

Full wwPDB X-ray Structure Validation Report (i)

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PDB ID	:	2VPM
Title	:	Trypanothione synthetase
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Deposited on	:	2008-03-03
Resolution	:	2.80 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 (i) symbol.

The types of validation reports are described at http://www.wwpdb.org/validation/2017/FAQs#types.

The following versions of software and data (see references (1)) were used in the production of this report:

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.36
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)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: $X\text{-}RAY \, DIFFRACTION$

The reported resolution of this entry is 2.80 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	$egin{array}{c} { m Whole \ archive} \ (\#{ m Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries,\ resolution\ range}({ m \AA}))$
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

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 of 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 <=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	А	653	80%	11%	8%
1	В	653	3% 	11%	9%



2VPM

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 9867 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 TRYPANOTHIONE SYNTHETASE.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace	
1	А	600	Total 4865	C 3116	N 830	O 897	S 22	0	0	0
1	В	596	Total 4832	C 3095	N 825	0 891	S 21	0	0	0

• Molecule 2 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	А	5	Total Cl 5 5	0	0
2	В	3	Total Cl 3 3	0	0

• Molecule 3 is BROMIDE ION (three-letter code: BR) (formula: Br).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	А	1	Total Br 1 1	0	0

• Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	А	73	Total O 73 73	0	0
4	В	88	Total O 88 88	0	0



3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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: TRYPANOTHIONE SYNTHETASE



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants	70.52Å 167.08Å 84.92Å	Deperitor
a, b, c, α , β , γ	90.00° 94.11° 90.00°	Depositor
Bosolution(A)	84.82 - 2.80	Depositor
Resolution (A)	20.79 - 2.80	EDS
% Data completeness	100.0 (84.82-2.80)	Depositor
(in resolution range)	$100.0\ (20.79-2.80)$	EDS
R_{merge}	0.07	Depositor
R _{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	$2.10 (at 2.79 \text{\AA})$	Xtriage
Refinement program	REFMAC 5.3.0027	Depositor
B B.	0.203 , 0.253	Depositor
II, II, <i>free</i>	0.207 , (Not available)	DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor $(Å^2)$	55.8	Xtriage
Anisotropy	0.057	Xtriage
Bulk solvent $k_{sol}(e/Å^3), B_{sol}(Å^2)$	0.31 , 50.1	EDS
L-test for twinning ²	$ < L >=0.47, < L^2>=0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	9867	wwPDB-VP
Average B, all atoms $(Å^2)$	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 4.30% of the height of the origin peak. No significant pseudotranslation is detected.

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.



¹Intensities estimated from amplitudes.

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: BR, CL

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).

Mal	Chain	Bond	lengths	Bond angles		
1VIOI	Chain	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	А	0.49	0/5004	0.60	0/6789	
1	В	0.49	0/4971	0.61	1/6746~(0.0%)	
All	All	0.49	0/9975	0.61	1/13535~(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	А	0	3
1	В	0	2
All	All	0	5

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
1	В	503	GLU	N-CA-C	-5.20	96.95	111.00

There are no chirality outliers.

All (5)) planarity	outliers are	e listed	below:
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Mol	Chain	Res	Type	Group
1	А	343	PHE	Peptide
1	А	502	PHE	Peptide
1	А	56	LYS	Peptide
1	В	502	PHE	Peptide
1	В	56	LYS	Peptide



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	А	4865	0	4669	44	0
1	В	4832	0	4632	43	0
2	А	5	0	0	0	0
2	В	3	0	0	0	0
3	А	1	0	0	0	0
4	А	73	0	0	1	0
4	В	88	0	0	3	0
All	All	9867	0	9301	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 4.

All (84) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom 1	Atom 2	Interatomic	Clash
Atom-1	Atom-2	$distance ({ m \AA})$	overlap (Å)
1:A:293:GLN:HE22	1:A:522:ASN:HD21	1.15	0.94
1:A:225:PHE:O	1:A:262:VAL:HG13	1.72	0.90
1:B:101:VAL:HG11	1:B:189:LEU:HG	1.67	0.74
1:A:222:ARG:NE	1:A:614:GLU:OE2	2.26	0.69
1:B:351:SER:O	1:B:352:THR:HB	1.92	0.67
1:A:623:ASP:N	1:A:623:ASP:OD1	2.28	0.67
1:A:293:GLN:HE22	1:A:522:ASN:ND2	1.90	0.66
1:B:204:VAL:HG13	4:B:2042:HOH:O	1.94	0.66
1:A:5:GLN:HE22	1:B:586:PHE:H	1.42	0.65
1:A:220:LEU:HD21	1:A:612:ILE:HD11	1.79	0.64
1:B:403:ASP:OD1	1:B:404:GLU:N	2.32	0.62
1:A:407:GLU:HA	1:A:410:TYR:CE1	2.34	0.62
1:A:39:HIS:CD2	1:A:387:MET:HG3	2.35	0.61
1:B:33:TYR:HB2	1:B:57:TYR:CG	2.38	0.59
1:A:344:GLU:CD	1:A:513:LYS:HZ1	2.07	0.58
1:A:36:LYS:HE3	1:A:605:ASP:OD2	2.05	0.56
1:A:543:HIS:CE1	1:A:584:GLN:HE21	2.24	0.55
1:A:162:LEU:HD22	1:A:175:ASP:HA	1.89	0.55
1:B:220:LEU:HD21	1:B:612:ILE:HD11	1.87	0.55
1:A:326:SER:OG	1:A:350:ALA:HB2	2.07	0.54



		Interatomic	Clash
Atom-1	Atom-2	distance (Å)	overlap (Å)
1:A:347:ALA:HB1	1:A:507:LYS:HE2	1.89	0.54
1:A:39:HIS:CG	1:A:387:MET:HG3	2.43	0.54
1:B:0:HIS:CG	1:B:1:MET:H	2.26	0.54
1:B:162:LEU:HD22	1:B:175:ASP:HA	1.90	0.53
1:A:225:PHE:CE1	1:A:226:LEU:HD22	2.44	0.53
1:B:150:ARG:NH1	1:B:177:ILE:HD12	2.24	0.52
1:A:51:ILE:CG1	1:A:206:LEU:HD22	2.40	0.52
1:B:295:ILE:HD12	1:B:318:TRP:HB2	1.92	0.52
1:A:5:GLN:HE22	1:B:586:PHE:N	2.09	0.51
1:A:590:LYS:HE3	1:A:595:TYR:CZ	2.46	0.50
1:A:5:GLN:NE2	1:B:586:PHE:H	2.08	0.50
1:B:235:LEU:HD22	1:B:242:GLU:HG2	1.93	0.50
1:A:239:ASN:HD21	1:A:417:GLN:HE22	1.59	0.50
1:A:444:VAL:HG12	1:A:450:ARG:HA	1.93	0.50
1:B:503:GLU:OE1	1:B:507:LYS:NZ	2.41	0.50
1:A:138:GLY:HA3	1:A:141:TYR:CE1	2.46	0.50
1:B:593:ASP:O	1:B:616:LYS:HB3	2.11	0.50
1:A:51:ILE:HG12	1:A:206:LEU:HD22	1.93	0.49
1:B:95:THR:O	1:B:96:THR:HB	2.11	0.49
1:B:439:ASP:N	1:B:439:ASP:OD1	2.44	0.49
1:B:549:PRO:HA	1:B:580:MET:HB3	1.95	0.49
1:A:58:GLN:HE21	1:A:148:ASN:ND2	2.11	0.49
1:A:443:VAL:HG12	1:A:451:VAL:HG21	1.95	0.49
1:B:221:LEU:HD21	4:B:2057:HOH:O	2.11	0.49
1:B:613:ARG:NH1	1:B:624:SER:O	2.45	0.49
1:B:222:ARG:NE	1:B:614:GLU:OE2	2.36	0.48
1:A:613:ARG:NH1	1:A:624:SER:O	2.47	0.48
1:B:294:VAL:HG12	1:B:301:LEU:HD23	1.96	0.48
1:A:291:THR:HG23	1:A:509:ILE:CG2	2.43	0.48
1:A:293:GLN:NE2	1:A:522:ASN:HD21	1.97	0.47
1:B:0:HIS:CD2	1:B:1:MET:H	2.32	0.47
1:B:597:ILE:HG12	1:B:619:ILE:HD12	1.95	0.47
1:B:101:VAL:CG1	1:B:189:LEU:HG	2.39	0.47
1:A:291:THR:HG23	1:A:509:ILE:HG21	1.97	0.46
1:A:331:PHE:HB2	1:A:598:ILE:HB	1.97	0.46
1:A:474:ARG:NH1	1:A:478:TRP:CZ3	2.83	0.46
1:A:39:HIS:HB2	1:A:387:MET:HE2	1.98	0.46
1:B:235:LEU:HD12	1:B:246:VAL:HG22	1.98	0.46
1:B:33:TYR:HB2	1:B:57:TYR:CD1	2.52	0.45
1:B:331:PHE:HB2	1:B:598:ILE:HB	1.98	0.45
1:A:550:ILE:HA	4:A:2068:HOH:O	2.17	0.44



A + 1	A + a	Interatomic	Clash
Atom-1	Atom-2	distance (\AA)	overlap (Å)
1:A:17:ILE:HB	1:A:18:PRO:HD2	2.00	0.44
1:A:58:GLN:HE21	1:A:148:ASN:HD21	1.66	0.44
1:A:352:THR:HA	1:A:355:GLU:OE1	2.17	0.44
1:B:536:LEU:HD12	1:B:537:THR:H	1.83	0.43
1:A:84:ILE:O	1:A:189:LEU:HD11	2.19	0.43
1:A:313:ARG:HA	1:A:497:TRP:CZ3	2.54	0.43
1:B:294:VAL:HG21	1:B:304:PHE:CZ	2.54	0.43
1:B:177:ILE:HG22	1:B:178:ASP:OD1	2.19	0.42
1:A:413:LEU:HD21	1:A:428:LEU:HD22	2.01	0.42
1:A:240:PRO:HB2	1:A:413:LEU:HD13	2.01	0.42
1:B:273:HIS:O	1:B:277:VAL:HG23	2.19	0.42
1:B:402:VAL:HG12	1:B:457:THR:OG1	2.20	0.42
1:B:332:ALA:HB2	1:B:343:PHE:CE2	2.55	0.41
1:B:340:VAL:O	1:B:528:ALA:HA	2.20	0.41
1:A:443:VAL:HG12	1:A:451:VAL:CG2	2.50	0.41
1:B:101:VAL:CG1	1:B:102:LEU:N	2.84	0.41
1:B:513:LYS:HB3	1:B:581:ILE:HD13	2.02	0.41
1:A:344:GLU:HG3	1:A:513:LYS:HZ2	1.86	0.41
1:B:416:MET:HB2	1:B:416:MET:HE2	1.96	0.41
1:B:325:ILE:O	1:B:603:ILE:HA	2.21	0.41
1:B:169:GLY:HA2	4:B:2039:HOH:O	2.21	0.40
1:B:510:PRO:HA	1:B:515:ILE:HD13	2.03	0.40
1:B:70:ARG:NH2	1:B:196:ASN:ND2	2.70	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	А	594/653~(91%)	571 (96%)	22 (4%)	1 (0%)	47 78
1	В	590/653~(90%)	566 (96%)	20 (3%)	4 (1%)	22 53



Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	1184/1306~(91%)	1137 (96%)	42 (4%)	5 (0%)	34 66

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	А	124	ASP
1	В	57	TYR
1	В	124	ASP
1	В	232	ALA
1	В	621	GLY

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	А	513/557~(92%)	499~(97%)	14 (3%)	44 78
1	В	509/557~(91%)	492 (97%)	17 (3%)	38 72
All	All	1022/1114 (92%)	991~(97%)	31 (3%)	41 75

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	А	10	SER
1	А	58	GLN
1	А	77	ASP
1	А	98	SER
1	А	126	ASN
1	А	149	TYR
1	А	226	LEU
1	А	234	TRP
1	А	238	ASN
1	А	251	MET
1	А	456	LYS
1	А	587	GLU
1	А	618	VAL



Mol	Chain	Res	Type
1	А	623	ASP
1	В	16	HIS
1	В	70	ARG
1	В	96	THR
1	В	135	THR
1	В	180	ASP
1	В	189	LEU
1	В	194	ARG
1	В	214	GLU
1	В	226	LEU
1	В	295	ILE
1	В	349	SER
1	В	352	THR
1	В	361	GLN
1	В	410	TYR
1	В	439	ASP
1	В	456	LYS
1	В	581	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	А	5	GLN
1	А	35	ASN
1	А	126	ASN
1	А	148	ASN
1	А	152	HIS
1	А	293	GLN
1	А	417	GLN
1	А	584	GLN
1	В	0	HIS
1	В	543	HIS
1	В	584	GLN

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 monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 9 ligands modelled in this entry, 9 are monoatomic - leaving 0 for Mogul analysis.

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.

No monomer is involved in short contacts.

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^{th} 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	А	600/653~(91%)	-0.09	27 (4%) 33 23	38, 48, 68, 78	27 (4%)
1	В	596/653~(91%)	-0.10	21 (3%) 44 34	39, 49, 64, 74	23 (3%)
All	All	1196/1306 (91%)	-0.10	48 (4%) 38 28	38, 49, 66, 78	50 (4%)

All (48) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ	
1	А	636	LEU	5.2	
1	В	233	ASN	5.1	
1	В	232	ALA	4.3	
1	А	233	ASN	3.7	
1	А	232	ALA	3.7	
1	А	477	ASN	3.7	
1	А	543	HIS	3.6	
1	А	228	THR	3.5	
1	А	476	GLU	3.4	
1	В	372	GLN	3.2	
1	А	179	ALA	3.2	
1	В	199	GLU	3.2	
1	А	439	ASP	3.2	
1	В	375	ARG	3.0	
1	В	476	GLU	3.0	
1	А	251	MET	3.0	
1	А	634	ASP	2.9	
1	В	178	ASP	2.8	
1	В	477	ASN	2.8	
1	А	635	LYS	2.7	
1	А	178	ASP	2.7	
1	А	317	SER	2.6	
1	A	633	THR	2.5	
1	А	238	ASN	2.5	



Mol	Chain	Res	Type	RSRZ
1	А	155	GLU	2.5
1	А	15	GLY	2.5
1	В	238	ASN	2.4
1	А	180	ASP	2.3
1	А	415	CYS	2.3
1	В	180	ASP	2.2
1	В	179	ALA	2.2
1	А	375	ARG	2.2
1	В	337	THR	2.2
1	А	14	PRO	2.2
1	А	177	ILE	2.2
1	А	475	GLY	2.2
1	В	228	THR	2.2
1	А	478	TRP	2.1
1	В	229	GLU	2.1
1	В	633	THR	2.1
1	В	230	SER	2.1
1	В	472	GLU	2.1
1	В	475	GLY	2.1
1	В	231	LYS	2.1
1	В	80	TRP	2.1
1	А	199	GLU	2.1
1	А	230	SER	2.0
1	В	14	PRO	2.0

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 monosaccharides 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^{th} 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 ²)	Q<0.9
3	BR	А	1642	1/1	0.97	0.11	70,70,70,70	0
2	CL	А	1638	1/1	0.98	0.04	45,45,45,45	0
2	CL	А	1639	1/1	0.98	0.04	21,21,21,21	0
2	CL	А	1640	1/1	0.98	0.04	33,33,33,33	0
2	CL	А	1641	1/1	0.98	0.06	20,20,20,20	0
2	CL	В	1635	1/1	0.98	0.05	45,45,45,45	0
2	CL	В	1636	1/1	0.98	0.04	$35,\!35,\!35,\!35$	0
2	CL	A	1637	1/1	0.98	0.04	17,17,17,17	0
2	CL	В	1634	1/1	0.99	0.06	19,19,19,19	0

6.5 Other polymers (i)

There are no such residues in this entry.

