

Full wwPDB X-ray Structure Validation Report (i)

Jul 25, 2023 – 12:21 AM EDT

PDB ID	:	1QBZ
Title	:	THE CRYSTAL STRUCTURE OF THE SIV GP41 ECTODOMAIN AT 1.47
		А
Authors	:	Yang, ZN.; Mueser, T.C.; Kaufman, J.; Stahl, S.J.; Wingfield, P.T.; Hyde,
		C.C.
Deposited on		
Resolution	:	1.47 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 (1)) were used in the production of this report:

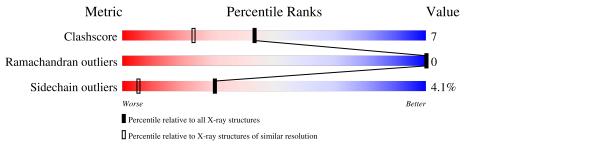
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.34

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: $X\hbox{-}RAY\,DIFFRACTION$

The reported resolution of this entry is 1.47 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	Similar resolution		
Metric	$(\# {\rm Entries})$	$(\# { m Entries}, { m resolution} { m range}({ m \AA}))$		
Clashscore	141614	4955 (1.50-1.46)		
Ramachandran outliers	138981	4846 (1.50-1.46)		
Sidechain outliers	138945	4844 (1.50-1.46)		

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%

Note EDS was not executed.

Mol	Chain	Length	Quality of ch	nain	
1	А	123	58%	24%	• 17%
1	В	123	81%		10% • 7%
1	С	123	76%		18% •••



2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 3229 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.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	А	102	Total	С	Ν	0	\mathbf{S}	0	14	0
	Л	102	894	564	157	172	1	0		
1	В	114	Total	С	Ν	0	S	0	13	0
	D	114	1016	643	182	189	2	0	15	U
1	C	120	Total	С	Ν	0	S	0	0	0
I C	120	1035	654	184	196	1	0	9	0	

• Molecule 1 is a protein called PROTEIN (SIV GP41 ECTODOMAIN).

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
А	86	ALA	CYS	engineered mutation	GB AAA47708
В	86	ALA	CYS	engineered mutation	GB AAA47708
С	86	ALA	CYS	engineered mutation	GB AAA47708
А	92	ALA	CYS	engineered mutation	GB AAA47708
В	92	ALA	CYS	engineered mutation	GB AAA47708
С	92	ALA	CYS	engineered mutation	GB AAA47708

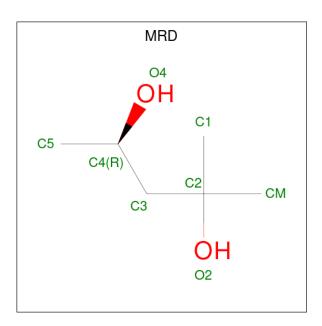
• Molecule 2 is MERCURY (II) ION (three-letter code: HG) (formula: Hg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	А	1	Total Hg 1 1	0	0
2	В	1	Total Hg 1 1	0	0

• Molecule 3 is (4R)-2-METHYLPENTANE-2,4-DIOL (three-letter code: MRD) (formula: $C_6H_{14}O_2$).







Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	А	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 8 & 6 & 2 \end{array}$	0	0
3	С	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 8 & 6 & 2 \end{array}$	0	0

• Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	В	1	Total Cl 1 1	0	0
4	С	2	Total Cl 2 2	0	0

• Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	А	66	Total O 66 66	0	1
5	В	105	Total O 105 105	0	0
5	С	92	TotalO9292	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 (SIV GP41 ECTODOMAIN)





4 Data and refinement statistics (i)

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

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants	91.12Å 47.21Å 77.47Å	Depositor
a, b, c, α , β , γ	90.00° 100.83° 90.00°	Depositor
Resolution (Å)	15.00 - 1.47	Depositor
% Data completeness	5.0 (15.00-1.47)	Depositor
(in resolution range)	5.6 (15.66-1.47)	Depositor
R_{merge}	0.06	Depositor
R _{sym}	(Not available)	Depositor
Refinement program	SHELXL-97	Depositor
R, R_{free}	0.155 , 0.218	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3229	wwPDB-VP
Average B, all atoms $(Å^2)$	24.0	wwPDB-VP



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: CL, MRD, HG

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		
	Unam	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	А	0.55	0/968	1.35	11/1304~(0.8%)	
1	В	0.57	0/1095	1.28	6/1478~(0.4%)	
1	С	0.57	0/1095	1.36	9/1482~(0.6%)	
All	All	0.56	0/3158	1.33	26/4264~(0.6%)	

There are no bond length outliers.

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
1	С	30	ARG	NE-CZ-NH2	10.47	125.54	120.30
1	С	30	ARG	NE-CZ-NH1	-10.41	115.09	120.30
1	В	49	ARG	NE-CZ-NH2	9.96	125.28	120.30
1	А	49	ARG	NE-CZ-NH2	-9.49	115.56	120.30
1	С	67	ARG	NE-CZ-NH2	-9.11	115.75	120.30
1	В	121[A]	PHE	CB-CG-CD1	-8.33	114.97	120.80
1	В	121[B]	PHE	CB-CG-CD1	-8.33	114.97	120.80
1	С	55	ARG	CD-NE-CZ	8.27	135.18	123.60
1	В	121[A]	PHE	CB-CG-CD2	8.23	126.56	120.80
1	В	121[B]	PHE	CB-CG-CD2	8.23	126.56	120.80
1	С	117[A]	ARG	NE-CZ-NH2	7.70	124.15	120.30
1	С	117[B]	ARG	NE-CZ-NH2	7.70	124.15	120.30
1	А	120	ASP	CB-CG-OD2	7.42	124.98	118.30
1	А	117	ARG	NE-CZ-NH2	7.02	123.81	120.30
1	А	117	ARG	CD-NE-CZ	6.77	133.08	123.60
1	А	132	GLU	CA-CB-CG	6.44	127.57	113.40
1	А	67	ARG	NE-CZ-NH2	6.11	123.35	120.30
1	А	55[A]	ARG	CD-NE-CZ	5.96	131.95	123.60
1	А	55[B]	ARG	CD-NE-CZ	5.96	131.95	123.60
1	А	138	GLU	OE1-CD-OE2	-5.91	116.21	123.30
1	А	81	LEU	CB-CG-CD2	5.80	120.86	111.00

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Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$Observed(^{o})$	$Ideal(^{o})$		
1	С	146	LYS	CA-CB-CG	5.53	125.56	113.40		
1	В	30	ARG	NE-CZ-NH1	5.21	122.90	120.30		
1	А	120	ASP	CB-CG-OD1	-5.17	113.64	118.30		
1	С	67	ARG	NH1-CZ-NH2	5.07	124.97	119.40		
1	С	114	GLU	OE1-CD-OE2	-5.01	117.29	123.30		

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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	А	894	0	894	16	0
1	В	1016	0	1017	8	0
1	С	1035	0	1026	18	0
2	А	1	0	0	0	0
2	В	1	0	0	0	0
3	А	8	0	14	0	0
3	С	8	0	14	0	0
4	В	1	0	0	0	0
4	С	2	0	0	0	0
5	А	66	0	0	6	0
5	В	105	0	0	3	0
5	С	92	0	0	2	0
All	All	3229	0	2965	40	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 (40) 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:C:30:ARG:HE	1:C:148:ASN:HD22	1.31	0.77
1:A:42[B]:GLN:HG2	5:B:8101:HOH:O	1.87	0.73
1:A:131[B]:GLU:HG3	5:A:8066:HOH:O	1.90	0.70

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Continuea from previo	1.0	Interatomic	Clash
Atom-1	Atom-2	distance (Å)	overlap (Å)
1:A:38:GLN:O	1:A:42[A]:GLN:HG3	1.94	0.67
5:A:8052:HOH:O	1:C:66:THR:HG22	2.00	0.61
1:A:142:TYR:O	1:A:145[A]:GLN:HB2	2.02	0.59
1:B:42[B]:GLN:HG2	5:B:8043:HOH:O	2.02	0.58
1:C:30:ARG:HE	1:C:148:ASN:ND2	2.02	0.58
1:C:65[B]:GLN:HG3	1:C:112:TRP:CD2	2.41	0.56
1:C:91:VAL:O	1:C:91:VAL:HG12	2.07	0.55
1:C:65[B]:GLN:HG3	1:C:112:TRP:CE3	2.44	0.53
1:A:144:LEU:O	1:A:147:LEU:HB2	2.09	0.52
1:B:30:ARG:HG3	1:B:147:LEU:CD2	2.39	0.52
1:B:65:GLN:HG3	1:B:112:TRP:CE3	2.44	0.52
1:B:65:GLN:HG3	1:B:112:TRP:CD2	2.47	0.49
1:A:139[A]:LYS:HD2	5:A:8025:HOH:O	2.12	0.48
1:A:36:ILE:O	1:A:40[A]:GLN:HG2	2.14	0.48
1:A:65[A]:GLN:HG3	1:A:112:TRP:CD2	2.49	0.47
1:A:82:ASN:ND2	5:A:8040:HOH:O	2.48	0.47
5:B:8027:HOH:O	1:C:50:GLN:NE2	2.47	0.46
1:C:30:ARG:HH11	1:C:148:ASN:ND2	2.14	0.46
1:C:85:GLY:O	1:C:89:ARG:N	2.48	0.46
1:C:104:THR:HA	5:C:8083:HOH:O	2.14	0.46
1:A:108:ASN:ND2	5:A:8022:HOH:O	2.49	0.46
1:C:114:GLU:OE1	1:C:117[A]:ARG:NH2	2.49	0.46
1:A:74:TYR:OH	1:A:78[A]:GLN:NE2	2.49	0.45
1:C:131[A]:GLU:HG2	5:C:8084:HOH:O	2.16	0.45
1:B:116:GLU:OE2	1:B:120:ASP:OD1	2.35	0.44
1:A:77[A]:ASP:OD2	1:B:96:VAL:HG21	2.18	0.43
1:A:137:GLN:O	1:A:141:MET:HG3	2.18	0.43
1:C:124:GLU:HA	1:C:124:GLU:OE1	2.19	0.43
1:C:49[A]:ARG:HD3	1:C:52[A]:GLU:OE2	2.20	0.42
1:A:104:THR:HG22	1:A:105:PRO:N	2.35	0.42
1:C:30:ARG:NE	1:C:148:ASN:HD22	2.08	0.42
1:A:42[B]:GLN:NE2	1:B:132[B]:GLU:HG3	2.35	0.41
1:C:58:VAL:O	1:C:62:LYS:HG3	2.20	0.41
1:C:145:GLN:O	1:C:149:SER:HB2	2.20	0.41
1:C:58:VAL:CG1	1:C:62:LYS:HE3	2.51	0.40
1:A:48[B]:LYS:HG2	5:A:8021:HOH:O	2.20	0.40

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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	Percent	tiles
1	А	112/123~(91%)	111 (99%)	1 (1%)	0	100	100
1	В	124/123~(101%)	123~(99%)	1 (1%)	0	100	100
1	С	125/123~(102%)	124 (99%)	1 (1%)	0	100	100
All	All	361/369~(98%)	358~(99%)	3 (1%)	0	100	100

There are no Ramachandran outliers to report.

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 side chain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	sed Rotameric Outliers		Percentiles		
1	А	102/109~(94%)	100~(98%)	2(2%)	55 24		
1	В	116/109~(106%)	113~(97%)	3~(3%)	46 15		
1	С	115/109~(106%)	107~(93%)	8 (7%)	15 1		
All	All	333/327~(102%)	320~(96%)	13 (4%)	30 6		

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	А	104	THR
1	А	109	ASN
1	В	28	GLN
1	В	29	SER
1	В	65	GLN

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Contr	nued fron	i previo	us page
Mol	Chain	Res	Type
1	С	30	ARG
1	С	39	GLN
1	С	65[A]	GLN
1	С	65[B]	GLN
1	С	90	GLN
1	С	102	SER
1	С	147	LEU
1	С	149	SER

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Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	А	51	GLN
1	А	80	GLN
1	А	108	ASN
1	А	109	ASN
1	А	125	ASN
1	В	38	GLN
1	В	78	GLN
1	В	108	ASN
1	В	148	ASN
1	С	39	GLN
1	С	40	GLN
1	С	50	GLN
1	С	90	GLN
1	С	108	ASN
1	С	125	ASN
1	С	148	ASN

5.3.3RNA (i)

There are no RNA molecules in this entry.

Non-standard residues in protein, DNA, RNA chains (i) 5.4

There are no non-standard protein/DNA/RNA residues in this entry.

Carbohydrates (i) 5.5

There are no monosaccharides in this entry.



5.6 Ligand geometry (i)

Of 7 ligands modelled in this entry, 5 are monoatomic - leaving 2 for Mogul analysis.

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	Chain	Dec	Link	В	ond leng	gths	В	ond ang	gles
IVIOI	туре	Unam	Res	LIIIK	Counts	RMSZ	# Z >2	Counts	RMSZ	# Z >2
3	MRD	С	7002	-	7,7,7	0.29	0	$9,\!10,\!10$	0.49	0
3	MRD	А	7001	-	7,7,7	0.32	0	9,10,10	1.05	1 (11%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	MRD	С	7002	-	-	2/5/5/5	-
3	MRD	А	7001	-	-	1/5/5/5	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Ζ	$Observed(^{o})$	$Ideal(^{o})$
3	А	7001	MRD	O4-C4-C3	-2.17	102.59	111.36

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	А	7001	MRD	C2-C3-C4-O4
3	С	7002	MRD	C1-C2-C3-C4
3	С	7002	MRD	C2-C3-C4-O4

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)

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.

