



# Full wwPDB X-ray Structure Validation Report ⓘ

May 13, 2020 – 02:32 am BST

PDB ID : 6D9G  
Title : X-ray Structure of the FAB Fragment of 15B8, a Murine Monoclonal Antibody  
Specific for the Human Serotonin Transporter  
Authors : Coleman, J.A.; Yang, D.; Gouaux, E.  
Deposited on : 2018-04-29  
Resolution : 2.30 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](mailto: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.

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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)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

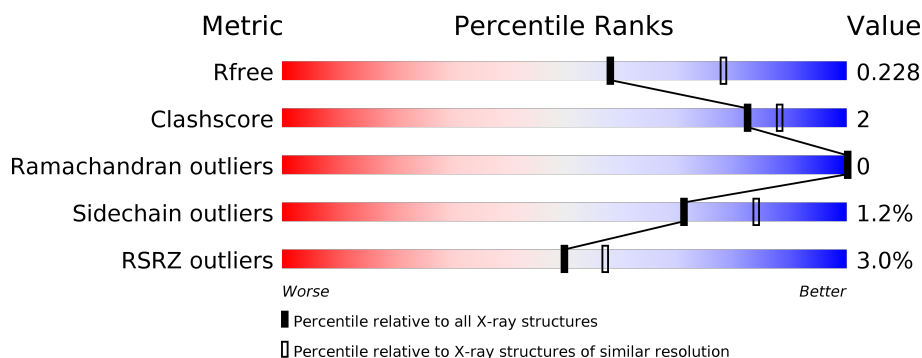
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.30 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))
$R_{free}$	130704	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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 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  $\leq 5\%$ . 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	A	229	<div> <div>2%</div> <div> <div></div> <div>93%</div> <div>6%</div> <div>• • •</div> </div> </div>
1	C	229	<div> <div>3%</div> <div> <div></div> <div>90%</div> <div>6%</div> <div>• •</div> </div> </div>
2	B	216	<div> <div></div> <div> <div></div> <div>96%</div> <div>•</div> </div> </div>
2	D	216	<div> <div>6%</div> <div> <div></div> <div>96%</div> <div>•</div> </div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 13396 atoms, of which 6404 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 Antibody heavy chain Fab.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	222	Total	C	H	N	O	S	0	0	0
			3285	1049	1627	276	325	8			
1	C	221	Total	C	H	N	O	S	0	0	0
			3280	1047	1626	275	324	8			

- Molecule 2 is a protein called Antibody light chain Fab.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	B	216	Total	C	H	N	O	S	0	0	0
			3239	1034	1581	283	335	6			
2	D	216	Total	C	H	N	O	S	0	0	0
			3228	1034	1570	283	335	6			

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	109	Total	O	0	0
			109	109		
3	B	121	Total	O	0	0
			121	121		
3	C	67	Total	O	0	0
			67	67		
3	D	67	Total	O	0	0
			67	67		

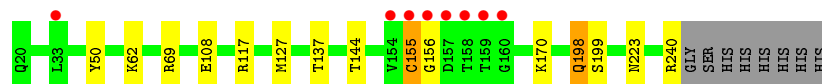
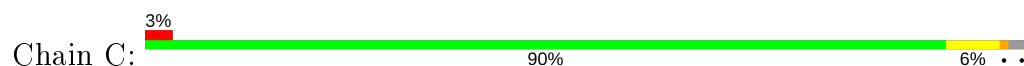
### 3 Residue-property plots [i](#)

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: Antibody heavy chain Fab



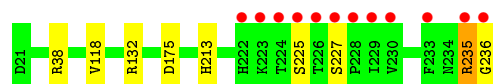
- Molecule 1: Antibody heavy chain Fab



- Molecule 2: Antibody light chain Fab



- Molecule 2: Antibody light chain Fab



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 2 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	82.87Å 85.49Å 141.70Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.93 – 2.30 73.20 – 2.17	Depositor EDS
% Data completeness (in resolution range)	98.4 (19.93-2.30) 91.7 (73.20-2.17)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.17 (at 2.16Å)	Xtriage
Refinement program	PHENIX dev_2597	Depositor
R, $R_{free}$	0.195 , 0.228 0.198 , 0.228	Depositor DCC
$R_{free}$ test set	2487 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	36.0	Xtriage
Anisotropy	0.227	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.41 , 43.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.003 for k,h,-l	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	13396	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.29% 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	A	0.25	0/1700	0.46	0/2313
1	C	0.25	0/1696	0.47	0/2308
2	B	0.26	0/1697	0.45	0/2304
2	D	0.26	0/1697	0.47	0/2304
All	All	0.25	0/6790	0.46	0/9229

There are no bond length outliers.

There are no bond angle outliers.

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	A	1658	1627	1629	6	0
1	C	1654	1626	1626	9	0
2	B	1658	1581	1581	7	0
2	D	1658	1570	1581	11	2
3	A	109	0	0	2	1
3	B	121	0	0	7	1
3	C	67	0	0	5	1
3	D	67	0	0	9	1
All	All	6992	6404	6417	32	4

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 2.

All (32) 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:240:ARG:NH1	3:C:302:HOH:O	1.88	1.03
1:C:144:THR:OG1	3:C:301:HOH:O	1.82	0.95
2:D:236:GLY:CA	3:D:302:HOH:O	2.14	0.94
2:D:236:GLY:N	3:D:302:HOH:O	1.99	0.94
2:B:175:ASP:OD2	3:B:301:HOH:O	1.87	0.93
2:B:189:ASP:OD1	3:B:302:HOH:O	1.88	0.91
1:C:137:THR:OG1	3:C:303:HOH:O	1.89	0.90
2:B:134:ASP:OD2	3:B:303:HOH:O	1.92	0.88
2:D:132:ARG:O	3:D:301:HOH:O	1.93	0.85
1:A:20:GLN:N	3:A:301:HOH:O	2.14	0.79
1:A:160:GLY:O	1:A:212:SER:OG	2.00	0.79
1:C:170:LYS:NZ	3:C:304:HOH:O	2.17	0.76
2:D:213:HIS:HD2	3:D:303:HOH:O	1.72	0.70
2:D:175:ASP:OD2	3:D:303:HOH:O	2.09	0.70
1:A:179:LEU:O	3:A:302:HOH:O	2.14	0.66
1:C:198:GLN:O	1:C:199:SER:OG	2.14	0.65
2:D:235:ARG:NH2	3:D:306:HOH:O	2.31	0.64
2:D:236:GLY:HA2	3:D:302:HOH:O	1.91	0.61
2:D:132:ARG:NH2	3:D:307:HOH:O	2.35	0.58
1:C:50:TYR:O	3:C:305:HOH:O	2.18	0.56
1:A:158:THR:O	1:A:160:GLY:N	2.40	0.53
2:B:147:GLU:OE2	3:B:305:HOH:O	2.19	0.52
1:A:127:MET:HG3	3:B:321:HOH:O	2.14	0.47
2:D:235:ARG:HG2	2:D:236:GLY:H	1.80	0.47
1:C:69:ARG:NH2	2:D:118:VAL:O	2.47	0.46
1:A:157:ASP:O	1:A:158:THR:OG1	2.31	0.46
2:B:116:LYS:NZ	3:B:311:HOH:O	2.48	0.44
2:B:204:THR:HG23	3:B:315:HOH:O	2.18	0.44
1:C:117:ARG:O	1:C:127:MET:HA	2.19	0.41
1:C:155:CYS:SG	1:C:156:GLY:N	2.93	0.41
2:B:74:ALA:O	2:B:75:ALA:HB3	2.20	0.41
2:D:38:ARG:NH2	3:D:304:HOH:O	2.23	0.40

All (4) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:225:SER:O	2:D:227:SER:N[2_555]	1.83	0.37
2:D:225:SER:O	2:D:227:SER:H[2_555]	1.35	0.25

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:318:HOH:O	3:D:360:HOH:O[3_445]	1.98	0.22
3:A:337:HOH:O	3:B:369:HOH:O[3_545]	2.07	0.13

## 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	A	220/229 (96%)	217 (99%)	3 (1%)	0	100	100
1	C	219/229 (96%)	210 (96%)	9 (4%)	0	100	100
2	B	214/216 (99%)	207 (97%)	7 (3%)	0	100	100
2	D	214/216 (99%)	201 (94%)	13 (6%)	0	100	100
All	All	867/890 (97%)	835 (96%)	32 (4%)	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 sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	186/193 (96%)	183 (98%)	3 (2%)	62	78
1	C	186/193 (96%)	181 (97%)	5 (3%)	44	61
2	B	186/186 (100%)	186 (100%)	0	100	100
2	D	186/186 (100%)	185 (100%)	1 (0%)	88	95
All	All	744/758 (98%)	735 (99%)	9 (1%)	71	84



All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	20	GLN
1	A	157	ASP
1	A	223	ASN
1	C	62	LYS
1	C	108	GLU
1	C	155	CYS
1	C	198	GLN
1	C	223	ASN
2	D	235	ARG

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 ⓘ

### 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<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	A	222/229 (96%)	0.03	5 (2%) 60 67	27, 39, 62, 127	0
1	C	221/229 (96%)	0.20	8 (3%) 42 49	30, 49, 86, 167	0
2	B	216/216 (100%)	-0.16	1 (0%) 91 94	26, 37, 55, 91	0
2	D	216/216 (100%)	0.31	12 (5%) 24 30	30, 45, 86, 319	0
All	All	875/890 (98%)	0.09	26 (2%) 50 57	26, 42, 77, 319	0

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	227	SER	13.5
1	C	157	ASP	12.2
2	D	224	THR	12.1
2	D	228	PRO	11.6
1	C	156	GLY	11.0
1	A	156	GLY	10.0
1	C	158	THR	7.5
2	D	229	ILE	7.3
1	C	155	CYS	5.8
2	D	225	SER	5.7
2	D	226	THR	5.7
1	A	157	ASP	5.3
1	A	155	CYS	5.2
2	D	223	LYS	4.9
1	A	159	THR	4.1
1	C	154	VAL	4.0
2	D	233	PHE	3.5
2	D	235	ARG	3.4
2	D	230	VAL	3.1
2	D	222	HIS	2.8
1	C	159	THR	2.4

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Mol	Chain	Res	Type	RSRZ
1	C	160	GLY	2.3
2	B	236	GLY	2.3
1	C	33	LEU	2.3
2	D	236	GLY	2.1
1	A	160	GLY	2.1

## 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.