



Full wwPDB EM Validation Report ⓘ

Mar 9, 2025 – 12:08 AM JST

PDB ID : 9IJ3
EMDB ID : EMD-60614
Title : Cryo-EM Structure of MILI-piRNA-target (26-nt)
Authors : Li, Z.Q.; Xu, Q.K.; Wu, J.P.; Shen, E.Z.
Deposited on : 2024-06-21
Resolution : 2.60 Å(reported)

This is a Full wwPDB EM 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/EMValidationReportHelp>

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:

EMDB validation analysis : **FAILED**
Mogul : 1.8.5 (274361), CSD as541be (2020)
MolProbity : 4.02b-467
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
MapQ : **FAILED**
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.41.2

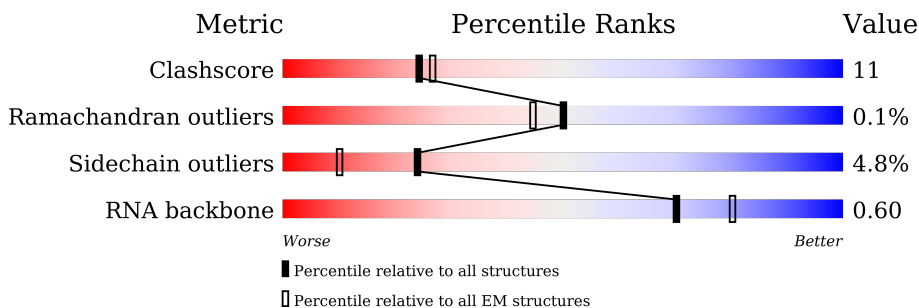
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 2.60 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)	EM structures (#Entries)
Clashscore	210492	15764
Ramachandran outliers	207382	16835
Sidechain outliers	206894	16415
RNA backbone	6643	2191

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the 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 $\leq 5\%$.

Mol	Chain	Length	Quality of chain
1	A	971	<div> <div>56%</div> <div>19%</div> <div>•</div> <div>24%</div> </div>
2	B	26	<div> <div>35%</div> <div>42%</div> <div>23%</div> </div>
3	C	25	<div> <div>64%</div> <div>32%</div> <div>•</div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7005 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, 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 Piwi-like protein 2.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	738	Total	C	N	O	S	0	0
			5921	3799	1023	1056	43		

- Molecule 2 is a RNA chain called RNA (26-MER).

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	26	Total	C	N	O	P	0	0
			550	247	95	182	26		

- Molecule 3 is a RNA chain called RNA (25-MER).

Mol	Chain	Residues	Atoms					AltConf	Trace
3	C	25	Total	C	N	O	P	0	0
			533	238	93	177	25		

- Molecule 4 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		AltConf
4	A	1	Total	Mn	0
			1	1	

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.

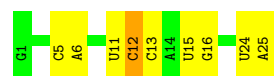
• Molecule 1: Piwi-like protein 2



• Molecule 2: RNA (26-MER)



Chain C: 64% 32% .



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	139074	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	50	Depositor
Minimum defocus (nm)	1500	Depositor
Maximum defocus (nm)	2000	Depositor
Magnification	Not provided	
Image detector	TFS FALCON 4i (4k x 4k)	Depositor

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: MN, OMC

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	0.29	0/6054	0.62	6/8200 (0.1%)
2	B	0.58	1/590 (0.2%)	1.07	4/914 (0.4%)
3	C	0.28	0/595	0.88	0/925
All	All	0.32	1/7239 (0.0%)	0.70	10/10039 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1	U	OP3-P	-10.65	1.48	1.61

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	357	ILE	CG1-CB-CG2	-7.25	95.46	111.40
1	A	619	LEU	CA-CB-CG	6.51	130.28	115.30
1	A	429	ASP	CB-CG-OD2	6.50	124.15	118.30
2	B	11	C	C5-C6-N1	6.14	124.07	121.00
1	A	451	LEU	CA-CB-CG	6.09	129.30	115.30
2	B	9	A	O5'-P-OP2	6.01	117.91	110.70
1	A	613	MET	CB-CG-SD	5.77	129.71	112.40
2	B	24	C	C2-N1-C1'	5.77	125.14	118.80
2	B	11	C	C6-N1-C2	-5.59	118.06	120.30
1	A	277	LEU	CA-CB-CG	5.17	127.20	115.30

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	5921	0	6010	127	0
2	B	550	0	282	16	0
3	C	533	0	268	10	0
4	A	1	0	0	0	0
All	All	7005	0	6560	142	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 (142) 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:252:LYS:HE3	3:C:11:U:OP1	1.56	1.04
1:A:252:LYS:CE	3:C:11:U:OP1	2.12	0.97
1:A:516:THR:HG23	1:A:520:ASN:OD1	1.71	0.89
1:A:307:THR:HG23	1:A:308:LYS:N	2.00	0.76
1:A:229:ILE:HG13	1:A:376:ALA:H	1.54	0.72
1:A:610:LYS:O	1:A:613:MET:HB2	1.90	0.72
1:A:699:VAL:HA	1:A:702:ILE:HG22	1.75	0.69
1:A:526:HIS:HD1	1:A:859:TYR:HH	1.41	0.68
1:A:252:LYS:NZ	3:C:11:U:P	2.68	0.66
1:A:600:PRO:HB3	1:A:633:ARG:HH21	1.61	0.66
1:A:307:THR:HG23	1:A:308:LYS:HG3	1.76	0.66
1:A:851:GLN:HB2	1:A:903:HIS:HB3	1.81	0.63
1:A:384:ARG:HG3	1:A:498:MET:HB2	1.81	0.63
1:A:249:VAL:O	1:A:255:ARG:NH1	2.32	0.62
1:A:523:PRO:HB3	1:A:859:TYR:HB3	1.81	0.62
1:A:338:GLY:O	1:A:422:ASN:ND2	2.34	0.61
2:B:14:G:H2'	2:B:15:G:H8	1.65	0.60
1:A:240:GLN:O	1:A:307:THR:HG22	2.01	0.60
1:A:287:VAL:HB	1:A:304:ILE:HB	1.82	0.59
1:A:354:ARG:HH22	1:A:384:ARG:HH21	1.51	0.59
1:A:516:THR:HG22	1:A:516:THR:O	2.02	0.57
1:A:307:THR:CG2	1:A:308:LYS:N	2.67	0.57
1:A:671:ILE:HD13	1:A:678:LEU:HD12	1.86	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:931:ASN:HB2	1:A:971:LEU:HD22	1.85	0.57
1:A:434:LYS:HB3	1:A:469:LEU:HD12	1.87	0.56
1:A:951:LEU:O	1:A:956:LEU:N	2.38	0.56
1:A:697:ILE:HD13	2:B:2:U:H5'	1.88	0.55
1:A:584:LEU:HD12	1:A:871:PRO:HB2	1.88	0.55
1:A:888:TYR:HB3	1:A:901:PRO:HB2	1.88	0.55
1:A:515:LEU:HG	1:A:519:ILE:HD12	1.89	0.55
1:A:952:SER:HA	1:A:956:LEU:HB2	1.89	0.55
1:A:252:LYS:NZ	3:C:11:U:OP1	2.39	0.55
1:A:931:ASN:ND2	1:A:971:LEU:O	2.40	0.55
1:A:244:THR:HB	1:A:303:LYS:HB3	1.88	0.54
1:A:712:ALA:HA	1:A:715:ILE:HD12	1.90	0.54
1:A:252:LYS:HZ1	3:C:11:U:P	2.31	0.54
1:A:695:GLN:NE2	2:B:1:U:OP2	2.40	0.54
2:B:1:U:H5'	2:B:2:U:H5''	1.90	0.54
1:A:889:LEU:HB3	1:A:904:TYR:HE2	1.73	0.53
1:A:429:ASP:OD2	1:A:430:VAL:N	2.41	0.53
1:A:886:ASP:N	1:A:886:ASP:OD1	2.42	0.53
1:A:326:ARG:NH2	1:A:339:ARG:O	2.42	0.53
1:A:672:MET:HA	1:A:699:VAL:HG22	1.90	0.52
1:A:733:PRO:O	1:A:735:LYS:NZ	2.43	0.52
1:A:646:ARG:NH2	1:A:648:GLU:OE2	2.43	0.51
1:A:717:LEU:HB3	1:A:727:LEU:HD11	1.92	0.51
1:A:240:GLN:HB2	1:A:308:LYS:HB2	1.93	0.51
1:A:328:VAL:HG13	1:A:543:ALA:HB1	1.91	0.51
1:A:224:LEU:HD21	1:A:873:THR:HG21	1.92	0.51
1:A:514:ASP:HA	1:A:517:GLN:HG2	1.92	0.51
1:A:816:SER:OG	1:A:819:GLN:HG3	2.11	0.50
1:A:668:VAL:HG22	1:A:695:GLN:HB3	1.92	0.50
1:A:810:VAL:HB	1:A:847:VAL:HA	1.93	0.50
1:A:516:THR:O	1:A:516:THR:CG2	2.59	0.50
3:C:15:U:H2'	3:C:16:G:C8	2.48	0.49
1:A:944:TYR:HE1	1:A:971:LEU:HD21	1.78	0.49
1:A:859:TYR:HB2	1:A:866:PHE:HD2	1.78	0.49
1:A:225:GLY:HA3	1:A:866:PHE:HE2	1.76	0.49
1:A:361:TYR:CZ	1:A:533:LEU:HB2	2.48	0.48
1:A:661:GLU:HG2	1:A:664:ILE:HG12	1.93	0.48
2:B:14:G:H2'	2:B:15:G:C8	2.48	0.48
1:A:255:ARG:HB3	1:A:272:PHE:HE2	1.79	0.48
1:A:353:HIS:HB2	1:A:355:LEU:HD22	1.96	0.48
1:A:680:GLY:HA2	1:A:955:ILE:HG23	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:10:A:H2'	2:B:11:C:C6	2.49	0.48
2:B:7:U:H2'	2:B:8:C:H6	1.79	0.47
1:A:228:LEU:HB2	1:A:879:ILE:HG22	1.96	0.47
1:A:359:PRO:HB3	1:A:380:HIS:CE1	2.50	0.47
1:A:904:TYR:OH	1:A:937:ARG:O	2.32	0.47
1:A:642:LEU:HD13	1:A:650:TYR:CZ	2.50	0.47
1:A:382:VAL:HG11	1:A:501:ILE:HD13	1.97	0.47
1:A:951:LEU:HA	1:A:955:ILE:HB	1.96	0.47
2:B:11:C:H2'	2:B:12:A:C8	2.50	0.46
1:A:607:PHE:HB3	1:A:642:LEU:HD21	1.98	0.46
2:B:17:A:H2'	2:B:18:A:C8	2.50	0.46
1:A:307:THR:HG23	1:A:308:LYS:H	1.79	0.46
1:A:541:GLU:O	1:A:545:ASN:HB2	2.16	0.46
1:A:251:CYS:SG	1:A:253:SER:OG	2.66	0.46
1:A:815:VAL:HG23	1:A:820:LEU:HD12	1.98	0.46
1:A:526:HIS:ND1	1:A:859:TYR:OH	2.38	0.45
1:A:648:GLU:HB2	1:A:652:ARG:HH21	1.81	0.45
1:A:415:ILE:HG22	1:A:426:ARG:HA	1.98	0.45
1:A:645:ASP:O	1:A:675:ARG:NH1	2.45	0.45
1:A:852:LYS:N	3:C:16:G:OP1	2.34	0.45
1:A:617:ARG:HA	1:A:617:ARG:HD2	1.79	0.45
1:A:685:LEU:HD12	1:A:689:GLN:HB2	1.98	0.45
1:A:229:ILE:HB	1:A:374:LEU:O	2.18	0.45
1:A:604:TRP:HZ3	1:A:606:LEU:HD13	1.82	0.45
1:A:728:TRP:O	1:A:930:TRP:NE1	2.48	0.45
1:A:272:PHE:HD1	1:A:277:LEU:HB2	1.81	0.44
1:A:361:TYR:CG	1:A:533:LEU:HD22	2.52	0.44
1:A:850:VAL:HG22	1:A:904:TYR:HD1	1.82	0.44
1:A:616:ALA:HA	1:A:619:LEU:HD13	1.99	0.44
1:A:604:TRP:CD2	1:A:637:PRO:HB3	2.52	0.44
1:A:226:LEU:HD11	1:A:229:ILE:HG23	2.00	0.44
1:A:235:ASN:HB2	1:A:551:GLY:HA3	2.00	0.44
1:A:790:LEU:HD22	1:A:831:LEU:HG	2.00	0.44
1:A:383:ILE:HG12	1:A:419:ARG:HH22	1.83	0.43
1:A:860:LEU:HD12	1:A:869:PRO:HA	1.99	0.43
2:B:16:A:H2'	2:B:17:A:C8	2.53	0.43
2:B:17:A:H2'	2:B:18:A:H8	1.83	0.43
1:A:427:ILE:HD12	1:A:470:LEU:HB3	2.01	0.43
1:A:740:ILE:HD12	1:A:808:ILE:HG23	2.00	0.43
1:A:289:LEU:HB2	1:A:302:ILE:HB	1.99	0.43
1:A:322:ASN:ND2	3:C:12:C:O3'	2.34	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:606:LEU:HD12	1:A:668:VAL:HB	2.00	0.43
1:A:226:LEU:HD22	1:A:859:TYR:CZ	2.53	0.43
1:A:228:LEU:HD23	1:A:228:LEU:HA	1.86	0.43
1:A:584:LEU:HD13	1:A:935:THR:HG21	2.00	0.43
1:A:388:VAL:O	1:A:392:MET:HG2	2.18	0.43
1:A:425:TYR:CE2	1:A:489:ILE:HG21	2.54	0.43
1:A:742:MET:HA	1:A:757:GLY:O	2.18	0.43
1:A:544:SER:O	1:A:548:THR:HG23	2.19	0.42
2:B:7:U:H2'	2:B:8:C:C6	2.55	0.42
1:A:234:HIS:CG	1:A:552:LEU:HD22	2.53	0.42
1:A:273:ASP:OD2	1:A:273:ASP:N	2.52	0.42
1:A:486:LYS:HE2	1:A:486:LYS:HB3	1.88	0.42
1:A:415:ILE:HD12	1:A:424:THR:HB	2.02	0.42
1:A:684:LYS:HB2	1:A:684:LYS:HE2	1.87	0.42
1:A:228:LEU:HG	1:A:879:ILE:HB	2.02	0.42
1:A:758:PHE:HB3	1:A:772:ARG:O	2.20	0.42
1:A:671:ILE:HG21	1:A:678:LEU:HG	2.02	0.41
1:A:852:LYS:HB2	1:A:852:LYS:HE2	1.82	0.41
1:A:951:LEU:HB3	1:A:956:LEU:HG	2.01	0.41
1:A:591:THR:HG21	1:A:713:GLN:HG2	2.02	0.41
3:C:24:U:H2'	3:C:25:A:C8	2.56	0.41
1:A:342:TYR:HA	1:A:360:GLY:HA3	2.03	0.41
1:A:224:LEU:HA	1:A:860:LEU:HA	2.02	0.41
1:A:828:ILE:HG21	1:A:908:LEU:HD22	2.03	0.41
1:A:378:VAL:H	1:A:526:HIS:CE1	2.39	0.41
1:A:470:LEU:HB2	1:A:491:LEU:HB2	2.02	0.41
1:A:734:LEU:HD23	1:A:734:LEU:HA	1.91	0.41
2:B:13:U:H2'	2:B:14:G:C8	2.56	0.41
2:B:16:A:H2'	2:B:17:A:H8	1.86	0.41
1:A:698:ASN:ND2	2:B:1:U:O2'	2.54	0.40
3:C:5:C:H2'	3:C:6:A:H8	1.86	0.40
1:A:240:GLN:HG3	1:A:310:LEU:HD22	2.03	0.40
1:A:325:PHE:CE2	1:A:329:MET:HE2	2.56	0.40
1:A:415:ILE:HG12	2:B:10:A:H5''	2.02	0.40
1:A:451:LEU:O	1:A:455:SER:HB3	2.22	0.40
1:A:588:LYS:HB3	1:A:588:LYS:HE2	1.80	0.40
1:A:816:SER:OG	1:A:819:GLN:CG	2.69	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 PDB entries followed by that with respect to all EM entries.

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	730/971 (75%)	665 (91%)	64 (9%)	1 (0%)	48 71

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	226	LEU

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 PDB entries followed by that with respect to all EM entries.

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	666/863 (77%)	634 (95%)	32 (5%)	21 44

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	261	ASP
1	A	268	ASN
1	A	320	PHE
1	A	348	MET
1	A	354	ARG
1	A	392	MET
1	A	400	LYS
1	A	403	PHE
1	A	422	ASN
1	A	437	LYS

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Mol	Chain	Res	Type
1	A	465	ASP
1	A	466	ASP
1	A	477	ARG
1	A	520	ASN
1	A	575	ASN
1	A	586	TRP
1	A	633	ARG
1	A	639	TRP
1	A	644	ASP
1	A	698	ASN
1	A	736	GLN
1	A	742	MET
1	A	746	HIS
1	A	747	ASP
1	A	763	ASN
1	A	812	ARG
1	A	838	PHE
1	A	839	ASP
1	A	859	TYR
1	A	877	HIS
1	A	918	MET
1	A	924	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	776	GLN
1	A	946	HIS

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	B	25/26 (96%)	6 (24%)	0
3	C	24/25 (96%)	2 (8%)	0
All	All	49/51 (96%)	8 (16%)	0

All (8) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	B	2	U

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Mol	Chain	Res	Type
2	B	3	A
2	B	7	U
2	B	8	C
2	B	19	C
2	B	24	C
3	C	12	C
3	C	13	C

There are no RNA pucker outliers to report.

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

1 non-standard protein/DNA/RNA residue is 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	OMC	B	26	3,2	19,22,23	0.48	0	26,31,34	0.67	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
2	OMC	B	26	3,2	-	0/9/27/28	0/2/2/2

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 1 ligands modelled in this entry, 1 is monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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.