wwPDB X-ray Structure Validation Summary Report (i)

Oct 12, 2024 – 05:51 PM EDT

PDB ID	:	1FVC
Title	:	X-RAY STRUCTURES OF THE ANTIGEN-BINDING DOMAINS FROM THREE VARIANTS OF HUMANIZED ANTI-P185-HER2 ANTIBODY 4D5 AND COMPARISON WITH MOLECULAR MODELING
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Deposited on	:	1992-10-20
Resolution	:	2.20 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 (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:

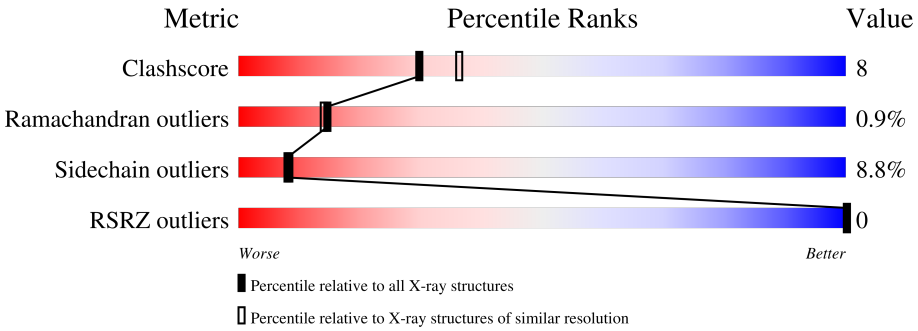
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.20.1
EDS	:	3.0
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.003 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.39

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 2.20 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	180529	6634 (2.20-2.20)
Ramachandran outliers	177936	6560 (2.20-2.20)
Sidechain outliers	177891	6561 (2.20-2.20)
RSRZ outliers	164620	5791 (2.20-2.20)

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	109	74% 22%
1	C	109	66% 28%
2	B	120	63% 28% 6%
2	D	120	65% 26% 9%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 3647 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 IGG1-KAPPA 4D5 FV (LIGHT CHAIN).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	109	Total	C	N	O	S	0	0	0
			835	526	139	167	3			
1	C	108	Total	C	N	O	S	4	0	0
			828	522	138	165	3			

There are 28 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	28	ASP	SER	conflict	GB 185985
A	29	VAL	ILE	conflict	GB 185985
A	30	ASN	SER	conflict	GB 185985
A	31	THR	SER	conflict	GB 185985
A	32	ALA	TYR	conflict	GB 185985
A	33	VAL	LEU	conflict	GB 185985
A	34	ALA	ASN	conflict	GB 185985
A	50	SER	ALA	conflict	GB 185985
A	53	PHE	SER	conflict	GB 185985
A	55	TYR	GLN	conflict	GB 185985
A	66	ARG	GLY	conflict	GB 185985
A	91	HIS	SER	conflict	GB 185985
A	93	THR	SER	conflict	GB 185985
A	?	-	TRP	deletion	GB 185985
C	28	ASP	SER	conflict	GB 185985
C	29	VAL	ILE	conflict	GB 185985
C	30	ASN	SER	conflict	GB 185985
C	31	THR	SER	conflict	GB 185985
C	32	ALA	TYR	conflict	GB 185985
C	33	VAL	LEU	conflict	GB 185985
C	34	ALA	ASN	conflict	GB 185985
C	50	SER	ALA	conflict	GB 185985
C	53	PHE	SER	conflict	GB 185985
C	55	TYR	GLN	conflict	GB 185985
C	66	ARG	GLY	conflict	GB 185985

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Chain	Residue	Modelled	Actual	Comment	Reference
C	91	HIS	SER	conflict	GB 185985
C	93	THR	SER	conflict	GB 185985
C	?	-	TRP	deletion	GB 185985

- Molecule 2 is a protein called IGG1-KAPPA 4D5 FV (HEAVY CHAIN).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	120	Total	C	N	O	S	4	0	0
			929	586	160	179	4			
2	D	120	Total	C	N	O	S	0	0	0
			929	586	160	179	4			

- Molecule 3 is water.

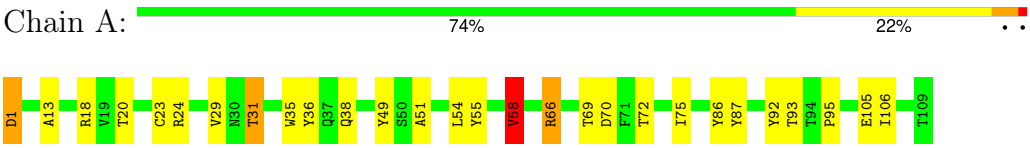
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	47	Total	O	0	0
			47	47		
3	B	23	Total	O	0	0
			23	23		
3	C	31	Total	O	0	0
			31	31		
3	D	25	Total	O	0	0
			25	25		

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.

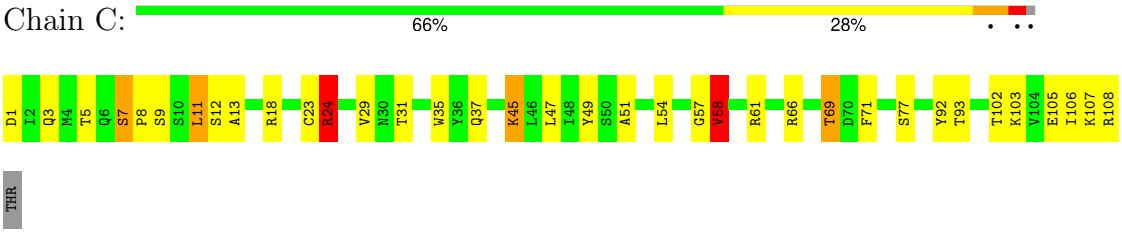
• Molecule 1: IGG1-KAPPA 4D5 FV (LIGHT CHAIN)

Chain A:



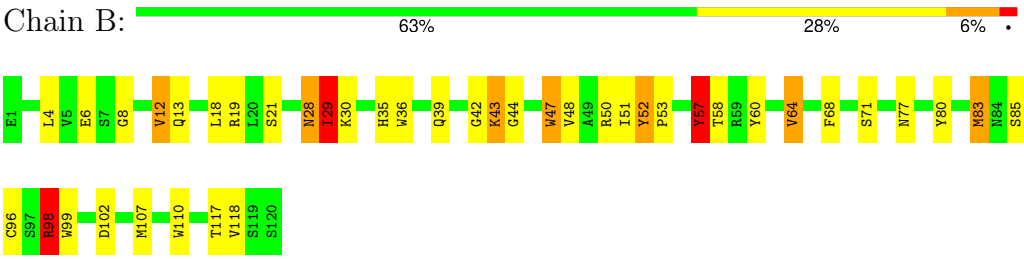
• Molecule 1: IGG1-KAPPA 4D5 FV (LIGHT CHAIN)

Chain C:



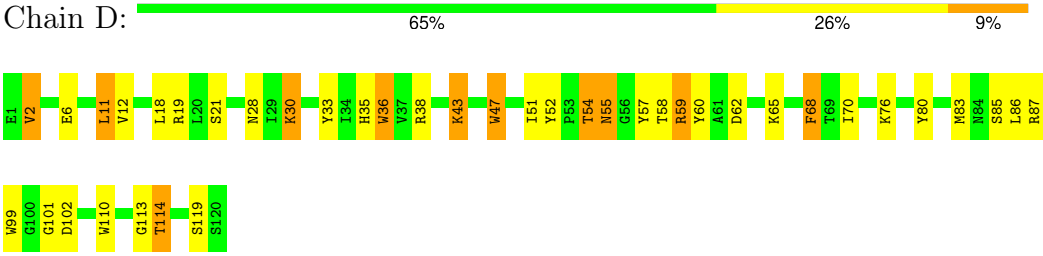
• Molecule 2: IGG1-KAPPA 4D5 FV (HEAVY CHAIN)

Chain B:



• Molecule 2: IGG1-KAPPA 4D5 FV (HEAVY CHAIN)

Chain D:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	37.60Å 63.40Å 90.20Å 90.00° 98.20° 90.00°	Depositor
Resolution (Å)	10.00 – 2.20 10.00 – 2.20	Depositor EDS
% Data completeness (in resolution range)	(Not available) (10.00-2.20) 79.6 (10.00-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$	-	Xtriage
Refinement program	X-PLOR	Depositor
R, R_{free}	0.183 , (Not available) 0.174 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å^2)	25.4	Xtriage
Anisotropy	0.303	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.29 , 77.8	EDS
L-test for twinning[1]	$< \|L\| > = 0.45$, $< L^2 > = 0.27$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	3647	wwPDB-VP
Average B, all atoms (Å^2)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 11.00% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]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

5.1 Standard geometry

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	0.89	1/855 (0.1%)	1.71	21/1164 (1.8%)
1	C	0.86	0/848	1.70	18/1154 (1.6%)
2	B	0.89	0/952	1.84	27/1291 (2.1%)
2	D	0.96	0/952	1.80	24/1291 (1.9%)
All	All	0.90	1/3607 (0.0%)	1.77	90/4900 (1.8%)

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
2	B	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	58	VAL	CB-CG1	-5.04	1.42	1.52

The worst 5 of 90 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	11	LEU	CA-CB-CG	10.58	139.63	115.30
1	A	35	TRP	CD1-CG-CD2	10.45	114.66	106.30
2	D	57	TYR	CB-CG-CD2	-9.61	115.23	121.00
2	D	47	TRP	CD1-CG-CD2	9.04	113.53	106.30
1	A	35	TRP	CG-CD2-CE3	8.91	141.91	133.90

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	57	TYR	Sidechain

5.2 Too-close contacts

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	835	0	810	9	0
1	C	828	0	803	12	0
2	B	929	0	885	21	0
2	D	929	0	885	18	0
3	A	47	0	0	1	0
3	B	23	0	0	0	0
3	C	31	0	0	0	0
3	D	25	0	0	0	0
All	All	3647	0	3383	58	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 8.

The worst 5 of 58 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:D:51:ILE:HG13	2:D:58:THR:HG22	1.63	0.80
2:D:35:HIS:HD2	2:D:47:TRP:HE1	1.30	0.77
2:B:35:HIS:HD2	2:B:47:TRP:HE1	1.32	0.75
1:A:38:GLN:HE22	2:B:39:GLN:HE22	1.42	0.68
2:D:11:LEU:HD21	2:D:119:SER:HB3	1.78	0.64

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	107/109 (98%)	102 (95%)	4 (4%)	1 (1%)	14 14
1	C	106/109 (97%)	98 (92%)	7 (7%)	1 (1%)	14 14
2	B	118/120 (98%)	112 (95%)	6 (5%)	0	100 100
2	D	118/120 (98%)	110 (93%)	6 (5%)	2 (2%)	7 5
All	All	449/458 (98%)	422 (94%)	23 (5%)	4 (1%)	14 14

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	51	ALA
2	D	2	VAL
2	D	102	ASP
1	A	51	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	94/95 (99%)	86 (92%)	8 (8%)	8 9
1	C	93/95 (98%)	82 (88%)	11 (12%)	4 4
2	B	95/95 (100%)	88 (93%)	7 (7%)	11 12
2	D	95/95 (100%)	88 (93%)	7 (7%)	11 12
All	All	377/380 (99%)	344 (91%)	33 (9%)	8 8

5 of 33 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	D	43	LYS
2	D	54	THR
2	D	114	THR
2	B	85	SER

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Mol	Chain	Res	Type
2	B	71	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 6 such sidechains are listed below:

Mol	Chain	Res	Type
2	D	3	GLN
2	D	35	HIS
2	D	55	ASN
2	B	35	HIS
2	B	28	ASN

5.3.3 RNA

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains

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

5.5 Carbohydrates

There are no oligosaccharides in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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^{th} 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	109/109 (100%)	-0.94	0	100	100	7, 18, 39, 70	0
1	C	108/109 (99%)	-0.91	0	100	100	7, 19, 36, 47	1 (0%)
2	B	120/120 (100%)	-0.66	0	100	100	10, 25, 48, 72	1 (0%)
2	D	120/120 (100%)	-0.72	0	100	100	8, 26, 51, 65	0
All	All	457/458 (99%)	-0.80	0	100	100	7, 22, 46, 72	2 (0%)

There are no RSRZ outliers to report.

6.2 Non-standard residues in protein, DNA, RNA chains

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

6.3 Carbohydrates

There are no monosaccharides in this entry.

6.4 Ligands

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.