



Full wwPDB X-ray Structure Validation Report ⓘ

May 13, 2020 – 05:19 am BST

PDB ID : 3U3D
Title : Plasmodium falciparum Sir2A preferentially hydrolyzes medium and long chain fatty acyl lysine
Authors : Zhou, Y.; Hao, Q.
Deposited on : 2011-10-05
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 ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) 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 : 2.11
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

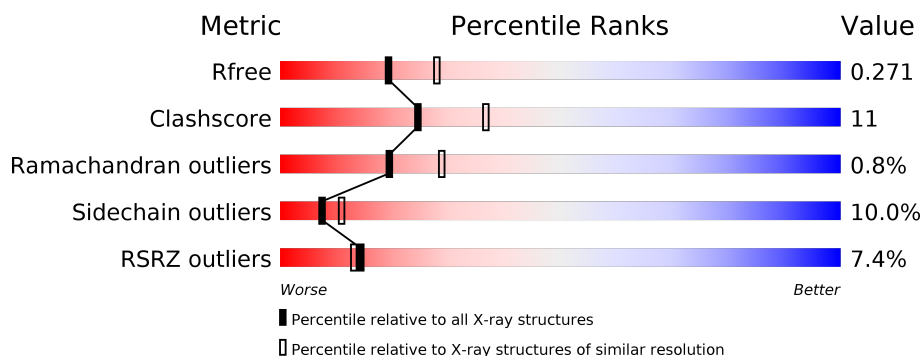
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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))
R_{free}	130704	3907 (2.40-2.40)
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)
RSRZ outliers	127900	3811 (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 $\leq 5\%$. 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	290	<div> <div>6%</div> <div>68%</div> <div>19%</div> <div>•</div> <div>9%</div> </div>
2	B	12	<div> <div>17%</div> <div>33%</div> <div>17%</div> <div>8%</div> <div>42%</div> </div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 2151 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 Transcriptional regulatory protein sir2 homologue.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	263	Total	C	N	O	S	0	1	0
			2055	1310	345	387	13			

There are 17 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-16	GLY	-	EXPRESSION TAG	UNP Q8IE47
A	-15	SER	-	EXPRESSION TAG	UNP Q8IE47
A	-14	HIS	-	EXPRESSION TAG	UNP Q8IE47
A	-13	MET	-	EXPRESSION TAG	UNP Q8IE47
A	-12	ALA	-	EXPRESSION TAG	UNP Q8IE47
A	-11	SER	-	EXPRESSION TAG	UNP Q8IE47
A	-10	MET	-	EXPRESSION TAG	UNP Q8IE47
A	-9	THR	-	EXPRESSION TAG	UNP Q8IE47
A	-8	GLY	-	EXPRESSION TAG	UNP Q8IE47
A	-7	GLY	-	EXPRESSION TAG	UNP Q8IE47
A	-6	GLN	-	EXPRESSION TAG	UNP Q8IE47
A	-5	GLN	-	EXPRESSION TAG	UNP Q8IE47
A	-4	MET	-	EXPRESSION TAG	UNP Q8IE47
A	-3	GLY	-	EXPRESSION TAG	UNP Q8IE47
A	-2	ARG	-	EXPRESSION TAG	UNP Q8IE47
A	-1	GLY	-	EXPRESSION TAG	UNP Q8IE47
A	0	SER	-	EXPRESSION TAG	UNP Q8IE47

- Molecule 2 is a protein called histone 3 myristoyl lysine 9 peptide.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	B	7	Total	C	N	O	0	0	0
			64	42	11	11			

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



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

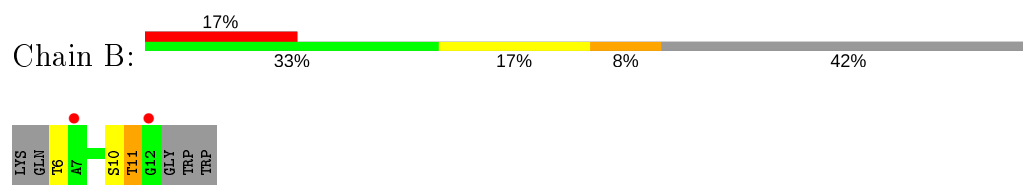
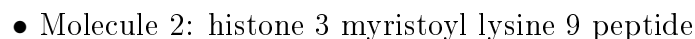
- Molecule 4 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Zn	0	0
			1	1		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	25	Total	O	0	0
			25	25		

- Molecule 1: Transcriptional regulatory protein sir2 homologue



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	32.51Å 103.33Å 105.51Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.40 36.91 – 2.40	Depositor EDS
% Data completeness (in resolution range)	98.0 (50.00-2.40) 98.0 (36.91-2.40)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.49 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.224 , 0.276 0.222 , 0.271	Depositor DCC
R_{free} test set	720 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	48.8	Xtriage
Anisotropy	0.347	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 51.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.009 for -h,l,k	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	2151	wwPDB-VP
Average B, all atoms (Å ²)	59.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.54% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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 ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, ZN, MYK

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	1.01	4/2095 (0.2%)	0.94	3/2825 (0.1%)
2	B	1.00	0/38	1.30	0/48
All	All	1.01	4/2133 (0.2%)	0.95	3/2873 (0.1%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	141	CYS	CB-SG	-9.84	1.65	1.82
1	A	73	TYR	CE1-CZ	5.71	1.46	1.38
1	A	73	TYR	CD1-CE1	5.36	1.47	1.39
1	A	181	PHE	CD2-CE2	5.21	1.49	1.39

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	168	CYS	CA-CB-SG	7.70	127.86	114.00
1	A	34	LEU	CB-CG-CD2	5.83	120.92	111.00
1	A	83	ASP	CB-CG-OD1	5.58	123.32	118.30

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	2055	0	2113	44	0
2	B	64	0	76	3	0
3	A	6	0	8	2	0
4	A	1	0	0	0	0
5	A	25	0	0	0	0
All	All	2151	0	2197	47	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 (47) 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:210:SER:N	1:A:211:THR:HA	1.70	1.01
1:A:268:LYS:O	1:A:268:LYS:HG3	1.71	0.88
1:A:49[A]:ARG:HB2	1:A:49[A]:ARG:NH2	1.91	0.85
1:A:77:ILE:HD11	1:A:178:ILE:HD11	1.60	0.83
1:A:43:SER:O	1:A:44:ASN:HB2	1.78	0.79
1:A:209:SER:C	1:A:211:THR:HA	2.05	0.78
1:A:268:LYS:O	1:A:268:LYS:CG	2.34	0.76
1:A:210:SER:N	1:A:211:THR:CA	2.53	0.66
2:B:6:THR:O	2:B:6:THR:HG23	1.97	0.65
1:A:49[A]:ARG:NH2	1:A:49[A]:ARG:CB	2.62	0.63
1:A:49[A]:ARG:CZ	1:A:49[A]:ARG:HB2	2.27	0.62
1:A:267:GLU:OE1	1:A:267:GLU:HA	2.01	0.61
1:A:48:PHE:HD1	1:A:51:SER:HB3	1.66	0.60
1:A:252:PHE:O	1:A:255:LEU:HB2	2.01	0.59
1:A:49[A]:ARG:HH21	1:A:49[A]:ARG:CB	2.15	0.59
1:A:77:ILE:CD1	1:A:178:ILE:HD11	2.30	0.59
1:A:48:PHE:CD1	1:A:51:SER:HB3	2.38	0.58
1:A:116:VAL:HG22	3:A:274:GOL:H11	1.87	0.56
2:B:6:THR:CG2	2:B:6:THR:O	2.54	0.55
1:A:49[A]:ARG:CA	1:A:49[A]:ARG:HH21	2.20	0.55
1:A:267:GLU:HA	1:A:269:ASN:HB2	1.91	0.53
1:A:49[A]:ARG:HD3	1:A:56:TRP:CD2	2.44	0.52
1:A:233:ASN:O	1:A:249:CYS:HA	2.10	0.51
2:B:10:SER:OG	2:B:11:THR:N	2.41	0.51
1:A:46:PRO:O	1:A:49[A]:ARG:NH2	2.44	0.50
1:A:49[B]:ARG:HA	1:A:56:TRP:CD1	2.46	0.50
1:A:236:LYS:HG2	1:A:240:THR:HG21	1.92	0.50
1:A:92:ILE:HD11	1:A:123:SER:N	2.28	0.49
1:A:209:SER:HB3	1:A:210:SER:HA	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:49[A]:ARG:HA	1:A:56:TRP:CD1	2.48	0.48
1:A:49[A]:ARG:HA	1:A:49[A]:ARG:HH21	1.76	0.48
1:A:92:ILE:HD11	1:A:122:ALA:C	2.33	0.48
1:A:46:PRO:HG2	1:A:54:SER:CB	2.44	0.48
1:A:197:ILE:HD12	1:A:218:LEU:HD22	1.96	0.47
1:A:48:PHE:O	1:A:49[A]:ARG:HB3	2.14	0.47
1:A:114:GLN:HG2	1:A:132:HIS:CE1	2.49	0.47
1:A:116:VAL:HG22	3:A:274:GOL:C1	2.44	0.46
1:A:17:THR:HG23	1:A:20:GLU:H	1.80	0.46
1:A:103:GLU:OE1	1:A:126:THR:HB	2.16	0.44
1:A:63:ILE:O	1:A:69:GLY:HA3	2.17	0.44
1:A:202:LEU:HD21	1:A:204:LEU:HD21	1.99	0.43
1:A:209:SER:CB	1:A:210:SER:HA	2.49	0.43
1:A:48:PHE:CD1	1:A:51:SER:CB	3.03	0.41
1:A:60:ASP:HA	1:A:61:PRO:HD2	1.82	0.41
1:A:105:LEU:HD23	1:A:105:LEU:HA	1.85	0.41
1:A:263:LYS:O	1:A:266:SER:HB3	2.21	0.41
1:A:48:PHE:O	1:A:49[B]:ARG:HB3	2.21	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	262/290 (90%)	243 (93%)	17 (6%)	2 (1%)	19	29
2	B	4/12 (33%)	4 (100%)	0	0	100	100
All	All	266/302 (88%)	247 (93%)	17 (6%)	2 (1%)	19	29

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	44	ASN
1	A	268	LYS

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	237/256 (93%)	213 (90%)	24 (10%)	7	11
2	B	4/8 (50%)	3 (75%)	1 (25%)	0	0
All	All	241/264 (91%)	216 (90%)	25 (10%)	7	10

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	14	GLN
1	A	17	THR
1	A	26	LYS
1	A	29	LYS
1	A	34	LEU
1	A	49[A]	ARG
1	A	49[B]	ARG
1	A	104	SER
1	A	148	LYS
1	A	163	GLN
1	A	170	CYS
1	A	173	ILE
1	A	189	LEU
1	A	192	GLU
1	A	200	CYS
1	A	211	THR
1	A	212	VAL
1	A	228	LYS
1	A	233	ASN
1	A	235	SER
1	A	236	LYS
1	A	255	LEU

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Mol	Chain	Res	Type
1	A	270	LYS
1	A	272	ILE
2	B	11	THR

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

Mol	Chain	Res	Type
1	A	30	HIS
1	A	233	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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	MYK	B	9	2	22,23,24	0.72	1 (4%)	19,24,26	1.09	1 (5%)

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	MYK	B	9	2	-	9/22/23/25	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	9	MYK	CB-CA	-2.04	1.50	1.53

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	9	MYK	CD-CG-CB	-2.16	106.00	113.62

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	9	MYK	CI-CT-CY-CX
2	B	9	MYK	CU-CQ-CS-CW
2	B	9	MYK	CT-CI-CK-CL
2	B	9	MYK	CQ-CS-CW-CV
2	B	9	MYK	CL-CM-CP-CR
2	B	9	MYK	CK-CI-CT-CY
2	B	9	MYK	CS-CQ-CU-CR
2	B	9	MYK	OX-CX-CY-CT
2	B	9	MYK	NZ-CX-CY-CT

There are no ring outliers.

No monomer is involved in short contacts.

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	GOL	A	274	-	5,5,5	0.44	0	5,5,5	0.47	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	GOL	A	274	-	-	0/4/4/4	-

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.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	274	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

6.1 Protein, DNA and RNA chains

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, 95th 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 ²)	Q<0.9
1	A	263/290 (90%)	0.28	18 (6%) 17 15	24, 57, 105, 179	6 (2%)
2	B	6/12 (50%)	1.21	2 (33%) 0 0	50, 60, 70, 74	0
All	All	269/302 (89%)	0.30	20 (7%) 14 13	24, 57, 105, 179	6 (2%)

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	49[A]	ARG	4.9
1	A	266	SER	4.6
1	A	12	ASP	4.6
1	A	210	SER	4.5
1	A	269	ASN	4.3
1	A	50	GLY	4.1
1	A	53	ASN	4.0
1	A	51	SER	3.8
1	A	48	PHE	3.7
2	B	12	GLY	3.7
1	A	11	LYS	3.4
1	A	271	LYS	3.4
1	A	52	SER	3.3
1	A	263	LYS	3.1
1	A	267	GLU	3.1
1	A	264	GLY	2.9
2	B	7	ALA	2.9
1	A	243	MET	2.4
1	A	92	ILE	2.1
1	A	224	LYS	2.1

6.2 Non-standard residues in protein, DNA, RNA chains [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. The B-factors column lists the minimum, median, 95th 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	B-factors(\AA^2)	Q<0.9
2	MYK	B	9	24/25	0.83	0.23	36,47,62,80	0

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. The B-factors column lists the minimum, median, 95th 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	B-factors(\AA^2)	Q<0.9
3	GOL	A	274	6/6	0.71	0.32	73,75,79,80	0
4	ZN	A	275	1/1	0.98	0.07	39,39,39,39	0

6.5 Other polymers [i](#)

There are no such residues in this entry.