



wwPDB X-ray Structure Validation Summary Report ⓘ

Aug 5, 2026 – 03:39 PM JST

PDB ID : 6JY3 / pdb_00006jy3
Title : Monomeric Form of Bovine Heart Cytochrome c Oxidase in the Fully Oxidized State
Authors : Shinzawa-Itoh, K.; Muramoto, K.
Deposited on : 2019-04-26
Resolution : 1.85 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	:	1.8.5 (274361), CSD as541be (2020)
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

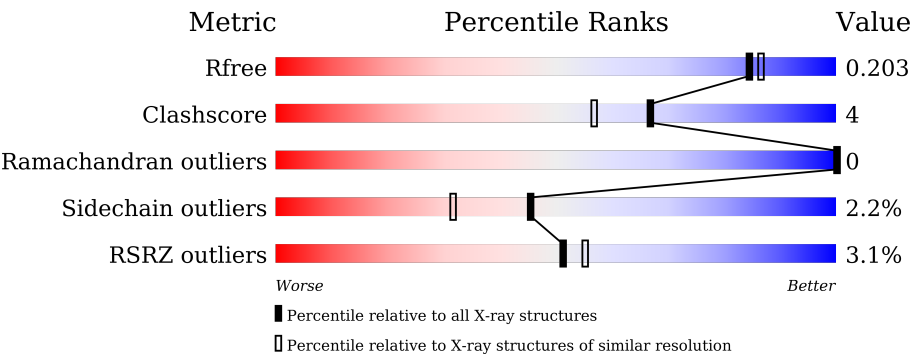
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.85 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	3428 (1.86-1.86)
Clashscore	190562	3579 (1.86-1.86)
Ramachandran outliers	187476	3553 (1.86-1.86)
Sidechain outliers	187428	3553 (1.86-1.86)
RSRZ outliers	180081	3429 (1.86-1.86)

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	514	0% 93%6%
2	B	227	4% 85%12% .
3	C	261	0% 90%8% .
4	D	147	3% 86%7%7%
5	E	109	3% 87%6%6%
6	F	98	5% 82%7%.7%

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Mol	Chain	Length	Quality of chain
7	G	85	8% 73% 11% • 15%
8	H	85	4% 84% 5% 12%
9	I	73	7% 85% 10% • •
10	J	59	7% 88% 5% 7%
11	K	56	4% 80% 7% 12%
12	L	47	2% 79% 13% • 6%
13	M	46	80% • • 13%

2 Entry composition

There are 26 unique types of molecules in this entry. The entry contains 15245 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 Cytochrome c oxidase subunit 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	514	Total	C	N	O	S	0	0	0
			4027	2691	623	678	35			

- Molecule 2 is a protein called Cytochrome c oxidase subunit 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	227	Total	C	N	O	S	0	0	0
			1824	1185	281	340	18			

- Molecule 3 is a protein called Cytochrome c oxidase subunit 3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	254	Total	C	N	O	S	0	0	0
			2061	1379	327	343	12			

- Molecule 4 is a protein called Cytochrome c oxidase subunit 4 isoform 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	136	Total	C	N	O	S	0	0	0
			1133	740	186	203	4			

- Molecule 5 is a protein called Cytochrome c oxidase subunit 5A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	E	102	Total	C	N	O	S	0	0	0
			825	528	139	156	2			

- Molecule 6 is a protein called Cytochrome c oxidase subunit 5B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	F	91	Total	C	N	O	S	0	0	0
			694	432	122	135	5			

- Molecule 7 is a protein called Cytochrome c oxidase subunit 6A2, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
7	G	72	Total	C	N	O	S	0	0	0
			595	387	113	94	1			

- Molecule 8 is a protein called Cytochrome c oxidase subunit 6B1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
8	H	75	Total	C	N	O	S	0	0	0
			628	395	114	114	5			

- Molecule 9 is a protein called Cytochrome c oxidase subunit 6C.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
9	I	70	Total	C	N	O	S	0	0	0
			575	375	103	93	4			

- Molecule 10 is a protein called Cytochrome c oxidase subunit 7A1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
10	J	55	Total	C	N	O	S	0	0	0
			434	280	72	79	3			

- Molecule 11 is a protein called Cytochrome c oxidase subunit 7B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
11	K	49	Total	C	N	O	S	0	0	0
			384	250	65	67	2			

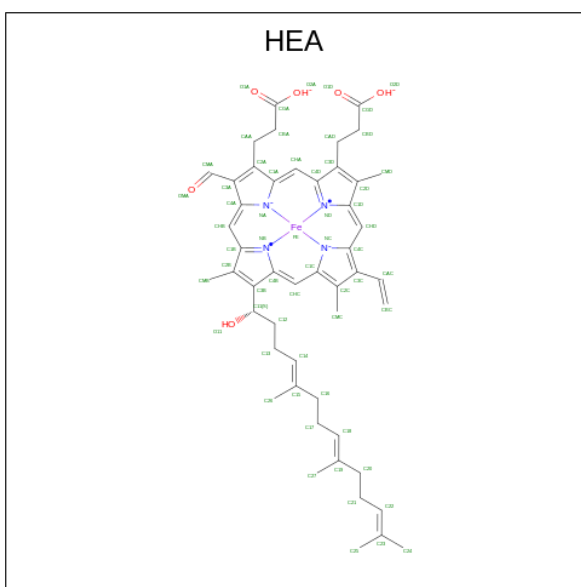
- Molecule 12 is a protein called Cytochrome c oxidase subunit 7C, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
12	L	44	Total	C	N	O	S	0	0	0
			360	242	59	57	2			

- Molecule 13 is a protein called Cytochrome c oxidase subunit 8B.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
13	M	40	Total	C	N	O	0	0	0
			311	208	48	55			

- Molecule 14 is HEME-A (CCD ID: HEA) (formula: C₄₉H₅₆FeN₄O₆).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
14	A	1	Total 60	C 49	Fe 1	N 4	O 6	0	0
14	A	1	Total 60	C 49	Fe 1	N 4	O 6	0	0

- Molecule 15 is COPPER (II) ION (CCD ID: CU) (formula: Cu).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
15	A	1	Total	Cu	0	0
			1	1		

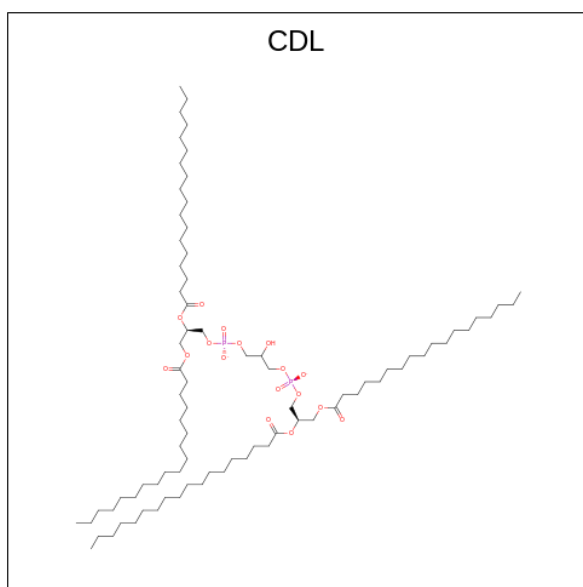
- Molecule 16 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
16	A	1	Total	Mg	0	0
			1	1		

- Molecule 17 is SODIUM ION (CCD ID: NA) (formula: Na).

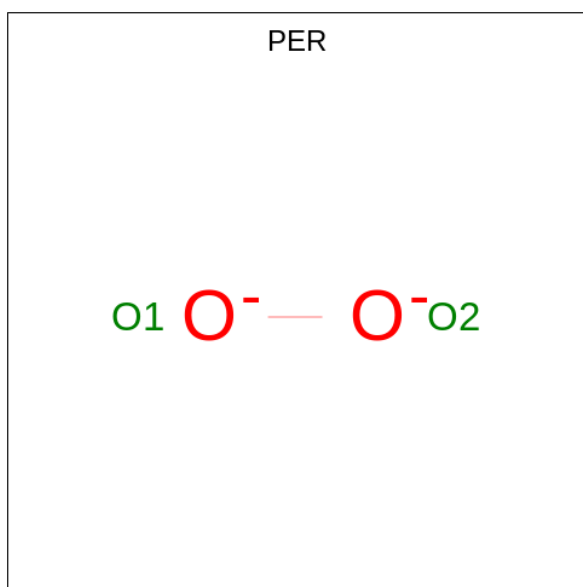
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
17	A	1	Total	Na	0	0
			1	1		
17	C	1	Total	Na	0	0
			1	1		

- Molecule 18 is CARDIOLIPIN (CCD ID: CDL) (formula: C₈₁H₁₅₆O₁₇P₂) (labeled as "Ligand of Interest" by depositor).



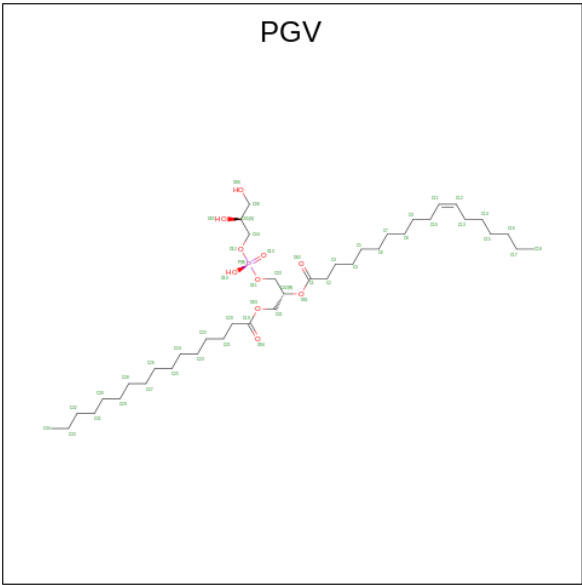
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
18	A	1	Total	C	O	P	0	0
			94	75	17	2		
18	B	1	Total	C	O	P	0	0
			64	45	17	2		
18	C	1	Total	C	O	P	0	0
			87	68	17	2		

- Molecule 19 is PEROXIDE ION (CCD ID: PER) (formula: O₂).



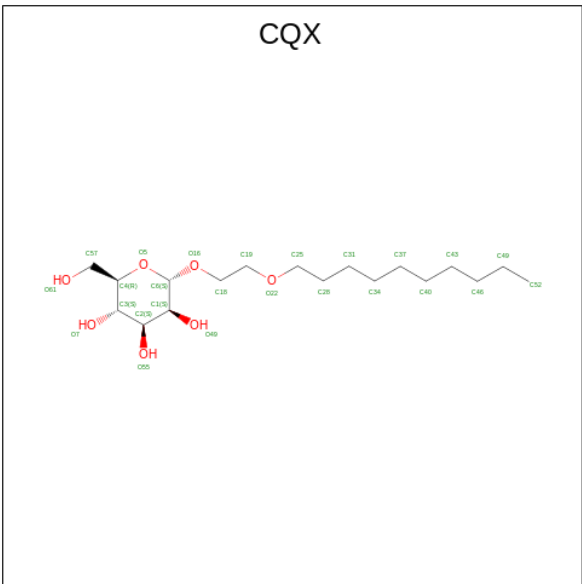
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
19	A	1	Total	O	0	0
			2	2		

- Molecule 20 is (1R)-2-{{[[[(2S)-2,3-DIHYDROXYPROPYL]OXY}(HYDROXY)PHOSPHORYL]OXY}-1-[(PALMITOYLOXY)METHYL]ETHYL (11E)-OCTADEC-11-ENOATE (CCD ID: PGV) (formula: C₄₀H₇₇O₁₀P).



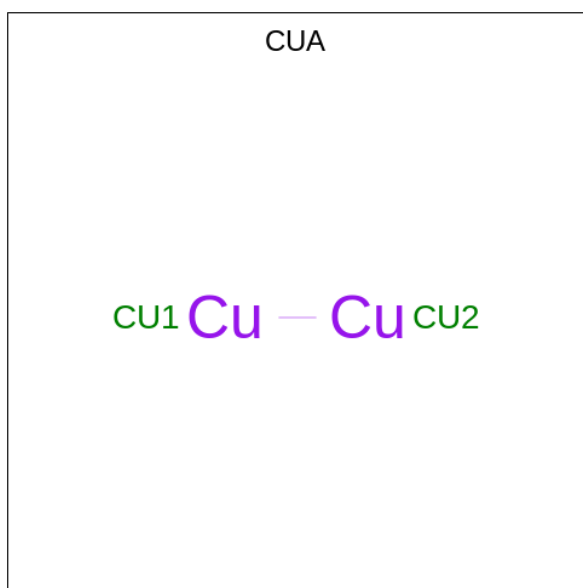
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
20	A	1	Total	C	O	P	0	0
			51	40	10	1		
20	C	1	Total	C	O	P	0	0
			41	30	10	1		

- Molecule 21 is (2S,3S,4S,5S,6R)-2-(2-decoxyethoxy)-6-(hydroxymethyl)oxane-3,4,5-triol (CCD ID: CQX) (formula: C₁₈H₃₆O₇) (labeled as "Ligand of Interest" by depositor).



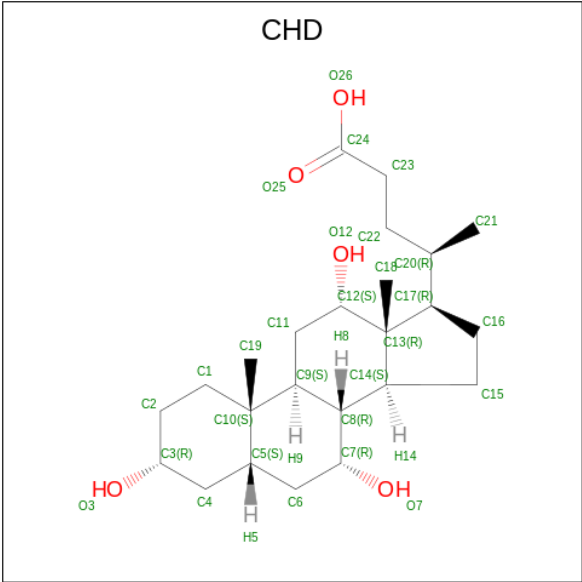
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
21	A	1	Total C O 25 18 7	0	0
21	A	1	Total C O 25 18 7	0	0
21	A	1	Total C O 25 18 7	0	0
21	B	1	Total C O 25 18 7	0	0
21	B	1	Total C O 16 9 7	0	0
21	C	1	Total C O 25 18 7	0	0
21	C	1	Total C O 25 18 7	0	0
21	G	1	Total C O 25 18 7	0	0

- Molecule 22 is DINUCLEAR COPPER ION (CCD ID: CUA) (formula: Cu₂).



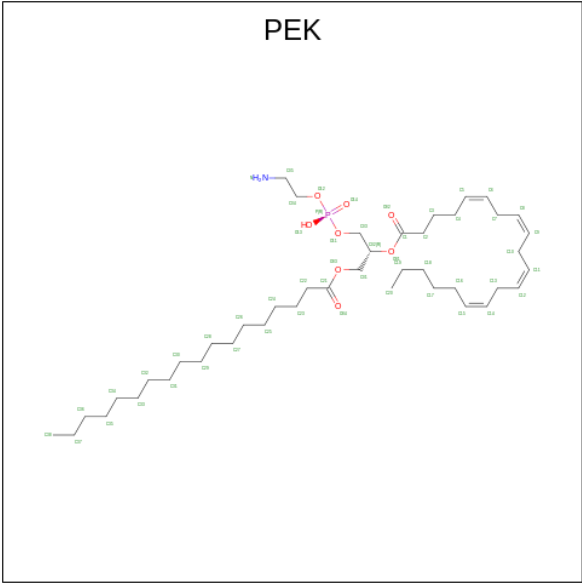
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
22	B	1	Total Cu 2 2	0	0

- Molecule 23 is CHOLIC ACID (CCD ID: CHD) (formula: C₂₄H₄₀O₅) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
23	C	1	Total	C	O	0	0
			29	24	5		

- Molecule 24 is (1S)-2-{[(2-AMINOETHOXY)(HYDROXY)PHOSPHORYL]OXY}-1-[(STEAROYLOXY)METHYL]ETHYL (5E,8E,11E,14E)-ICOSA-5,8,11,14-TETRAENOATE (CCD ID: PEK) (formula: C₄₃H₇₈NO₈P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
24	C	1	Total	C	N	O	P	0	0
			43	33	1	8	1		

- Molecule 25 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
25	F	1	Total Zn 1 1	0	0

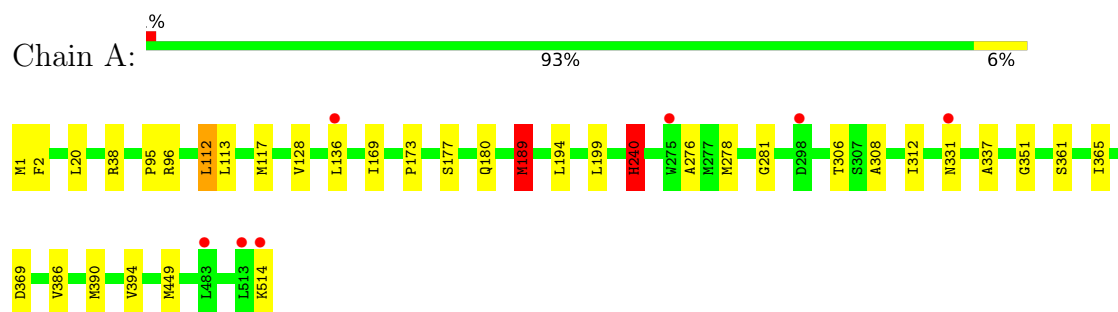
- Molecule 26 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
26	A	204	Total O 205 205	0	1
26	B	124	Total O 125 125	0	1
26	C	79	Total O 79 79	0	0
26	D	41	Total O 41 41	0	0
26	E	31	Total O 31 31	0	0
26	F	46	Total O 47 47	0	1
26	G	37	Total O 37 37	0	0
26	H	40	Total O 40 40	0	0
26	I	17	Total O 17 17	0	0
26	J	9	Total O 9 9	0	0
26	K	12	Total O 12 12	0	0
26	L	13	Total O 13 13	0	0
26	M	9	Total O 9 9	0	0

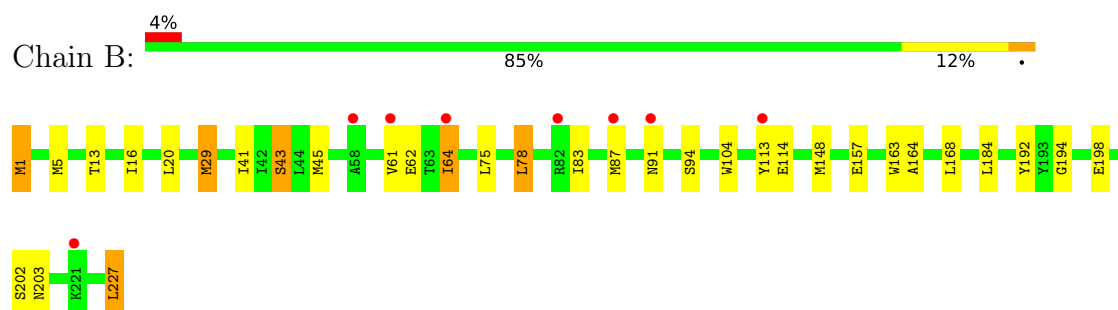
3 Residue-property plots [i](#)

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.

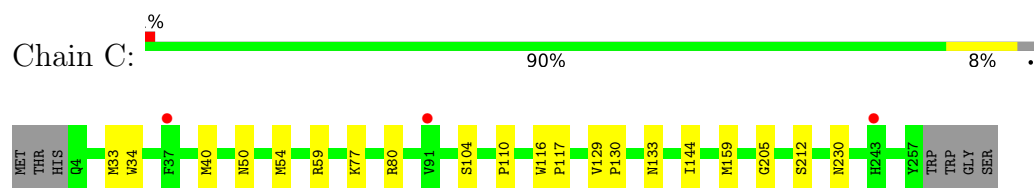
- Molecule 1: Cytochrome c oxidase subunit 1



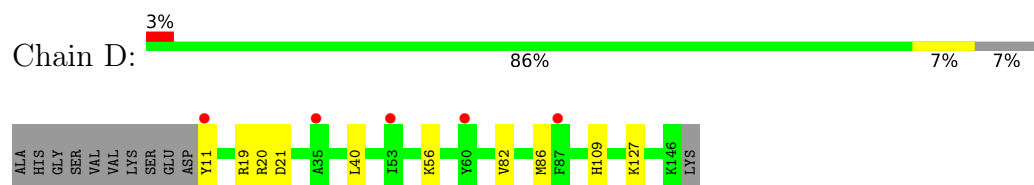
- Molecule 2: Cytochrome c oxidase subunit 2



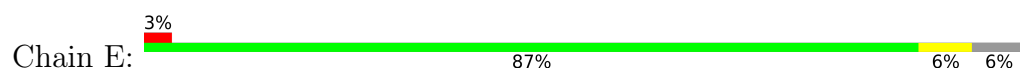
- Molecule 3: Cytochrome c oxidase subunit 3



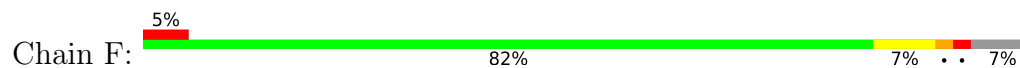
- Molecule 4: Cytochrome c oxidase subunit 4 isoform 1



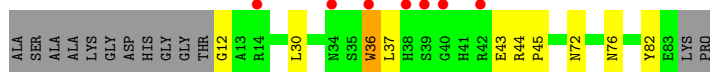
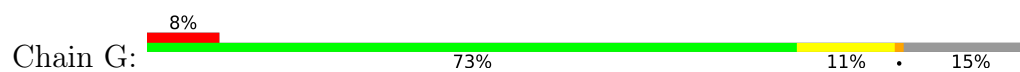
- Molecule 5: Cytochrome c oxidase subunit 5A



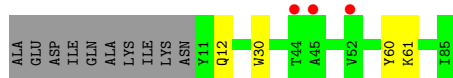
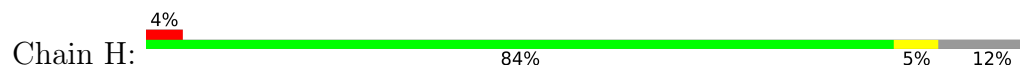
• Molecule 6: Cytochrome c oxidase subunit 5B



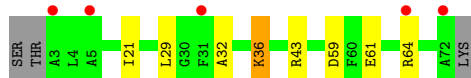
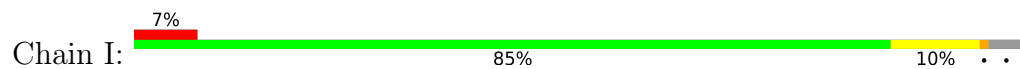
• Molecule 7: Cytochrome c oxidase subunit 6A2, mitochondrial



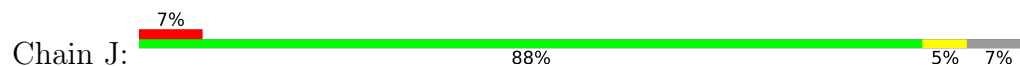
• Molecule 8: Cytochrome c oxidase subunit 6B1



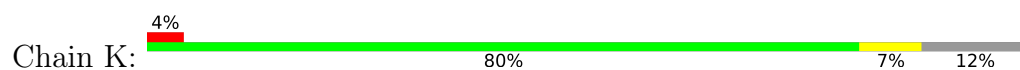
• Molecule 9: Cytochrome c oxidase subunit 6C



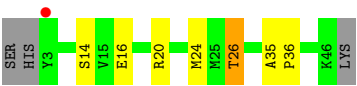
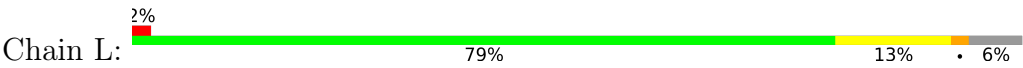
• Molecule 10: Cytochrome c oxidase subunit 7A1



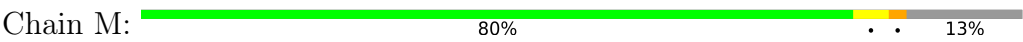
• Molecule 11: Cytochrome c oxidase subunit 7B



- Molecule 12: Cytochrome c oxidase subunit 7C, mitochondrial



- Molecule 13: Cytochrome c oxidase subunit 8B



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	149.53Å 152.13Å 174.08Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.00 – 1.85 40.00 – 1.85	Depositor EDS
% Data completeness (in resolution range)	99.0 (40.00-1.85) 99.0 (40.00-1.85)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.86 (at 1.80Å)	Xtriage
Refinement program	REFMAC 5.8.0049	Depositor
R, R_{free}	0.151 , 0.192 0.166 , 0.203	Depositor DCC
R_{free} test set	17846 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å ²)	30.9	Xtriage
Anisotropy	0.472	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 67.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.005 for k,h,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	15245	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.14% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, FME, CHD, HEA, PER, CQX, NA, MG, CDL, CUA, PEK, PGV, CU

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# $ Z > 5$	RMSZ	# $ Z > 5$
1	A	1.38	20/4156 (0.5%)	1.10	4/5678 (0.1%)
2	B	1.36	4/1860 (0.2%)	1.16	5/2534 (0.2%)
3	C	1.27	5/2143 (0.2%)	1.03	1/2931 (0.0%)
4	D	1.14	1/1167 (0.1%)	1.11	2/1577 (0.1%)
5	E	1.13	1/843 (0.1%)	1.02	0/1145
6	F	1.11	0/709	1.09	2/963 (0.2%)
7	G	1.25	1/621 (0.2%)	1.14	1/848 (0.1%)
8	H	1.20	0/648	1.07	0/877
9	I	1.16	0/588	1.06	2/781 (0.3%)
10	J	1.16	0/443	1.00	1/598 (0.2%)
11	K	1.15	0/398	1.15	2/546 (0.4%)
12	L	1.35	1/372 (0.3%)	1.16	1/500 (0.2%)
13	M	1.20	0/321	1.05	0/440
All	All	1.28	33/14269 (0.2%)	1.09	21/19418 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
6	F	0	1
All	All	0	2

The worst 5 of 33 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	278	MET	SD-CE	-13.69	1.45	1.79
7	G	36	TRP	CB-CG	10.06	1.81	1.50
3	C	104	SER	CA-CB	8.96	1.68	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	104	SER	CB-OG	8.51	1.59	1.42
1	A	189	MET	SD-CE	-7.95	1.59	1.79

The worst 5 of 21 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	20	ARG	NE-CZ-NH2	10.31	128.48	119.20
2	B	29	MET	CG-SD-CE	-8.72	81.71	100.90
1	A	189	MET	CB-CG-SD	-8.03	88.61	112.70
7	G	36	TRP	CB-CA-C	7.43	124.94	110.67
11	K	49	THR	CA-C-N	-6.82	112.76	119.78

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	240	HIS	Sidechain
6	F	65	ASP	Peptide

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	4027	0	4001	25	0
2	B	1824	0	1833	20	0
3	C	2061	0	1992	9	0
4	D	1133	0	1119	6	0
5	E	825	0	823	3	0
6	F	694	0	677	8	0
7	G	595	0	569	11	0
8	H	628	0	580	1	0
9	I	575	0	584	5	0
10	J	434	0	432	4	0
11	K	384	0	366	2	0
12	L	360	0	360	5	0
13	M	311	0	321	4	0
14	A	120	0	108	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
15	A	1	0	0	0	0
16	A	1	0	0	0	0
17	A	1	0	0	0	0
17	C	1	0	0	0	0
18	A	94	0	141	10	0
18	B	64	0	72	0	0
18	C	87	0	124	9	0
19	A	2	0	0	1	0
20	A	51	0	76	1	0
20	C	41	0	55	0	0
21	A	75	0	0	2	0
21	B	41	0	0	0	0
21	C	50	0	0	1	0
21	G	25	0	0	0	0
22	B	2	0	0	0	0
23	C	29	0	39	1	0
24	C	43	0	58	2	0
25	F	1	0	0	0	0
26	A	205	0	0	5	0
26	B	125	0	0	2	0
26	C	79	0	0	0	0
26	D	41	0	0	1	0
26	E	31	0	0	0	0
26	F	47	0	0	0	0
26	G	37	0	0	3	0
26	H	40	0	0	0	0
26	I	17	0	0	1	0
26	J	9	0	0	1	0
26	K	12	0	0	0	0
26	L	13	0	0	0	0
26	M	9	0	0	2	0
All	All	15245	0	14330	106	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 4.

The worst 5 of 106 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)
7:G:36:TRP:CB	7:G:36:TRP:CG	1.81	1.59
19:A:607:PER:O1	19:A:607:PER:O2	1.55	1.22

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:16:ILE:HG23	2:B:87:MET:HE2	1.36	1.04
1:A:112:LEU:HG	26:A:892:HOH:O	1.58	1.02
18:C:305:CDL:O1	10:J:8:LYS:HD2	1.65	0.95

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	512/514 (100%)	498 (97%)	14 (3%)	0	100	100
2	B	225/227 (99%)	220 (98%)	5 (2%)	0	100	100
3	C	252/261 (97%)	249 (99%)	3 (1%)	0	100	100
4	D	134/147 (91%)	128 (96%)	6 (4%)	0	100	100
5	E	100/109 (92%)	99 (99%)	1 (1%)	0	100	100
6	F	89/98 (91%)	87 (98%)	2 (2%)	0	100	100
7	G	70/85 (82%)	66 (94%)	4 (6%)	0	100	100
8	H	73/85 (86%)	72 (99%)	1 (1%)	0	100	100
9	I	68/73 (93%)	68 (100%)	0	0	100	100
10	J	53/59 (90%)	53 (100%)	0	0	100	100
11	K	47/56 (84%)	46 (98%)	1 (2%)	0	100	100
12	L	42/47 (89%)	41 (98%)	1 (2%)	0	100	100
13	M	38/46 (83%)	38 (100%)	0	0	100	100
All	All	1703/1807 (94%)	1665 (98%)	38 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	426/426 (100%)	421 (99%)	5 (1%)	63	55
2	B	210/210 (100%)	203 (97%)	7 (3%)	33	18
3	C	220/226 (97%)	217 (99%)	3 (1%)	59	49
4	D	120/129 (93%)	118 (98%)	2 (2%)	53	42
5	E	89/95 (94%)	87 (98%)	2 (2%)	45	32
6	F	76/81 (94%)	72 (95%)	4 (5%)	20	7
7	G	62/69 (90%)	60 (97%)	2 (3%)	34	19
8	H	67/75 (89%)	64 (96%)	3 (4%)	24	10
9	I	55/58 (95%)	53 (96%)	2 (4%)	31	16
10	J	46/50 (92%)	46 (100%)	0	100	100
11	K	39/46 (85%)	39 (100%)	0	100	100
12	L	37/40 (92%)	36 (97%)	1 (3%)	39	24
13	M	34/38 (90%)	33 (97%)	1 (3%)	37	22
All	All	1481/1543 (96%)	1449 (98%)	32 (2%)	45	32

5 of 32 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
9	I	29	LEU
9	I	36	LYS
3	C	144	ILE
2	B	227	LEU
12	L	16	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 23 such sidechains are listed below:

Mol	Chain	Res	Type
6	F	80	GLN
8	H	22	ASN

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Mol	Chain	Res	Type
8	H	12	GLN
8	H	25	GLN
2	B	181	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	FME	A	1	1	8,9,10	0.95	0	7,9,11	1.32	1 (14%)
2	FME	B	1	2	8,9,10	2.18	4 (50%)	7,9,11	6.04	4 (57%)

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
1	FME	A	1	1	-	5/7/9/11	-
2	FME	B	1	2	-	1/7/9/11	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1	FME	O1-CN	-3.54	1.12	1.22
2	B	1	FME	CN-N	3.52	1.45	1.33
2	B	1	FME	CG-SD	-2.19	1.69	1.81
2	B	1	FME	CA-N	2.14	1.49	1.46

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1	FME	CA-N-CN	-12.03	104.31	122.82
2	B	1	FME	C-CA-N	8.75	125.52	109.73
2	B	1	FME	CG-CB-CA	-4.70	99.89	112.95
1	A	1	FME	CA-N-CN	-2.38	119.16	122.82
2	B	1	FME	O1-CN-N	-2.29	119.23	125.27

There are no chirality outliers.

5 of 6 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	1	FME	N-CA-CB-CG
1	A	1	FME	CA-CB-CG-SD
2	B	1	FME	O1-CN-N-CA
1	A	1	FME	CB-CG-SD-CE
1	A	1	FME	C-CA-CB-CG

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	1	FME	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 24 ligands modelled in this entry, 5 are monoatomic - leaving 19 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
20	PGV	A	608	-	50,50,50	0.92	3 (6%)	53,56,56	1.15	3 (5%)
21	CQX	C	306	-	25,25,25	1.22	2 (8%)	30,30,30	2.04	5 (16%)
18	CDL	B	301	-	63,63,99	1.53	5 (7%)	69,75,111	1.57	14 (20%)
21	CQX	G	101	-	25,25,25	0.78	1 (4%)	30,30,30	0.89	0
20	PGV	C	304	-	40,40,50	1.05	3 (7%)	43,46,56	1.35	4 (9%)
21	CQX	B	304	-	16,16,25	0.69	0	21,21,30	1.34	3 (14%)
18	CDL	C	305	-	86,86,99	1.19	6 (6%)	92,98,111	1.38	13 (14%)
21	CQX	A	611	-	25,25,25	0.87	2 (8%)	30,30,30	1.65	7 (23%)
21	CQX	B	303	-	25,25,25	0.78	1 (4%)	30,30,30	1.99	11 (36%)
19	PER	A	607	14,15	1,1,1	1.26	0	-	-	-
23	CHD	C	301	-	32,32,32	0.95	2 (6%)	51,51,51	1.44	9 (17%)
24	PEK	C	303	-	42,42,52	0.80	2 (4%)	45,47,57	0.96	2 (4%)
14	HEA	A	601	1	66,67,67	1.97	17 (25%)	78,103,103	2.02	23 (29%)
21	CQX	A	609	-	25,25,25	0.88	1 (4%)	30,30,30	2.24	11 (36%)
14	HEA	A	602	19,1	66,67,67	1.78	16 (24%)	78,103,103	2.20	20 (25%)
21	CQX	A	610	-	25,25,25	1.06	2 (8%)	30,30,30	1.74	6 (20%)
22	CUA	B	302	2	0,1,1	-	-	-	-	-
21	CQX	C	307	-	25,25,25	0.57	0	30,30,30	0.82	1 (3%)
18	CDL	A	606	-	93,93,99	0.98	4 (4%)	99,105,111	1.55	15 (15%)

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
21	CQX	A	611	-	-	4/16/36/36	0/1/1/1
21	CQX	C	307	-	-	7/16/36/36	0/1/1/1
21	CQX	B	303	-	-	7/16/36/36	0/1/1/1
14	HEA	A	601	1	-	7/36/76/76	-
20	PGV	A	608	-	-	7/55/55/55	-
21	CQX	C	306	-	-	7/16/36/36	0/1/1/1
18	CDL	A	606	-	-	50/104/104/110	-
21	CQX	A	609	-	-	6/16/36/36	0/1/1/1
24	PEK	C	303	-	-	8/46/46/56	-
14	HEA	A	602	19,1	-	5/36/76/76	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
18	CDL	B	301	-	-	40/74/74/110	-
21	CQX	A	610	-	-	2/16/36/36	0/1/1/1
20	PGV	C	304	-	-	9/45/45/55	-
21	CQX	G	101	-	-	1/16/36/36	0/1/1/1
23	CHD	C	301	-	-	2/9/74/74	0/4/4/4
21	CQX	B	304	-	-	0/7/27/36	0/1/1/1
18	CDL	C	305	-	-	51/97/97/110	-

The worst 5 of 67 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
18	B	301	CDL	OA6-CA5	6.94	1.53	1.34
14	A	601	HEA	C4B-NB	-6.36	1.29	1.40
18	B	301	CDL	OB6-CB5	4.75	1.47	1.34
14	A	602	HEA	C4B-NB	-4.71	1.32	1.40
18	C	305	CDL	OB6-CB5	4.68	1.47	1.34

The worst 5 of 147 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
21	C	306	CQX	C18-O16-C6	6.78	125.09	113.84
14	A	602	HEA	C4A-NA-C1A	-6.61	98.88	105.35
14	A	602	HEA	C4C-NC-C1C	-6.21	99.27	105.35
14	A	602	HEA	CHB-C4A-NA	-6.11	117.83	124.44
14	A	601	HEA	C3C-C4C-NC	6.02	113.95	110.25

There are no chirality outliers.

5 of 213 torsion outliers are listed below:

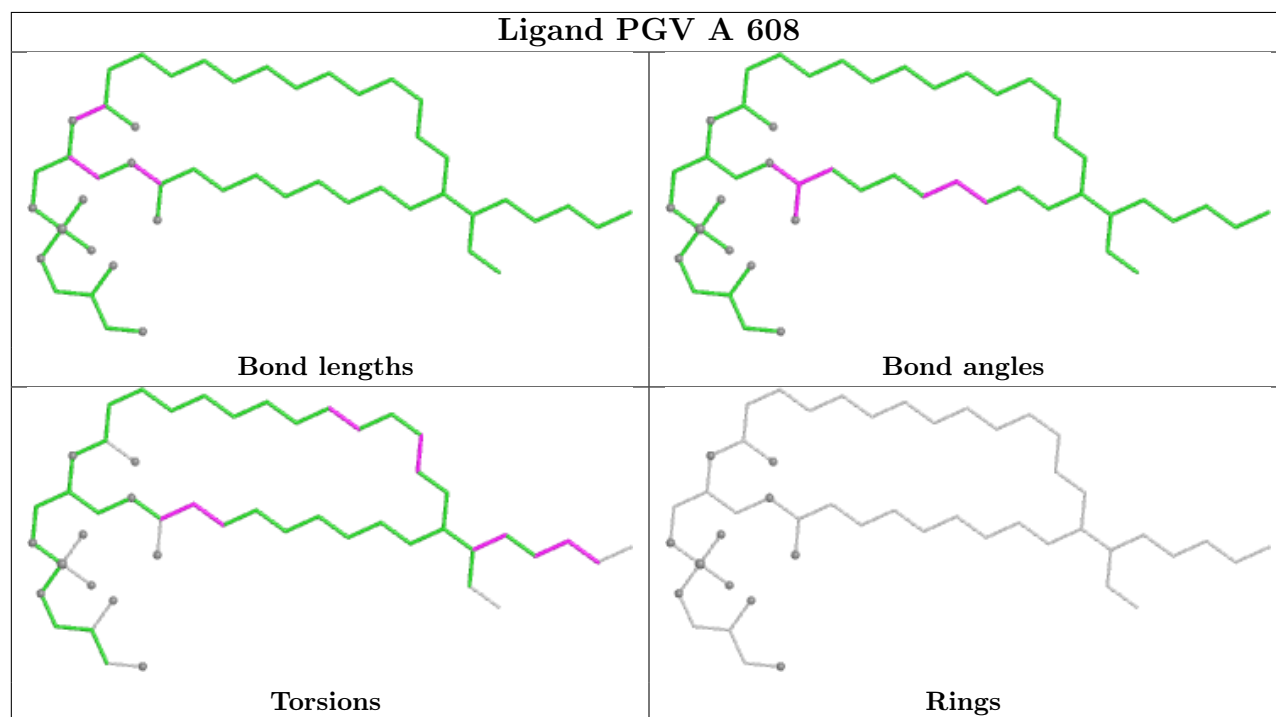
Mol	Chain	Res	Type	Atoms
18	A	606	CDL	CA2-C1-CB2-OB2
18	A	606	CDL	CA2-OA2-PA1-OA3
18	A	606	CDL	CA3-OA5-PA1-OA3
18	A	606	CDL	C11-CA5-OA6-CA4
18	A	606	CDL	CB2-OB2-PB2-OB5

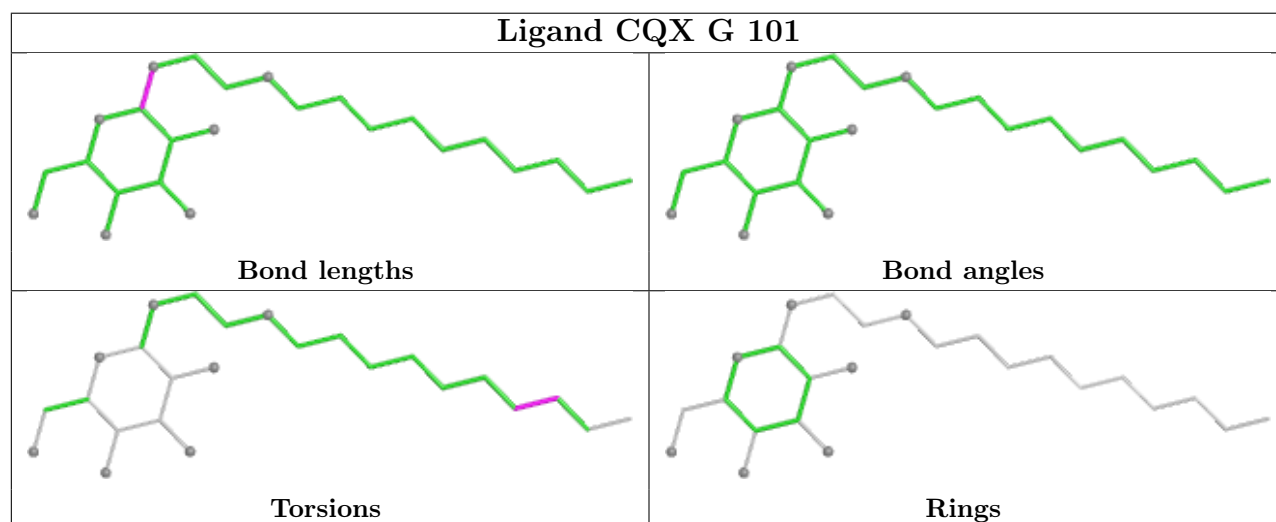
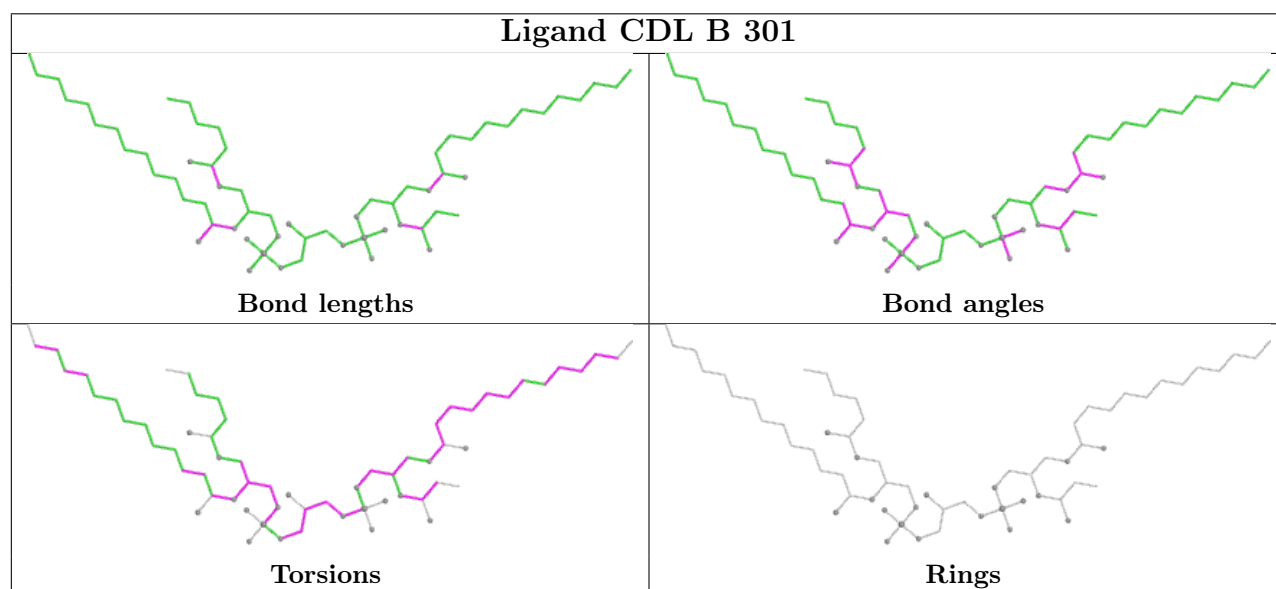
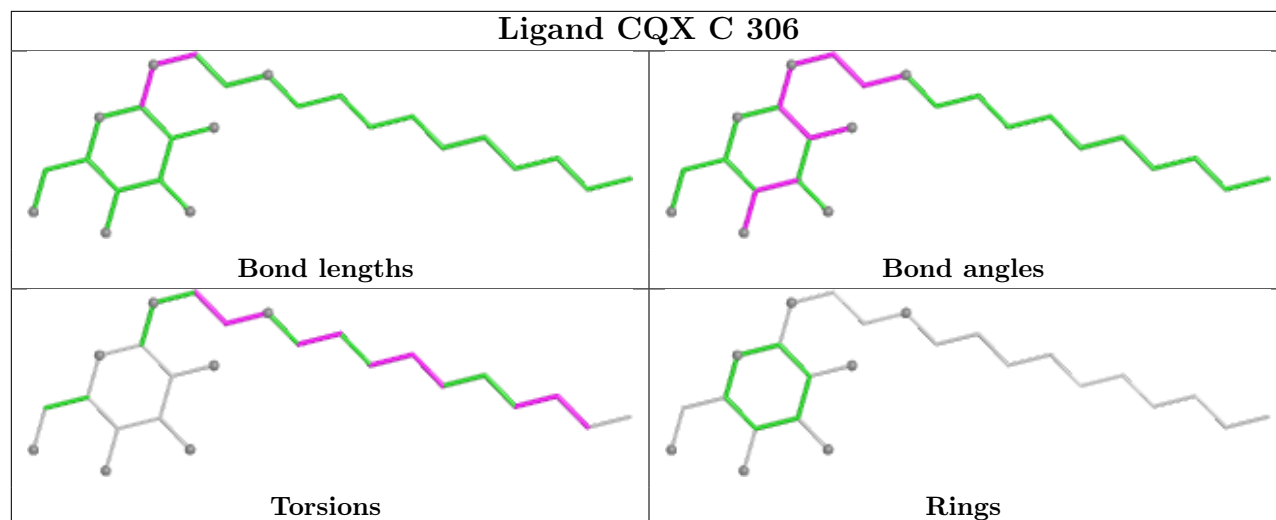
There are no ring outliers.

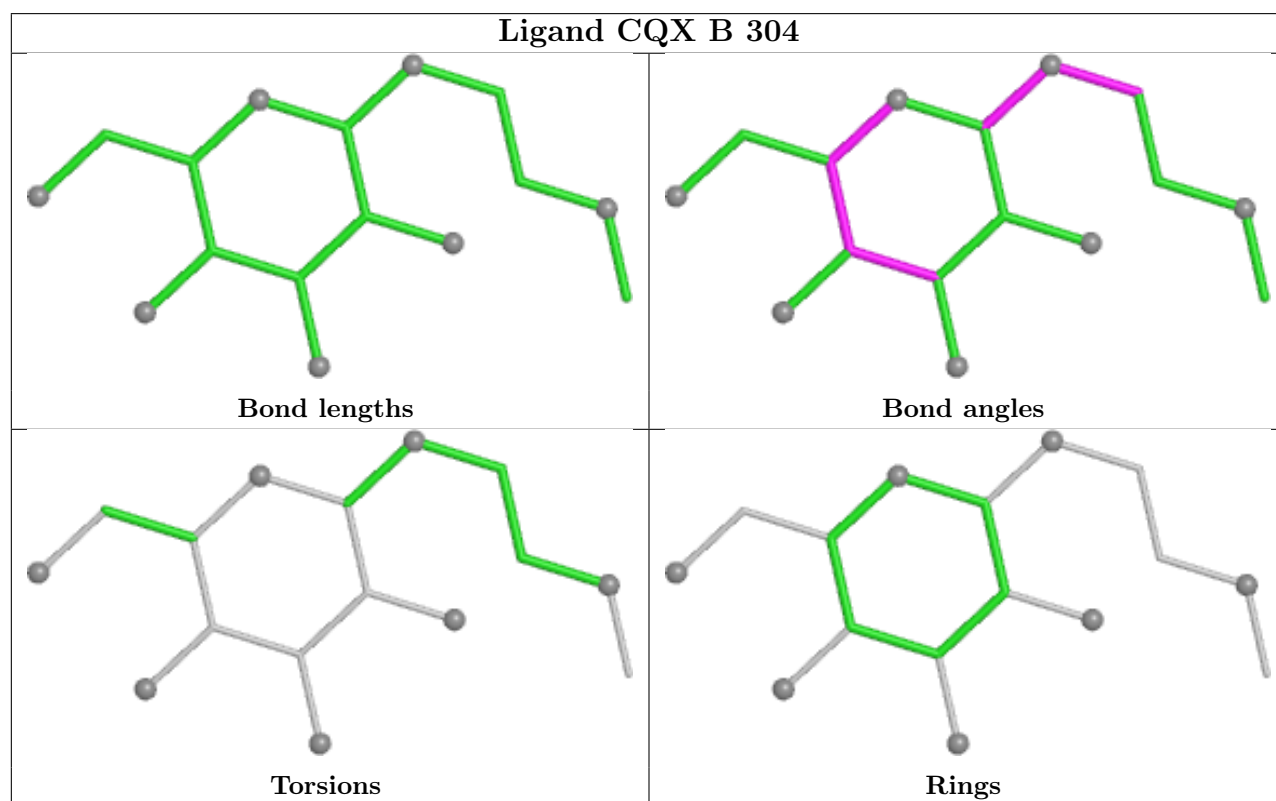
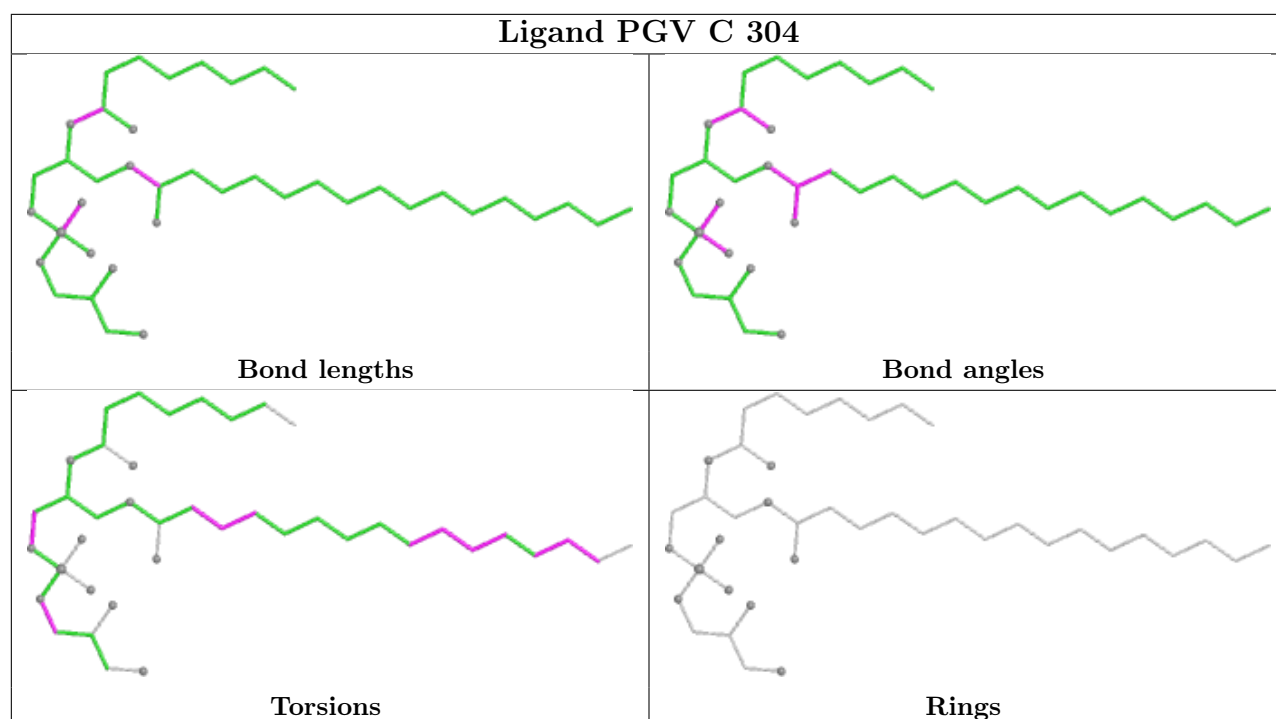
10 monomers are involved in 30 short contacts:

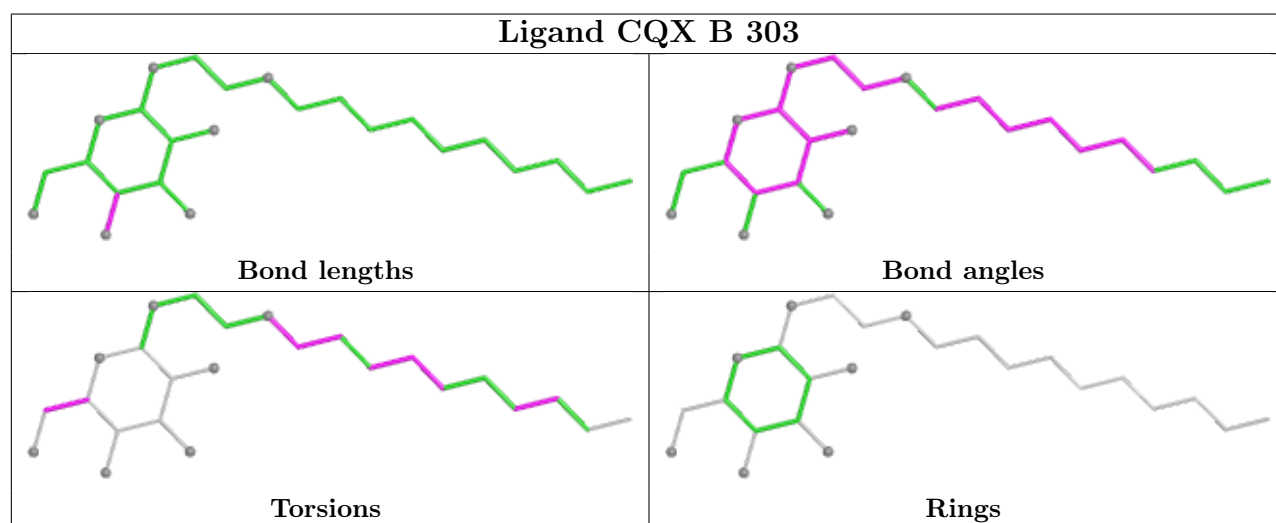
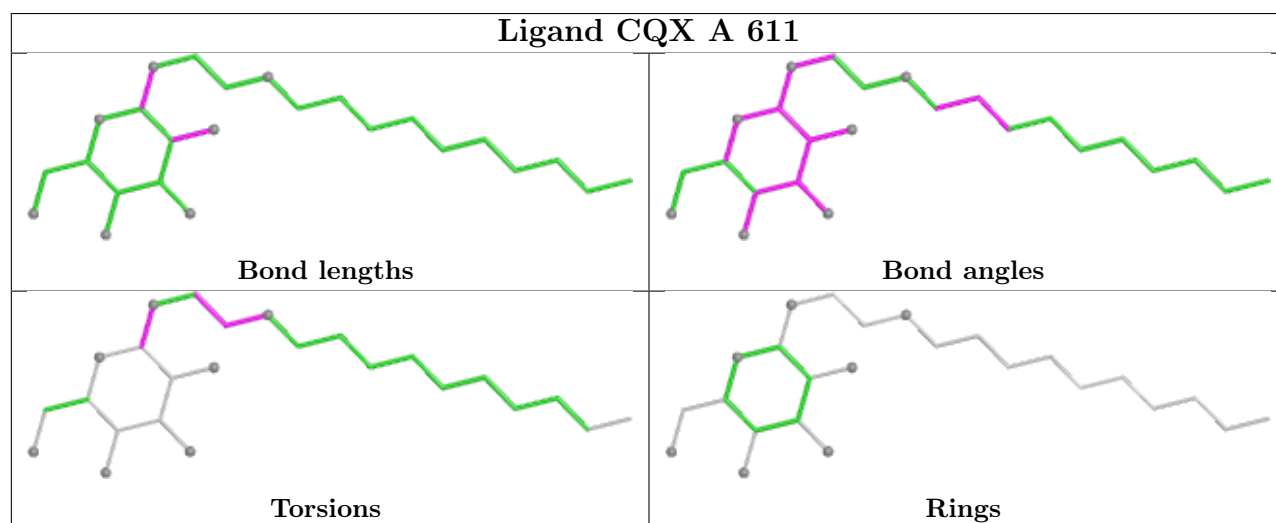
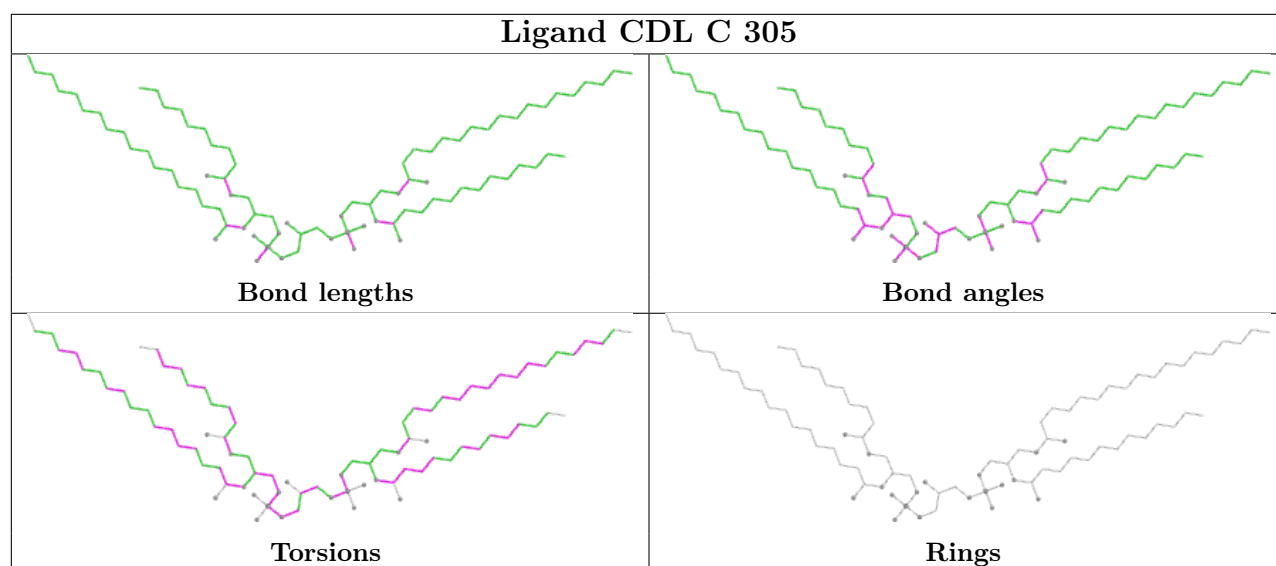
Mol	Chain	Res	Type	Clashes	Symm-Clashes
20	A	608	PGV	1	0
21	C	306	CQX	1	0
18	C	305	CDL	9	0
19	A	607	PER	1	0
23	C	301	CHD	1	0
24	C	303	PEK	2	0
14	A	601	HEA	3	0
14	A	602	HEA	1	0
21	A	610	CQX	2	0
18	A	606	CDL	10	0

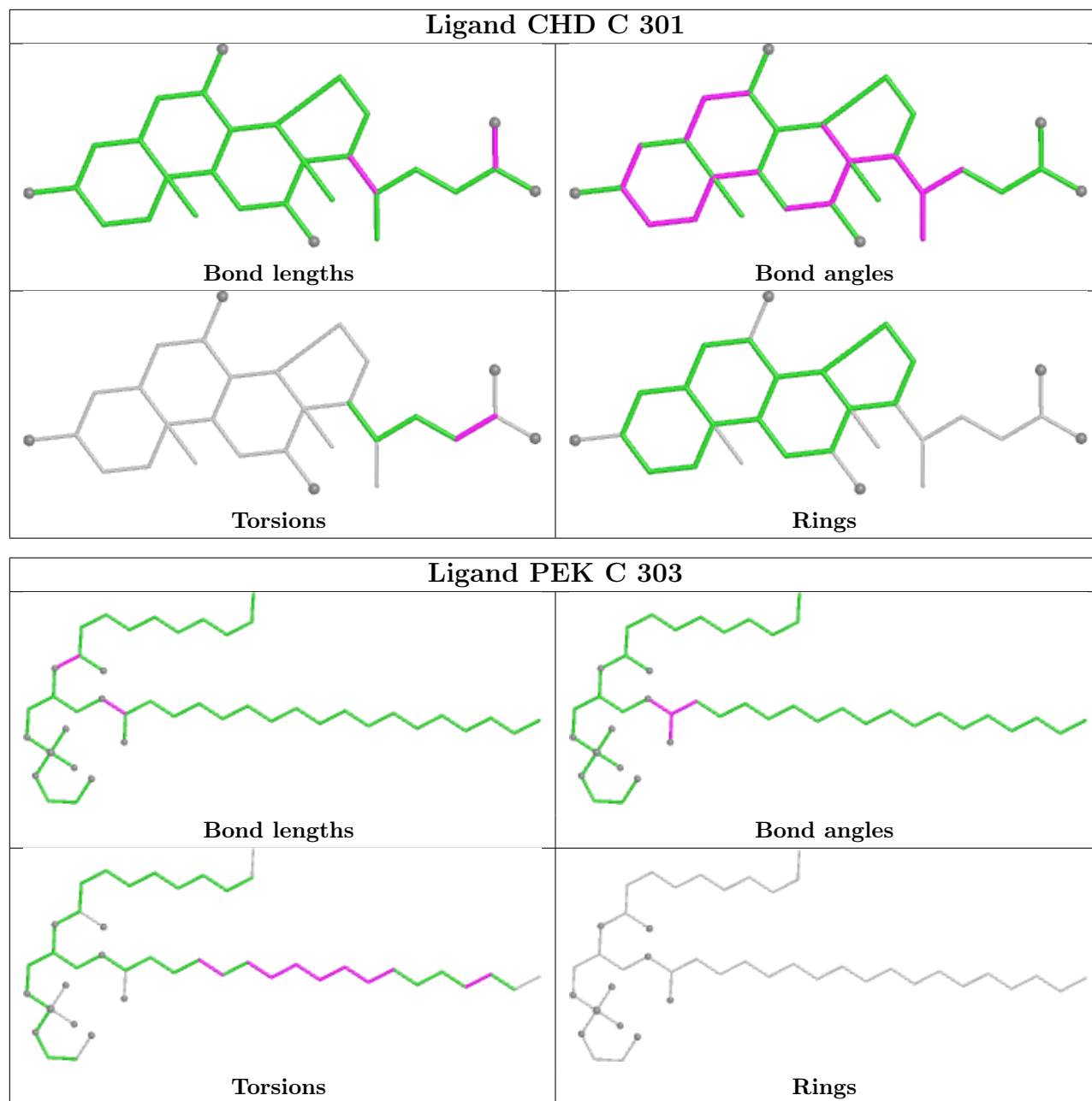
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

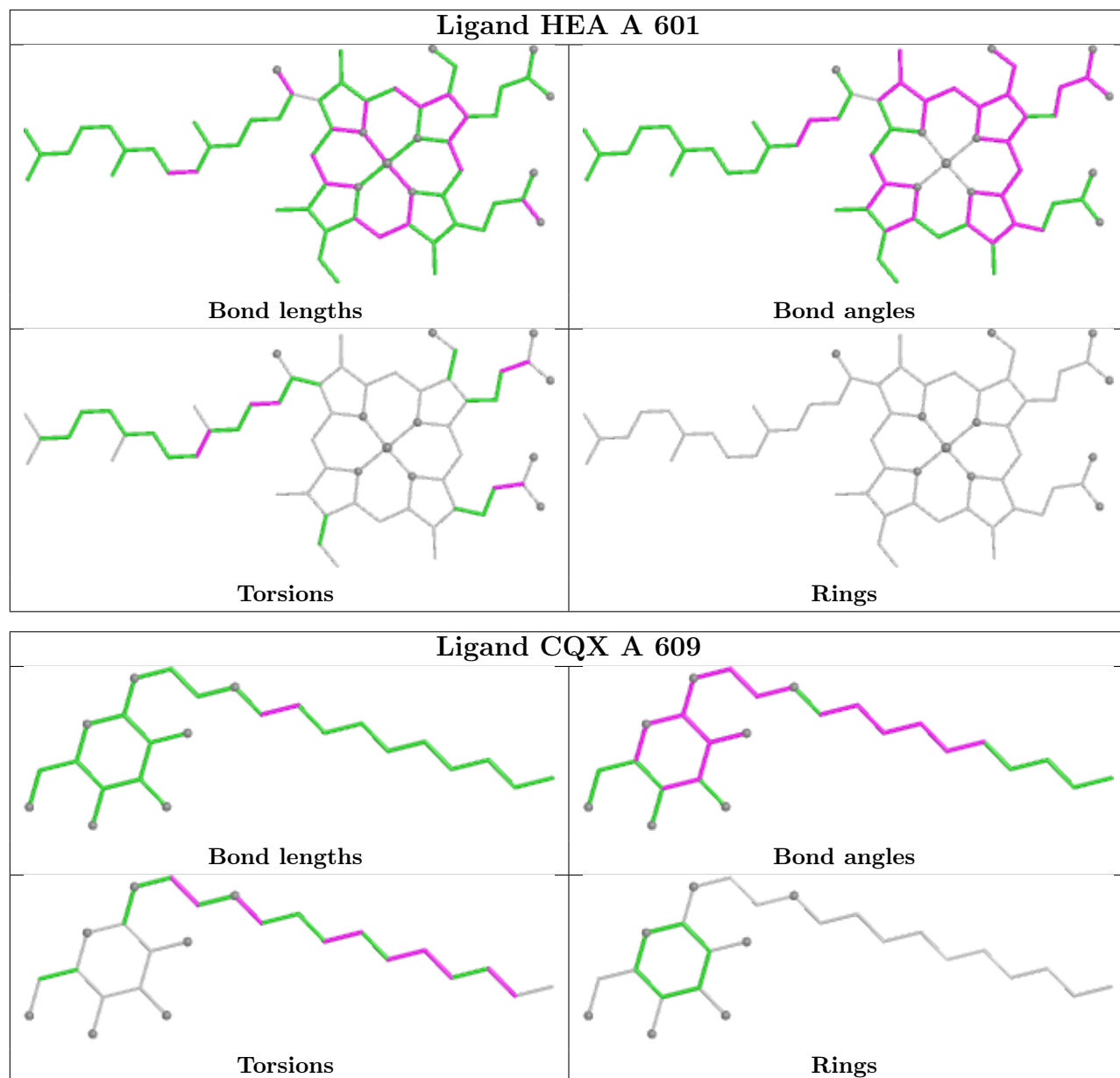


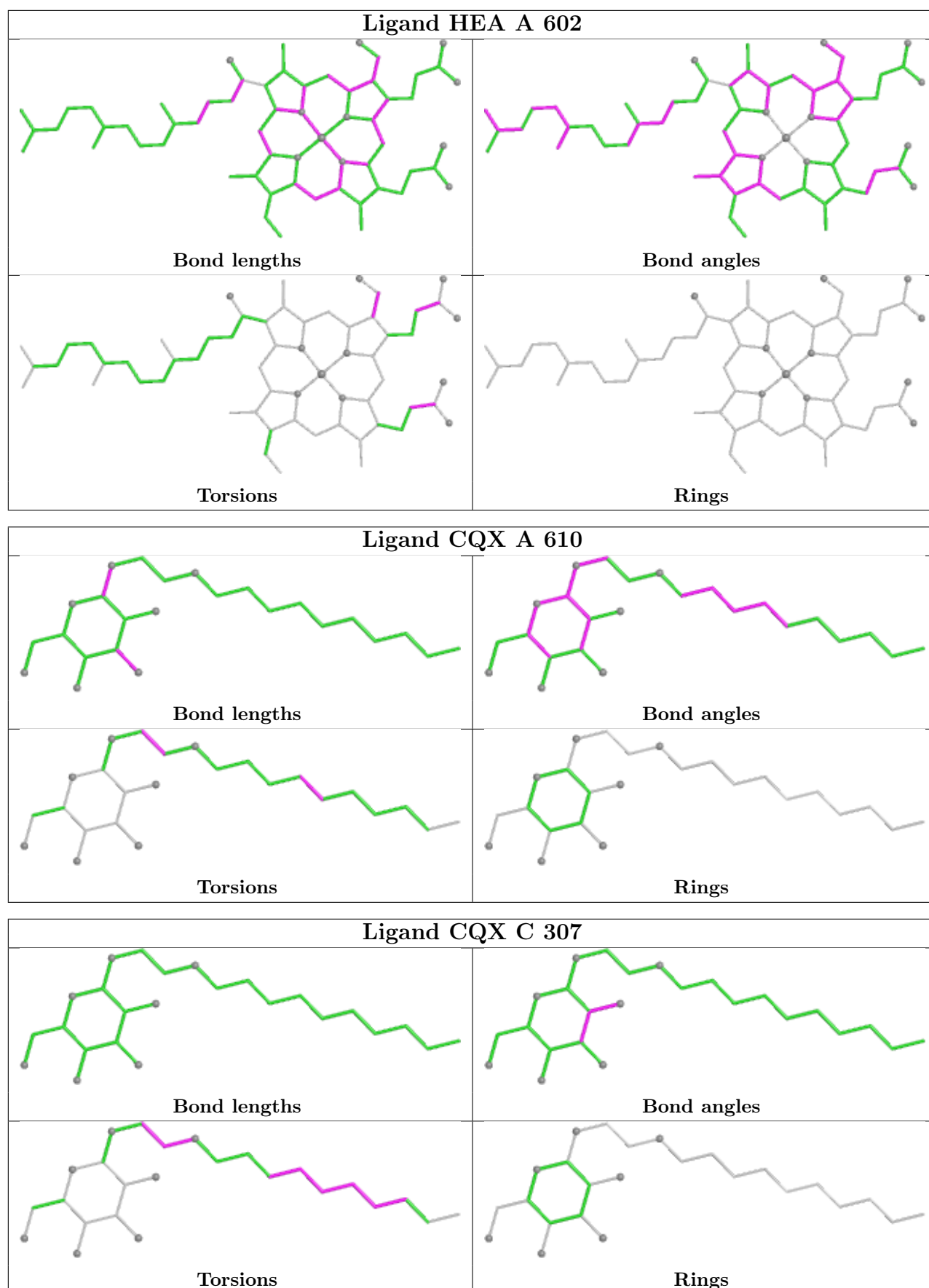


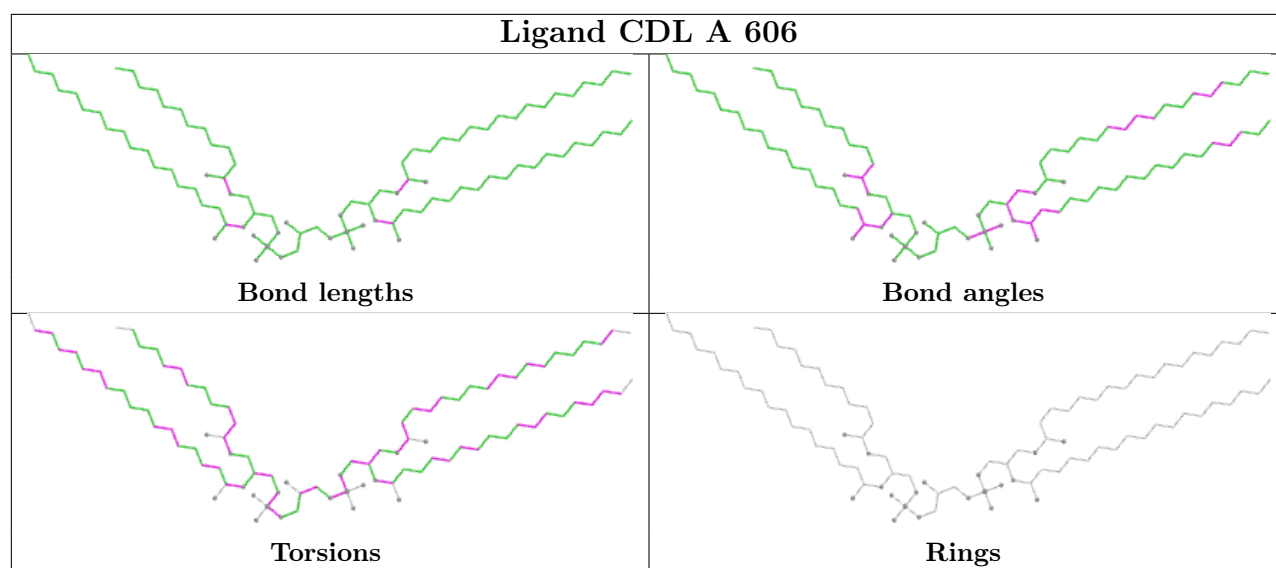












5.7 Other polymers [i](#)

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	513/514 (99%)	-0.25	7 (1%) 73 77	25, 32, 42, 96	0
2	B	226/227 (99%)	0.14	8 (3%) 47 51	29, 38, 63, 95	0
3	C	254/261 (97%)	0.03	3 (1%) 76 80	28, 37, 51, 80	0
4	D	136/147 (92%)	0.48	5 (3%) 45 49	37, 50, 77, 84	0
5	E	102/109 (93%)	0.42	3 (2%) 53 57	39, 52, 71, 86	0
6	F	91/98 (92%)	0.37	5 (5%) 30 33	34, 49, 72, 93	0
7	G	72/85 (84%)	0.64	7 (9%) 13 13	35, 45, 89, 102	0
8	H	75/85 (88%)	0.31	3 (4%) 42 46	33, 43, 76, 84	0
9	I	70/73 (95%)	0.38	5 (7%) 22 24	36, 46, 68, 81	0
10	J	55/59 (93%)	0.63	4 (7%) 21 23	38, 48, 77, 86	0
11	K	49/56 (87%)	0.66	2 (4%) 41 45	42, 52, 69, 83	0
12	L	44/47 (93%)	0.11	1 (2%) 61 64	34, 41, 58, 67	0
13	M	40/46 (86%)	0.51	0 100 100	39, 47, 62, 71	0
All	All	1727/1807 (95%)	0.14	53 (3%) 51 55	25, 39, 68, 102	0

The worst 5 of 53 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
7	G	36	TRP	8.8
2	B	113	TYR	6.3
11	K	6	ALA	4.5
4	D	11	TYR	4.2
1	A	136	LEU	3.6

6.2 Non-standard residues in protein, DNA, RNA chains [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
1	FME	A	1	10/11	0.90	0.21	47,60,99,117	0
2	FME	B	1	10/11	0.97	0.07	28,42,55,64	0

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.

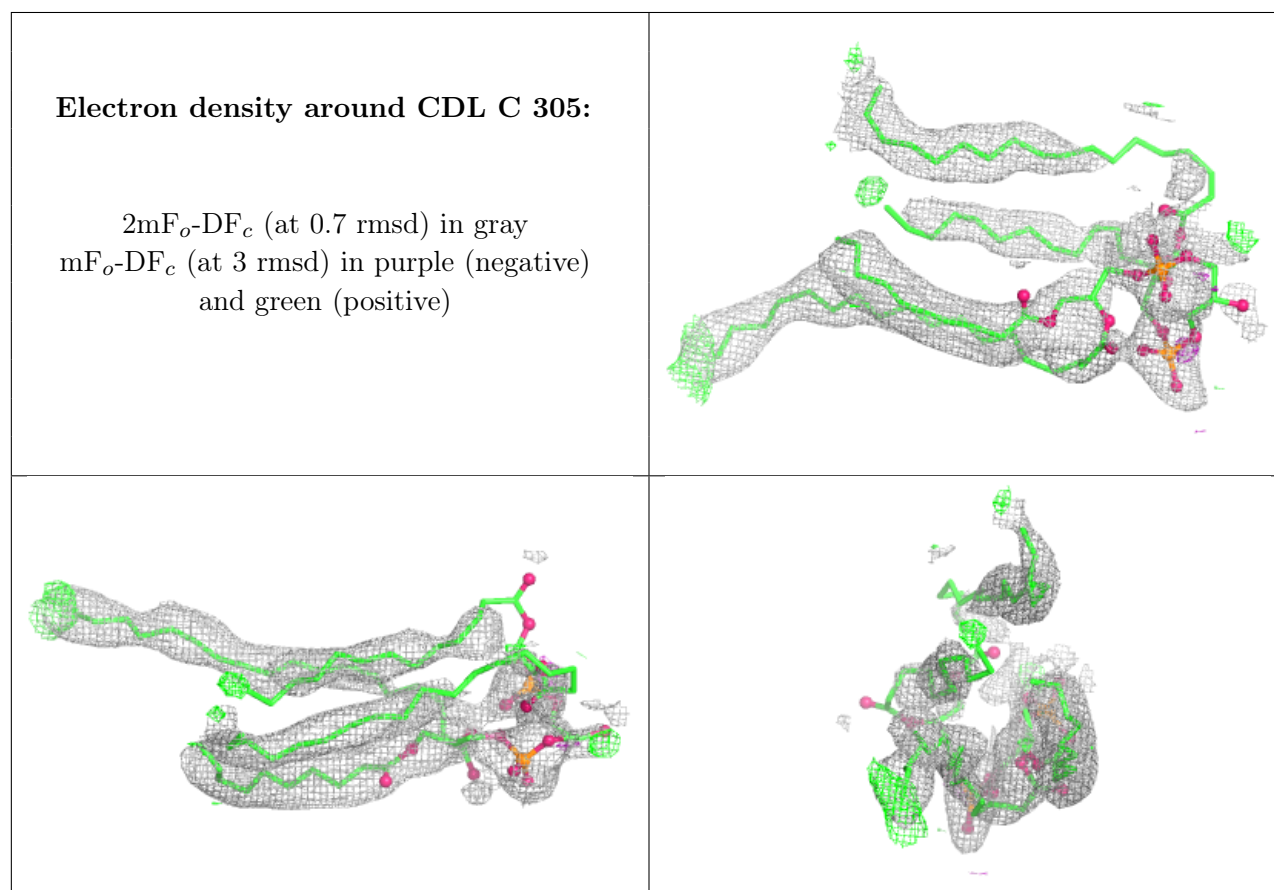
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
17	NA	C	302	1/1	0.75	0.30	64,64,64,64	0
18	CDL	C	305	87/100	0.89	0.18	44,87,139,152	0
21	CQX	A	611	25/25	0.89	0.17	49,63,92,102	0
18	CDL	A	606	94/100	0.90	0.17	46,79,133,139	0
18	CDL	B	301	64/100	0.92	0.15	52,88,124,138	0
21	CQX	B	304	16/25	0.92	0.12	46,54,81,83	0
21	CQX	B	303	25/25	0.94	0.11	40,53,73,75	0
21	CQX	A	609	25/25	0.94	0.11	40,60,74,99	0
21	CQX	C	307	25/25	0.94	0.14	40,57,89,98	0
21	CQX	C	306	25/25	0.95	0.12	38,49,93,95	0
21	CQX	A	610	25/25	0.96	0.09	47,57,72,80	0
21	CQX	G	101	25/25	0.96	0.08	43,49,59,61	0
23	CHD	C	301	29/29	0.96	0.08	29,34,39,48	0
19	PER	A	607	2/2	0.97	0.09	25,25,25,33	0
20	PGV	A	608	51/51	0.97	0.09	28,38,78,102	0
20	PGV	C	304	41/51	0.97	0.09	32,39,68,84	0
24	PEK	C	303	43/53	0.97	0.09	31,42,69,97	0
16	MG	A	604	1/1	0.99	0.03	31,31,31,31	0
14	HEA	A	601	60/60	0.99	0.06	24,29,49,60	0
14	HEA	A	602	60/60	0.99	0.05	21,26,34,45	0

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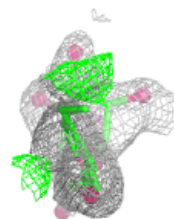
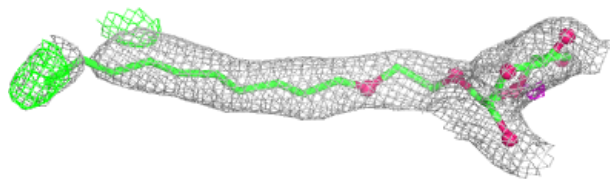
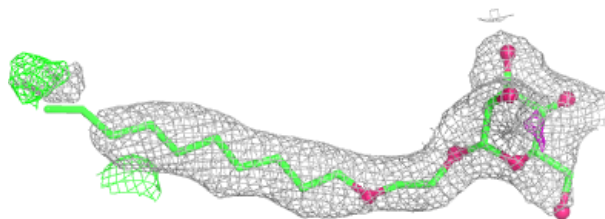
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
22	CUA	B	302	2/2	1.00	0.01	30,30,30,31	0
17	NA	A	605	1/1	1.00	0.07	35,35,35,35	0
15	CU	A	603	1/1	1.00	0.01	26,26,26,26	0
25	ZN	F	101	1/1	1.00	0.04	43,43,43,43	0

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.



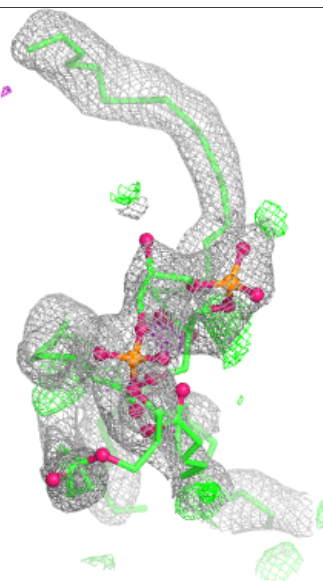
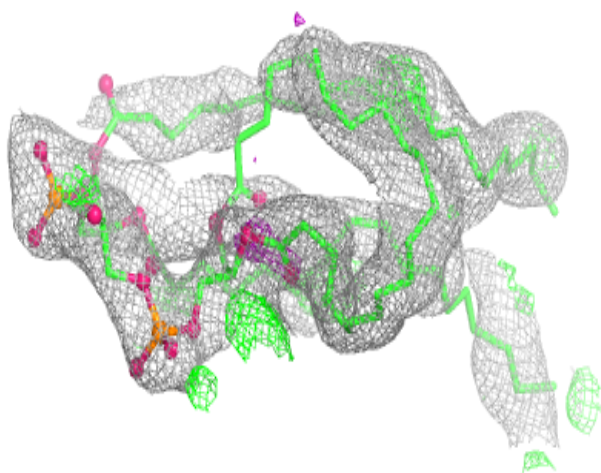
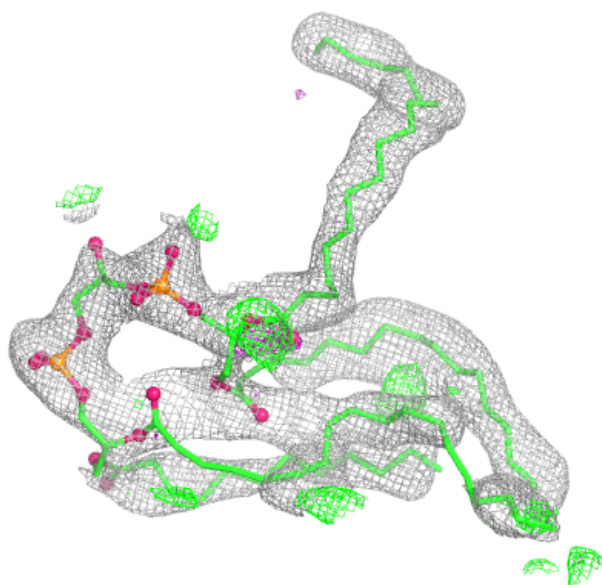
Electron density around CQX A 611:

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



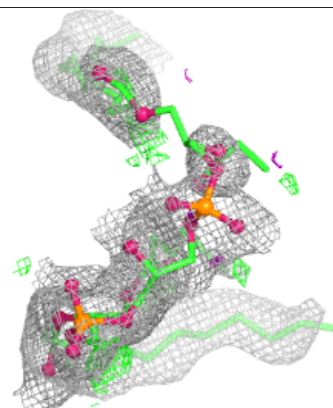
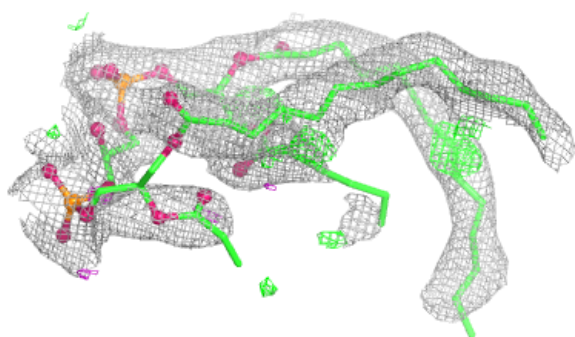
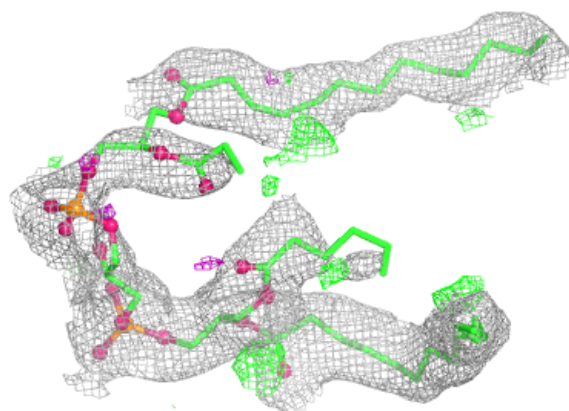
Electron density around CDL A 606:

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

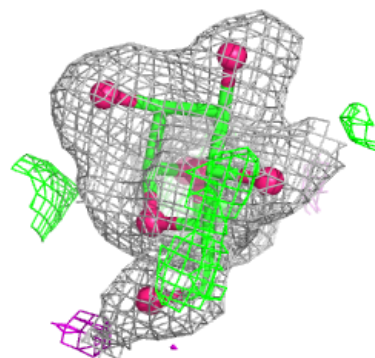
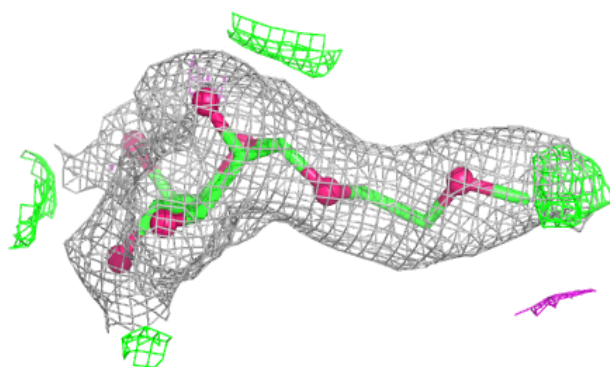
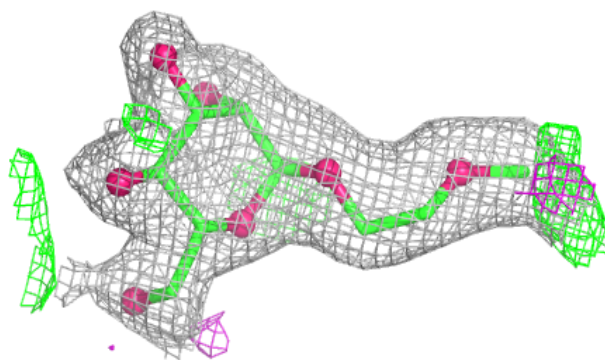


Electron density around CDL B 301:

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and green (positive)

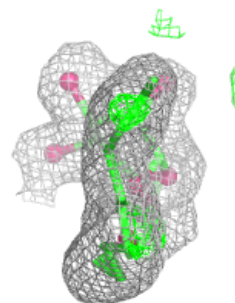
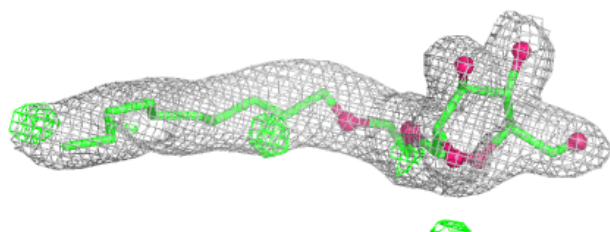
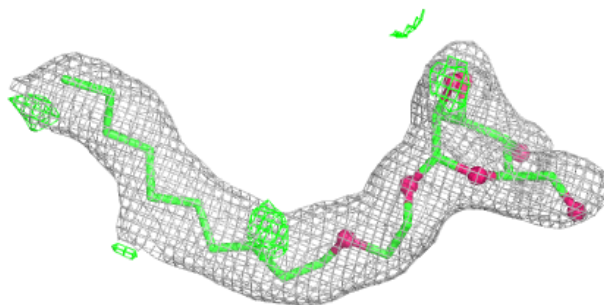
**Electron density around CQX B 304:**

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

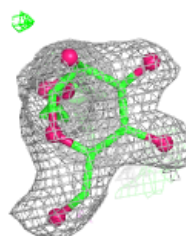
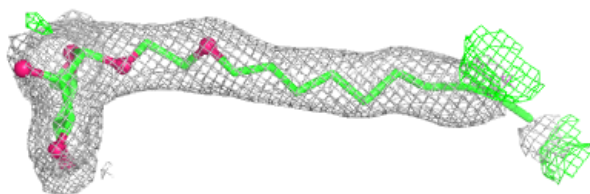
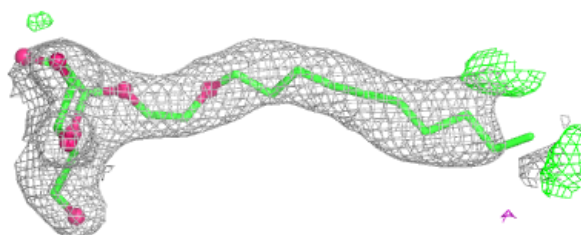


Electron density around CQX B 303:

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

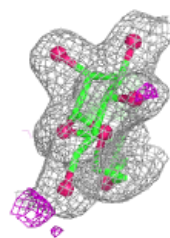
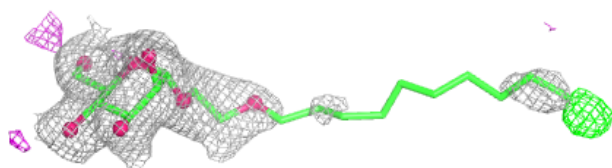
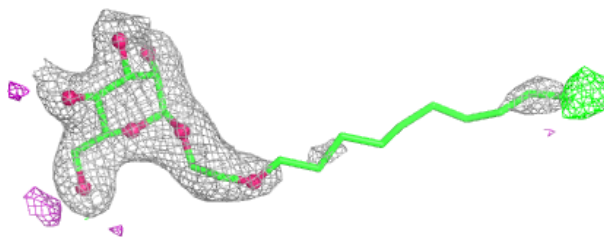
**Electron density around CQX A 609:**

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and green (positive)

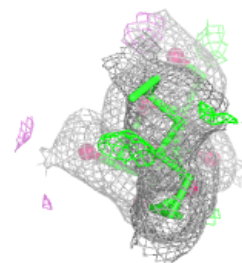
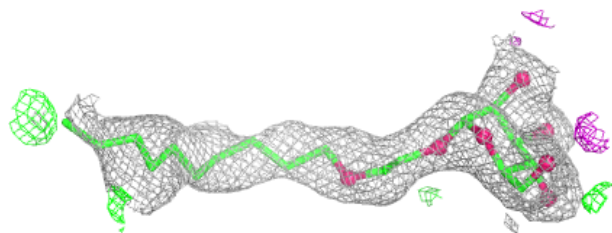
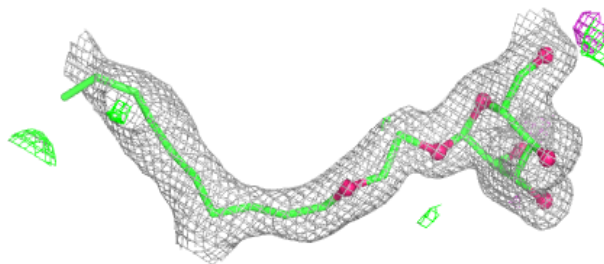


Electron density around CQX C 307:

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

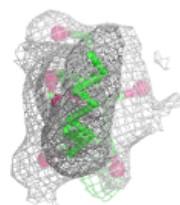
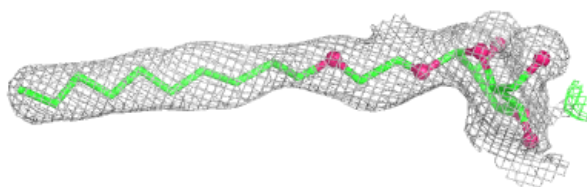
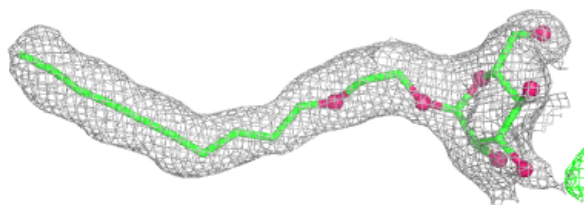
**Electron density around CQX C 306:**

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

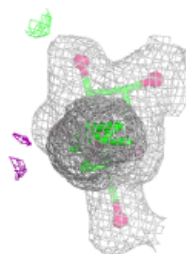
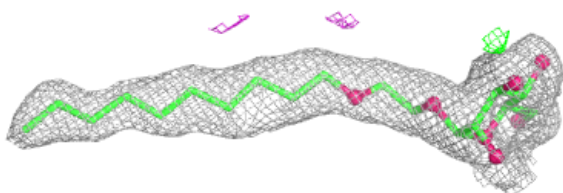
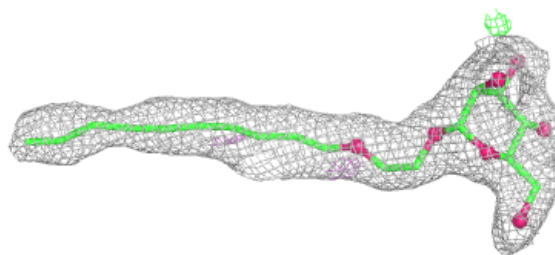


Electron density around CQX A 610:

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

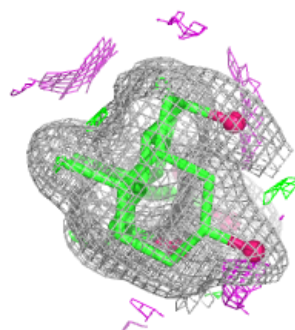
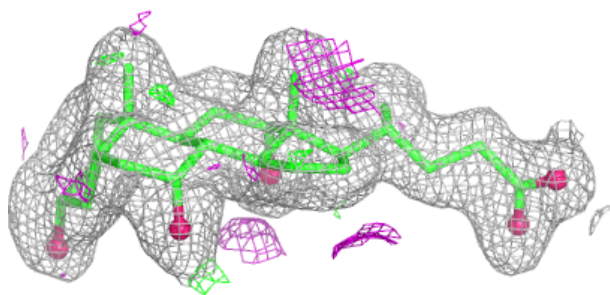
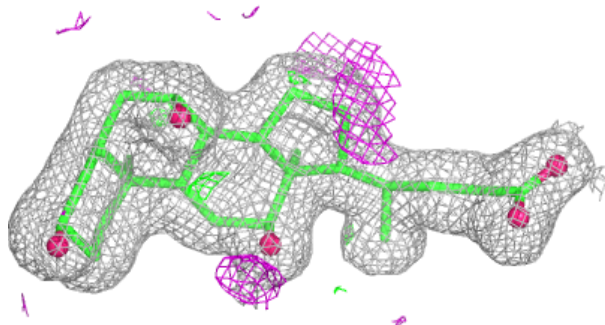
**Electron density around CQX G 101:**

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

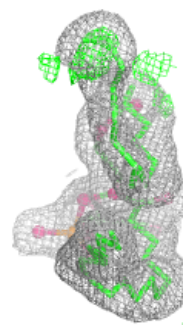
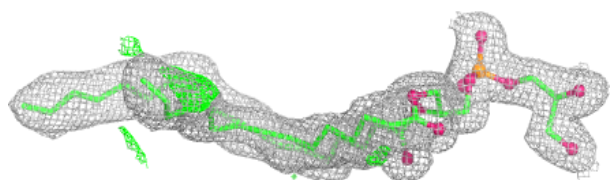
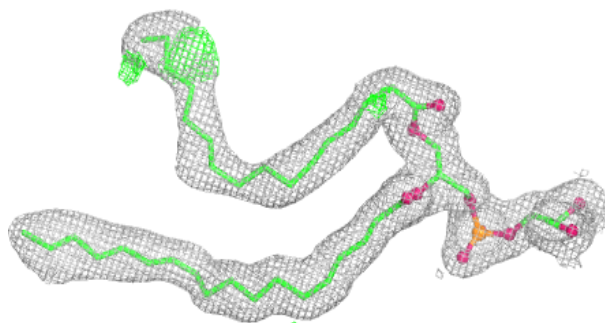


Electron density around CHD C 301:

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

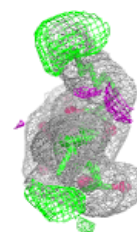
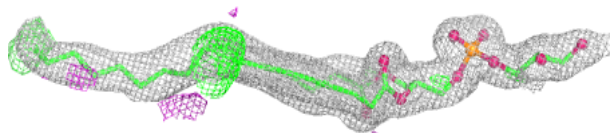
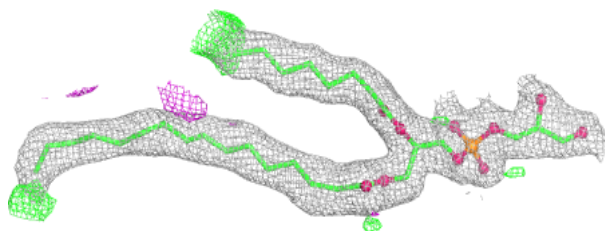
**Electron density around PGV A 608:**

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

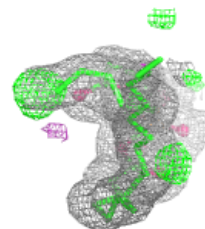
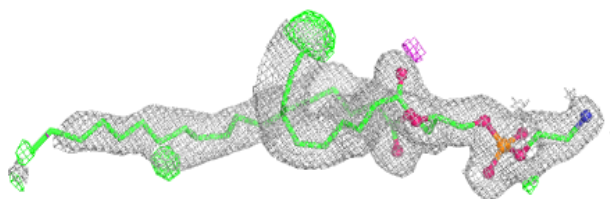
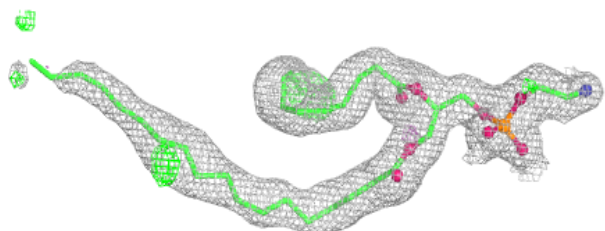


Electron density around PGV C 304:

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

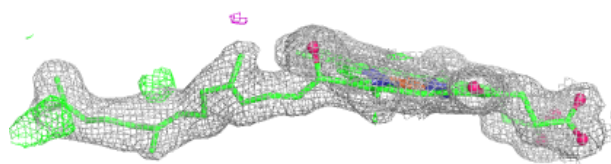
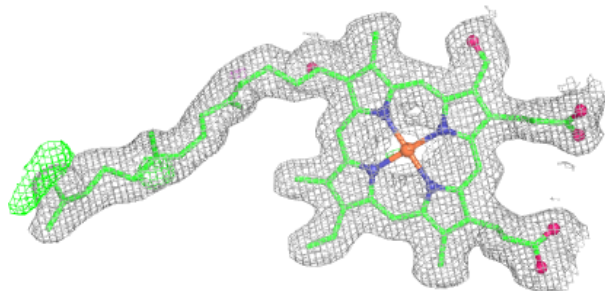
**Electron density around PEK C 303:**

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

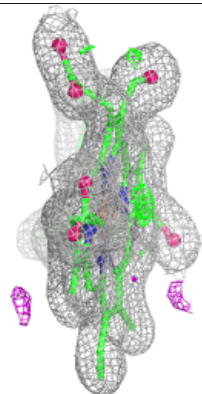
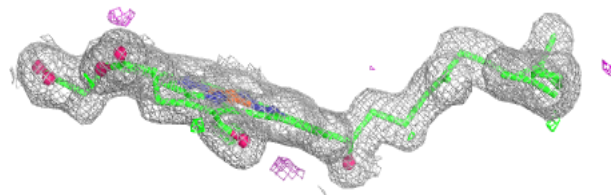
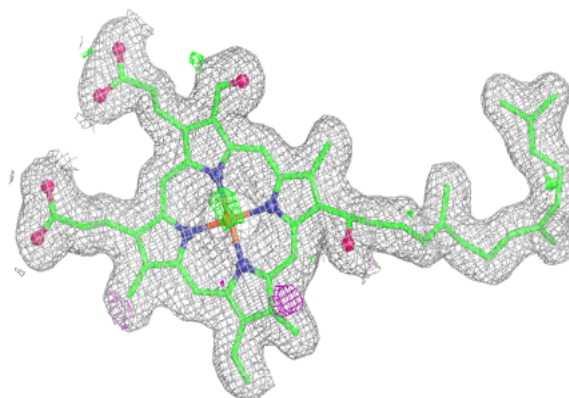


Electron density around HEA A 601:

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

**Electron density around HEA A 602:**

$2mF_o-DF_c$ (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.