



wwPDB X-ray Structure Validation Summary Report ⓘ

Dec 3, 2023 – 12:05 pm GMT

PDB ID : 2J8C
Title : X-ray high resolution structure of the photosynthetic reaction center from Rb. sphaeroides at pH 8 in the neutral state
Authors : Koepke, J.; Diehm, R.; Fritzsich, G.
Deposited on : 2006-10-24
Resolution : 1.87 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity : 4.02b-467
Mogul : 1.8.4, CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.36
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.36

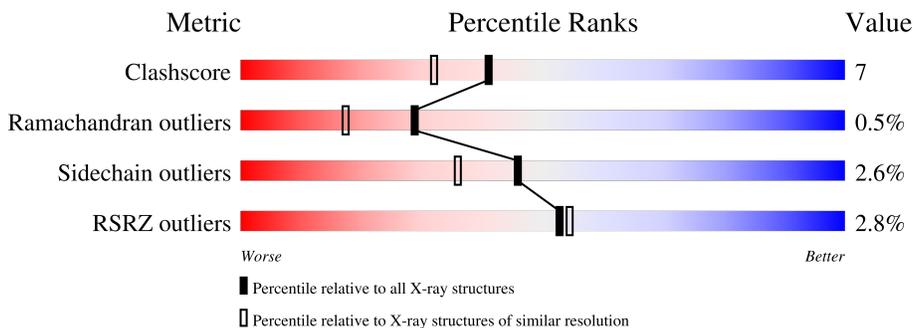
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.87 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))
Clashscore	141614	10282 (1.90-1.86)
Ramachandran outliers	138981	10152 (1.90-1.86)
Sidechain outliers	138945	10152 (1.90-1.86)
RSRZ outliers	127900	9303 (1.90-1.86)

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	H	260	
2	L	281	
3	M	307	

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
10	HTO	L	1291	-	-	-	X
14	PC1	M	1312	X	-	-	-
4	GOL	H	1251	-	-	X	-
5	BCL	L	1282	X	-	-	-
5	BCL	L	1288	X	-	-	-
5	BCL	M	1303	X	-	-	-
5	BCL	M	1304	X	-	-	-

2 Entry composition [i](#)

There are 16 unique types of molecules in this entry. The entry contains 7774 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 REACTION CENTER PROTEIN H CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	H	241	1850	1183	321	337	9	0	5	1

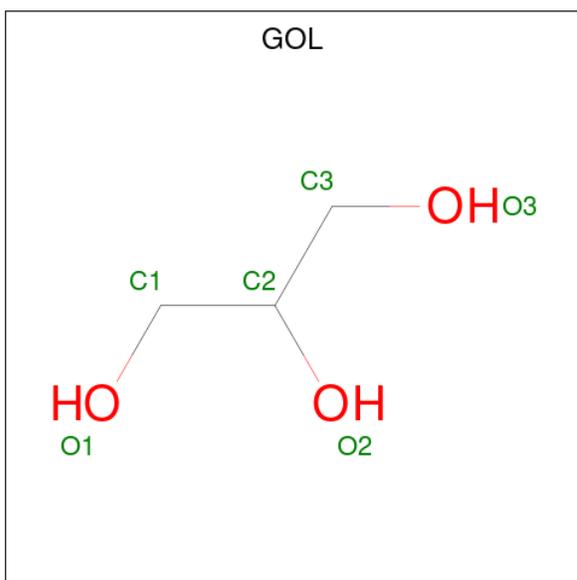
- Molecule 2 is a protein called REACTION CENTER PROTEIN L CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	L	281	2239	1513	355	363	8	0	2	0

- Molecule 3 is a protein called REACTION CENTER PROTEIN M CHAIN.

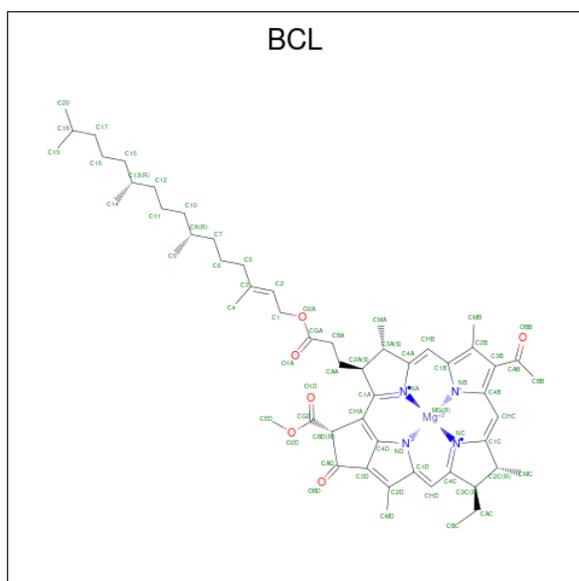
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	M	303	2411	1607	396	398	10	0	1	1

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	H	1	Total	C	O	0	0
			6	3	3		
4	H	1	Total	C	O	0	0
			6	3	3		
4	H	1	Total	C	O	0	0
			6	3	3		
4	H	1	Total	C	O	0	0
			6	3	3		
4	L	1	Total	C	O	0	0
			6	3	3		
4	L	1	Total	C	O	0	0
			6	3	3		
4	L	1	Total	C	O	0	0
			6	3	3		

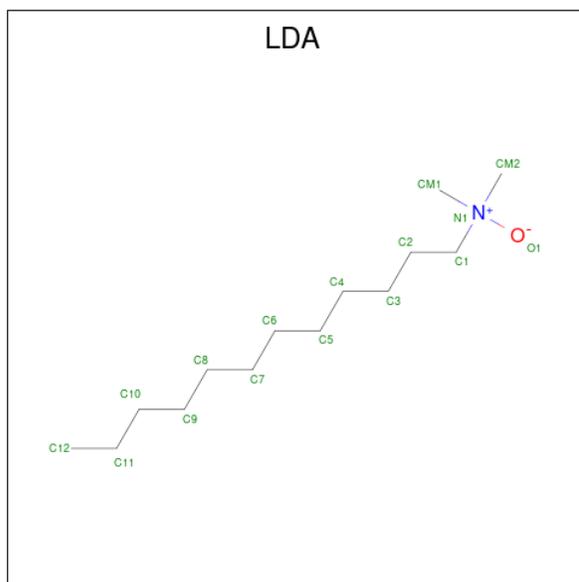
- Molecule 5 is BACTERIOCHLOROPHYLL A (three-letter code: BCL) (formula: $C_{55}H_{74}MgN_4O_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
5	L	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
5	L	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
5	M	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
5	M	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		

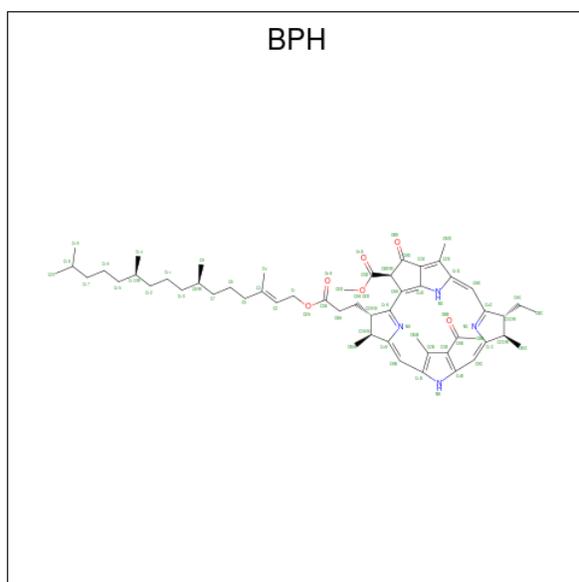
- Molecule 6 is LAURYL DIMETHYLAMINE-N-OXIDE (three-letter code: LDA) (formula:

C₁₄H₃₁NO).



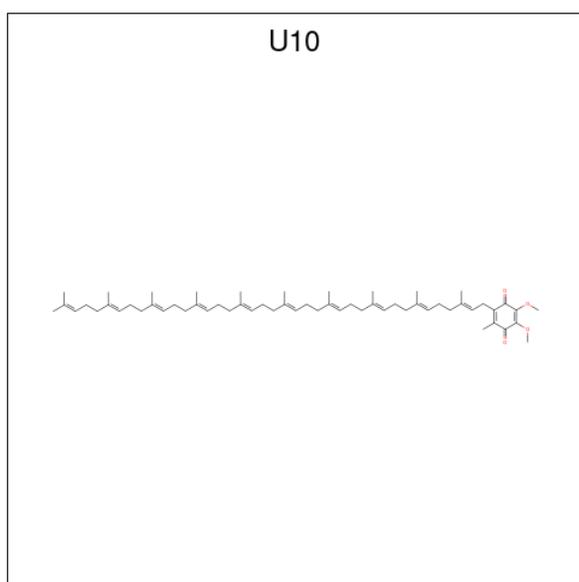
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	
6	L	1	Total	C	N	O	0	0
			16	14	1	1		
6	L	1	Total	C	N	O	0	0
			16	14	1	1		
6	L	1	Total	C	N	O	0	0
			16	14	1	1		
6	M	1	Total	C	N	O	0	0
			16	14	1	1		
6	M	1	Total	C	N	O	0	0
			16	14	1	1		

- Molecule 7 is BACTERIOPHEOPHYTIN A (three-letter code: BPH) (formula: C₅₅H₇₆N₄O₆).



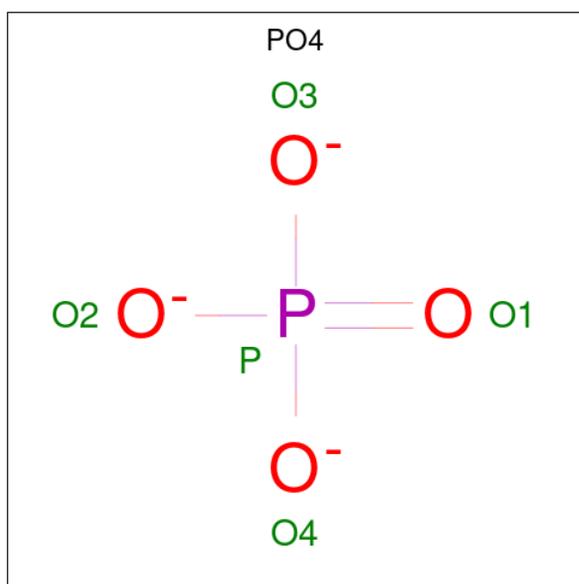
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
7	L	1	65	55	4	6	0	0
7	M	1	65	55	4	6	0	0

- Molecule 8 is UBIQUINONE-10 (three-letter code: U10) (formula: $C_{59}H_{90}O_4$).



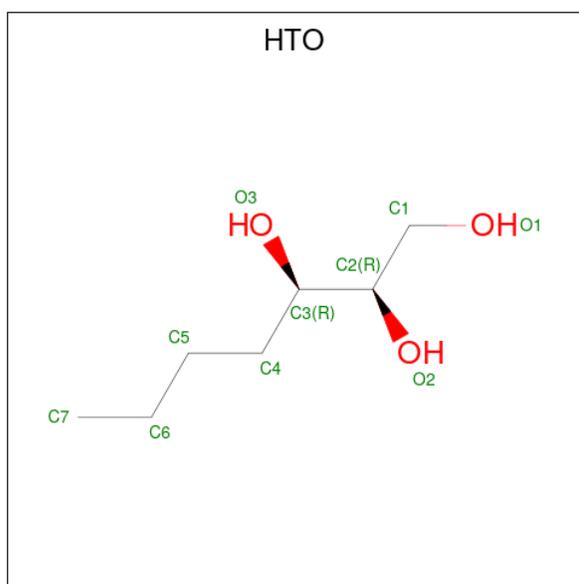
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
8	L	1	46	38	8	0	1
8	M	1	48	44	4	0	0

- Molecule 9 is PHOSPHATE ION (three-letter code: PO4) (formula: O_4P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	L	1	Total	O	P	0	0
			5	4	1		

- Molecule 10 is HEPTANE-1,2,3-TRIOL (three-letter code: HTO) (formula: $C_7H_{16}O_3$).

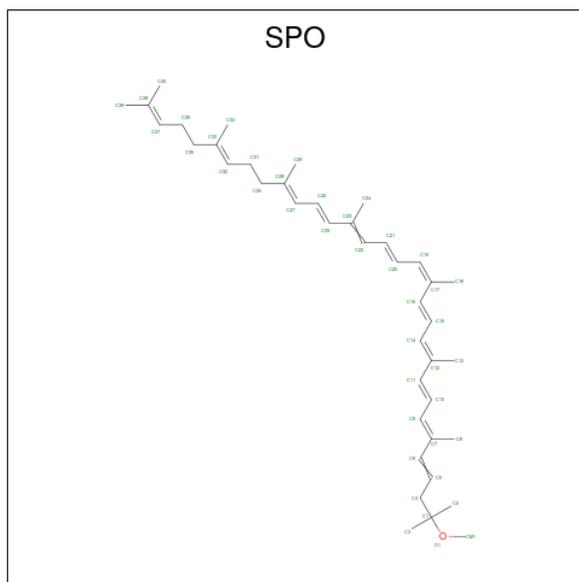


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
10	L	1	Total	C	O	0	0
			10	7	3		
10	L	1	Total	C	O	0	0
			10	7	3		

- Molecule 11 is FE (III) ION (three-letter code: FE) (formula: Fe).

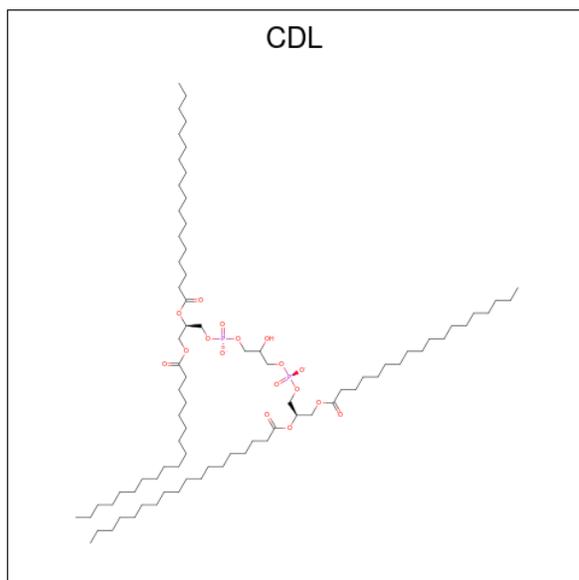
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
11	M	1	Total	Fe	0	0
			1	1		

- Molecule 12 is SPHEROIDENE (three-letter code: SPO) (formula: C₄₁H₆₀O).



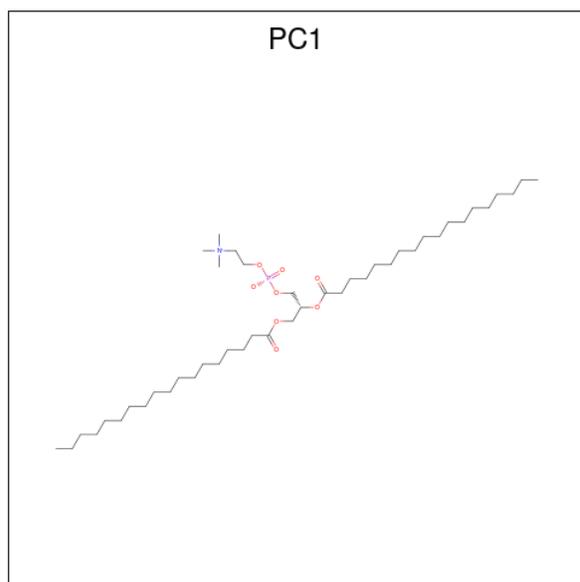
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
12	M	1	Total	C	O	0	0
			42	41	1		

- Molecule 13 is CARDIOLIPIN (three-letter code: CDL) (formula: C₈₁H₁₅₆O₁₇P₂).



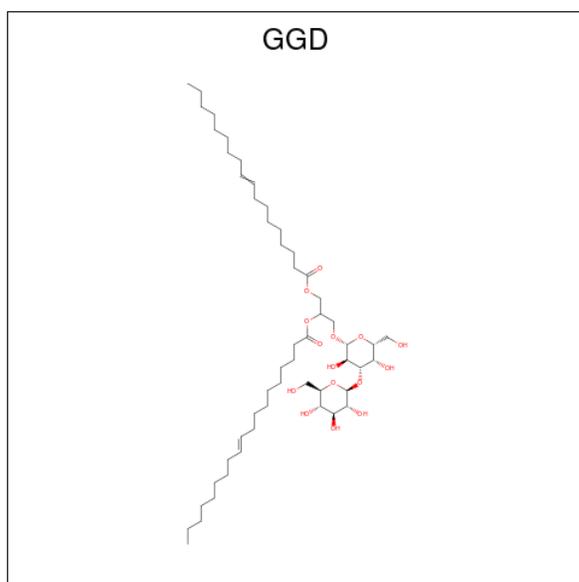
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	O	P		
13	M	1	81	62	17	2	0	0

- Molecule 14 is 1,2-DIACYL-SN-GLYCERO-3-PHOSPHOCHOLINE (three-letter code: PC1) (formula: $C_{44}H_{88}NO_8P$).



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

- Molecule 15 is NONADEC-10-ENOIC ACID 2-[3,4-DIHYDROXY-6-HYDROXYMETHYL-5-(3,4,5-TRIHYDROXY-6-HYDROXYMETHYL-TETRAHYDRO-PYRAN-2-YLOXY)-TETRAHYDRO-PYRAN-2-YLOXY]-1-OCTADEC-9-ENOYLOXYMETHYL-ETHYL ESTER (three-letter code: GGD) (formula: $C_{52}H_{94}O_{15}$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
15	M	1	Total	C	O	0	0
			57	42	15		

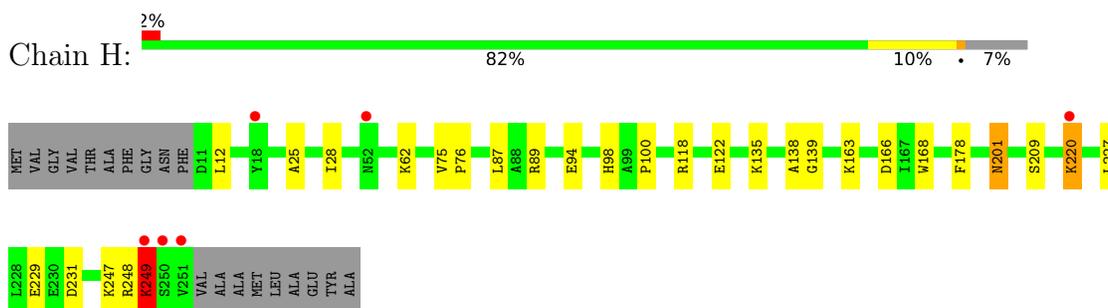
- Molecule 16 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
16	H	172	Total	O	0	0
			172	172		
16	L	116	Total	O	0	0
			116	116		
16	M	127	Total	O	0	0
			127	127		

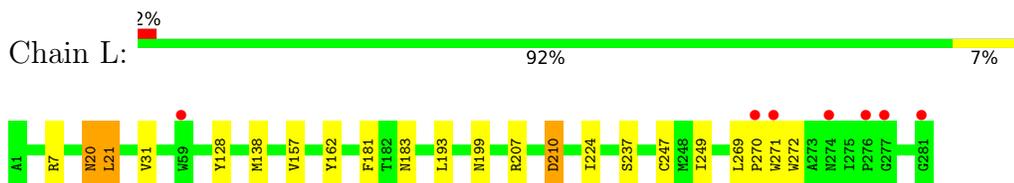
3 Residue-property plots [i](#)

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.

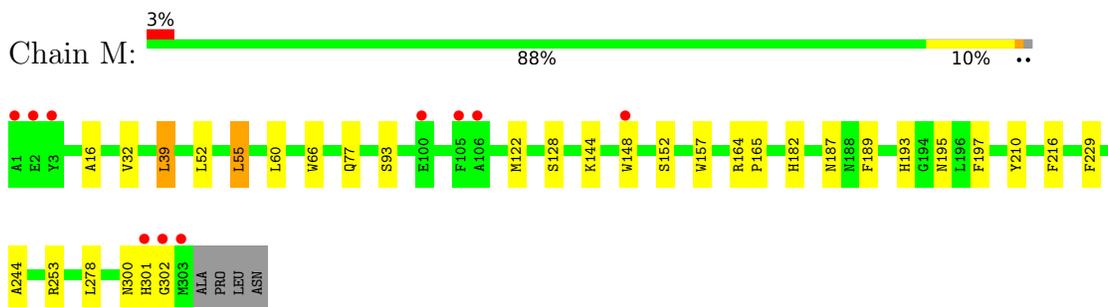
- Molecule 1: REACTION CENTER PROTEIN H CHAIN



- Molecule 2: REACTION CENTER PROTEIN L CHAIN



- Molecule 3: REACTION CENTER PROTEIN M CHAIN



4 Data and refinement statistics i

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	138.69Å 138.69Å 184.61Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	119.52 – 1.87 29.81 – 1.87	Depositor EDS
% Data completeness (in resolution range)	87.0 (119.52-1.87) 87.0 (29.81-1.87)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.90 (at 1.87Å)	Xtrriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.178 , 0.196 0.179 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	23.7	Xtrriage
Anisotropy	0.074	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 70.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	0.018 for -h,-k,l	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7774	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.69% 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: GOL, PC1, FE, PO4, U10, BCL, CDL, LDA, BPH, GGD, SPO, HTO

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	H	0.63	0/1930	0.70	1/2621 (0.0%)
2	L	0.66	0/2339	0.61	1/3203 (0.0%)
3	M	0.65	0/2508	0.64	0/3424
All	All	0.65	0/6777	0.65	2/9248 (0.0%)

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	H	0	1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	139	GLY	N-CA-C	-5.27	99.92	113.10
2	L	210	ASP	CB-CG-OD1	5.25	123.03	118.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	H	138	ALA	Peptide

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	H	1850	0	1873	29	0
2	L	2239	0	2185	20	0
3	M	2411	0	2319	23	0
4	H	24	0	32	7	0
4	L	18	0	24	2	0
5	L	132	0	147	6	0
5	M	132	0	148	16	0
6	L	48	0	93	0	0
6	M	32	0	62	2	0
7	L	65	0	76	4	0
7	M	65	0	76	12	0
8	L	46	0	46	4	0
8	M	48	0	63	1	0
9	L	5	0	0	0	0
10	L	20	0	32	0	0
11	M	1	0	0	0	0
12	M	42	0	60	4	0
13	M	81	0	90	8	0
14	M	43	0	60	0	0
15	M	57	0	67	2	0
16	H	172	0	0	4	1
16	L	116	0	0	1	0
16	M	127	0	0	1	0
All	All	7774	0	7453	105	1

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 7.

The worst 5 of 105 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:248:ARG:HA	1:H:249[A]:LYS:HB2	1.33	1.06
7:M:1308:BPH:HHC	7:M:1308:BPH:HBB3	1.39	1.01
4:H:1251:GOL:H32	16:H:2002:HOH:O	1.67	0.93
7:L:1286:BPH:HHC	7:L:1286:BPH:HBB3	1.58	0.85

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:L:1282:BCL:HBB2	5:L:1282:BCL:HHC	1.59	0.85

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
16:H:2039:HOH:O	16:H:2039:HOH:O[4_555]	1.66	0.54

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	H	245/260 (94%)	241 (98%)	2 (1%)	2 (1%)	19	9
2	L	281/281 (100%)	275 (98%)	5 (2%)	1 (0%)	34	22
3	M	302/307 (98%)	293 (97%)	7 (2%)	2 (1%)	22	11
All	All	828/848 (98%)	809 (98%)	14 (2%)	5 (1%)	29	14

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	H	249[A]	LYS
1	H	249[B]	LYS
3	M	301	HIS
3	M	195	ASN
2	L	31	VAL

5.3.2 Protein sidechains [i](#)

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

resolution.

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	H	200/208 (96%)	190 (95%)	10 (5%)	24	13
2	L	221/220 (100%)	215 (97%)	6 (3%)	44	34
3	M	237/240 (99%)	232 (98%)	5 (2%)	53	45
All	All	658/668 (98%)	637 (97%)	21 (3%)	46	27

5 of 21 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	L	247	CYS
3	M	52	LEU
3	M	216	PHE
3	M	55	LEU
3	M	39	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 10 such sidechains are listed below:

Mol	Chain	Res	Type
3	M	77	GLN
3	M	187	ASN
3	M	193	HIS
2	L	159	ASN
2	L	183	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry

Of 29 ligands modelled in this entry, 1 is monoatomic - leaving 28 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
12	SPO	M	1310	-	40,41,41	3.83	12 (30%)	47,50,50	2.27	20 (42%)
7	BPH	L	1286	-	51,70,70	2.51	9 (17%)	52,101,101	1.99	13 (25%)
4	GOL	H	1251	-	5,5,5	0.48	0	5,5,5	0.81	0
5	BCL	L	1288	2	64,74,74	1.92	12 (18%)	78,115,115	1.98	21 (26%)
8	U10	L	1287[B]	-	23,23,63	2.70	8 (34%)	28,31,79	1.27	3 (10%)
4	GOL	H	1254	-	5,5,5	0.37	0	5,5,5	0.25	0
6	LDA	L	1285	-	12,15,15	2.07	1 (8%)	14,17,17	0.58	0
5	BCL	M	1304	3	64,74,74	2.02	13 (20%)	78,115,115	1.94	19 (24%)
7	BPH	M	1308	-	51,70,70	2.65	9 (17%)	52,101,101	2.21	14 (26%)
9	PO4	L	1289	-	4,4,4	1.04	0	6,6,6	0.75	0
6	LDA	L	1283	-	12,15,15	2.02	1 (8%)	14,17,17	0.45	0
5	BCL	L	1282	2	64,74,74	1.88	13 (20%)	78,115,115	1.97	20 (25%)
8	U10	L	1287[A]	-	23,23,63	2.49	8 (34%)	28,31,79	1.89	7 (25%)
13	CDL	M	1311	-	79,79,99	1.89	16 (20%)	84,90,111	2.71	14 (16%)
6	LDA	M	1305	-	12,15,15	2.03	1 (8%)	14,17,17	0.48	0
8	U10	M	1309	-	48,48,63	2.54	11 (22%)	58,61,79	1.59	11 (18%)
14	PC1	M	1312	-	42,42,53	1.74	8 (19%)	48,50,61	1.32	6 (12%)
10	HTO	L	1291	-	9,9,9	0.33	0	10,10,10	0.87	1 (10%)
10	HTO	L	1290	-	9,9,9	0.29	0	10,10,10	0.79	1 (10%)
4	GOL	L	1293	-	5,5,5	0.41	0	5,5,5	0.27	0
4	GOL	L	1292	-	5,5,5	0.46	0	5,5,5	0.86	0
15	GGD	M	1313	-	58,58,68	2.14	12 (20%)	72,72,82	1.79	18 (25%)
4	GOL	L	1294	-	5,5,5	0.36	0	5,5,5	0.18	0
6	LDA	L	1284	-	12,15,15	2.03	1 (8%)	14,17,17	0.45	0
4	GOL	H	1252	-	5,5,5	0.43	0	5,5,5	0.39	0
4	GOL	H	1253	-	5,5,5	0.36	0	5,5,5	0.19	0
6	LDA	M	1306	-	12,15,15	2.01	1 (8%)	14,17,17	0.54	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	BCL	M	1303	3	64,74,74	1.90	11 (17%)	78,115,115	1.73	11 (14%)

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
12	SPO	M	1310	-	-	6/47/47/47	-
7	BPH	L	1286	-	-	6/37/105/105	0/5/6/6
4	GOL	H	1251	-	-	2/4/4/4	-
5	BCL	L	1288	2	2/2/21/25	10/37/137/137	-
8	U10	L	1287[B]	-	-	8/15/39/87	0/1/1/1
4	GOL	H	1254	-	-	0/4/4/4	-
6	LDA	L	1285	-	-	10/13/13/13	-
5	BCL	M	1304	3	2/2/21/25	10/37/137/137	-
7	BPH	M	1308	-	-	17/37/105/105	0/5/6/6
6	LDA	L	1283	-	-	7/13/13/13	-
5	BCL	L	1282	2	2/2/21/25	9/37/137/137	-
8	U10	L	1287[A]	-	-	7/15/39/87	0/1/1/1
13	CDL	M	1311	-	-	50/88/88/110	-
6	LDA	M	1305	-	-	5/13/13/13	-
14	PC1	M	1312	-	1/1/5/5	25/46/46/57	-
8	U10	M	1309	-	-	6/45/69/87	0/1/1/1
10	HTO	L	1291	-	-	9/10/10/10	-
10	HTO	L	1290	-	-	10/10/10/10	-
4	GOL	L	1293	-	-	2/4/4/4	-
4	GOL	L	1292	-	-	2/4/4/4	-
15	GGD	M	1313	-	-	19/47/87/97	0/2/2/2
4	GOL	L	1294	-	-	1/4/4/4	-
6	LDA	L	1284	-	-	7/13/13/13	-
4	GOL	H	1252	-	-	2/4/4/4	-
4	GOL	H	1253	-	-	2/4/4/4	-
6	LDA	M	1306	-	-	7/13/13/13	-
5	BCL	M	1303	3	2/2/21/25	18/37/137/137	-

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

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
12	M	1310	SPO	C27-C28	12.95	1.47	1.34
7	M	1308	BPH	OBD-CAD	11.84	1.38	1.22
7	L	1286	BPH	OBD-CAD	10.36	1.36	1.22
5	L	1282	BCL	OBD-CAD	9.22	1.38	1.22
5	M	1304	BCL	OBD-CAD	9.10	1.38	1.22

The worst 5 of 179 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
13	M	1311	CDL	C13-C12-C11	12.96	159.78	113.19
13	M	1311	CDL	C17-C16-C15	11.70	173.84	114.42
7	M	1308	BPH	O2D-CGD-CBD	10.24	123.98	111.00
13	M	1311	CDL	C20-C19-C18	9.40	162.17	114.42
7	L	1286	BPH	O2D-CGD-CBD	8.13	121.29	111.00

5 of 9 chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
5	L	1282	BCL	C13
5	L	1282	BCL	C8
5	L	1288	BCL	C13
5	L	1288	BCL	C8
5	M	1303	BCL	C13

5 of 257 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	H	1252	GOL	C1-C2-C3-O3
4	L	1292	GOL	C1-C2-C3-O3
4	L	1293	GOL	O1-C1-C2-C3
5	M	1303	BCL	C1-C2-C3-C4
5	M	1303	BCL	C1-C2-C3-C5

There are no ring outliers.

17 monomers are involved in 62 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
12	M	1310	SPO	4	0
7	L	1286	BPH	4	0
4	H	1251	GOL	6	0
5	L	1288	BCL	2	0
5	M	1304	BCL	11	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	M	1308	BPH	12	0
5	L	1282	BCL	4	0
8	L	1287[A]	U10	4	0
13	M	1311	CDL	8	0
6	M	1305	LDA	1	0
8	M	1309	U10	1	0
4	L	1292	GOL	1	0
15	M	1313	GGD	2	0
4	L	1294	GOL	1	0
4	H	1252	GOL	1	0
6	M	1306	LDA	1	0
5	M	1303	BCL	7	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

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

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

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	H	241/260 (92%)	-0.33	6 (2%) 57 59	17, 24, 33, 61	3 (1%)
2	L	281/281 (100%)	-0.49	7 (2%) 57 59	16, 21, 37, 50	0
3	M	303/307 (98%)	-0.24	10 (3%) 46 47	14, 25, 42, 56	6 (1%)
All	All	825/848 (97%)	-0.35	23 (2%) 53 54	14, 23, 41, 61	9 (1%)

The worst 5 of 23 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	H	251	VAL	9.9
1	H	250	SER	8.3
3	M	1	ALA	5.6
3	M	301	HIS	5.3
2	L	281	GLY	4.7

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

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

6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

6.4 Ligands [i](#)

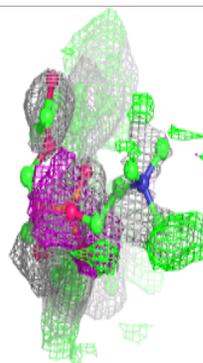
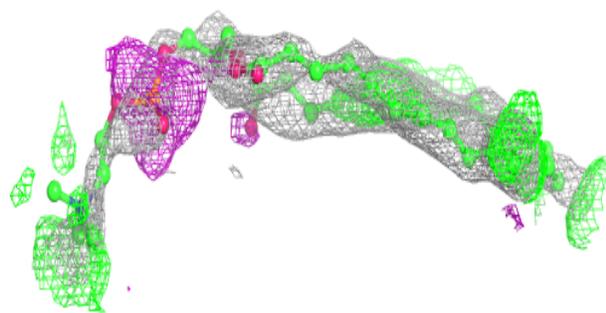
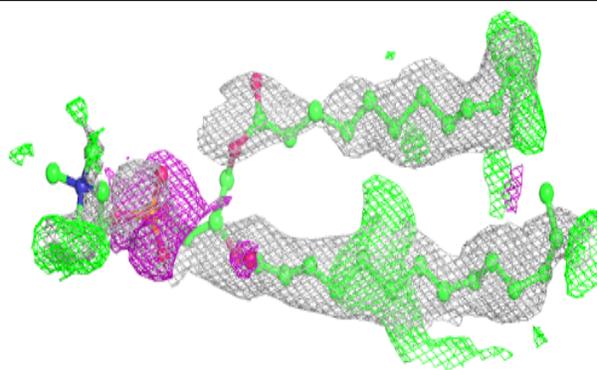
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
14	PC1	M	1312	43/54	0.33	0.36	66,82,95,96	0
10	HTO	L	1291	10/10	0.35	0.56	78,80,81,82	0
6	LDA	L	1284	16/16	0.43	0.34	72,73,77,77	0
15	GGD	M	1313	57/67	0.43	0.36	47,73,90,90	0
6	LDA	L	1285	16/16	0.56	0.34	72,76,80,81	0
6	LDA	L	1283	16/16	0.56	0.30	41,64,73,74	0
4	GOL	H	1253	6/6	0.59	0.28	65,66,67,68	0
4	GOL	L	1294	6/6	0.60	0.18	80,81,81,81	0
4	GOL	L	1292	6/6	0.64	0.21	29,42,44,45	0
8	U10	L	1287[B]	23/63	0.70	0.34	34,42,44,44	23
8	U10	L	1287[A]	23/63	0.70	0.34	24,40,48,49	23
13	CDL	M	1311	81/100	0.71	0.38	42,69,83,83	0
10	HTO	L	1290	10/10	0.72	0.28	60,62,63,64	0
4	GOL	H	1251	6/6	0.79	0.27	32,46,47,49	0
6	LDA	M	1306	16/16	0.79	0.23	54,58,63,64	0
4	GOL	H	1254	6/6	0.79	0.19	63,64,64,65	0
4	GOL	H	1252	6/6	0.80	0.26	63,63,63,64	0
6	LDA	M	1305	16/16	0.80	0.20	31,45,51,51	0
12	SPO	M	1310	42/42	0.88	0.15	21,26,48,51	0
8	U10	M	1309	48/63	0.92	0.13	15,25,45,48	0
4	GOL	L	1293	6/6	0.93	0.15	66,68,68,69	0
7	BPH	M	1308	65/65	0.94	0.12	17,22,69,71	0
5	BCL	M	1304	66/66	0.96	0.11	12,19,38,46	0
5	BCL	M	1303	66/66	0.96	0.12	16,21,58,58	0
5	BCL	L	1288	66/66	0.97	0.10	14,17,31,37	0
5	BCL	L	1282	66/66	0.97	0.11	13,18,41,43	0
7	BPH	L	1286	65/65	0.98	0.08	10,16,26,28	0
9	PO4	L	1289	5/5	0.99	0.13	33,35,36,36	0
11	FE	M	1307	1/1	1.00	0.04	15,15,15,15	0

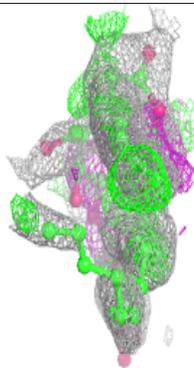
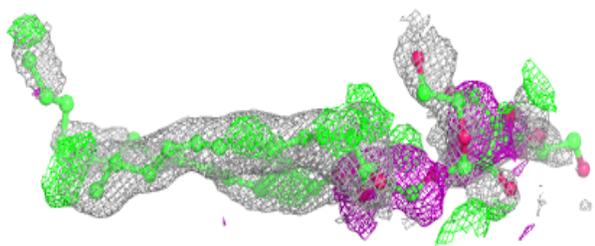
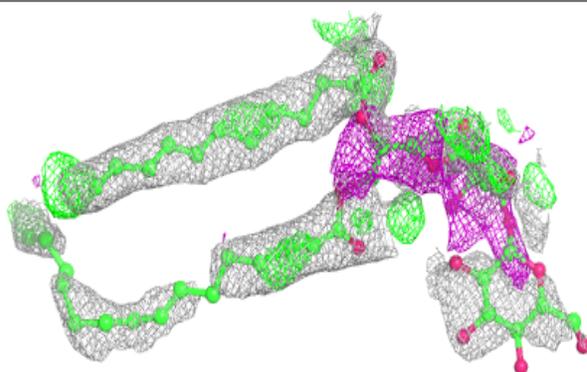
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around PC1 M 1312:

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

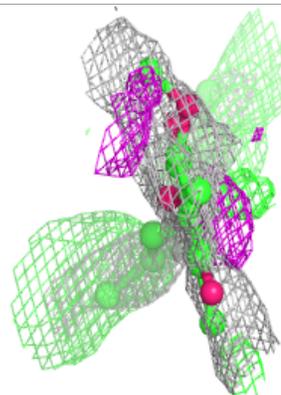
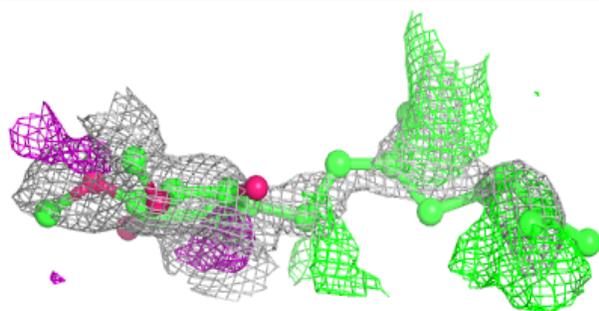
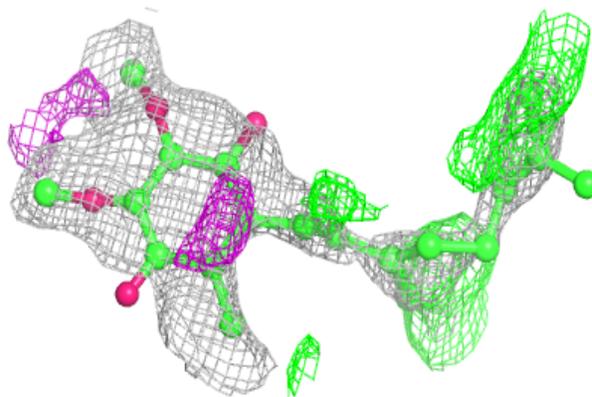
**Electron density around GGD M 1313:**

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

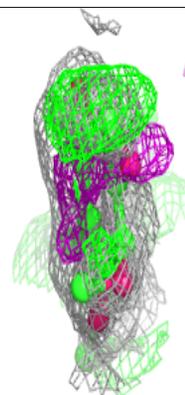
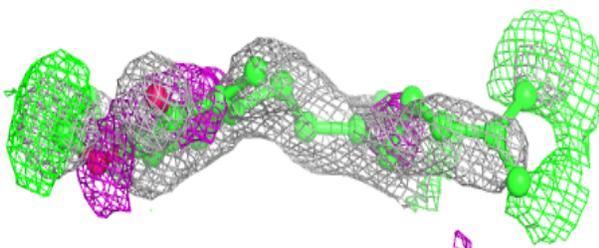
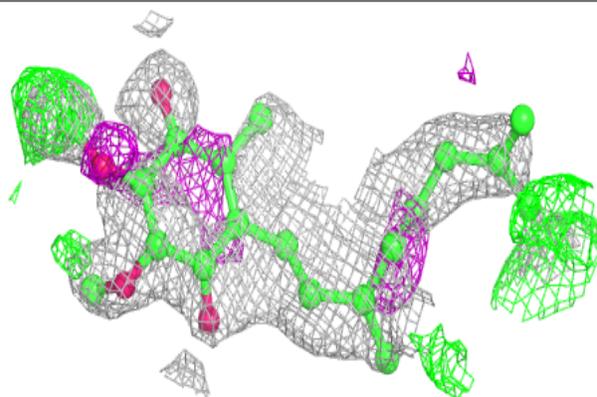


Electron density around U10 L 1287 (B):

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

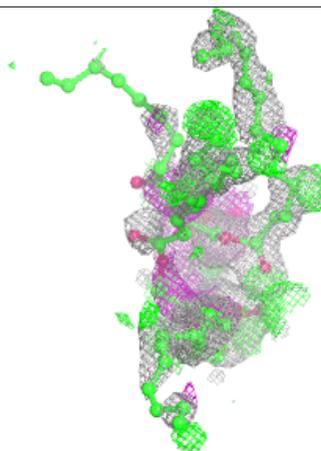
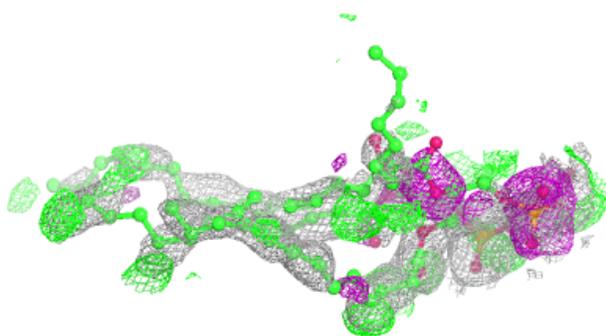
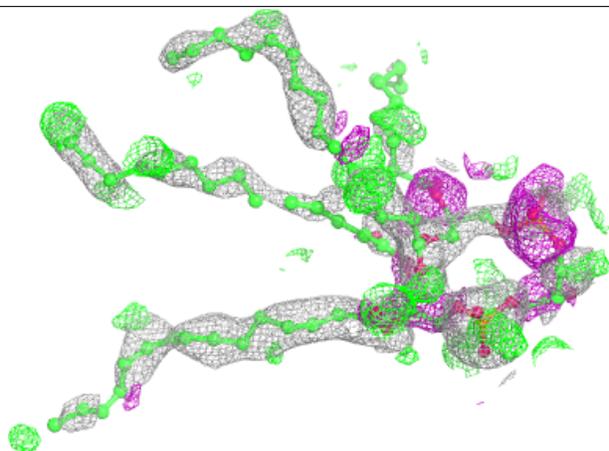
**Electron density around U10 L 1287 (A):**

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

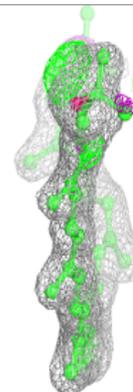
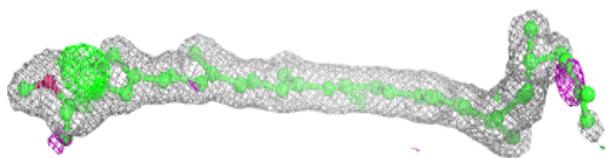
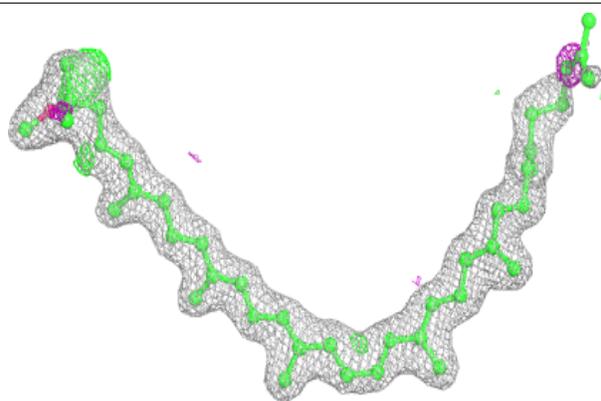


Electron density around CDL M 1311:

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and green (positive)

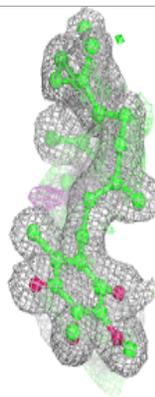
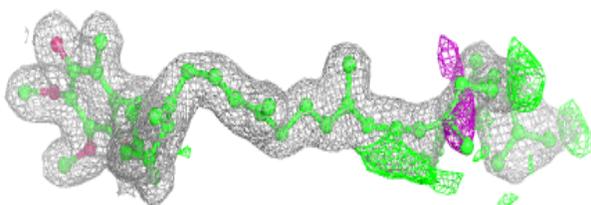
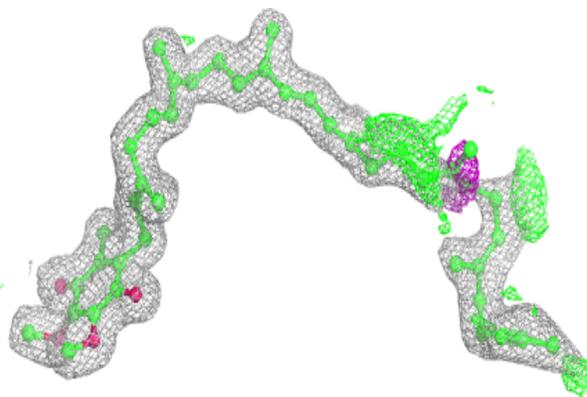
**Electron density around SPO M 1310:**

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and green (positive)

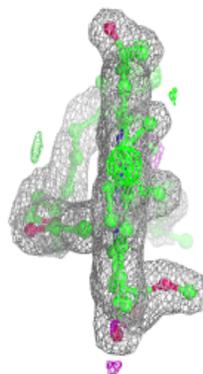
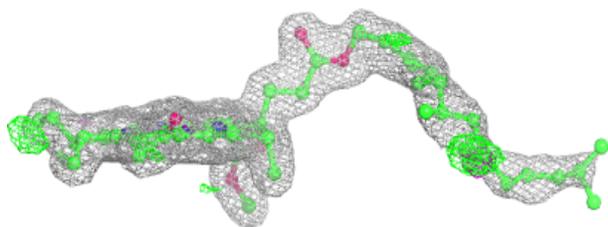
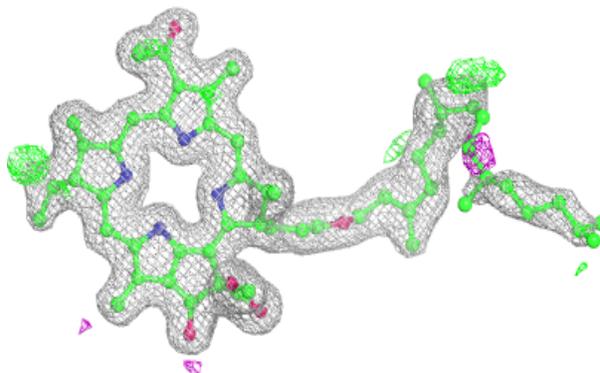


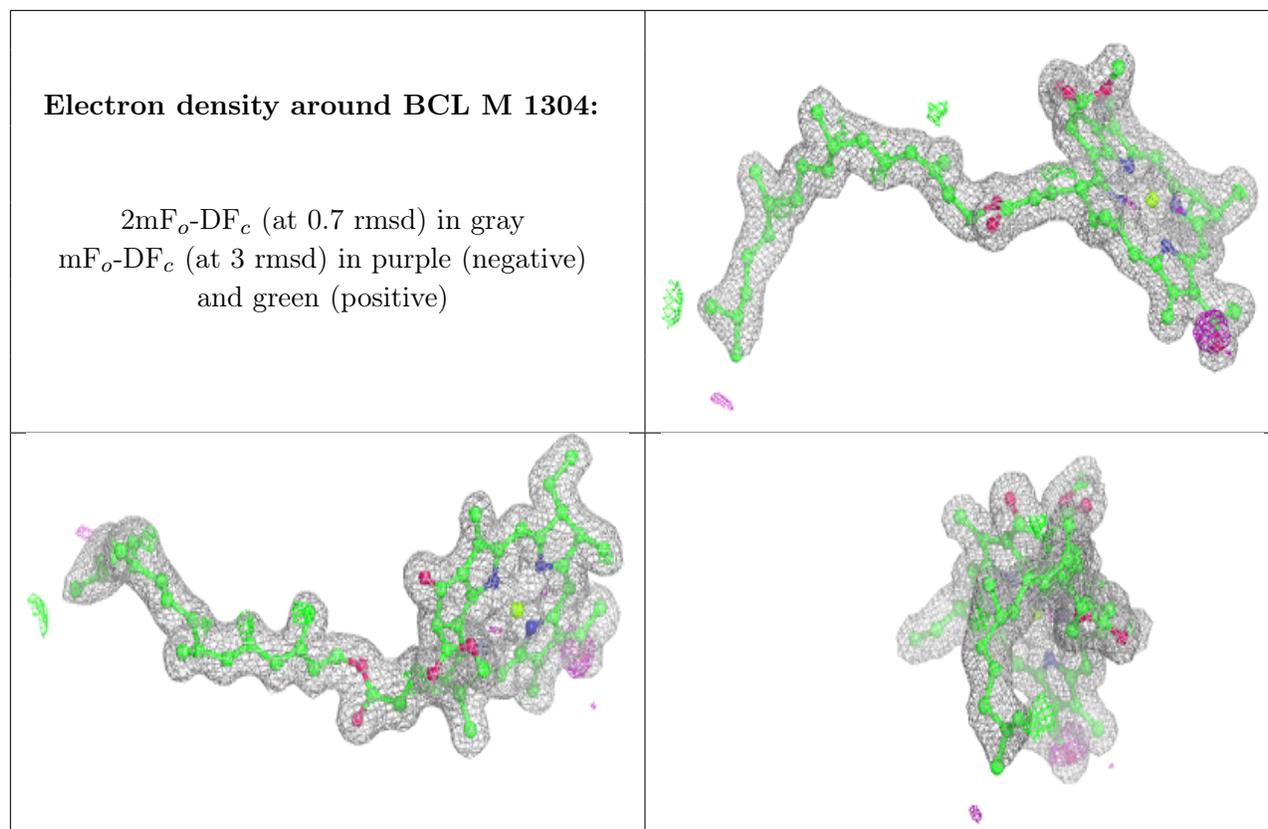
Electron density around U10 M 1309:

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

**Electron density around BPH M 1308:**

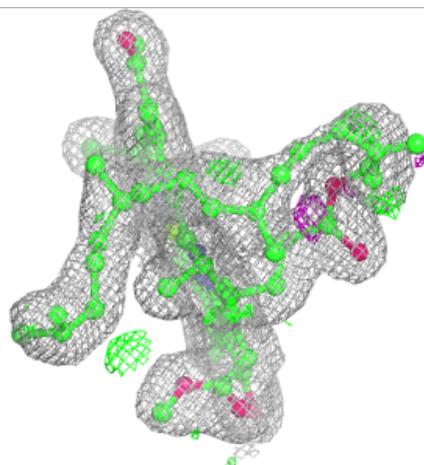
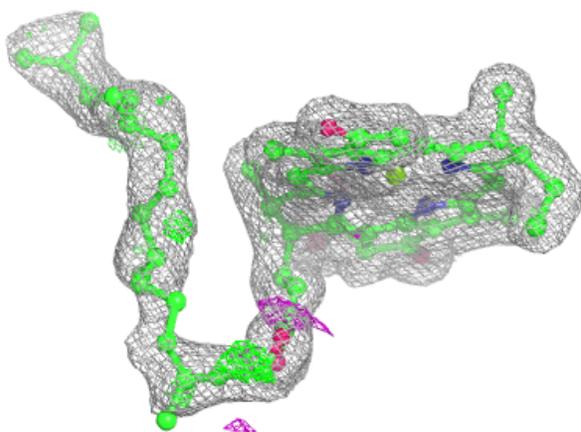
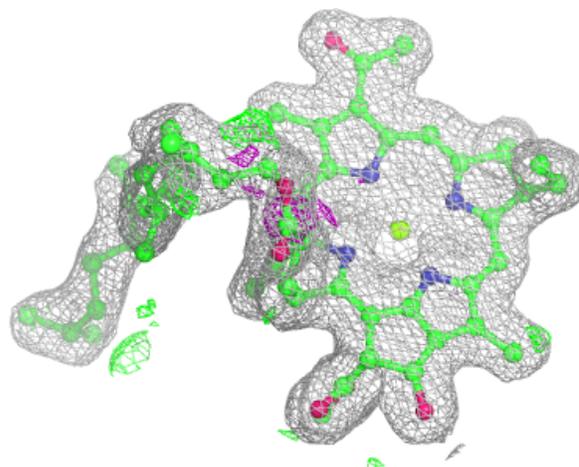
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

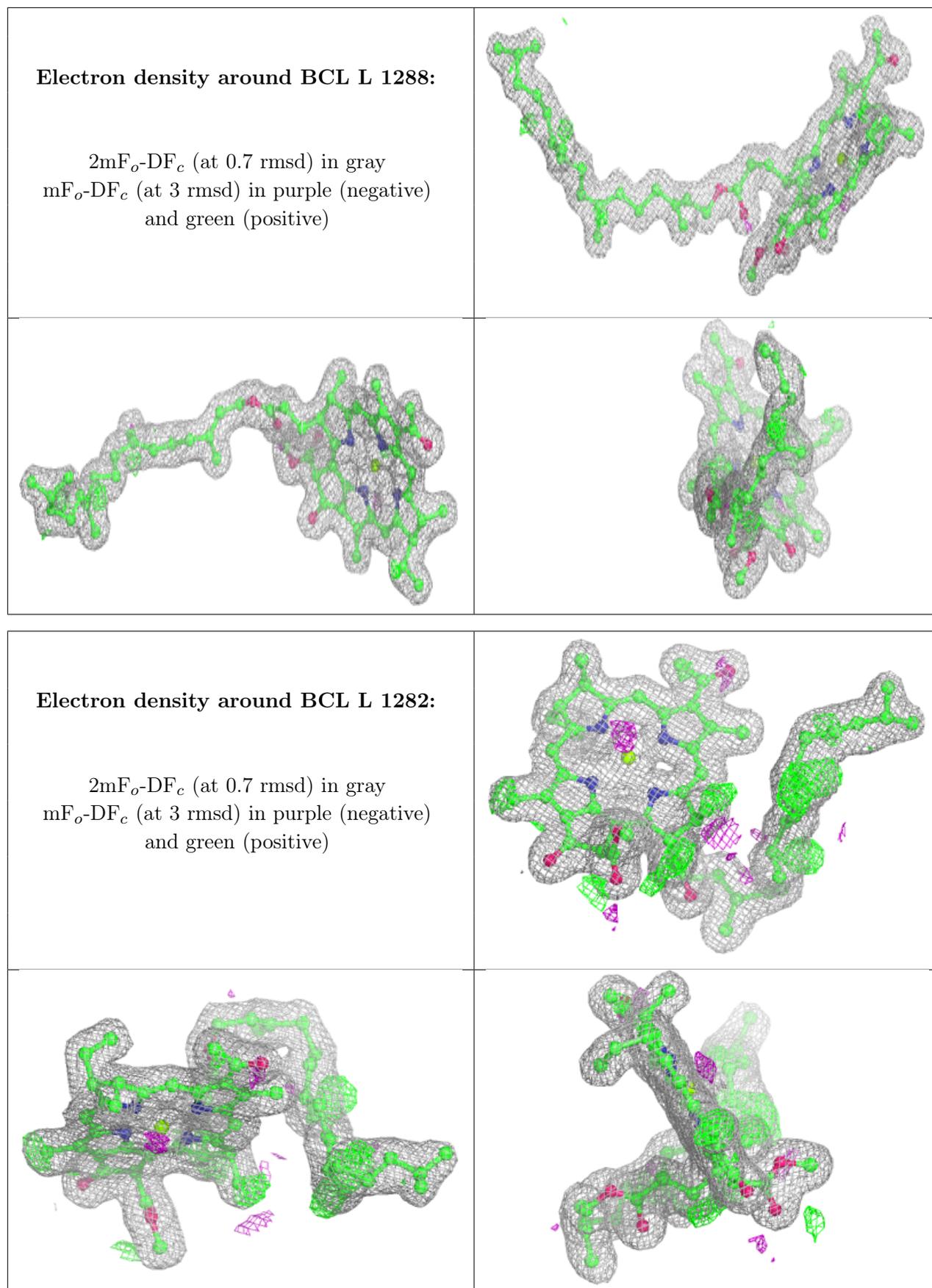


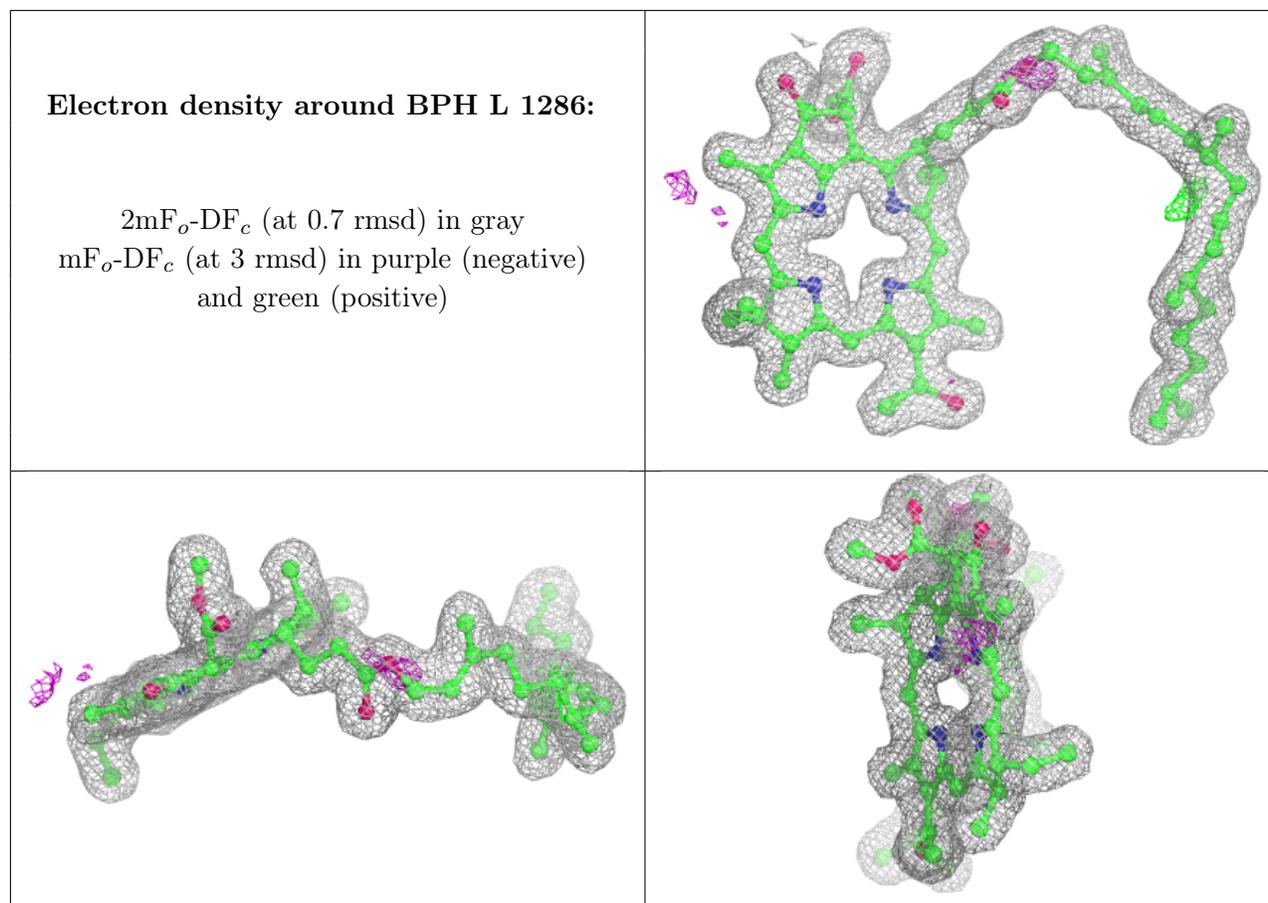


Electron density around BCL M 1303:

$2mF_o-DF_c$ (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.