



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 9, 2024 – 01:29 pm GMT

PDB ID : 5NIV
Title : Crystal structure of 5D3 Fab
Authors : Manolaridis, I.; Locher, K.P.
Deposited on : 2017-03-27
Resolution : 1.50 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

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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	:	3.0
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.003 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.39

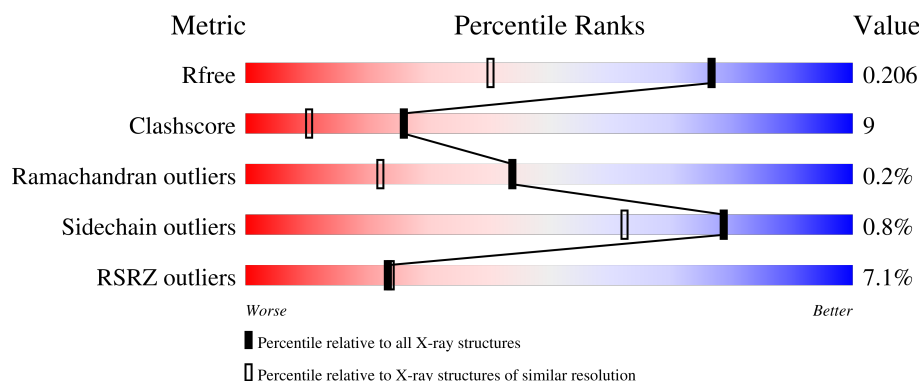
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.50 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}	164625	3717 (1.50-1.50)
Clashscore	180529	4048 (1.50-1.50)
Ramachandran outliers	177936	3970 (1.50-1.50)
Sidechain outliers	177891	3967 (1.50-1.50)
RSRZ outliers	164620	3718 (1.50-1.50)

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	A	212	
2	B	221	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3848 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 Light chain of 5D3 Fab.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	212	Total	C	N	O	S	0	0	0
			1645	1029	278	333	5			

- Molecule 2 is a protein called Heavy chain of 5D3 Fab.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	213	Total	C	N	O	S	0	0	0
			1628	1038	266	318	6			

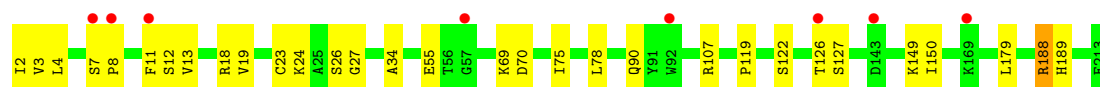
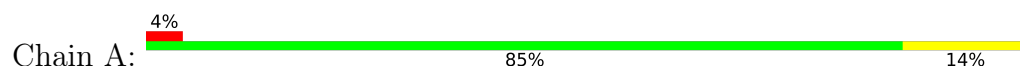
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	272	Total	O	0	0
			272	272		
3	B	303	Total	O	0	0
			303	303		

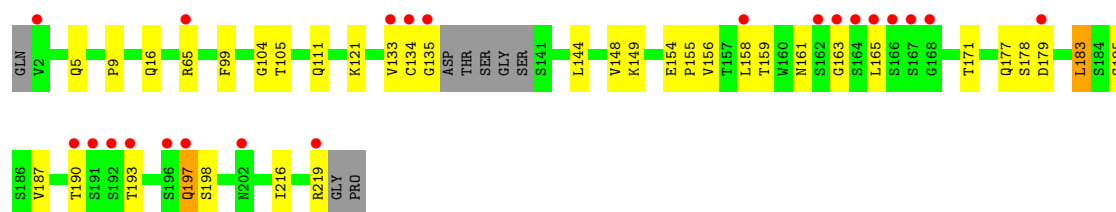
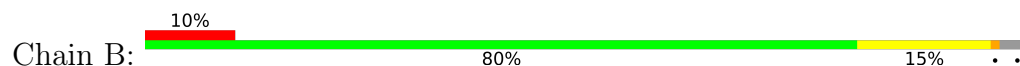
3 Residue-property plots [i](#)

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: Light chain of 5D3 Fab



- Molecule 2: Heavy chain of 5D3 Fab



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	139.20Å 45.19Å 83.28Å 90.00° 122.14° 90.00°	Depositor
Resolution (Å)	42.20 – 1.50 42.20 – 1.50	Depositor EDS
% Data completeness (in resolution range)	93.7 (42.20-1.50) 94.0 (42.20-1.50)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.09 (at 1.50Å)	Xtriage
Refinement program	PHENIX (1.10.1_2155: ???)	Depositor
R, R_{free}	0.181 , 0.205 0.179 , 0.206	Depositor DCC
R_{free} test set	3329 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	15.0	Xtriage
Anisotropy	0.064	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 35.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	3848	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.24% 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	A	0.35	0/1684	0.60	0/2289
2	B	0.35	0/1673	0.64	1/2289 (0.0%)
All	All	0.35	0/3357	0.62	1/4578 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	183	LEU	CB-CG-CD2	-5.89	100.98	111.00

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	1645	0	1578	27	0
2	B	1628	0	1590	36	0
3	A	272	0	0	11	1
3	B	303	0	0	12	2
All	All	3848	0	3168	60	3

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 9.

All (60) 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:B:111:GLN:OE1	3:B:301:HOH:O	1.95	0.82
1:A:2:ILE:N	3:A:302:HOH:O	2.12	0.81
2:B:155:PRO:HD2	3:B:505:HOH:O	1.81	0.79
2:B:133:VAL:O	2:B:135:GLY:N	2.15	0.78
1:A:18:ARG:NH1	3:A:303:HOH:O	2.16	0.78
2:B:177:GLN:OE1	3:B:303:HOH:O	2.03	0.76
3:A:406:HOH:O	2:B:133:VAL:HG12	1.88	0.74
1:A:3:VAL:HG22	1:A:26:SER:HB3	1.70	0.73
2:B:197:GLN:NE2	3:B:304:HOH:O	2.22	0.72
1:A:119:PRO:HG2	2:B:133:VAL:HB	1.72	0.71
1:A:19:VAL:HG12	1:A:75:ILE:HB	1.76	0.67
1:A:127:SER:HB3	3:A:320:HOH:O	1.96	0.65
1:A:24:LYS:HD2	1:A:70:ASP:OD1	2.01	0.61
1:A:7:SER:O	3:A:304:HOH:O	2.17	0.59
1:A:55:GLU:OE1	3:A:305:HOH:O	2.17	0.58
2:B:158:LEU:HD11	2:B:183:LEU:HD11	1.86	0.57
1:A:107:ARG:NH1	3:A:310:HOH:O	2.38	0.57
1:A:7:SER:N	3:A:304:HOH:O	2.37	0.56
2:B:171:THR:OG1	3:B:302:HOH:O	2.00	0.56
2:B:156:VAL:CG2	2:B:183:LEU:HD21	2.36	0.55
2:B:9:PRO:HG2	3:B:440:HOH:O	2.07	0.55
1:A:23:CYS:O	1:A:24:LYS:HD3	2.06	0.54
2:B:65:ARG:HG2	2:B:65:ARG:HH11	1.72	0.54
2:B:5:GLN:OE1	2:B:111:GLN:NE2	2.41	0.53
2:B:178:SER:O	2:B:179:ASP:HB3	2.08	0.53
2:B:158:LEU:HD22	2:B:185:SER:CB	2.39	0.53
1:A:18:ARG:NH2	3:A:306:HOH:O	2.21	0.53
2:B:121:LYS:NZ	3:B:308:HOH:O	2.28	0.52
1:A:4:LEU:HD11	1:A:90:GLN:N	2.25	0.52
2:B:149:LYS:HE2	3:B:303:HOH:O	2.10	0.52
1:A:4:LEU:HD11	1:A:90:GLN:H	1.74	0.52
2:B:154:GLU:HB3	2:B:155:PRO:HG3	1.94	0.50
1:A:122:SER:O	1:A:126:THR:HG23	2.13	0.49
2:B:144:LEU:HD13	2:B:216:ILE:HG21	1.94	0.49
1:A:149:LYS:NZ	3:A:301:HOH:O	1.97	0.49
2:B:158:LEU:HD22	2:B:185:SER:HB2	1.93	0.49
1:A:188:ARG:HG2	1:A:189:HIS:CE1	2.48	0.48
2:B:99:PHE:CE1	2:B:104:GLY:HA2	2.48	0.48
1:A:7:SER:C	3:A:304:HOH:O	2.51	0.47
2:B:65:ARG:HG2	2:B:65:ARG:NH1	2.30	0.47
1:A:150:ILE:HD11	1:A:179:LEU:HD21	1.96	0.47
2:B:148:VAL:HG21	2:B:158:LEU:HD21	1.97	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:154:GLU:HG3	3:B:367:HOH:O	2.16	0.46
1:A:24:LYS:HD3	1:A:24:LYS:HA	1.69	0.45
2:B:161:ASN:ND2	3:B:304:HOH:O	2.25	0.45
1:A:34:ALA:HB2	2:B:105:THR:HG22	1.99	0.45
1:A:13:VAL:HB	1:A:78:LEU:HD22	1.99	0.44
2:B:158:LEU:N	2:B:158:LEU:HD12	2.33	0.44
2:B:190:THR:O	2:B:193:THR:HG22	2.17	0.44
1:A:27:GLY:HA2	1:A:69:LYS:HE2	1.99	0.43
1:A:119:PRO:CG	2:B:133:VAL:HB	2.45	0.43
2:B:158:LEU:HD22	2:B:185:SER:HB3	2.01	0.43
2:B:158:LEU:HD11	2:B:183:LEU:CD1	2.48	0.42
2:B:155:PRO:HB3	3:B:452:HOH:O	2.19	0.41
1:A:12:SER:HB3	1:A:107:ARG:NH2	2.36	0.41
2:B:16:GLN:OE1	3:B:305:HOH:O	2.22	0.41
2:B:159:THR:HG22	2:B:163:GLY:N	2.36	0.41
2:B:165:LEU:HD13	2:B:187:VAL:HG21	2.03	0.41
1:A:8:PRO:HD2	1:A:11:PHE:CZ	2.56	0.41
2:B:197:GLN:HG2	2:B:198:SER:N	2.36	0.40

All (3) 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)
3:B:556:HOH:O	3:B:601:HOH:O[4_546]	2.06	0.14
3:A:361:HOH:O	3:A:457:HOH:O[4_555]	2.10	0.10
3:B:557:HOH:O	3:B:565:HOH:O[4_556]	2.18	0.02

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	A	210/212 (99%)	208 (99%)	2 (1%)	0	100 100

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	B	209/221 (95%)	201 (96%)	7 (3%)	1 (0%)	25	8
All	All	419/433 (97%)	409 (98%)	9 (2%)	1 (0%)	44	22

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	134	CYS

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	187/187 (100%)	186 (100%)	1 (0%)	86	75
2	B	187/193 (97%)	185 (99%)	2 (1%)	70	48
All	All	374/380 (98%)	371 (99%)	3 (1%)	79	62

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

Mol	Chain	Res	Type
1	A	188	ARG
2	B	197	GLN
2	B	219	ARG

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

Mol	Chain	Res	Type
1	A	79	GLN
2	B	3	GLN
2	B	202	ASN

5.3.3 RNA ⓘ

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 oligosaccharides 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, 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	A	212/212 (100%)	0.17	8 (3%) 44 48	10, 17, 27, 39	0
2	B	213/221 (96%)	0.44	22 (10%) 13 13	10, 16, 38, 50	0
All	All	425/433 (98%)	0.31	30 (7%) 23 24	10, 16, 33, 50	0

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	165	LEU	7.8
2	B	133	VAL	7.1
2	B	135	GLY	5.4
2	B	166	SER	4.7
2	B	167	SER	4.7
2	B	2	VAL	4.4
2	B	163	GLY	3.9
2	B	162	SER	3.9
2	B	134	CYS	3.9
2	B	219	ARG	3.6
2	B	164	SER	3.5
2	B	197	GLN	3.5
1	A	8	PRO	3.4
2	B	179	ASP	3.4
2	B	190	THR	3.4
2	B	158	LEU	3.3
1	A	7	SER	3.1
2	B	65	ARG	2.9
1	A	126	THR	2.7
2	B	196	SER	2.6
2	B	168	GLY	2.6
2	B	191	SER	2.6
2	B	193	THR	2.5
2	B	192	SER	2.5

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	11	PHE	2.4
1	A	92	TRP	2.4
1	A	57	GLY	2.3
2	B	202	ASN	2.3
1	A	169	LYS	2.2
1	A	143	ASP	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 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.