



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 7, 2022 – 10:20 AM EST

PDB ID : 7RXL
Title : Fab1488 in complex with the C-terminal alpha-TSR domain of *P. falciparum*
Authors : Pholcharee, T.; Oyen, D.; Wilson, I.A.
Deposited on : 2021-08-23
Resolution : 1.82 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 ⓘ 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 : 2.27
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) : 2.27

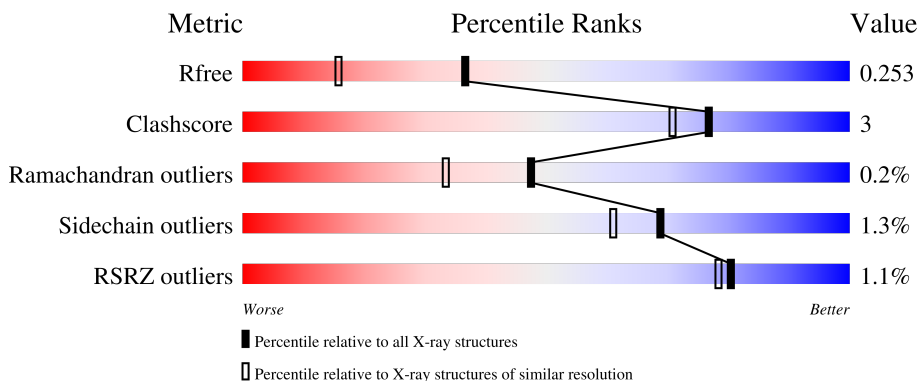
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.82 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.



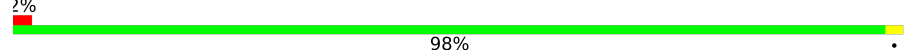
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	7484 (1.84-1.80)
Clashscore	141614	8401 (1.84-1.80)
Ramachandran outliers	138981	8290 (1.84-1.80)
Sidechain outliers	138945	8290 (1.84-1.80)
RSRZ outliers	127900	7371 (1.84-1.80)

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 $\leq 5\%$. 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	D	212	 91% 7%
1	L	212	 85% 12%
2	C	224	 92% 7%
2	H	224	 89% 8%
3	E	66	 98%

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Mol	Chain	Length	Quality of chain
3	F	66	 A horizontal bar chart representing the quality of the chain. The bar is divided into three segments: a small red segment at the beginning labeled '2%', a large green segment in the middle labeled '98%', and a very small yellow segment at the end. A small black dot is visible at the far right end of the bar.

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 14692 atoms, of which 7040 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 Fab1488 light chain.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
1	L	208	3019	966	1475	257	315	6	0	2	0
1	D	207	3004	965	1469	252	312	6	0	2	0

- Molecule 2 is a protein called Fab1488 heavy chain.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
2	H	218	3170	1027	1550	271	316	6	1	1	0
2	C	222	3234	1043	1586	277	322	6	0	0	0

- Molecule 3 is a protein called Circumsporozoite protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
3	E	66	976	309	477	85	100	5	0	0	0
3	F	66	982	309	483	86	99	5	0	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	L	67	Total O 67 67	0	0
4	H	60	Total O 60 60	0	0
4	D	60	Total O 60 60	0	0
4	C	86	Total O 86 86	0	0

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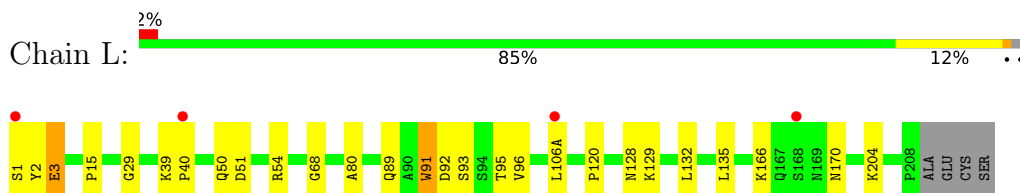
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	E	21	Total 21	O 21	0	0
4	F	13	Total 13	O 13	0	0

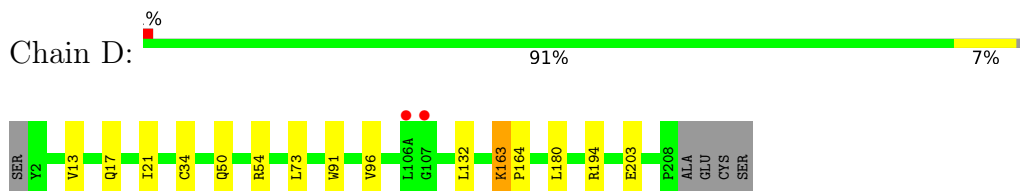
3 Residue-property plots [i](#)

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.

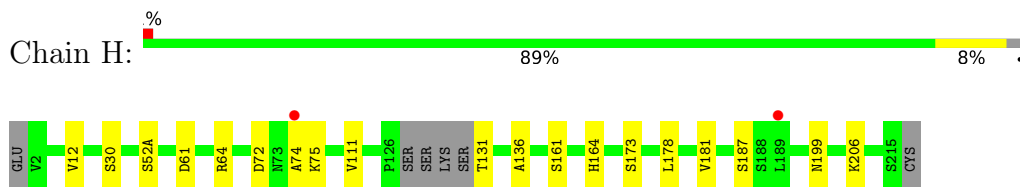
- Molecule 1: Fab1488 light chain



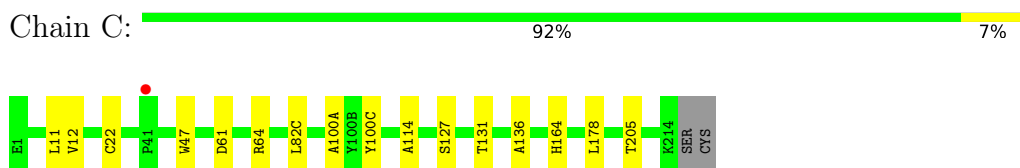
- Molecule 1: Fab1488 light chain



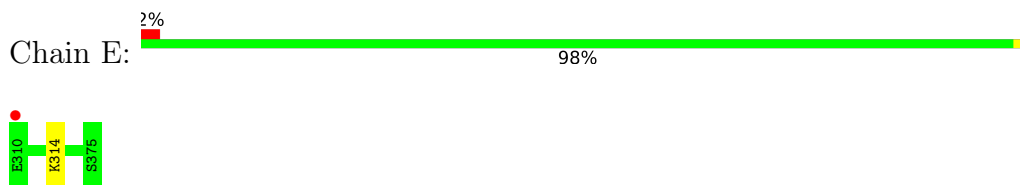
- Molecule 2: Fab1488 heavy chain



- Molecule 2: Fab1488 heavy chain



- Molecule 3: Circumsporozoite protein



- Molecule 3: Circumsporozoite protein



4 Data and refinement statistics i

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	40.69Å 69.34Å 93.21Å 85.06° 78.14° 79.18°	Depositor
Resolution (Å)	39.21 – 1.82 39.21 – 1.82	Depositor EDS
% Data completeness (in resolution range)	94.6 (39.21-1.82) 94.7 (39.21-1.82)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.20 (at 1.82Å)	Xtrriage
Refinement program	PHENIX (1.14_3260: ???)	Depositor
R, R_{free}	0.206 , 0.255 0.205 , 0.253	Depositor DCC
R_{free} test set	3911 reflections (4.74%)	wwPDB-VP
Wilson B-factor (Å ²)	24.6	Xtrriage
Anisotropy	0.847	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.43 , 43.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	0.013 for -h,-k,-h+l	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	14692	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.64% 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 [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	D	0.40	0/1574	0.53	0/2156
1	L	0.39	0/1582	0.52	0/2167
2	C	0.42	0/1690	0.57	0/2304
2	H	0.40	0/1666	0.56	0/2272
3	E	0.33	0/507	0.46	0/684
3	F	0.40	0/507	0.49	0/683
All	All	0.40	0/7526	0.54	0/10266

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	D	1535	1469	1457	10	0
1	L	1544	1475	1470	16	0
2	C	1648	1586	1584	11	0
2	H	1620	1550	1537	10	0
3	E	499	477	473	1	0
3	F	499	483	481	1	0
4	C	86	0	0	0	0
4	D	60	0	0	0	0
4	E	21	0	0	0	0
4	F	13	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	H	60	0	0	1	0
4	L	67	0	0	0	0
All	All	7652	7040	7002	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 3.

The worst 5 of 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:12:VAL:HG11	2:C:82(C):LEU:HD13	1.78	0.65
2:C:127:SER:O	2:C:131:THR:HG23	1.97	0.65
1:D:132:LEU:HD11	1:D:180:LEU:HD11	1.81	0.62
2:H:72:ASP:OD1	2:H:74:ALA:HB3	1.99	0.62
2:C:131:THR:HG22	2:C:136:ALA:HB2	1.82	0.61

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	D	207/212 (98%)	201 (97%)	6 (3%)	0	100	100
1	L	208/212 (98%)	203 (98%)	4 (2%)	1 (0%)	29	15
2	C	220/224 (98%)	214 (97%)	6 (3%)	0	100	100
2	H	215/224 (96%)	207 (96%)	7 (3%)	1 (0%)	29	15
3	E	64/66 (97%)	63 (98%)	1 (2%)	0	100	100
3	F	64/66 (97%)	63 (98%)	1 (2%)	0	100	100
All	All	978/1004 (97%)	951 (97%)	25 (3%)	2 (0%)	47	33

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	L	3	GLU
2	H	173	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	D	168/179 (94%)	166 (99%)	2 (1%)	71	64
1	L	170/179 (95%)	167 (98%)	3 (2%)	59	48
2	C	179/187 (96%)	177 (99%)	2 (1%)	73	67
2	H	174/187 (93%)	171 (98%)	3 (2%)	60	50
3	E	55/61 (90%)	55 (100%)	0	100	100
3	F	56/61 (92%)	56 (100%)	0	100	100
All	All	802/854 (94%)	792 (99%)	10 (1%)	69	64

5 of 10 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	163	LYS
2	C	22	CYS
2	C	164	HIS
2	H	161	SER
2	H	164	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	L	69	ASN
1	L	108	GLN
1	D	126	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 monosaccharides 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	D	207/212 (97%)	0.12	2 (0%) 82 80	21, 34, 51, 66	0
1	L	208/212 (98%)	0.18	4 (1%) 66 63	23, 36, 50, 62	0
2	C	222/224 (99%)	-0.00	1 (0%) 91 89	20, 31, 43, 61	0
2	H	218/224 (97%)	0.06	2 (0%) 84 82	22, 32, 56, 69	0
3	E	66/66 (100%)	0.09	1 (1%) 73 70	28, 39, 54, 72	0
3	F	66/66 (100%)	0.08	1 (1%) 73 70	28, 38, 55, 63	0
All	All	987/1004 (98%)	0.09	11 (1%) 80 78	20, 34, 52, 72	0

The worst 5 of 11 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	L	168	SER	3.7
3	E	310	GLU	3.0
2	H	74	ALA	3.0
2	H	189	LEU	2.8
1	L	106(A)	LEU	2.8

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 monosaccharides 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.