



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 05:01 am GMT

PDB ID : 4CZA  
Title : Structure of the sodium proton antiporter PaNhaP from *Pyrococcus abyssi* with bound thallium ion.  
Authors : Woehlert, D.; Kuhlbrandt, W.; Yildiz, O.  
Deposited on : 2014-04-16  
Resolution : 3.20 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](mailto:validation@mail.wwpdb.org)

A user guide is available at

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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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 : 1.7.2 (RC1), CSD as538be (2017)  
Xtriage (Phenix) : 1.9-1692  
EDS : trunk28620  
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : recalc28949

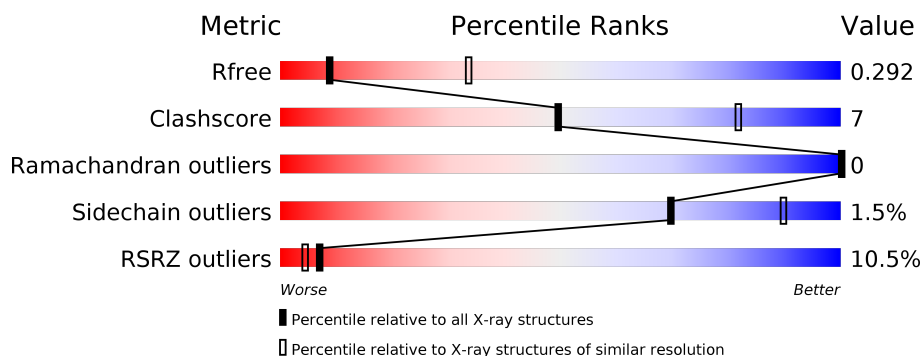
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 3.20 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))
$R_{free}$	100719	1015 (3.22-3.18)
Clashscore	112137	1009 (3.20-3.20)
Ramachandran outliers	110173	1118 (3.22-3.18)
Sidechain outliers	110143	1117 (3.22-3.18)
RSRZ outliers	101464	1020 (3.22-3.18)

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. 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	420	<div> <div>7%</div> <div>83%</div> <div>16%</div> <div>.</div> </div>
1	B	420	<div> <div>13%</div> <div>80%</div> <div>20%</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	TAM	A	1421	-	-	-	X
5	UND	B	1423	-	-	-	X
6	ACT	B	1424	-	-	-	X

## 2 Entry composition [i](#)

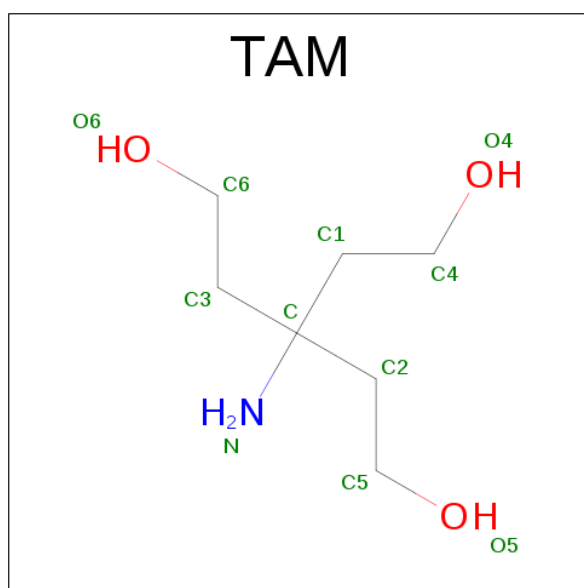
There are 7 unique types of molecules in this entry. The entry contains 6651 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 NA<sup>+</sup>/H<sup>+</sup> ANTIPORTER, PUTATIVE.

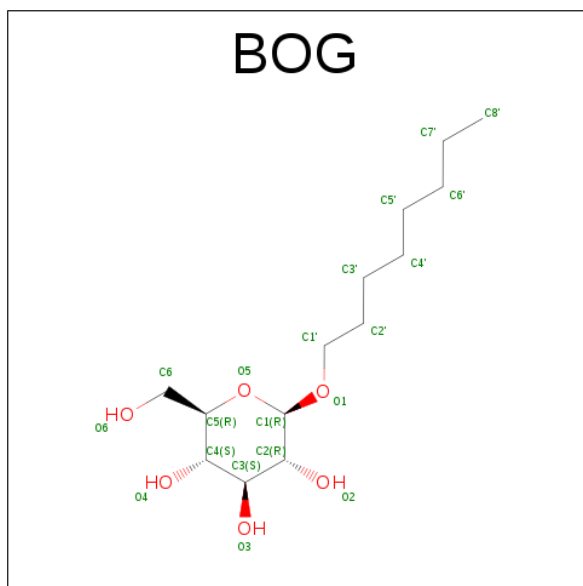
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	420	Total	C	N	O	Se	0	0	0
			3284	2210	514	553	7			
1	B	419	Total	C	N	O	Se	0	0	0
			3276	2206	513	550	7			

- Molecule 2 is TRIS(HYDROXYETHYL)AMINOMETHANE (three-letter code: TAM) (formula: C<sub>7</sub>H<sub>17</sub>NO<sub>3</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			11	7	1	3		
2	B	1	Total	C	N	O	0	0
			11	7	1	3		
2	B	1	Total	C	N	O	0	0
			11	7	1	3		
2	B	1	Total	C	N	O	0	0
			11	7	1	3		

- Molecule 3 is SUGAR (B-OCTYLGLUCOSIDE) (three-letter code: BOG) (formula:  $C_{14}H_{28}O_6$ ).

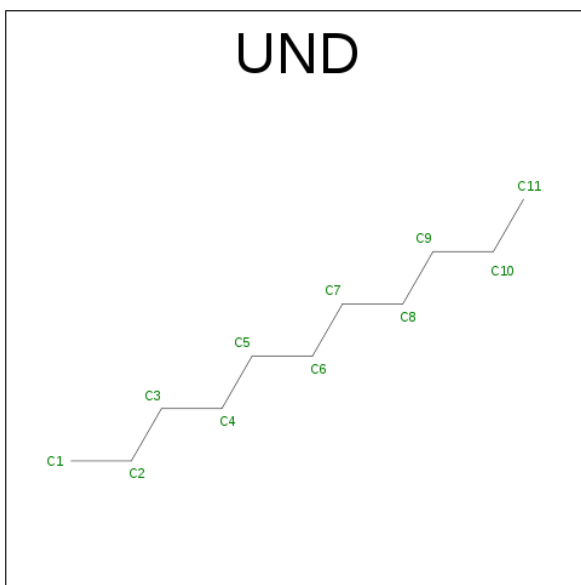


Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	C O	0	0
			20	14 6		

- Molecule 4 is THALLIUM (I) ION (three-letter code: TL) (formula: Tl).

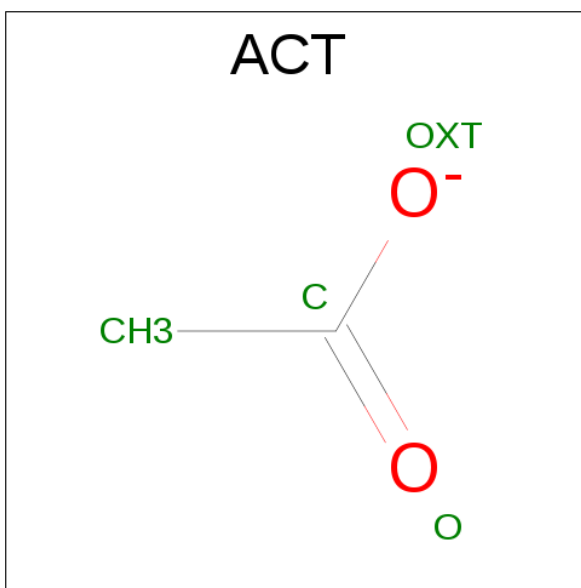
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Tl	0	0
			1	1		
4	A	1	Total	Tl	0	0
			1	1		

- Molecule 5 is UNDECANE (three-letter code: UND) (formula:  $C_{11}H_{24}$ ).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	1	Total	C	0	0
			11	11		

- Molecule 6 is ACETATE ION (three-letter code: ACT) (formula:  $\text{C}_2\text{H}_3\text{O}_2$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	B	1	Total	C	O	0	0
			4	2	2		

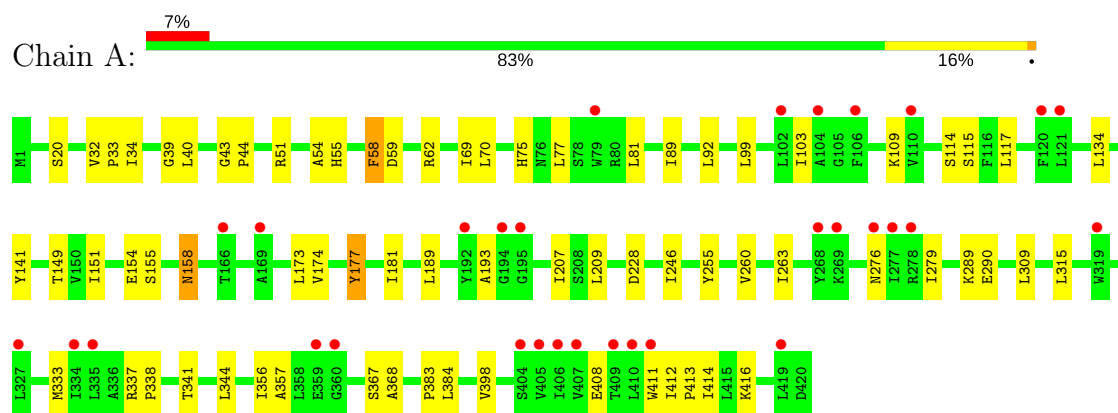
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	4	Total 4	O 4	0	0
7	B	6	Total 6	O 6	0	0

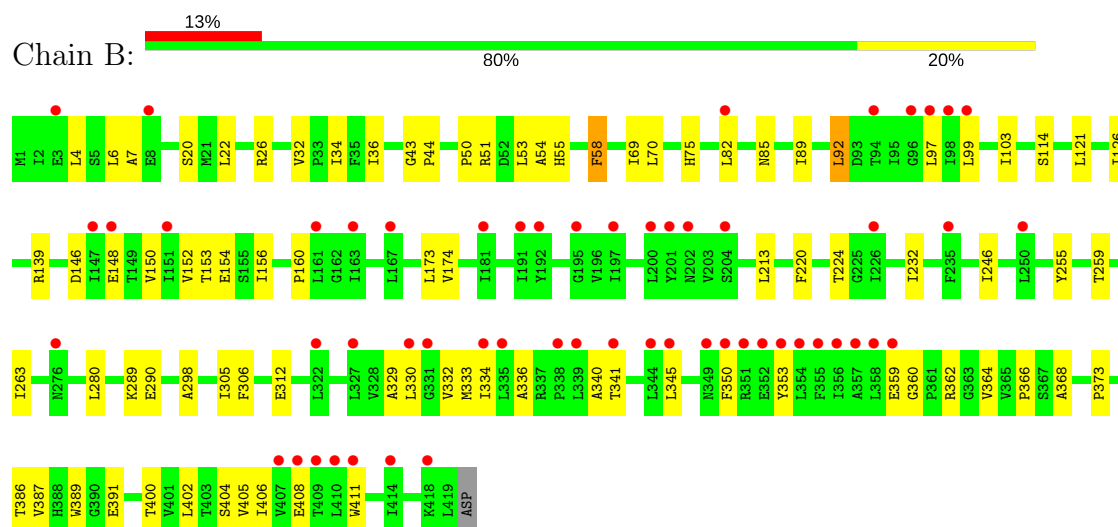
### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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 ( $\text{RSRZ} > 2$ ). 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: NA<sup>+</sup>/H<sup>+</sup> ANTIPTORTER, PUTATIVE



- Molecule 1: NA<sup>+</sup>/H<sup>+</sup> ANTIPTORTER, PUTATIVE





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	54.06Å 107.40Å 99.80Å 90.00° 96.37° 90.00°	Depositor
Resolution (Å)	49.60 – 3.20 49.60 – 3.20	Depositor EDS
% Data completeness (in resolution range)	99.5 (49.60-3.20) 98.9 (49.60-3.20)	Depositor EDS
$R_{merge}$	0.04	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.01 (at 3.19Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, $R_{free}$	0.248 , 0.297 0.250 , 0.292	Depositor DCC
$R_{free}$ test set	950 reflections (5.12%)	DCC
Wilson B-factor (Å <sup>2</sup> )	117.4	Xtriage
Anisotropy	0.564	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.30 , 70.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.87	EDS
Total number of atoms	6651	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	81.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.73% of the height of the origin peak. No significant pseudotranslation is detected.*

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

<sup>2</sup>Theoretical 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

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: TAM, TL, BOG, UND, ACT

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	0.22	0/3351	0.40	0/4549
1	B	0.23	0/3343	0.40	0/4538
All	All	0.23	0/6694	0.40	0/9087

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-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 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	3284	0	3522	42	0
1	B	3276	0	3518	47	0
2	A	11	0	17	1	0
2	B	33	0	51	1	0
3	A	20	0	28	1	0
4	A	1	0	0	1	0
4	B	1	0	0	0	0
5	B	11	0	24	0	0
6	B	4	0	3	0	0
7	A	4	0	0	0	0
7	B	6	0	0	0	0
All	All	6651	0	7163	90	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 7.

All (90) 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)
1:A:69:ILE:HG23	2:A:1421:TAM:HN1	1.54	0.72
1:A:155:SER:OG	4:A:1423:TL:TL	2.19	0.65
1:B:82:LEU:HD23	1:B:153:THR:HG21	1.83	0.60
1:A:40:LEU:HD11	1:A:315:LEU:HD21	1.84	0.60
1:B:75:HIS:NE2	1:B:290:GLU:OE2	2.30	0.60
1:A:228:ASP:OD1	1:A:289:LYS:NZ	2.35	0.60
1:A:337:ARG:O	1:A:341:THR:OG1	2.15	0.60
1:B:85:ASN:ND2	1:B:146:ASP:OD1	2.36	0.59
1:B:70:LEU:HD12	1:B:255:TYR:HB3	1.87	0.57
1:B:345:LEU:HD11	1:B:350:PHE:HD1	1.70	0.56
1:A:134:LEU:HD22	1:A:412:ILE:HD11	1.86	0.56
1:B:4:LEU:HD13	1:B:6:LEU:H	1.70	0.56
1:B:213:LEU:HD11	1:B:246:ILE:HD11	1.87	0.56
1:A:207:ILE:HA	1:A:260:VAL:HG21	1.89	0.55
2:B:1420:TAM:N	2:B:1420:TAM:O6	2.40	0.55
1:B:50:PRO:HD2	1:B:53:LEU:HD12	1.90	0.54
1:B:360:GLY:O	1:B:362:ARG:NH1	2.36	0.53
1:A:20:SER:HB3	1:A:34:ILE:HG22	1.90	0.53
1:A:70:LEU:HD12	1:A:255:TYR:HB3	1.90	0.53
1:B:332:VAL:HA	1:B:336:ALA:HB3	1.89	0.53
1:A:62:ARG:NH2	1:A:368:ALA:O	2.38	0.52
3:A:1422:BOG:H1'1	1:B:289:LYS:HE2	1.91	0.52
1:B:126:ILE:HG12	1:B:362:ARG:HG3	1.91	0.52
1:A:92:LEU:HD21	1:A:344:LEU:HD22	1.92	0.52
1:B:92:LEU:HD11	1:B:154:GLU:HG3	1.92	0.52
1:A:77:LEU:HD22	1:A:263:ILE:HG23	1.92	0.51
1:A:338:PRO:HA	1:A:357:ALA:HB1	1.93	0.51
1:B:329:ALA:HA	1:B:400:THR:HG23	1.93	0.51
1:B:51:ARG:O	1:B:55:HIS:ND1	2.28	0.50
1:B:69:ILE:HD13	1:B:298:ALA:HB1	1.93	0.50
1:B:139:ARG:NH1	1:B:148:GLU:OE1	2.45	0.50
1:A:141:TYR:HB3	1:A:416:LYS:HE2	1.94	0.50
1:A:413:PRO:HG2	1:A:414:ILE:HD12	1.92	0.50
1:A:209:LEU:HD21	1:A:246:ILE:HG22	1.94	0.50
1:B:54:ALA:O	1:B:58:PHE:HB2	2.12	0.49
1:B:306:PHE:HE2	1:B:364:VAL:HG22	1.78	0.49
1:A:81:LEU:HB3	1:A:149:THR:HG21	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:51:ARG:HG2	1:A:55:HIS:CE1	2.47	0.48
1:A:51:ARG:O	1:A:55:HIS:ND1	2.45	0.48
1:A:75:HIS:NE2	1:A:290:GLU:OE2	2.44	0.48
1:B:330:LEU:HA	1:B:333:MSE:HG2	1.95	0.48
1:A:276:ASN:HB2	1:A:279:ILE:HG12	1.95	0.48
1:A:181:ILE:HD12	1:A:181:ILE:H	1.77	0.48
1:A:189:LEU:HB3	1:A:193:ALA:HB3	1.96	0.47
1:B:20:SER:HB3	1:B:34:ILE:HG22	1.96	0.47
1:B:51:ARG:NH1	1:B:312:GLU:O	2.48	0.46
1:B:152:VAL:O	1:B:156:ILE:HG12	2.14	0.46
1:A:32:VAL:HB	1:A:33:PRO:HD3	1.98	0.46
1:B:220:PHE:O	1:B:224:THR:HG22	2.17	0.45
1:B:70:LEU:HD13	1:B:160:PRO:HB3	1.98	0.45
1:A:109:LYS:HA	1:A:114:SER:HB3	1.97	0.45
1:B:121:LEU:HD21	1:B:373:PRO:HB3	1.99	0.45
1:B:362:ARG:HB3	1:B:366:PRO:HG2	1.98	0.45
1:A:39:GLY:O	1:A:43:GLY:N	2.51	0.44
1:A:309:LEU:HD22	1:A:367:SER:HB3	1.99	0.44
1:A:151:ILE:HD11	1:A:356:ILE:HG23	1.99	0.44
1:B:402:LEU:O	1:B:406:ILE:HG22	2.16	0.44
1:A:99:LEU:O	1:A:103:ILE:HG22	2.17	0.44
1:B:386:THR:OG1	1:B:387:VAL:N	2.50	0.44
1:B:173:LEU:HD12	1:B:174:VAL:N	2.33	0.44
1:A:333:MSE:HE3	1:A:408:GLU:HG3	2.00	0.43
1:B:4:LEU:HD12	1:B:7:ALA:H	1.82	0.43
1:A:59:ASP:HA	1:A:62:ARG:HD2	2.00	0.43
1:A:51:ARG:HG2	1:A:55:HIS:HE1	1.84	0.43
1:B:22:LEU:O	1:B:26:ARG:HG2	2.19	0.43
1:A:383:PRO:HG2	1:A:384:LEU:HD12	2.01	0.43
1:A:309:LEU:HD11	1:A:398:VAL:HG22	2.01	0.42
1:A:43:GLY:HA3	1:A:44:PRO:HD3	1.89	0.42
1:B:333:MSE:HG3	1:B:334:ILE:HG12	2.00	0.42
1:A:174:VAL:HG23	1:A:177:TYR:HB2	2.00	0.42
1:B:51:ARG:NH2	1:B:391:GLU:OE2	2.52	0.42
1:B:359:GLU:O	1:B:408:GLU:HG2	2.19	0.42
1:A:54:ALA:O	1:A:58:PHE:HB2	2.20	0.42
1:A:154:GLU:OE2	1:A:337:ARG:NH1	2.43	0.42
1:B:305:ILE:HD13	1:B:368:ALA:HB2	2.02	0.42
1:B:32:VAL:HG11	1:B:405:VAL:HG21	2.02	0.42
1:B:114:SER:HB3	1:B:389:TRP:CH2	2.54	0.41
1:A:117:LEU:HB3	1:A:173:LEU:HD21	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:414:ILE:H	1:A:414:ILE:HD12	1.86	0.41
1:B:36:ILE:HG21	1:B:402:LEU:HD21	2.02	0.41
1:B:89:ILE:HD11	1:B:150:VAL:HG12	2.01	0.41
1:B:259:THR:O	1:B:263:ILE:HG13	2.20	0.41
1:B:92:LEU:O	1:B:97:LEU:N	2.53	0.41
1:B:232:ILE:HG13	1:B:290:GLU:HB2	2.02	0.41
1:B:404:SER:O	1:B:408:GLU:HB2	2.21	0.41
1:B:43:GLY:HA3	1:B:44:PRO:HD3	1.93	0.41
1:A:155:SER:HA	1:A:158:ASN:ND2	2.35	0.41
1:A:114:SER:OG	1:A:115:SER:N	2.54	0.41
1:B:99:LEU:HD23	1:B:340:ALA:HA	2.04	0.40
1:B:99:LEU:O	1:B:103:ILE:HG12	2.21	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	Percentiles	
1	A	418/420 (100%)	398 (95%)	20 (5%)	0	100	100
1	B	417/420 (99%)	394 (94%)	23 (6%)	0	100	100
All	All	835/840 (99%)	792 (95%)	43 (5%)	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	Percentiles	
1	A	358/351 (102%)	353 (99%)	5 (1%)	71	90
1	B	357/351 (102%)	351 (98%)	6 (2%)	66	88
All	All	715/702 (102%)	704 (98%)	11 (2%)	70	90

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	58	PHE
1	A	89	ILE
1	A	158	ASN
1	A	177	TYR
1	A	411	TRP
1	B	58	PHE
1	B	92	LEU
1	B	280	LEU
1	B	341	THR
1	B	353	TYR
1	B	411	TRP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 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 carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 9 ligands modelled in this entry, 2 are monoatomic - leaving 7 for Mogul analysis.

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	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
2	TAM	A	1421	-	7,10,10	1.21	0	9,12,12	0.49	0
3	BOG	A	1422	-	20,20,20	0.49	0	25,25,25	0.86	1 (4%)
2	TAM	B	1420	-	7,10,10	1.19	0	9,12,12	0.63	0
2	TAM	B	1421	-	7,10,10	1.22	0	9,12,12	0.63	0
2	TAM	B	1422	-	7,10,10	1.22	0	9,12,12	0.66	0
5	UND	B	1423	-	10,10,10	0.10	0	9,9,9	0.82	0
6	ACT	B	1424	-	1,3,3	2.24	1 (100%)	0,3,3	0.00	-

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	TAM	A	1421	-	-	0/12/12/12	0/0/0/0
3	BOG	A	1422	-	-	0/11/31/31	0/1/1/1
2	TAM	B	1420	-	-	0/12/12/12	0/0/0/0
2	TAM	B	1421	-	-	0/12/12/12	0/0/0/0
2	TAM	B	1422	-	-	0/12/12/12	0/0/0/0
5	UND	B	1423	-	-	0/8/8/8	0/0/0/0
6	ACT	B	1424	-	-	0/0/0/0	0/0/0/0

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	B	1424	ACT	CH3-C	2.24	1.51	1.48

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1422	BOG	C1-O5-C5	-2.80	108.44	113.72

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1421	TAM	1	0
3	A	1422	BOG	1	0
2	B	1420	TAM	1	0

## 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, 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	413/420 (98%)	0.22	31 (7%) 15 8	26, 72, 113, 138	0
1	B	412/420 (98%)	0.50	56 (13%) 3 2	36, 81, 133, 162	0
All	All	825/840 (98%)	0.36	87 (10%) 7 4	26, 77, 128, 162	0

All (87) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	358	LEU	6.9
1	B	354	LEU	6.5
1	A	278	ARG	4.8
1	A	192	TYR	4.7
1	A	277	ILE	4.7
1	B	351	ARG	4.6
1	B	94	THR	4.6
1	B	407	VAL	4.5
1	B	339	LEU	4.4
1	A	407	VAL	4.3
1	B	161	LEU	4.3
1	B	355	PHE	4.2
1	B	411	TRP	4.1
1	B	335	LEU	4.1
1	B	167	LEU	4.0
1	A	79	TRP	3.9
1	B	197	ILE	3.9
1	B	350	PHE	3.9
1	B	414	ILE	3.7
1	A	359	GLU	3.7
1	A	410	LEU	3.7
1	B	356	ILE	3.7
1	A	269	LYS	3.6
1	B	98	ILE	3.6

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Mol	Chain	Res	Type	RSRZ
1	B	359	GLU	3.6
1	B	334	ILE	3.6
1	B	235	PHE	3.6
1	B	327	LEU	3.6
1	B	357	ALA	3.5
1	B	200	LEU	3.4
1	A	335	LEU	3.4
1	B	82	LEU	3.4
1	A	106	PHE	3.4
1	A	319	TRP	3.4
1	B	97	LEU	3.4
1	B	322	LEU	3.4
1	B	3	GLU	3.2
1	B	192	TYR	3.2
1	B	410	LEU	3.1
1	A	409	THR	3.1
1	B	349	ASN	3.1
1	B	353	TYR	3.0
1	A	195	GLY	2.9
1	B	338	PRO	2.9
1	B	418	LYS	2.8
1	B	409	THR	2.8
1	A	268	TYR	2.8
1	A	405	VAL	2.7
1	B	195	GLY	2.7
1	A	406	ILE	2.7
1	A	102	LEU	2.7
1	B	344	LEU	2.7
1	B	201	TYR	2.7
1	B	8	GLU	2.6
1	B	276	ASN	2.6
1	B	341	THR	2.6
1	B	181	ILE	2.6
1	A	276	ASN	2.6
1	A	334	ILE	2.6
1	B	250	LEU	2.5
1	A	110	VAL	2.5
1	A	120	PHE	2.5
1	A	169	ALA	2.5
1	B	352	GLU	2.5
1	A	419	LEU	2.5
1	B	345	LEU	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	147	ILE	2.4
1	A	327	LEU	2.4
1	A	360	GLY	2.4
1	B	202	ASN	2.4
1	A	404	SER	2.4
1	A	166	THR	2.4
1	A	411	TRP	2.3
1	B	96	GLY	2.3
1	B	151	ILE	2.2
1	A	194	GLY	2.2
1	A	121	LEU	2.2
1	B	204	SER	2.1
1	B	191	ILE	2.1
1	B	408	GLU	2.1
1	B	148	GLU	2.1
1	B	99	LEU	2.1
1	B	330	LEU	2.1
1	B	163	ILE	2.1
1	B	226	ILE	2.0
1	A	104	ALA	2.0
1	B	331	GLY	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 carbohydrates 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. 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	RSCC	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
2	TAM	A	1421	11/11	0.84	0.38	3.35	23,51,65,66	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
6	ACT	B	1424	4/4	0.78	0.45	2.94	71,76,82,85	0
3	BOG	A	1422	20/20	0.81	0.23	1.91	58,68,80,81	0
5	UND	B	1423	11/11	0.46	0.62	1.76	76,89,101,103	0
2	TAM	B	1421	11/11	0.86	0.31	0.82	86,98,103,110	0
2	TAM	B	1420	11/11	0.68	0.19	-0.04	81,101,111,111	0
2	TAM	B	1422	11/11	0.78	0.18	-0.65	88,109,120,123	0
4	TL	B	1425	1/1	0.89	0.05	-2.47	84,84,84,84	1
4	TL	A	1423	1/1	0.88	0.08	-3.17	56,56,56,56	1

## 6.5 Other polymers [i](#)

There are no such residues in this entry.