

Full wwPDB X-ray Structure Validation Report (i)

Oct 23, 2021 – 02:12 PM EDT

PDB ID : 6PRN

Title : E1M, K50A, R52A MUTANT OF RH. BLASTICA PORIN

Authors: Maveyraud, L.; Schmid, B.; Schulz, G.E.

Deposited on : 1998-06-12

Resolution : 2.04 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)

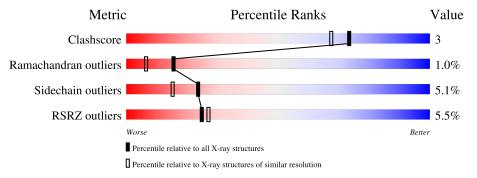
 $\begin{tabular}{lll} Validation Pipeline (wwPDB-VP) & : & 2.23.2 \end{tabular}$

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: X-RAY DIFFRACTION

The reported resolution of this entry is 2.04 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	Similar resolution
	$(\# \mathrm{Entries})$	$(\# ext{Entries}, ext{ resolution range}(ext{Å}))$
Clashscore	141614	1773 (2.04-2.04)
Ramachandran outliers	138981	1752 (2.04-2.04)
Sidechain outliers	138945	1752 (2.04-2.04)
RSRZ outliers	127900	1672 (2.04-2.04)

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% 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		
			6%		
1	A	289	91%	7%	•



2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 2387 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.

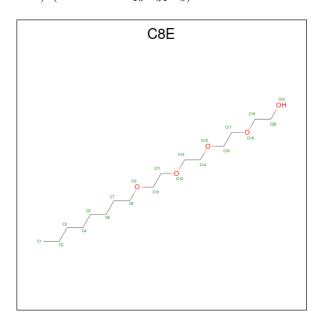
• Molecule 1 is a protein called PORIN.

Mol	Chain	Residues		Ato	oms			ZeroOcc	AltConf	Trace
1	A	289	Total 2167	C 1362	N 342	O 458	S 5	0	4	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	50	ALA	LYS	engineered mutation	UNP P39767
A	52	ALA	ARG	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	A	1	Total C O 21 16 5	0	0
2	A	1	Total C O 21 16 5	0	0



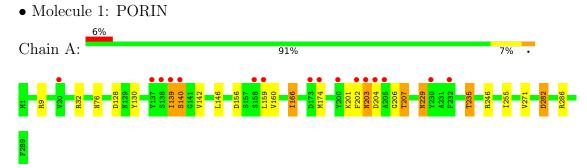
• Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	178	Total 178	O 178	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.





4 Data and refinement statistics (i)

Property	Value	Source	
Space group	H 3	Depositor	
Cell constants	104.36Å 104.36Å 124.25Å	Donositon	
a, b, c, α , β , γ	90.00° 90.00° 120.00°	Depositor	
Resolution (Å)	19.68 - 2.04	Depositor	
, ,	19.51 - 2.04	EDS	
% Data completeness	99.8 (19.68-2.04)	Depositor	
(in resolution range)	99.8 (19.51-2.04)	EDS	
R_{merge}	0.04	Depositor	
R_{sym}	0.04	Depositor	
$< I/\sigma(I) > 1$	4.15 (at 2.04Å)	Xtriage	
Refinement program	REFMAC	Depositor	
R, R_{free}	0.163 , 0.186	Depositor	
	0.165 , (Not available)	DCC	
R_{free} test set	No test flags present.	wwPDB-VP	
Wilson B-factor (\mathring{A}^2)	29.4	Xtriage	
Anisotropy	0.182	Xtriage	
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.35 , 71.8	EDS	
L-test for twinning ²	$< L > = 0.50, < L^2> = 0.34$	Xtriage	
Estimated twinning fraction	$\begin{array}{c} 0.004 \; \text{for} \; -2/3*\text{h-}1/3*\text{k+}2/3*\text{l}, -1/3*\text{h-}2/3*\text{k-}\\ 2/3*\text{l}, 2/3*\text{h-}2/3*\text{k+}1/3*\text{l}\\ 0.000 \; \text{for} \; -\text{h}, 1/3*\text{h-}1/3*\text{k+}2/3*\text{l}, 2/3*\text{h+}4/3*\\ & & & & & & & & & & & \\ k+1/3*\text{l}\\ 0.000 \; \text{for} \; -\text{l}/3*\text{h+}1/3*\text{k-}2/3*\text{l}, -\text{k}, -4/3*\text{h-}2/3\\ & & & & & & & & \\ k+1/3*\text{l}\\ 0.000 \; \text{for} \; -\text{h}, 2/3*\text{h+}1/3*\text{k-}2/3*\text{l}, -2/3*\text{h-}4/3*\\ & & & & & & & \\ k-1/3*\text{l}\\ 0.000 \; \text{for} \; 1/3*\text{h+}2/3*\text{k+}2/3*\text{l}, -\text{k}, 4/3*\text{h+}2/3\\ & & & & & & \\ *\text{k-}1/3*\text{l}\\ 0.013 \; \text{for} \; -\text{l}/3*\text{h-}2/3*\text{k-}2/3*\text{l}, -2/3*\text{h-}1/3*\text{k+}\\ & & & & & \\ 2/3*\text{l}, -2/3*\text{h+}2/3*\text{k-}1/3*\text{l}\\ & & & & & \\ 0.023 \; \text{for} \; \text{h}, -\text{h-k}, -\text{l} \end{array}$	Xtriage	
F_o, F_c correlation	0.96	EDS	
Total number of atoms	2387	wwPDB-VP	
Average B, all atoms (Å ²)	34.0	wwPDB-VP	

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 5.48% of the height of the origin peak. No significant pseudotranslation is detected.

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



¹Intensities 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: 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	Bond	lengths	Bond angles	
IVIOI	Chain	RMSZ	# Z > 5	RMSZ	# Z > 5
1	A	0.83	0/2226	1.24	7/3032 (0.2%)

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$Observed(^o)$	$\operatorname{Ideal}({}^{o})$
1	A	9	ARG	NE-CZ-NH2	-10.25	115.18	120.30
1	A	286	ARG	NE-CZ-NH2	7.46	124.03	120.30
1	A	246	ARG	NE-CZ-NH2	7.16	123.88	120.30
1	A	128	ASP	CB-CG-OD1	6.03	123.72	118.30
1	A	166	ILE	CB-CA-C	-5.81	99.98	111.60
1	A	32	ARG	NE-CZ-NH2	5.25	122.93	120.30
1	A	282	ASP	CB-CG-OD2	5.24	123.01	118.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2167	0	1986	14	0
2	A	42	0	68	1	0
3	A	178	0	0	4	0
All	All	2387	0	2054	14	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 (14) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$\begin{array}{c} {\rm Interatomic} \\ {\rm distance} \ ({\rm \AA}) \end{array}$	Clash overlap (Å)
1:A:207:THR:HG23	1:A:229:ASN:HD21	1.40	0.86
1:A:139:ILE:HG12	1:A:142:VAL:HB	1.56	0.84
1:A:139:ILE:HG23	3:A:458:HOH:O	1.86	0.72
1:A:207:THR:HG23	1:A:229:ASN:ND2	2.05	0.70
1:A:207:THR:CG2	1:A:229:ASN:HD21	2.17	0.51
1:A:202:PHE:O	1:A:203:ASN:HB2	2.12	0.48
1:A:271:VAL:HG23	2:A:291:C8E:H101	1.96	0.48
1:A:235:THR:HG23	3:A:434:HOH:O	2.15	0.46
1:A:201:LYS:HD2	1:A:206:GLY:O	2.17	0.45
1:A:156:ASP:HB3	1:A:159:LEU:HD12	1.99	0.45
1:A:139:ILE:HG13	3:A:468:HOH:O	2.18	0.43
1:A:139:ILE:HG13	1:A:140:SER:H	1.85	0.42
1:A:282:ASP:CG	3:A:464:HOH:O	2.57	0.42
1:A:139:ILE:HD11	1:A:142:VAL:HG21	2.03	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	290/289 (100%)	279 (96%)	8 (3%)	3 (1%)	15 6

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	203	ASN
1	A	139	ILE

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Mol	Chain	Res	Type
1	A	140	SER

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.

\mathbf{Mol}	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	221/217 (102%)	210 (95%)	11 (5%)	24 16	

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	76	ASN
1	A	130	TYR
1	A	146	LEU
1	A	160	VAL
1	A	166	ILE
1	A	174	MET
1	A	204	ASP
1	A	207	THR
1	A	229	ASN
1	A	235	THR
1	A	255	ILE

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

Mol	Chain	Res	Type
1	A	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)

2 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	hain Res	es Link	Bo	Bond lengths			Bond angles		
MIOI	туре	Chain	nes	LIIIK	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z >2	
2	C8E	A	291	-	20,20,20	0.83	0	19,19,19	2.22	8 (42%)	
2	C8E	A	290	-	20,20,20	0.92	0	19,19,19	2.08	8 (42%)	

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.

Mo	l Type	Chain	Res	Link	Chirals	Torsions	Rings
2	C8E	A	291	-	-	11/18/18/18	-
2	C8E	A	290	-	-	10/18/18/18	-

There are no bond length outliers.

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$Observed(^o)$	$\operatorname{Ideal}({}^{o})$
2	A	291	C8E	O15-C14-C13	5.48	135.09	110.39
2	A	290	C8E	O15-C14-C13	4.30	129.79	110.39
2	A	290	C8E	O18-C17-C16	3.80	127.54	110.39
2	A	291	C8E	O12-C13-C14	3.67	126.94	110.39

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Mol	Chain	Res	Type	Atoms	Z	$Observed(^o)$	$Ideal(^{o})$
2	A	291	C8E	O18-C19-C20	3.59	125.84	110.07
2	A	290	C8E	O18-C19-C20	3.50	125.44	110.07
2	A	290	C8E	C7-C6-C5	-3.08	98.80	114.42
2	A	291	C8E	C7-C6-C5	-2.76	100.41	114.42
2	A	291	C8E	O18-C17-C16	2.39	121.19	110.39
2	A	290	C8E	O12-C13-C14	2.38	121.12	110.39
2	A	291	C8E	O9-C8-C7	2.23	122.00	110.26
2	A	291	C8E	O15-C16-C17	2.22	120.41	110.39
2	A	290	C8E	O9-C8-C7	2.20	121.80	110.26
2	A	290	C8E	O15-C16-C17	2.08	119.78	110.39
2	A	290	C8E	C5-C4-C3	-2.01	104.20	114.42
2	A	291	C8E	C5-C4-C3	-2.01	104.24	114.42

There are no chirality outliers.

All (21) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	291	C8E	O12-C13-C14-O15
2	A	290	C8E	C16-C17-O18-C19
2	A	290	C8E	C2-C3-C4-C5
2	A	290	C8E	C3-C4-C5-C6
2	A	291	C8E	C2-C3-C4-C5
2	A	291	C8E	O18-C19-C20-O21
2	A	290	C8E	O12-C13-C14-O15
2	A	291	C8E	C3-C4-C5-C6
2	A	291	C8E	C1-C2-C3-C4
2	A	291	C8E	C7-C8-O9-C10
2	A	291	C8E	C4-C5-C6-C7
2	A	291	C8E	C20-C19-O18-C17
2	A	290	C8E	C17-C16-O15-C14
2	A	290	C8E	C1-C2-C3-C4
2	A	290	C8E	C20-C19-O18-C17
2	A	290	C8E	C14-C13-O12-C11
2	A	290	C8E	C7-C8-O9-C10
2	A	291	C8E	C11-C10-O9-C8
2	A	291	C8E	C14-C13-O12-C11
2	A	290	C8E	O15-C16-C17-O18
2	A	291	C8E	O15-C16-C17-O18

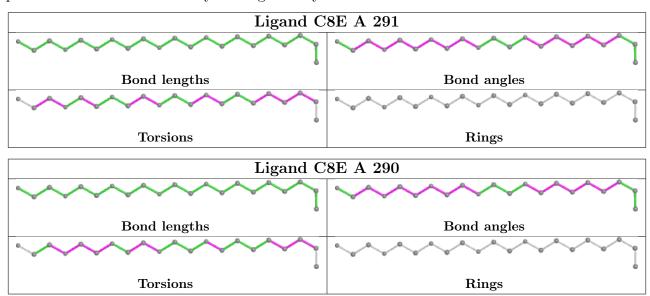
There are no ring outliers.

1 monomer is involved in 1 short contact:



Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	291	C8E	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 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)

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^{th} 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	$\langle { m RSRZ} \rangle$	$\# \mathrm{RSRZ}{>}2$	$OWAB(Å^2)$	Q<0.9
1	A	289/289 (100%)	-0.24	16 (5%) 25 27	17, 29, 54, 70	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	139	ILE	6.4
1	A	140	SER	5.9
1	A	173	ASP	5.0
1	A	204	ASP	4.3
1	A	232	PHE	3.6
1	A	20	VAL	3.5
1	A	202	PHE	3.4
1	A	203	ASN	3.1
1	A	174	MET	2.6
1	A	138	SER	2.6
1	A	158	SER	2.6
1	A	230	TYR	2.5
1	A	137	TYR	2.3
1	A	200	TYR	2.2
1	A	205	ALA	2.1
1	A	159	LEU	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 monosaccharides in this entry.

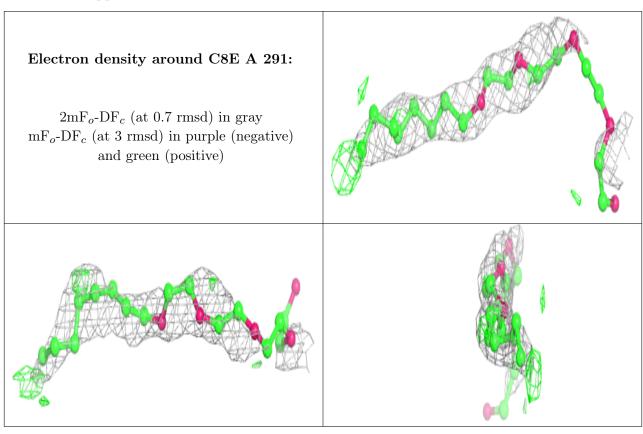


6.4 Ligands (i)

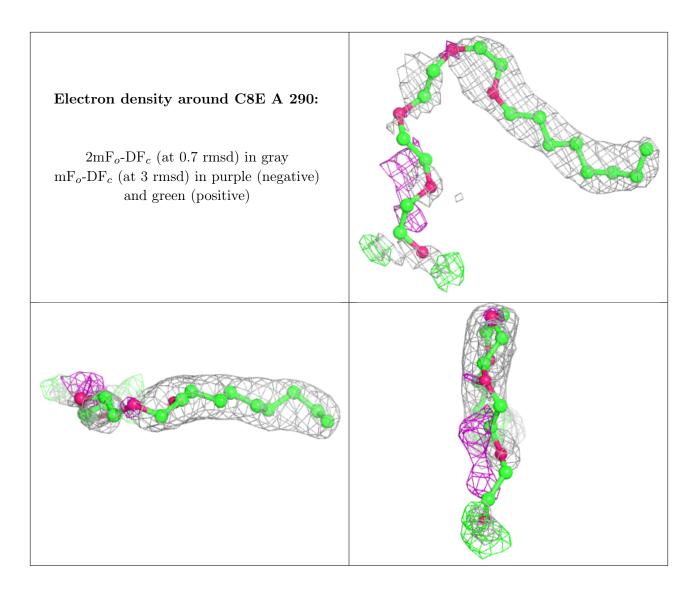
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9'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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q < 0.9
2	C8E	A	291	21/21	0.66	0.39	80,89,107,109	0
2	C8E	A	290	21/21	0.70	0.33	42,65,76,77	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.







6.5 Other polymers (i)

There are no such residues in this entry.

