Full wwPDB X-ray Structure Validation Report (i)

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PDB ID	:	1XWI
Title	:	Crystal Structure of VPS4B
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Deposited on	:	2004-11-01
Resolution	:	2.80 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 (i) 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 (i)) were used in the production of this report:

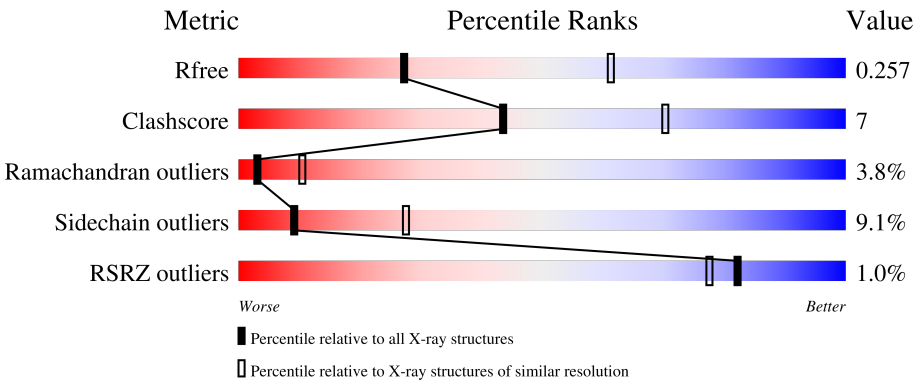
MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	2.36.2
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

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 <=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	322	% 72% 23% • •

2 Entry composition (i)

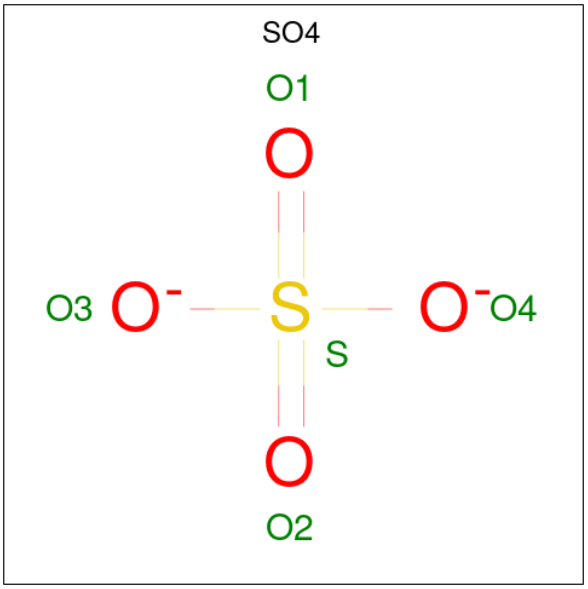
There are 3 unique types of molecules in this entry. The entry contains 2527 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 SKD1 protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	322	2508	1591	438	470	9	158	0	0

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



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

- Molecule 3 is water.

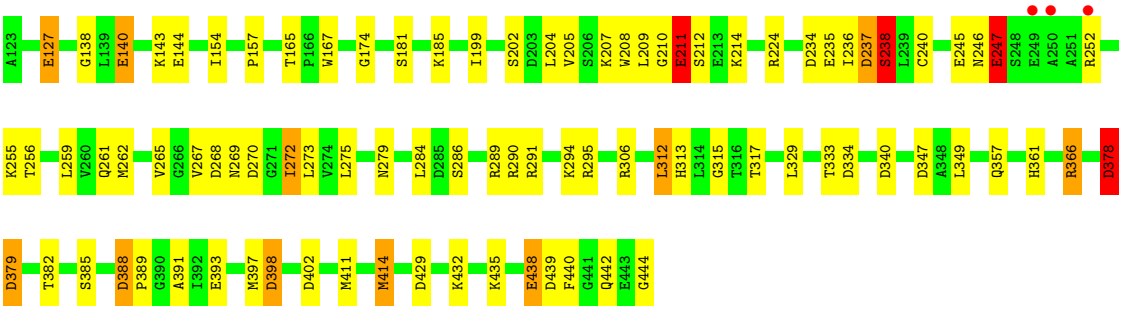
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	A	9	9	9	0	0

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: SKD1 protein

Chain A:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 65	Depositor
Cell constants a, b, c, α, β, γ	88.06Å 88.06Å 112.57Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	76.70 – 2.80 63.14 – 2.80	Depositor EDS
% Data completeness (in resolution range)	98.8 (76.70-2.80) 98.8 (63.14-2.80)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	0.06	Depositor
$< I/\sigma(I) >$ [1]	7.32 (at 2.81Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.192 , 0.260 0.193 , 0.257	Depositor DCC
R_{free} test set	1190 reflections (9.83%)	wwPDB-VP
Wilson B-factor (Å^2)	52.9	Xtriage
Anisotropy	0.013	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 37.1	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.055 for h,-h-k,-l	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	2527	wwPDB-VP
Average B, all atoms (Å^2)	53.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.64% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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: 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	1.21	5/2560 (0.2%)	1.70	26/3464 (0.8%)

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 outliers	#Planarity outliers
1	A	0	5

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	247	GLU	C-N	-39.79	0.42	1.34
1	A	211	GLU	C-N	-8.81	1.13	1.34
1	A	199	ILE	C-N	-8.53	1.14	1.34
1	A	205	VAL	N-CA	6.05	1.58	1.46
1	A	204	LEU	CA-C	5.27	1.66	1.52

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	238	SER	O-C-N	-44.21	51.96	122.70
1	A	247	GLU	C-N-CA	-40.23	21.12	121.70
1	A	247	GLU	CA-C-N	-36.54	36.80	117.20
1	A	211	GLU	O-C-N	-23.81	84.61	122.70
1	A	247	GLU	O-C-N	-14.19	100.00	122.70
1	A	199	ILE	O-C-N	-14.18	100.01	122.70
1	A	199	ILE	CA-C-N	10.08	139.38	117.20
1	A	211	GLU	CA-C-N	8.29	135.44	117.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	347	ASP	CB-CG-OD2	8.14	125.63	118.30
1	A	340	ASP	CB-CG-OD2	7.33	124.89	118.30
1	A	204	LEU	CA-C-N	6.53	131.57	117.20
1	A	209	LEU	CA-C-N	-6.47	103.26	116.20
1	A	245	GLU	CB-CA-C	6.41	123.23	110.40
1	A	237	ASP	CB-CG-OD2	6.24	123.91	118.30
1	A	246	ASN	C-N-CA	-5.72	107.39	121.70
1	A	378	ASP	CB-CG-OD2	5.64	123.38	118.30
1	A	205	VAL	CB-CA-C	-5.61	100.75	111.40
1	A	275	LEU	CA-CB-CG	5.56	128.08	115.30
1	A	379	ASP	CB-CG-OD2	5.37	123.13	118.30
1	A	402	ASP	CB-CG-OD2	5.35	123.11	118.30
1	A	388	ASP	CB-CG-OD2	5.25	123.03	118.30
1	A	209	LEU	CB-CA-C	5.22	120.12	110.20
1	A	429	ASP	CB-CG-OD2	5.21	122.99	118.30
1	A	334	ASP	CB-CG-OD1	5.19	122.97	118.30
1	A	398	ASP	CB-CG-OD2	5.13	122.92	118.30
1	A	209	LEU	C-N-CA	5.12	133.05	122.30

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	211	GLU	Peptide,Mainchain
1	A	238	SER	Mainchain
1	A	247	GLU	Peptide,Mainchain

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	2508	0	2540	32	0
2	A	10	0	0	0	0
3	A	9	0	0	0	0
All	All	2527	0	2540	32	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 7.

All (32) 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:259:LEU:HD23	1:A:291:ARG:HD2	1.61	0.82
1:A:329:LEU:HD21	1:A:411:MET:HE1	1.65	0.78
1:A:379:ASP:N	1:A:379:ASP:OD1	2.16	0.76
1:A:411:MET:O	1:A:411:MET:HE2	1.93	0.68
1:A:411:MET:CE	1:A:414:MET:HB2	2.23	0.68
1:A:411:MET:HE3	1:A:414:MET:HB2	1.77	0.64
1:A:317:THR:HG21	1:A:349:LEU:HD22	1.80	0.63
1:A:286:SER:O	1:A:290:ARG:HG3	1.99	0.62
1:A:127:GLU:OE1	1:A:185:LYS:HE3	2.02	0.60
1:A:289:ARG:NH1	1:A:444:GLY:O	2.36	0.58
1:A:255:LYS:O	1:A:259:LEU:HD12	2.06	0.55
1:A:366:ARG:HD3	1:A:378:ASP:OD1	2.07	0.55
1:A:167:TRP:CE2	1:A:294:LYS:HE3	2.44	0.52
1:A:312:LEU:O	1:A:315:GLY:N	2.43	0.52
1:A:295:ARG:NH2	1:A:444:GLY:H	2.09	0.50
1:A:306:ARG:HD2	1:A:333:THR:OG1	2.11	0.49
1:A:236:ILE:HG23	1:A:284:LEU:HD11	1.95	0.49
1:A:154:ILE:O	1:A:157:PRO:HD3	2.14	0.48
1:A:388:ASP:HB2	1:A:391:ALA:HB2	1.95	0.48
1:A:224:ARG:HA	1:A:272:ILE:HD11	1.94	0.48
1:A:414:MET:HA	1:A:414:MET:HE3	1.96	0.48
1:A:295:ARG:HH22	1:A:444:GLY:H	1.62	0.47
1:A:361:HIS:CD2	1:A:393:GLU:HB2	2.50	0.47
1:A:268:ASP:OD2	1:A:270:ASP:HB3	2.15	0.47
1:A:261:GLN:O	1:A:269:ASN:ND2	2.48	0.46
1:A:140:GLU:H	1:A:140:GLU:CD	2.21	0.44
1:A:438:GLU:O	1:A:440:PHE:N	2.50	0.44
1:A:174:GLY:O	1:A:279:ASN:HA	2.17	0.44
1:A:138:GLY:O	1:A:143:LYS:HE3	2.19	0.43
1:A:234:ASP:O	1:A:235:GLU:C	2.57	0.42
1:A:235:GLU:HG2	1:A:238:SER:HB3	2.01	0.41
1:A:237:ASP:HA	1:A:284:LEU:HD12	2.03	0.41

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	320/322 (99%)	290 (91%)	18 (6%)	12 (4%)	3 10

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	240	CYS
1	A	312	LEU
1	A	202	SER
1	A	210	GLY
1	A	212	SER
1	A	265	VAL
1	A	313	HIS
1	A	247	GLU
1	A	439	ASP
1	A	208	TRP
1	A	438	GLU
1	A	389	PRO

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	276/276 (100%)	251 (91%)	25 (9%)	9 27

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	127	GLU
1	A	140	GLU
1	A	144	GLU
1	A	165	THR
1	A	181	SER
1	A	207	LYS
1	A	211	GLU
1	A	214	LYS
1	A	252	ARG
1	A	256	THR
1	A	262	MET
1	A	267	VAL
1	A	272	ILE
1	A	273	LEU
1	A	357	GLN
1	A	366	ARG
1	A	378	ASP
1	A	382	THR
1	A	385	SER
1	A	397	MET
1	A	398	ASP
1	A	414	MET
1	A	432	LYS
1	A	435	LYS
1	A	442	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	318	GLN
1	A	357	GLN
1	A	420	ASN
1	A	428	HIS
1	A	442	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 monosaccharides in this entry.

5.6 Ligand geometry (i)

2 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	601	-	4,4,4	0.36	0	6,6,6	0.44	0
2	SO4	A	602	-	4,4,4	0.17	0	6,6,6	0.24	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)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	A	3

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	199:ILE	C	200:SER	N	1.14
1	A	211:GLU	C	212:SER	N	1.13
1	A	247:GLU	C	248:SER	N	0.42

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	301/322 (93%)	-0.06	3 (0%) 82 77	30, 46, 87, 105	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	249	GLU	2.9
1	A	252	ARG	2.2
1	A	250	ALA	2.1

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 monosaccharides 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	SO4	A	601	5/5	0.97	0.14	49,49,51,53	0
2	SO4	A	602	5/5	0.99	0.16	38,38,40,41	0

6.5 Other polymers (i)

There are no such residues in this entry.