



wwPDB X-ray Structure Validation Summary Report ⓘ

Jun 17, 2024 – 04:27 PM EDT

PDB ID : 2YCE
Title : Structure of an Archaeal fructose-1,6-bisphosphate aldolase with the catalytic
Lys covalently bound to the carbinolamine intermediate of the substrate.
Authors : Lorentzen, E.; Siebers, B.; Hensel, R.; Pohl, E.
Deposited on : 2011-03-14
Resolution : 1.93 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

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	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	2.37.1
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)	:	2.37.1

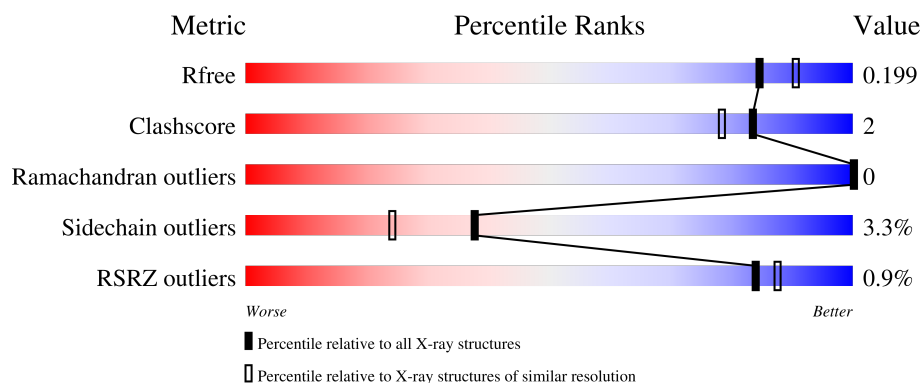
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION





The reported resolution of this entry is 1.93 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	4310 (1.96-1.92)
Clashscore	141614	1023 (1.94-1.94)
Ramachandran outliers	138981	1007 (1.94-1.94)
Sidechain outliers	138945	1007 (1.94-1.94)
RSRZ outliers	127900	4250 (1.96-1.92)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	263	
1	B	263	
1	C	263	
1	D	263	
1	E	263	

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Mol	Chain	Length	Quality of chain
1	F	263	<div><div><div></div><div></div><div></div></div><div>%</div><div><div></div><div></div><div></div></div><div>88%</div><div>7%</div><div>5%</div></div>
1	G	263	<div><div><div></div><div></div><div></div></div><div>%</div><div><div></div><div></div><div></div></div><div>87%</div><div>7%</div><div>5%</div></div>
1	H	263	<div><div><div></div><div></div><div></div></div><div>2%</div><div><div></div><div></div><div></div></div><div>84%</div><div>10%</div><div>5%</div></div>
1	I	263	<div><div><div></div><div></div><div></div></div><div></div><div><div></div><div></div><div></div></div><div>88%</div><div>7%</div><div>5%</div></div>
1	J	263	<div><div><div></div><div></div><div></div></div><div>%</div><div><div></div><div></div><div></div></div><div>86%</div><div>9%</div><div>5%</div></div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 21227 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 FRUCTOSE-BISPHOSPHATE ALDOLASE CLASS 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	250	Total	C	N	O	S	0	5	0
			1962	1264	334	358	6			
1	B	251	Total	C	N	O	S	0	2	0
			1950	1256	332	357	5			
1	C	250	Total	C	N	O	S	0	4	0
			1953	1261	331	356	5			
1	D	250	Total	C	N	O	S	0	1	0
			1943	1252	331	355	5			
1	E	255	Total	C	N	O	S	0	4	0
			1983	1278	340	359	6			
1	F	250	Total	C	N	O	S	0	3	0
			1951	1256	334	355	6			
1	G	250	Total	C	N	O	S	0	1	0
			1938	1249	328	355	6			
1	H	250	Total	C	N	O	S	0	2	0
			1945	1253	331	356	5			
1	I	250	Total	C	N	O	S	0	1	0
			1937	1248	328	356	5			
1	J	250	Total	C	N	O	S	0	2	0
			1947	1255	331	355	6			

There are 20 discrepancies between the modelled and reference sequences:

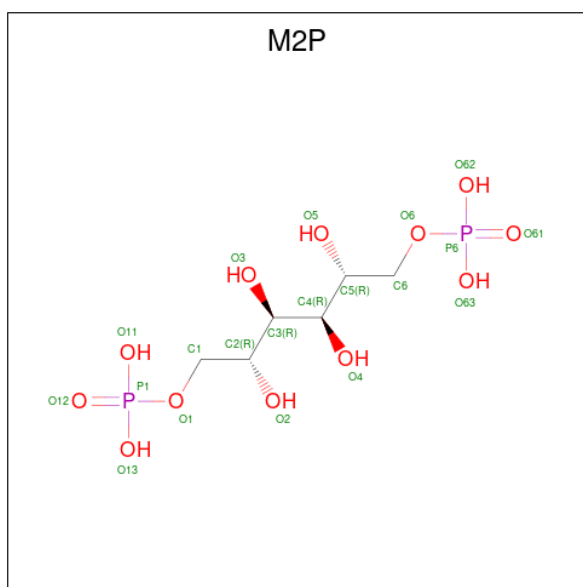
Chain	Residue	Modelled	Actual	Comment	Reference
A	146	PHE	TYR	engineered mutation	UNP P58315
A	173	SER	ALA	engineered mutation	UNP P58315
B	146	PHE	TYR	engineered mutation	UNP P58315
B	173	SER	ALA	engineered mutation	UNP P58315
C	146	PHE	TYR	engineered mutation	UNP P58315
C	173	SER	ALA	engineered mutation	UNP P58315
D	146	PHE	TYR	engineered mutation	UNP P58315
D	173	SER	ALA	engineered mutation	UNP P58315
E	146	PHE	TYR	engineered mutation	UNP P58315

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Chain	Residue	Modelled	Actual	Comment	Reference
E	173	SER	ALA	engineered mutation	UNP P58315
F	146	PHE	TYR	engineered mutation	UNP P58315
F	173	SER	ALA	engineered mutation	UNP P58315
G	146	PHE	TYR	engineered mutation	UNP P58315
G	173	SER	ALA	engineered mutation	UNP P58315
H	146	PHE	TYR	engineered mutation	UNP P58315
H	173	SER	ALA	engineered mutation	UNP P58315
I	146	PHE	TYR	engineered mutation	UNP P58315
I	173	SER	ALA	engineered mutation	UNP P58315
J	146	PHE	TYR	engineered mutation	UNP P58315
J	173	SER	ALA	engineered mutation	UNP P58315

- Molecule 2 is D-MANNITOL-1,6-DIPHOSPHATE (three-letter code: M2P) (formula: $C_6H_{16}O_{12}P_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	O	P	0	0
			20	6	12	2		
2	B	1	Total	C	O	P	0	0
			20	6	12	2		
2	C	1	Total	C	O	P	0	0
			20	6	12	2		
2	D	1	Total	C	O	P	0	0
			20	6	12	2		
2	E	1	Total	C	O	P	0	0
			20	6	12	2		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	F	1	Total	C	O	P	0	0
			20	6	12	2		
2	G	1	Total	C	O	P	0	0
			20	6	12	2		
2	H	1	Total	C	O	P	0	0
			20	6	12	2		
2	I	1	Total	C	O	P	0	0
			20	6	12	2		
2	J	1	Total	C	O	P	0	0
			20	6	12	2		


- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	158	Total	O	0	0
			158	158		
3	B	173	Total	O	0	0
			173	173		
3	C	136	Total	O	0	0
			136	136		
3	D	126	Total	O	0	0
			126	126		
3	E	173	Total	O	0	0
			173	173		
3	F	122	Total	O	0	0
			122	122		
3	G	148	Total	O	0	0
			148	148		
3	H	150	Total	O	0	0
			150	150		
3	I	158	Total	O	0	0
			158	158		
3	J	174	Total	O	0	0
			174	174		

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: FRUCTOSE-BISPHOSPHATE ALDOLASE CLASS 1

Chain A: 




• Molecule 1: FRUCTOSE-BISPHOSPHATE ALDOLASE CLASS 1

Chain B: 




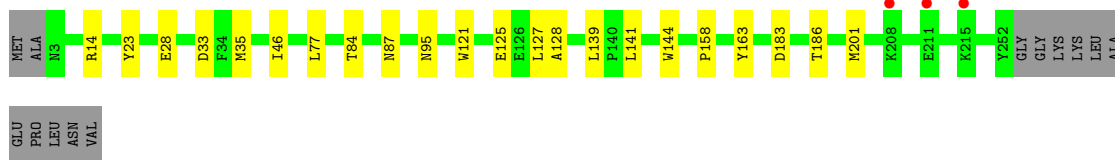
• Molecule 1: FRUCTOSE-BISPHOSPHATE ALDOLASE CLASS 1

Chain C: 




• Molecule 1: FRUCTOSE-BISPHOSPHATE ALDOLASE CLASS 1

Chain D: 



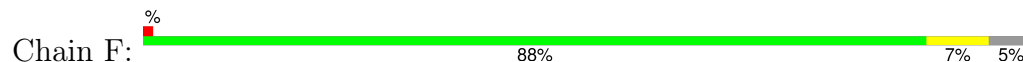
• Molecule 1: FRUCTOSE-BISPHOSPHATE ALDOLASE CLASS 1

Chain E: 



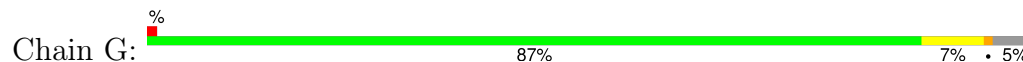
ALA
GLU
PRO
LEU
ASN
VAL

• Molecule 1: FRUCTOSE-BISPHOSPHATE ALDOLASE CLASS 1



MET ALA N3 K7 Y23 E28 P31 I46 R51 L77 T84 N95 T112 A128 L139 W144 P158 Y163 K177 D183 T186 M201 A229 E249 L250 V251 Y252 GLY GLY LYS LYS LEU LEU ALA GLU PRO LEU ASN VAL

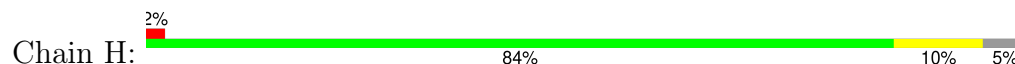
• Molecule 1: FRUCTOSE-BISPHOSPHATE ALDOLASE CLASS 1



MET ALA N3 R14 E28 D33 E44 L77 T84 N95 Y98 A128 D133 L139 W144 P158 Y163 K177 T186 K192 V193 M201 K208 E218 E222 R238 Y252 GLY GLY LYS LYS LEU LEU ALA GLU PRO LEU ASN

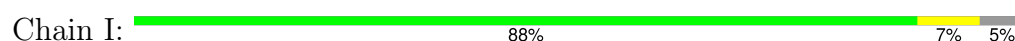
VAL

• Molecule 1: FRUCTOSE-BISPHOSPHATE ALDOLASE CLASS 1



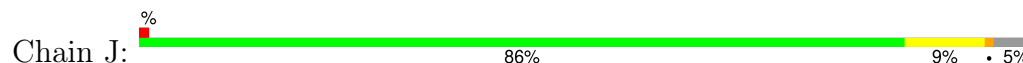
MET ALA N3 R14 Y23 E28 P31 M35 I46 L47 R48 R51 S72 L77 T84 N95 G110 W121 E125 E126 L127 A128 L139 P140 L141 V142 V143 W144 P158 Y163 K177 G182 K185 T186 M201 K208 E211 R215 Q236 R237 R238 Y252 GLY GLY LYS LYS LEU ALA PRO LEU ASN VAL

• Molecule 1: FRUCTOSE-BISPHOSPHATE ALDOLASE CLASS 1



MET ALA N3 Y23 E28 M35 I46 L77 T84 V93 A94 N95 W121 K122 L127 A128 L139 P140 L141 W144 P158 Y163 T186 M201 E249 Y252 GLY GLY LYS LYS LEU ALA PRO LEU ASN VAL

• Molecule 1: FRUCTOSE-BISPHOSPHATE ALDOLASE CLASS 1



MET ALA N3 K7 Y23 E28 D33 I46 L77 T84 N95 V98 G110 Y111 T112 A128 R132 D133 L139 V142 V143 W144 P158 Y163 K177 T186 V193 M201 K208 E211 K215 A229 Y252 GLY GLY LYS LYS LEU ALA PRO LEU ASN VAL

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	82.90Å 159.20Å 101.40Å 90.00° 107.80° 90.00°	Depositor
Resolution (Å)	39.80 – 1.93 39.80 – 1.93	Depositor EDS
% Data completeness (in resolution range)	100.0 (39.80-1.93) 97.2 (39.80-1.93)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.35 (at 1.94Å)	Xtriage
Refinement program	REFMAC 5.5.0110	Depositor
R, R_{free}	0.156 , 0.198 0.157 , 0.199	Depositor DCC
R_{free} test set	2723 reflections (1.50%)	wwPDB-VP
Wilson B-factor (Å ²)	19.5	Xtriage
Anisotropy	0.036	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 42.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	21227	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.45% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: M2P

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	0.84	1/2023 (0.0%)	0.77	0/2735
1	B	0.82	0/1999	0.74	1/2704 (0.0%)
1	C	0.79	1/2009 (0.0%)	0.71	0/2720
1	D	0.79	0/1989	0.71	0/2691
1	E	0.85	2/2041 (0.1%)	0.74	1/2758 (0.0%)
1	F	0.77	0/2008	0.71	0/2715
1	G	0.81	0/1986	0.75	1/2687 (0.0%)
1	H	0.81	0/1995	0.73	0/2699
1	I	0.82	0/1984	0.74	0/2685
1	J	0.84	0/1997	0.76	1/2701 (0.0%)
All	All	0.81	4/20031 (0.0%)	0.74	4/27095 (0.0%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	89	GLU	CB-CG	6.83	1.65	1.52
1	A	89	GLU	CB-CG	5.68	1.62	1.52
1	E	89	GLU	CG-CD	5.65	1.60	1.51
1	C	89	GLU	CB-CG	5.09	1.61	1.52

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	J	132	ARG	NE-CZ-NH1	7.43	124.02	120.30
1	G	14	ARG	NE-CZ-NH2	-6.32	117.14	120.30
1	E	89	GLU	OE1-CD-OE2	-5.89	116.23	123.30
1	B	14	ARG	NE-CZ-NH1	5.34	122.97	120.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1962	0	1981	9	0
1	B	1950	0	1968	9	0
1	C	1953	0	1979	9	0
1	D	1943	0	1961	10	0
1	E	1983	0	2010	14	0
1	F	1951	0	1971	9	0
1	G	1938	0	1952	10	0
1	H	1945	0	1961	16	0
1	I	1937	0	1948	9	0
1	J	1947	0	1965	14	0
2	A	20	0	11	0	0
2	B	20	0	11	0	0
2	C	20	0	11	0	0
2	D	20	0	11	0	0
2	E	20	0	11	0	0
2	F	20	0	11	0	0
2	G	20	0	11	0	0
2	H	20	0	11	0	0
2	I	20	0	11	0	0
2	J	20	0	11	0	0
3	A	158	0	0	3	0
3	B	173	0	0	2	0
3	C	136	0	0	3	0
3	D	126	0	0	4	0
3	E	173	0	0	2	0
3	F	122	0	0	1	0
3	G	148	0	0	2	0
3	H	150	0	0	3	0
3	I	158	0	0	1	0
3	J	174	0	0	3	0
All	All	21227	0	19806	95	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 95 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:C:7:LYS:HE2	3:C:2076:HOH:O	1.81	0.81
1:E:7:LYS:HE2	3:E:2091:HOH:O	1.83	0.79
1:D:33:ASP:HA	3:D:2025:HOH:O	1.86	0.75
1:A:7:LYS:HE2	3:A:2077:HOH:O	1.87	0.74
1:D:121:TRP:HD1	3:D:2075:HOH:O	1.77	0.67

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	253/263 (96%)	247 (98%)	6 (2%)	0	100	100
1	B	251/263 (95%)	244 (97%)	7 (3%)	0	100	100
1	C	252/263 (96%)	246 (98%)	6 (2%)	0	100	100
1	D	249/263 (95%)	243 (98%)	6 (2%)	0	100	100
1	E	257/263 (98%)	251 (98%)	6 (2%)	0	100	100
1	F	251/263 (95%)	243 (97%)	8 (3%)	0	100	100
1	G	249/263 (95%)	243 (98%)	6 (2%)	0	100	100
1	H	250/263 (95%)	243 (97%)	7 (3%)	0	100	100
1	I	249/263 (95%)	244 (98%)	5 (2%)	0	100	100
1	J	250/263 (95%)	245 (98%)	5 (2%)	0	100	100
All	All	2511/2630 (96%)	2449 (98%)	62 (2%)	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	203/207 (98%)	196 (97%)	7 (3%)	37	22
1	B	200/207 (97%)	194 (97%)	6 (3%)	41	27
1	C	202/207 (98%)	196 (97%)	6 (3%)	41	27
1	D	199/207 (96%)	191 (96%)	8 (4%)	31	16
1	E	203/207 (98%)	196 (97%)	7 (3%)	37	22
1	F	201/207 (97%)	195 (97%)	6 (3%)	41	27
1	G	199/207 (96%)	192 (96%)	7 (4%)	36	21
1	H	200/207 (97%)	192 (96%)	8 (4%)	31	16
1	I	199/207 (96%)	193 (97%)	6 (3%)	41	27
1	J	200/207 (97%)	194 (97%)	6 (3%)	41	27
All	All	2006/2070 (97%)	1939 (97%)	67 (3%)	38	24

5 of 67 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	I	139	LEU
1	I	163	TYR
1	J	201	MET
1	D	163	TYR
1	D	144	TRP

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

Mol	Chain	Res	Type
1	D	87	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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](#)

10 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	M2P	D	270	1	19,19,19	1.76	6 (31%)	26,28,28	1.06	3 (11%)
2	M2P	G	270	1	19,19,19	1.67	6 (31%)	26,28,28	1.14	2 (7%)
2	M2P	H	270	1	19,19,19	1.70	4 (21%)	26,28,28	1.05	1 (3%)
2	M2P	I	270	1	19,19,19	1.66	4 (21%)	26,28,28	0.74	0
2	M2P	J	270	1	19,19,19	1.32	2 (10%)	26,28,28	0.92	0
2	M2P	A	270	1	19,19,19	1.37	2 (10%)	26,28,28	1.09	0
2	M2P	B	270	1	19,19,19	1.76	3 (15%)	26,28,28	0.92	0
2	M2P	C	270	1	19,19,19	1.46	4 (21%)	26,28,28	1.29	5 (19%)
2	M2P	E	270	1	19,19,19	1.23	2 (10%)	26,28,28	1.13	2 (7%)
2	M2P	F	270	1	19,19,19	1.41	2 (10%)	26,28,28	1.05	1 (3%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	M2P	D	270	1	-	8/24/24/24	-
2	M2P	G	270	1	-	7/24/24/24	-
2	M2P	H	270	1	-	9/24/24/24	-
2	M2P	I	270	1	-	8/24/24/24	-
2	M2P	J	270	1	-	9/24/24/24	-
2	M2P	A	270	1	-	5/24/24/24	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	M2P	B	270	1	-	6/24/24/24	-
2	M2P	C	270	1	-	9/24/24/24	-
2	M2P	E	270	1	-	9/24/24/24	-
2	M2P	F	270	1	-	5/24/24/24	-

The worst 5 of 35 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	270	M2P	C1-C2	4.74	1.58	1.51
2	G	270	M2P	C1-C2	3.64	1.56	1.51
2	D	270	M2P	P1-O12	3.50	1.61	1.50
2	I	270	M2P	C1-C2	3.48	1.56	1.51
2	F	270	M2P	P1-O12	3.37	1.61	1.50

The worst 5 of 14 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	270	M2P	O6-P6-O61	2.67	113.66	106.44
2	C	270	M2P	O62-P6-O6	-2.60	99.90	106.67
2	C	270	M2P	O63-P6-O6	2.56	113.34	106.67
2	D	270	M2P	O4-C4-C3	2.39	115.03	109.46
2	E	270	M2P	O62-P6-O6	-2.38	100.47	106.67

There are no chirality outliers.

5 of 75 torsion outliers are listed below:

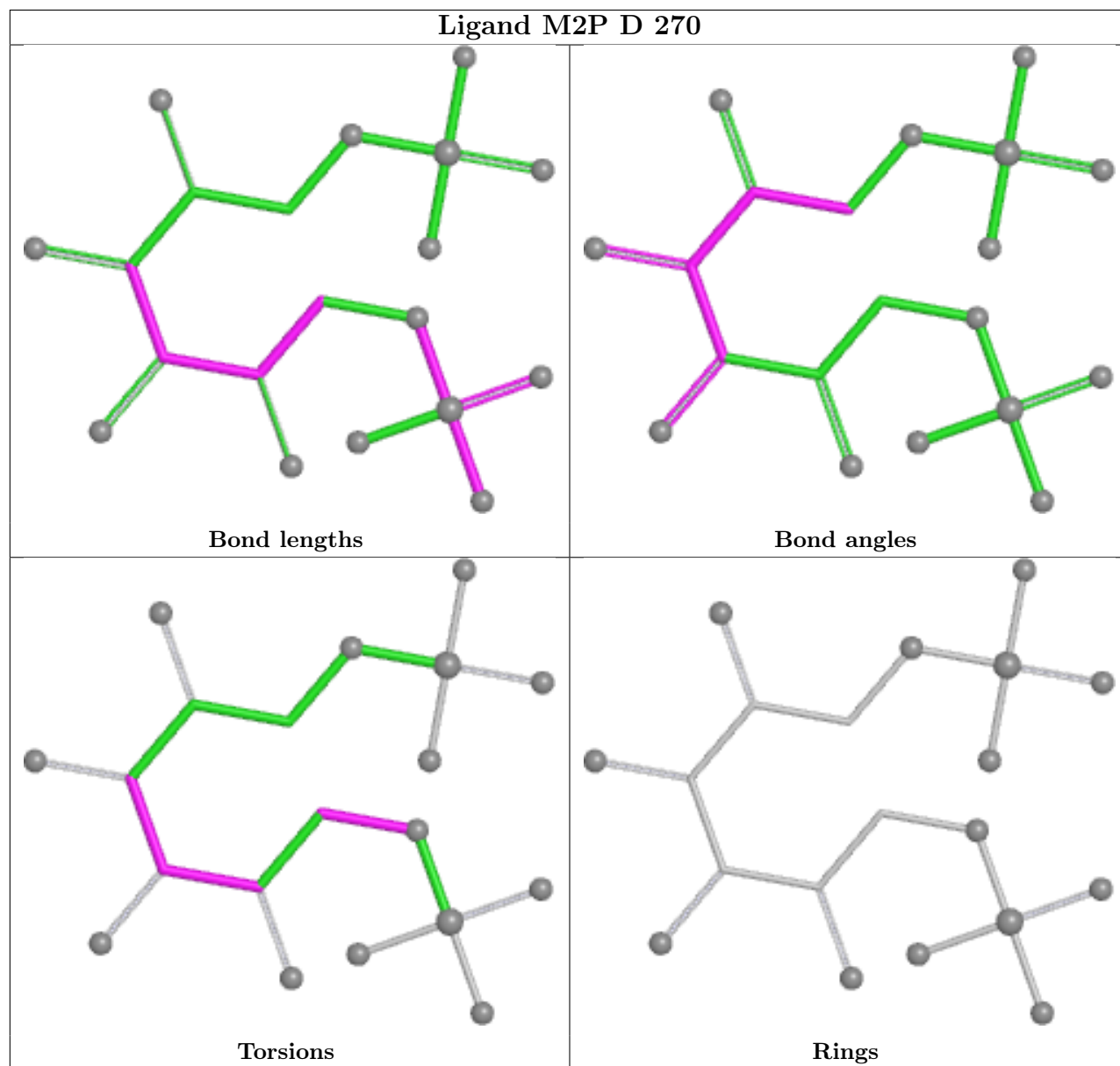
Mol	Chain	Res	Type	Atoms
2	A	270	M2P	C2-C3-C4-C5
2	A	270	M2P	O3-C3-C4-C5
2	A	270	M2P	C2-C3-C4-O4
2	A	270	M2P	O3-C3-C4-O4
2	B	270	M2P	C2-C3-C4-C5

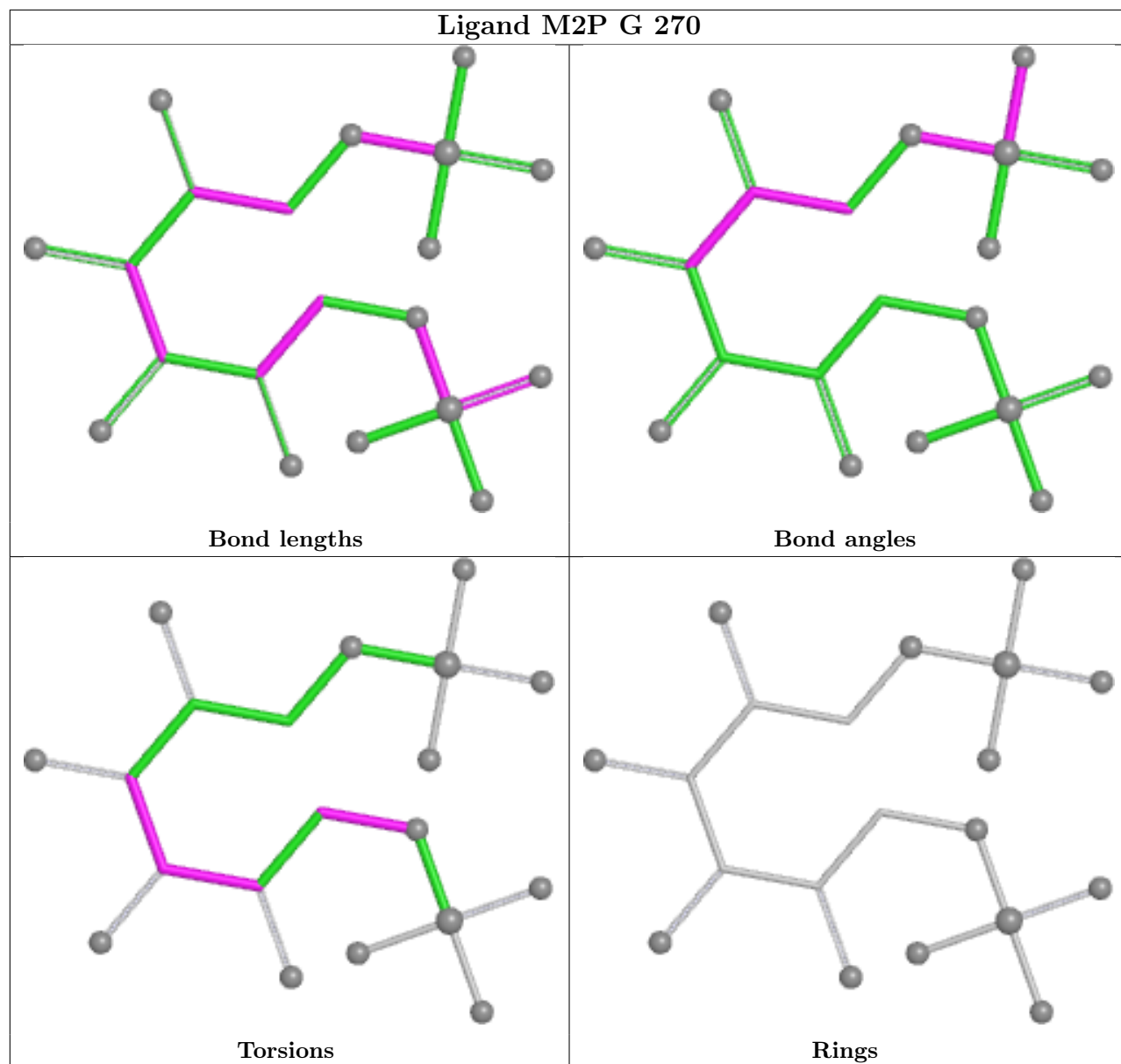
There are no ring outliers.

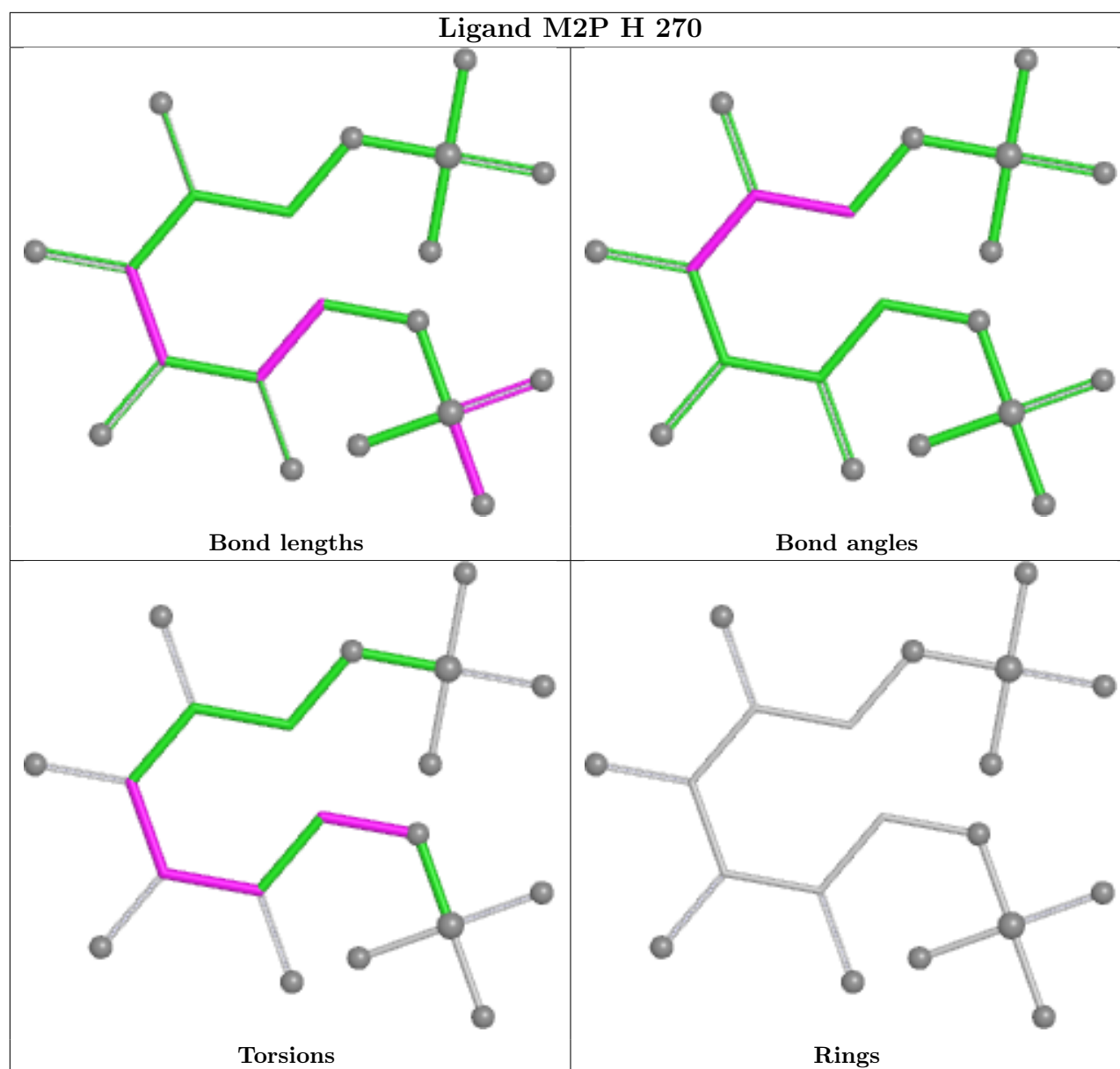
No monomer is involved in short contacts.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier.

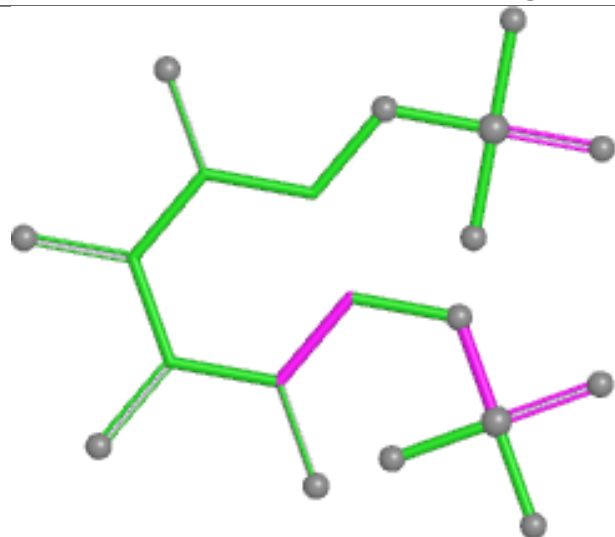
Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



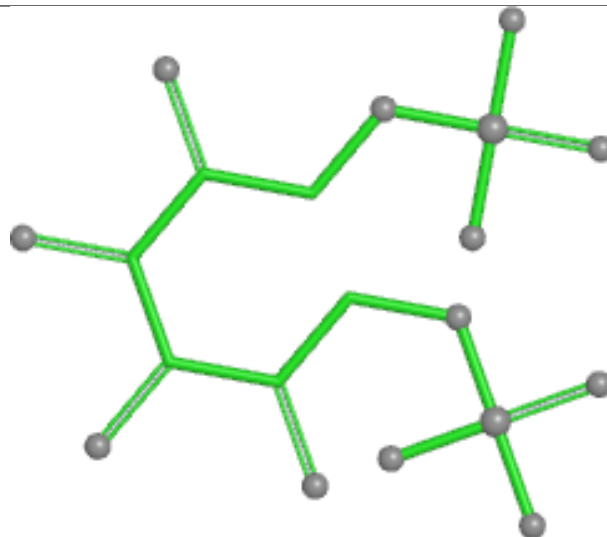




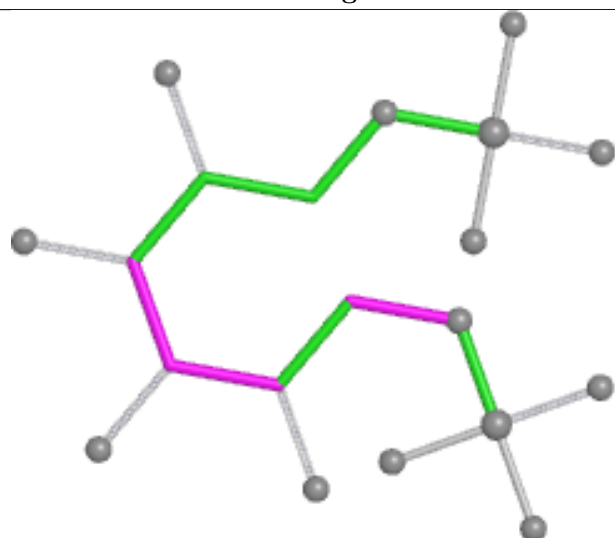
Ligand M2P I 270



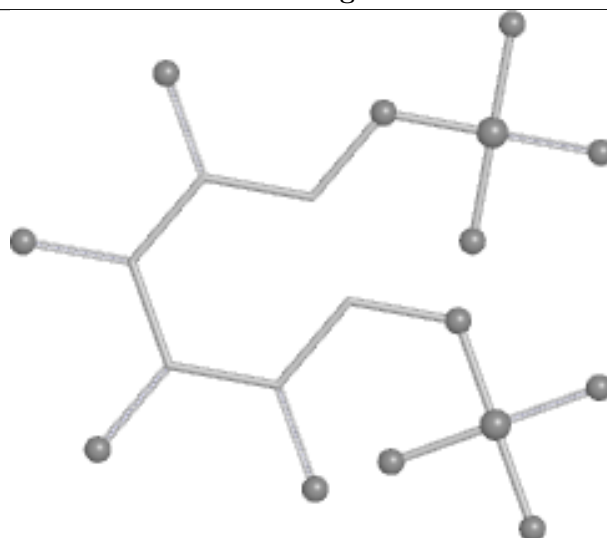
Bond lengths



Bond angles

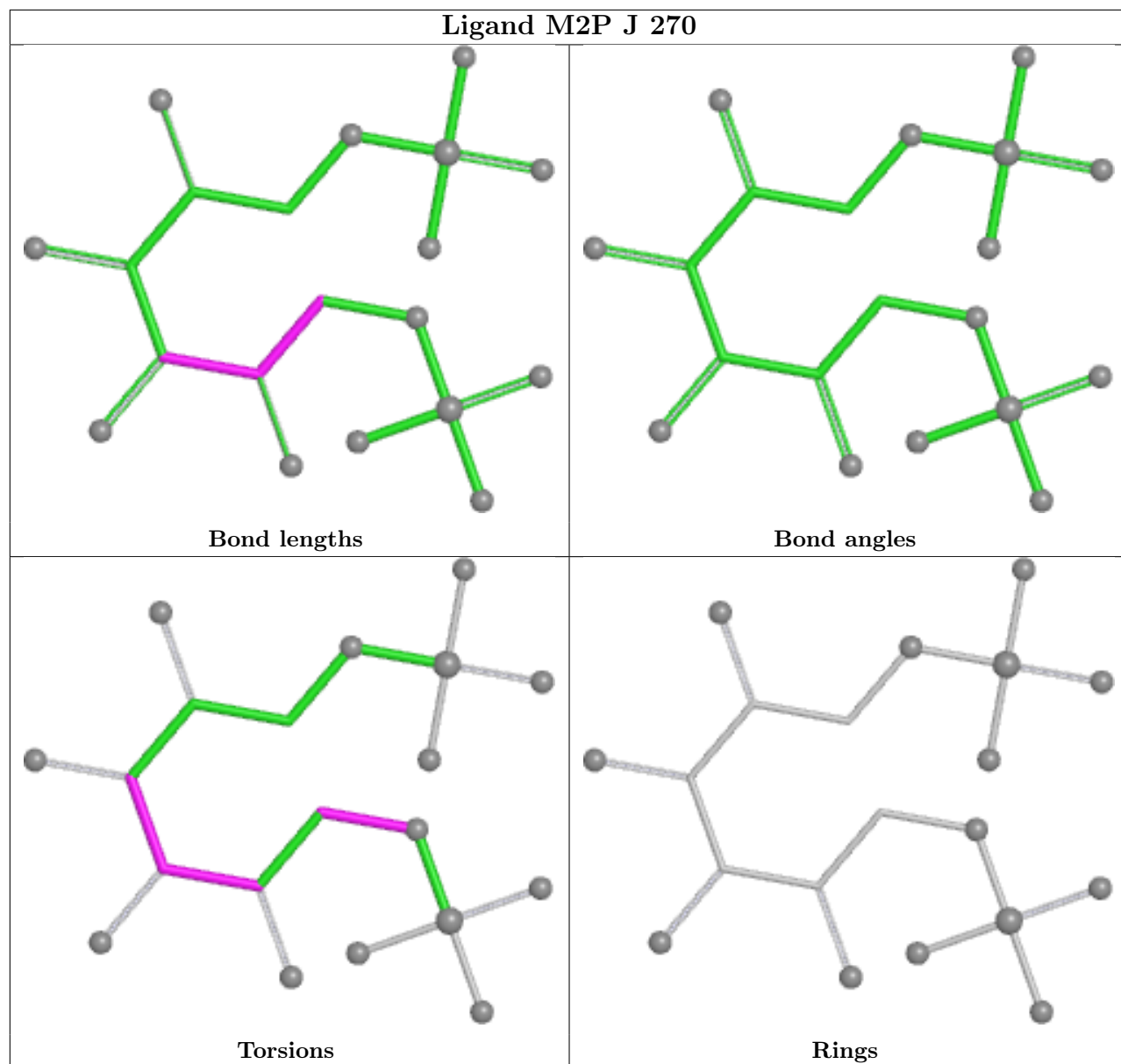


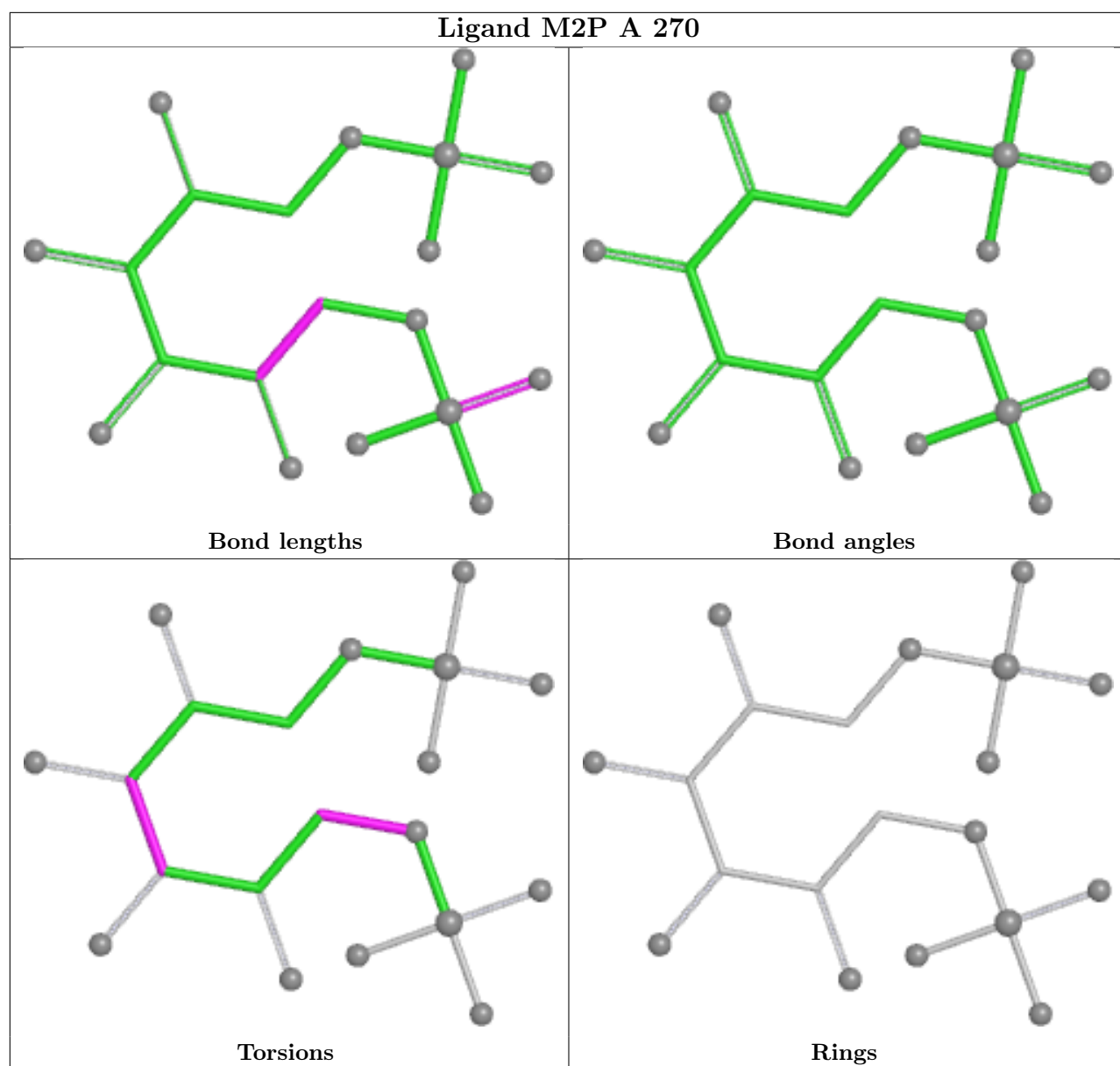
Torsions

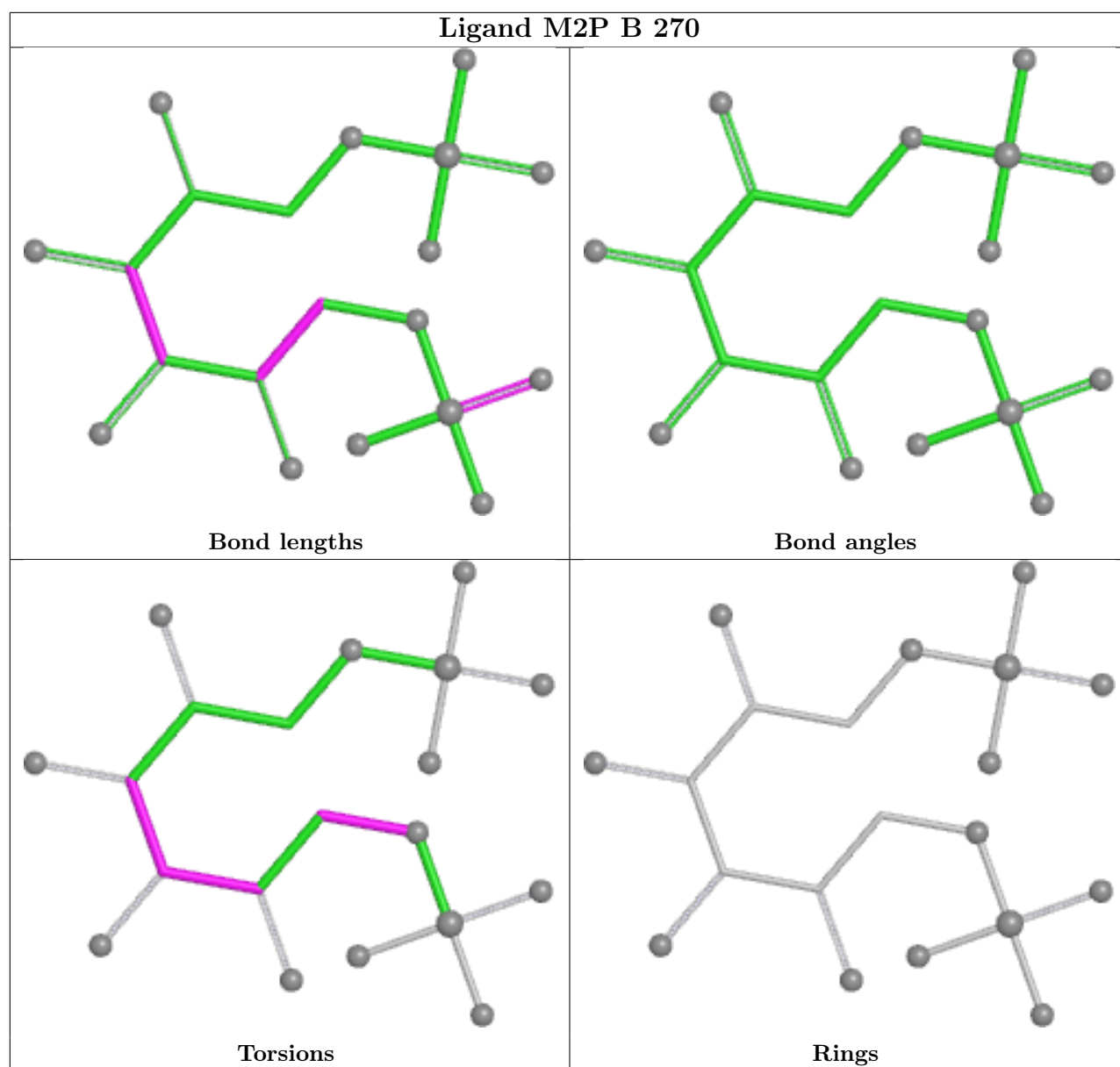


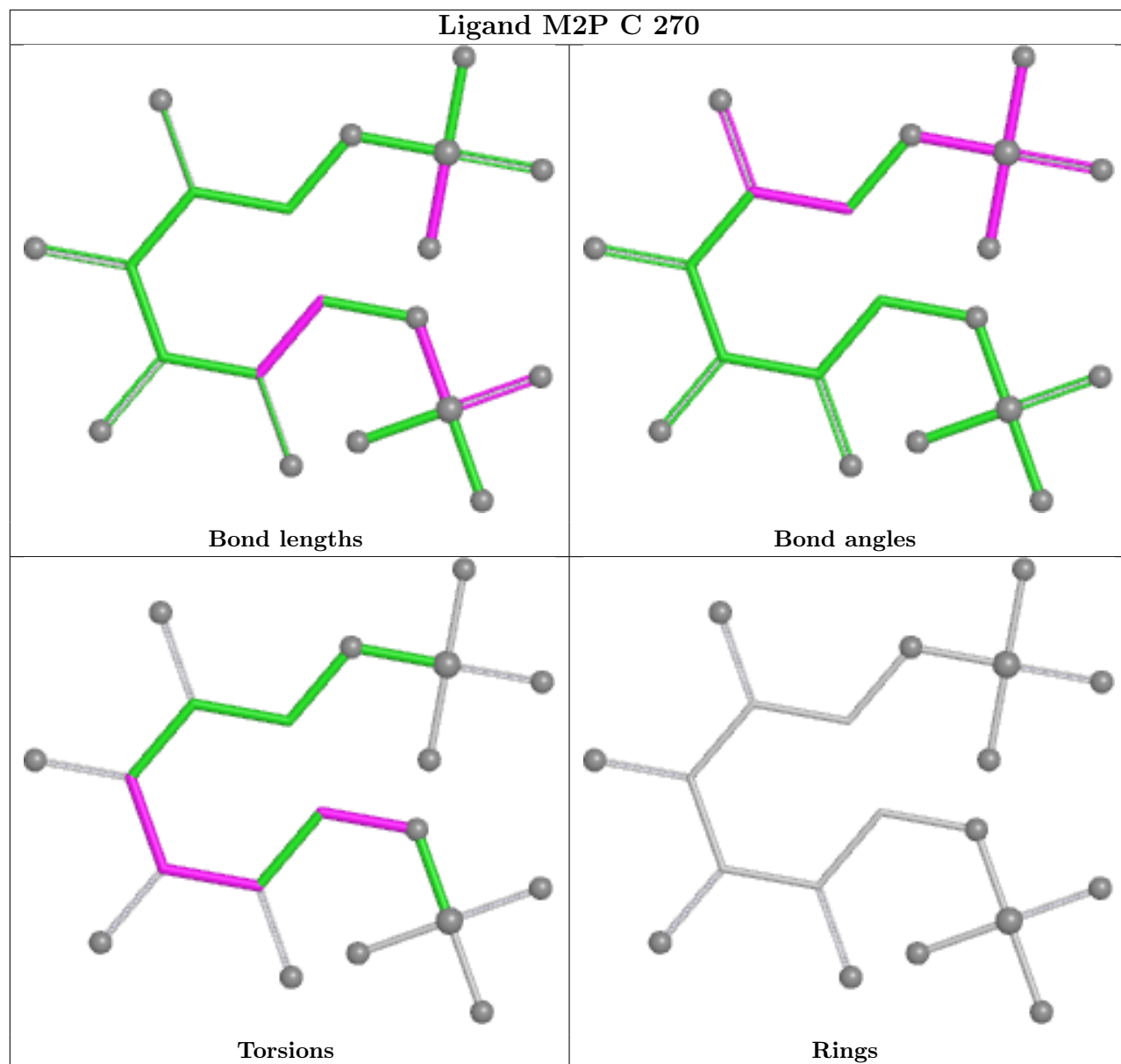
Rings

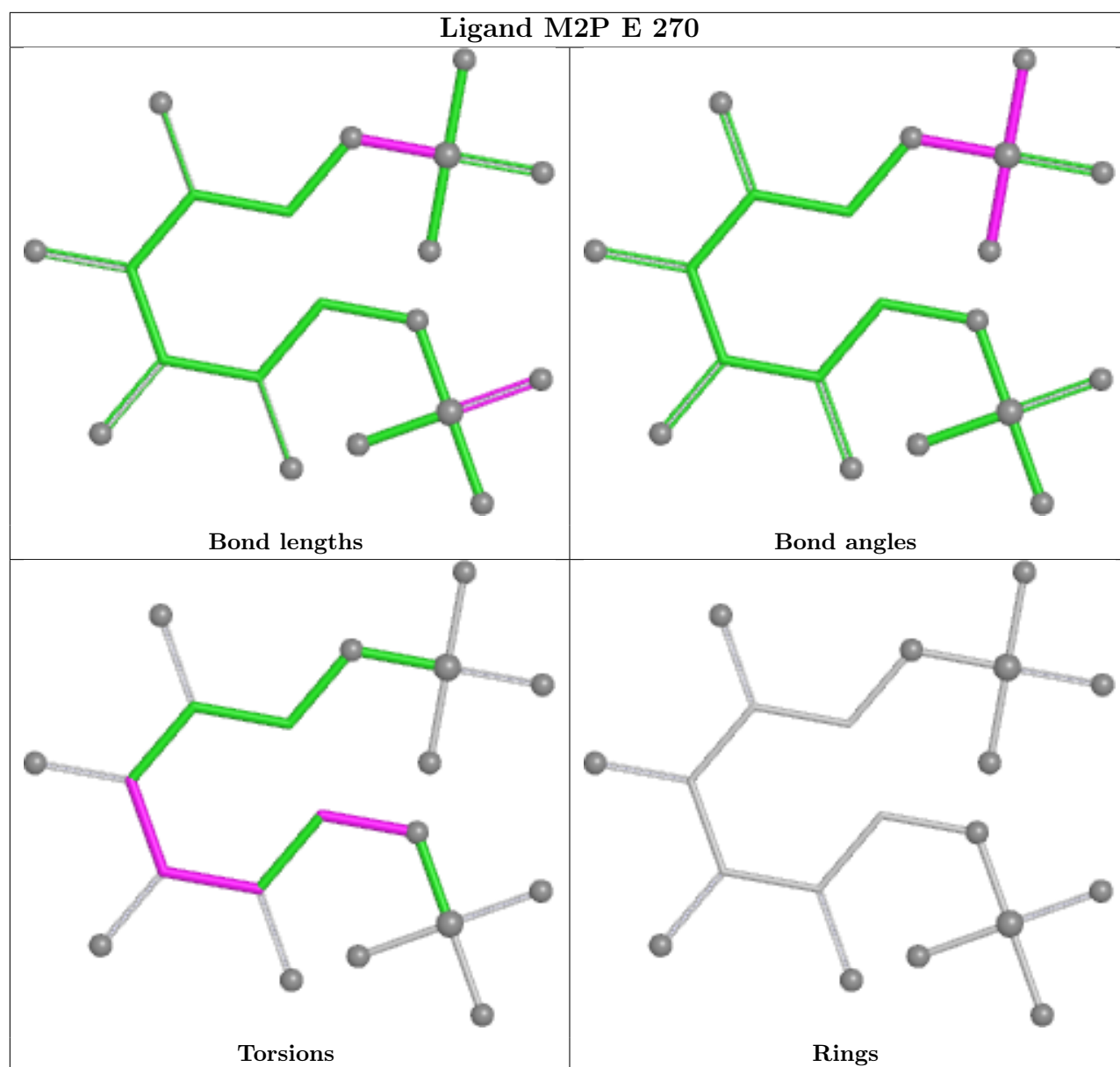
Ligand M2P J 270

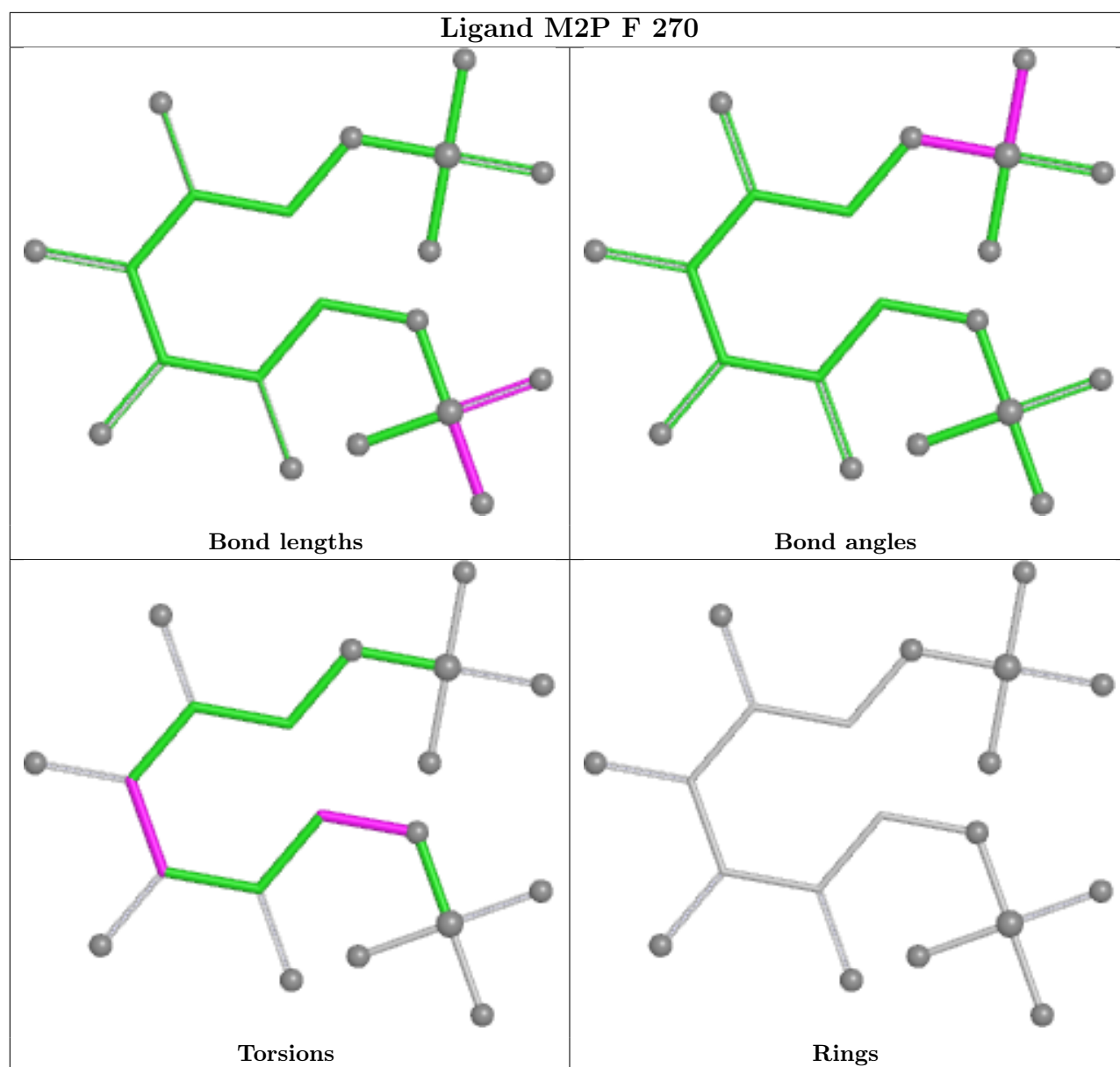












5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	250/263 (95%)	-0.32	1 (0%) 92 95	11, 19, 37, 50	0
1	B	251/263 (95%)	-0.46	1 (0%) 92 95	11, 18, 31, 39	0
1	C	250/263 (95%)	-0.41	0 100 100	12, 23, 46, 59	0
1	D	250/263 (95%)	-0.48	3 (1%) 79 83	11, 21, 41, 50	0
1	E	255/263 (96%)	-0.30	4 (1%) 72 77	11, 19, 38, 60	0
1	F	250/263 (95%)	-0.43	3 (1%) 79 83	13, 23, 43, 52	0
1	G	250/263 (95%)	-0.48	2 (0%) 86 89	12, 21, 37, 49	0
1	H	250/263 (95%)	-0.24	5 (2%) 65 71	12, 21, 38, 54	0
1	I	250/263 (95%)	-0.44	1 (0%) 92 95	11, 19, 36, 44	0
1	J	250/263 (95%)	-0.48	2 (0%) 86 89	12, 18, 33, 46	0
All	All	2506/2630 (95%)	-0.40	22 (0%) 84 87	11, 20, 38, 60	0

The worst 5 of 22 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	208	LYS	3.4
1	E	256	LYS	2.9
1	E	238	ARG	2.8
1	E	257	LEU	2.7
1	H	238	ARG	2.6

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](#)

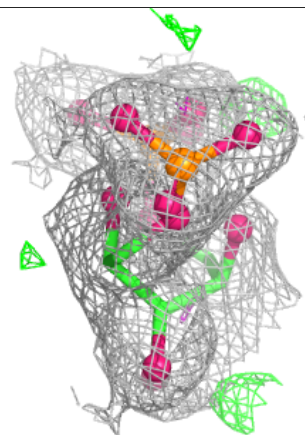
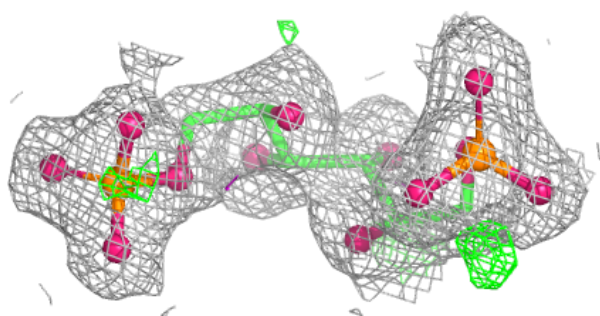
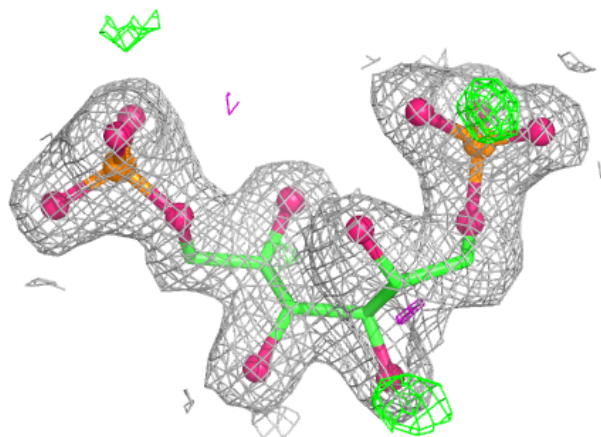
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	M2P	D	270	20/20	0.97	0.08	20,26,30,31	0
2	M2P	F	270	20/20	0.97	0.07	21,27,31,31	0
2	M2P	G	270	20/20	0.97	0.09	22,25,29,31	0
2	M2P	H	270	20/20	0.97	0.08	20,26,29,31	0
2	M2P	I	270	20/20	0.97	0.07	18,24,28,30	0
2	M2P	B	270	20/20	0.98	0.09	17,24,28,29	0
2	M2P	C	270	20/20	0.98	0.07	21,25,29,30	0
2	M2P	A	270	20/20	0.98	0.07	20,24,30,31	0
2	M2P	E	270	20/20	0.98	0.07	19,25,29,32	0
2	M2P	J	270	20/20	0.98	0.08	17,24,28,30	0

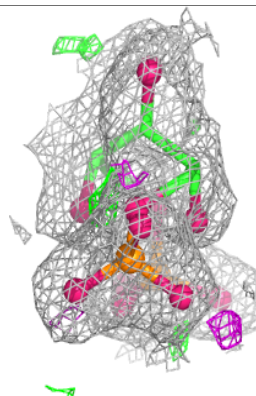
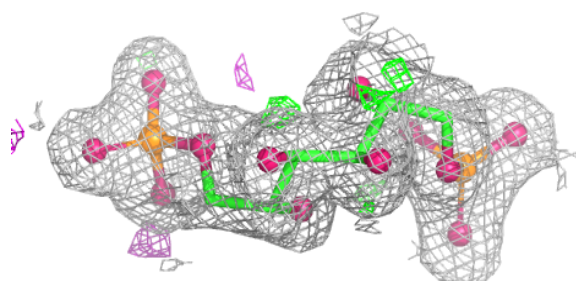
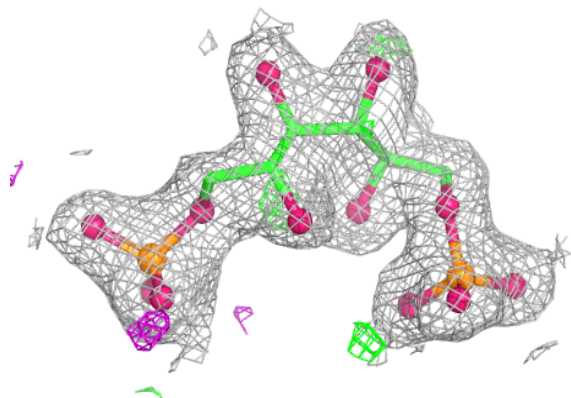
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around M2P D 270:

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

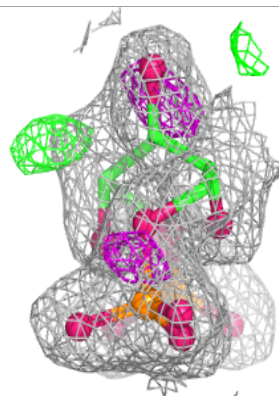
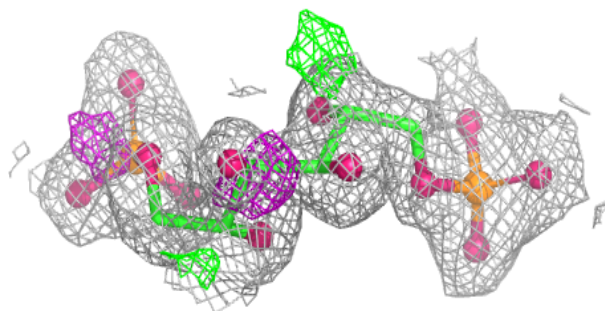
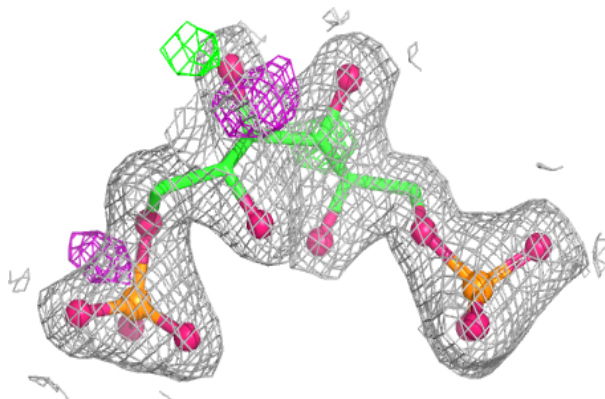
**Electron density around M2P F 270:**

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

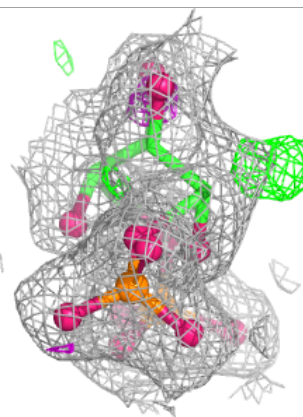
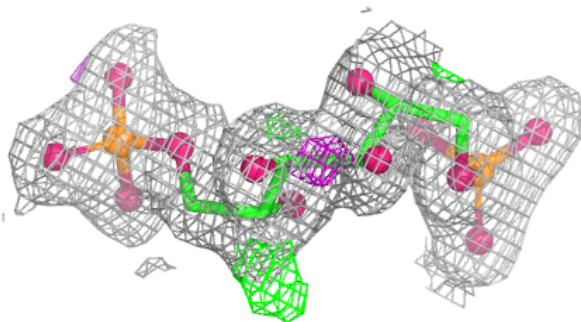
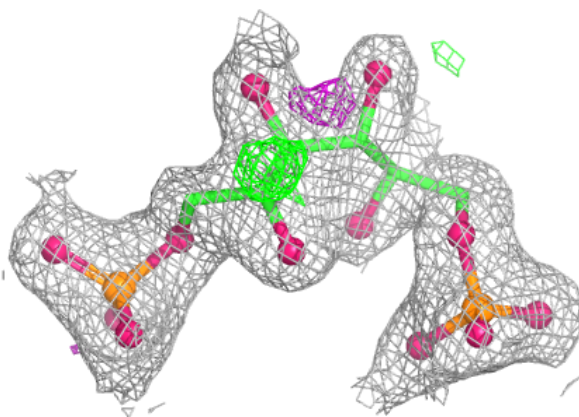


Electron density around M2P G 270:

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

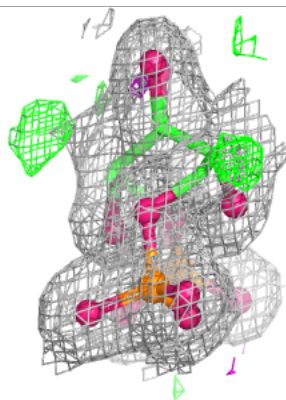
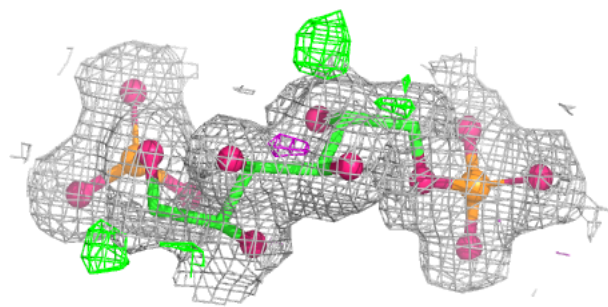
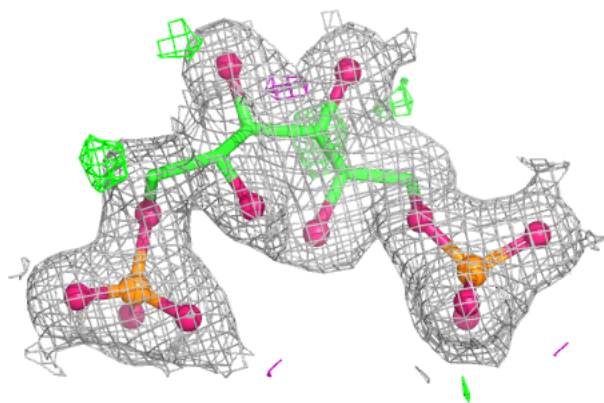
**Electron density around M2P H 270:**

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

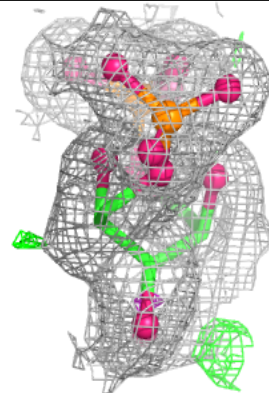
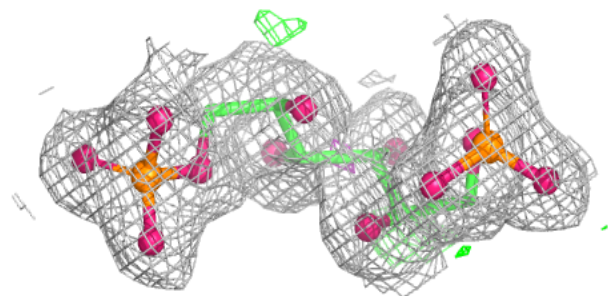
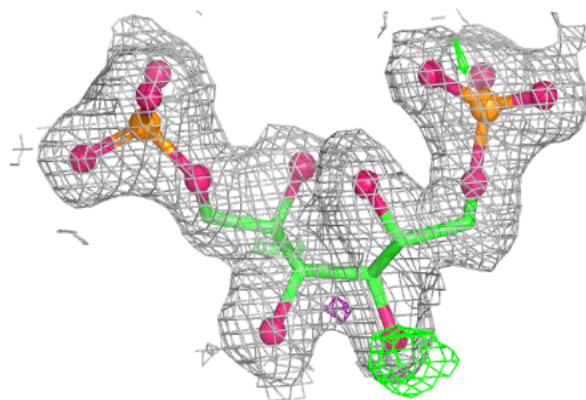


Electron density around M2P I 270:

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

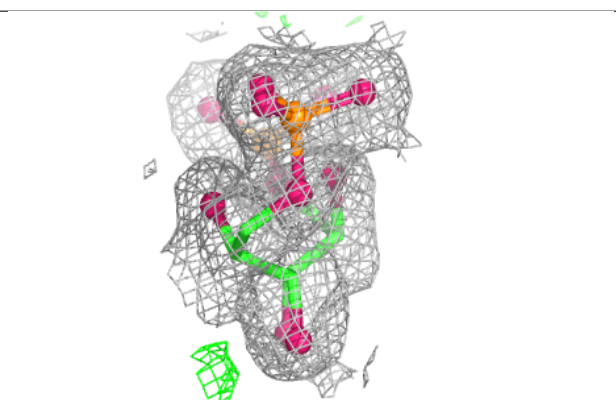
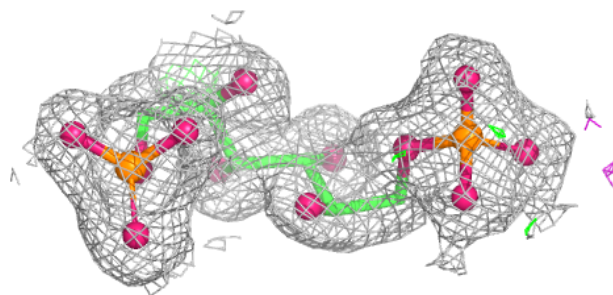
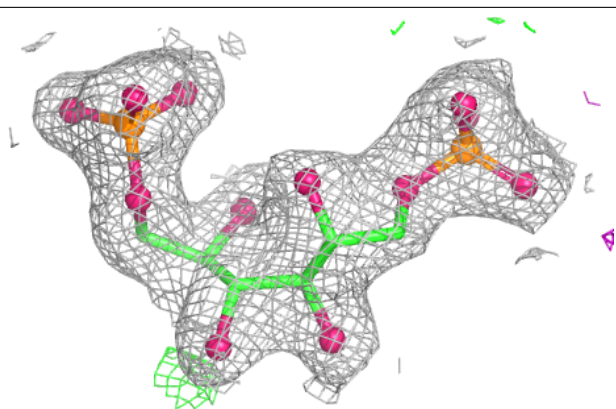
**Electron density around M2P B 270:**

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

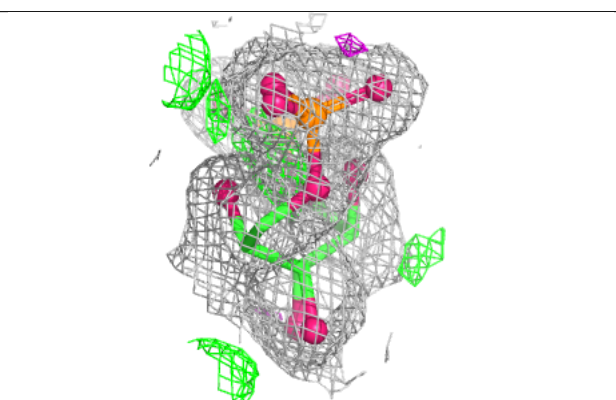
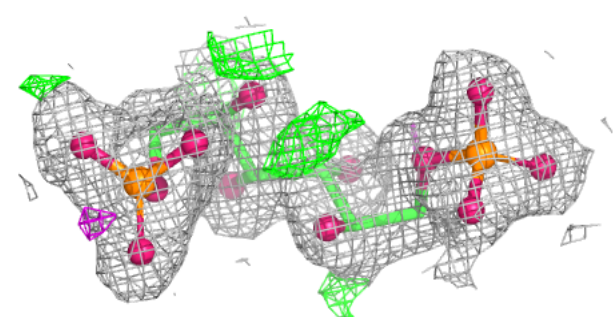
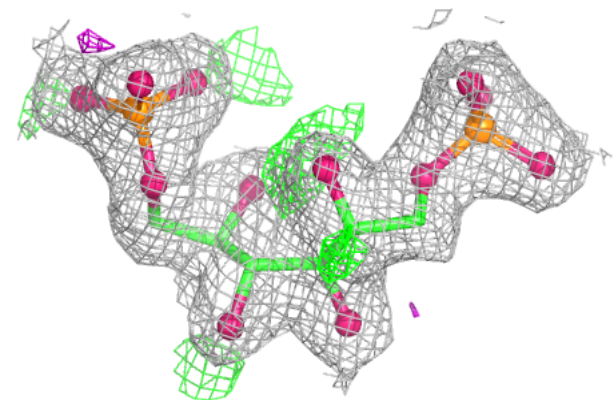


Electron density around M2P C 270:

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

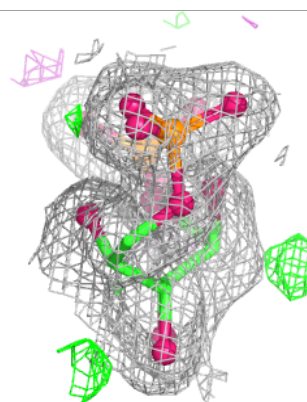
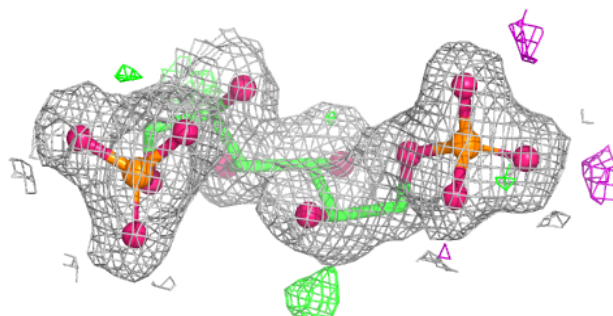
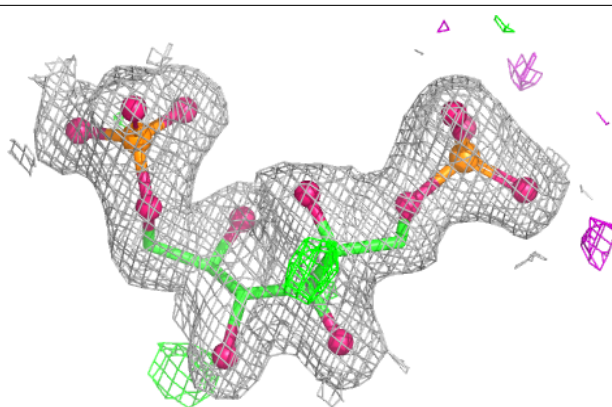
**Electron density around M2P A 270:**

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

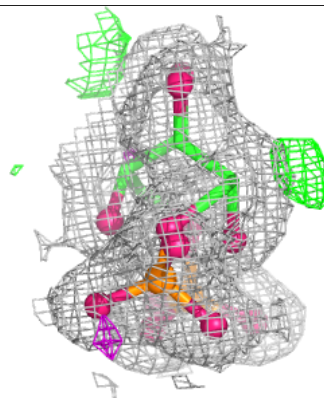
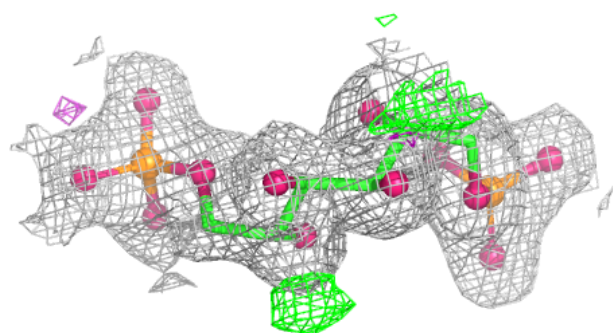
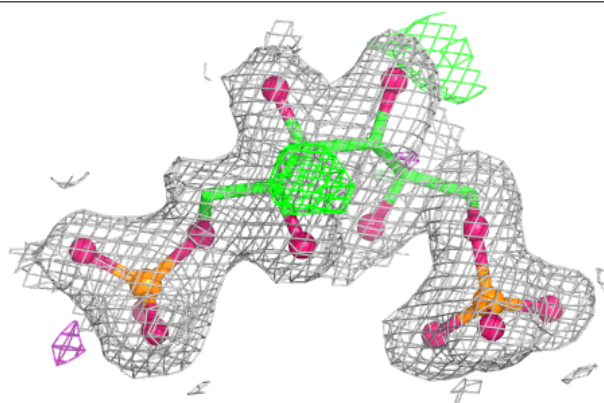


Electron density around M2P E 270:

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

**Electron density around M2P J 270:**

$2mF_o-DF_c$ (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.