



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 12, 2019 – 11:44 AM EST

PDB ID : 6DDM  
Title : Crystal Structure Analysis of the Epitope of an Anti-MICA Antibody  
Authors : Matsumoto, M.L.  
Deposited on : 2018-05-10  
Resolution : 1.30 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 : rb-20031633  
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) : rb-20031633

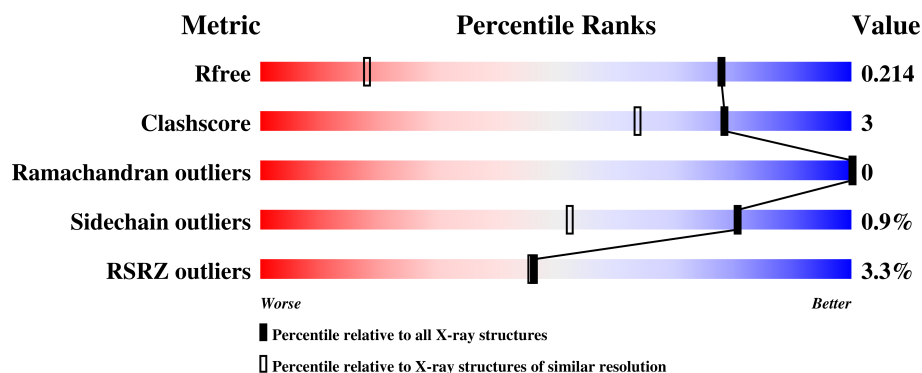
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.30 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))
$R_{free}$	111664	1286 (1.32-1.28)
Clashscore	122126	1332 (1.32-1.28)
Ramachandran outliers	120053	1282 (1.32-1.28)
Sidechain outliers	120020	1282 (1.32-1.28)
RSRZ outliers	108989	1250 (1.32-1.28)

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	A	215	<div> <div>2%</div> <div>93% 7%</div> </div>
2	C	94	<div> <div>3%</div> <div>86% 6% • 6%</div> </div>
3	B	220	<div> <div>5%</div> <div>95% • •</div> </div>

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4811 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 Anti-MICA Fab fragment light chain clone 1D5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	215	Total	C	N	O	S	0	20	0
			1753	1090	286	367	10			

- Molecule 2 is a protein called MHC class I polypeptide-related sequence A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	88	Total	C	N	O	S	0	10	0
			757	467	135	147	8			

- Molecule 3 is a protein called Anti-MICA Fab fragment heavy chain clone 1D5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	B	216	Total	C	N	O	S	0	14	0
			1676	1056	272	341	7			

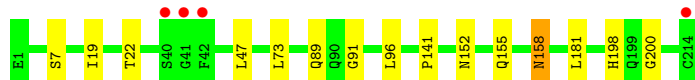
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	261	Total	O	0	1
			262	262		
4	C	112	Total	O	0	0
			112	112		
4	B	245	Total	O	0	6
			251	251		

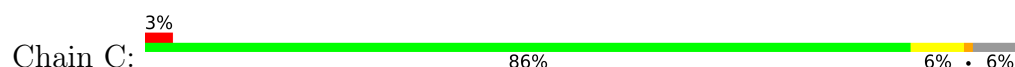
### 3 Residue-property plots [i](#)

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.

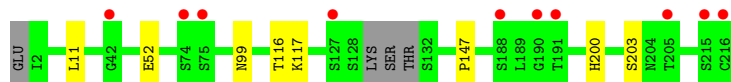
- Molecule 1: Anti-MICA Fab fragment light chain clone 1D5



- Molecule 2: MHC class I polypeptide-related sequence A



- Molecule 3: Anti-MICA Fab fragment heavy chain clone 1D5



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	58.69Å 50.18Å 88.63Å 90.00° 90.85° 90.00°	Depositor
Resolution (Å)	35.00 – 1.30 34.91 – 1.30	Depositor EDS
% Data completeness (in resolution range)	99.6 (35.00-1.30) 99.4 (34.91-1.30)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.64 (at 1.30Å)	Xtriage
Refinement program	REFMAC 5.8.0155	Depositor
R, $R_{free}$	0.188 , 0.205 0.197 , 0.214	Depositor DCC
$R_{free}$ test set	6152 reflections (4.87%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	14.7	Xtriage
Anisotropy	0.648	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.40 , 45.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.020 for h,-k,-l	Xtriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	4811	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.91% 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	0.37	0/1788	0.61	0/2433
2	C	0.38	0/778	0.62	0/1064
3	B	0.35	0/1717	0.58	0/2345
All	All	0.36	0/4283	0.60	0/5842

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	1753	0	1672	15	0
2	C	757	0	679	6	0
3	B	1676	0	1579	5	0
4	A	262	0	0	0	0
4	B	251	0	0	1	0
4	C	112	0	0	0	0
All	All	4811	0	3930	25	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 3.

All (25) 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:A:19[B]:ILE:C	1:A:19[B]:ILE:HD12	2.05	0.77
1:A:181[B]:LEU:N	1:A:181[B]:LEU:HD12	2.09	0.66
1:A:19[B]:ILE:O	1:A:19[B]:ILE:HD12	1.97	0.65
1:A:155:GLN:HB3	1:A:158:ASN:HD21	1.64	0.62
2:C:239:TRP:H	2:C:251:GLN:HE22	1.48	0.61
1:A:181[B]:LEU:HD12	1:A:181[B]:LEU:H	1.64	0.61
3:B:147:PRO:O	3:B:200:HIS:HE1	1.86	0.59
1:A:158:ASN:H	1:A:158:ASN:HD22	1.53	0.55
2:C:207:PRO:HG2	2:C:284[B]:MET:HE1	1.89	0.55
2:C:234:ASN:H	2:C:234:ASN:HD22	1.56	0.52
3:B:200:HIS:HD2	3:B:203:SER:OG	1.93	0.52
2:C:239:TRP:H	2:C:251:GLN:NE2	2.09	0.51
1:A:19[B]:ILE:C	1:A:19[B]:ILE:CD1	2.74	0.51
3:B:117:LYS:NZ	4:B:301:HOH:O	2.44	0.50
1:A:198:HIS:CD2	1:A:200:GLY:H	2.30	0.50
3:B:11:LEU:HD23	3:B:116[B]:THR:HG23	1.97	0.47
1:A:19[B]:ILE:O	1:A:19[B]:ILE:CD1	2.64	0.45
1:A:141:PRO:O	1:A:198:HIS:HE1	2.00	0.44
1:A:7[A]:SER:HG	1:A:22[A]:THR:HG1	1.66	0.44
2:C:284[B]:MET:HE2	2:C:284[B]:MET:HB3	1.91	0.43
1:A:181[B]:LEU:N	1:A:181[B]:LEU:CD1	2.80	0.43
2:C:211:VAL:HG22	2:C:282[B]:CYS:SG	2.60	0.42
1:A:89:GLN:HE22	3:B:99:ASN:ND2	2.18	0.41
1:A:91:GLY:HA2	1:A:96:LEU:HD22	2.01	0.41
1:A:47:LEU:HD23	1:A:73:LEU:HD11	2.04	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	A	233/215 (108%)	231 (99%)	2 (1%)	0	<b>100</b> <b>100</b>

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	C	93/94 (99%)	93 (100%)	0	0	100	100
3	B	225/220 (102%)	221 (98%)	4 (2%)	0	100	100
All	All	551/529 (104%)	545 (99%)	6 (1%)	0	100	100

There are no Ramachandran outliers to report.

### 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	A	202/186 (109%)	200 (99%)	2 (1%)	78	48
2	C	86/84 (102%)	85 (99%)	1 (1%)	74	40
3	B	191/190 (100%)	190 (100%)	1 (0%)	90	68
All	All	479/460 (104%)	475 (99%)	4 (1%)	81	57

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	152	ASN
1	A	158	ASN
2	C	234	ASN
3	B	52	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	124	GLN
1	A	152	ASN
1	A	155	GLN
1	A	158	ASN
1	A	198	HIS
1	A	210	ASN
2	C	234	ASN

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Mol	Chain	Res	Type
2	C	251	GLN
2	C	261	ASN
3	B	6	GLN
3	B	82(A)	ASN
3	B	99	ASN
3	B	200	HIS

### 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	A	215/215 (100%)	0.05	4 (1%) 66 67	12, 18, 28, 36	0
2	C	88/94 (93%)	0.03	3 (3%) 45 44	12, 17, 25, 29	0
3	B	216/220 (98%)	0.25	10 (4%) 32 32	13, 19, 31, 43	0
All	All	519/529 (98%)	0.13	17 (3%) 46 46	12, 19, 29, 43	0

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	B	216	CYS	6.8
1	A	214	CYS	5.0
3	B	42	GLY	3.3
1	A	40	SER	3.3
3	B	215	SER	2.9
3	B	190	GLY	2.8
2	C	282[A]	CYS	2.8
3	B	74	SER	2.8
1	A	42	PHE	2.7
3	B	205	THR	2.7
2	C	214	SER	2.6
3	B	75	SER	2.4
3	B	188[A]	SER	2.4
3	B	191	THR	2.4
2	C	271	ARG	2.3
1	A	41	GLY	2.3
3	B	127	SER	2.3

### 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.