



wwPDB X-ray Structure Validation Summary Report ⓘ

May 29, 2020 – 06:45 am BST

PDB ID : 4ZKT
Title : Crystal structure of the progenitor M complex of Clostridium botulinum type E neurotoxin
Authors : Eswaramoorthy, S.; Swaminathan, S.
Deposited on : 2015-04-30
Resolution : 3.05 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 ⓘ symbol.

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

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.11
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)	:	2.11

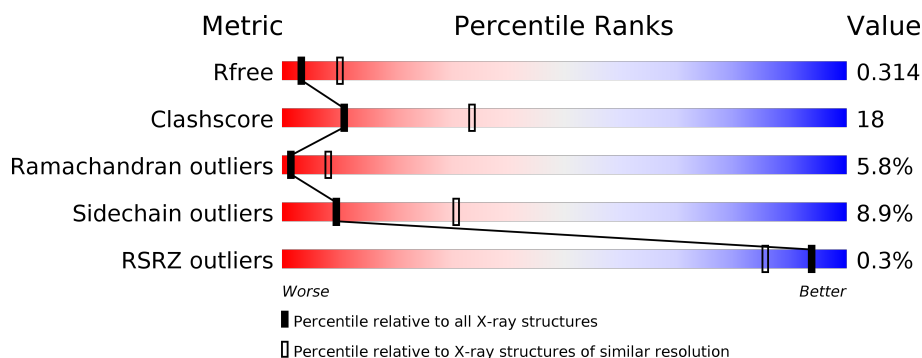
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.05 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}	130704	1754 (3.10-3.02)
Clashscore	141614	1864 (3.10-3.02)
Ramachandran outliers	138981	1794 (3.10-3.02)
Sidechain outliers	138945	1793 (3.10-3.02)
RSRZ outliers	127900	1713 (3.10-3.02)

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 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 $\leq 5\%$. 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	1252	
1	C	1252	
1	E	1252	
2	B	1163	
2	D	1163	
2	F	1163	

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 56952 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 Bontoxilysin A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	1235	Total	C	N	O	S	0	0	0
			9978	6351	1665	1940	22			
1	C	1235	Total	C	N	O	S	0	0	0
			9978	6351	1665	1940	22			
1	E	1235	Total	C	N	O	S	0	0	0
			9978	6351	1665	1940	22			

- Molecule 2 is a protein called Botulinum neurotoxin type E, nontoxic-nonhemagglutinin component, NTNH.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	1114	Total	C	N	O	S	0	0	0
			9005	5804	1451	1726	24			
2	D	1114	Total	C	N	O	S	0	0	0
			9005	5804	1451	1726	24			
2	F	1114	Total	C	N	O	S	0	0	0
			9005	5804	1451	1726	24			

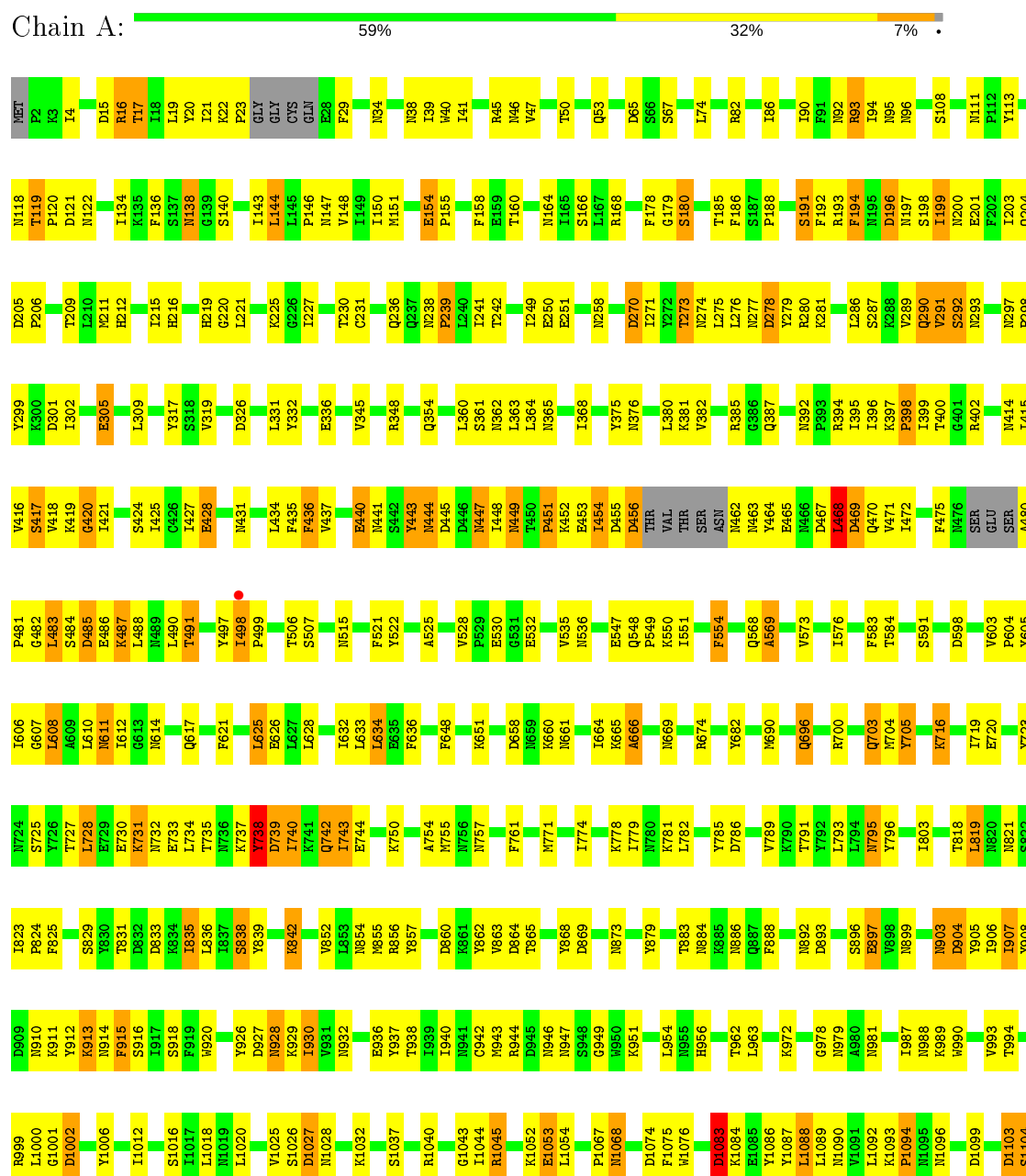
- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

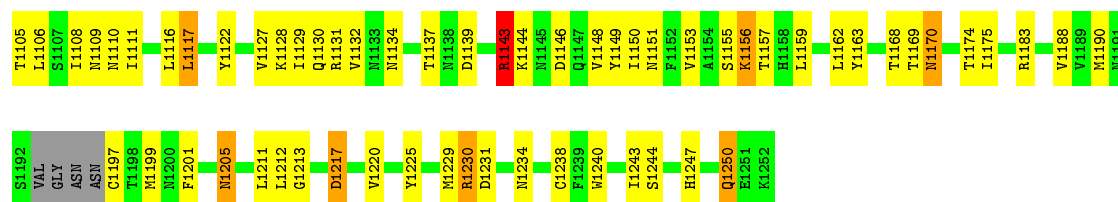
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Zn	0	0
			1	1		
3	C	1	Total	Zn	0	0
			1	1		
3	E	1	Total	Zn	0	0
			1	1		

3 Residue-property plots

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.

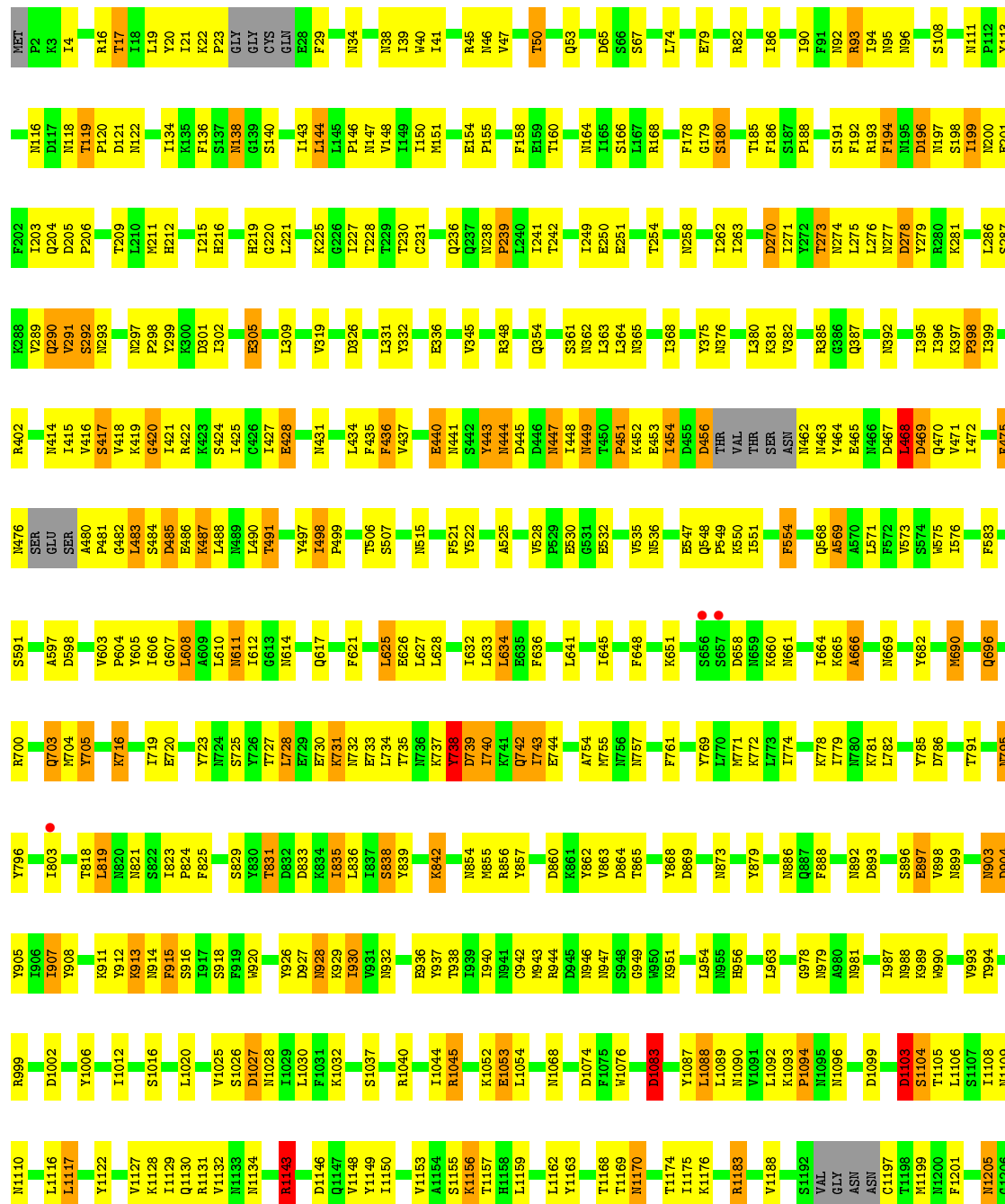
• Molecule 1: Bontoxilysin A

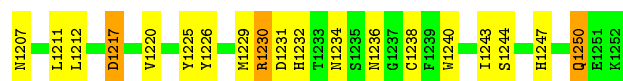




• Molecule 1: Bontoxilysin A

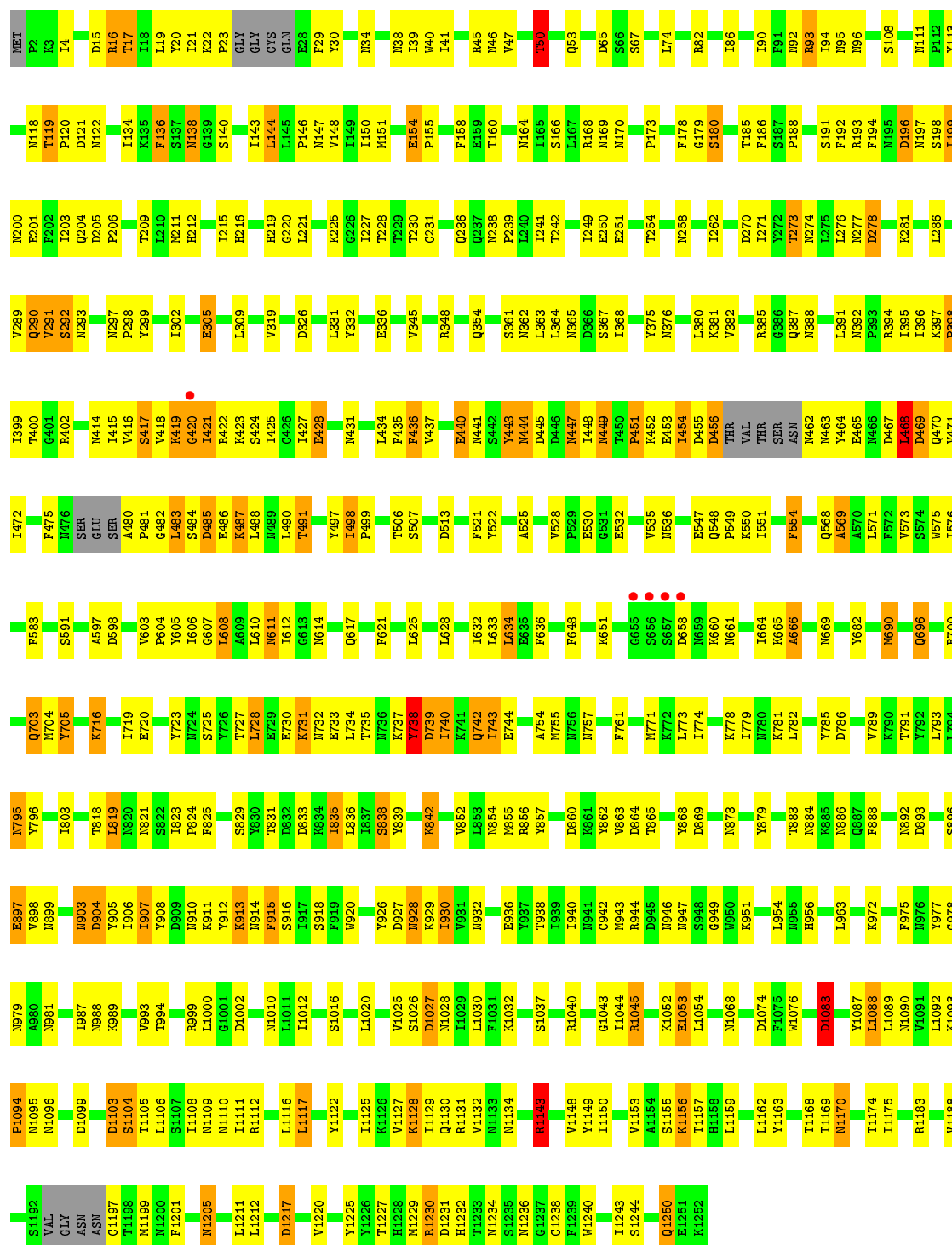
Chain C:  60%  32% 7% •





• Molecule 1: Bontoxilysin A

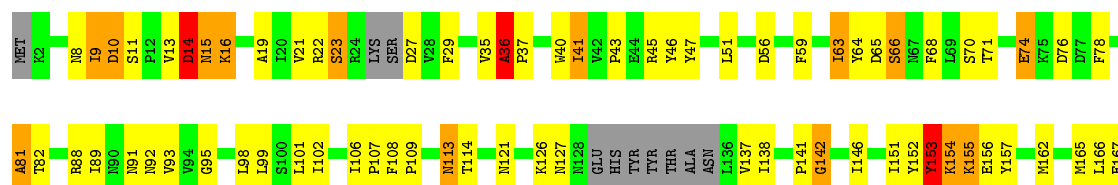
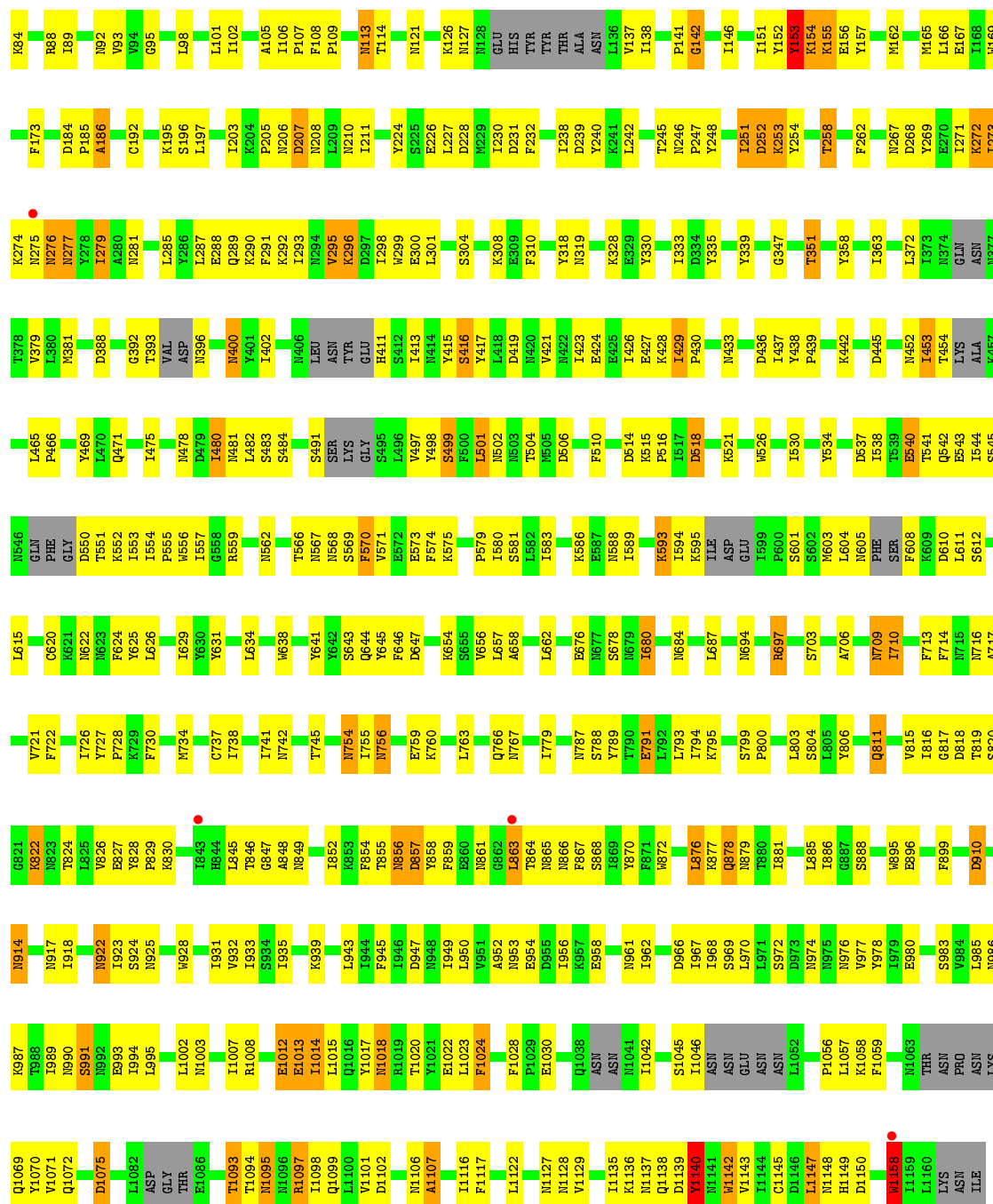
Chain E: 59% 32% 7% .



• Molecule 2: Botulinum neurotoxin type E, nontoxic-nonhemagglutinin component, NTNH

[illegible]

MET	K2	I9	D10	S11	P12	V13	D14	N15	K16	A19	L20	V21	R22	S23	R24	LVS	SER	D27	V28	F29	V35	A36	P37	W40	I41	V42	P43	E44	R45	Y46	Y47	L51	F59	I63	Y64	D65	S66	N67	F68	L69	S70	T71	E74	K75	D76	D77	F78	A81	T82	E83
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4 Data and refinement statistics

Property	Value	Source
Space group	P 31	Depositor
Cell constants a, b, c, α , β , γ	192.60Å 192.60Å 286.54Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	49.30 – 3.05 49.25 – 3.05	Depositor EDS
% Data completeness (in resolution range)	59.7 (49.30-3.05) 59.8 (49.25-3.05)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.46 (at 3.07Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.243 , 0.321 0.247 , 0.314	Depositor DCC
R_{free} test set	1334 reflections (0.99%)	wwPDB-VP
Wilson B-factor (Å ²)	110.6	Xtriage
Anisotropy	0.154	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.25 , 110.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.36$, $\langle L^2 \rangle = 0.19$	Xtriage
Estimated twinning fraction	0.135 for -h,-k,l 0.196 for h,-h-k,-l 0.137 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	56952	wwPDB-VP
Average B, all atoms (Å ²)	134.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.18% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ZN

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	0.58	0/10176	0.74	2/13800 (0.0%)
1	C	0.56	0/10176	0.73	1/13800 (0.0%)
1	E	0.54	0/10176	0.72	2/13800 (0.0%)
2	B	0.64	2/9182 (0.0%)	0.75	0/12454
2	D	0.60	1/9182 (0.0%)	0.74	0/12454
2	F	0.58	1/9182 (0.0%)	0.73	0/12454
All	All	0.58	4/58074 (0.0%)	0.73	5/78762 (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
2	B	0	1
2	D	0	1
2	F	0	2
All	All	0	4

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1158	TRP	CB-CG	-7.00	1.37	1.50
2	D	1158	TRP	CB-CG	-5.84	1.39	1.50
2	F	1158	TRP	CB-CG	-5.69	1.40	1.50
2	B	116	ASP	C-O	5.60	1.33	1.23

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1143	ARG	NE-CZ-NH2	7.34	123.97	120.30
1	E	1143	ARG	NE-CZ-NH2	7.21	123.90	120.30
1	A	394	ARG	NE-CZ-NH2	-6.68	116.96	120.30
1	C	1143	ARG	NE-CZ-NH2	6.28	123.44	120.30
1	E	394	ARG	NE-CZ-NH2	-5.63	117.48	120.30

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	36	ALA	Peptide
2	D	36	ALA	Peptide
2	F	1138	GLN	Peptide
2	F	36	ALA	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	9978	0	9735	338	1
1	C	9978	0	9735	348	1
1	E	9978	0	9735	352	0
2	B	9005	0	8666	361	0
2	D	9005	0	8666	344	0
2	F	9005	0	8666	356	0
3	A	1	0	0	0	0
3	C	1	0	0	0	0
3	E	1	0	0	0	0
All	All	56952	0	55203	2034	1

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 18.

The worst 5 of 2034 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)
2:D:885:LEU:HG	2:D:886:ILE:HD12	1.33	1.09

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:885:LEU:HG	2:F:886:ILE:HD12	1.32	1.09
2:B:436:ASP:HA	2:B:437:ILE:HB	1.34	1.08
2:B:885:LEU:HG	2:B:886:ILE:HD12	1.29	1.08
2:D:436:ASP:HA	2:D:437:ILE:HB	1.35	1.05

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:280:ARG:NH2	1:C:420:GLY:O[3_564]	2.16	0.04

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	1225/1252 (98%)	999 (82%)	159 (13%)	67 (6%)	2	9
1	C	1225/1252 (98%)	995 (81%)	166 (14%)	64 (5%)	2	10
1	E	1225/1252 (98%)	997 (81%)	162 (13%)	66 (5%)	2	9
2	B	1084/1163 (93%)	832 (77%)	183 (17%)	69 (6%)	1	7
2	D	1084/1163 (93%)	835 (77%)	179 (16%)	70 (6%)	1	6
2	F	1084/1163 (93%)	838 (77%)	177 (16%)	69 (6%)	1	7
All	All	6927/7245 (96%)	5496 (79%)	1026 (15%)	405 (6%)	1	8

5 of 405 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	119	THR
1	A	199	ILE
1	A	291	VAL
1	A	419	LYS

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Mol	Chain	Res	Type
1	A	449	ASN

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	1125/1155 (97%)	1013 (90%)	112 (10%)	7	25
1	C	1125/1155 (97%)	1011 (90%)	114 (10%)	7	25
1	E	1125/1155 (97%)	1014 (90%)	111 (10%)	8	25
2	B	990/1103 (90%)	915 (92%)	75 (8%)	13	38
2	D	990/1103 (90%)	915 (92%)	75 (8%)	13	38
2	F	990/1103 (90%)	914 (92%)	76 (8%)	13	38
All	All	6345/6774 (94%)	5782 (91%)	563 (9%)	9	31

5 of 563 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	348	ARG
1	E	1027	ASP
2	F	570	PHE
1	E	440	GLU
1	E	696	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 110 such sidechains are listed below:

Mol	Chain	Res	Type
1	E	956	HIS
2	B	636	GLN
2	F	677	ASN
1	E	1060	GLN
2	B	171	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 3 ligands modelled in this entry, 3 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, 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	1235/1252 (98%)	-0.69	1 (0%) 95 91	53, 120, 189, 296	0
1	C	1235/1252 (98%)	-0.66	3 (0%) 95 89	46, 129, 207, 365	0
1	E	1235/1252 (98%)	-0.67	5 (0%) 92 82	62, 143, 214, 391	0
2	B	1114/1163 (95%)	-0.69	1 (0%) 95 91	37, 109, 179, 274	0
2	D	1114/1163 (95%)	-0.68	4 (0%) 92 82	48, 129, 194, 278	0
2	F	1114/1163 (95%)	-0.61	10 (0%) 84 66	51, 143, 220, 328	0
All	All	7047/7245 (97%)	-0.67	24 (0%) 94 85	37, 129, 203, 391	0

The worst 5 of 24 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	657	SER	9.4
1	E	655	GLY	7.5
1	E	657	SER	7.4
2	B	275	ASN	6.8
1	E	656	SER	4.7

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, 95th 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(\AA^2)	Q<0.9
3	ZN	E	1301	1/1	0.98	0.14	99,99,99,99	0
3	ZN	C	1301	1/1	0.98	0.17	83,83,83,83	0
3	ZN	A	1301	1/1	0.99	0.18	86,86,86,86	0

6.5 Other polymers [i](#)

There are no such residues in this entry.