



# wwPDB X-ray Structure Validation Summary Report ⓘ

Apr 14, 2022 – 02:24 PM EDT

PDB ID : 7T0J  
Title : Crystal structure of S25-2 Fab Unliganded 3  
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Deposited on : 2021-11-29  
Resolution : 2.15 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](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	:	2.27
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

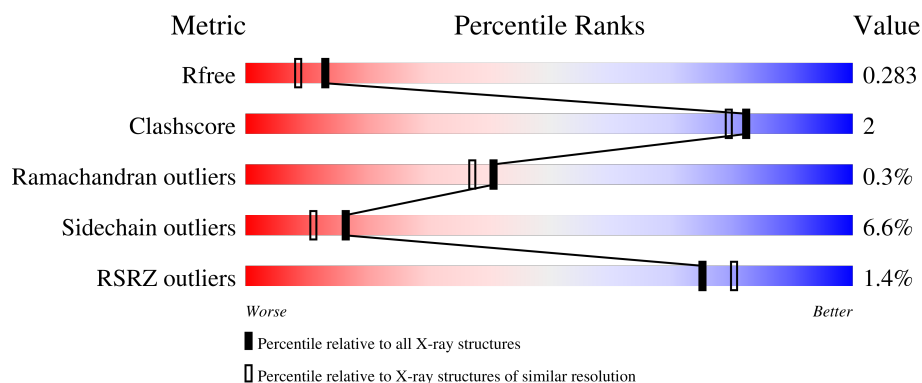
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.15 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	1479 (2.16-2.16)
Clashscore	141614	1585 (2.16-2.16)
Ramachandran outliers	138981	1560 (2.16-2.16)
Sidechain outliers	138945	1559 (2.16-2.16)
RSRZ outliers	127900	1456 (2.16-2.16)

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  $\geq 3$ , 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	219	<div> <div>5%</div> <div>92%</div> <div>6%</div> </div>
1	C	219	<div> <div>89%</div> <div>10%</div> </div>
1	E	219	<div> <div>86%</div> <div>13%</div> </div>
1	L	219	<div> <div>85%</div> <div>13%</div> </div>
2	B	223	<div> <div>%</div> <div>82%</div> <div>13%</div> </div>

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Mol	Chain	Length	Quality of chain
2	D	223	<div><div>%</div><div><div></div><div>87%</div><div>9%</div><div></div></div><div></div></div>
2	F	223	<div><div>3%</div><div><div></div><div>88%</div><div>9%</div><div></div></div><div></div></div>
2	H	223	<div><div></div><div><div></div><div>93%</div><div>6%</div><div></div></div><div></div></div>

## 2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 13697 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 S25-2 Fab light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	218	Total	C	N	O	S	0	0	0
			1694	1053	289	345	7			
1	A	218	Total	C	N	O	S	0	0	0
			1681	1047	285	342	7			
1	C	219	Total	C	N	O	S	0	0	0
			1700	1056	290	346	8			
1	E	218	Total	C	N	O	S	0	0	0
			1688	1050	286	345	7			

- Molecule 2 is a protein called S25-2 Fab heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	223	Total	C	N	O	S	0	0	0
			1702	1078	284	333	7			
2	B	214	Total	C	N	O	S	0	0	0
			1628	1032	270	319	7			
2	D	214	Total	C	N	O	S	0	0	0
			1642	1045	270	320	7			
2	F	219	Total	C	N	O	S	0	0	0
			1678	1064	279	328	7			

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	L	25	Total	O	0	0
			25	25		
3	H	64	Total	O	0	0
			64	64		
3	A	14	Total	O	0	0
			14	14		
3	B	13	Total	O	0	0
			13	13		

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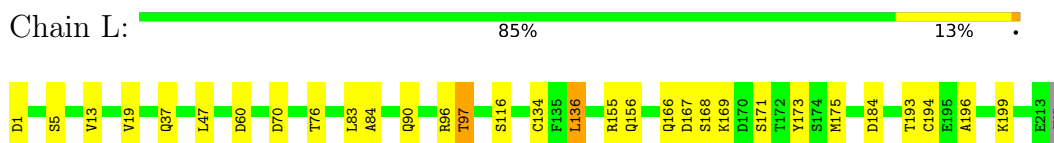
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	C	53	Total 53	O 53	0	0
3	D	47	Total 47	O 47	0	0
3	E	29	Total 29	O 29	0	0
3	F	39	Total 39	O 39	0	0

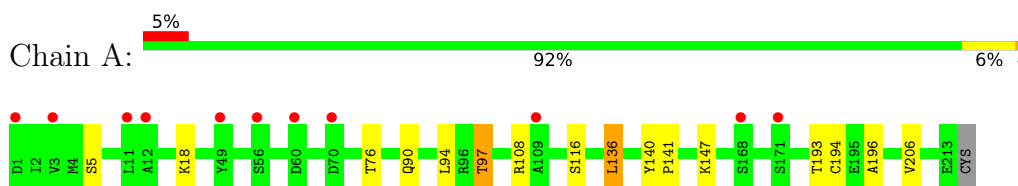
### 3 Residue-property plots [i](#)

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.

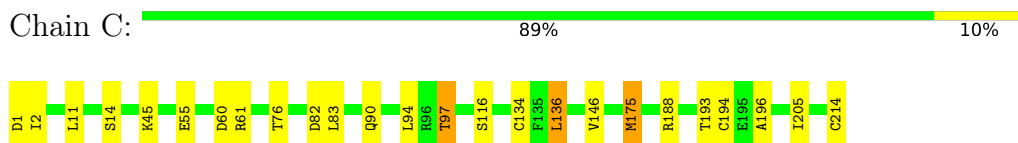
- Molecule 1: S25-2 Fab light chain



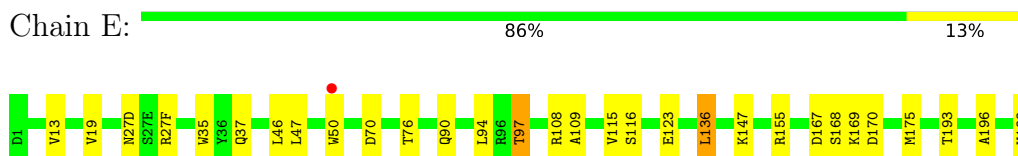
- Molecule 1: S25-2 Fab light chain



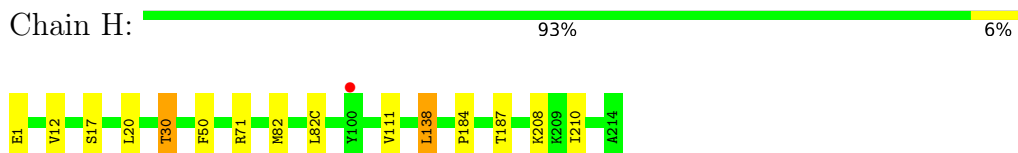
- Molecule 1: S25-2 Fab light chain



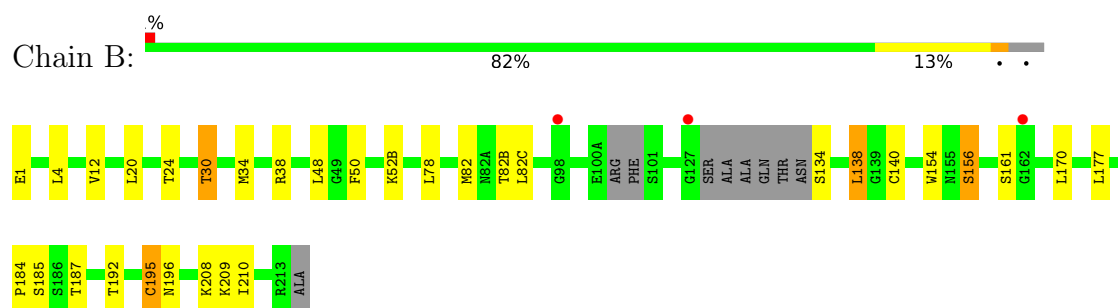
- Molecule 1: S25-2 Fab light chain



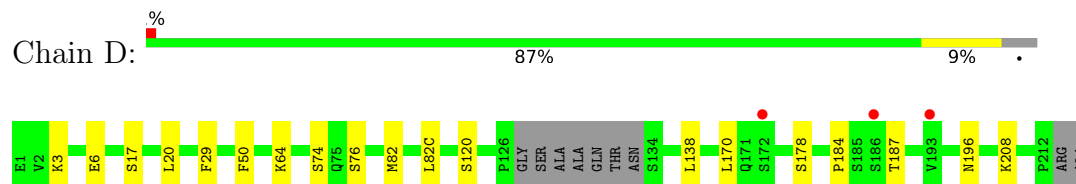
- Molecule 2: S25-2 Fab heavy chain



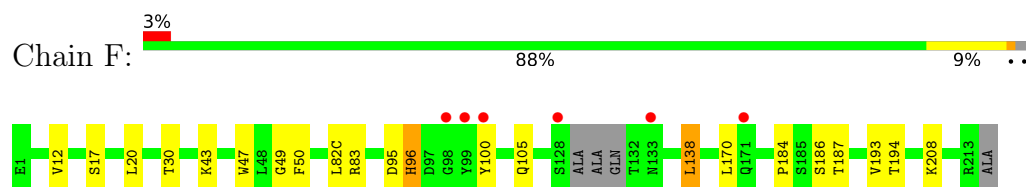
- Molecule 2: S25-2 Fab heavy chain



- Molecule 2: S25-2 Fab heavy chain



- Molecule 2: S25-2 Fab heavy chain



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	73.73Å 135.33Å 92.67Å 90.00° 94.35° 90.00°	Depositor
Resolution (Å)	40.00 – 2.15 38.45 – 2.15	Depositor EDS
% Data completeness (in resolution range)	99.7 (40.00-2.15) 99.7 (38.45-2.15)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.12 (at 2.14Å)	Xtriage
Refinement program	REFMAC 5.8.0155	Depositor
R, $R_{free}$	0.230 , 0.281 0.234 , 0.283	Depositor DCC
$R_{free}$ test set	4992 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	36.2	Xtriage
Anisotropy	0.505	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.29 , 14.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	13697	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.88% 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	0.53	0/1716	0.72	0/2327
1	C	0.63	0/1735	0.75	0/2351
1	E	0.58	0/1723	0.76	0/2336
1	L	0.54	0/1729	0.75	1/2343 (0.0%)
2	B	0.54	0/1670	0.73	0/2279
2	D	0.59	0/1687	0.75	0/2304
2	F	0.59	0/1723	0.78	0/2352
2	H	0.62	0/1748	0.78	1/2388 (0.0%)
All	All	0.58	0/13731	0.75	2/18680 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	L	96	ARG	NE-CZ-NH1	5.81	123.21	120.30
2	H	71	ARG	NE-CZ-NH1	5.77	123.18	120.30

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	1681	0	1625	5	0
1	C	1700	0	1649	9	0
1	E	1688	0	1633	12	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	L	1694	0	1644	7	0
2	B	1628	0	1581	16	0
2	D	1642	0	1595	4	0
2	F	1678	0	1629	8	0
2	H	1702	0	1653	6	0
3	A	14	0	0	0	0
3	B	13	0	0	0	0
3	C	53	0	0	2	0
3	D	47	0	0	0	0
3	E	29	0	0	0	0
3	F	39	0	0	1	0
3	H	64	0	0	2	0
3	L	25	0	0	0	0
All	All	13697	0	13009	65	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.

The worst 5 of 65 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:13:VAL:HG11	1:L:19:VAL:HG12	1.69	0.72
1:C:90:GLN:HE21	1:C:97:THR:HG22	1.60	0.67
1:C:90:GLN:HE21	1:C:97:THR:CG2	2.08	0.66
1:L:90:GLN:HE21	1:L:97:THR:HG22	1.64	0.63
1:E:108:ARG:NH1	1:E:109:ALA:O	2.33	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	A	216/219 (99%)	205 (95%)	11 (5%)	0	100	100
1	C	217/219 (99%)	210 (97%)	7 (3%)	0	100	100
1	E	216/219 (99%)	207 (96%)	7 (3%)	2 (1%)	17	11
1	L	216/219 (99%)	209 (97%)	5 (2%)	2 (1%)	17	11
2	B	208/223 (93%)	198 (95%)	9 (4%)	1 (0%)	29	22
2	D	210/223 (94%)	208 (99%)	2 (1%)	0	100	100
2	F	215/223 (96%)	207 (96%)	7 (3%)	1 (0%)	29	22
2	H	221/223 (99%)	217 (98%)	4 (2%)	0	100	100
All	All	1719/1768 (97%)	1661 (97%)	52 (3%)	6 (0%)	41	37

5 of 6 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	27(F)	ARG
2	F	96	HIS
1	L	84	ALA
1	L	169	LYS
2	B	156	SER

### 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	191/195 (98%)	181 (95%)	10 (5%)	23	19
1	C	195/195 (100%)	178 (91%)	17 (9%)	10	6
1	E	193/195 (99%)	179 (93%)	14 (7%)	14	9
1	L	194/195 (100%)	176 (91%)	18 (9%)	9	5
2	B	183/190 (96%)	172 (94%)	11 (6%)	19	14
2	D	185/190 (97%)	173 (94%)	12 (6%)	17	12
2	F	189/190 (100%)	177 (94%)	12 (6%)	18	13
2	H	190/190 (100%)	183 (96%)	7 (4%)	34	32

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	1520/1540 (99%)	1419 (93%)	101 (7%)	16 11

5 of 101 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	134	CYS
2	D	170	LEU
2	F	194	THR
1	C	175	MET
2	D	20	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 8 such sidechains are listed below:

Mol	Chain	Res	Type
2	F	164	HIS
2	F	105	GLN
2	D	96	HIS
1	C	93	ASN
1	E	79	GLN

### 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 ⓘ

There are no monosaccharides in this entry.

## 5.6 Ligand geometry ⓘ

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	218/219 (99%)	0.38	11 (5%) 28 37	35, 58, 79, 92	0
1	C	219/219 (100%)	-0.26	0 100 100	25, 37, 58, 70	0
1	E	218/219 (99%)	-0.02	1 (0%) 91 93	27, 46, 68, 86	0
1	L	218/219 (99%)	-0.02	0 100 100	27, 46, 73, 91	0
2	B	214/223 (95%)	-0.04	3 (1%) 75 80	28, 44, 69, 85	0
2	D	214/223 (95%)	-0.12	3 (1%) 75 80	23, 38, 70, 91	0
2	F	219/223 (98%)	-0.13	6 (2%) 54 63	25, 36, 62, 86	0
2	H	223/223 (100%)	-0.21	1 (0%) 92 94	20, 33, 54, 80	0
All	All	1743/1768 (98%)	-0.06	25 (1%) 75 80	20, 42, 70, 92	0

The worst 5 of 25 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	162	GLY	4.0
2	F	171	GLN	3.6
2	B	98	GLY	3.2
1	A	49	TYR	3.1
1	A	109	ALA	2.9

### 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.