



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 10:33 PM GMT

PDB ID : 1U6H  
Title : Vinculin head (0-258) in complex with the talin vinculin binding site 2 (849-879)  
Authors : Fillingham, I.; Gingras, A.R.; Papagrigoriou, E.; Patel, B.; Emsley, J.; Roberts, G.C.K.; Critchley, D.R.; Barsukov, I.L.  
Deposited on : 2004-07-30  
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  
Mogul : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : **NOT EXECUTED**  
EDS : **NOT EXECUTED**  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

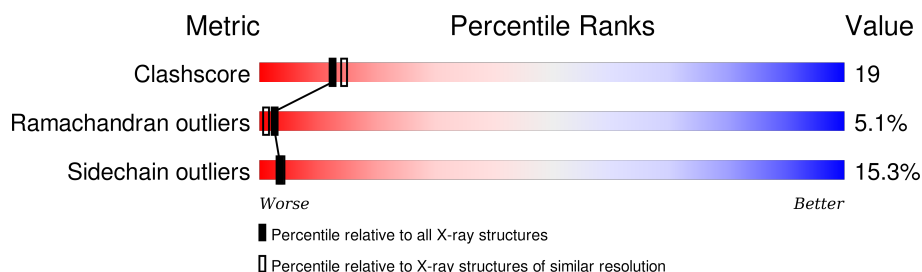
# 1 Overall quality at a glance

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))
Clashscore	102246	4595 (2.40-2.36)
Ramachandran outliers	100387	4520 (2.40-2.36)
Sidechain outliers	100360	4522 (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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	280	
2	B	31	

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2172 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 Vinculin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	252	1947	1225	332	376	14	0	0	0

There are 21 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-21	MET	-	EXPRESSION TAG	UNP P12003
A	-20	GLY	-	EXPRESSION TAG	UNP P12003
A	-19	SER	-	EXPRESSION TAG	UNP P12003
A	-18	SER	-	EXPRESSION TAG	UNP P12003
A	-17	HIS	-	EXPRESSION TAG	UNP P12003
A	-16	HIS	-	EXPRESSION TAG	UNP P12003
A	-15	HIS	-	EXPRESSION TAG	UNP P12003
A	-14	HIS	-	EXPRESSION TAG	UNP P12003
A	-13	HIS	-	EXPRESSION TAG	UNP P12003
A	-12	HIS	-	EXPRESSION TAG	UNP P12003
A	-11	SER	-	EXPRESSION TAG	UNP P12003
A	-10	SER	-	EXPRESSION TAG	UNP P12003
A	-9	GLY	-	EXPRESSION TAG	UNP P12003
A	-8	LEU	-	EXPRESSION TAG	UNP P12003
A	-7	VAL	-	EXPRESSION TAG	UNP P12003
A	-6	PRO	-	EXPRESSION TAG	UNP P12003
A	-5	ARG	-	EXPRESSION TAG	UNP P12003
A	-4	GLY	-	EXPRESSION TAG	UNP P12003
A	-3	SER	-	EXPRESSION TAG	UNP P12003
A	-2	HIS	-	EXPRESSION TAG	UNP P12003
A	-1	MET	-	EXPRESSION TAG	UNP P12003

- Molecule 2 is a protein called Talin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	24	Total 168	C 106	N 31	O 30	S 1	0	0	0

- Molecule 3 is water.

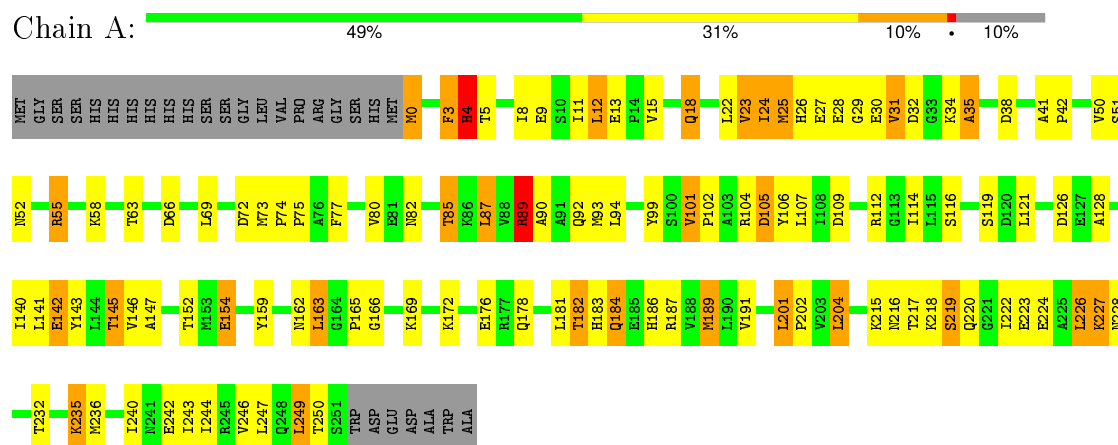
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	52	Total 52	O 52	0	0
3	B	5	Total 5	O 5	0	0

### 3 Residue-property plots

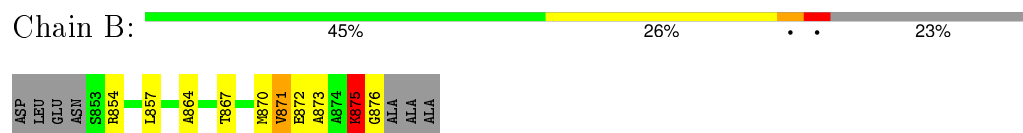
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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 was not executed.

#### • Molecule 1: Vinculin



#### • Molecule 2: Talin



## 4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	52.12Å 69.15Å 96.16Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.38	Depositor
% Data completeness (in resolution range)	99.7 (30.00-2.38)	Depositor
$R_{merge}$	0.12	Depositor
$R_{sym}$	(Not available)	Depositor
Refinement program	REFMAC 5.2.0003	Depositor
R, $R_{free}$	0.242 , 0.322	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2172	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	48.0	wwPDB-VP

## 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	0.91	0/1970	1.10	13/2669 (0.5%)
2	B	0.95	0/167	1.06	0/221
All	All	0.91	0/2137	1.10	13/2890 (0.4%)

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

There are no bond length outliers.

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	126	ASP	CB-CG-OD2	8.93	126.34	118.30
1	A	105	ASP	CB-CG-OD2	6.92	124.53	118.30
1	A	109	ASP	CB-CG-OD2	6.68	124.31	118.30
1	A	72	ASP	CB-CG-OD2	6.67	124.30	118.30
1	A	66	ASP	CB-CG-OD2	6.41	124.07	118.30
1	A	93	MET	CG-SD-CE	6.26	110.22	100.20
1	A	141	LEU	CA-CB-CG	5.87	128.80	115.30
1	A	227	LYS	CD-CE-NZ	-5.75	98.47	111.70
1	A	226	LEU	CA-CB-CG	5.56	128.08	115.30
1	A	89	ARG	CG-CD-NE	5.40	123.14	111.80
1	A	4	HIS	N-CA-C	5.26	125.22	111.00
1	A	38	ASP	CB-CG-OD2	5.25	123.02	118.30
1	A	204	LEU	CA-CB-CG	5.05	126.91	115.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	3	PHE	Peptide

## 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	1947	0	2009	79	0
2	B	168	0	188	9	0
3	A	52	0	0	7	0
3	B	5	0	0	0	0
All	All	2172	0	2197	84	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 19.

All (84) 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:184:GLN:HG2	3:A:301:HOH:O	1.42	1.17
1:A:9:GLU:O	1:A:13:GLU:HB2	1.82	0.79
1:A:89:ARG:HG2	1:A:89:ARG:HH11	1.47	0.79
1:A:178:GLN:HG3	1:A:191:VAL:HG23	1.66	0.76
1:A:89:ARG:CG	1:A:89:ARG:HH11	1.99	0.75
3:A:290:HOH:O	2:B:876:GLY:HA3	1.90	0.71
1:A:23:VAL:HG21	1:A:112:ARG:HH11	1.55	0.70
1:A:172:LYS:O	1:A:176:GLU:HG3	1.94	0.68
1:A:119:SER:HB3	3:A:268:HOH:O	1.93	0.67
1:A:215:LYS:O	1:A:216:ASN:ND2	2.28	0.65
1:A:217:THR:O	1:A:219:SER:N	2.30	0.64
1:A:73:MET:HB3	1:A:74:PRO:HD3	1.80	0.63
1:A:152:THR:CG2	1:A:154:GLU:HG3	2.27	0.63
1:A:215:LYS:C	1:A:216:ASN:HD22	2.03	0.61
1:A:25:MET:O	1:A:30:GLU:HB3	2.02	0.60
1:A:18:GLN:NE2	2:B:872:GLU:HG2	2.16	0.60
1:A:4:HIS:HD2	1:A:8:ILE:HD12	1.66	0.60
1:A:63:THR:HG22	1:A:63:THR:O	2.02	0.59
1:A:104:ARG:HG3	1:A:105:ASP:N	2.18	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:178:GLN:HG3	1:A:191:VAL:CG2	2.33	0.59
1:A:89:ARG:HD3	1:A:92:GLN:NE2	2.18	0.58
1:A:11:ILE:HG22	1:A:12:LEU:HD13	1.84	0.58
1:A:152:THR:HG21	1:A:154:GLU:HG3	1.85	0.58
1:A:18:GLN:HE21	2:B:872:GLU:HG2	1.69	0.57
1:A:140:ILE:HD11	1:A:166:GLY:O	2.04	0.56
1:A:3:PHE:HB3	1:A:8:ILE:HG22	1.87	0.56
1:A:186:HIS:ND1	3:A:279:HOH:O	2.33	0.56
1:A:236:MET:O	1:A:240:ILE:HG13	2.06	0.56
2:B:875:LYS:HD2	2:B:875:LYS:N	2.19	0.55
1:A:23:VAL:HG12	1:A:28:GLU:HG3	1.89	0.55
1:A:22:LEU:N	3:A:290:HOH:O	2.39	0.54
1:A:246:VAL:HA	1:A:249:LEU:HD22	1.90	0.54
1:A:30:GLU:HG3	1:A:31:VAL:N	2.23	0.53
1:A:201:LEU:HB3	1:A:202:PRO:HD3	1.90	0.53
1:A:142:GLU:O	1:A:145:THR:HB	2.08	0.53
1:A:9:GLU:O	1:A:13:GLU:CB	2.56	0.52
1:A:82:ASN:O	1:A:85:THR:HG22	2.11	0.51
1:A:101:VAL:N	1:A:102:PRO:CD	2.74	0.50
1:A:58:LYS:HG2	1:A:77:PHE:CE1	2.47	0.50
1:A:163:LEU:O	1:A:163:LEU:HG	2.12	0.50
1:A:69:LEU:HD13	1:A:128:ALA:HB2	1.95	0.49
1:A:15:VAL:CG2	2:B:864:ALA:HB1	2.43	0.49
2:B:867:THR:O	2:B:871:VAL:HG12	2.12	0.49
1:A:89:ARG:CD	1:A:92:GLN:NE2	2.75	0.49
1:A:232:THR:O	1:A:236:MET:HG3	2.13	0.48
1:A:178:GLN:HE21	1:A:187:ARG:HD2	1.78	0.48
1:A:143:TYR:OH	1:A:162:ASN:HB2	2.14	0.48
1:A:23:VAL:CG2	1:A:112:ARG:HH11	2.25	0.47
1:A:140:ILE:HD11	1:A:166:GLY:C	2.34	0.47
1:A:89:ARG:HG2	1:A:89:ARG:NH1	2.25	0.47
1:A:74:PRO:HB2	1:A:75:PRO:HD3	1.97	0.47
1:A:89:ARG:HD3	1:A:92:GLN:HE22	1.80	0.46
1:A:26:HIS:O	1:A:29:GLY:O	2.34	0.46
1:A:35:ALA:HB2	1:A:99:TYR:CE2	2.51	0.46
1:A:0:MET:O	1:A:0:MET:HG3	2.14	0.46
1:A:178:GLN:O	1:A:187:ARG:CD	2.64	0.45
1:A:30:GLU:HG3	1:A:31:VAL:H	1.81	0.45
1:A:31:VAL:O	1:A:34:LYS:CB	2.64	0.45
1:A:152:THR:HG22	1:A:154:GLU:H	1.82	0.45
1:A:119:SER:CB	3:A:268:HOH:O	2.60	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:4:HIS:CD2	1:A:5:THR:HG23	2.52	0.45
1:A:73:MET:HB3	1:A:74:PRO:CD	2.47	0.44
2:B:873:ALA:C	2:B:875:LYS:H	2.21	0.44
1:A:181:LEU:HB2	1:A:187:ARG:HG2	2.00	0.44
1:A:147:ALA:HB2	1:A:159:TYR:OH	2.19	0.43
1:A:152:THR:HG21	1:A:154:GLU:CG	2.49	0.43
1:A:63:THR:CG2	1:A:63:THR:O	2.67	0.43
1:A:4:HIS:HE1	3:A:267:HOH:O	2.02	0.43
1:A:189:MET:CE	1:A:246:VAL:HG22	2.49	0.43
1:A:73:MET:HG2	1:A:77:PHE:CZ	2.55	0.42
1:A:114:ILE:HG21	2:B:867:THR:OG1	2.19	0.42
1:A:4:HIS:CE1	1:A:250:THR:HG1	2.38	0.42
1:A:235:LYS:HA	1:A:235:LYS:HD2	1.81	0.41
2:B:870:MET:O	2:B:870:MET:HG2	2.18	0.41
1:A:52:ASN:ND2	1:A:55:ARG:HE	2.18	0.41
1:A:242:GLU:O	1:A:243:ILE:C	2.58	0.41
1:A:34:LYS:O	1:A:35:ALA:HB3	2.20	0.41
1:A:244:ILE:HA	1:A:247:LEU:HD12	2.02	0.41
1:A:41:ALA:N	1:A:42:PRO:HD2	2.35	0.41
1:A:4:HIS:CE1	1:A:250:THR:OG1	2.74	0.41
1:A:87:LEU:O	1:A:90:ALA:HB3	2.21	0.41
1:A:50:VAL:HG13	1:A:80:VAL:HG12	2.03	0.40
1:A:182:THR:HG22	1:A:183:HIS:ND1	2.36	0.40
1:A:121:LEU:C	1:A:121:LEU:HD23	2.41	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	250/280 (89%)	216 (86%)	21 (8%)	13 (5%)	<b>2</b> <b>1</b>

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	B	22/31 (71%)	19 (86%)	2 (9%)	1 (4%)	3	2
All	All	272/311 (88%)	235 (86%)	23 (8%)	14 (5%)	2	1

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	4	HIS
1	A	23	VAL
1	A	24	ILE
1	A	25	MET
1	A	218	LYS
1	A	27	GLU
1	A	220	GLN
2	B	875	LYS
1	A	31	VAL
1	A	35	ALA
1	A	55	ARG
1	A	163	LEU
1	A	219	SER
1	A	165	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	220/247 (89%)	188 (86%)	32 (14%)	4	4
2	B	15/20 (75%)	11 (73%)	4 (27%)	0	0
All	All	235/267 (88%)	199 (85%)	36 (15%)	3	4

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	0	MET
1	A	12	LEU

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Mol	Chain	Res	Type
1	A	18	GLN
1	A	24	ILE
1	A	32	ASP
1	A	51	SER
1	A	85	THR
1	A	87	LEU
1	A	89	ARG
1	A	94	LEU
1	A	101	VAL
1	A	106	TYR
1	A	107	LEU
1	A	116	SER
1	A	142	GLU
1	A	145	THR
1	A	146	VAL
1	A	154	GLU
1	A	169	LYS
1	A	182	THR
1	A	184	GLN
1	A	189	MET
1	A	201	LEU
1	A	204	LEU
1	A	222	ILE
1	A	223	GLU
1	A	224	GLU
1	A	226	LEU
1	A	227	LYS
1	A	228	ASN
1	A	235	LYS
1	A	249	LEU
2	B	854	ARG
2	B	857	LEU
2	B	871	VAL
2	B	875	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	18	GLN
1	A	26	HIS
1	A	47	GLN
1	A	52	ASN

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Mol	Chain	Res	Type
1	A	62	GLN
1	A	92	GLN
1	A	178	GLN
1	A	186	HIS
1	A	192	ASN
1	A	195	ASN
1	A	216	ASN
1	A	228	ASN
1	A	230	ASN
1	A	248	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](#)

There are no ligands in this entry.

## 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 ⓘ

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

### 6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

### 6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

### 6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.