



# wwPDB X-ray Structure Validation Summary Report ⓘ

Sep 10, 2017 – 11:08 PM EDT

PDB ID : 4XMK  
Title : Crystal structure of Fab of HIV-1 gp120 V3-specific human monoclonal anti-body 2424 in complex with JR-FL V3 peptide  
Authors : Kong, X.-P.; Pan, R.  
Deposited on : unknown  
Resolution : 3.18 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](mailto:validation@mail.wwpdb.org)

A user guide is available at

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20029824
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	rb-20029824

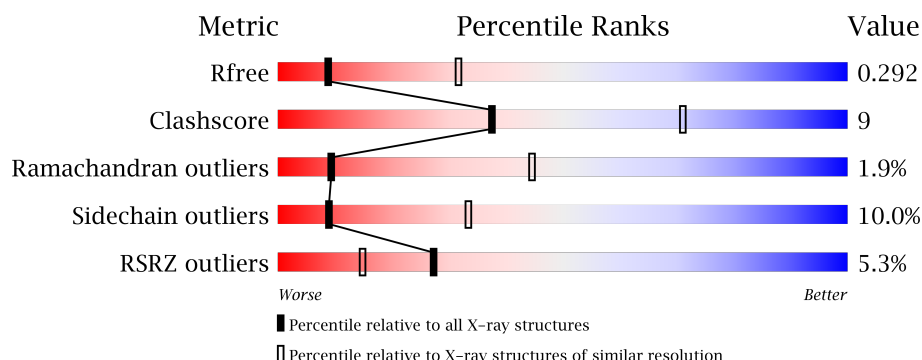
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.18 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))
$R_{free}$	100719	1123 (3.20-3.16)
Clashscore	112137	1255 (3.20-3.16)
Ramachandran outliers	110173	1233 (3.20-3.16)
Sidechain outliers	110143	1232 (3.20-3.16)
RSRZ outliers	101464	1128 (3.20-3.16)

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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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  $\leq 5\%$ . 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	L	215	
1	M	215	
1	N	215	
2	H	223	
2	I	223	

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Mol	Chain	Length	Quality of chain
2	J	223	<div><div></div><div>13%</div><div>73%</div><div>22%</div><div></div><div></div></div>
3	P	11	<div><div></div><div>9%</div><div>64%</div><div>27%</div><div>9%</div><div></div></div>
3	Q	11	<div><div></div><div>82%</div><div>18%</div><div></div><div></div></div>
3	R	11	<div><div></div><div>45%</div><div>55%</div><div>45%</div><div></div></div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 10085 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 Light chain of HIV-1 gp120 V3-specific human monoclonal antibody 2424.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	215	Total	C	N	O	S	0	0	0
			1654	1038	277	333	6			
1	M	215	Total	C	N	O	S	0	0	0
			1654	1038	277	333	6			
1	N	213	Total	C	N	O	S	0	0	0
			1635	1028	272	329	6			

- Molecule 2 is a protein called Heavy chain of HIV-1 gp120 V3-specific human monoclonal antibody 2424.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	217	Total	C	N	O	S	0	0	0
			1629	1032	270	320	7			
2	I	217	Total	C	N	O	S	0	0	0
			1629	1032	270	320	7			
2	J	216	Total	C	N	O	S	0	0	0
			1623	1029	269	318	7			

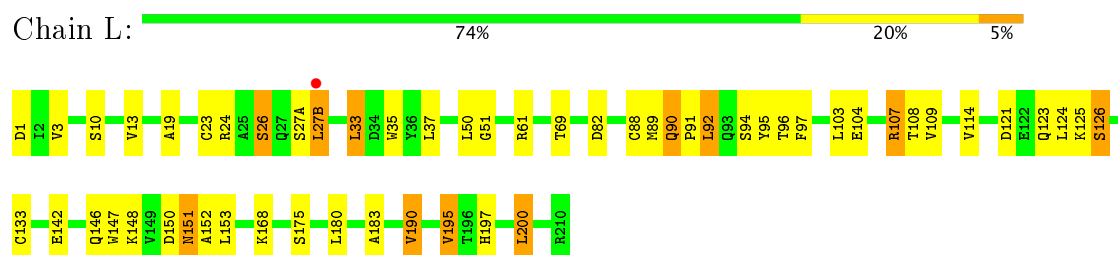
- Molecule 3 is a protein called HIV-1 JR-FL gp120 V3 peptide.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	P	11	Total	C	N	O	0	0	0
			87	58	16	13			
3	Q	11	Total	C	N	O	0	0	0
			87	58	16	13			
3	R	11	Total	C	N	O	0	0	0
			87	58	16	13			

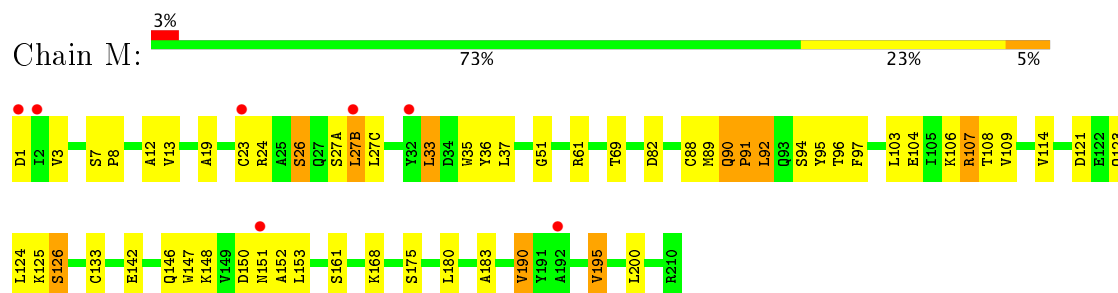
### 3 Residue-property plots [i](#)

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 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.

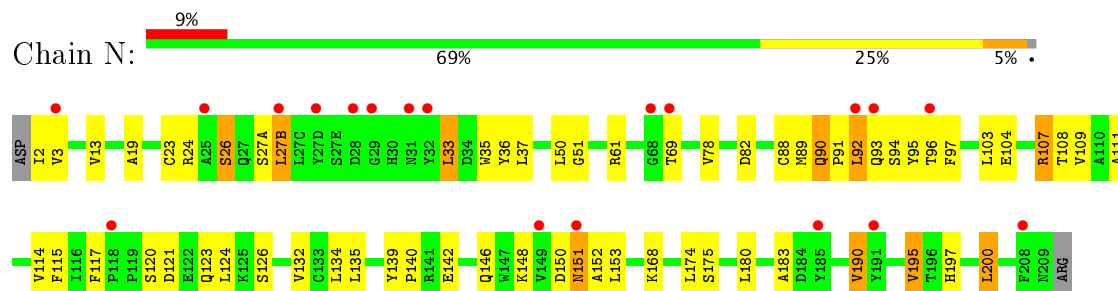
- Molecule 1: Light chain of HIV-1 gp120 V3-specific human monoclonal antibody 2424



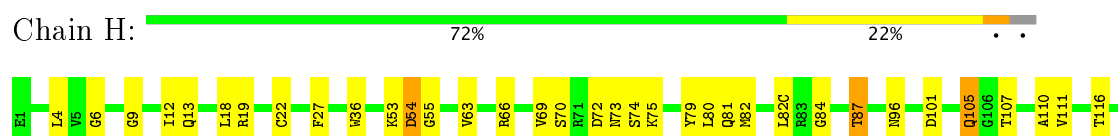
- Molecule 1: Light chain of HIV-1 gp120 V3-specific human monoclonal antibody 2424



- Molecule 1: Light chain of HIV-1 gp120 V3-specific human monoclonal antibody 2424

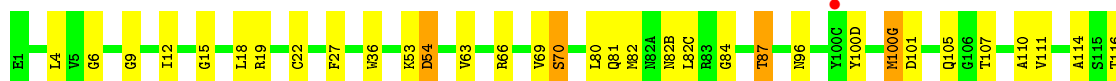


- Molecule 2: Heavy chain of HIV-1 gp120 V3-specific human monoclonal antibody 2424

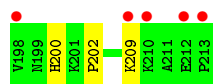
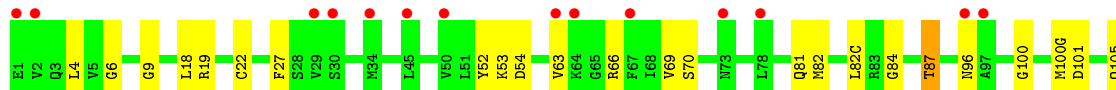




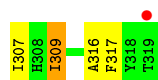
- Molecule 2: Heavy chain of HIV-1 gp120 V3-specific human monoclonal antibody 2424



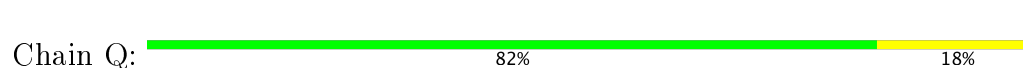
- Molecule 2: Heavy chain of HIV-1 gp120 V3-specific human monoclonal antibody 2424



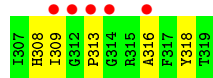
- Molecule 3: HIV-1 JR-FL gp120 V3 peptide



- Molecule 3: HIV-1 JR-FL gp120 V3 peptide



- Molecule 3: HIV-1 JR-FL gp120 V3 peptide



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 2 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	94.95Å 122.25Å 140.81Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.28 – 3.18 48.28 – 3.18	Depositor EDS
% Data completeness (in resolution range)	100.0 (48.28-3.18) 91.6 (48.28-3.18)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.85 (at 3.19Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.9_1692)	Depositor
R, $R_{free}$	0.225 , 0.283 0.231 , 0.292	Depositor DCC
$R_{free}$ test set	1342 reflections (5.04%)	DCC
Wilson B-factor (Å <sup>2</sup> )	61.5	Xtriage
Anisotropy	0.465	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 62.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.46$ , $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.89	EDS
Total number of atoms	10085	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	69.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.84% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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 [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	L	0.49	0/1691	0.77	3/2298 (0.1%)
1	M	0.48	0/1691	0.65	2/2298 (0.1%)
1	N	0.43	0/1672	0.63	2/2273 (0.1%)
2	H	0.52	0/1669	0.64	0/2272
2	I	0.51	0/1669	0.63	0/2272
2	J	0.44	0/1663	0.59	0/2264
3	P	0.32	0/90	0.47	0/121
3	Q	0.41	0/90	0.49	0/121
3	R	0.36	0/90	0.46	0/121
All	All	0.48	0/10325	0.65	7/14040 (0.0%)

There are no bond length outliers.

The worst 5 of 7 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	L	24	ARG	NE-CZ-NH2	15.54	128.07	120.30
1	L	24	ARG	NE-CZ-NH1	-15.34	112.63	120.30
1	M	24	ARG	NE-CZ-NH2	-8.96	115.82	120.30
1	N	24	ARG	NE-CZ-NH1	8.74	124.67	120.30
1	M	24	ARG	NE-CZ-NH1	8.70	124.65	120.30

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	L	1654	0	1609	23	0
1	M	1654	0	1609	27	0
1	N	1635	0	1589	35	0
2	H	1629	0	1575	36	0
2	I	1629	0	1575	42	0
2	J	1623	0	1570	36	0
3	P	87	0	84	2	0
3	Q	87	0	84	1	0
3	R	87	0	84	5	0
All	All	10085	0	9779	174	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 9.

The worst 5 of 174 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:H:84:GLY:O	2:H:87:THR:HG23	1.89	0.72
2:J:84:GLY:O	2:J:87:THR:HG23	1.90	0.71
1:N:3:VAL:H	1:N:26:SER:HB3	1.57	0.69
2:H:9:GLY:HA2	2:H:18:LEU:HD21	1.74	0.69
1:N:150:ASP:HA	1:N:190:VAL:HG23	1.75	0.68

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	L	213/215 (99%)	197 (92%)	9 (4%)	7 (3%)	4	28
1	M	213/215 (99%)	198 (93%)	8 (4%)	7 (3%)	4	28
1	N	211/215 (98%)	197 (93%)	8 (4%)	6 (3%)	6	33

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	H	213/223 (96%)	192 (90%)	18 (8%)	3 (1%)	13	52
2	I	213/223 (96%)	194 (91%)	17 (8%)	2 (1%)	20	63
2	J	212/223 (95%)	193 (91%)	19 (9%)	0	100	100
3	P	9/11 (82%)	8 (89%)	1 (11%)	0	100	100
3	Q	9/11 (82%)	8 (89%)	1 (11%)	0	100	100
3	R	9/11 (82%)	8 (89%)	1 (11%)	0	100	100
All	All	1302/1347 (97%)	1195 (92%)	82 (6%)	25 (2%)	9	44

5 of 25 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	L	126	SER
1	L	151	ASN
1	M	126	SER
1	N	126	SER
1	M	151	ASN

### 5.3.2 Protein sidechains ⓘ

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	L	188/188 (100%)	166 (88%)	22 (12%)	6	27
1	M	188/188 (100%)	167 (89%)	21 (11%)	7	29
1	N	186/188 (99%)	166 (89%)	20 (11%)	7	30
2	H	179/184 (97%)	163 (91%)	16 (9%)	11	40
2	I	179/184 (97%)	164 (92%)	15 (8%)	13	44
2	J	178/184 (97%)	164 (92%)	14 (8%)	14	48
3	P	8/8 (100%)	6 (75%)	2 (25%)	1	2
3	Q	8/8 (100%)	7 (88%)	1 (12%)	5	24
3	R	8/8 (100%)	7 (88%)	1 (12%)	5	24
All	All	1122/1140 (98%)	1010 (90%)	112 (10%)	9	33

5 of 112 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	M	107	ARG
2	I	70	SER
2	J	161	SER
1	M	121	ASP
1	M	175	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 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 [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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	L	215/215 (100%)	0.15	1 (0%) 90 85	38, 60, 80, 97	0
1	M	215/215 (100%)	0.38	7 (3%) 47 30	43, 66, 88, 95	0
1	N	213/215 (99%)	0.73	19 (8%) 10 6	48, 83, 113, 120	0
2	H	217/223 (97%)	0.11	0 100 100	32, 49, 69, 84	0
2	I	217/223 (97%)	0.44	9 (4%) 38 23	41, 59, 93, 109	0
2	J	216/223 (96%)	0.97	28 (12%) 4 2	51, 90, 116, 128	0
3	P	11/11 (100%)	0.77	1 (9%) 10 5	62, 74, 84, 102	0
3	Q	11/11 (100%)	0.57	0 100 100	68, 78, 85, 88	0
3	R	11/11 (100%)	1.99	5 (45%) 0 0	88, 100, 105, 110	0
All	All	1326/1347 (98%)	0.48	70 (5%) 27 15	32, 66, 106, 128	0

The worst 5 of 70 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	I	188	SER	5.2
3	P	319	THR	4.2
1	N	32	TYR	3.9
2	J	1	GLU	3.8
1	N	185	TYR	3.8

### 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 carbohydrates in this entry.

## 6.4 Ligands [i](#)

There are no ligands in this entry.

## 6.5 Other polymers [i](#)

There are no such residues in this entry.