



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 9, 2018 – 08:33 am GMT

PDB ID : 1J06
Title : Crystal structure of mouse acetylcholinesterase in the apo form
Authors : Bourne, Y.; Taylor, P.; Radic, Z.; Marchot, P.
Deposited on : 2002-11-07
Resolution : 2.35 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
Mogul : 1.7.3 (157068), CSD as539be (2018)
Xtriage (Phenix) : 1.13
EDS : trunk30967
Percentile statistics : 20171227.v01 (using entries in the PDB archive December 27th 2017)
Refmac : 5.8.0158
CCP4 : 7.0 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk30967

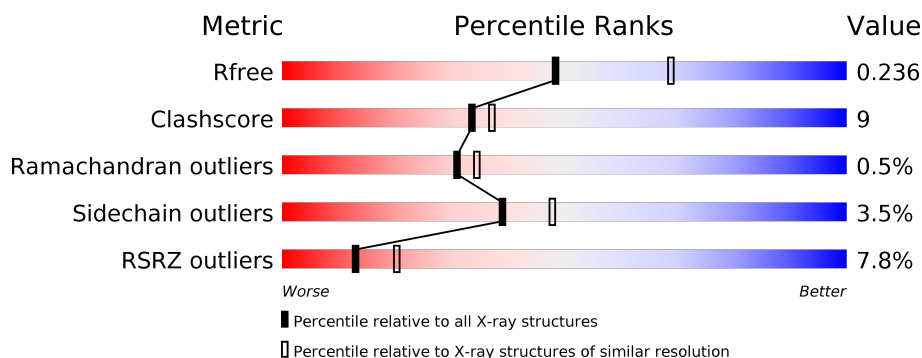
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.35 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}	111664	1015 (2.36-2.36)
Clashscore	122126	1081 (2.36-2.36)
Ramachandran outliers	120053	1066 (2.36-2.36)
Sidechain outliers	120020	1067 (2.36-2.36)
RSRZ outliers	108989	1002 (2.36-2.36)

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 ≥ 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	543	<div> <div>7%</div> <div>79%</div> <div>19%</div> <div>..</div> </div>
1	B	543	<div> <div>8%</div> <div>77%</div> <div>19%</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
6	AE3	A	902	-	-	-	X

2 Entry composition [i](#)

There are 8 unique types of molecules in this entry. The entry contains 8804 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 acetylcholinesterase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	535	Total	C	N	O	S	0	0	0
			4177	2679	725	759	14			
1	B	533	Total	C	N	O	S	0	0	0
			4159	2670	718	757	14			

- Molecule 2 is N-ACETYL-D-GLUCOSAMINE (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



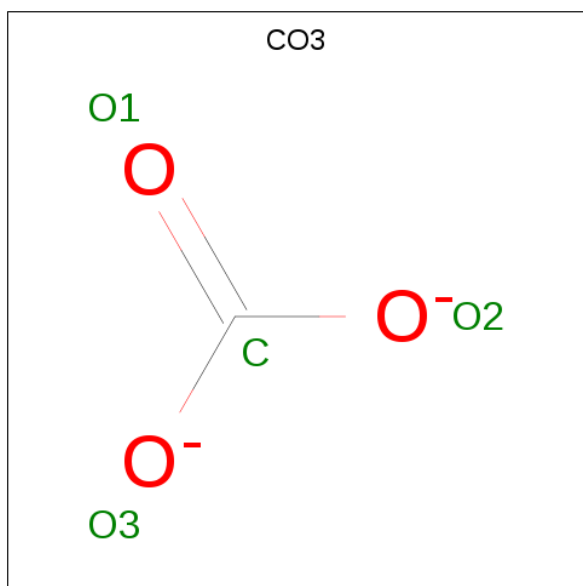
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	B	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 3 is ALPHA-L-FUCOSE (three-letter code: FUC) (formula: $C_6H_{12}O_5$).



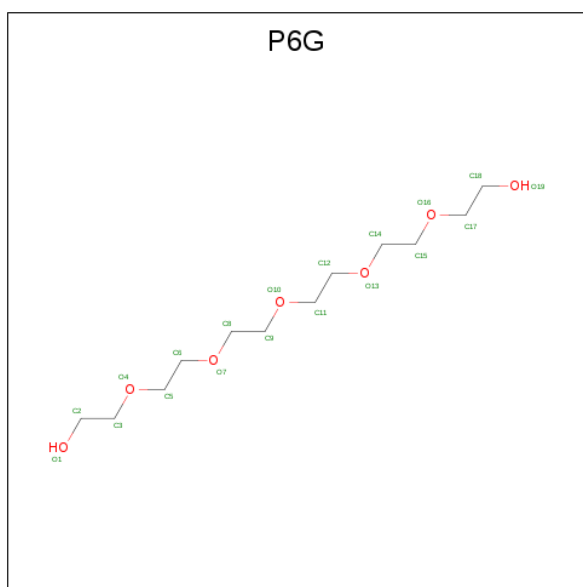
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			10	6	4		

- Molecule 4 is CARBONATE ION (three-letter code: CO3) (formula: CO_3).



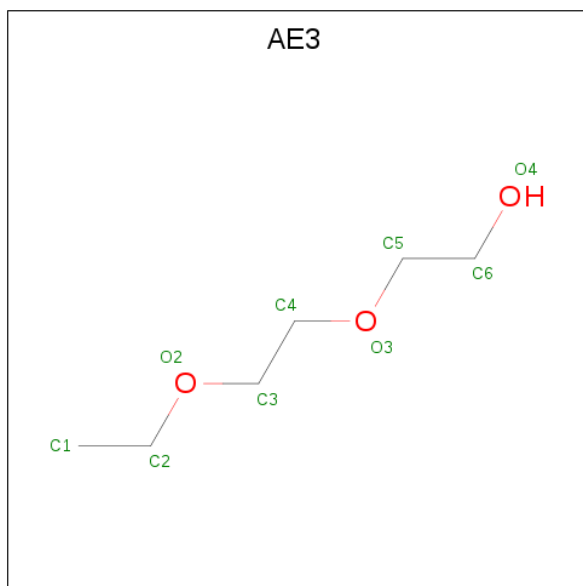
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			4	1	3		
4	B	1	Total	C	O	0	0
			4	1	3		

- Molecule 5 is HEXAETHYLENE GLYCOL (three-letter code: P6G) (formula: $\text{C}_{12}\text{H}_{26}\text{O}_7$).



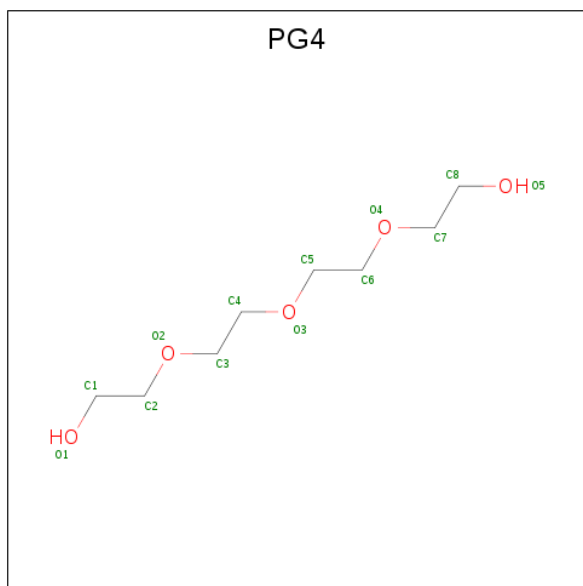
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			19	12	7		

- Molecule 6 is 2-(2-ETHOXYETHOXY)ETHANOL (three-letter code: AE3) (formula: $C_6H_{14}O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			9	6	3		
6	B	1	Total	C	O	0	0
			9	6	3		

- Molecule 7 is TETRAETHYLENE GLYCOL (three-letter code: PG4) (formula: $C_8H_{18}O_5$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	B	1	Total	C	O	0	0
			11	8	3		

- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	210	Total	O	0	0
			210	210		
8	B	150	Total	O	0	0
			150	150		

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	79.20Å 111.97Å 226.83Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.35 39.84 – 2.35	Depositor EDS
% Data completeness (in resolution range)	(Not available) (20.00-2.35) 99.0 (39.84-2.35)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.25 (at 2.34Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.209 , 0.232 0.210 , 0.236	Depositor DCC
R_{free} test set	1692 reflections (2.02%)	wwPDB-VP
Wilson B-factor (Å ²)	41.9	Xtriage
Anisotropy	0.847	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 40.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	8804	wwPDB-VP
Average B, all atoms (Å ²)	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.88% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: AE3, NAG, CO3, PG4, FUC, P6G

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	0.69	0/4300	0.80	0/5875
1	B	0.61	0/4282	0.78	1/5851 (0.0%)
All	All	0.65	0/8582	0.79	1/11726 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	1
All	All	0	2

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	B	161	LEU	N-CA-C	-5.82	95.28	111.00

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	449	TYR	Sidechain
1	B	70	TYR	Sidechain

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	4177	0	4063	72	0
1	B	4159	0	4046	85	0
2	A	28	0	25	0	0
2	B	14	0	13	0	0
3	A	10	0	10	0	0
4	A	4	0	0	1	0
4	B	4	0	0	1	0
5	A	19	0	26	1	0
6	A	9	0	14	0	0
6	B	9	0	14	2	0
7	B	11	0	12	0	0
8	A	210	0	0	6	0
8	B	150	0	0	2	0
All	All	8804	0	8223	154	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.

The worst 5 of 154 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:48:MET:HE3	1:A:166:GLU:HA	1.29	1.09
1:B:48:MET:HE1	1:B:166:GLU:HA	1.32	1.09
1:A:197:VAL:H	1:A:223:HIS:HD2	1.14	0.94
1:B:197:VAL:H	1:B:223:HIS:HD2	1.03	0.93
1:A:485:ARG:HG3	8:A:1107:HOH:O	1.70	0.92

There are no symmetry-related clashes.

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	531/543 (98%)	506 (95%)	25 (5%)	0	100	100
1	B	529/543 (97%)	507 (96%)	17 (3%)	5 (1%)	19	19
All	All	1060/1086 (98%)	1013 (96%)	42 (4%)	5 (0%)	31	34

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	542	ALA
1	B	496	LYS
1	B	494	ASP
1	B	497	SER
1	B	108	PRO

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	439/443 (99%)	423 (96%)	16 (4%)	38	47
1	B	438/443 (99%)	423 (97%)	15 (3%)	40	49
All	All	877/886 (99%)	846 (96%)	31 (4%)	39	48

5 of 31 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	424	ARG
1	B	136	ARG
1	B	496	LYS
1	A	525	ARG
1	B	203	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 6 such

sidechains are listed below:

Mol	Chain	Res	Type
1	A	527	GLN
1	B	499	GLN
1	B	223	HIS
1	A	291	GLN
1	B	291	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](#)

10 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	$\# Z > 2$
2	NAG	A	544	1,3	14,14,15	0.97	1 (7%)	17,19,21	1.18	3 (17%)
3	FUC	A	545	2	9,10,11	0.74	0	13,14,16	0.91	0
2	NAG	A	701	1	14,14,15	0.78	0	17,19,21	0.82	1 (5%)
5	P6G	A	901	-	18,18,18	2.01	6 (33%)	17,17,17	1.14	1 (5%)
6	AE3	A	902	-	8,8,8	1.53	2 (25%)	7,7,7	1.36	1 (14%)
4	CO3	A	951	-	0,3,3	0.00	-	0,3,3	0.00	-
2	NAG	B	601	1	14,14,15	1.02	1 (7%)	17,19,21	0.85	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	AE3	B	903	-	8,8,8	1.65	2 (25%)	7,7,7	1.49	1 (14%)
7	PG4	B	904	-	10,10,12	1.82	2 (20%)	9,9,11	0.83	0
4	CO3	B	952	-	0,3,3	0.00	-	0,3,3	0.00	-

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	NAG	A	544	1,3	-	0/6/23/26	0/1/1/1
3	FUC	A	545	2	-	0/0/17/20	0/1/1/1
2	NAG	A	701	1	-	0/6/23/26	0/1/1/1
5	P6G	A	901	-	-	0/16/16/16	0/0/0/0
6	AE3	A	902	-	-	0/6/6/6	0/0/0/0
4	CO3	A	951	-	-	0/0/0/0	0/0/0/0
2	NAG	B	601	1	-	0/6/23/26	0/1/1/1
6	AE3	B	903	-	-	0/6/6/6	0/0/0/0
7	PG4	B	904	-	-	0/8/8/10	0/0/0/0
4	CO3	B	952	-	-	0/0/0/0	0/0/0/0

The worst 5 of 14 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	601	NAG	C1-C2	2.14	1.55	1.52
2	A	544	NAG	C1-C2	2.43	1.55	1.52
6	A	902	AE3	O2-C3	2.50	1.53	1.42
7	B	904	PG4	O3-C5	2.68	1.54	1.42
5	A	901	P6G	O7-C6	2.75	1.54	1.42

The worst 5 of 7 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	701	NAG	C2-N2-C7	-2.36	119.50	122.94
2	A	544	NAG	C4-C3-C2	-2.32	107.62	111.02
2	A	544	NAG	C2-N2-C7	-2.29	119.60	122.94
2	A	544	NAG	C1-O5-C5	2.19	115.20	112.19
6	A	902	AE3	O4-C6-C5	3.04	129.76	111.91

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	901	P6G	1	0
4	A	951	CO3	1	0
6	B	903	AE3	2	0
4	B	952	CO3	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	535/543 (98%)	0.12	39 (7%) 15 23	30, 44, 66, 98	0
1	B	533/543 (98%)	0.31	44 (8%) 11 17	35, 49, 71, 99	0
All	All	1068/1086 (98%)	0.22	83 (7%) 13 20	30, 46, 69, 99	0

The worst 5 of 83 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	543	THR	6.4
1	B	497	SER	5.6
1	B	118	ILE	5.2
1	B	540	LEU	4.5
1	B	117	TRP	4.4

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(\AA^2)	Q<0.9
2	NAG	B	601	14/15	0.62	0.39	87,89,91,92	0
6	AE3	B	903	9/9	0.68	0.39	83,87,89,89	0
6	AE3	A	902	9/9	0.80	0.46	81,82,83,83	0
2	NAG	A	544	14/15	0.82	0.40	76,79,85,89	0
7	PG4	B	904	11/13	0.83	0.18	87,88,90,90	0
2	NAG	A	701	14/15	0.85	0.33	81,85,88,89	0
3	FUC	A	545	10/11	0.86	0.53	91,93,95,95	0
4	CO3	B	952	4/4	0.91	0.34	63,64,64,65	0
4	CO3	A	951	4/4	0.91	0.31	50,50,51,52	0
5	P6G	A	901	19/19	0.92	0.17	61,71,75,76	0

6.5 Other polymers [i](#)

There are no such residues in this entry.