



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 17, 2026 – 11:04 PM UTC

PDB ID : 1A47 / pdb_00001a47
Title : CGTASE FROM THERMOANAEROBACTERIUM THERMOSULFURIGENES EM1 IN COMPLEX WITH A MALTOHEXAOSE INHIBITOR
Authors : Uitdehaag, J.C.M.; Kalk, K.H.; Rozeboom, H.J.; Dijkstra, B.W.
Deposited on : 1998-02-11
Resolution : 2.56 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

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

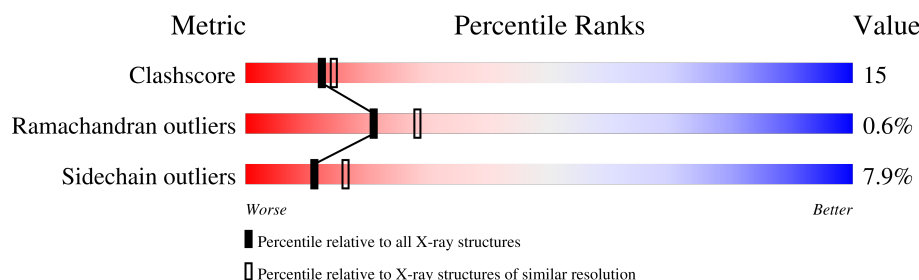
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.56 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	190562	1897 (2.58-2.54)
Ramachandran outliers	187476	1875 (2.58-2.54)
Sidechain outliers	187428	1875 (2.58-2.54)

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 of 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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	683	68% 28% .
2	B	3	100%
3	C	3	33% 67%
4	D	2	50% 50%

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 5620 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 CYCLODEXTRIN GLYCOSYLTRANSFERASE.

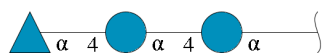
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	683	Total	C	N	O	S	0	0	0
			5333	3377	885	1056	15			

- Molecule 2 is an oligosaccharide called alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose.



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	B	3	Total	C	O	0	0	0
			34	18	16			

- Molecule 3 is an oligosaccharide called alpha-D-quinovopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose.



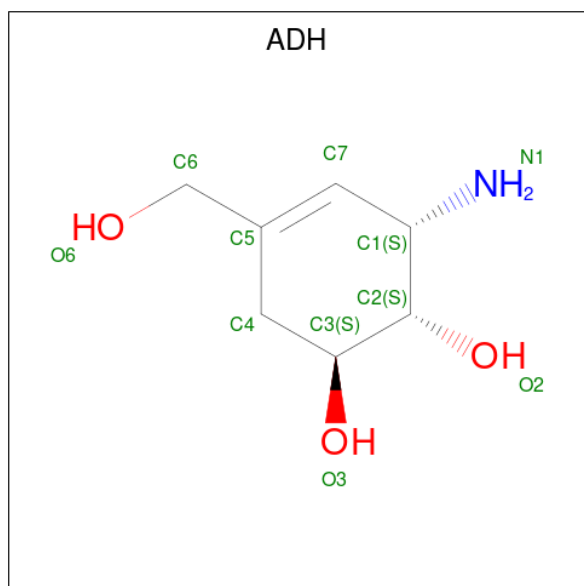
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
3	C	3	Total	C	O	0	0	0
			32	18	14			

- Molecule 4 is an oligosaccharide called alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose.



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
4	D	2	Total	C	O	0	0	0
			23	12	11			

- Molecule 5 is 1-AMINO-2,3-DIHYDROXY-5-HYDROXYMETHYL CYCLOHEX-5-ENE (CCD ID: ADH) (formula: $C_7H_{13}NO_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	0	0
			11	7	1	3		

- Molecule 6 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	2	Total	Ca	0	0
			2	2		

- Molecule 7 is water.

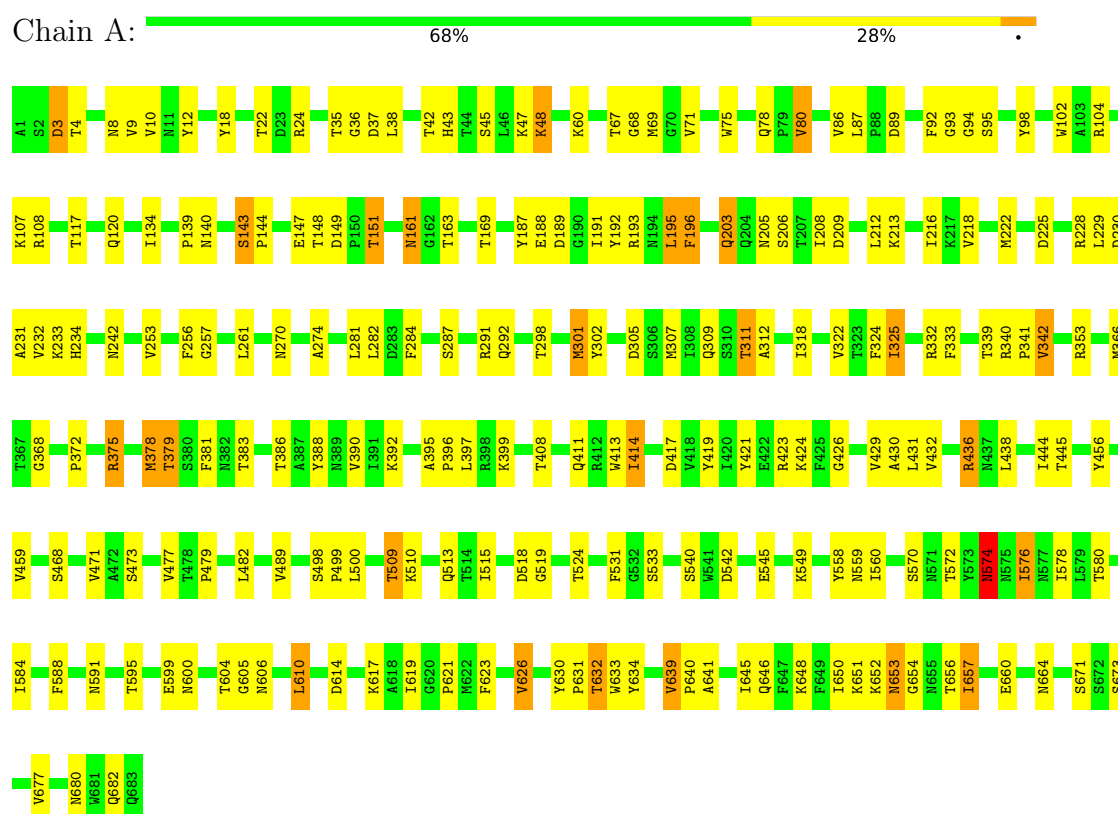
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	185	Total	O	0	0
			185	185		

3 Residue-property plots

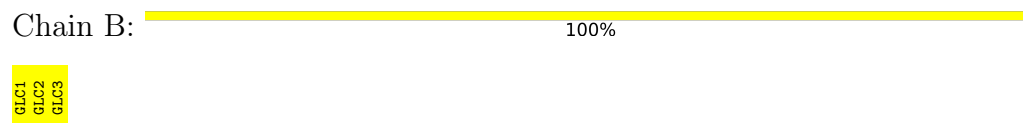
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded 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. 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: CYCLODEXTRIN GLYCOSYLTRANSFERASE



• Molecule 2: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose



• Molecule 3: alpha-D-quinovopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose





- Molecule 4: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose

Chain D: 50% 50%



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	74.00Å 97.79Å 116.26Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	8.00 – 2.56	Depositor
% Data completeness (in resolution range)	84.0 (8.00-2.56)	Depositor
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
Refinement program	TNT 5D	Depositor
R, R_{free}	0.186 , 0.250	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	5620	wwPDB-VP
Average B, all atoms (Å ²)	18.0	wwPDB-VP

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: GLC, CA, G6D, ADH

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	0.67	0/5470	0.94	3/7466 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	605	GLY	CA-C-N	6.93	129.88	120.38
1	A	605	GLY	C-N-CA	6.93	129.88	120.38
1	A	574	ASN	N-CA-C	5.12	118.10	110.52

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	5333	0	5056	159	0
2	B	34	0	30	0	0
3	C	32	0	21	1	0
4	D	23	0	20	0	0
5	A	11	0	10	3	0
6	A	2	0	0	0	0
7	A	185	0	0	4	0
All	All	5620	0	5137	162	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 15.

All (162) 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:311:THR:HG21	1:A:353:ARG:HH21	1.21	1.00
5:A:694:ADH:N1	5:A:694:ADH:C1	2.30	0.93
1:A:75:TRP:CZ2	1:A:228:ARG:HD3	2.06	0.91
1:A:301:MET:CG	1:A:419:TYR:HB2	2.02	0.88
1:A:311:THR:HG21	1:A:353:ARG:NH2	1.89	0.88
1:A:3:ASP:HB3	1:A:518:ASP:CG	2.02	0.85
1:A:3:ASP:HA	1:A:8:ASN:HD22	1.41	0.84
1:A:648:LYS:HD2	1:A:660:GLU:HB3	1.60	0.83
1:A:652:LYS:HG2	1:A:657:ILE:HD13	1.64	0.79
1:A:445:THR:HG22	1:A:479:PRO:HB3	1.63	0.79
1:A:301:MET:HG2	1:A:419:TYR:HB2	1.64	0.78
1:A:444:ILE:HD13	1:A:482:LEU:HB2	1.63	0.78
1:A:301:MET:HG3	1:A:419:TYR:HB2	1.66	0.77
1:A:386:THR:O	1:A:390:VAL:HG23	1.86	0.75
1:A:9:VAL:HG13	1:A:225:ASP:O	1.88	0.73
1:A:206:SER:OG	1:A:673:SER:HB2	1.89	0.73
1:A:417:ASP:HA	1:A:436:ARG:HD2	1.70	0.73
1:A:287:SER:O	1:A:291:ARG:HG3	1.89	0.72
1:A:9:VAL:HG22	1:A:225:ASP:HB3	1.72	0.72
1:A:307:MET:O	1:A:311:THR:HB	1.90	0.72
1:A:498:SER:HB2	1:A:499:PRO:HD2	1.71	0.70
1:A:209:ASP:O	1:A:213:LYS:HG3	1.91	0.70
1:A:379:THR:HA	7:A:868:HOH:O	1.93	0.69
1:A:558:TYR:CD1	1:A:578:ILE:HD12	2.29	0.68
1:A:499:PRO:HB2	1:A:570:SER:HB3	1.76	0.68
1:A:397:LEU:HD11	1:A:459:VAL:HG11	1.76	0.68
1:A:432:VAL:HG13	1:A:489:VAL:HG22	1.75	0.68
1:A:584:ILE:CG1	1:A:641:ALA:HB2	2.25	0.67
1:A:9:VAL:HG11	1:A:253:VAL:HG12	1.77	0.66
1:A:660:GLU:HG2	1:A:682:GLN:HB2	1.76	0.66
1:A:444:ILE:CD1	1:A:482:LEU:HB2	2.26	0.64
1:A:652:LYS:HD2	1:A:657:ILE:HD11	1.79	0.64
1:A:595:THR:HB	1:A:599:GLU:HB3	1.82	0.62
1:A:632:THR:HG22	7:A:864:HOH:O	1.99	0.61
1:A:417:ASP:O	1:A:436:ARG:HG3	2.00	0.61
1:A:584:ILE:HG13	1:A:641:ALA:HB2	1.82	0.61
1:A:189:ASP:OD1	1:A:193:ARG:HD2	2.01	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:3:ASP:HA	1:A:8:ASN:ND2	2.13	0.60
1:A:274:ALA:HB2	1:A:281:LEU:HD12	1.81	0.60
1:A:456:TYR:O	1:A:468:SER:HB2	2.01	0.60
1:A:630:TYR:CG	1:A:631:PRO:HA	2.37	0.60
1:A:187:TYR:O	1:A:191:ILE:HG13	2.02	0.60
1:A:614:ASP:OD2	1:A:617:LYS:HD3	2.02	0.60
1:A:89:ASP:O	1:A:93:GLY:HA2	2.02	0.59
1:A:648:LYS:CD	1:A:660:GLU:HB3	2.32	0.59
1:A:231:ALA:HB1	1:A:234:HIS:HD2	1.68	0.58
1:A:38:LEU:HD23	1:A:86:VAL:HG22	1.84	0.58
1:A:228:ARG:HG3	1:A:256:PHE:CE2	2.38	0.58
1:A:595:THR:HG22	1:A:651:LYS:HE2	1.86	0.58
1:A:47:LYS:HG2	1:A:372:PRO:HG3	1.85	0.57
1:A:9:VAL:HG12	1:A:253:VAL:HA	1.87	0.57
1:A:584:ILE:HG12	1:A:641:ALA:HB2	1.87	0.57
1:A:395:ALA:HB3	1:A:396:PRO:HD3	1.87	0.56
1:A:139:PRO:HB3	1:A:212:LEU:HD22	1.88	0.56
5:A:694:ADH:N1	5:A:694:ADH:C7	2.69	0.55
1:A:3:ASP:HB3	1:A:518:ASP:OD1	2.07	0.55
1:A:424:LYS:HE3	1:A:426:GLY:O	2.07	0.54
1:A:120:GLN:HG3	1:A:222:MET:HE1	1.88	0.54
1:A:378:MET:HE3	1:A:381:PHE:CE2	2.43	0.54
1:A:574:ASN:ND2	1:A:574:ASN:H	2.06	0.54
1:A:499:PRO:HB2	1:A:570:SER:CB	2.38	0.53
1:A:515:ILE:HD11	1:A:531:PHE:CZ	2.43	0.53
1:A:342:VAL:HG22	7:A:715:HOH:O	2.09	0.53
1:A:509:THR:HG22	1:A:513:GLN:OE1	2.09	0.53
1:A:230:ASP:OD1	5:A:694:ADH:H1	2.09	0.53
1:A:340:ARG:N	1:A:341:PRO:HD2	2.24	0.52
1:A:652:LYS:CG	1:A:657:ILE:HD13	2.37	0.51
1:A:102:TRP:HB3	1:A:143:SER:HB3	1.92	0.51
1:A:292:GLN:HG2	1:A:298:THR:OG1	2.10	0.51
1:A:149:ASP:OD2	1:A:151:THR:HG22	2.12	0.49
1:A:192:TYR:HB2	1:A:626:VAL:HG11	1.94	0.49
1:A:261:LEU:HB2	1:A:284:PHE:HB3	1.93	0.49
1:A:339:THR:O	1:A:342:VAL:HG23	2.12	0.49
1:A:282:LEU:HD23	1:A:322:VAL:HB	1.95	0.49
1:A:652:LYS:HD2	1:A:657:ILE:CD1	2.43	0.49
1:A:588:PHE:O	1:A:634:TYR:HA	2.14	0.48
1:A:383:THR:HG22	1:A:388:TYR:CE2	2.49	0.47
1:A:35:THR:HG22	1:A:36:GLY:N	2.28	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:188:GLU:HB2	1:A:623:PHE:CD1	2.49	0.47
1:A:24:ARG:HG3	1:A:375:ARG:O	2.15	0.47
1:A:80:VAL:HA	1:A:108:ARG:O	2.15	0.47
1:A:408:THR:O	1:A:423:ARG:HA	2.15	0.47
1:A:87:LEU:HB2	1:A:95:SER:HB3	1.97	0.47
1:A:325:ILE:HB	1:A:333:PHE:CD2	2.51	0.46
1:A:471:VAL:HG22	1:A:477:VAL:HG22	1.98	0.46
1:A:580:THR:HB	1:A:640:PRO:HG3	1.98	0.46
1:A:311:THR:HG22	1:A:312:ALA:N	2.30	0.45
1:A:143:SER:HB2	1:A:144:PRO:HD2	1.97	0.45
1:A:429:VAL:HG12	1:A:430:ALA:N	2.31	0.45
3:C:2:GLC:H62	3:C:3:G6D:O5	2.16	0.45
1:A:35:THR:CG2	1:A:36:GLY:N	2.80	0.45
1:A:195:LEU:HD22	1:A:196:PHE:CD2	2.52	0.45
1:A:213:LYS:O	1:A:216:ILE:HG22	2.17	0.45
1:A:332:ARG:HG3	1:A:368:GLY:O	2.17	0.45
1:A:107:LYS:C	1:A:108:ARG:HG2	2.42	0.45
1:A:606:ASN:HB2	1:A:646:GLN:H	1.80	0.45
1:A:78:GLN:HG2	1:A:80:VAL:HG22	1.97	0.45
1:A:149:ASP:C	1:A:151:THR:H	2.25	0.45
1:A:218:VAL:O	1:A:222:MET:HG2	2.17	0.45
1:A:302:TYR:HA	1:A:414:ILE:HD11	1.99	0.44
1:A:498:SER:HB2	1:A:499:PRO:CD	2.42	0.44
1:A:510:LYS:HE2	1:A:513:GLN:NE2	2.32	0.44
1:A:340:ARG:N	1:A:341:PRO:CD	2.80	0.44
1:A:305:ASP:O	1:A:309:GLN:HG2	2.17	0.44
1:A:340:ARG:HB3	1:A:341:PRO:HD3	2.00	0.44
1:A:108:ARG:HD3	1:A:161:ASN:ND2	2.32	0.44
1:A:140:ASN:HB3	7:A:700:HOH:O	2.17	0.44
1:A:232:VAL:HG11	1:A:257:GLY:HA3	1.99	0.44
1:A:395:ALA:N	1:A:396:PRO:HD2	2.33	0.44
1:A:559:ASN:OD1	1:A:574:ASN:HB3	2.18	0.44
1:A:604:THR:HG22	1:A:648:LYS:O	2.18	0.43
1:A:24:ARG:NH2	1:A:98:TYR:CD2	2.86	0.43
1:A:383:THR:HG22	1:A:388:TYR:CZ	2.52	0.43
1:A:287:SER:HB2	1:A:324:PHE:O	2.17	0.43
1:A:633:TRP:CD1	1:A:633:TRP:N	2.86	0.43
1:A:231:ALA:HB1	1:A:234:HIS:CD2	2.51	0.43
1:A:560:ILE:O	1:A:572:THR:HA	2.18	0.43
1:A:9:VAL:CG1	1:A:253:VAL:HG12	2.47	0.43
1:A:423:ARG:HB2	1:A:430:ALA:HB3	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:47:LYS:HE3	1:A:92:PHE:CD1	2.53	0.43
1:A:42:THR:O	1:A:43:HIS:HB2	2.19	0.43
1:A:591:ASN:HB2	1:A:680:ASN:OD1	2.18	0.43
1:A:37:ASP:O	1:A:48:LYS:HE3	2.19	0.42
1:A:233:LYS:O	1:A:233:LYS:HG2	2.18	0.42
1:A:68:GLY:O	1:A:392:LYS:HG3	2.20	0.42
1:A:411:GLN:HE21	1:A:413:TRP:N	2.18	0.42
1:A:652:LYS:CD	1:A:657:ILE:CD1	2.98	0.42
1:A:203:GLN:HE21	1:A:203:GLN:HB3	1.68	0.42
1:A:639:VAL:HB	1:A:640:PRO:HD2	2.01	0.42
1:A:120:GLN:HG3	1:A:222:MET:CE	2.49	0.42
1:A:195:LEU:O	1:A:196:PHE:HB2	2.20	0.42
1:A:38:LEU:HD23	1:A:38:LEU:HA	1.75	0.42
1:A:147:GLU:O	1:A:169:THR:HG23	2.20	0.42
1:A:610:LEU:HA	1:A:610:LEU:HD12	1.78	0.42
1:A:89:ASP:HB3	1:A:94:GLY:H	1.84	0.42
1:A:143:SER:CB	1:A:144:PRO:HD2	2.50	0.42
1:A:584:ILE:HG13	1:A:641:ALA:CB	2.49	0.42
1:A:4:THR:HB	1:A:399:LYS:HD3	2.01	0.41
1:A:92:PHE:CZ	1:A:372:PRO:HD2	2.55	0.41
1:A:630:TYR:CD2	1:A:631:PRO:HA	2.55	0.41
1:A:600:ASN:HB3	1:A:621:PRO:HB3	2.01	0.41
1:A:648:LYS:HE2	1:A:664:ASN:OD1	2.20	0.41
1:A:281:LEU:HD23	1:A:281:LEU:HA	1.94	0.41
1:A:108:ARG:HD3	1:A:161:ASN:HD22	1.86	0.41
1:A:531:PHE:CD1	1:A:531:PHE:N	2.88	0.41
1:A:75:TRP:CD1	1:A:75:TRP:C	2.98	0.41
1:A:149:ASP:CG	1:A:151:THR:HG22	2.45	0.41
1:A:421:TYR:O	1:A:431:LEU:HD12	2.21	0.41
1:A:518:ASP:OD1	1:A:545:GLU:HG3	2.20	0.41
1:A:18:TYR:HB2	1:A:71:VAL:HG21	2.03	0.41
1:A:69:MET:O	1:A:392:LYS:HA	2.21	0.41
1:A:120:GLN:HE21	1:A:222:MET:HE1	1.85	0.41
1:A:341:PRO:HD3	1:A:438:LEU:HD11	2.01	0.41
1:A:366:MET:HE2	1:A:366:MET:HB3	1.96	0.41
1:A:653:ASN:CG	1:A:654:GLY:N	2.79	0.41
1:A:576:ILE:HD12	1:A:576:ILE:HA	1.74	0.41
1:A:500:LEU:O	1:A:519:GLY:HA3	2.21	0.40
1:A:513:GLN:O	1:A:549:LYS:HA	2.22	0.40
1:A:149:ASP:OD1	1:A:151:THR:HB	2.22	0.40
1:A:12:TYR:CE2	1:A:134:ILE:HD11	2.57	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:104:ARG:HG2	1:A:104:ARG:HH11	1.86	0.40
1:A:205:ASN:HB3	1:A:208:ILE:HG12	2.03	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	681/683 (100%)	633 (93%)	44 (6%)	4 (1%)	21	28

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	22	THR
1	A	626	VAL
1	A	196	PHE
1	A	379	THR

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	583/583 (100%)	537 (92%)	46 (8%)	11	16

All (46) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	ASP
1	A	10	VAL
1	A	45	SER
1	A	48	LYS
1	A	60	LYS
1	A	67	THR
1	A	80	VAL
1	A	117	THR
1	A	143	SER
1	A	148	THR
1	A	151	THR
1	A	161	ASN
1	A	163	THR
1	A	195	LEU
1	A	203	GLN
1	A	229	LEU
1	A	242	ASN
1	A	270	ASN
1	A	301	MET
1	A	311	THR
1	A	318	ILE
1	A	325	ILE
1	A	342	VAL
1	A	375	ARG
1	A	378	MET
1	A	414	ILE
1	A	436	ARG
1	A	473	SER
1	A	509	THR
1	A	524	THR
1	A	533	SER
1	A	540	SER
1	A	542	ASP
1	A	574	ASN
1	A	576	ILE
1	A	610	LEU
1	A	619	ILE
1	A	632	THR
1	A	639	VAL
1	A	645	ILE
1	A	650	ILE
1	A	653	ASN
1	A	656	THR

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Mol	Chain	Res	Type
1	A	657	ILE
1	A	671	SER
1	A	677	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	8	ASN
1	A	43	HIS
1	A	62	ASN
1	A	120	GLN
1	A	161	ASN
1	A	203	GLN
1	A	270	ASN
1	A	382	ASN
1	A	410	GLN
1	A	411	GLN
1	A	428	ASN
1	A	497	ASN
1	A	559	ASN
1	A	574	ASN
1	A	600	ASN
1	A	655	ASN

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 ⓘ

8 monosaccharides 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	GLC	B	1	2	12,12,12	0.38	0	17,17,17	1.03	2 (11%)
2	GLC	B	2	2	11,11,12	0.45	0	15,15,17	1.12	1 (6%)
2	GLC	B	3	2	11,11,12	0.37	0	15,15,17	1.26	1 (6%)
3	GLC	C	1	3	12,12,12	0.35	0	17,17,17	0.94	1 (5%)
3	GLC	C	2	3	11,11,12	0.43	0	15,15,17	1.31	2 (13%)
3	G6D	C	3	3,5	9,9,11	0.41	0	11,12,16	1.41	2 (18%)
4	GLC	D	1	5,4	12,12,12	0.41	0	17,17,17	0.62	0
4	GLC	D	2	4	11,11,12	0.45	0	15,15,17	1.75	3 (20%)

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	GLC	B	1	2	-	2/2/22/22	0/1/1/1
2	GLC	B	2	2	-	0/2/19/22	0/1/1/1
2	GLC	B	3	2	-	2/2/19/22	0/1/1/1
3	GLC	C	1	3	-	0/2/22/22	0/1/1/1
3	GLC	C	2	3	-	0/2/19/22	0/1/1/1
3	G6D	C	3	3,5	-	-	0/1/1/1
4	GLC	D	1	5,4	-	0/2/22/22	0/1/1/1
4	GLC	D	2	4	-	1/2/19/22	0/1/1/1

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	2	GLC	C1-O5-C5	4.97	118.84	112.19
2	B	3	GLC	C1-O5-C5	3.75	117.22	112.19
2	B	2	GLC	C1-O5-C5	3.33	116.65	112.19
3	C	2	GLC	C1-O5-C5	3.28	116.58	112.19
3	C	3	G6D	O5-C5-C4	-2.68	106.79	109.39
4	D	2	GLC	C3-C4-C5	2.60	114.95	110.23
2	B	1	GLC	C4-C3-C2	2.46	115.15	110.83
4	D	2	GLC	O5-C5-C4	2.32	116.47	110.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	2	GLC	O5-C1-C2	2.18	116.00	110.79
3	C	3	G6D	C1-O5-C5	2.15	114.95	111.80
3	C	1	GLC	C3-C4-C5	2.12	114.08	110.23
2	B	1	GLC	C1-C2-C3	2.06	114.55	110.36

There are no chirality outliers.

All (5) torsion outliers are listed below:

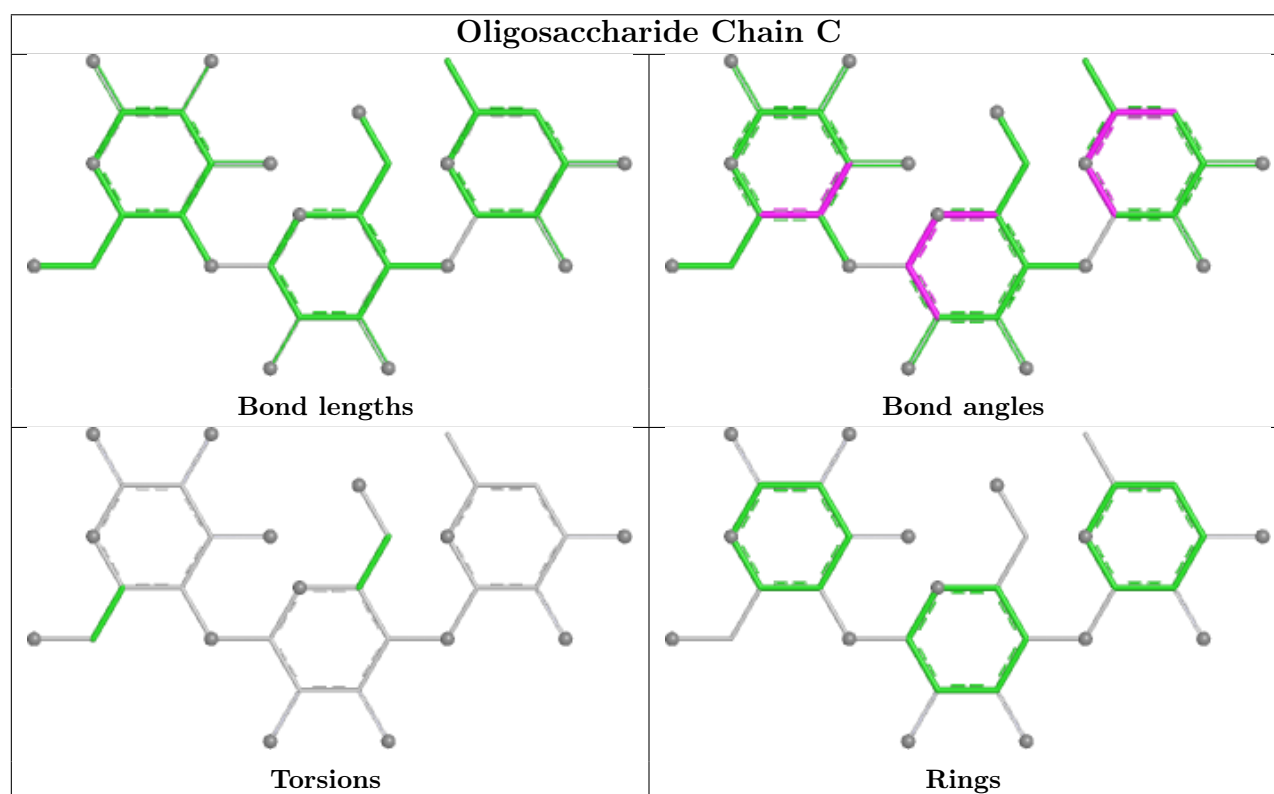
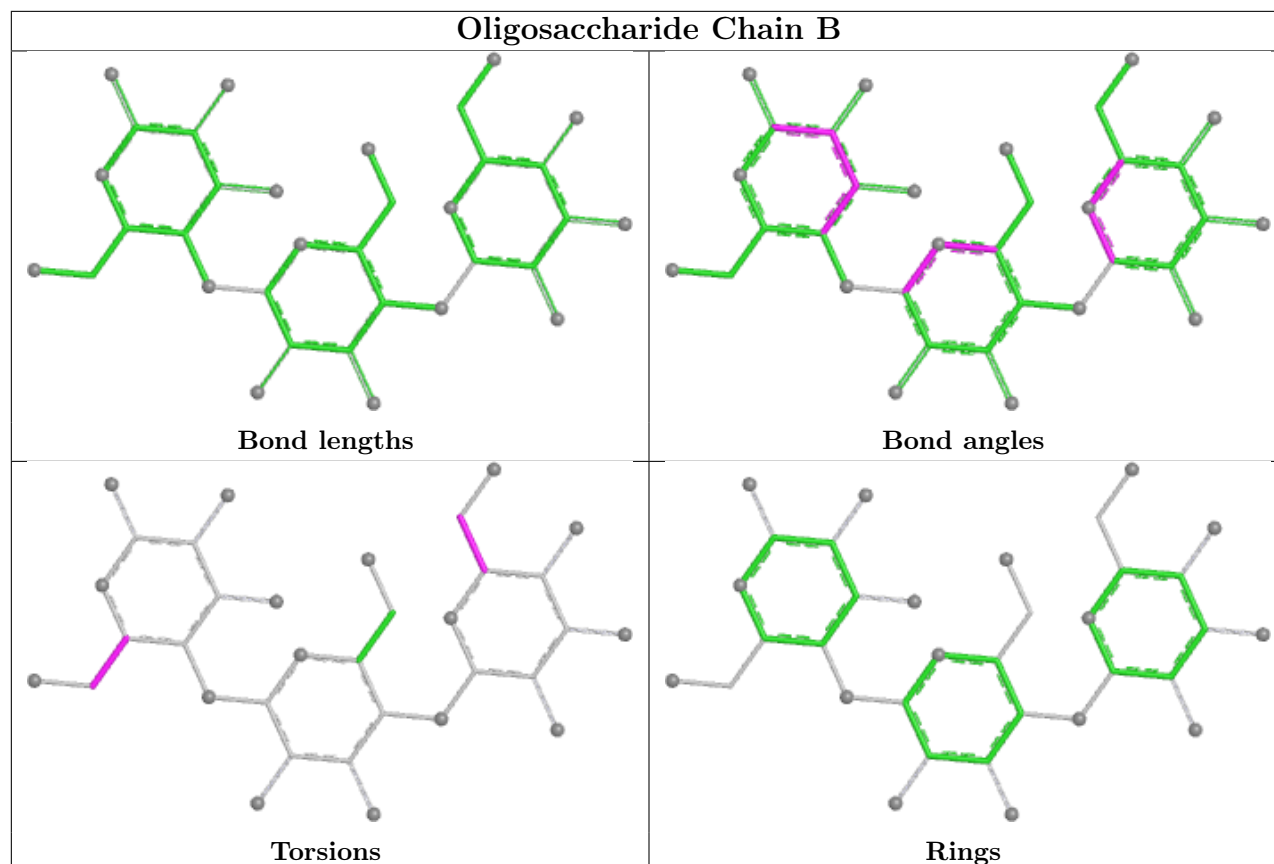
Mol	Chain	Res	Type	Atoms
2	B	1	GLC	C4-C5-C6-O6
2	B	3	GLC	O5-C5-C6-O6
2	B	1	GLC	O5-C5-C6-O6
2	B	3	GLC	C4-C5-C6-O6
4	D	2	GLC	O5-C5-C6-O6

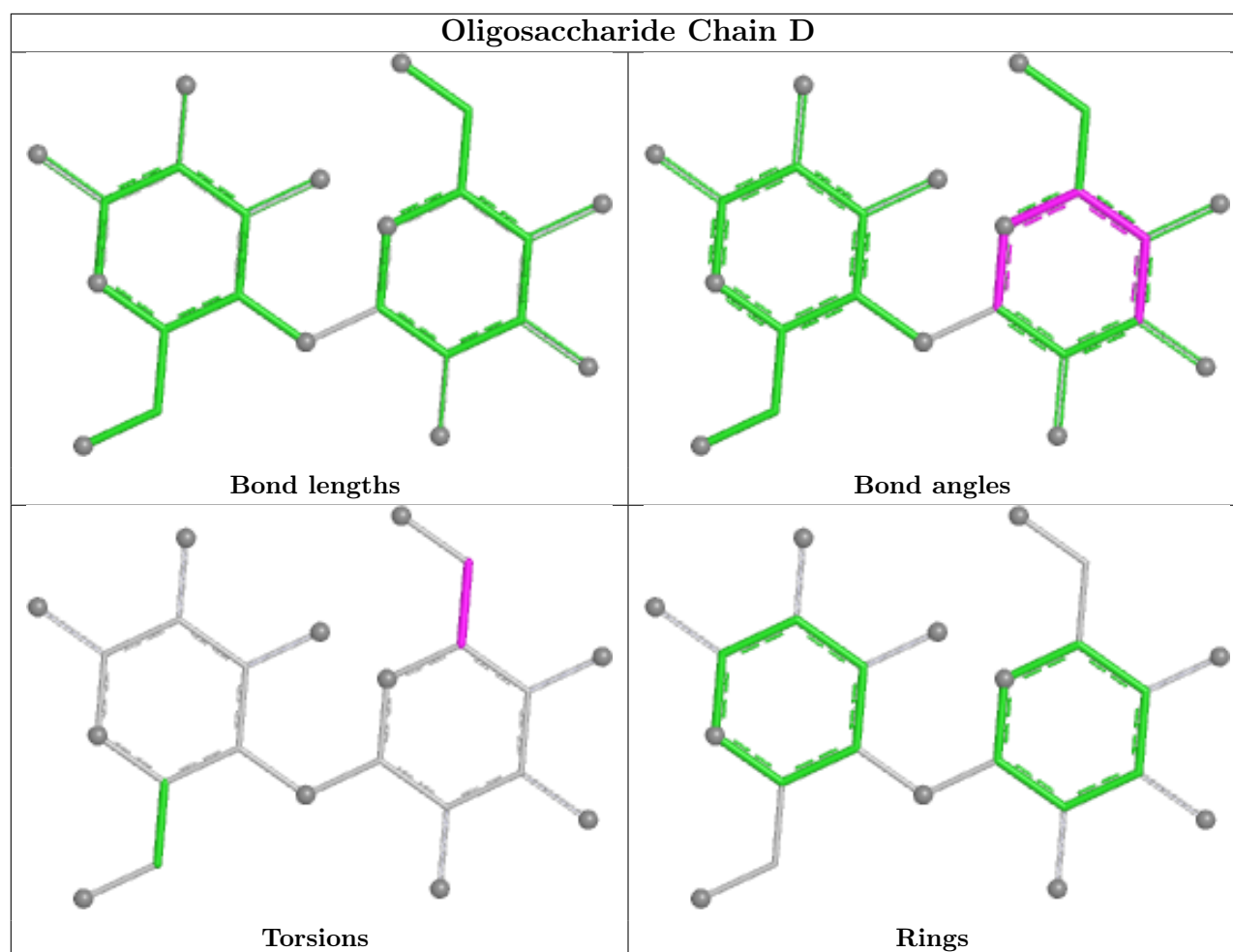
There are no ring outliers.

2 monomers are involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	2	GLC	1	0
3	C	3	G6D	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.





5.6 Ligand geometry [i](#)

Of 3 ligands modelled in this entry, 2 are monoatomic - leaving 1 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	$\# Z > 2$
5	ADH	A	694	3,4	11,11,11	8.89	4 (36%)	8,15,15	1.32	1 (12%)

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
5	ADH	A	694	3,4	-	1/2/18/18	0/1/1/1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	694	ADH	C1-N1	28.53	2.30	1.47
5	A	694	ADH	C4-C5	-5.45	1.40	1.50
5	A	694	ADH	C7-C5	3.08	1.42	1.33
5	A	694	ADH	C1-C7	-3.01	1.43	1.50

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	694	ADH	C2-C1-N1	3.46	118.44	111.40

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	694	ADH	C7-C5-C6-O6

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	694	ADH	3	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

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.