



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 05:09 AM UTC

PDB ID : 4D2D / pdb\_00004d2d  
Title : Structure of a tri peptide bound POT family peptide transporter  
Authors : Lyons, J.A.; Parker, J.L.; Solcan, N.; Brinth, A.; Li, D.; Shah, S.T.A.; Caffrey, M.; Newstead, S.  
Deposited on : 2014-05-09  
Resolution : 2.52 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.

The types of validation reports are described at

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

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

MolProbity : 4-5-2 with Phenix2.0  
Mogul : 2022.3.0, CSD as543be (2022)  
Xtrriage (Phenix) : 2.0  
EDS : 3.0  
Buster-report : wwPDB partial adaption of 1.1.7 (2018)  
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)  
CCP4 : 9.0.010 (Gargrove)  
Density-Fitness : 1.0.12  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.49

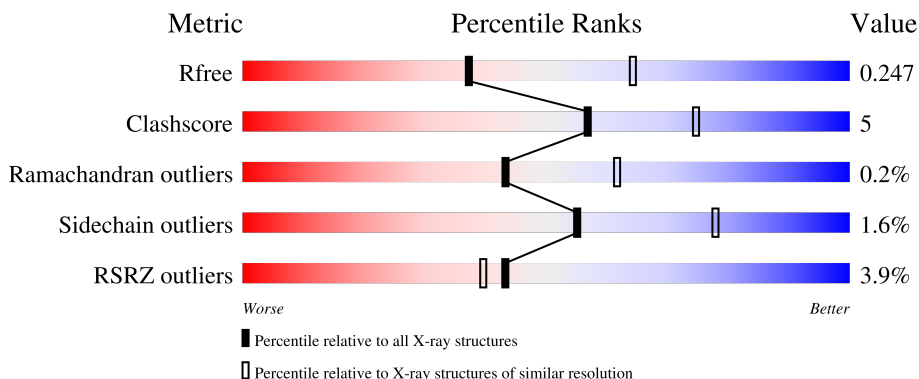
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.52 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	180053	7383 (2.54-2.50)
Clashscore	190562	8079 (2.54-2.50)
Ramachandran outliers	187476	7944 (2.54-2.50)
Sidechain outliers	187428	7946 (2.54-2.50)
RSRZ outliers	180081	7387 (2.54-2.50)

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	491	
2	B	3	

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 3780 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+ SYMPORTER.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	463	3582	2407	556	602	17	0	0	0

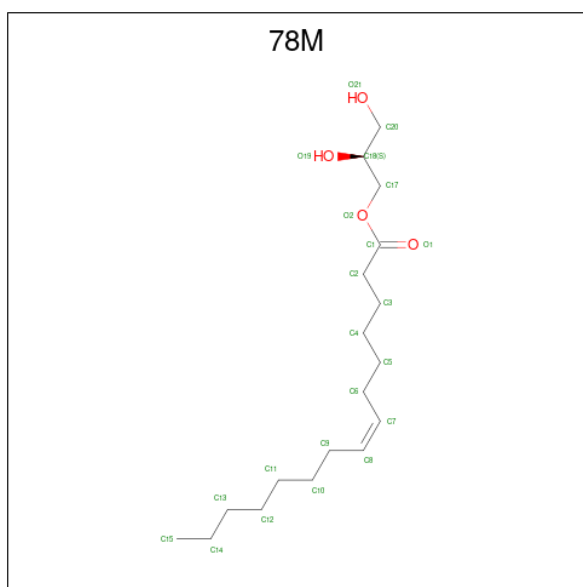
There are 8 discrepancies between the modelled and reference sequences:

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 a protein called ALANINE-TRIPETIDE.

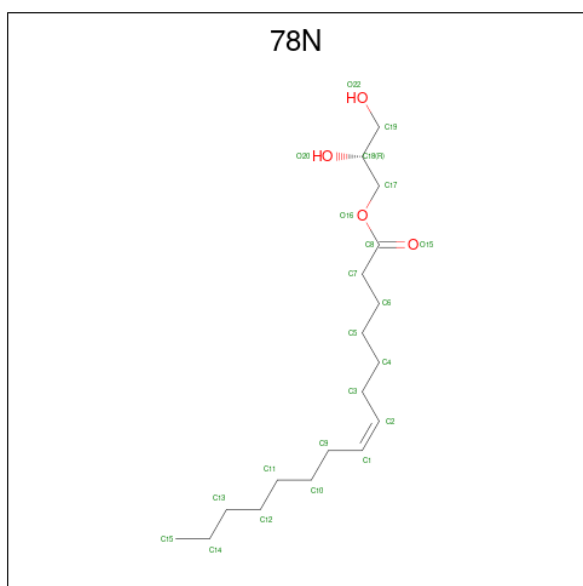
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
2	B	3	16	9	3	4	0	0	0

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



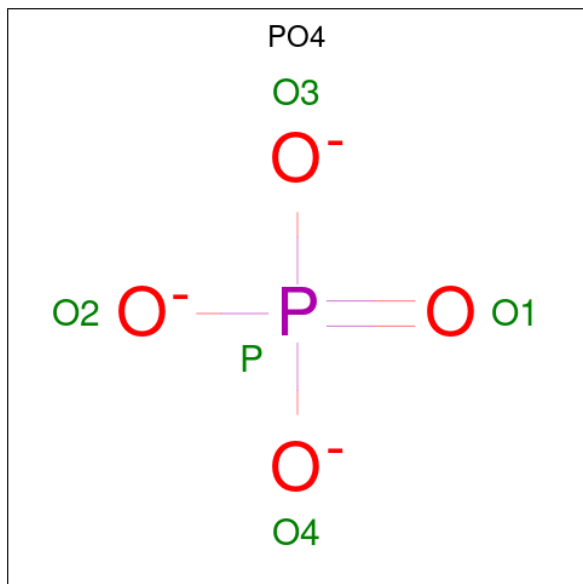
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	C O	0	0
			22	18 4		
3	A	1	Total	C O	0	0
			22	18 4		
3	A	1	Total	C O	0	0
			22	18 4		

- Molecule 4 is (2R)-2,3-DIHYDROXYPROPYL(7Z)-PENTADEC-7-ENOATE (CCD ID: 78N) (formula:  $C_{18}H_{34}O_4$ ).



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

- Molecule 5 is PHOSPHATE ION (CCD ID: PO4) (formula: O<sub>4</sub>P).



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

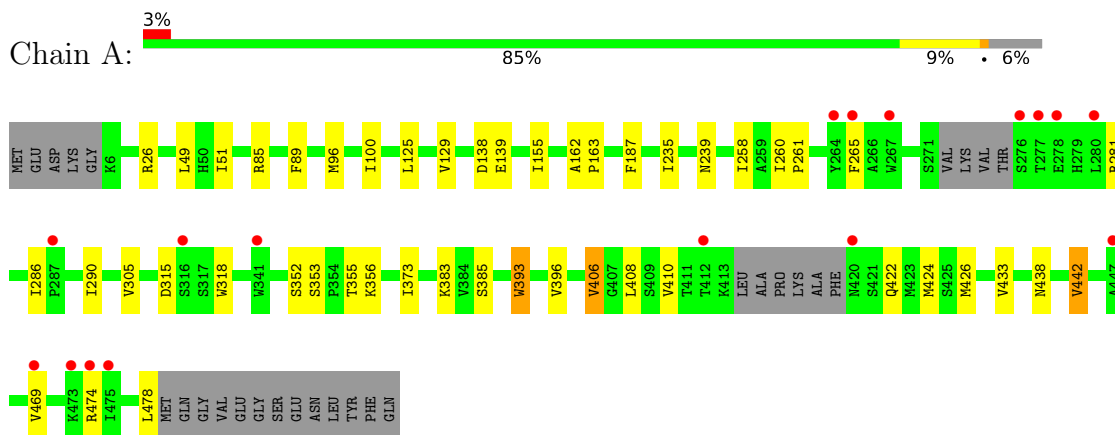
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	45	Total	O	0	0
			45	45		

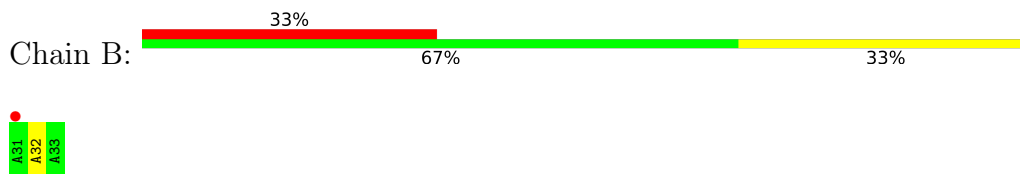
### 3 Residue-property plots [i](#)

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: DI-OR TRIPEPTIDE\ :H+ SYMPORTER



- Molecule 2: ALANINE-TRIPEPTIDE



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	103.38Å 110.70Å 110.62Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	75.56 – 2.52 75.56 – 2.52	Depositor EDS
% Data completeness (in resolution range)	98.2 (75.56-2.52) 98.2 (75.56-2.52)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.25 (at 2.51Å)	Xtrriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, $R_{free}$	0.198 , 0.242 0.205 , 0.247	Depositor DCC
$R_{free}$ test set	1102 reflections (5.16%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	44.3	Xtrriage
Anisotropy	0.383	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 45.7	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	3780	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	50.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.16% 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: 78N, PO4, 78M

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	0.34	0/3688	0.77	2/5028 (0.0%)
2	B	0.21	0/15	0.23	0/18
All	All	0.34	0/3703	0.77	2/5046 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	85	ARG	CA-C-N	-5.30	114.21	119.56
1	A	85	ARG	C-N-CA	-5.30	114.21	119.56

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	3582	0	3668	28	0
2	B	16	0	14	0	0
3	A	66	0	102	7	0
4	A	66	0	102	6	0
5	A	5	0	0	0	0
6	A	45	0	0	0	0
All	All	3780	0	3886	35	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 5.

All (35) 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:281:ARG:HG2	1:A:478:LEU:HB3	1.77	0.66
4:A:1481:78N:H121	4:A:1484:78N:H52C	1.80	0.64
1:A:353:SER:HB3	1:A:410:VAL:HG22	1.79	0.63
1:A:385:SER:HB3	3:A:1483:78M:H172	1.82	0.61
1:A:408:LEU:HD11	1:A:424:MET:HG3	1.87	0.57
1:A:49:LEU:HG	1:A:51:ILE:HG12	1.87	0.56
1:A:96:MET:O	1:A:100:ILE:HG12	2.06	0.56
1:A:438:ASN:O	1:A:442:VAL:HB	2.08	0.54
1:A:26:ARG:HG3	1:A:155:ILE:HG23	1.93	0.50
1:A:315:ASP:OD1	1:A:383:LYS:HB2	2.11	0.50
1:A:422:GLN:O	1:A:426:MET:HG3	2.12	0.49
1:A:265:PHE:CE2	1:A:286:ILE:HG12	2.47	0.49
1:A:89:PHE:CE2	3:A:1479:78M:H52C	2.47	0.49
1:A:474:ARG:HH11	1:A:474:ARG:HG3	1.77	0.49
1:A:352:SER:O	1:A:355:THR:OG1	2.31	0.48
1:A:406:VAL:O	1:A:410:VAL:HG23	2.14	0.48
1:A:260:ILE:HB	1:A:261:PRO:HD3	1.95	0.48
1:A:305:VAL:HG23	1:A:442:VAL:HG22	1.95	0.48
3:A:1482:78M:H91C	3:A:1482:78M:H61C	1.69	0.47
4:A:1480:78N:H92C	4:A:1480:78N:H31C	1.75	0.47
1:A:286:ILE:O	1:A:290:ILE:HG12	2.15	0.46
3:A:1483:78M:H51C	4:A:1484:78N:O15	2.16	0.46
1:A:261:PRO:HG2	1:A:433:VAL:HG21	1.99	0.45
3:A:1483:78M:H121	3:A:1483:78M:H92C	1.65	0.45
1:A:235:ILE:O	1:A:239:ASN:ND2	2.39	0.45
1:A:162:ALA:HB3	1:A:163:PRO:HD3	2.01	0.43
1:A:125:LEU:O	1:A:129:VAL:HG23	2.19	0.42
1:A:258:ILE:O	1:A:261:PRO:HD2	2.20	0.42
1:A:353:SER:HA	1:A:356:LYS:HD2	2.01	0.42
3:A:1483:78M:H22C	4:A:1484:78N:H191	2.01	0.42
1:A:187:PHE:CE1	3:A:1482:78M:H153	2.55	0.41
4:A:1480:78N:H152	4:A:1484:78N:H51C	2.02	0.41
1:A:318:TRP:CD1	1:A:318:TRP:H	2.38	0.41
1:A:393:TRP:HA	1:A:396:VAL:HG12	2.02	0.41
1:A:373:ILE:HD11	4:A:1480:78N:H91C	2.03	0.41

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	457/491 (93%)	450 (98%)	7 (2%)	0	100	100
2	B	1/3 (33%)	0	0	1 (100%)	0	0
All	All	458/494 (93%)	450 (98%)	7 (2%)	1 (0%)	43	62

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	32	ALA

### 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	375/398 (94%)	369 (98%)	6 (2%)	55	78

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	138	ASP
1	A	139	GLU
1	A	393	TRP
1	A	406	VAL
1	A	442	VAL
1	A	469	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such

sidechains are listed below:

Mol	Chain	Res	Type
1	A	279	HIS
1	A	328	ASN
1	A	349	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 oligosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

7 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
4	78N	A	1480	-	21,21,21	0.91	1 (4%)	22,22,22	1.10	2 (9%)
4	78N	A	1481	-	21,21,21	0.91	1 (4%)	22,22,22	1.19	2 (9%)
3	78M	A	1483	-	21,21,21	0.92	1 (4%)	22,22,22	1.17	2 (9%)
3	78M	A	1482	-	21,21,21	0.91	1 (4%)	22,22,22	1.15	2 (9%)
4	78N	A	1484	-	21,21,21	0.90	1 (4%)	22,22,22	1.23	2 (9%)
3	78M	A	1479	-	21,21,21	0.91	1 (4%)	22,22,22	1.21	2 (9%)
5	PO4	A	1485	-	4,4,4	0.95	0	6,6,6	0.49	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
4	78N	A	1480	-	-	9/21/21/21	-
4	78N	A	1481	-	-	5/21/21/21	-
3	78M	A	1483	-	-	12/21/21/21	-
3	78M	A	1482	-	-	13/21/21/21	-
4	78N	A	1484	-	-	9/21/21/21	-
3	78M	A	1479	-	-	11/21/21/21	-

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1480	78N	O16-C17	-2.67	1.39	1.45
4	A	1481	78N	O16-C17	-2.66	1.39	1.45
3	A	1479	78M	O2-C17	-2.63	1.39	1.45
3	A	1483	78M	O2-C17	-2.60	1.39	1.45
3	A	1482	78M	O2-C17	-2.58	1.39	1.45
4	A	1484	78N	O16-C17	-2.57	1.39	1.45

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1484	78N	O16-C8-C7	3.10	121.29	111.83
3	A	1479	78M	O2-C17-C18	2.98	119.87	105.85
4	A	1481	78N	O16-C8-C7	2.93	120.77	111.83
3	A	1483	78M	O2-C1-C2	2.78	120.32	111.83
4	A	1484	78N	O16-C17-C18	2.76	118.85	105.85
3	A	1483	78M	O2-C17-C18	2.74	118.71	105.85
3	A	1479	78M	O2-C1-C2	2.72	120.12	111.83
4	A	1480	78N	O16-C8-C7	2.71	120.09	111.83
4	A	1481	78N	O16-C17-C18	2.67	118.42	105.85
3	A	1482	78M	O2-C17-C18	2.63	118.22	105.85
3	A	1482	78M	O2-C1-C2	2.50	119.46	111.83
4	A	1480	78N	O16-C17-C18	2.44	117.31	105.85

There are no chirality outliers.

All (59) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1479	78M	O2-C17-C18-C20
3	A	1479	78M	O2-C17-C18-O19
3	A	1482	78M	C17-C18-C20-O21
3	A	1482	78M	O2-C17-C18-C20
4	A	1484	78N	O16-C17-C18-O20
3	A	1483	78M	O1-C1-O2-C17
3	A	1483	78M	C2-C1-O2-C17
3	A	1483	78M	C9-C10-C11-C12
4	A	1481	78N	O16-C17-C18-C19
3	A	1482	78M	O2-C17-C18-O19
3	A	1482	78M	O19-C18-C20-O21
4	A	1484	78N	C5-C6-C7-C8
4	A	1481	78N	O16-C17-C18-O20
4	A	1480	78N	C7-C8-O16-C17
4	A	1484	78N	C11-C12-C13-C14
3	A	1483	78M	C17-C18-C20-O21
3	A	1483	78M	O19-C18-C20-O21
3	A	1483	78M	C3-C4-C5-C6
4	A	1480	78N	C4-C5-C6-C7
4	A	1480	78N	C10-C11-C12-C13
4	A	1480	78N	O15-C8-O16-C17
3	A	1479	78M	C11-C12-C13-C14
4	A	1484	78N	C9-C10-C11-C12
4	A	1481	78N	C5-C6-C7-C8
4	A	1480	78N	C3-C4-C5-C6
4	A	1484	78N	O16-C17-C18-C19
4	A	1484	78N	C10-C11-C12-C13
4	A	1480	78N	C12-C13-C14-C15
3	A	1483	78M	C2-C3-C4-C5
3	A	1479	78M	C11-C10-C9-C8
3	A	1482	78M	C11-C10-C9-C8
3	A	1483	78M	C7-C8-C9-C10
3	A	1482	78M	C18-C17-O2-C1
3	A	1483	78M	C11-C10-C9-C8
3	A	1482	78M	C4-C5-C6-C7
3	A	1482	78M	C1-C2-C3-C4
3	A	1479	78M	C10-C11-C12-C13
4	A	1481	78N	C3-C4-C5-C6
4	A	1480	78N	C2-C3-C4-C5
3	A	1482	78M	C9-C10-C11-C12
3	A	1482	78M	C10-C11-C12-C13
4	A	1480	78N	C11-C10-C9-C1
3	A	1482	78M	C2-C3-C4-C5

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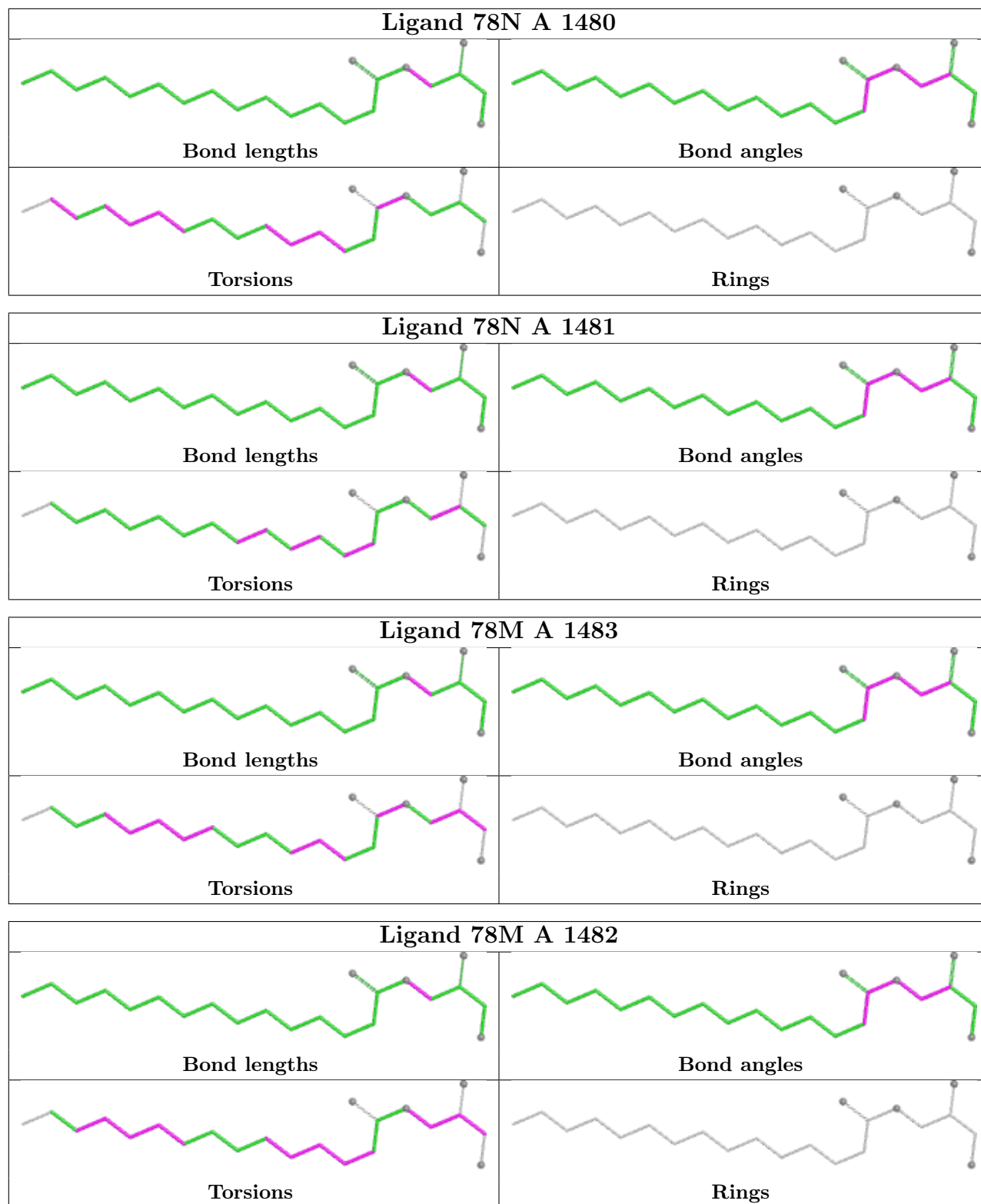
Mol	Chain	Res	Type	Atoms
3	A	1482	78M	C3-C4-C5-C6
3	A	1483	78M	O2-C17-C18-C20
3	A	1479	78M	C3-C4-C5-C6
3	A	1483	78M	C10-C11-C12-C13
4	A	1480	78N	C9-C10-C11-C12
3	A	1479	78M	O2-C1-C2-C3
3	A	1479	78M	C12-C13-C14-C15
4	A	1481	78N	C1-C2-C3-C4
3	A	1483	78M	O2-C17-C18-O19
4	A	1484	78N	C3-C4-C5-C6
3	A	1482	78M	C11-C12-C13-C14
3	A	1479	78M	C4-C5-C6-C7
3	A	1479	78M	C5-C6-C7-C8
4	A	1484	78N	C1-C2-C3-C4
4	A	1484	78N	O20-C18-C19-O22
3	A	1479	78M	C7-C8-C9-C10

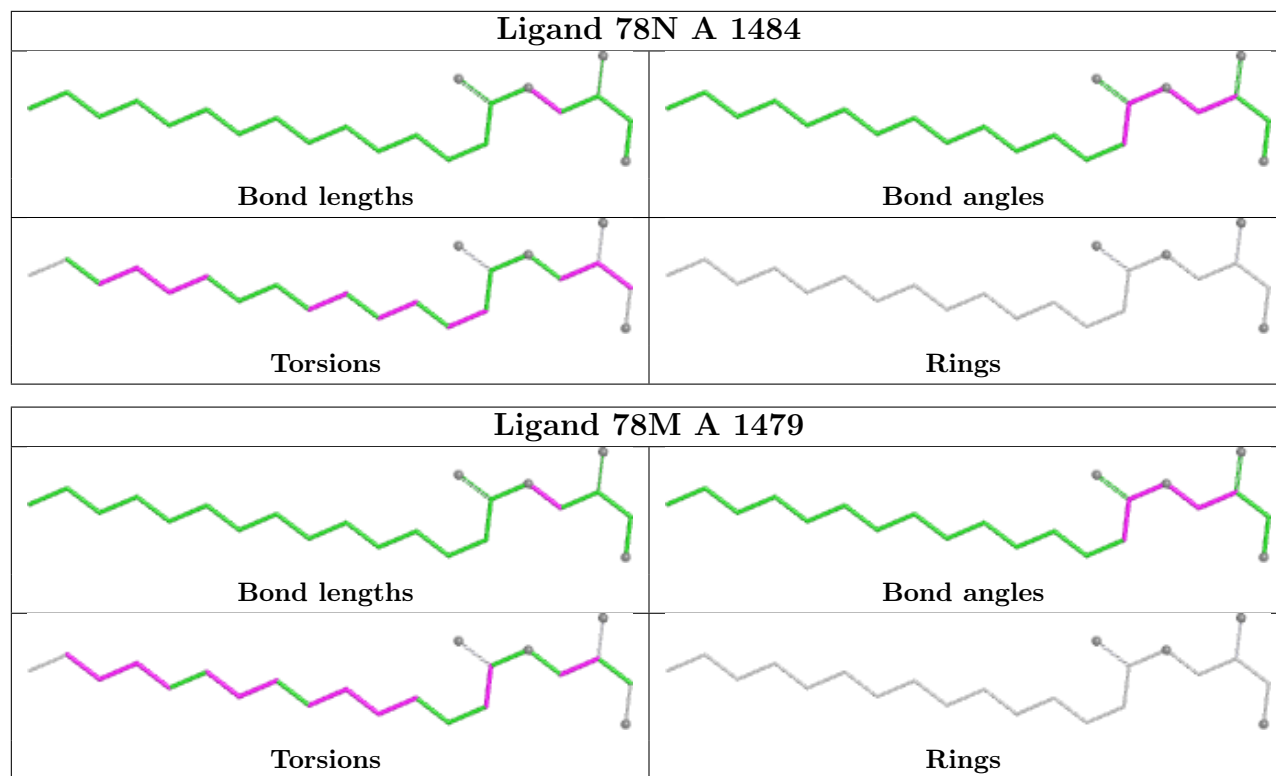
There are no ring outliers.

6 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1480	78N	3	0
4	A	1481	78N	1	0
3	A	1483	78M	4	0
3	A	1482	78M	2	0
4	A	1484	78N	4	0
3	A	1479	78M	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 [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	463/491 (94%)	0.32	17 (3%) 45 41	32, 44, 79, 101	0
2	B	3/3 (100%)	2.16	1 (33%) 1 1	63, 63, 75, 82	0
All	All	466/494 (94%)	0.33	18 (3%) 43 39	32, 44, 79, 101	0

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	31	ALA	3.4
1	A	276	SER	3.1
1	A	277	THR	3.0
1	A	267	TRP	2.9
1	A	316	SER	2.8
1	A	473	LYS	2.8
1	A	287	PRO	2.5
1	A	280	LEU	2.5
1	A	469	VAL	2.5
1	A	264	TYR	2.3
1	A	475	ILE	2.2
1	A	474	ARG	2.2
1	A	278	GLU	2.1
1	A	447	ALA	2.1
1	A	341	TRP	2.1
1	A	420	ASN	2.1
1	A	265	PHE	2.0
1	A	412	THR	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](#)

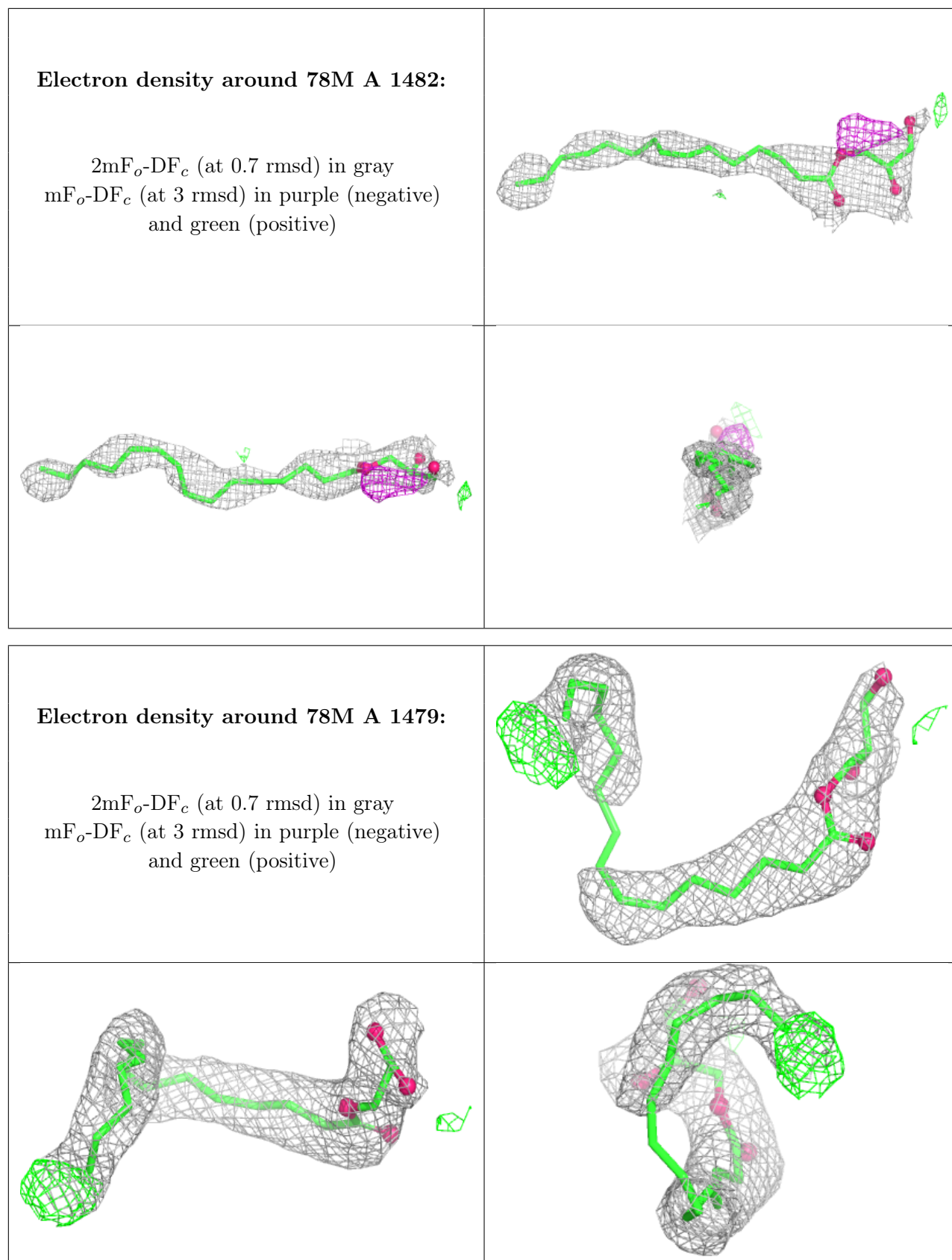
There are no oligosaccharides in this entry.

### 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 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.

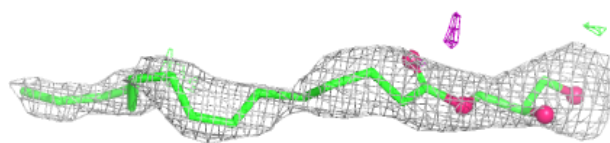
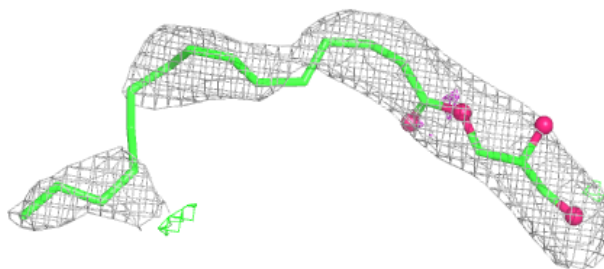
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
3	78M	A	1482	22/22	0.83	0.18	44,61,72,84	0
3	78M	A	1479	22/22	0.84	0.19	33,55,66,83	0
4	78N	A	1484	22/22	0.85	0.19	56,66,77,80	0
4	78N	A	1481	22/22	0.88	0.16	49,65,74,83	0
4	78N	A	1480	22/22	0.88	0.14	42,58,65,79	0
3	78M	A	1483	22/22	0.89	0.13	43,56,67,80	0
5	PO4	A	1485	5/5	0.92	0.12	58,67,74,75	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.

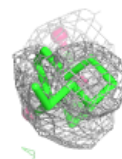
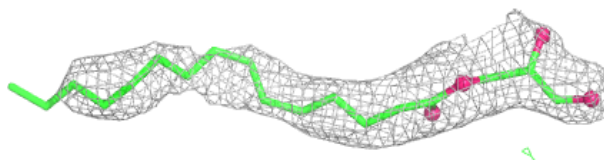


**Electron density around 78N A 1484:**

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

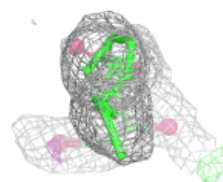
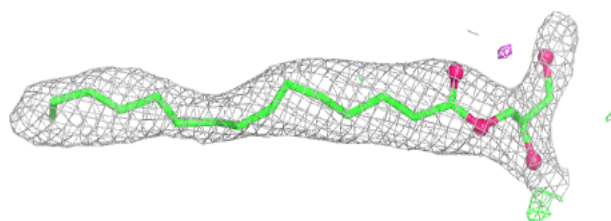
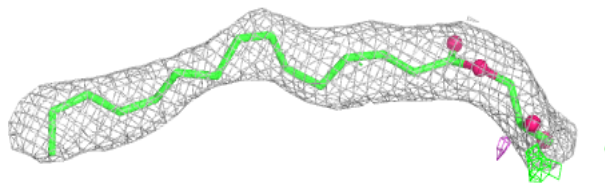
**Electron density around 78N A 1481:**

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

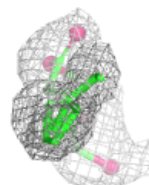
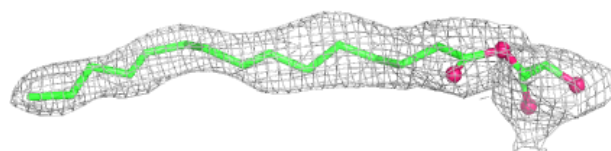
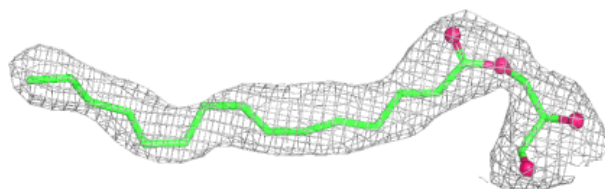


**Electron density around 78N A 1480:**

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

**Electron density around 78M A 1483:**

$2mF_o-DF_c$  (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.