



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 7, 2026 – 02:03 AM UTC

PDB ID : 1VCB / pdb_00001vcb
Title : THE VHL-ELONGINC-ELONGINB STRUCTURE
Authors : Stebbins, C.E.; Kaelin, W.G.; Pavletich, N.P.
Deposited on : 1999-03-13
Resolution : 2.70 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) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

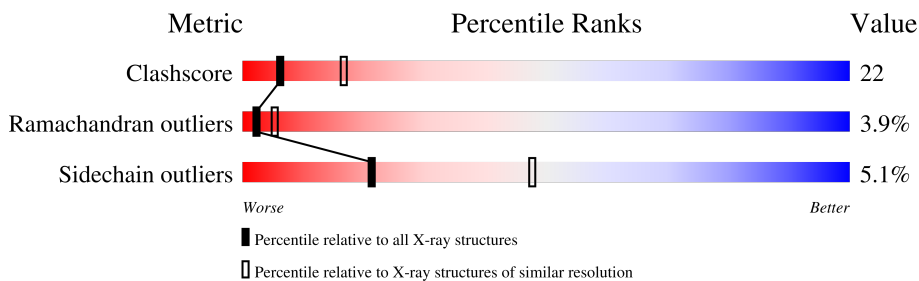
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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))
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)





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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	118	42% 32% 8% 17%
1	D	118	42% 34% 8% 17%
1	G	118	43% 33% 7% 17%
1	J	118	41% 31% 11% 17%
2	B	112	45% 29% • 21%
2	E	112	45% 30% • 21%
2	H	112	44% 31% • 21%
2	K	112	43% 31% • 21%

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Mol	Chain	Length	Quality of chain
3	C	160	 53% 28% 7% • 11%
3	F	160	 54% 28% 6% • 11%
3	I	160	 54% 28% 6% • 11%
3	L	160	 54% 28% 6% • 11%

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 10858 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 PROTEIN (ELONGIN B).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	98	755	477	128	146	4	0	0	0
1	D	98	755	477	128	146	4	0	0	0
1	G	98	755	477	128	146	4	0	0	0
1	J	98	755	477	128	146	4	0	0	0

- Molecule 2 is a protein called PROTEIN (ELONGIN C).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	88	692	447	111	128	6	0	0	0
2	E	88	692	447	111	128	6	0	0	0
2	H	88	692	447	111	128	6	0	0	0
2	K	88	692	447	111	128	6	0	0	0

- Molecule 3 is a protein called PROTEIN (VHL).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	C	142	1154	734	211	207	2	0	0	0
3	F	142	1154	734	211	207	2	0	0	0
3	I	142	1154	734	211	207	2	0	0	0
3	L	142	1154	734	211	207	2	0	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	32	Total O 32 32	0	0
4	B	28	Total O 28 28	0	0
4	C	59	Total O 59 59	0	0
4	D	27	Total O 27 27	0	0
4	E	28	Total O 28 28	0	0
4	F	45	Total O 45 45	0	0
4	G	22	Total O 22 22	0	0
4	H	19	Total O 19 19	0	0
4	I	55	Total O 55 55	0	0
4	J	41	Total O 41 41	0	0
4	K	29	Total O 29 29	0	0
4	L	69	Total O 69 69	0	0

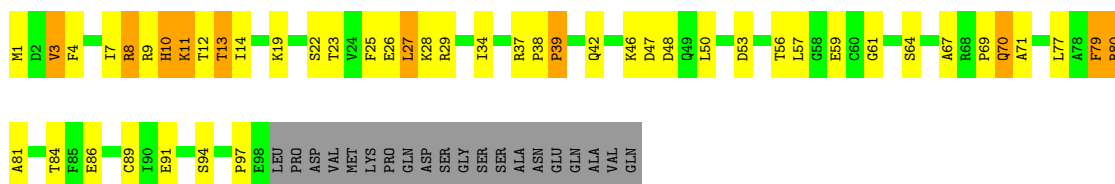
3 Residue-property plots [i](#)

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. 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. 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.

Note EDS was not executed.

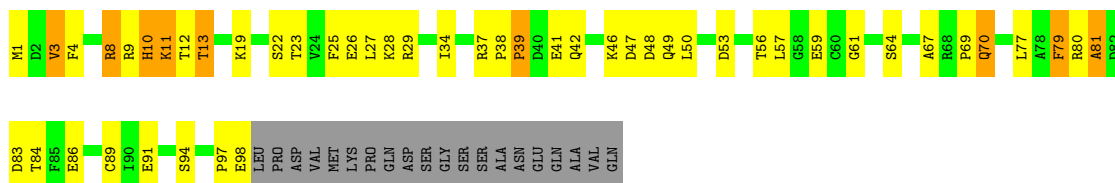
- Molecule 1: PROTEIN (ELONGIN B)

Chain A: 



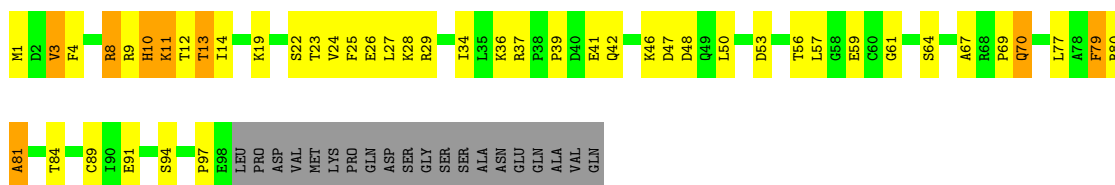
- Molecule 1: PROTEIN (ELONGIN B)

Chain D: 



- Molecule 1: PROTEIN (ELONGIN B)

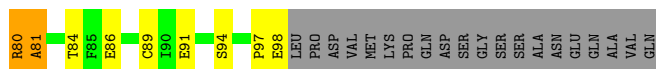
Chain G: 



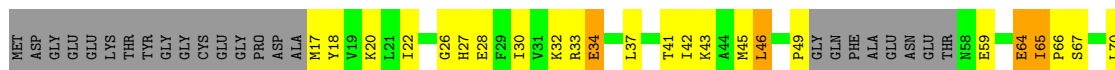
- Molecule 1: PROTEIN (ELONGIN B)

Chain J: 

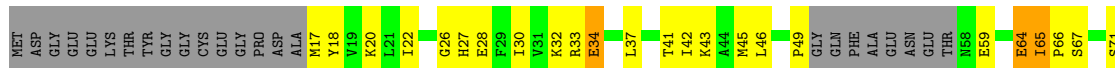




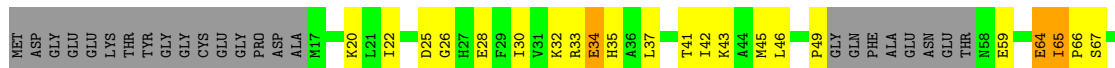
- Molecule 2: PROTEIN (ELONGIN C)



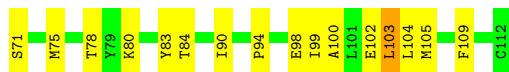
- Molecule 2: PROTEIN (ELONGIN C)



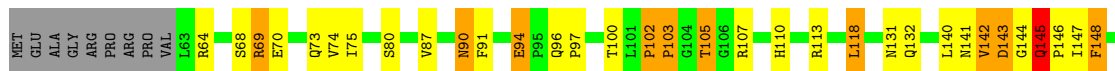
- Molecule 2: PROTEIN (ELONGIN C)



- Molecule 2: PROTEIN (ELONGIN C)



- Molecule 3: PROTEIN (VHL)





- Molecule 3: PROTEIN (VHL)

Chain F: 54% 28% 6% • 11%



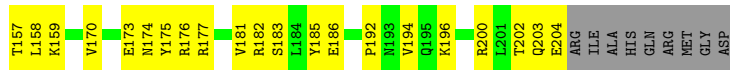
- Molecule 3: PROTEIN (VHL)

Chain I: 54% 28% 6% • 11%



- Molecule 3: PROTEIN (VHL)

Chain L: 54% 28% 6% • 11%



4 Data and refinement statistics

Xtrriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 41 2 2	Depositor
Cell constants a, b, c, α , β , γ	93.50Å 93.50Å 362.30Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.70	Depositor
% Data completeness (in resolution range)	84.6 (20.00-2.70)	Depositor
R_{merge}	0.07	Depositor
R_{sym}	7.00	Depositor
Refinement program	CNS 0.3	Depositor
R, R_{free}	0.238 , 0.289	Depositor
Estimated twinning fraction	No twinning to report.	Xtrriage
Total number of atoms	10858	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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	0.83	0/769	1.18	6/1040 (0.6%)
1	D	0.77	0/769	1.16	4/1040 (0.4%)
1	G	0.78	0/769	1.17	4/1040 (0.4%)
1	J	0.84	0/769	1.20	5/1040 (0.5%)
2	B	0.82	0/706	1.18	8/952 (0.8%)
2	E	0.78	0/706	1.18	7/952 (0.7%)
2	H	0.79	0/706	1.19	7/952 (0.7%)
2	K	0.83	0/706	1.18	8/952 (0.8%)
3	C	0.79	0/1184	1.25	14/1616 (0.9%)
3	F	0.77	0/1184	1.25	9/1616 (0.6%)
3	I	0.80	0/1184	1.28	11/1616 (0.7%)
3	L	0.84	0/1184	1.28	11/1616 (0.7%)
All	All	0.80	0/10636	1.22	94/14432 (0.7%)

There are no bond length outliers.

The worst 5 of 94 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	26	GLY	N-CA-C	8.72	125.07	114.69
2	K	26	GLY	N-CA-C	8.56	124.88	114.69
2	E	26	GLY	N-CA-C	8.39	124.68	114.69
3	I	145	GLN	N-CA-C	8.28	128.12	109.81
3	F	145	GLN	N-CA-C	8.22	127.98	109.81

There are no chirality outliers.

There are no planarity outliers.

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	755	0	744	38	0
1	D	755	0	744	34	0
1	G	755	0	744	35	0
1	J	755	0	744	39	0
2	B	692	0	685	24	0
2	E	692	0	685	26	0
2	H	692	0	685	29	0
2	K	692	0	685	38	0
3	C	1154	0	1144	64	1
3	F	1154	0	1144	56	1
3	I	1154	0	1144	58	0
3	L	1154	0	1144	59	0
4	A	32	0	0	1	0
4	B	28	0	0	0	0
4	C	59	0	0	6	0
4	D	27	0	0	2	0
4	E	28	0	0	3	0
4	F	45	0	0	5	0
4	G	22	0	0	0	0
4	H	19	0	0	4	0
4	I	55	0	0	7	0
4	J	41	0	0	2	0
4	K	29	0	0	6	0
4	L	69	0	0	11	0
All	All	10858	0	10292	464	1

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 22.

The worst 5 of 464 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)
3:I:142:VAL:HA	4:I:264:HOH:O	1.59	1.02
3:F:196:LYS:HE3	3:F:200:ARG:HH22	1.21	1.00
3:C:196:LYS:HE3	3:C:200:ARG:HH22	1.25	0.99
3:L:196:LYS:HE3	3:L:200:ARG:HH22	1.25	0.99
3:I:196:LYS:HE3	3:I:200:ARG:HH22	1.24	0.98

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the sym-

metry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:70:GLU:OE2	3:F:142:VAL:O[6_565]	2.17	0.03

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	96/118 (81%)	85 (88%)	5 (5%)	6 (6%)	1	2
1	D	96/118 (81%)	85 (88%)	5 (5%)	6 (6%)	1	2
1	G	96/118 (81%)	85 (88%)	5 (5%)	6 (6%)	1	2
1	J	96/118 (81%)	85 (88%)	5 (5%)	6 (6%)	1	2
2	B	84/112 (75%)	82 (98%)	2 (2%)	0	100	100
2	E	84/112 (75%)	81 (96%)	3 (4%)	0	100	100
2	H	84/112 (75%)	81 (96%)	3 (4%)	0	100	100
2	K	84/112 (75%)	81 (96%)	3 (4%)	0	100	100
3	C	140/160 (88%)	127 (91%)	6 (4%)	7 (5%)	1	3
3	F	140/160 (88%)	128 (91%)	5 (4%)	7 (5%)	1	3
3	I	140/160 (88%)	127 (91%)	7 (5%)	6 (4%)	2	4
3	L	140/160 (88%)	126 (90%)	8 (6%)	6 (4%)	2	4
All	All	1280/1560 (82%)	1173 (92%)	57 (4%)	50 (4%)	2	5

5 of 50 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	81	ALA
3	C	142	VAL
3	C	143	ASP
3	C	145	GLN
1	D	81	ALA

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	81/103 (79%)	75 (93%)	6 (7%)	13	32
1	D	81/103 (79%)	75 (93%)	6 (7%)	13	32
1	G	81/103 (79%)	75 (93%)	6 (7%)	13	32
1	J	81/103 (79%)	75 (93%)	6 (7%)	13	32
2	B	76/96 (79%)	71 (93%)	5 (7%)	15	36
2	E	76/96 (79%)	72 (95%)	4 (5%)	20	46
2	H	76/96 (79%)	72 (95%)	4 (5%)	20	46
2	K	76/96 (79%)	72 (95%)	4 (5%)	20	46
3	C	130/147 (88%)	127 (98%)	3 (2%)	44	73
3	F	130/147 (88%)	125 (96%)	5 (4%)	29	58
3	I	130/147 (88%)	125 (96%)	5 (4%)	29	58
3	L	130/147 (88%)	126 (97%)	4 (3%)	35	65
All	All	1148/1384 (83%)	1090 (95%)	58 (5%)	21	48

5 of 58 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	F	186	GLU
3	L	118	LEU
2	H	34	GLU
3	L	77	CYS
1	J	70	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 28 such sidechains are listed below:

Mol	Chain	Res	Type
1	G	10	HIS
3	L	174	ASN
2	H	35	HIS
1	J	70	GLN

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Mol	Chain	Res	Type
1	G	70	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](#)

There are no ligands in this entry.

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](#)

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates [i](#)

EDS was not executed - this section is therefore empty.

6.4 Ligands [i](#)

EDS was not executed - this section is therefore empty.

6.5 Other polymers [i](#)

EDS was not executed - this section is therefore empty.