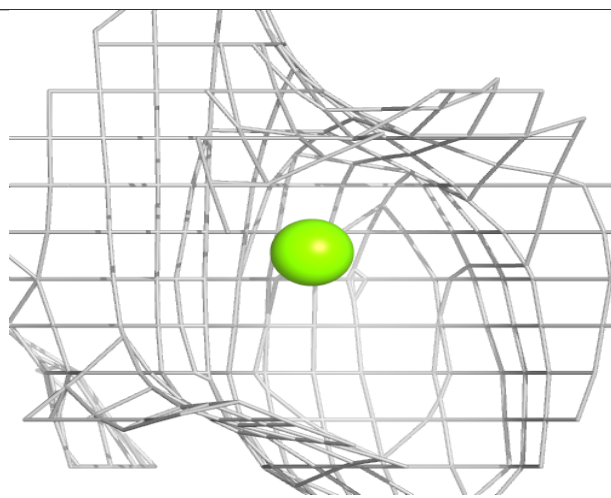
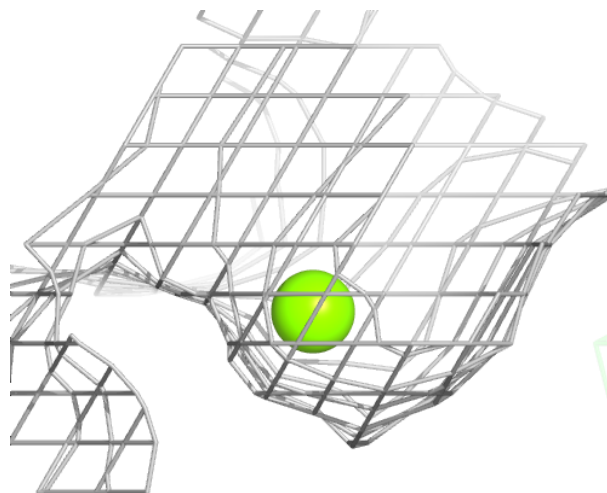
**Electron density around MG C 401:**

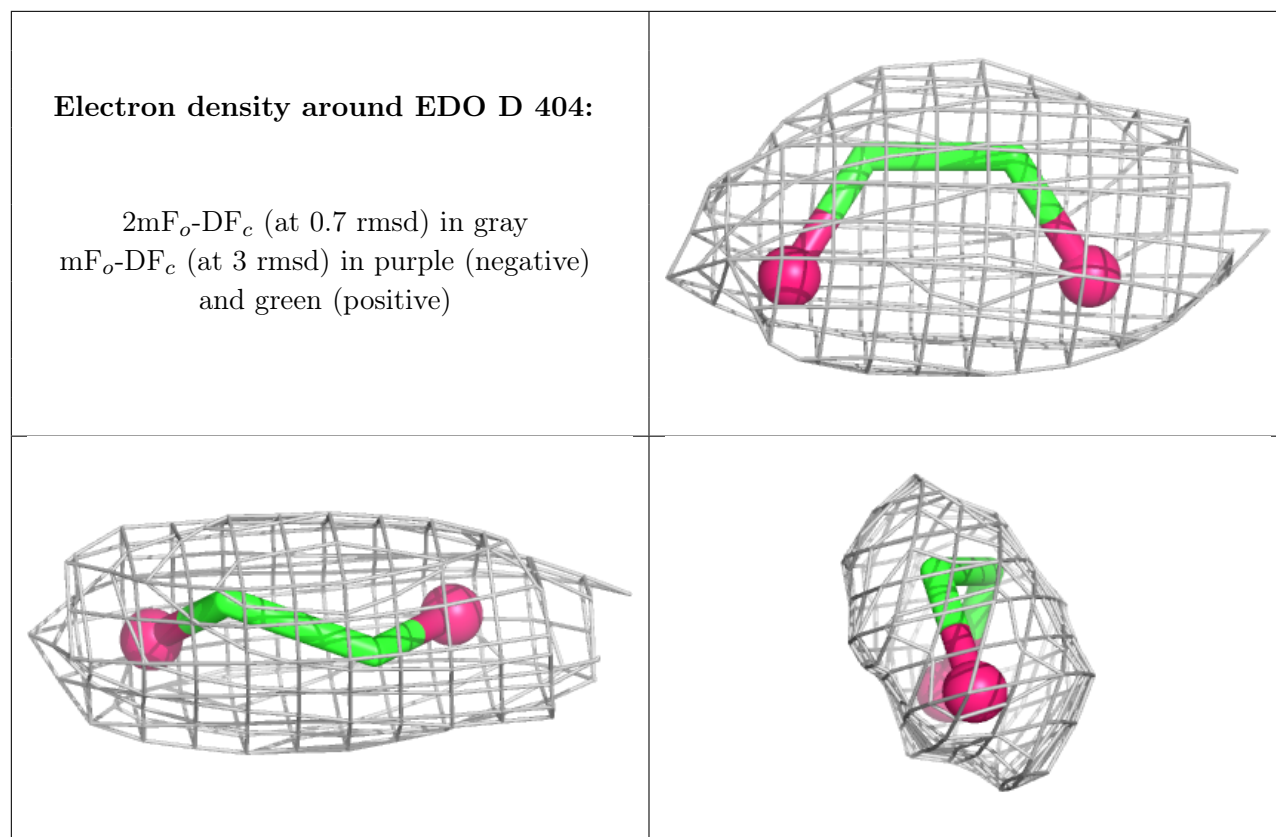
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



**Electron density around MG F 401:**

$2mF_o-DF_c$  (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.