



# Full wwPDB NMR Structure Validation Report ⓘ

Sep 17, 2024 – 02:03 PM EDT

PDB ID : 1PCO  
Title : SOLUTION STRUCTURE OF PORCINE PANCREATIC PROCOLIPASE  
AS DETERMINED FROM 1H HOMONUCLEAR TWO-AND THREE-  
DIMENSIONAL NMR  
Authors : Breg, J.N.; Sarda, L.; Cozzone, P.J.; Rugani, N.; Boelens, R.; Kaptein, R.  
Deposited on : 1994-06-08

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We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

A user guide is available at

<https://www.wwpdb.org/validation/2017/NMRValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4.02b-467
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
wwPDB-RCI	:	v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV	:	Wang et al. (2010)
wwPDB-ShiftChecker	:	v1.2
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.38.3

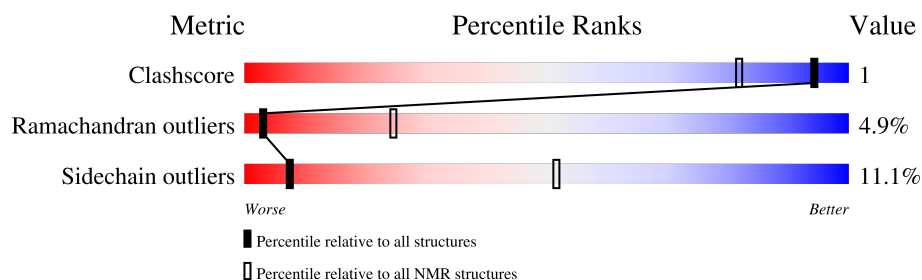
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*SOLUTION NMR*

The overall completeness of chemical shifts assignment was not calculated.

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)	NMR archive (#Entries)
Clashscore	210492	14027
Ramachandran outliers	207382	12486
Sidechain outliers	206894	12463

The table below summarises the geometric issues observed across the polymeric chains and their fit to the experimental data. The red, orange, yellow and green segments indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. A cyan segment indicates the fraction of residues that are not part of the well-defined cores, and 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  $\leq 5\%$

Mol	Chain	Length	Quality of chain
1	A	93	<div> <div>48%</div> <div>25%</div> <div>27%</div> </div>

## 2 Ensemble composition and analysis

This entry contains 25 models. Model 6 is the overall representative, medoid model (most similar to other models).

The following residues are included in the computation of the global validation metrics.

Well-defined (core) protein residues			
Well-defined core	Residue range (total)	Backbone RMSD (Å)	Medoid model
1	A:10-A:70, A:84-A:90 (68)	1.25	6

Ill-defined regions of proteins are excluded from the global statistics.

Ligands and non-protein polymers are included in the analysis.

NmrClust was unable to cluster the ensemble.

Error message: Inconsistent models

### 3 Entry composition [i](#)

There is only 1 type of molecule in this entry. The entry contains 700 atoms, of which 0 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called PORCINE PANCREATIC PROCOLIPASE B.

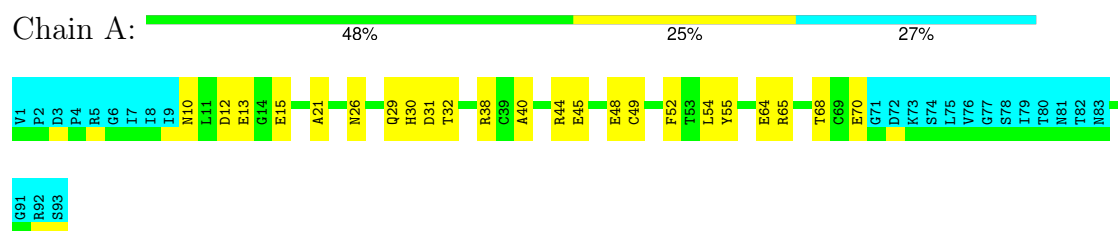
Mol	Chain	Residues	Atoms					Trace
1	A	93	Total	C	N	O	S	0
			700	425	125	140	10	

## 4 Residue-property plots

### 4.1 Average score per residue in the NMR ensemble

These plots are provided for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic is the same as shown in the summary in section 1 of this report. The second graphic shows the sequence where residues are colour-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. Stretches of 2 or more consecutive residues without any outliers are shown as green connectors. Residues which are classified as ill-defined in the NMR ensemble, are shown in cyan with an underline colour-coded according to the previous scheme. Residues which were present in the experimental sample, but not modelled in the final structure are shown in grey.

- Molecule 1: PORCINE PANCREATIC PROCOLIPASE B

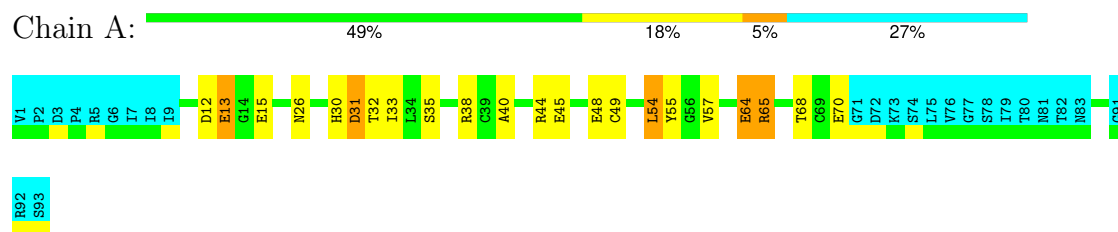


### 4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

#### 4.2.1 Score per residue for model 1

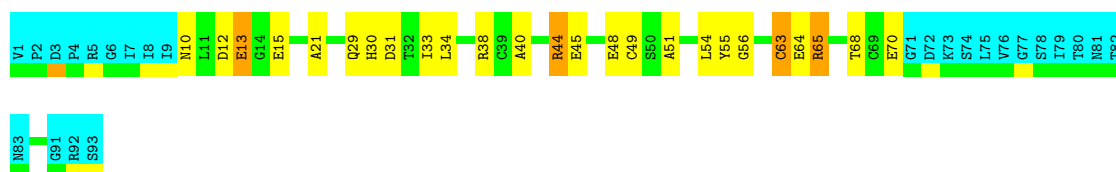
- Molecule 1: PORCINE PANCREATIC PROCOLIPASE B



#### 4.2.2 Score per residue for model 2

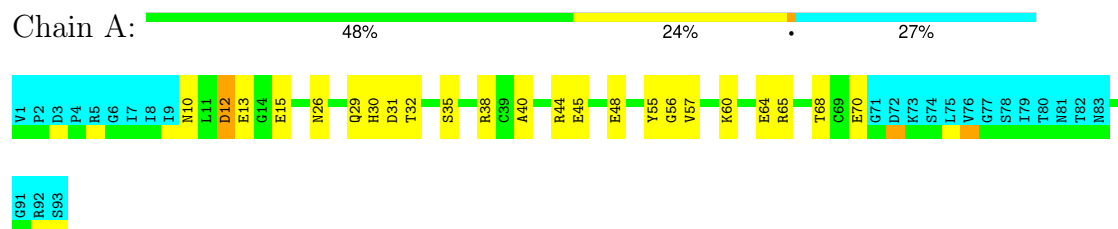
- Molecule 1: PORCINE PANCREATIC PROCOLIPASE B





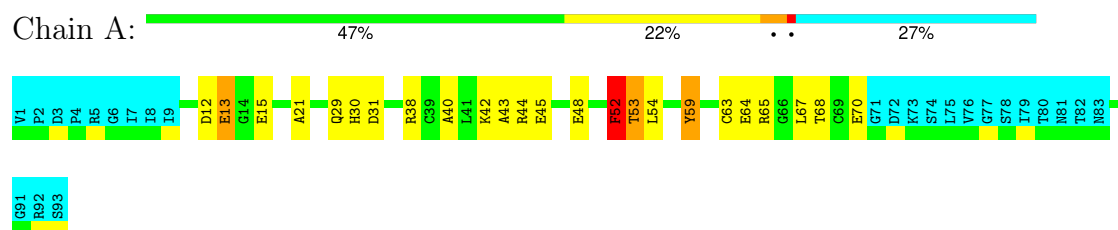
#### 4.2.3 Score per residue for model 3

- Molecule 1: PORCINE PANCREATIC PROCOLIPASE B



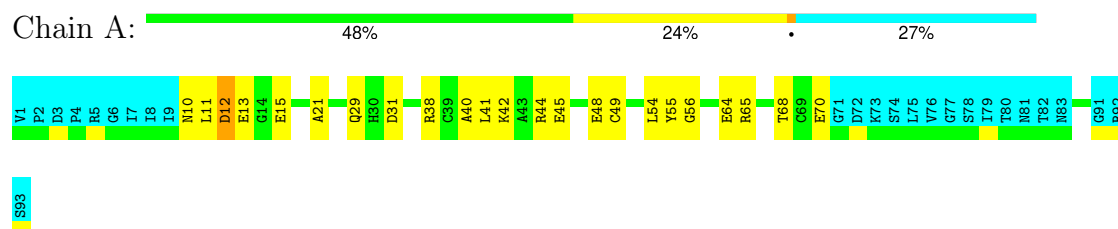
#### 4.2.4 Score per residue for model 4

- Molecule 1: PORCINE PANCREATIC PROCOLIPASE B



#### 4.2.5 Score per residue for model 5

- Molecule 1: PORCINE PANCREATIC PROCOLIPASE B



#### 4.2.6 Score per residue for model 6 (medoid)

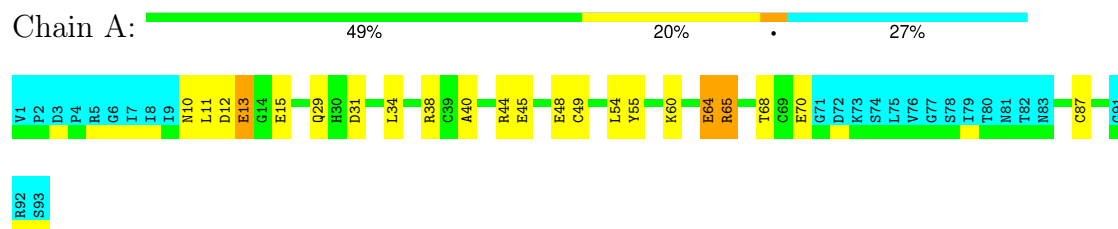
- Molecule 1: PORCINE PANCREATIC PROCOLIPASE B





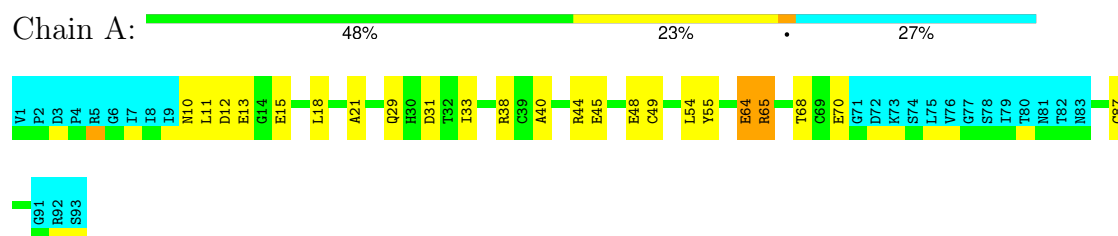
#### 4.2.7 Score per residue for model 7

- Molecule 1: PORCINE PANCREATIC PROCOLIPASE B



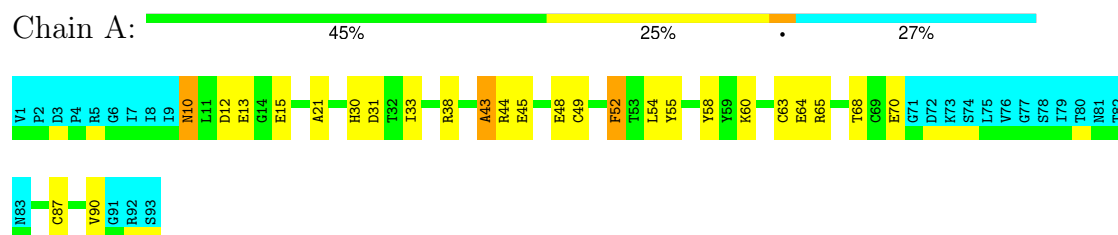
#### 4.2.8 Score per residue for model 8

- Molecule 1: PORCINE PANCREATIC PROCOLIPASE B



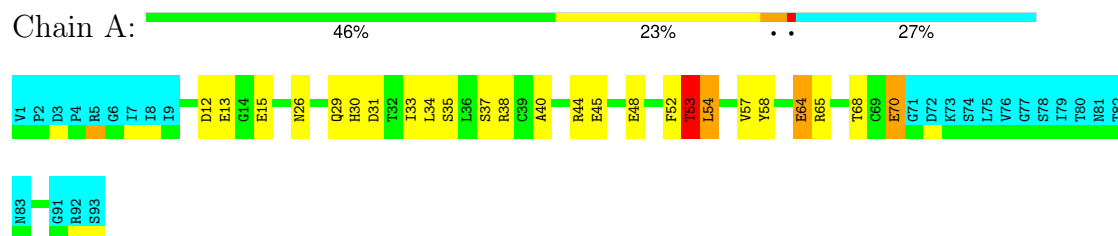
#### 4.2.9 Score per residue for model 9

- Molecule 1: PORCINE PANCREATIC PROCOLIPASE B



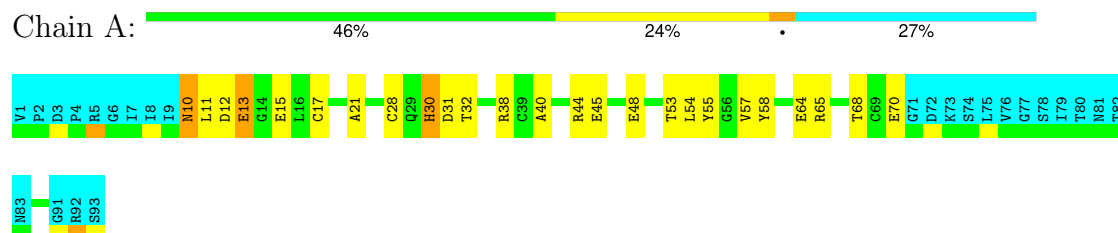
#### 4.2.10 Score per residue for model 10

- Molecule 1: PORCINE PANCREATIC PROCOLIPASE B



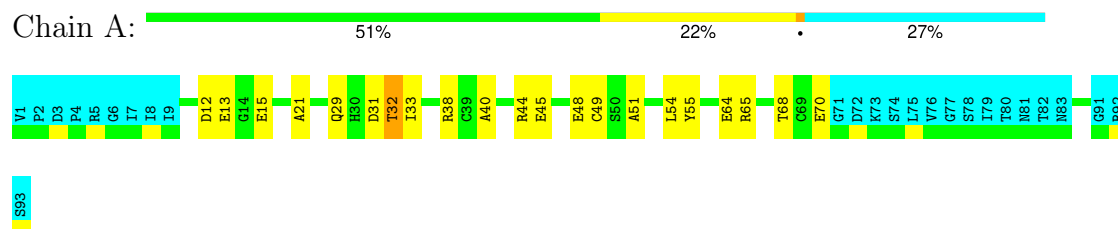
#### 4.2.11 Score per residue for model 11

- Molecule 1: PORCINE PANCREATIC PROCOLIPASE B



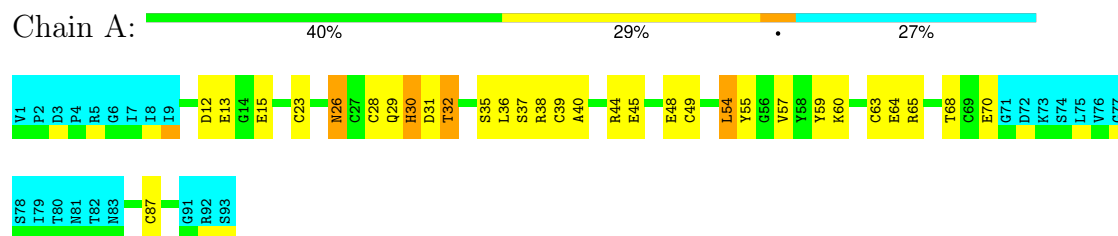
#### 4.2.12 Score per residue for model 12

- Molecule 1: PORCINE PANCREATIC PROCOLIPASE B



#### 4.2.13 Score per residue for model 13

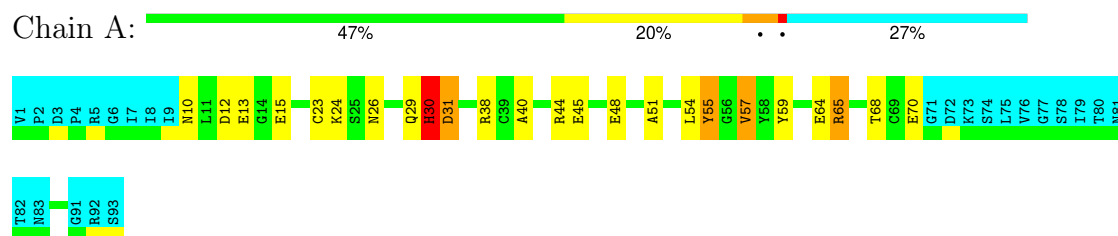
- Molecule 1: PORCINE PANCREATIC PROCOLIPASE B





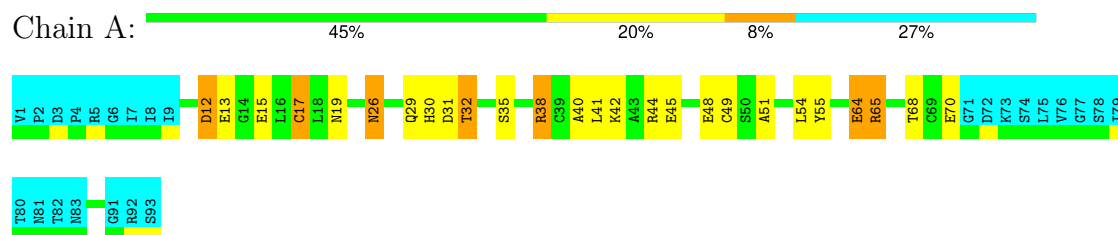
#### 4.2.14 Score per residue for model 14

- Molecule 1: PORCINE PANCREATIC PROCOLIPASE B



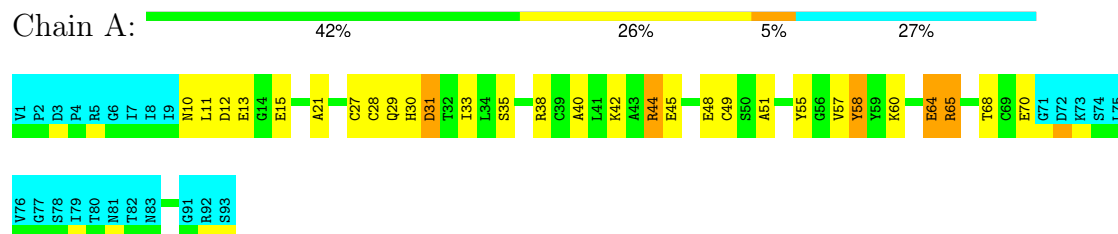
#### 4.2.15 Score per residue for model 15

- Molecule 1: PORCINE PANCREATIC PROCOLIPASE B



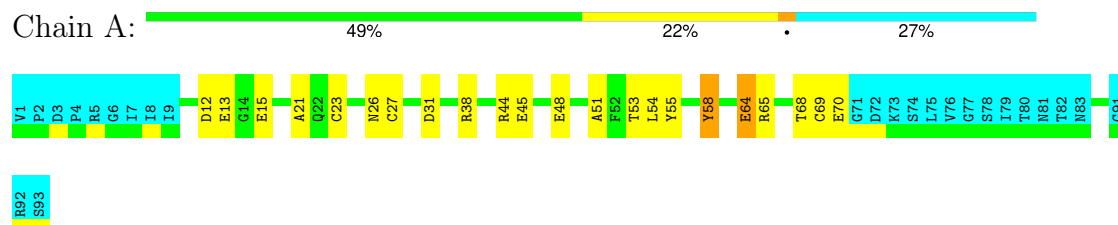
#### 4.2.16 Score per residue for model 16

- Molecule 1: PORCINE PANCREATIC PROCOLIPASE B



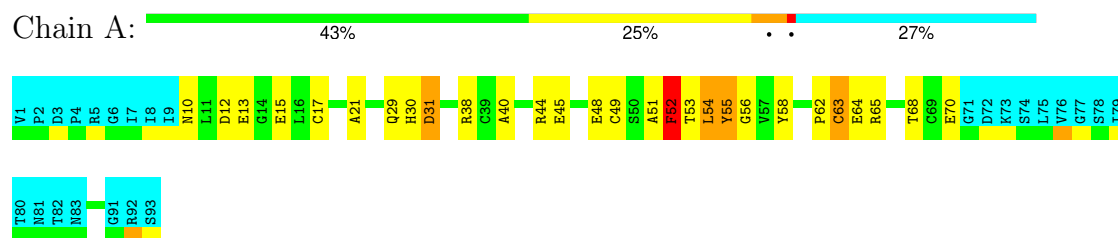
#### 4.2.17 Score per residue for model 17

- Molecule 1: PORCINE PANCREATIC PROCOLIPASE B



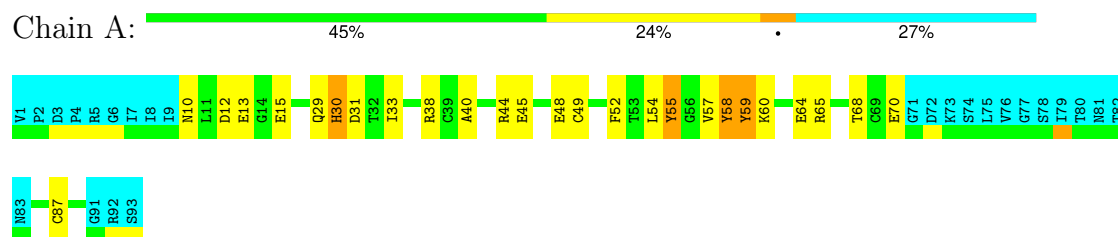
## 4.2.18 Score per residue for model 18

- Molecule 1: PORCINE PANCREATIC PROCOLIPASE B



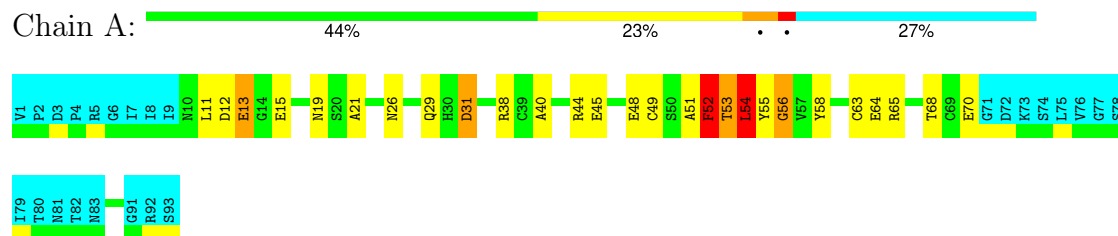
## 4.2.19 Score per residue for model 19

- Molecule 1: PORCINE PANCREATIC PROCOLIPASE B



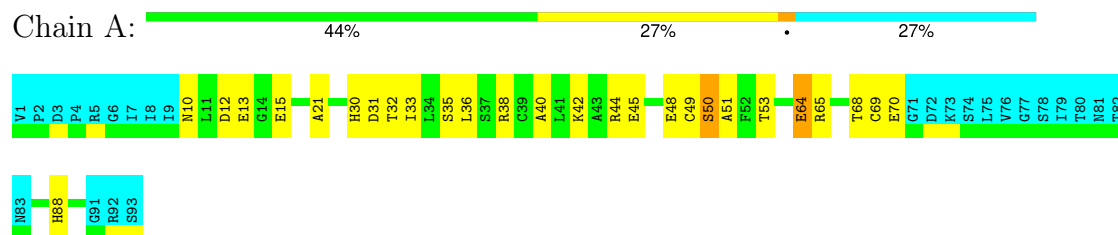
## 4.2.20 Score per residue for model 20

- Molecule 1: PORCINE PANCREATIC PROCOLIPASE B



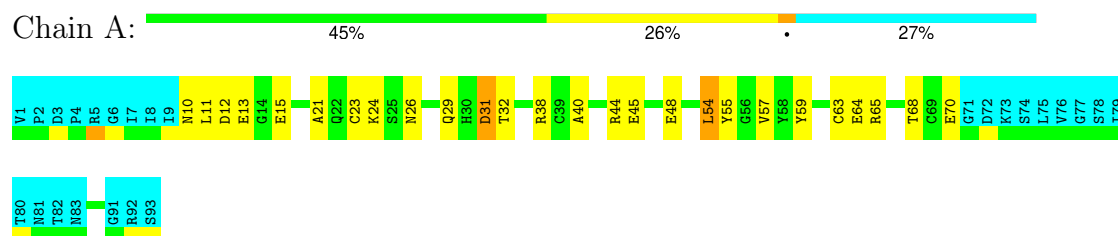
## 4.2.21 Score per residue for model 21

- Molecule 1: PORCINE PANCREATIC PROCOLIPASE B



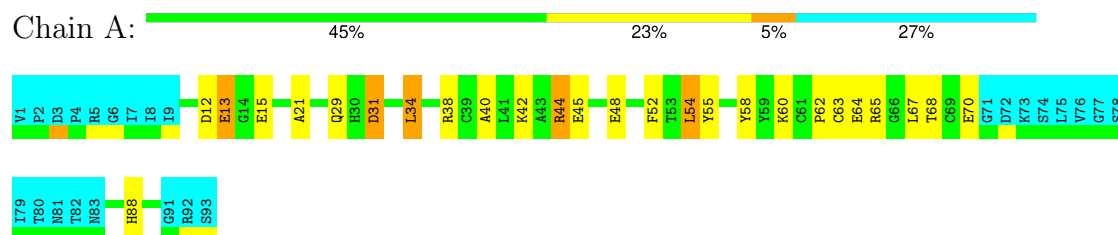
#### 4.2.22 Score per residue for model 22

- Molecule 1: PORCINE PANCREATIC PROCOLIPASE B



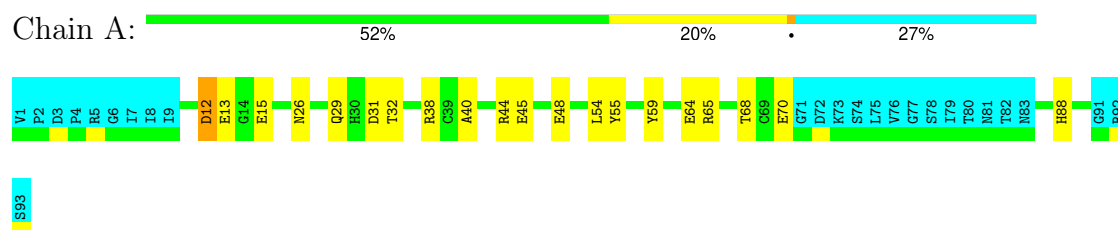
#### 4.2.23 Score per residue for model 23

- Molecule 1: PORCINE PANCREATIC PROCOLIPASE B



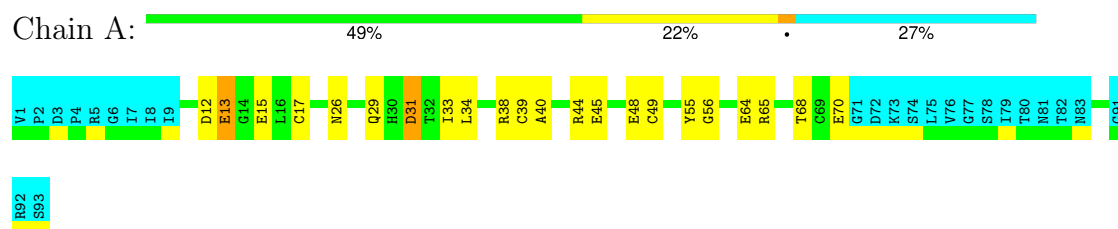
#### 4.2.24 Score per residue for model 24

- Molecule 1: PORCINE PANCREATIC PROCOLIPASE B



#### 4.2.25 Score per residue for model 25

- Molecule 1: PORCINE PANCREATIC PROCOLIPASE B



## 5 Refinement protocol and experimental data overview ⓘ

Of the ? calculated structures, 25 were deposited, based on the following criterion: ?.

The authors did not provide any information on software used for structure solution, optimization or refinement.

No chemical shift data was provided.

## 6 Model quality

### 6.1 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 (average) 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	1.61±0.01	6±1/528 ( 1.2± 0.1%)	1.70±0.05	13±3/712 ( 1.8± 0.4%)
All	All	1.61	159/13200 ( 1.2%)	1.70	323/17800 ( 1.8%)

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	Planarity
1	A	0.0±0.0	1.6±1.0
All	All	0	41

All unique bond outliers are listed below. They are sorted according to the Z-score of the worst occurrence in the ensemble.

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)	Models	
								Worst	Total
1	A	70	GLU	CD-OE2	10.45	1.37	1.25	20	11
1	A	70	GLU	CD-OE1	10.28	1.36	1.25	4	14
1	A	13	GLU	CD-OE1	10.26	1.36	1.25	7	7
1	A	64	GLU	CD-OE1	10.24	1.36	1.25	9	16
1	A	48	GLU	CD-OE2	10.12	1.36	1.25	8	11
1	A	13	GLU	CD-OE2	10.04	1.36	1.25	5	18
1	A	15	GLU	CD-OE1	10.01	1.36	1.25	23	7
1	A	48	GLU	CD-OE1	9.96	1.36	1.25	11	14
1	A	45	GLU	CD-OE2	9.93	1.36	1.25	12	12
1	A	15	GLU	CD-OE2	9.90	1.36	1.25	6	18
1	A	45	GLU	CD-OE1	9.86	1.36	1.25	2	13
1	A	64	GLU	CD-OE2	9.80	1.36	1.25	25	9
1	A	31	ASP	CG-OD1	5.11	1.37	1.25	23	1
1	A	12	ASP	CG-OD2	5.08	1.37	1.25	1	5
1	A	12	ASP	CG-OD1	5.08	1.37	1.25	8	3

All unique angle outliers are listed below. They are sorted according to the Z-score of the worst occurrence in the ensemble.

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
1	A	30	HIS	N-CA-C	11.79	142.84	111.00	14	2
1	A	65	ARG	NE-CZ-NH1	8.86	124.73	120.30	22	25
1	A	44	ARG	NE-CZ-NH1	8.82	124.71	120.30	17	25
1	A	38	ARG	NE-CZ-NH1	8.73	124.67	120.30	20	25
1	A	21	ALA	N-CA-CB	-8.72	97.89	110.10	20	14
1	A	54	LEU	CB-CA-C	8.44	126.23	110.20	20	10
1	A	31	ASP	CB-CG-OD1	-8.14	110.97	118.30	8	24
1	A	30	HIS	N-CA-CB	-8.03	96.14	110.60	14	2
1	A	31	ASP	CB-CG-OD2	-7.49	111.56	118.30	15	22
1	A	58	TYR	CB-CG-CD2	-7.34	116.60	121.00	16	6
1	A	12	ASP	CB-CG-OD2	-7.05	111.95	118.30	1	24
1	A	12	ASP	CB-CG-OD1	-6.96	112.04	118.30	8	24
1	A	51	ALA	N-CA-CB	-6.85	100.50	110.10	14	5
1	A	30	HIS	CA-C-N	6.56	131.63	117.20	14	1
1	A	40	ALA	N-CA-CB	-6.54	100.94	110.10	11	23
1	A	57	VAL	CG1-CB-CG2	-6.53	100.46	110.90	14	5
1	A	59	TYR	CB-CG-CD2	-6.21	117.27	121.00	19	2
1	A	38	ARG	NE-CZ-NH2	-6.09	117.25	120.30	22	22
1	A	35	SER	N-CA-CB	-6.06	101.40	110.50	13	1
1	A	23	CYS	CA-CB-SG	-6.06	103.09	114.00	13	1
1	A	15	GLU	N-CA-CB	-6.00	99.80	110.60	5	1
1	A	31	ASP	CA-CB-CG	5.97	126.53	113.40	14	1
1	A	32	THR	N-CA-CB	5.88	121.48	110.30	12	1
1	A	53	THR	O-C-N	-5.84	113.35	122.70	10	1
1	A	44	ARG	NE-CZ-NH2	-5.80	117.40	120.30	16	11
1	A	55	TYR	CB-CG-CD1	-5.80	117.52	121.00	14	1
1	A	40	ALA	CB-CA-C	5.73	118.70	110.10	14	1
1	A	65	ARG	NE-CZ-NH2	-5.56	117.52	120.30	11	10
1	A	27	CYS	CA-CB-SG	-5.55	104.01	114.00	16	2
1	A	53	THR	N-CA-CB	5.48	120.72	110.30	20	1
1	A	52	PHE	CB-CG-CD1	-5.47	116.97	120.80	20	3
1	A	43	ALA	N-CA-CB	-5.41	102.53	110.10	9	1
1	A	64	GLU	N-CA-C	5.40	125.57	111.00	23	1
1	A	28	CYS	CA-CB-SG	-5.39	104.30	114.00	11	2
1	A	57	VAL	CA-CB-CG2	5.29	118.84	110.90	14	1
1	A	30	HIS	CG-ND1-CE1	-5.29	98.83	105.70	11	7
1	A	67	LEU	CB-CA-C	5.28	120.24	110.20	23	1
1	A	23	CYS	CB-CA-C	5.27	120.95	110.40	13	1
1	A	52	PHE	N-CA-C	5.22	125.11	111.00	18	1
1	A	31	ASP	N-CA-C	5.19	125.00	111.00	14	1

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
1	A	11	LEU	CB-CA-C	5.17	120.03	110.20	22	2
1	A	53	THR	N-CA-C	-5.17	97.04	111.00	10	1
1	A	54	LEU	N-CA-CB	5.14	120.67	110.40	10	1
1	A	52	PHE	CB-CG-CD2	-5.06	117.26	120.80	20	1
1	A	88	HIS	CG-ND1-CE1	-5.06	99.12	105.70	21	4
1	A	30	HIS	ND1-CE1-NE2	5.06	121.03	109.90	11	1
1	A	23	CYS	N-CA-CB	-5.06	101.50	110.60	13	1

There are no chirality outliers.

All unique planar outliers are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type	Group	Models (Total)
1	A	55	TYR	Sidechain	18
1	A	58	TYR	Sidechain	6
1	A	52	PHE	Sidechain	5
1	A	56	GLY	Peptide	4
1	A	59	TYR	Sidechain	4
1	A	53	THR	Mainchain	2
1	A	30	HIS	Sidechain	2

## 6.2 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 each 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 averaged over the ensemble.

Mol	Chain	Non-H	H(model)	H(added)	Clashes
1	A	520	0	485	1±1
All	All	13000	0	12125	27

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 1.

All unique clashes are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:49:CYS:SG	1:A:87:CYS:SG	0.67	2.92	9	2
1:A:30:HIS:HA	1:A:57:VAL:CG1	0.61	2.25	14	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:17:CYS:SG	1:A:19:ASN:O	0.54	2.65	15	1
1:A:54:LEU:HD23	1:A:56:GLY:H	0.53	1.64	20	1
1:A:49:CYS:HB3	1:A:87:CYS:SG	0.52	2.44	7	3
1:A:30:HIS:HA	1:A:57:VAL:HG12	0.50	1.83	14	1
1:A:37:SER:HB3	1:A:57:VAL:HG11	0.50	1.83	10	2
1:A:49:CYS:HB2	1:A:87:CYS:SG	0.50	2.46	19	1
1:A:30:HIS:HB2	1:A:57:VAL:HG12	0.49	1.83	10	5
1:A:49:CYS:SG	1:A:63:CYS:SG	0.48	3.12	18	3
1:A:51:ALA:HB2	1:A:69:CYS:SG	0.46	2.50	17	1
1:A:17:CYS:SG	1:A:39:CYS:SG	0.45	3.14	25	1
1:A:49:CYS:SG	1:A:62:PRO:HA	0.44	2.52	18	1
1:A:50:SER:O	1:A:69:CYS:SG	0.43	2.77	21	1
1:A:28:CYS:HA	1:A:39:CYS:SG	0.43	2.53	13	1
1:A:43:ALA:HB3	1:A:63:CYS:SG	0.41	2.56	9	2

## 6.3 Torsion angles [i](#)

### 6.3.1 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 PDB entries followed by that with respect to all NMR entries. 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	68/93 (73%)	53±3 (78±4%)	11±3 (17±4%)	3±2 (5±2%)	<b>3</b>	24
All	All	1700/2325 (73%)	1330 (78%)	286 (17%)	84 (5%)	<b>3</b>	24

All 20 unique Ramachandran outliers are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type	Models (Total)
1	A	32	THR	9
1	A	33	ILE	9
1	A	10	ASN	9
1	A	26	ASN	8
1	A	53	THR	7
1	A	35	SER	6
1	A	34	LEU	6
1	A	63	CYS	5

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Mol	Chain	Res	Type	Models (Total)
1	A	31	ASP	5
1	A	55	TYR	3
1	A	23	CYS	3
1	A	51	ALA	3
1	A	56	GLY	2
1	A	54	LEU	2
1	A	36	LEU	2
1	A	90	VAL	1
1	A	24	LYS	1
1	A	30	HIS	1
1	A	50	SER	1
1	A	62	PRO	1

### 6.3.2 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 PDB entries followed by that with respect to all NMR entries. 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	60/81 (74%)	53±2 (89±4%)	7±2 (11±4%)	7	51
All	All	1500/2025 (74%)	1334 (89%)	166 (11%)	7	51

All 31 unique residues with a non-rotameric sidechain are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type	Models (Total)
1	A	68	THR	25
1	A	29	GLN	20
1	A	54	LEU	15
1	A	13	GLU	9
1	A	49	CYS	8
1	A	64	GLU	8
1	A	60	LYS	8
1	A	65	ARG	7
1	A	42	LYS	7
1	A	10	ASN	6
1	A	52	PHE	6
1	A	26	ASN	5
1	A	30	HIS	5

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Mol	Chain	Res	Type	Models (Total)
1	A	11	LEU	5
1	A	12	ASP	4
1	A	31	ASP	3
1	A	44	ARG	3
1	A	17	CYS	3
1	A	32	THR	3
1	A	55	TYR	3
1	A	59	TYR	2
1	A	41	LEU	2
1	A	67	LEU	1
1	A	18	LEU	1
1	A	70	GLU	1
1	A	38	ARG	1
1	A	33	ILE	1
1	A	19	ASN	1
1	A	24	LYS	1
1	A	34	LEU	1
1	A	63	CYS	1

### 6.3.3 RNA [i](#)

There are no RNA molecules in this entry.

## 6.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

### 6.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

### 6.6 Ligand geometry [i](#)

There are no ligands in this entry.

### 6.7 Other polymers [i](#)

There are no such molecules in this entry.

## 6.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

## 7 Chemical shift validation

No chemical shift data were provided