

Full wwPDB X-ray Structure Validation Report (i)

Oct 23, 2021 – 03:08 PM EDT

PDB ID	:	7PRN
Title	:	E1M, D97A, E99A MUTANT OF RH. BLASTICA PORIN
Authors	:	Maveyraud, L.; Schmid, B.; Schulz, G.E.
Deposited on	:	1998-06-12
Resolution	:	2.25 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* A user guide is available at https://www.wwpdb.org/validation/2017/XrayValidationReportHelp with specific help available everywhere you see the (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

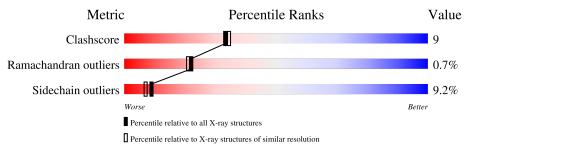
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.23.2
buster-report	:	1.1.7(2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.23.2

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: $X\text{-}RAY \, DIFFRACTION$

The reported resolution of this entry is 2.25 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	$egin{array}{c} { m Whole \ archive} \ (\#{ m Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries,\ resolution\ range}({ m \AA}))$
Clashscore	141614	1487 (2.26-2.26)
Ramachandran outliers	138981	1449 (2.26-2.26)
Sidechain outliers	138945	1450 (2.26-2.26)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=5%

Mol	Chain	Length	Quality of chain		
1	А	289	74%	22%	••



7PRN

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 2238 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.

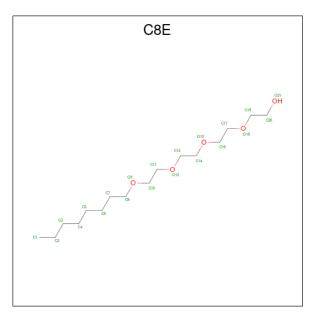
• Molecule 1 is a protein called PORIN.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace	
1	А	289	Total 2156	C 1353	N 346	0 453	${f S}$ 4	0	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
А	97	ALA	ASP	engineered mutation	UNP P39767
А	99	ALA	GLU	engineered mutation	UNP P39767

• Molecule 2 is (HYDROXYETHYLOXY)TRI(ETHYLOXY)OCTANE (three-letter code: C8E) (formula: $C_{16}H_{34}O_5$).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf	
2	А	1	Total 21	C 16	O 5	0	0

• Molecule 3 is water.

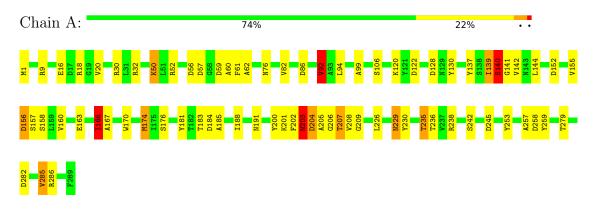


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	А	61	$\begin{array}{cc} \text{Total} & \text{O} \\ 61 & 61 \end{array}$	0	0



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. 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. 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: PORIN



4 Data and refinement statistics (i)

Property	Value	Source
Space group	H 3	Depositor
Cell constants	104.40Å 104.40Å 124.25Å	Depositor
a, b, c, α , β , γ	90.00° 90.00° 120.00°	-
Resolution (Å)	32.90 - 2.25	Depositor
	32.95 - 2.25	EDS
% Data completeness	99.8 (32.90-2.25)	Depositor
(in resolution range)	99.7 (32.95-2.25)	EDS
R _{merge}	0.04	Depositor
R_{sym}	0.04	Depositor
$< I/\sigma(I) > 1$	9.80 (at 2.24Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.215 , 0.222	Depositor
	0.217 , (Not available)	DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor $(Å^2)$	29.7	Xtriage
Anisotropy	0.076	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.36 , 42.5	EDS
L-test for $twinning^2$	$< L >=0.40, < L^2>=0.23$	Xtriage
Estimated twinning fraction	$\begin{array}{c} 0.084 \ {\rm for} \ -2/3 ^{\rm *h-1}/3 ^{\rm *k} + 2/3 ^{\rm *l}, -1/3 ^{\rm *h-2}/3 ^{\rm *k} + 2/3 ^{\rm *l}, 2/3 ^{\rm *h-2}/3 ^{\rm *k} + 1/3 ^{\rm *l} \\ 0.069 \ {\rm for} \ -h, 1/3 ^{\rm *h-1}/3 ^{\rm *k} + 2/3 ^{\rm *l}, 2/3 ^{\rm *h} + 4/3 ^{\rm *k} \\ {\rm k} + 1/3 ^{\rm *l} \\ 0.075 \ {\rm for} \ -1/3 ^{\rm *h+1}/3 ^{\rm *k} - 2/3 ^{\rm *l}, -{\rm k}, -4/3 ^{\rm *h-2}/3 \\ {\rm *k} + 1/3 ^{\rm *l} \\ 0.069 \ {\rm for} \ -h, 2/3 ^{\rm *h} + 1/3 ^{\rm *k} - 2/3 ^{\rm *l}, -2/3 ^{\rm *h-4}/3 ^{\rm *l} \\ {\rm 0.077 \ for} \ 1/3 ^{\rm *h+2}/3 ^{\rm *k} + 2/3 ^{\rm *l}, -2/3 ^{\rm *h-4}/3 ^{\rm *h+2}/3 \\ {\rm *k} - 1/3 ^{\rm *l} \\ 0.079 \ {\rm for} \ -1/3 ^{\rm *h-2}/3 ^{\rm *k} - 2/3 ^{\rm *l}, -2/3 ^{\rm *h-1}/3 ^{\rm *k} + 2/3 ^{\rm *l}, -2/3 ^{\rm *h-2}/3 ^{\rm *k}, -1/3 ^{\rm *l} \\ 0.315 \ {\rm for} \ h, -h \cdot {\rm k}, -1 \end{array}$	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	2238	wwPDB-VP
Average B, all atoms $(Å^2)$	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 5.69% of the height of the origin peak. No significant pseudotranslation is detected.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.



 $^{^1 \}mathrm{Intensities}$ estimated from amplitudes.

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: $\rm C8E$

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).

Mal	Chain	Bo	nd lengths	Bo	ond angles
Mol	Chain	RMSZ	# Z > 5	RMSZ	# Z > 5
1	А	1.10	1/2199~(0.0%)	1.58	27/2994~(0.9%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	А	242	SER	CB-OG	5.43	1.49	1.42

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
1	А	9	ARG	NE-CZ-NH2	-13.85	113.38	120.30
1	А	286	ARG	NE-CZ-NH2	7.44	124.02	120.30
1	А	92	VAL	CB-CA-C	-7.29	97.55	111.40
1	А	52	ARG	NE-CZ-NH2	7.23	123.92	120.30
1	А	57	ASP	CB-CG-OD2	-7.21	111.81	118.30
1	А	56	ASP	CB-CG-OD2	7.05	124.64	118.30
1	А	152	ASP	CB-CG-OD1	7.02	124.62	118.30
1	А	32	ARG	NE-CZ-NH2	6.87	123.74	120.30
1	А	18	ARG	NE-CZ-NH2	6.79	123.69	120.30
1	А	230	TYR	CB-CG-CD2	-6.77	116.94	121.00
1	А	30	ARG	NE-CZ-NH1	6.57	123.59	120.30
1	А	30	ARG	NE-CZ-NH2	-6.57	117.02	120.30
1	А	184	ASP	CB-CG-OD1	6.08	123.78	118.30
1	А	230	TYR	CB-CG-CD1	5.95	124.57	121.00
1	А	238	ARG	NE-CZ-NH2	5.92	123.26	120.30
1	А	259	TYR	CB-CG-CD1	-5.76	117.55	121.00
1	А	166	ILE	CB-CA-C	-5.68	100.24	111.60
1	А	245	ASP	CB-CG-OD1	5.62	123.36	118.30
1	А	128	ASP	CB-CG-OD1	5.42	123.18	118.30
1	А	174	MET	CG-SD-CE	5.37	108.80	100.20

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Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
1	А	130	TYR	CA-CB-CG	5.35	123.57	113.40
1	А	253	TYR	CB-CG-CD2	-5.35	117.79	121.00
1	А	285	VAL	CB-CA-C	-5.30	101.33	111.40
1	А	137	TYR	CB-CG-CD2	-5.29	117.82	121.00
1	А	152	ASP	CB-CG-OD2	-5.09	113.72	118.30
1	А	82	VAL	CB-CA-C	-5.06	101.78	111.40
1	А	156	ASP	CB-CG-OD1	5.05	122.84	118.30

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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	А	2156	0	1984	36	0
2	А	21	0	34	3	0
3	А	61	0	0	3	0
All	All	2238	0	2018	36	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 9.

All (36) 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:203:ASN:HD22	1:A:204:ASP:H	1.06	1.00
1:A:92:VAL:HG13	1:A:167:ALA:HB1	1.48	0.92
1:A:139:ILE:HG23	1:A:142:VAL:HB	1.55	0.88
1:A:92:VAL:CG1	1:A:167:ALA:HB1	2.11	0.80
1:A:92:VAL:HG13	1:A:167:ALA:CB	2.22	0.68
1:A:201:LYS:HG3	1:A:203:ASN:O	1.96	0.66
1:A:142:VAL:HG22	1:A:170:TRP:HD1	1.63	0.64
1:A:61:PHE:HZ	2:A:290:C8E:H142	1.62	0.64
1:A:203:ASN:ND2	1:A:204:ASP:H	1.88	0.63
1:A:207:THR:HG23	1:A:229:ASN:HD21	1.63	0.63

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		Interatomic	Clash
Atom-1	Atom-2	distance (\AA)	overlap (Å)
1:A:181:TYR:CE2	1:A:183:THR:HG23	2.33	0.63
1:A:203:ASN:HD22	1:A:204:ASP:N	1.90	0.58
1:A:188:ILE:HG22	1:A:191:ASN:HB2	1.89	0.54
1:A:86:ASP:HA	3:A:298:HOH:O	2.09	0.53
1:A:99:ALA:HA	1:A:106:SER:HB2	1.91	0.52
1:A:207:THR:HG23	1:A:229:ASN:ND2	2.24	0.52
1:A:61:PHE:CZ	2:A:290:C8E:H142	2.45	0.51
1:A:50:LYS:NZ	3:A:327:HOH:O	2.44	0.50
1:A:209:GLY:O	1:A:226:LEU:HD12	2.13	0.48
1:A:155:VAL:O	1:A:156:ASP:C	2.52	0.48
1:A:200:TYR:CE2	1:A:202:PHE:HB3	2.50	0.46
1:A:202:PHE:CE1	1:A:208:VAL:HG13	2.50	0.46
1:A:163:GLU:HG3	1:A:185:ALA:HA	1.97	0.45
1:A:16:GLU:HB2	1:A:279:THR:HB	1.98	0.45
1:A:166:ILE:HD12	1:A:181:TYR:HB3	1.98	0.45
1:A:181:TYR:HE2	1:A:183:THR:HG23	1.82	0.44
1:A:59:ASP:OD2	1:A:120:LYS:NZ	2.50	0.43
1:A:94:LEU:HD11	1:A:176:SER:HB2	2.01	0.43
1:A:140:SER:HB3	1:A:141:GLY:H	1.77	0.42
1:A:60:ALA:CB	2:A:290:C8E:H172	2.50	0.42
1:A:122:ASP:OD1	1:A:122:ASP:C	2.58	0.42
1:A:236:THR:O	1:A:257:ALA:HA	2.20	0.41
1:A:62:ALA:HA	3:A:306:HOH:O	2.20	0.41
1:A:203:ASN:HB3	1:A:205:ALA:H	1.86	0.41
1:A:203:ASN:HB3	1:A:206:GLY:H	1.86	0.41
1:A:235:THR:HA	1:A:258:ASP:O	2.22	0.40

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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	А	287/289~(99%)	274 (96%)	11 (4%)	2(1%)	22 21	

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	А	140	SER
1	А	203	ASN

5.3.2 Protein sidechains (i)

In the following table, the Percentiles column shows the percent side chain 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.

Ν	Лоl	Chain	Analysed	Rotameric	Outliers	Percentiles
	1	А	217/217~(100%)	197~(91%)	20~(9%)	9 7

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	А	1	MET
1	А	20	VAL
1	А	50	LYS
1	А	76	ASN
1	А	92	VAL
1	А	139	ILE
1	А	140	SER
1	А	144	LEU
1	А	157	SER
1	А	158	SER
1	А	160	VAL
1	А	166	ILE
1	А	174	MET
1	А	203	ASN
1	А	204	ASP
1	А	207	THR
1	А	229	ASN
1	А	235	THR
1	А	282	ASP

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Mol	Chain	Res	Type
1	А	285	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	А	118	ASN
1	А	203	ASN
1	А	229	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)

1 ligand is 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	Dec	Tink	Bo	ond leng	ths	В	ond ang	les
IVIOI	Type	Unam	Res	Link	Counts	RMSZ	# Z >2	Counts	RMSZ	# Z >2
2	C8E	А	290	-	20,20,20	0.93	0	$19,\!19,\!19$	2.18	6 (31%)

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	C8E	А	290	-	-	11/18/18/18	-

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Ζ	$Observed(^{o})$	$Ideal(^{o})$
2	А	290	C8E	O15-C14-C13	4.90	132.47	110.39
2	А	290	C8E	O18-C17-C16	4.20	129.34	110.39
2	А	290	C8E	O18-C19-C20	3.68	126.25	110.07
2	А	290	C8E	C7-C6-C5	-3.32	97.55	114.42
2	А	290	C8E	O12-C13-C14	2.21	120.35	110.39
2	А	290	C8E	O15-C16-C17	2.17	120.17	110.39

There are no chirality outliers.

Mol	Chain	Res	Type	Atoms
2	А	290	C8E	C16-C17-O18-C19
2	А	290	C8E	O12-C13-C14-O15
2	А	290	C8E	C3-C4-C5-C6
2	А	290	C8E	C2-C3-C4-C5
2	А	290	C8E	C1-C2-C3-C4
2	А	290	C8E	C20-C19-O18-C17
2	А	290	C8E	C7-C8-O9-C10
2	А	290	C8E	O18-C19-C20-O21
2	А	290	C8E	C14-C13-O12-C11
2	А	290	C8E	O15-C16-C17-O18
2	А	290	C8E	C4-C5-C6-C7

All (11) torsion outliers are listed below:

There are no ring outliers.

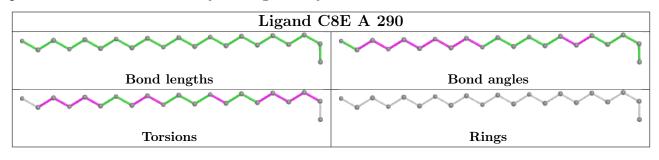
1 monomer is involved in 3 short contacts:

Mo	Chain	Res	Type	Clashes	Symm-Clashes
2	А	290	C8E	3	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 then 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)

Unable to reproduce the depositors R factor - this section is therefore empty.

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

Unable to reproduce the depositors R factor - this section is therefore empty.

6.3 Carbohydrates (i)

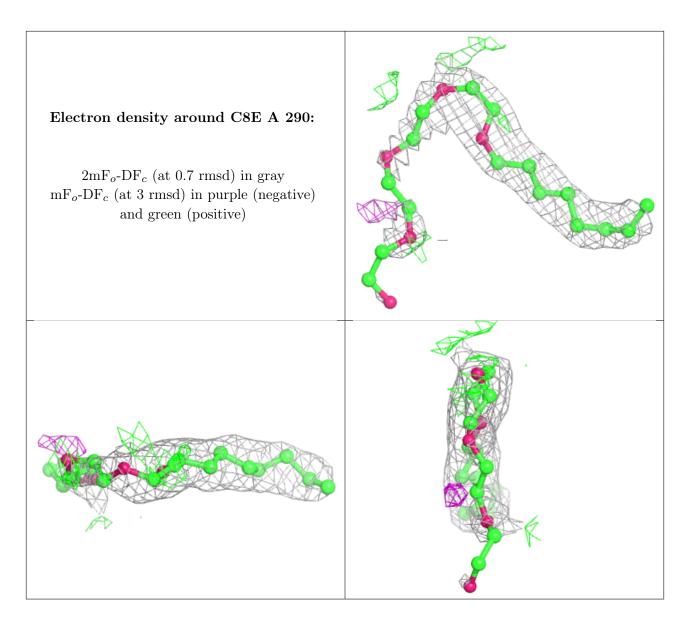
Unable to reproduce the depositors R factor - this section is therefore empty.

6.4 Ligands (i)

Unable to reproduce the depositors R factor - this section is therefore empty.

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.





6.5 Other polymers (i)

Unable to reproduce the depositors R factor - this section is therefore empty.

