



Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 11:04 AM GMT

PDB ID : 3D6N
Title : Crystal Structure of Aquifex Dihydroorotase Activated by Aspartate Transcarbamoylase
Authors : Edwards, B.F.P.
Deposited on : 2008-05-20
Resolution : 2.30 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

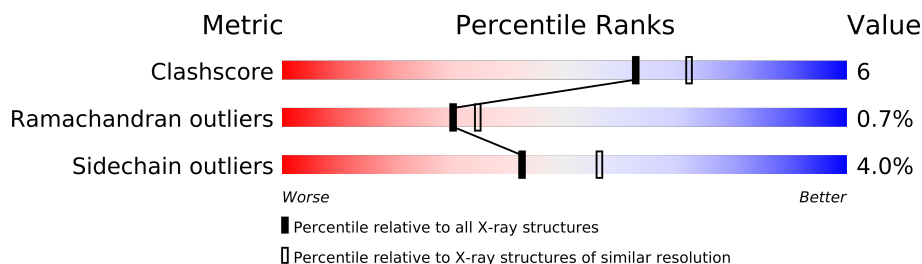
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : **FAILED**
Percentile statistics : 21963
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.30 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))
Clashscore	79885	3679 (2.30-2.30)
Ramachandran outliers	78287	3642 (2.30-2.30)
Sidechain outliers	78261	3641 (2.30-2.30)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Note EDS failed to run properly.

Mol	Chain	Length	Quality of chain
1	A	422	
2	B	291	

2 Entry composition i

There are 5 unique types of molecules in this entry. The entry contains 6081 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Dihydroorotase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	422	Total	C	N	O	S	0	0	0
			3256	2071	554	615	16			

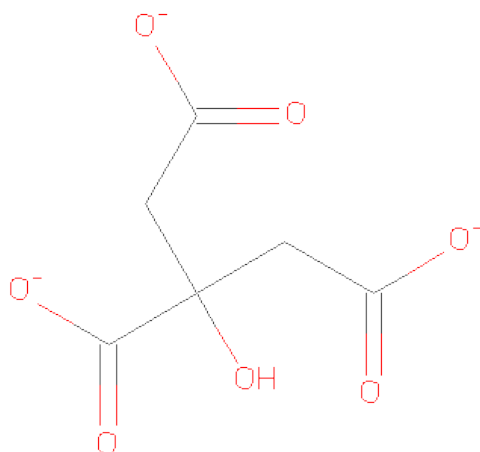
- Molecule 2 is a protein called Aspartate carbamoyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	291	Total	C	N	O	S	0	0	0
			2376	1541	395	436	4			

- 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 CITRATE ANION (three-letter code: FLC) (formula: C₆H₅O₇).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total 13	C 6	O 7	0	0
4	B	1	Total 13	C 6	O 7	0	0

- Molecule 5 is water.

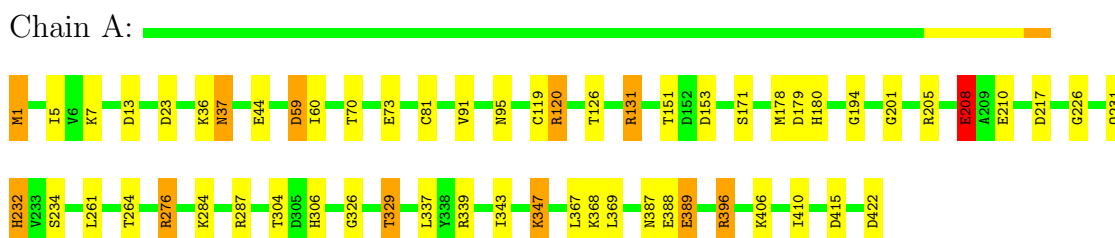
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	209	Total 209	O 209	0	0
5	B	213	Total 213	O 213	0	0

3 Residue-property plots

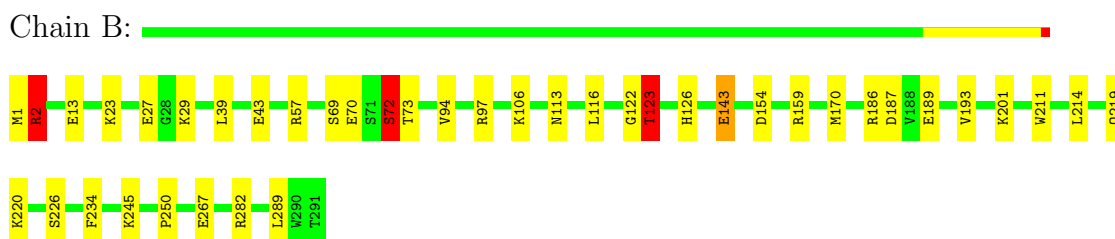
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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.

Note EDS failed to run properly.

- Molecule 1: Dihydroorotase



- Molecule 2: Aspartate carbamoyltransferase



4 Data and refinement statistics

EDS failed to run properly - this section will therefore be incomplete.

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, α , β , γ	158.44Å 158.44Å 233.55Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	65.00 – 2.30	Depositor
% Data completeness (in resolution range)	100.0 (65.00-2.30)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	0.80	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.02 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.164 , 0.204	Depositor
Wilson B-factor (Å ²)	32.3	Xtriage
Anisotropy	0.052	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	1 of 50158 reflections (0.002%)	Xtriage
Total number of atoms	6081	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 2.72% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, FLC

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	0.85	3/3310 (0.1%)	0.85	6/4472 (0.1%)
2	B	1.10	4/2431 (0.2%)	0.95	8/3274 (0.2%)
All	All	0.97	7/5741 (0.1%)	0.89	14/7746 (0.2%)

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
2	B	1	1

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	143	GLU	CB-CG	9.72	1.70	1.52
2	B	143	GLU	CG-CD	8.23	1.64	1.51
1	A	119	CYS	CB-SG	-7.00	1.70	1.82
1	A	81	CYS	CB-SG	-6.58	1.71	1.82
2	B	13	GLU	CG-CD	5.79	1.60	1.51
2	B	245	LYS	CE-NZ	5.46	1.62	1.49
1	A	208	GLU	CB-CG	5.06	1.61	1.52

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	131	ARG	NE-CZ-NH2	-9.04	115.78	120.30
1	A	205	ARG	NE-CZ-NH2	-8.94	115.83	120.30
2	B	123	THR	OG1-CB-CG2	6.23	124.33	110.00
2	B	123	THR	N-CA-CB	-6.19	98.54	110.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	2	ARG	NE-CZ-NH1	6.12	123.36	120.30
1	A	205	ARG	NE-CZ-NH1	5.98	123.29	120.30
2	B	57	ARG	NE-CZ-NH2	-5.79	117.41	120.30
1	A	231	GLN	C-N-CA	-5.75	107.34	121.70
1	A	59	ASP	CB-CG-OD1	5.62	123.35	118.30
2	B	214	LEU	CA-CB-CG	-5.52	102.59	115.30
2	B	282	ARG	NE-CZ-NH2	-5.44	117.58	120.30
1	A	131	ARG	NE-CZ-NH1	5.37	122.99	120.30
2	B	143	GLU	OE1-CD-OE2	-5.26	116.98	123.30
2	B	154	ASP	CB-CG-OD1	5.18	122.96	118.30

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	B	123	THR	CB

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	72	SER	Peptide

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3256	0	3362	35	0
2	B	2376	0	2393	32	0
3	A	1	0	0	0	0
4	A	13	0	5	1	0
4	B	13	0	5	0	0
5	A	209	0	0	9	1
5	B	213	0	0	8	1
All	All	6081	0	5765	64	2

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 6.

All (64) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:120:ARG:HD3	5:A:579:HOH:O	1.36	1.24
2:B:69:SER:HB3	2:B:72:SER:HA	1.20	1.09
2:B:70:GLU:OE1	5:B:491:HOH:O	1.81	0.97
2:B:1:MET:HE1	5:B:425:HOH:O	1.65	0.95
2:B:2:ARG:HG3	2:B:289:LEU:HD22	1.58	0.85
2:B:69:SER:CB	2:B:72:SER:HA	2.05	0.84
2:B:211:TRP:HZ3	2:B:234:PHE:O	1.64	0.79
1:A:201:GLY:O	2:B:123:THR:CG2	2.33	0.77
2:B:2:ARG:HG3	2:B:289:LEU:CD2	2.17	0.74
1:A:59:ASP:OD2	1:A:304:THR:HG23	1.91	0.71
1:A:232:HIS:NE2	5:A:577:HOH:O	2.15	0.70
2:B:69:SER:HB3	2:B:72:SER:CA	2.12	0.70
1:A:339:ARG:NH2	1:A:406:LYS:O	2.26	0.69
2:B:126:HIS:HD2	5:B:427:HOH:O	1.76	0.69
2:B:2:ARG:HH11	2:B:2:ARG:HG2	1.58	0.67
2:B:70:GLU:HA	5:B:494:HOH:O	1.97	0.65
1:A:179:ASP:OD2	1:A:217:ASP:OD2	2.15	0.65
2:B:2:ARG:CG	2:B:2:ARG:HH11	2.12	0.62
1:A:264:THR:HG21	1:A:284:LYS:HE3	1.83	0.61
2:B:23:LYS:O	2:B:27:GLU:HG3	1.99	0.61
1:A:178:MET:HE3	5:A:468:HOH:O	2.01	0.60
1:A:178:MET:CE	5:A:468:HOH:O	2.51	0.59
1:A:276:ARG:NH2	5:A:547:HOH:O	2.27	0.58
1:A:70:THR:HA	1:A:73:GLU:O	2.05	0.57
2:B:2:ARG:CG	2:B:289:LEU:CD2	2.83	0.56
1:A:201:GLY:O	2:B:123:THR:HG21	2.07	0.54
4:A:424:FLC:OA1	4:A:424:FLC:CGC	2.56	0.53
2:B:2:ARG:NH1	2:B:2:ARG:HG2	2.22	0.53
1:A:387:ASN:OD1	1:A:389:GLU:HG2	2.09	0.52
2:B:219:GLN:HG3	2:B:226:SER:HA	1.92	0.52
1:A:388:GLU:OE1	1:A:396:ARG:NH2	2.43	0.52
1:A:343:ILE:HB	1:A:347:LYS:HG2	1.91	0.51
2:B:211:TRP:CZ3	2:B:234:PHE:O	2.55	0.50
2:B:43:GLU:HB3	2:B:97:ARG:HG2	1.92	0.50
2:B:122:GLY:O	2:B:159:ARG:HD3	2.11	0.50
1:A:120:ARG:NH2	5:A:562:HOH:O	2.45	0.50
1:A:261:LEU:HD13	1:A:337:LEU:HD11	1.95	0.49
1:A:7:LYS:HA	1:A:23:ASP:OD1	2.13	0.49
1:A:208:GLU:OE2	2:B:186:ARG:NH2	2.43	0.48
1:A:329:THR:HG23	5:A:530:HOH:O	2.12	0.48
1:A:276:ARG:NE	5:A:547:HOH:O	2.36	0.48
1:A:1:MET:SD	1:A:44:GLU:HG3	2.53	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:2:ARG:HG2	2:B:289:LEU:HD23	1.96	0.47
1:A:210:GLU:OE1	1:A:234:SER:OG	2.28	0.47
1:A:59:ASP:OD2	1:A:304:THR:CG2	2.61	0.47
1:A:60:ILE:HA	1:A:91:VAL:HB	1.97	0.47
1:A:95:ASN:HB3	1:A:153:ASP:O	2.15	0.47
1:A:13:ASP:HB2	1:A:367:LEU:HD12	1.96	0.47
2:B:2:ARG:CG	2:B:289:LEU:HD23	2.46	0.46
1:A:410:ILE:HA	1:A:422:ASP:HB2	1.96	0.46
1:A:194:GLY:HA2	2:B:187:ASP:OD2	2.15	0.46
2:B:267:GLU:HB2	5:B:416:HOH:O	2.16	0.45
1:A:180:HIS:ND1	5:A:577:HOH:O	2.36	0.45
2:B:126:HIS:CD2	5:B:427:HOH:O	2.61	0.43
1:A:126:THR:HG21	1:A:131:ARG:HG3	1.99	0.43
1:A:208:GLU:HG3	1:A:208:GLU:H	1.32	0.43
2:B:94:VAL:HB	2:B:116:LEU:HD23	1.99	0.43
2:B:170:MET:HE2	2:B:170:MET:HB3	1.99	0.42
1:A:284:LYS:HE2	1:A:287:ARG:HH11	1.84	0.42
1:A:126:THR:HA	1:A:151:THR:O	2.20	0.42
2:B:201:LYS:HZ3	2:B:201:LYS:HG3	1.75	0.41
2:B:113:ASN:HB2	5:B:482:HOH:O	2.21	0.41
2:B:1:MET:CE	5:B:425:HOH:O	2.45	0.40
1:A:171:SER:HB3	1:A:226:GLY:HA2	2.04	0.40

All (2) 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	Distance(Å)	Clash(Å)
5:B:487:HOH:O	5:B:492:HOH:O[17_555]	2.11	0.09
5:A:539:HOH:O	5:A:543:HOH:O[5_556]	2.18	0.02

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	420/422 (100%)	398 (95%)	19 (4%)	3 (1%)	30	34

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	B	289/291 (99%)	276 (96%)	11 (4%)	2 (1%)	30	34
All	All	709/713 (99%)	674 (95%)	30 (4%)	5 (1%)	30	34

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	72	SER
1	A	37	ASN
1	A	306	HIS
2	B	250	PRO
1	A	326	GLY

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	362/362 (100%)	347 (96%)	15 (4%)	41	55
2	B	262/262 (100%)	252 (96%)	10 (4%)	44	59
All	All	624/624 (100%)	599 (96%)	25 (4%)	42	56

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	5	ILE
1	A	36	LYS
1	A	37	ASN
1	A	120	ARG
1	A	208	GLU
1	A	232	HIS
1	A	276	ARG
1	A	329	THR
1	A	347	LYS
1	A	368	LYS
1	A	369	LEU
1	A	389	GLU

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Mol	Chain	Res	Type
1	A	396	ARG
1	A	415	ASP
2	B	2	ARG
2	B	29	LYS
2	B	39	LEU
2	B	73	THR
2	B	106	LYS
2	B	123	THR
2	B	143	GLU
2	B	189	GLU
2	B	193	VAL
2	B	220	LYS

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

Mol	Chain	Res	Type
2	B	126	HIS

5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 3 ligands modelled in this entry, 1 is monoatomic - leaving 2 for Mogul analysis.

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	$\# Z > 2$
4	FLC	A	424	3	5,12,12	5.98	2 (40%)	7,17,17	1.17	0
4	FLC	B	292	-	5,12,12	4.93	2 (40%)	7,17,17	1.26	1 (14%)

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
4	FLC	A	424	3	-	0/6/16/16	0/0/0/0
4	FLC	B	292	-	-	0/6/16/16	0/0/0/0

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	424	FLC	CA-CAC	10.11	1.55	1.49
4	A	424	FLC	CG-CGC	8.48	1.54	1.49
4	B	292	FLC	CG-CGC	7.78	1.54	1.49
4	B	292	FLC	CA-CAC	7.75	1.54	1.49

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	292	FLC	CB-CG-CGC	-2.79	110.84	115.01

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

EDS failed to run properly - this section will therefore be empty.

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

EDS failed to run properly - this section will therefore be empty.

6.3 Carbohydrates ⓘ

EDS failed to run properly - this section will therefore be empty.

6.4 Ligands ⓘ

EDS failed to run properly - this section will therefore be empty.

6.5 Other polymers ⓘ

EDS failed to run properly - this section will therefore be empty.