



# Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 03:56 PM GMT

PDB ID : 3A8N  
Title : Crystal structure of the Tiam1 PHCCEX domain  
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Deposited on : 2009-10-07  
Resolution : 4.50 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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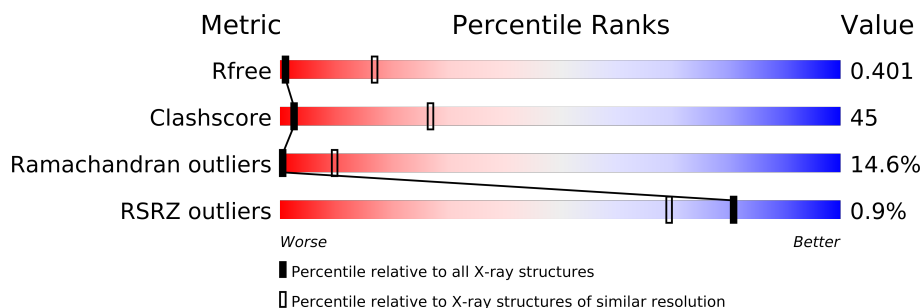
The following versions of software and data (see [references](#)) were used in the production of this report:

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 4.50 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))
$R_{free}$	66092	1029 (5.50-3.50)
Clashscore	79885	1300 (5.50-3.50)
Ramachandran outliers	78287	1222 (5.50-3.50)
RSRZ outliers	66119	1028 (5.50-3.50)

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 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	279	

## 2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 1142 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 T-lymphoma invasion and metastasis-inducing protein 1.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
1	A	230	Total	C	N	O	0	0	0
			1142	682	230	230			

There are 5 discrepancies between the modelled and reference sequences:

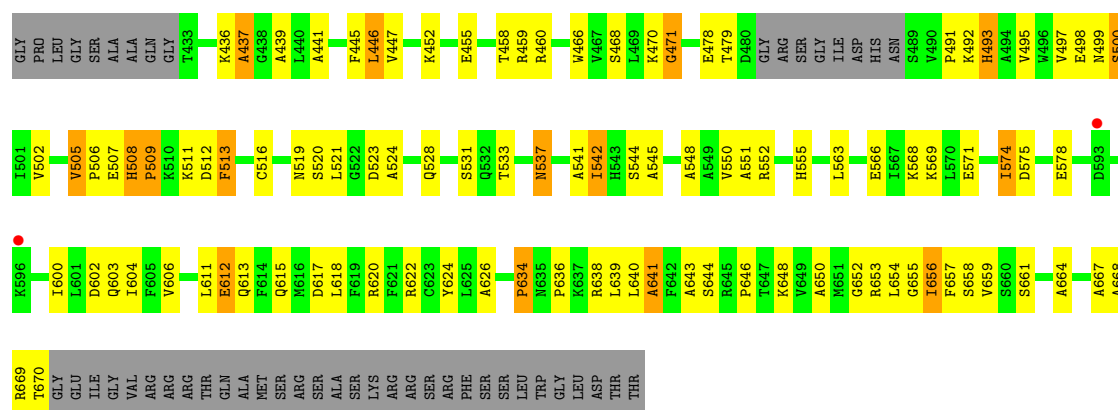
Chain	Residue	Modelled	Actual	Comment	Reference
A	424	GLY	-	EXPRESSION TAG	UNP Q60610
A	425	PRO	-	EXPRESSION TAG	UNP Q60610
A	426	LEU	-	EXPRESSION TAG	UNP Q60610
A	427	GLY	-	EXPRESSION TAG	UNP Q60610
A	428	SER	-	EXPRESSION TAG	UNP Q60610

### 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: T-lymphoma invasion and metastasis-inducing protein 1

Chain A: 



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 64 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	113.47Å 113.47Å 113.84Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	49.25 – 4.50 49.25 – 4.50	Depositor EDS
% Data completeness (in resolution range)	99.0 (49.25-4.50) 98.9 (49.25-4.50)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.25 (at 4.45Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, $R_{free}$	0.306 , 0.396 0.317 , 0.401	Depositor DCC
$R_{free}$ test set	262 reflections (10.25%)	DCC
Wilson B-factor (Å <sup>2</sup> )	153.9	Xtriage
Anisotropy	0.337	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.29 , 553.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 2817 reflections	Xtriage
$F_o, F_c$ correlation	0.83	EDS
Total number of atoms	1142	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	107.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.81% 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

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	0.44	0/1140	0.80	5/1588 (0.3%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	491	PRO	N-CA-CB	5.58	109.99	103.30
1	A	646	PRO	N-CA-CB	5.57	109.98	103.30
1	A	634	PRO	N-CA-CB	5.51	109.92	103.30
1	A	478	GLU	N-CA-C	5.28	125.26	111.00
1	A	636	PRO	N-CA-CB	5.22	109.57	103.30

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	1142	0	520	75	0
All	All	1142	0	520	75	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 45.

All (75) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:439:ALA:HA	1:A:466:TRP:HA	1.69	0.75
1:A:541:ALA:O	1:A:544:SER:N	2.18	0.73
1:A:437:ALA:CB	1:A:468:SER:HA	2.20	0.72
1:A:502:VAL:HA	1:A:516:CYS:O	1.90	0.72
1:A:446:LEU:O	1:A:524:ALA:N	2.24	0.71
1:A:656:ILE:O	1:A:658:SER:N	2.24	0.70
1:A:604:ILE:C	1:A:606:VAL:H	1.96	0.69
1:A:446:LEU:N	1:A:524:ALA:HB3	2.10	0.67
1:A:437:ALA:HB2	1:A:468:SER:HA	1.79	0.65
1:A:604:ILE:C	1:A:606:VAL:N	2.47	0.64
1:A:658:SER:H	1:A:661:SER:CB	2.11	0.63
1:A:545:ALA:O	1:A:548:ALA:HB3	2.00	0.62
1:A:498:GLU:C	1:A:500:SER:H	2.03	0.61
1:A:640:LEU:O	1:A:643:ALA:N	2.33	0.60
1:A:471:GLY:O	1:A:545:ALA:HB1	2.02	0.60
1:A:622:ARG:C	1:A:624:TYR:H	2.05	0.59
1:A:505:VAL:O	1:A:507:GLU:N	2.35	0.59
1:A:436:LYS:O	1:A:437:ALA:HB2	2.00	0.59
1:A:650:ALA:C	1:A:652:GLY:H	2.06	0.58
1:A:667:ALA:C	1:A:669:ARG:N	2.55	0.58
1:A:650:ALA:C	1:A:652:GLY:N	2.57	0.57
1:A:667:ALA:C	1:A:669:ARG:H	2.07	0.57
1:A:667:ALA:O	1:A:669:ARG:N	2.38	0.56
1:A:505:VAL:C	1:A:507:GLU:H	2.09	0.56
1:A:669:ARG:O	1:A:670:THR:C	2.45	0.55
1:A:602:ASP:C	1:A:604:ILE:H	2.09	0.55
1:A:498:GLU:O	1:A:500:SER:N	2.41	0.54
1:A:650:ALA:O	1:A:653:ARG:N	2.38	0.54
1:A:541:ALA:O	1:A:542:ILE:C	2.47	0.53
1:A:445:PHE:O	1:A:446:LEU:CB	2.59	0.51
1:A:571:GLU:O	1:A:574:ILE:N	2.44	0.51
1:A:441:ALA:HB3	1:A:528:GLN:CB	2.40	0.51
1:A:655:GLY:O	1:A:656:ILE:O	2.29	0.50
1:A:505:VAL:C	1:A:507:GLU:N	2.65	0.50
1:A:563:LEU:O	1:A:566:GLU:N	2.45	0.50
1:A:661:SER:O	1:A:664:ALA:HB3	2.11	0.50
1:A:655:GLY:O	1:A:656:ILE:C	2.49	0.49
1:A:615:GLN:C	1:A:617:ASP:N	2.66	0.49
1:A:624:TYR:C	1:A:626:ALA:H	2.16	0.49
1:A:441:ALA:HB3	1:A:528:GLN:H	1.78	0.48
1:A:615:GLN:O	1:A:618:LEU:N	2.47	0.47
1:A:519:ASN:O	1:A:521:LEU:N	2.48	0.47
1:A:519:ASN:C	1:A:521:LEU:N	2.67	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:512:ASP:O	1:A:513:PHE:O	2.33	0.46
1:A:658:SER:O	1:A:659:VAL:C	2.52	0.46
1:A:446:LEU:HA	1:A:455:GLU:O	2.15	0.46
1:A:638:ARG:O	1:A:641:ALA:HB3	2.16	0.46
1:A:658:SER:O	1:A:661:SER:N	2.49	0.46
1:A:602:ASP:O	1:A:604:ILE:N	2.49	0.45
1:A:615:GLN:C	1:A:617:ASP:H	2.20	0.45
1:A:622:ARG:C	1:A:624:TYR:N	2.66	0.45
1:A:612:GLU:O	1:A:613:GLN:C	2.55	0.45
1:A:537:ASN:O	1:A:541:ALA:N	2.37	0.45
1:A:548:ALA:O	1:A:551:ALA:N	2.49	0.45
1:A:644:SER:O	1:A:648:LYS:N	2.49	0.45
1:A:611:LEU:O	1:A:612:GLU:C	2.55	0.45
1:A:492:LYS:O	1:A:493:HIS:CB	2.65	0.44
1:A:618:LEU:C	1:A:620:ARG:N	2.72	0.44
1:A:519:ASN:C	1:A:521:LEU:H	2.21	0.43
1:A:600:ILE:C	1:A:602:ASP:N	2.69	0.43
1:A:498:GLU:C	1:A:500:SER:N	2.71	0.43
1:A:541:ALA:O	1:A:544:SER:CB	2.67	0.42
1:A:658:SER:O	1:A:661:SER:CB	2.67	0.42
1:A:658:SER:N	1:A:661:SER:CB	2.81	0.42
1:A:602:ASP:C	1:A:604:ILE:N	2.72	0.42
1:A:447:VAL:HA	1:A:523:ASP:HA	2.02	0.42
1:A:470:LYS:O	1:A:471:GLY:C	2.58	0.41
1:A:552:ARG:O	1:A:555:HIS:N	2.44	0.41
1:A:508:HIS:O	1:A:509:PRO:O	2.38	0.41
1:A:604:ILE:O	1:A:606:VAL:N	2.55	0.40
1:A:548:ALA:C	1:A:550:VAL:N	2.74	0.40
1:A:574:ILE:O	1:A:578:GLU:N	2.29	0.40
1:A:618:LEU:O	1:A:620:ARG:N	2.55	0.40
1:A:531:SER:C	1:A:533:THR:N	2.75	0.40
1:A:458:THR:O	1:A:460:ARG:N	2.54	0.40

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	226/279 (81%)	143 (63%)	50 (22%)	33 (15%)	0	10

All (33) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	446	LEU
1	A	459	ARG
1	A	479	THR
1	A	493	HIS
1	A	500	SER
1	A	509	PRO
1	A	537	ASN
1	A	656	ILE
1	A	471	GLY
1	A	497	VAL
1	A	499	ASN
1	A	513	PHE
1	A	568	LYS
1	A	574	ILE
1	A	575	ASP
1	A	603	GLN
1	A	654	LEU
1	A	657	PHE
1	A	452	LYS
1	A	495	VAL
1	A	511	LYS
1	A	569	LYS
1	A	668	ALA
1	A	437	ALA
1	A	506	PRO
1	A	639	LEU
1	A	641	ALA
1	A	505	VAL
1	A	520	SER
1	A	612	GLU
1	A	542	ILE
1	A	634	PRO
1	A	508	HIS

### 5.3.2 Protein sidechains ⓘ

There are no protein residues with a non-rotameric sidechain to report in this entry.

### 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 ⓘ

There are no ligands in this entry.

### 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	230/279 (82%)	-0.23	2 (0%) 81 69	4, 99, 179, 203	0

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	596	LYS	3.4
1	A	593	ASP	2.1

### 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 ⓘ

There are no ligands in this entry.

### 6.5 Other polymers ⓘ

There are no such residues in this entry.