



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 05:19 AM UTC

PDB ID : 5IIG / pdb_00005iig
Title : Structure of the SPX-TTM domain fragment of the yeast inorganic polyphosphate polymerase Vtc4 (form A).
Authors : Wild, R.; Hothorn, M.
Deposited on : 2016-03-01
Resolution : 2.99 Å(reported)

This is a Full wwPDB X-ray Structure Validation 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
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (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

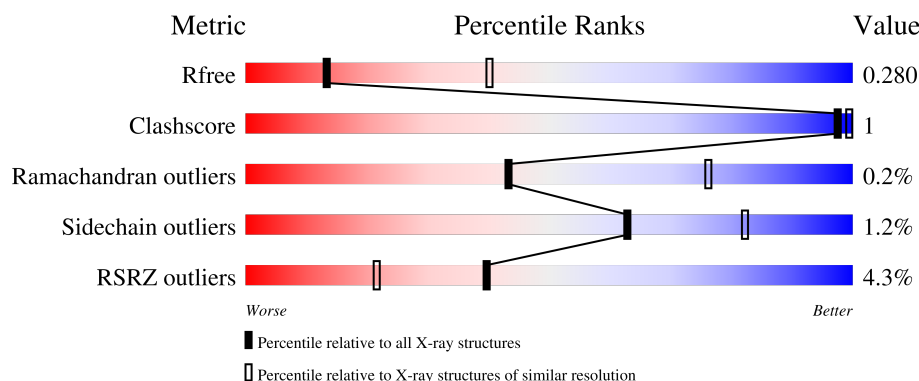
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.99 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	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

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	485	4% 79% 16% • •

2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 3871 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.

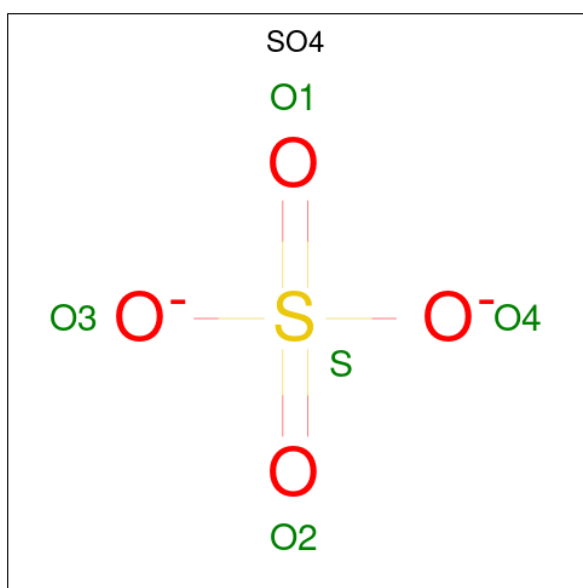
- Molecule 1 is a protein called Vacuolar transporter chaperone 4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	467	Total	C	N	O	S	0	0	0
			3851	2466	664	715	6			

There are 7 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	expression tag	UNP P47075
A	-1	ALA	-	expression tag	UNP P47075
A	0	MET	-	expression tag	UNP P47075
A	1	GLY	-	expression tag	UNP P47075
A	426	ASN	GLU	engineered mutation	UNP P47075
A	481	GLY	-	expression tag	UNP P47075
A	482	SER	-	expression tag	UNP P47075

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).

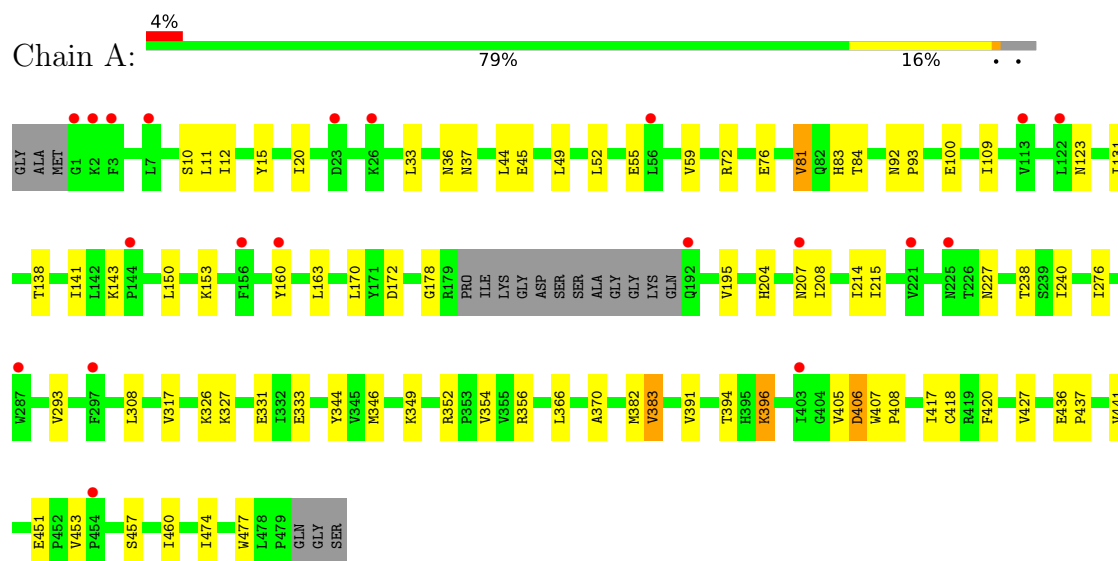


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		

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: Vacuolar transporter chaperone 4



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 2 2	Depositor
Cell constants a, b, c, α , β , γ	145.39Å 145.39Å 71.89Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.22 – 2.99 48.22 – 2.99	Depositor EDS
% Data completeness (in resolution range)	99.8 (48.22-2.99) 99.8 (48.22-2.99)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.58 (at 3.01Å)	Xtriage
Refinement program	BUSTER 2.10.1	Depositor
R, R_{free}	0.244 , 0.273 0.252 , 0.280	Depositor DCC
R_{free} test set	797 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å ²)	105.3	Xtriage
Anisotropy	0.087	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 68.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	3871	wwPDB-VP
Average B, all atoms (Å ²)	113.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.35% 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: SO4

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	1.96	79/3938 (2.0%)	1.19	0/5323

All (79) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	153	LYS	CA-C	8.84	1.64	1.52
1	A	405	VAL	CA-C	7.98	1.62	1.52
1	A	143	LYS	C-N	7.68	1.43	1.34
1	A	352	ARG	C-N	7.53	1.42	1.33
1	A	436	GLU	C-N	7.28	1.42	1.33
1	A	451	GLU	C-N	7.14	1.41	1.33
1	A	352	ARG	CA-C	6.94	1.59	1.52
1	A	474	ILE	C-N	6.67	1.41	1.33
1	A	437	PRO	C-N	6.60	1.42	1.33
1	A	93	PRO	C-N	6.57	1.42	1.33
1	A	12	ILE	CA-CB	6.57	1.60	1.53
1	A	240	ILE	CA-CB	6.44	1.61	1.53
1	A	227	ASN	CA-C	6.39	1.61	1.52
1	A	143	LYS	CA-C	6.29	1.59	1.52
1	A	453	VAL	C-N	6.28	1.41	1.33
1	A	406	ASP	CA-C	6.15	1.61	1.52
1	A	37	ASN	CA-C	6.08	1.59	1.53
1	A	327	LYS	C-N	6.07	1.41	1.33
1	A	407	TRP	CA-C	6.06	1.60	1.53
1	A	204	HIS	C-N	6.05	1.42	1.33
1	A	92	ASN	C-N	6.00	1.40	1.33
1	A	366	LEU	C-N	5.95	1.40	1.33
1	A	317	VAL	CA-CB	5.94	1.61	1.54
1	A	153	LYS	C-N	5.85	1.40	1.33
1	A	20	ILE	CA-CB	5.81	1.60	1.54
1	A	33	LEU	C-N	5.77	1.41	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	215	ILE	C-N	5.77	1.41	1.33
1	A	405	VAL	CA-CB	5.71	1.61	1.54
1	A	370	ALA	C-N	5.71	1.41	1.33
1	A	36	ASN	C-N	5.69	1.40	1.33
1	A	36	ASN	CA-C	5.67	1.61	1.52
1	A	349	LYS	C-N	5.64	1.41	1.33
1	A	141	ILE	CG1-CD1	5.63	1.73	1.51
1	A	123	ASN	C-N	5.57	1.41	1.33
1	A	474	ILE	CA-C	5.57	1.58	1.52
1	A	141	ILE	CA-CB	5.55	1.60	1.54
1	A	109	ILE	C-N	5.51	1.40	1.33
1	A	131	ILE	CG1-CD1	5.51	1.73	1.51
1	A	195	VAL	CA-CB	5.50	1.61	1.54
1	A	408	PRO	C-N	5.46	1.40	1.33
1	A	420	PHE	C-N	5.46	1.40	1.34
1	A	163	LEU	C-N	5.45	1.40	1.33
1	A	72	ARG	C-N	5.42	1.41	1.33
1	A	208	ILE	CA-CB	5.41	1.60	1.54
1	A	293	VAL	CA-CB	5.40	1.59	1.53
1	A	83	HIS	C-N	5.39	1.41	1.33
1	A	214	ILE	CA-CB	5.39	1.60	1.54
1	A	276	ILE	CA-CB	5.35	1.60	1.54
1	A	45	GLU	C-N	5.34	1.40	1.33
1	A	208	ILE	CA-C	5.32	1.59	1.52
1	A	453	VAL	CA-CB	5.32	1.59	1.54
1	A	10	SER	CA-C	5.32	1.61	1.52
1	A	383	VAL	CA-CB	5.30	1.61	1.54
1	A	207	ASN	C-N	5.27	1.40	1.33
1	A	160	TYR	C-N	5.24	1.41	1.34
1	A	396	LYS	CG-CD	5.20	1.68	1.52
1	A	15	TYR	CA-C	5.19	1.59	1.53
1	A	59	VAL	CA-CB	5.18	1.60	1.54
1	A	76	GLU	C-N	5.17	1.40	1.33
1	A	333	GLU	C-N	5.15	1.41	1.33
1	A	441	VAL	CA-CB	5.14	1.61	1.54
1	A	172	ASP	C-N	5.14	1.40	1.33
1	A	153	LYS	CA-CB	5.13	1.61	1.53
1	A	417	ILE	CA-CB	5.13	1.61	1.54
1	A	427	VAL	CA-CB	5.13	1.59	1.53
1	A	394	THR	C-N	5.12	1.41	1.33
1	A	150	LEU	C-N	5.10	1.40	1.33
1	A	331	GLU	C-N	5.10	1.40	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	354	VAL	CA-CB	5.09	1.59	1.55
1	A	100	GLU	C-N	5.09	1.40	1.33
1	A	477	TRP	CA-C	5.07	1.59	1.52
1	A	344	TYR	C-N	5.07	1.40	1.33
1	A	436	GLU	CA-C	5.06	1.59	1.52
1	A	178	GLY	C-N	5.05	1.40	1.33
1	A	84	THR	C-N	5.04	1.40	1.33
1	A	55	GLU	C-N	5.02	1.40	1.33
1	A	11	LEU	C-N	5.02	1.38	1.33
1	A	81	VAL	C-N	5.01	1.40	1.34
1	A	391	VAL	CA-CB	5.01	1.60	1.53

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	3851	0	3786	6	0
2	A	20	0	0	0	0
All	All	3871	0	3786	6	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 1.

All (6) 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:382:MET:HB2	1:A:418:CYS:HB3	1.94	0.49
1:A:308:LEU:HB3	1:A:346:MET:HE2	2.00	0.43
1:A:49:LEU:HA	1:A:52:LEU:HD12	2.00	0.43
1:A:356:ARG:HB3	1:A:383:VAL:HB	2.00	0.43
1:A:81:VAL:HG21	1:A:170:LEU:HD22	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:457:SER:HB3	1:A:460:ILE:HB	2.04	0.40

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	463/485 (96%)	449 (97%)	13 (3%)	1 (0%)	43 76

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	406	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	419/444 (94%)	414 (99%)	5 (1%)	63 82

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

Mol	Chain	Res	Type
1	A	44	LEU
1	A	138	THR

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Mol	Chain	Res	Type
1	A	238	THR
1	A	326	LYS
1	A	396	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	42	GLN
1	A	147	GLN
1	A	283	HIS
1	A	316	GLN
1	A	334	ASN
1	A	395	HIS
1	A	430	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

4 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	SO4	A	502	-	4,4,4	0.40	0	6,6,6	0.12	0
2	SO4	A	501	-	4,4,4	0.26	0	6,6,6	0.09	0
2	SO4	A	504	-	4,4,4	0.24	0	6,6,6	0.08	0
2	SO4	A	503	-	4,4,4	0.26	0	6,6,6	0.08	0

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	467/485 (96%)	0.41	20 (4%) 40 21	81, 108, 156, 202	0

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1	GLY	4.0
1	A	2	LYS	3.2
1	A	156	PHE	3.1
1	A	7	LEU	2.9
1	A	454	PRO	2.8
1	A	3	PHE	2.8
1	A	297	PHE	2.8
1	A	56	LEU	2.7
1	A	287	TRP	2.4
1	A	144	PRO	2.3
1	A	160	TYR	2.3
1	A	192	GLN	2.2
1	A	23	ASP	2.2
1	A	403	ILE	2.1
1	A	113	VAL	2.1
1	A	221	VAL	2.1
1	A	225	ASN	2.1
1	A	122	LEU	2.1
1	A	26	LYS	2.0
1	A	207	ASN	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($\text{\AA}^2$)	Q<0.9
2	SO4	A	503	5/5	0.75	0.06	177,177,177,177	0
2	SO4	A	504	5/5	0.83	0.09	169,170,170,170	0
2	SO4	A	502	5/5	0.84	0.12	109,109,109,110	0
2	SO4	A	501	5/5	0.86	0.09	140,140,140,140	0

6.5 Other polymers [i](#)

There are no such residues in this entry.