



Full wwPDB/EMDataBank EM Map/Model Validation Report ⓘ

Aug 15, 2018 – 09:46 AM EDT

PDB ID : 5YHQ
EMDB ID: : EMD-6829
Title : Cryo-EM Structure of CVA6 VLP
Authors : Chen, J.; Zhang, C.; Huang, Z.; Cong, Y.
Deposited on : 2017-09-29
Resolution : 3.00 Å(reported)

This is a Full wwPDB/EMDataBank EM Map/Model Validation Report
for a publicly released PDB/EMDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

MolProbity : 4.02b-467
Percentile statistics : 20171227.v01 (using entries in the PDB archive December 27th 2017)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20031172

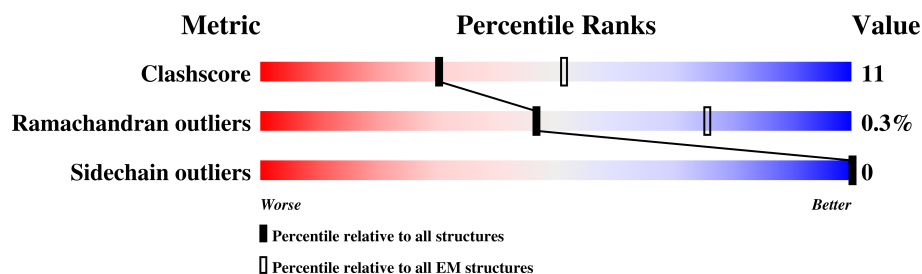
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.00 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)	EM structures (#Entries)
Clashscore	136327	1886
Ramachandran outliers	132723	1663
Sidechain outliers	132532	1531

The table below summarises the geometric issues observed across the polymeric chains. The red, orange, yellow and green segments on the bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	305	
2	C	240	
3	B	325	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5335 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, 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 Capsid protein VP1.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	226	Total	C	N	O	S	0	0
			1800	1144	308	338	10		

- Molecule 2 is a protein called Capsid protein VP3.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	C	228	Total	C	N	O	S	0	0
			1758	1123	289	337	9		

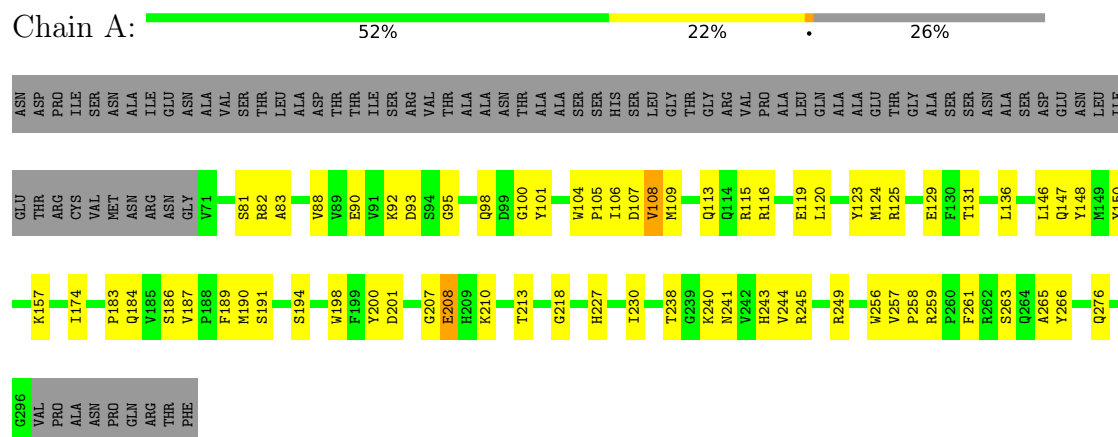
- Molecule 3 is a protein called capsid protein VP0.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	B	228	Total	C	N	O	S	0	0
			1777	1147	297	326	7		

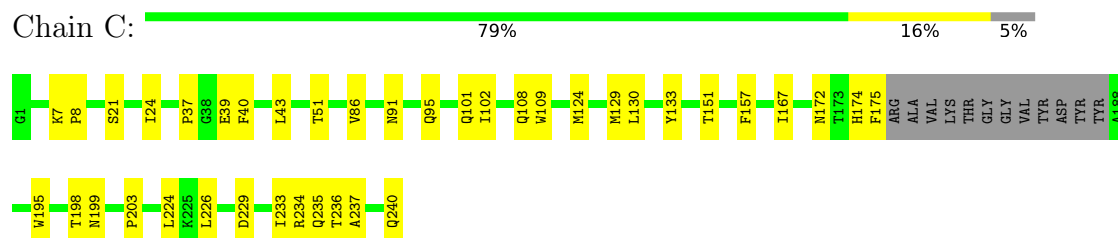
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA 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.

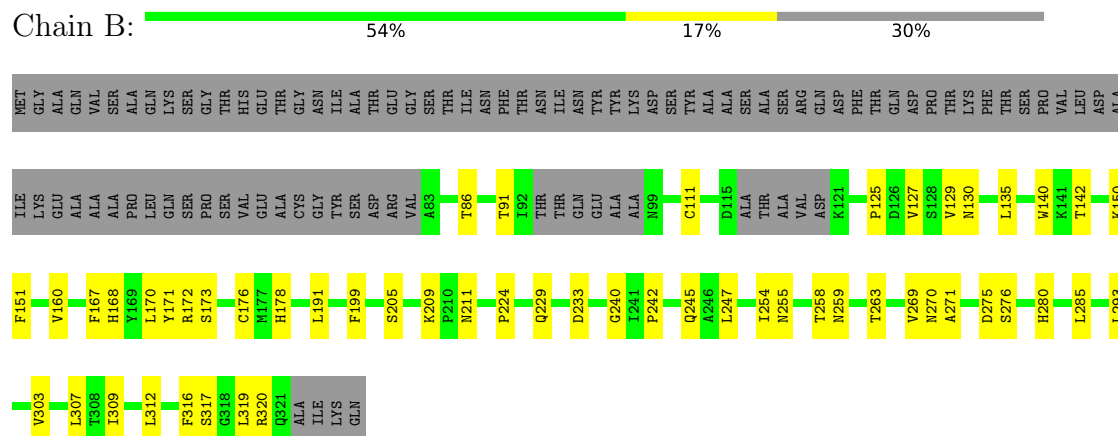
• Molecule 1: Capsid protein VP1



• Molecule 2: Capsid protein VP3



• Molecule 3: capsid protein VP0



4 Experimental information

Property	Value	Source
Reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	Depositor
Number of particles used	11596	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	42	Depositor
Minimum defocus (nm)	Not provided	Depositor
Maximum defocus (nm)	Not provided	Depositor
Magnification	Not provided	Depositor
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor

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	$\# Z > 2$	RMSZ	$\# Z > 2$
1	A	0.52	0/1855	0.57	0/2527
2	C	0.44	0/1803	0.54	0/2470
3	B	0.47	0/1832	0.56	1/2508 (0.0%)
All	All	0.48	0/5490	0.56	1/7505 (0.0%)

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	2

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	233	ASP	N-CA-C	-5.88	95.11	111.00

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	207	GLY	Peptide
1	A	208	GLU	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	1800	0	1720	53	0
2	C	1758	0	1731	40	0
3	B	1777	0	1722	38	0
All	All	5335	0	5173	113	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 11.

All (113) 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)
2:C:226:LEU:HD23	2:C:226:LEU:O	1.74	0.85
2:C:151:THR:O	2:C:151:THR:HG22	1.82	0.79
3:B:275:ASP:OD2	3:B:280:HIS:ND1	2.15	0.78
1:A:257:VAL:HG12	1:A:257:VAL:O	1.84	0.77
1:A:208:GLU:HG2	1:A:210:LYS:HE3	1.66	0.77
1:A:83:ALA:HB2	1:A:249:ARG:HB2	1.68	0.74
2:C:234:ARG:HG3	2:C:235:GLN:H	1.52	0.74
3:B:167:PHE:O	3:B:320:ARG:HG2	1.89	0.72
2:C:133:TYR:HB3	2:C:151:THR:HG22	1.71	0.72
3:B:258:THR:HG22	3:B:259:ASN:H	1.54	0.72
1:A:90:GLU:HG2	1:A:243:HIS:HD1	1.54	0.72
2:C:86:VAL:HG11	2:C:167:ILE:HG21	1.71	0.72
2:C:133:TYR:HB3	2:C:151:THR:CG2	2.21	0.70
2:C:91:ASN:OD1	2:C:95:GLN:NE2	2.26	0.69
3:B:209:LYS:HG2	3:B:211:ASN:H	1.59	0.68
2:C:43:LEU:HB3	2:C:102:ILE:HD11	1.78	0.66
1:A:124:MET:HG3	1:A:198:TRP:HZ3	1.61	0.66
2:C:108:GLN:HA	2:C:172:ASN:HD21	1.63	0.64
3:B:125:PRO:HB2	3:B:129:VAL:HG22	1.80	0.63
1:A:106:ILE:O	1:A:106:ILE:HG22	1.98	0.63
1:A:113:GLN:NE2	2:C:229:ASP:HB3	2.13	0.62
2:C:37:PRO:HG2	3:B:269:VAL:HG22	1.81	0.62
1:A:150:TYR:OH	1:A:190:MET:SD	2.58	0.62
1:A:148:TYR:HB2	1:A:174:ILE:HB	1.81	0.61
2:C:235:GLN:HB3	2:C:240:GLN:HE22	1.66	0.61
1:A:119:GLU:OE1	1:A:266:TYR:OH	2.18	0.61
3:B:270:ASN:OD1	3:B:271:ALA:N	2.31	0.60
1:A:116:ARG:HD2	2:C:233:ILE:HD13	1.84	0.59
1:A:265:ALA:HA	2:C:235:GLN:NE2	2.17	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:127:VAL:HA	3:B:130:ASN:HB2	1.83	0.59
2:C:51:THR:HA	3:B:247:LEU:HD12	1.84	0.58
1:A:131:THR:OG1	1:A:184:GLN:OE1	2.20	0.57
1:A:129:GLU:OE1	1:A:249:ARG:NH2	2.37	0.57
1:A:93:ASP:HB3	1:A:238:THR:O	2.04	0.57
1:A:125:ARG:NH2	1:A:194:SER:O	2.38	0.57
1:A:113:GLN:HE22	2:C:229:ASP:HB3	1.71	0.56
1:A:124:MET:HG3	1:A:198:TRP:CZ3	2.40	0.56
3:B:176:CYS:HB2	3:B:312:LEU:HD11	1.88	0.55
1:A:201:ASP:O	3:B:280:HIS:HD2	1.90	0.55
3:B:171:TYR:HH	3:B:173:SER:HG	1.55	0.55
1:A:276:GLN:HG3	3:B:205:SER:O	2.07	0.54
1:A:146:LEU:HD13	1:A:230:ILE:HD11	1.88	0.54
2:C:198:THR:OG1	2:C:199:ASN:N	2.41	0.54
1:A:123:TYR:HB2	1:A:256:TRP:HB2	1.90	0.54
1:A:120:LEU:O	1:A:259:ARG:N	2.33	0.53
2:C:235:GLN:O	2:C:237:ALA:N	2.41	0.53
1:A:148:TYR:HE1	1:A:230:ILE:HD12	1.73	0.53
2:C:133:TYR:HD2	2:C:151:THR:HG21	1.73	0.53
3:B:224:PRO:HB3	3:B:229:GLN:HB2	1.90	0.53
3:B:171:TYR:HD1	3:B:316:PHE:HE1	1.56	0.53
1:A:148:TYR:CE1	1:A:230:ILE:HD12	2.44	0.52
3:B:151:PHE:HE1	3:B:309:ILE:HG21	1.74	0.52
3:B:151:PHE:CE2	3:B:285:LEU:HD13	2.45	0.52
2:C:234:ARG:HG3	2:C:235:GLN:N	2.24	0.51
2:C:226:LEU:CD2	2:C:226:LEU:O	2.53	0.50
3:B:275:ASP:OD1	3:B:276:SER:N	2.37	0.49
3:B:178:HIS:ND1	3:B:263:THR:OG1	2.39	0.49
3:B:142:THR:HG22	3:B:293:LEU:O	2.12	0.49
3:B:168:HIS:CE1	3:B:319:LEU:HD13	2.47	0.49
3:B:111:CYS:SG	3:B:172:ARG:NE	2.86	0.49
1:A:148:TYR:CE2	1:A:183:PRO:HG2	2.48	0.49
3:B:170:LEU:HB2	3:B:317:SER:HB3	1.94	0.48
2:C:129:MET:HG2	2:C:157:PHE:CZ	2.48	0.48
3:B:151:PHE:HZ	3:B:285:LEU:HD22	1.79	0.48
1:A:256:TRP:CD1	2:C:39:GLU:HB3	2.49	0.48
2:C:7:LYS:HB3	2:C:8:PRO:HD2	1.94	0.48
3:B:86:THR:OG1	3:B:91:THR:HG22	2.15	0.47
3:B:135:LEU:HD21	3:B:307:LEU:HD23	1.97	0.47
1:A:95:GLY:O	1:A:98:GLN:HG3	2.14	0.47
1:A:157:LYS:HE2	1:A:227:HIS:CE1	2.50	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:258:PRO:HG3	2:C:40:PHE:CD2	2.50	0.46
2:C:174:HIS:CG	2:C:175:PHE:H	2.33	0.46
2:C:101:GLN:HG3	3:B:245:GLN:HE22	1.80	0.46
3:B:129:VAL:HG12	3:B:160:VAL:HG12	1.98	0.46
3:B:209:LYS:HB3	3:B:211:ASN:OD1	2.16	0.46
2:C:133:TYR:HB3	2:C:151:THR:HG21	1.98	0.46
3:B:171:TYR:OH	3:B:173:SER:OG	2.26	0.46
1:A:257:VAL:O	1:A:257:VAL:CG1	2.57	0.45
2:C:109:TRP:HB3	2:C:224:LEU:HD23	1.98	0.45
1:A:113:GLN:HG3	2:C:233:ILE:HG21	1.98	0.45
1:A:88:VAL:HG12	1:A:245:ARG:HG2	1.98	0.45
1:A:256:TRP:NE1	2:C:39:GLU:HB3	2.31	0.45
2:C:43:LEU:HD22	2:C:102:ILE:HD11	1.99	0.45
1:A:263:SER:HB3	3:B:242:PRO:HD3	2.00	0.44
3:B:151:PHE:CZ	3:B:285:LEU:HD22	2.52	0.44
2:C:124:MET:HB3	2:C:203:PRO:HG3	1.99	0.44
1:A:116:ARG:HH21	1:A:261:PHE:HB3	1.82	0.44
1:A:210:LYS:HD3	1:A:213:THR:O	2.17	0.44
2:C:86:VAL:O	2:C:86:VAL:HG12	2.18	0.43
1:A:200:TYR:O	1:A:218:GLY:HA2	2.19	0.43
1:A:189:PHE:CZ	1:A:191:SER:HB3	2.53	0.43
1:A:230:ILE:HD13	1:A:244:VAL:HG21	2.01	0.43
1:A:104:TRP:CD1	1:A:105:PRO:HD2	2.54	0.43
2:C:233:ILE:HG13	2:C:233:ILE:O	2.19	0.43
3:B:150:LYS:HE2	3:B:199:PHE:HD2	1.84	0.42
1:A:129:GLU:OE2	2:C:21:SER:OG	2.37	0.42
1:A:92:LYS:HA	1:A:241:ASN:HD22	1.83	0.42
1:A:186:SER:OG	2:C:21:SER:HB2	2.19	0.42
2:C:130:LEU:HB3	2:C:195:TRP:HB2	2.01	0.42
1:A:109:MET:HE2	1:A:115:ARG:HG3	2.02	0.42
1:A:136:LEU:HD11	1:A:240:LYS:HB3	2.02	0.42
2:C:133:TYR:CD2	2:C:151:THR:HG21	2.55	0.41
1:A:81:SER:O	1:A:82:ARG:HG2	2.21	0.41
1:A:100:GLY:O	1:A:101:TYR:HB3	2.21	0.41
1:A:107:ASP:O	1:A:108:VAL:HB	2.20	0.41
3:B:191:LEU:HB2	3:B:254:ILE:HB	2.02	0.41
1:A:147:GLN:O	1:A:230:ILE:HA	2.21	0.41
1:A:120:LEU:HD13	1:A:261:PHE:CE1	2.56	0.41
3:B:140:TRP:HB3	3:B:303:VAL:HG13	2.03	0.40
1:A:263:SER:OG	3:B:240:GLY:O	2.24	0.40
1:A:187:VAL:HG22	2:C:24:ILE:HD11	2.02	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:255:ASN:O	3:B:258:THR:O	2.38	0.40
3:B:258:THR:HG22	3:B:259:ASN:N	2.29	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 PDB entries followed by that with respect to all EM 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	224/305 (73%)	202 (90%)	21 (9%)	1 (0%)	36	76
2	C	224/240 (93%)	216 (96%)	7 (3%)	1 (0%)	36	76
3	B	222/325 (68%)	212 (96%)	10 (4%)	0	100	100
All	All	670/870 (77%)	630 (94%)	38 (6%)	2 (0%)	47	80

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	108	VAL
2	C	236	THR

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 PDB entries followed by that with respect to all EM 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	196/258 (76%)	196 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	C	197/206 (96%)	197 (100%)	0	100	100
3	B	196/272 (72%)	196 (100%)	0	100	100
All	All	589/736 (80%)	589 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	113	GLN
1	A	137	ASN
1	A	219	GLN
2	C	152	HIS
2	C	235	GLN
3	B	164	ASN
3	B	168	HIS
3	B	245	GLN
3	B	313	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 carbohydrates 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 ⓘ

There are no chain breaks in this entry.