



# wwPDB X-ray Structure Validation Summary Report ⓘ

Jan 31, 2024 – 02:25 PM EST

PDB ID : 8FMG  
Title : Structure of CBASS Cap5 from *Pseudomonas syringae* as an activated tetramer with the cyclic dinucleotide 3'2'-c-diAMP ligand (3 tetramers in the AU)  
Authors : Rechkoblit, O.; Kreitler, D.F.; Aggarwal, A.K.  
Deposited on : 2022-12-23  
Resolution : 1.79 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>.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.36  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.36

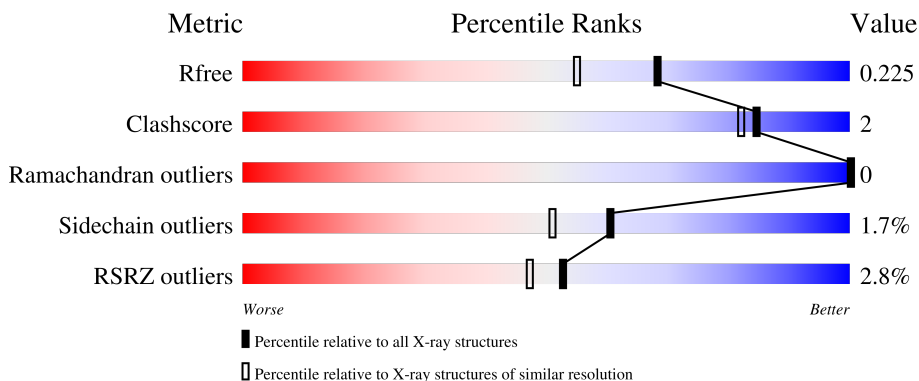
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.79 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	130704	5950 (1.80-1.80)
Clashscore	141614	6793 (1.80-1.80)
Ramachandran outliers	138981	6697 (1.80-1.80)
Sidechain outliers	138945	6696 (1.80-1.80)
RSRZ outliers	127900	5850 (1.80-1.80)

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	388	 2% 89% 5% 5%
1	B	388	 4% 83% 7% 9%
1	C	388	 3% 86% 8% 5%
1	D	388	 2% 85% 6% 9%

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Mol	Chain	Length	Quality of chain
1	E	388	<p>3% 89% 6% 5%</p>
1	F	388	<p>3% 84% 6% 9%</p>
1	G	388	<p>1% 86% 9% 5%</p>
1	H	388	<p>3% 87% 6% 8%</p>
1	K	388	<p>2% 92% 5% 1%</p>
1	L	388	<p>1% 85% 6% 8%</p>
1	M	388	<p>3% 90% 5% 5%</p>
1	N	388	<p>4% 86% 5% 9%</p>

## 2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 69460 atoms, of which 32951 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 SAVED domain-containing protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
1	A	369	5714	1817	2826	520	540	11	0	4	0
1	B	354	5368	1718	2656	482	502	10	0	0	0
1	C	368	5600	1784	2764	512	530	10	0	0	0
1	D	353	5340	1713	2635	481	501	10	0	0	0
1	E	369	5597	1791	2754	510	532	10	0	0	0
1	F	353	5412	1730	2677	485	510	10	0	1	0
1	G	368	5640	1793	2795	514	528	10	0	0	0
1	H	356	5431	1737	2685	486	513	10	0	0	0
1	K	369	5580	1787	2745	508	530	10	0	1	0
1	L	357	5480	1748	2713	491	518	10	0	1	0
1	M	369	5637	1794	2788	514	531	10	0	0	0
1	N	354	5382	1727	2661	484	500	10	0	2	0

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Zn	0	0
			1	1		
2	B	1	Total	Zn	0	0
			1	1		

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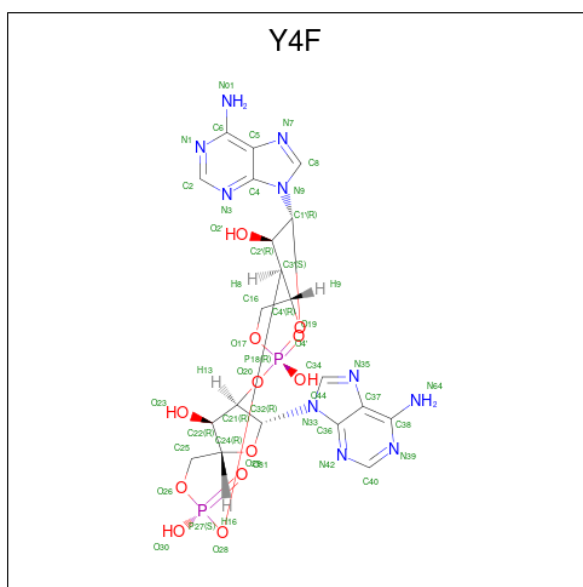
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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	C	1	Total Zn 1 1	0	0
2	D	1	Total Zn 1 1	0	0
2	E	1	Total Zn 1 1	0	0
2	F	1	Total Zn 1 1	0	0
2	G	1	Total Zn 1 1	0	0
2	H	1	Total Zn 1 1	0	0
2	K	1	Total Zn 1 1	0	0
2	L	1	Total Zn 1 1	0	0
2	M	1	Total Zn 1 1	0	0
2	N	1	Total Zn 1 1	0	0

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	2	Total Mg 2 2	0	0
3	C	2	Total Mg 2 2	0	0
3	E	2	Total Mg 2 2	0	0
3	G	2	Total Mg 2 2	0	0
3	H	1	Total Mg 1 1	0	0
3	K	2	Total Mg 2 2	0	0
3	M	2	Total Mg 2 2	0	0

- Molecule 4 is Cyclic (adenosine-(2'-5')-monophosphate-adenosine-(3'-5')-monophosphate (three-letter code: Y4F) (formula: C<sub>20</sub>H<sub>24</sub>N<sub>10</sub>O<sub>12</sub>P<sub>2</sub>) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	
			Total	C	H	N	O			P
4	A	1	Total	C	H	N	O	P	0	0
			65	20	21	10	12	2		
4	B	1	Total	C	H	N	O	P	0	0
			65	20	21	10	12	2		
4	C	1	Total	C	H	N	O	P	0	0
			65	20	21	10	12	2		
4	D	1	Total	C	H	N	O	P	0	0
			65	20	21	10	12	2		
4	E	1	Total	C	H	N	O	P	0	0
			65	20	21	10	12	2		
4	F	1	Total	C	H	N	O	P	0	0
			65	20	21	10	12	2		
4	G	1	Total	C	H	N	O	P	0	0
			65	20	21	10	12	2		
4	H	1	Total	C	H	N	O	P	0	0
			65	20	21	10	12	2		
4	K	1	Total	C	H	N	O	P	0	0
			65	20	21	10	12	2		
4	L	1	Total	C	H	N	O	P	0	0
			65	20	21	10	12	2		
4	M	1	Total	C	H	N	O	P	0	0
			65	20	21	10	12	2		
4	N	1	Total	C	H	N	O	P	0	0
			65	20	21	10	12	2		

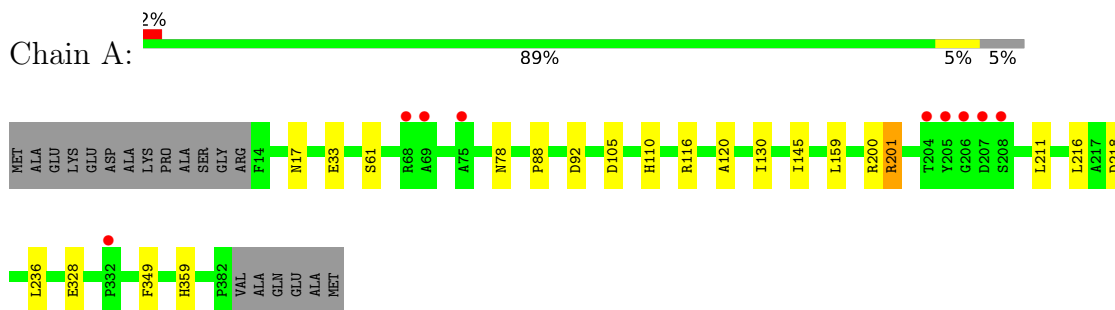
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	230	Total O 230 230	0	0
5	B	218	Total O 218 218	0	0
5	C	206	Total O 206 206	0	0
5	D	141	Total O 141 141	0	0
5	E	252	Total O 252 252	0	0
5	F	162	Total O 162 162	0	0
5	G	232	Total O 232 232	0	0
5	H	212	Total O 212 212	0	0
5	K	233	Total O 233 233	0	0
5	L	223	Total O 223 223	0	0
5	M	202	Total O 202 202	0	0
5	N	163	Total O 163 163	0	0

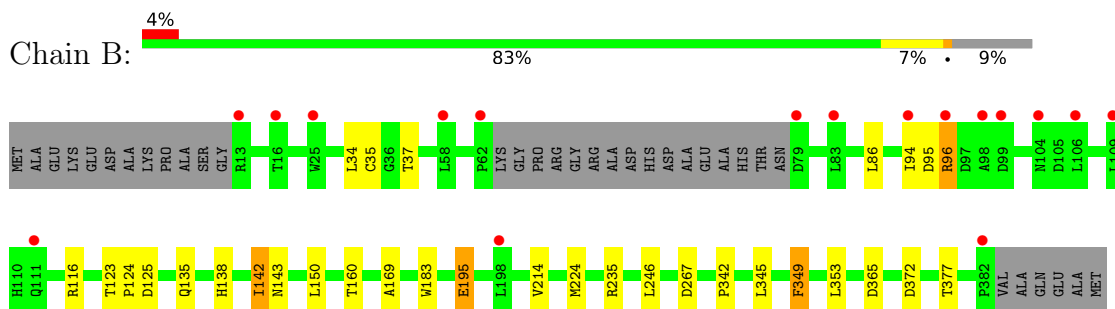
### 3 Residue-property plots [i](#)

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.

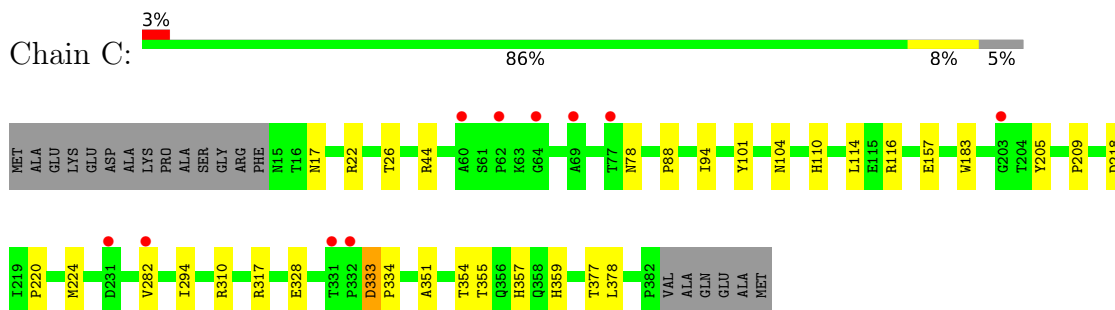
- Molecule 1: SAVED domain-containing protein



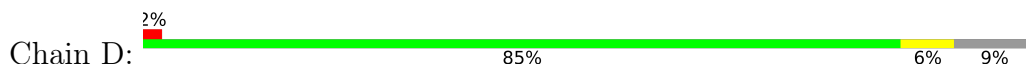
- Molecule 1: SAVED domain-containing protein



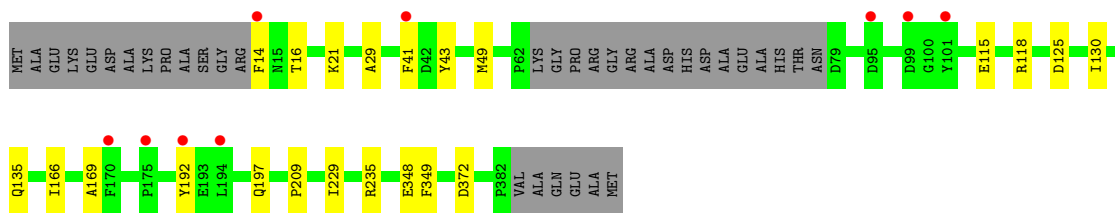
- Molecule 1: SAVED domain-containing protein



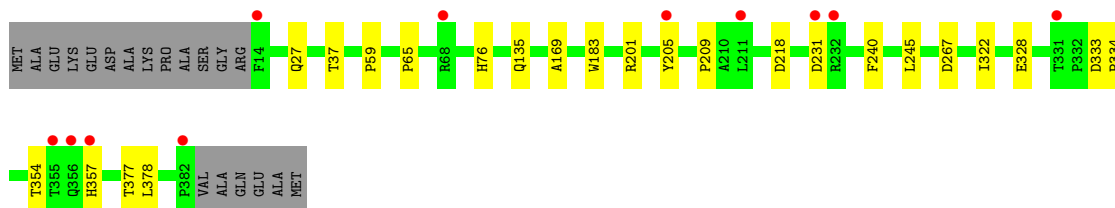
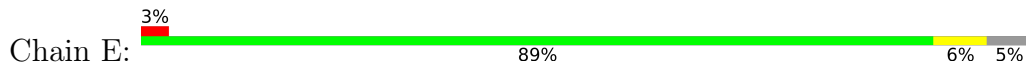
- Molecule 1: SAVED domain-containing protein



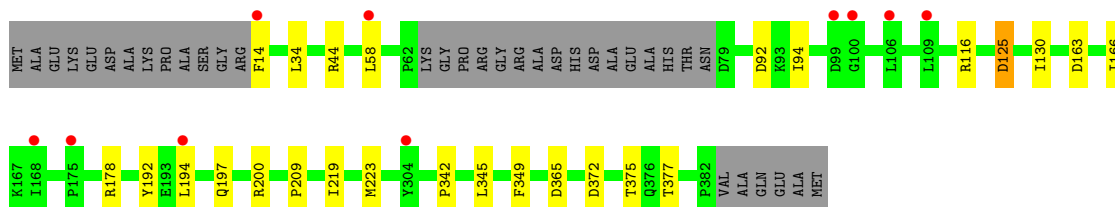
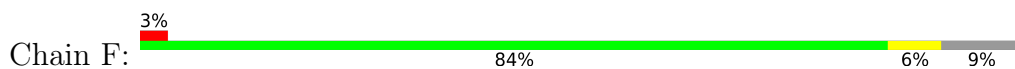




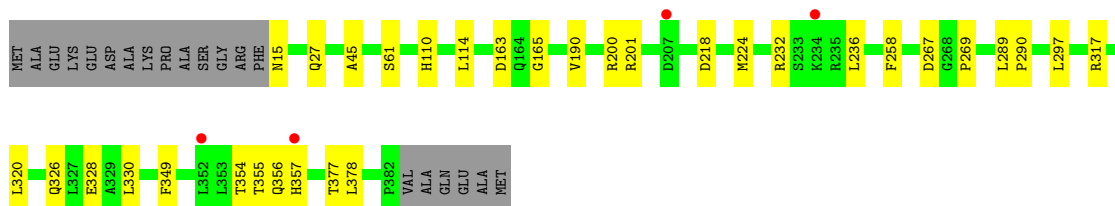
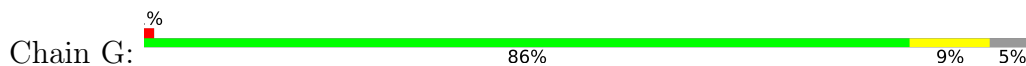
- Molecule 1: SAVED domain-containing protein



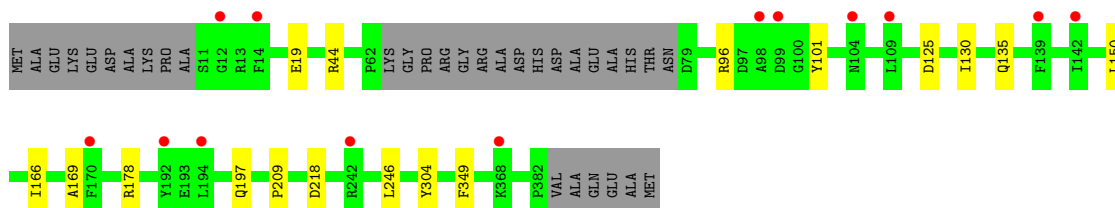
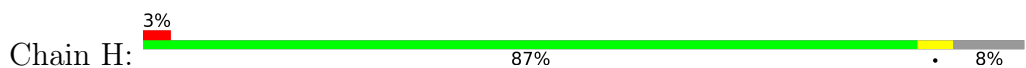
- Molecule 1: SAVED domain-containing protein



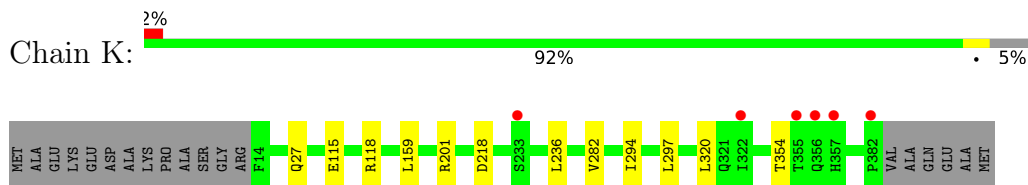
- Molecule 1: SAVED domain-containing protein



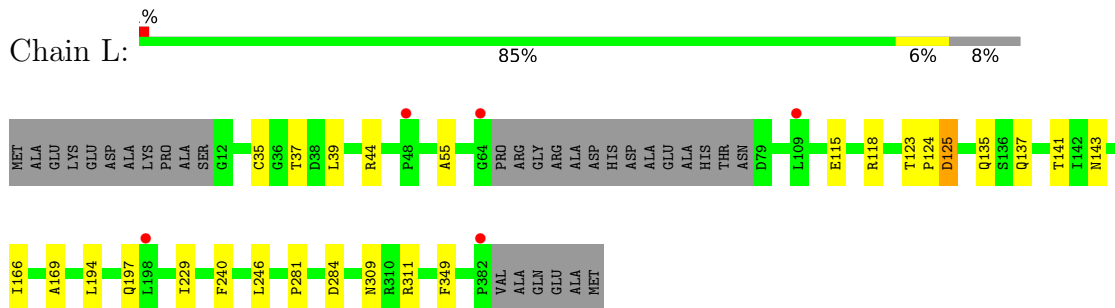
- Molecule 1: SAVED domain-containing protein



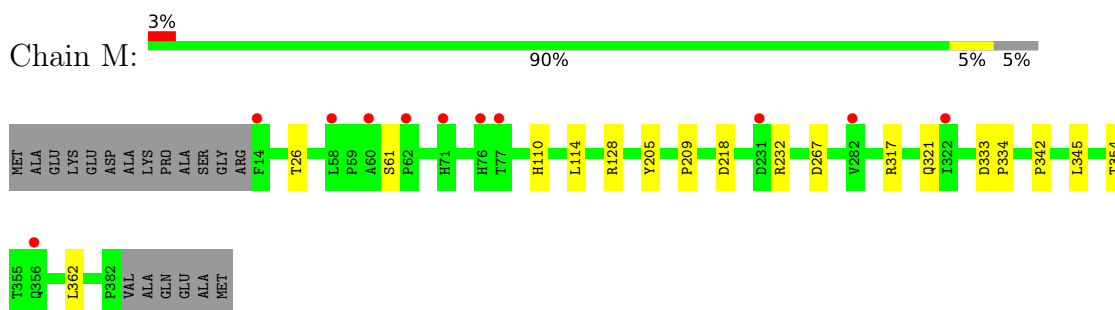
- Molecule 1: SAVED domain-containing protein



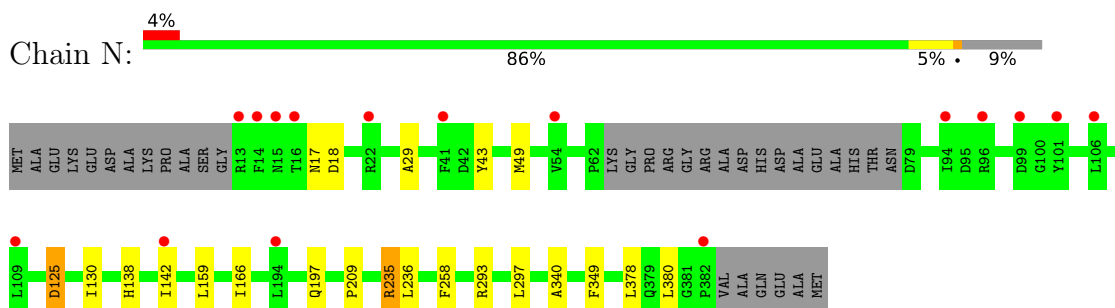
- Molecule 1: SAVED domain-containing protein



- Molecule 1: SAVED domain-containing protein



- Molecule 1: SAVED domain-containing protein



## 4 Data and refinement statistics i

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	145.19Å 80.52Å 406.84Å 90.00° 90.62° 90.00°	Depositor
Resolution (Å)	58.12 – 1.79 406.82 – 1.79	Depositor EDS
% Data completeness (in resolution range)	61.0 (58.12-1.79) 61.0 (406.82-1.79)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.54 (at 1.80Å)	Xtrriage
Refinement program	PHENIX 1.20.1_4487, PHENIX 1.20.1_4487	Depositor
R, $R_{free}$	0.179 , 0.225 0.180 , 0.225	Depositor DCC
$R_{free}$ test set	12884 reflections (4.82%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	27.0	Xtrriage
Anisotropy	0.067	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.41 , 45.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	0.006 for -1/2*h-3/2*k,-1/2*h+1/2*k,-l 0.006 for -1/2*h+3/2*k,1/2*h+1/2*k,-l 0.025 for 1/2*h-3/2*k,-1/2*h-1/2*k,-l 0.020 for 1/2*h+3/2*k,1/2*h-1/2*k,-l 0.007 for -h,-k,l	Xtrriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	69460	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	36.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.14% 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: Y4F, ZN, 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	0.30	0/2958	0.56	0/4031
1	B	0.29	0/2777	0.55	0/3788
1	C	0.28	0/2905	0.54	0/3961
1	D	0.27	0/2771	0.53	0/3782
1	E	0.28	0/2913	0.56	0/3975
1	F	0.28	0/2801	0.56	0/3820
1	G	0.28	0/2914	0.54	0/3971
1	H	0.28	0/2812	0.54	0/3832
1	K	0.29	0/2904	0.55	0/3963
1	L	0.29	0/2833	0.55	0/3861
1	M	0.28	0/2918	0.54	0/3978
1	N	0.28	0/2787	0.54	0/3804
All	All	0.28	0/34293	0.55	0/46766

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 [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	2888	2826	2821	14	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	2712	2656	2654	20	0
1	C	2836	2764	2763	21	0
1	D	2705	2635	2633	12	0
1	E	2843	2754	2753	15	0
1	F	2735	2677	2674	14	0
1	G	2845	2795	2794	15	0
1	H	2746	2685	2683	10	0
1	K	2835	2745	2743	8	0
1	L	2767	2713	2710	16	0
1	M	2849	2788	2787	9	0
1	N	2721	2661	2657	13	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
2	E	1	0	0	0	0
2	F	1	0	0	0	0
2	G	1	0	0	0	0
2	H	1	0	0	0	0
2	K	1	0	0	0	0
2	L	1	0	0	0	0
2	M	1	0	0	0	0
2	N	1	0	0	0	0
3	A	2	0	0	0	0
3	C	2	0	0	0	0
3	E	2	0	0	0	0
3	G	2	0	0	0	0
3	H	1	0	0	0	0
3	K	2	0	0	0	0
3	M	2	0	0	0	0
4	A	44	21	0	0	0
4	B	44	21	0	0	0
4	C	44	21	0	0	0
4	D	44	21	0	0	0
4	E	44	21	0	0	0
4	F	44	21	0	0	0
4	G	44	21	0	0	0
4	H	44	21	0	0	0
4	K	44	21	0	0	0
4	L	44	21	0	1	0
4	M	44	21	0	0	0
4	N	44	21	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	A	230	0	0	4	0
5	B	218	0	0	3	0
5	C	206	0	0	3	0
5	D	141	0	0	0	0
5	E	252	0	0	3	0
5	F	162	0	0	3	0
5	G	232	0	0	3	0
5	H	212	0	0	1	0
5	K	233	0	0	1	0
5	L	223	0	0	1	0
5	M	202	0	0	1	0
5	N	163	0	0	0	0
All	All	36509	32951	32672	153	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 2.

The worst 5 of 153 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:F:166:ILE:HD11	1:F:197:GLN:HG3	1.60	0.83
1:D:166:ILE:HD11	1:D:197:GLN:HG3	1.73	0.70
1:A:159:LEU:HD11	1:A:236:LEU:HD22	1.76	0.67
1:B:365:ASP:OD2	5:B:501:HOH:O	2.14	0.65
1:N:166:ILE:HD11	1:N:197:GLN:HG3	1.79	0.64

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	371/388 (96%)	356 (96%)	15 (4%)	0	100	100
1	B	350/388 (90%)	337 (96%)	13 (4%)	0	100	100
1	C	366/388 (94%)	355 (97%)	11 (3%)	0	100	100
1	D	349/388 (90%)	337 (97%)	12 (3%)	0	100	100
1	E	367/388 (95%)	351 (96%)	16 (4%)	0	100	100
1	F	350/388 (90%)	342 (98%)	8 (2%)	0	100	100
1	G	366/388 (94%)	353 (96%)	13 (4%)	0	100	100
1	H	352/388 (91%)	340 (97%)	12 (3%)	0	100	100
1	K	368/388 (95%)	350 (95%)	18 (5%)	0	100	100
1	L	354/388 (91%)	348 (98%)	6 (2%)	0	100	100
1	M	367/388 (95%)	355 (97%)	12 (3%)	0	100	100
1	N	352/388 (91%)	338 (96%)	14 (4%)	0	100	100
All	All	4312/4656 (93%)	4162 (96%)	150 (4%)	0	100	100

There are no Ramachandran outliers to report.

### 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	304/317 (96%)	300 (99%)	4 (1%)	69	62
1	B	283/317 (89%)	274 (97%)	9 (3%)	39	25
1	C	296/317 (93%)	292 (99%)	4 (1%)	67	59
1	D	281/317 (89%)	275 (98%)	6 (2%)	53	42
1	E	295/317 (93%)	292 (99%)	3 (1%)	76	71
1	F	288/317 (91%)	281 (98%)	7 (2%)	49	36
1	G	298/317 (94%)	290 (97%)	8 (3%)	44	31
1	H	288/317 (91%)	286 (99%)	2 (1%)	84	81
1	K	292/317 (92%)	290 (99%)	2 (1%)	84	81

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	L	292/317 (92%)	289 (99%)	3 (1%)	76	71
1	M	298/317 (94%)	292 (98%)	6 (2%)	55	44
1	N	282/317 (89%)	278 (99%)	4 (1%)	67	59
All	All	3497/3804 (92%)	3439 (98%)	58 (2%)	60	51

5 of 58 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	F	92	ASP
1	N	235	ARG
1	G	218	ASP
1	N	125	ASP
1	M	128	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	78	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 monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

Of 37 ligands modelled in this entry, 25 are monoatomic - leaving 12 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
4	Y4F	G	404	-	43,50,50	0.75	1 (2%)	50,78,78	1.03	2 (4%)
4	Y4F	D	402	-	43,50,50	0.76	1 (2%)	50,78,78	0.99	3 (6%)
4	Y4F	L	402	-	43,50,50	0.72	1 (2%)	50,78,78	1.11	3 (6%)
4	Y4F	K	404	-	43,50,50	0.73	1 (2%)	50,78,78	1.03	2 (4%)
4	Y4F	B	402	-	43,50,50	0.77	1 (2%)	50,78,78	0.97	2 (4%)
4	Y4F	H	403	-	43,50,50	0.74	1 (2%)	50,78,78	1.04	3 (6%)
4	Y4F	A	404	-	43,50,50	0.77	1 (2%)	50,78,78	1.12	3 (6%)
4	Y4F	N	402	-	43,50,50	0.77	1 (2%)	50,78,78	1.02	3 (6%)
4	Y4F	F	402	-	43,50,50	0.78	1 (2%)	50,78,78	1.05	3 (6%)
4	Y4F	E	404	-	43,50,50	0.76	1 (2%)	50,78,78	1.05	3 (6%)
4	Y4F	C	404	-	43,50,50	0.74	1 (2%)	50,78,78	0.98	2 (4%)
4	Y4F	M	404	-	43,50,50	0.75	1 (2%)	50,78,78	0.98	2 (4%)

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
4	Y4F	G	404	-	-	5/22/62/62	0/6/7/7
4	Y4F	D	402	-	-	2/22/62/62	0/6/7/7
4	Y4F	L	402	-	-	4/22/62/62	0/6/7/7
4	Y4F	K	404	-	-	5/22/62/62	0/6/7/7
4	Y4F	B	402	-	-	3/22/62/62	0/6/7/7
4	Y4F	H	403	-	-	5/22/62/62	0/6/7/7
4	Y4F	A	404	-	-	4/22/62/62	0/6/7/7
4	Y4F	N	402	-	-	7/22/62/62	0/6/7/7
4	Y4F	F	402	-	-	4/22/62/62	0/6/7/7
4	Y4F	E	404	-	-	4/22/62/62	0/6/7/7
4	Y4F	C	404	-	-	5/22/62/62	0/6/7/7
4	Y4F	M	404	-	-	5/22/62/62	0/6/7/7

The worst 5 of 12 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	G	404	Y4F	C40-N42	3.06	1.37	1.32
4	M	404	Y4F	C40-N42	3.04	1.37	1.32
4	E	404	Y4F	C40-N42	3.03	1.37	1.32
4	F	402	Y4F	C40-N42	3.01	1.36	1.32
4	A	404	Y4F	C40-N42	2.99	1.36	1.32

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

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	G	404	Y4F	N42-C40-N39	-4.19	122.13	128.68
4	E	404	Y4F	N42-C40-N39	-4.15	122.19	128.68
4	L	402	Y4F	N42-C40-N39	-4.13	122.22	128.68
4	N	402	Y4F	N42-C40-N39	-4.12	122.24	128.68
4	H	403	Y4F	N42-C40-N39	-4.08	122.30	128.68

There are no chirality outliers.

5 of 53 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	404	Y4F	C3'-O28-P27-O30
4	B	402	Y4F	C3'-O28-P27-O29
4	B	402	Y4F	C3'-O28-P27-O30
4	C	404	Y4F	C3'-O28-P27-O30
4	E	404	Y4F	C3'-O28-P27-O30

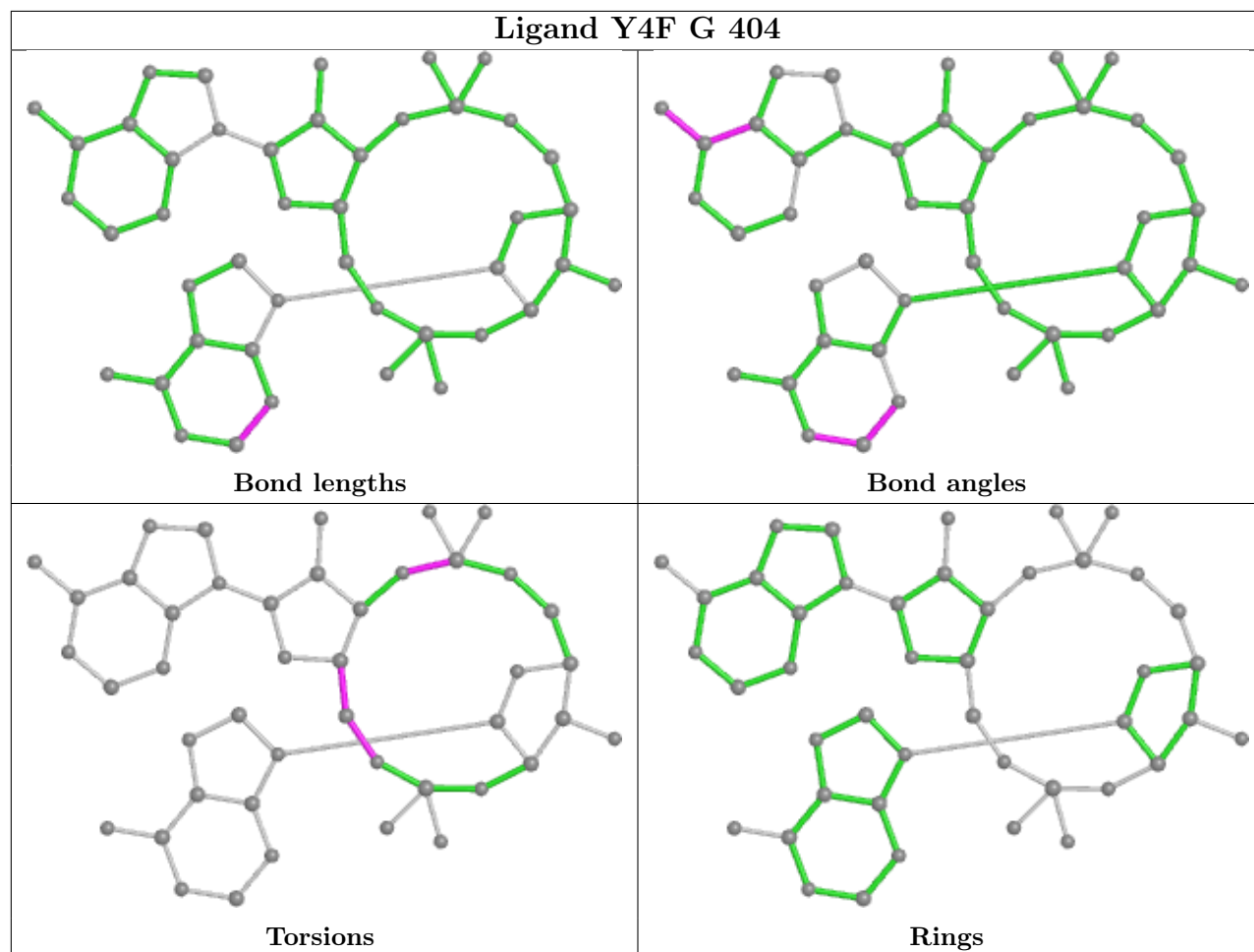
There are no ring outliers.

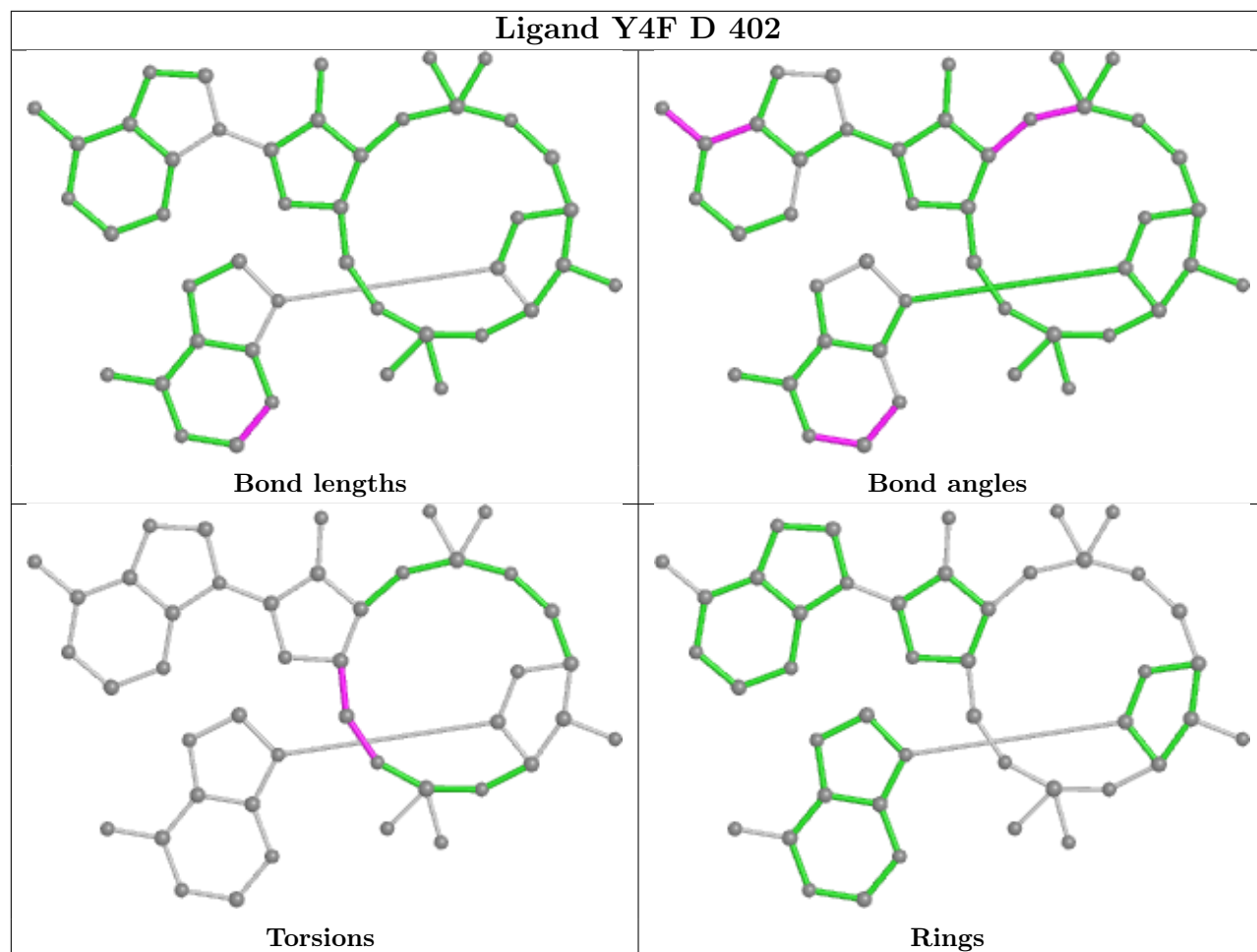
2 monomers are involved in 2 short contacts:

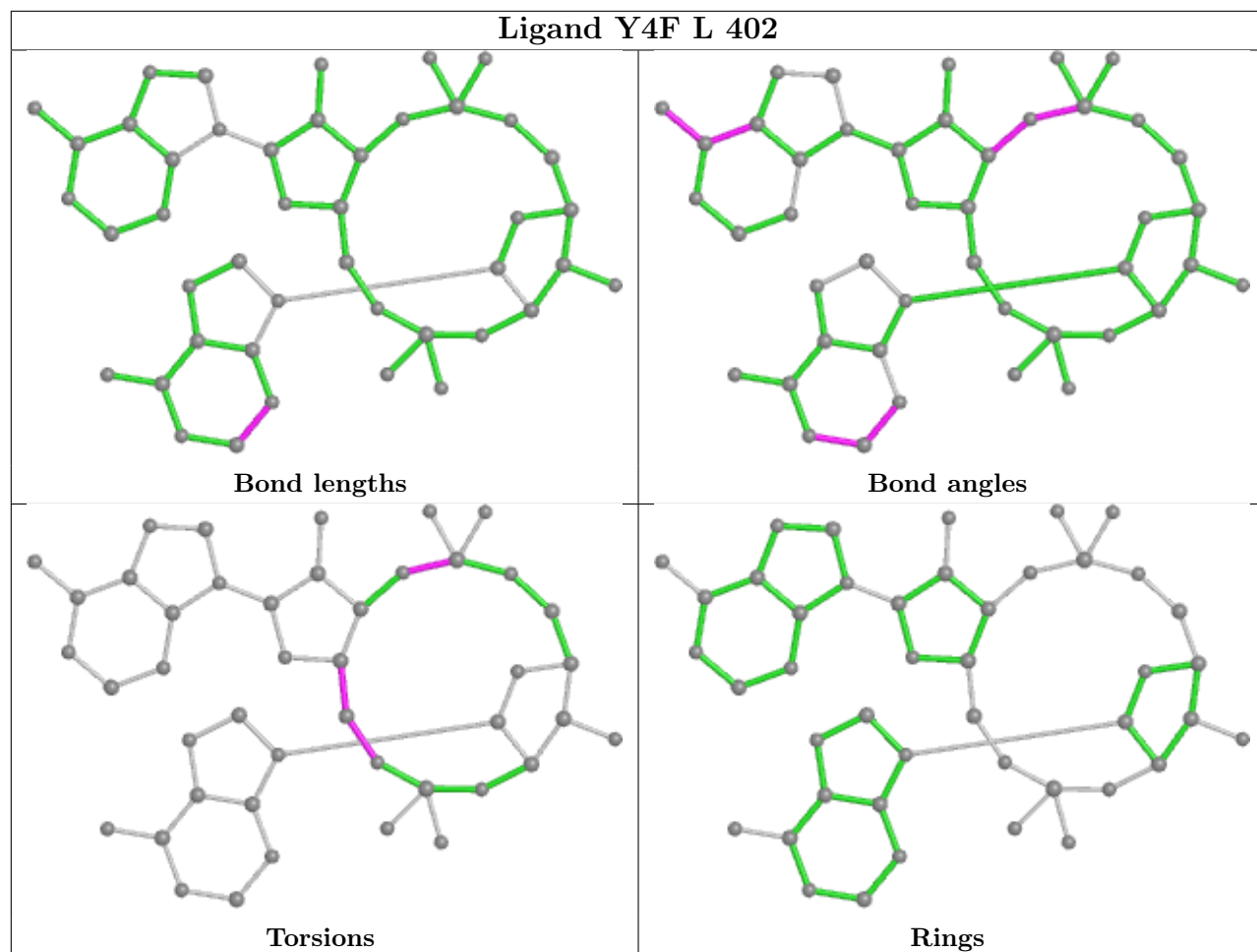
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	L	402	Y4F	1	0
4	N	402	Y4F	1	0

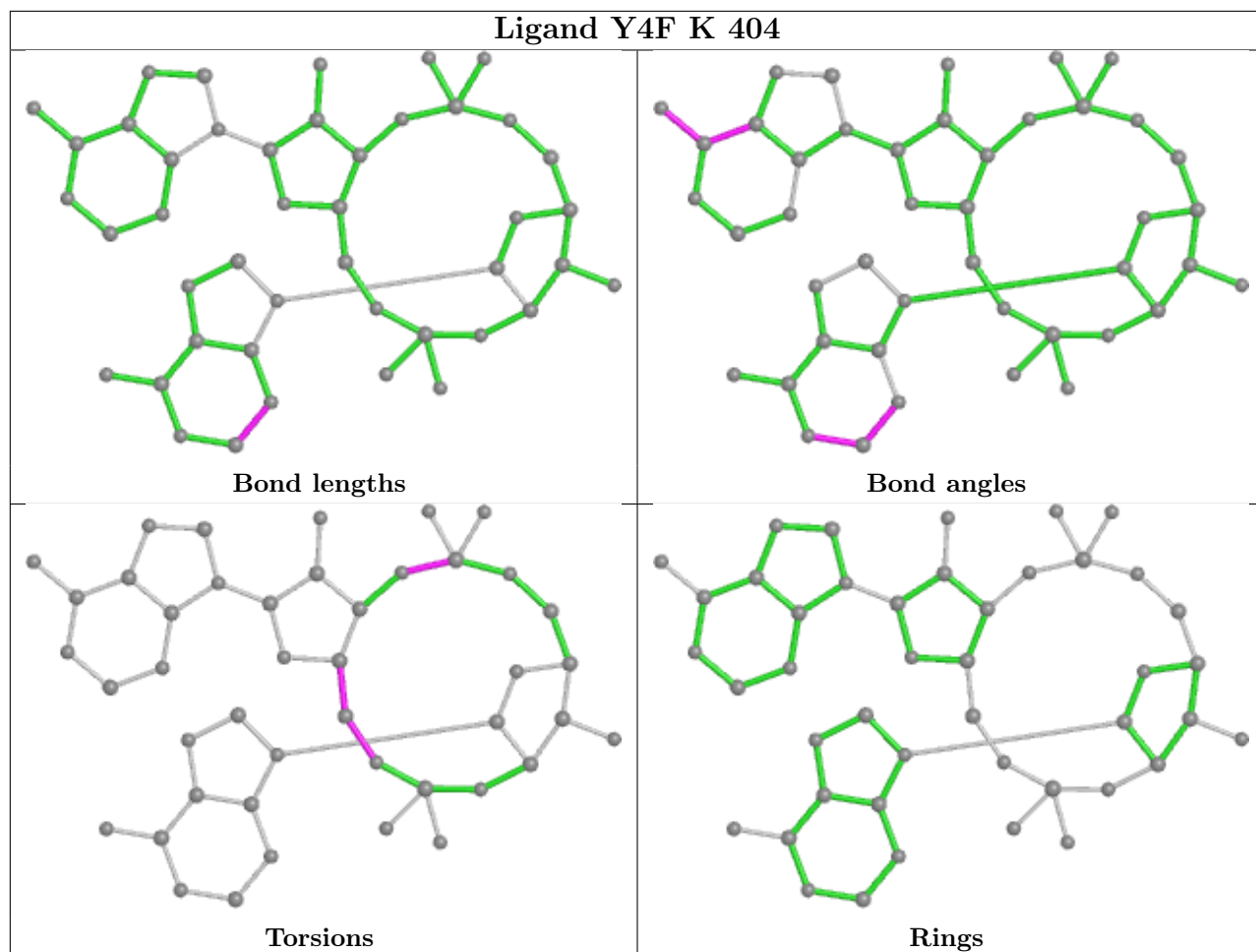
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

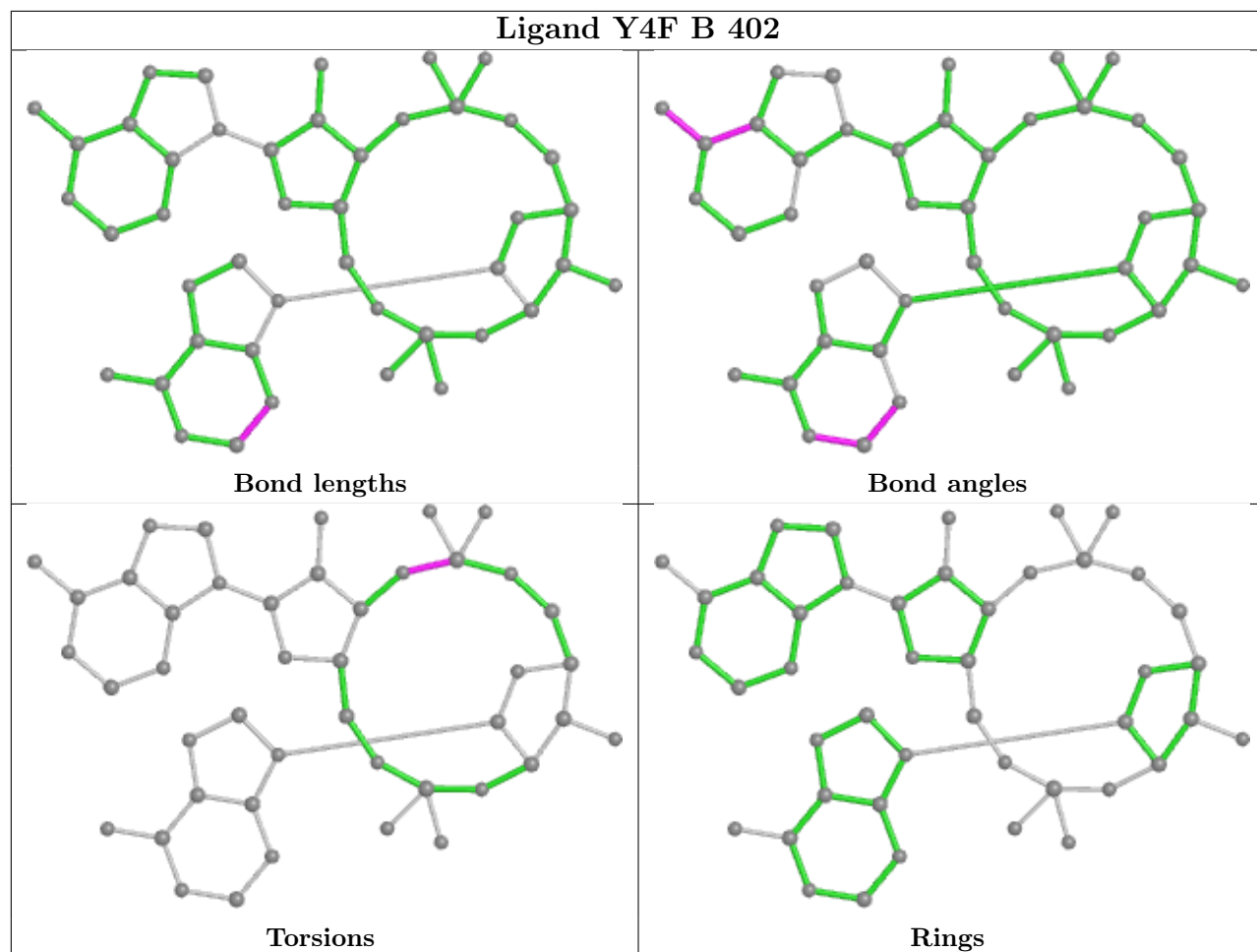
any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

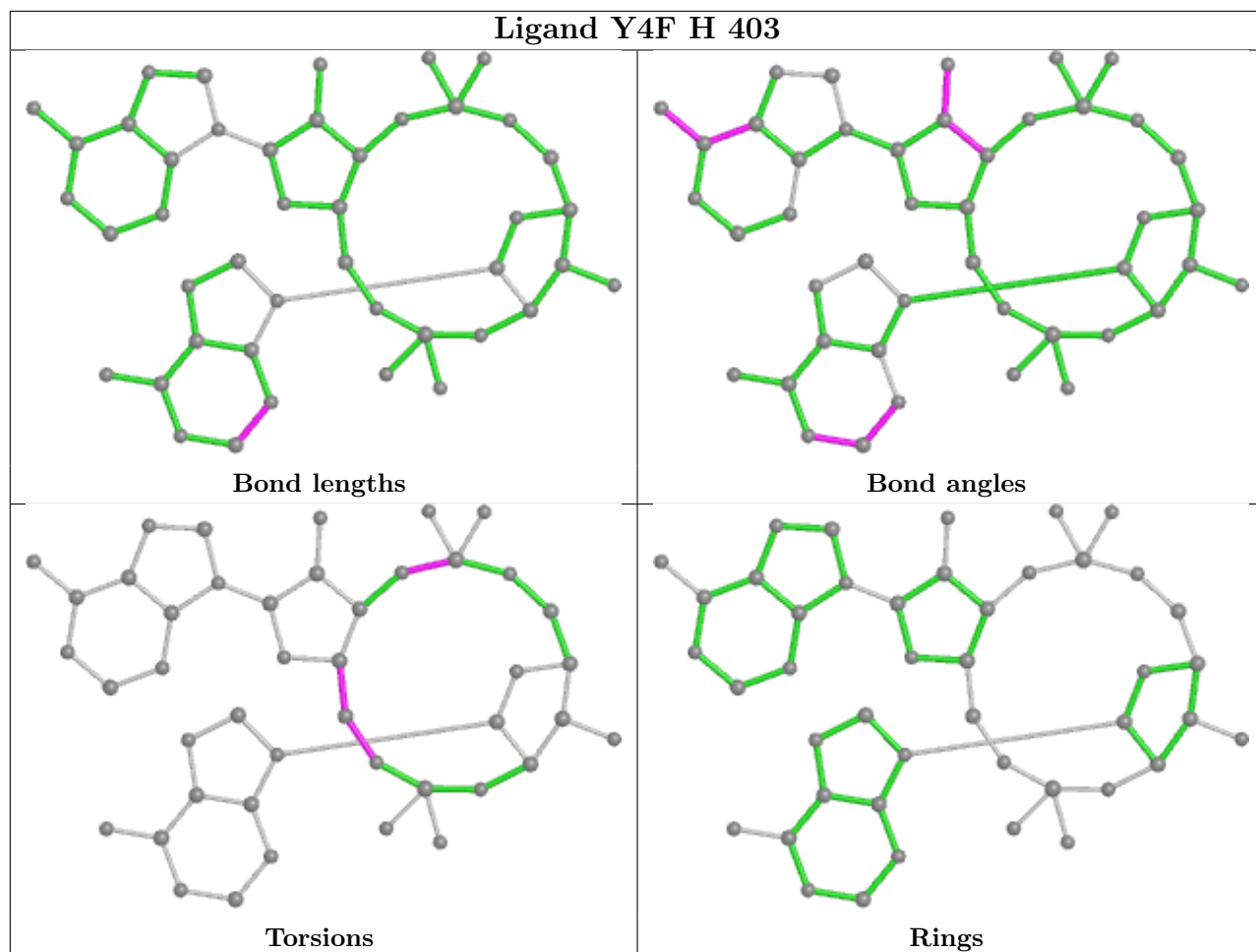




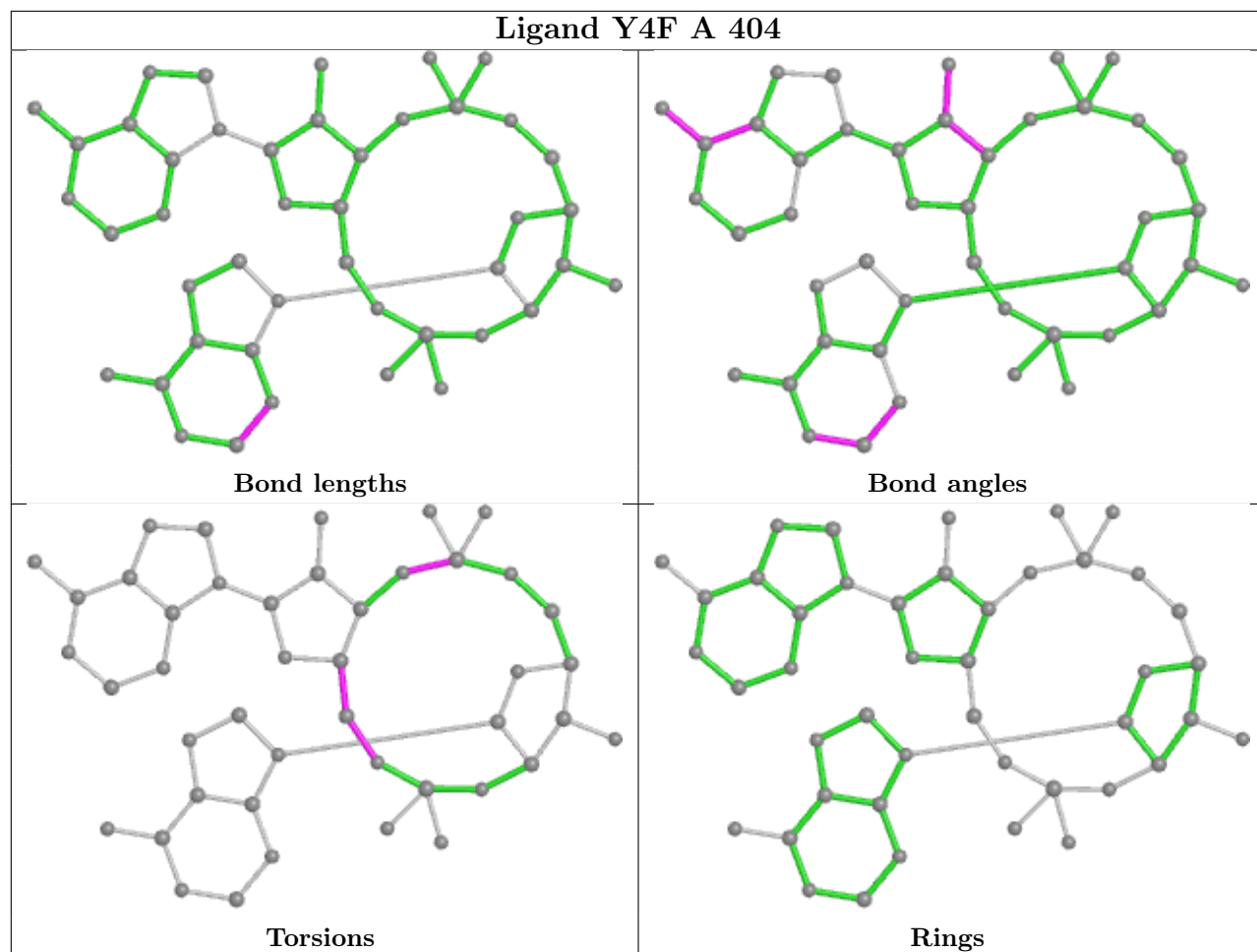


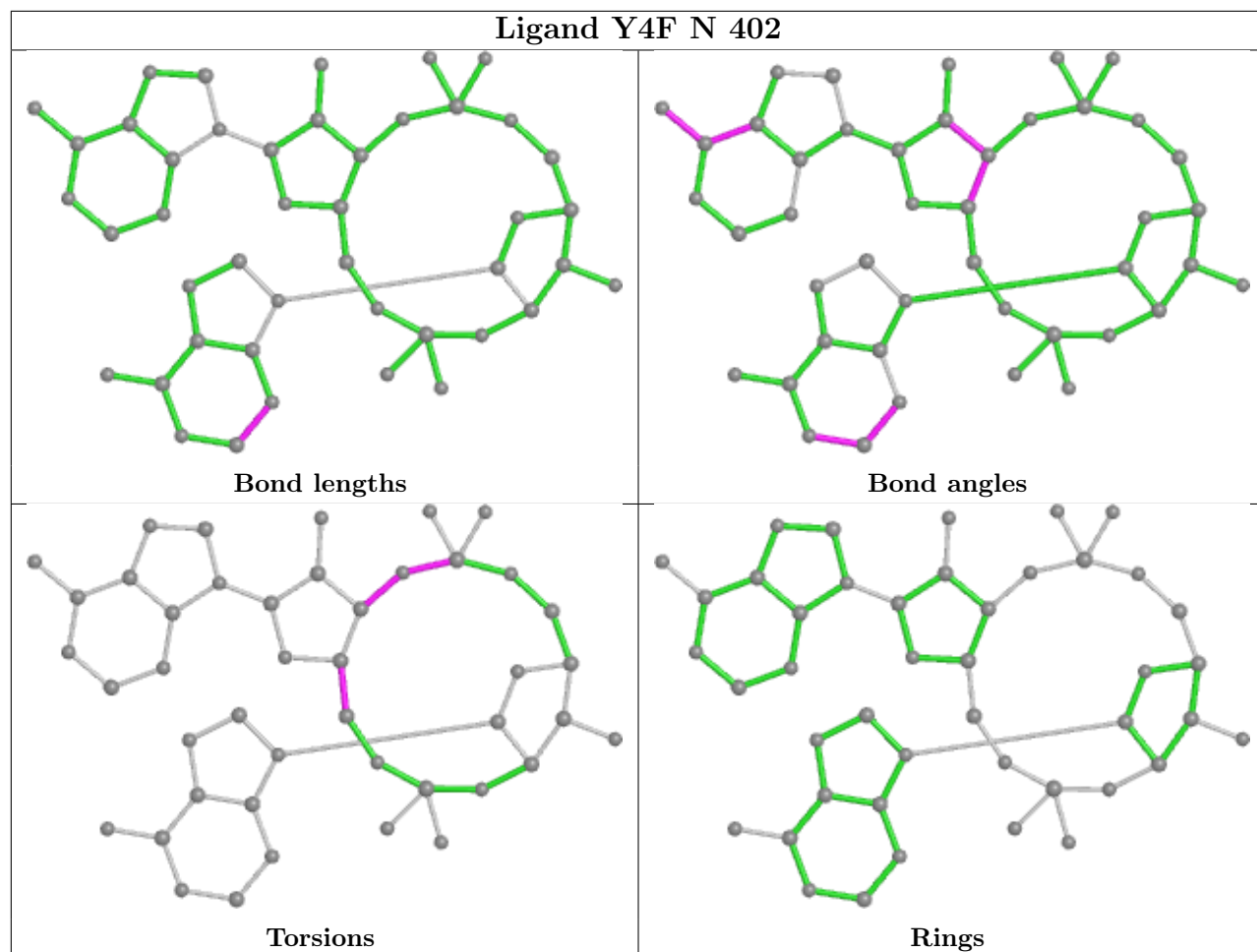


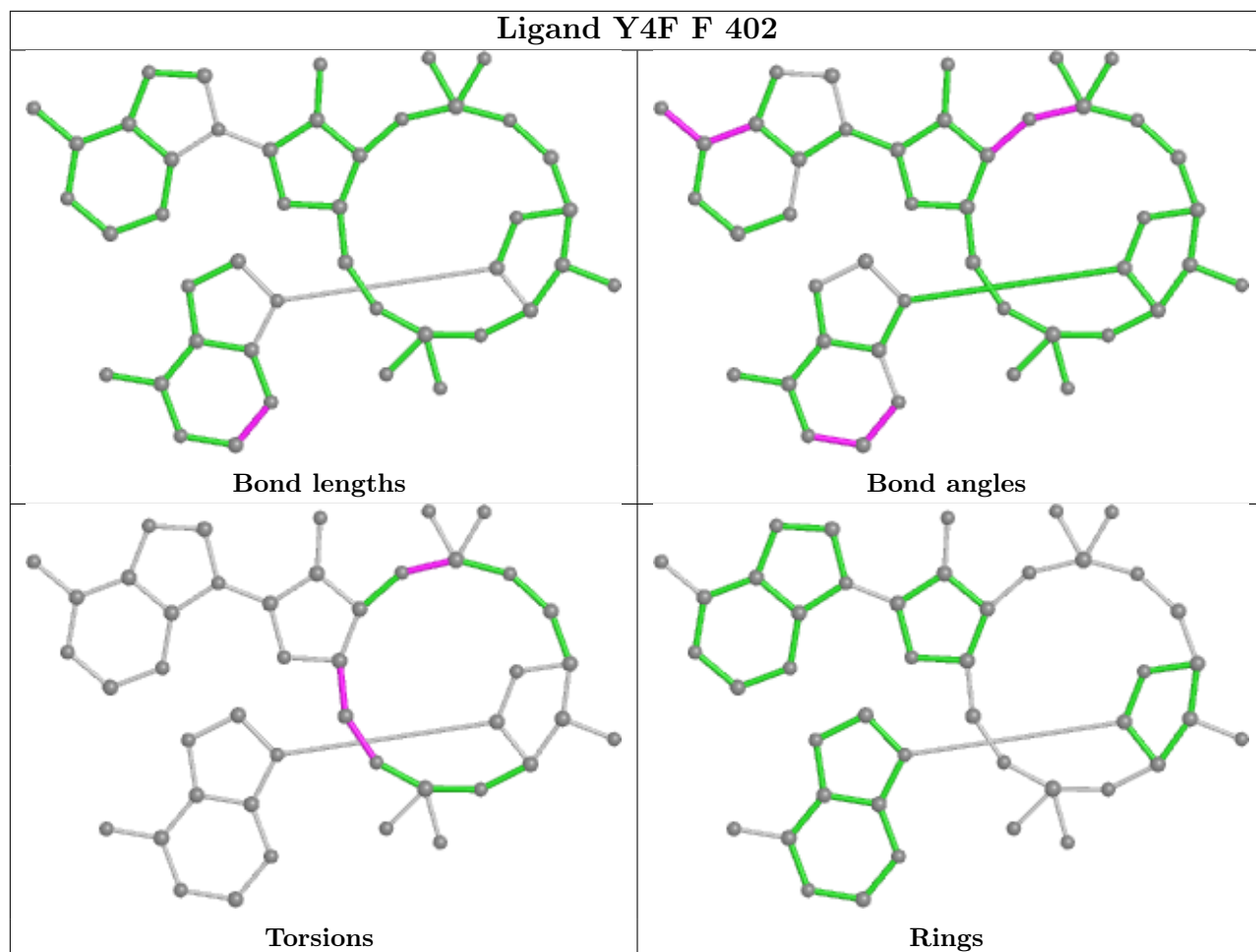


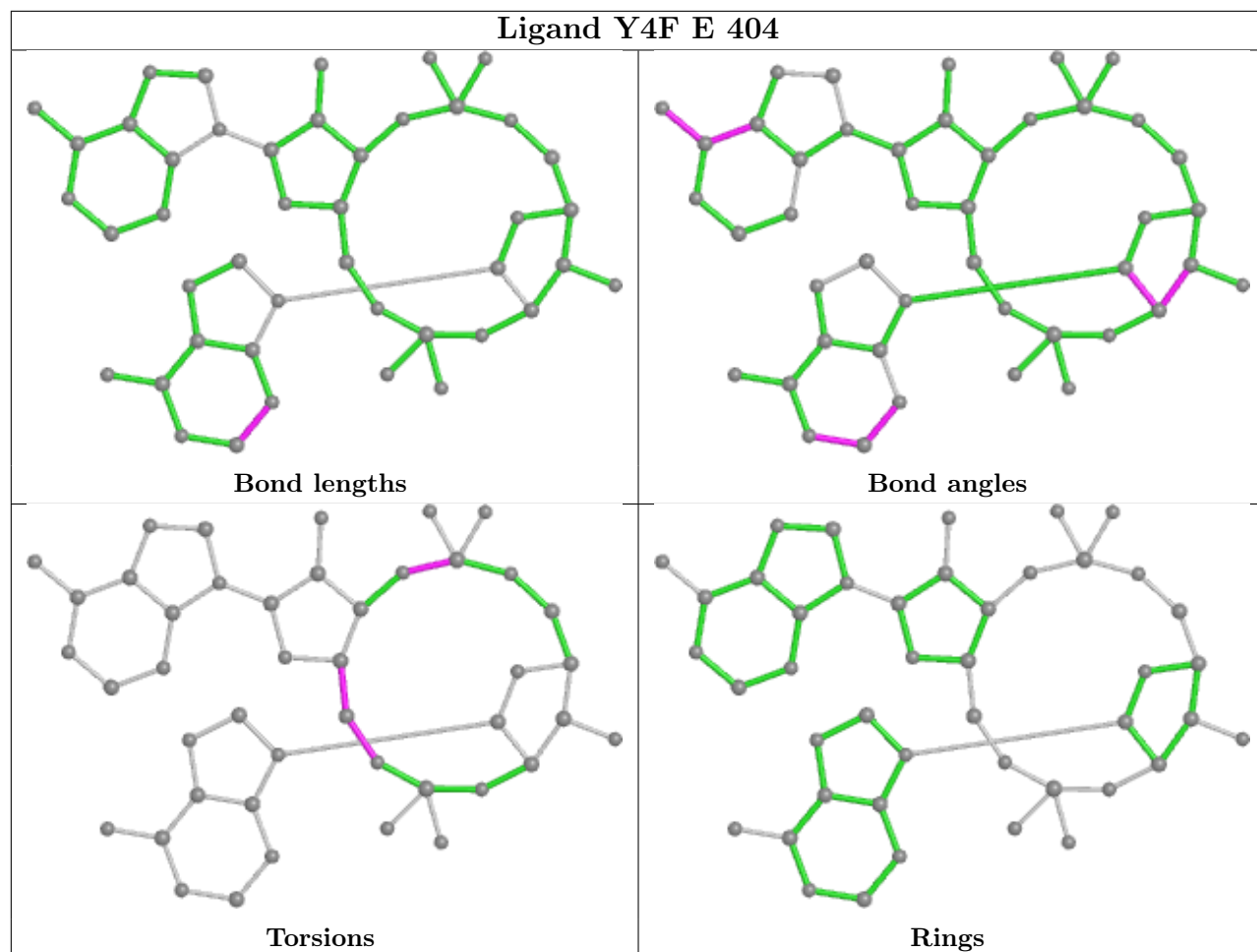


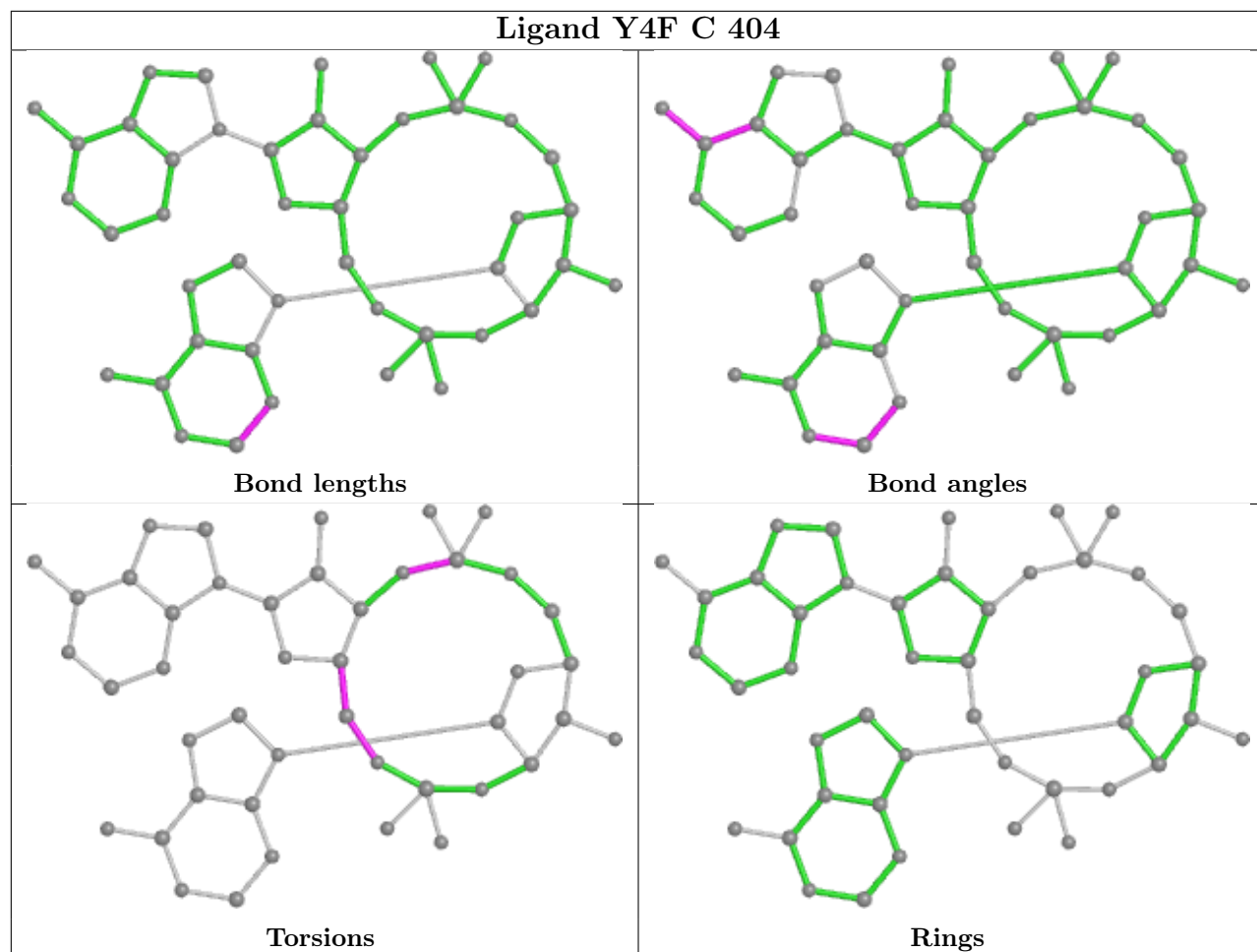


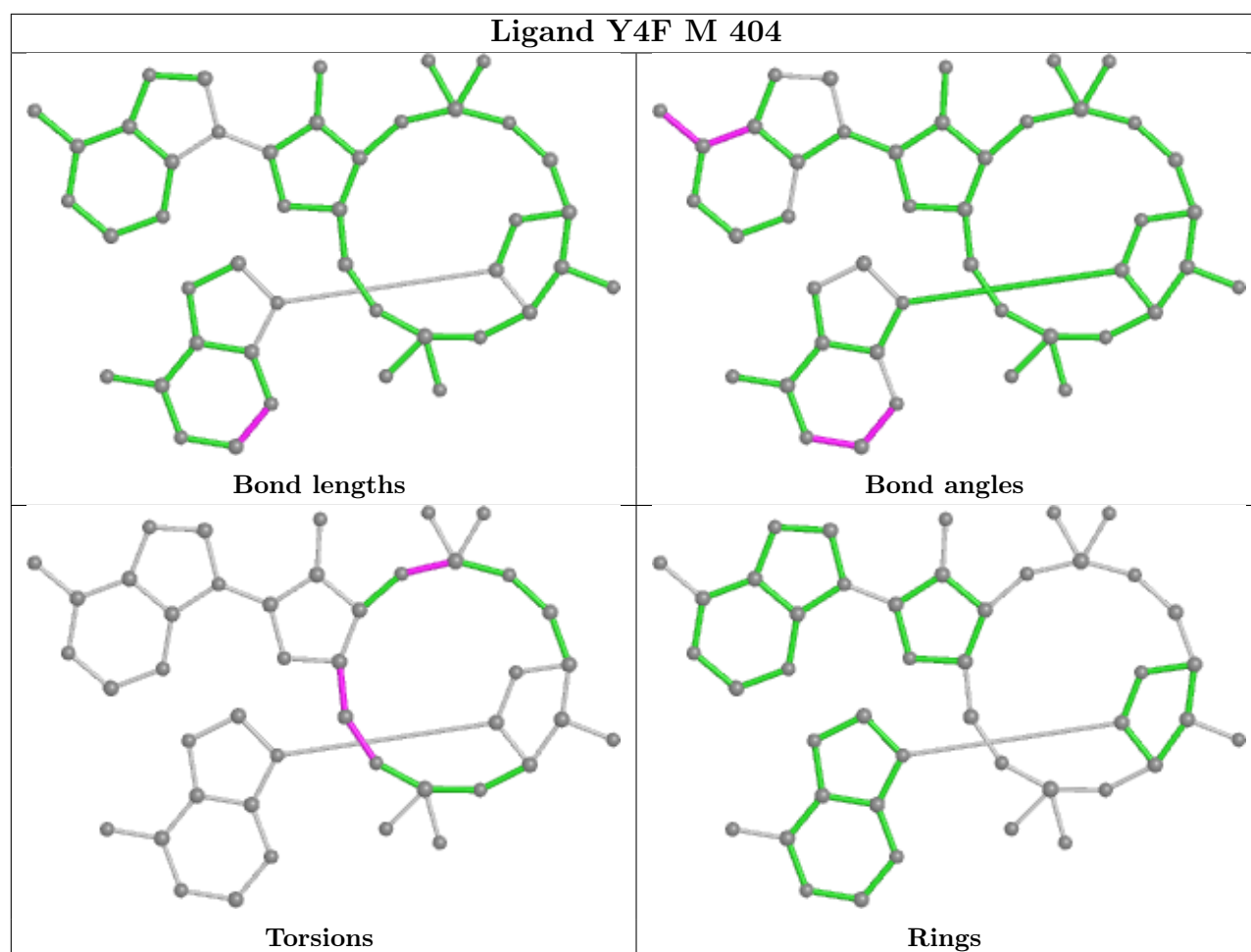












## 5.7 Other polymers [i](#)

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	369/388 (95%)	0.15	9 (2%) 59 54	16, 27, 60, 125	0
1	B	354/388 (91%)	0.26	17 (4%) 30 25	16, 28, 72, 98	0
1	C	368/388 (94%)	0.17	10 (2%) 54 49	18, 30, 66, 94	0
1	D	353/388 (90%)	0.26	9 (2%) 57 52	22, 35, 68, 94	0
1	E	369/388 (95%)	0.12	11 (2%) 50 44	19, 29, 60, 106	0
1	F	353/388 (90%)	0.20	10 (2%) 53 47	20, 32, 62, 93	0
1	G	368/388 (94%)	0.08	4 (1%) 80 78	18, 28, 56, 91	0
1	H	356/388 (91%)	0.15	13 (3%) 41 36	18, 28, 62, 89	0
1	K	369/388 (95%)	0.09	6 (1%) 72 68	19, 29, 64, 92	0
1	L	357/388 (92%)	0.05	5 (1%) 75 72	17, 27, 50, 76	0
1	M	369/388 (95%)	0.18	11 (2%) 50 44	20, 31, 64, 97	0
1	N	354/388 (91%)	0.23	16 (4%) 33 27	21, 33, 70, 96	0
All	All	4339/4656 (93%)	0.16	121 (2%) 53 47	16, 30, 64, 125	0

The worst 5 of 121 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	94	ILE	7.8
1	D	14	PHE	6.8
1	A	207	ASP	6.8
1	H	142	ILE	5.6
1	A	208	SER	5.5

### 6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

### 6.3 Carbohydrates [i](#)

There are no monosaccharides 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
3	MG	C	402	1/1	0.90	0.06	38,38,38,38	0
3	MG	E	403	1/1	0.90	0.13	66,66,66,66	0
3	MG	E	402	1/1	0.91	0.08	23,23,23,23	0
3	MG	M	402	1/1	0.91	0.07	30,30,30,30	0
4	Y4F	L	402	44/44	0.92	0.15	35,56,70,77	0
4	Y4F	F	402	44/44	0.93	0.15	35,49,73,79	0
3	MG	G	402	1/1	0.94	0.06	27,27,27,27	0
4	Y4F	D	402	44/44	0.94	0.12	31,45,62,69	0
3	MG	M	403	1/1	0.95	0.07	45,45,45,45	0
3	MG	K	402	1/1	0.96	0.05	28,28,28,28	0
3	MG	A	402	1/1	0.96	0.05	28,28,28,28	0
4	Y4F	H	403	44/44	0.96	0.08	28,36,45,49	0
3	MG	H	402	1/1	0.96	0.04	33,33,33,33	0
4	Y4F	N	402	44/44	0.96	0.12	24,35,52,63	0
4	Y4F	B	402	44/44	0.97	0.10	20,30,49,53	0
3	MG	C	403	1/1	0.97	0.03	44,44,44,44	0
3	MG	G	403	1/1	0.97	0.10	43,43,43,43	0
4	Y4F	E	404	44/44	0.98	0.09	14,21,30,36	0
4	Y4F	A	404	44/44	0.98	0.10	11,17,21,26	0
4	Y4F	G	404	44/44	0.98	0.10	13,20,26,27	0
3	MG	A	403	1/1	0.98	0.18	44,44,44,44	0
4	Y4F	K	404	44/44	0.98	0.09	14,20,25,29	0
4	Y4F	C	404	44/44	0.98	0.09	15,22,31,37	0
3	MG	K	403	1/1	0.98	0.16	43,43,43,43	0
2	ZN	G	401	1/1	0.99	0.19	29,29,29,29	0
2	ZN	H	401	1/1	0.99	0.11	30,30,30,30	0
2	ZN	L	401	1/1	0.99	0.11	32,32,32,32	0
2	ZN	N	401	1/1	0.99	0.12	35,35,35,35	0
2	ZN	B	401	1/1	0.99	0.11	42,42,42,42	0
2	ZN	D	401	1/1	0.99	0.12	35,35,35,35	0
4	Y4F	M	404	44/44	0.99	0.09	17,23,31,34	0
2	ZN	F	401	1/1	0.99	0.16	33,33,33,33	0

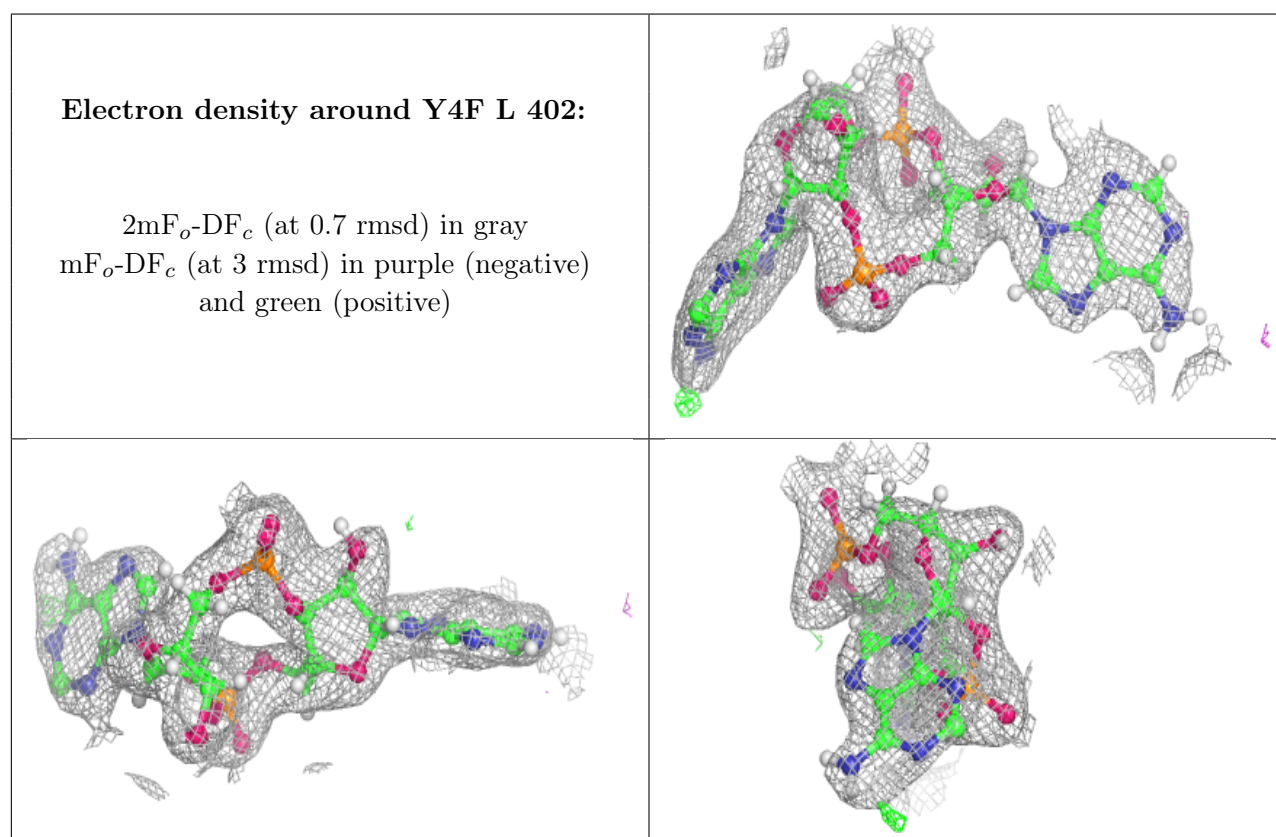
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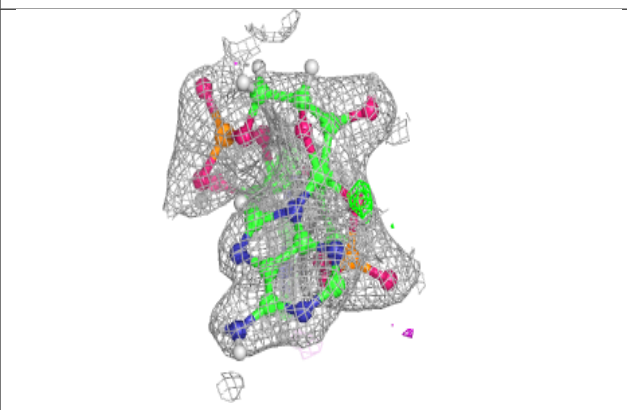
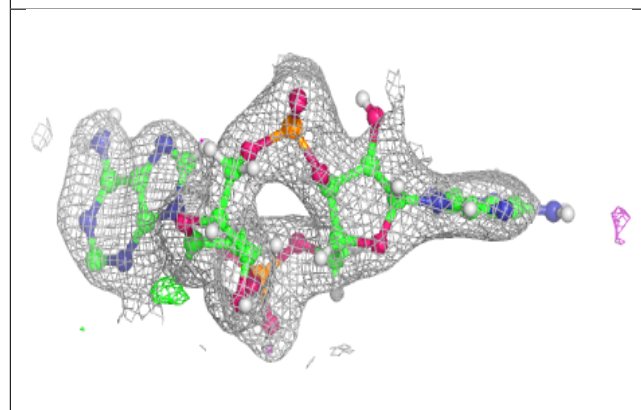
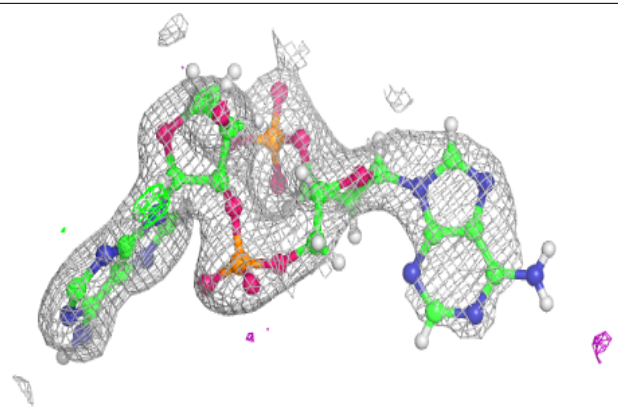
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	ZN	E	401	1/1	1.00	0.13	24,24,24,24	0
2	ZN	K	401	1/1	1.00	0.10	20,20,20,20	0
2	ZN	C	401	1/1	1.00	0.14	26,26,26,26	0
2	ZN	M	401	1/1	1.00	0.12	23,23,23,23	0
2	ZN	A	401	1/1	1.00	0.14	26,26,26,26	0

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.

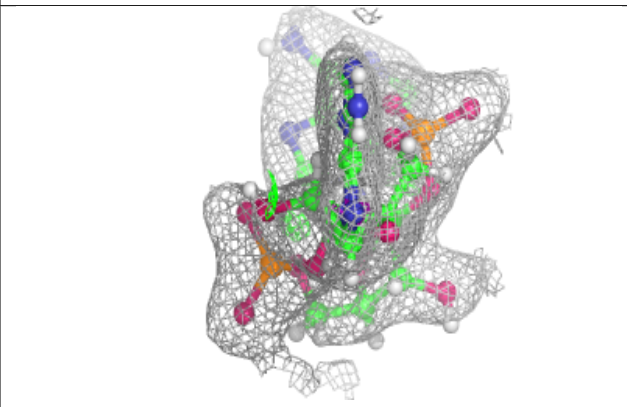
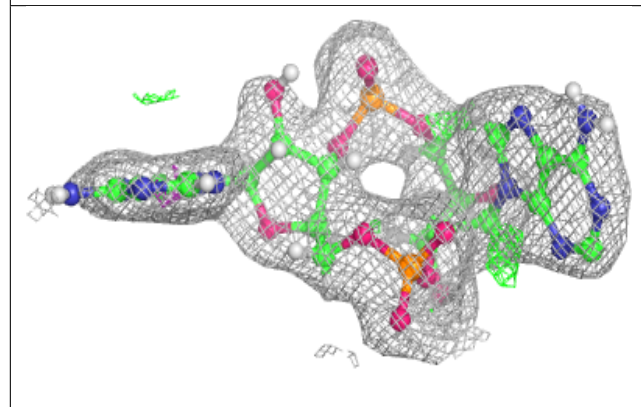
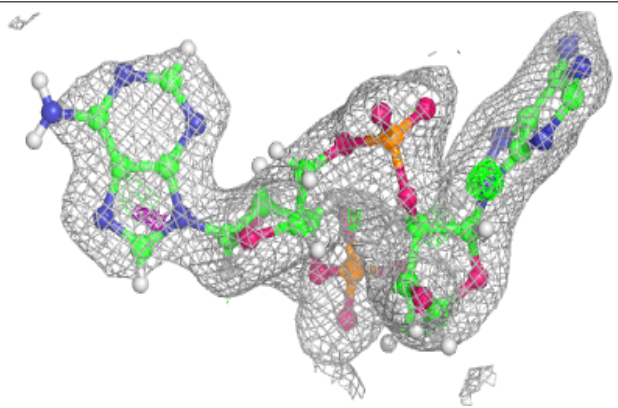


**Electron density around Y4F F 402:**

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

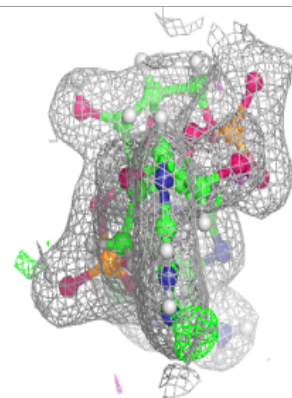
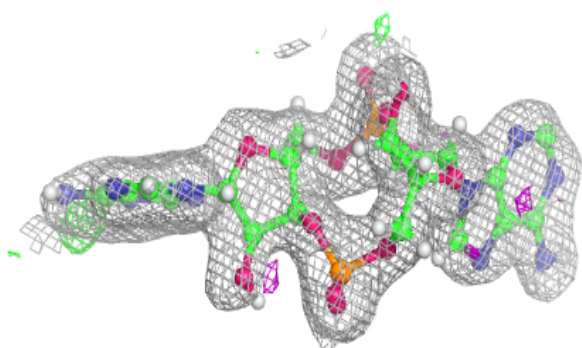
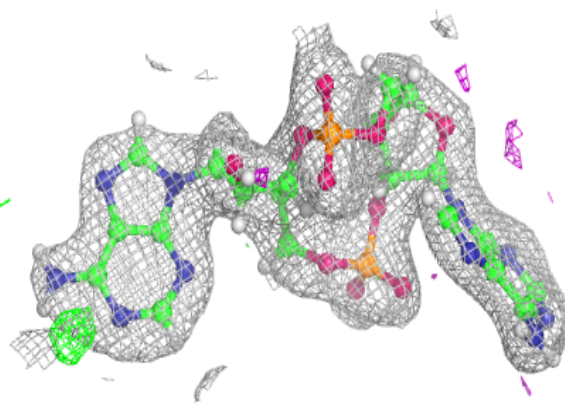
**Electron density around Y4F D 402:**

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

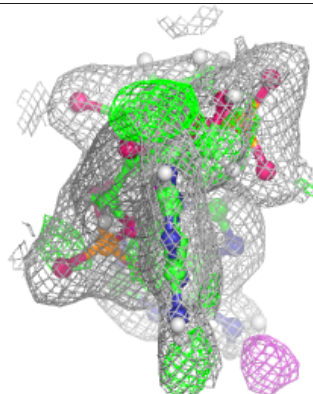
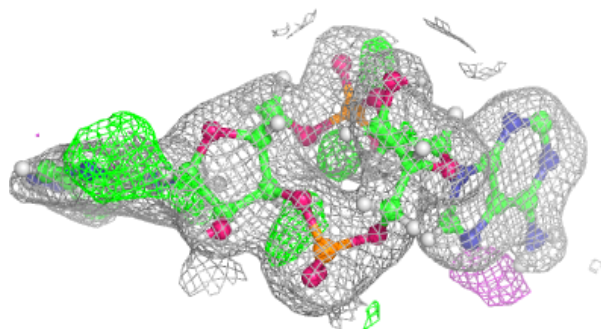
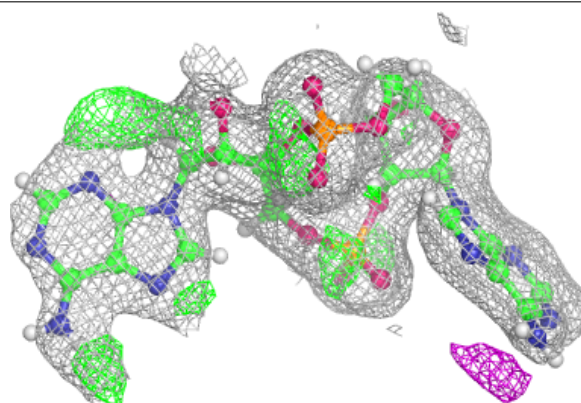


**Electron density around Y4F H 403:**

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

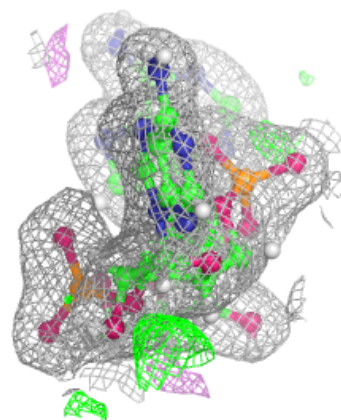
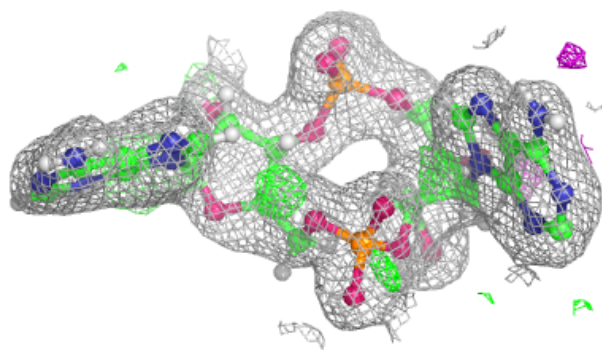
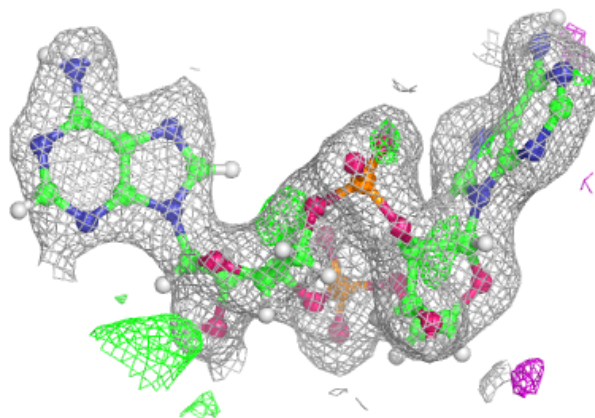
**Electron density around Y4F N 402:**

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



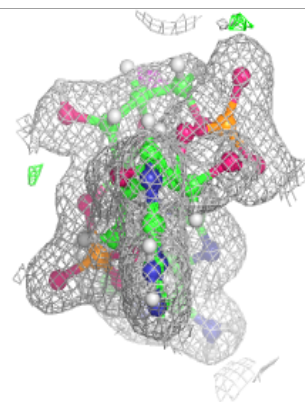
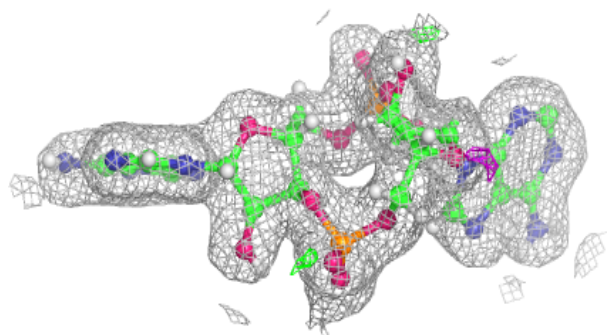
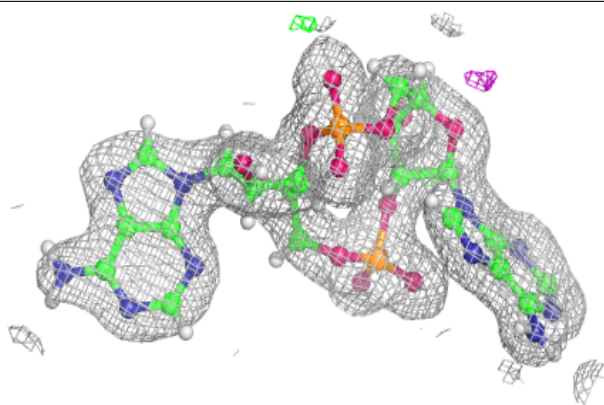
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and green (positive)

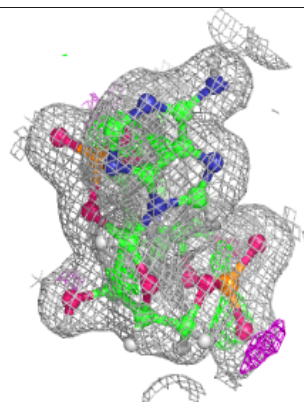
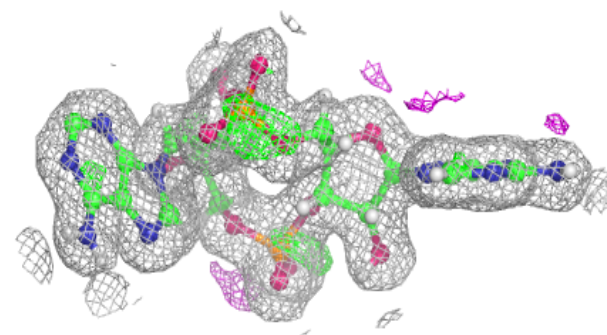
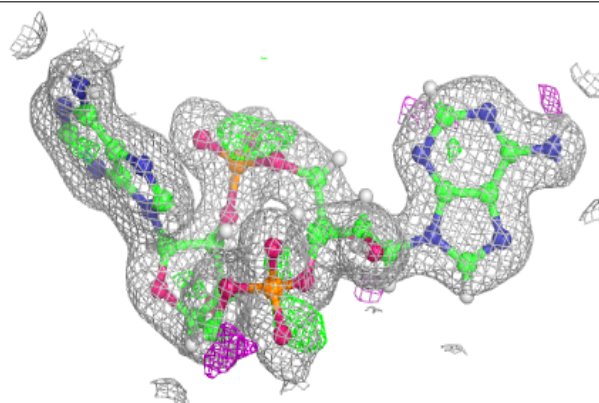


**Electron density around Y4F E 404:**

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

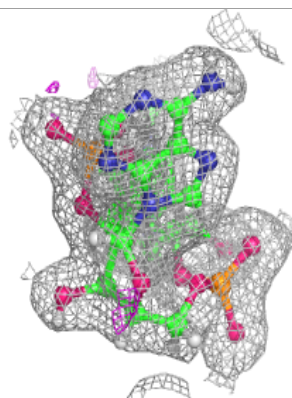
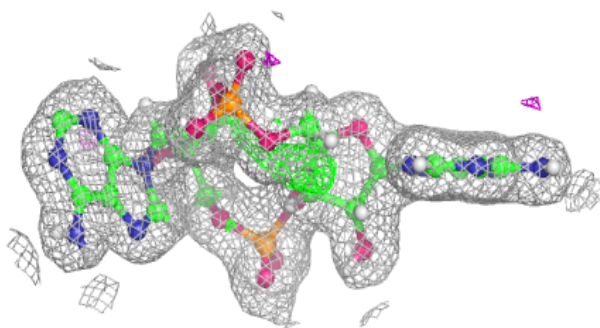
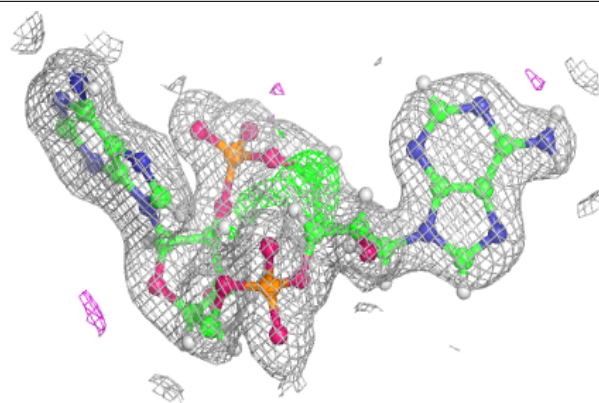
**Electron density around Y4F A 404:**

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

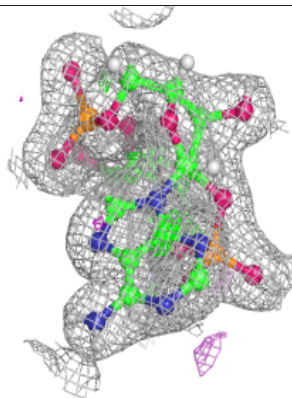
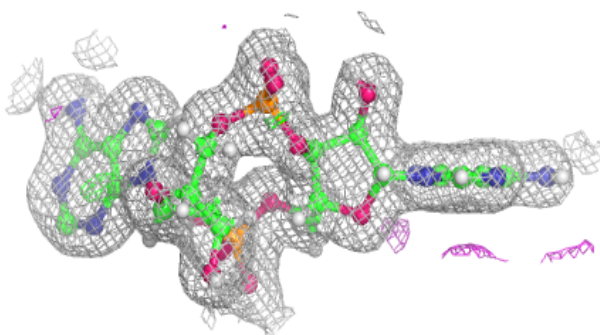
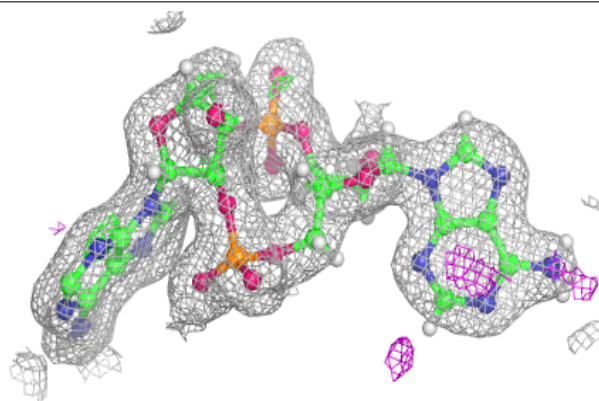


**Electron density around Y4F G 404:**

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

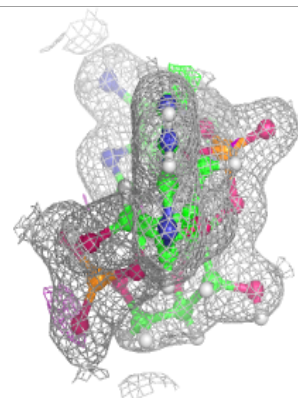
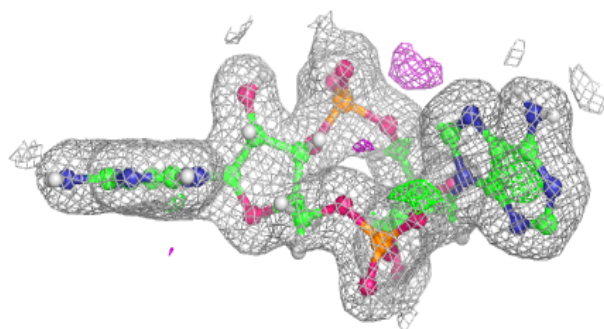
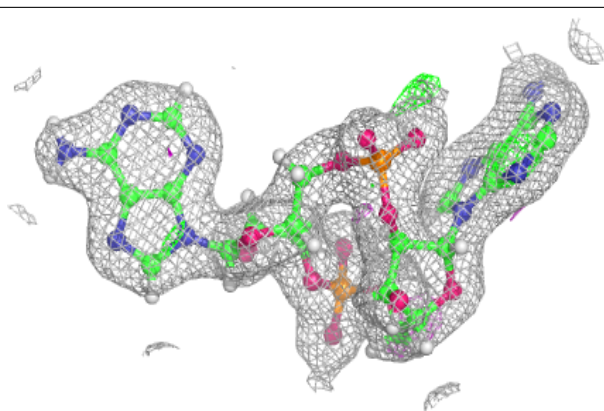
**Electron density around Y4F K 404:**

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

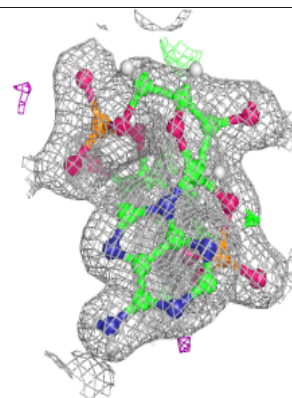
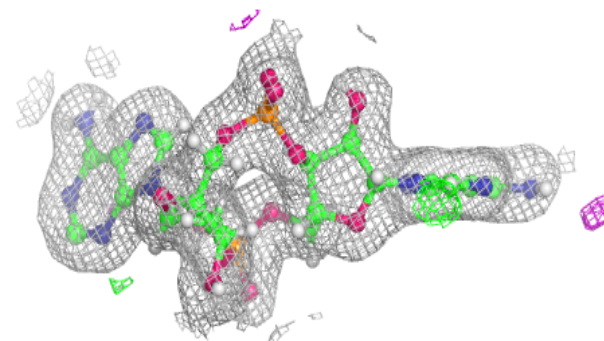
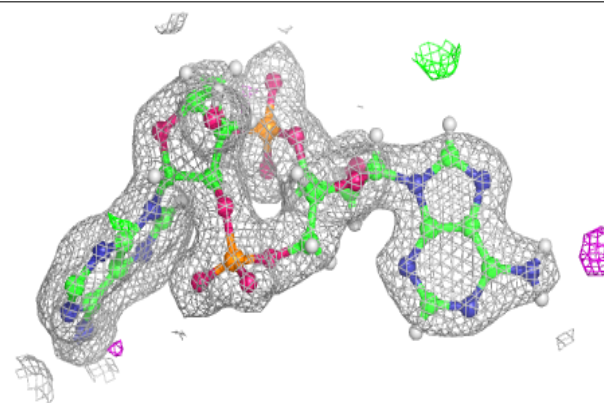


**Electron density around Y4F C 404:**

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

**Electron density around Y4F M 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 [i](#)

There are no such residues in this entry.