



Full wwPDB X-ray Structure Validation Report ⓘ

May 27, 2024 – 09:19 PM EDT

PDB ID : 6MFC
Title : GphF GNAT-like decarboxylase
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Deposited on : 2018-09-10
Resolution : 2.59 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 ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtrriage (Phenix) : 1.13
EDS : 2.36.2
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) : 2.36.2

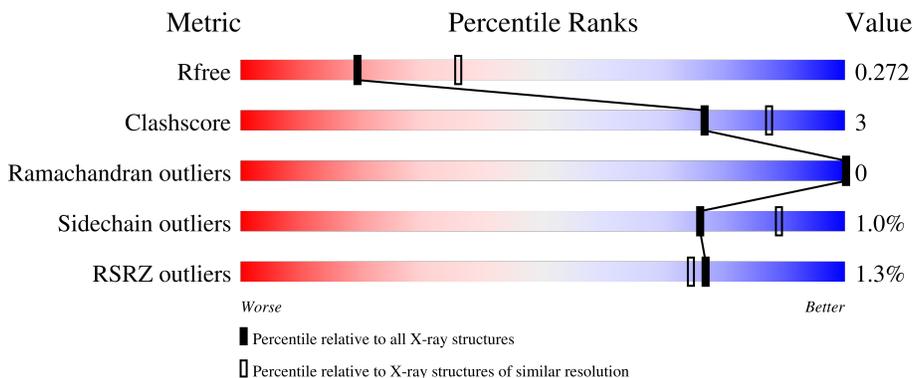
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.59 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	3676 (2.60-2.56)
Clashscore	141614	4049 (2.60-2.56)
Ramachandran outliers	138981	3979 (2.60-2.56)
Sidechain outliers	138945	3979 (2.60-2.56)
RSRZ outliers	127900	3614 (2.60-2.56)

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 $\leq 5\%$. 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	232	 2% 77% 7% 16%
1	B	232	 77% 7% 16%

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 3011 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 GphF.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	B	195	1467	915	291	257	4	0	0	0
1	A	195	1467	915	291	257	4	0	0	0

There are 48 discrepancies between the modelled and reference sequences:

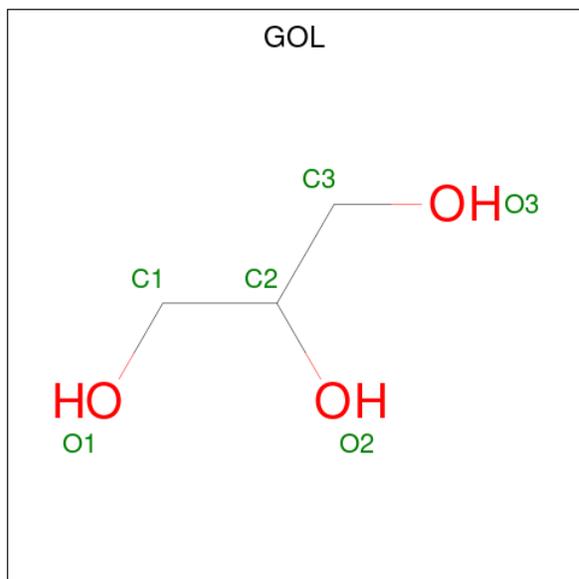
Chain	Residue	Modelled	Actual	Comment	Reference
B	474	MET	-	initiating methionine	UNP U6BSB2
B	475	HIS	-	expression tag	UNP U6BSB2
B	476	HIS	-	expression tag	UNP U6BSB2
B	477	HIS	-	expression tag	UNP U6BSB2
B	478	HIS	-	expression tag	UNP U6BSB2
B	479	HIS	-	expression tag	UNP U6BSB2
B	480	HIS	-	expression tag	UNP U6BSB2
B	481	SER	-	expression tag	UNP U6BSB2
B	482	SER	-	expression tag	UNP U6BSB2
B	483	GLY	-	expression tag	UNP U6BSB2
B	484	VAL	-	expression tag	UNP U6BSB2
B	485	ASP	-	expression tag	UNP U6BSB2
B	486	LEU	-	expression tag	UNP U6BSB2
B	487	GLY	-	expression tag	UNP U6BSB2
B	488	THR	-	expression tag	UNP U6BSB2
B	489	GLU	-	expression tag	UNP U6BSB2
B	490	ASN	-	expression tag	UNP U6BSB2
B	491	LEU	-	expression tag	UNP U6BSB2
B	492	TYR	-	expression tag	UNP U6BSB2
B	493	PHE	-	expression tag	UNP U6BSB2
B	494	GLN	-	expression tag	UNP U6BSB2
B	495	SER	-	expression tag	UNP U6BSB2
B	496	ASN	-	expression tag	UNP U6BSB2
B	497	ALA	-	expression tag	UNP U6BSB2
A	474	MET	-	initiating methionine	UNP U6BSB2

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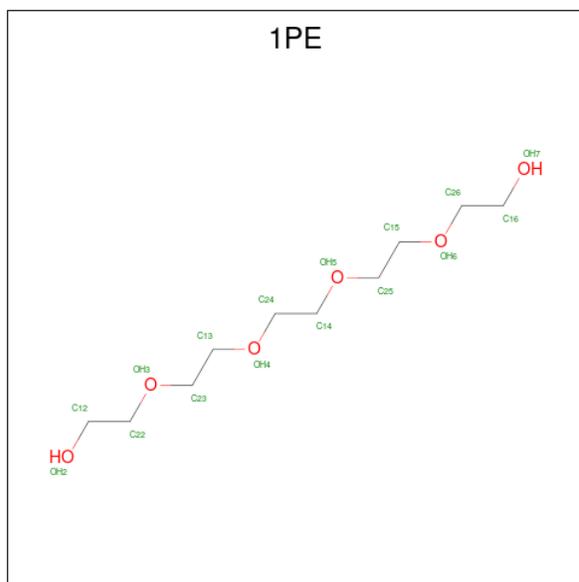
Chain	Residue	Modelled	Actual	Comment	Reference
A	475	HIS	-	expression tag	UNP U6BSB2
A	476	HIS	-	expression tag	UNP U6BSB2
A	477	HIS	-	expression tag	UNP U6BSB2
A	478	HIS	-	expression tag	UNP U6BSB2
A	479	HIS	-	expression tag	UNP U6BSB2
A	480	HIS	-	expression tag	UNP U6BSB2
A	481	SER	-	expression tag	UNP U6BSB2
A	482	SER	-	expression tag	UNP U6BSB2
A	483	GLY	-	expression tag	UNP U6BSB2
A	484	VAL	-	expression tag	UNP U6BSB2
A	485	ASP	-	expression tag	UNP U6BSB2
A	486	LEU	-	expression tag	UNP U6BSB2
A	487	GLY	-	expression tag	UNP U6BSB2
A	488	THR	-	expression tag	UNP U6BSB2
A	489	GLU	-	expression tag	UNP U6BSB2
A	490	ASN	-	expression tag	UNP U6BSB2
A	491	LEU	-	expression tag	UNP U6BSB2
A	492	TYR	-	expression tag	UNP U6BSB2
A	493	PHE	-	expression tag	UNP U6BSB2
A	494	GLN	-	expression tag	UNP U6BSB2
A	495	SER	-	expression tag	UNP U6BSB2
A	496	ASN	-	expression tag	UNP U6BSB2
A	497	ALA	-	expression tag	UNP U6BSB2

- Molecule 2 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



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

- Molecule 3 is PENTAETHYLENE GLYCOL (three-letter code: 1PE) (formula: C₁₀H₂₂O₆).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	1	Total C O 7 5 2	0	0
3	A	1	Total C O 10 6 4	0	0

- Molecule 4 is water.

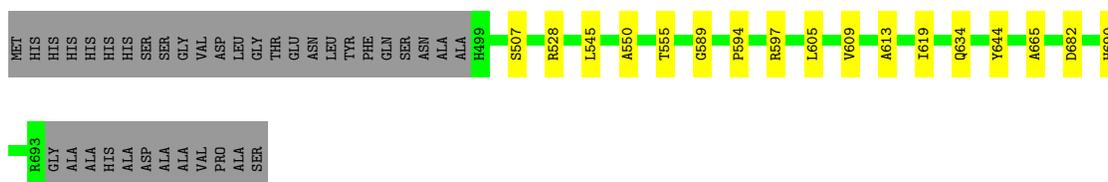
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	24	Total O 24 24	0	0
4	A	18	Total O 18 18	0	0

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.

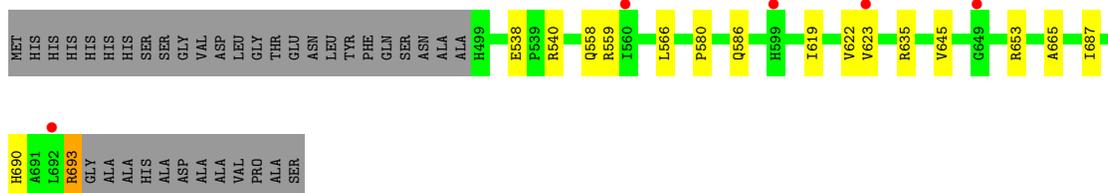
- Molecule 1: GphF

Chain B: 



- Molecule 1: GphF

Chain A: 



4 Data and refinement statistics i

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	138.33Å 145.70Å 78.01Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	43.96 – 2.59 43.96 – 2.59	Depositor EDS
% Data completeness (in resolution range)	99.6 (43.96-2.59) 99.6 (43.96-2.59)	Depositor EDS
R_{merge}	0.18	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.42 (at 2.58Å)	Xtrriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.234 , 0.272 0.234 , 0.272	Depositor DCC
R_{free} test set	1242 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	77.7	Xtrriage
Anisotropy	0.399	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 58.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	0.000 for -k,-h,-l	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	3011	wwPDB-VP
Average B, all atoms (Å ²)	91.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.94% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 1PE, GOL

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	0/1500	0.42	0/2036
1	B	0.24	0/1500	0.41	0/2036
All	All	0.24	0/3000	0.41	0/4072

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	1467	0	1463	11	0
1	B	1467	0	1463	9	0
2	A	6	0	8	0	0
2	B	12	0	16	0	0
3	A	10	0	13	0	0
3	B	7	0	6	1	0
4	A	18	0	0	1	0
4	B	24	0	0	0	0
All	All	3011	0	2969	18	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 (18) 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:619:ILE:O	1:A:693:ARG:NH1	2.18	0.77
1:B:555:THR:HG21	1:B:609:VAL:HG21	1.74	0.68
1:A:559:ARG:NH1	1:A:580:PRO:O	2.32	0.63
1:A:645:VAL:O	1:A:653:ARG:NH2	2.31	0.63
1:B:665:ALA:HB2	1:B:690:HIS:CE1	2.44	0.52
1:A:665:ALA:HB2	1:A:690:HIS:CE1	2.44	0.52
1:B:550:ALA:HB2	1:A:693:ARG:NH2	2.25	0.51
1:B:589:GLY:HA2	3:B:803:1PE:H242	1.94	0.50
1:B:545:LEU:HB2	1:B:605:LEU:HD21	1.93	0.50
1:A:622:VAL:HB	1:A:690:HIS:HB2	1.98	0.45
1:A:538:GLU:HG3	1:A:558:GLN:HA	1.99	0.45
1:B:528:ARG:HD2	1:A:540:ARG:NH1	2.32	0.43
1:B:594:PRO:HA	1:B:597:ARG:HE	1.83	0.43
1:A:635:ARG:NH2	4:A:902:HOH:O	2.51	0.42
1:B:634:GLN:HG3	1:B:644:TYR:CE1	2.54	0.42
1:B:613:ALA:HB1	1:B:619:ILE:HG21	2.01	0.42
1:A:586:GLN:HA	1:A:623:VAL:O	2.19	0.42
1:A:566:LEU:HD12	1:A:687:ILE:HG23	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	193/232 (83%)	189 (98%)	4 (2%)	0	100	100
1	B	193/232 (83%)	187 (97%)	6 (3%)	0	100	100
All	All	386/464 (83%)	376 (97%)	10 (3%)	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	143/169 (85%)	142 (99%)	1 (1%)	84	93
1	B	143/169 (85%)	141 (99%)	2 (1%)	67	84
All	All	286/338 (85%)	283 (99%)	3 (1%)	76	89

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	507	SER
1	B	682	ASP
1	A	693	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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](#)

5 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
3	1PE	A	802	-	9,9,15	0.51	0	8,8,14	0.25	0
2	GOL	A	801	-	5,5,5	0.88	0	5,5,5	1.02	0
3	1PE	B	803	-	6,6,15	0.51	0	5,5,14	0.25	0
2	GOL	B	802	-	5,5,5	0.90	0	5,5,5	1.03	0
2	GOL	B	801	-	5,5,5	0.93	0	5,5,5	0.99	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
3	1PE	A	802	-	-	3/7/7/13	-
2	GOL	A	801	-	-	2/4/4/4	-
3	1PE	B	803	-	-	3/4/4/13	-
2	GOL	B	802	-	-	0/4/4/4	-
2	GOL	B	801	-	-	4/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	801	GOL	O1-C1-C2-O2
2	A	801	GOL	O1-C1-C2-C3
2	B	801	GOL	O1-C1-C2-C3
2	B	801	GOL	C1-C2-C3-O3
2	B	801	GOL	O1-C1-C2-O2
3	B	803	1PE	C24-C14-OH5-C25
3	A	802	1PE	C13-C23-OH3-C22
3	B	803	1PE	C14-C24-OH4-C13
3	A	802	1PE	C12-C22-OH3-C23
2	B	801	GOL	O2-C2-C3-O3

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Mol	Chain	Res	Type	Atoms
3	A	802	1PE	OH4-C13-C23-OH3
3	B	803	1PE	C23-C13-OH4-C24

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	803	1PE	1	0

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, 95th 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 ²)	Q<0.9
1	A	195/232 (84%)	0.53	5 (2%) 56 52	70, 88, 122, 140	0
1	B	195/232 (84%)	0.39	0 100 100	65, 88, 123, 160	0
All	All	390/464 (84%)	0.46	5 (1%) 77 74	65, 88, 123, 160	0

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	599	HIS	2.3
1	A	560	ILE	2.2
1	A	623	VAL	2.1
1	A	692	LEU	2.1
1	A	649	GLY	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, 95th 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(\AA^2)	Q<0.9
2	GOL	A	801	6/6	0.78	0.29	91,106,108,109	0
2	GOL	B	801	6/6	0.79	0.23	104,111,113,116	0
2	GOL	B	802	6/6	0.85	0.53	121,127,133,135	0
3	1PE	B	803	7/16	0.85	0.35	107,110,139,141	0
3	1PE	A	802	10/16	0.86	0.27	94,102,105,106	0

6.5 Other polymers [i](#)

There are no such residues in this entry.