



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2018 – 06:37 am GMT

PDB ID : 1A7Q
Title : FV FRAGMENT OF MOUSE MONOCLONAL ANTIBODY D1.3 (BALB/C, IGG1, K) HIGH AFFINITY EXPRESSED VARIANT CONTAINING SER26L->GLY, ILE29L->THR, GLU81L->ASP, THR97L->SER, PRO240H->LEU, ASP258H->ALA, LYS281H->GLU, ASN283H->ASP AND LEU312H->VAL
Authors : Marks, C.; Henrick, K.; Winter, G.
Deposited on : 1998-03-16
Resolution : 2.00 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 ⓘ symbol.

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

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	trunk30967
Percentile statistics	:	20171227.v01 (using entries in the PDB archive December 27th 2017)
Refmac	:	5.8.0158
CCP4	:	7.0 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk30967

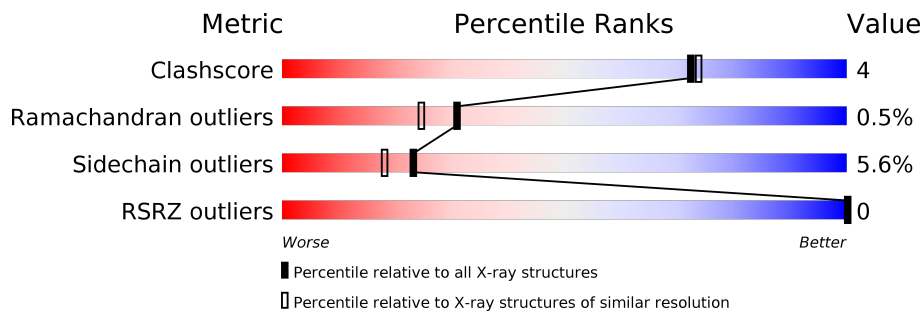
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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)	Similar resolution (#Entries, resolution range(Å))
Clashscore	122126	8267 (2.00-2.00)
Ramachandran outliers	120053	8166 (2.00-2.00)
Sidechain outliers	120020	8165 (2.00-2.00)
RSRZ outliers	108989	7011 (2.00-2.00)

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 ≥ 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	106	
2	H	116	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 1728 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 D1.3 FV (LIGHT CHAIN).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	106	Total	C	N	O	S	0	0	0
			791	503	128	158	2			

There are 11 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
L	3	VAL	GLU	CONFLICT	GB 545862
L	26	GLY	SER	SOMATIC VARIANT	GB 545862
L	29	THR	ILE	SOMATIC VARIANT	GB 545862
L	50	TYR	LYS	CONFLICT	GB 545862
L	51	THR	ALA	CONFLICT	GB 545862
L	52	THR	GLN	CONFLICT	GB 545862
L	56	ALA	ASP	CONFLICT	GB 545862
L	81	ASP	GLU	VARIANT	GB 545862
L	?	-	PRO	DELETION	GB 545862
L	96	ARG	TRP	CONFLICT	GB 545862
L	97	SER	THR	SOMATIC VARIANT	GB 545862

- Molecule 2 is a protein called IGG1-KAPPA D1.3 FV (HEAVY CHAIN).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	116	Total	C	N	O	S	0	0	0
			890	558	153	175	4			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
H	240	LEU	PRO	SOMATIC VARIANT	GB 896294
H	258	ALA	ASP	SOMATIC VARIANT	GB 896294
H	281	GLU	LYS	SOMATIC VARIANT	GB 896294
H	283	ASP	ASN	SOMATIC VARIANT	GB 896294

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	L	23	Total 23	O 23	0	0
3	H	24	Total 24	O 24	0	0

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 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.

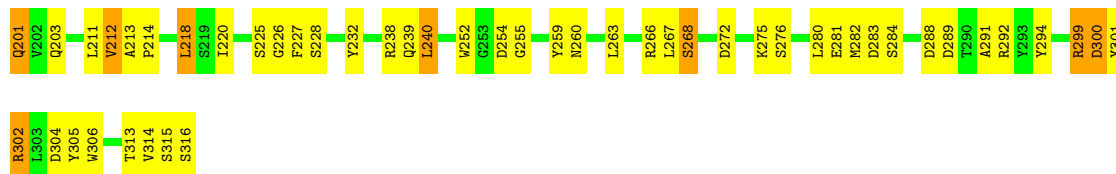
- Molecule 1: IGG1-KAPPA D1.3 FV (LIGHT CHAIN)

Chain L: 



- Molecule 2: IGG1-KAPPA D1.3 FV (HEAVY CHAIN)

Chain H: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	90.31Å 90.31Å 56.85Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	6.00 – 2.00 13.69 – 2.00	Depositor EDS
% Data completeness (in resolution range)	(Not available) (6.00-2.00) 97.5 (13.69-2.00)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.81 (at 2.01Å)	Xtriage
Refinement program	CCP4	Depositor
R, R_{free}	0.178 , (Not available) 0.177 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	24.7	Xtriage
Anisotropy	0.087	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 68.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	1728	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.09% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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

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	L	1.20	0/812	2.19	31/1108 (2.8%)
2	H	1.19	0/909	2.71	64/1236 (5.2%)
All	All	1.19	0/1721	2.48	95/2344 (4.1%)

There are no bond length outliers.

All (95) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	302	ARG	NE-CZ-NH2	-22.58	109.01	120.30
2	H	266	ARG	NE-CZ-NH2	-20.29	110.16	120.30
2	H	302	ARG	NE-CZ-NH1	19.53	130.06	120.30
2	H	266	ARG	NE-CZ-NH1	15.16	127.88	120.30
1	L	96	ARG	NE-CZ-NH2	-14.07	113.27	120.30
2	H	238	ARG	NE-CZ-NH2	-13.45	113.58	120.30
2	H	304	ASP	CB-CG-OD1	13.36	130.32	118.30
2	H	268	SER	N-CA-CB	13.31	130.47	110.50
1	L	96	ARG	NE-CZ-NH1	12.75	126.68	120.30
2	H	218	LEU	CA-CB-CG	11.18	141.01	115.30
2	H	299	ARG	NE-CZ-NH2	-10.32	115.14	120.30
2	H	238	ARG	NE-CZ-NH1	10.15	125.37	120.30
2	H	284	SER	N-CA-CB	-10.05	95.42	110.50
2	H	305	TYR	CB-CG-CD1	-9.97	115.02	121.00
1	L	82	ASP	CB-CG-OD1	9.80	127.12	118.30
2	H	272	ASP	CB-CG-OD2	-9.61	109.65	118.30
2	H	232	TYR	CB-CG-CD2	-9.53	115.28	121.00
2	H	268	SER	CB-CA-C	-9.16	92.70	110.10
1	L	1	ASP	CB-CG-OD2	-9.14	110.07	118.30
2	H	201	GLN	N-CA-CB	-8.55	95.21	110.60
2	H	283	ASP	CB-CG-OD2	-8.39	110.75	118.30
2	H	306	TRP	CE3-CZ3-CH2	-8.32	112.05	121.20
2	H	259	TYR	CB-CG-CD1	-8.20	116.08	121.00
2	H	304	ASP	CB-CG-OD2	-8.12	110.99	118.30
1	L	10	SER	CB-CA-C	-7.73	95.42	110.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	300	ASP	CB-CG-OD2	7.67	125.20	118.30
2	H	315	SER	N-CA-CB	-7.67	99.00	110.50
1	L	83	PHE	CB-CA-C	7.50	125.40	110.40
2	H	201	GLN	O-C-N	-7.41	110.84	122.70
2	H	299	ARG	CD-NE-CZ	-7.32	113.35	123.60
2	H	232	TYR	CB-CG-CD1	7.22	125.33	121.00
2	H	254	ASP	CB-CG-OD2	-7.22	111.80	118.30
2	H	292	ARG	NE-CZ-NH2	-7.21	116.69	120.30
1	L	3	VAL	CA-CB-CG2	7.20	121.70	110.90
1	L	86	TYR	CB-CG-CD2	-7.10	116.74	121.00
1	L	71	TYR	CB-CG-CD1	6.84	125.10	121.00
1	L	50	TYR	CB-CG-CD2	-6.78	116.93	121.00
2	H	240	LEU	CA-CB-CG	6.62	130.52	115.30
2	H	299	ARG	NH1-CZ-NH2	6.56	126.62	119.40
1	L	11	LEU	CA-CB-CG	6.54	130.35	115.30
2	H	281	GLU	OE1-CD-OE2	6.52	131.12	123.30
1	L	4	LEU	O-C-N	6.51	133.12	122.70
2	H	294	TYR	CB-CG-CD2	-6.41	117.15	121.00
1	L	1	ASP	CB-CG-OD1	6.36	124.02	118.30
2	H	302	ARG	CD-NE-CZ	-6.18	114.94	123.60
2	H	213	ALA	N-CA-CB	-6.18	101.45	110.10
2	H	263	LEU	N-CA-CB	-6.14	98.11	110.40
2	H	301	TYR	CB-CG-CD1	-6.04	117.38	121.00
2	H	212	VAL	CG1-CB-CG2	-6.02	101.26	110.90
1	L	53	THR	CA-CB-OG1	-6.00	96.40	109.00
2	H	227	PHE	CA-C-O	6.00	132.70	120.10
2	H	201	GLN	CA-CB-CG	-5.96	100.28	113.40
2	H	211	LEU	CA-CB-CG	5.95	128.98	115.30
1	L	17	GLU	OE1-CD-OE2	5.94	130.42	123.30
1	L	103	LYS	CA-CB-CG	5.80	126.16	113.40
1	L	79	GLN	CB-CA-C	5.75	121.89	110.40
2	H	275	LYS	CB-CA-C	5.74	121.89	110.40
1	L	86	TYR	CB-CG-CD1	5.73	124.44	121.00
2	H	306	TRP	CZ3-CH2-CZ2	5.72	128.46	121.60
2	H	201	GLN	CB-CA-C	5.65	121.70	110.40
2	H	201	GLN	CA-C-N	5.64	129.61	117.20
1	L	84	GLY	O-C-N	5.63	131.70	122.70
2	H	288	ASP	CB-CG-OD1	-5.61	113.25	118.30
2	H	268	SER	O-C-N	5.60	131.66	122.70
2	H	289	ASP	CB-CG-OD1	-5.56	113.29	118.30
1	L	52	THR	CA-C-O	-5.53	108.48	120.10
1	L	18	THR	CA-CB-OG1	-5.48	97.50	109.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	304	ASP	CA-C-O	-5.46	108.64	120.10
2	H	283	ASP	CA-C-N	-5.45	105.21	117.20
2	H	239	GLN	OE1-CD-NE2	5.44	134.42	121.90
1	L	33	LEU	CD1-CG-CD2	5.43	126.80	110.50
1	L	61	ARG	N-CA-CB	5.38	120.29	110.60
2	H	228	SER	N-CA-CB	-5.36	102.46	110.50
2	H	313	THR	CA-CB-OG1	-5.35	97.76	109.00
2	H	304	ASP	O-C-N	5.32	131.22	122.70
1	L	71	TYR	CB-CG-CD2	-5.31	117.81	121.00
2	H	282	MET	CA-C-O	5.31	131.25	120.10
1	L	21	ILE	CB-CA-C	5.31	122.21	111.60
1	L	82	ASP	OD1-CG-OD2	-5.26	113.30	123.30
2	H	260	ASN	N-CA-CB	-5.26	101.14	110.60
2	H	255	GLY	CA-C-O	-5.24	111.16	120.60
2	H	203	GLN	CB-CG-CD	5.23	125.21	111.60
1	L	52	THR	CA-C-N	5.21	128.67	117.20
2	H	203	GLN	CB-CA-C	5.21	120.81	110.40
2	H	254	ASP	CB-CG-OD1	5.20	122.98	118.30
1	L	42	LYS	CB-CA-C	5.15	120.71	110.40
1	L	77	SER	CA-C-O	5.15	130.91	120.10
1	L	36	TYR	CB-CA-C	-5.14	100.11	110.40
2	H	288	ASP	CB-CA-C	5.11	120.62	110.40
2	H	252	TRP	CD2-CE3-CZ3	-5.11	112.16	118.80
2	H	225	SER	CB-CA-C	-5.08	100.44	110.10
2	H	272	ASP	CB-CG-OD1	5.02	122.82	118.30
2	H	268	SER	CA-C-N	-5.01	106.17	117.20
1	L	80	PRO	C-N-CA	5.01	134.22	121.70
2	H	252	TRP	CG-CD2-CE3	-5.00	129.40	133.90

There are no chirality outliers.

There are no planarity outliers.

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	L	791	0	735	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	H	890	0	842	9	0
3	H	24	0	0	0	0
3	L	23	0	0	0	0
All	All	1728	0	1577	12	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 4.

All (12) 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:L:48:VAL:HG22	1:L:54:LEU:HD23	1.68	0.74
2:H:220:ILE:HD12	2:H:280:LEU:HD23	1.86	0.57
2:H:240:LEU:HD22	2:H:291:ALA:HB2	1.89	0.55
2:H:300:ASP:HB2	2:H:302:ARG:NH1	2.28	0.48
2:H:214:PRO:HD2	2:H:316:SER:HB3	1.99	0.44
2:H:299:ARG:HH11	2:H:299:ARG:HD3	1.48	0.43
2:H:299:ARG:NH1	2:H:302:ARG:NH2	2.66	0.43
2:H:212:VAL:O	2:H:314:VAL:HA	2.20	0.42
1:L:43:SER:HA	1:L:44:PRO:HD3	1.84	0.42
2:H:267:LEU:HD22	2:H:280:LEU:HD11	2.03	0.41
1:L:15:VAL:HG21	1:L:80:PRO:HG3	2.02	0.41
2:H:201:GLN:O	2:H:226:GLY:HA3	2.22	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 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	104/106 (98%)	101 (97%)	2 (2%)	1 (1%)	17	10
2	H	114/116 (98%)	110 (96%)	4 (4%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	218/222 (98%)	211 (97%)	6 (3%)	1 (0%)	31	25

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	L	77	SER

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	L	82/88 (93%)	75 (92%)	7 (8%)	12	7
2	H	95/99 (96%)	92 (97%)	3 (3%)	42	41
All	All	177/187 (95%)	167 (94%)	10 (6%)	23	18

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	L	3	VAL
1	L	14	SER
1	L	33	LEU
1	L	70	GLN
1	L	76	ASN
1	L	81	ASP
1	L	96	ARG
2	H	218	LEU
2	H	268	SER
2	H	276	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	L	45	GLN
1	L	89	GLN

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Mol	Chain	Res	Type
2	H	256	ASN
2	H	308	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 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, 95th 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	L	106/106 (100%)	-0.45	0 100 100	13, 22, 32, 38	11 (10%)
2	H	116/116 (100%)	-0.56	0 100 100	12, 20, 29, 32	11 (9%)
All	All	222/222 (100%)	-0.51	0 100 100	12, 21, 31, 38	22 (9%)

There are no RSRZ outliers to report.

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.