



wwPDB X-ray Structure Validation Summary Report ⓘ

Aug 5, 2026 – 10:15 PM EDT

PDB ID : 3L75 / pdb_00003L75
Title : Cytochrome BC1 complex from chicken with fenamidone bound
Authors : Huang, L.; Berry, E.A.
Deposited on : 2009-12-28
Resolution : 2.79 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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

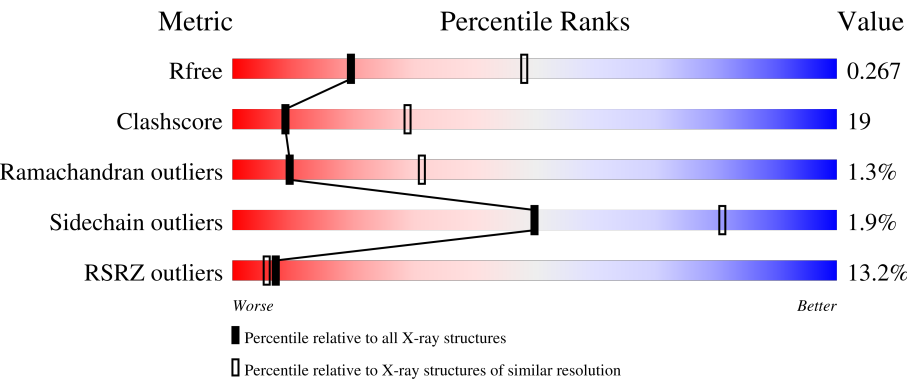
MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.015 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.50

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.79 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.



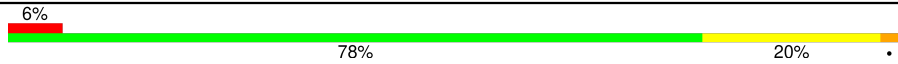
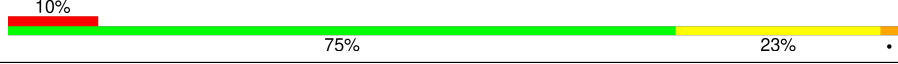
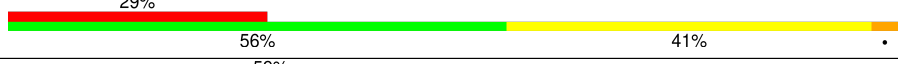
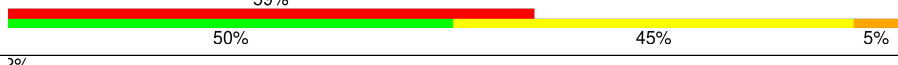
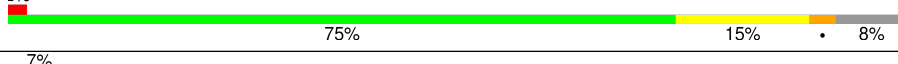
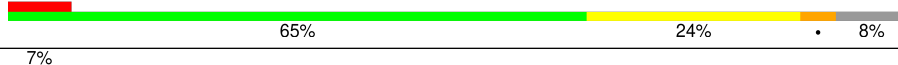
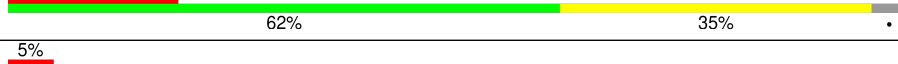
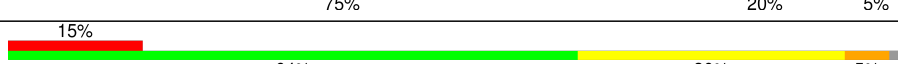
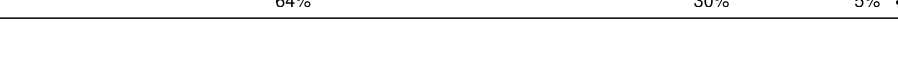
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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 ≥ 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	446	
1	N	446	
2	B	441	
2	O	441	
3	C	380	

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Mol	Chain	Length	Quality of chain
3	P	380	
4	D	241	
4	Q	241	
5	E	196	
5	R	196	
6	F	110	
6	S	110	
7	G	81	
7	T	81	
8	H	77	
8	U	77	
9	I	47	
9	V	47	
10	J	61	
10	W	61	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
11	PEE	N	3008	-	X	-	-
12	UNL	P	3048	-	-	-	X
16	AZI	C	2011	-	X	-	-
16	AZI	P	3011	-	X	-	-
17	BOG	D	2091	-	-	-	X
18	HEC	D	501	X	-	-	-
18	HEC	Q	501	X	-	-	-
20	FES	R	501	-	-	X	-

2 Entry composition

There are 21 unique types of molecules in this entry. The entry contains 32691 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 MITOCHONDRIAL UBIQUINOL-CYTOCHROME-C REDUCTASE COMPLEX CORE PROTEIN I.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	443	Total	C	N	O	S	0	0	1
			3440	2155	606	658	21			
1	N	442	Total	C	N	O	S	0	0	0
			3437	2154	605	657	21			

- Molecule 2 is a protein called MITOCHONDRIAL UBIQUINOL-CYTOCHROME-C REDUCTASE COMPLEX CORE PROTEIN 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	421	Total	C	N	O	S	0	0	0
			3141	1974	545	613	9			
2	O	422	Total	C	N	O	S	0	0	0
			3147	1977	546	614	10			

- Molecule 3 is a protein called CYTOCHROME B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	380	Total	C	N	O	S	0	0	0
			3017	2022	478	505	12			
3	P	379	Total	C	N	O	S	0	0	0
			3012	2019	477	504	12			

- Molecule 4 is a protein called MITOCHONDRIAL CYTOCHROME C1, HEME PROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	241	Total	C	N	O	S	0	0	0
			1898	1212	327	347	12			
4	Q	241	Total	C	N	O	S	0	0	0
			1898	1212	327	347	12			

- Molecule 5 is a protein called CYTOCHROME B-C1 COMPLEX SUBUNIT 5, RIESKE IRONSULFUR PROTEIN, MITOCHONDRIAL.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	E	196	Total	C	N	O	S	0	0	0
			1513	952	263	292	6			
5	R	196	Total	C	N	O	S	0	0	0
			1512	952	262	292	6			

- Molecule 6 is a protein called MITOCHONDRIAL UBIQUINOL-CYTOCHROME C REDUCTASE 14 KDA PROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	F	101	Total	C	N	O	S	0	0	0
			891	570	159	159	3			
6	S	101	Total	C	N	O	S	0	0	0
			891	570	159	159	3			

- Molecule 7 is a protein called MITOCHONDRIAL UBIQUINOL-CYTOCHROME C REDUCTASE UBIQUINONE-BINDING PROTEIN QP-C.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
7	G	81	Total	C	N	O	0	0	0
			676	439	120	117			
7	T	78	Total	C	N	O	0	0	0
			654	428	116	110			

- Molecule 8 is a protein called MITOCHONDRIAL UBIQUINOL-CYTOCHROME C REDUCTASE 11 KDA PROTEIN, COMPLEX III SUBUNIT VIII.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
8	H	70	Total	C	N	O	S	0	0	0
			574	350	105	114	5			
8	U	67	Total	C	N	O	S	0	0	0
			553	338	103	107	5			

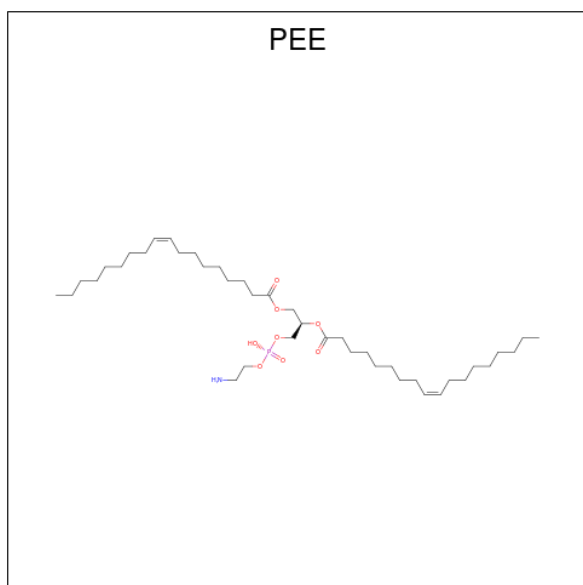
- Molecule 9 is a protein called CYTOCHROME B-C1 COMPLEX SUBUNIT RIESKE, MITOCHONDRIAL.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
9	I	46	Total	C	N	O	S	0	0	0
			288	172	58	56	2			
9	V	44	Total	C	N	O	S	0	0	1
			278	167	56	53	2			

- Molecule 10 is a protein called MITOCHONDRIAL UBIQUINOL-CYTOCHROME C REDUCTASE 7.2 KDA PROTEIN.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
10	J	61	Total	C	N	O	0	0	0
			497	321	87	89			
10	W	60	Total	C	N	O	0	0	1
			479	311	86	82			

- Molecule 11 is 1,2-dioleoyl-sn-glycero-3-phosphoethanolamine (CCD ID: PEE) (formula: $C_{41}H_{78}NO_8P$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
11	A	1	Total	C	O	P		0	0
			21	12	8	1			
11	C	1	Total	C	N	O	P	0	0
			50	40	1	8	1		
11	C	1	Total	C	N	O	P	0	0
			48	38	1	8	1		
11	N	1	Total	O	P			0	0
			5	4	1				
11	P	1	Total	C	N	O	P	0	0
			50	40	1	8	1		
11	P	1	Total	C	N	O	P	0	0
			48	38	1	8	1		

- Molecule 12 is UNKNOWN LIGAND (CCD ID: UNL) (formula:).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
12	A	1	Total 1 1	0	0

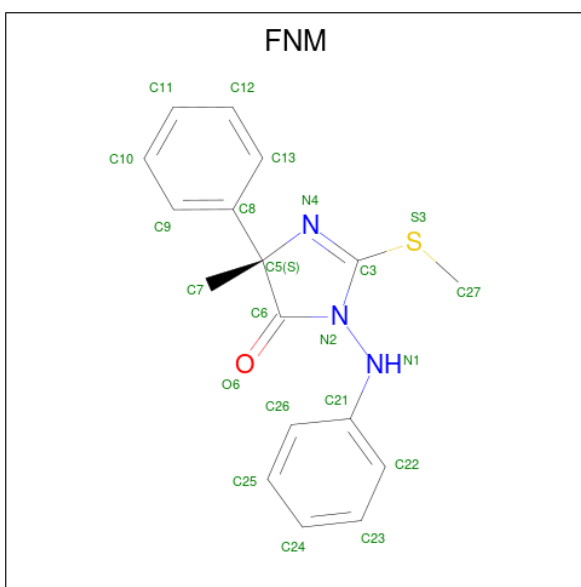
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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
12	C	3	Total O 5 5	0	0
12	D	1	Total O 2 2	0	0
12	P	5	Total O 7 7	0	0
12	R	1	Total O 1 1	0	0

- # HEM

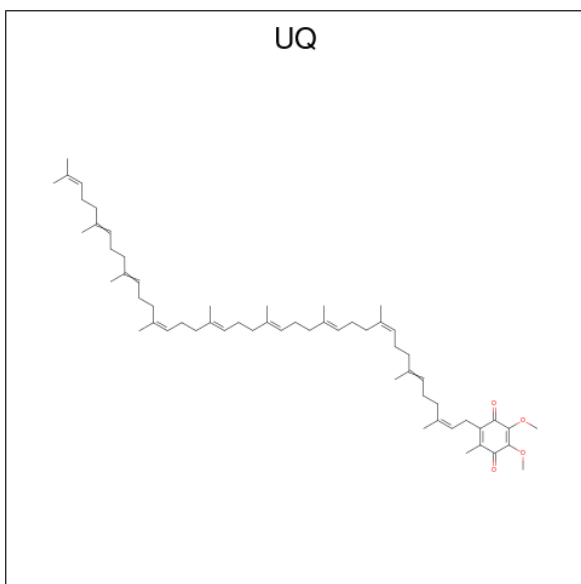
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
13	C	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
13	C	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
13	P	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
13	P	1	Total 43	C 34	Fe 1	N 4	O 4	0	0

- 



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
14	C	1	Total	C	N	O	S	0	0
			22	17	3	1	1		
14	P	1	Total	C	N	O	S	0	0
			22	17	3	1	1		

- Molecule 15 is Coenzyme Q10, (2Z,6E,10Z,14E,18E,22E,26Z)-isomer (CCD ID: UQ) (formula: C₅₉H₉₀O₄).



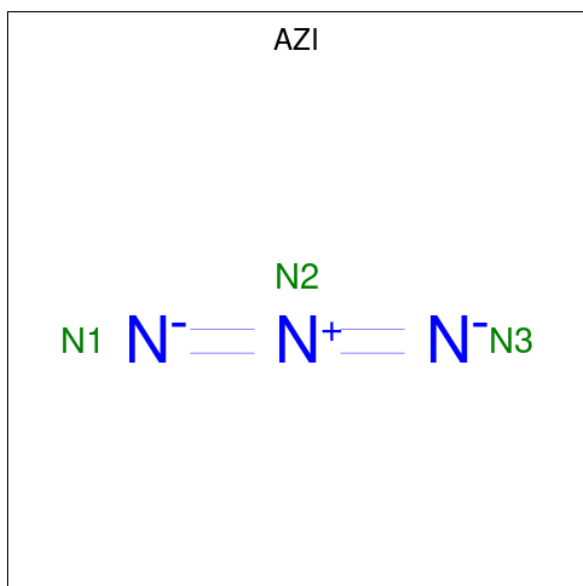
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
15	C	1	Total	C	O	0	0
			19	15	4		

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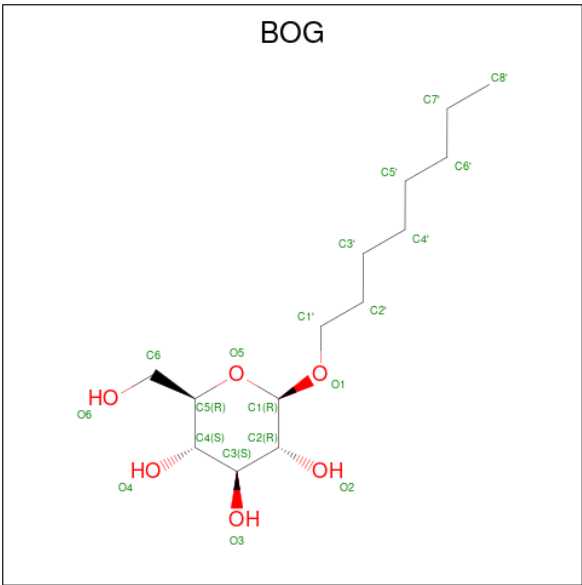
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
15	P	1	Total	C	O	0	0
			19	15	4		

- Molecule 16 is AZIDE ION (CCD ID: AZI) (formula: N₃).



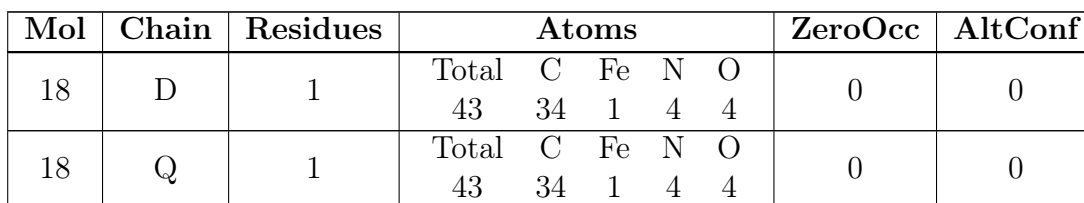
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
16	C	1	Total	N	0	0
			3	3		
16	P	1	Total	N	0	0
			3	3		

- Molecule 17 is octyl beta-D-glucopyranoside (CCD ID: BOG) (formula: C₁₄H₂₈O₆).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
17	C	1	Total	C	O	0	0
			12	10	2		
17	D	1	Total	C	O	0	0
			20	14	6		
17	D	1	Total	C	O	0	0
			20	14	6		
17	P	1	Total	C	O	0	0
			19	13	6		
17	Q	1	Total	C	O	0	0
			20	14	6		
17	Q	1	Total	C	O	0	0
			20	14	6		

- Molecule 18 is HEME C (CCD ID: HEC) (formula: C₃₄H₃₆FeN₄O₄).



- # CDL
-
- The chemical structure of CDL (1,3-bis(sn-3'-phosphatidyl)-sn-glycerol) is shown. It consists of a glycerol backbone with three phosphate groups and three long-chain fatty acid tails. The structure is labeled "CDL".

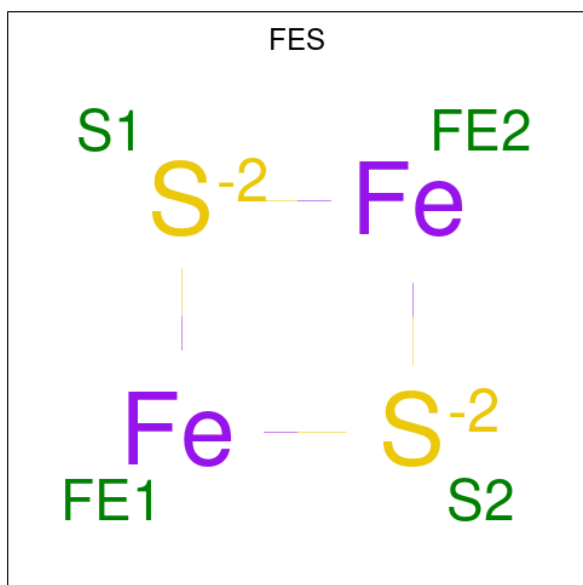
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
19	D	1	Total 42	C 23	O 17	P 2	0	0
19	G	1	Total 40	C 21	O 17	P 2	0	0



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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
19	Q	1	Total	C	O	P	0	0
			42	23	17	2		
19	T	1	Total	C	O	P	0	0
			40	21	17	2		

- Molecule 20 is FE2/S2 (INORGANIC) CLUSTER (CCD ID: FES) (formula: Fe_2S_2).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
20	E	1	Total	Fe	S	0	0
			4	2	2		
20	R	1	Total	Fe	S	0	0
			4	2	2		

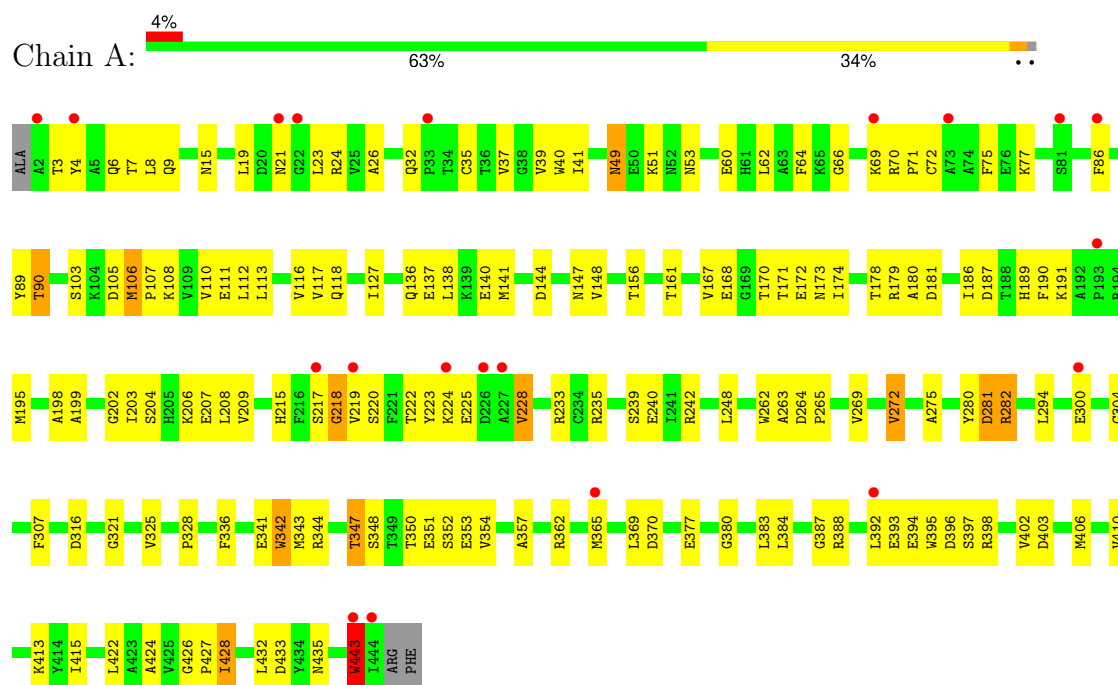
- Molecule 21 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
21	A	2	Total	O	0	0
			2	2		
21	C	9	Total	O	0	0
			9	9		
21	E	3	Total	O	0	0
			3	3		
21	P	10	Total	O	0	0
			10	10		
21	R	4	Total	O	0	0
			4	4		

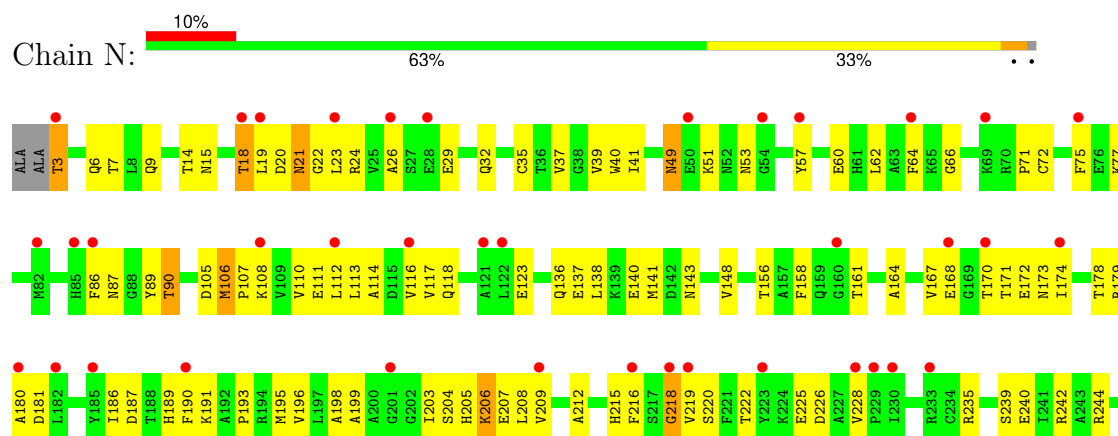
3 Residue-property plots

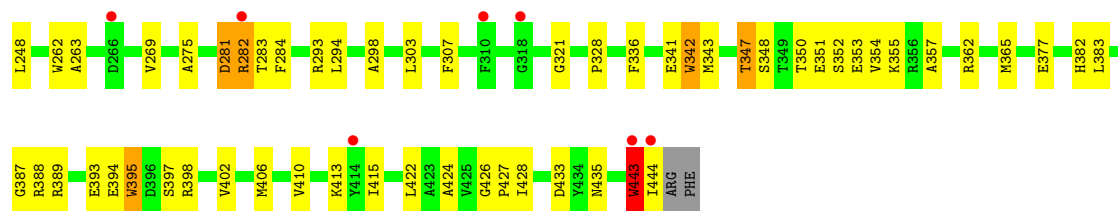
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: MITOCHONDRIAL UBIQUINOL-CYTOCHROME-C REDUCTASE COMPLEX CORE PROTEIN I



• Molecule 1: MITOCHONDRIAL UBIQUINOL-CYTOCHROME-C REDUCTASE COMPLEX CORE PROTEIN I

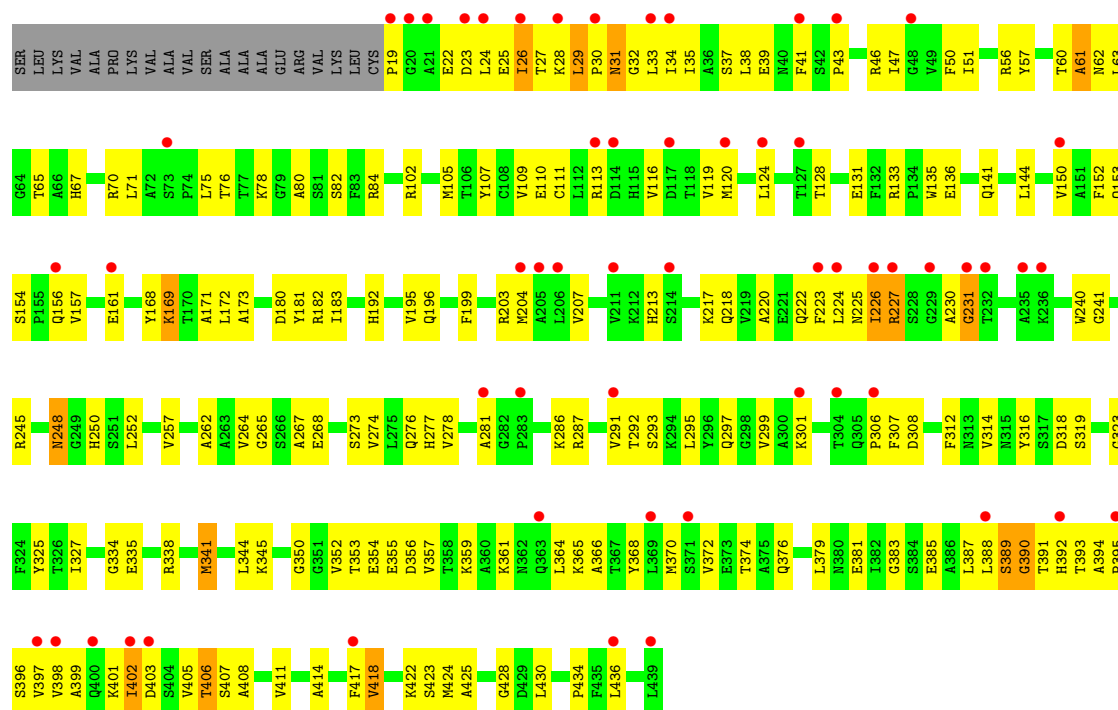




• Molecule 2: MITOCHONDRIAL UBIQUINOL-CYTOCHROME-C REDUCTASE COMPLEX CORE PROTEIN 2



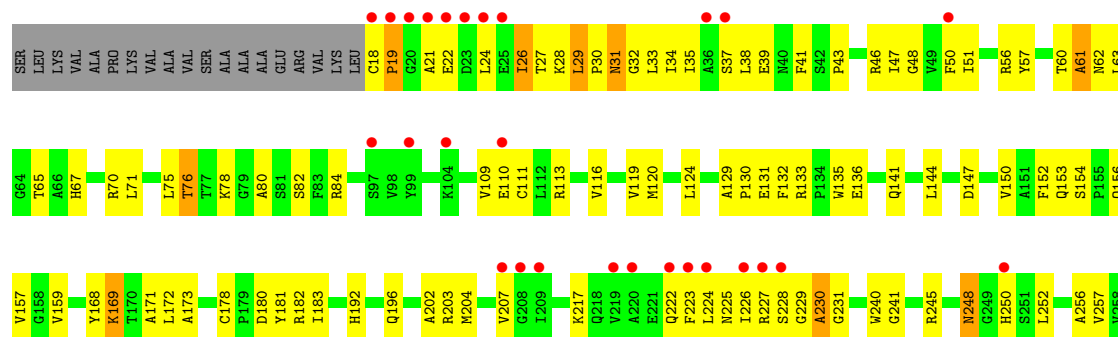
Chain B:

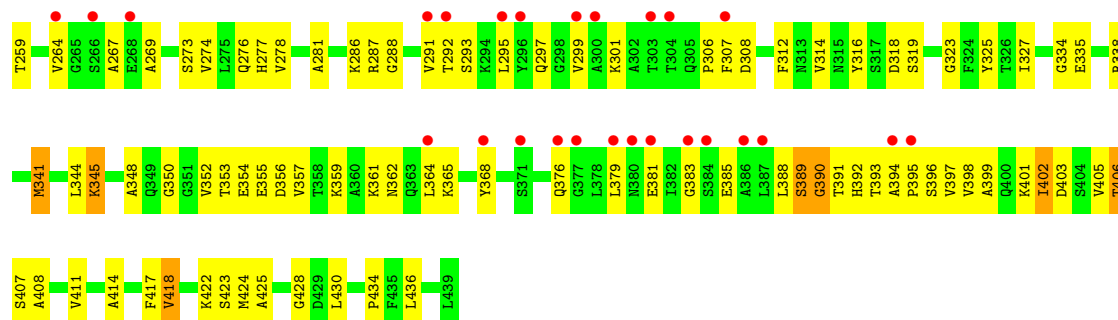


• Molecule 2: MITOCHONDRIAL UBIQUINOL-CYTOCHROME-C REDUCTASE COMPLEX CORE PROTEIN 2

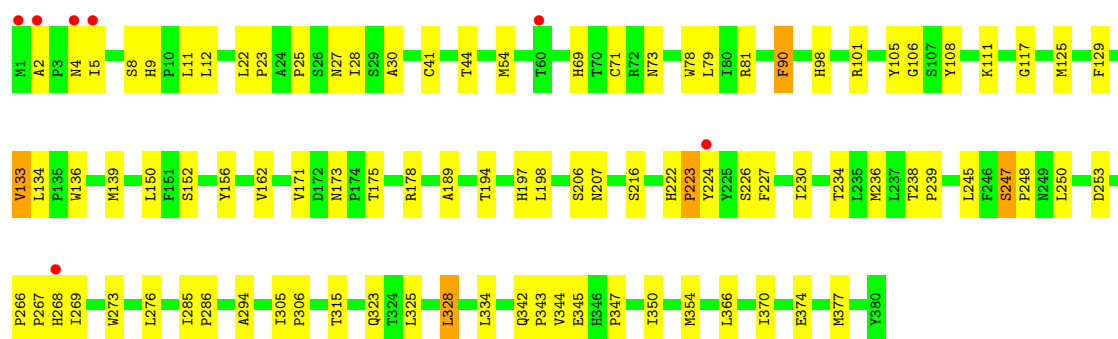
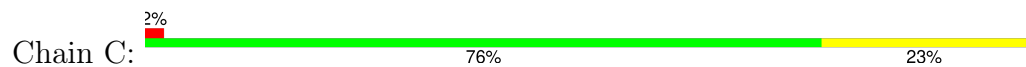


Chain O:

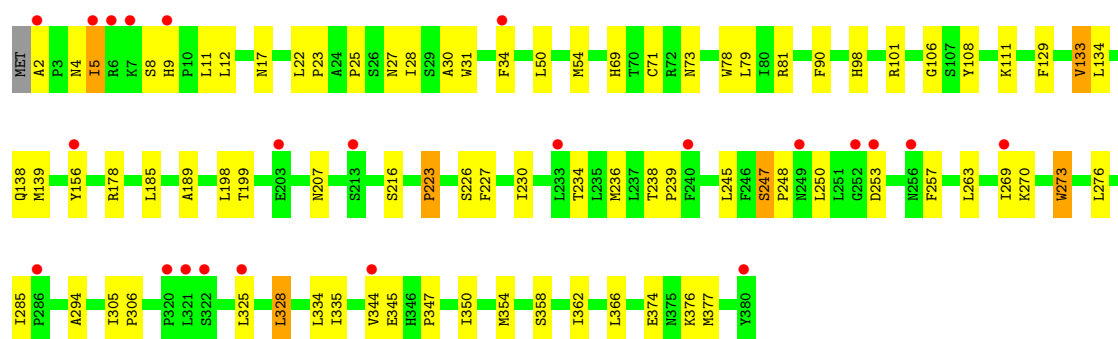
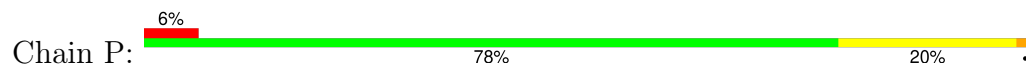




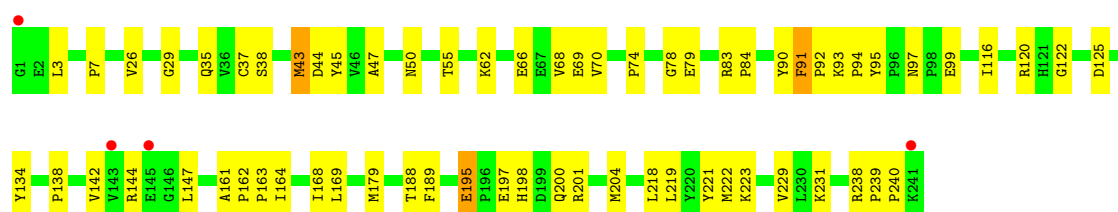
• Molecule 3: CYTOCHROME B



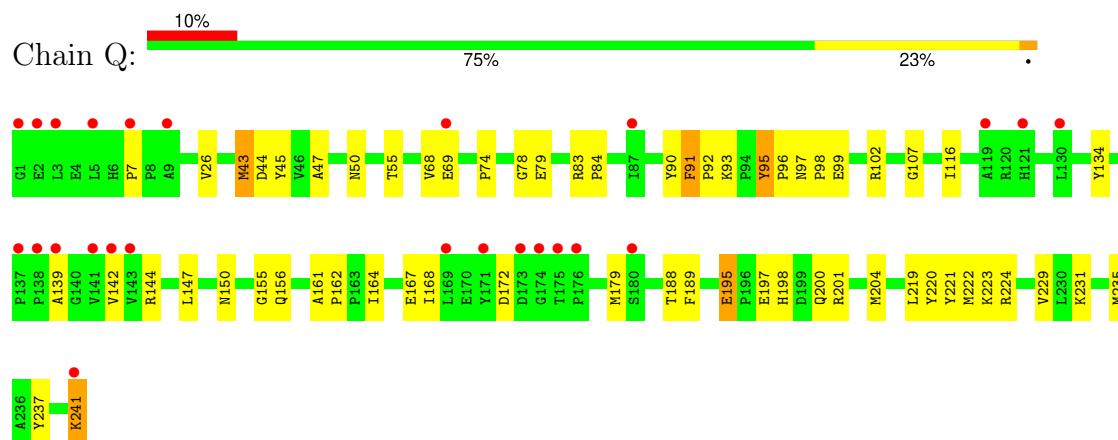
• Molecule 3: CYTOCHROME B



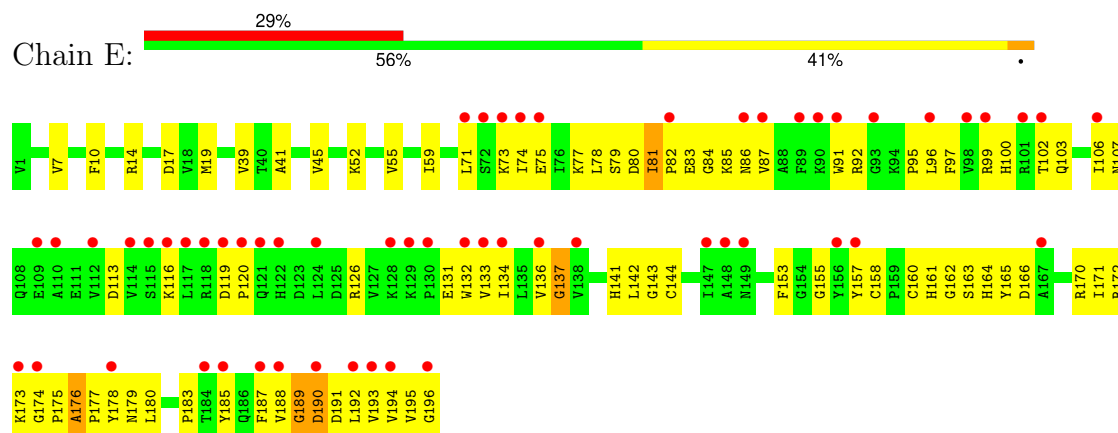
• Molecule 4: MITOCHONDRIAL CYTOCHROME C1, HEME PROTEIN



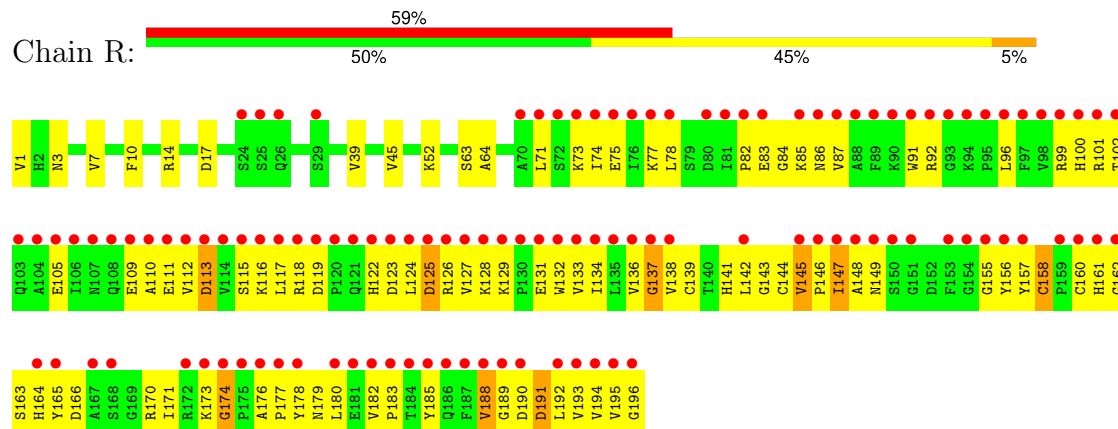
• Molecule 4: MITOCHONDRIAL CYTOCHROME C1, HEME PROTEIN



• Molecule 5: CYTOCHROME B-C1 COMPLEX SUBUNIT 5, RIESKE IRONSULFUR PROTEIN, MITOCHONDRIAL



• Molecule 5: CYTOCHROME B-C1 COMPLEX SUBUNIT 5, RIESKE IRONSULFUR PROTEIN, MITOCHONDRIAL

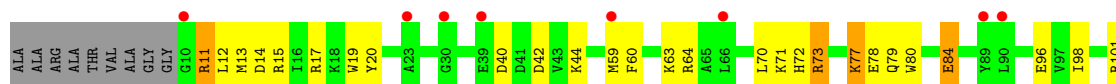


• Molecule 6: MITOCHONDRIAL UBIQUINOL-CYTOCHROME C REDUCTASE 14 KDA PROTEIN

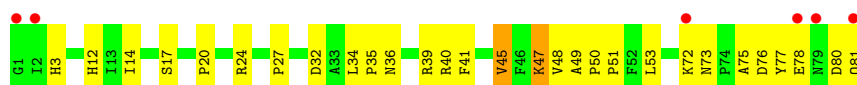




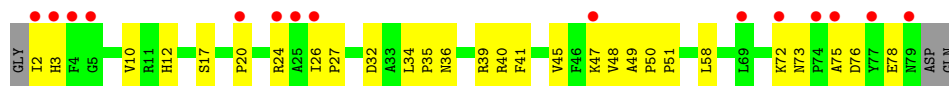
- Molecule 6: MITOCHONDRIAL UBIQUINOL-CYTOCHROME C REDUCTASE 14 KDA PROTEIN



- Molecule 7: MITOCHONDRIAL UBIQUINOL-CYTOCHROME C REDUCTASE UBIQUINONE-BINDING PROTEIN QP-C



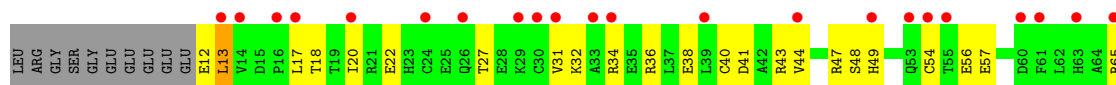
- Molecule 7: MITOCHONDRIAL UBIQUINOL-CYTOCHROME C REDUCTASE UBIQUINONE-BINDING PROTEIN QP-C

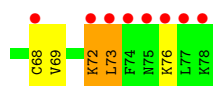


- Molecule 8: MITOCHONDRIAL UBIQUINOL-CYTOCHROME C REDUCTASE 11 KDA PROTEIN, COMPLEX III SUBUNIT VIII

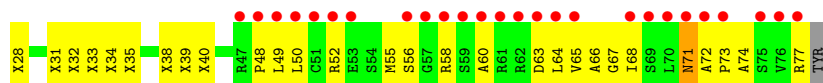


- Molecule 8: MITOCHONDRIAL UBIQUINOL-CYTOCHROME C REDUCTASE 11 KDA PROTEIN, COMPLEX III SUBUNIT VIII

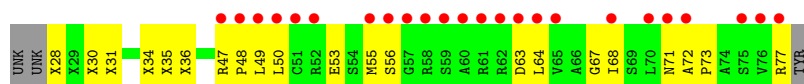




● Molecule 9: CYTOCHROME B-C1 COMPLEX SUBUNIT RIESKE, MITOCHONDRIAL



● Molecule 9: CYTOCHROME B-C1 COMPLEX SUBUNIT RIESKE, MITOCHONDRIAL



● Molecule 10: MITOCHONDRIAL UBIQUINOL-CYTOCHROME C REDUCTASE 7.2 KDA PROTEIN



● Molecule 10: MITOCHONDRIAL UBIQUINOL-CYTOCHROME C REDUCTASE 7.2 KDA PROTEIN



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	172.28Å 182.14Å 241.20Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	58.78 – 2.79 58.78 – 2.79	Depositor EDS
% Data completeness (in resolution range)	97.6 (58.78-2.79) 97.6 (58.78-2.79)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.03 (at 2.77Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.248 , 0.275 0.236 , 0.267	Depositor DCC
R_{free} test set	3613 reflections (1.96%)	wwPDB-VP
Wilson B-factor (Å ²)	64.6	Xtriage
Anisotropy	0.413	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 45.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	32691	wwPDB-VP
Average B, all atoms (Å ²)	69.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.81% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CDL, UQ, PEE, BOG, AZI, FES, FNM, UNL, HEC, HEM

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	0.52	0/3511	0.97	13/4757 (0.3%)
1	N	0.49	0/3508	0.98	12/4753 (0.3%)
2	B	0.45	0/3196	0.95	8/4334 (0.2%)
2	O	0.48	0/3202	0.99	11/4343 (0.3%)
3	C	0.62	0/3119	1.01	16/4270 (0.4%)
3	P	0.55	0/3114	0.99	14/4263 (0.3%)
4	D	0.58	0/1956	1.01	9/2658 (0.3%)
4	Q	0.48	0/1956	0.98	10/2658 (0.4%)
5	E	0.47	0/1547	0.94	5/2103 (0.2%)
5	R	0.45	0/1545	0.90	7/2098 (0.3%)
6	F	0.59	0/911	0.90	1/1219 (0.1%)
6	S	0.48	0/911	0.89	1/1219 (0.1%)
7	G	0.57	0/698	1.03	3/946 (0.3%)
7	T	0.45	0/676	1.01	3/918 (0.3%)
8	H	0.51	0/582	0.88	0/779
8	U	0.38	0/561	0.88	1/751 (0.1%)
9	I	0.52	0/218	1.00	1/293 (0.3%)
9	V	0.53	0/218	0.94	2/293 (0.7%)
10	J	0.51	0/508	0.89	1/682 (0.1%)
10	W	0.47	0/490	0.87	1/660 (0.2%)
All	All	0.51	0/32427	0.97	119/43997 (0.3%)

There are no bond length outliers.

The worst 5 of 119 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	O	227	ARG	N-CA-C	9.96	125.81	110.36
3	C	111	LYS	N-CA-C	9.20	122.46	111.33
2	O	18	CYS	CA-C-N	8.91	128.97	119.89
2	O	18	CYS	C-N-CA	8.91	128.97	119.89

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	415	ILE	CB-CA-C	-8.12	104.81	111.71

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	3440	0	3353	141	0
1	N	3437	0	3349	148	0
2	B	3141	0	3142	200	0
2	O	3147	0	3146	187	0
3	C	3017	0	3063	78	0
3	P	3012	0	3058	70	0
4	D	1898	0	1846	52	0
4	Q	1898	0	1846	51	0
5	E	1513	0	1478	70	0
5	R	1512	0	1476	100	0
6	F	891	0	893	23	0
6	S	891	0	893	30	0
7	G	676	0	659	26	0
7	T	654	0	641	25	0
8	H	574	0	548	18	0
8	U	553	0	535	24	0
9	I	288	0	254	45	0
9	V	278	0	252	32	0
10	J	497	0	490	11	0
10	W	479	0	478	14	0
11	A	21	0	13	1	0
11	C	98	0	147	0	0
11	N	5	0	0	0	0
11	P	98	0	147	1	0
12	A	1	0	0	0	0
12	C	5	0	0	0	0
12	D	2	0	0	0	0
12	P	7	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
12	R	1	0	0	0	0
13	C	86	0	60	5	0
13	P	86	0	60	5	0
14	C	22	0	17	5	0
14	P	22	0	17	4	0
15	C	19	0	17	3	0
15	P	19	0	17	2	0
16	C	3	0	0	0	0
16	P	3	0	0	0	0
17	C	12	0	18	0	0
17	D	40	0	56	3	0
17	P	19	0	24	1	0
17	Q	40	0	56	1	0
18	D	43	0	30	1	0
18	Q	43	0	30	0	0
19	D	42	0	28	2	0
19	G	40	0	24	1	0
19	Q	42	0	28	3	0
19	T	40	0	24	1	0
20	E	4	0	0	1	0
20	R	4	0	0	2	0
21	A	2	0	0	0	0
21	C	9	0	0	1	0
21	E	3	0	0	0	0
21	P	10	0	0	1	0
21	R	4	0	0	0	0
All	All	32691	0	32213	1210	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 19.

The worst 5 of 1210 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)
2:O:353:THR:HG22	2:O:355:GLU:H	1.16	1.11
5:R:83:GLU:HG2	5:R:102:THR:HG22	1.33	1.07
2:B:353:THR:HG22	2:B:355:GLU:H	1.20	1.06
2:B:157:VAL:HG23	9:I:64:LEU:HD21	1.38	1.03
1:N:178:THR:HG22	1:N:180:ALA:H	1.22	1.02

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	441/446 (99%)	412 (93%)	23 (5%)	6 (1%)	9	30
1	N	440/446 (99%)	411 (93%)	22 (5%)	7 (2%)	7	27
2	B	419/441 (95%)	374 (89%)	35 (8%)	10 (2%)	4	17
2	O	420/441 (95%)	372 (89%)	41 (10%)	7 (2%)	7	25
3	C	378/380 (100%)	364 (96%)	14 (4%)	0	100	100
3	P	377/380 (99%)	360 (96%)	17 (4%)	0	100	100
4	D	239/241 (99%)	228 (95%)	11 (5%)	0	100	100
4	Q	239/241 (99%)	224 (94%)	15 (6%)	0	100	100
5	E	194/196 (99%)	170 (88%)	18 (9%)	6 (3%)	3	12
5	R	192/196 (98%)	156 (81%)	28 (15%)	8 (4%)	2	7
6	F	99/110 (90%)	93 (94%)	5 (5%)	1 (1%)	12	38
6	S	99/110 (90%)	91 (92%)	6 (6%)	2 (2%)	6	21
7	G	79/81 (98%)	69 (87%)	10 (13%)	0	100	100
7	T	76/81 (94%)	68 (90%)	8 (10%)	0	100	100
8	H	68/77 (88%)	62 (91%)	6 (9%)	0	100	100
8	U	65/77 (84%)	60 (92%)	4 (6%)	1 (2%)	8	28
9	I	29/47 (62%)	26 (90%)	2 (7%)	1 (3%)	3	11
9	V	29/47 (62%)	24 (83%)	5 (17%)	0	100	100
10	J	59/61 (97%)	55 (93%)	3 (5%)	1 (2%)	7	25
10	W	58/61 (95%)	54 (93%)	2 (3%)	2 (3%)	3	11
All	All	4000/4160 (96%)	3673 (92%)	275 (7%)	52 (1%)	9	31

5 of 52 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	282	ARG

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Mol	Chain	Res	Type
2	B	26	ILE
2	B	226	ILE
2	B	227	ARG
2	B	389	SER

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	365/368 (99%)	357 (98%)	8 (2%)	45	78
1	N	365/368 (99%)	353 (97%)	12 (3%)	33	69
2	B	332/347 (96%)	326 (98%)	6 (2%)	51	82
2	O	333/347 (96%)	325 (98%)	8 (2%)	43	77
3	C	328/329 (100%)	323 (98%)	5 (2%)	57	84
3	P	328/329 (100%)	324 (99%)	4 (1%)	63	87
4	D	200/200 (100%)	198 (99%)	2 (1%)	68	88
4	Q	200/200 (100%)	197 (98%)	3 (2%)	57	84
5	E	166/166 (100%)	166 (100%)	0	100	100
5	R	165/166 (99%)	164 (99%)	1 (1%)	78	92
6	F	93/96 (97%)	89 (96%)	4 (4%)	26	60
6	S	93/96 (97%)	88 (95%)	5 (5%)	20	51
7	G	71/71 (100%)	70 (99%)	1 (1%)	59	85
7	T	69/71 (97%)	68 (99%)	1 (1%)	59	85
8	H	65/71 (92%)	64 (98%)	1 (2%)	57	84
8	U	63/71 (89%)	62 (98%)	1 (2%)	55	83
9	I	23/26 (88%)	22 (96%)	1 (4%)	26	60
9	V	23/26 (88%)	23 (100%)	0	100	100
10	J	49/49 (100%)	48 (98%)	1 (2%)	48	80
10	W	47/49 (96%)	46 (98%)	1 (2%)	47	79

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	3378/3446 (98%)	3313 (98%)	65 (2%)	50 81

5 of 65 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
6	S	14	ASP
6	S	70	LEU
6	F	84	GLU
6	F	73	ARG
6	S	73	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 86 such sidechains are listed below:

Mol	Chain	Res	Type
2	O	248	ASN
4	Q	35	GLN
2	O	329	GLN
3	P	82	ASN
4	Q	200	GLN

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 ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 41 ligands modelled in this entry, 11 are unknown - leaving 30 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
11	PEE	P	3007	-	47,47,50	1.38	5 (10%)	50,52,55	0.77	2 (4%)
19	CDL	G	2004	-	39,39,99	1.21	3 (7%)	45,51,111	1.10	3 (6%)
16	AZI	P	3011	-	2,2,2	3.11	2 (100%)	0,1,1	-	-
17	BOG	P	2010	-	18,18,20	1.10	3 (16%)	22,22,25	0.57	0
17	BOG	C	3010	-	11,11,20	1.08	2 (18%)	10,11,25	0.97	1 (10%)
14	FNM	P	3001	-	24,24,24	1.84	4 (16%)	24,34,34	1.73	2 (8%)
13	HEM	C	501	3	50,50,50	1.39	4 (8%)	67,82,82	1.44	10 (14%)
17	BOG	D	2091	-	20,20,20	1.12	2 (10%)	25,25,25	1.00	1 (4%)
13	HEM	C	502	3	50,50,50	1.57	8 (16%)	67,82,82	1.46	11 (16%)
14	FNM	C	2001	-	24,24,24	1.81	3 (12%)	24,34,34	1.89	3 (12%)
17	BOG	Q	3091	-	20,20,20	1.14	2 (10%)	25,25,25	0.90	1 (4%)
11	PEE	A	2008	-	20,20,50	1.71	4 (20%)	23,25,55	0.65	0
19	CDL	T	3004	-	39,39,99	1.22	3 (7%)	45,51,111	1.10	2 (4%)
18	HEC	D	501	4	50,50,50	1.23	6 (12%)	68,82,82	1.62	15 (22%)
17	BOG	D	2009	-	20,20,20	0.98	1 (5%)	25,25,25	0.93	2 (8%)
13	HEM	P	501	3	50,50,50	1.44	5 (10%)	67,82,82	1.21	4 (5%)
15	UQ	P	3002	-	19,19,63	2.46	10 (52%)	24,26,79	1.05	2 (8%)
13	HEM	P	502	3	50,50,50	1.53	4 (8%)	67,82,82	1.24	7 (10%)
17	BOG	Q	3009	-	20,20,20	0.94	1 (5%)	25,25,25	0.85	1 (4%)
11	PEE	P	3005	-	49,49,50	1.48	9 (18%)	52,54,55	0.83	3 (5%)
20	FES	E	501	5	0,4,4	-	-	-	-	-
11	PEE	C	2005	-	49,49,50	1.48	9 (18%)	52,54,55	0.83	2 (3%)
16	AZI	C	2011	-	2,2,2	2.74	2 (100%)	0,1,1	-	-
11	PEE	N	3008	-	4,4,50	3.65	4 (100%)	6,6,55	0.63	0
19	CDL	Q	3003	-	41,41,99	1.20	2 (4%)	47,53,111	1.07	4 (8%)
11	PEE	C	2007	-	47,47,50	1.32	6 (12%)	50,52,55	0.80	3 (6%)
20	FES	R	501	5	0,4,4	-	-	-	-	-
19	CDL	D	2003	-	41,41,99	1.15	1 (2%)	47,53,111	1.07	4 (8%)
18	HEC	Q	501	4	50,50,50	1.09	2 (4%)	68,82,82	1.52	13 (19%)
15	UQ	C	2002	-	19,19,63	2.63	10 (52%)	24,26,79	1.10	2 (8%)

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
11	PEE	P	3007	-	-	24/51/51/54	-
19	CDL	G	2004	-	-	21/49/49/110	-
17	BOG	P	2010	-	-	1/6/26/31	0/1/1/1
17	BOG	C	3010	-	-	5/9/9/31	-
14	FNM	P	3001	-	-	2/12/31/31	0/3/3/3
13	HEM	C	501	3	-	2/14/54/54	-
17	BOG	D	2091	-	-	6/11/31/31	0/1/1/1
13	HEM	C	502	3	-	6/14/54/54	-
14	FNM	C	2001	-	-	2/12/31/31	0/3/3/3
17	BOG	Q	3091	-	-	4/11/31/31	0/1/1/1
11	PEE	A	2008	-	-	10/24/24/54	-
19	CDL	T	3004	-	-	22/49/49/110	-
18	HEC	D	501	4	1/1/3/7	9/14/54/54	-
17	BOG	D	2009	-	-	4/11/31/31	0/1/1/1
13	HEM	P	501	3	-	2/14/54/54	-
15	UQ	P	3002	-	-	2/11/35/87	0/1/1/1
13	HEM	P	502	3	-	6/14/54/54	-
17	BOG	Q	3009	-	-	6/11/31/31	0/1/1/1
11	PEE	P	3005	-	-	29/53/53/54	-
20	FES	E	501	5	-	-	0/1/1/1
11	PEE	C	2005	-	-	28/53/53/54	-
19	CDL	Q	3003	-	-	24/51/51/110	-
11	PEE	C	2007	-	-	24/51/51/54	-
20	FES	R	501	5	-	-	0/1/1/1
19	CDL	D	2003	-	-	26/51/51/110	-
18	HEC	Q	501	4	1/1/3/7	8/14/54/54	-
15	UQ	C	2002	-	-	3/11/35/87	0/1/1/1

The worst 5 of 117 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
14	C	2001	FNM	C3-N4	6.99	1.34	1.28
14	P	3001	FNM	C3-N4	6.60	1.33	1.28
13	P	502	HEM	CBB-CAB	5.30	1.55	1.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
15	C	2002	UQ	C6-C5	5.09	1.44	1.35
15	C	2002	UQ	C7-C6	5.03	1.60	1.51

The worst 5 of 98 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
14	C	2001	FNM	C27-S3-C3	7.71	110.67	100.43
14	P	3001	FNM	C27-S3-C3	6.94	109.64	100.43
13	C	502	HEM	CAD-C3D-C4D	5.47	134.23	124.70
18	D	501	HEC	CAD-C3D-C4D	4.96	133.34	124.70
13	P	502	HEM	CAD-C3D-C4D	4.33	132.25	124.70

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
18	D	501	HEC	NA
18	Q	501	HEC	NA

5 of 276 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
11	A	2008	PEE	C1-O3P-P-O1P
11	A	2008	PEE	C1-O3P-P-O4P
11	C	2005	PEE	C11-C10-O2-C2
11	C	2005	PEE	C4-O4P-P-O3P
11	C	2005	PEE	C4-O4P-P-O2P

There are no ring outliers.

21 monomers are involved in 42 short contacts:

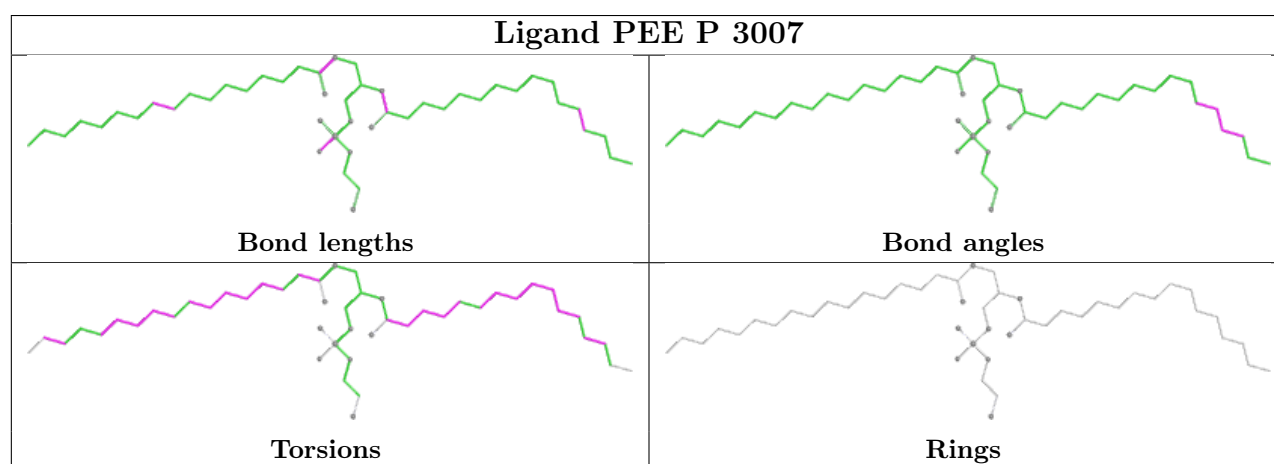
Mol	Chain	Res	Type	Clashes	Symm-Clashes
11	P	3007	PEE	1	0
19	G	2004	CDL	1	0
17	P	2010	BOG	1	0
14	P	3001	FNM	4	0
13	C	501	HEM	1	0
17	D	2091	BOG	2	0
13	C	502	HEM	4	0
14	C	2001	FNM	5	0
17	Q	3091	BOG	1	0
11	A	2008	PEE	1	0

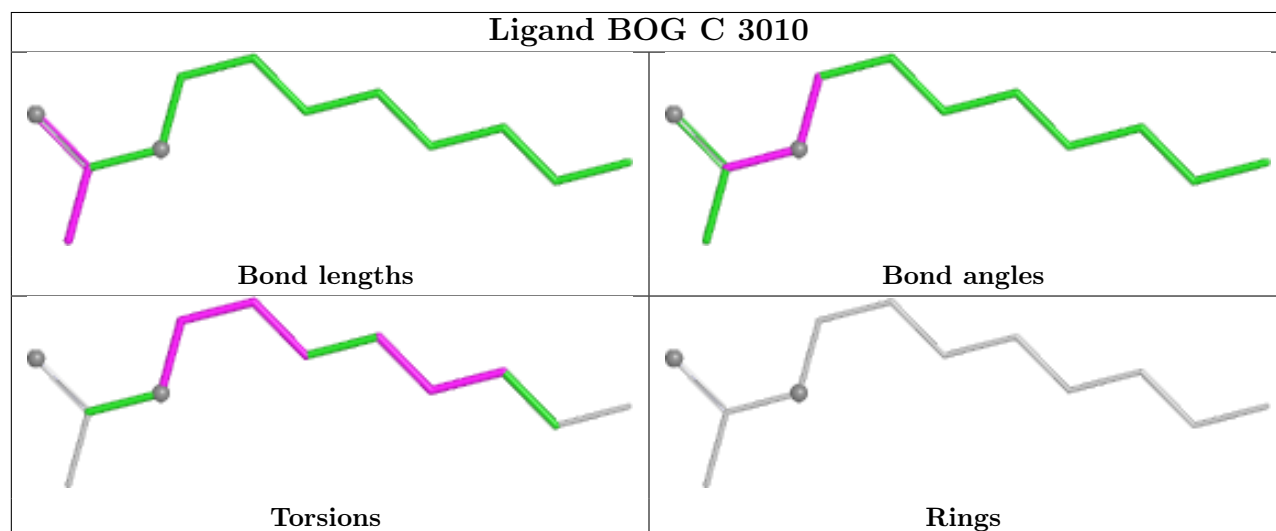
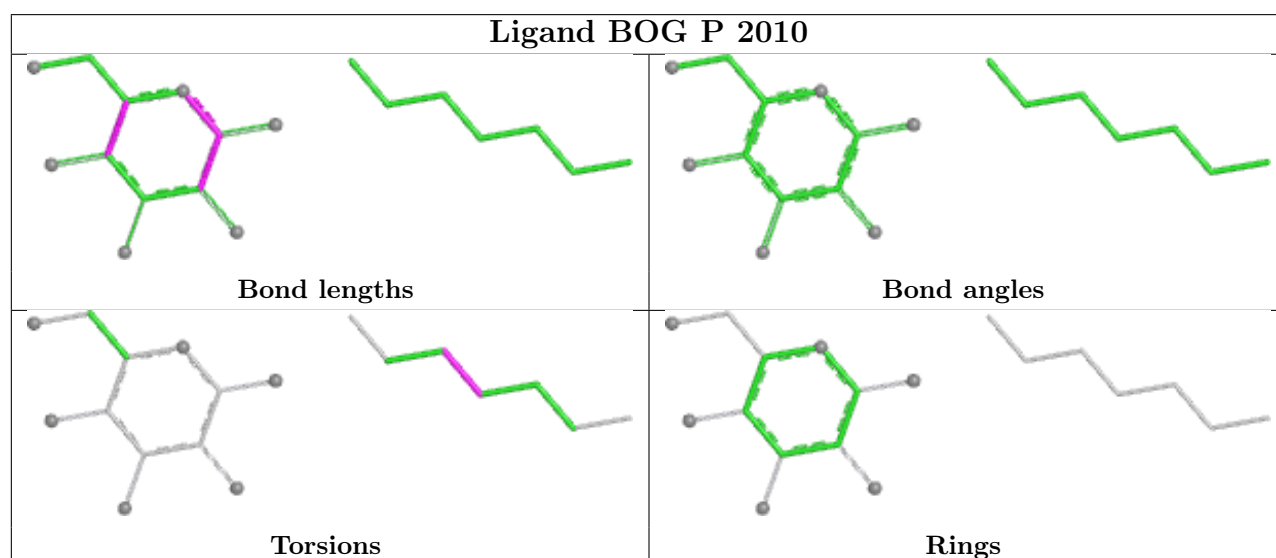
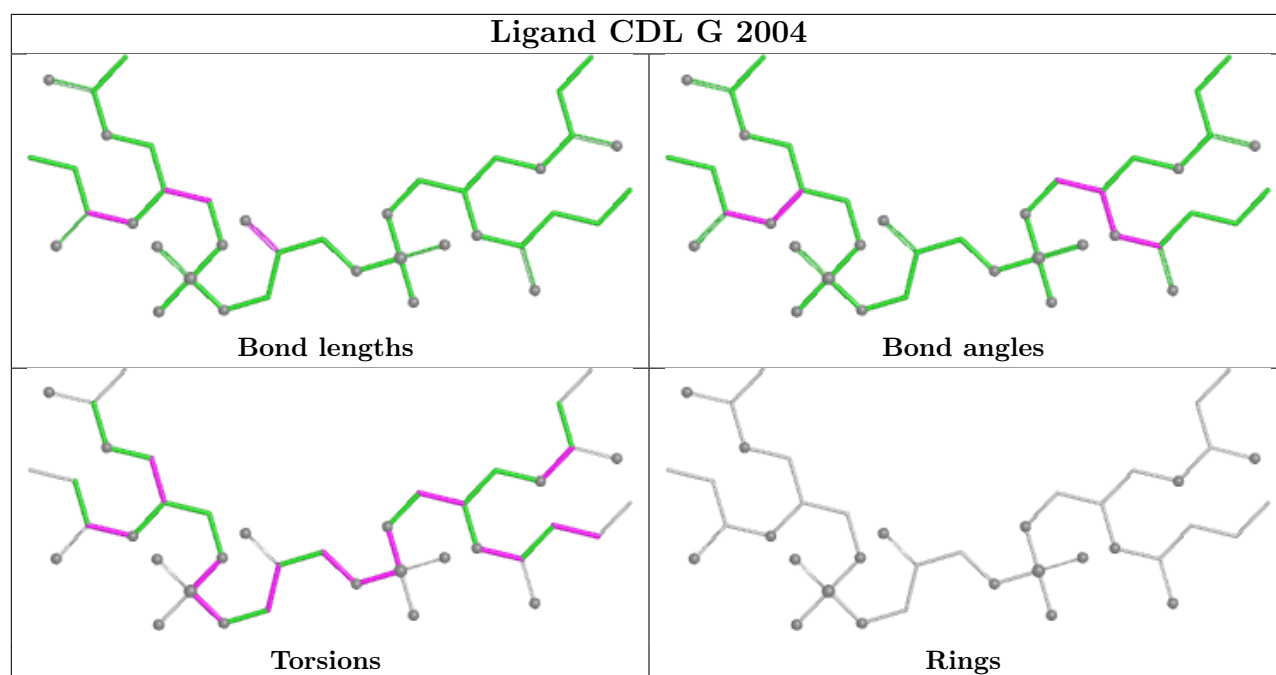
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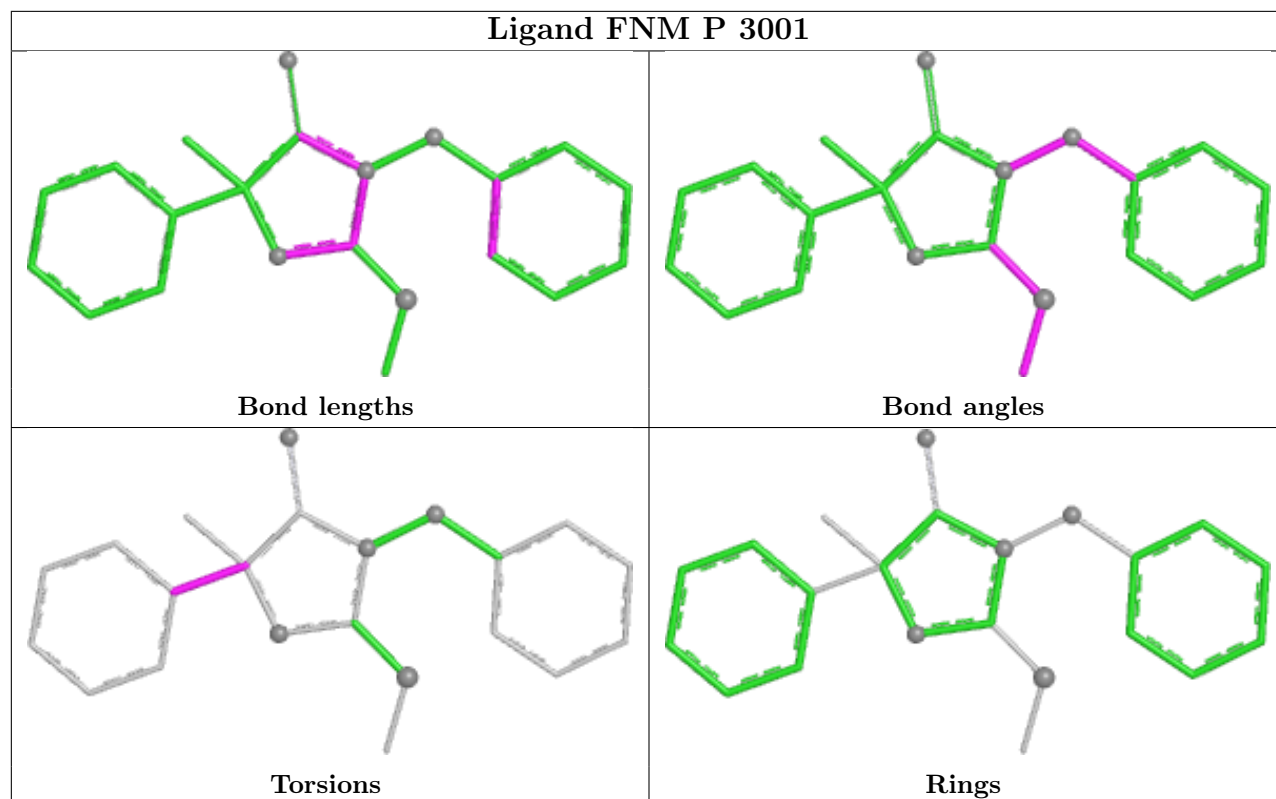
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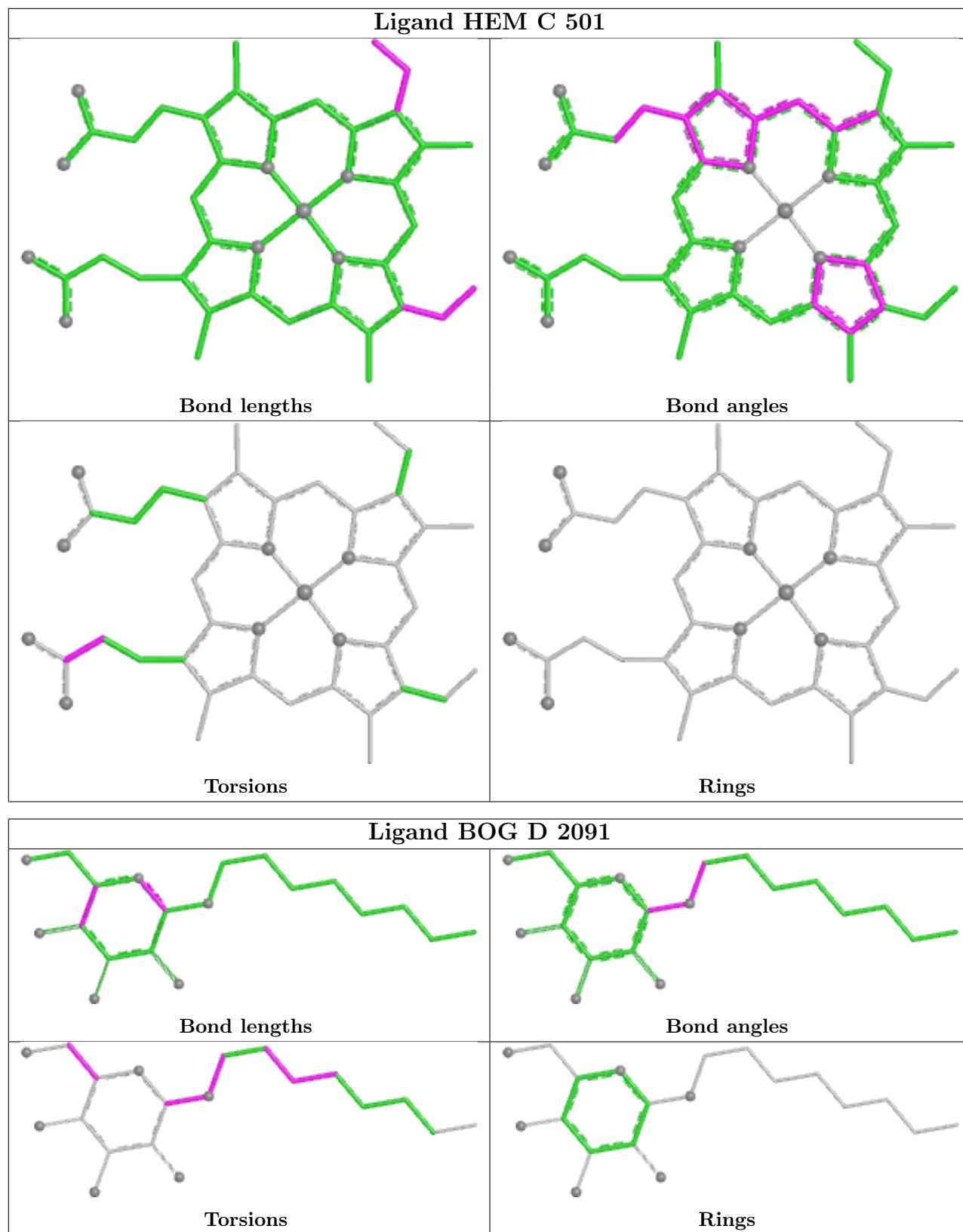
Mol	Chain	Res	Type	Clashes	Symm-Clashes
19	T	3004	CDL	1	0
18	D	501	HEC	1	0
17	D	2009	BOG	1	0
13	P	501	HEM	1	0
15	P	3002	UQ	2	0
13	P	502	HEM	4	0
20	E	501	FES	1	0
19	Q	3003	CDL	3	0
20	R	501	FES	2	0
19	D	2003	CDL	2	0
15	C	2002	UQ	3	0

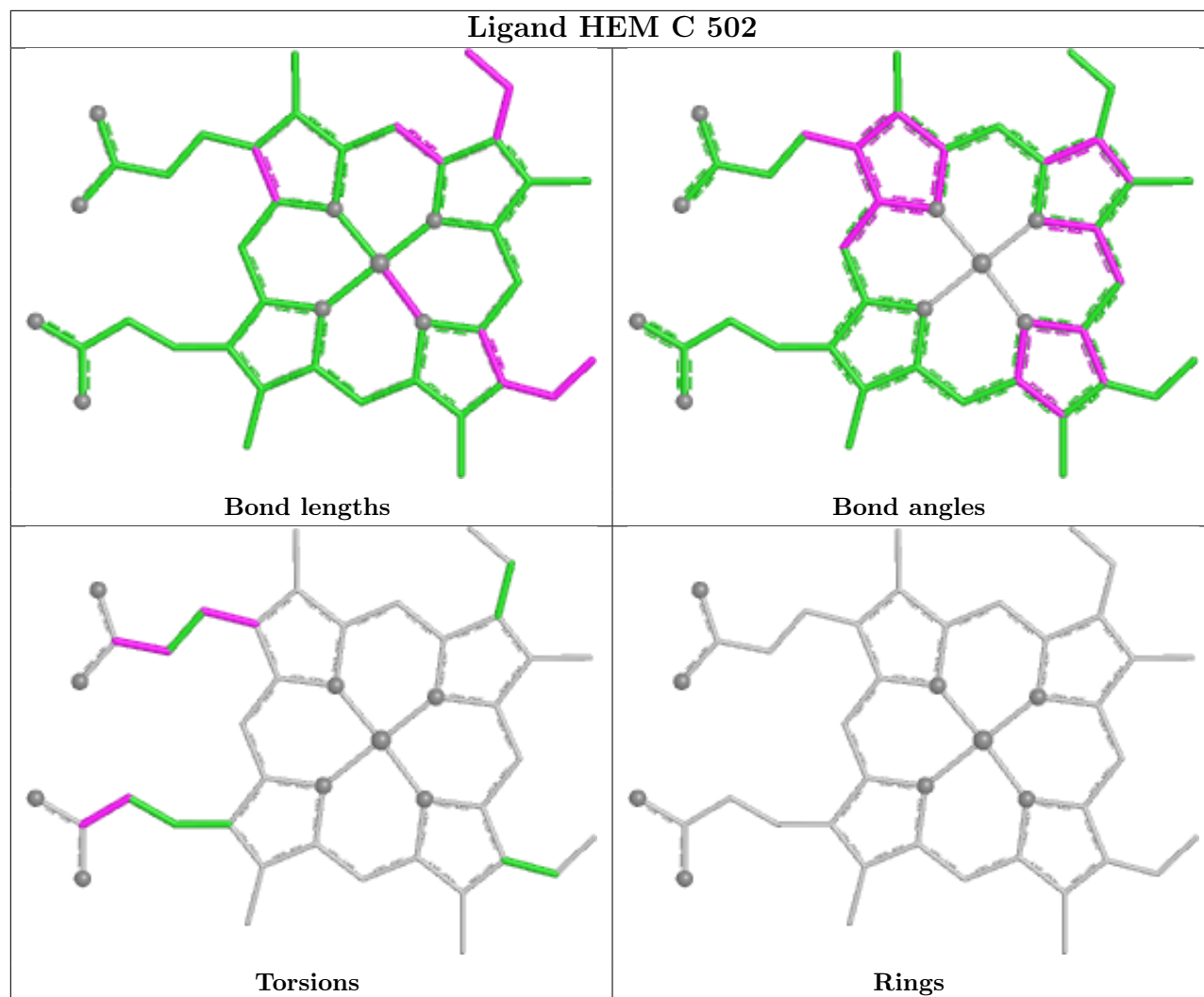
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

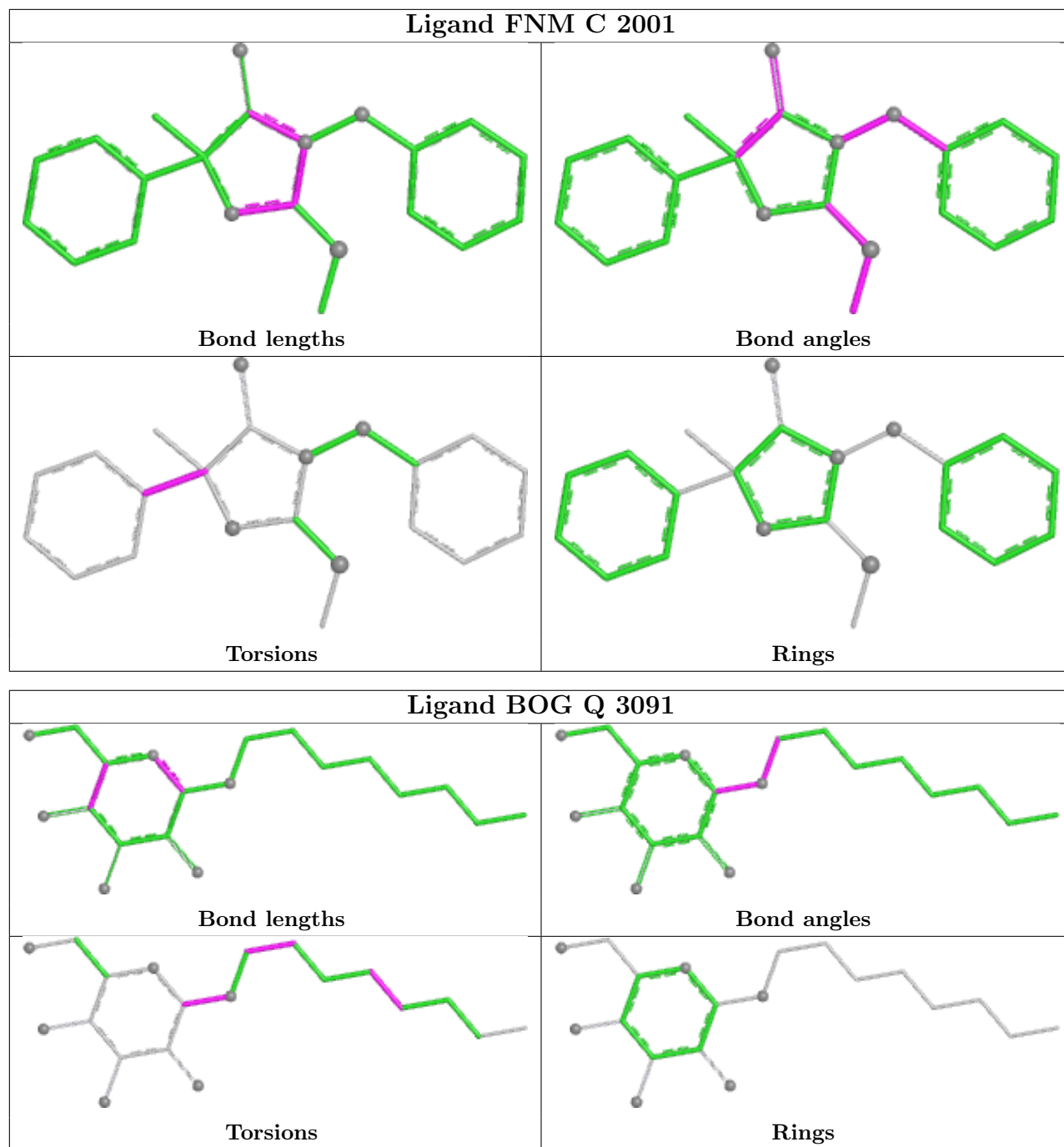


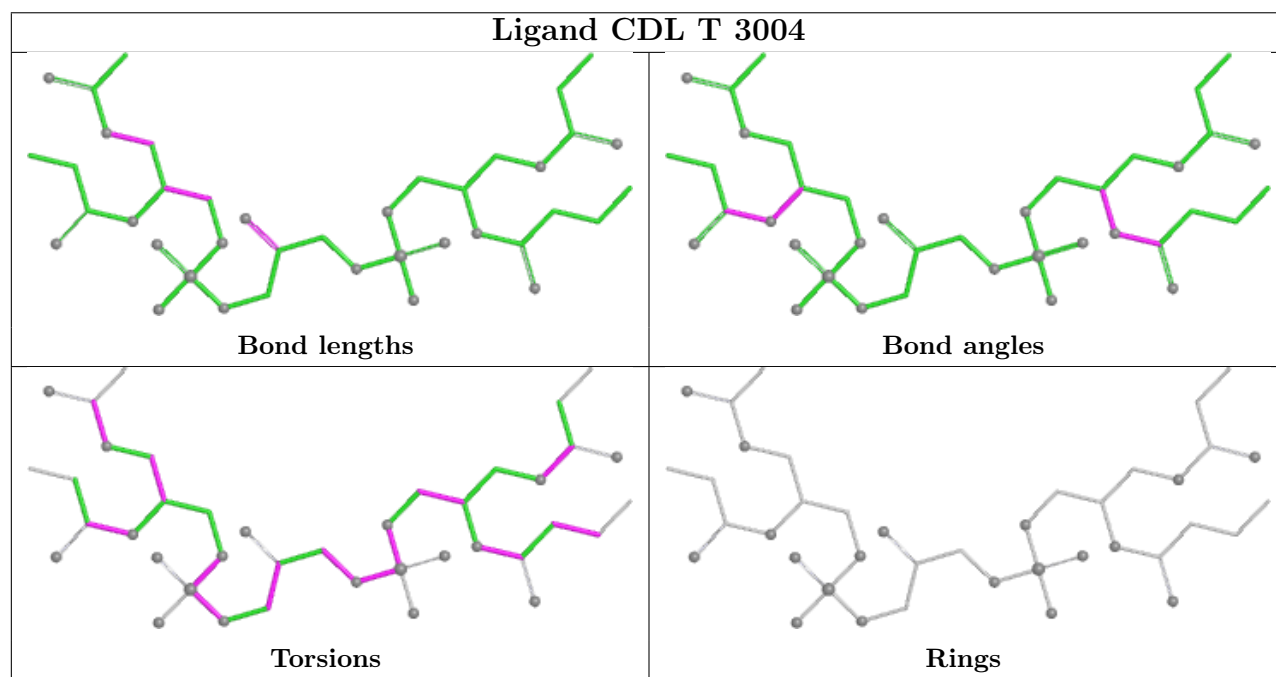
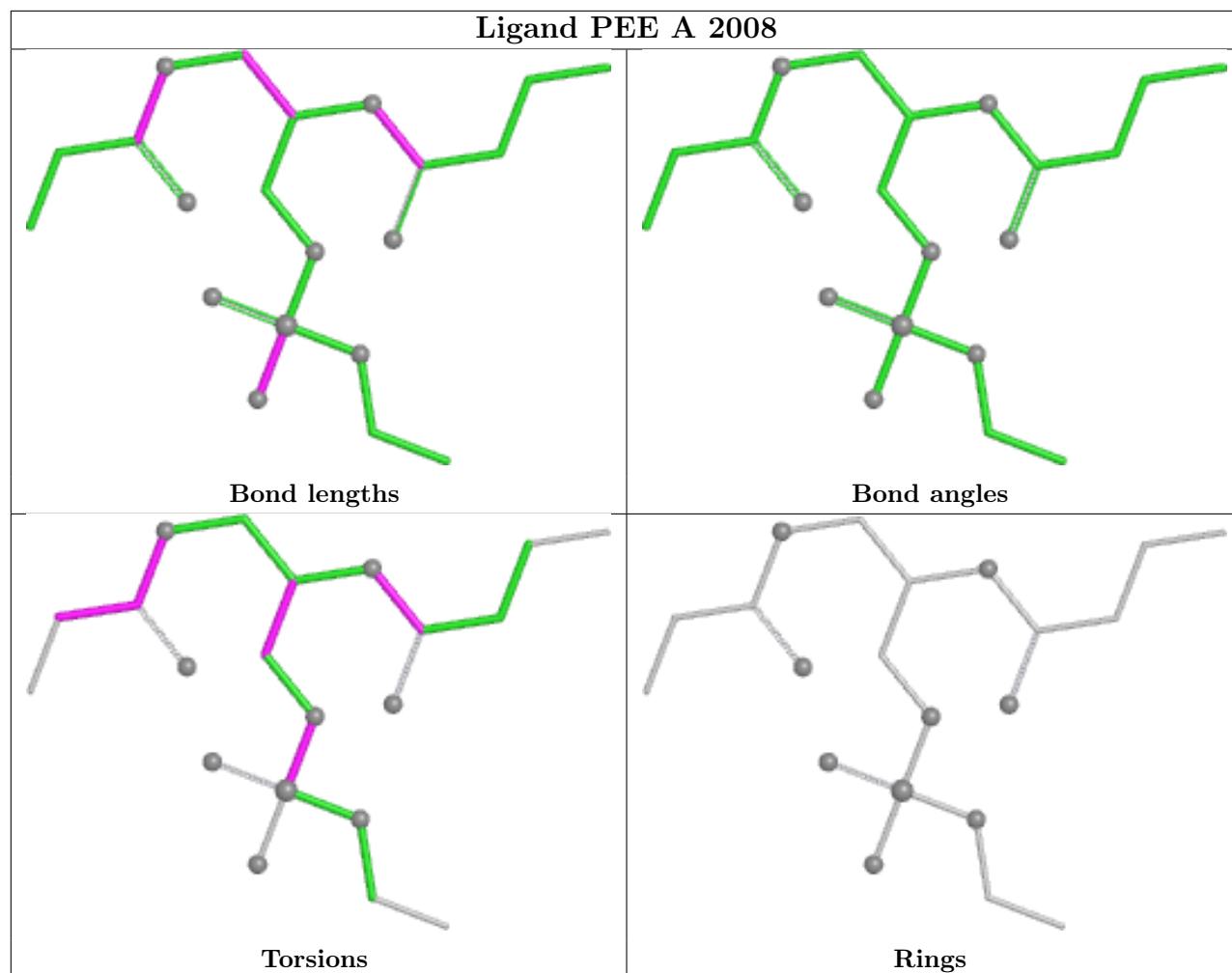


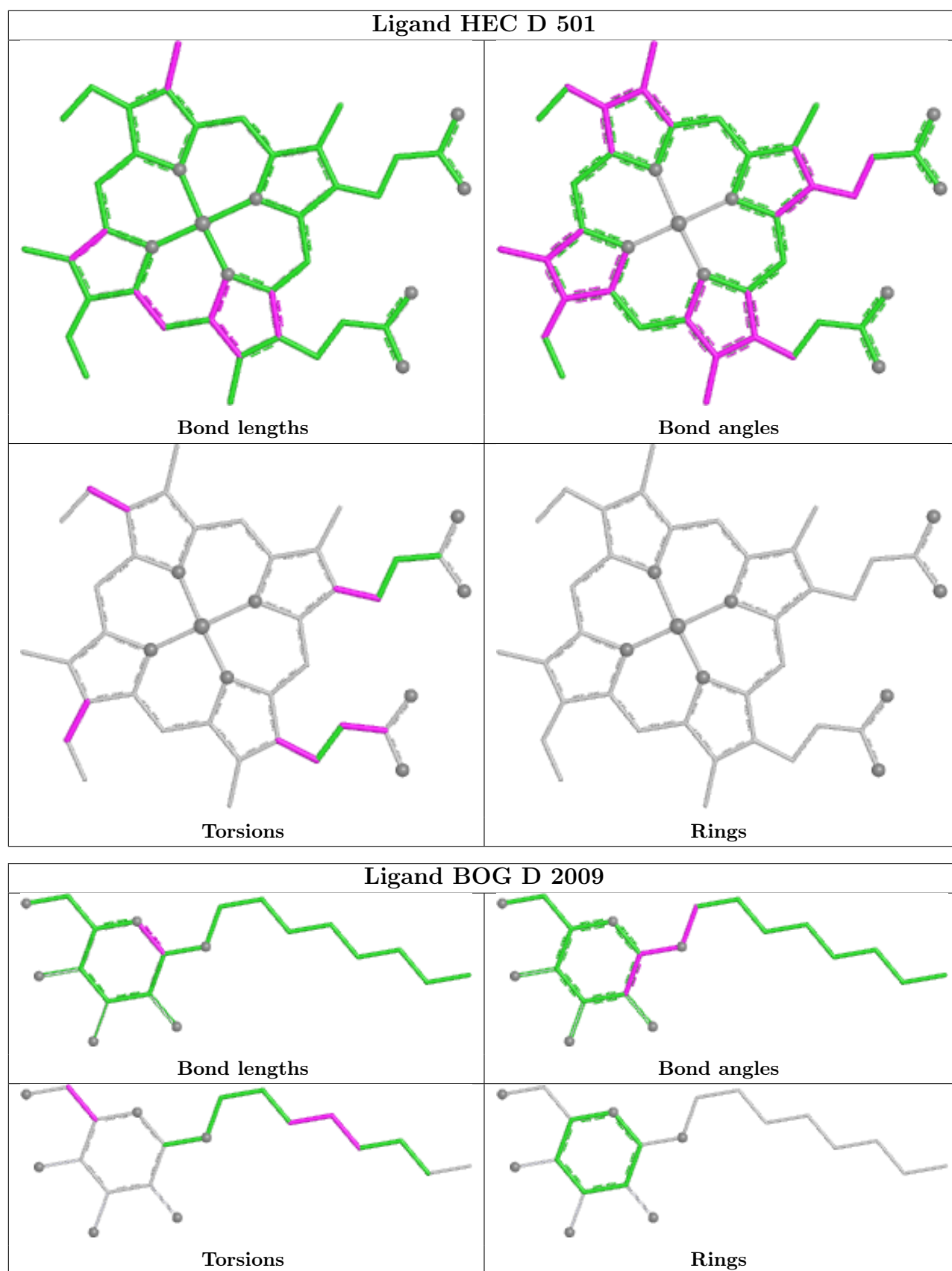


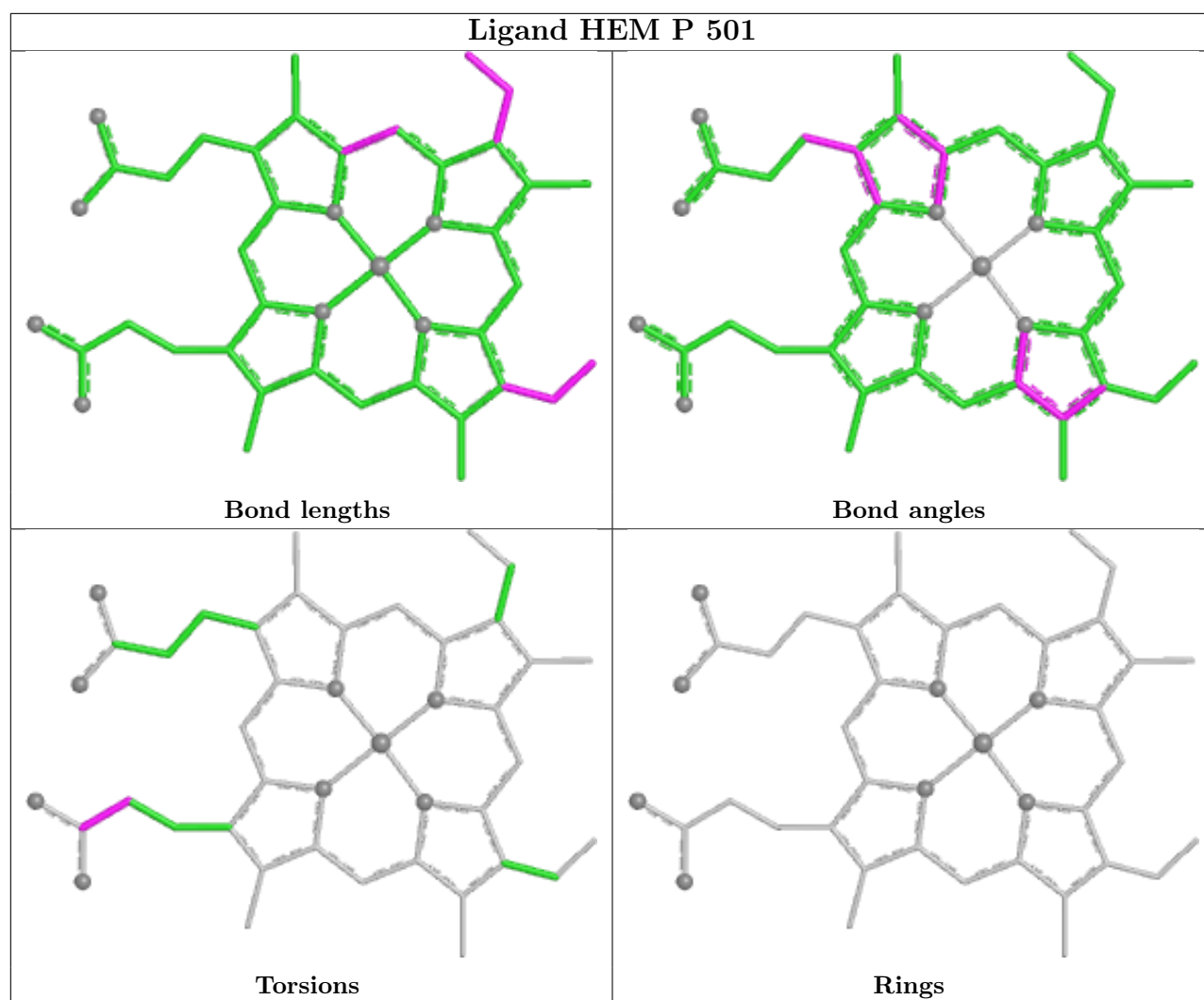


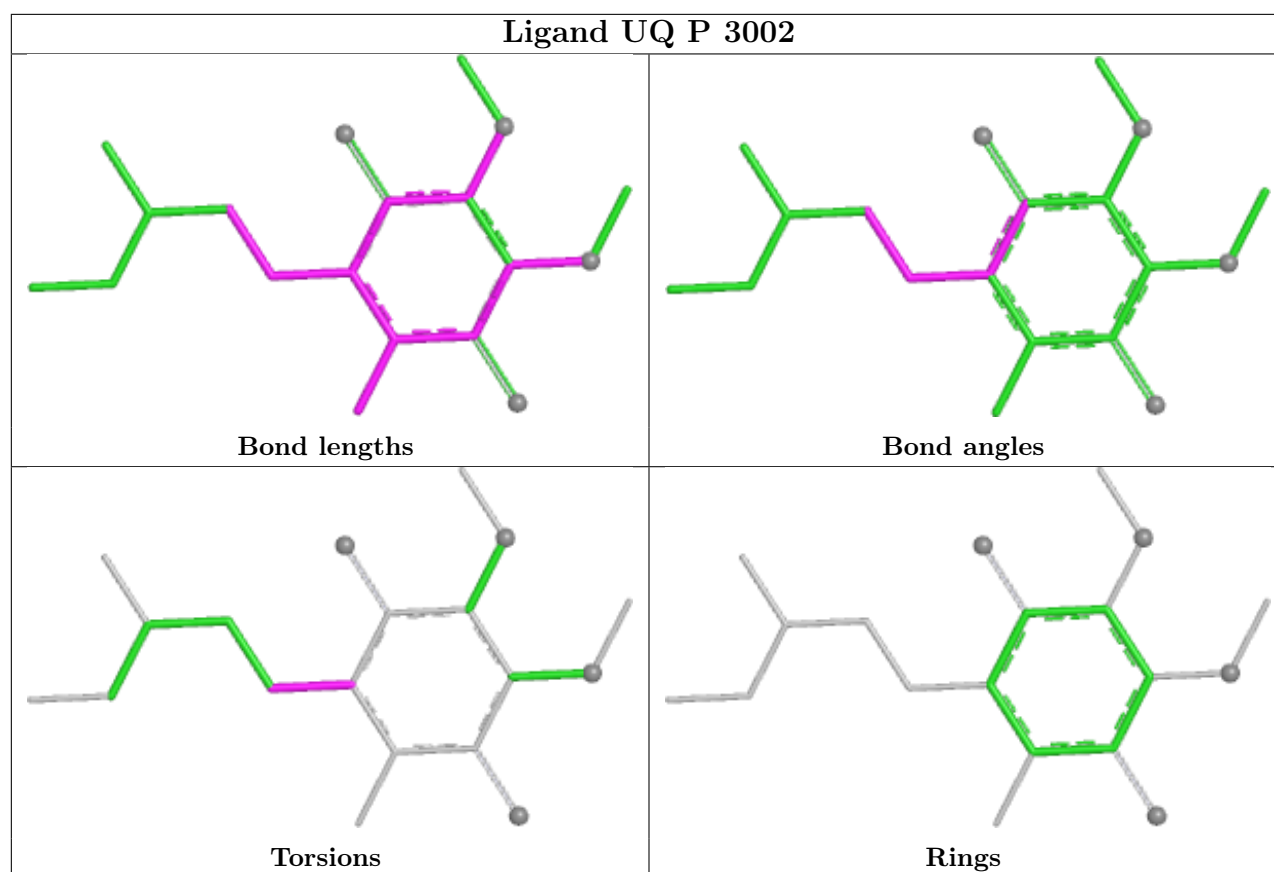


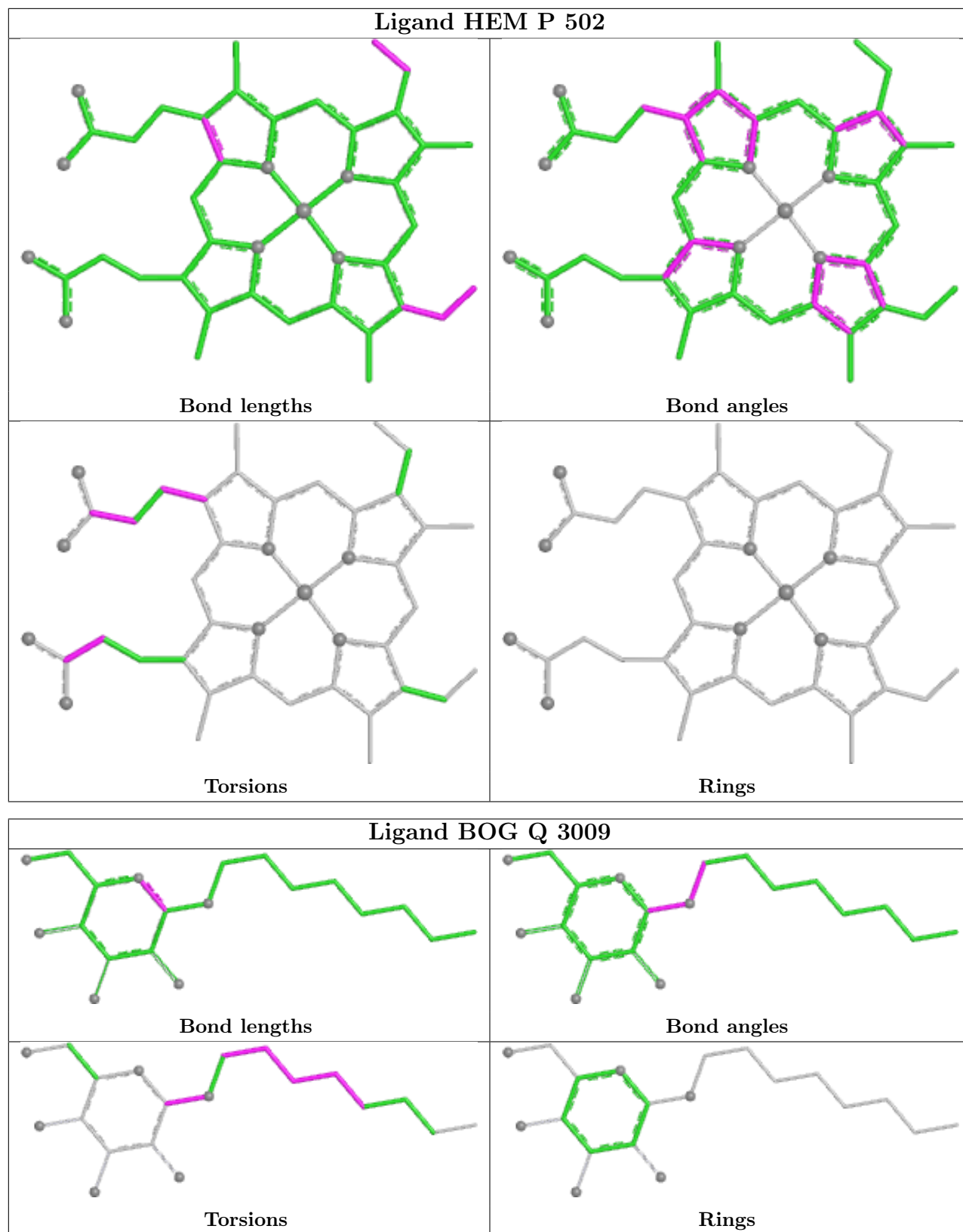


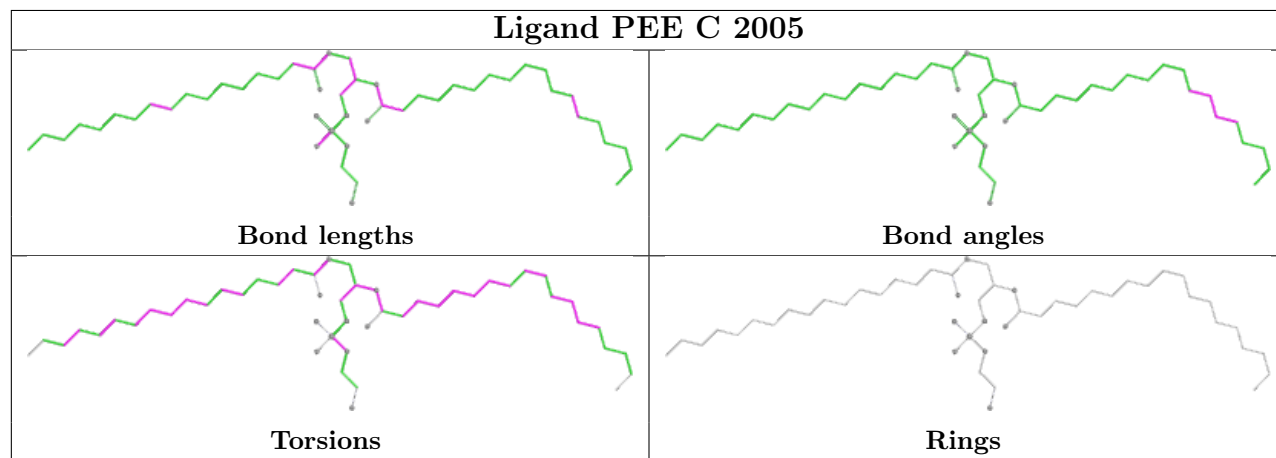
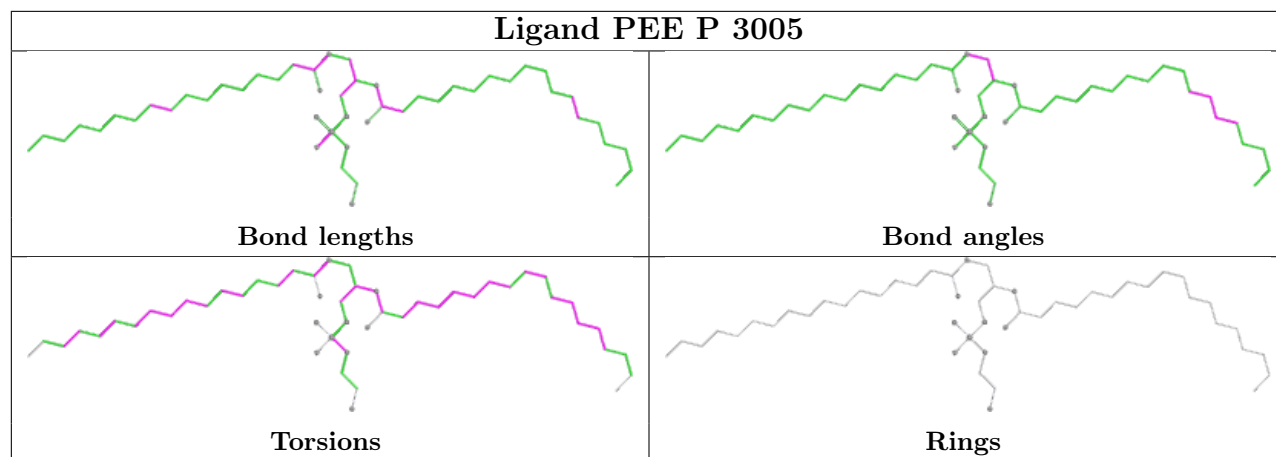


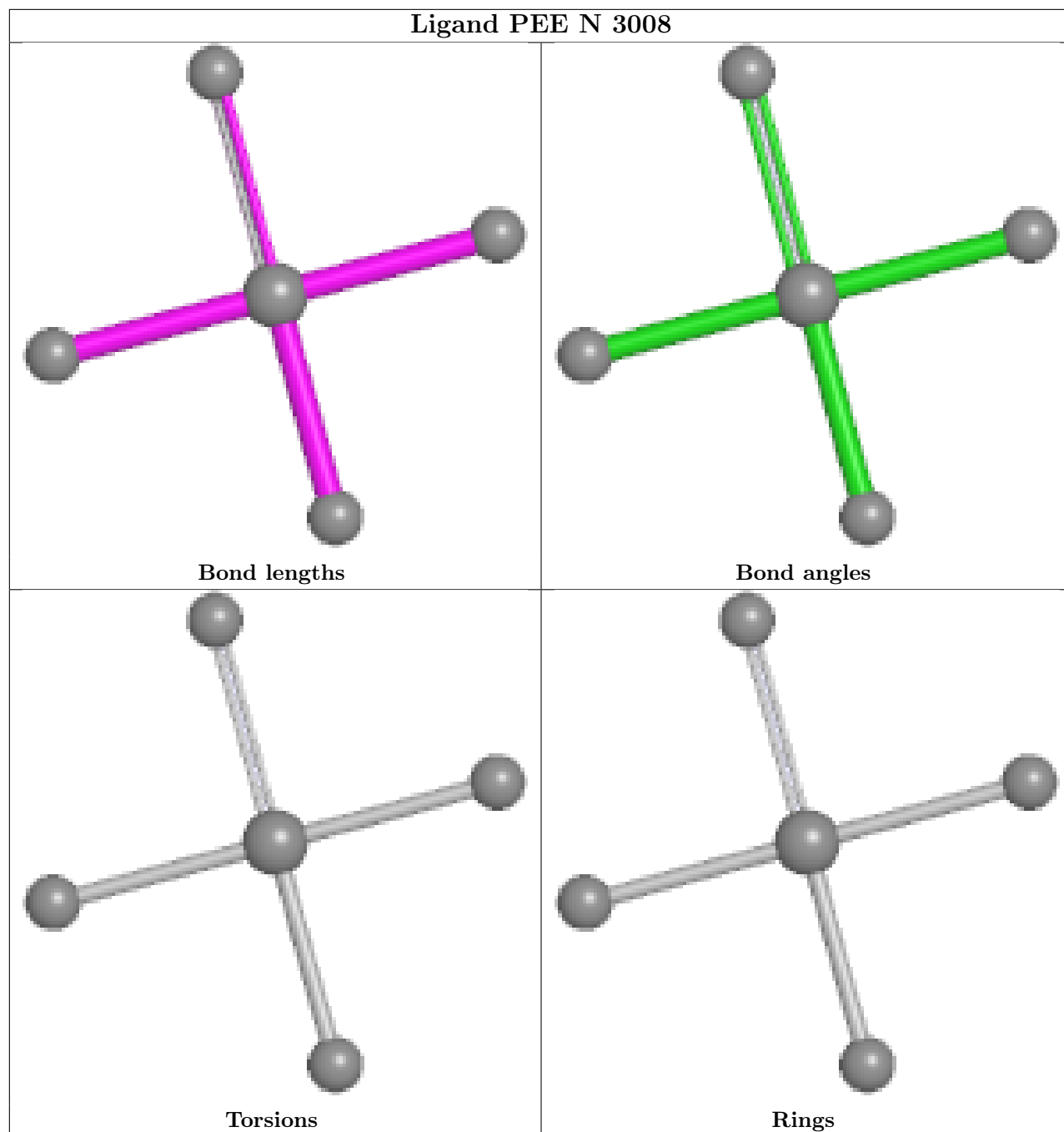


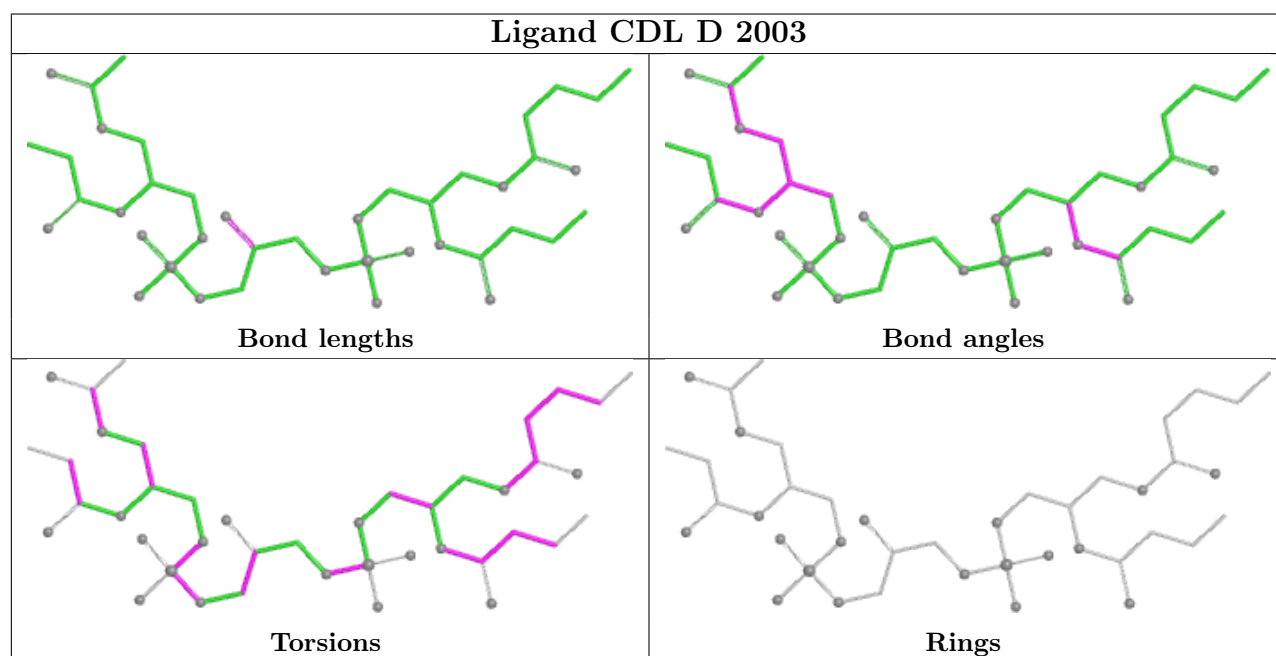
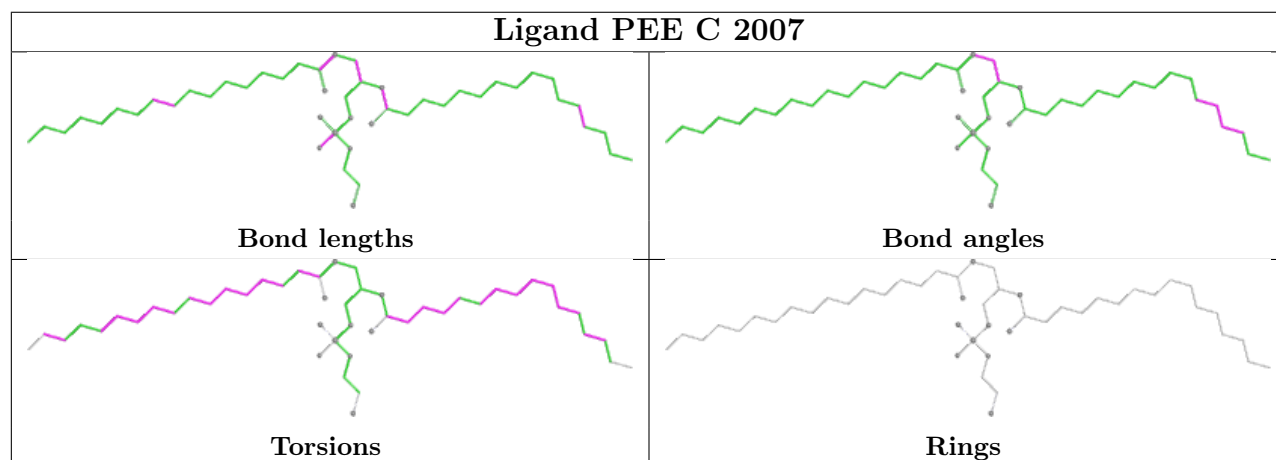
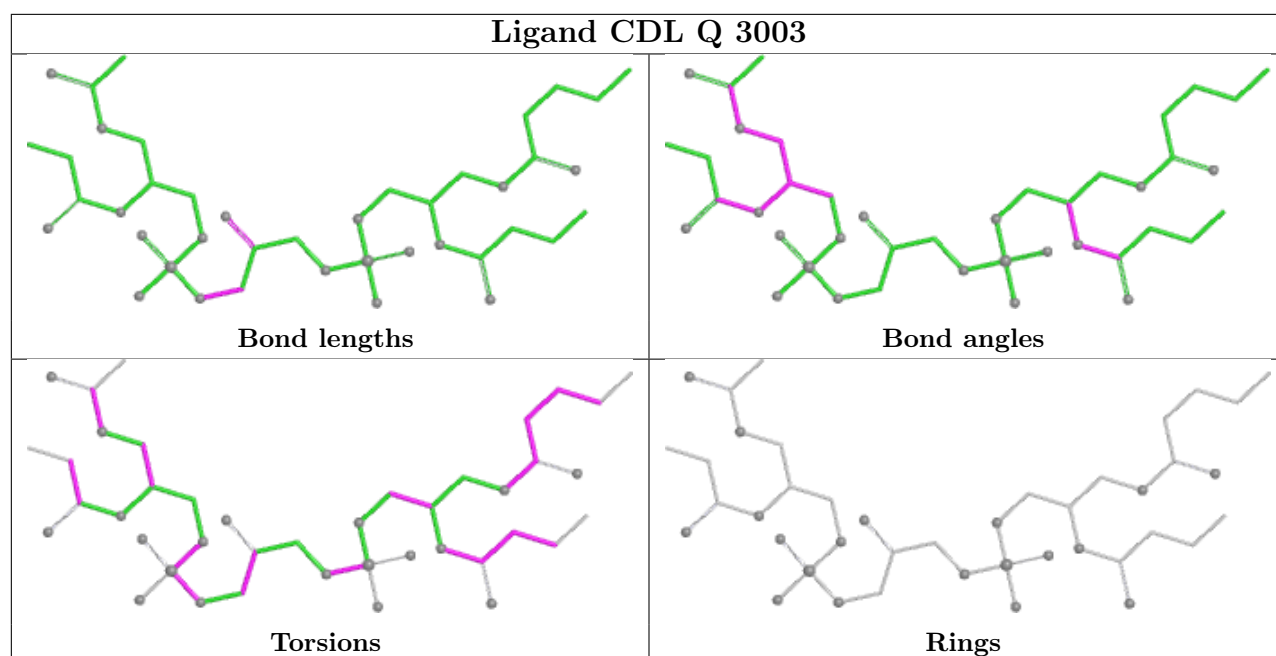




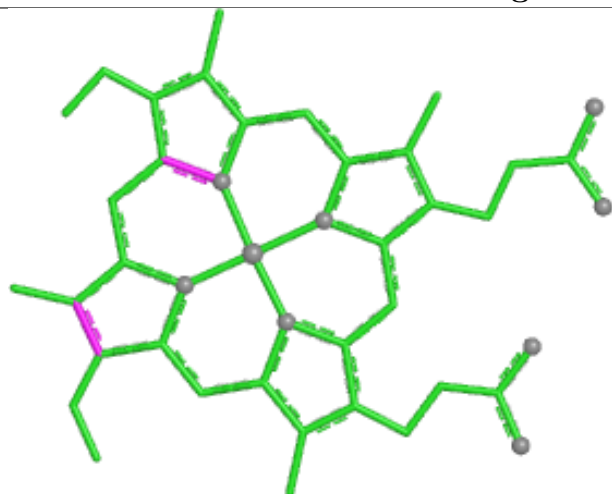




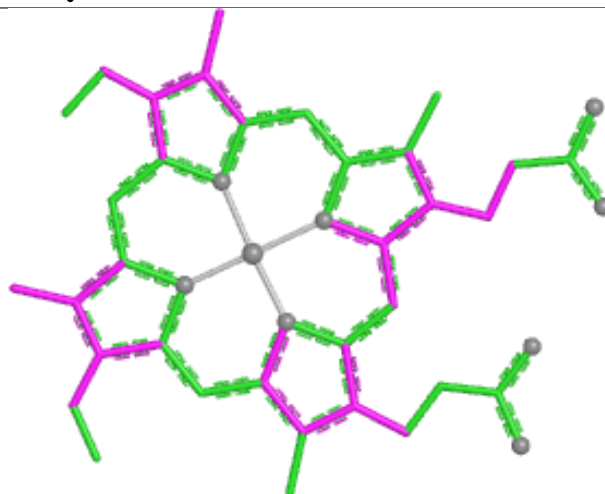




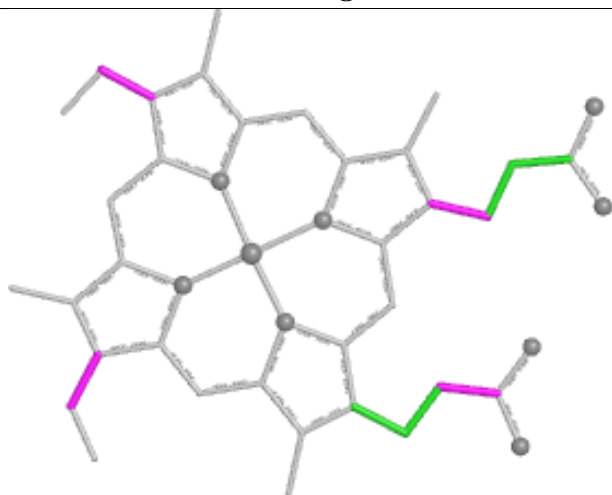
Ligand HEC Q 501



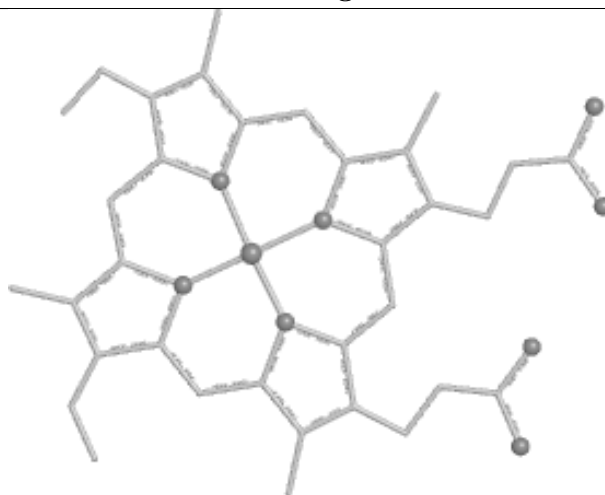
Bond lengths



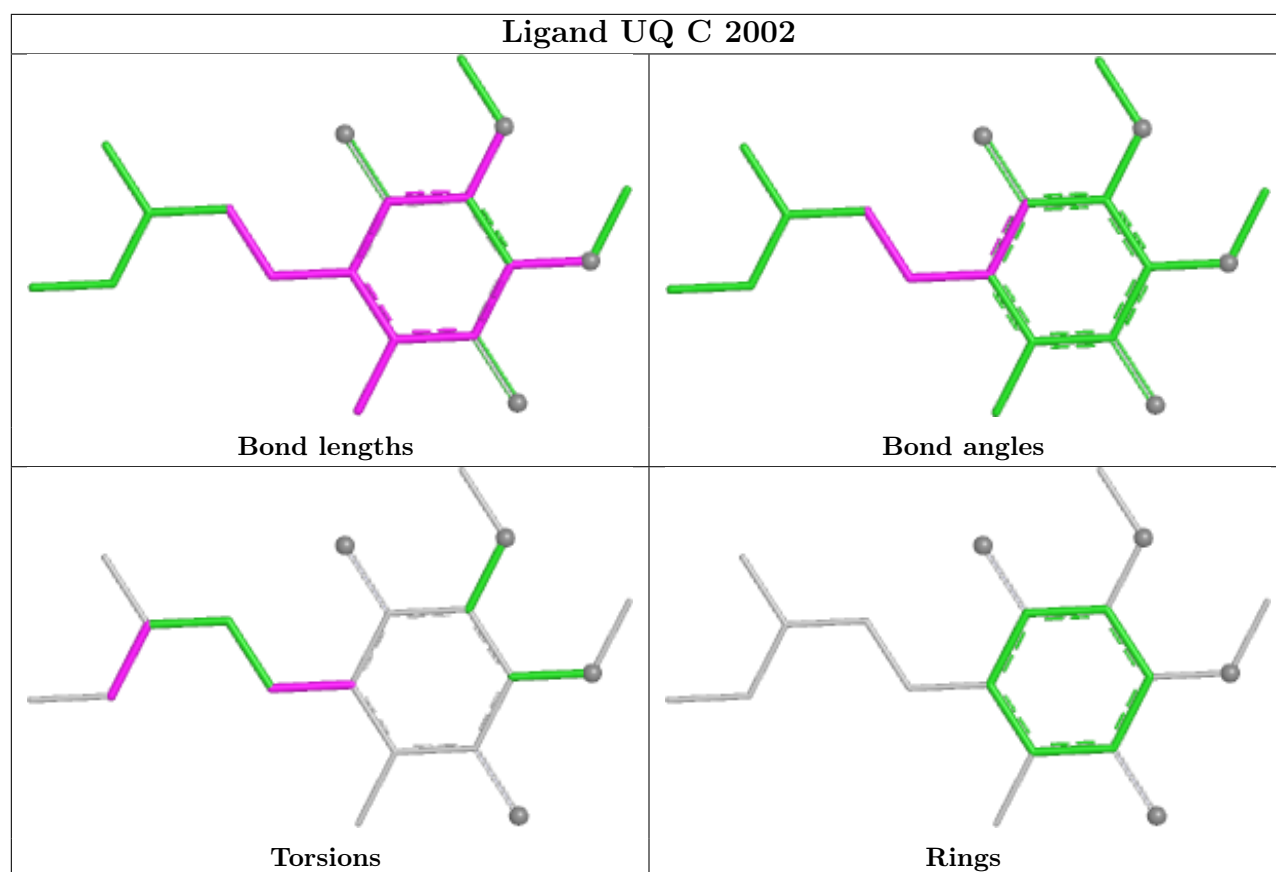
Bond angles



Torsions



Rings



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	443/446 (99%)	0.52	20 (4%) 38 30	29, 61, 87, 104	0
1	N	442/446 (99%)	0.94	45 (10%) 12 9	39, 69, 95, 104	0
2	B	421/441 (95%)	1.05	57 (13%) 7 5	50, 78, 116, 141	0
2	O	422/441 (95%)	0.95	53 (12%) 8 6	37, 74, 106, 120	0
3	C	380/380 (100%)	-0.11	7 (1%) 67 58	21, 39, 73, 109	0
3	P	379/380 (99%)	0.52	23 (6%) 27 20	30, 62, 84, 96	0
4	D	241/241 (100%)	-0.14	4 (1%) 69 60	30, 42, 78, 94	0
4	Q	241/241 (100%)	0.84	25 (10%) 11 8	50, 69, 99, 119	0
5	E	196/196 (100%)	1.24	57 (29%) 1 1	38, 80, 108, 113	0
5	R	196/196 (100%)	2.86	115 (58%) 0 0	42, 99, 140, 144	127 (64%)
6	F	101/110 (91%)	-0.17	2 (1%) 65 56	29, 42, 60, 96	0
6	S	101/110 (91%)	0.79	8 (7%) 18 13	55, 68, 111, 135	0
7	G	81/81 (100%)	0.44	6 (7%) 20 15	32, 52, 93, 108	0
7	T	78/81 (96%)	1.22	15 (19%) 3 2	48, 82, 144, 161	0
8	H	70/77 (90%)	0.41	4 (5%) 29 22	39, 59, 84, 123	0
8	U	67/77 (87%)	2.03	30 (44%) 0 0	96, 115, 133, 140	0
9	I	31/47 (65%)	3.40	26 (83%) 0 0	81, 105, 125, 126	0
9	V	31/47 (65%)	3.76	24 (77%) 0 0	69, 109, 142, 145	0
10	J	61/61 (100%)	0.33	2 (3%) 49 39	44, 60, 92, 130	0
10	W	60/61 (98%)	1.04	9 (15%) 5 4	56, 77, 99, 111	0
All	All	4042/4160 (97%)	0.81	532 (13%) 7 6	21, 65, 110, 161	127 (3%)

The worst 5 of 532 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
9	V	63	ASP	12.5
5	R	195	VAL	10.8
5	R	194	VAL	10.5
5	R	124	LEU	10.0
5	R	133	VAL	10.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 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($\text{\AA}^2$)	Q<0.9
17	BOG	Q	3091	20/20	0.53	0.34	178,186,187,187	0
12	UNL	P	3048	2/-	0.54	0.88	92,92,92,95	0
17	BOG	D	2091	20/20	0.72	0.50	177,189,190,190	0
12	UNL	D	2012	2/-	0.78	0.21	59,59,59,60	0
12	UNL	C	2048	2/-	0.78	0.37	62,62,62,63	0
17	BOG	P	2010	19/20	0.79	0.35	102,183,184,184	0
19	CDL	T	3004	40/100	0.79	0.23	94,97,106,107	0
19	CDL	Q	3003	42/100	0.81	0.26	119,124,137,138	0
19	CDL	D	2003	42/100	0.81	0.22	81,93,97,98	0
11	PEE	A	2008	21/51	0.82	0.24	95,113,116,118	0
12	UNL	P	3046	2/-	0.86	0.17	85,85,85,85	0
15	UQ	C	2002	19/63	0.86	0.27	85,87,89,89	0
12	UNL	P	3014	1/-	0.87	0.17	55,55,55,55	0
19	CDL	G	2004	40/100	0.87	0.17	60,70,85,86	0
11	PEE	C	2005	50/51	0.87	0.25	83,93,101,104	0
16	AZI	C	2011	3/3	0.87	0.25	68,68,69,71	0
15	UQ	P	3002	19/63	0.88	0.33	138,141,143,143	0
11	PEE	N	3008	5/51	0.88	0.11	88,88,89,89	0
11	PEE	P	3005	50/51	0.88	0.26	74,100,105,105	0

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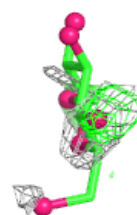
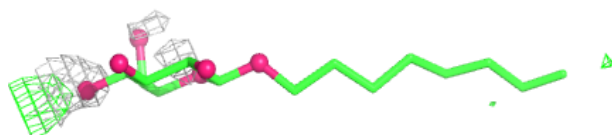
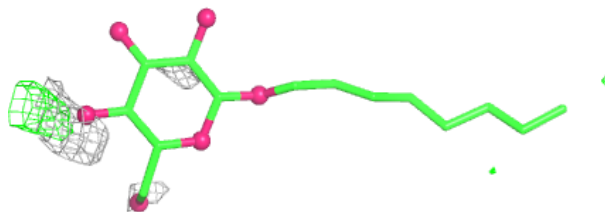
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
17	BOG	Q	3009	20/20	0.90	0.17	71,79,81,82	0
11	PEE	P	3007	48/51	0.90	0.24	76,90,105,106	0
17	BOG	C	3010	12/20	0.90	0.26	94,96,97,98	0
12	UNL	C	2047	1/-	0.91	0.09	34,34,34,34	0
17	BOG	D	2009	20/20	0.91	0.13	56,63,65,67	0
12	UNL	C	2046	2/-	0.92	0.22	87,87,87,91	0
12	UNL	P	3013	1/-	0.92	0.29	56,56,56,56	0
12	UNL	A	3015	1/-	0.92	0.15	48,48,48,48	0
12	UNL	P	3047	1/-	0.94	0.09	47,47,47,47	0
16	AZI	P	3011	3/3	0.94	0.20	74,74,76,77	0
14	FNM	P	3001	22/22	0.95	0.11	49,53,59,62	0
20	FES	R	501	4/4	0.95	0.17	106,106,106,107	4
18	HEC	Q	501	43/43	0.96	0.10	47,55,64,66	0
12	UNL	R	3012	1/-	0.96	0.19	38,38,38,38	0
11	PEE	C	2007	48/51	0.96	0.14	41,52,81,82	0
20	FES	E	501	4/4	0.97	0.08	83,84,85,86	0
14	FNM	C	2001	22/22	0.97	0.07	38,40,42,44	0
13	HEM	P	502	43/43	0.98	0.07	34,40,55,62	0
18	HEC	D	501	43/43	0.98	0.07	23,30,35,37	0
13	HEM	P	501	43/43	0.98	0.09	40,43,54,58	0
13	HEM	C	501	43/43	0.99	0.07	23,31,38,43	0
13	HEM	C	502	43/43	0.99	0.06	23,26,33,39	0

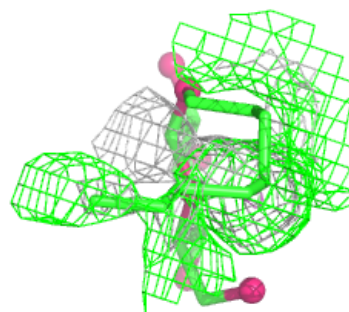
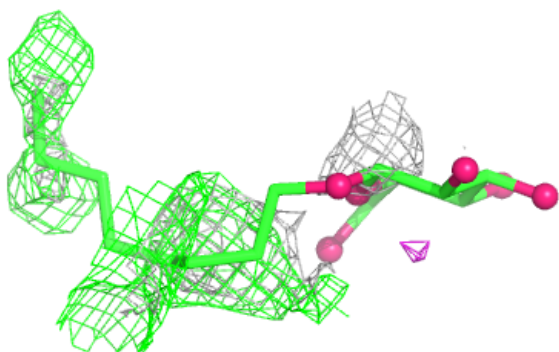
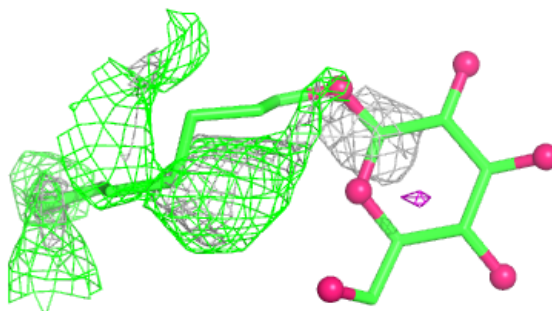
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around BOG Q 3091:

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

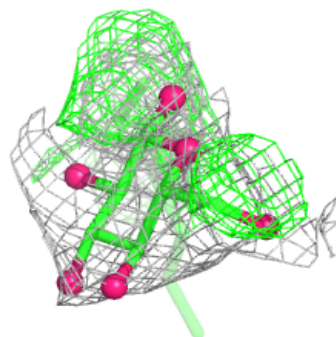
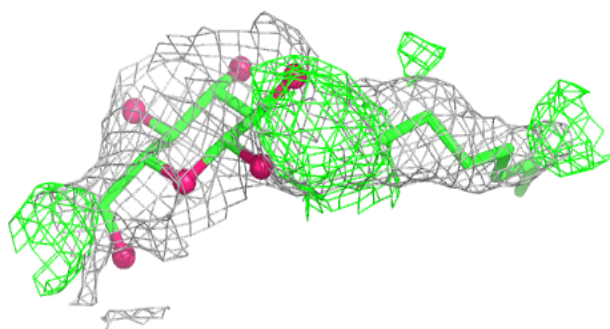
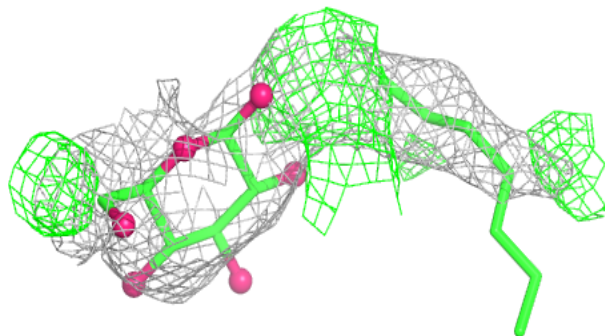
**Electron density around BOG D 2091:**

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



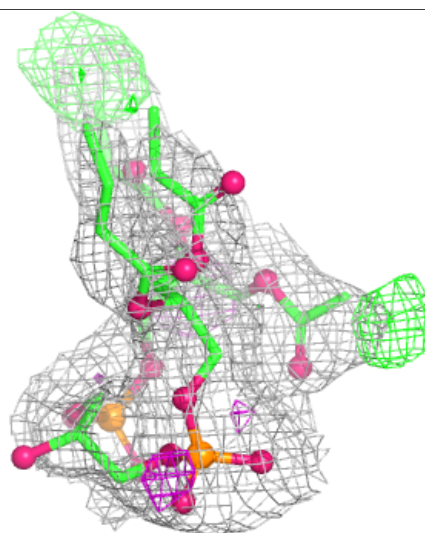
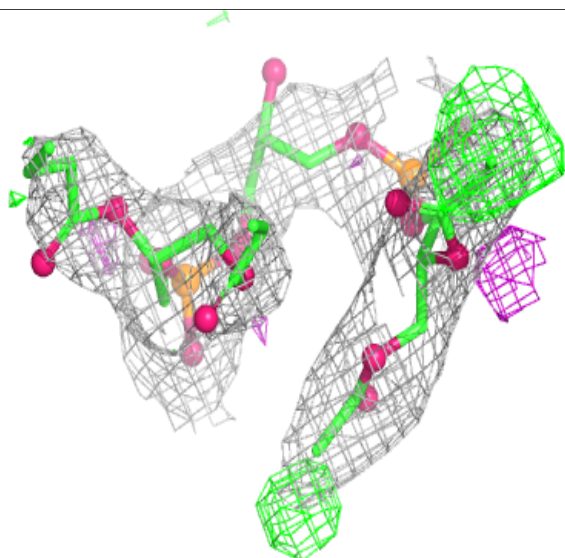
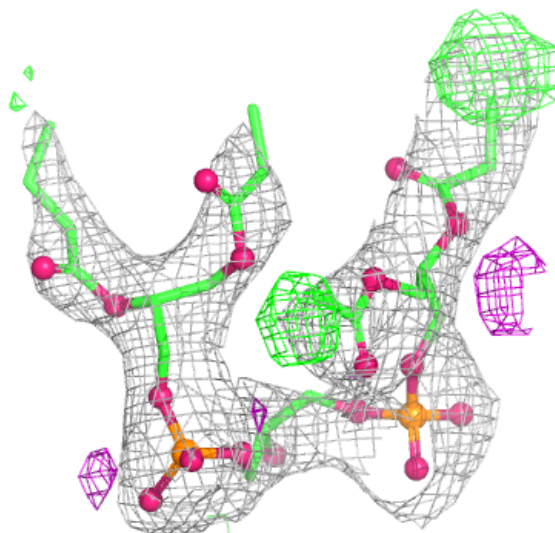
Electron density around BOG P 2010:

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



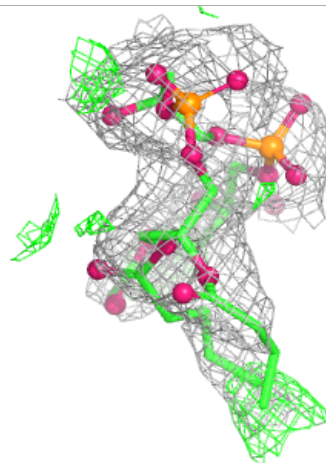
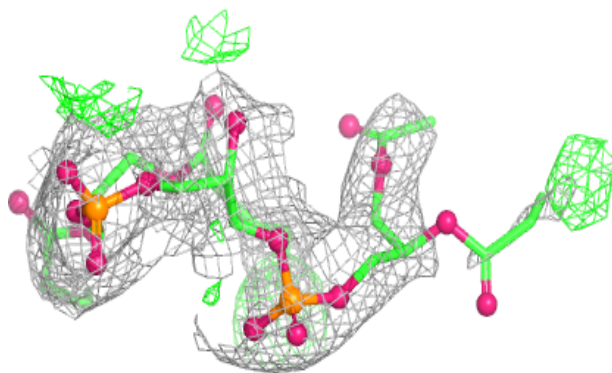
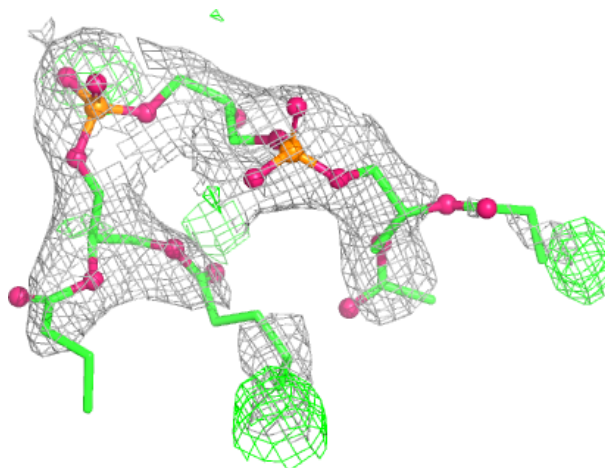
Electron density around CDL T 3004:

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



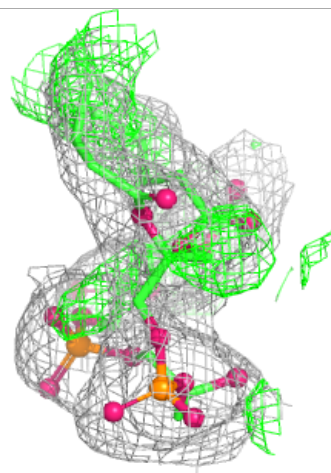
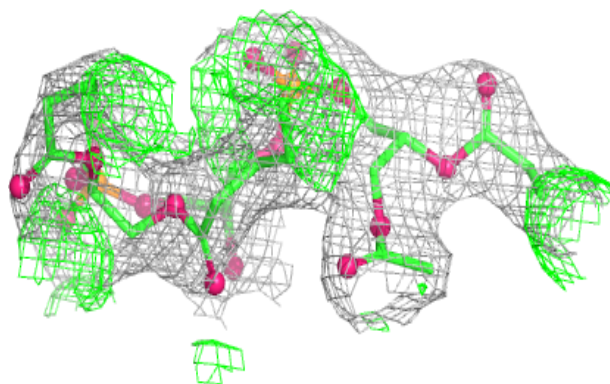
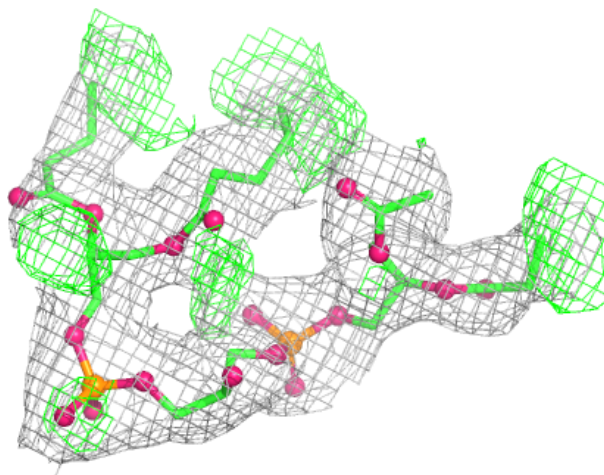
Electron density around CDL Q 3003:

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



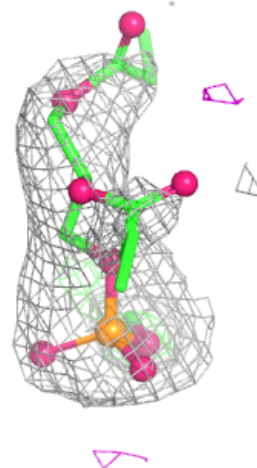
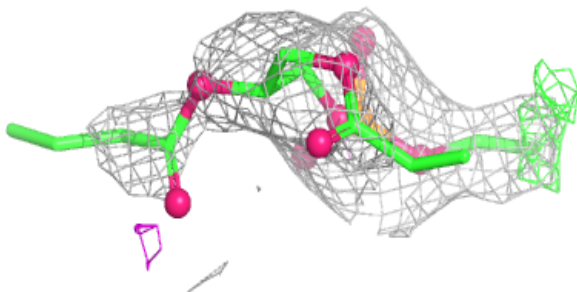
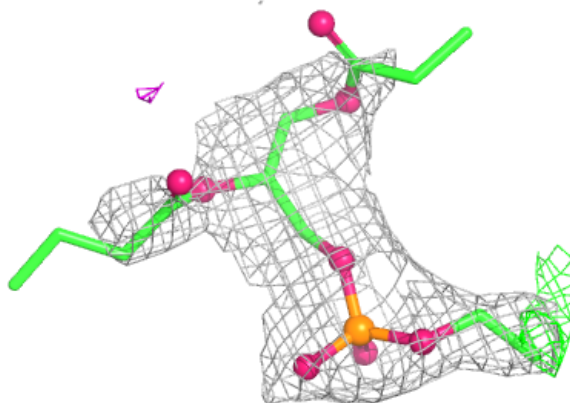
Electron density around CDL D 2003:

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



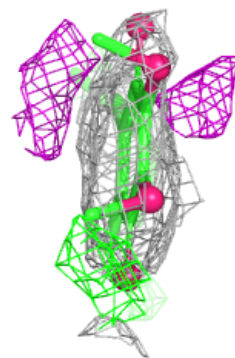
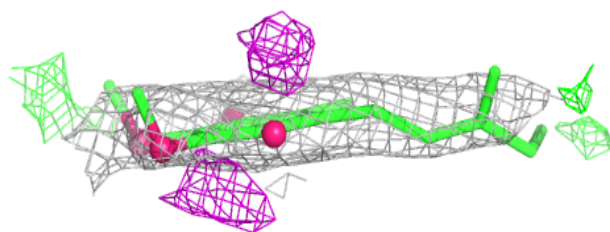
Electron density around PEE A 2008:

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



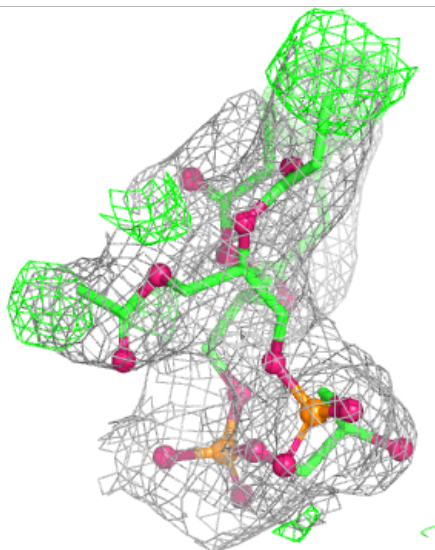
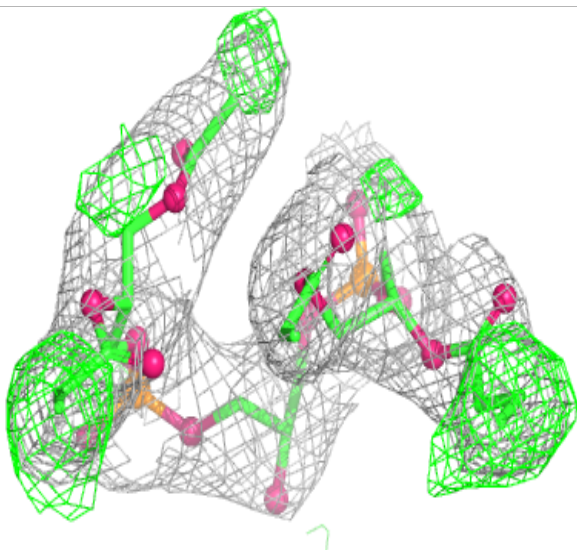
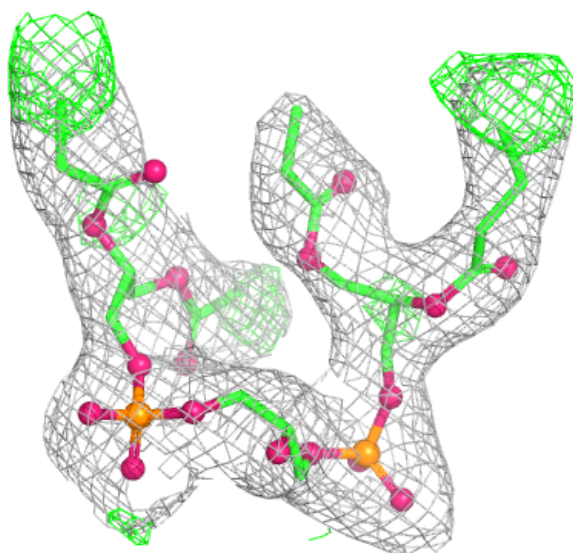
Electron density around UQ C 2002:

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



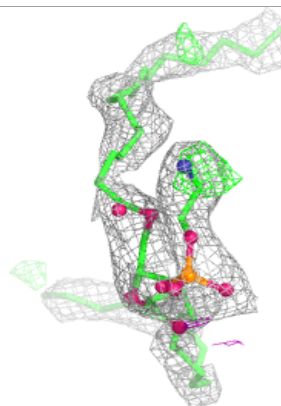
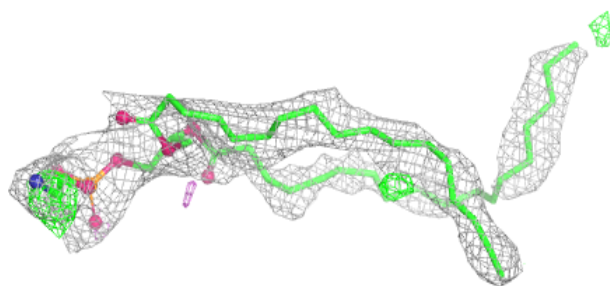
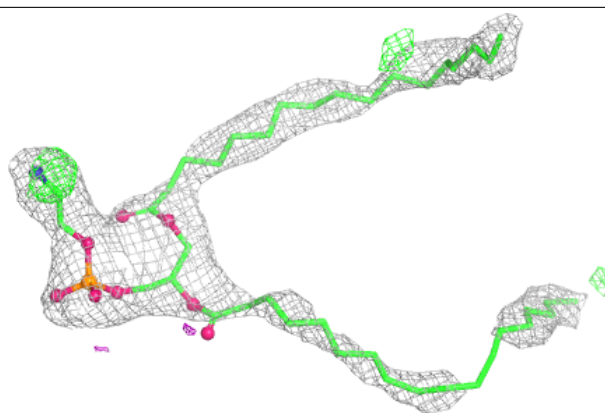
Electron density around CDL G 2004:

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

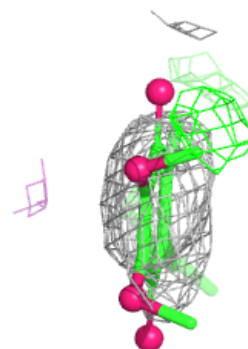
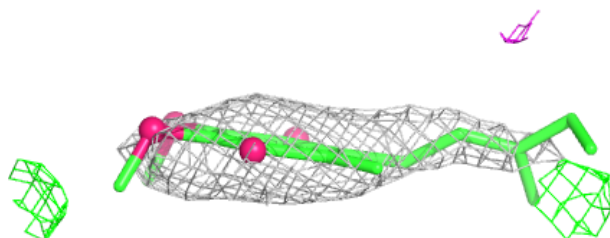
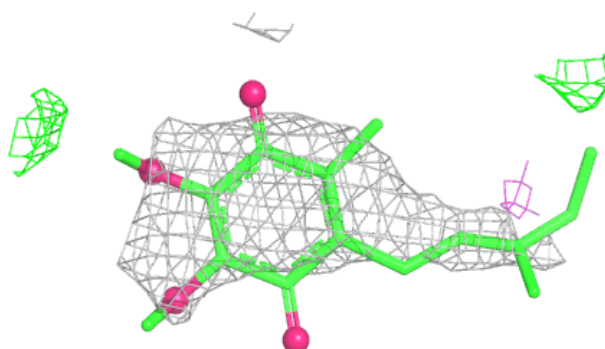


Electron density around PEE C 2005:

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

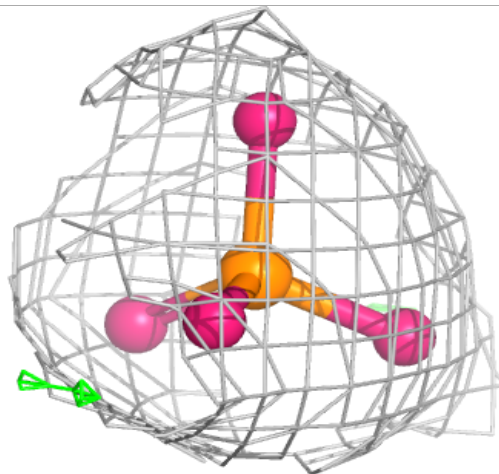
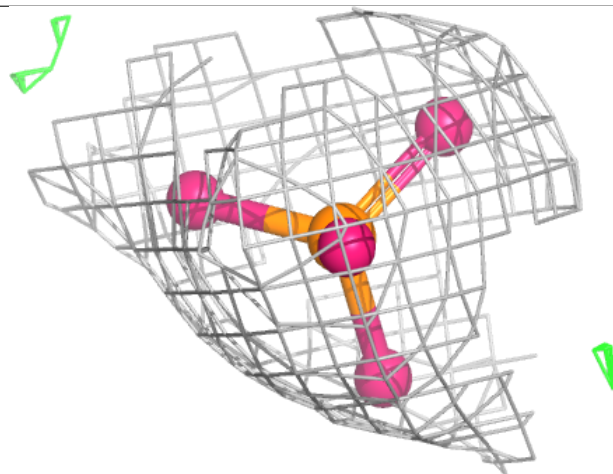
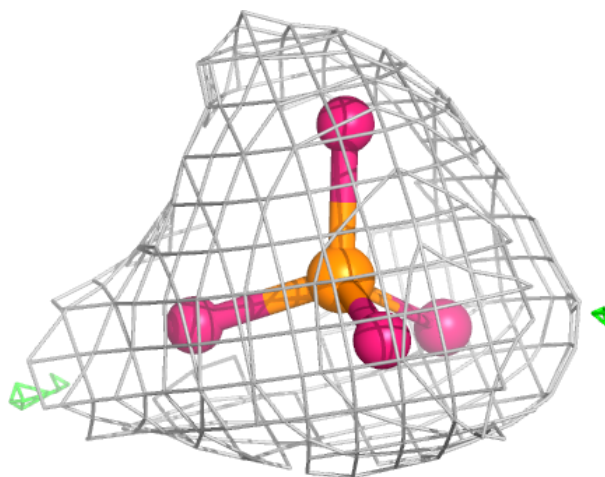
**Electron density around UQ P 3002:**

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



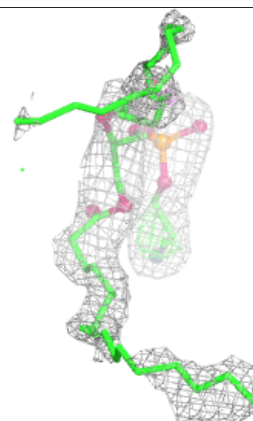
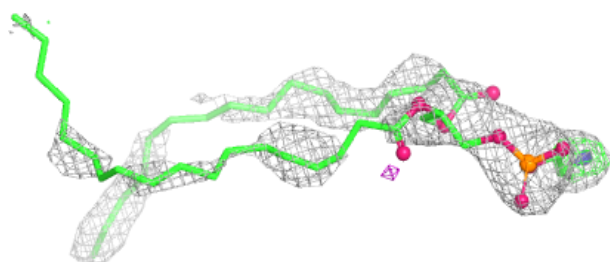
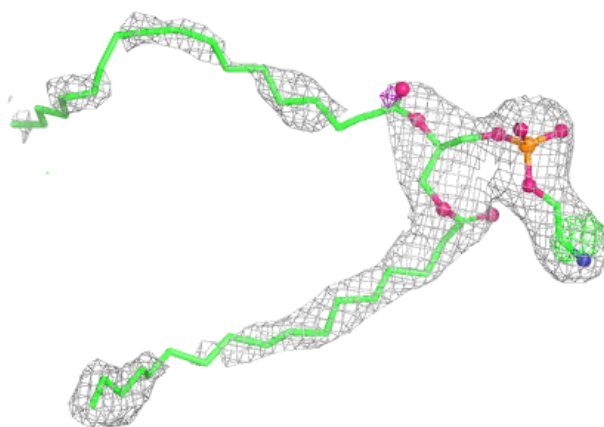
Electron density around PEE N 3008:

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

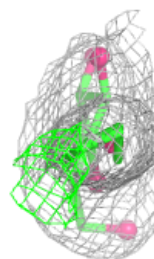
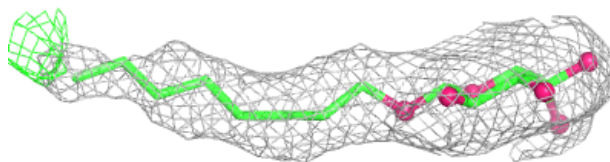
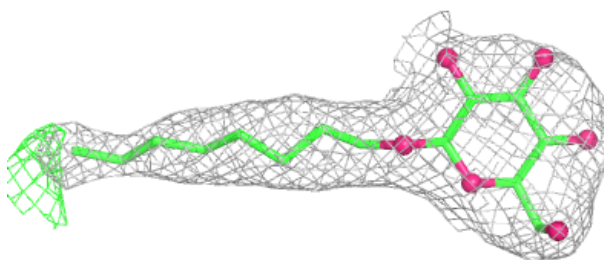


Electron density around PEE P 3005:

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

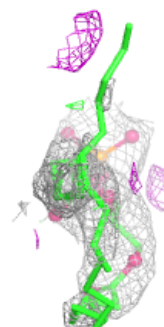
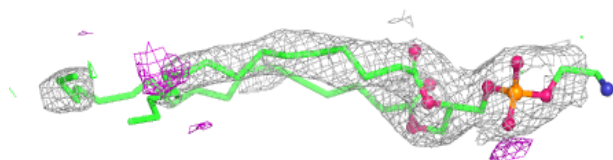
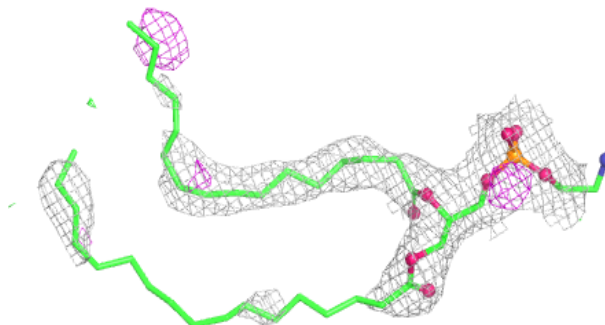
**Electron density around BOG Q 3009:**

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

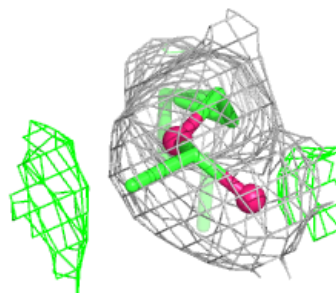
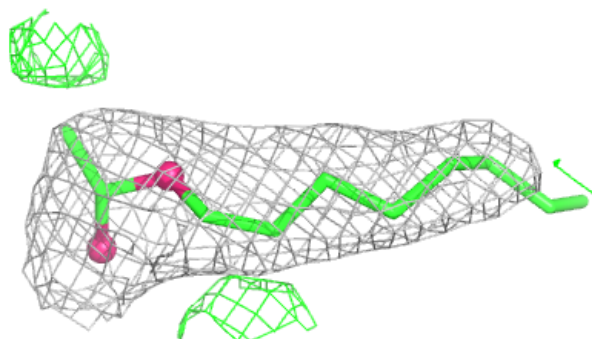
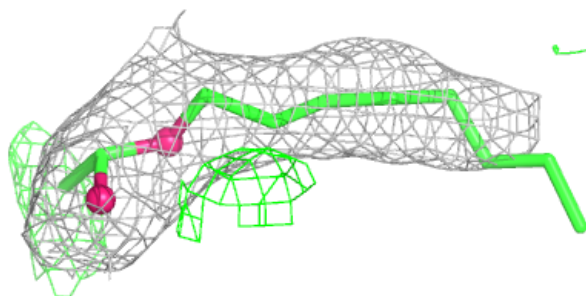


Electron density around PEE P 3007:

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

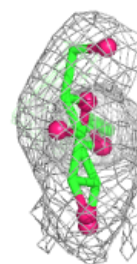
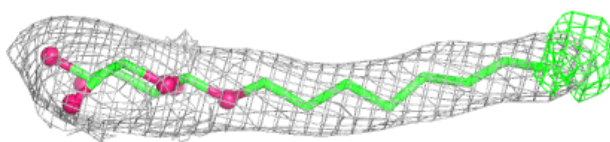
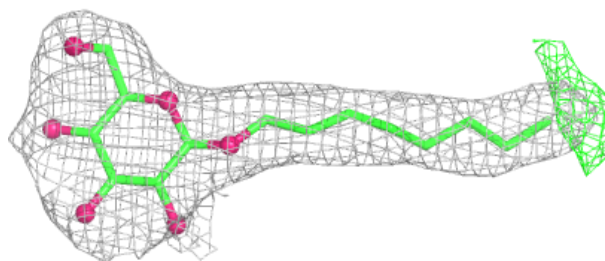
**Electron density around BOG C 3010:**

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

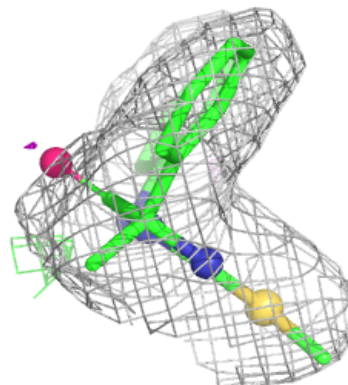
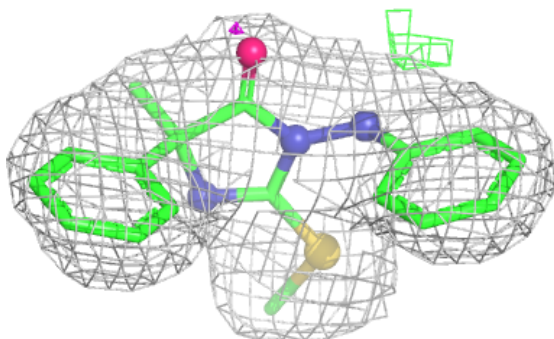
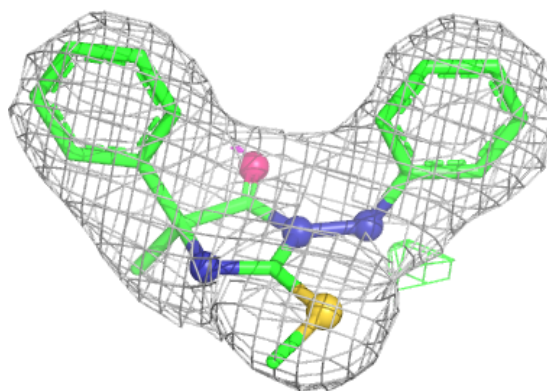


Electron density around BOG D 2009:

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

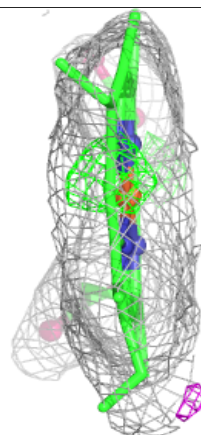
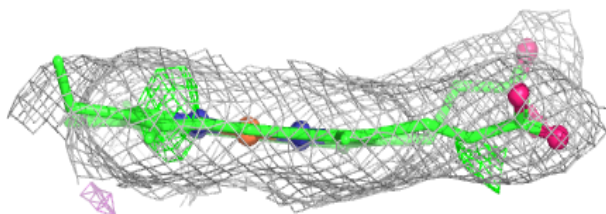
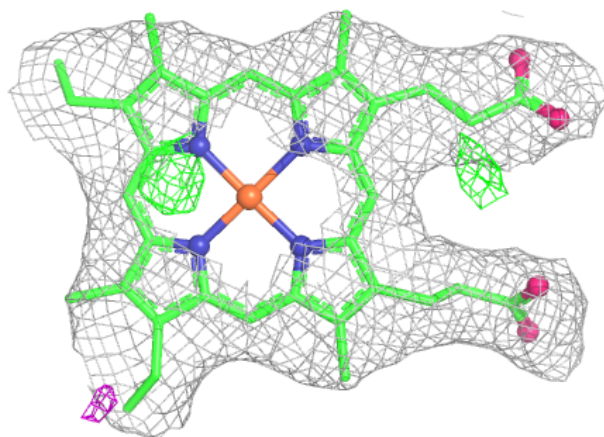
**Electron density around FNM P 3001:**

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

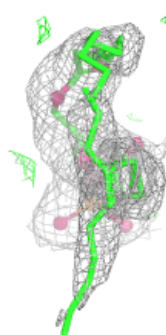
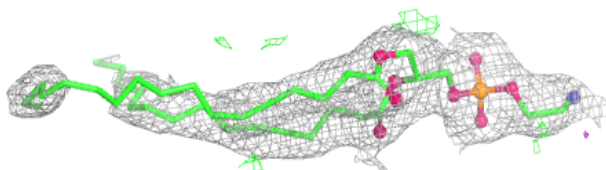
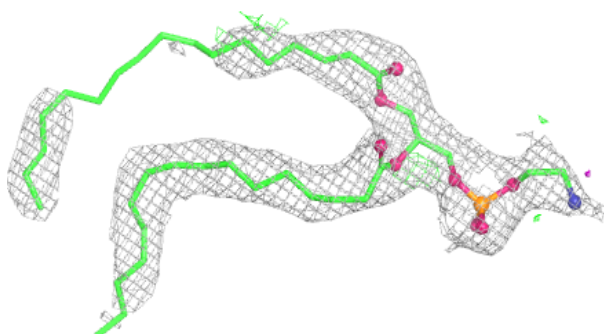


Electron density around HEC Q 501:

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

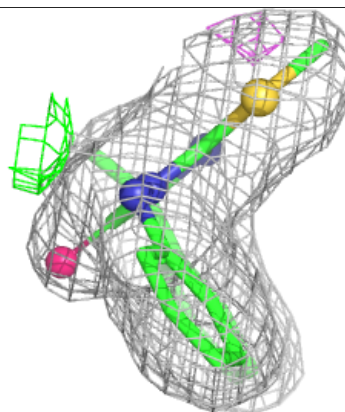
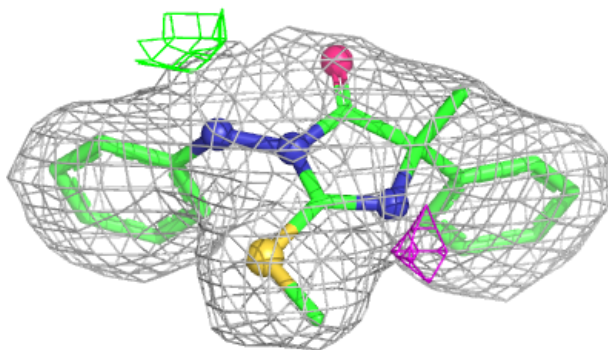
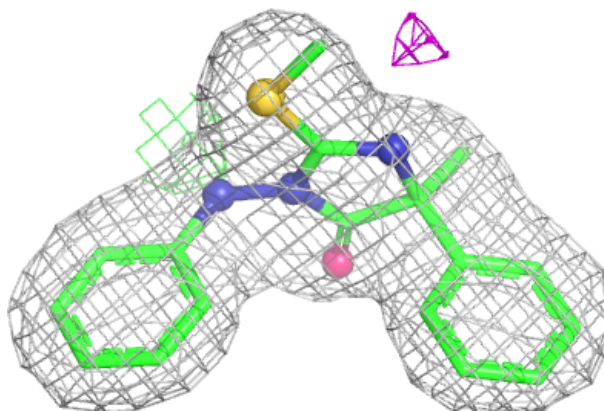
**Electron density around PEE C 2007:**

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



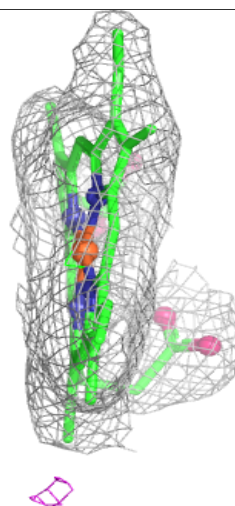
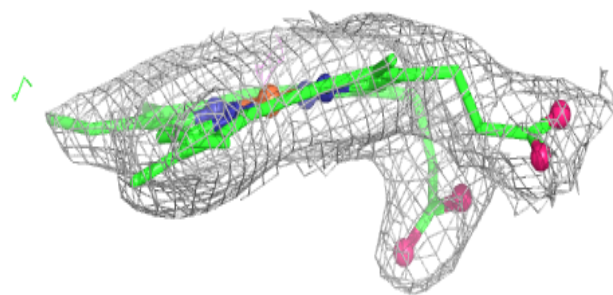
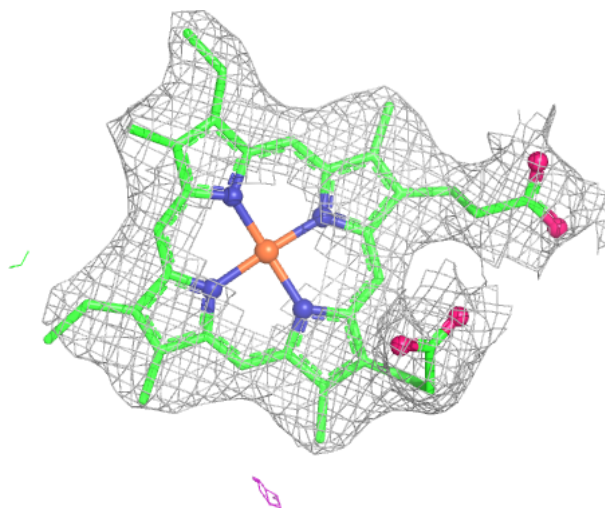
Electron density around FNM C 2001:

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



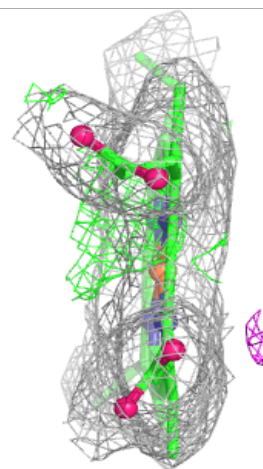
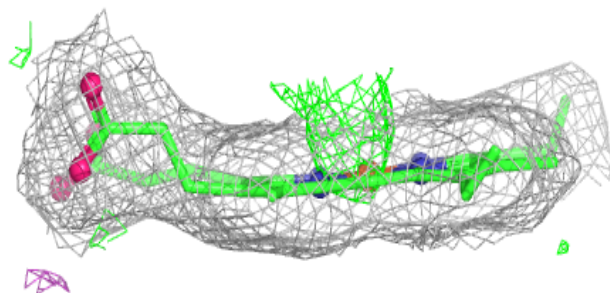
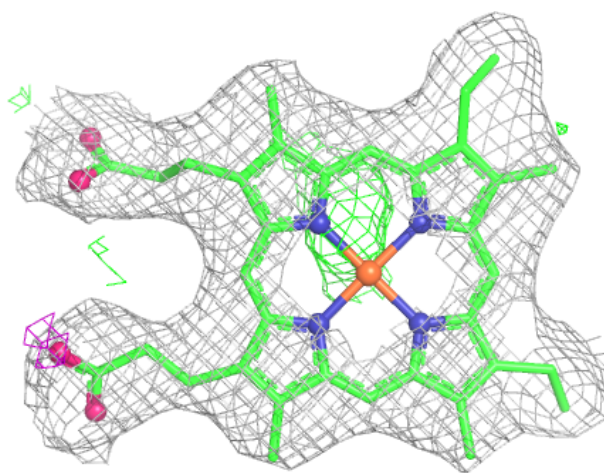
Electron density around HEM P 502:

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



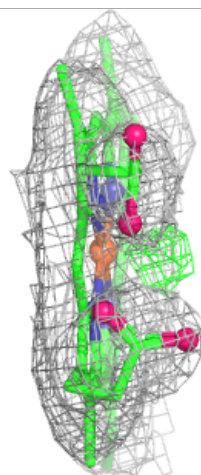
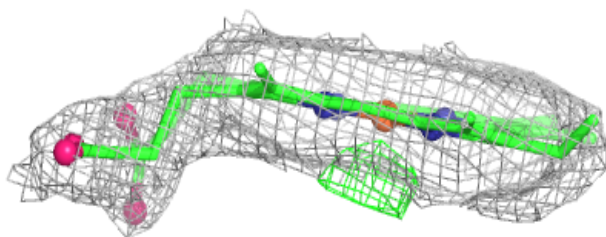
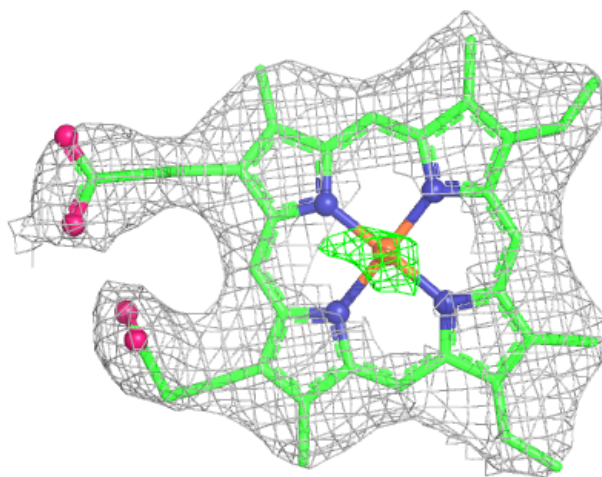
Electron density around HEC D 501:

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



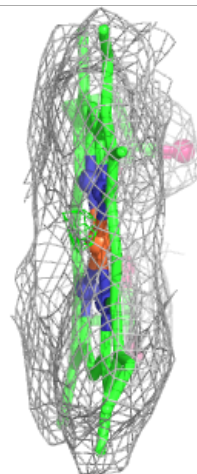
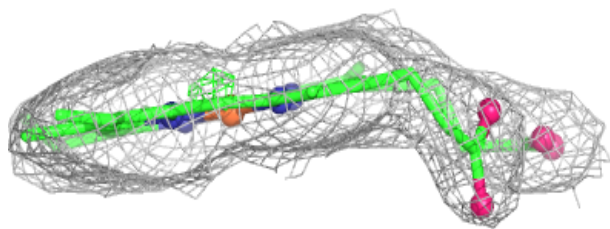
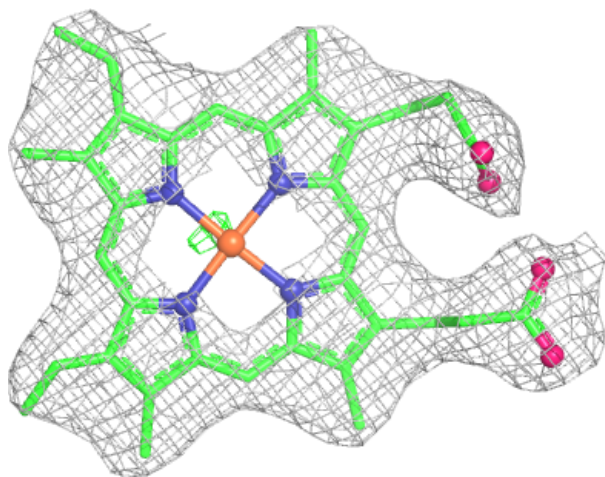
Electron density around HEM P 501:

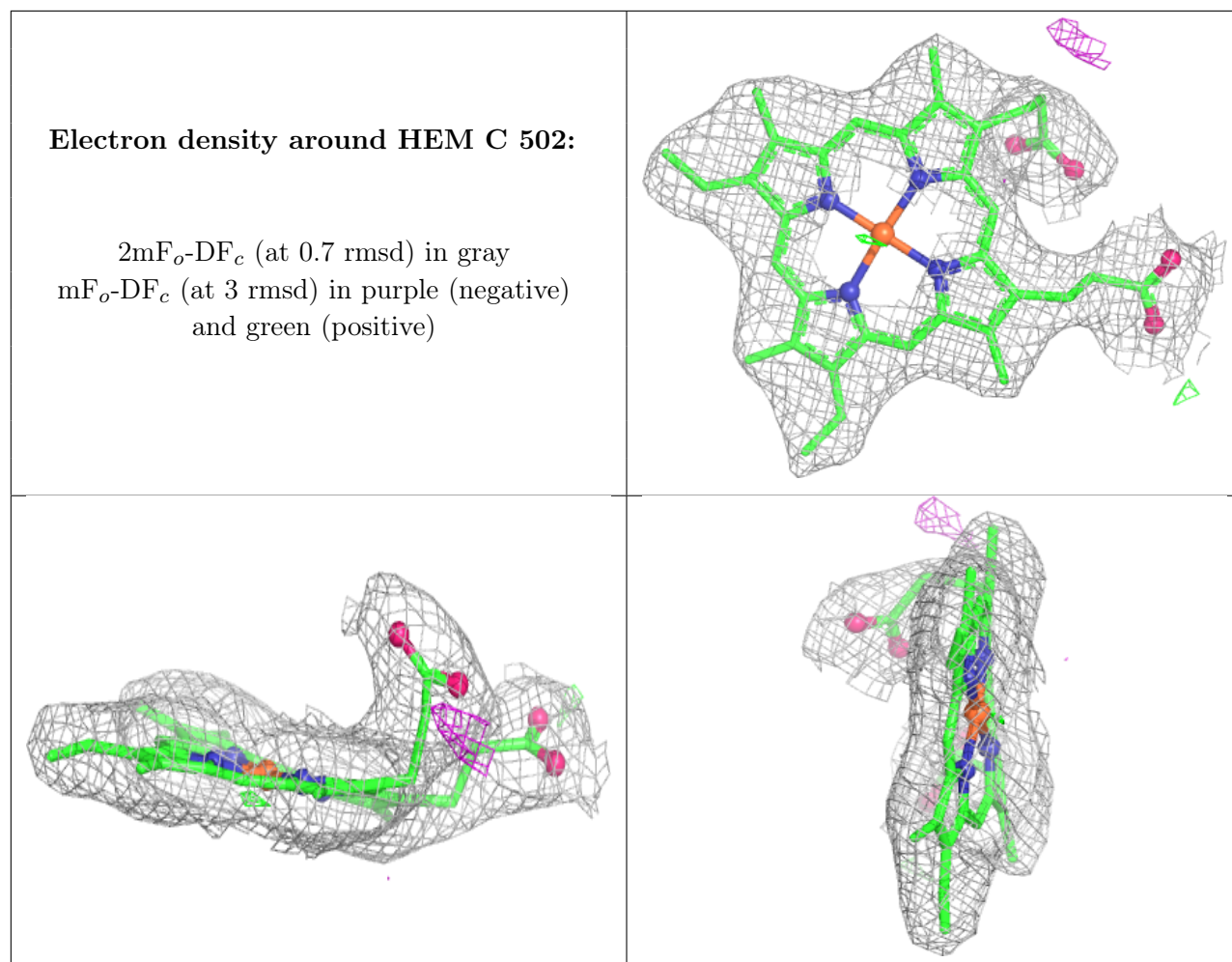
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around HEM C 501:

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





6.5 Other polymers ⓘ

There are no such residues in this entry.