



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 6, 2026 – 08:30 AM UTC

PDB ID : 2ZWP / pdb_00002zwp
Title : Crystal structure of Ca3 site mutant of Pro-S324A
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Deposited on : 2008-12-17
Resolution : 2.40 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 types of validation reports are described at

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

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

MolProbity : 4-5-2 with Phenix2.0
Xtrriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

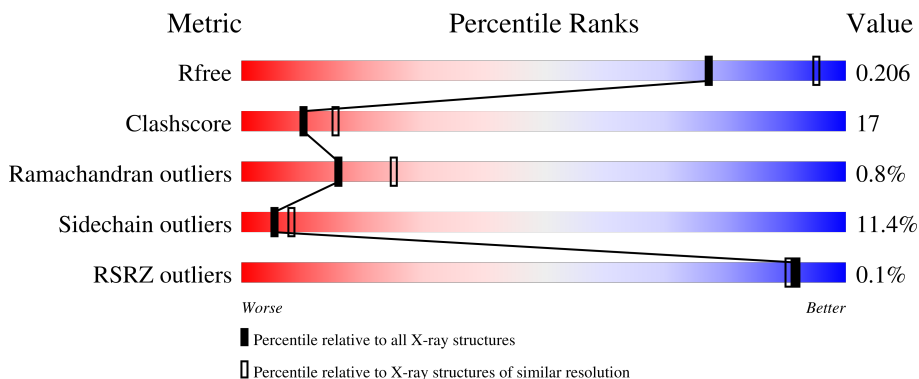
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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 $\leq 5\%$. 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	398	 64% 26% 5% . .
1	B	398	 66% 24% 6% . .

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6103 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 Tk-subtilisin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	389	2853	1811	476	559	7	0	0	0
1	B	383	2821	1790	471	553	7	0	0	0

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	225	ALA	ASP	engineered mutation	UNP P58502
A	324	ALA	SER	engineered mutation	UNP P58502
B	225	ALA	ASP	engineered mutation	UNP P58502
B	324	ALA	SER	engineered mutation	UNP P58502

- Molecule 2 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	6	Total 6	Ca 6	0	0
2	B	5	Total 5	Ca 5	0	0

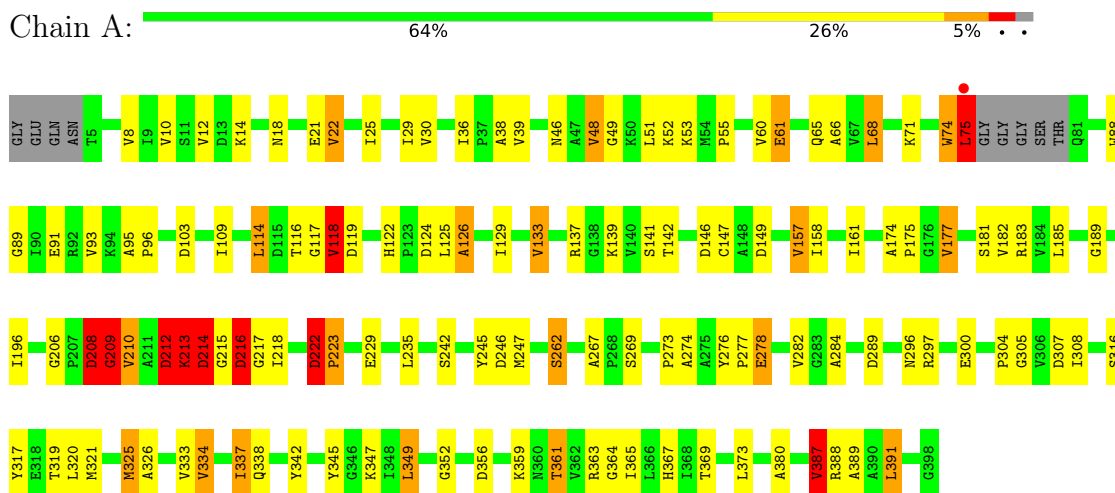
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	218	Total 218	O 218	0	0
3	B	200	Total 200	O 200	0	0

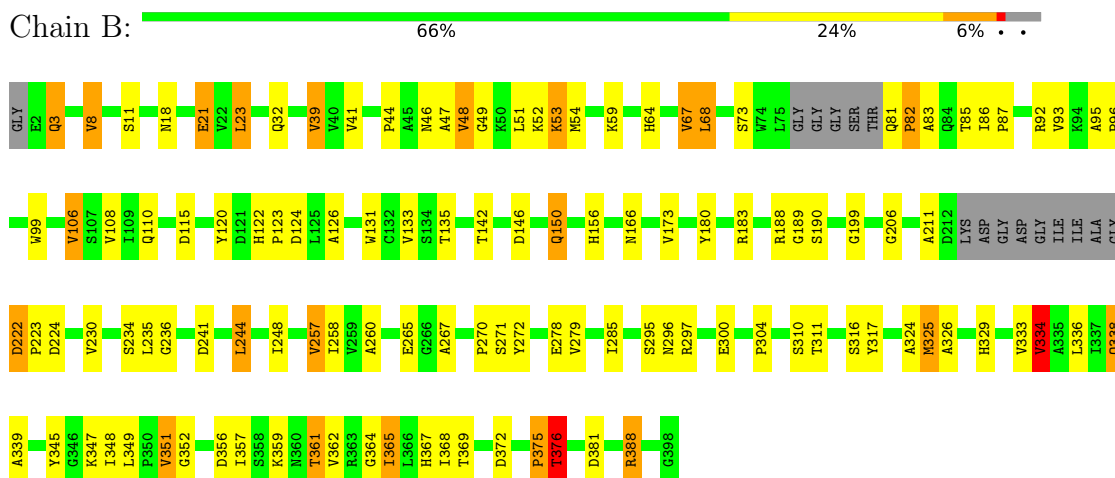
3 Residue-property plots [i](#)

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: Tk-subtilisin



- Molecule 1: Tk-subtilisin



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	50.58Å 87.74Å 155.49Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	43.85 – 2.40 43.85 – 2.40	Depositor EDS
% Data completeness (in resolution range)	94.3 (43.85-2.40) 94.2 (43.85-2.40)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.80 (at 2.39Å)	Xtrriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.157 , 0.202 0.157 , 0.206	Depositor DCC
R_{free} test set	1327 reflections (4.76%)	wwPDB-VP
Wilson B-factor (Å ²)	27.7	Xtrriage
Anisotropy	0.097	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 49.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6103	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.38% 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 [i](#)

5.1 Standard geometry [i](#)

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

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	1.27	6/2913 (0.2%)	1.40	28/3986 (0.7%)
1	B	1.28	10/2880 (0.3%)	1.36	23/3941 (0.6%)
All	All	1.28	16/5793 (0.3%)	1.38	51/7927 (0.6%)

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
1	A	1	5
1	B	0	2
All	All	1	7

The worst 5 of 16 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	48	VAL	CA-CB	6.72	1.63	1.54
1	A	214	ASP	N-CA	6.55	1.54	1.46
1	B	348	ILE	CA-CB	6.05	1.64	1.54
1	A	177	VAL	CA-CB	6.03	1.61	1.54
1	B	351	VAL	CA-CB	5.88	1.61	1.54

The worst 5 of 51 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	222	ASP	CA-C-N	16.45	140.40	119.84
1	B	222	ASP	C-N-CA	16.45	140.40	119.84
1	A	213	LYS	N-CA-C	-15.01	94.94	114.31
1	A	223	PRO	N-CA-C	-10.79	99.19	114.18

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	74	TRP	N-CA-C	10.05	123.29	111.02

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	216	ASP	CA

5 of 7 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	208	ASP	Peptide
1	A	209	GLY	Peptide
1	A	212	ASP	Peptide
1	A	213	LYS	Peptide
1	A	222	ASP	Peptide

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	A	2853	0	2817	112	0
1	B	2821	0	2779	81	0
2	A	6	0	0	0	0
2	B	5	0	0	0	0
3	A	218	0	0	1	0
3	B	200	0	0	12	0
All	All	6103	0	5596	193	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 17.

The worst 5 of 193 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:A:320:LEU:CB	1:A:325:MET:HE3	1.73	1.19
1:A:320:LEU:HB2	1:A:325:MET:HE3	1.14	1.14

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:361:THR:HG21	3:B:562:HOH:O	1.54	1.07
1:A:215:GLY:CA	1:A:216:ASP:HB3	1.86	1.04
1:A:361:THR:HG22	1:A:364:GLY:H	1.23	1.02

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	385/398 (97%)	360 (94%)	21 (6%)	4 (1%)	12	20
1	B	377/398 (95%)	348 (92%)	27 (7%)	2 (0%)	24	37
All	All	762/796 (96%)	708 (93%)	48 (6%)	6 (1%)	16	25

5 of 6 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	214	ASP
1	A	210	VAL
1	A	216	ASP
1	B	82	PRO
1	A	119	ASP

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	Rotameric	Outliers	Percentiles	
1	A	300/306 (98%)	266 (89%)	34 (11%)	5	8
1	B	298/306 (97%)	264 (89%)	34 (11%)	5	8
All	All	598/612 (98%)	530 (89%)	68 (11%)	5	8

5 of 68 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	325	MET
1	B	338	GLN
1	B	375	PRO
1	A	246	ASP
1	A	218	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 18 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	296	ASN
1	B	367	HIS
1	B	338	GLN
1	B	18	ASN
1	B	290	ASN

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 11 ligands modelled in this entry, 11 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

There are no such residues in this entry.

5.8 Polymer linkage issues

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	389/398 (97%)	-0.61	1 (0%) 90 88	12, 21, 39, 49	0
1	B	383/398 (96%)	-0.57	0 100 100	13, 22, 37, 49	0
All	All	772/796 (96%)	-0.59	1 (0%) 92 90	12, 22, 39, 49	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	75	LEU	2.0

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 oligosaccharides 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(Å ²)	Q<0.9
2	CA	A	401	1/1	0.97	0.04	26,26,26,26	0
2	CA	A	403	1/1	0.98	0.03	49,49,49,49	0
2	CA	B	399	1/1	0.98	0.03	29,29,29,29	0
2	CA	B	400	1/1	0.98	0.03	33,33,33,33	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	CA	B	401	1/1	0.98	0.05	29,29,29,29	0
2	CA	B	402	1/1	0.98	0.02	17,17,17,17	0
2	CA	A	399	1/1	0.99	0.03	21,21,21,21	0
2	CA	A	402	1/1	0.99	0.02	31,31,31,31	0
2	CA	A	400	1/1	0.99	0.04	24,24,24,24	0
2	CA	A	404	1/1	0.99	0.02	28,28,28,28	0
2	CA	B	403	1/1	0.99	0.01	19,19,19,19	0

6.5 Other polymers [\(i\)](#)

There are no such residues in this entry.