



# Full wwPDB X-ray Structure Validation Report ⓘ

Sep 19, 2019 – 11:00 AM EDT

PDB ID : 4D2C  
Title : Structure of a di peptide bound POT family peptide transporter  
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Deposited on : 2014-05-09  
Resolution : 2.47 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 : 1.8.0 (224370), CSD as540be (2019)  
Xtriage (Phenix) : 1.13  
EDS : 2.4  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20171227.v01 (using entries in the PDB archive December 27th 2017)  
Refmac : 5.8.0158  
CCP4 : 7.0 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.4

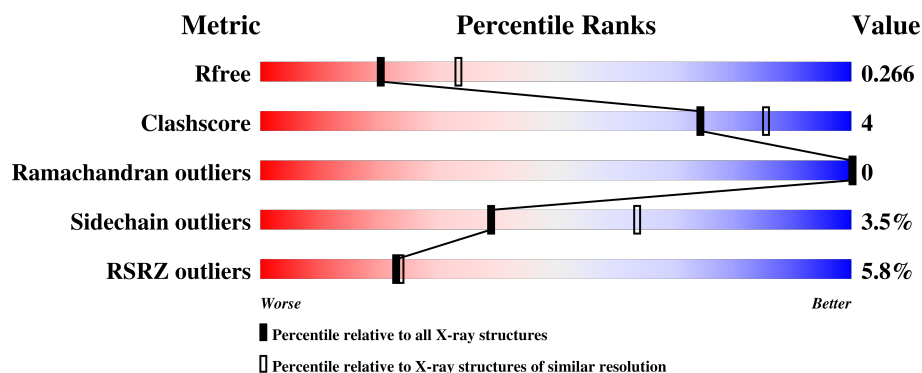
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.47 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))
$R_{free}$	111664	5140 (2.50-2.46)
Clashscore	122126	5860 (2.50-2.46)
Ramachandran outliers	120053	5763 (2.50-2.46)
Sidechain outliers	120020	5765 (2.50-2.46)
RSRZ outliers	108989	5026 (2.50-2.46)

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. 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	491	<div> <div>5%</div> <div>78%</div> <div>8%</div> <div>13%</div> </div>

## 2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 3513 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.

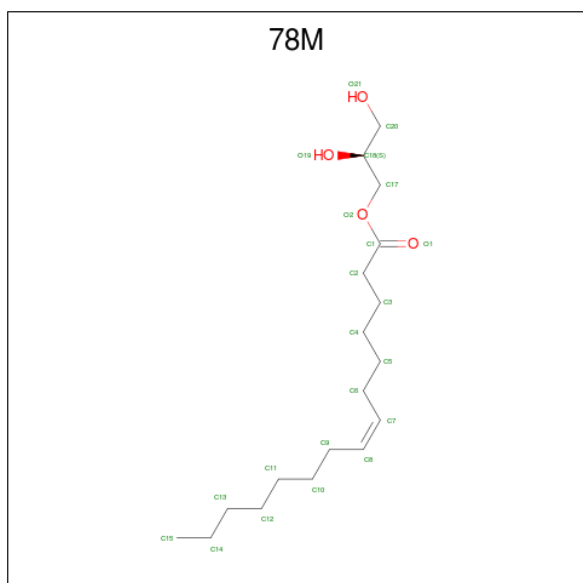
- Molecule 1 is a protein called Di-or tripeptide:H<sup>+</sup> symporter.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	429	3321	2248	505	552	16	0	0	0

There are 8 discrepancies between the modelled and reference sequences:

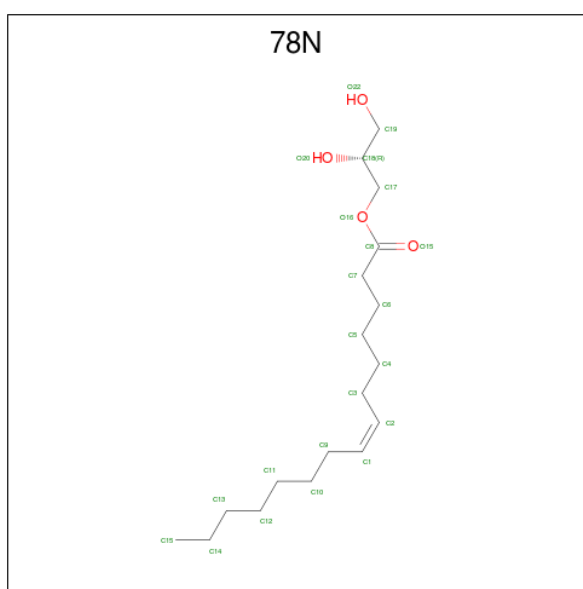
Chain	Residue	Modelled	Actual	Comment	Reference
A	484	GLY	-	expression tag	UNP Q5M4H8
A	485	SER	-	expression tag	UNP Q5M4H8
A	486	GLU	-	expression tag	UNP Q5M4H8
A	487	ASN	-	expression tag	UNP Q5M4H8
A	488	LEU	-	expression tag	UNP Q5M4H8
A	489	TYR	-	expression tag	UNP Q5M4H8
A	490	PHE	-	expression tag	UNP Q5M4H8
A	491	GLN	-	expression tag	UNP Q5M4H8

- Molecule 2 is (2S)-2,3-DIHYDROXYPROPYL(7Z)-PENTADEC-7-ENOATE (three-letter code: 78M) (formula: C<sub>18</sub>H<sub>34</sub>O<sub>4</sub>).



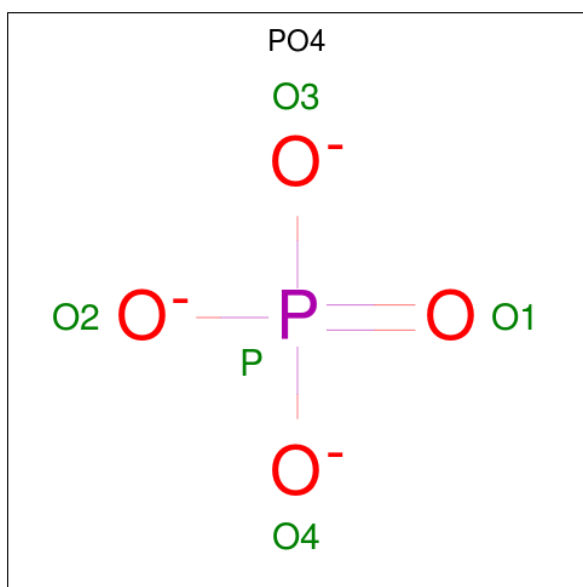
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			22	18	4		
2	A	1	Total	C	O	0	0
			22	18	4		
2	A	1	Total	C	O	0	0
			22	18	4		
2	A	1	Total	C	O	0	0
			22	18	4		

- Molecule 3 is (2R)-2,3-DIHYDROXYPROPYL(7Z)-PENTADEC-7-ENOATE (three-letter code: 78N) (formula: C<sub>18</sub>H<sub>34</sub>O<sub>4</sub>).



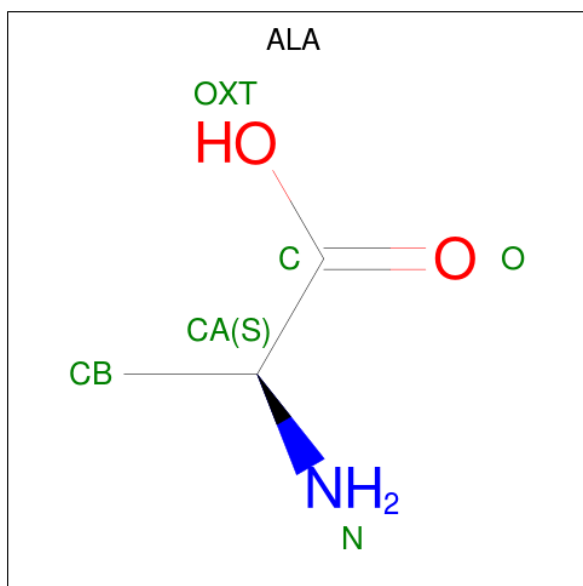
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			22	18	4		
3	A	1	Total	C	O	0	0
			22	18	4		

- Molecule 4 is PHOSPHATE ION (three-letter code: PO4) (formula: O<sub>4</sub>P).



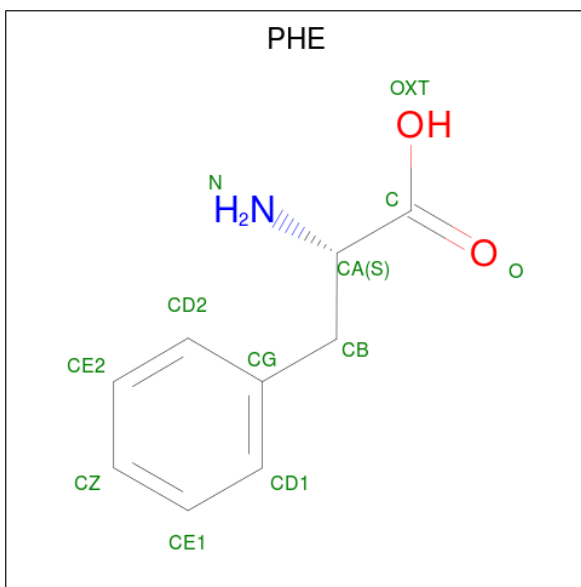
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	P	0	0
			5	4	1		

- Molecule 5 is ALANINE (three-letter code: ALA) (formula:  $C_3H_7NO_2$ ).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	0	0
			5	3	1	1		

- Molecule 6 is PHENYLALANINE (three-letter code: PHE) (formula:  $C_9H_{11}NO_2$ ).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	1	Total	C	N	O	0	0
			12	9	1	2		

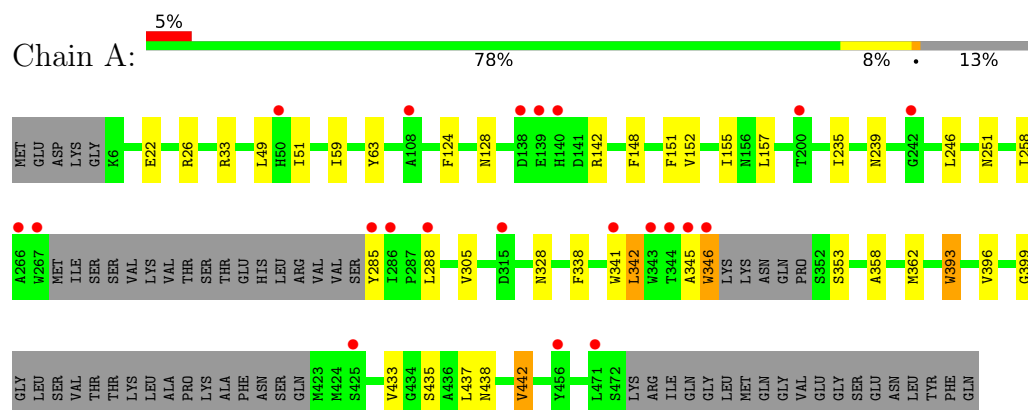
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	38	Total	O	0	0
			38	38		

### 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: Di- or tripeptide:H<sup>+</sup> symporter



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	102.06Å 110.33Å 110.74Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	51.03 – 2.47 51.03 – 2.47	Depositor EDS
% Data completeness (in resolution range)	97.2 (51.03-2.47) 97.2 (51.03-2.47)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.10 (at 2.48Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, $R_{free}$	0.222 , 0.266 0.222 , 0.266	Depositor DCC
$R_{free}$ test set	1159 reflections (5.24%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	47.0	Xtriage
Anisotropy	0.220	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 57.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	3513	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.05% 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

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 78M, PO4, 78N

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	0.24	0/3424	0.39	0/4673

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

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	3321	0	3385	22	0
2	A	88	0	136	3	0
3	A	44	0	68	2	0
4	A	5	0	0	0	0
5	A	5	0	4	1	0
6	A	12	0	9	0	0
7	A	38	0	0	1	0
All	All	3513	0	3602	25	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 4.

All (25) 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:A:251:ASN:HD22	2:A:502:78M:H172	1.66	0.61
1:A:235:ILE:O	1:A:239:ASN:ND2	2.31	0.57
1:A:341:TRP:O	1:A:345:ALA:N	2.34	0.56
1:A:59:ILE:HD11	1:A:246:LEU:HD21	1.88	0.56
1:A:285:TYR:HA	1:A:288:LEU:HD12	1.88	0.56
1:A:124:PHE:O	1:A:128:ASN:ND2	2.36	0.55
1:A:258:ILE:HG12	1:A:433:VAL:HG13	1.90	0.54
1:A:261:PRO:HG2	1:A:433:VAL:HG21	1.90	0.53
1:A:305:VAL:HG23	1:A:442:VAL:HG22	1.90	0.53
1:A:399:GLY:O	1:A:403:ILE:HG12	2.12	0.50
1:A:33:ARG:NH1	7:A:606:HOH:O	2.44	0.50
1:A:342:LEU:O	1:A:346:TRP:HD1	1.97	0.47
3:A:503:78N:H91C	2:A:504:78M:H8	1.99	0.45
1:A:328:ASN:OD1	5:A:508:ALA:N	2.50	0.44
3:A:506:78N:H32C	3:A:506:78N:H92C	1.86	0.44
1:A:358:ALA:O	1:A:362:MET:HG3	2.19	0.43
1:A:148:PHE:O	1:A:152:VAL:HG23	2.19	0.43
1:A:438:ASN:O	1:A:442:VAL:HB	2.20	0.42
1:A:338:PHE:O	1:A:341:TRP:HB2	2.21	0.41
1:A:258:ILE:O	1:A:261:PRO:HD2	2.20	0.41
1:A:26:ARG:HG3	1:A:155:ILE:HG23	2.01	0.41
1:A:393:TRP:HA	1:A:396:VAL:HG12	2.01	0.41
1:A:22:GLU:HG3	1:A:151:PHE:CZ	2.56	0.41
1:A:49:LEU:HG	1:A:51:ILE:HG12	2.02	0.40
2:A:501:78M:H62C	2:A:501:78M:H92C	1.89	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles ⓘ

### 5.3.1 Protein backbone ⓘ

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	421/491 (86%)	417 (99%)	4 (1%)	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	343/398 (86%)	331 (96%)	12 (4%)	39 64

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	63	TYR
1	A	142	ARG
1	A	157	LEU
1	A	264	TYR
1	A	342	LEU
1	A	346	TRP
1	A	353	SER
1	A	393	TRP
1	A	404	SER
1	A	435	SER
1	A	437	LEU
1	A	442	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	328	ASN
1	A	440	GLN

### 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 carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

9 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
2	78M	A	501	-	21,21,21	0.89	1 (4%)	22,22,22	1.10	2 (9%)
2	78M	A	502	-	21,21,21	0.90	1 (4%)	22,22,22	1.11	2 (9%)
3	78N	A	503	-	21,21,21	0.90	1 (4%)	22,22,22	1.10	2 (9%)
2	78M	A	504	-	21,21,21	0.89	1 (4%)	22,22,22	1.13	2 (9%)
2	78M	A	505	-	21,21,21	0.88	1 (4%)	22,22,22	1.02	2 (9%)
3	78N	A	506	-	21,21,21	0.88	1 (4%)	22,22,22	1.16	2 (9%)
4	PO4	A	507	-	4,4,4	0.83	0	6,6,6	0.33	0

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	78M	A	501	-	-	11/21/21/21	-
2	78M	A	502	-	-	11/21/21/21	-
3	78N	A	503	-	-	11/21/21/21	-
2	78M	A	504	-	-	9/21/21/21	-
2	78M	A	505	-	-	14/21/21/21	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	78N	A	506	-	-	15/21/21/21	-

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	78M	O2-C17	-2.70	1.38	1.45
2	A	504	78M	O2-C17	-2.67	1.39	1.45
3	A	503	78N	O16-C17	-2.62	1.39	1.45
2	A	505	78M	O2-C17	-2.61	1.39	1.45
3	A	506	78N	O16-C17	-2.56	1.39	1.45
2	A	502	78M	O2-C17	-2.55	1.39	1.45

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	506	78N	O16-C8-C7	2.79	120.90	111.93
2	A	504	78M	O2-C1-C2	2.68	120.53	111.93
3	A	503	78N	O16-C8-C7	2.67	120.51	111.93
2	A	502	78M	O2-C1-C2	2.63	120.38	111.93
2	A	501	78M	O2-C1-C2	2.56	120.14	111.93
2	A	504	78M	O2-C17-C18	2.55	118.07	105.77
2	A	502	78M	O2-C17-C18	2.55	118.07	105.77
3	A	506	78N	O16-C17-C18	2.40	117.35	105.77
2	A	501	78M	O2-C17-C18	2.33	117.01	105.77
3	A	503	78N	O16-C17-C18	2.30	116.87	105.77
2	A	505	78M	O2-C17-C18	2.19	116.35	105.77
2	A	505	78M	O2-C1-C2	2.19	118.96	111.93

There are no chirality outliers.

All (71) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	501	78M	C17-C18-C20-O21
3	A	506	78N	O16-C17-C18-C19
3	A	503	78N	C17-C18-C19-O22
2	A	505	78M	C17-C18-C20-O21
2	A	504	78M	O2-C17-C18-O19
3	A	506	78N	O15-C8-O16-C17
3	A	506	78N	C7-C8-O16-C17
2	A	501	78M	C2-C1-O2-C17
2	A	501	78M	O1-C1-O2-C17

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Mol	Chain	Res	Type	Atoms
2	A	504	78M	C2-C1-O2-C17
2	A	505	78M	O2-C17-C18-C20
2	A	505	78M	O2-C17-C18-O19
3	A	506	78N	O20-C18-C19-O22
3	A	503	78N	O20-C18-C19-O22
2	A	504	78M	O1-C1-O2-C17
2	A	501	78M	O2-C17-C18-O19
3	A	506	78N	O16-C17-C18-O20
2	A	502	78M	O2-C17-C18-O19
2	A	501	78M	C3-C4-C5-C6
2	A	501	78M	C2-C3-C4-C5
2	A	501	78M	O2-C17-C18-C20
2	A	502	78M	O2-C17-C18-C20
2	A	504	78M	O2-C17-C18-C20
3	A	506	78N	C3-C4-C5-C6
3	A	506	78N	C4-C5-C6-C7
3	A	506	78N	C17-C18-C19-O22
2	A	504	78M	C17-C18-C20-O21
2	A	501	78M	C11-C12-C13-C14
2	A	502	78M	C11-C12-C13-C14
2	A	505	78M	C9-C10-C11-C12
3	A	506	78N	C2-C3-C4-C5
2	A	505	78M	C11-C10-C9-C8
3	A	503	78N	C7-C8-O16-C17
2	A	502	78M	C2-C3-C4-C5
2	A	501	78M	O19-C18-C20-O21
2	A	504	78M	O19-C18-C20-O21
3	A	503	78N	C11-C10-C9-C1
2	A	502	78M	C3-C4-C5-C6
2	A	504	78M	C2-C3-C4-C5
3	A	503	78N	O15-C8-O16-C17
3	A	506	78N	C11-C10-C9-C1
2	A	502	78M	C4-C5-C6-C7
2	A	504	78M	C10-C11-C12-C13
2	A	504	78M	C9-C10-C11-C12
3	A	506	78N	C11-C12-C13-C14
2	A	505	78M	O19-C18-C20-O21
2	A	502	78M	C12-C13-C14-C15
3	A	506	78N	C10-C11-C12-C13
2	A	505	78M	C4-C5-C6-C7
3	A	503	78N	C11-C12-C13-C14
2	A	505	78M	C11-C12-C13-C14

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Mol	Chain	Res	Type	Atoms
3	A	503	78N	C2-C3-C4-C5
2	A	502	78M	C1-C2-C3-C4
2	A	502	78M	C11-C10-C9-C8
3	A	503	78N	C4-C5-C6-C7
3	A	506	78N	C9-C10-C11-C12
2	A	502	78M	C10-C11-C12-C13
2	A	505	78M	O2-C1-C2-C3
2	A	505	78M	C5-C6-C7-C8
2	A	505	78M	C2-C1-O2-C17
2	A	505	78M	O1-C1-O2-C17
3	A	503	78N	C12-C13-C14-C15
3	A	503	78N	C1-C2-C3-C4
2	A	505	78M	C2-C3-C4-C5
2	A	501	78M	C5-C6-C7-C8
3	A	506	78N	C2-C1-C9-C10
3	A	503	78N	C2-C1-C9-C10
2	A	505	78M	C10-C11-C12-C13
2	A	502	78M	C5-C6-C7-C8
3	A	506	78N	C5-C6-C7-C8
2	A	501	78M	C1-C2-C3-C4

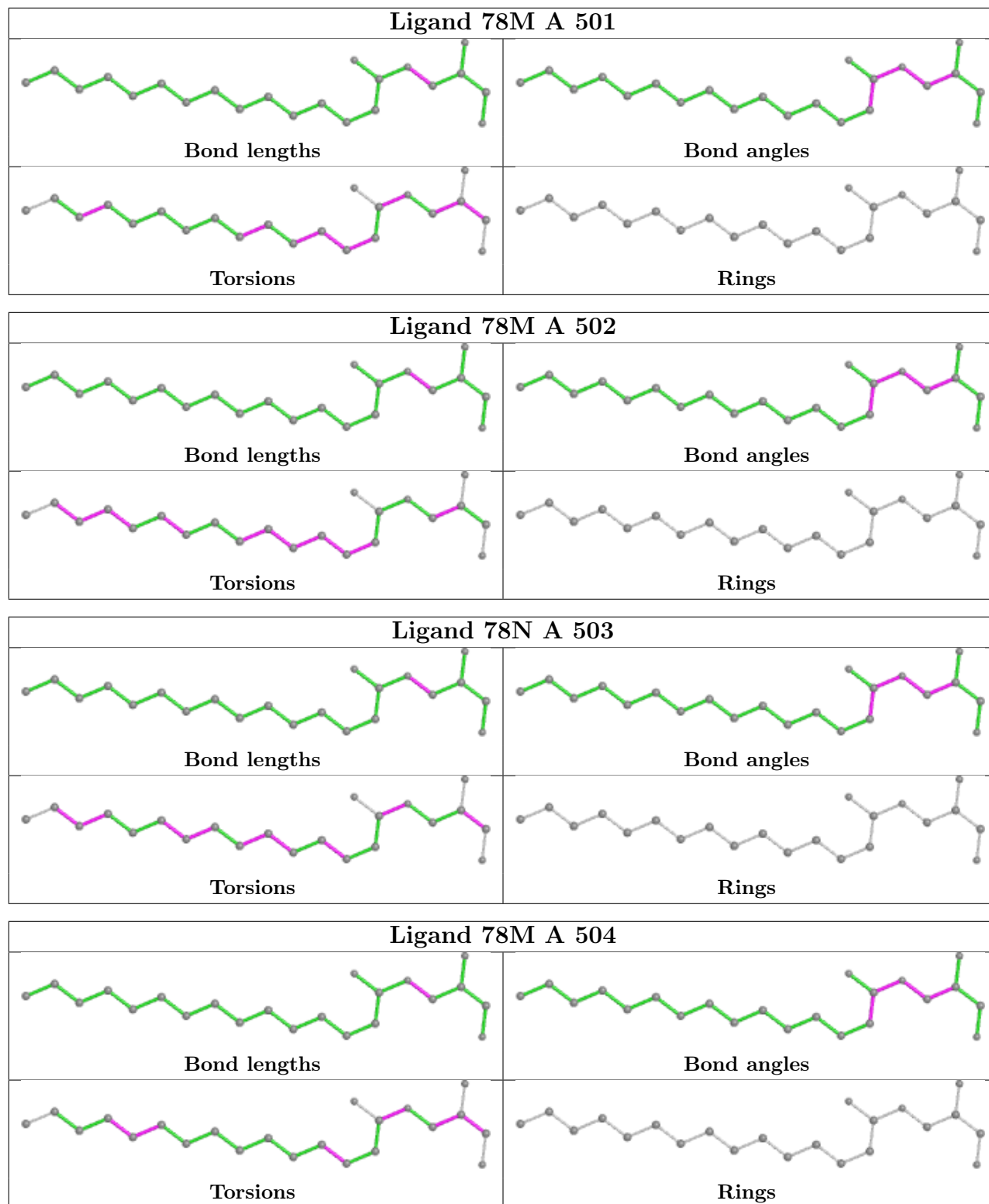
There are no ring outliers.

5 monomers are involved in 4 short contacts:

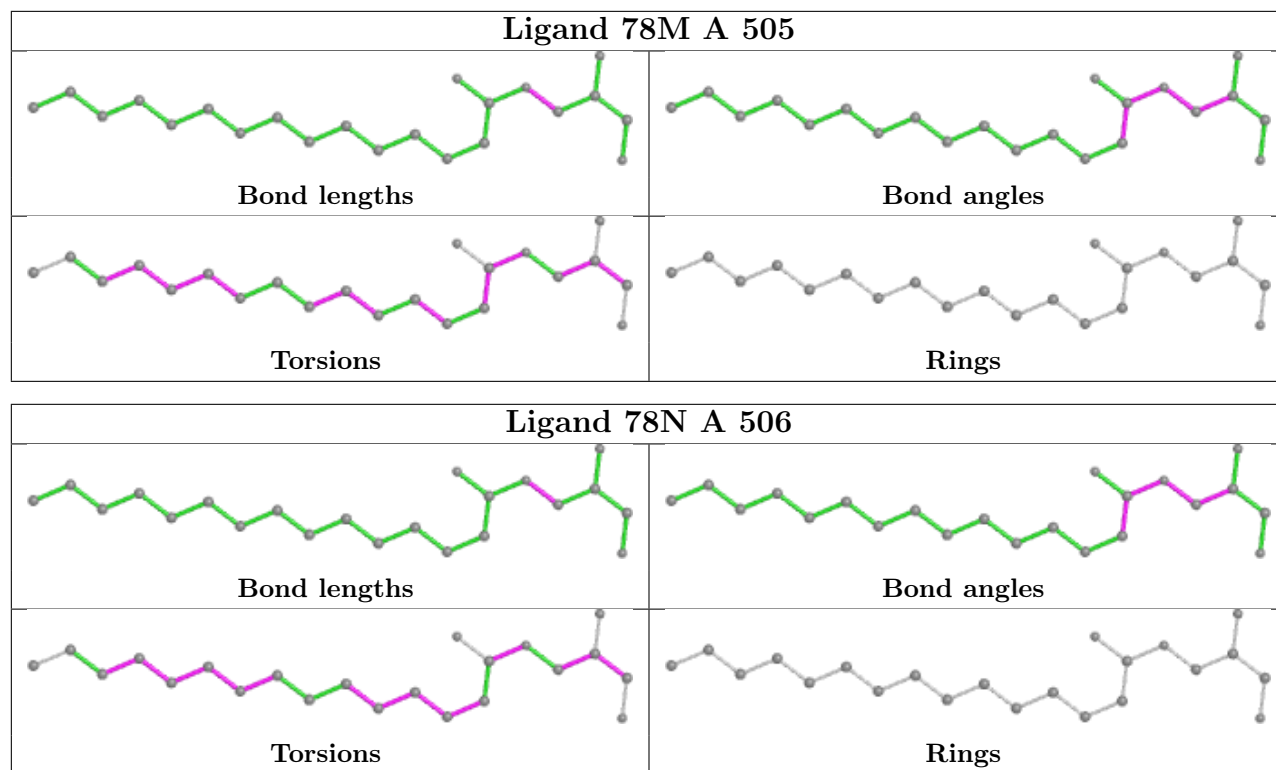
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	501	78M	1	0
2	A	502	78M	1	0
3	A	503	78N	1	0
2	A	504	78M	1	0
3	A	506	78N	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient

equivalents in the CSD to analyse the geometry.







## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	429/491 (87%)	0.12	25 (5%) 23 24	19, 33, 75, 101	0

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	346	TRP	6.0
1	A	344	THR	5.5
1	A	140	HIS	4.3
1	A	139	GLU	3.9
1	A	343	TRP	3.5
1	A	341	TRP	3.5
1	A	405	PRO	3.5
1	A	242	GLY	3.4
1	A	345	ALA	3.3
1	A	267	TRP	3.3
1	A	285	TYR	3.1
1	A	288	LEU	2.9
1	A	264	TYR	2.8
1	A	138	ASP	2.7
1	A	286	ILE	2.6
1	A	200	THR	2.5
1	A	108	ALA	2.5
1	A	50	HIS	2.3
1	A	265	PHE	2.3
1	A	315	ASP	2.3
1	A	425	SER	2.3
1	A	404	SER	2.2
1	A	471	LEU	2.2
1	A	266	ALA	2.1
1	A	456	TYR	2.0

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

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

## 6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 6.4 Ligands [i](#)

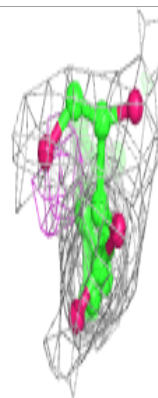
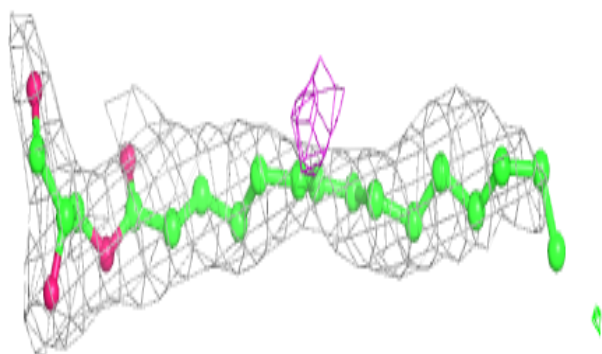
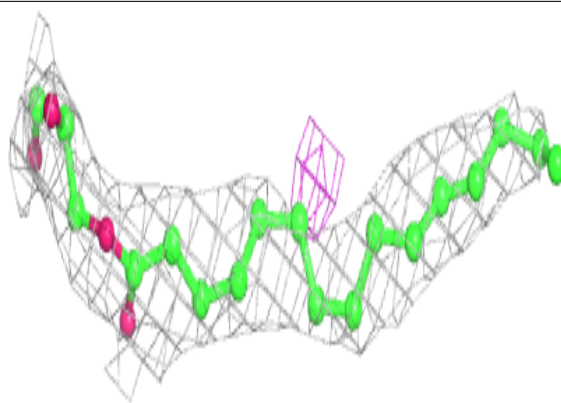
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
3	78N	A	503	22/22	0.79	0.29	28,49,60,68	0
3	78N	A	506	22/22	0.82	0.21	44,59,66,68	0
2	78M	A	501	22/22	0.82	0.26	32,44,62,74	0
2	78M	A	505	22/22	0.85	0.25	32,46,67,71	0
2	78M	A	502	22/22	0.87	0.35	38,63,74,78	0
2	78M	A	504	22/22	0.87	0.31	38,56,67,84	0
6	PHE	A	509	12/12	0.88	0.24	53,62,79,81	0
5	ALA	A	508	5/6	0.93	0.15	46,46,50,55	0
4	PO4	A	507	5/5	0.95	0.28	43,53,72,73	0

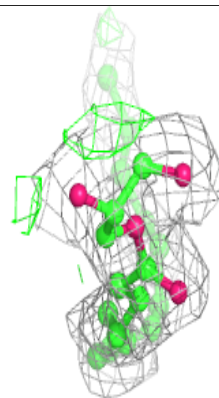
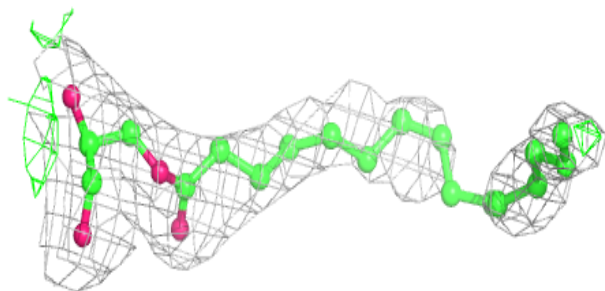
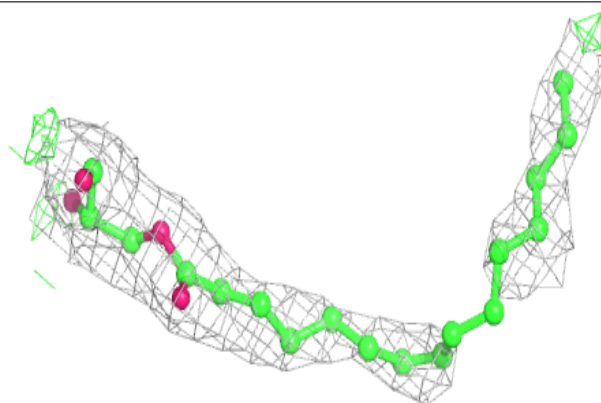
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

**Electron density around 78N A 503:**

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

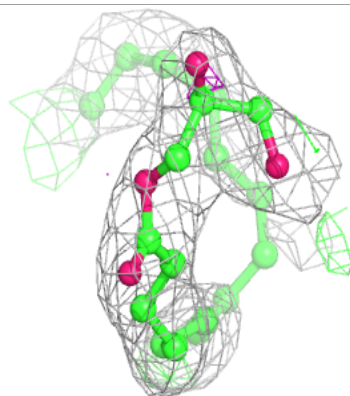
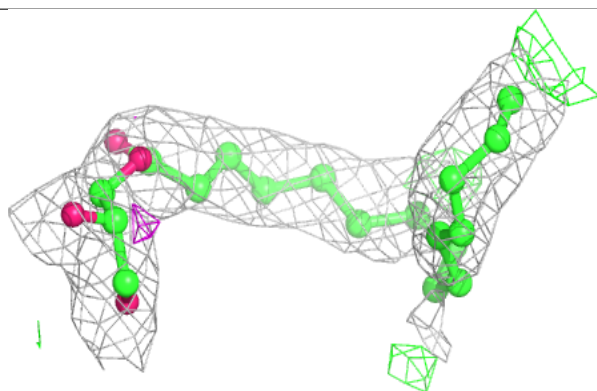
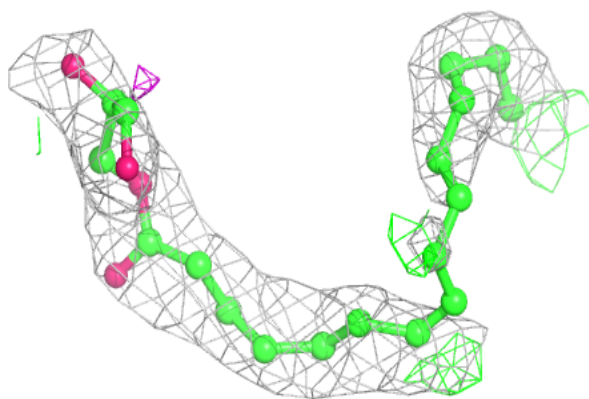
**Electron density around 78N A 506:**

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

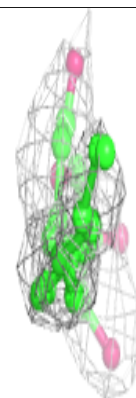
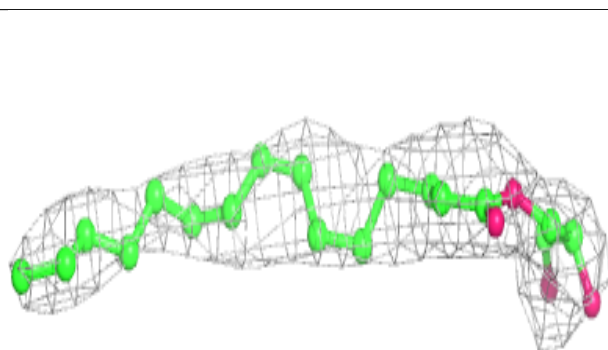
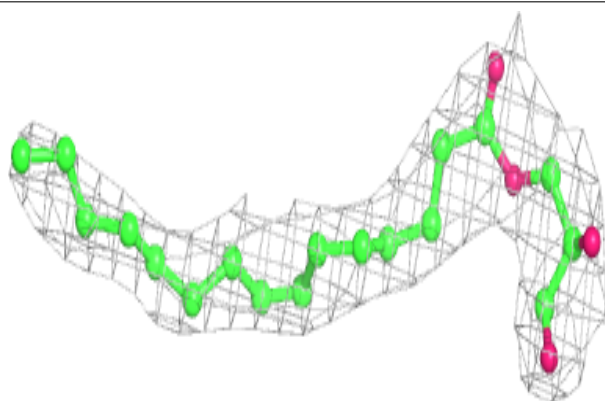


**Electron density around 78M A 501:**

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

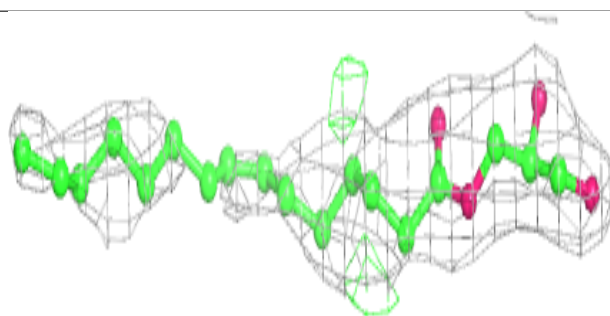
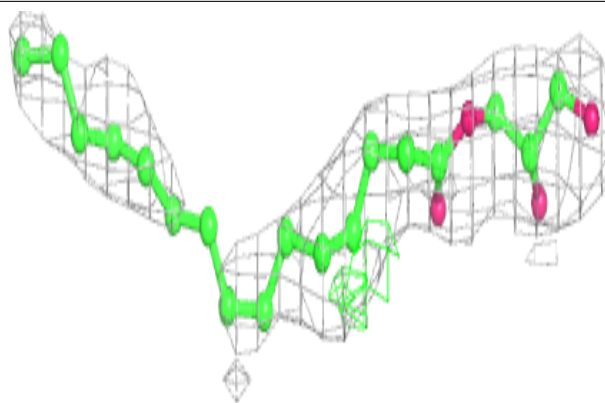
**Electron density around 78M A 505:**

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

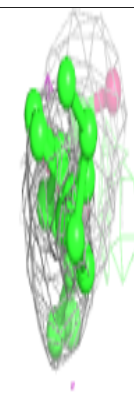
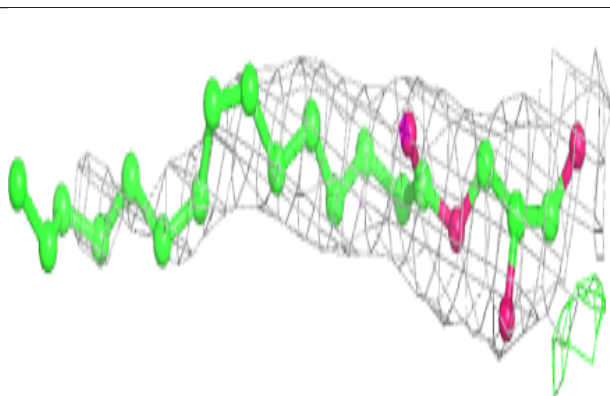
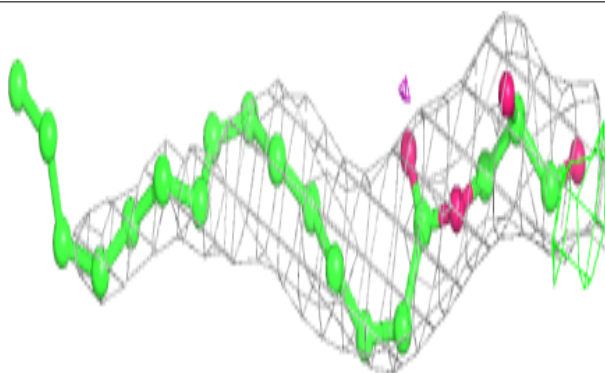


**Electron density around 78M A 502:**

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

**Electron density around 78M A 504:**

$2mF_o-DF_c$  (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.