

wwPDB X-ray Structure Validation Summary Report (i)

Aug 28, 2023 – 10:20 AM EDT

PDB ID : 3K48

Title : Crystal structure of APRIL bound to a peptide

Authors : Hymowitz, S.G. Deposited on : 2009-10-05

Resolution : 2.80 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 (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:

 $\begin{array}{ccc} \text{MolProbity} & : & 4.02\text{b-}467 \\ \text{Xtriage (Phenix)} & : & 1.13 \end{array}$

EDS: 2.35

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

Refmac : 5.8.0158

CCP4 : 7.0.044 (Gargrove) roteins) : Engh & Huber (2001)

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

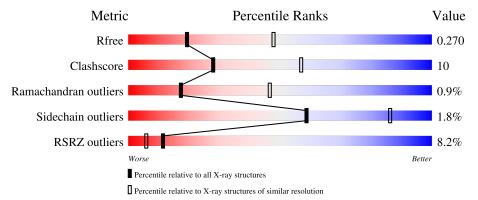
Validation Pipeline (wwPDB-VP) : 2.35

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: $X\text{-}RAY\ DIFFRACTION$

The reported resolution of this entry is 2.80 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	$\begin{array}{c} \text{Whole archive} \\ (\#\text{Entries}) \end{array}$	Similar resolution $(\# \text{Entries, resolution range}(\text{\AA}))$
R_{free}	130704	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

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 <=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		
			9%		
1	A	140	79%	17%	•••
			9%		
1	В	140	81%	16%	• •
			6%		
1	D	140	83%	14%	• •
			19%		
2	R	16	69%	5%	6%
2	S	16	69% 12%	19%	

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Mol	Chain	Length	Quality of chain		
		1.0			
2	(T)	16	75%	6%	19%



2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 3613 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 Tumor necrosis factor ligand superfamily member 13.

Mol	Chain	Residues		At	oms			ZeroOcc	AltConf	Trace
1	A	137		С		О	S	0	0	0
_	11	131	1080	683	198	194	5	U		
1	D	137	Total	\mathbf{C}	N	O	\mathbf{S}	0	0	0
1	Б	137	1080	683	198	194	5	0	U	U
1	D	137	Total	С	N	О	S	0	0	0
1	ש	197	1080	683	198	194	5		U	U

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	102	GLY	-	expression tag	UNP Q9D777
A	103	SER	-	expression tag	UNP Q9D777
В	102	GLY	-	expression tag	UNP Q9D777
В	103	SER	-	expression tag	UNP Q9D777
D	102	GLY	-	expression tag	UNP Q9D777
D	103	SER	-	expression tag	UNP Q9D777

• Molecule 2 is a protein called peptide.

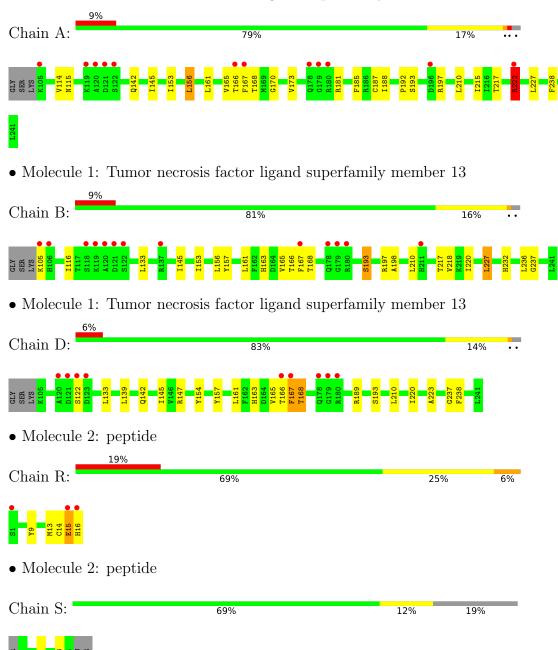
Mol	Chain	Residues		Ato	oms			ZeroOcc	AltConf	Trace
2	R	16	Total	С	N O S	0	0	0		
2	Ιι	10	141	91	23	24	3	U	U	
2	C	13	Total	С	N	О	S	0	0	0
2	b	10	116	77	18	18	3	U	0	0
2	Т	13	Total	С	N	О	S	0	0	0
	1	10	116	77	18	18	3	U	U	U



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: Tumor necrosis factor ligand superfamily member 13





• Molecule 2: peptide

Chain T: 75% 6% 19%





4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants	100.81Å 84.77Å 55.56Å	Donositor
a, b, c, α , β , γ	90.00° 99.45° 90.00°	Depositor
Resolution (Å)	30.00 - 2.80	Depositor
rtesolution (A)	29.19 - 2.80	EDS
% Data completeness	99.6 (30.00-2.80)	Depositor
(in resolution range)	99.6 (29.19-2.80)	EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.11	Depositor
$< I/\sigma(I) > 1$	2.25 (at 2.80Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
D D.	0.205 , 0.267	Depositor
R, R_{free}	0.209 , 0.270	DCC
R_{free} test set	1143 reflections (10.03%)	wwPDB-VP
Wilson B-factor (Å ²)	35.9	Xtriage
Anisotropy	0.072	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.35, 40.4	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	0.91	EDS
Total number of atoms	3613	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 7.88% 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)

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		
IVIOI	Chain	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	A	0.48	0/1104	0.68	1/1495 (0.1%)	
1	В	0.48	0/1104	0.63	0/1495	
1	D	0.49	0/1104	0.64	0/1495	
2	R	0.60	0/149	0.81	0/203	
2	S	0.62	0/123	0.62	0/168	
2	Т	0.56	0/123	0.63	0/168	
All	All	0.50	0/3707	0.65	1/5024 (0.0%)	

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^o)$	$\operatorname{Ideal}({}^{o})$
1	A	222	ARG	NE-CZ-NH1	5.49	123.05	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	1080	0	1080	31	0
1	В	1080	0	1080	25	0
1	D	1080	0	1080	19	0
2	R	141	0	113	6	0
2	S	116	0	92	3	0
2	Т	116	0	92	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	3613	0	3537	71	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 10.

The worst 5 of 71 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$\begin{array}{c} {\rm Interatomic} \\ {\rm distance} \ ({\rm \AA}) \end{array}$	$egin{aligned} ext{Clash} \ ext{overlap } (ext{Å}) \end{aligned}$
1:A:145:ILE:HG22	1:A:217:THR:HG22	1.48	0.96
1:A:197:ARG:HD2	1:D:167:PHE:HA	1.50	0.93
1:D:161:LEU:HD21	2:T:13:MET:CE	2.13	0.79
1:D:161:LEU:HD21	2:T:13:MET:HE2	1.66	0.77
1:D:165:VAL:HG22	1:D:193:SER:HA	1.65	0.76

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
1	A	135/140 (96%)	128 (95%)	6 (4%)	1 (1%)	22	53
1	В	135/140 (96%)	129 (96%)	5 (4%)	1 (1%)	22	53
1	D	135/140 (96%)	126 (93%)	8 (6%)	1 (1%)	22	53
2	R	14/16 (88%)	11 (79%)	2 (14%)	1 (7%)	1	2
2	S	11/16 (69%)	9 (82%)	2 (18%)	0	100	100
2	Т	11/16 (69%)	9 (82%)	2 (18%)	0	100	100
All	All	441/468 (94%)	412 (93%)	25 (6%)	4 (1%)	17	46

All (4) Ramachandran outliers are listed below:



Mol	Chain	Res	Type
1	В	167	PHE
1	D	167	PHE
1	A	167	PHE
2	R	15	GLU

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	A	119/122 (98%)	117 (98%)	2 (2%)	60 87
1	В	119/122 (98%)	116 (98%)	3 (2%)	47 80
1	D	119/122 (98%)	117 (98%)	2 (2%)	60 87
2	R	15/15 (100%)	15 (100%)	0	100 100
2	S	12/15 (80%)	12 (100%)	0	100 100
2	Т	12/15 (80%)	12 (100%)	0	100 100
All	All	396/411 (96%)	389 (98%)	7 (2%)	59 86

5 of 7 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	В	193	SER
1	В	227	LEU
1	D	189	ARG
1	D	168	THR
1	В	105	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	В	142	GLN
1	В	163	HIS
1	В	178	GLN
1	D	142	GLN
1	D	178	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, 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></rsrz>	# RSRZ > 2	$OWAB(A^2)$	Q < 0.9
1	A	137/140 (97%)	0.09	12 (8%) 10 5	6, 16, 34, 43	0
1	В	137/140 (97%)	0.05	13 (9%) 8 4	5, 16, 34, 43	0
1	D	137/140 (97%)	-0.00	9 (6%) 18 11	6, 16, 34, 43	0
2	R	16/16 (100%)	0.60	3 (18%) 1 1	22, 24, 36, 39	0
2	S	13/16 (81%)	0.06	0 100 100	22, 23, 29, 30	0
2	Т	13/16 (81%)	0.02	0 100 100	22, 23, 28, 30	0
All	All	453/468 (96%)	0.07	37 (8%) 11 6	5, 17, 34, 43	0

The worst 5 of 37 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	179	GLY	6.5
1	A	179	GLY	4.1
2	R	16	HIS	4.0
1	В	118	SER	4.0
1	A	120	ALA	3.9

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.

