



# Full wwPDB X-ray Structure Validation Report ⓘ

Jun 4, 2019 – 09:46 PM EDT

PDB ID : 5T6C  
Title : Crystal structure of *Aspergillus fumigatus* N-myristoyl transferase in complex with myristoyl CoA and a dichloro-methylpyridinyl-methoxy-phenyl-pyridine piperazine ligand  
Authors : Robinson, D.A.; Wyatt, P.G.  
Deposited on : 2016-09-01  
Resolution : 1.90 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

<https://www.wwpdb.org/validation/2017/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 : 1.8.0 (224370), CSD as540be (2019)  
Xtriage (Phenix) : 1.13  
EDS : rb-20031633  
Percentile statistics : 20171227.v01 (using entries in the PDB archive December 27th 2017)  
Refmac : 5.8.0158  
CCP4 : 7.0 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : rb-20031633

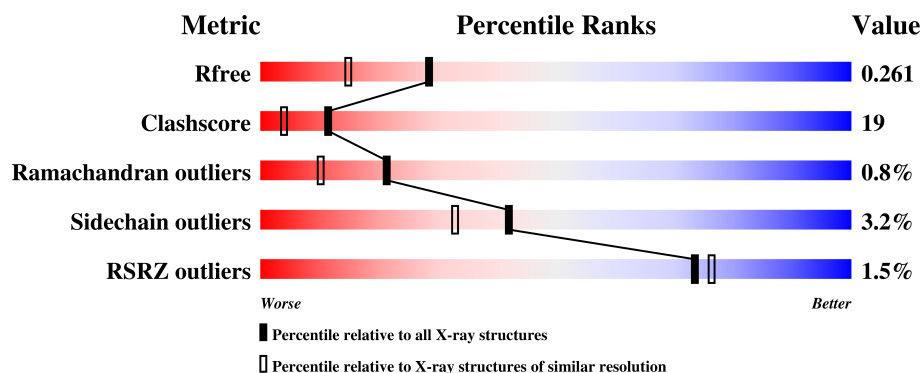
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.90 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))
$R_{free}$	111664	5502 (1.90-1.90)
Clashscore	122126	6115 (1.90-1.90)
Ramachandran outliers	120053	6048 (1.90-1.90)
Sidechain outliers	120020	6048 (1.90-1.90)
RSRZ outliers	108989	5379 (1.90-1.90)

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	411	<div> <div></div> <div>72%</div> <div>20%</div> <div>• 5%</div> </div>

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 3779 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.

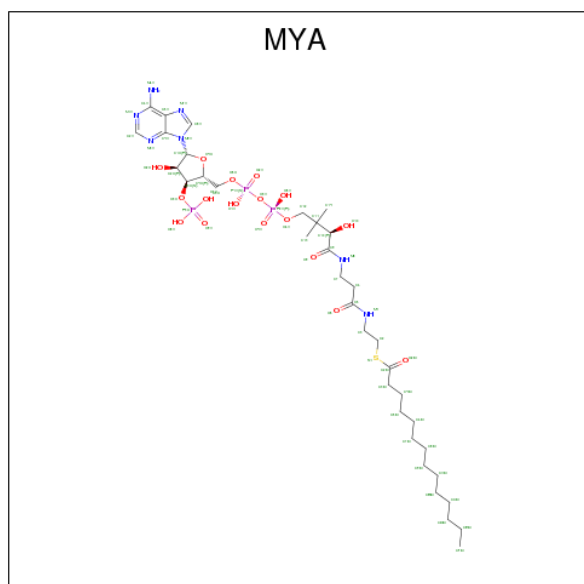
- Molecule 1 is a protein called Glycylpeptide N-tetradecanoyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	392	3185	2059	542	572	12	0	0	0

There are 4 discrepancies between the modelled and reference sequences:

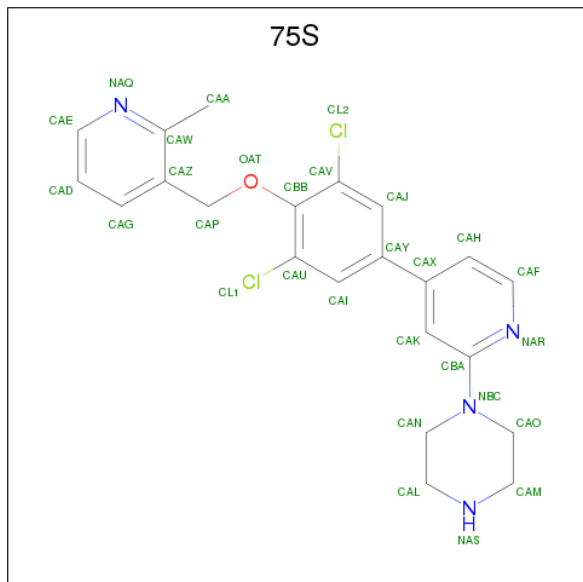
Chain	Residue	Modelled	Actual	Comment	Reference
A	82	GLY	-	expression tag	UNP Q9UVX3
A	83	PRO	-	expression tag	UNP Q9UVX3
A	84	ARG	-	expression tag	UNP Q9UVX3
A	85	SER	-	expression tag	UNP Q9UVX3

- Molecule 2 is TETRADECANOYL-COA (three-letter code: MYA) (formula:  $C_{35}H_{62}N_7O_{17}P_3S$ ).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
			Total	C	N	O	P	S		
2	A	1	63	35	7	17	3	1	0	0

- Molecule 3 is 1-(4-{3,5-dichloro-4-[(2-methylpyridin-3-yl)methoxy]phenyl}pyridin-2-yl)piperazine (three-letter code: 75S) (formula: C<sub>22</sub>H<sub>22</sub>Cl<sub>2</sub>N<sub>4</sub>O).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	Cl	N	O	0	0
			29	22	2	4	1		

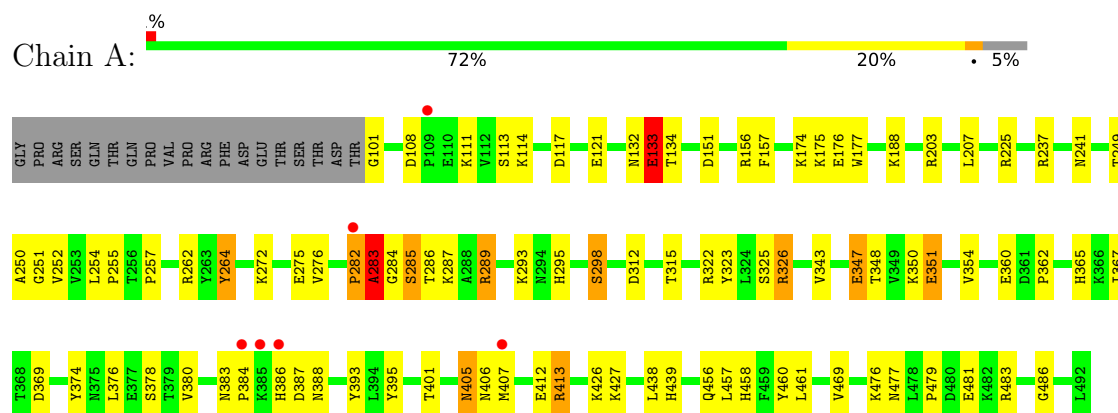
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	502	Total	O	0	0
			502	502		

### 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: Glycylpeptide N-tetradecanoyltransferase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	51.14Å 58.60Å 152.15Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	38.00 – 1.90 38.53 – 1.90	Depositor EDS
% Data completeness (in resolution range)	97.9 (38.00-1.90) 97.9 (38.53-1.90)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.27 (at 1.89Å)	Xtriage
Refinement program	REFMAC 5.8.0049	Depositor
R, $R_{free}$	0.223 , 0.256 0.230 , 0.261	Depositor DCC
$R_{free}$ test set	1802 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	12.3	Xtriage
Anisotropy	0.070	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 37.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.91	EDS
Total number of atoms	3779	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 5.27% 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: 75S, MYA

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	1.10	5/3272 (0.2%)	1.07	19/4448 (0.4%)

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	1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	133	GLU	CD-OE2	6.95	1.33	1.25
1	A	133	GLU	CD-OE1	5.61	1.31	1.25
1	A	264	TYR	CG-CD2	5.52	1.46	1.39
1	A	412	GLU	CD-OE1	5.33	1.31	1.25
1	A	481	GLU	CD-OE1	5.13	1.31	1.25

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	413	ARG	NE-CZ-NH1	12.47	126.53	120.30
1	A	483	ARG	NE-CZ-NH1	8.96	124.78	120.30
1	A	413	ARG	NE-CZ-NH2	-8.28	116.16	120.30
1	A	203	ARG	NE-CZ-NH2	8.00	124.30	120.30
1	A	289	ARG	NE-CZ-NH2	-7.35	116.63	120.30
1	A	483	ARG	NE-CZ-NH2	-6.76	116.92	120.30
1	A	326	ARG	NE-CZ-NH2	6.56	123.58	120.30
1	A	207	LEU	CB-CG-CD1	-6.43	100.06	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	283	ALA	N-CA-C	-6.17	94.33	111.00
1	A	156	ARG	NE-CZ-NH1	6.13	123.36	120.30
1	A	156	ARG	NE-CZ-NH2	-5.98	117.31	120.30
1	A	413	ARG	CB-CG-CD	5.83	126.76	111.60
1	A	262	ARG	NE-CZ-NH2	-5.57	117.52	120.30
1	A	262	ARG	NE-CZ-NH1	5.50	123.05	120.30
1	A	151	ASP	CB-CG-OD1	5.41	123.17	118.30
1	A	117	ASP	CB-CG-OD1	5.31	123.08	118.30
1	A	322	ARG	NE-CZ-NH2	-5.28	117.66	120.30
1	A	322	ARG	NE-CZ-NH1	5.26	122.93	120.30
1	A	108	ASP	CB-CG-OD1	5.18	122.96	118.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	282	PRO	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	3185	0	3166	121	0
2	A	63	0	58	2	0
3	A	29	0	0	3	0
4	A	502	0	0	96	0
All	All	3779	0	3224	123	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 19.

All (123) 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)
3:A:502:75S:OAT	3:A:502:75S:CAP	1.65	1.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:376:LEU:HG	4:A:1008:HOH:O	1.22	1.27
1:A:286:THR:HB	4:A:957:HOH:O	1.20	1.26
1:A:401:THR:HB	4:A:604:HOH:O	1.22	1.26
1:A:188:LYS:HE3	4:A:832:HOH:O	1.28	1.25
1:A:323:TYR:CE1	4:A:608:HOH:O	1.94	1.18
1:A:257:PRO:HB3	4:A:909:HOH:O	1.40	1.17
1:A:283:ALA:HB2	4:A:1056:HOH:O	1.41	1.17
1:A:354:VAL:HG23	4:A:965:HOH:O	1.02	1.16
1:A:407:MET:HG3	4:A:1028:HOH:O	1.42	1.15
1:A:250:ALA:HA	4:A:700:HOH:O	1.46	1.14
1:A:388:ASN:HB2	4:A:621:HOH:O	1.47	1.11
1:A:323:TYR:HE1	4:A:608:HOH:O	1.32	1.01
1:A:250:ALA:N	4:A:601:HOH:O	1.92	1.01
1:A:387:ASP:HB3	4:A:992:HOH:O	1.61	1.00
1:A:351:GLU:HB2	4:A:725:HOH:O	1.62	0.98
1:A:461:LEU:HD22	4:A:845:HOH:O	1.66	0.96
1:A:395:TYR:CE2	4:A:752:HOH:O	2.19	0.95
1:A:476:LYS:HE3	4:A:695:HOH:O	1.67	0.94
1:A:348:THR:HB	4:A:1011:HOH:O	1.67	0.94
1:A:121:GLU:HG2	4:A:805:HOH:O	1.65	0.93
1:A:350:LYS:HD3	4:A:804:HOH:O	1.70	0.91
1:A:286:THR:HG21	4:A:1035:HOH:O	1.71	0.90
1:A:401:THR:CB	4:A:604:HOH:O	1.88	0.89
1:A:157:PHE:CD2	1:A:376:LEU:HD21	2.07	0.88
3:A:502:75S:CBB	3:A:502:75S:CAP	2.51	0.88
1:A:293:LYS:HE2	4:A:787:HOH:O	1.72	0.87
1:A:250:ALA:HB3	4:A:601:HOH:O	1.76	0.85
1:A:250:ALA:CB	4:A:602:HOH:O	2.25	0.84
1:A:254:LEU:HA	4:A:734:HOH:O	1.77	0.83
1:A:157:PHE:HD2	1:A:376:LEU:HD21	1.42	0.82
1:A:325:SER:HB2	4:A:723:HOH:O	1.80	0.81
1:A:362:PRO:HB2	4:A:628:HOH:O	1.82	0.80
1:A:406:ASN:CB	4:A:840:HOH:O	2.31	0.78
1:A:376:LEU:CG	4:A:1008:HOH:O	1.96	0.76
1:A:362:PRO:CD	4:A:628:HOH:O	2.33	0.76
1:A:250:ALA:CB	4:A:601:HOH:O	2.34	0.76
1:A:401:THR:HG23	4:A:856:HOH:O	1.87	0.74
1:A:395:TYR:CD2	4:A:752:HOH:O	2.37	0.74
1:A:252:VAL:N	4:A:602:HOH:O	2.08	0.74
1:A:362:PRO:CB	4:A:628:HOH:O	2.34	0.73
1:A:387:ASP:CB	4:A:992:HOH:O	2.27	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:378:SER:HB2	4:A:967:HOH:O	1.91	0.71
1:A:350:LYS:CE	4:A:804:HOH:O	2.37	0.71
1:A:458:HIS:HE1	4:A:662:HOH:O	1.72	0.71
1:A:286:THR:HG23	1:A:289:ARG:H	1.56	0.70
1:A:407:MET:CB	4:A:1028:HOH:O	2.41	0.69
1:A:250:ALA:HB1	4:A:602:HOH:O	1.91	0.68
1:A:298:SER:O	4:A:603:HOH:O	2.11	0.67
1:A:362:PRO:HD2	4:A:628:HOH:O	1.93	0.67
1:A:257:PRO:CG	4:A:909:HOH:O	2.42	0.67
1:A:133:GLU:HG3	4:A:946:HOH:O	1.95	0.67
1:A:257:PRO:CB	4:A:909:HOH:O	2.15	0.65
1:A:456:GLN:HE22	1:A:477:ASN:HD21	1.42	0.65
1:A:460:TYR:HE1	4:A:601:HOH:O	1.80	0.64
1:A:175:LYS:HE2	4:A:1053:HOH:O	1.97	0.64
1:A:315:THR:HB	4:A:639:HOH:O	1.98	0.63
1:A:132:ASN:OD1	1:A:134:THR:OG1	2.17	0.61
1:A:252:VAL:O	4:A:602:HOH:O	2.16	0.60
1:A:282:PRO:O	1:A:285:SER:HB2	2.01	0.60
1:A:295:HIS:CD2	4:A:618:HOH:O	2.56	0.58
1:A:376:LEU:CD2	4:A:1008:HOH:O	2.36	0.58
1:A:250:ALA:HB3	4:A:602:HOH:O	1.91	0.58
1:A:250:ALA:CA	4:A:700:HOH:O	2.22	0.58
1:A:250:ALA:CA	4:A:601:HOH:O	2.44	0.57
1:A:350:LYS:CD	4:A:804:HOH:O	2.35	0.57
1:A:282:PRO:HA	1:A:283:ALA:HB2	1.86	0.57
1:A:476:LYS:CE	4:A:695:HOH:O	2.38	0.56
1:A:312:ASP:HA	4:A:639:HOH:O	2.05	0.55
1:A:133:GLU:HB2	4:A:646:HOH:O	2.06	0.55
1:A:350:LYS:HE3	4:A:804:HOH:O	2.02	0.55
1:A:295:HIS:ND1	4:A:610:HOH:O	2.33	0.55
1:A:362:PRO:CG	4:A:628:HOH:O	2.56	0.52
1:A:426:LYS:NZ	4:A:618:HOH:O	2.41	0.52
1:A:405:ASN:HA	4:A:656:HOH:O	2.08	0.52
1:A:469:VAL:HG13	1:A:486:GLY:CA	2.39	0.52
1:A:283:ALA:CA	4:A:912:HOH:O	2.57	0.52
1:A:380:VAL:HG11	1:A:383:ASN:HB3	1.91	0.51
1:A:286:THR:CB	4:A:957:HOH:O	2.06	0.51
1:A:174:LYS:HE2	4:A:996:HOH:O	2.10	0.50
1:A:407:MET:CG	4:A:1028:HOH:O	2.18	0.50
1:A:134:THR:HG21	4:A:990:HOH:O	2.11	0.50
1:A:348:THR:HG21	4:A:1075:HOH:O	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:380:VAL:CG1	1:A:383:ASN:HB3	2.42	0.50
1:A:225:ARG:HH22	2:A:501:MYA:P3X	2.35	0.49
1:A:177:TRP:HE1	1:A:241:ASN:HD22	1.60	0.49
1:A:174:LYS:CE	4:A:996:HOH:O	2.60	0.49
1:A:388:ASN:CB	4:A:621:HOH:O	2.27	0.49
1:A:132:ASN:ND2	4:A:624:HOH:O	2.46	0.49
1:A:113:SER:HB2	4:A:1010:HOH:O	2.12	0.48
1:A:388:ASN:CA	4:A:621:HOH:O	2.60	0.47
1:A:272:LYS:HD3	4:A:769:HOH:O	2.14	0.47
1:A:347:GLU:HB2	4:A:916:HOH:O	2.15	0.47
1:A:407:MET:HA	4:A:1028:HOH:O	2.14	0.47
1:A:284:GLY:C	4:A:607:HOH:O	2.53	0.47
1:A:326:ARG:NE	4:A:606:HOH:O	2.30	0.46
1:A:114:LYS:HE2	4:A:779:HOH:O	2.15	0.46
1:A:360:GLU:OE2	1:A:365:HIS:HD2	2.00	0.45
1:A:176:GLU:O	1:A:237:ARG:HD3	2.17	0.45
1:A:251:GLY:HA3	3:A:502:75S:CAF	2.47	0.45
1:A:275:GLU:OE2	1:A:386:HIS:NE2	2.50	0.45
1:A:439:HIS:HD2	4:A:905:HOH:O	2.00	0.45
1:A:101:GLY:HA3	4:A:722:HOH:O	2.17	0.44
1:A:257:PRO:HG3	4:A:909:HOH:O	2.12	0.44
1:A:283:ALA:HA	4:A:912:HOH:O	2.16	0.44
1:A:374:TYR:CE1	1:A:393:TYR:HB2	2.53	0.44
1:A:293:LYS:NZ	4:A:630:HOH:O	2.50	0.43
1:A:362:PRO:N	4:A:628:HOH:O	2.49	0.43
1:A:264:TYR:OH	1:A:479:PRO:HD3	2.19	0.43
1:A:254:LEU:HB3	1:A:255:PRO:HD2	2.00	0.43
1:A:323:TYR:CZ	4:A:608:HOH:O	2.38	0.43
1:A:293:LYS:HG3	4:A:1091:HOH:O	2.18	0.43
1:A:249:THR:HB	1:A:457:LEU:HG	1.99	0.43
1:A:133:GLU:CB	4:A:646:HOH:O	2.65	0.42
1:A:351:GLU:CB	4:A:725:HOH:O	2.39	0.42
1:A:458:HIS:HB3	4:A:909:HOH:O	2.19	0.42
1:A:315:THR:CB	4:A:639:HOH:O	2.63	0.41
1:A:225:ARG:NH2	2:A:501:MYA:O7A	2.51	0.41
1:A:406:ASN:C	4:A:840:HOH:O	2.59	0.41
1:A:286:THR:OG1	1:A:287:LYS:N	2.54	0.41
1:A:111:LYS:HB3	1:A:111:LYS:HE3	1.86	0.40
1:A:276:VAL:HG13	1:A:386:HIS:CD2	2.56	0.40
1:A:376:LEU:HD23	1:A:376:LEU:C	2.41	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	390/411 (95%)	374 (96%)	13 (3%)	3 (1%)	21	10

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	283	ALA
1	A	405	ASN
1	A	384	PRO

### 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 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	345/367 (94%)	334 (97%)	11 (3%)	42	33

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	133	GLU
1	A	285	SER
1	A	298	SER
1	A	343	VAL
1	A	347	GLU
1	A	351	GLU
1	A	367	ILE
1	A	369	ASP

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Mol	Chain	Res	Type
1	A	413	ARG
1	A	427	LYS
1	A	438	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	137	GLN
1	A	161	GLN
1	A	241	ASN
1	A	365	HIS
1	A	419	ASN
1	A	429	HIS
1	A	439	HIS
1	A	447	GLN
1	A	458	HIS
1	A	477	ASN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

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	$\# Z  > 2$
2	MYA	A	501	-	53,65,65	1.23	7 (13%)	66,91,91	1.49	10 (15%)
3	75S	A	502	-	32,32,32	1.95	8 (25%)	43,44,44	3.01	12 (27%)

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	MYA	A	501	-	-	0/59/80/80	0/3/3/3
3	75S	A	502	-	-	0/13/21/21	0/4/4/4

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	502	75S	CAN-NBC	-2.57	1.42	1.46
2	A	501	MYA	C2-S1	-2.49	1.78	1.81
2	A	501	MYA	O2X-C2X	-2.41	1.37	1.43
3	A	502	75S	CAY-CAX	-2.31	1.43	1.49
2	A	501	MYA	C2X-C3X	-2.29	1.47	1.52
2	A	501	MYA	C5A-N7A	-2.19	1.31	1.39
3	A	502	75S	CBB-CAV	-2.06	1.36	1.40
3	A	502	75S	CAN-CAL	2.04	1.56	1.51
2	A	501	MYA	P3X-O8A	2.04	1.63	1.54
3	A	502	75S	CAJ-CAV	2.65	1.43	1.38
3	A	502	75S	CAI-CAU	2.70	1.43	1.38
2	A	501	MYA	O10-C10	2.85	1.47	1.42
3	A	502	75S	CAO-NBC	3.51	1.52	1.46
2	A	501	MYA	P3X-O9A	4.13	1.64	1.50
3	A	502	75S	OAT-CAP	6.98	1.65	1.43

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	MYA	N3A-C2A-N1A	-5.53	119.77	128.68
3	A	502	75S	CAG-CAZ-CAW	-3.57	114.94	117.30
3	A	502	75S	CAD-CAE-NAQ	-3.33	117.93	123.42
3	A	502	75S	CAM-CAO-NBC	-3.01	103.91	110.50
3	A	502	75S	CAO-NBC-CBA	-2.96	113.51	120.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	MYA	C2X-C3X-C4X	-2.80	98.25	103.22
2	A	501	MYA	O3X-P3X-O9A	-2.40	100.12	109.39
3	A	502	75S	CAO-CAM-NAS	-2.37	104.93	111.14
3	A	502	75S	CAK-CBA-NAR	-2.21	118.85	122.69
3	A	502	75S	CAN-NBC-CBA	2.02	125.11	120.40
2	A	501	MYA	O1A-P1A-O2A	2.04	122.43	112.21
2	A	501	MYA	C2A-N1A-C6A	2.13	122.46	118.77
3	A	502	75S	CBB-CAV-CL2	2.16	121.31	118.42
2	A	501	MYA	O4X-C4X-C3X	2.18	109.52	104.84
2	A	501	MYA	C2-S1-C2M	2.24	104.42	100.16
2	A	501	MYA	C3X-C2X-C1X	2.25	104.87	99.89
3	A	502	75S	CAG-CAD-CAE	2.26	122.28	118.92
3	A	502	75S	CAK-CAX-CAY	2.30	124.73	120.87
2	A	501	MYA	O7A-P3X-O8A	2.63	117.80	107.57
3	A	502	75S	OAT-CAP-CAZ	3.80	120.61	109.36
2	A	501	MYA	O2M-C2M-C3M	5.09	118.12	109.07
3	A	502	75S	CAX-CAK-CBA	16.56	124.25	118.30

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	501	MYA	2	0
3	A	502	75S	3	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	392/411 (95%)	0.12	6 (1%) 73 76	9, 14, 31, 58	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	385	LYS	4.5
1	A	384	PRO	4.1
1	A	282	PRO	2.6
1	A	109	PRO	2.3
1	A	386	HIS	2.2
1	A	407	MET	2.2

### 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. 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	B-factors(Å <sup>2</sup> )	Q<0.9
3	75S	A	502	29/29	0.93	0.12	12,15,19,21	0
2	MYA	A	501	63/63	0.95	0.10	8,11,29,37	0



## 6.5 Other polymers [i](#)

There are no such residues in this entry.