



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 01:00 am GMT

PDB ID : 1T2E  
Title : Plasmodium falciparum lactate dehydrogenase S245A, A327P mutant complexed with NADH and oxamate  
Authors : Cameron, A.; Read, J.; Tranter, R.; Winter, V.J.; Sessions, R.B.; Brady, R.L.; Vivas, L.; Easton, A.; Kendrick, H.; Croft, S.L.; Barros, D.; Lavandera, J.L.; Martin, J.J.; Risco, F.; Garcia-Ochoa, S.; Gamo, F.J.; Sanz, L.; Leon, L.; Ruiz, J.R.; Gabarro, R.; Mallo, A.; De Las Heras, F.G.  
Deposited on : 2004-04-21  
Resolution : 1.85 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	:	1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix)	:	1.9-1692
EDS	:	trunk28620
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
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)	:	recalc28949

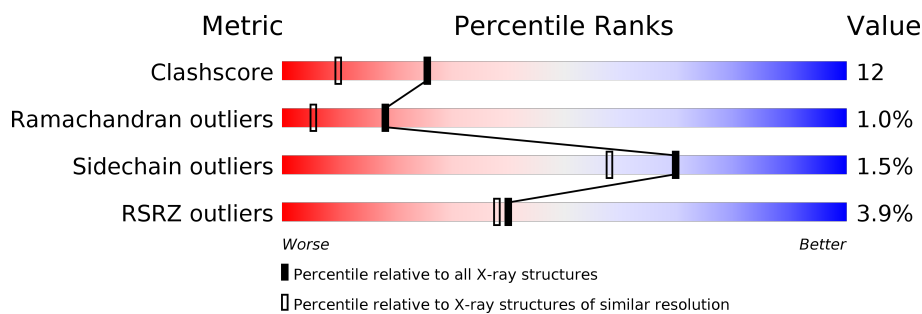
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.85 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	112137	2083 (1.86-1.86)
Ramachandran outliers	110173	2060 (1.86-1.86)
Sidechain outliers	110143	2060 (1.86-1.86)
RSRZ outliers	101464	1932 (1.86-1.86)

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	322	

## 2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 2702 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.

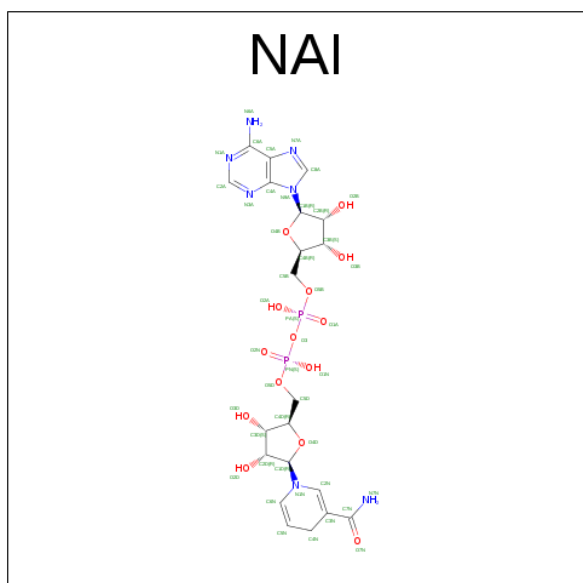
- Molecule 1 is a protein called L-lactate dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	317	2404	1532	413	446	13	39	0	0

There are 8 discrepancies between the modelled and reference sequences:

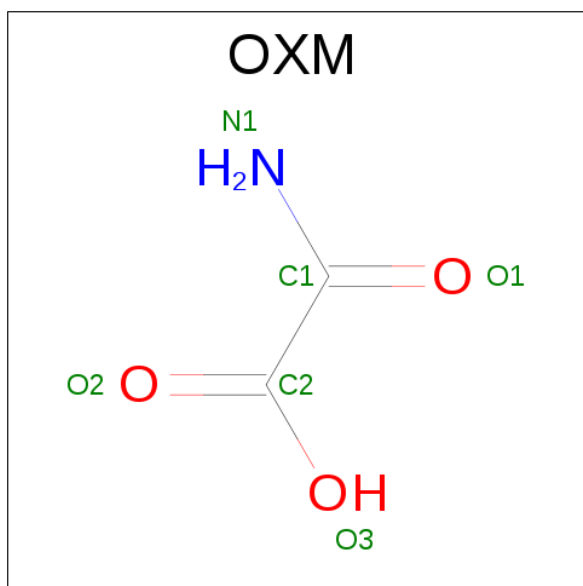
Chain	Residue	Modelled	Actual	Comment	Reference
A	245	ALA	SER	ENGINEERED	UNP Q27743
A	327	PRO	ALA	ENGINEERED	UNP Q27743
A	330	HIS	-	EXPRESSION TAG	UNP Q27743
A	331	HIS	-	EXPRESSION TAG	UNP Q27743
A	332	HIS	-	EXPRESSION TAG	UNP Q27743
A	333	HIS	-	EXPRESSION TAG	UNP Q27743
A	334	HIS	-	EXPRESSION TAG	UNP Q27743
A	335	HIS	-	EXPRESSION TAG	UNP Q27743

- Molecule 2 is 1,4-DIHYDRONICOTINAMIDE ADENINE DINUCLEOTIDE (three-letter code: NAI) (formula: C<sub>21</sub>H<sub>29</sub>N<sub>7</sub>O<sub>14</sub>P<sub>2</sub>).



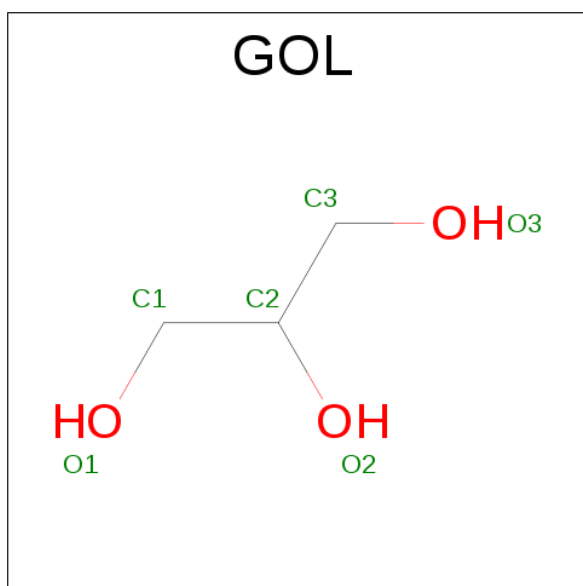
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			44	21	7	14	2		

- Molecule 3 is OXAMIC ACID (three-letter code: OXM) (formula:  $C_2H_3NO_3$ ).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			6	2	1	3		

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula:  $C_3H_8O_3$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			6	3	3		

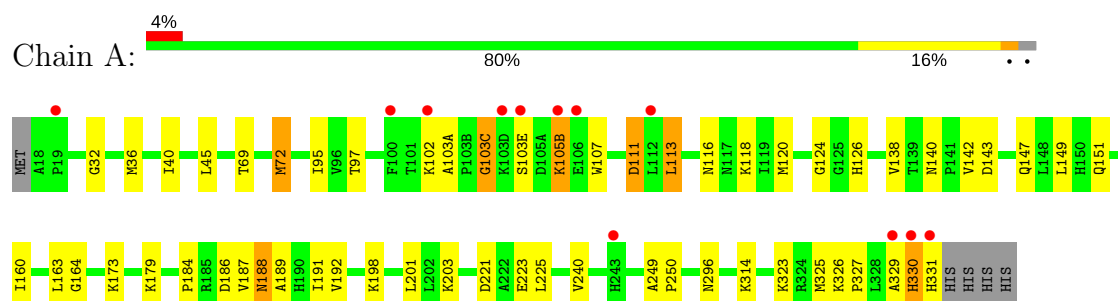
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	242	Total	O	0	0
			242	242		

### 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 the various outlier classes 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 ( $RSRZ > 2$ ). 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.

- Molecule 1: L-lactate dehydrogenase



## 4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	79.07Å 85.34Å 89.71Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 1.85 19.27 – 1.85	Depositor EDS
% Data completeness (in resolution range)	94.0 (20.00-1.85) 94.1 (19.27-1.85)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.06	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.42 (at 1.85Å)	Xtriage
Refinement program	REFMAC 5	Depositor
R, $R_{free}$	0.156 , 0.191 0.159 , (Not available)	Depositor DCC
$R_{free}$ test set	No test flags present.	DCC
Wilson B-factor (Å <sup>2</sup> )	14.0	Xtriage
Anisotropy	0.662	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.40 , 52.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.011 for -h,-l,-k	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	2702	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	17.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.07% 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, 0.333 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: GOL, NAI, OXM

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	1.08	5/2444 (0.2%)	1.07	5/3312 (0.2%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	103(A)	ALA	C-N	-18.00	1.00	1.34
1	A	113	LEU	C-N	-14.13	1.07	1.34
1	A	116	ASN	C-N	-5.74	1.20	1.34
1	A	192	VAL	CB-CG1	5.35	1.64	1.52
1	A	72	MET	SD-CE	-5.20	1.48	1.77

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	221	ASP	CB-CG-OD2	6.63	124.26	118.30
1	A	143	ASP	CB-CG-OD2	6.60	124.24	118.30
1	A	186	ASP	CB-CG-OD2	5.92	123.62	118.30
1	A	103(C)	GLY	C-N-CA	-5.26	108.55	121.70
1	A	111	ASP	CB-CG-OD2	5.08	122.87	118.30

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	2404	0	2483	55	2
2	A	44	0	20	6	0
3	A	6	0	2	0	0
4	A	6	0	8	2	0
5	A	242	0	0	8	5
All	All	2702	0	2513	58	5

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 12.

All (58) 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)
2:A:336:NAI:O5D	2:A:336:NAI:C5D	1.66	1.43
1:A:103(E):SER:OG	1:A:105(B):LYS:HD2	1.23	1.34
1:A:103(E):SER:OG	1:A:105(B):LYS:CD	2.01	1.08
1:A:296:ASN:ND2	5:A:576:HOH:O	2.02	0.93
2:A:336:NAI:PN	2:A:336:NAI:C5D	2.64	0.85
1:A:103(E):SER:CB	1:A:105(B):LYS:NZ	2.42	0.82
1:A:103(E):SER:HB3	1:A:105(B):LYS:NZ	1.96	0.81
1:A:103(E):SER:CB	1:A:105(B):LYS:CE	2.60	0.79
1:A:113:LEU:HD23	1:A:113:LEU:C	2.03	0.78
1:A:103(E):SER:HB2	1:A:105(B):LYS:HE3	1.70	0.72
1:A:147:GLN:HG2	1:A:151:GLN:HE21	1.57	0.70
1:A:103(E):SER:OG	1:A:105(B):LYS:CE	2.41	0.68
1:A:69:THR:HA	1:A:72:MET:HE1	1.74	0.68
1:A:329:ALA:C	1:A:331:HIS:H	1.98	0.66
1:A:103(E):SER:HB3	1:A:105(B):LYS:HZ1	1.60	0.66
1:A:326:LYS:HB3	1:A:327:PRO:HD3	1.79	0.65
1:A:138:VAL:O	2:A:336:NAI:H2N	1.98	0.63
1:A:102:LYS:HZ2	1:A:240:VAL:HG21	1.63	0.62
1:A:330:HIS:HD2	5:A:509:HOH:O	1.81	0.62
1:A:102:LYS:NZ	1:A:107:TRP:CD1	2.70	0.59
1:A:102:LYS:NZ	1:A:240:VAL:HG21	2.17	0.58
1:A:124:GLY:CA	1:A:149:LEU:HD11	2.36	0.56
1:A:330:HIS:HE1	5:A:518:HOH:O	1.88	0.56
1:A:103(E):SER:CB	1:A:105(B):LYS:HZ2	2.17	0.56
1:A:188:ASN:ND2	4:A:338:GOL:O2	2.32	0.55
1:A:124:GLY:HA2	1:A:149:LEU:HD11	1.87	0.55
1:A:140:ASN:HD22	1:A:142:VAL:H	1.55	0.54
1:A:173:LYS:HG2	1:A:187:VAL:O	2.07	0.54
1:A:102:LYS:HD3	1:A:107:TRP:CD1	2.42	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:102:LYS:HD3	1:A:107:TRP:HD1	1.73	0.52
2:A:336:NAI:O1N	5:A:566:HOH:O	2.19	0.51
1:A:249:ALA:N	1:A:250:PRO:CD	2.73	0.51
1:A:69:THR:HA	1:A:72:MET:CE	2.38	0.50
1:A:191:ILE:HG12	1:A:201:LEU:HG	1.93	0.50
1:A:113:LEU:CD2	1:A:113:LEU:C	2.78	0.49
1:A:179:LYS:NZ	1:A:223:GLU:OE1	2.44	0.49
1:A:329:ALA:C	1:A:331:HIS:N	2.65	0.49
1:A:103(E):SER:OG	1:A:105(B):LYS:NZ	2.45	0.49
1:A:126:HIS:HE1	5:A:579:HOH:O	1.96	0.48
1:A:124:GLY:CA	1:A:149:LEU:CD1	2.91	0.48
1:A:97:THR:HB	2:A:336:NAI:H51N	1.96	0.47
1:A:179:LYS:CE	1:A:223:GLU:OE1	2.62	0.47
1:A:40:ILE:HG23	1:A:45:LEU:HB2	1.96	0.46
1:A:203:LYS:HE3	1:A:225:LEU:HD23	1.99	0.45
1:A:102:LYS:HZ2	1:A:240:VAL:CG2	2.28	0.45
1:A:113:LEU:CD1	1:A:325:MET:HB3	2.47	0.45
1:A:314:LYS:HD3	1:A:314:LYS:HA	1.86	0.44
1:A:36:MET:SD	1:A:95:ILE:HG21	2.58	0.43
1:A:323:LYS:HE2	5:A:523:HOH:O	2.17	0.42
1:A:111:ASP:HA	5:A:563:HOH:O	2.18	0.42
1:A:160:ILE:HG23	1:A:160:ILE:O	2.20	0.42
1:A:163:LEU:HD23	2:A:336:NAI:C7N	2.49	0.41
1:A:120:MET:O	1:A:149:LEU:HD13	2.20	0.41
1:A:32:GLY:HA2	1:A:36:MET:HE2	2.02	0.41
1:A:188:ASN:HD22	1:A:189:ALA:H	1.67	0.41
1:A:147:GLN:HG2	5:A:518:HOH:O	2.20	0.40
1:A:188:ASN:HD21	4:A:338:GOL:C2	2.33	0.40
1:A:329:ALA:O	1:A:331:HIS:N	2.50	0.40

All (5) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:A:560:HOH:O	5:A:571:HOH:O[6_555]	0.88	1.32
5:A:574:HOH:O	5:A:574:HOH:O[2_665]	1.12	1.08
5:A:455:HOH:O	5:A:551:HOH:O[4_565]	1.23	0.97
1:A:118:LYS:NZ	5:A:554:HOH:O[6_555]	1.84	0.36
1:A:118:LYS:CE	5:A:554:HOH:O[6_555]	2.19	0.01

## 5.3 Torsion angles

### 5.3.1 Protein backbone

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	315/322 (98%)	304 (96%)	8 (2%)	3 (1%)	18 5

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	330	HIS
1	A	103(C)	GLY
1	A	164	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	264/269 (98%)	260 (98%)	4 (2%)	70 58

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	105(B)	LYS
1	A	184	PRO
1	A	188	ASN
1	A	198	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	132(B)	ASN
1	A	140	ASN
1	A	151	GLN
1	A	188	ASN
1	A	210(B)	GLN
1	A	234	ASN
1	A	296	ASN
1	A	330	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 ⓘ

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	$\# Z  > 2$
2	NAI	A	336	-	40,48,48	4.12	23 (57%)	41,73,73	3.57	19 (46%)
3	OXM	A	337	-	2,5,5	1.49	1 (50%)	2,6,6	1.27	0
4	GOL	A	338	-	5,5,5	0.45	0	5,5,5	0.82	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	NAI	A	336	-	-	0/25/72/72	0/5/5/5
3	OXM	A	337	-	-	0/0/4/4	0/0/0/0
4	GOL	A	338	-	-	0/4/4/4	0/0/0/0

All (24) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	336	NAI	C2B-C1B	-9.92	1.37	1.53
2	A	336	NAI	O2B-C2B	-7.29	1.26	1.43
2	A	336	NAI	O3B-C3B	-6.03	1.29	1.43
2	A	336	NAI	C3B-C4B	-5.83	1.37	1.53
2	A	336	NAI	C2B-C3B	-5.54	1.38	1.53
2	A	336	NAI	C7N-N7N	-4.03	1.21	1.33
2	A	336	NAI	C5D-C4D	-3.81	1.39	1.51
2	A	336	NAI	O2D-C2D	-3.78	1.34	1.43
2	A	336	NAI	PN-O5D	-3.40	1.44	1.59
2	A	336	NAI	PN-O1N	-3.15	1.39	1.55
2	A	336	NAI	PN-O2N	-2.61	1.41	1.50
2	A	336	NAI	PA-O2A	-2.42	1.43	1.55
3	A	337	OXM	O1-C1	2.04	1.28	1.24
2	A	336	NAI	C5A-C4A	2.84	1.46	1.40
2	A	336	NAI	O4B-C1B	3.10	1.45	1.41
2	A	336	NAI	C2D-C1D	3.34	1.63	1.53
2	A	336	NAI	C2D-C3D	3.75	1.63	1.53
2	A	336	NAI	C3D-C4D	5.09	1.66	1.53
2	A	336	NAI	O5D-C5D	5.56	1.66	1.44
2	A	336	NAI	C4N-C5N	5.90	1.61	1.49
2	A	336	NAI	C1D-N1N	5.91	1.63	1.46
2	A	336	NAI	O7N-C7N	6.47	1.40	1.24
2	A	336	NAI	C6N-N1N	7.15	1.57	1.37
2	A	336	NAI	C6N-C5N	8.32	1.48	1.33

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	336	NAI	C4B-O4B-C1B	-8.87	100.33	109.77
2	A	336	NAI	C3D-C2D-C1D	-7.91	86.23	101.43
2	A	336	NAI	O4D-C1D-N1N	-7.57	92.83	108.07
2	A	336	NAI	C1D-N1N-C2N	-7.14	108.97	121.09
2	A	336	NAI	C2D-C1D-N1N	-5.51	99.15	113.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	336	NAI	N3A-C2A-N1A	-5.46	124.10	128.86
2	A	336	NAI	C3N-C2N-N1N	-4.91	115.96	123.08
2	A	336	NAI	O4D-C1D-C2D	-3.99	97.80	106.64
2	A	336	NAI	O4D-C4D-C3D	-3.27	98.67	105.17
2	A	336	NAI	C2A-N1A-C6A	2.17	122.57	118.77
2	A	336	NAI	O4B-C4B-C3B	2.88	110.90	105.17
2	A	336	NAI	O5D-C5D-C4D	3.03	119.73	109.00
2	A	336	NAI	O2A-PA-O1A	3.41	129.92	112.28
2	A	336	NAI	O5D-PN-O2N	3.43	123.09	109.25
2	A	336	NAI	O7N-C7N-N7N	3.90	132.40	122.92
2	A	336	NAI	O3B-C3B-C4B	4.08	123.00	111.09
2	A	336	NAI	O2B-C2B-C1B	4.24	124.88	111.61
2	A	336	NAI	C1D-N1N-C6N	4.27	130.04	120.77
2	A	336	NAI	C5B-C4B-C3B	4.43	132.19	115.29

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	336	NAI	6	0
4	A	338	GOL	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 [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	311/322 (96%)	-0.31	12 (3%) 40 38	7, 13, 32, 51	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	100	PHE	3.7
1	A	331	HIS	3.6
1	A	105(B)	LYS	3.4
1	A	19	PRO	3.0
1	A	103(D)	LYS	2.9
1	A	330	HIS	2.9
1	A	243	HIS	2.7
1	A	329	ALA	2.7
1	A	106	GLU	2.4
1	A	112	LEU	2.2
1	A	102	LYS	2.2
1	A	103(E)	SER	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(Å <sup>2</sup> )	Q<0.9
4	GOL	A	338	6/6	0.97	0.10	1.35	14,19,20,20	0
3	OXM	A	337	6/6	0.95	0.11	0.61	18,20,21,23	0
2	NAI	A	336	44/44	0.97	0.07	-0.38	8,14,17,19	0

## 6.5 Other polymers

There are no such residues in this entry.