



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2022 – 04:18 PM EST

PDB ID : 5SC9
Title : Structure of liver pyruvate kinase in complex with anthraquinone derivative 29
Authors : Lulla, A.; Foller, A.; Nain-Perez, A.; Grotli, M.; Brear, P.; Hyvonen, M.
Deposited on : 2021-12-01
Resolution : 1.69 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The 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 : 2.27
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) : 2.27

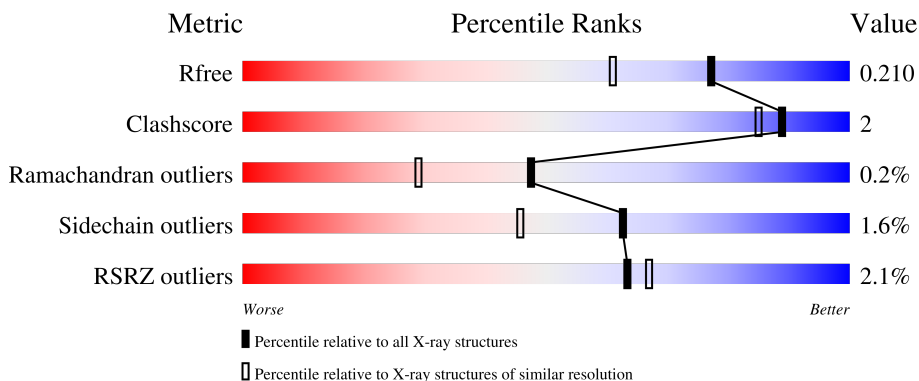
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.69 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.



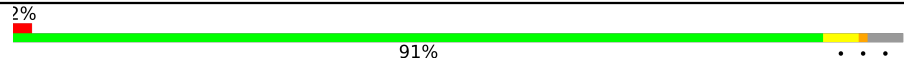

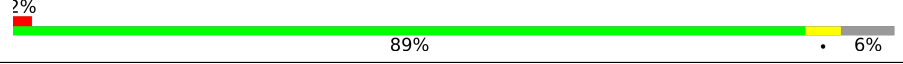
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	6780 (1.70-1.66)
Clashscore	141614	7310 (1.70-1.66)
Ramachandran outliers	138981	7173 (1.70-1.66)
Sidechain outliers	138945	7172 (1.70-1.66)
RSRZ outliers	127900	6661 (1.70-1.66)

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	447	 2% 86% 6% 7%
1	B	447	 2% 91% 5%
1	C	447	 2% 89% 7%
1	D	447	 2% 90% 6%
1	E	447	 2% 89% 7%

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Mol	Chain	Length	Quality of chain
1	F	447	 2% 91% 5% 6% . . .
1	G	447	 % 88% 5% 6%
1	H	447	 2% 89% 5% 6% .

2 Entry composition i

There are 6 unique types of molecules in this entry. The entry contains 29743 atoms, of which 119 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 Pyruvate kinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	414	3232	2032	581	599	20	0	15	0
1	B	431	3325	2093	596	616	20	3	9	0
1	C	417	3215	2020	579	597	19	0	9	0
1	D	419	3231	2031	585	596	19	0	8	0
1	E	414	3247	2040	583	604	20	0	17	0
1	F	431	3325	2093	596	616	20	0	9	0
1	G	419	3237	2038	581	599	19	0	9	0
1	H	421	3256	2047	590	600	19	0	10	0

There are 48 discrepancies between the modelled and reference sequences:

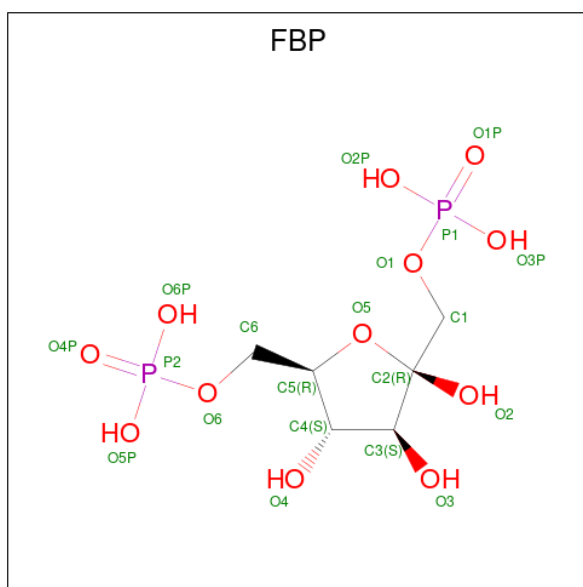
Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	expression tag	UNP Q16716
A	0	SER	-	expression tag	UNP Q16716
A	12	ASP	SER	conflict	UNP Q16716
A	228	GLY	-	linker	UNP Q16716
A	229	SER	-	linker	UNP Q16716
A	230	GLY	-	linker	UNP Q16716
B	-1	GLY	-	expression tag	UNP Q16716
B	0	SER	-	expression tag	UNP Q16716
B	12	ASP	SER	conflict	UNP Q16716
B	228	GLY	-	linker	UNP Q16716
B	229	SER	-	linker	UNP Q16716
B	230	GLY	-	linker	UNP Q16716
C	-1	GLY	-	expression tag	UNP Q16716

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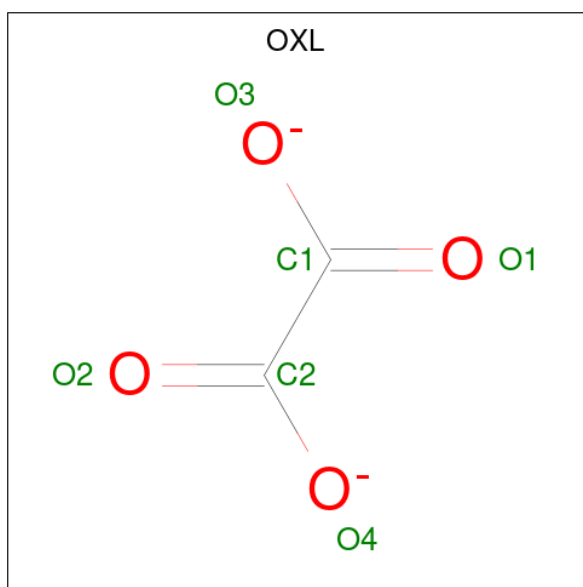
Chain	Residue	Modelled	Actual	Comment	Reference
C	0	SER	-	expression tag	UNP Q16716
C	12	ASP	SER	conflict	UNP Q16716
C	228	GLY	-	linker	UNP Q16716
C	229	SER	-	linker	UNP Q16716
C	230	GLY	-	linker	UNP Q16716
D	-1	GLY	-	expression tag	UNP Q16716
D	0	SER	-	expression tag	UNP Q16716
D	12	ASP	SER	conflict	UNP Q16716
D	228	GLY	-	linker	UNP Q16716
D	229	SER	-	linker	UNP Q16716
D	230	GLY	-	linker	UNP Q16716
E	-1	GLY	-	expression tag	UNP Q16716
E	0	SER	-	expression tag	UNP Q16716
E	12	ASP	SER	conflict	UNP Q16716
E	228	GLY	-	linker	UNP Q16716
E	229	SER	-	linker	UNP Q16716
E	230	GLY	-	linker	UNP Q16716
F	-1	GLY	-	expression tag	UNP Q16716
F	0	SER	-	expression tag	UNP Q16716
F	12	ASP	SER	conflict	UNP Q16716
F	228	GLY	-	linker	UNP Q16716
F	229	SER	-	linker	UNP Q16716
F	230	GLY	-	linker	UNP Q16716
G	-1	GLY	-	expression tag	UNP Q16716
G	0	SER	-	expression tag	UNP Q16716
G	12	ASP	SER	conflict	UNP Q16716
G	228	GLY	-	linker	UNP Q16716
G	229	SER	-	linker	UNP Q16716
G	230	GLY	-	linker	UNP Q16716
H	-1	GLY	-	expression tag	UNP Q16716
H	0	SER	-	expression tag	UNP Q16716
H	12	ASP	SER	conflict	UNP Q16716
H	130	GLY	-	linker	UNP Q16716
H	229	SER	-	linker	UNP Q16716
H	230	GLY	-	linker	UNP Q16716

- Molecule 2 is 1,6-di-O-phosphono-beta-D-fructofuranose (three-letter code: FBP) (formula: $C_6H_{14}O_{12}P_2$).



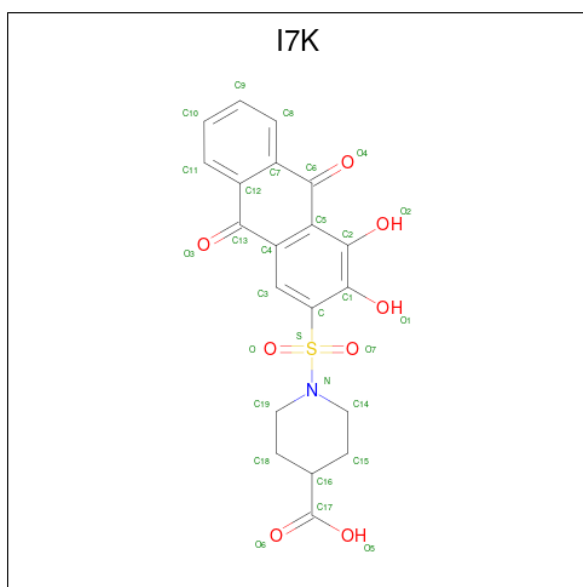
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	O	P	0	0
			20	6	12	2		
2	B	1	Total	C	O	P	0	0
			20	6	12	2		
2	C	1	Total	C	O	P	0	0
			20	6	12	2		
2	D	1	Total	C	O	P	0	0
			20	6	12	2		
2	E	1	Total	C	O	P	0	0
			20	6	12	2		
2	F	1	Total	C	O	P	0	0
			20	6	12	2		
2	G	1	Total	C	O	P	0	0
			20	6	12	2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

- Molecule 3 is OXALATE ION (three-letter code: OXL) (formula: C_2O_4).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 6 2 4	0	0
3	B	1	Total C O 6 2 4	0	0
3	C	1	Total C O 6 2 4	0	0
3	D	1	Total C O 6 2 4	0	0
3	E	1	Total C O 6 2 4	0	0
3	F	1	Total C O 6 2 4	0	0
3	G	1	Total C O 6 2 4	0	0
3	H	1	Total C O 6 2 4	0	0

- Molecule 4 is 1-(3,4-dihydroxy-9,10-dioxo-9,10-dihydroanthracene-2-sulfonyl)piperidine-4-carboxylic acid (three-letter code: I7K) (formula: C₂₀H₁₇NO₈S) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	
			Total	C	H	N	O			S
4	A	1	47	20	17	1	8	1	17	0
4	C	1	47	20	17	1	8	1	17	0
4	D	1	47	20	17	1	8	1	17	0
4	E	1	47	20	17	1	8	1	17	0
4	F	1	47	20	17	1	8	1	17	0
4	G	1	47	20	17	1	8	1	17	0
4	H	1	47	20	17	1	8	1	17	0

- Molecule 5 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Mg		
5	A	1	1	1	0	0
5	B	1	1	1	0	0
5	C	2	2	2	0	0
5	D	1	1	1	0	0
5	E	1	1	1	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	F	2	Total 2	Mg 2	0	0
5	G	1	Total 1	Mg 1	0	0
5	H	1	Total 1	Mg 1	0	0

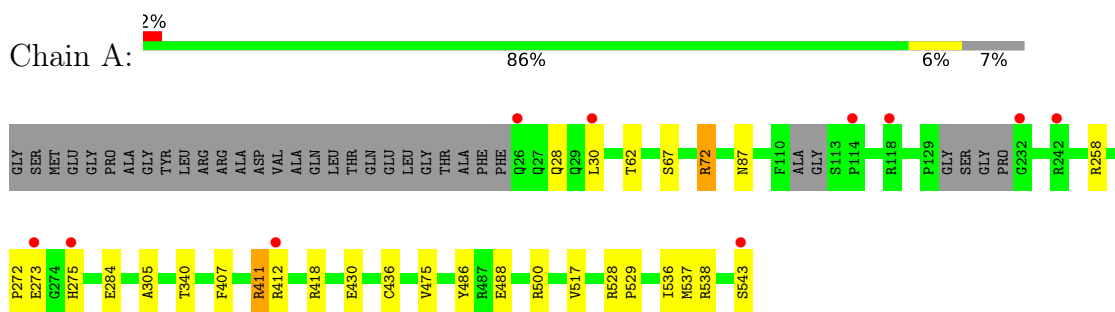
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	353	Total 353	O 353	0	0
6	B	295	Total 295	O 295	0	0
6	C	398	Total 398	O 398	0	0
6	D	431	Total 431	O 431	0	0
6	E	338	Total 338	O 338	0	0
6	F	398	Total 398	O 398	0	0
6	G	437	Total 437	O 437	0	0
6	H	478	Total 478	O 478	0	0

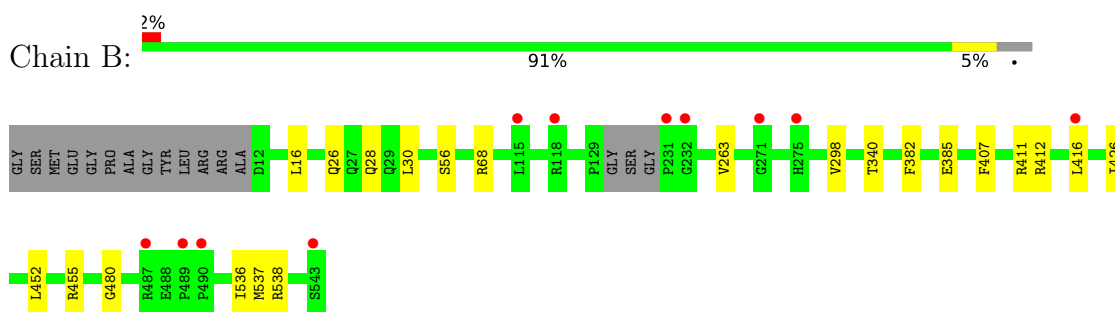
3 Residue-property plots [i](#)

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.

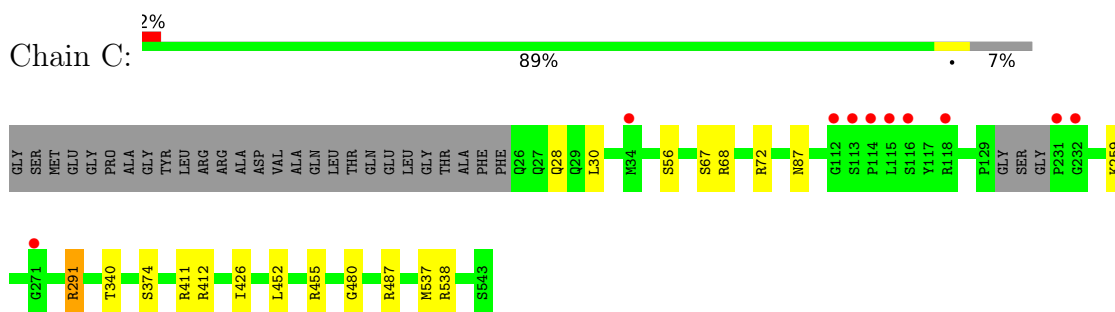
- Molecule 1: Pyruvate kinase



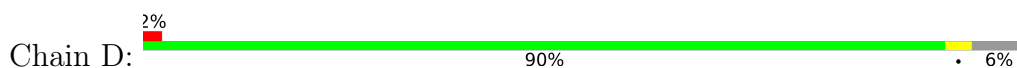
- Molecule 1: Pyruvate kinase

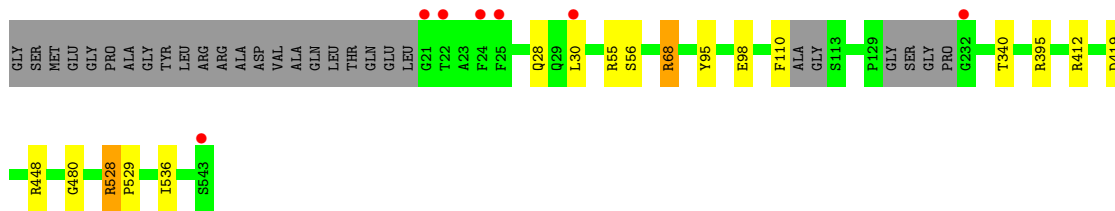


- Molecule 1: Pyruvate kinase

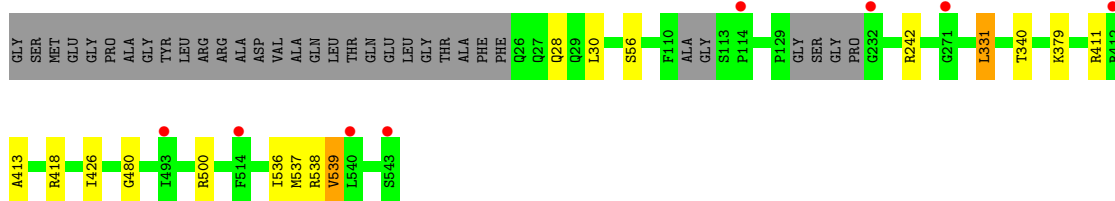


- Molecule 1: Pyruvate kinase

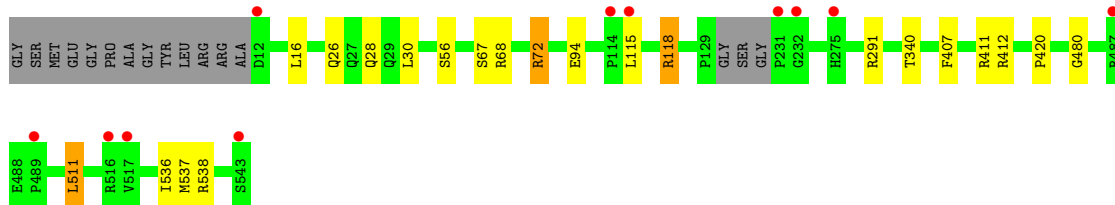
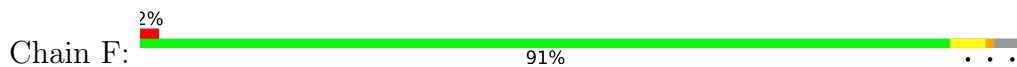




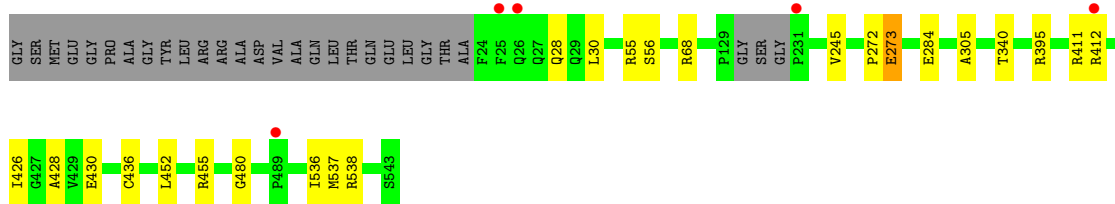
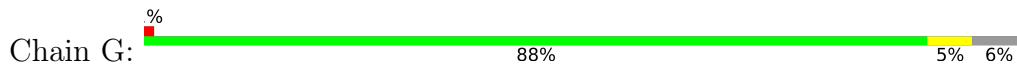
● Molecule 1: Pyruvate kinase



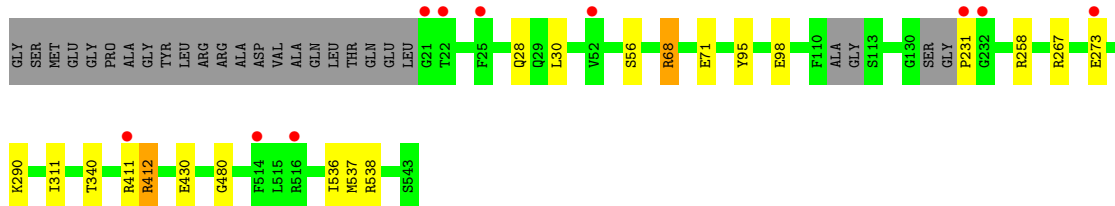
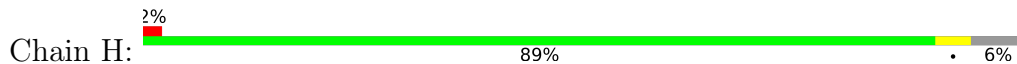
● Molecule 1: Pyruvate kinase



● Molecule 1: Pyruvate kinase



● Molecule 1: Pyruvate kinase



4 Data and refinement statistics i

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	208.27Å 112.68Å 187.90Å 90.00° 91.34° 90.00°	Depositor
Resolution (Å)	187.85 – 1.69 187.85 – 1.69	Depositor EDS
% Data completeness (in resolution range)	87.4 (187.85-1.69) 87.5 (187.85-1.69)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.71 (at 1.69Å)	Xtrriage
Refinement program	BUSTER 2.10.4 (16-JUL-2021)	Depositor
R, R_{free}	0.195 , 0.216 0.189 , 0.210	Depositor DCC
R_{free} test set	21129 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å ²)	21.8	Xtrriage
Anisotropy	0.003	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 54.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	0.000 for -h,-k,l	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	29743	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 41.24 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 2.4315e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: I7K, MG, FBP, OXL

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	0.39	0/3327	0.53	0/4493
1	B	0.38	0/3406	0.53	0/4604
1	C	0.43	0/3295	0.56	0/4454
1	D	0.45	0/3308	0.56	0/4469
1	E	0.36	0/3349	0.51	0/4523
1	F	0.41	0/3406	0.53	0/4604
1	G	0.44	0/3319	0.55	0/4486
1	H	0.47	0/3340	0.56	0/4511
All	All	0.42	0/26750	0.54	0/36144

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	3232	0	3317	20	0
1	B	3325	0	3400	14	0
1	C	3215	0	3283	11	0
1	D	3231	0	3296	12	0
1	E	3247	0	3325	9	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	F	3325	0	3400	15	0
1	G	3237	0	3301	12	0
1	H	3256	0	3326	19	0
2	A	20	0	10	0	0
2	B	20	0	10	0	0
2	C	20	0	10	0	0
2	D	20	0	10	0	0
2	E	20	0	10	0	0
2	F	20	0	10	0	0
2	G	20	0	10	1	0
2	H	20	0	10	0	0
3	A	6	0	0	0	0
3	B	6	0	0	0	0
3	C	6	0	0	0	0
3	D	6	0	0	0	0
3	E	6	0	0	0	0
3	F	6	0	0	0	0
3	G	6	0	0	0	0
3	H	6	0	0	0	0
4	A	30	17	0	3	0
4	C	30	17	0	2	0
4	D	30	17	0	0	0
4	E	30	17	0	0	0
4	F	30	17	0	1	0
4	G	30	17	0	0	0
4	H	30	17	0	0	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
5	C	2	0	0	0	0
5	D	1	0	0	0	0
5	E	1	0	0	0	0
5	F	2	0	0	0	0
5	G	1	0	0	0	0
5	H	1	0	0	0	0
6	A	353	0	0	1	0
6	B	295	0	0	0	0
6	C	398	0	0	1	0
6	D	431	0	0	2	0
6	E	338	0	0	0	0
6	F	398	0	0	0	0
6	G	437	0	0	2	0
6	H	478	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	29624	119	26728	100	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.

All (100) 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:H:311:ILE:CD1	1:H:311:ILE:CG1	1.79	1.55
1:D:110:PHE:C	6:D:715:HOH:O	1.77	1.23
2:G:601:FBP:H5	6:G:712:HOH:O	0.86	1.04
1:A:272:PRO:HA	1:A:275:HIS:NE2	1.92	0.85
1:G:538:ARG:HG2	1:H:536:ILE:HG12	1.59	0.84
1:A:500:ARG:HG2	6:A:907:HOH:O	1.75	0.84
1:H:68:ARG:NH2	1:H:95:TYR:O	2.12	0.81
1:D:528:ARG:HD2	1:D:529:PRO:O	1.81	0.81
1:F:115:LEU:HD21	1:F:511:LEU:CB	2.09	0.81
1:F:115:LEU:HD21	1:F:511:LEU:HB3	1.63	0.79
1:G:536:ILE:HG12	1:H:538[A]:ARG:HG2	1.65	0.79
1:A:418[B]:ARG:HG3	1:B:16:LEU:HD11	1.66	0.78
1:A:536:ILE:HG12	1:B:538:ARG:HG2	1.67	0.77
1:E:418[B]:ARG:HG3	1:F:16:LEU:HD11	1.68	0.76
1:C:538:ARG:HG2	1:D:536:ILE:HG12	1.72	0.72
1:F:115:LEU:CD2	1:F:511:LEU:HB3	2.20	0.71
1:D:68:ARG:HH22	1:D:98:GLU:HB2	1.55	0.70
1:A:528:ARG:HD2	1:A:529:PRO:O	1.93	0.68
1:B:407:PHE:CE2	1:B:411:ARG:NH1	2.62	0.68
1:E:411:ARG:HG3	1:E:426:ILE:HD11	1.75	0.68
1:A:62:THR:HG21	4:A:603:I7K:O	1.94	0.67
1:D:68:ARG:NH2	1:D:95:TYR:O	2.27	0.66
1:F:115:LEU:HD21	1:F:511:LEU:HB2	1.77	0.66
1:D:536:ILE:HD11	6:D:915:HOH:O	1.95	0.65
1:H:267:ARG:HG3	6:H:958:HOH:O	1.95	0.65
1:F:407:PHE:CE2	1:F:411:ARG:NH1	2.65	0.65
1:H:68:ARG:HH22	1:H:98:GLU:HB2	1.60	0.65
1:E:418[A]:ARG:HD3	1:F:16:LEU:HD11	1.78	0.63
1:E:536:ILE:HG12	1:F:538:ARG:HG2	1.80	0.63
1:E:538:ARG:HG2	1:F:536:ILE:HG12	1.81	0.62
1:G:411:ARG:HG3	1:G:426:ILE:HD11	1.80	0.62
1:A:87:ASN:ND2	4:A:603:I7K:C14	2.63	0.62
1:C:538:ARG:HD3	6:C:860:HOH:O	1.98	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:87:ASN:HD22	4:A:603:I7K:C14	2.13	0.61
1:D:68:ARG:NH2	1:D:98:GLU:HB2	2.16	0.60
1:C:411:ARG:HG2	1:C:426:ILE:HD11	1.82	0.60
1:A:407:PHE:CE2	1:A:411:ARG:NH1	2.70	0.59
1:D:528:ARG:CD	1:D:529:PRO:O	2.51	0.58
1:H:311:ILE:CD1	1:H:311:ILE:CB	2.75	0.56
1:H:411:ARG:HG3	1:H:412:ARG:N	2.22	0.55
1:C:259:LYS:HB3	1:C:291:ARG:HH11	1.72	0.54
1:B:411:ARG:HG2	1:B:426:ILE:HD11	1.87	0.54
1:B:263:VAL:HG21	1:B:298:VAL:HG23	1.90	0.54
6:G:1027:HOH:O	1:H:411:ARG:HD2	2.08	0.54
4:F:603:I7K:O1	4:F:603:I7K:C14	2.56	0.53
1:H:71[B]:GLU:CD	1:H:71[B]:GLU:H	2.13	0.52
1:A:517:VAL:HG22	1:A:543:SER:HB2	1.92	0.52
1:F:28:GLN:HG3	1:F:30:LEU:HG	1.91	0.52
1:G:245:VAL:CG1	1:G:273:GLU:HG2	2.40	0.51
1:E:28:GLN:HG3	1:E:30:LEU:HG	1.93	0.50
1:A:272:PRO:HA	1:A:275:HIS:CE1	2.46	0.49
1:B:28:GLN:HG3	1:B:30:LEU:HG	1.93	0.49
1:H:28:GLN:HG3	1:H:30:LEU:HG	1.94	0.49
1:D:28:GLN:HG3	1:D:30:LEU:HG	1.93	0.49
1:G:28:GLN:HG3	1:G:30:LEU:HG	1.94	0.49
1:F:67:SER:HA	1:F:72:ARG:HG2	1.95	0.49
1:C:28:GLN:HG3	1:C:30:LEU:HG	1.94	0.49
1:A:538:ARG:HG3	1:B:536:ILE:HG12	1.95	0.49
1:A:28:GLN:HG3	1:A:30:LEU:HG	1.94	0.48
1:A:418[A]:ARG:HD3	1:B:16:LEU:HD11	1.94	0.48
1:A:411:ARG:NH1	1:A:430:GLU:OE2	2.46	0.48
1:H:56:SER:HB2	1:H:480:GLY:HA2	1.96	0.48
1:A:284:GLU:HG2	1:A:305:ALA:HB3	1.95	0.47
1:B:407:PHE:CD2	1:B:411:ARG:NH1	2.82	0.47
1:A:517:VAL:HG13	1:A:543:SER:HB3	1.97	0.46
1:G:56:SER:HB2	1:G:480:GLY:HA2	1.96	0.46
1:G:272:PRO:HD2	1:G:273:GLU:OE1	2.16	0.46
1:G:430[B]:GLU:OE2	1:H:430[B]:GLU:OE1	2.34	0.46
1:A:67:SER:HA	1:A:72:ARG:HG2	1.98	0.45
1:A:436:CYS:SG	1:B:416:LEU:HD13	2.57	0.44
1:C:87:ASN:HB3	4:C:603:I7K:O2	2.19	0.43
1:C:411:ARG:CG	1:C:426:ILE:HD11	2.49	0.43
1:D:55:ARG:HB2	1:D:395:ARG:HG3	2.01	0.43
1:H:231:PRO:O	1:H:258:ARG:NH2	2.51	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:486:TYR:CZ	1:A:488[B]:GLU:HB2	2.54	0.42
1:B:56:SER:HB2	1:B:480:GLY:HA2	2.01	0.42
1:D:56:SER:HB2	1:D:480:GLY:HA2	2.01	0.42
1:H:290:LYS:HA	1:H:290:LYS:HD2	1.86	0.42
1:F:56:SER:HB2	1:F:480:GLY:HA2	2.02	0.42
1:G:428:ALA:HB2	1:G:537:MET:HG3	2.02	0.42
1:H:56:SER:HB2	1:H:480:GLY:CA	2.50	0.42
1:E:331:LEU:HD11	1:E:413:ALA:HB1	2.02	0.41
1:D:419:ASP:OD2	1:D:448:ARG:NH2	2.53	0.41
1:G:284:GLU:HG2	1:G:305:ALA:HB3	2.02	0.41
1:E:56:SER:HB2	1:E:480:GLY:HA2	2.02	0.41
1:H:311:ILE:HD13	1:H:311:ILE:HG21	2.02	0.41
1:B:382:PHE:HB3	1:B:385:GLU:HB2	2.02	0.41
1:C:56:SER:HB2	1:C:480:GLY:HA2	2.02	0.41
1:H:311:ILE:CD1	1:H:311:ILE:HG21	2.50	0.41
1:C:67:SER:HA	1:C:72:ARG:HG2	2.02	0.41
1:C:374:SER:HB3	4:C:603:I7K:C14	2.51	0.41
1:G:55:ARG:HB2	1:G:395:ARG:HG3	2.02	0.41
1:G:452:LEU:O	1:G:455:ARG:HG2	2.21	0.41
1:B:263:VAL:HG11	1:B:298:VAL:HG23	2.03	0.40
1:C:452:LEU:O	1:C:455:ARG:HG2	2.21	0.40
1:B:452:LEU:O	1:B:455:ARG:HG2	2.22	0.40
1:E:539:VAL:HG22	1:F:420:PRO:HB3	2.03	0.40
1:F:16:LEU:HD12	1:F:16:LEU:HA	1.91	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [\(i\)](#)

5.3.1 Protein backbone [\(i\)](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	423/447 (95%)	417 (99%)	5 (1%)	1 (0%)	47 29

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	B	436/447 (98%)	431 (99%)	4 (1%)	1 (0%)	47 29
1	C	422/447 (94%)	414 (98%)	7 (2%)	1 (0%)	47 29
1	D	421/447 (94%)	416 (99%)	4 (1%)	1 (0%)	47 29
1	E	425/447 (95%)	420 (99%)	4 (1%)	1 (0%)	47 29
1	F	436/447 (98%)	431 (99%)	4 (1%)	1 (0%)	47 29
1	G	424/447 (95%)	417 (98%)	6 (1%)	1 (0%)	47 29
1	H	425/447 (95%)	419 (99%)	5 (1%)	1 (0%)	47 29
All	All	3412/3576 (95%)	3365 (99%)	39 (1%)	8 (0%)	47 29

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	340	THR
1	B	340	THR
1	C	340	THR
1	D	340	THR
1	E	340	THR
1	F	340	THR
1	G	340	THR
1	H	340	THR

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	346/352 (98%)	338 (98%)	8 (2%)	50 30
1	B	352/352 (100%)	347 (99%)	5 (1%)	67 51
1	C	341/352 (97%)	336 (98%)	5 (2%)	65 48
1	D	342/352 (97%)	339 (99%)	3 (1%)	78 69
1	E	348/352 (99%)	341 (98%)	7 (2%)	55 36
1	F	352/352 (100%)	342 (97%)	10 (3%)	43 22

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	G	343/352 (97%)	339 (99%)	4 (1%)	71	57
1	H	345/352 (98%)	341 (99%)	4 (1%)	71	57
All	All	2769/2816 (98%)	2723 (98%)	46 (2%)	62	43

All (46) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	72	ARG
1	A	258	ARG
1	A	273	GLU
1	A	411	ARG
1	A	412	ARG
1	A	475	VAL
1	A	537[A]	MET
1	A	537[B]	MET
1	B	26	GLN
1	B	68	ARG
1	B	412	ARG
1	B	537[A]	MET
1	B	537[B]	MET
1	C	68	ARG
1	C	291	ARG
1	C	412	ARG
1	C	487	ARG
1	C	537	MET
1	D	68	ARG
1	D	412	ARG
1	D	528	ARG
1	E	242	ARG
1	E	331	LEU
1	E	379	LYS
1	E	500	ARG
1	E	537[A]	MET
1	E	537[B]	MET
1	E	539	VAL
1	F	26	GLN
1	F	68	ARG
1	F	72	ARG
1	F	94	GLU
1	F	118	ARG
1	F	291	ARG

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Mol	Chain	Res	Type
1	F	412	ARG
1	F	511	LEU
1	F	537[A]	MET
1	F	537[B]	MET
1	G	68	ARG
1	G	273	GLU
1	G	412	ARG
1	G	436	CYS
1	H	68	ARG
1	H	273	GLU
1	H	412	ARG
1	H	537	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	470	GLN
1	B	236	GLN
1	E	236	GLN
1	G	470	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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 33 ligands modelled in this entry, 10 are monoatomic - leaving 23 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
2	FBP	G	601	-	18,20,20	0.65	0	23,32,32	0.93	2 (8%)
3	OXL	F	602	5	0,5,5	-	-	0,6,6	-	-
4	I7K	F	603	-	30,33,33	0.25	0	46,51,51	0.79	2 (4%)
3	OXL	A	602	5	0,5,5	-	-	0,6,6	-	-
2	FBP	H	601	-	18,20,20	0.75	0	23,32,32	0.82	1 (4%)
4	I7K	D	603	-	30,33,33	0.26	0	46,51,51	0.53	1 (2%)
2	FBP	D	601	-	18,20,20	0.63	1 (5%)	23,32,32	0.76	0
2	FBP	E	601	-	18,20,20	0.64	0	23,32,32	0.70	1 (4%)
4	I7K	H	603	-	30,33,33	0.44	0	46,51,51	0.65	2 (4%)
4	I7K	C	603	-	30,33,33	0.28	0	46,51,51	0.82	3 (6%)
2	FBP	B	601	-	18,20,20	0.83	1 (5%)	23,32,32	0.75	1 (4%)
3	OXL	B	602	5	0,5,5	-	-	0,6,6	-	-
3	OXL	H	602	5	0,5,5	-	-	0,6,6	-	-
3	OXL	D	602	5	0,5,5	-	-	0,6,6	-	-
3	OXL	E	602	5	0,5,5	-	-	0,6,6	-	-
4	I7K	A	603	-	30,33,33	0.30	0	46,51,51	0.43	0
2	FBP	C	601	-	18,20,20	0.71	1 (5%)	23,32,32	0.81	0
4	I7K	G	603	-	30,33,33	0.34	0	46,51,51	0.50	0
3	OXL	G	602	5	0,5,5	-	-	0,6,6	-	-
3	OXL	C	602	5	0,5,5	-	-	0,6,6	-	-
4	I7K	E	603	-	30,33,33	0.27	0	46,51,51	0.70	0
2	FBP	A	601	-	18,20,20	0.60	0	23,32,32	0.63	0
2	FBP	F	601	-	18,20,20	0.41	0	23,32,32	0.59	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
2	FBP	G	601	-	-	3/13/32/32	0/1/1/1
3	OXL	F	602	5	-	0/0/4/4	-
4	I7K	F	603	-	-	3/12/42/42	0/4/4/4
3	OXL	A	602	5	-	0/0/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	FBP	H	601	-	-	2/13/32/32	0/1/1/1
4	I7K	D	603	-	-	4/12/42/42	0/4/4/4
2	FBP	D	601	-	-	2/13/32/32	0/1/1/1
2	FBP	E	601	-	-	2/13/32/32	0/1/1/1
4	I7K	H	603	-	-	3/12/42/42	0/4/4/4
4	I7K	C	603	-	-	4/12/42/42	0/4/4/4
2	FBP	B	601	-	-	2/13/32/32	0/1/1/1
3	OXL	B	602	5	-	0/0/4/4	-
3	OXL	H	602	5	-	0/0/4/4	-
3	OXL	D	602	5	-	0/0/4/4	-
3	OXL	E	602	5	-	0/0/4/4	-
4	I7K	A	603	-	-	7/12/42/42	0/4/4/4
2	FBP	C	601	-	-	2/13/32/32	0/1/1/1
4	I7K	G	603	-	-	4/12/42/42	0/4/4/4
3	OXL	G	602	5	-	0/0/4/4	-
3	OXL	C	602	5	-	0/0/4/4	-
4	I7K	E	603	-	-	3/12/42/42	0/4/4/4
2	FBP	A	601	-	-	2/13/32/32	0/1/1/1
2	FBP	F	601	-	-	2/13/32/32	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	601	FBP	P1-O3P	-2.43	1.45	1.54
2	D	601	FBP	P2-O5P	-2.09	1.46	1.54
2	B	601	FBP	P2-O5P	-2.08	1.46	1.54

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	F	603	I7K	C3-C-C1	-3.79	117.77	121.14
4	C	603	I7K	C3-C-C1	-3.45	118.07	121.14
4	H	603	I7K	C19-N-S	2.85	122.23	117.05
4	C	603	I7K	C15-C14-N	2.56	112.92	109.43
4	F	603	I7K	C15-C14-N	2.26	112.51	109.43
2	B	601	FBP	O3P-P1-O2P	2.25	116.24	107.64
2	G	601	FBP	O3P-P1-O1	2.25	112.71	106.73
2	E	601	FBP	O3P-P1-O2P	2.24	116.20	107.64
4	D	603	I7K	C14-N-S	2.21	121.07	117.05

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	601	FBP	O5P-P2-O6	2.16	112.48	106.73
4	H	603	I7K	C14-N-S	2.12	120.91	117.05
2	G	601	FBP	O5P-P2-O6	2.03	112.13	106.73
4	C	603	I7K	C18-C19-N	2.01	112.16	109.43

There are no chirality outliers.

All (45) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	E	601	FBP	C4-C5-C6-O6
2	G	601	FBP	O1-C1-C2-O2
2	G	601	FBP	O1-C1-C2-C3
2	G	601	FBP	O1-C1-C2-O5
4	A	603	I7K	C19-N-S-O
4	A	603	I7K	C19-N-S-C
4	C	603	I7K	C19-N-S-O7
4	E	603	I7K	C14-N-S-O
4	C	603	I7K	C19-N-S-C
2	A	601	FBP	C4-C5-C6-O6
2	B	601	FBP	C4-C5-C6-O6
2	D	601	FBP	C4-C5-C6-O6
2	F	601	FBP	C4-C5-C6-O6
2	H	601	FBP	C4-C5-C6-O6
4	F	603	I7K	C19-N-S-O7
4	E	603	I7K	C14-N-S-C
4	F	603	I7K	C19-N-S-C
4	A	603	I7K	C19-N-S-O7
2	C	601	FBP	C4-C5-C6-O6
4	D	603	I7K	C14-N-S-O7
4	D	603	I7K	C14-N-S-O
2	A	601	FBP	O5-C5-C6-O6
2	B	601	FBP	O5-C5-C6-O6
2	C	601	FBP	O5-C5-C6-O6
2	D	601	FBP	O5-C5-C6-O6
2	E	601	FBP	O5-C5-C6-O6
2	F	601	FBP	O5-C5-C6-O6
2	H	601	FBP	O5-C5-C6-O6
4	G	603	I7K	C14-N-S-O7
4	G	603	I7K	C14-N-S-O
4	H	603	I7K	C19-N-S-O7
4	H	603	I7K	C19-N-S-O
4	D	603	I7K	C19-N-S-O

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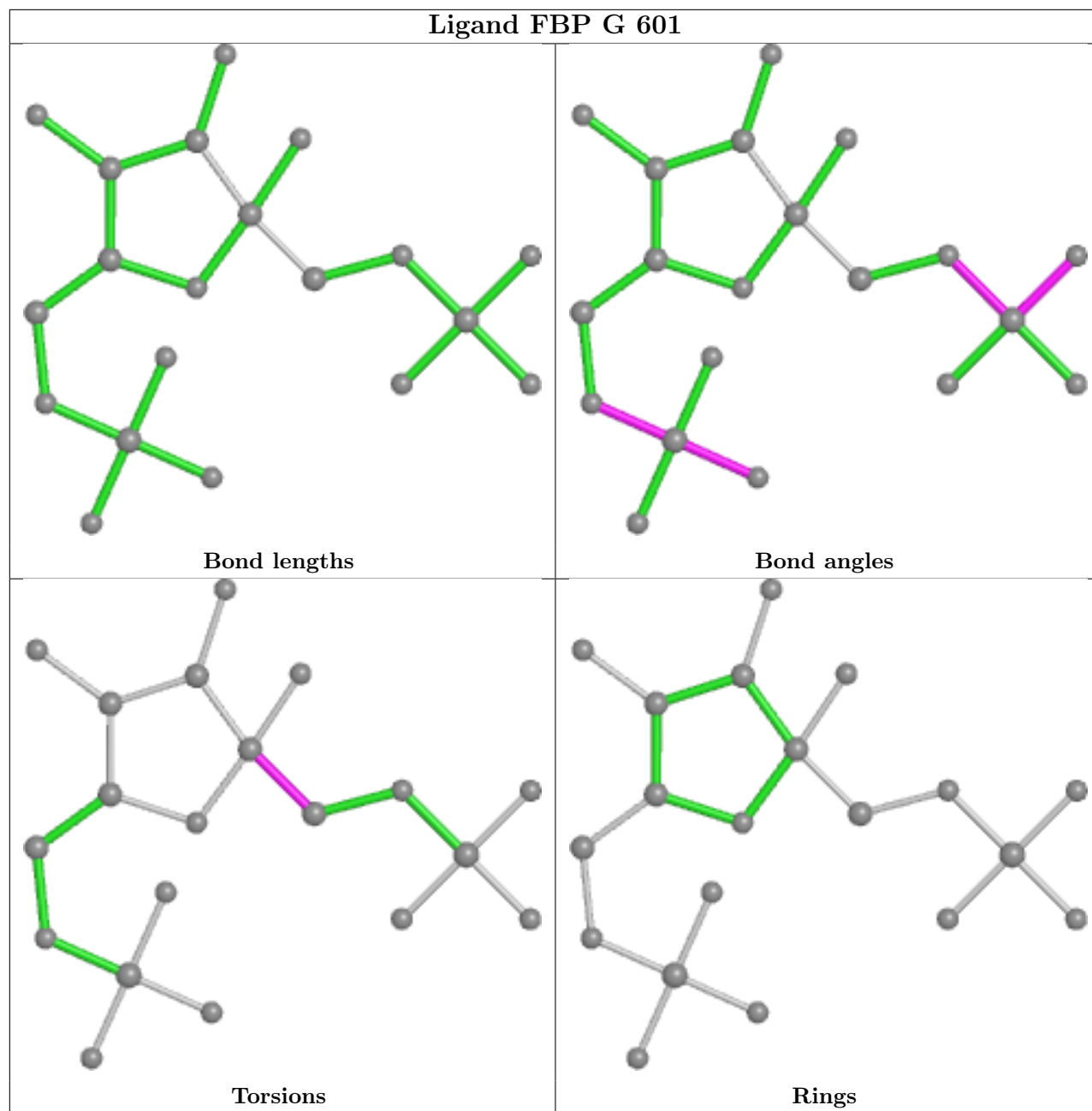
Mol	Chain	Res	Type	Atoms
4	A	603	I7K	C14-N-S-O
4	A	603	I7K	C14-N-S-O7
4	G	603	I7K	C19-N-S-O
4	D	603	I7K	C14-N-S-C
4	C	603	I7K	C19-N-S-O
4	A	603	I7K	C1-C-S-O7
4	H	603	I7K	C19-N-S-C
4	A	603	I7K	C14-N-S-C
4	G	603	I7K	C14-N-S-C
4	C	603	I7K	C1-C-S-O7
4	F	603	I7K	C1-C-S-O7
4	E	603	I7K	C1-C-S-N

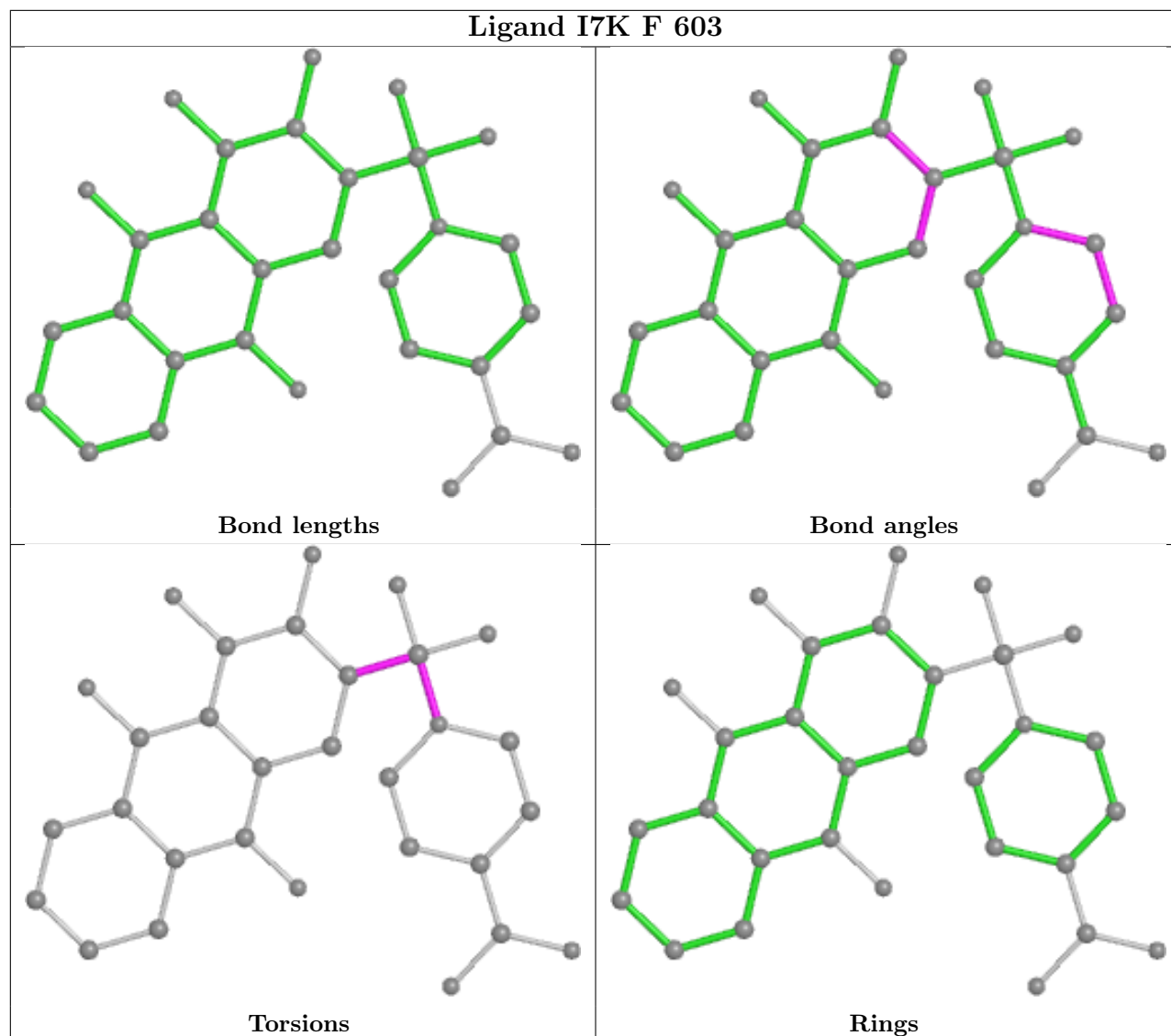
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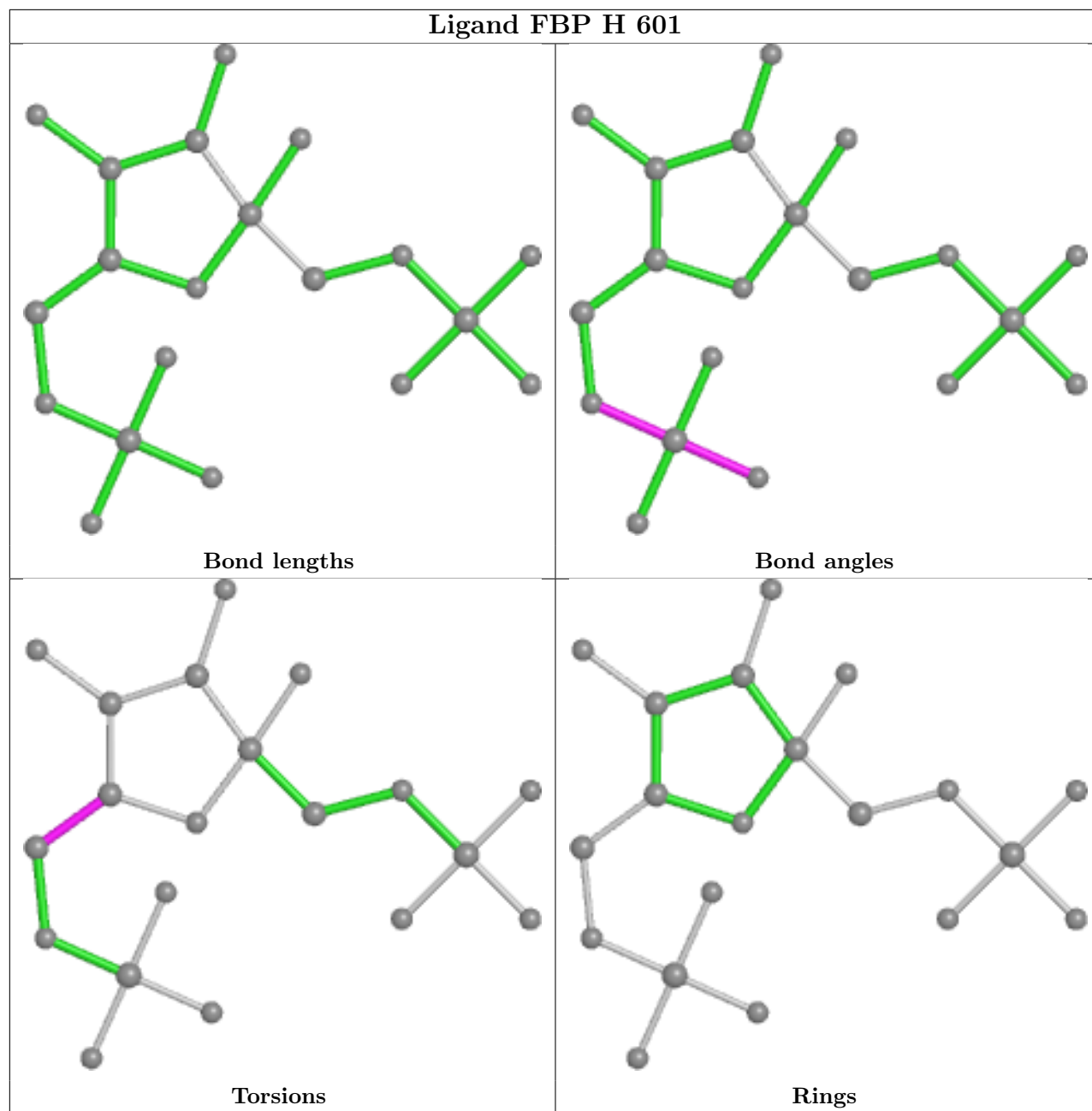
4 monomers are involved in 7 short contacts:

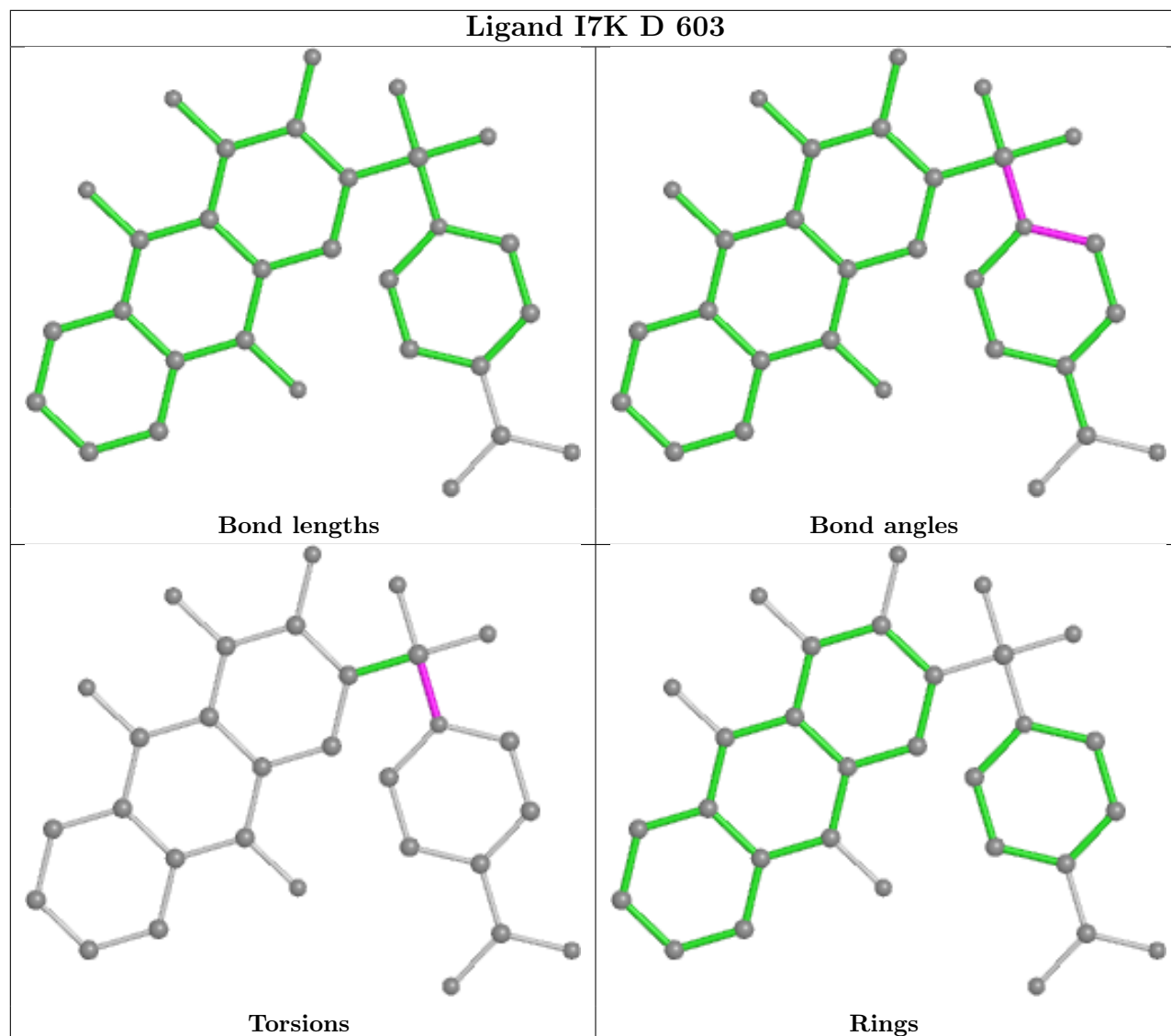
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	G	601	FBP	1	0
4	F	603	I7K	1	0
4	C	603	I7K	2	0
4	A	603	I7K	3	0

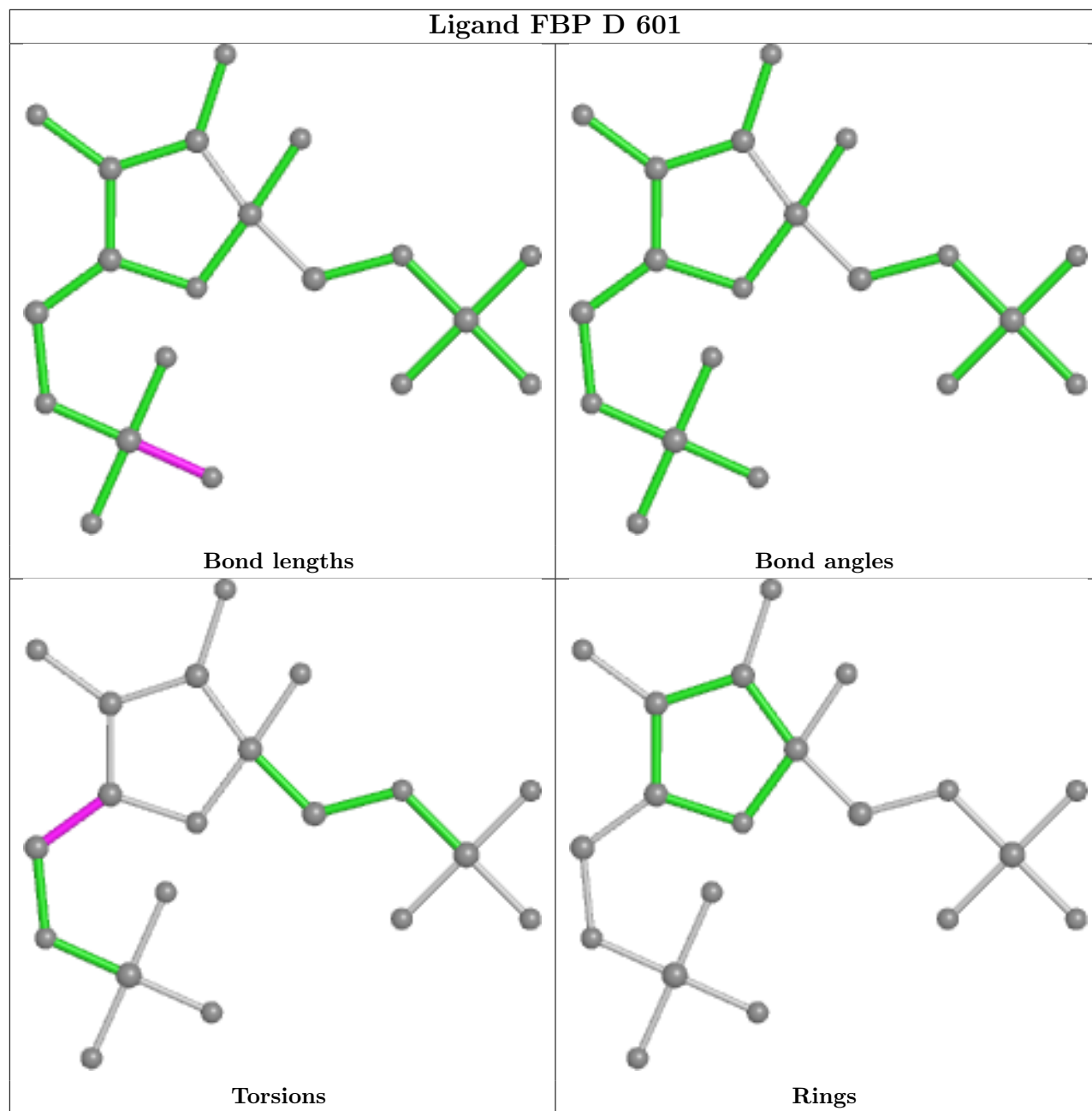
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

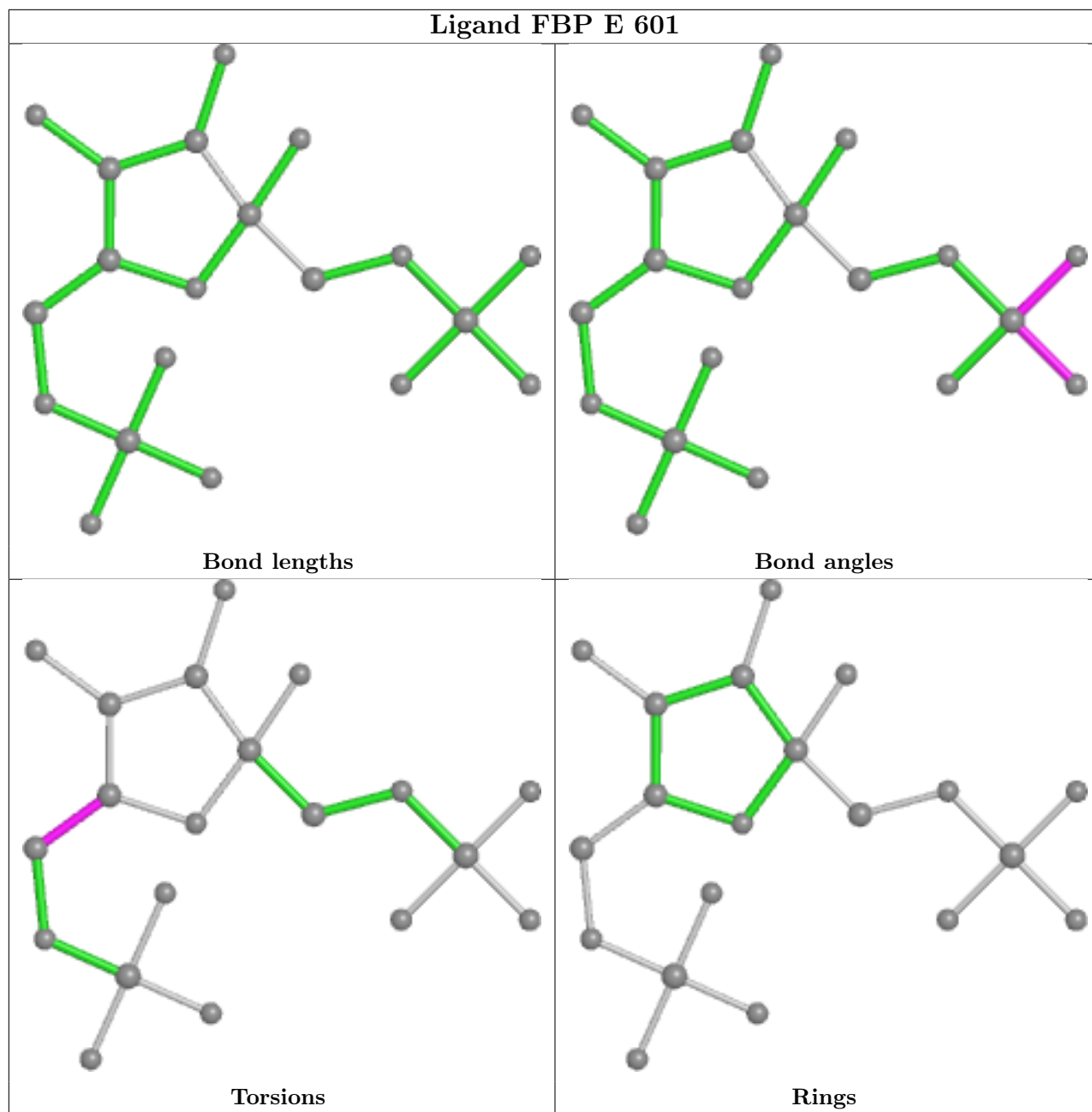


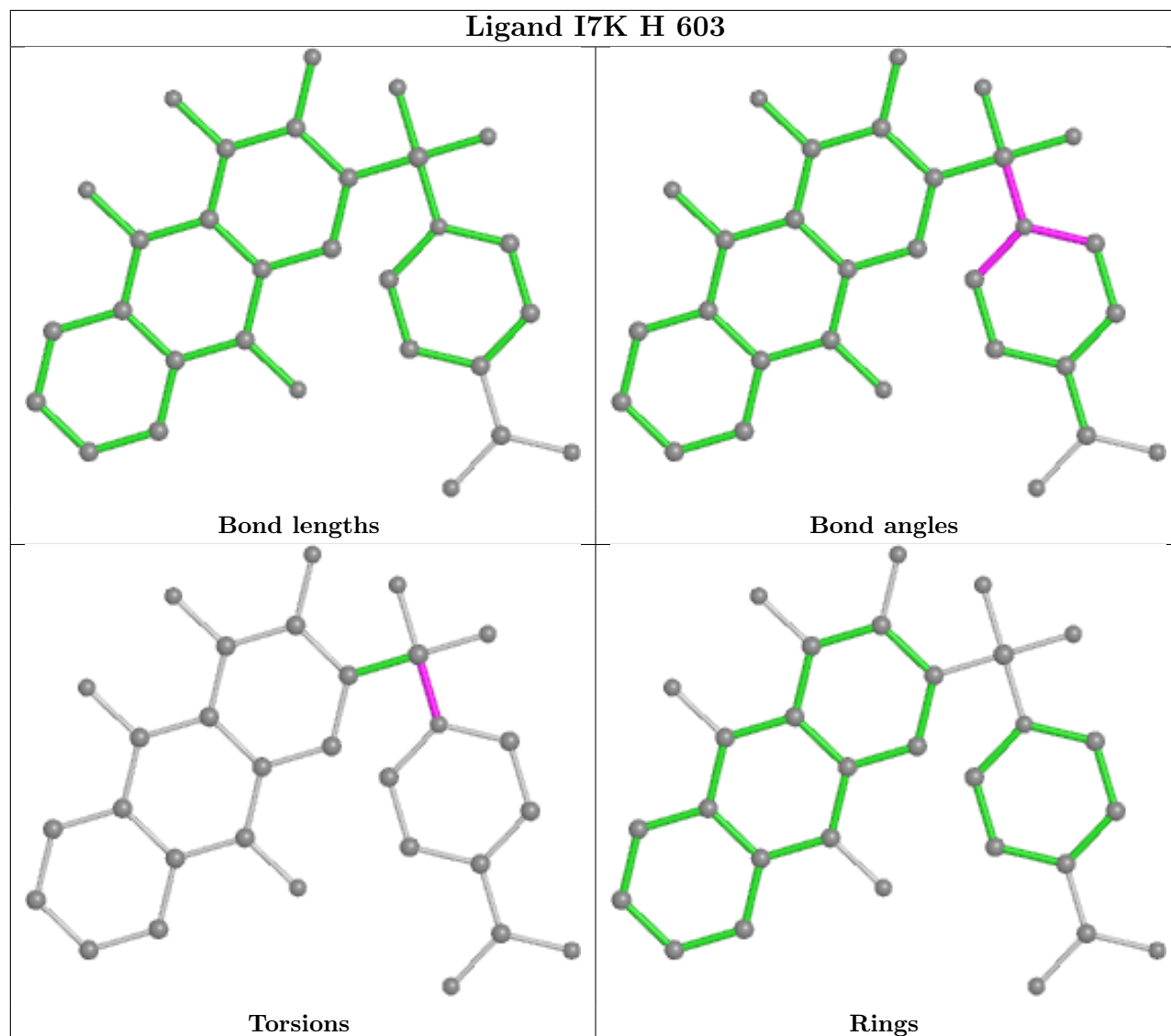


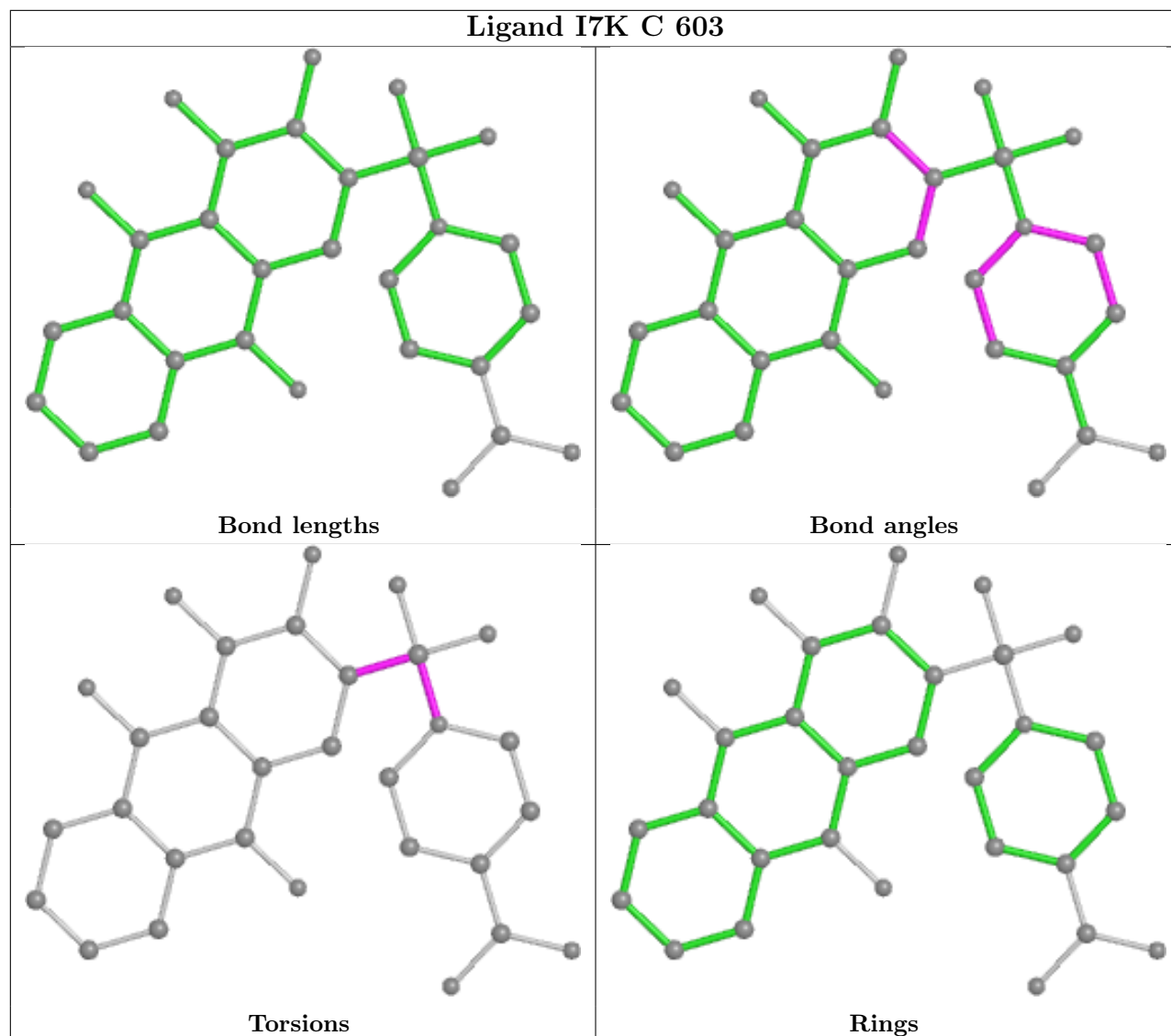


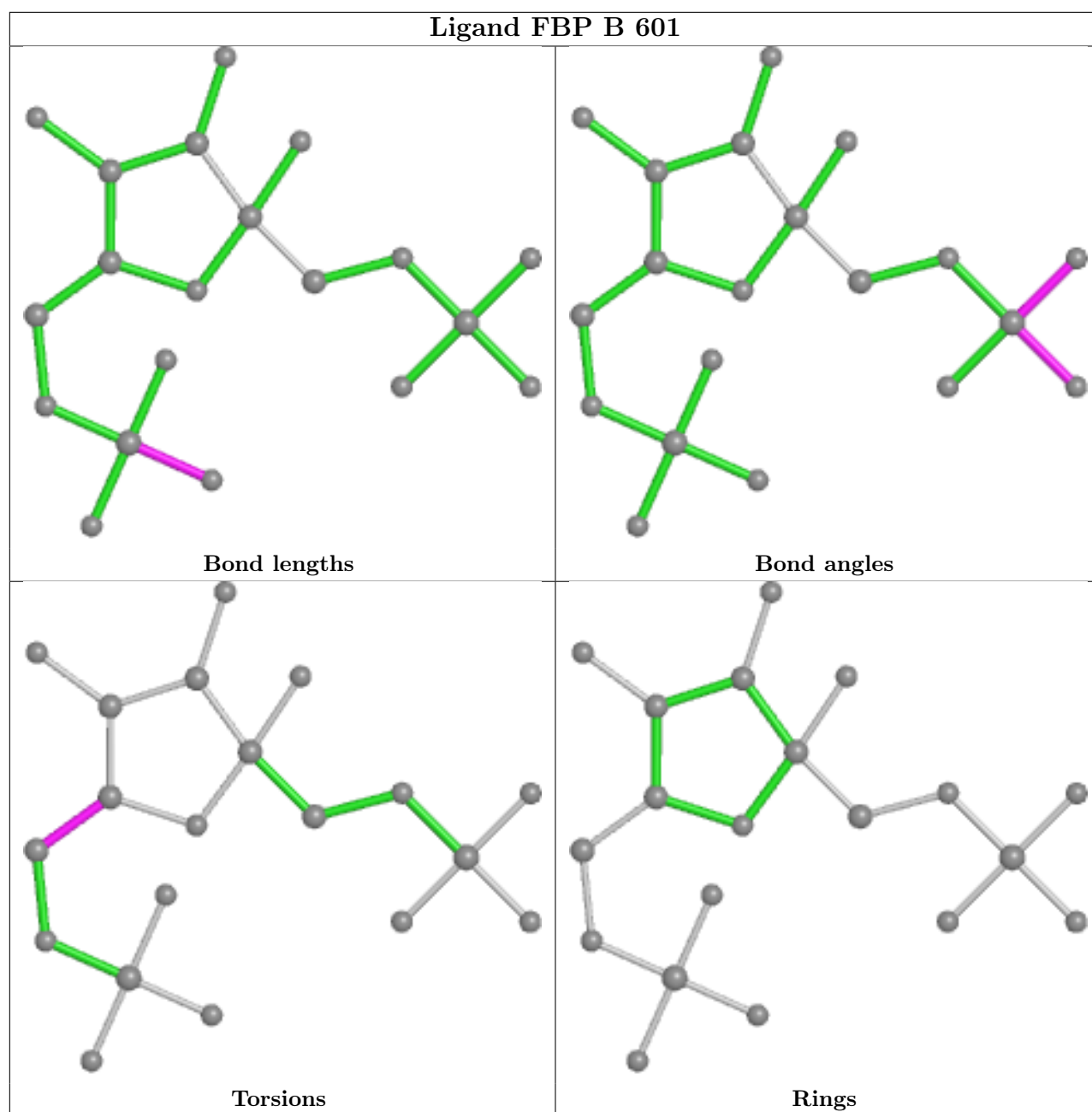


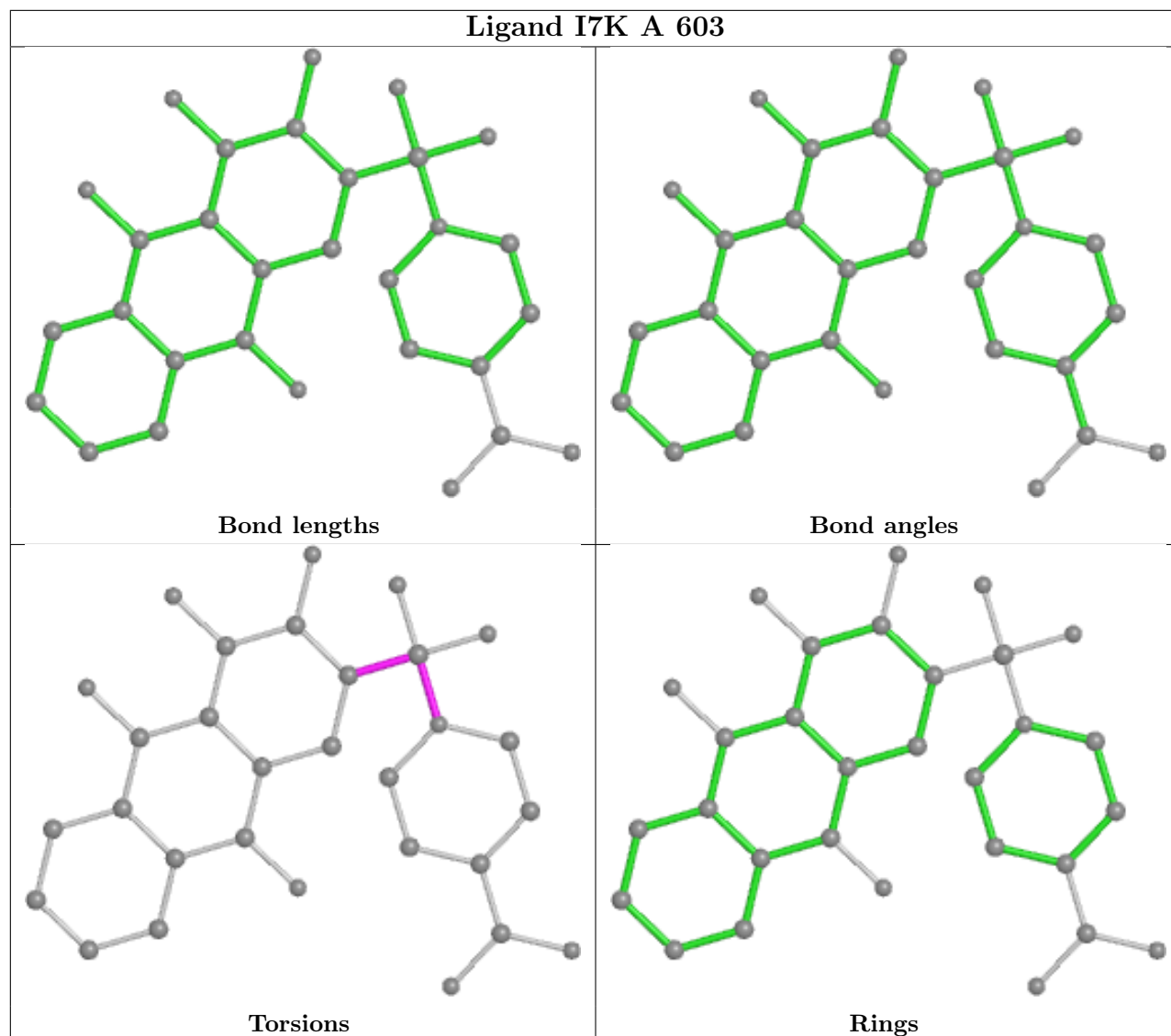


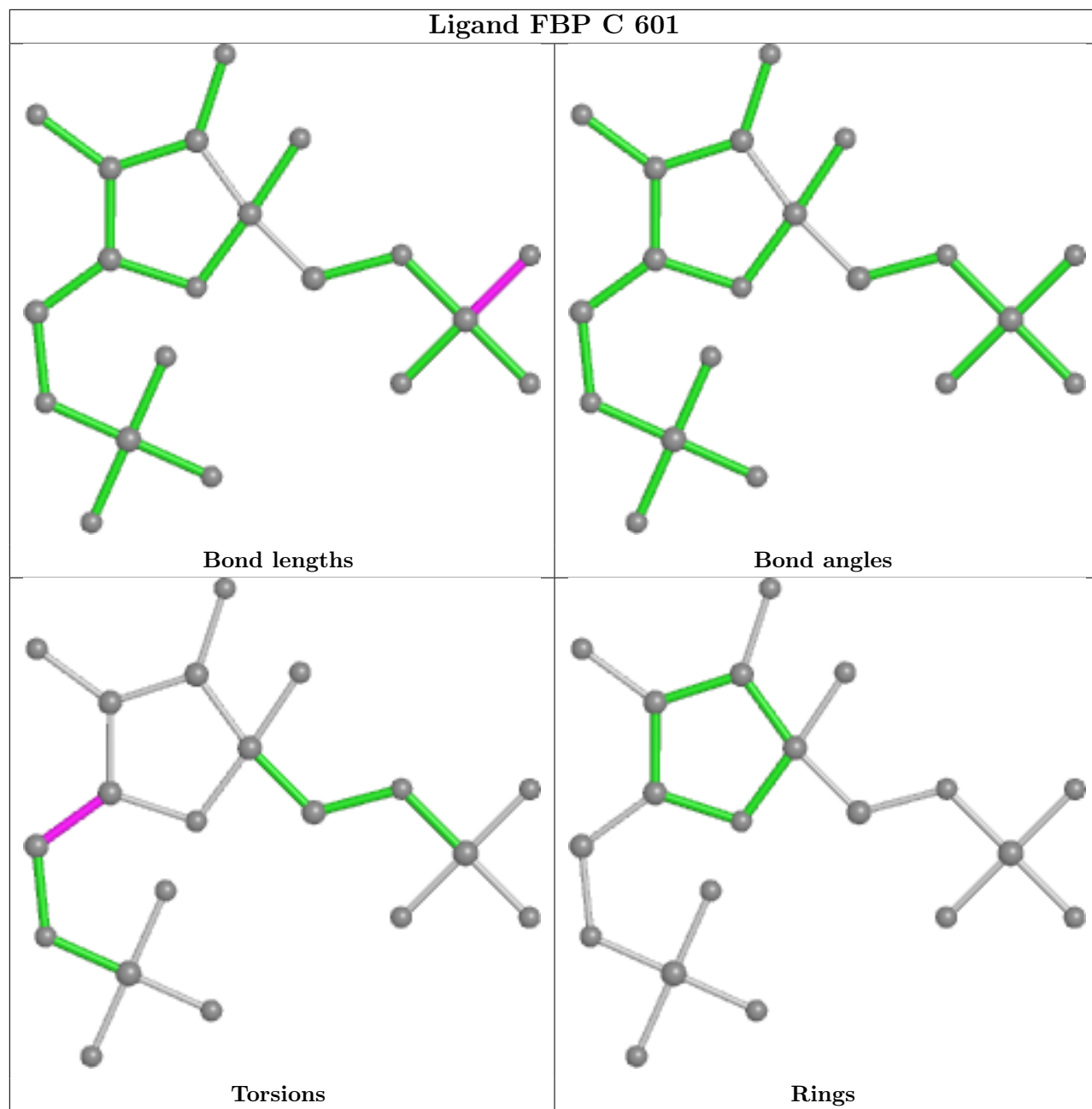


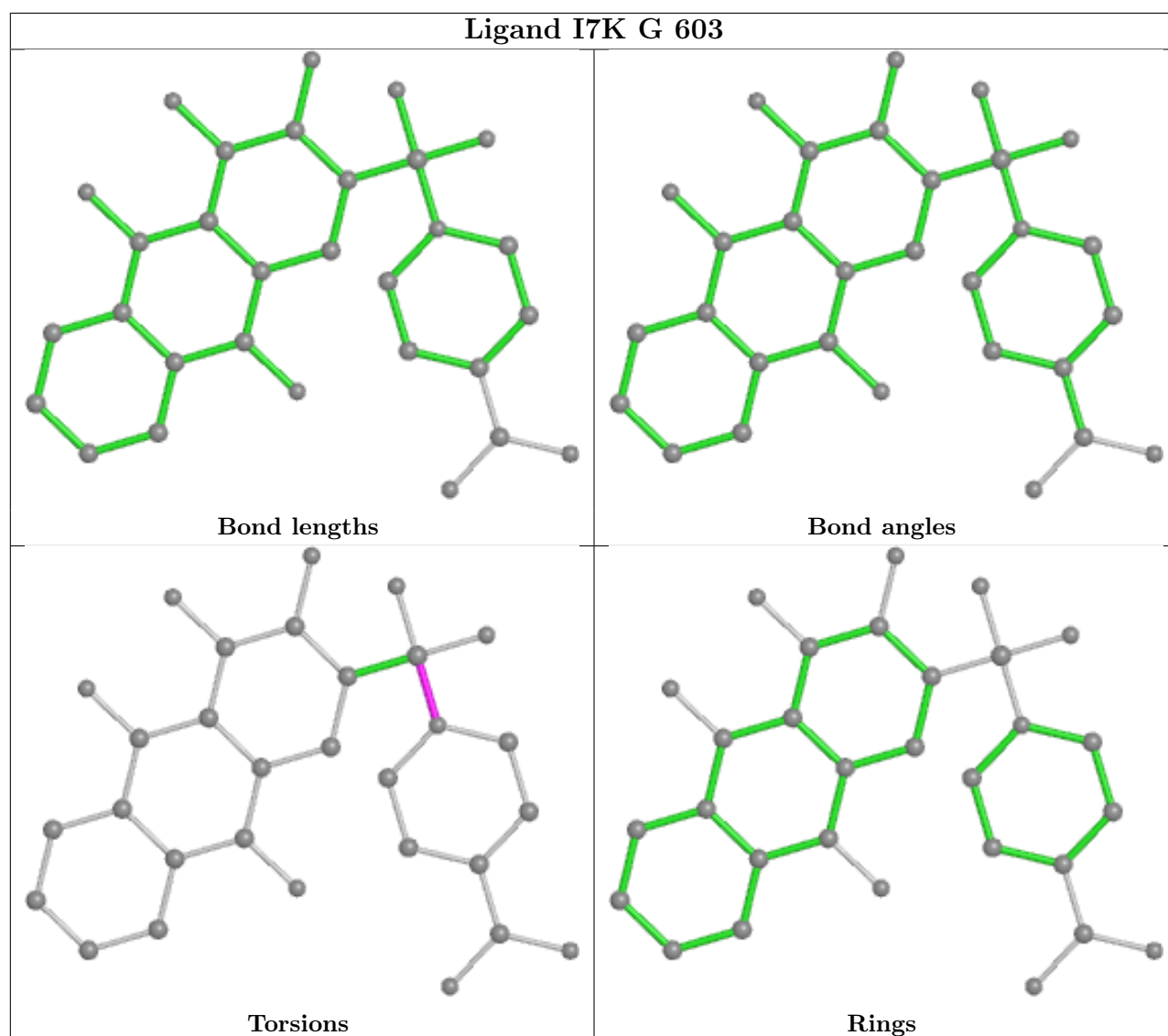


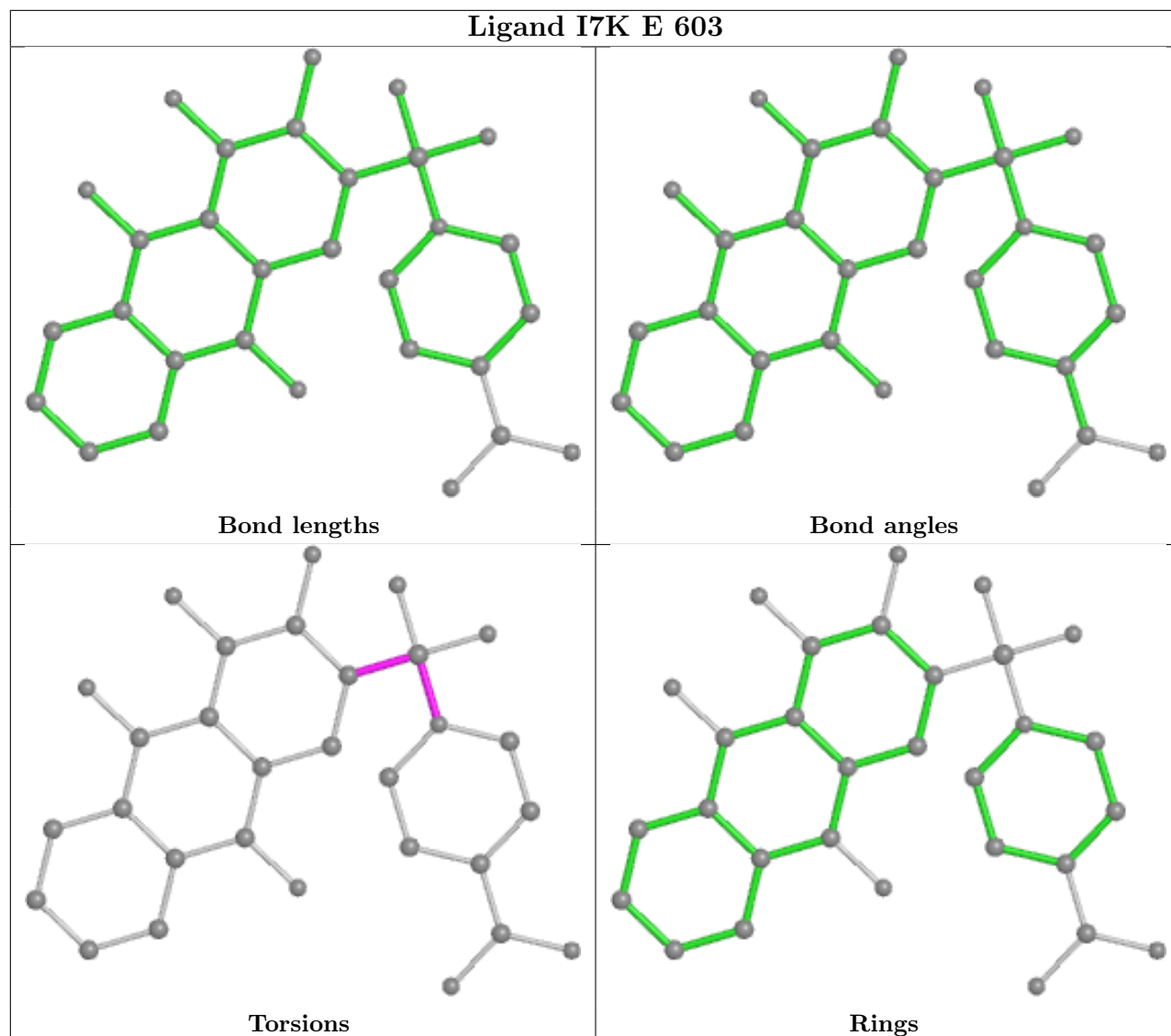


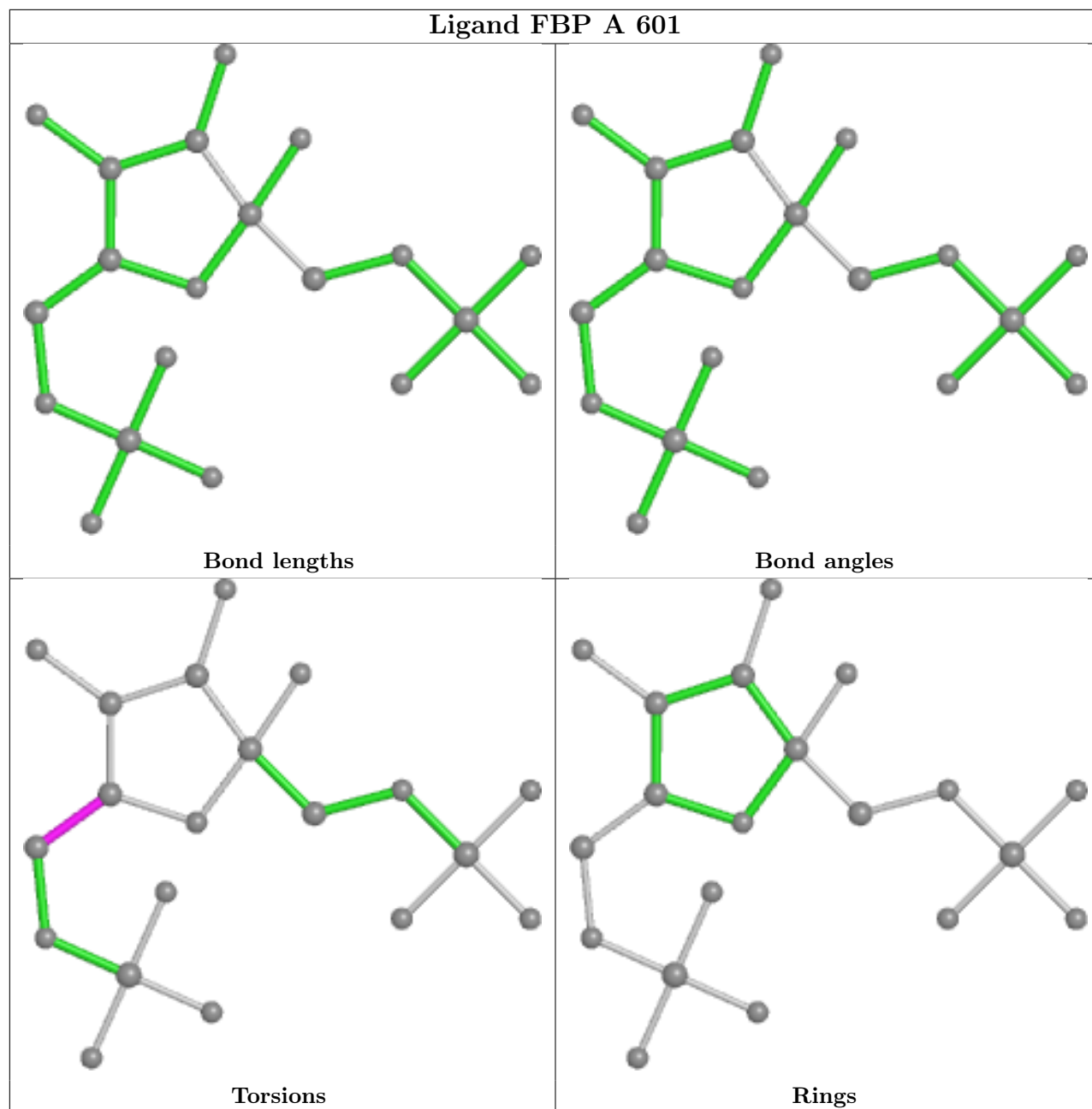


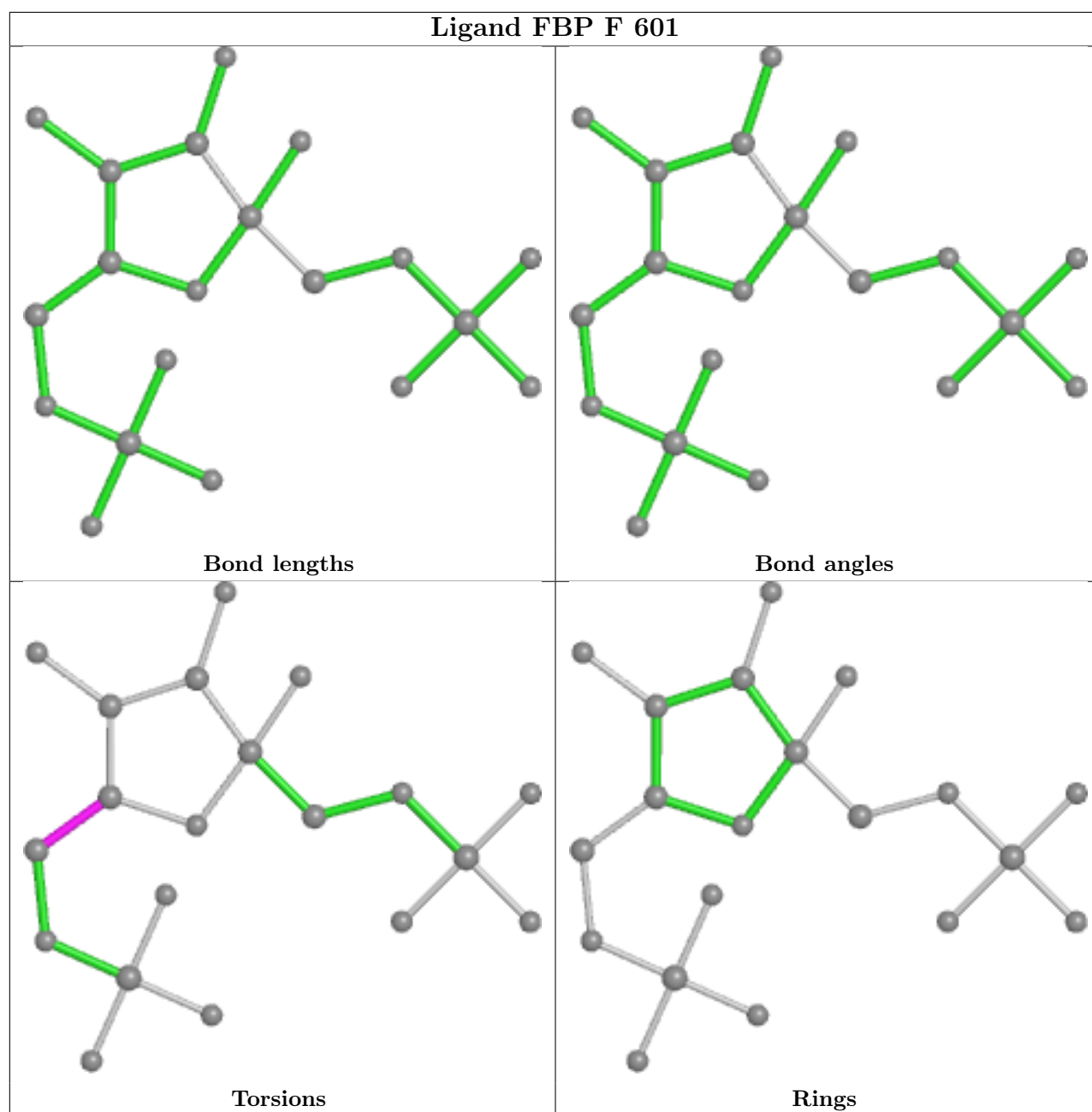












5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	414/447 (92%)	-0.02	10 (2%) 59 62	16, 26, 45, 61	0
1	B	431/447 (96%)	0.09	11 (2%) 56 58	15, 28, 49, 72	0
1	C	417/447 (93%)	-0.14	10 (2%) 59 62	11, 21, 39, 55	0
1	D	419/447 (93%)	-0.10	7 (1%) 70 74	12, 19, 38, 75	0
1	E	414/447 (92%)	-0.02	8 (1%) 66 70	16, 27, 48, 61	0
1	F	431/447 (96%)	-0.06	11 (2%) 56 58	13, 23, 44, 60	0
1	G	419/447 (93%)	-0.15	5 (1%) 79 82	12, 20, 36, 50	1 (0%)
1	H	421/447 (94%)	-0.16	10 (2%) 59 62	11, 17, 35, 57	0
All	All	3366/3576 (94%)	-0.07	72 (2%) 63 67	11, 23, 44, 75	1 (0%)

All (72) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	25	PHE	10.2
1	D	22	THR	8.2
1	D	24	PHE	7.9
1	D	21	GLY	5.4
1	A	543	SER	5.2
1	H	21	GLY	5.1
1	E	232	GLY	5.0
1	G	25	PHE	5.0
1	F	489	PRO	4.9
1	C	271	GLY	4.7
1	B	232	GLY	4.3
1	H	22	THR	4.3
1	C	231	PRO	4.3
1	H	25	PHE	4.2
1	F	115	LEU	4.2
1	A	232	GLY	3.9

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Mol	Chain	Res	Type	RSRZ
1	C	232	GLY	3.8
1	H	232	GLY	3.5
1	B	489	PRO	3.4
1	H	516	ARG	3.2
1	F	516	ARG	3.2
1	D	232	GLY	3.1
1	E	540	LEU	3.1
1	D	543	SER	3.1
1	B	115	LEU	3.1
1	F	114	PRO	2.9
1	G	231	PRO	2.9
1	H	231	PRO	2.9
1	B	275	HIS	2.8
1	C	34	MET	2.8
1	F	275	HIS	2.8
1	A	114	PRO	2.8
1	A	242	ARG	2.7
1	B	543	SER	2.7
1	H	52	VAL	2.7
1	G	26	GLN	2.6
1	F	232	GLY	2.6
1	F	231	PRO	2.5
1	C	112	GLY	2.5
1	B	231	PRO	2.4
1	E	114	PRO	2.4
1	A	30	LEU	2.4
1	E	514	PHE	2.4
1	E	412	ARG	2.4
1	C	113	SER	2.4
1	D	30	LEU	2.4
1	F	543	SER	2.4
1	F	487	ARG	2.3
1	B	490	PRO	2.3
1	H	411	ARG	2.3
1	E	543	SER	2.3
1	A	26	GLN	2.3
1	C	114	PRO	2.3
1	E	493	ILE	2.3
1	C	116[A]	SER	2.3
1	G	489	PRO	2.3
1	H	514	PHE	2.2
1	A	275	HIS	2.2

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Mol	Chain	Res	Type	RSRZ
1	C	118	ARG	2.2
1	A	412	ARG	2.2
1	F	517	VAL	2.1
1	B	416	LEU	2.1
1	H	273	GLU	2.1
1	B	118	ARG	2.1
1	A	118	ARG	2.1
1	B	487	ARG	2.1
1	B	271	GLY	2.0
1	F	12	ASP	2.0
1	A	273	GLU	2.0
1	G	412	ARG	2.0
1	C	115	LEU	2.0
1	E	271	GLY	2.0

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

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

6.3 Carbohydrates [i](#)

There are no 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(Å ²)	Q<0.9
4	I7K	C	603	30/30	0.70	0.23	54,56,58,58	17
4	I7K	A	603	30/30	0.74	0.30	70,70,72,72	17
4	I7K	F	603	30/30	0.74	0.29	51,53,55,55	17
4	I7K	D	603	30/30	0.84	0.25	28,31,36,38	17
3	OXL	B	602	6/6	0.85	0.16	31,32,34,34	0
3	OXL	E	602	6/6	0.85	0.15	29,31,32,33	0
3	OXL	A	602	6/6	0.92	0.15	28,29,30,30	0
4	I7K	E	603	30/30	0.92	0.20	40,42,47,48	17
3	OXL	F	602	6/6	0.92	0.12	24,26,27,29	0

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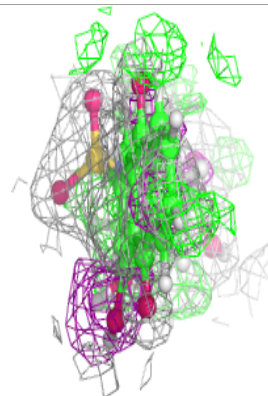
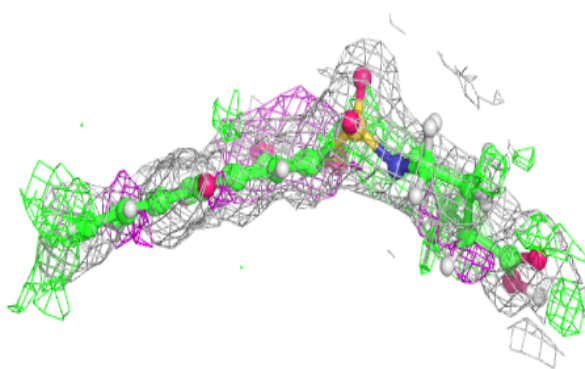
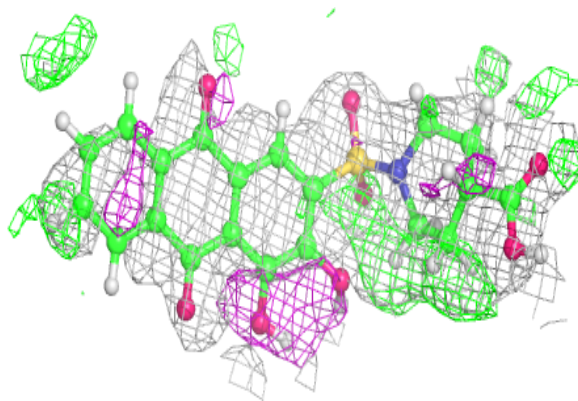
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	I7K	G	603	30/30	0.92	0.16	26,31,34,34	17
3	OXL	C	602	6/6	0.94	0.11	28,28,29,30	0
3	OXL	D	602	6/6	0.94	0.08	23,24,25,25	0
3	OXL	G	602	6/6	0.95	0.07	22,23,24,25	0
4	I7K	H	603	30/30	0.95	0.11	21,22,25,26	17
2	FBP	G	601	20/20	0.96	0.10	13,15,17,24	0
3	OXL	H	602	6/6	0.97	0.08	17,19,20,22	0
2	FBP	A	601	20/20	0.97	0.08	19,22,23,23	0
5	MG	F	605	1/1	0.97	0.11	23,23,23,23	0
2	FBP	F	601	20/20	0.98	0.08	20,23,26,26	0
2	FBP	B	601	20/20	0.98	0.08	22,23,26,27	0
5	MG	A	604	1/1	0.98	0.04	30,30,30,30	0
5	MG	B	603	1/1	0.98	0.05	29,29,29,29	0
5	MG	E	604	1/1	0.98	0.04	28,28,28,28	0
2	FBP	E	601	20/20	0.98	0.08	20,21,25,25	0
2	FBP	C	601	20/20	0.99	0.09	14,15,17,18	0
5	MG	C	604	1/1	0.99	0.04	24,24,24,24	0
5	MG	C	605	1/1	0.99	0.10	20,20,20,20	0
5	MG	D	604	1/1	0.99	0.05	22,22,22,22	0
2	FBP	H	601	20/20	0.99	0.10	11,13,15,16	0
5	MG	F	604	1/1	0.99	0.06	24,24,24,24	0
2	FBP	D	601	20/20	0.99	0.09	13,15,17,17	0
5	MG	H	604	1/1	0.99	0.05	20,20,20,20	0
5	MG	G	604	1/1	1.00	0.06	21,21,21,21	0

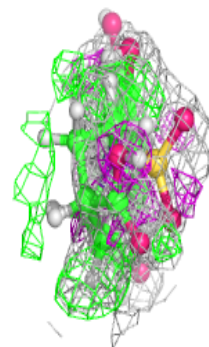
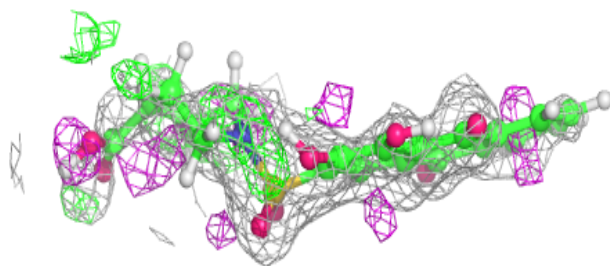
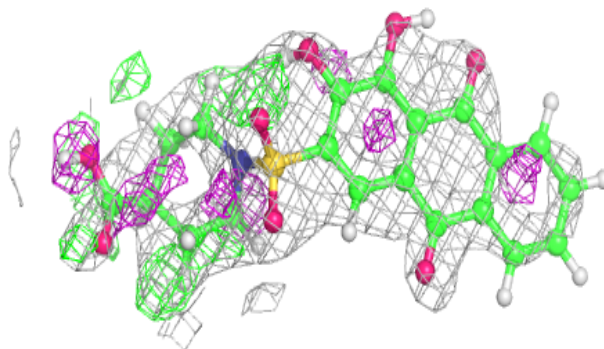
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around I7K C 603:

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

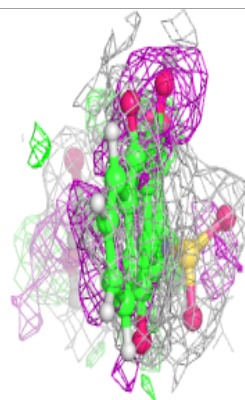
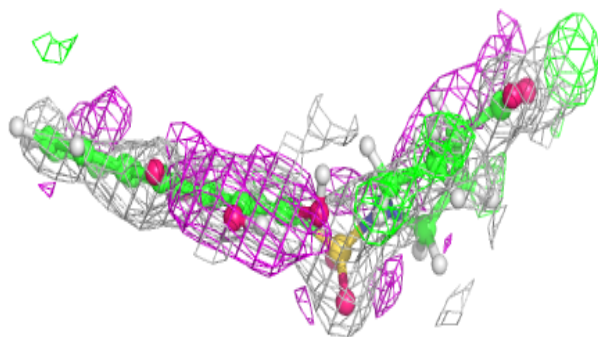
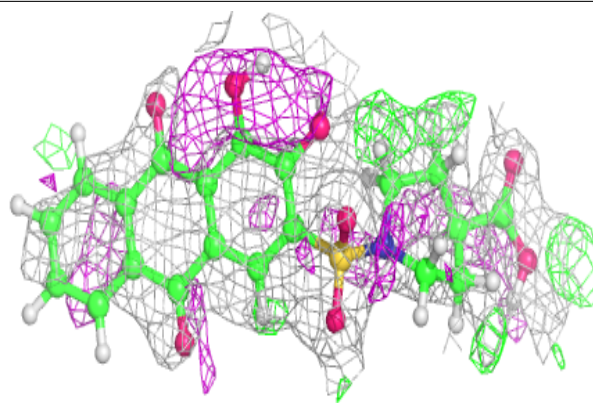
**Electron density around I7K A 603:**

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

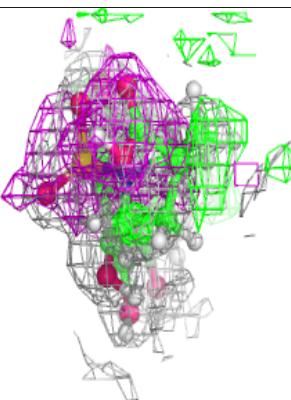
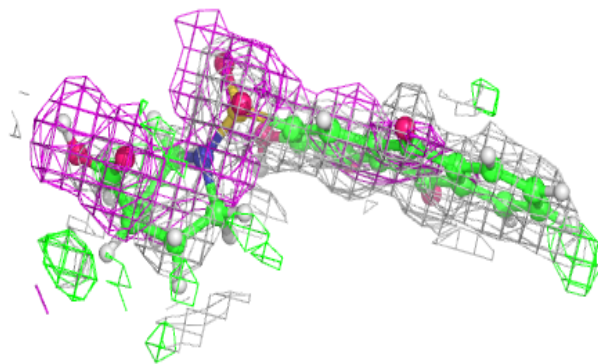
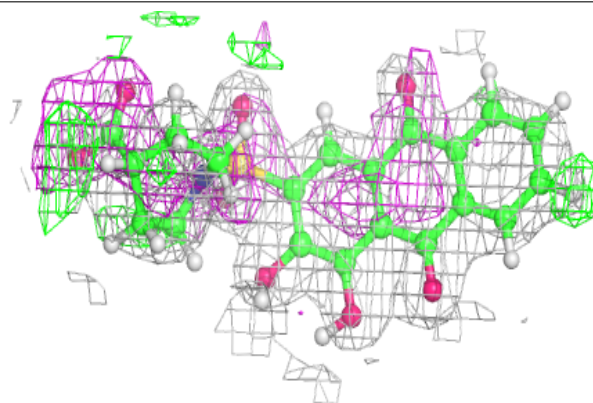


Electron density around I7K F 603:

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

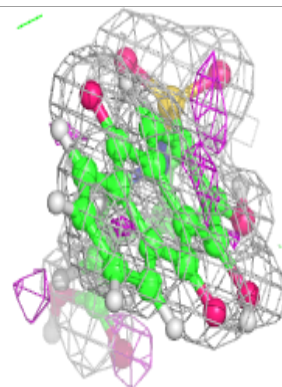
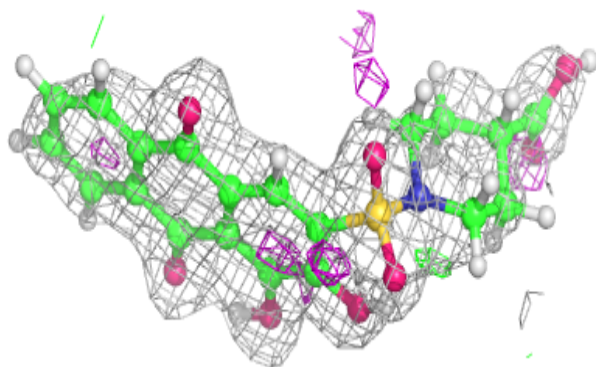
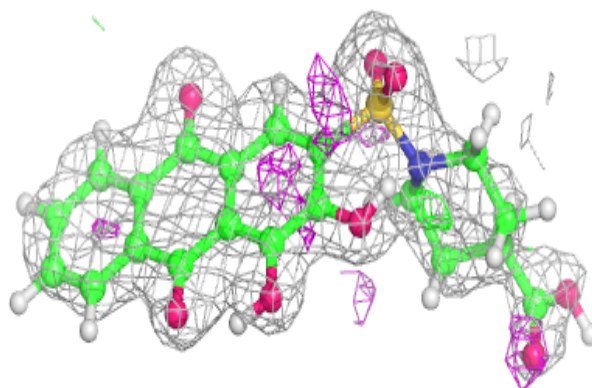
**Electron density around I7K D 603:**

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

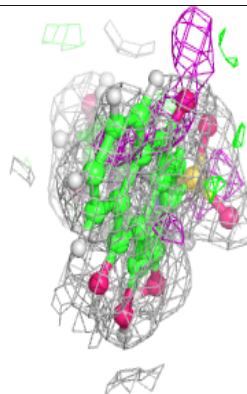
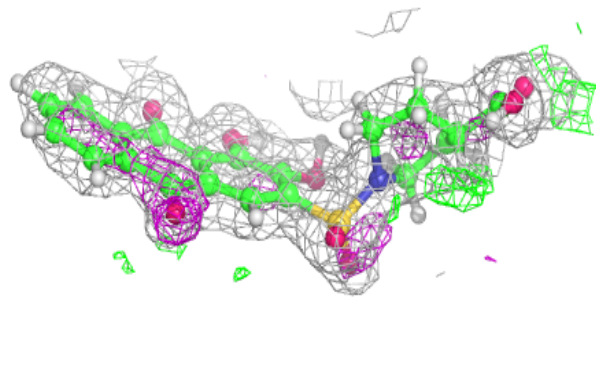
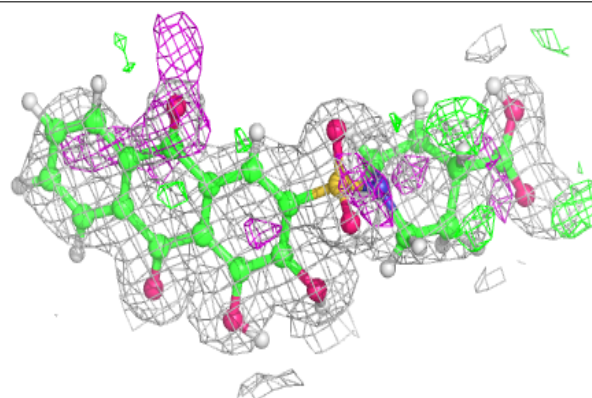


Electron density around I7K E 603:

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

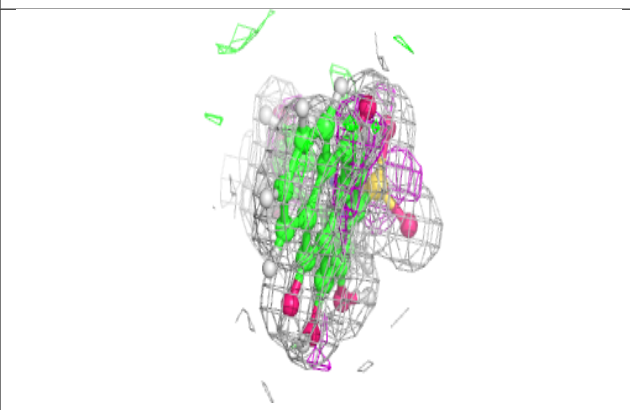
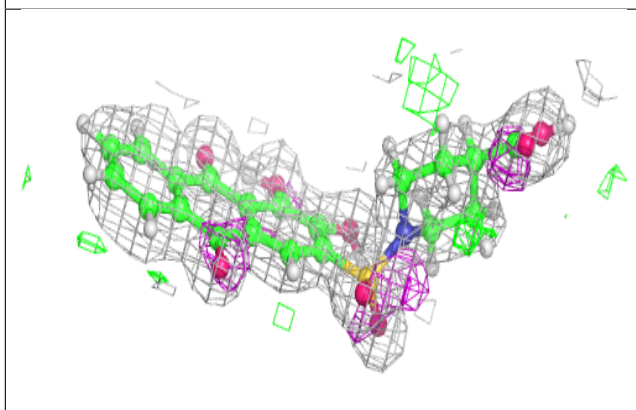
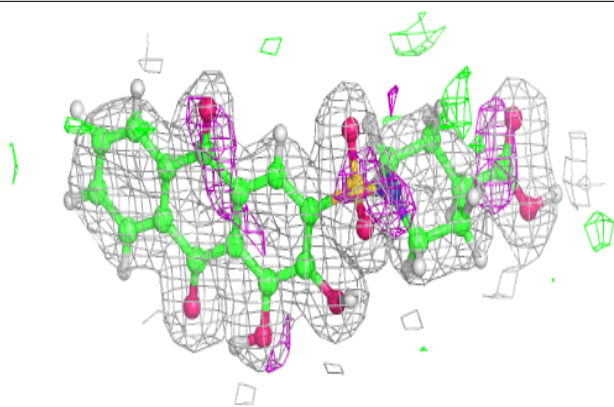
**Electron density around I7K G 603:**

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and green (positive)

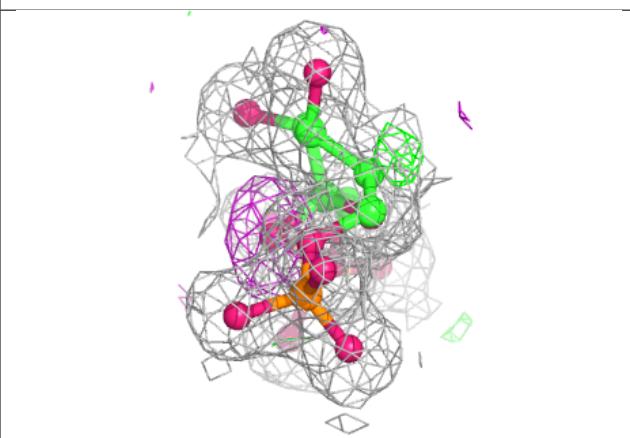
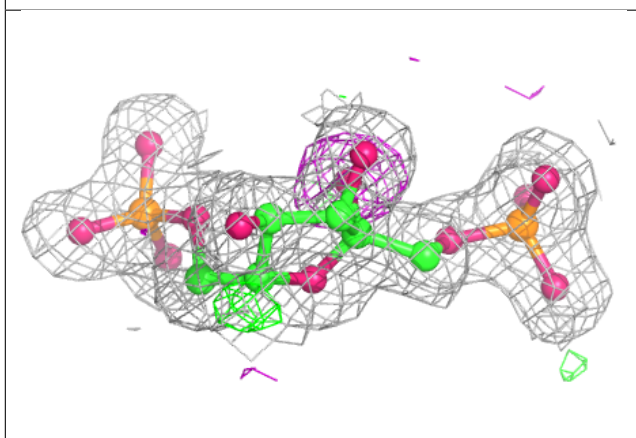
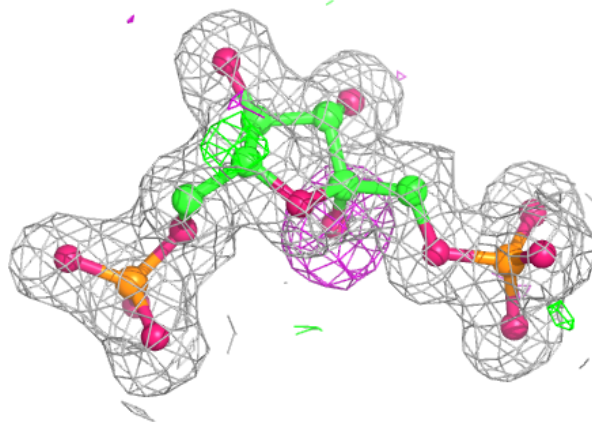


Electron density around I7K H 603:

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

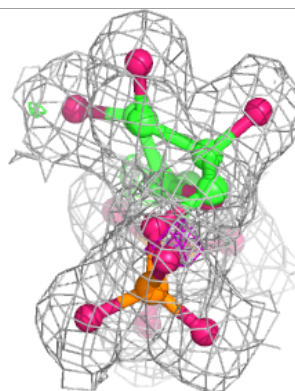
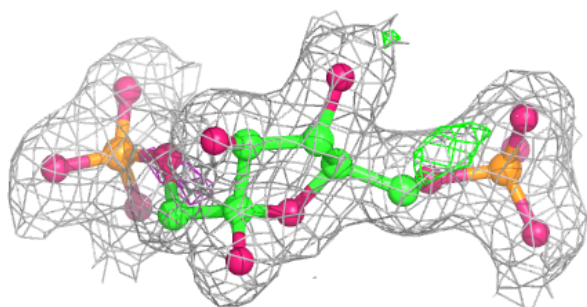
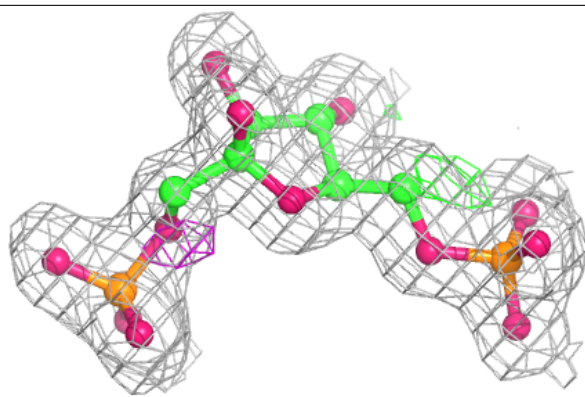
**Electron density around FBP G 601:**

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and green (positive)

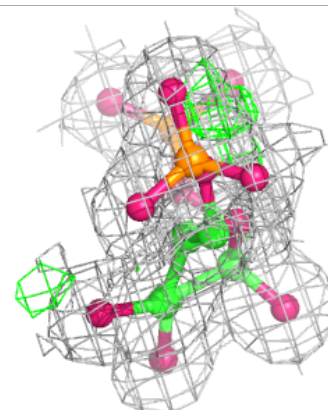
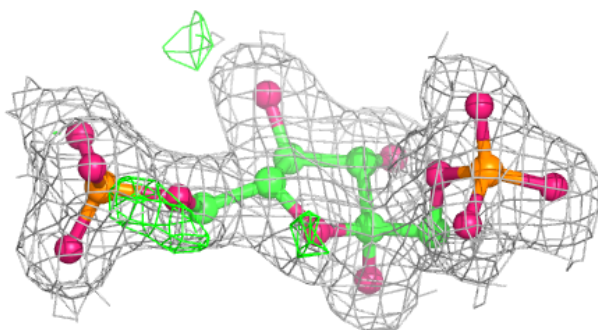
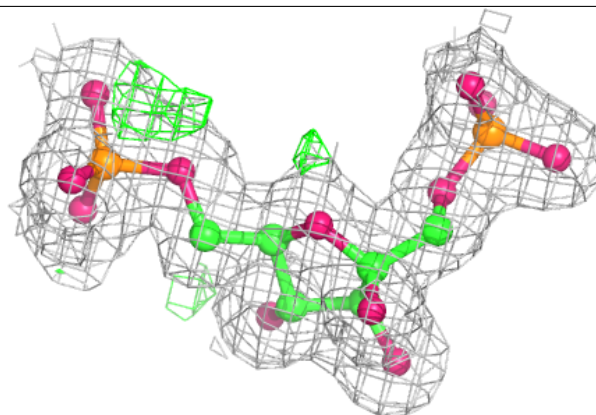


Electron density around FBP A 601:

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

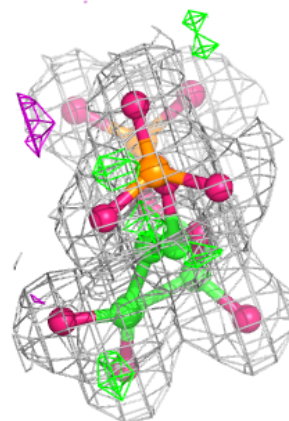
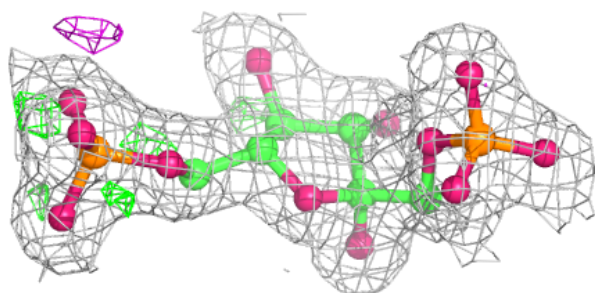
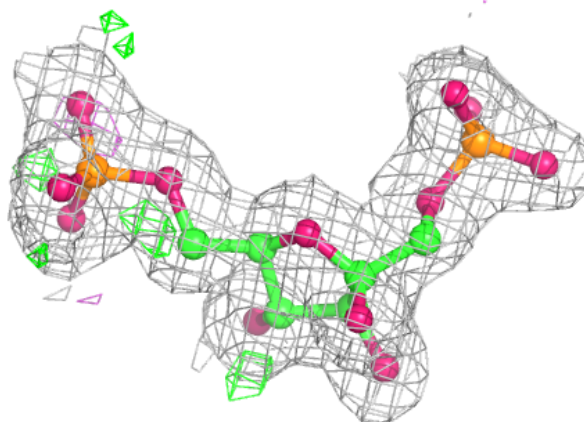
**Electron density around FBP F 601:**

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

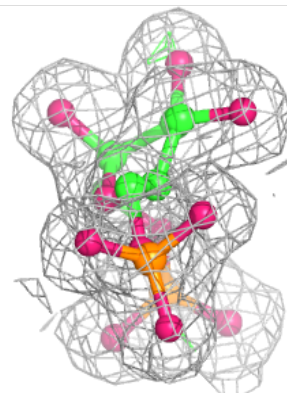
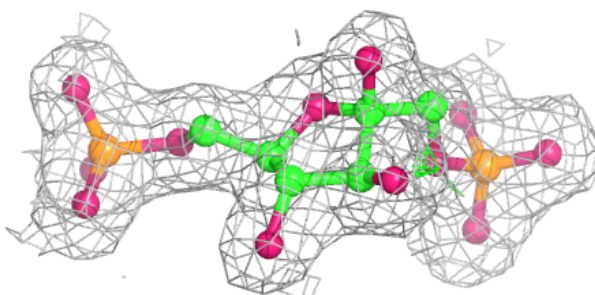
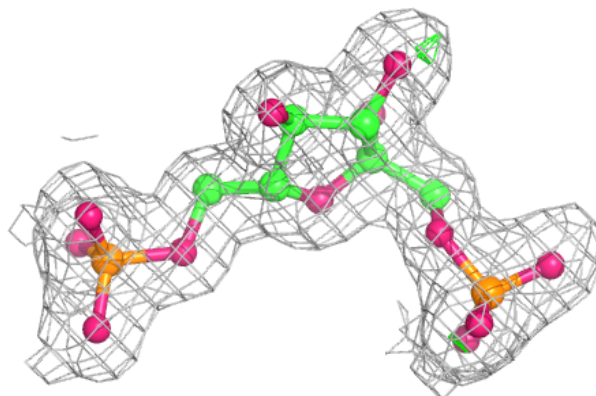


Electron density around FBP B 601:

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

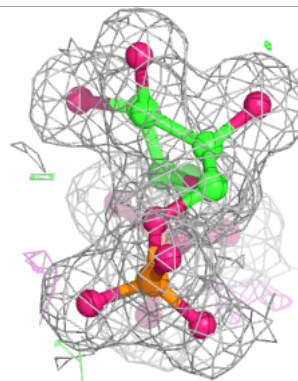
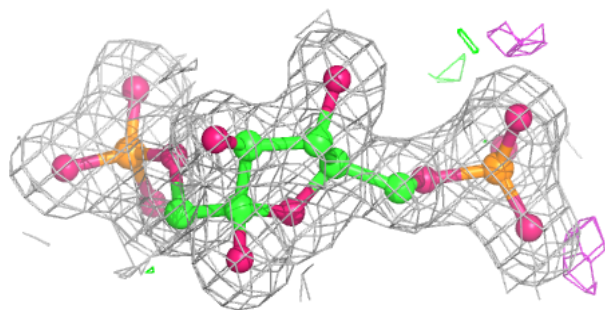
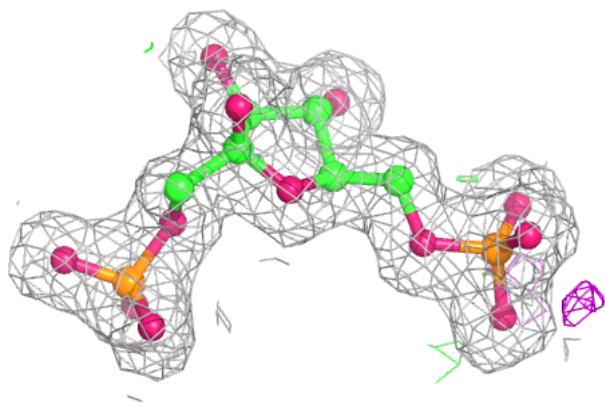
**Electron density around FBP E 601:**

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

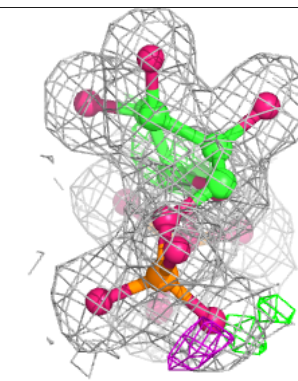
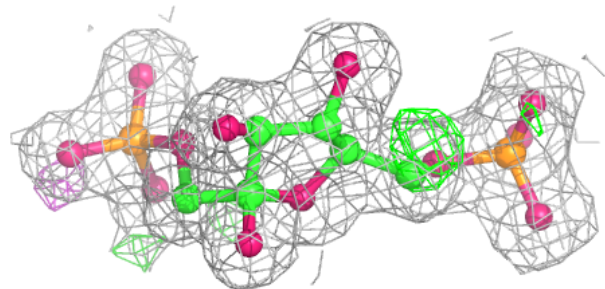
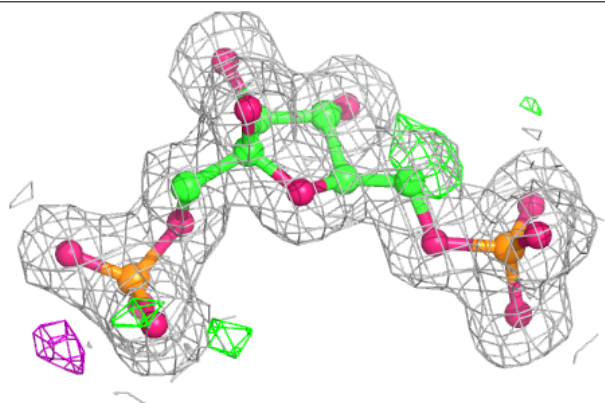


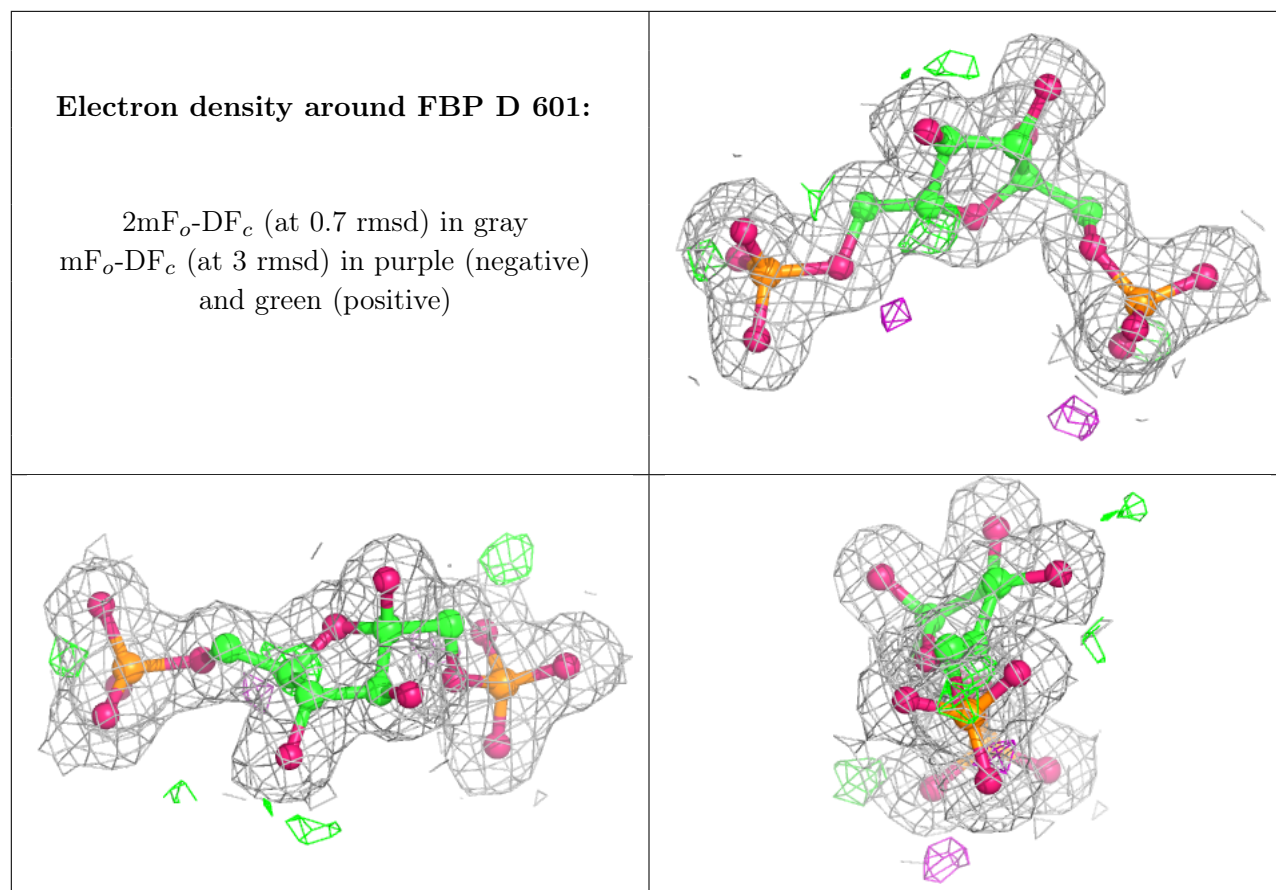
Electron density around FBP C 601:

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

**Electron density around FBP H 601:**

$2mF_o-DF_c$ (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.