



# Full wwPDB X-ray Structure Validation Report ⓘ

Dec 6, 2021 – 01:02 pm GMT

PDB ID : 7PCW  
Title : X-RAY STRUCTURE OF THE HALOALKANE DEHALOGENASE  
HALOTAG7-M175W LABELED WITH A CHLOROALKANE-TETRAM  
ETHYLRHODAMINE FLUOROPHORE SUBSTRATE  
Authors : Tarnawski, M.; Frei, M.; Hiblot, J.; Johnsson, K.  
Deposited on : 2021-08-04  
Resolution : 2.30 Å(reported)

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

We welcome your comments at [validation@mail.wwpdb.org](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 : 1.8.4 (270009), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.23.2  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0267  
CCP4 : 7.1.010 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.23.2

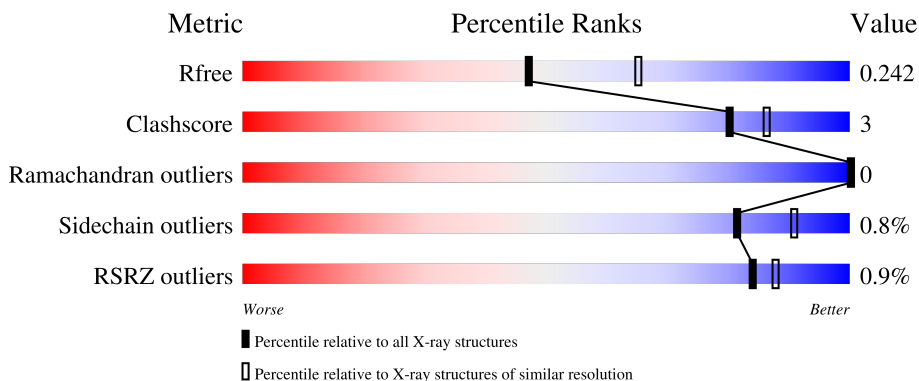
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

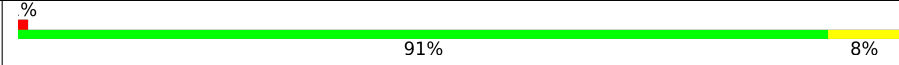
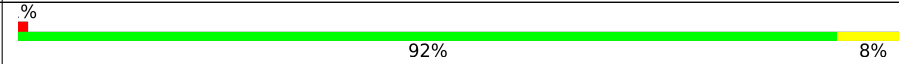
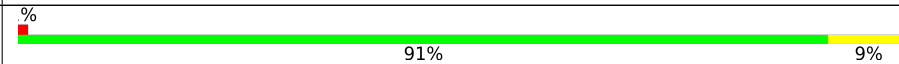
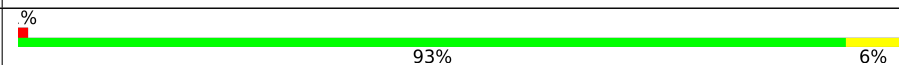
The reported resolution of this entry is 2.30 Å.

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	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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  $\geq 3$ , 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	293	 91% 8%
1	B	293	 92% 8%
1	C	293	 91% 9%
1	D	293	 93% 6%

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 9711 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 Haloalkane dehalogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	292	2347	1528	395	416	8	0	0	0
1	B	293	2356	1534	396	418	8	0	0	0
1	C	293	2356	1534	396	418	8	0	0	0
1	D	292	2347	1528	395	416	8	0	0	0

There are 96 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	3	GLY	-	expression tag	UNP P0A3G3
A	47	VAL	LEU	engineered mutation	UNP P0A3G3
A	58	THR	SER	engineered mutation	UNP P0A3G3
A	78	GLY	ASP	engineered mutation	UNP P0A3G3
A	87	PHE	TYR	engineered mutation	UNP P0A3G3
A	88	MET	LEU	engineered mutation	UNP P0A3G3
A	128	PHE	CYS	engineered mutation	UNP P0A3G3
A	155	THR	ALA	engineered mutation	UNP P0A3G3
A	160	LYS	GLU	engineered mutation	UNP P0A3G3
A	167	VAL	ALA	engineered mutation	UNP P0A3G3
A	172	THR	ALA	engineered mutation	UNP P0A3G3
A	175	TRP	LYS	engineered mutation	UNP P0A3G3
A	176	GLY	CYS	engineered mutation	UNP P0A3G3
A	195	ASN	LYS	engineered mutation	UNP P0A3G3
A	224	GLU	ALA	engineered mutation	UNP P0A3G3
A	227	ASP	ASN	engineered mutation	UNP P0A3G3
A	257	LYS	GLU	engineered mutation	UNP P0A3G3
A	264	ALA	THR	engineered mutation	UNP P0A3G3
A	272	ASN	HIS	engineered mutation	UNP P0A3G3
A	273	LEU	TYR	engineered mutation	UNP P0A3G3
A	291	SER	PRO	engineered mutation	UNP P0A3G3

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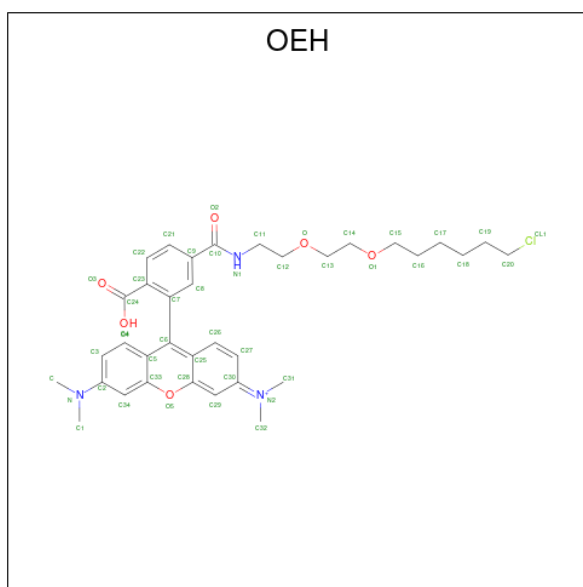
Chain	Residue	Modelled	Actual	Comment	Reference
A	292	THR	ALA	engineered mutation	UNP P0A3G3
A	294	GLU	-	expression tag	UNP P0A3G3
A	295	ILE	-	expression tag	UNP P0A3G3
B	3	GLY	-	expression tag	UNP P0A3G3
B	47	VAL	LEU	engineered mutation	UNP P0A3G3
B	58	THR	SER	engineered mutation	UNP P0A3G3
B	78	GLY	ASP	engineered mutation	UNP P0A3G3
B	87	PHE	TYR	engineered mutation	UNP P0A3G3
B	88	MET	LEU	engineered mutation	UNP P0A3G3
B	128	PHE	CYS	engineered mutation	UNP P0A3G3
B	155	THR	ALA	engineered mutation	UNP P0A3G3
B	160	LYS	GLU	engineered mutation	UNP P0A3G3
B	167	VAL	ALA	engineered mutation	UNP P0A3G3
B	172	THR	ALA	engineered mutation	UNP P0A3G3
B	175	TRP	LYS	engineered mutation	UNP P0A3G3
B	176	GLY	CYS	engineered mutation	UNP P0A3G3
B	195	ASN	LYS	engineered mutation	UNP P0A3G3
B	224	GLU	ALA	engineered mutation	UNP P0A3G3
B	227	ASP	ASN	engineered mutation	UNP P0A3G3
B	257	LYS	GLU	engineered mutation	UNP P0A3G3
B	264	ALA	THR	engineered mutation	UNP P0A3G3
B	272	ASN	HIS	engineered mutation	UNP P0A3G3
B	273	LEU	TYR	engineered mutation	UNP P0A3G3
B	291	SER	PRO	engineered mutation	UNP P0A3G3
B	292	THR	ALA	engineered mutation	UNP P0A3G3
B	294	GLU	-	expression tag	UNP P0A3G3
B	295	ILE	-	expression tag	UNP P0A3G3
C	3	GLY	-	expression tag	UNP P0A3G3
C	47	VAL	LEU	engineered mutation	UNP P0A3G3
C	58	THR	SER	engineered mutation	UNP P0A3G3
C	78	GLY	ASP	engineered mutation	UNP P0A3G3
C	87	PHE	TYR	engineered mutation	UNP P0A3G3
C	88	MET	LEU	engineered mutation	UNP P0A3G3
C	128	PHE	CYS	engineered mutation	UNP P0A3G3
C	155	THR	ALA	engineered mutation	UNP P0A3G3
C	160	LYS	GLU	engineered mutation	UNP P0A3G3
C	167	VAL	ALA	engineered mutation	UNP P0A3G3
C	172	THR	ALA	engineered mutation	UNP P0A3G3
C	175	TRP	LYS	engineered mutation	UNP P0A3G3
C	176	GLY	CYS	engineered mutation	UNP P0A3G3
C	195	ASN	LYS	engineered mutation	UNP P0A3G3
C	224	GLU	ALA	engineered mutation	UNP P0A3G3

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Chain	Residue	Modelled	Actual	Comment	Reference
C	227	ASP	ASN	engineered mutation	UNP P0A3G3
C	257	LYS	GLU	engineered mutation	UNP P0A3G3
C	264	ALA	THR	engineered mutation	UNP P0A3G3
C	272	ASN	HIS	engineered mutation	UNP P0A3G3
C	273	LEU	TYR	engineered mutation	UNP P0A3G3
C	291	SER	PRO	engineered mutation	UNP P0A3G3
C	292	THR	ALA	engineered mutation	UNP P0A3G3
C	294	GLU	-	expression tag	UNP P0A3G3
C	295	ILE	-	expression tag	UNP P0A3G3
D	3	GLY	-	expression tag	UNP P0A3G3
D	47	VAL	LEU	engineered mutation	UNP P0A3G3
D	58	THR	SER	engineered mutation	UNP P0A3G3
D	78	GLY	ASP	engineered mutation	UNP P0A3G3
D	87	PHE	TYR	engineered mutation	UNP P0A3G3
D	88	MET	LEU	engineered mutation	UNP P0A3G3
D	128	PHE	CYS	engineered mutation	UNP P0A3G3
D	155	THR	ALA	engineered mutation	UNP P0A3G3
D	160	LYS	GLU	engineered mutation	UNP P0A3G3
D	167	VAL	ALA	engineered mutation	UNP P0A3G3
D	172	THR	ALA	engineered mutation	UNP P0A3G3
D	175	TRP	LYS	engineered mutation	UNP P0A3G3
D	176	GLY	CYS	engineered mutation	UNP P0A3G3
D	195	ASN	LYS	engineered mutation	UNP P0A3G3
D	224	GLU	ALA	engineered mutation	UNP P0A3G3
D	227	ASP	ASN	engineered mutation	UNP P0A3G3
D	257	LYS	GLU	engineered mutation	UNP P0A3G3
D	264	ALA	THR	engineered mutation	UNP P0A3G3
D	272	ASN	HIS	engineered mutation	UNP P0A3G3
D	273	LEU	TYR	engineered mutation	UNP P0A3G3
D	291	SER	PRO	engineered mutation	UNP P0A3G3
D	292	THR	ALA	engineered mutation	UNP P0A3G3
D	294	GLU	-	expression tag	UNP P0A3G3
D	295	ILE	-	expression tag	UNP P0A3G3

- Molecule 2 is [9-[2-carboxy-5-[2-[2-(6-chloranylhexoxy)ethoxy]ethylcarbamoyl]phenyl]-6-(dimethylamino)xanthen-3-ylidene]-dimethyl-azanium (three-letter code: OEH) (formula: C<sub>35</sub>H<sub>43</sub>ClN<sub>3</sub>O<sub>6</sub>) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
2	A	1	44	35	3	6	0	0
2	B	1	44	35	3	6	0	0
2	C	1	44	35	3	6	0	0
2	D	1	44	35	3	6	0	0

- Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Cl		
3	A	1	1	1	0	0
3	B	1	1	1	0	0
3	C	1	1	1	0	0
3	D	1	1	1	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	A	34	34	34	0	0

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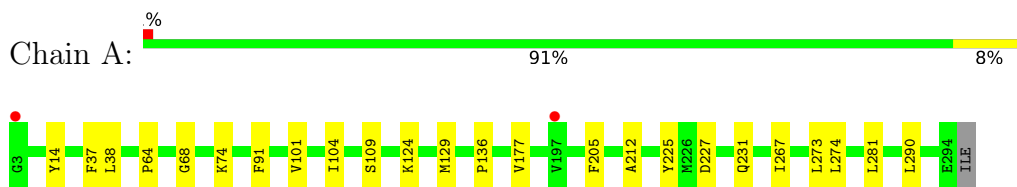
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<b>Mol</b>	<b>Chain</b>	<b>Residues</b>	<b>Atoms</b>		<b>ZeroOcc</b>	<b>AltConf</b>
4	B	36	Total 36	O 36	0	0
4	C	32	Total 32	O 32	0	0
4	D	23	Total 23	O 23	0	0

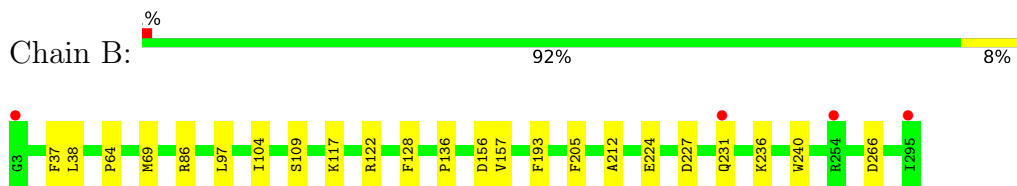
### 3 Residue-property plots [i](#)

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.

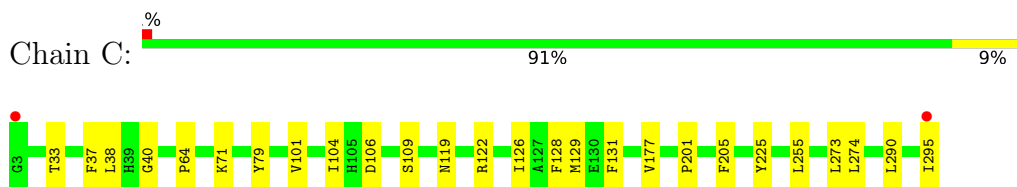
- Molecule 1: Haloalkane dehalogenase



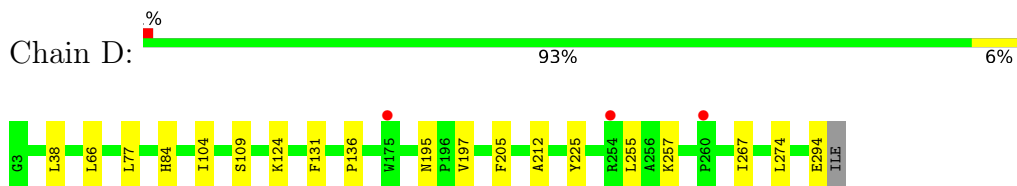
- Molecule 1: Haloalkane dehalogenase



- Molecule 1: Haloalkane dehalogenase



- Molecule 1: Haloalkane dehalogenase





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	44.32Å 76.56Å 85.76Å 68.40° 79.81° 79.02°	Depositor
Resolution (Å)	40.36 – 2.30 45.43 – 2.30	Depositor EDS
% Data completeness (in resolution range)	95.8 (40.36-2.30) 95.7 (45.43-2.30)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.89 (at 2.29Å)	Xtrriage
Refinement program	PHENIX 1.19.2_4158	Depositor
R, $R_{free}$	0.191 , 0.243 0.190 , 0.242	Depositor DCC
$R_{free}$ test set	2174 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	28.1	Xtrriage
Anisotropy	0.610	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	(Not available) , (Not available)	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	9711	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	29.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 16.91% of the height of the origin peak. No significant pseudotranslation is detected.*

<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: CL, OEH

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	0.25	0/2428	0.48	0/3322
1	B	0.24	0/2437	0.49	0/3333
1	C	0.25	0/2437	0.48	0/3333
1	D	0.24	0/2428	0.48	0/3322
All	All	0.24	0/9730	0.48	0/13310

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2347	0	2281	14	0
1	B	2356	0	2292	12	0
1	C	2356	0	2292	14	0
1	D	2347	0	2281	11	0
2	A	44	0	0	0	0
2	B	44	0	0	0	0
2	C	44	0	0	0	0
2	D	44	0	0	0	0
3	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	34	0	0	0	0
4	B	36	0	0	1	0
4	C	32	0	0	1	0
4	D	23	0	0	0	0
All	All	9711	0	9146	50	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 3.

All (50) 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:C:71:LYS:NZ	4:C:401:HOH:O	2.23	0.72
1:C:129:MET:HB2	1:C:274:LEU:HD12	1.85	0.58
1:B:117:LYS:HE3	1:B:224:GLU:HB3	1.85	0.58
1:C:38:LEU:HB2	1:C:104:ILE:HG22	1.86	0.57
1:C:119:ASN:HB3	1:C:122:ARG:HD3	1.88	0.55
1:A:129:MET:HB2	1:A:274:LEU:HD12	1.89	0.54
1:A:38:LEU:HB2	1:A:104:ILE:HG22	1.90	0.53
1:D:124:LYS:NZ	1:D:294:GLU:O	2.40	0.52
1:B:227:ASP:O	1:B:231:GLN:HG2	2.10	0.52
1:B:86:ARG:NH1	4:B:401:HOH:O	2.43	0.52
1:C:104:ILE:HD12	1:C:109:SER:HA	1.93	0.51
1:D:104:ILE:HD12	1:D:109:SER:HA	1.93	0.50
1:B:69:MET:HG2	1:B:193:PHE:HZ	1.77	0.49
1:B:156:ASP:OD1	1:B:157:VAL:N	2.40	0.49
1:D:195:ASN:ND2	1:D:197:VAL:HG22	2.28	0.49
1:D:38:LEU:HB2	1:D:104:ILE:HG22	1.96	0.48
1:C:104:ILE:HB	1:C:109:SER:HA	1.95	0.48
1:A:68:GLY:HA2	1:A:74:LYS:HG2	1.95	0.48
1:A:104:ILE:HB	1:A:109:SER:HA	1.96	0.48
1:B:97:LEU:O	1:B:122:ARG:NH2	2.43	0.48
1:C:177:VAL:HG22	1:C:273:LEU:HB2	1.96	0.47
1:A:104:ILE:HD12	1:A:109:SER:HA	1.96	0.47
1:D:104:ILE:HB	1:D:109:SER:HA	1.96	0.47
1:C:33:THR:HB	1:C:295:ILE:HB	1.96	0.47
1:B:37:PHE:O	1:B:64:PRO:HD2	2.15	0.46
1:A:64:PRO:HB3	1:A:91:PHE:CZ	2.51	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:66:LEU:HD12	1:D:84:HIS:ND1	2.31	0.46
1:B:104:ILE:HB	1:B:109:SER:HA	1.98	0.46
1:C:131:PHE:HZ	1:C:255:LEU:HD13	1.82	0.45
1:B:136:PRO:O	1:B:212:ALA:HA	2.16	0.45
1:A:101:VAL:HG11	1:A:290:LEU:HD21	1.98	0.44
1:A:267:ILE:HA	1:A:281:LEU:HD23	1.99	0.44
1:B:128:PHE:HE1	1:B:236:LYS:HB3	1.82	0.44
1:C:40:GLY:HA3	1:C:106:ASP:HB3	1.98	0.44
1:C:79:TYR:CG	1:C:201:PRO:HG3	2.52	0.44
1:B:38:LEU:HB2	1:B:104:ILE:HG22	1.99	0.44
1:A:136:PRO:O	1:A:212:ALA:HA	2.18	0.44
1:C:126:ILE:HG22	1:C:128:PHE:HD1	1.83	0.43
1:A:177:VAL:HG22	1:A:273:LEU:HB2	2.00	0.43
1:D:131:PHE:HZ	1:D:255:LEU:HD13	1.84	0.42
1:D:136:PRO:O	1:D:212:ALA:HA	2.19	0.42
1:D:195:ASN:HD21	1:D:197:VAL:HG13	1.84	0.42
1:A:227:ASP:O	1:A:231:GLN:HG2	2.20	0.42
1:B:240:TRP:CZ2	1:B:266:ASP:HB2	2.55	0.42
1:A:124:LYS:HA	1:A:124:LYS:HD3	1.85	0.42
1:C:37:PHE:O	1:C:64:PRO:HD2	2.20	0.42
1:A:37:PHE:O	1:A:64:PRO:HD2	2.20	0.42
1:C:101:VAL:HG11	1:C:290:LEU:HD21	2.02	0.41
1:A:14:TYR:OH	1:D:257:LYS:HB3	2.20	0.41
1:D:267:ILE:HD13	1:D:274:LEU:HB3	2.02	0.40

There are no symmetry-related clashes.

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

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

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	290/293 (99%)	281 (97%)	9 (3%)	0	<a href="#">100</a> <a href="#">100</a>

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	291/293 (99%)	280 (96%)	11 (4%)	0	100	100
1	C	291/293 (99%)	281 (97%)	10 (3%)	0	100	100
1	D	290/293 (99%)	278 (96%)	12 (4%)	0	100	100
All	All	1162/1172 (99%)	1120 (96%)	42 (4%)	0	100	100

There are no Ramachandran outliers to report.

### 5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	251/252 (100%)	249 (99%)	2 (1%)	81	91
1	B	252/252 (100%)	251 (100%)	1 (0%)	91	96
1	C	252/252 (100%)	250 (99%)	2 (1%)	81	91
1	D	251/252 (100%)	248 (99%)	3 (1%)	71	84
All	All	1006/1008 (100%)	998 (99%)	8 (1%)	81	91

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	205	PHE
1	A	225	TYR
1	B	205	PHE
1	C	205	PHE
1	C	225	TYR
1	D	77	LEU
1	D	205	PHE
1	D	225	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	D	195	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 [i](#)

Of 8 ligands modelled in this entry, 4 are monoatomic - leaving 4 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

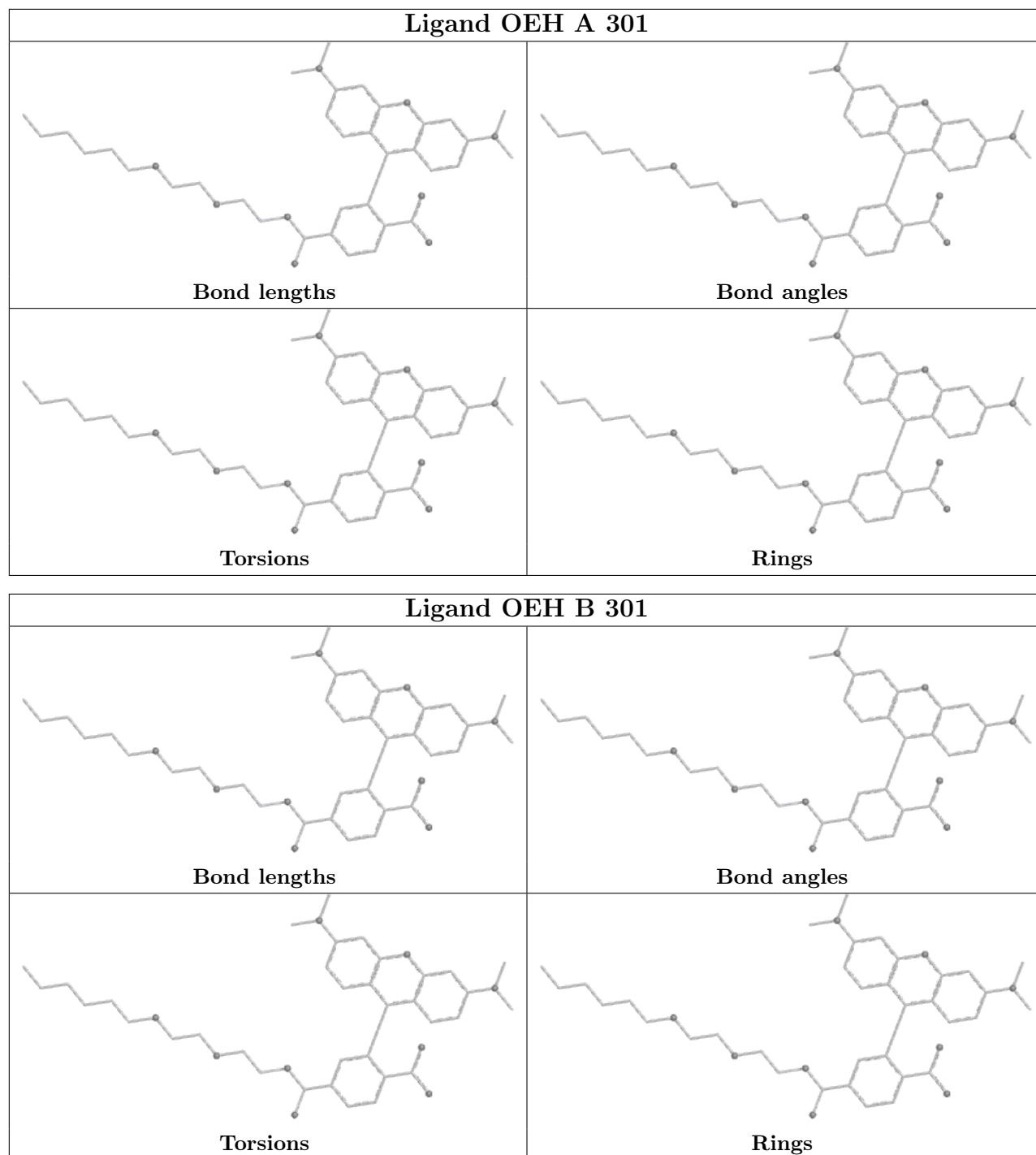
There are no chirality outliers.

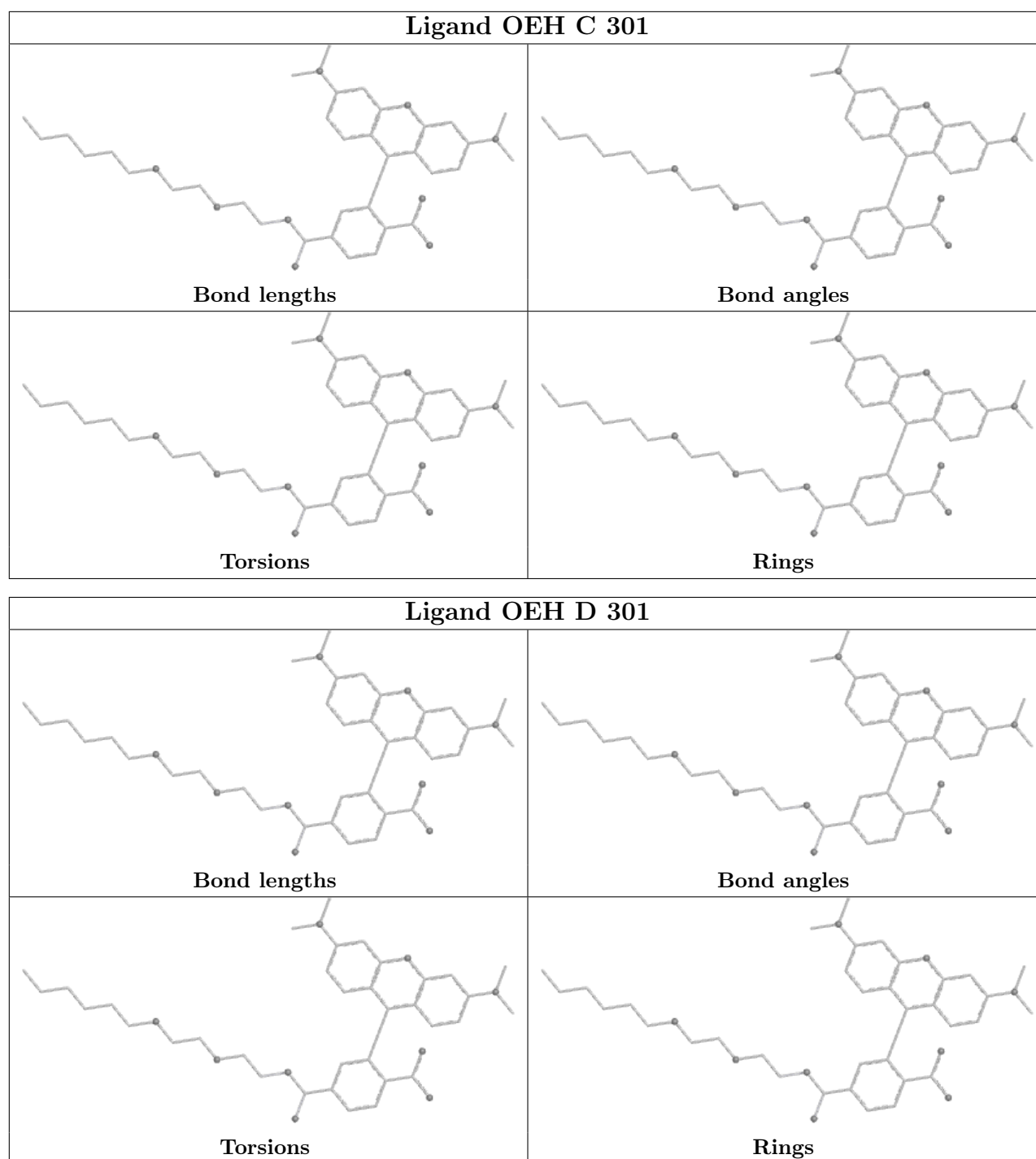
There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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 [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, 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	292/293 (99%)	0.01	2 (0%) 87 91	20, 28, 41, 60	0
1	B	293/293 (100%)	0.03	4 (1%) 75 80	20, 26, 39, 68	0
1	C	293/293 (100%)	0.03	2 (0%) 87 91	21, 28, 39, 59	0
1	D	292/293 (99%)	0.16	3 (1%) 82 86	22, 31, 43, 72	0
All	All	1170/1172 (99%)	0.06	11 (0%) 84 88	20, 28, 41, 72	0

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	295	ILE	6.2
1	C	3	GLY	4.5
1	A	197	VAL	4.0
1	A	3	GLY	3.5
1	B	3	GLY	2.9
1	B	295	ILE	2.7
1	B	231	GLN	2.6
1	D	175	TRP	2.4
1	B	254	ARG	2.4
1	D	260	PRO	2.3
1	D	254	ARG	2.1

### 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](#)

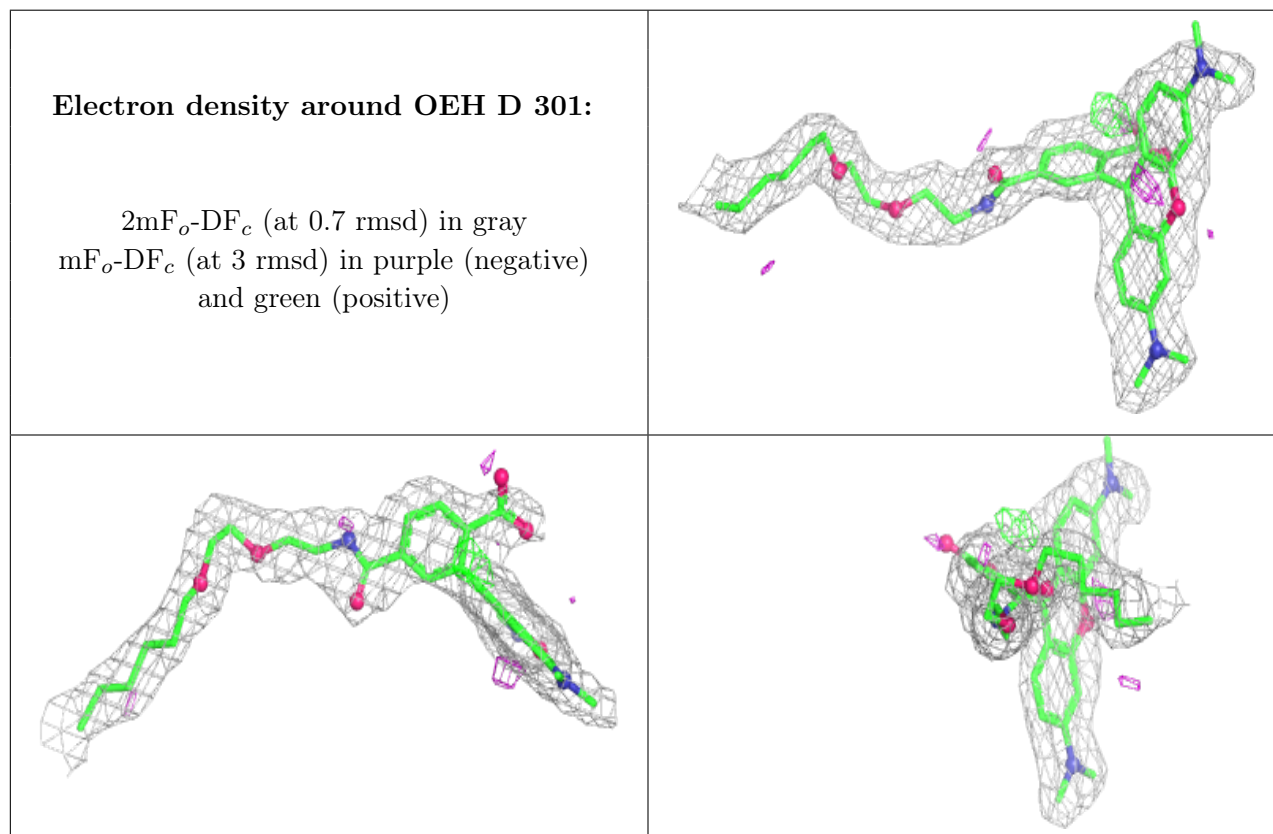
There are no monosaccharides in this entry.

## 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, 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.

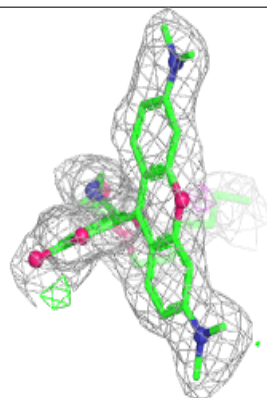
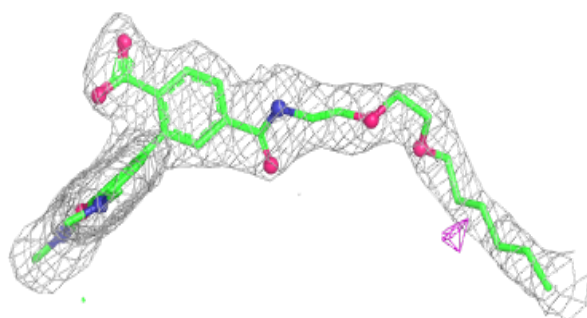
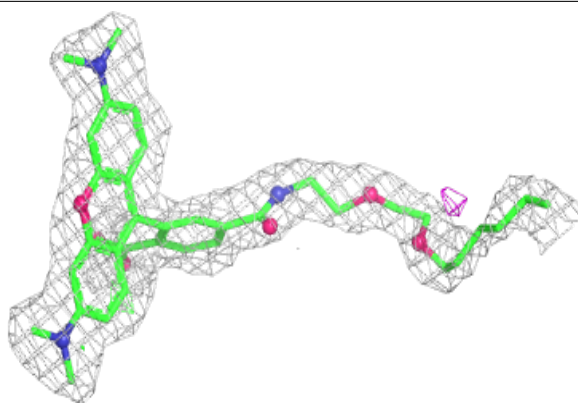
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
2	OEH	D	301	44/45	0.89	0.18	22,34,45,62	0
2	OEH	C	301	44/45	0.91	0.15	16,26,37,44	0
2	OEH	B	301	44/45	0.92	0.14	19,26,31,43	0
2	OEH	A	301	44/45	0.93	0.13	19,28,35,43	0
3	CL	A	302	1/1	0.97	0.12	22,22,22,22	0
3	CL	B	302	1/1	0.98	0.12	20,20,20,20	0
3	CL	D	302	1/1	0.98	0.12	25,25,25,25	0
3	CL	C	302	1/1	1.00	0.11	21,21,21,21	0

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

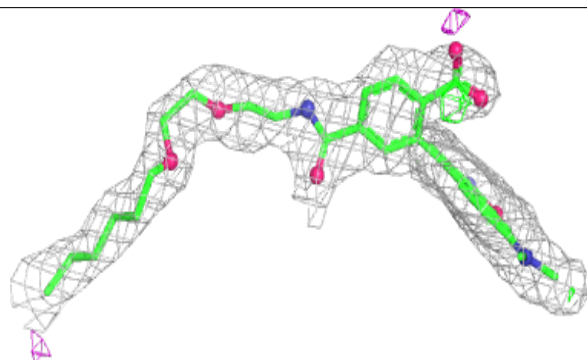
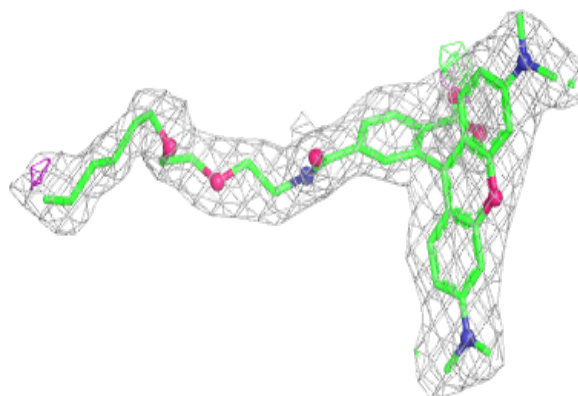


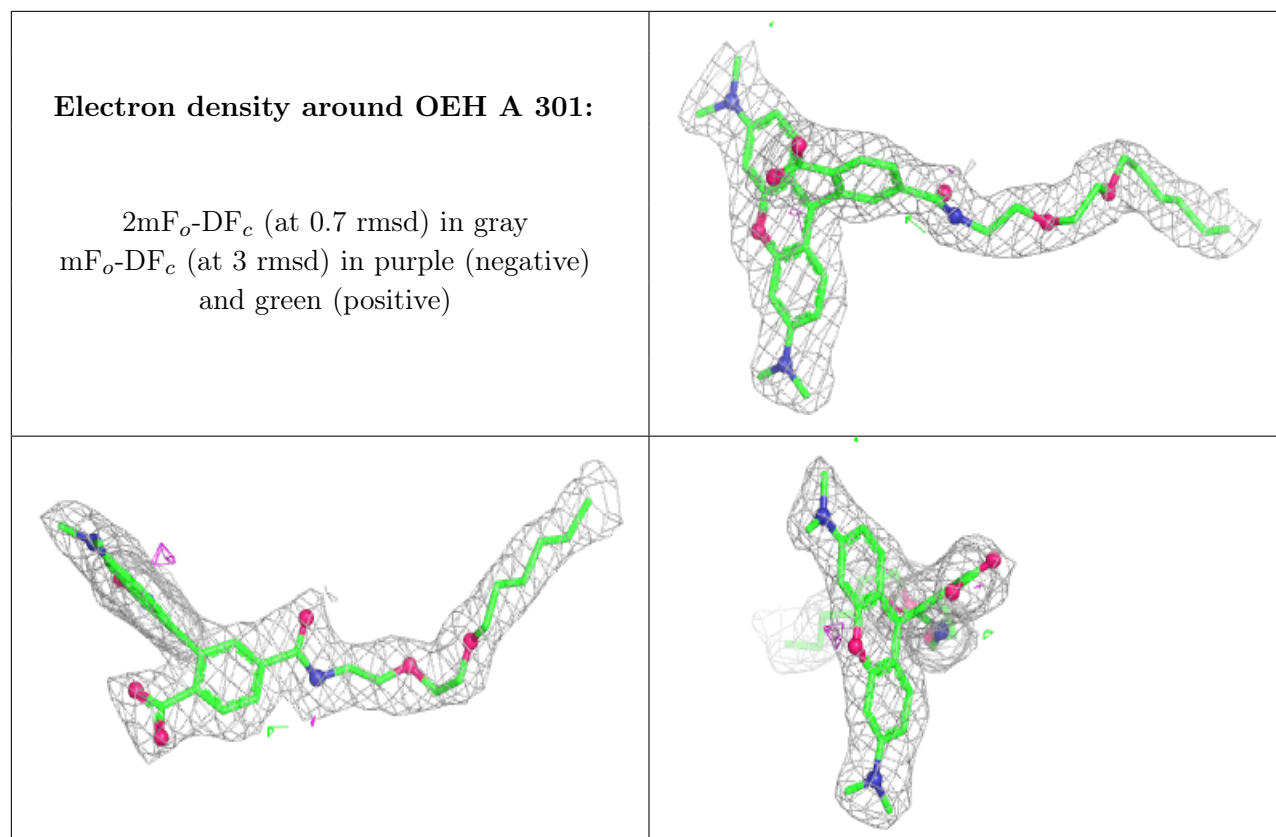
**Electron density around OEH C 301:**

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

**Electron density around OEH B 301:**

$2mF_o-DF_c$  (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.