



Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 08:42 AM GMT

PDB ID : 3OSG
Title : The structure of protozoan parasite *Trichomonas vaginalis* Myb2 in complex with MRE-1-12 DNA
Authors : Jiang, I.; Tsai, C.K.; Chen, S.C.; Wang, S.H.; Amiraslanov, I.; Chang, C.F.; Wu, W.J.; Tai, J.H.; Liaw, Y.C.; Huang, T.H.
Deposited on : 2010-09-09
Resolution : 2.00 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

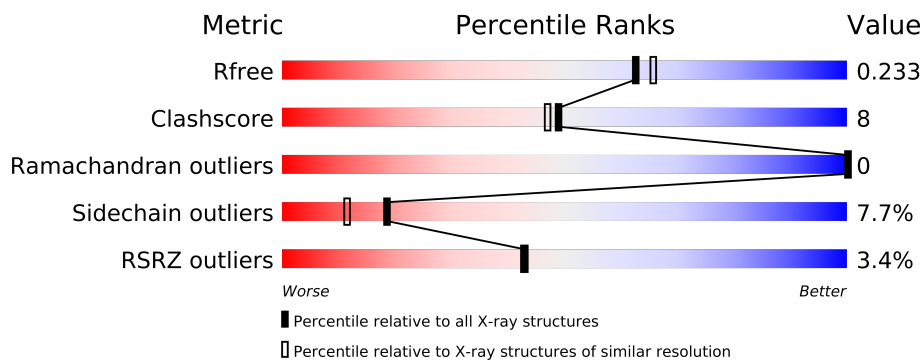
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.00 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}	66092	4888 (2.00-2.00)
Clashscore	79885	6188 (2.00-2.00)
Ramachandran outliers	78287	6102 (2.00-2.00)
Sidechain outliers	78261	6100 (2.00-2.00)
RSRZ outliers	66119	4890 (2.00-2.00)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	126	
1	D	126	
2	B	12	
2	E	12	
3	C	12	
3	F	12	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 3034 atoms, of which 0 are hydrogen and 0 are deuterium.

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 MYB21.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	110	Total	C	N	O	S	0	0	0
			906	577	169	157	3			
1	D	103	Total	C	N	O	S	0	0	0
			856	546	159	148	3			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	39	MET	-	EXPRESSION TAG	UNP Q58HP3
A	157	LEU	-	EXPRESSION TAG	UNP Q58HP3
A	158	GLU	-	EXPRESSION TAG	UNP Q58HP3
A	159	HIS	-	EXPRESSION TAG	UNP Q58HP3
A	160	HIS	-	EXPRESSION TAG	UNP Q58HP3
A	161	HIS	-	EXPRESSION TAG	UNP Q58HP3
A	162	HIS	-	EXPRESSION TAG	UNP Q58HP3
A	163	HIS	-	EXPRESSION TAG	UNP Q58HP3
A	164	HIS	-	EXPRESSION TAG	UNP Q58HP3
D	39	MET	-	EXPRESSION TAG	UNP Q58HP3
D	157	LEU	-	EXPRESSION TAG	UNP Q58HP3
D	158	GLU	-	EXPRESSION TAG	UNP Q58HP3
D	159	HIS	-	EXPRESSION TAG	UNP Q58HP3
D	160	HIS	-	EXPRESSION TAG	UNP Q58HP3
D	161	HIS	-	EXPRESSION TAG	UNP Q58HP3
D	162	HIS	-	EXPRESSION TAG	UNP Q58HP3
D	163	HIS	-	EXPRESSION TAG	UNP Q58HP3
D	164	HIS	-	EXPRESSION TAG	UNP Q58HP3

- Molecule 2 is a DNA chain called 5'-D(*AP*AP*AP*TP*AP*TP*CP*GP*TP*TP*AP*T)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	12	Total	C	N	O	P	0	0	0
			243	119	43	70	11			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	12	Total 243	C 119	N 43	O 70	P 11	0	0	0

- Molecule 3 is a DNA chain called 5'-D(*AP*TP*AP*AP*CP*GP*AP*TP*AP*TP*TP*T)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	12	Total 243	C 119	N 43	O 70	P 11	0	0	0
3	F	12	Total 243	C 119	N 43	O 70	P 11	0	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	83	Total 83	O 83	0	0
4	B	44	Total 44	O 44	0	0
4	C	42	Total 42	O 42	0	0
4	D	82	Total 82	O 82	0	0
4	E	24	Total 24	O 24	0	0
4	F	25	Total 25	O 25	0	0



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	40.12Å 127.53Å 41.47Å 90.00° 100.23° 90.00°	Depositor
Resolution (Å)	24.19 – 2.00 24.19 – 2.00	Depositor EDS
% Data completeness (in resolution range)	93.8 (24.19-2.00) 93.9 (24.19-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.59 (at 1.99Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.6_289)	Depositor
R, R_{free}	0.191 , 0.233 0.186 , 0.233	Depositor DCC
R_{free} test set	1293 reflections (4.97%)	DCC
Wilson B-factor (Å ²)	28.5	Xtriage
Anisotropy	0.531	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 46.3	EDS
Estimated twinning fraction	0.041 for l,-k,h	Xtriage
L-test for twinning	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 26881 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	3034	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.08% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality i

5.1 Standard geometry i

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	0.41	0/928	0.58	0/1252
1	D	0.41	0/879	0.56	0/1187
2	B	0.83	0/272	1.61	6/418 (1.4%)
2	E	0.69	0/272	1.48	7/418 (1.7%)
3	C	0.84	0/272	1.60	7/418 (1.7%)
3	F	0.74	0/272	1.61	7/418 (1.7%)
All	All	0.58	0/2895	1.10	27/4111 (0.7%)

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

There are no bond length outliers.

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	F	10	DT	C1'-O4'-C4'	-7.38	102.72	110.10
2	B	10	DT	O4'-C1'-N1	-7.09	103.03	108.00
2	B	3	DA	O4'-C1'-N9	-6.95	103.14	108.00
3	C	9	DA	O4'-C1'-N9	-6.78	103.26	108.00
3	F	2	DT	O4'-C1'-C2'	-6.16	100.97	105.90
2	B	6	DT	O4'-C1'-N1	-6.00	103.80	108.00
2	E	6	DT	N3-C4-O4	5.94	123.47	119.90
3	F	9	DA	P-O3'-C3'	5.94	126.83	119.70
3	F	8	DT	O4'-C1'-N1	-5.84	103.91	108.00
3	F	5	DC	O4'-C1'-N1	5.82	112.08	108.00
3	F	2	DT	C3'-C2'-C1'	-5.76	95.58	102.50
3	F	2	DT	C1'-O4'-C4'	-5.75	104.35	110.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	5	DC	P-O3'-C3'	5.65	126.48	119.70
2	E	10	DT	C5-C4-O4	-5.57	121.00	124.90
3	C	7	DA	P-O3'-C3'	5.51	126.31	119.70
2	E	6	DT	C5-C4-O4	-5.50	121.05	124.90
3	C	3	DA	O4'-C1'-N9	-5.49	104.16	108.00
2	E	12	DT	O4'-C1'-N1	5.37	111.76	108.00
2	B	2	DA	O4'-C1'-N9	-5.29	104.29	108.00
3	C	6	DG	O4'-C1'-N9	-5.29	104.30	108.00
2	B	12	DT	C5-C4-O4	-5.20	121.26	124.90
2	E	1	DA	O4'-C4'-C3'	-5.15	102.44	104.50
3	C	2	DT	C3'-C2'-C1'	-5.13	96.34	102.50
3	C	7	DA	C1'-O4'-C4'	-5.12	104.98	110.10
2	B	8	DG	O4'-C1'-N9	-5.12	104.42	108.00
2	E	10	DT	O4'-C1'-N1	-5.07	104.45	108.00
2	E	10	DT	N3-C4-O4	5.07	122.94	119.90

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	40	VAL	Peptide

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	906	0	909	18	0
1	D	856	0	855	18	0
2	B	243	0	139	2	0
2	E	243	0	139	2	0
3	C	243	0	139	1	0
3	F	243	0	139	1	0
4	A	83	0	0	4	0
4	B	44	0	0	0	0
4	C	42	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	D	82	0	0	0	0
4	E	24	0	0	1	0
4	F	25	0	0	0	0
All	All	3034	0	2320	39	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 8.

All (39) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:129:PHE:HB3	1:A:132:ARG:HG3	1.55	0.87
1:D:48:LYS:NZ	1:D:48:LYS:HB3	1.95	0.81
1:A:42:VAL:O	1:A:43:ASN:HB2	1.93	0.68
3:F:8:DT:H2''	3:F:9:DA:C8	2.30	0.67
1:D:70:ASP:OD2	1:D:73:MET:HE1	1.96	0.66
1:D:48:LYS:HZ3	1:D:48:LYS:HB3	1.60	0.66
1:A:78:PHE:O	4:A:234:HOH:O	2.15	0.64
1:A:109:ALA:O	1:A:112:VAL:HG22	2.02	0.60
1:A:133:THR:H	1:A:136:HIS:HD2	1.52	0.56
1:A:133:THR:H	1:A:136:HIS:CD2	2.24	0.55
1:D:141:TRP:O	1:D:145:SER:HB2	2.08	0.54
1:D:111:LEU:HD23	1:D:144:ILE:HD12	1.90	0.52
1:D:150:ILE:HG22	1:D:150:ILE:O	2.10	0.52
1:D:124:ILE:HA	1:D:127:LYS:HD2	1.91	0.52
1:D:120:ARG:CZ	1:D:122:TRP:HZ2	2.23	0.51
1:A:107:GLU:OE2	1:A:132:ARG:HD2	2.09	0.51
1:A:43:ASN:N	1:A:45:LYS:O	2.41	0.51
1:D:49:LYS:HE3	2:E:6:DT:O2	2.10	0.51
1:D:133:THR:H	1:D:136:HIS:CD2	2.29	0.50
1:A:51:LYS:HD3	2:B:6:DT:H5'	1.95	0.49
1:D:108:ASP:OD2	1:D:144:ILE:HD11	2.14	0.47
1:A:75:ALA:O	4:A:234:HOH:O	2.21	0.46
1:A:45:LYS:HA	1:A:45:LYS:HD2	1.68	0.45
1:D:49:LYS:HG2	4:E:26:HOH:O	2.16	0.45
1:A:150:ILE:HG22	1:A:150:ILE:O	2.17	0.45
2:E:11:DA:H1'	2:E:12:DT:H5''	1.98	0.45
1:D:48:LYS:HZ2	1:D:48:LYS:HB3	1.75	0.45
1:A:40:VAL:HG22	1:A:51:LYS:O	2.18	0.44
2:B:11:DA:C2	3:C:3:DA:C2	3.06	0.44
1:A:43:ASN:OD1	4:A:227:HOH:O	2.21	0.44
1:A:112:VAL:O	1:A:116:GLN:HG3	2.18	0.43
1:D:112:VAL:O	1:D:115:ILE:HG22	2.18	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:60:LEU:HD13	1:A:94:LEU:HD12	2.02	0.41
1:D:137:ILE:HG22	1:D:138:LYS:N	2.36	0.41
1:D:50:GLN:NE2	1:D:80:ASN:OD1	2.53	0.41
1:D:48:LYS:CB	1:D:48:LYS:NZ	2.75	0.41
1:D:120:ARG:NH2	1:D:122:TRP:HZ2	2.19	0.41
1:A:91:LYS:HD3	4:A:197:HOH:O	2.21	0.40
1:A:113:GLN:HA	1:A:116:GLN:HE21	1.86	0.40

There are no symmetry-related clashes.

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	106/126 (84%)	102 (96%)	4 (4%)	0	100	100
1	D	101/126 (80%)	97 (96%)	4 (4%)	0	100	100
All	All	207/252 (82%)	199 (96%)	8 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	94/110 (86%)	86 (92%)	8 (8%)	15	9
1	D	89/110 (81%)	83 (93%)	6 (7%)	23	16
All	All	183/220 (83%)	169 (92%)	14 (8%)	18	11

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	40	VAL
1	A	41	GLN
1	A	42	VAL
1	A	43	ASN
1	A	45	LYS
1	A	60	LEU
1	A	84	ARG
1	A	132	ARG
1	D	48	LYS
1	D	50	GLN
1	D	60	LEU
1	D	73	MET
1	D	84	ARG
1	D	145	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	43	ASN
1	A	66	GLN
1	A	113	GLN
1	A	116	GLN
1	A	136	HIS
1	D	50	GLN
1	D	67	HIS
1	D	113	GLN
1	D	136	HIS

5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

There are no ligands in this entry.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

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	110/126 (87%)	0.31	3 (2%) 52 52	21, 34, 47, 54	0
1	D	103/126 (81%)	0.36	6 (5%) 22 22	19, 34, 48, 61	0
2	B	12/12 (100%)	-0.49	0 100 100	20, 26, 32, 33	0
2	E	12/12 (100%)	-0.28	0 100 100	25, 35, 40, 41	0
3	C	12/12 (100%)	-0.63	0 100 100	19, 24, 29, 30	0
3	F	12/12 (100%)	-0.44	0 100 100	31, 34, 38, 39	0
All	All	261/300 (87%)	0.19	9 (3%) 43 43	19, 34, 47, 61	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	150	ILE	4.9
1	A	150	ILE	3.7
1	D	149	GLY	3.2
1	A	135	ILE	3.0
1	D	135	ILE	2.7
1	A	40	VAL	2.7
1	D	102	PRO	2.3
1	D	137	ILE	2.3
1	D	100	HIS	2.2

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.