



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 03:17 PM GMT

PDB ID : 4C2X  
Title : Human N-myristoyltransferase isoform 2 (NMT2)  
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Deposited on : 2013-08-20  
Resolution : 2.33 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) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

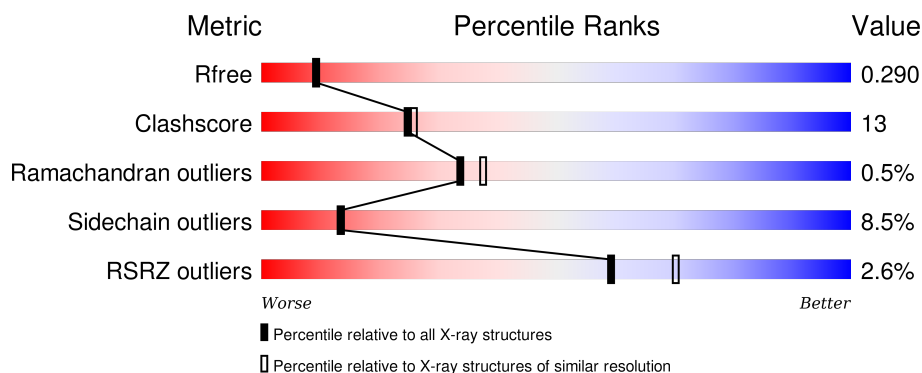
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*


The reported resolution of this entry is 2.33 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))
$R_{free}$	91344	1406 (2.36-2.32)
Clashscore	102246	1509 (2.36-2.32)
Ramachandran outliers	100387	1490 (2.36-2.32)
Sidechain outliers	100360	1491 (2.36-2.32)
RSRZ outliers	91569	1412 (2.36-2.32)

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.

Mol	Chain	Length	Quality of chain
1	A	410	

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	MG	A	999	-	-	-	X

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 3345 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 GLYCYLPEPTIDE N-TETRADECANOYLTRANSFERASE 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	388	Total	C	N	O	S	0	4	0
			3212	2081	544	572	15			

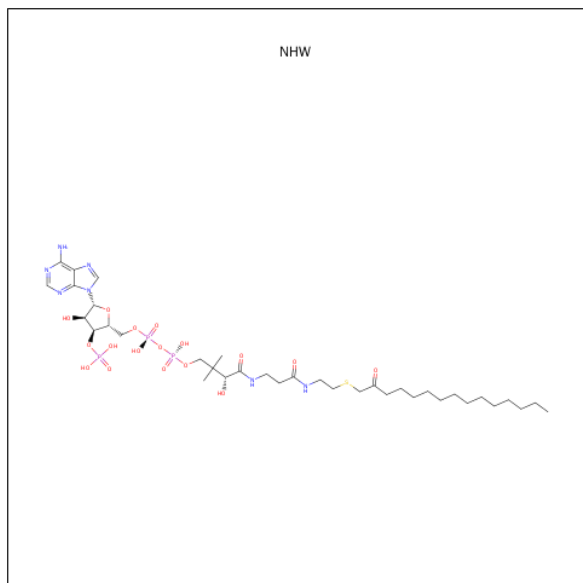
There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	87	MET	-	EXPRESSION TAG	UNP O60551
A	88	GLY	-	EXPRESSION TAG	UNP O60551
A	89	SER	-	EXPRESSION TAG	UNP O60551
A	90	SER	-	EXPRESSION TAG	UNP O60551
A	91	HIS	-	EXPRESSION TAG	UNP O60551
A	92	HIS	-	EXPRESSION TAG	UNP O60551
A	93	HIS	-	EXPRESSION TAG	UNP O60551
A	94	HIS	-	EXPRESSION TAG	UNP O60551
A	95	HIS	-	EXPRESSION TAG	UNP O60551
A	96	HIS	-	EXPRESSION TAG	UNP O60551
A	97	SER	-	EXPRESSION TAG	UNP O60551
A	98	SER	-	EXPRESSION TAG	UNP O60551
A	99	GLY	-	EXPRESSION TAG	UNP O60551
A	100	LEU	-	EXPRESSION TAG	UNP O60551
A	101	GLU	-	EXPRESSION TAG	UNP O60551
A	102	VAL	-	EXPRESSION TAG	UNP O60551
A	103	LEU	-	EXPRESSION TAG	UNP O60551
A	104	PHE	-	EXPRESSION TAG	UNP O60551
A	105	GLN	-	EXPRESSION TAG	UNP O60551
A	106	GLY	-	EXPRESSION TAG	UNP O60551
A	107	PRO	-	EXPRESSION TAG	UNP O60551
A	108	HIS	-	EXPRESSION TAG	UNP O60551
A	109	MET	-	EXPRESSION TAG	UNP O60551
A	304	LYS	ARG	VARIANT	UNP O60551

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Mg	0	0
			1	1		

- Molecule 3 is 2-OXOPENTADECYL-COA (three-letter code: NHW) (formula:  $C_{36}H_{64}N_7O_{17}P_3S$ ).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
			Total	C	N	O	P	S		
3	A	1	64	36	7	17	3	1	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	68	Total	O	0	0
			68	68		



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	49.70 Å 72.32 Å 114.69 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	61.17 – 2.33 57.34 – 2.33	Depositor EDS
% Data completeness (in resolution range)	99.8 (61.17-2.33) 99.8 (57.34-2.33)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.50 (at 2.32 Å)	Xtriage
Refinement program	REFMAC 5.8.0033	Depositor
R, $R_{free}$	0.218 , 0.288 0.221 , 0.290	Depositor DCC
$R_{free}$ test set	936 reflections (5.40%)	DCC
Wilson B-factor (Å <sup>2</sup> )	51.0	Xtriage
Anisotropy	0.275	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 41.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 18367 reflections	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	3345	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.66% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

## 5 Model quality

### 5.1 Standard geometry

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

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	0.62	0/3312	0.79	2/4493 (0.0%)

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	2

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
1	A	234	ARG	NE-CZ-NH2	-6.80	116.90	120.30
1	A	234	ARG	NE-CZ-NH1	5.94	123.27	120.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	129	ASP	Peptide
1	A	312	SER	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	3212	0	3227	83	0
2	A	1	0	0	0	0
3	A	64	0	60	6	0
4	A	68	0	0	0	0
All	All	3345	0	3287	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 13.

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:234:ARG:NH2	1:A:360:PHE:O	1.88	1.06
1:A:116:ARG:HA	1:A:116:ARG:NE	1.65	1.06
1:A:116:ARG:HE	1:A:116:ARG:HA	1.26	0.93
1:A:132:ILE:HD13	1:A:482:ARG:HD2	1.62	0.80
1:A:183:ASP:O	1:A:186:ASN:ND2	2.17	0.78
1:A:322:ARG:HG2	1:A:322:ARG:HH11	1.52	0.73
1:A:171:GLU:OE1	1:A:215:ARG:NH2	2.16	0.72
1:A:312:SER:HA	1:A:313:HIS:ND1	2.06	0.70
1:A:341:MET:CE	1:A:376:LEU:HD13	2.23	0.68
1:A:410:HIS:NE2	1:A:412:ALA:HB3	2.10	0.66
1:A:307:VAL:HA	1:A:312:SER:O	1.96	0.65
1:A:347:LYS:O	1:A:351:GLU:HG3	1.97	0.65
1:A:435:MET:HG3	1:A:461:PHE:CE2	2.32	0.65
1:A:295:ARG:HD3	1:A:473:ASN:OD1	1.97	0.64
1:A:248:LEU:HG	3:A:1000:NHW:H14	1.82	0.60
1:A:414:LYS:HD3	1:A:414:LYS:H	1.67	0.59
1:A:398:LEU:HD23	1:A:434:LEU:HD22	1.84	0.59
1:A:410:HIS:CD2	1:A:412:ALA:H	2.21	0.58
1:A:240:LYS:HE2	1:A:481:TRP:CZ3	2.39	0.58
1:A:441:LEU:O	1:A:445:LYS:HG3	2.04	0.57
1:A:341:MET:HE2	1:A:376:LEU:HD13	1.87	0.57
1:A:379:GLU:HG2	1:A:380:HIS:CE1	2.41	0.56
1:A:240:LYS:HE2	1:A:481:TRP:HZ3	1.71	0.55
1:A:157:MET:HE1	1:A:220:LYS:CD	2.37	0.55
1:A:251:HIS:CE1	1:A:253:LYS:HG2	2.41	0.55
1:A:113:ALA:O	1:A:114:LYS:HD3	2.07	0.55
1:A:295:ARG:CD	1:A:473:ASN:OD1	2.54	0.55
1:A:341:MET:HE1	1:A:376:LEU:HD13	1.87	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:395:THR:HB	1:A:428:GLU:HB2	1.89	0.53
1:A:116:ARG:HD3	1:A:121:ASP:OD2	2.08	0.53
1:A:157:MET:HE1	1:A:220:LYS:HD2	1.89	0.53
1:A:410:HIS:CD2	1:A:412:ALA:N	2.76	0.53
1:A:171:GLU:CD	1:A:215:ARG:HH22	2.08	0.52
1:A:266:GLU:HA	1:A:266:GLU:OE1	2.11	0.51
1:A:350:ARG:HE	1:A:368:GLU:HG3	1.76	0.51
1:A:322:ARG:CG	1:A:322:ARG:HH11	2.23	0.50
1:A:414:LYS:HD3	1:A:414:LYS:N	2.26	0.50
1:A:260:ALA:HB2	3:A:1000:NHW:H12A	1.93	0.50
1:A:312:SER:HA	1:A:313:HIS:CG	2.47	0.49
1:A:260:ALA:HB3	1:A:261:PRO:HD3	1.95	0.49
1:A:366:MET:HA	1:A:370:GLU:OE1	2.12	0.49
1:A:250:VAL:HG23	3:A:1000:NHW:H14B	1.94	0.49
1:A:152:LEU:HD12	1:A:157:MET:HA	1.95	0.49
1:A:367:ASP:C	1:A:367:ASP:OD1	2.52	0.48
1:A:353:ILE:HG12	1:A:357:LEU:CD1	2.44	0.48
1:A:254:LEU:HD13	1:A:259:VAL:HG21	1.96	0.48
1:A:395:THR:O	1:A:427:THR:OG1	2.21	0.47
1:A:356:TYR:O	1:A:359:GLN:HG2	2.14	0.47
1:A:316:ARG:HD2	1:A:317:ASN:HA	1.96	0.46
1:A:475[A]:GLN:HB2	1:A:477:TYR:CE2	2.51	0.46
3:A:1000:NHW:H6	3:A:1000:NHW:O9	2.16	0.45
1:A:350:ARG:HD3	1:A:368:GLU:HG3	1.98	0.45
1:A:313:HIS:CD2	1:A:313:HIS:O	2.69	0.45
1:A:209:GLN:CD	1:A:209:GLN:H	2.19	0.45
1:A:320:LEU:O	1:A:324:MET:HG2	2.17	0.45
1:A:389:SER:O	1:A:391:ASN:N	2.50	0.45
1:A:260:ALA:CB	3:A:1000:NHW:H12A	2.47	0.44
1:A:117:TYR:O	1:A:121:ASP:HB2	2.17	0.44
1:A:241:LYS:HD2	1:A:241:LYS:HA	1.62	0.44
1:A:125:VAL:HA	1:A:138:ILE:CD1	2.48	0.43
1:A:399:SER:OG	1:A:422:PHE:HB3	2.18	0.43
1:A:290:PRO:HB3	1:A:475[A]:GLN:HE21	1.83	0.43
1:A:157:MET:CE	1:A:220:LYS:HD2	2.49	0.43
1:A:316:ARG:HA	1:A:317:ASN:HA	1.44	0.43
1:A:308:GLU:O	1:A:410:HIS:CE1	2.72	0.43
1:A:215:ARG:NH1	1:A:220:LYS:O	2.49	0.42
1:A:350:ARG:HE	1:A:368:GLU:CG	2.33	0.42
1:A:110:ASP:OD1	1:A:111:GLU:N	2.52	0.42
1:A:127:LYS:HE2	1:A:127:LYS:HB3	1.87	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:207:LEU:HB3	1:A:209:GLN:NE2	2.35	0.42
1:A:159:ASP:OD2	1:A:161:LEU:HD23	2.20	0.42
1:A:350:ARG:NE	1:A:368:GLU:HG3	2.35	0.41
1:A:299:ARG:HB2	1:A:467:PHE:CE2	2.55	0.41
1:A:188:PHE:HA	1:A:406:THR:O	2.20	0.41
1:A:322:ARG:CG	1:A:322:ARG:NH1	2.83	0.41
1:A:226:ILE:HD12	1:A:248:LEU:HD13	2.03	0.41
1:A:185:ASP:N	1:A:185:ASP:OD1	2.54	0.41
1:A:261:PRO:HD3	3:A:1000:NHW:O5X	2.21	0.41
1:A:426:HIS:CD2	1:A:431:LEU:HD13	2.56	0.41
1:A:489:GLU:H	1:A:489:GLU:HG3	1.54	0.41
1:A:316:ARG:HA	1:A:316:ARG:HD2	1.86	0.41
1:A:228:ALA:HB2	1:A:245:ILE:HD13	2.02	0.41
1:A:177:ASN:O	1:A:189:ARG:HD2	2.20	0.40
1:A:312:SER:HA	1:A:313:HIS:HD1	1.84	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	390/410 (95%)	357 (92%)	31 (8%)	2 (0%)	<a href="#">34</a> <a href="#">37</a>

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	111	GLU
1	A	392	GLY

### 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	358/373 (96%)	328 (92%)	30 (8%)	14	14

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	111	GLU
1	A	114	LYS
1	A	116	ARG
1	A	128	LEU
1	A	130	GLU
1	A	157	MET
1	A	164	SER
1	A	185	ASP
1	A	198	LEU
1	A	208	LEU
1	A	209	GLN
1	A	221	LYS
1	A	241	LYS
1	A	254	LEU
1	A	256	SER
1	A	299	ARG
1	A	324	MET
1	A	336	SER
1	A	348	SER
1	A	376	LEU
1	A	393	LYS
1	A	405	SER
1	A	407	VAL
1	A	409	HIS
1	A	414	LYS
1	A	425	ILE
1	A	428	GLU
1	A	432	LEU
1	A	455	LEU
1	A	489	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	135	HIS
1	A	186	ASN
1	A	317	ASN
1	A	361	HIS
1	A	410	HIS
1	A	426	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 ⓘ

Of 2 ligands modelled in this entry, 1 is 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	$\# Z  > 2$
3	NHW	A	1000	2	56,66,66	1.02	3 (5%)	68,92,92	1.78	10 (14%)

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	NHW	A	1000	2	-	0/61/81/81	0/3/3/3

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1000	NHW	CP-S1	2.27	1.86	1.81
3	A	1000	NHW	O4X-C1X	2.42	1.44	1.41
3	A	1000	NHW	P3X-O3X	2.68	1.68	1.60

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1000	NHW	N3A-C2A-N1A	-9.82	121.37	128.89
3	A	1000	NHW	C6-C7-N8	-3.46	104.29	111.88
3	A	1000	NHW	C4X-O4X-C1X	-3.30	106.09	109.72
3	A	1000	NHW	C2X-C1X-N9A	-2.90	109.86	114.29
3	A	1000	NHW	P2A-O3A-P1A	-2.78	124.91	132.73
3	A	1000	NHW	O4X-C4X-C5X	-2.18	101.53	109.32
3	A	1000	NHW	C7-C6-C5	-2.15	108.77	112.31
3	A	1000	NHW	C2-S1-CP	-2.00	98.56	101.89
3	A	1000	NHW	O7A-P3X-O9A	2.13	117.44	110.58
3	A	1000	NHW	C13-C11-C10	3.09	114.98	109.34

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1000	NHW	6	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](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	388/410 (94%)	0.04	10 (2%) 59 70	31, 53, 85, 124	4 (1%)

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	412	ALA	7.7
1	A	113	ALA	3.8
1	A	114	LYS	3.6
1	A	144	ASN	3.2
1	A	112	ALA	2.7
1	A	316	ARG	2.6
1	A	129	ASP	2.5
1	A	413	HIS	2.4
1	A	325	LYS	2.2
1	A	185	ASP	2.1

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

There are no non-standard protein/DNA/RNA residues in this entry.

### 6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

### 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors

of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
2	MG	A	999	1/1	0.94	0.15	3.28	65,65,65,65	0
3	NHW	A	1000	64/64	0.96	0.14	0.27	30,45,52,58	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.