Table S1| Related to Figure 1; Data collection and Model Refinement

Data collection	
Particles	426,107
Pixel size (Å)	1.00
Defocus range	-1.4 to -2.4
Voltage (kV)	300
Electron dose (e ⁻ /A ²)	45
Refinement	
CC map_model	0.735
Model content	
Total number of atoms	10,138
No. of protein chains	1
No. of sugar modifications	8
Model quality	
RMSD	
Bond lengths (Å) / Bond angles (°)	0.01/0.9
Ramachandran plot statistics	
Most favored (%)	90.25
Outliers (%)	0.15
Rotamer outliers (%)	0.81
C-beta deviations	0
Clashescore	19.75
B_iso mean	45.07

Table S2 | Related to Figures 1-2; Structural homologs of TEN2 domains

Domain structure (description)

Monobody (5dc4, light Furin-P-domain (1p8j, TcA (409y, receptor ECR2 (Ig-like domain) chain Immunoglobulin) catalytic site stabilization binding domain D in of peptidase activity) ABC toxins) ECR3 (β-propeller) Monooxigenase (3fvz, Glycine Olfactomedin-domain TcC (409x, TcA binding alpha-amidating lyase) (2rmf, latrophilin3 ECR) domain in ABC toxins) (Yen C) ECR4 (YD-barrel) TOM40(5080, Translocase of the Outer Mitochondrial membrane) Tcd B (Yen B) TcB-C/Yenb-C (409x/4igl, ABC toxin encapsulation core) **HNH** endonuclease (5h0m, (tox-HNH), DNase **Barrel** exit point Furin cleavage site ECR5 (TOX-GHH)

Structural homologs (PDB ID, function)

TEN2 domains are presented to the left. Domain names and numbers are indicated below the structure. Structural homologs are highlighted to the right. ABC toxin homologs are indicated by a yellow background. Homologs with protease activity, suspected to be involved in TCAP processing, in blue. Additional homologs in green. PDB ID and protein functions are indicated in parenthesis.