



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 30, 2017 – 08:44 PM EDT

PDB ID : 5EHH  
Title : Structure of human DPP3 in complex with endomorphin-2.  
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eroux, P.; Gruber, K.  
Deposited on : unknown  
Resolution : 2.38 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

<http://wwpdb.org/validation/2016/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  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20029824  
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : rb-20029824

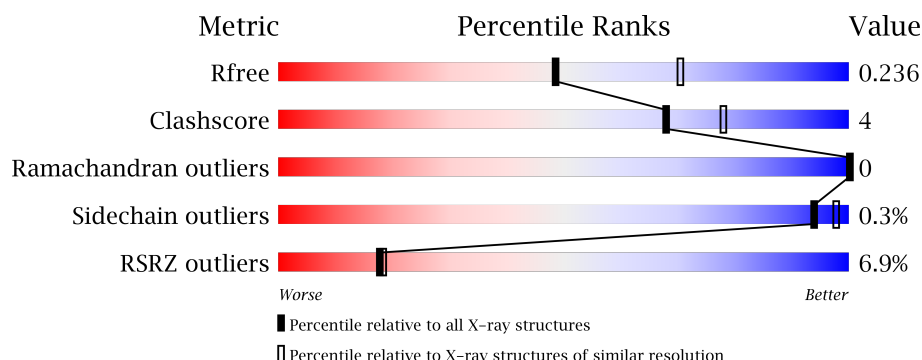
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.38 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))
$R_{free}$	100719	4388 (2.40-2.36)
Clashscore	112137	4984 (2.40-2.36)
Ramachandran outliers	110173	4907 (2.40-2.36)
Sidechain outliers	110143	4909 (2.40-2.36)
RSRZ outliers	101464	4423 (2.40-2.36)

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	726	<div> <div>6%</div> <div> <div></div> <div>90%</div> <div>10%</div> </div> </div>
2	B	5	<div> <div>60%</div> <div> <div></div> <div>60%</div> <div>40%</div> </div> </div>

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 6061 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 Dipeptidyl peptidase 3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	724	Total	C	N	O	S	0	0	0
			5745	3662	971	1100	12			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	19	SER	CYS	engineered mutation	UNP Q9NY33
A	207	CYS	GLU	engineered mutation	UNP Q9NY33
A	451	ALA	GLU	engineered mutation	UNP Q9NY33
A	491	CYS	SER	engineered mutation	UNP Q9NY33
A	519	SER	CYS	engineered mutation	UNP Q9NY33
A	654	SER	CYS	engineered mutation	UNP Q9NY33

- Molecule 2 is a protein called Endomorphin-2.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	B	5	Total	C	N	O	0	0	1
			42	32	5	5			

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Zn	0	0
			1	1		

- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	2	Total	Mg	0	0
			2	2		

- Molecule 5 is POTASSIUM ION (three-letter code: K) (formula: K).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total 1	K 1	0	0

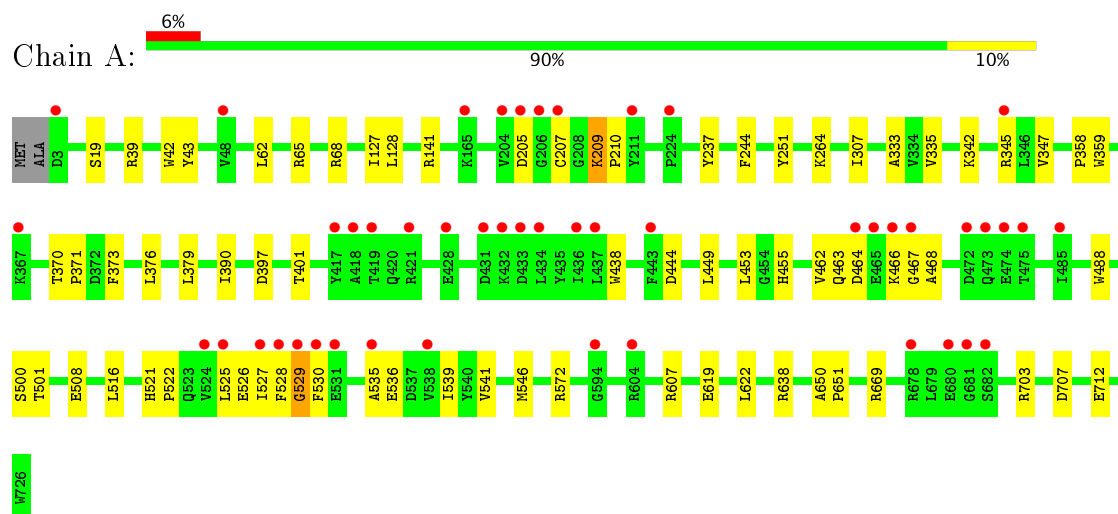
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	270	Total 270	O 270	0	0

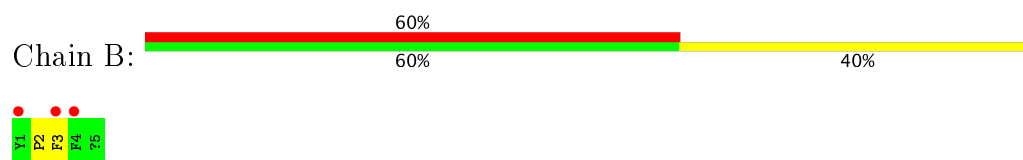
### 3 Residue-property plots [i](#)

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: Dipeptidyl peptidase 3



- Molecule 2: Endomorphin-2



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	120.03 Å   105.46 Å   64.72 Å 90.00°   93.49°   90.00°	Depositor
Resolution (Å)	49.09 – 2.38 49.09 – 2.38	Depositor EDS
% Data completeness (in resolution range)	98.9 (49.09-2.38) 98.9 (49.09-2.38)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	0.09	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.06 (at 2.37 Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, $R_{free}$	0.194   ,   0.237 0.194   ,   0.236	Depositor DCC
$R_{free}$ test set	1600 reflections (5.00%)	DCC
Wilson B-factor (Å <sup>2</sup> )	40.6	Xtriage
Anisotropy	0.366	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 36.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	6061	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.59% 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: K, MG, ZN, NH2

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	0.32	1/5877 (0.0%)	0.40	2/7962 (0.0%)
2	B	0.28	0/44	0.37	0/59
All	All	0.32	1/5921 (0.0%)	0.40	2/8021 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	210	PRO	N-CD	5.14	1.55	1.47

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	529	GLY	N-CA-C	5.90	127.86	113.10
1	A	209	LYS	C-N-CD	5.44	139.82	128.40

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	5745	0	5617	50	1
2	B	42	0	36	2	0
3	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	2	0	0	0	0
5	A	1	0	0	0	0
6	A	270	0	0	4	0
All	All	6061	0	5653	50	1

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 (50) 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:463:GLN:HE21	1:A:467:GLY:HA3	1.23	1.01
1:A:464:ASP:OD1	1:A:466:LYS:O	1.89	0.90
1:A:463:GLN:HE21	1:A:467:GLY:CA	1.92	0.83
1:A:516:LEU:HD11	1:A:546:MET:HG2	1.70	0.74
1:A:463:GLN:NE2	1:A:467:GLY:HA3	2.04	0.69
1:A:345:ARG:NH1	1:A:527:ILE:O	2.27	0.67
1:A:525:LEU:HD13	1:A:535:ALA:HB1	1.76	0.66
1:A:358:PRO:HB3	1:A:619:GLU:HG3	1.82	0.61
1:A:462:VAL:HG12	1:A:463:GLN:N	2.17	0.60
1:A:669:ARG:NH2	6:A:906:HOH:O	2.35	0.59
1:A:370:THR:O	1:A:370:THR:HG23	2.02	0.59
1:A:438:TRP:HB2	1:A:541:VAL:HG21	1.83	0.58
1:A:501:THR:OG1	1:A:638:ARG:NH2	2.39	0.56
1:A:397:ASP:O	1:A:401:THR:OG1	2.20	0.56
1:A:19:SER:HB3	1:A:379:LEU:HD23	1.88	0.55
1:A:463:GLN:CG	1:A:467:GLY:HA2	2.37	0.55
1:A:342:LYS:NZ	1:A:444:ASP:OD2	2.39	0.55
1:A:43:TYR:OH	1:A:707:ASP:OD2	2.22	0.54
1:A:68:ARG:NH1	1:A:712:GLU:OE2	2.42	0.53
1:A:65:ARG:NH1	6:A:911:HOH:O	2.42	0.53
1:A:390:ILE:HA	2:B:2:PRO:HA	1.91	0.52
1:A:463:GLN:OE1	1:A:488:TRP:NE1	2.42	0.52
1:A:463:GLN:HG2	1:A:467:GLY:HA2	1.93	0.51
1:A:39:ARG:NH1	6:A:917:HOH:O	2.44	0.50
1:A:264:LYS:NZ	6:A:912:HOH:O	2.44	0.50
1:A:205:ASP:OD2	1:A:209:LYS:HB3	2.13	0.49
1:A:525:LEU:HD11	1:A:539:ILE:HD11	1.96	0.47
1:A:205:ASP:OD1	1:A:209:LYS:N	2.43	0.47
1:A:536:GLU:HG2	1:A:607:ARG:HH21	1.79	0.47
1:A:462:VAL:CG1	1:A:463:GLN:N	2.78	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:345:ARG:HB3	1:A:527:ILE:HG23	1.96	0.47
1:A:500:SER:OG	1:A:501:THR:N	2.47	0.47
1:A:237:TYR:N	1:A:244:PHE:O	2.46	0.47
1:A:347:VAL:HG13	1:A:371:PRO:HD2	1.95	0.46
1:A:455:HIS:CE1	1:A:508:GLU:OE1	2.68	0.45
1:A:62:LEU:HD21	1:A:127:ILE:HD13	1.98	0.45
1:A:397:ASP:OD1	1:A:397:ASP:N	2.49	0.45
1:A:449:LEU:HD22	1:A:453:LEU:HD12	1.99	0.45
1:A:526:GLU:C	1:A:529:GLY:H	2.22	0.43
1:A:528:PHE:CB	1:A:530:PHE:HE2	2.32	0.43
1:A:650:ALA:HA	1:A:651:PRO:HD3	1.92	0.42
1:A:572:ARG:HH21	2:B:3:PHE:HE1	1.67	0.42
1:A:522:PRO:HA	1:A:525:LEU:HD12	2.00	0.42
1:A:42:TRP:CD2	1:A:703:ARG:HD3	2.56	0.41
1:A:466:LYS:O	1:A:468:ALA:N	2.53	0.41
1:A:307:ILE:HA	1:A:335:VAL:HG12	2.03	0.41
1:A:521:HIS:HA	1:A:522:PRO:HD3	1.93	0.41
1:A:333:ALA:HB3	1:A:376:LEU:HD13	2.02	0.41
1:A:128:LEU:HD22	1:A:141:ARG:HG2	2.03	0.40
1:A:359:TRP:HZ3	1:A:622:LEU:HG	1.86	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:207:CYS:SG	1:A:207:CYS:SG 2_857	1.69	0.51

## 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	722/726 (99%)	703 (97%)	19 (3%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	B	3/5 (60%)	2 (67%)	1 (33%)	0	100	100
All	All	725/731 (99%)	705 (97%)	20 (3%)	0	100	100

There are no Ramachandran outliers to report.

### 5.3.2 Protein sidechains ⓘ

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	608/609 (100%)	606 (100%)	2 (0%)	94	98
2	B	4/4 (100%)	4 (100%)	0	100	100
All	All	612/613 (100%)	610 (100%)	2 (0%)	94	98

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	251	TYR
1	A	373	PHE

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

Mol	Chain	Res	Type
1	A	463	GLN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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](#)

Of 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 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	724/726 (99%)	0.24	47 (6%) 20 20	17, 33, 66, 100	0
2	B	4/5 (80%)	2.86	3 (75%) 0 0	38, 41, 41, 46	4 (100%)
All	All	728/731 (99%)	0.26	50 (6%) 18 18	17, 33, 66, 100	4 (0%)

All (50) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	467	GLY	8.2
1	A	206	GLY	5.5
2	B	3	PHE	5.4
1	A	465	GLU	4.6
1	A	535	ALA	4.4
1	A	485	ILE	3.7
1	A	207	CYS	3.7
1	A	419	THR	3.7
1	A	529	GLY	3.7
1	A	434	LEU	3.6
1	A	475	THR	3.2
1	A	525	LEU	3.0
1	A	466	LYS	3.0
1	A	472	ASP	3.0
1	A	681	GLY	2.9
1	A	417	TYR	2.9
1	A	205	ASP	2.8
1	A	204	VAL	2.7
1	A	528	PHE	2.7
1	A	211	TYR	2.7
1	A	524	VAL	2.7
1	A	224	PRO	2.6
1	A	474	GLU	2.6
1	A	680	GLU	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	594	GLY	2.6
1	A	530	PHE	2.5
1	A	531	GLU	2.5
1	A	428	GLU	2.5
1	A	604	ARG	2.5
1	A	527	ILE	2.5
1	A	682	SER	2.4
2	B	1	TYR	2.4
1	A	473	GLN	2.4
1	A	436	ILE	2.4
1	A	3	ASP	2.4
1	A	538	VAL	2.4
1	A	421	ARG	2.3
1	A	464	ASP	2.3
1	A	418	ALA	2.3
1	A	48	VAL	2.3
1	A	165	LYS	2.2
1	A	678	ARG	2.1
1	A	443	PHE	2.1
1	A	367	LYS	2.1
1	A	431	ASP	2.1
1	A	433	ASP	2.1
2	B	4	PHE	2.1
1	A	345	ARG	2.0
1	A	437	LEU	2.0
1	A	432	LYS	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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
4	MG	A	804	1/1	0.98	0.09	-0.70	17,17,17,17	0
3	ZN	A	801	1/1	0.97	0.05	-1.56	49,49,49,49	0
5	K	A	803	1/1	0.99	0.09	-2.39	31,31,31,31	0
4	MG	A	802	1/1	0.98	0.21	-	21,21,21,21	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.