



# wwPDB X-ray Structure Validation Summary Report ⓘ

May 24, 2020 – 01:45 am BST

PDB ID : 3NZ2  
Title : Crystal Structure of Hexapeptide-Repeat containing-Acetyltransferase VCA0836 Complexed with Acetyl Co Enzyme A from *Vibrio cholerae* O1 bio-var eltor  
Authors : Kim, Y.; Maltseva, N.; Hasseman, J.; Anderson, W.F.; Joachimiak, A.; Center for Structural Genomics of Infectious Diseases (CSGID)  
Deposited on : 2010-07-15  
Resolution : 2.35 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](mailto: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.

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

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.11  
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) : 2.11

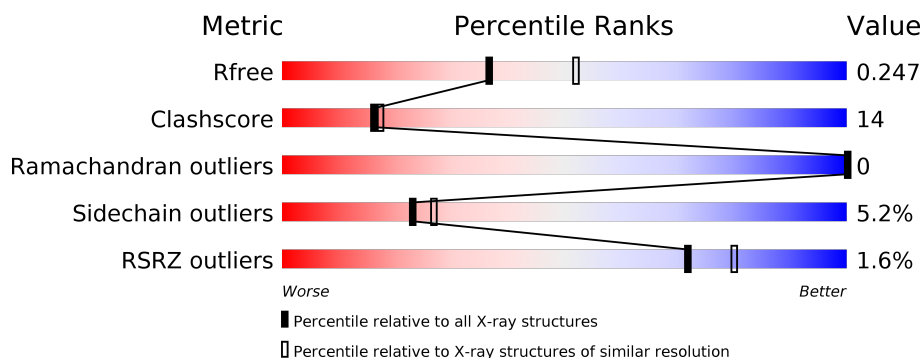
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.35 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	130704	1164 (2.36-2.36)
Clashscore	141614	1232 (2.36-2.36)
Ramachandran outliers	138981	1211 (2.36-2.36)
Sidechain outliers	138945	1212 (2.36-2.36)
RSRZ outliers	127900	1150 (2.36-2.36)

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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	195	<div> <div>75%</div> <div>17%</div> <div>• 7%</div> </div>
1	B	195	<div>2%</div> <div>71%</div> <div>23%</div> <div>6%</div>

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Mol	Chain	Length	Quality of chain
1	G	195	 71% 20% • 7%
1	H	195	 69% 24% • 6%
1	I	195	 67% 25% • 6%
1	J	195	 72% 22% • 5%
1	K	195	 68% 24% • 6%
1	L	195	 69% 23% • 6%

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
3	ACY	K	198	-	-	X	-
4	GOL	E	196	-	-	X	-

## 2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 18822 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 Hexapeptide-repeat containing-acetyltransferase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	182	Total	C	N	O	S	Se	0	1	0
			1420	891	257	265	3	4			
1	B	184	Total	C	N	O	S	Se	0	1	0
			1437	902	260	267	3	5			
1	C	183	Total	C	N	O	S	Se	0	0	0
			1417	889	254	267	3	4			
1	D	183	Total	C	N	O	S	Se	0	1	0
			1431	900	256	268	3	4			
1	E	183	Total	C	N	O	S	Se	0	1	0
			1428	895	258	268	3	4			
1	F	183	Total	C	N	O	S	Se	0	0	0
			1417	889	254	267	3	4			
1	G	182	Total	C	N	O	S	Se	0	1	0
			1418	890	254	267	3	4			
1	H	183	Total	C	N	O	S	Se	0	1	0
			1428	895	258	268	3	4			
1	I	184	Total	C	N	O	S	Se	0	1	0
			1437	901	260	269	3	4			
1	J	185	Total	C	N	O	S	Se	0	0	0
			1431	899	257	267	3	5			
1	K	183	Total	C	N	O	S	Se	0	0	0
			1417	889	254	267	3	4			
1	L	183	Total	C	N	O	S	Se	0	1	0
			1428	895	258	268	3	4			

There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	EXPRESSION TAG	UNP Q9KLB0
A	-1	ASN	-	EXPRESSION TAG	UNP Q9KLB0
A	0	ALA	-	EXPRESSION TAG	UNP Q9KLB0
A	1	MSE	-	EXPRESSION TAG	UNP Q9KLB0
B	-2	SER	-	EXPRESSION TAG	UNP Q9KLB0

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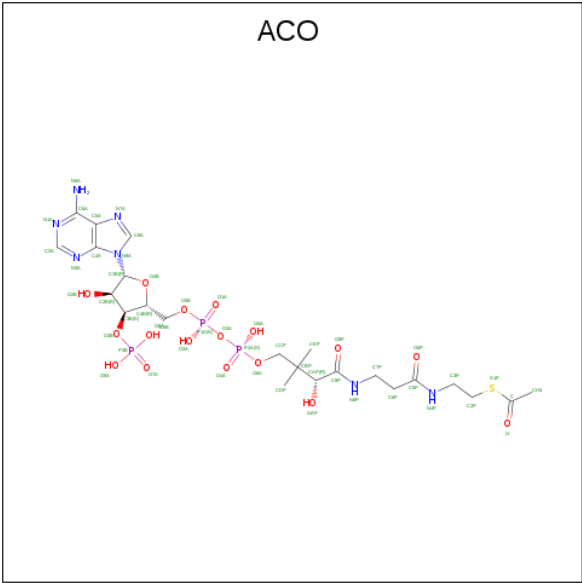
Chain	Residue	Modelled	Actual	Comment	Reference
B	-1	ASN	-	EXPRESSION TAG	UNP Q9KLB0
B	0	ALA	-	EXPRESSION TAG	UNP Q9KLB0
B	1	MSE	-	EXPRESSION TAG	UNP Q9KLB0
C	-2	SER	-	EXPRESSION TAG	UNP Q9KLB0
C	-1	ASN	-	EXPRESSION TAG	UNP Q9KLB0
C	0	ALA	-	EXPRESSION TAG	UNP Q9KLB0
C	1	MSE	-	EXPRESSION TAG	UNP Q9KLB0
D	-2	SER	-	EXPRESSION TAG	UNP Q9KLB0
D	-1	ASN	-	EXPRESSION TAG	UNP Q9KLB0
D	0	ALA	-	EXPRESSION TAG	UNP Q9KLB0
D	1	MSE	-	EXPRESSION TAG	UNP Q9KLB0
E	-2	SER	-	EXPRESSION TAG	UNP Q9KLB0
E	-1	ASN	-	EXPRESSION TAG	UNP Q9KLB0
E	0	ALA	-	EXPRESSION TAG	UNP Q9KLB0
E	1	MSE	-	EXPRESSION TAG	UNP Q9KLB0
F	-2	SER	-	EXPRESSION TAG	UNP Q9KLB0
F	-1	ASN	-	EXPRESSION TAG	UNP Q9KLB0
F	0	ALA	-	EXPRESSION TAG	UNP Q9KLB0
F	1	MSE	-	EXPRESSION TAG	UNP Q9KLB0
G	-2	SER	-	EXPRESSION TAG	UNP Q9KLB0
G	-1	ASN	-	EXPRESSION TAG	UNP Q9KLB0
G	0	ALA	-	EXPRESSION TAG	UNP Q9KLB0
G	1	MSE	-	EXPRESSION TAG	UNP Q9KLB0
H	-2	SER	-	EXPRESSION TAG	UNP Q9KLB0
H	-1	ASN	-	EXPRESSION TAG	UNP Q9KLB0
H	0	ALA	-	EXPRESSION TAG	UNP Q9KLB0
H	1	MSE	-	EXPRESSION TAG	UNP Q9KLB0
I	-2	SER	-	EXPRESSION TAG	UNP Q9KLB0
I	-1	ASN	-	EXPRESSION TAG	UNP Q9KLB0
I	0	ALA	-	EXPRESSION TAG	UNP Q9KLB0
I	1	MSE	-	EXPRESSION TAG	UNP Q9KLB0
J	-2	SER	-	EXPRESSION TAG	UNP Q9KLB0
J	-1	ASN	-	EXPRESSION TAG	UNP Q9KLB0
J	0	ALA	-	EXPRESSION TAG	UNP Q9KLB0
J	1	MSE	-	EXPRESSION TAG	UNP Q9KLB0
K	-2	SER	-	EXPRESSION TAG	UNP Q9KLB0
K	-1	ASN	-	EXPRESSION TAG	UNP Q9KLB0
K	0	ALA	-	EXPRESSION TAG	UNP Q9KLB0
K	1	MSE	-	EXPRESSION TAG	UNP Q9KLB0
L	-2	SER	-	EXPRESSION TAG	UNP Q9KLB0
L	-1	ASN	-	EXPRESSION TAG	UNP Q9KLB0
L	0	ALA	-	EXPRESSION TAG	UNP Q9KLB0

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Chain	Residue	Modelled	Actual	Comment	Reference
L	1	MSE	-	EXPRESSION TAG	UNP Q9KLB0

- Molecule 2 is ACETYL COENZYME \*A (three-letter code: ACO) (formula: C<sub>23</sub>H<sub>38</sub>N<sub>7</sub>O<sub>17</sub>P<sub>3</sub>S).



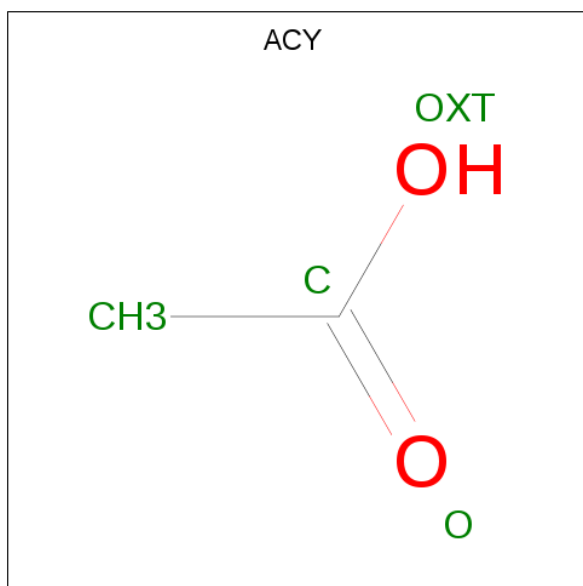
Mol	Chain	Residues	Atoms							ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	S		0	0
			51	23	7	17	3	1			
2	A	1	Total	C	N	O	P	S		0	0
			51	23	7	17	3	1			
2	C	1	Total	C	N	O	P	S		0	0
			51	23	7	17	3	1			
2	D	1	Total	C	N	O	P	S		0	0
			51	23	7	17	3	1			
2	E	1	Total	C	N	O	P	S		0	0
			51	23	7	17	3	1			
2	F	1	Total	C	N	O	P	S		0	0
			51	23	7	17	3	1			
2	H	1	Total	C	N	O	P	S		0	0
			51	23	7	17	3	1			
2	I	1	Total	C	N	O	P	S		0	0
			51	23	7	17	3	1			
2	I	1	Total	C	N	O	P	S		0	0
			51	23	7	17	3	1			
2	J	1	Total	C	N	O	P	S		0	0
			51	23	7	17	3	1			
2	L	1	Total	C	N	O	P	S		0	0
			51	23	7	17	3	1			

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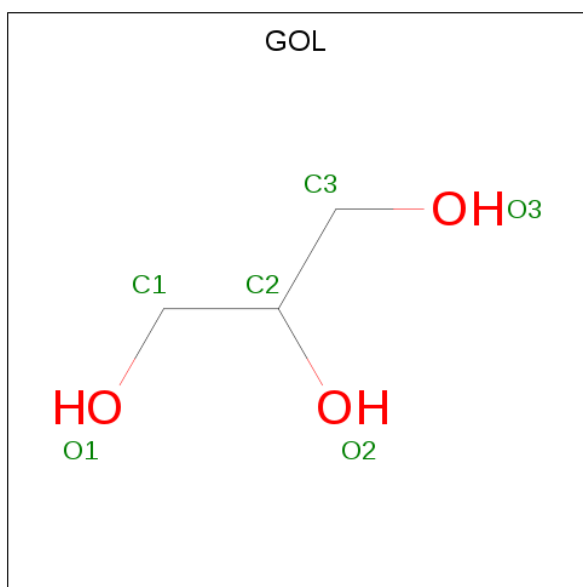
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	L	1	Total	C	N	O	P	S	0	0
			51	23	7	17	3	1		

- Molecule 3 is ACETIC ACID (three-letter code: ACY) (formula: C<sub>2</sub>H<sub>4</sub>O<sub>2</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		
3	B	1	Total	C	O	0	0
			4	2	2		
3	D	1	Total	C	O	0	0
			4	2	2		
3	F	1	Total	C	O	0	0
			4	2	2		
3	G	1	Total	C	O	0	0
			4	2	2		
3	I	1	Total	C	O	0	0
			4	2	2		
3	K	1	Total	C	O	0	0
			4	2	2		
3	K	1	Total	C	O	0	0
			4	2	2		

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	O	0	0
			6	3	3		
4	E	1	Total	C	O	0	0
			6	3	3		
4	F	1	Total	C	O	0	0
			6	3	3		
4	L	1	Total	C	O	0	0
			6	3	3		

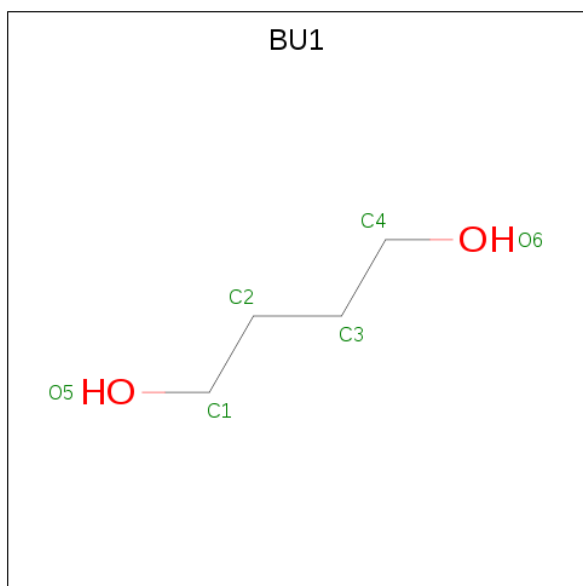
- Molecule 5 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	H	1	Total	Cl	0	0
			1	1		
5	B	1	Total	Cl	0	0
			1	1		
5	C	1	Total	Cl	0	0
			1	1		
5	J	1	Total	Cl	0	0
			1	1		
5	E	1	Total	Cl	0	0
			1	1		

- Molecule 6 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	C	1	Total	Mg	0	1
			2	2		

- Molecule 7 is 1,4-BUTANEDIOL (three-letter code: BU1) (formula: C<sub>4</sub>H<sub>10</sub>O<sub>2</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	G	1	Total	C	O	0	0
			6	4	2		
7	K	1	Total	C	O	0	0
			6	4	2		

- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	98	Total	O	0	0
			98	98		
8	B	88	Total	O	0	0
			88	88		
8	C	80	Total	O	0	0
			80	80		
8	D	66	Total	O	0	0
			66	66		
8	E	97	Total	O	0	0
			97	97		
8	F	93	Total	O	0	0
			93	93		
8	G	69	Total	O	0	0
			69	69		

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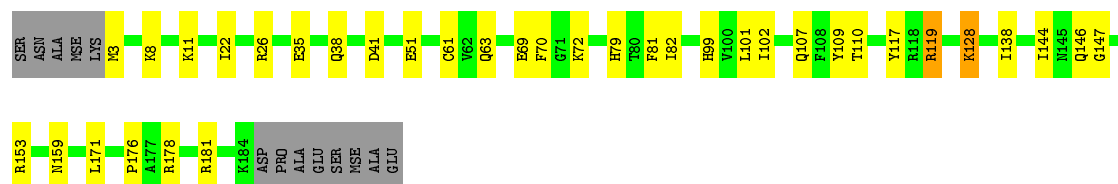
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	H	75	Total 75	O 75	0	0
8	I	80	Total 80	O 80	0	0
8	J	87	Total 87	O 87	0	0
8	K	92	Total 92	O 92	0	0
8	L	97	Total 97	O 97	0	0

### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

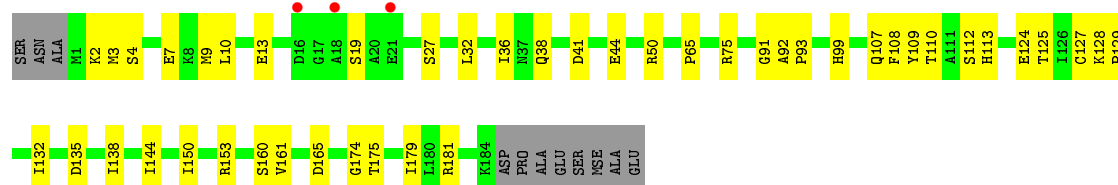
- Molecule 1: Hexapeptide-repeat containing-acetyltransferase

Chain A: 



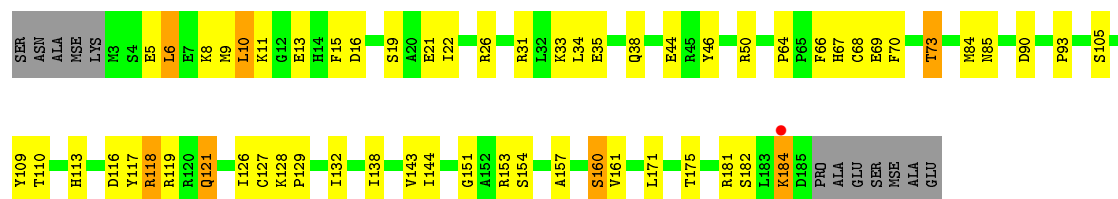
- Molecule 1: Hexapeptide-repeat containing-acetyltransferase

Chain B: 



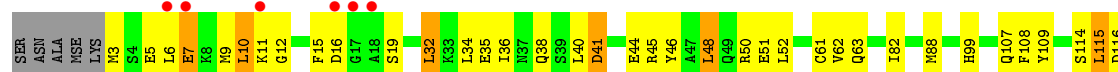
- Molecule 1: Hexapeptide-repeat containing-acetyltransferase

Chain C: 



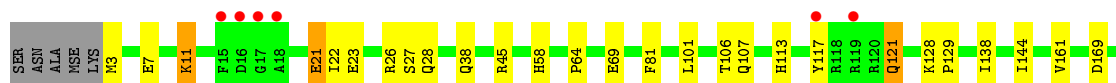
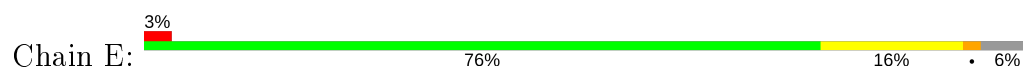
- Molecule 1: Hexapeptide-repeat containing-acetyltransferase

Chain D: 





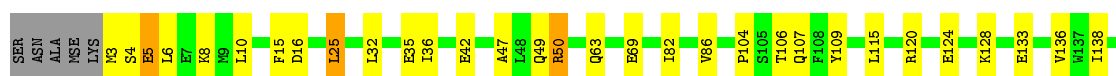
- Molecule 1: Hexapeptide-repeat containing-acetyltransferase



- Molecule 1: Hexapeptide-repeat containing-acetyltransferase



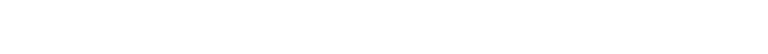
- Molecule 1: Hexapeptide-repeat containing-acetyltransferase

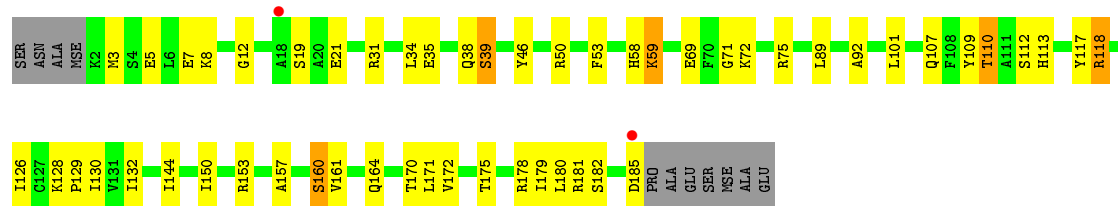


- Molecule 1: Hexapeptide-repeat containing-acetyltransferase

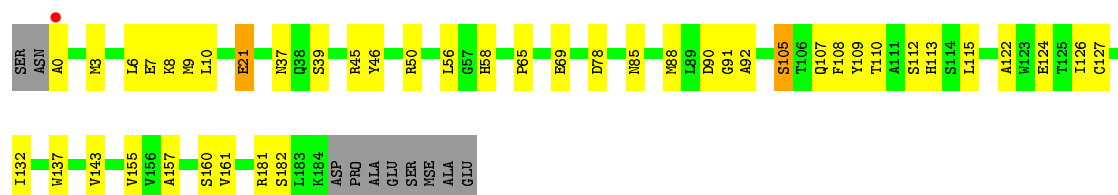
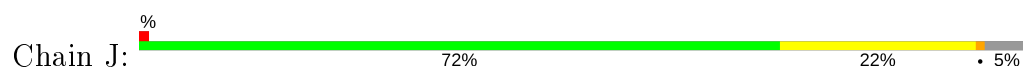


- Molecule 1: Hexapeptide-repeat containing-acetyltransferase

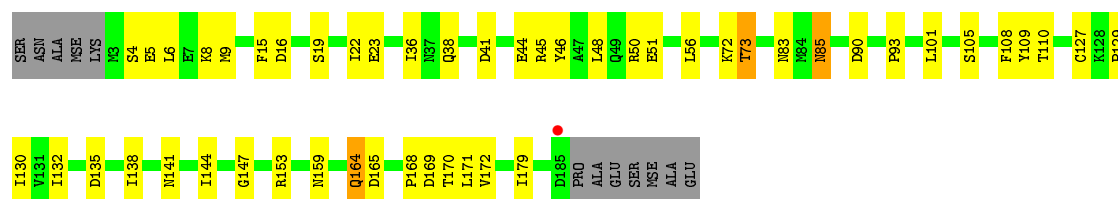




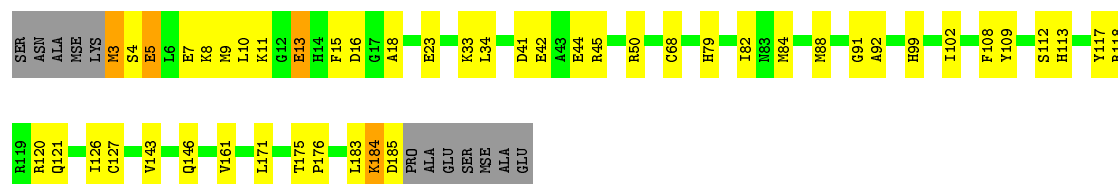
• Molecule 1: Hexapeptide-repeat containing-acetyltransferase



• Molecule 1: Hexapeptide-repeat containing-acetyltransferase



• Molecule 1: Hexapeptide-repeat containing-acetyltransferase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	77.32Å 135.80Å 120.53Å 90.00° 91.77° 90.00°	Depositor
Resolution (Å)	46.63 – 2.35 46.63 – 2.35	Depositor EDS
% Data completeness (in resolution range)	96.9 (46.63-2.35) 97.0 (46.63-2.35)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.18	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.16 (at 2.34Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.6.2_432)	Depositor
R, $R_{free}$	0.188 , 0.253 0.182 , 0.247	Depositor DCC
$R_{free}$ test set	5025 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	25.0	Xtriage
Anisotropy	0.767	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 57.7	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.056 for h,-k,-l	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	18822	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 46.23 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.1800e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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, MG, CL, BU1, ACO, ACY

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	0.61	0/1444	0.69	0/1950
1	B	0.64	0/1460	0.70	0/1968
1	C	0.60	0/1441	0.68	0/1947
1	D	0.58	0/1457	0.67	0/1970
1	E	0.66	0/1452	0.67	0/1961
1	F	0.59	0/1441	0.65	0/1947
1	G	0.61	0/1442	0.69	0/1948
1	H	0.61	0/1452	0.74	0/1961
1	I	0.64	0/1460	0.69	0/1969
1	J	0.63	0/1453	0.74	1/1958 (0.1%)
1	K	0.59	0/1441	0.67	0/1947
1	L	0.64	0/1452	0.70	0/1961
All	All	0.62	0/17395	0.69	1/23487 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	J	90	ASP	CB-CG-OD1	5.20	122.98	118.30

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1420	0	1416	32	0
1	B	1437	0	1441	35	0
1	C	1417	0	1407	57	0
1	D	1431	0	1417	74	0
1	E	1428	0	1420	44	0
1	F	1417	0	1408	37	0
1	G	1418	0	1409	33	0
1	H	1428	0	1420	44	0
1	I	1437	0	1433	42	0
1	J	1431	0	1434	33	0
1	K	1417	0	1408	42	0
1	L	1428	0	1420	43	0
2	A	102	0	68	17	0
2	C	51	0	34	5	0
2	D	51	0	34	2	0
2	E	51	0	34	4	0
2	F	51	0	34	11	0
2	H	51	0	34	2	0
2	I	102	0	68	8	0
2	J	51	0	34	6	0
2	L	102	0	68	6	1
3	A	8	0	6	1	0
3	B	4	0	3	0	0
3	D	4	0	3	0	0
3	F	4	0	3	0	0
3	G	4	0	3	1	0
3	I	4	0	3	0	0
3	K	8	0	6	2	0
4	B	6	0	8	1	0
4	E	6	0	8	4	0
4	F	6	0	8	1	0
4	L	6	0	8	3	0
5	B	1	0	0	0	0
5	C	1	0	0	0	0
5	E	1	0	0	0	0
5	H	1	0	0	0	0
5	J	1	0	0	0	0
6	C	2	0	0	0	0
7	G	6	0	10	2	0
7	K	6	0	10	0	0
8	A	98	0	0	4	0
8	B	88	0	0	9	0
8	C	80	0	0	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
8	D	66	0	0	5	0
8	E	97	0	0	8	0
8	F	93	0	0	5	0
8	G	69	0	0	3	0
8	H	75	0	0	4	0
8	I	80	0	0	3	1
8	J	87	0	0	1	0
8	K	92	0	0	2	0
8	L	97	0	0	2	0
All	All	18822	0	17520	507	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 14.

The worst 5 of 507 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:E:3:MSE:HE2	1:E:7:GLU:OE2	1.26	1.28
2:F:195:ACO:HH33	8:F:646:HOH:O	1.39	1.20
1:E:45:ARG:HE	4:E:196:GOL:H11	0.93	1.09
1:I:92:ALA:H	1:I:110:THR:HG21	1.09	1.08
1:H:45[A]:ARG:HG2	1:H:45[A]:ARG:HH11	1.01	1.07

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)
2:L:195:ACO:O5A	8:I:624:HOH:O[1_655]	2.16	0.04

## 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	181/195 (93%)	178 (98%)	3 (2%)	0	100	100
1	B	183/195 (94%)	182 (100%)	1 (0%)	0	100	100
1	C	181/195 (93%)	179 (99%)	2 (1%)	0	100	100
1	D	182/195 (93%)	178 (98%)	4 (2%)	0	100	100
1	E	182/195 (93%)	181 (100%)	1 (0%)	0	100	100
1	F	181/195 (93%)	180 (99%)	1 (1%)	0	100	100
1	G	181/195 (93%)	179 (99%)	2 (1%)	0	100	100
1	H	182/195 (93%)	182 (100%)	0	0	100	100
1	I	183/195 (94%)	182 (100%)	1 (0%)	0	100	100
1	J	183/195 (94%)	179 (98%)	4 (2%)	0	100	100
1	K	181/195 (93%)	179 (99%)	2 (1%)	0	100	100
1	L	182/195 (93%)	180 (99%)	2 (1%)	0	100	100
All	All	2182/2340 (93%)	2159 (99%)	23 (1%)	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	154/157 (98%)	151 (98%)	3 (2%)	57	68
1	B	156/157 (99%)	151 (97%)	5 (3%)	39	47
1	C	154/157 (98%)	143 (93%)	11 (7%)	14	15
1	D	155/157 (99%)	142 (92%)	13 (8%)	11	10
1	E	155/157 (99%)	149 (96%)	6 (4%)	32	40
1	F	154/157 (98%)	148 (96%)	6 (4%)	32	40
1	G	154/157 (98%)	144 (94%)	10 (6%)	17	18
1	H	155/157 (99%)	142 (92%)	13 (8%)	11	10
1	I	156/157 (99%)	142 (91%)	14 (9%)	9	8

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	J	155/157 (99%)	149 (96%)	6 (4%)	32	40
1	K	154/157 (98%)	148 (96%)	6 (4%)	32	40
1	L	155/157 (99%)	149 (96%)	6 (4%)	32	40
All	All	1857/1884 (99%)	1758 (95%)	99 (5%)	23	26

5 of 99 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	G	3	MSE
1	H	5	GLU
1	K	85	ASN
1	G	10	LEU
1	G	50	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 40 such sidechains are listed below:

Mol	Chain	Res	Type
1	F	145	ASN
1	G	145	ASN
1	K	164	GLN
1	G	67	HIS
1	G	146	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 carbohydrates in this entry.

## 5.6 Ligand geometry

Of 34 ligands modelled in this entry, 7 are monoatomic - leaving 27 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	$\# Z  > 2$
3	ACY	K	197	-	1,3,3	1.63	0	0,3,3	0.00	-
3	ACY	G	196	-	1,3,3	3.03	1 (100%)	0,3,3	0.00	-
4	GOL	L	196	-	5,5,5	0.36	0	5,5,5	1.16	1 (20%)
2	ACO	A	193	-	45,53,53	1.81	10 (22%)	56,79,79	1.59	7 (12%)
4	GOL	E	196	-	5,5,5	0.53	0	5,5,5	1.20	1 (20%)
3	ACY	D	197	-	1,3,3	0.82	0	0,3,3	0.00	-
3	ACY	F	196	-	1,3,3	0.81	0	0,3,3	0.00	-
7	BU1	G	197	-	5,5,5	0.33	0	4,4,4	0.60	0
4	GOL	B	196	-	5,5,5	0.37	0	5,5,5	0.28	0
2	ACO	C	195	-	45,53,53	1.81	11 (24%)	56,79,79	1.44	6 (10%)
4	GOL	F	197	-	5,5,5	0.38	0	5,5,5	0.27	0
3	ACY	K	198	-	1,3,3	1.21	0	0,3,3	0.00	-
3	ACY	A	196	-	1,3,3	1.48	0	0,3,3	0.00	-
2	ACO	D	195	-	45,53,53	1.77	6 (13%)	56,79,79	1.51	8 (14%)
7	BU1	K	196	-	5,5,5	0.35	0	4,4,4	0.36	0
3	ACY	A	197	-	1,3,3	1.36	0	0,3,3	0.00	-
2	ACO	I	195	-	45,53,53	1.80	11 (24%)	56,79,79	1.47	11 (19%)
2	ACO	H	195	-	45,53,53	1.95	11 (24%)	56,79,79	1.53	9 (16%)
2	ACO	E	195	-	45,53,53	1.82	11 (24%)	56,79,79	1.21	5 (8%)
2	ACO	L	195	-	45,53,53	1.87	10 (22%)	56,79,79	1.37	10 (17%)
2	ACO	J	195	-	45,53,53	1.92	12 (26%)	56,79,79	1.64	10 (17%)
2	ACO	I	193	-	45,53,53	1.79	9 (20%)	56,79,79	1.38	5 (8%)
3	ACY	B	198	-	1,3,3	1.40	0	0,3,3	0.00	-
2	ACO	F	195	-	45,53,53	1.86	11 (24%)	56,79,79	1.66	9 (16%)
2	ACO	L	193	-	45,53,53	1.88	12 (26%)	56,79,79	1.43	9 (16%)
2	ACO	A	195	-	45,53,53	1.88	10 (22%)	56,79,79	1.45	7 (12%)
3	ACY	I	196	-	1,3,3	1.18	0	0,3,3	0.00	-

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	ACO	J	195	-	-	20/47/67/67	0/3/3/3
2	ACO	I	193	-	-	17/47/67/67	0/3/3/3
2	ACO	D	195	-	-	16/47/67/67	0/3/3/3
4	GOL	E	196	-	-	2/4/4/4	-
7	BU1	K	196	-	-	1/3/3/3	-
2	ACO	A	193	-	-	18/47/67/67	0/3/3/3
2	ACO	F	195	-	-	16/47/67/67	0/3/3/3
7	BU1	G	197	-	-	0/3/3/3	-
2	ACO	L	193	-	-	18/47/67/67	0/3/3/3
2	ACO	I	195	-	-	12/47/67/67	0/3/3/3
4	GOL	B	196	-	-	2/4/4/4	-
4	GOL	L	196	-	-	2/4/4/4	-
2	ACO	C	195	-	-	20/47/67/67	0/3/3/3
2	ACO	A	195	-	-	22/47/67/67	0/3/3/3
2	ACO	H	195	-	-	18/47/67/67	0/3/3/3
4	GOL	F	197	-	-	0/4/4/4	-
2	ACO	E	195	-	-	21/47/67/67	0/3/3/3
2	ACO	L	195	-	-	13/47/67/67	0/3/3/3

The worst 5 of 125 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	H	195	ACO	C9P-N8P	7.82	1.50	1.33
2	I	193	ACO	C9P-N8P	7.37	1.49	1.33
2	D	195	ACO	C9P-N8P	7.36	1.49	1.33
2	A	195	ACO	C9P-N8P	7.14	1.49	1.33
2	I	195	ACO	C9P-N8P	7.12	1.49	1.33

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

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	193	ACO	O6A-CCP-CBP	6.60	121.15	110.55
2	F	195	ACO	O6A-CCP-CBP	6.06	120.29	110.55
2	J	195	ACO	O6A-CCP-CBP	5.50	119.38	110.55
2	C	195	ACO	N3A-C2A-N1A	-4.83	121.14	128.68

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	L	193	ACO	N3A-C2A-N1A	-4.65	121.41	128.68

There are no chirality outliers.

5 of 218 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	193	ACO	C3B-O3B-P3B-O7A
2	A	193	ACO	C3B-C4B-C5B-O5B
2	A	193	ACO	O4B-C4B-C5B-O5B
2	A	193	ACO	C5B-O5B-P1A-O1A
2	A	193	ACO	C5B-O5B-P1A-O2A

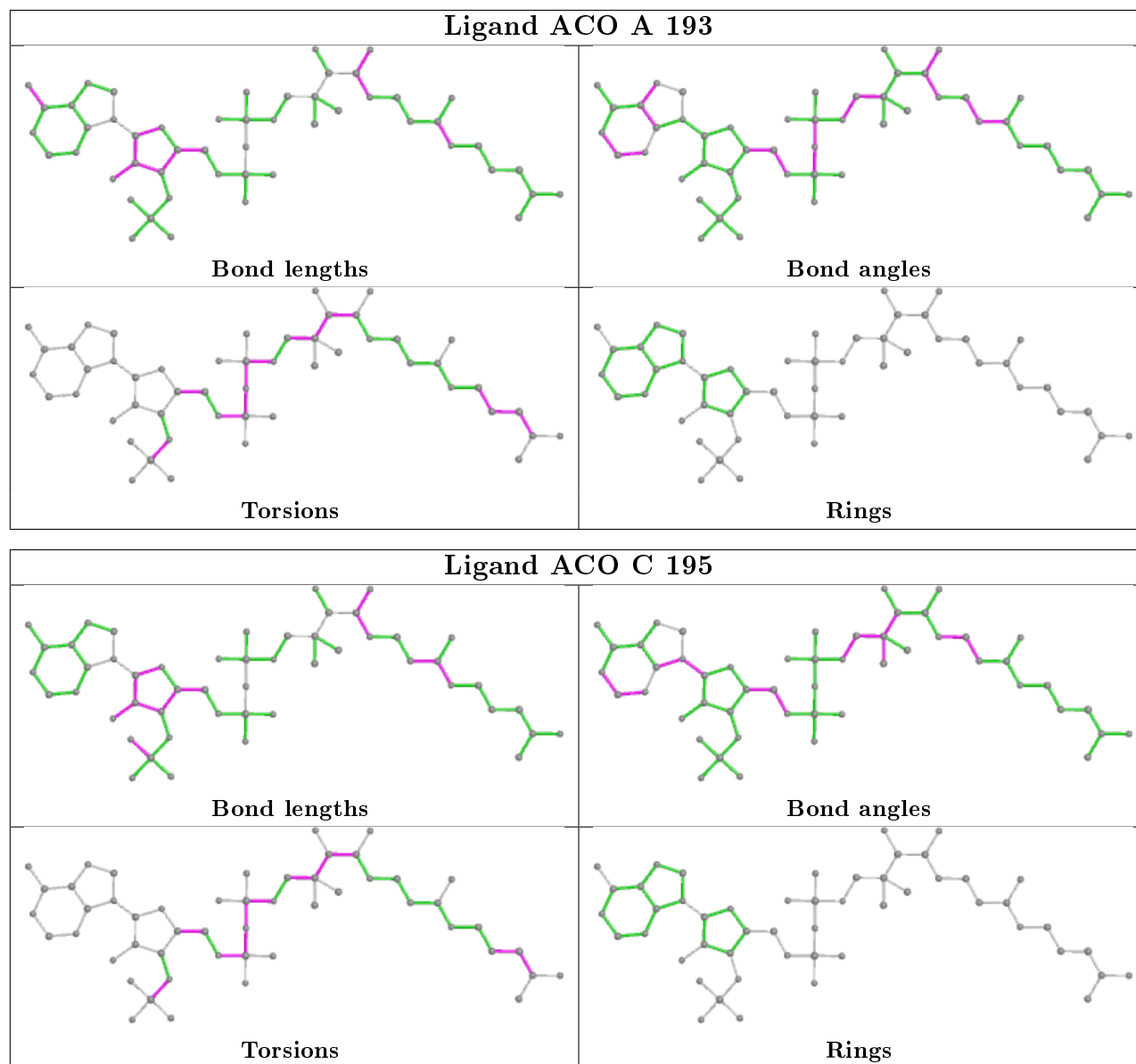
There are no ring outliers.

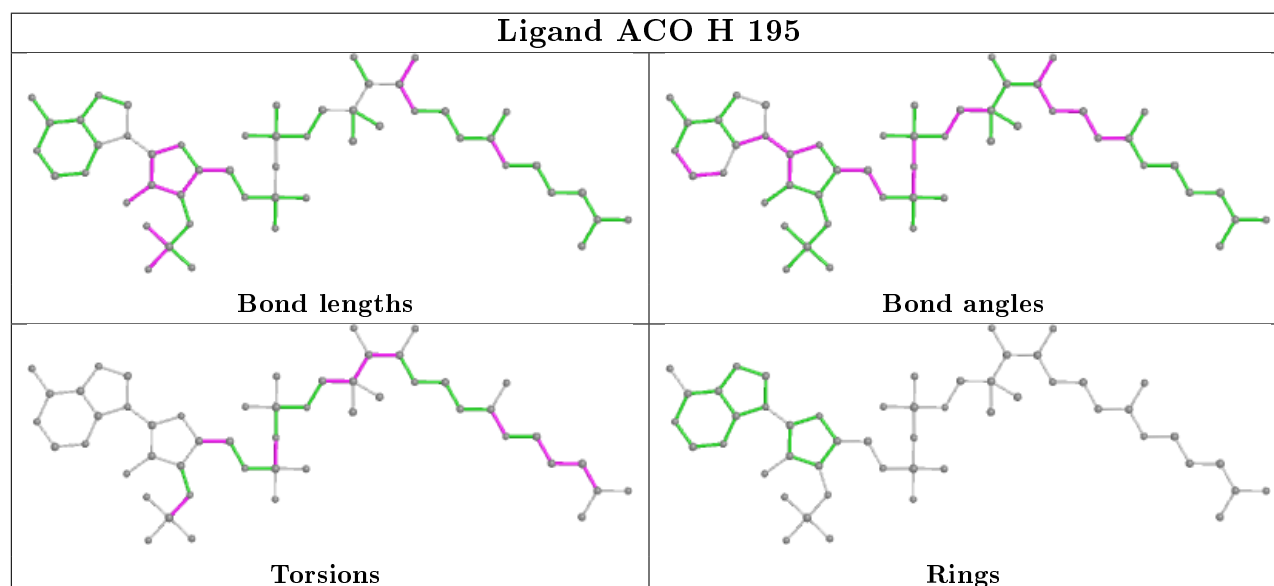
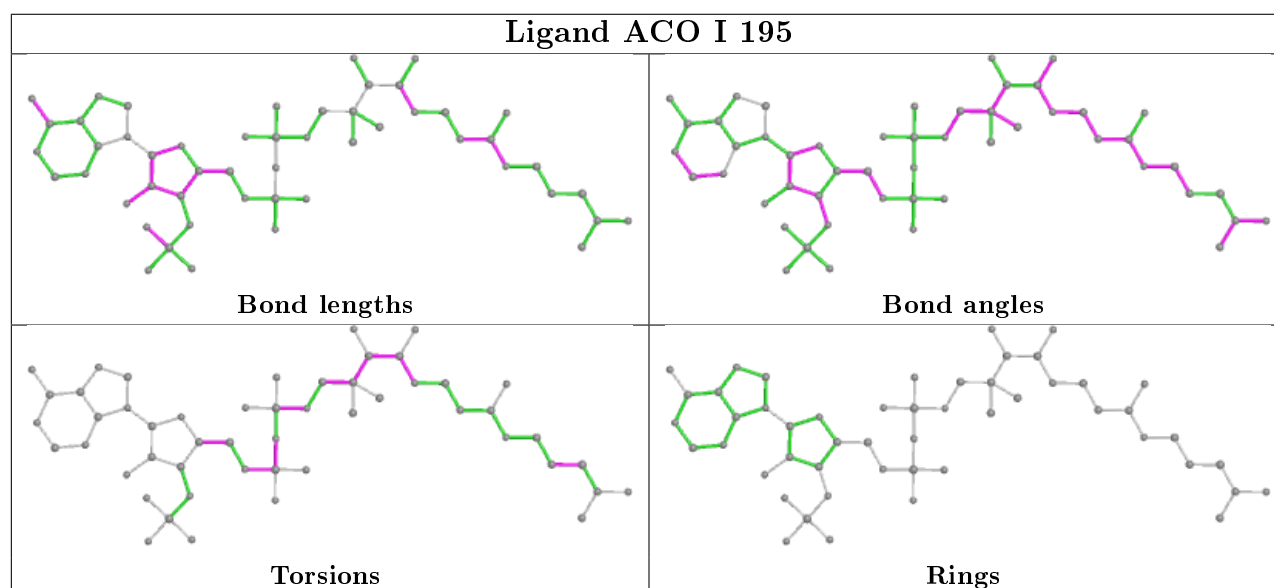
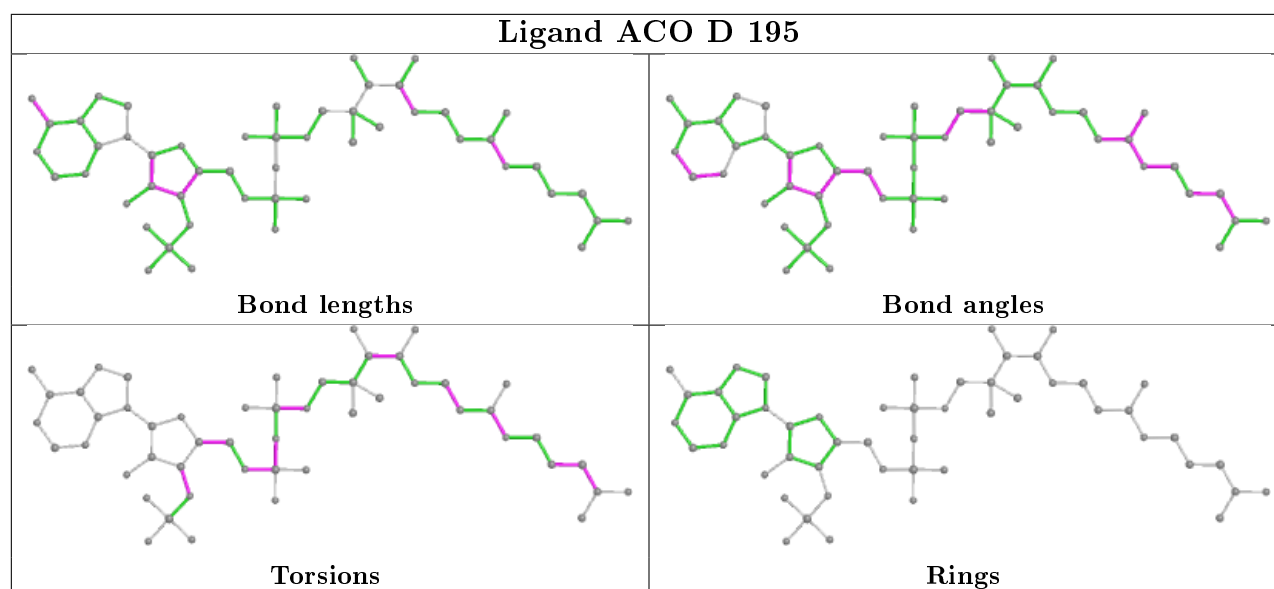
20 monomers are involved in 77 short contacts:

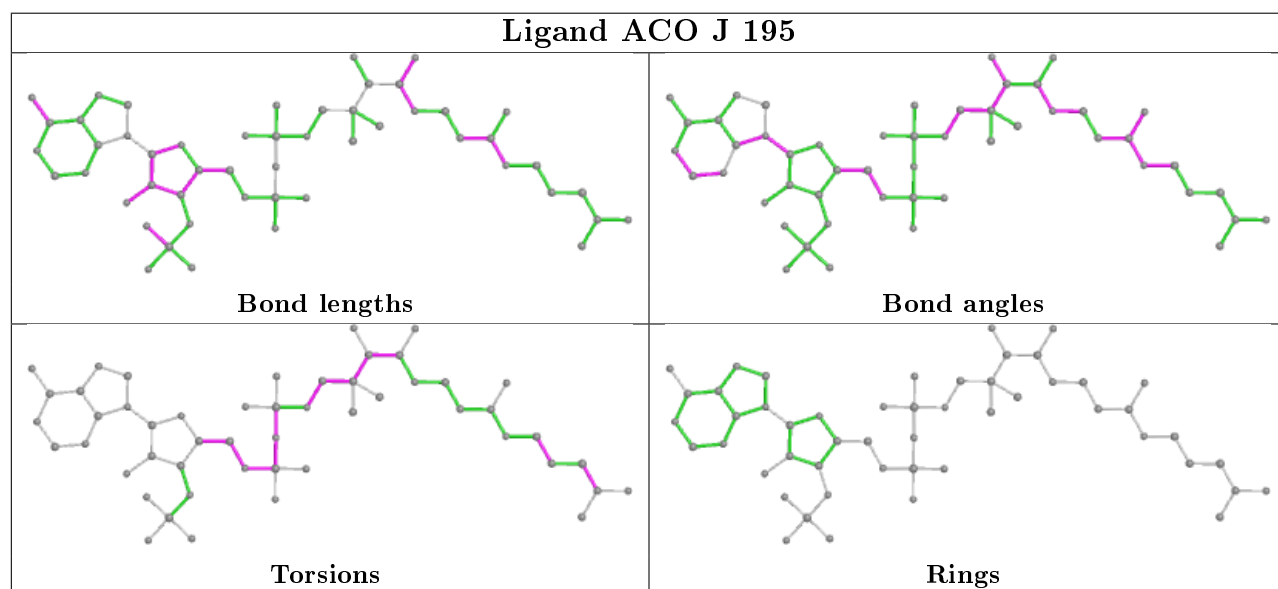
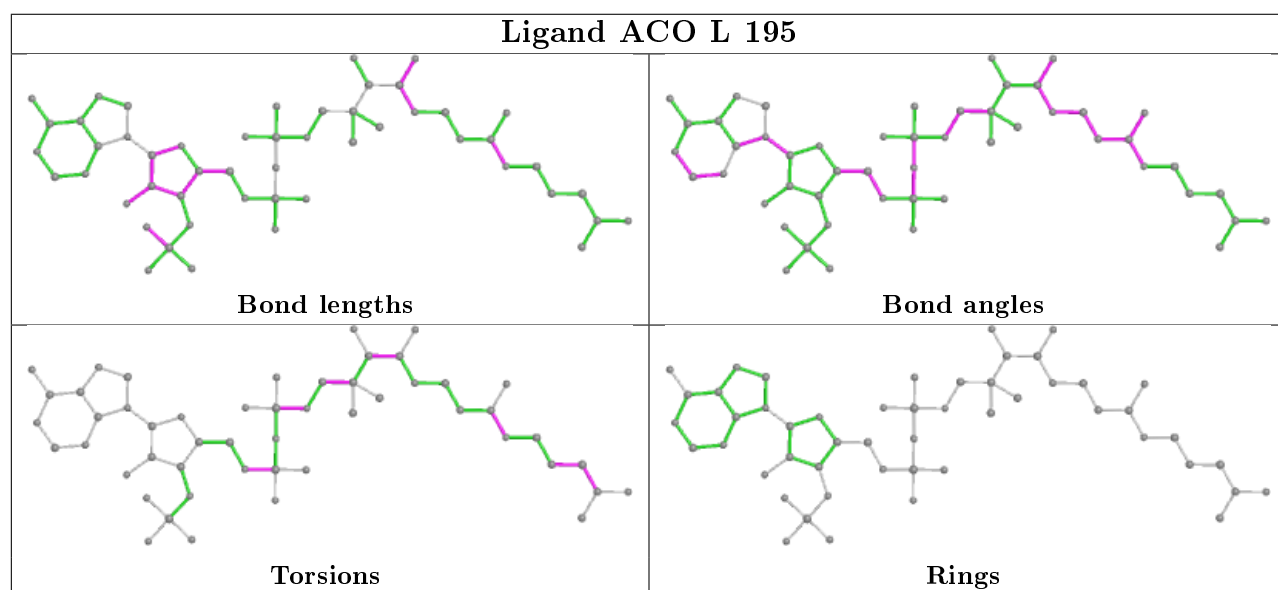
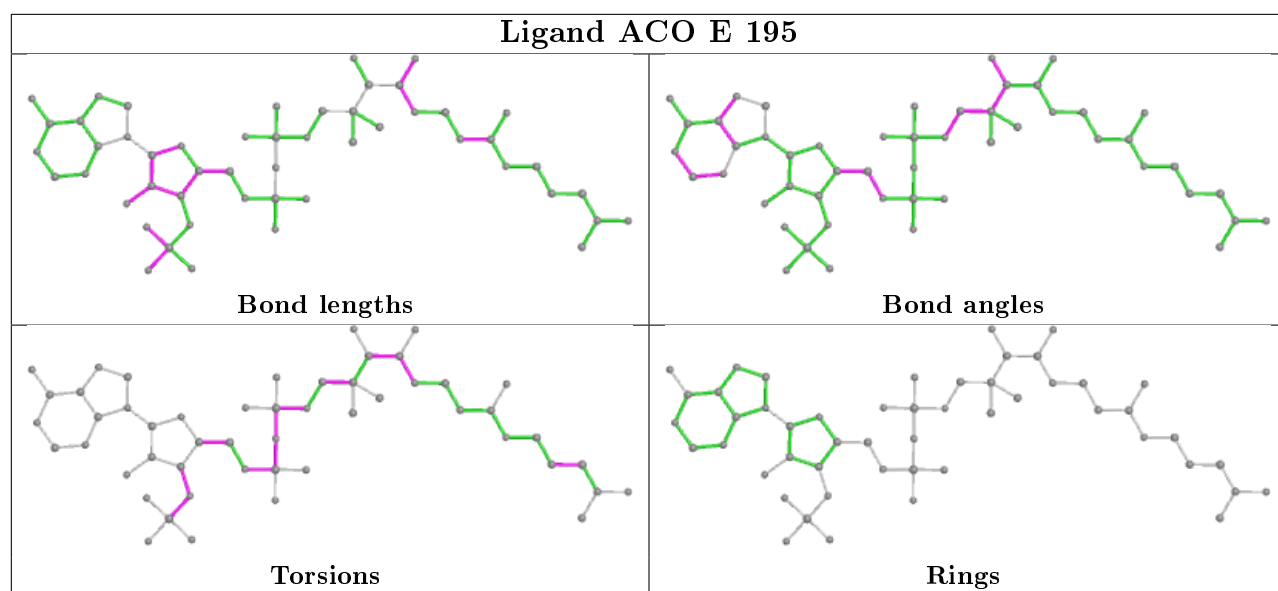
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	G	196	ACY	1	0
4	L	196	GOL	3	0
2	A	193	ACO	3	0
4	E	196	GOL	4	0
7	G	197	BU1	2	0
4	B	196	GOL	1	0
2	C	195	ACO	5	0
4	F	197	GOL	1	0
3	K	198	ACY	2	0
3	A	196	ACY	1	0
2	D	195	ACO	2	0
2	I	195	ACO	7	0
2	H	195	ACO	2	0
2	E	195	ACO	4	0
2	L	195	ACO	5	1
2	J	195	ACO	6	0
2	I	193	ACO	1	0
2	F	195	ACO	11	0
2	L	193	ACO	1	0
2	A	195	ACO	14	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.

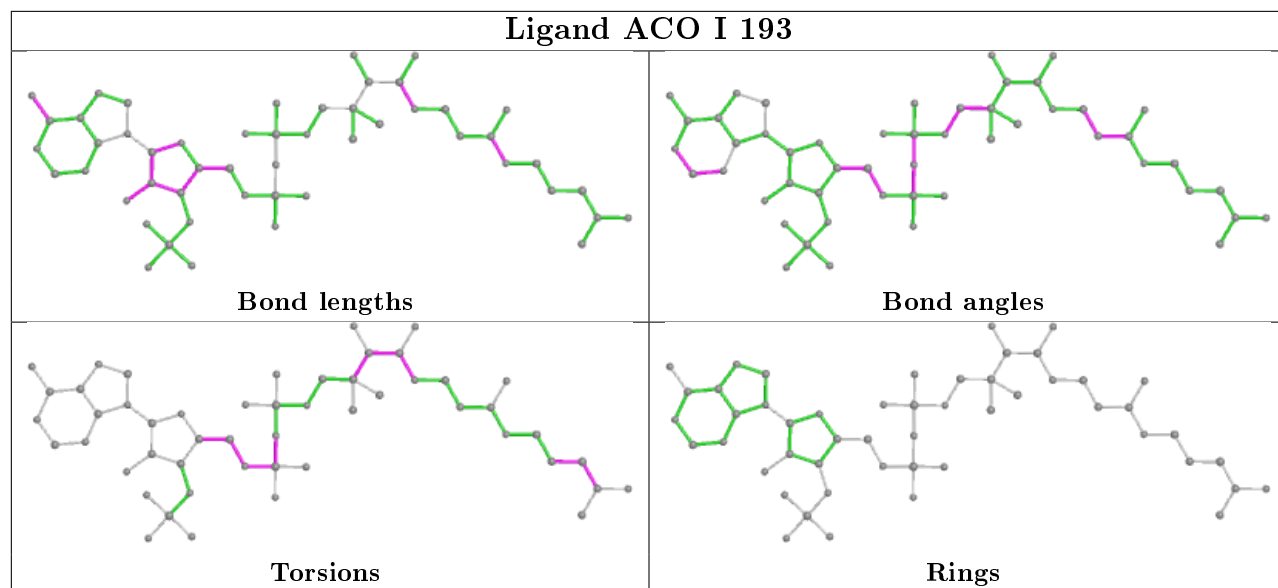
Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD 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.



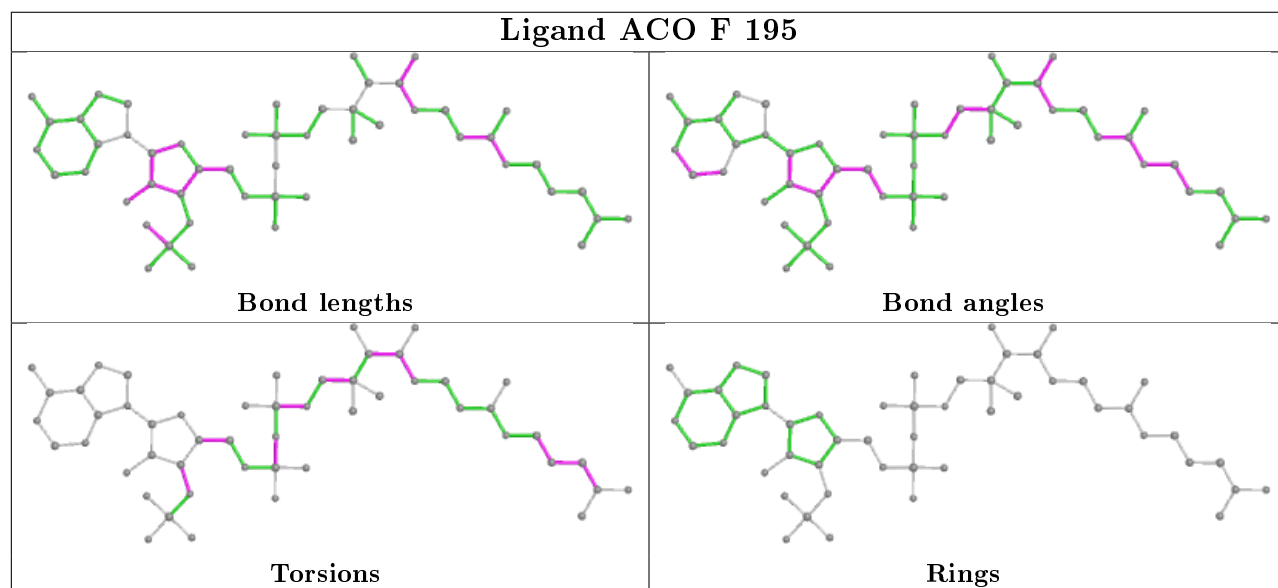




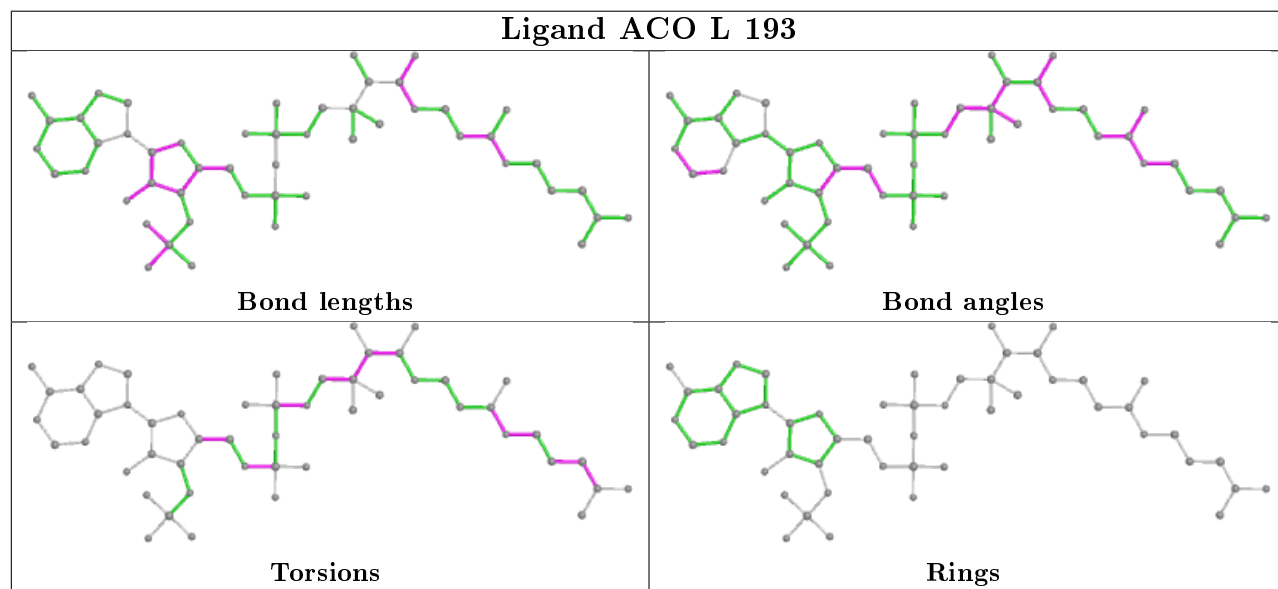
## Ligand ACO I 193

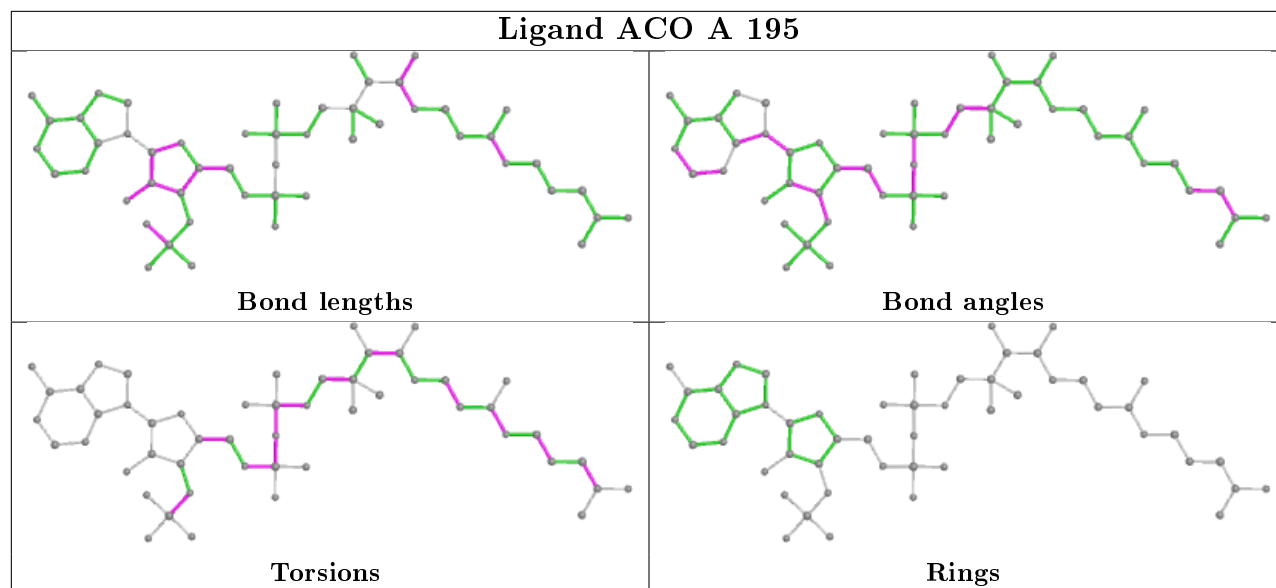


## Ligand ACO F 195



## Ligand ACO L 193





## 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	178/195 (91%)	-0.32	0 100 100	10, 22, 40, 62	0
1	B	179/195 (91%)	-0.14	3 (1%) 70 78	10, 23, 47, 88	0
1	C	179/195 (91%)	-0.18	1 (0%) 89 93	11, 26, 55, 87	0
1	D	179/195 (91%)	-0.19	7 (3%) 39 52	16, 28, 59, 82	0
1	E	179/195 (91%)	-0.05	6 (3%) 45 57	11, 22, 48, 69	0
1	F	179/195 (91%)	-0.10	7 (3%) 39 52	15, 27, 53, 84	0
1	G	178/195 (91%)	-0.32	1 (0%) 89 93	13, 25, 45, 64	0
1	H	179/195 (91%)	-0.08	6 (3%) 45 57	13, 28, 61, 86	0
1	I	180/195 (92%)	-0.14	2 (1%) 80 87	15, 25, 54, 74	0
1	J	180/195 (92%)	-0.29	1 (0%) 89 93	11, 24, 48, 79	0
1	K	179/195 (91%)	-0.24	1 (0%) 89 93	10, 24, 49, 75	0
1	L	179/195 (91%)	-0.21	0 100 100	9, 20, 47, 65	0
All	All	2148/2340 (91%)	-0.19	35 (1%) 72 80	9, 24, 51, 88	0

The worst 5 of 35 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	K	185	ASP	4.0
1	D	6	LEU	3.7
1	H	18	ALA	3.7
1	D	185	ASP	3.4
1	I	18	ALA	3.4

### 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

## 6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 6.4 Ligands ⓘ

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, 95<sup>th</sup> 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.

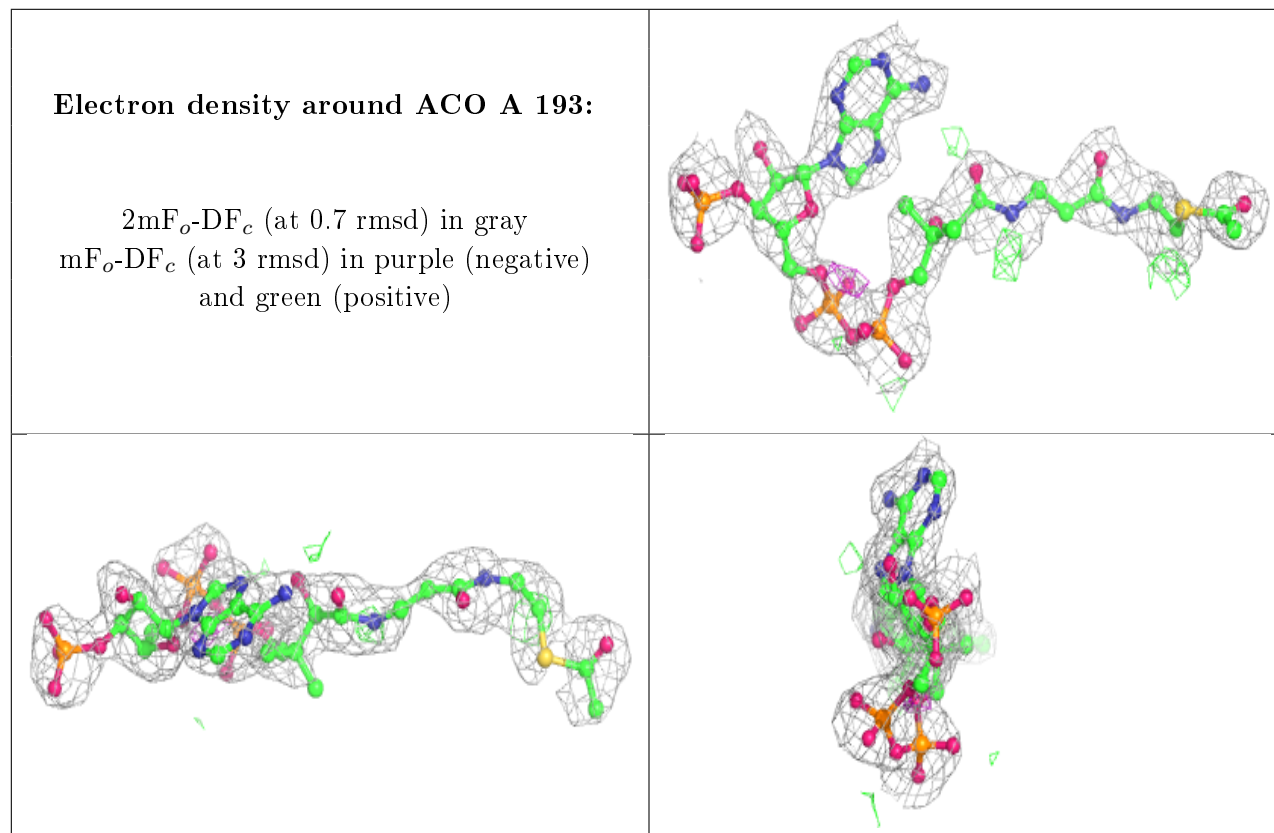
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
3	ACY	I	196	4/4	0.64	0.37	74,76,76,76	0
4	GOL	B	196	6/6	0.71	0.35	54,59,62,66	0
3	ACY	K	197	4/4	0.72	0.21	50,54,57,60	0
5	CL	B	197	1/1	0.79	0.16	73,73,73,73	0
3	ACY	F	196	4/4	0.81	0.21	54,59,59,59	0
3	ACY	B	198	4/4	0.81	0.30	48,64,67,68	0
6	MG	C	197[A]	1/1	0.81	0.23	38,38,38,38	1
6	MG	C	197[B]	1/1	0.81	0.23	35,35,35,35	1
4	GOL	F	197	6/6	0.83	0.25	50,57,61,64	0
7	BU1	G	197	6/6	0.83	0.20	46,50,52,52	0
2	ACO	A	193	51/51	0.83	0.19	26,59,88,114	0
3	ACY	D	197	4/4	0.83	0.22	43,45,51,54	0
2	ACO	H	195	51/51	0.85	0.17	25,51,128,131	0
2	ACO	I	195	51/51	0.85	0.18	24,67,88,93	0
2	ACO	A	195	51/51	0.87	0.18	23,60,89,152	0
3	ACY	A	196	4/4	0.87	0.21	35,36,38,42	0
2	ACO	I	193	51/51	0.88	0.18	22,64,86,104	0
3	ACY	K	198	4/4	0.88	0.19	44,49,49,50	0
2	ACO	F	195	51/51	0.88	0.16	23,52,125,132	0
2	ACO	C	195	51/51	0.89	0.18	23,40,64,70	0
4	GOL	E	196	6/6	0.90	0.18	28,30,34,37	0
5	CL	J	196	1/1	0.90	0.17	58,58,58,58	0
2	ACO	L	195	51/51	0.90	0.17	27,44,80,88	0
2	ACO	J	195	51/51	0.90	0.16	28,49,68,77	0
7	BU1	K	196	6/6	0.90	0.13	43,45,47,49	0
5	CL	C	196	1/1	0.91	0.08	61,61,61,61	0
2	ACO	D	195	51/51	0.91	0.15	5,48,85,156	0
3	ACY	G	196	4/4	0.92	0.14	24,29,33,33	0
4	GOL	L	196	6/6	0.93	0.17	21,34,35,37	0
5	CL	H	196	1/1	0.93	0.24	70,70,70,70	0
2	ACO	L	193	51/51	0.94	0.12	18,34,53,62	0
2	ACO	E	195	51/51	0.94	0.13	19,36,45,71	0

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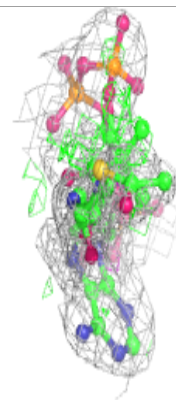
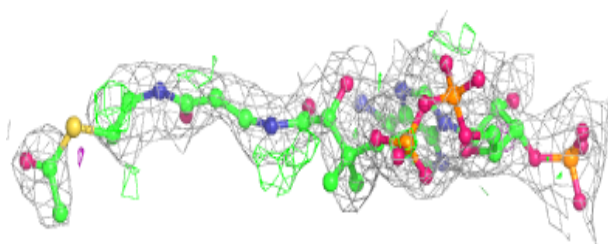
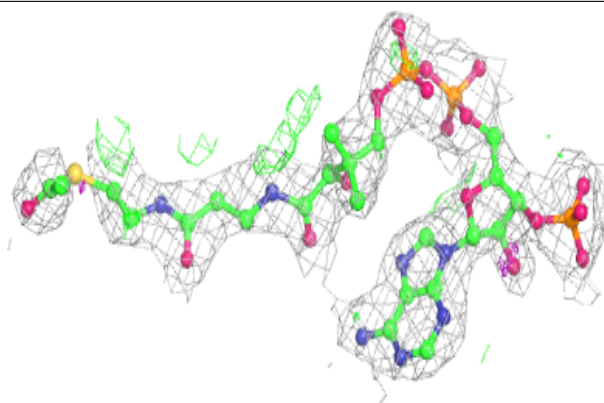
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
3	ACY	A	197	4/4	0.96	0.12	42,44,46,48	0
5	CL	E	193	1/1	0.96	0.12	51,51,51,51	0

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.

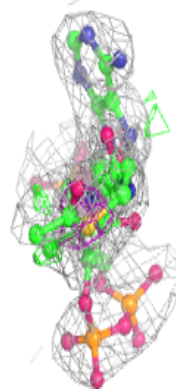
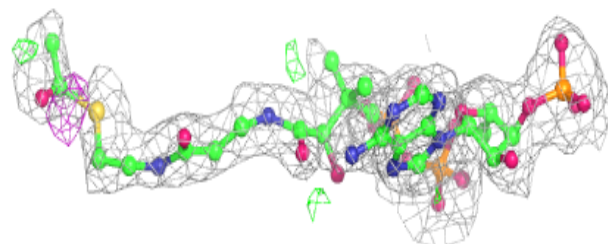
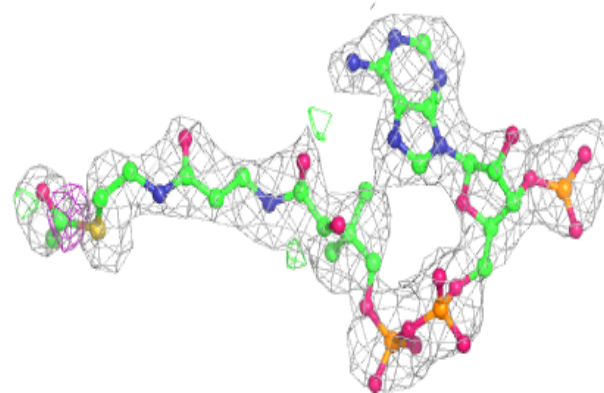


**Electron density around ACO H 195:**

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

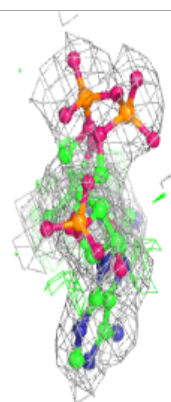
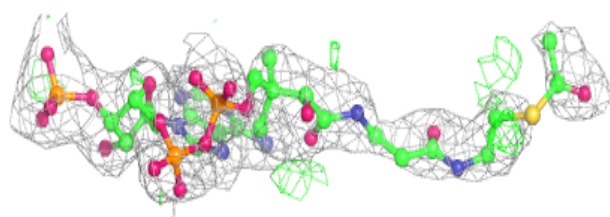
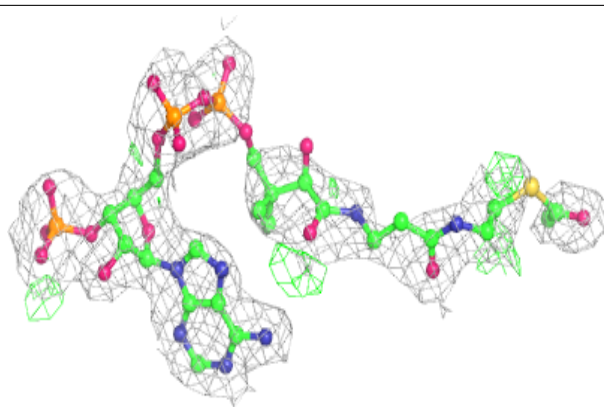
**Electron density around ACO I 195:**

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

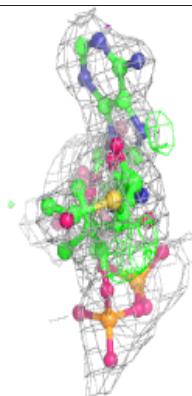
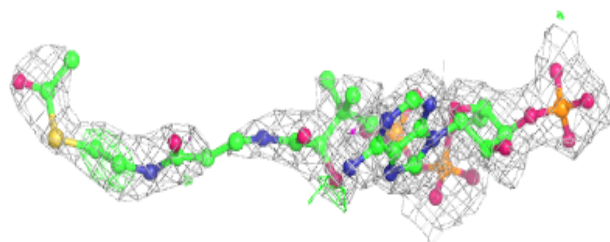
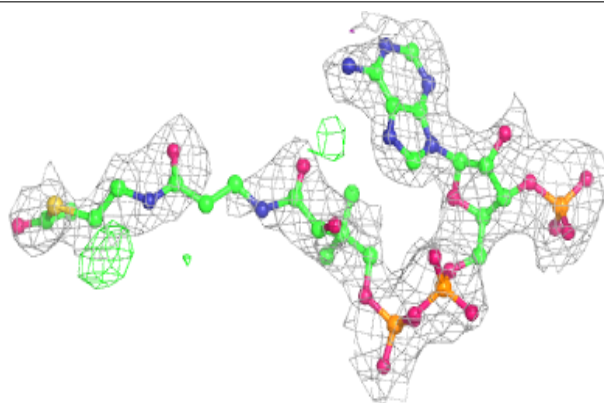


**Electron density around ACO A 195:**

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

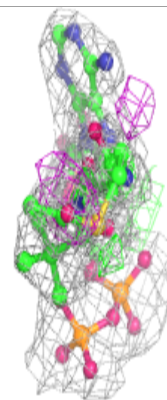
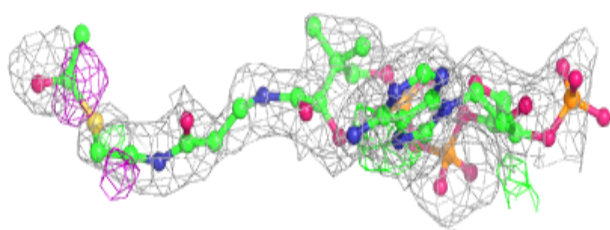
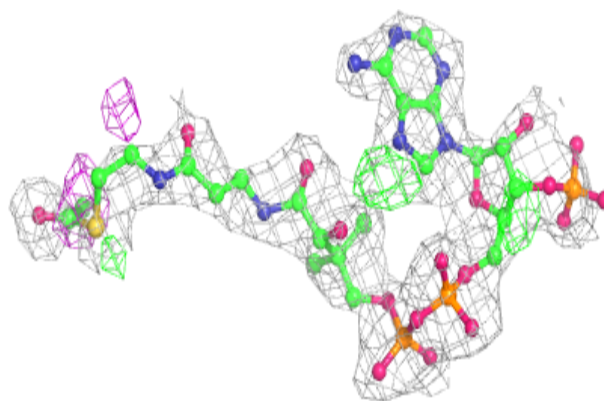
**Electron density around ACO I 193:**

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

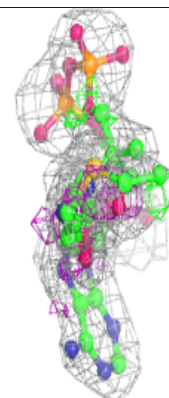
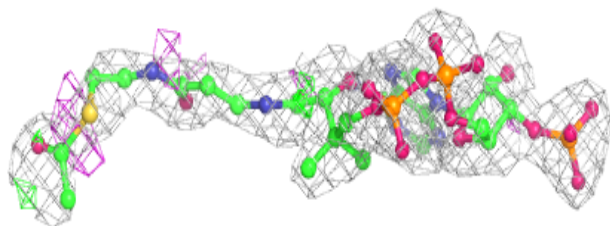
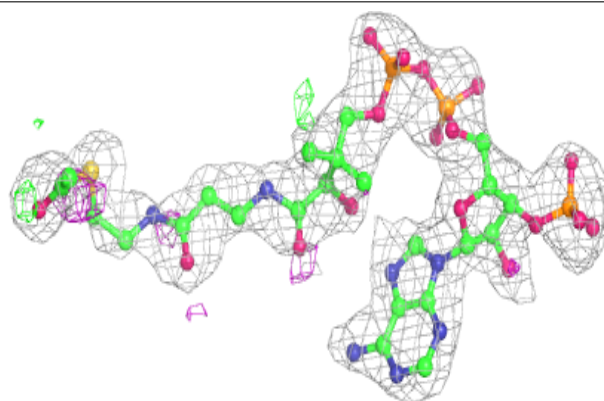


**Electron density around ACO F 195:**

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

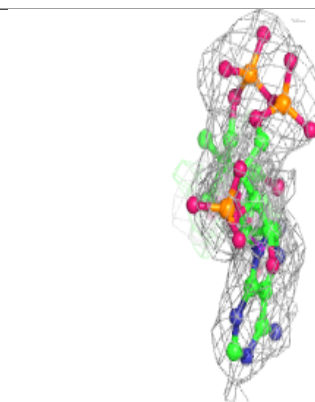
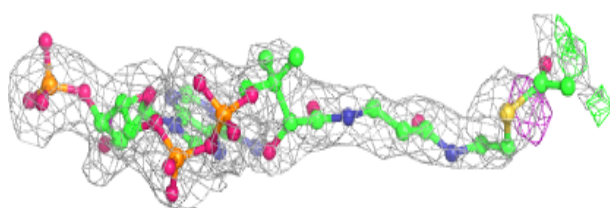
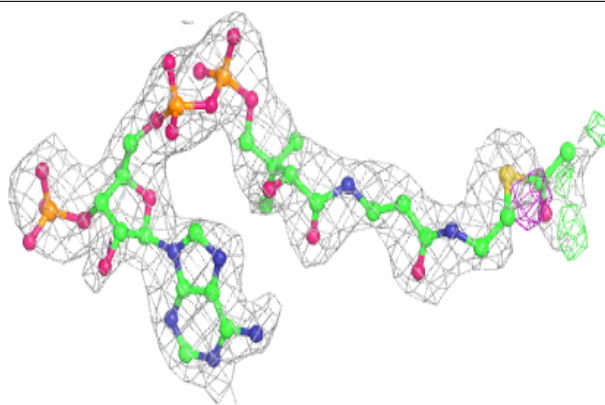
**Electron density around ACO C 195:**

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

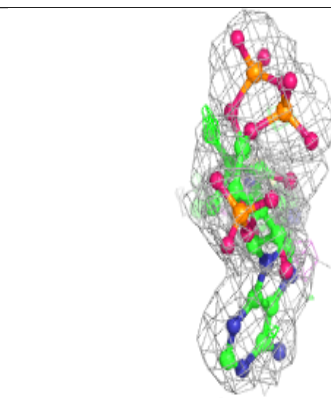
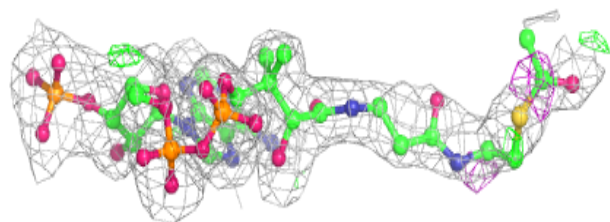
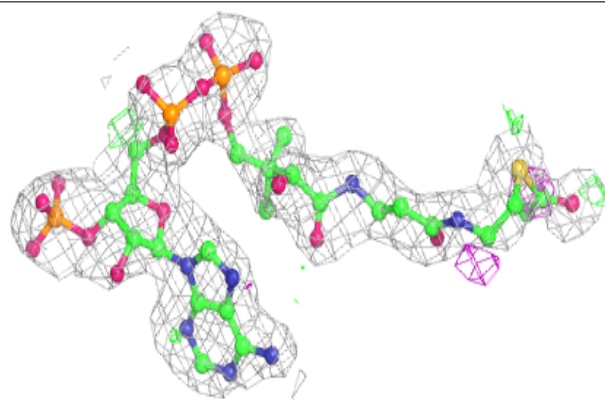


**Electron density around ACO L 195:**

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

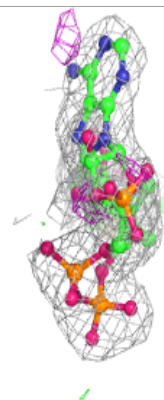
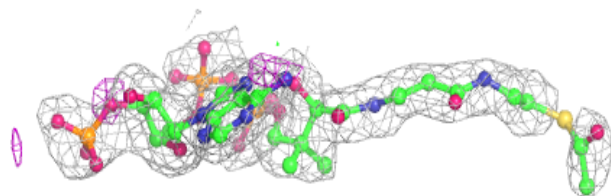
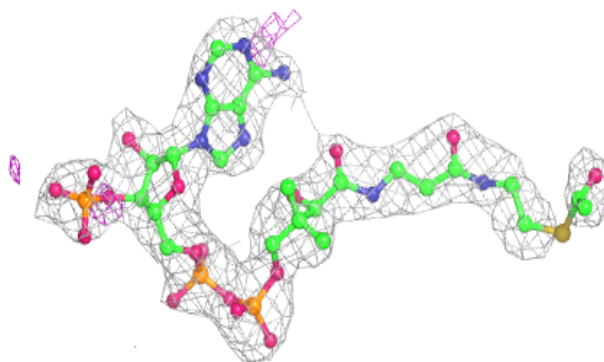
**Electron density around ACO J 195:**

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

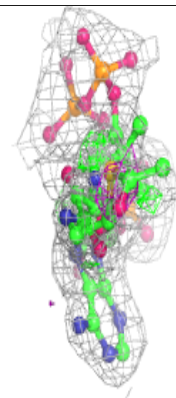
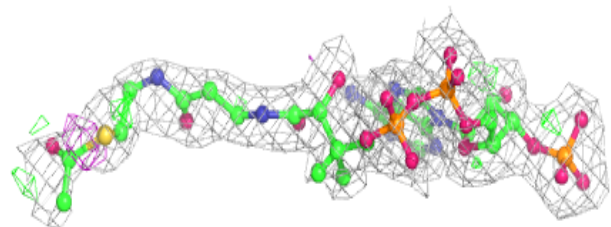
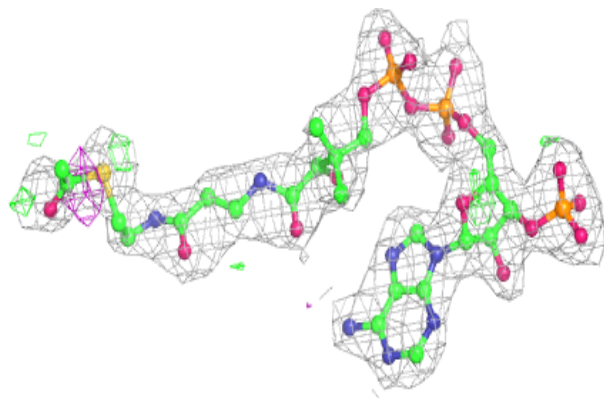


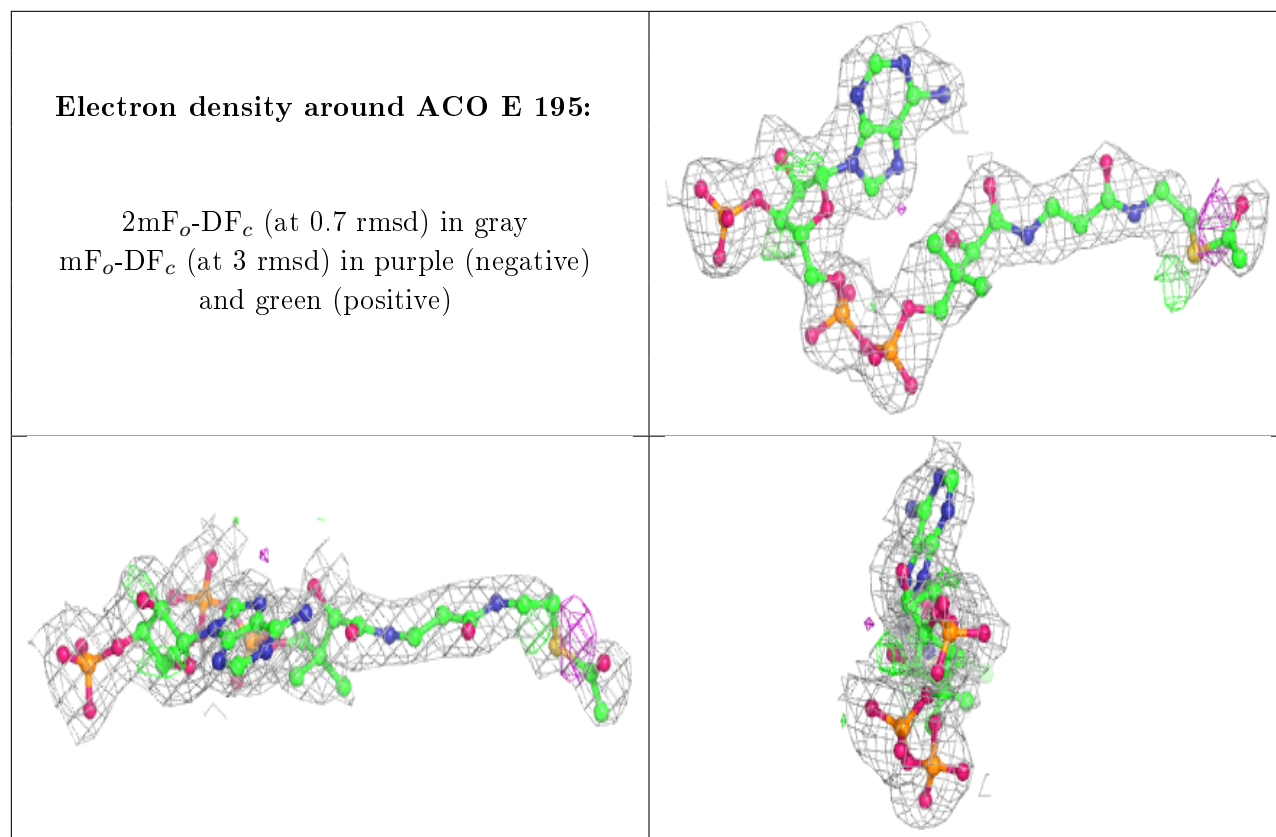
**Electron density around ACO D 195:**

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

**Electron density around ACO L 193:**

$2mF_o-DF_c$  (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.