



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 7, 2018 – 01:56 PM EST

PDB ID : 6AYZ  
Title : Crystal structure of Asf1-Fab 12E complex  
Authors : Bailey, L.J.; Kossiakoff, A.A.  
Deposited on : 2017-09-08  
Resolution : 2.10 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

<http://wwpdb.org/validation/2016/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) : 1.9-1692  
EDS : rb-20030736  
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-20030736

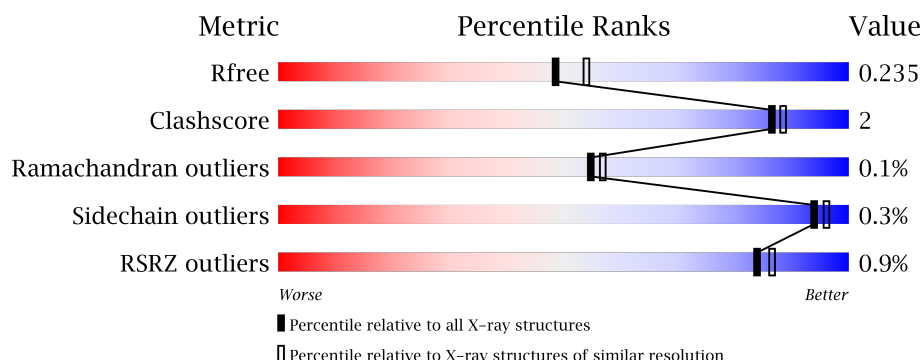
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.10 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	4243 (2.10-2.10)
Clashscore	112137	4788 (2.10-2.10)
Ramachandran outliers	110173	4740 (2.10-2.10)
Sidechain outliers	110143	4741 (2.10-2.10)
RSRZ outliers	101464	4275 (2.10-2.10)

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	B	224	<div> <div>95%</div> <div> <div></div> <div></div> <div></div> <div></div> </div> <div>..</div> </div>
1	R	224	<div> <div>94%</div> <div> <div></div> <div></div> <div></div> <div></div> </div> <div>5%</div> </div>
2	C	215	<div> <div>2%</div> <div> <div></div> <div></div> <div></div> <div></div> </div> <div>91%</div> <div>8%</div> <div>.</div> </div>
2	D	215	<div> <div>87%</div> <div> <div></div> <div></div> <div></div> <div></div> </div> <div>10%</div> <div>..</div> </div>
3	A	154	<div> <div>0%</div> <div> <div></div> <div></div> <div></div> <div></div> </div> <div>90%</div> <div>6%</div> <div>.</div> </div>

*Continued on next page...*

Continued from previous page...

Mol	Chain	Length	Quality of chain
3	M	154	<div><div><div>%</div><div><div></div></div><div>96%</div><div><div></div></div><div>...</div></div></div>

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 9637 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 Fab Heavy Chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	220	Total	C	N	O	S	0	0	0
			1623	1026	269	323	5			
1	R	221	Total	C	N	O	S	0	0	0
			1642	1039	271	326	6			

- Molecule 2 is a protein called Fab Light Chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	213	Total	C	N	O	S	0	0	0
			1615	1008	272	329	6			
2	D	211	Total	C	N	O	S	0	0	0
			1611	1008	272	326	5			

- Molecule 3 is a protein called Histone chaperone ASF1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	M	152	Total	C	N	O	S	0	0	0
			1210	778	196	234	2			
3	A	149	Total	C	N	O	S	0	1	0
			1197	768	196	231	2			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
M	1	GLY	-	expression tag	UNP P32447
A	1	GLY	-	expression tag	UNP P32447

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	152	Total	O	0	0
			152	152		

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	C	118	Total 118	O 118	0	0
4	D	130	Total 130	O 130	0	0
4	M	100	Total 100	O 100	0	0
4	R	180	Total 180	O 180	0	0
4	A	59	Total 59	O 59	0	0

### 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: Fab Heavy Chain

Chain B:  95%



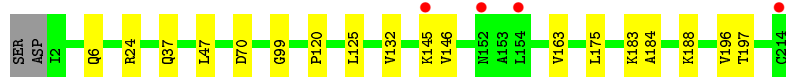
- Molecule 1: Fab Heavy Chain

Chain R:  94%




- Molecule 2: Fab Light Chain

Chain C:  91%



- Molecule 2: Fab Light Chain

Chain D:  87%

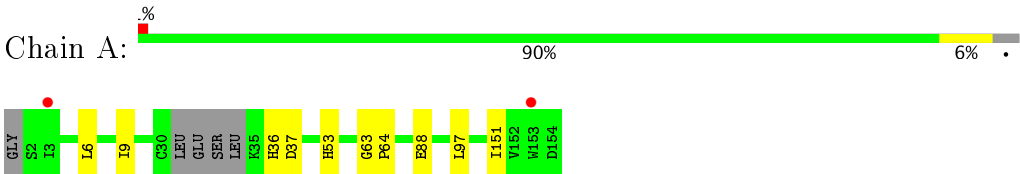


- Molecule 3: Histone chaperone ASF1

Chain M:  96%



- Molecule 3: Histone chaperone ASF1



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	86.81Å 77.44Å 101.73Å 90.00° 110.56° 90.00°	Depositor
Resolution (Å)	35.95 – 2.10 35.95 – 2.10	Depositor EDS
% Data completeness (in resolution range)	99.4 (35.95-2.10) 99.4 (35.95-2.10)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.21 (at 2.10Å)	Xtriage
Refinement program	PHENIX (1.10 _2152: ???)	Depositor
R, $R_{free}$	0.191 , 0.236 0.191 , 0.235	Depositor DCC
$R_{free}$ test set	3715 reflections (5.03%)	DCC
Wilson B-factor (Å <sup>2</sup> )	32.3	Xtriage
Anisotropy	0.421	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.29 , 37.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	9637	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.82% 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

### 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	B	0.28	0/1665	0.51	0/2276
1	R	0.29	0/1685	0.51	0/2302
2	C	0.29	0/1649	0.50	0/2240
2	D	0.28	0/1645	0.51	0/2233
3	A	0.27	0/1226	0.47	0/1674
3	M	0.28	0/1239	0.46	0/1693
All	All	0.28	0/9109	0.50	0/12418

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

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	B	1623	0	1564	5	0
1	R	1642	0	1585	6	0
2	C	1615	0	1559	10	0
2	D	1611	0	1575	14	0
3	A	1197	0	1157	6	0
3	M	1210	0	1181	3	0
4	A	59	0	0	0	0
4	B	152	0	0	1	0
4	C	118	0	0	2	0
4	D	130	0	0	2	0

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	M	100	0	0	0	0
4	R	180	0	0	0	0
All	All	9637	0	8621	43	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 2.

All (43) 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:C:120:PRO:HD3	2:C:132:VAL:HG22	1.73	0.69
2:D:2:ILE:HG12	2:D:95:LEU:HD12	1.75	0.69
3:A:6:LEU:HD21	3:A:9:ILE:HD11	1.84	0.58
2:C:37:GLN:HB2	2:C:47:LEU:HD11	1.84	0.58
1:B:121:LYS:NZ	4:B:303:HOH:O	2.41	0.54
1:R:136:SER:O	1:R:139:THR:HG22	2.08	0.53
3:A:36:HIS:O	3:A:37:ASP:HB2	2.09	0.53
2:D:105:GLU:OE2	2:D:142:ARG:NH2	2.40	0.52
3:A:63:GLY:HA2	3:A:64:PRO:C	2.29	0.52
3:A:9:ILE:HD13	3:A:97:LEU:HD23	1.92	0.51
2:D:2:ILE:HD11	2:D:95:LEU:HB2	1.91	0.51
2:C:125:LEU:O	2:C:183:LYS:HD3	2.11	0.51
1:B:12:VAL:HG11	1:B:86:LEU:HD13	1.92	0.51
3:A:37:ASP:OD1	3:A:63:GLY:HA3	2.12	0.50
2:C:6:GLN:N	4:C:301:HOH:O	2.45	0.49
2:C:163:VAL:HG22	2:C:175:LEU:HD12	1.96	0.48
2:C:184:ALA:O	2:C:188:LYS:HG2	2.14	0.48
1:R:214:LYS:HE3	1:R:216:GLU:OE1	2.13	0.47
4:D:305:HOH:O	3:M:49:ARG:NH1	2.46	0.46
2:C:146:VAL:HG22	2:C:196:VAL:HG22	1.99	0.45
1:B:33:SER:OG	1:B:99:GLY:HA3	2.17	0.45
2:D:184:ALA:O	2:D:188:LYS:HG3	2.16	0.45
2:D:154:LEU:HD23	2:D:154:LEU:HA	1.80	0.44
2:D:210:ASN:O	2:D:211:ARG:HB2	2.17	0.44
2:D:187:GLU:HA	2:D:212:GLY:C	2.38	0.44
3:A:53[B]:HIS:HE2	3:A:88:GLU:CD	2.21	0.44
2:C:145:LYS:HB2	2:C:197:THR:HB	1.98	0.44
2:C:99:GLY:HA2	4:C:301:HOH:O	2.18	0.44
2:D:125:LEU:O	2:D:183:LYS:HD2	2.17	0.43
1:R:33:SER:OG	1:R:99:GLY:HA3	2.18	0.43
3:M:145:ARG:HG3	1:R:102:TRP:CE2	2.53	0.43

*Continued on next page...*

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:121:LYS:HE3	1:B:179:LEU:HD13	2.00	0.42
2:C:24:ARG:HD3	2:C:70:ASP:OD1	2.19	0.42
1:R:131:SER:O	1:R:135:THR:HG23	2.20	0.42
1:R:188:VAL:HG11	1:R:198:TYR:CE1	2.55	0.42
1:B:121:LYS:HE3	1:B:148:ASP:HB3	2.02	0.41
2:D:169:LYS:HD3	4:D:421:HOH:O	2.19	0.41
2:D:83:PHE:CD1	2:D:106:ILE:HG12	2.55	0.41
2:D:33:VAL:HA	2:D:89:GLN:O	2.21	0.41
3:M:53:HIS:HE2	3:M:88:GLU:CD	2.24	0.41
2:D:35:TRP:CZ3	2:D:88:CYS:HB3	2.56	0.41
2:D:37:GLN:HB2	2:D:47:LEU:HD11	2.02	0.40
2:D:149:LYS:HG2	2:D:154:LEU:HD23	2.02	0.40

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	B	218/224 (97%)	213 (98%)	5 (2%)	0	100	100
1	R	219/224 (98%)	215 (98%)	4 (2%)	0	100	100
2	C	211/215 (98%)	205 (97%)	6 (3%)	0	100	100
2	D	209/215 (97%)	202 (97%)	6 (3%)	1 (0%)	32	28
3	A	146/154 (95%)	142 (97%)	4 (3%)	0	100	100
3	M	150/154 (97%)	149 (99%)	1 (1%)	0	100	100
All	All	1153/1186 (97%)	1126 (98%)	26 (2%)	1 (0%)	55	57

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	D	211	ARG

### 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	B	178/185 (96%)	178 (100%)	0	100	100
1	R	181/185 (98%)	181 (100%)	0	100	100
2	C	184/189 (97%)	184 (100%)	0	100	100
2	D	185/189 (98%)	184 (100%)	1 (0%)	91	94
3	A	136/141 (96%)	135 (99%)	1 (1%)	87	91
3	M	138/141 (98%)	137 (99%)	1 (1%)	87	91
All	All	1002/1030 (97%)	999 (100%)	3 (0%)	94	96

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	D	169	LYS
3	M	49	ARG
3	A	151	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
2	D	147	GLN
3	A	70	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 [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	B	220/224 (98%)	-0.38	1 (0%) 90 92	24, 35, 52, 77	0
1	R	221/224 (98%)	-0.38	1 (0%) 90 92	21, 33, 54, 73	0
2	C	213/215 (99%)	-0.08	4 (1%) 67 71	22, 39, 75, 127	0
2	D	211/215 (98%)	-0.23	1 (0%) 90 92	23, 35, 59, 106	0
3	A	149/154 (96%)	0.10	2 (1%) 77 81	26, 45, 73, 91	0
3	M	152/154 (98%)	-0.20	2 (1%) 77 81	23, 37, 55, 61	0
All	All	1166/1186 (98%)	-0.21	11 (0%) 84 86	21, 36, 64, 127	0

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	C	214	CYS	6.3
2	D	2	ILE	5.5
3	A	153	TRP	4.7
2	C	154	LEU	4.0
2	C	152	ASN	2.7
3	A	3	ILE	2.5
1	B	0	SER	2.5
3	M	31	LEU	2.4
3	M	67	VAL	2.2
2	C	145	LYS	2.2
1	R	221	ASP	2.2

### 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 [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.