



wwPDB X-ray Structure Validation Summary Report ⓘ

Jun 15, 2020 – 10:40 pm BST

PDB ID : 3BBB
Title : Crystal structure of the NM23-H2 transcription factor complex with dinucleotide d(AG)
Authors : Weichsel, A.; Montfort, W.R.
Deposited on : 2007-11-09
Resolution : 1.30 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.11
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)	:	2.11

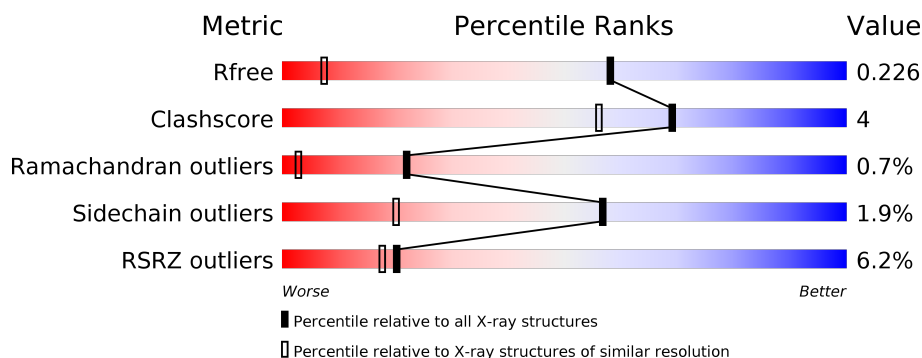
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.30 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	1058 (1.30-1.30)
Clashscore	141614	1101 (1.30-1.30)
Ramachandran outliers	138981	1058 (1.30-1.30)
Sidechain outliers	138945	1058 (1.30-1.30)
RSRZ outliers	127900	1029 (1.30-1.30)

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 ≥ 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	151	<div> <div>12%</div> <div> <div></div> <div>91%</div> <div>8%</div> <div>..</div> </div> </div>
1	B	151	<div> <div>%</div> <div> <div></div> <div>93%</div> <div>5%</div> <div>.</div> </div> </div>
1	C	151	<div> <div>7%</div> <div> <div></div> <div>86%</div> <div>12%</div> <div>.</div> </div> </div>
1	D	151	<div> <div>10%</div> <div> <div></div> <div>91%</div> <div>9%</div> <div>.</div> </div> </div>
1	E	151	<div> <div>3%</div> <div> <div></div> <div>86%</div> <div>12%</div> <div>.</div> </div> </div>
1	F	151	<div> <div>4%</div> <div> <div></div> <div>91%</div> <div>8%</div> <div>.</div> </div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	DG	C	154[A]	-	-	-	X
2	DG	E	154[A]	-	-	-	X
3	DA	C	153[A]	-	-	-	X
3	DA	E	153[A]	-	-	-	X

2 Entry composition [i](#)

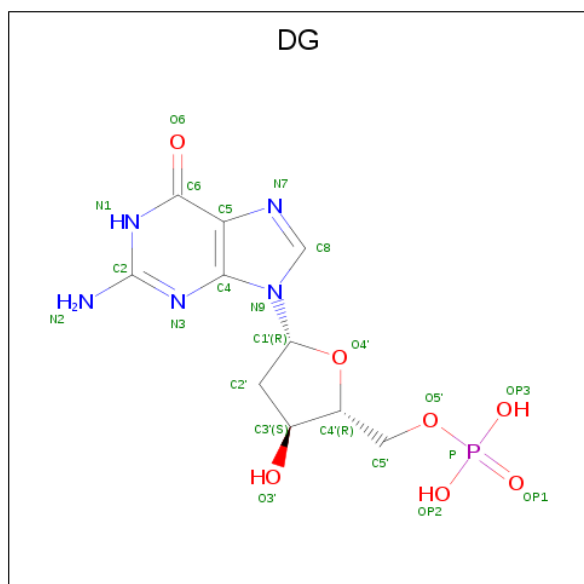
There are 4 unique types of molecules in this entry. The entry contains 8098 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Nucleoside diphosphate kinase B.

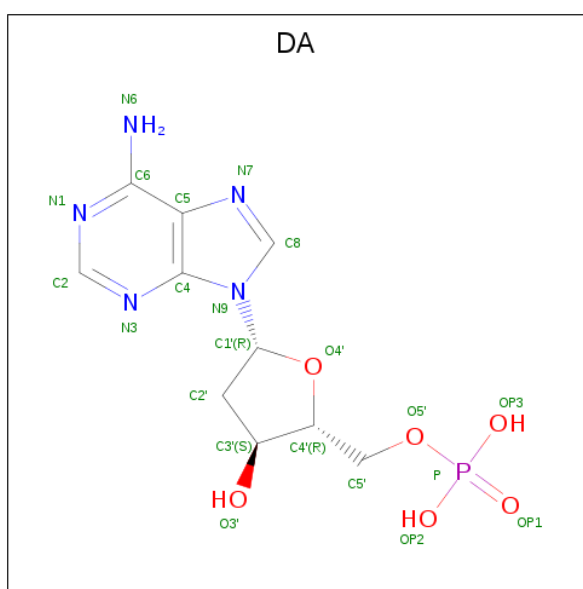
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	151	Total	C	N	O	S	0	2	0
			1218	781	212	219	6			
1	B	151	Total	C	N	O	S	0	6	0
			1244	798	215	224	7			
1	C	148	Total	C	N	O	S	0	3	0
			1211	775	213	217	6			
1	D	151	Total	C	N	O	S	0	1	0
			1213	778	211	218	6			
1	E	148	Total	C	N	O	S	0	3	0
			1209	777	209	217	6			
1	F	151	Total	C	N	O	S	0	4	0
			1235	796	212	220	7			

- Molecule 2 is 2'-DEOXYGUANOSINE-5'-MONOPHOSPHATE (three-letter code: DG) (formula: C₁₀H₁₄N₅O₇P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	B	1	Total	C	N	O	P	0	0
			23	10	5	7	1		
2	C	1	Total	C	N	O	P	0	1
			22	10	5	6	1		
2	E	1	Total	C	N	O	P	0	1
			22	10	5	6	1		
2	F	1	Total	C	N	O	P	0	0
			23	10	5	7	1		

- Molecule 3 is 2'-DEOXYADENOSINE-5'-MONOPHOSPHATE (three-letter code: DA) (formula: C₁₀H₁₄N₅O₆P).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	C	1	Total	C	N	O	0	1
			18	10	5	3		
3	E	1	Total	C	N	O	0	1
			18	10	5	3		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	88	Total	O	0	1
			88	88		
4	B	135	Total	O	0	2
			136	136		
4	C	94	Total	O	0	0
			94	94		

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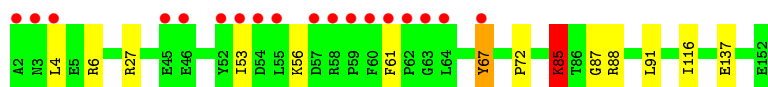
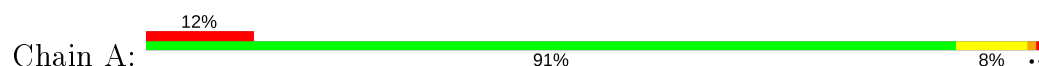
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	D	96	Total 96	O 96	0	1
4	E	121	Total 121	O 121	0	1
4	F	107	Total 107	O 107	0	1

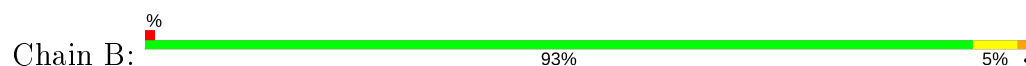
3 Residue-property plots [i](#)

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.

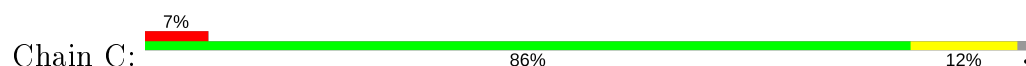
- Molecule 1: Nucleoside diphosphate kinase B



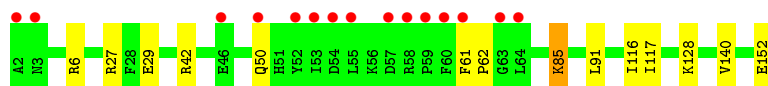
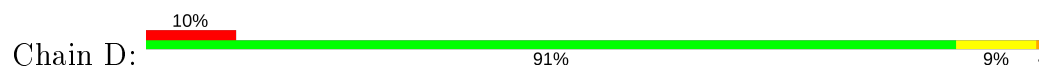
- Molecule 1: Nucleoside diphosphate kinase B



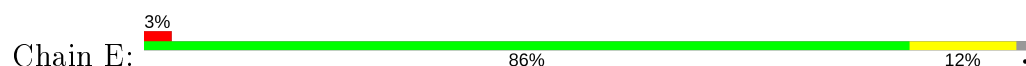
- Molecule 1: Nucleoside diphosphate kinase B



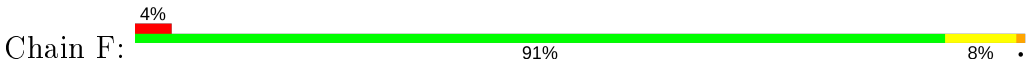
- Molecule 1: Nucleoside diphosphate kinase B



- Molecule 1: Nucleoside diphosphate kinase B



- Molecule 1: Nucleoside diphosphate kinase B



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	52.47Å 118.22Å 128.91Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	27.00 – 1.30 26.75 – 1.30	Depositor EDS
% Data completeness (in resolution range)	98.2 (27.00-1.30) 98.2 (26.75-1.30)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.84 (at 1.30Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.175 , 0.220 0.183 , 0.226	Depositor DCC
R_{free} test set	9678 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	12.9	Xtriage
Anisotropy	0.447	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 45.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	8098	wwPDB-VP
Average B, all atoms (Å ²)	17.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 83.62 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 2.0699e-07. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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

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	0.95	2/1252 (0.2%)	0.91	3/1685 (0.2%)
1	B	0.97	1/1287 (0.1%)	1.03	7/1730 (0.4%)
1	C	0.96	2/1242 (0.2%)	0.90	3/1670 (0.2%)
1	D	0.98	2/1244 (0.2%)	0.92	1/1674 (0.1%)
1	E	1.12	6/1244 (0.5%)	1.02	3/1674 (0.2%)
1	F	0.99	0/1273	0.93	0/1713
All	All	1.00	13/7542 (0.2%)	0.95	17/10146 (0.2%)

The worst 5 of 13 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	5	GLU	CB-CG	-8.53	1.35	1.52
1	A	137	GLU	CB-CG	-7.23	1.38	1.52
1	E	5	GLU	CD-OE2	-7.12	1.17	1.25
1	E	67[A]	TYR	CB-CG	6.42	1.61	1.51
1	E	67[B]	TYR	CB-CG	6.42	1.61	1.51

The worst 5 of 17 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	6	ARG	NE-CZ-NH1	-6.35	117.12	120.30
1	A	27	ARG	NE-CZ-NH2	-6.22	117.19	120.30
1	B	88	ARG	NE-CZ-NH1	6.17	123.38	120.30
1	C	88	ARG	NE-CZ-NH2	-5.90	117.35	120.30
1	D	42	ARG	NE-CZ-NH2	-5.88	117.36	120.30

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	1218	0	1227	5	0
1	B	1244	0	1259	9	0
1	C	1211	0	1216	11	0
1	D	1213	0	1221	8	0
1	E	1209	0	1213	15	0
1	F	1235	0	1249	8	0
2	B	23	0	12	3	0
2	C	22	0	12	3	0
2	E	22	0	12	5	0
2	F	23	0	12	3	0
3	C	18	0	12	4	0
3	E	18	0	12	1	0
4	A	88	0	0	0	0
4	B	136	0	0	5	0
4	C	94	0	0	3	0
4	D	96	0	0	0	0
4	E	121	0	0	3	0
4	F	107	0	0	1	0
All	All	8098	0	7457	55	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 4.

The worst 5 of 55 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:B:94:THR:HG21	2:B:153:DG:H2''	1.36	1.04
1:B:94:THR:CG2	2:B:153:DG:H2''	2.13	0.79
1:E:82[B]:ASN:HD21	1:E:85:LYS:HD3	1.48	0.79
3:C:153[A]:DA:H5'	2:C:154[A]:DG:C8	2.18	0.79
1:F:55:LEU:CD1	2:F:153:DG:H5'	2.15	0.76

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	151/151 (100%)	148 (98%)	2 (1%)	1 (1%)	22	3
1	B	155/151 (103%)	152 (98%)	2 (1%)	1 (1%)	25	4
1	C	149/151 (99%)	146 (98%)	2 (1%)	1 (1%)	22	3
1	D	150/151 (99%)	147 (98%)	2 (1%)	1 (1%)	22	3
1	E	149/151 (99%)	146 (98%)	2 (1%)	1 (1%)	22	3
1	F	153/151 (101%)	150 (98%)	2 (1%)	1 (1%)	22	3
All	All	907/906 (100%)	889 (98%)	12 (1%)	6 (1%)	22	3

5 of 6 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	116	ILE
1	B	116	ILE
1	C	116	ILE
1	D	116	ILE
1	E	116	ILE

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	132/130 (102%)	128 (97%)	4 (3%)	41	7
1	B	136/130 (105%)	133 (98%)	3 (2%)	52	15
1	C	131/130 (101%)	129 (98%)	2 (2%)	65	31

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	D	131/130 (101%)	129 (98%)	2 (2%)	65	31
1	E	131/130 (101%)	131 (100%)	0	100	100
1	F	134/130 (103%)	129 (96%)	5 (4%)	34	3
All	All	795/780 (102%)	779 (98%)	16 (2%)	57	17

5 of 16 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	58	ARG
1	C	141	ASP
1	F	67[A]	TYR
1	B	85	LYS
1	F	67[B]	TYR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	147	HIS
1	F	50	GLN

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

6 ligands are modelled in this entry.

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 [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	151/151 (100%)	0.35	18 (11%) 4 3	9, 16, 39, 52	0
1	B	151/151 (100%)	-0.04	2 (1%) 77 79	8, 13, 22, 37	0
1	C	148/151 (98%)	0.11	11 (7%) 14 11	9, 16, 28, 45	0
1	D	151/151 (100%)	0.36	15 (9%) 7 5	9, 15, 32, 37	0
1	E	148/151 (98%)	-0.02	4 (2%) 54 52	8, 14, 23, 30	0
1	F	151/151 (100%)	0.20	6 (3%) 38 35	9, 15, 27, 39	0
All	All	900/906 (99%)	0.16	56 (6%) 20 18	8, 14, 30, 52	0

The worst 5 of 56 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	2	ALA	11.1
1	B	2	ALA	11.0
1	D	2	ALA	6.3
1	C	59	PRO	6.1
1	A	2	ALA	5.9

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. 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(Å ²)	Q<0.9
3	DA	C	153[A]	18/22	0.16	0.62	77,78,78,78	18
2	DG	C	154[A]	22/23	0.34	0.53	77,78,79,79	22
3	DA	E	153[A]	18/22	0.39	0.61	79,79,82,82	18
2	DG	E	154[A]	22/23	0.57	0.53	80,80,82,83	22
2	DG	B	153	23/23	0.59	0.27	42,45,64,65	0
2	DG	F	153	23/23	0.65	0.27	33,38,51,52	0

6.5 Other polymers [i](#)

There are no such residues in this entry.