Full wwPDB X-ray Structure Validation Report (i)

Sep 13, 2020 – 09:52 AM BST

PDB ID	:	1N1I
Title	:	The structure of MSP-1(19) from Plasmodium knowlesi
Authors	:	Garman, S.C.; Simcoke, W.N.; Stowers, A.W.; Garboczi, D.N.
Deposited on	:	2002-10-17
Resolution	:	2.40 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 (i) symbol.

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

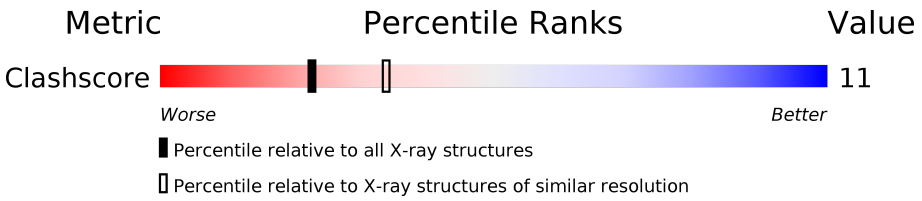
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	FAILED
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.14.4.dev1

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	141614	4398 (2.40-2.40)

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 <=5%

Note EDS failed to run properly.

Mol	Chain	Length	Quality of chain
1	A	105	72% 14% 13%
1	B	105	68% 21% 11%
1	C	105	72% 16% 11%
1	D	105	70% 12% 17%

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 3042 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 Merozoite surface protein-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	91	675	409	111	144	11	0	0	0
1	B	93	692	420	115	146	11	0	0	0
1	C	93	696	423	117	145	11	0	0	0
1	D	87	653	398	107	137	11	0	0	0

There are 52 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	GLU	GLN	SEE REMARK 999	UNP Q9GSQ9
A	2	ALA	THR	SEE REMARK 999	UNP Q9GSQ9
A	3	GLU	GLN	SEE REMARK 999	UNP Q9GSQ9
A	4	ALA	MET	SEE REMARK 999	UNP Q9GSQ9
A	5	SER	LEU	SEE REMARK 999	UNP Q9GSQ9
A	98	GLY	-	EXPRESSION TAG	UNP Q9GSQ9
A	99	PRO	-	EXPRESSION TAG	UNP Q9GSQ9
A	100	HIS	-	EXPRESSION TAG	UNP Q9GSQ9
A	101	HIS	-	EXPRESSION TAG	UNP Q9GSQ9
A	102	HIS	-	EXPRESSION TAG	UNP Q9GSQ9
A	103	HIS	-	EXPRESSION TAG	UNP Q9GSQ9
A	104	HIS	-	EXPRESSION TAG	UNP Q9GSQ9
A	105	HIS	-	EXPRESSION TAG	UNP Q9GSQ9
B	1	GLU	GLN	SEE REMARK 999	UNP Q9GSQ9
B	2	ALA	THR	SEE REMARK 999	UNP Q9GSQ9
B	3	GLU	GLN	SEE REMARK 999	UNP Q9GSQ9
B	4	ALA	MET	SEE REMARK 999	UNP Q9GSQ9
B	5	SER	LEU	SEE REMARK 999	UNP Q9GSQ9
B	98	GLY	-	EXPRESSION TAG	UNP Q9GSQ9
B	99	PRO	-	EXPRESSION TAG	UNP Q9GSQ9
B	100	HIS	-	EXPRESSION TAG	UNP Q9GSQ9

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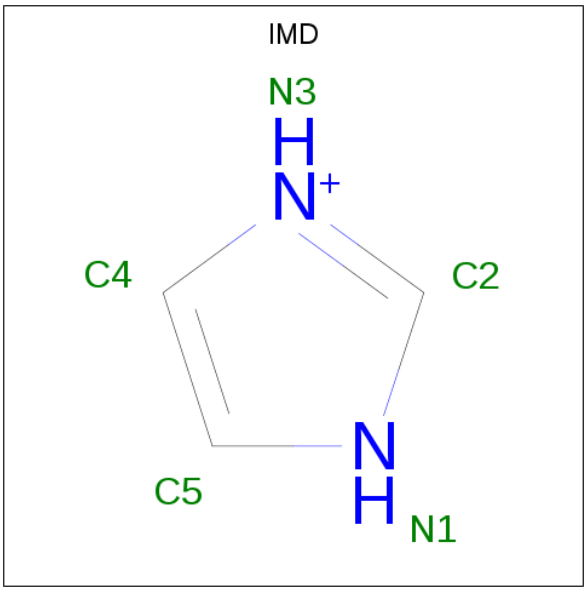
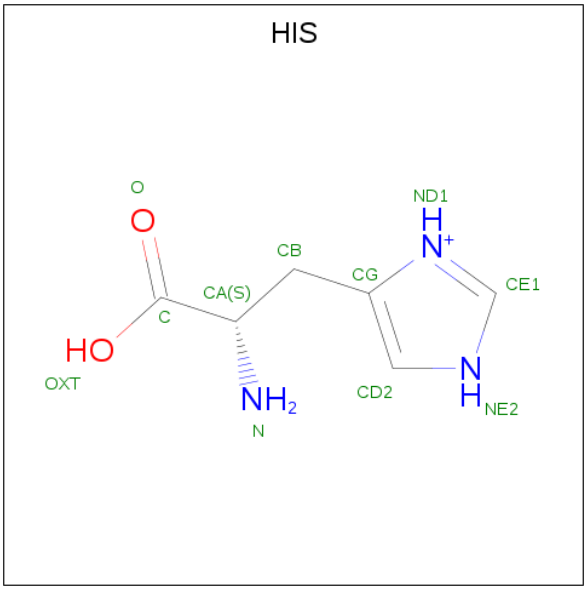
Chain	Residue	Modelled	Actual	Comment	Reference
B	101	HIS	-	EXPRESSION TAG	UNP Q9GSQ9
B	102	HIS	-	EXPRESSION TAG	UNP Q9GSQ9
B	103	HIS	-	EXPRESSION TAG	UNP Q9GSQ9
B	104	HIS	-	EXPRESSION TAG	UNP Q9GSQ9
B	105	HIS	-	EXPRESSION TAG	UNP Q9GSQ9
C	1	GLU	GLN	SEE REMARK 999	UNP Q9GSQ9
C	2	ALA	THR	SEE REMARK 999	UNP Q9GSQ9
C	3	GLU	GLN	SEE REMARK 999	UNP Q9GSQ9
C	4	ALA	MET	SEE REMARK 999	UNP Q9GSQ9
C	5	SER	LEU	SEE REMARK 999	UNP Q9GSQ9
C	98	GLY	-	EXPRESSION TAG	UNP Q9GSQ9
C	99	PRO	-	EXPRESSION TAG	UNP Q9GSQ9
C	100	HIS	-	EXPRESSION TAG	UNP Q9GSQ9
C	101	HIS	-	EXPRESSION TAG	UNP Q9GSQ9
C	102	HIS	-	EXPRESSION TAG	UNP Q9GSQ9
C	103	HIS	-	EXPRESSION TAG	UNP Q9GSQ9
C	104	HIS	-	EXPRESSION TAG	UNP Q9GSQ9
C	105	HIS	-	EXPRESSION TAG	UNP Q9GSQ9
D	1	GLU	GLN	SEE REMARK 999	UNP Q9GSQ9
D	2	ALA	THR	SEE REMARK 999	UNP Q9GSQ9
D	3	GLU	GLN	SEE REMARK 999	UNP Q9GSQ9
D	4	ALA	MET	SEE REMARK 999	UNP Q9GSQ9
D	5	SER	LEU	SEE REMARK 999	UNP Q9GSQ9
D	98	GLY	-	EXPRESSION TAG	UNP Q9GSQ9
D	99	PRO	-	EXPRESSION TAG	UNP Q9GSQ9
D	100	HIS	-	EXPRESSION TAG	UNP Q9GSQ9
D	101	HIS	-	EXPRESSION TAG	UNP Q9GSQ9
D	102	HIS	-	EXPRESSION TAG	UNP Q9GSQ9
D	103	HIS	-	EXPRESSION TAG	UNP Q9GSQ9
D	104	HIS	-	EXPRESSION TAG	UNP Q9GSQ9
D	105	HIS	-	EXPRESSION TAG	UNP Q9GSQ9

- Molecule 2 is IMIDAZOLE (three-letter code: IMD) (formula: $C_3H_5N_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N			
2	B	1	5	3	2		0	0
			Total	C	N			
2	D	1	5	3	2		0	0

- Molecule 3 is HISTIDINE (three-letter code: HIS) (formula: $C_6H_{10}N_3O_2$).



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

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	A	85	85	85	0	0
			Total	O		
4	B	80	80	80	0	0
			Total	O		
4	C	79	79	79	0	0
			Total	O		
4	D	62	62	62	0	0

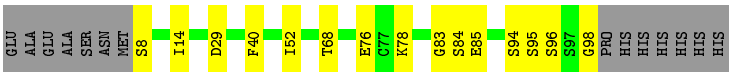
3 Residue-property plots (i)

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 failed to run properly.

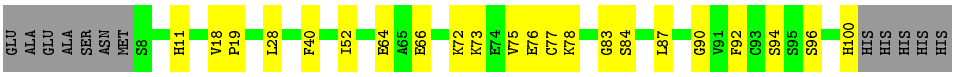
- Molecule 1: Merozoite surface protein-1

Chain A:



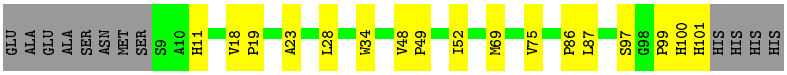
- Molecule 1: Merozoite surface protein-1

Chain B:



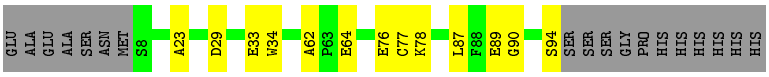
- Molecule 1: Merozoite surface protein-1

Chain C:



- Molecule 1: Merozoite surface protein-1

Chain D:



4 Data and refinement statistics (i)

EDS failed to run properly - this section is therefore incomplete.

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	33.88Å 106.46Å 62.69Å 90.00° 102.05° 90.00°	Depositor
Resolution (Å)	20.59 – 2.40	Depositor
% Data completeness (in resolution range)	97.4 (20.59-2.40)	Depositor
R_{merge}	0.08	Depositor
R_{sym}	0.08	Depositor
$< I/\sigma(I) >$ [1]	3.95 (at 2.41Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.234 , 0.264	Depositor
Wilson B-factor (Å^2)	45.5	Xtriage
Anisotropy	0.234	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3042	wwPDB-VP
Average B, all atoms (Å^2)	55.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.39% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: IMD

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	0.44	0/686	0.65	0/923
1	B	0.45	0/705	0.64	0/950
1	C	0.44	0/710	0.66	0/957
1	D	0.38	0/664	0.62	0/894
All	All	0.43	0/2765	0.64	0/3724

There are no bond length outliers.

There are no bond angle outliers.

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	675	0	616	12	0
1	B	692	0	630	14	0
1	C	696	0	632	20	0
1	D	653	0	598	12	0
2	B	5	0	5	0	0
2	D	5	0	5	0	0
3	C	10	0	6	3	0
4	A	85	0	0	6	0
4	B	80	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	C	79	0	0	7	0
4	D	62	0	0	3	0
All	All	3042	0	2492	60	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 11.

All (60) 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)
3:C:501:HIS:HA	4:C:3027:HOH:O	1.58	1.02
1:D:87:LEU:HD21	1:D:94:SER:HB2	1.55	0.87
1:C:101:HIS:HB3	4:C:3120:HOH:O	1.75	0.86
1:C:69:MET:SD	1:C:75:VAL:HG12	2.23	0.78
1:C:48:VAL:HG22	1:C:49:PRO:HD2	1.64	0.78
1:A:14:ILE:HG22	4:A:1007:HOH:O	1.86	0.76
1:B:66:GLU:HG2	4:B:2059:HOH:O	1.88	0.73
1:B:84:SER:HB2	1:B:94:SER:O	1.91	0.71
1:A:8:SER:HA	4:A:1129:HOH:O	1.92	0.70
1:A:98:GLY:HA3	1:D:33:GLU:HB2	1.73	0.69
3:C:501:HIS:HB2	4:C:3028:HOH:O	1.92	0.68
1:C:101:HIS:C	4:C:3120:HOH:O	2.34	0.67
1:A:40:PHE:O	4:A:1035:HOH:O	2.14	0.65
1:D:87:LEU:HD21	1:D:94:SER:CB	2.27	0.63
1:C:97:SER:OG	1:C:101:HIS:HB2	1.99	0.62
1:C:52:ILE:CD1	1:C:75:VAL:HG11	2.31	0.60
1:B:76:GLU:OE2	1:B:78:LYS:NZ	2.34	0.58
1:B:52:ILE:CD1	1:B:75:VAL:HG21	2.35	0.57
1:B:40:PHE:O	4:B:2035:HOH:O	2.17	0.57
1:A:95:SER:HB2	4:D:4009:HOH:O	2.04	0.57
1:B:52:ILE:HD12	1:B:75:VAL:HG21	1.87	0.56
1:D:87:LEU:N	1:D:87:LEU:HD22	2.20	0.56
1:A:84:SER:HB2	1:A:94:SER:O	2.06	0.56
1:C:48:VAL:HG22	1:C:49:PRO:CD	2.33	0.55
1:B:83:GLY:HA3	1:B:96:SER:HB2	1.90	0.54
3:C:501:HIS:HB3	4:C:3139:HOH:O	2.09	0.53
1:A:76:GLU:OE2	1:A:78:LYS:HE3	2.09	0.52
1:B:11:HIS:CE1	1:B:28:LEU:HD13	2.45	0.52
1:A:68:THR:HG23	4:A:1053:HOH:O	2.10	0.51
1:C:101:HIS:CA	4:C:3120:HOH:O	2.58	0.51
1:C:101:HIS:CB	4:C:3120:HOH:O	2.48	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:85:GLU:HB3	4:A:1108:HOH:O	2.11	0.49
1:C:52:ILE:HD12	1:C:75:VAL:HG11	1.94	0.49
1:D:62:ALA:HB1	1:D:64:GLU:OE1	2.13	0.48
1:C:52:ILE:CD1	1:C:75:VAL:CG1	2.92	0.48
1:D:89:GLU:HG2	4:D:4116:HOH:O	2.13	0.47
1:C:11:HIS:CE1	1:C:28:LEU:HD13	2.49	0.47
1:C:99:PRO:HG2	1:C:100:HIS:HD2	1.80	0.47
1:B:64:GLU:N	1:B:64:GLU:OE1	2.46	0.46
1:D:76:GLU:OE2	1:D:78:LYS:HE3	2.15	0.46
1:C:52:ILE:HD11	1:C:75:VAL:CG1	2.46	0.46
1:A:83:GLY:C	1:A:96:SER:HB3	2.36	0.46
1:A:52:ILE:HG21	4:A:1034:HOH:O	2.16	0.45
1:C:52:ILE:HD11	1:C:75:VAL:HG13	1.99	0.44
1:B:77:CYS:HB2	1:B:90:GLY:HA3	1.99	0.43
1:D:29:ASP:OD1	1:D:29:ASP:C	2.57	0.43
1:C:86:PRO:C	1:C:87:LEU:HD12	2.39	0.43
1:C:69:MET:SD	1:C:75:VAL:CG1	3.03	0.43
1:D:77:CYS:HB2	1:D:90:GLY:HA3	2.00	0.43
1:D:87:LEU:CD2	1:D:94:SER:HB2	2.39	0.42
1:B:72:LYS:O	1:B:73:LYS:HB2	2.20	0.42
1:B:18:VAL:HA	1:B:19:PRO:HD3	1.94	0.41
1:B:87:LEU:HB2	1:B:92:PHE:HB2	2.02	0.41
1:C:97:SER:HB3	1:C:101:HIS:O	2.20	0.41
1:D:23:ALA:O	1:D:34:TRP:HA	2.20	0.41
1:C:23:ALA:O	1:C:34:TRP:HA	2.20	0.41
1:D:94:SER:HB3	4:D:4104:HOH:O	2.19	0.41
1:B:100:HIS:C	1:B:100:HIS:CD2	2.94	0.40
1:A:29:ASP:C	1:A:29:ASP:OD1	2.59	0.40
1:C:18:VAL:HA	1:C:19:PRO:HD3	1.94	0.40

There are no symmetry-related clashes.

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

There are no protein backbone outliers to report in this entry.

5.3.2 Protein sidechains (i)

There are no protein residues with a non-rotameric sidechain to report in this entry.

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 monosaccharides in this entry.

5.6 Ligand geometry (i)

3 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	$\#\|Z\| > 2$
3	HIS	C	501	-	5,10,11	0.65	0	3,12,14	1.37	1 (33%)
2	IMD	B	401	-	3,5,5	0.57	0	4,5,5	0.55	0
2	IMD	D	601	-	3,5,5	0.58	0	4,5,5	0.48	0

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
3	HIS	C	501	-	-	0/5/6/8	0/1/1/1
2	IMD	B	401	-	-	-	0/1/1/1
2	IMD	D	601	-	-	-	0/1/1/1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	C	501	HIS	CD2-NE2-CE1	2.04	108.96	105.78

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	501	HIS	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 (i)

6.1 Protein, DNA and RNA chains (i)

EDS failed to run properly - this section is therefore empty.

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

EDS failed to run properly - this section is therefore empty.

6.3 Carbohydrates (i)

EDS failed to run properly - this section is therefore empty.

6.4 Ligands (i)

EDS failed to run properly - this section is therefore empty.

6.5 Other polymers (i)

EDS failed to run properly - this section is therefore empty.