wwPDB X-ray Structure Validation Summary Report (i)

Aug 23, 2026 – 02:23 PM JST

PDB ID	:	9KGA / pdb_00009kga
Title	:	Crystal structure of SpTx1 toxin
Authors	:	Chen, L.; Wang, M.; Fan, C.
Deposited on	:	2024-11-08
Resolution	:	1.90 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 (i)) were used in the production of this report:

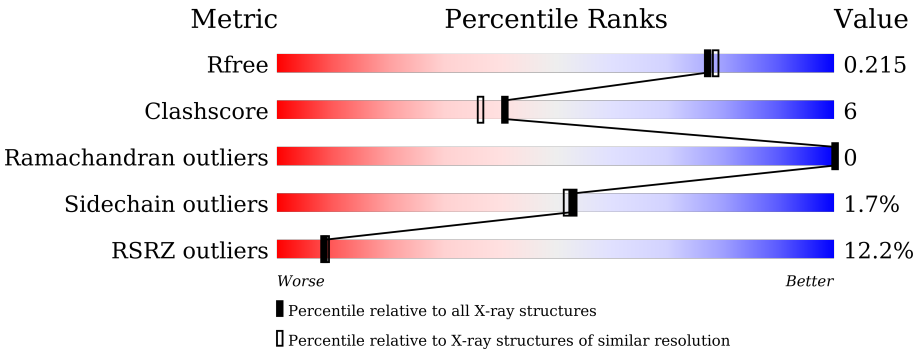
MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.015 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.50

1 Overall quality at a glance (i)

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}	180053	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (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 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	A	56	9% 84% 14% ·
1	B	56	2% 82% 16% ·
1	C	56	16% 84% 14% ·
1	D	56	23% 77% 18% 5%
1	E	56	7% 91% 7% ·
1	F	56	14% 77% 20% ·

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 3000 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 SpTx1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	55	Total	C	N	O	S	0	0	0
			453	289	78	82	4			
1	B	56	Total	C	N	O	S	0	0	0
			457	291	79	83	4			
1	C	56	Total	C	N	O	S	0	0	0
			453	288	78	83	4			
1	D	53	Total	C	N	O	S	0	0	0
			441	283	76	78	4			
1	E	55	Total	C	N	O	S	0	0	0
			451	288	78	81	4			
1	F	54	Total	C	N	O	S	0	0	0
			447	286	77	80	4			

- Molecule 2 is water.

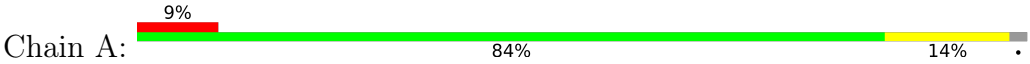
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	51	Total	O	0	0
			51	51		
2	B	61	Total	O	0	0
			61	61		
2	C	45	Total	O	0	0
			45	45		
2	D	37	Total	O	0	0
			37	37		
2	E	60	Total	O	0	0
			60	60		
2	F	44	Total	O	0	0
			44	44		

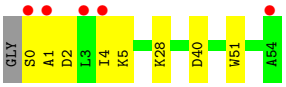
3 Residue-property plots (i)

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: SpTx1

Chain A:

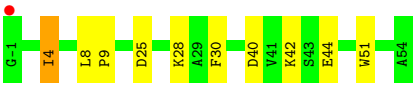




• Molecule 1: SpTx1

Chain B:

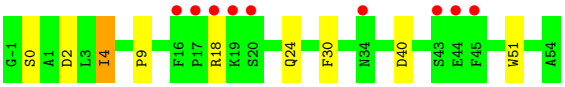




• Molecule 1: SpTx1

Chain C:

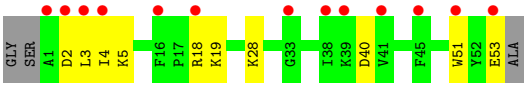




• Molecule 1: SpTx1

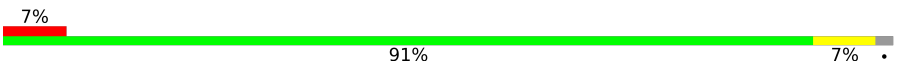
Chain D:

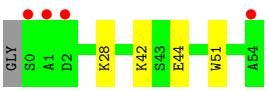




• Molecule 1: SpTx1

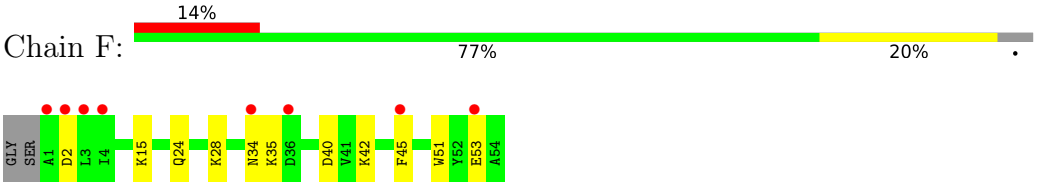
Chain E:





• Molecule 1: SpTx1

Chain F:


14%
77%
20%
GLY
SER
A1
D2
L3
I4
K15
Q24
K28
N34
K35
D36
D40
V41
K42
F45
W51
Y52
E53
A54

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α, β, γ	60.48Å 60.48Å 169.85Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	50.00 – 1.90 50.00 – 1.90	Depositor EDS
% Data completeness (in resolution range)	99.5 (50.00-1.90) 99.5 (50.00-1.90)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.28 (at 1.89Å)	Xtriage
Refinement program	PHENIX (1.19.2_4158: ???)	Depositor
R, R_{free}	0.203 , 0.214 0.204 , 0.215	Depositor DCC
R_{free} test set	1427 reflections (4.87%)	wwPDB-VP
Wilson B-factor (Å^2)	20.3	Xtriage
Anisotropy	0.033	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.41 , 55.6	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Estimated twinning fraction	0.037 for -h,-k,l	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	3000	wwPDB-VP
Average B, all atoms (Å^2)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 30.27 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.3448e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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

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	0.24	0/463	0.51	0/615
1	B	0.22	0/467	0.52	0/620
1	C	0.32	0/463	0.57	0/616
1	D	0.29	0/451	0.47	0/600
1	E	0.21	0/461	0.44	0/612
1	F	0.20	0/457	0.48	0/607
All	All	0.25	0/2762	0.50	0/3670

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

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	453	0	452	6	0
1	B	457	0	455	8	0
1	C	453	0	444	7	0
1	D	441	0	442	9	0
1	E	451	0	448	2	0
1	F	447	0	447	9	0
2	A	51	0	0	1	0
2	B	61	0	0	0	0
2	C	45	0	0	1	0
2	D	37	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	E	60	0	0	0	0
2	F	44	0	0	3	0
All	All	3000	0	2688	35	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 6.

The worst 5 of 35 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:D:19:LYS:NZ	2:D:101:HOH:O	2.22	0.68
1:A:5:LYS:NZ	2:A:101:HOH:O	2.30	0.64
1:F:2:ASP:O	1:F:53:GLU:HA	2.01	0.61
1:D:4:ILE:O	1:D:51:TRP:HA	2.03	0.58
1:A:0:SER:HB2	2:F:131:HOH:O	2.05	0.57

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	53/56 (95%)	53 (100%)	0	0	100	100
1	B	54/56 (96%)	54 (100%)	0	0	100	100
1	C	54/56 (96%)	54 (100%)	0	0	100	100
1	D	51/56 (91%)	51 (100%)	0	0	100	100
1	E	53/56 (95%)	52 (98%)	1 (2%)	0	100	100
1	F	52/56 (93%)	52 (100%)	0	0	100	100
All	All	317/336 (94%)	316 (100%)	1 (0%)	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	50/50 (100%)	49 (98%)	1 (2%)	48	46
1	B	50/50 (100%)	49 (98%)	1 (2%)	48	46
1	C	49/50 (98%)	48 (98%)	1 (2%)	48	46
1	D	49/50 (98%)	48 (98%)	1 (2%)	48	46
1	E	49/50 (98%)	48 (98%)	1 (2%)	48	46
1	F	49/50 (98%)	49 (100%)	0	100	100
All	All	296/300 (99%)	291 (98%)	5 (2%)	53	52

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	ILE
1	B	4	ILE
1	C	4	ILE
1	D	3	LEU
1	E	28	LYS

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

5.3.3 RNA

There are no RNA molecules 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 (i)

There are no oligosaccharides in this entry.

5.6 Ligand geometry (i)

There are no ligands in this entry.

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, 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	A	55/56 (98%)	0.35	5 (9%) 15 15	13, 24, 38, 58	0
1	B	56/56 (100%)	0.00	1 (1%) 67 71	12, 21, 32, 40	0
1	C	56/56 (100%)	0.95	9 (16%) 4 4	16, 28, 44, 49	0
1	D	53/56 (94%)	1.45	13 (24%) 2 1	17, 37, 54, 70	0
1	E	55/56 (98%)	0.22	4 (7%) 21 22	13, 23, 44, 57	0
1	F	54/56 (96%)	0.62	8 (14%) 5 6	14, 26, 47, 49	0
All	All	329/336 (97%)	0.59	40 (12%) 8 9	12, 26, 48, 70	0

The worst 5 of 40 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	1	ALA	6.9
1	D	2	ASP	5.8
1	D	4	ILE	4.7
1	C	16	PHE	4.6
1	A	0	SER	4.1

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 oligosaccharides in this entry.

6.4 Ligands

There are no ligands in this entry.

6.5 Other polymers (i)

There are no such residues in this entry.