

Full wwPDB X-ray Structure Validation Report (i)

Aug 15, 2023 – 03:05 PM EDT

PDB ID : 1XF2

> Title Structure of Fab DNA-1 complexed with dT3

Authors Schuermann, J.P.; Prewitt, S.P.; Deutscher, S.L.; Tanner, J.J.

2004-09-13 Deposited on

2.30 Å(reported) Resolution

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 (i) 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 (1)) were used in the production of this report:

MolProbity 4.02b-467

> 1.8.5 (274361), CSD as541be (2020) Mogul

Xtriage (Phenix) 1.13

EDS 2.35

20191225.v01 (using entries in the PDB archive December 25th 2019) Percentile statistics

> Refmac 5.8.0158

CCP4 7.0.044 (Gargrove)

Ideal geometry (proteins) Engh & Huber (2001) Ideal geometry (DNA, RNA) Parkinson et al. (1996)

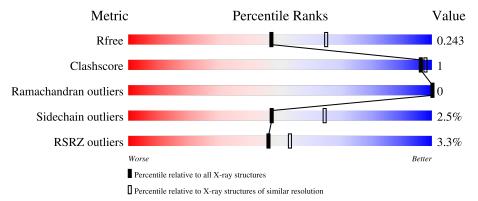
Validation Pipeline (wwPDB-VP) 2.35

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 2.30 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 $(\# \mathrm{Entries})$	$\begin{array}{c} {\rm Similar\ resolution} \\ (\#{\rm Entries,\ resolution\ range(\mathring{A})}) \end{array}$
R_{free}	130704	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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 <=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	Т	3	67%	33%			
2	A	214	94%	5%			
2	L	214	89%	10%			
3	В	230	89%	• 7%			
3	Н	230	90%	• 7%			



2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 6720 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 DNA chain called 5'-D(*TP*TP*T)-3'.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	Т	ર	Total	С	N	О	Р	0	0	0
1	1	3	41	20	4	15	2	U		

• Molecule 2 is a protein called antibody light chain Fab.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace	
2	L	213	Total 1637	C 1024	11	O 331	S 6	0	0	0
2	A	213	Total 1628	C 1020		O 330	S 6	0	0	0

• Molecule 3 is a protein called antibody heavy chain Fab.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
9	П	213	Total	С	N	О	S	0	0	0
3	Π	213	1596	1021	251	315	9	0	U	0
9	D	213	Total	Total C N O S	0	0	0			
3	D	213	1588	1018	252	309	9	0	U	U

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
Н	1	GLN	-	cloning artifact	GB 3399661
Н	2	VAL	-	cloning artifact	GB 3399661
Н	3	LYS	-	cloning artifact	GB 3399661
Н	4	LEU	-	cloning artifact	GB 3399661
Н	218	HIS	-	expression tag	GB 3399661
Н	219	HIS	-	expression tag	GB 3399661
Н	220	HIS	-	expression tag	GB 3399661
Н	221	HIS	-	expression tag	GB 3399661
Н	222	HIS	-	expression tag	GB 3399661
Н	223	HIS	-	expression tag	GB 3399661

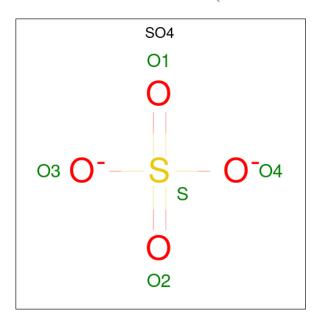
Continued on next page...



Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
В	1	GLN	-	cloning artifact	GB 3399661
В	2	VAL	-	cloning artifact	GB 3399661
В	3	LYS	-	cloning artifact	GB 3399661
В	4	LEU	-	cloning artifact	GB 3399661
В	218	HIS	-	expression tag	GB 3399661
В	219	HIS	_	expression tag	GB 3399661
В	220	HIS	-	expression tag	GB 3399661
В	221	HIS	-	expression tag	GB 3399661
В	222	HIS	-	expression tag	GB 3399661
В	223	HIS	-	expression tag	GB 3399661

 \bullet Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	L	1	Total O S 5 4 1	0	0
4	A	1	Total O S 5 4 1	0	0

• Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	Т	1	Total O 1 1	0	0
5	L	66	Total O 66 66	0	0

Continued on next page...



$Continued\ from\ previous\ page...$

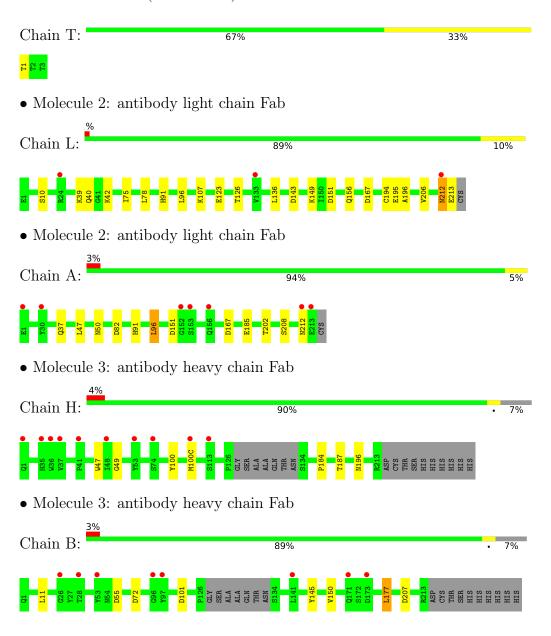
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	Н	65	Total O 65 65	0	0
5	A	40	Total O 40 40	0	0
5	В	48	Total O 48 48	0	0



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: 5'-D(*TP*TP*T)-3'





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants	170.68Å 170.68Å 144.67Å	Donogitor
a, b, c, α , β , γ	90.00° 90.00° 120.00°	Depositor
Resolution (Å)	29.50 - 2.30	Depositor
Resolution (A)	29.46 - 2.30	EDS
% Data completeness	99.8 (29.50-2.30)	Depositor
(in resolution range)	99.9 (29.46-2.30)	EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$< I/\sigma(I) > 1$	4.55 (at 2.31Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
D D	0.202 , 0.242	Depositor
R, R_{free}	0.205 , 0.243	DCC
R_{free} test set	2753 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	34.9	Xtriage
Anisotropy	0.276	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.37, 36.9	EDS
L-test for twinning ²	$ < L >=0.50, < L^2>=0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6720	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 3.01% of the height of the origin peak. No significant pseudotranslation is detected.

²Theoretical values of <|L|>, $<L^2>$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.



¹Intensities estimated from amplitudes.

5 Model quality (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: SO4

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).

Mal	Mol Chain		nd lengths	Bond angles		
IVIOI			# Z > 5	RMSZ	# Z > 5	
1	Т	1.17	0/44	1.90	$2/67 \ (3.0\%)$	
2	A	0.58	0/1666	0.75	3/2264 (0.1%)	
2	L	0.67	1/1675~(0.1%)	0.80	3/2274 (0.1%)	
3	В	0.58	0/1633	0.75	5/2234~(0.2%)	
3	Н	0.65	0/1641	0.76	0/2244	
All	All	0.63	$1/6659 \ (0.0\%)$	0.78	13/9083 (0.1%)	

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	Observed(A)	Ideal(A)
2	L	194	CYS	CB-SG	-5.38	1.73	1.81

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^o)$	$\operatorname{Ideal}({}^{o})$
1	Т	1	DT	O4'-C1'-N1	-8.90	101.77	108.00
2	L	167	ASP	CB-CG-OD2	7.17	124.75	118.30
2	L	151	ASP	CB-CG-OD2	6.69	124.32	118.30
3	В	72	ASP	CB-CG-OD2	6.29	123.96	118.30
1	Т	1	DT	N3-C2-O2	-6.18	118.59	122.30
2	A	167	ASP	CB-CG-OD2	5.84	123.56	118.30
2	A	82	ASP	CB-CG-OD2	5.50	123.25	118.30
2	L	143	ASP	CB-CG-OD2	5.45	123.21	118.30
3	В	177	LEU	CA-CB-CG	5.41	127.74	115.30
3	В	55	ASP	CB-CG-OD2	5.40	123.16	118.30
3	В	101	ASP	CB-CG-OD1	5.35	123.11	118.30
2	A	151	ASP	CB-CG-OD2	5.25	123.03	118.30
3	В	207	ASP	CB-CG-OD2	5.17	122.95	118.30

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	Τ	41	0	25	0	0
2	A	1628	0	1539	3	0
2	L	1637	0	1564	7	0
3	В	1588	0	1535	1	0
3	Н	1596	0	1542	3	0
4	A	5	0	0	0	0
4	L	5	0	0	0	0
5	A	40	0	0	0	0
5	В	48	0	0	0	0
5	Н	65	0	0	0	0
5	L	66	0	0	0	0
5	Т	1	0	0	0	0
All	All	6720	0	6205	13	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 1.

All (13) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$\begin{array}{c} {\rm Interatomic} \\ {\rm distance} \ ({\rm \AA}) \end{array}$	Clash overlap (Å)
2:A:91:HIS:HA	2:A:96:LEU:HD13	1.84	0.60
2:L:91:HIS:HB2	3:H:100:TYR:O	2.04	0.58
3:H:184:PRO:HB2	3:H:187:THR:HG23	1.89	0.53
2:A:91:HIS:HA	2:A:96:LEU:CD1	2.38	0.53
2:L:212:ASN:C	2:L:212:ASN:HD22	2.11	0.53
2:A:37:GLN:HB2	2:A:47:LEU:HD11	1.95	0.49
2:L:75:ILE:HG21	2:L:78:LEU:HD23	1.97	0.46
3:B:145:TYR:CE2	3:B:150:VAL:HG13	2.51	0.45
2:L:149:LYS:HE3	2:L:195:GLU:OE1	2.17	0.44
2:L:123:GLU:O	2:L:126:THR:OG1	2.29	0.43
2:L:39:LYS:HB2	2:L:42:LYS:HD2	2.02	0.41
2:L:136:LEU:CD2	2:L:196:ALA:HB2	2.51	0.41
3:H:47:TRP:CZ2	3:H:49:GLY:HA2	2.56	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	Perce	entiles
2	A	211/214~(99%)	202 (96%)	9 (4%)	0	100	100
2	L	211/214 (99%)	206 (98%)	5 (2%)	0	100	100
3	В	209/230 (91%)	203 (97%)	6 (3%)	0	100	100
3	Н	209/230 (91%)	206 (99%)	3 (1%)	0	100	100
All	All	840/888 (95%)	817 (97%)	23 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	Perce	ntiles
2	A	179/188~(95%)	173 (97%)	6 (3%)	37	51
2	L	183/188 (97%)	175 (96%)	8 (4%)	28	39
3	В	176/200 (88%)	174 (99%)	2 (1%)	73	86
3	Н	179/200 (90%)	177 (99%)	2 (1%)	73	86
All	All	717/776~(92%)	699 (98%)	18 (2%)	47	65

All (18) residues with a non-rotameric sidechain are listed below:



Mol	Chain	Res	Type
2	L	10	SER
2	L	40	GLN
2	L	96	LEU
2	L	107	LYS
2	L	156	GLN
2	L	206	VAL
2	L	212	ASN
2	L	213	GLU
3	Н	100(C)	MET
3	Н	196	ASN
2	A	50	ASN
2	A	96	LEU
2	A	185	GLU
2	A	202	THR
2	A	208	SER
2	A	212	ASN
3	В	11	LEU
3	В	177	LEU

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

Mol	Chain	Res	Type
2	L	40	GLN
2	L	90	HIS
2	L	137	ASN
2	L	145	ASN
2	L	156	GLN
2	L	161	ASN
2	L	212	ASN
2	A	40	GLN
2	A	90	HIS
2	A	137	ASN
2	A	161	ASN
2	A	212	ASN
3	В	171	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)

2 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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| > 2 is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Dec	Link	В	ond leng	gths	Е	ond ang	gles
MIOI	туре	Chain	Res	Lilik	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	SO4	L	215	-	4,4,4	0.14	0	6,6,6	0.40	0
4	SO4	A	215	-	4,4,4	0.13	0	6,6,6	0.34	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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^{th} 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 $>$	$\#\mathrm{RSRZ}{>}2$	$OWAB(A^2)$	Q<0.9
1	Т	3/3 (100%)	0.32	0 100 100	27, 27, 33, 42	3 (100%)
2	A	213/214 (99%)	0.07	7 (3%) 46 53	11, 22, 31, 41	0
2	L	213/214 (99%)	-0.09	3 (1%) 75 80	7, 17, 23, 35	0
3	В	213/230 (92%)	0.20	8 (3%) 40 47	7, 20, 26, 32	0
3	Н	213/230 (92%)	0.17	10 (4%) 31 38	5, 16, 23, 27	0
All	All	855/891 (95%)	0.09	28 (3%) 46 53	5, 18, 27, 42	3 (0%)

All (28) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	Н	53	TYR	4.6
3	В	97	TYR	3.8
2	A	152	GLY	3.7
3	В	171	GLN	3.4
3	В	96	GLY	3.4
2	A	212	ASN	3.2
3	Н	37	VAL	3.0
3	В	28	THR	3.0
2	A	156	GLN	3.0
2	A	1	GLU	2.9
3	Н	1	GLN	2.9
3	В	53	TYR	2.7
3	Н	48	ILE	2.6
2	L	24	ARG	2.6
2	A	213	GLU	2.6
3	Н	36	TRP	2.6
2	A	153	SER	2.3
3	Н	74	SER	2.2
3	Н	41	PRO	2.2
3	В	141	LEU	2.2

Continued on next page...



Continued from previous page...

Mol	Chain	Res	Type	RSRZ
2	L	212	ASN	2.1
3	Н	100(C)	MET	2.1
3	Н	35	HIS	2.1
3	В	26	GLY	2.1
2	A	30	TYR	2.1
3	В	173	ASP	2.0
2	L	133	VAL	2.0
3	Н	113	SER	2.0

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)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q<0.9
4	SO4	L	215	5/5	0.92	0.16	69,71,72,73	0
4	SO4	A	215	5/5	0.97	0.17	70,71,72,73	0

6.5 Other polymers (i)

There are no such residues in this entry.

