

Full wwPDB X-ray Structure Validation Report (i)

Oct 20, 2021 – 11:36 am BST

PDB ID	:	7PHX
Title	:	Tsetse thrombin inhibitor in complex with human alpha-thrombin - acid-stable
		sulfotyrosine analogue
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Deposited on		
Resolution	:	1.80 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 following versions of software and data (see references (1)) were used in the production of this report:

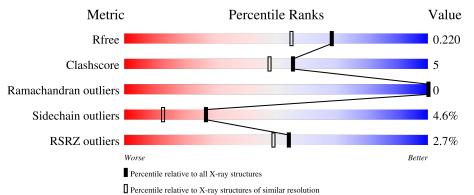
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.23.2
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0267
CCP4	:	7.1.010 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.23.2

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 1.80 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	$\begin{array}{c} \textbf{Whole archive} \\ (\#\textbf{Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries,\ resolution\ range}({ m \AA}))$
R_{free}	130704	5950(1.80-1.80)
Clashscore	141614	6793 (1.80-1.80)
Ramachandran outliers	138981	6697 (1.80-1.80)
Sidechain outliers	138945	6696 (1.80-1.80)
RSRZ outliers	127900	5850 (1.80-1.80)

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	L	36	6% 75% • •	19%
2	Н	259	^{2%} 82%	12% • 5%
3	Ι	32	^{3%} 69% • •	25%



$7\mathrm{PHX}$

2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 2590 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 Thrombin light chain.

Mol	Chain	Residues		Atc	\mathbf{ms}			ZeroOcc	AltConf	Trace
1	L	29	Total 234	C 146	N 38	O 49	S 1	0	0	0

• Molecule 2 is a protein called Thrombin heavy chain.

Mol	Chain	Residues		At	oms			ZeroOcc	AltConf	Trace
2	Н	247	Total 2027	C 1291	N 361	O 361	S 14	0	4	0

• Molecule 3 is a protein called Tsetse thrombin inhibitor.

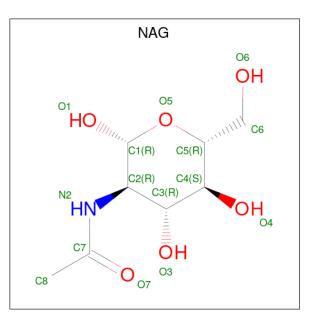
Mol	Chain	Residues		Atc	\mathbf{ms}			ZeroOcc	AltConf	Trace
3	Ι	24	Total 181	C 112	N 29	O 39	S 1	0	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
Ι	9	SMF	TYR	modified residue	UNP 097373
Ι	12	SMF	TYR	modified residue	UNP 097373

• Molecule 4 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).





Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	Н	1	Total C N O 14 8 1 5	0	0

• Molecule 5 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	Н	1	Total Na 1 1	0	0

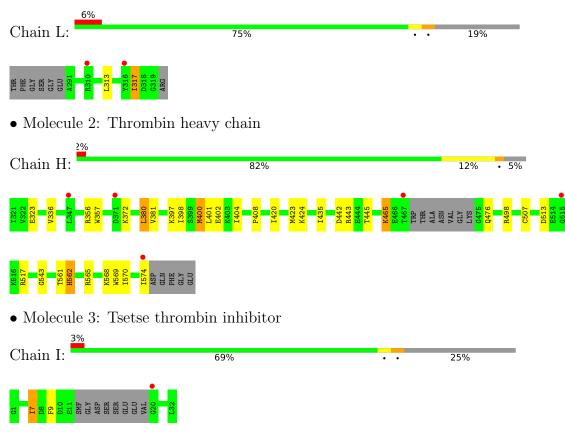
• Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	L	9	Total O 9 9	0	0
6	Н	111	Total O 111 111	0	0
6	Ι	13	Total O 13 13	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: Thrombin light chain



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants	43.70Å 80.79Å 85.89Å	Depositor
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	42.95 - 1.80	Depositor
Resolution (A)	42.95 - 1.80	EDS
% Data completeness	95.0(42.95-1.80)	Depositor
(in resolution range)	95.2(42.95-1.80)	EDS
R _{merge}	0.04	Depositor
R _{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	$2.50 (at 1.79 \text{\AA})$	Xtriage
Refinement program	PHENIX 1.19.2_4158, PHENIX 1.19.2_4158	Depositor
R, R_{free}	0.183 , 0.221	Depositor
$\mathbf{n}, \mathbf{n}_{free}$	0.182 , 0.220	DCC
R_{free} test set	1338 reflections (4.86%)	wwPDB-VP
Wilson B-factor $(Å^2)$	30.7	Xtriage
Anisotropy	0.846	Xtriage
Bulk solvent $k_{sol}(e/Å^3), B_{sol}(Å^2)$	(Not available), (Not available)	EDS
L-test for twinning ²	$< L >=0.45, < L^2>=0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	2590	wwPDB-VP
Average B, all atoms $(Å^2)$	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 9.69% 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: NAG, SMF, NA

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		
	Unam	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	L	0.56	0/236	0.68	0/314	
2	Н	0.42	0/2078	0.64	0/2808	
3	Ι	0.34	0/167	0.58	0/223	
All	All	0.43	0/2481	0.64	0/3345	

There are no bond length outliers.

There are no bond angle outliers.

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	L	234	0	229	1	0
2	Н	2027	0	2008	24	0
3	Ι	181	0	166	1	0
4	Н	14	0	13	0	0
5	Н	1	0	0	0	0
6	Н	111	0	0	2	0
6	Ι	13	0	0	0	0
6	L	9	0	0	0	0
All	All	2590	0	2416	25	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 5.

All (25) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

A 4 1	A + 9	Interatomic	Clash
Atom-1	Atom-2	distance (\AA)	overlap (Å)
2:H:498:ARG:HG3	6:H:701:HOH:O	1.99	0.61
2:H:380:LEU:HD11	2:H:400:MET:HG3	1.85	0.58
2:H:336:VAL:HG13	2:H:381[B]:VAL:HG13	1.88	0.56
2:H:404:ILE:HG12	2:H:423:MET:HG2	1.89	0.55
2:H:397:LYS:HG3	2:H:435:ILE:CD1	2.38	0.53
2:H:323:GLU:HG3	2:H:517:ARG:HG3	1.89	0.53
2:H:562:HIS:CD2	2:H:565:ARG:HB2	2.44	0.52
2:H:356:ARG:HG3	2:H:357:TRP:CD1	2.48	0.48
2:H:420:ILE:HD11	2:H:570:ILE:HD11	1.98	0.46
2:H:543:GLY:HA2	2:H:561:THR:O	2.16	0.45
2:H:465:LYS:HE2	2:H:465:LYS:HB2	1.88	0.44
2:H:356:ARG:HD3	2:H:424:LYS:HE3	2.00	0.44
2:H:357:TRP:CE2	2:H:574:ILE:HG12	2.53	0.44
1:L:317:ILE:O	1:L:317:ILE:HG13	2.18	0.44
2:H:408:PRO:HD3	2:H:569:TRP:NE1	2.33	0.43
2:H:380:LEU:CD1	2:H:400:MET:HG3	2.49	0.42
2:H:336:VAL:CG1	2:H:381[B]:VAL:HG13	2.50	0.42
2:H:442:ASP:OD1	2:H:445:THR:N	2.38	0.42
2:H:372:LYS:HG2	6:H:773:HOH:O	2.19	0.42
2:H:402:GLU:N	2:H:424:LYS:O	2.44	0.42
2:H:380:LEU:HD12	2:H:380:LEU:HA	1.71	0.41
2:H:443:ARG:NH1	3:I:7:ILE:O	2.54	0.41
2:H:408:PRO:HD3	2:H:569:TRP:HE1	1.86	0.41
2:H:357:TRP:CZ2	2:H:424:LYS:HD2	2.55	0.41
2:H:401:LEU:HD22	2:H:423:MET:HB3	2.02	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	L	27/36~(75%)	26~(96%)	1 (4%)	0	100 100
2	Н	247/259~(95%)	241 (98%)	6~(2%)	0	100 100
3	Ι	19/32~(59%)	19 (100%)	0	0	100 100
All	All	293/327~(90%)	286 (98%)	7 (2%)	0	100 100

There are no Ramachandran outliers to report.

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	ed Rotameric Out		Percentiles
1	L	26/31~(84%)	24~(92%)	2 (8%)	13 4
2	Н	220/225~(98%)	211~(96%)	9~(4%)	30 16
3	Ι	17/23~(74%)	16 (94%)	1 (6%)	19 7
All	All	263/279~(94%)	251 (95%)	12 (5%)	27 13

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	L	313	LEU
1	L	317	ILE
2	Н	380	LEU
2	Н	398	ILE
2	Н	400	MET
2	Н	465	LYS
2	Н	476	GLN
2	Н	507	CYS
2	Н	513	ASP
2	Н	562	HIS
2	Н	568	LYS
3	Ι	7	ILE

Sometimes side chains can be flipped to improve hydrogen bonding and reduce clashes. There are no such side chains identified.



5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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	Dog	Link	Bond lengths			Bond angles		
	Moi Type	Ullalli	nes	LIIIK	Counts	RMSZ	# Z >2	Counts	RMSZ	# Z >2
3	SMF	Ι	9	3	$15,\!16,\!17$	2.03	2 (13%)	18,22,24	0.80	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.

Mo	bl	Type	Chain	Res	Link	Chirals	Torsions	Rings
3		SMF	Ι	9	3	-	0/10/11/13	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Ζ	Observed(Å)	Ideal(Å)
3	Ι	9	SMF	O2-S	6.41	1.63	1.45
3	Ι	9	SMF	CH-S	-3.94	1.67	1.77

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.



5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

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).

М		Type	Chain	Res Link		Bond lengths			Bond angles		
	Mol Type	Unann	nes	LIIIK	Counts	RMSZ	# Z >2	Counts	RMSZ	# Z > 2	
4	:	NAG	Н	601	2	14,14,15	1.29	1 (7%)	$17,\!19,\!21$	1.35	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
4	NAG	Н	601	2	-	1/6/23/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	Н	601	NAG	O5-C1	4.57	1.51	1.43

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
4	Н	601	NAG	C1-O5-C5	5.34	119.42	112.19

There are no chirality outliers.

All (1) torsion outliers are listed below:

	Mol	Chain	Res	Type	Atoms
ſ	4	Н	601	NAG	O5-C5-C6-O6



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	$\langle RSRZ \rangle$	#RSRZ>2	$OWAB(Å^2)$	Q<0.9
1	L	29/36~(80%)	0.35	2 (6%) 16 13	38, 51, 77, 91	0
2	Н	247/259~(95%)	0.18	5 (2%) 65 61	27, 43, 70, 88	0
3	Ι	23/32~(71%)	0.57	1 (4%) 35 29	46, 53, 72, 96	0
All	All	299/327~(91%)	0.23	8 (2%) 54 49	27, 45, 72, 96	0

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	Ι	20	GLY	3.8
2	Н	371	ASP	3.4
2	Н	467	THR	3.2
2	Н	574	ILE	2.5
2	Н	347	LEU	2.5
1	L	316	TYR	2.0
1	L	310	ARG	2.0
2	Н	515	GLY	2.0

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, 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	$\mathbf{B} ext{-factors}(\mathbf{A}^2)$	Q < 0.9
3	SMF	Ι	9	16/17	0.86	0.13	57,66,85,94	0



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	$\mathbf{B} ext{-factors}(\mathrm{\AA}^2)$	Q<0.9
4	NAG	Н	601	14/15	0.80	0.37	69,82,92,96	0
5	NA	Н	602	1/1	0.94	0.13	43,43,43,43	0

6.5 Other polymers (i)

There are no such residues in this entry.

