



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 27, 2014 – 11:18 PM GMT

PDB ID : 2V4B  
Title : CRYSTAL STRUCTURE OF HUMAN ADAMTS-1 CATALYTIC DOMAIN  
AND CYSTEINE-RICH DOMAIN (APO-FORM)  
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Deposited on : 2007-06-28  
Resolution : 2.00 Å(reported)

This is a full wwPDB validation 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 <http://wwpdb.org/ValidationPDFNotes.html>

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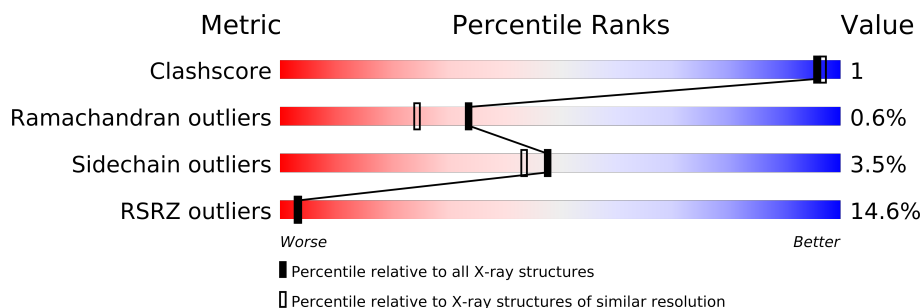
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance



The reported resolution of this entry is 2.00 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))
Clashscore	79885	6188 (2.00-2.00)
Ramachandran outliers	78287	6102 (2.00-2.00)
Sidechain outliers	78261	6100 (2.00-2.00)
RSRZ outliers	66119	4890 (2.00-2.00)

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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	300	
1	B	300	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
2	ZN	A	1552	-	X
4	NI	A	1556	-	X

## 2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 4556 atoms, of which 0 are hydrogen and 0 are deuterium.

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 ADAMTS-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	280	Total	C	N	O	S	0	0	0
			2151	1337	374	415	25			
1	B	281	Total	C	N	O	S	0	0	0
			2164	1343	379	417	25			

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Zn	0	0
			1	1		
2	A	1	Total	Zn	0	0
			1	1		

- Molecule 3 is CADMIUM ION (three-letter code: CD) (formula: Cd).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	2	Total	Cd	0	0
			2	2		
3	A	2	Total	Cd	0	0
			2	2		

- Molecule 4 is NICKEL (II) ION (three-letter code: NI) (formula: Ni).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	5	Total	Ni	0	0
			5	5		
4	A	9	Total	Ni	0	0
			9	9		

- Molecule 5 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	1	Total 1	Mg 1	0	0
5	A	2	Total 2	Mg 2	0	0

- Molecule 6 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	2	Total 2	Na 2	0	0
6	A	2	Total 2	Na 2	0	0

- Molecule 7 is water.

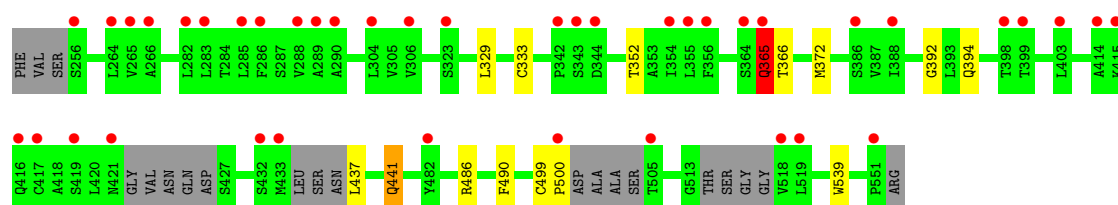
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	133	Total 133	O 133	0	0
7	B	81	Total 81	O 81	0	0

### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

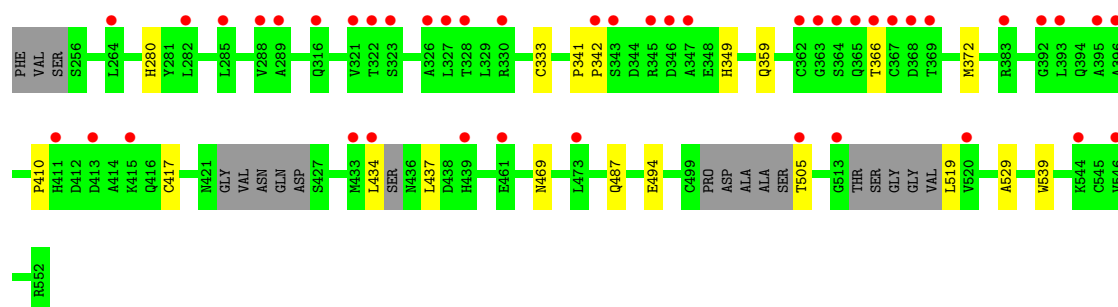
#### • Molecule 1: ADAMTS-1

Chain A: 



#### • Molecule 1: ADAMTS-1

Chain B: 



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	51.58Å 64.40Å 113.52Å 90.00° 90.91° 90.00°	Depositor
Resolution (Å)	40.26 – 2.00 40.25 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.9 (40.26-2.00) 99.9 (40.25-2.00)	Depositor EDS
$R_{merge}$	0.11	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.75 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, $R_{free}$	0.220 , 0.263 0.234 , (Not available)	Depositor DCC
$R_{free}$ test set	No test flags present.	DCC
Wilson B-factor (Å <sup>2</sup> )	27.5	Xtriage
Anisotropy	0.458	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.41 , 46.7	EDS
Estimated twinning fraction	0.053 for h,-k,-l	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.46$ , $\langle L^2 \rangle = 0.29$	Xtriage
Outliers	1 of 50441 reflections (0.002%)	Xtriage
$F_o, F_c$ correlation	0.91	EDS
Total number of atoms	4556	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.53% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: NI, ZN, CD, MG, NA

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	0.50	0/2202	0.59	0/2988
1	B	0.49	0/2214	0.59	0/3002
All	All	0.50	0/4416	0.59	0/5990

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Close contacts

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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2151	0	2021	6	0
1	B	2164	0	2035	4	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
4	A	9	0	0	0	0
4	B	5	0	0	0	0
5	A	2	0	0	0	0
5	B	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	A	2	0	0	0	0
6	B	2	0	0	0	0
7	A	133	0	0	0	0
7	B	81	0	0	0	0
All	All	4556	0	4056	9	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 1.

All (9) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:499:CYS:HA	1:A:500:PRO:C	2.02	0.80
1:B:487:GLN:HE22	1:B:529:ALA:H	1.45	0.64
1:B:341:PRO:O	1:B:349:HIS:HD2	1.91	0.54
1:A:441:GLN:HA	1:A:441:GLN:HE21	1.74	0.52
1:A:365:GLN:HG2	1:A:366:THR:H	1.78	0.47
1:A:392:GLY:HA2	1:A:490:PHE:HB3	1.97	0.47
1:A:329:LEU:HD22	1:A:372:MET:HG3	1.98	0.44
1:A:486:ARG:HD3	1:B:280:HIS:HB2	2.02	0.41
1:B:410:PRO:HG2	1:B:417:CYS:SG	2.60	0.41

There are no symmetry-related clashes.

## 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	270/300 (90%)	267 (99%)	2 (1%)	1 (0%)	43	36
1	B	271/300 (90%)	264 (97%)	5 (2%)	2 (1%)	30	20
All	All	541/600 (90%)	531 (98%)	7 (1%)	3 (1%)	33	24

All (3) Ramachandran outliers are listed below:



Mol	Chain	Res	Type
1	B	366	THR
1	A	365	GLN
1	B	342	PRO

### 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	243/259 (94%)	236 (97%)	7 (3%)	55	52
1	B	244/259 (94%)	234 (96%)	10 (4%)	41	35
All	All	487/518 (94%)	470 (96%)	17 (4%)	48	43

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	333	CYS
1	A	352	THR
1	A	365	GLN
1	A	394	GLN
1	A	437	LEU
1	A	441	GLN
1	A	539	TRP
1	B	333	CYS
1	B	359	GLN
1	B	372	MET
1	B	434	LEU
1	B	437	LEU
1	B	469	ASN
1	B	494	GLU
1	B	505	THR
1	B	519	LEU
1	B	539	TRP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	441	GLN

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Mol	Chain	Res	Type
1	A	457	ASN
1	B	349	HIS
1	B	365	GLN
1	B	408	ASN
1	B	441	GLN
1	B	487	GLN

### 5.3.3 RNA ⓘ

There are no RNA chains 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 carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

Of 27 ligands modelled in this entry, 27 are monoatomic - leaving 0 for Mogul analysis.

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.

## 5.7 Other polymers ⓘ

There are no such residues in this entry.

## 5.8 Polymer linkage issues

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, 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	280/300 (93%)	0.86	41 (14%) 3 3	23, 35, 48, 57	0
1	B	281/300 (93%)	1.06	44 (15%) 3 3	26, 39, 56, 62	0
All	All	561/600 (93%)	0.96	85 (15%) 3 3	23, 37, 53, 62	0

All (85) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	364	SER	6.7
1	B	365	GLN	6.6
1	B	363	GLY	5.3
1	A	551	PRO	5.2
1	A	285	LEU	4.3
1	A	518	VAL	4.3
1	B	366	THR	4.2
1	A	354	ILE	4.1
1	B	369	THR	4.0
1	A	264	LEU	3.9
1	B	326	ALA	3.7
1	B	367	CYS	3.6
1	A	364	SER	3.5
1	B	415	LYS	3.4
1	B	345	ARG	3.3
1	B	546	VAL	3.3
1	B	342	PRO	3.3
1	A	433	MET	3.3
1	A	519	LEU	3.3
1	B	520	VAL	3.2
1	B	323	SER	3.2
1	B	396	ALA	3.2
1	A	415	LYS	3.1
1	B	322	THR	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	432	SER	3.1
1	B	362	CYS	3.0
1	B	413	ASP	3.0
1	B	346	ASP	3.0
1	B	330	ARG	2.9
1	B	321	VAL	2.9
1	B	285	LEU	2.9
1	B	439	HIS	2.9
1	A	289	ALA	2.9
1	A	421	ASN	2.8
1	B	316	GLN	2.8
1	A	265	VAL	2.7
1	B	288	VAL	2.7
1	B	395	ALA	2.7
1	A	417	CYS	2.6
1	A	304	LEU	2.6
1	A	355	LEU	2.6
1	A	342	PRO	2.6
1	B	505	THR	2.6
1	B	328	THR	2.6
1	B	368	ASP	2.6
1	A	288	VAL	2.6
1	B	513	GLY	2.6
1	B	393	LEU	2.5
1	A	256	SER	2.5
1	B	343	SER	2.5
1	A	416	GLN	2.5
1	A	323	SER	2.5
1	A	344	ASP	2.5
1	A	398	THR	2.4
1	A	414	ALA	2.4
1	B	433	MET	2.4
1	A	505	THR	2.4
1	A	286	PHE	2.4
1	B	327	LEU	2.3
1	A	399	THR	2.3
1	A	290	ALA	2.3
1	A	403	LEU	2.3
1	A	266	ALA	2.3
1	B	461	GLU	2.3
1	B	347	ALA	2.2
1	A	282	LEU	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	388	ILE	2.2
1	A	482	TYR	2.2
1	A	365	GLN	2.2
1	B	434	LEU	2.2
1	B	289	ALA	2.2
1	A	343	SER	2.2
1	A	419	SER	2.2
1	A	500	PRO	2.1
1	A	306	VAL	2.1
1	B	411	HIS	2.1
1	A	356	PHE	2.1
1	B	282	LEU	2.1
1	B	383	ARG	2.1
1	B	392	GLY	2.1
1	A	283	LEU	2.1
1	B	544	LYS	2.1
1	B	264	LEU	2.0
1	B	473	LEU	2.0
1	A	386	SER	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

## 6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 6.4 Ligands ⓘ

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> 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	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
2	ZN	A	1552	1/1	0.18	4.90	25,25,25,25	0
4	NI	A	1556	1/1	0.22	2.27	26,26,26,26	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
5	MG	A	1565	1/1	0.17	1.32	52,52,52,52	0
5	MG	A	1560	1/1	0.18	1.20	39,39,39,39	0
5	MG	B	1558	1/1	0.18	0.00	51,51,51,51	0
4	NI	A	1558	1/1	0.15	-0.34	37,37,37,37	0
4	NI	A	1557	1/1	0.12	-0.52	28,28,28,28	0
4	NI	A	1563	1/1	0.14	-0.62	48,48,48,48	0
2	ZN	B	1553	1/1	0.10	-1.40	26,26,26,26	0
4	NI	A	1562	1/1	0.08	-1.81	53,53,53,53	0
4	NI	A	1564	1/1	0.09	-1.87	68,68,68,68	0
6	NA	B	1562	1/1	0.09	-1.90	45,45,45,45	0
4	NI	A	1555	1/1	0.10	-1.90	40,40,40,40	0
3	CD	A	1553	1/1	0.06	-2.16	38,38,38,38	0
4	NI	B	1559	1/1	0.08	-2.28	30,30,30,30	0
6	NA	B	1563	1/1	0.08	-2.34	45,45,45,45	0
3	CD	B	1556	1/1	0.03	-2.40	41,41,41,41	0
4	NI	B	1555	1/1	0.09	-2.46	49,49,49,49	0
4	NI	B	1557	1/1	0.07	-2.48	29,29,29,29	0
3	CD	A	1554	1/1	0.06	-3.05	35,35,35,35	0
4	NI	B	1560	1/1	0.05	-4.46	64,64,64,64	0
3	CD	B	1554	1/1	0.12	-4.68	43,43,43,43	0
6	NA	A	1566	1/1	0.06	-5.26	29,29,29,29	0
6	NA	A	1567	1/1	0.09	-7.22	40,40,40,40	0
4	NI	A	1559	1/1	0.05	-8.61	59,59,59,59	0
4	NI	B	1561	1/1	0.11	-13.93	44,44,44,44	0
4	NI	A	1561	1/1	0.06	-15.98	34,34,34,34	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.