



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 08:50 PM GMT

PDB ID : 1MBB
Title : OXIDOREDUCTASE
Authors : Benson, T.E.; Lees, W.J.; Walsh, C.T.; Hogle, J.M.
Deposited on : 1995-11-07
Resolution : 2.30 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
<http://wwpdb.org/validation/2016/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 (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

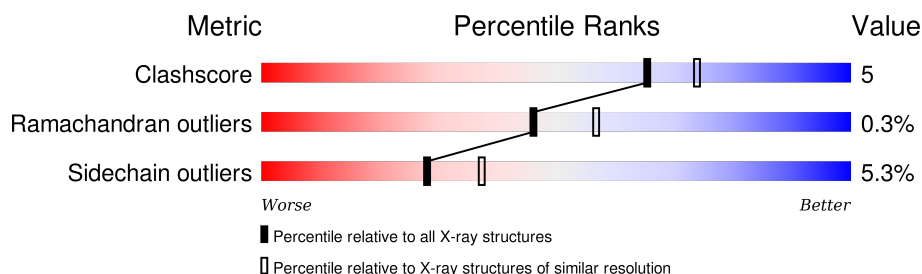
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (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. 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	342	

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
2	EEB	A	402	X	-	-	-

2 Entry composition [i](#)

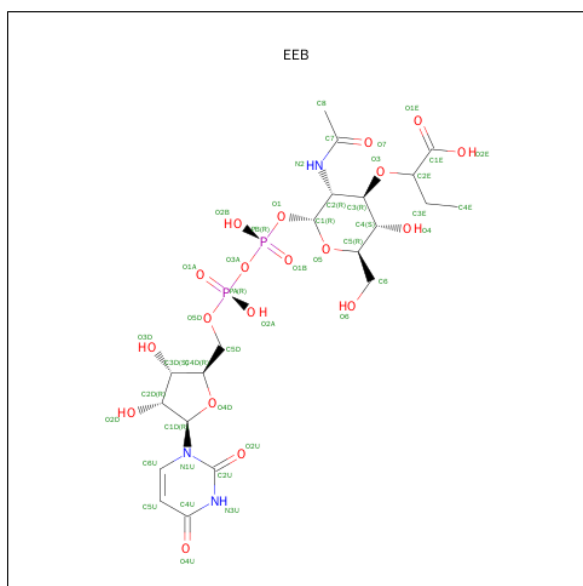
There are 4 unique types of molecules in this entry. The entry contains 2871 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 URIDINE DIPHOSPHO-N-ACETYLENOLPYRUVYLGLUCOSAMINE REDUCTASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	340	Total	C	N	O	S	0	0	0
			2654	1685	465	495	9			

- Molecule 2 is SUGAR (URIDINE-DIPHOSPHATE-3(N-ACETYLGLUCOSAMINYL)BUTYRIC ACID) (three-letter code: EEB) (formula: $C_{21}H_{33}N_3O_{19}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			45	21	3	19	2		

- Molecule 3 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

- Molecule 4 is water.

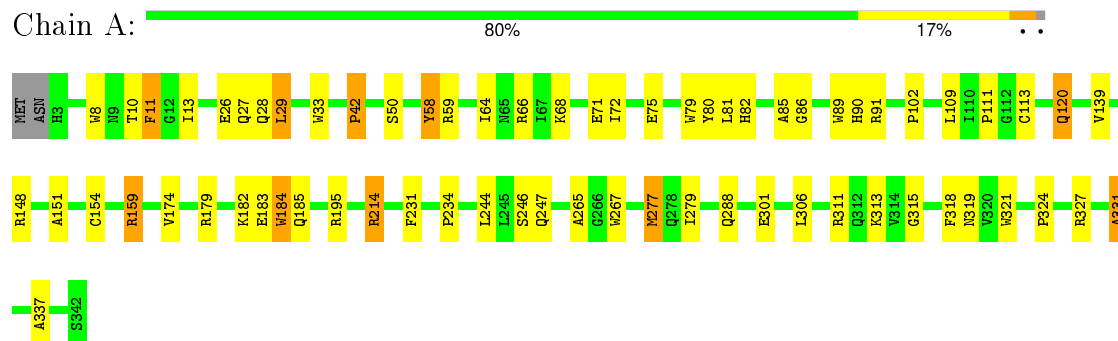
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	119	Total	O	0	0
			119	119		

3 Residue-property plots [i](#)

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 ($\text{RSRZ} > 2$). 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: URIDINE DIPHOSPHO-N-ACETYLENOLPYRUVYLGLUCOSAMINE REDUCTASE



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	50.00 Å 50.00 Å 264.30 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	12.00 – 2.30	Depositor
% Data completeness (in resolution range)	94.9 (12.00-2.30)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR	Depositor
R, R_{free}	0.238 , 0.324	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2871	wwPDB-VP
Average B, all atoms (Å ²)	14.0	wwPDB-VP

5 Model quality ⓘ

5.1 Standard geometry ⓘ

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

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	0.80	0/2716	1.55	36/3695 (1.0%)

There are no bond length outliers.

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	79	TRP	CD1-CG-CD2	9.72	114.08	106.30
1	A	8	TRP	CD1-CG-CD2	8.48	113.08	106.30
1	A	321	TRP	CD1-CG-CD2	8.30	112.94	106.30
1	A	33	TRP	CD1-CG-CD2	8.11	112.78	106.30
1	A	89	TRP	CD1-CG-CD2	8.04	112.73	106.30
1	A	79	TRP	CE2-CD2-CG	-7.88	101.00	107.30
1	A	184	TRP	CD1-CG-CD2	7.79	112.53	106.30
1	A	8	TRP	CE2-CD2-CG	-7.76	101.09	107.30
1	A	33	TRP	CE2-CD2-CG	-7.69	101.15	107.30
1	A	89	TRP	CE2-CD2-CG	-7.68	101.16	107.30
1	A	159	ARG	NE-CZ-NH1	7.43	124.02	120.30
1	A	321	TRP	CE2-CD2-CG	-7.26	101.49	107.30
1	A	184	TRP	CE2-CD2-CG	-7.09	101.63	107.30
1	A	66	ARG	NE-CZ-NH1	7.07	123.83	120.30
1	A	59	ARG	NE-CZ-NH1	6.88	123.74	120.30
1	A	267	TRP	CE2-CD2-CG	-6.86	101.81	107.30
1	A	267	TRP	CD1-CG-CD2	6.78	111.73	106.30
1	A	58	TYR	CB-CG-CD2	-6.73	116.96	121.00
1	A	91	ARG	NE-CZ-NH1	6.62	123.61	120.30
1	A	148	ARG	NE-CZ-NH1	6.61	123.61	120.30
1	A	277	MET	CA-CB-CG	-6.40	102.41	113.30
1	A	33	TRP	CB-CG-CD1	-6.31	118.80	127.00
1	A	79	TRP	CG-CD1-NE1	-6.31	103.79	110.10
1	A	29	LEU	CA-CB-CG	6.30	129.78	115.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	159	ARG	NE-CZ-NH2	-6.07	117.26	120.30
1	A	179	ARG	NE-CZ-NH1	5.97	123.28	120.30
1	A	214	ARG	NE-CZ-NH1	5.94	123.27	120.30
1	A	331	ALA	N-CA-C	5.60	126.12	111.00
1	A	33	TRP	CG-CD2-CE3	5.58	138.92	133.90
1	A	79	TRP	CG-CD2-CE3	5.56	138.91	133.90
1	A	311	ARG	NE-CZ-NH2	-5.56	117.52	120.30
1	A	195	ARG	NE-CZ-NH1	5.52	123.06	120.30
1	A	66	ARG	NE-CZ-NH2	-5.34	117.63	120.30
1	A	327	ARG	NE-CZ-NH1	5.33	122.97	120.30
1	A	184	TRP	CG-CD2-CE3	5.10	138.49	133.90
1	A	184	TRP	CG-CD1-NE1	-5.06	105.04	110.10

There are no chirality outliers.

There are no planarity outliers.

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	2654	0	2614	24	0
2	A	45	0	27	2	0
3	A	53	0	31	3	0
4	A	119	0	0	1	0
All	All	2871	0	2672	27	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 5.

All (27) 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:72:ILE:HG12	1:A:81:LEU:HD22	1.81	0.62
2:A:402:EEB:HE42	3:A:401:FAD:C5X	2.34	0.57
1:A:324:PRO:HG2	1:A:337:ALA:HB1	1.88	0.55
3:A:401:FAD:H8A	3:A:401:FAD:O5B	2.06	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:71:GLU:HB2	1:A:82:HIS:HB3	1.91	0.52
1:A:234:PRO:HD3	1:A:265:ALA:HB2	1.91	0.52
1:A:75:GLU:HG3	1:A:80:TYR:CE1	2.45	0.52
1:A:244:LEU:HD21	1:A:318:PHE:HB3	1.92	0.51
1:A:279:ILE:HD12	1:A:306:LEU:HA	1.91	0.51
1:A:120:GLN:O	1:A:159:ARG:HA	2.11	0.51
1:A:214:ARG:NH2	3:A:401:FAD:O4	2.44	0.50
1:A:85:ALA:O	1:A:113:CYS:HB3	2.12	0.50
1:A:10:THR:OG1	1:A:50:SER:HA	2.12	0.49
1:A:231:PHE:HA	1:A:324:PRO:HA	1.95	0.49
1:A:68:LYS:HE2	1:A:86:GLY:O	2.14	0.48
1:A:11:PHE:HB3	1:A:13:ILE:HG12	1.96	0.47
1:A:151:ALA:O	1:A:154:CYS:HB2	2.16	0.45
1:A:26:GLU:HG3	1:A:174:VAL:HG11	1.98	0.45
1:A:315:GLY:O	1:A:319:ASN:HA	2.18	0.44
1:A:75:GLU:O	1:A:182:LYS:NZ	2.50	0.44
1:A:214:ARG:NH1	4:A:563:HOH:O	2.50	0.43
1:A:90:HIS:HD2	1:A:109:LEU:H	1.65	0.43
1:A:313:LYS:HD3	1:A:313:LYS:HA	1.76	0.42
1:A:109:LEU:O	1:A:111:PRO:HD3	2.20	0.42
2:A:402:EEB:H3	2:A:402:EEB:HE31	1.76	0.41
1:A:102:PRO:HB2	1:A:184:TRP:CD1	2.56	0.41
1:A:42:PRO:HB2	1:A:58:TYR:OH	2.20	0.41

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	338/342 (99%)	320 (95%)	17 (5%)	1 (0%)	46 57

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	331	ALA

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	282/284 (99%)	267 (95%)	15 (5%)	28 37

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	PHE
1	A	27	GLN
1	A	28	GLN
1	A	29	LEU
1	A	42	PRO
1	A	64	ILE
1	A	120	GLN
1	A	139	VAL
1	A	183	GLU
1	A	185	GLN
1	A	246	SER
1	A	247	GLN
1	A	277	MET
1	A	288	GLN
1	A	301	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	82	HIS
1	A	90	HIS
1	A	98	GLN
1	A	185	GLN
1	A	205	GLN
1	A	309	HIS

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 ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

2 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	FAD	A	401	-	48,58,58	1.72	8 (16%)	54,89,89	2.91	14 (25%)
2	EEB	A	402	-	35,47,47	2.26	11 (31%)	46,70,70	3.84	15 (32%)

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	FAD	A	401	-	-	0/30/50/50	0/6/6/6
2	EEB	A	402	-	1/1/14/15	0/28/73/73	0/3/3/3

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	402	EEB	C3E-C2E	-4.59	1.37	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	402	EEB	O3-C2E	-4.58	1.36	1.44
2	A	402	EEB	PB-O1	-3.22	1.51	1.60
3	A	401	FAD	C5A-C4A	-3.02	1.33	1.40
3	A	401	FAD	C9A-C5X	-2.70	1.37	1.42
2	A	402	EEB	O2D-C2D	-2.46	1.37	1.43
2	A	402	EEB	C3D-C2D	-2.32	1.47	1.53
2	A	402	EEB	O7-C7	-2.26	1.18	1.23
2	A	402	EEB	O3-C3	-2.12	1.38	1.43
2	A	402	EEB	C8-C7	2.02	1.54	1.50
2	A	402	EEB	C5D-C4D	2.11	1.58	1.51
3	A	401	FAD	C4A-N3A	2.31	1.39	1.35
3	A	401	FAD	C9A-N10	2.61	1.42	1.38
2	A	402	EEB	C4-C3	3.15	1.61	1.52
3	A	401	FAD	C4X-N5	3.17	1.38	1.33
3	A	401	FAD	C5A-N7A	3.24	1.50	1.39
3	A	401	FAD	C4-C4X	5.48	1.52	1.41
3	A	401	FAD	C8A-N7A	5.72	1.45	1.34
2	A	402	EEB	C6U-N1U	8.45	1.47	1.35

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	401	FAD	C4X-C10-N10	-11.64	113.66	120.52
3	A	401	FAD	C4-C4X-C10	-7.83	114.93	119.94
2	A	402	EEB	PB-O3A-PA	-5.25	117.98	132.73
2	A	402	EEB	C8-C7-N2	-4.40	107.68	116.11
3	A	401	FAD	C4X-C4-N3	-3.80	118.39	123.59
3	A	401	FAD	C8M-C8-C7	-3.50	113.05	120.73
2	A	402	EEB	C1-C2-N2	-3.47	104.47	111.01
2	A	402	EEB	O5D-PA-O1A	-2.75	98.94	109.62
2	A	402	EEB	C5D-C4D-C3D	-2.51	105.23	115.21
3	A	401	FAD	C9A-C5X-N5	-2.48	118.69	122.36
3	A	401	FAD	C1B-N9A-C4A	-2.31	123.46	126.94
3	A	401	FAD	C2B-C1B-N9A	2.62	118.30	114.29
2	A	402	EEB	O5-C1-O1	2.81	115.07	111.36
3	A	401	FAD	C8M-C8-C9	2.82	127.95	120.28
3	A	401	FAD	C4-C4X-N5	3.01	122.38	118.72
3	A	401	FAD	C1'-N10-C9A	3.05	122.28	118.86
3	A	401	FAD	C5X-C9A-N10	3.22	120.06	117.62
3	A	401	FAD	O3P-P-O5'	3.38	111.91	102.94
2	A	402	EEB	O3D-C3D-C4D	3.59	121.82	111.05
3	A	401	FAD	C4X-N5-C5X	3.82	121.16	116.76

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	402	EEB	O1-C1-C2	3.93	115.66	108.42
2	A	402	EEB	O7-C7-N2	3.95	129.91	121.86
2	A	402	EEB	O3-C3-C4	4.69	119.27	107.17
2	A	402	EEB	C4E-C3E-C2E	7.96	125.95	113.19
2	A	402	EEB	C2-N2-C7	8.18	144.11	123.10
2	A	402	EEB	O3A-PA-O5D	9.28	127.56	102.94
3	A	401	FAD	C4-N3-C2	10.64	124.45	115.25
2	A	402	EEB	C4U-N3U-C2U	11.22	125.25	114.14
2	A	402	EEB	C3-C2-N2	12.46	132.79	111.07

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	A	402	EEB	C2E

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	401	FAD	3	0
2	A	402	EEB	2	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 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.