



# Full wwPDB X-ray Structure Validation Report ⓘ

May 13, 2020 – 07:59 pm BST

PDB ID : 5N9W  
Title : Structure of adenylation domain THR1 involved in the biosynthesis of 4-chlorothreonine in *Streptomyces* SP.OH-5093, apo structure  
Authors : Savino, C.; Vallone, B.; Scaglione, A.; Parisi, G.; Montemiglio, L.C.; Fullone, M.R.; Grgurina, I.  
Deposited on : 2017-02-27  
Resolution : 2.46 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

<https://www.wwpdb.org/validation/2017/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 : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.11  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
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) : 2.11

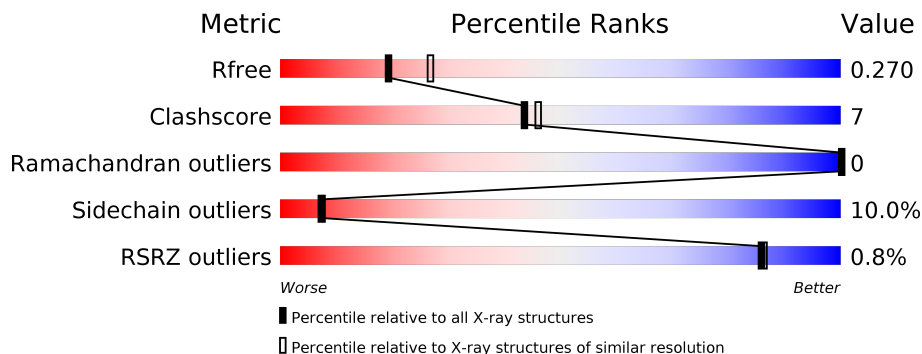
# 1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.46 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))
$R_{free}$	130704	1544 (2.48-2.44)
Clashscore	141614	1613 (2.48-2.44)
Ramachandran outliers	138981	1598 (2.48-2.44)
Sidechain outliers	138945	1598 (2.48-2.44)
RSRZ outliers	127900	1523 (2.48-2.44)

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 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  $\leq 5\%$ . 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	529	 % 69% 18% • 10%
1	B	529	 % 74% 14% • 10%

## 2 Entry composition [i](#)

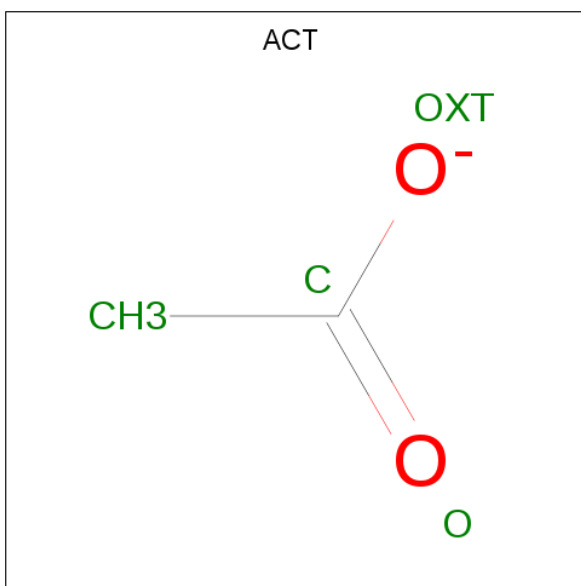
There are 3 unique types of molecules in this entry. The entry contains 7467 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Adenylation domain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	477	Total 3691	C 2335	N 668	O 679	S 9	0	1	0
1	B	477	Total 3674	C 2327	N 665	O 673	S 9	0	1	0

- Molecule 2 is ACETATE ION (three-letter code: ACT) (formula: C<sub>2</sub>H<sub>3</sub>O<sub>2</sub>).



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

- Molecule 3 is water.

<b>Mol</b>	<b>Chain</b>	<b>Residues</b>	<b>Atoms</b>		<b>ZeroOcc</b>	<b>AltConf</b>
3	A	44	Total 44	O 44	0	0
3	B	46	Total 46	O 46	0	0



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	177.63Å 52.81Å 108.86Å 90.00° 105.81° 90.00°	Depositor
Resolution (Å)	46.95 – 2.46 46.95 – 2.46	Depositor EDS
% Data completeness (in resolution range)	99.7 (46.95-2.46) 99.7 (46.95-2.46)	Depositor EDS
$R_{merge}$	0.14	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.64 (at 2.45Å)	Xtriage
Refinement program	PHENIX dev_1647	Depositor
R, $R_{free}$	0.197 , 0.270 0.200 , 0.270	Depositor DCC
$R_{free}$ test set	1793 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	37.9	Xtriage
Anisotropy	0.580	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 47.4	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	0.94	EDS
Total number of atoms	7467	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 13.22% 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 [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section:  
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	0.51	1/3778 (0.0%)	0.62	3/5141 (0.1%)
1	B	0.47	0/3761	0.61	0/5117
All	All	0.49	1/7539 (0.0%)	0.62	3/10258 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	123	PRO	N-CD	8.67	1.59	1.47

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	123	PRO	CA-N-CD	-6.81	101.97	111.50
1	A	123	PRO	N-CD-CG	-5.60	94.80	103.20
1	A	123	PRO	N-CA-C	5.25	125.76	112.10

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	3691	0	3659	52	1
1	B	3674	0	3650	46	1

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	4	0	3	0	0
2	B	8	0	6	0	0
3	A	44	0	0	2	0
3	B	46	0	0	1	0
All	All	7467	0	7318	98	1

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 (98) 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:B:155:ARG:NH2	3:B:701:HOH:O	2.13	0.80
1:A:42:GLU:OE2	1:A:143:ARG:NH2	2.14	0.78
1:B:294:ARG:NH1	1:B:330:LEU:O	2.20	0.74
1:B:174:PRO:HB3	1:B:512:ILE:HG12	1.75	0.68
1:A:244:ARG:NH1	1:A:271:GLU:OE1	2.27	0.68
1:A:247:ARG:HG3	1:A:248:ASP:N	2.09	0.67
1:B:19:THR:HG22	1:B:21:VAL:H	1.61	0.65
1:B:75:MET:HE1	1:B:80:ASN:HB2	1.80	0.63
1:B:257:THR:HG22	1:B:259:THR:H	1.64	0.62
1:A:364:LEU:O	1:A:404:SER:OG	2.16	0.62
1:A:202:ARG:HG3	1:A:227:GLU:HB3	1.83	0.61
1:B:204:LEU:HD11	1:B:231:LEU:HG	1.82	0.61
1:A:61:HIS:O	1:A:66:ARG:NH2	2.34	0.60
1:A:380:GLU:OE1	1:A:381:LEU:N	2.25	0.59
1:A:102:ASN:OD1	1:A:106:ARG:NH1	2.35	0.59
1:A:442:VAL:HG13	1:A:485:LEU:HD11	1.86	0.58
1:B:427:LYS:HD2	1:B:432:ARG:NH2	2.19	0.57
1:A:513:ASN:N	1:A:513:ASN:OD1	2.27	0.56
1:A:75:MET:CE	1:A:120:VAL:HG12	2.35	0.56
1:A:304:ARG:NH1	3:A:702:HOH:O	2.34	0.55
1:A:490:LYS:HA	1:A:498:GLN:NE2	2.20	0.55
1:A:438:VAL:HB	1:A:453:VAL:HG21	1.89	0.55
1:A:439:GLU:OE2	1:A:518:ARG:NH2	2.40	0.55
1:A:357:ALA:N	1:A:360:GLU:OE1	2.34	0.54
1:B:289:THR:O	1:B:292:VAL:HG12	2.06	0.54
1:A:423:ASP:N	1:A:423:ASP:OD1	2.41	0.54
1:A:114:SER:HA	1:A:376:LEU:HD21	1.89	0.53
1:B:287:LYS:HG3	1:B:418:TYR:HD2	1.74	0.53
1:A:439:GLU:HG3	1:A:453:VAL:HG13	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:38:HIS:CG	1:B:43:LYS:HD2	2.44	0.52
1:A:101:ARG:HH22	1:A:512:ILE:HD12	1.73	0.52
1:A:435:LEU:HD22	1:A:453:VAL:HG23	1.92	0.52
1:B:117:VAL:O	1:B:132:PRO:HG2	2.10	0.52
1:A:456:HIS:CE1	1:A:526:ASP:OD1	2.63	0.51
1:B:364:LEU:O	1:B:404:SER:HB3	2.11	0.51
1:B:19:THR:HG21	1:B:156:PRO:HG2	1.93	0.51
1:B:75:MET:HE1	1:B:80:ASN:CB	2.40	0.51
1:B:259:THR:OG1	1:B:286:GLU:OE2	2.22	0.49
1:A:503:LEU:HD11	1:A:528:ARG:HD2	1.95	0.49
1:B:247:ARG:HD2	1:B:248:ASP:OD1	2.12	0.49
1:A:173:ARG:HH11	1:A:173:ARG:HG2	1.78	0.49
1:A:221:ALA:HA	1:A:226:ALA:HB3	1.94	0.49
1:B:427:LYS:HD2	1:B:432:ARG:HH21	1.77	0.49
1:A:122:ASP:OD2	1:A:122:ASP:C	2.50	0.48
1:A:167:THR:HG22	1:A:513:ASN:HD21	1.78	0.48
1:B:257:THR:HG22	1:B:259:THR:N	2.26	0.48
1:B:459:ALA:HB3	1:B:462:ASP:HB3	1.96	0.47
1:A:26:ARG:HB3	1:A:26:ARG:HE	1.53	0.47
1:A:440:ALA:O	1:A:444:THR:HG23	2.15	0.47
1:A:358:THR:HG23	1:A:410:ARG:O	2.14	0.47
1:A:399:PRO:HB2	1:A:401:TYR:CE2	2.50	0.47
1:A:426:VAL:HG11	1:A:455:VAL:HG21	1.97	0.47
1:B:404:SER:OG	1:B:406:ASP:HB2	2.15	0.47
1:B:386:PHE:HA	1:B:401:TYR:O	2.14	0.46
1:A:339:ARG:HD3	1:A:340:PRO:HD2	1.98	0.46
1:A:373:GLU:O	1:A:400:ARG:NH2	2.48	0.46
1:A:352:ASP:OD1	1:A:354:ARG:NH1	2.49	0.46
1:B:231:LEU:HD21	1:B:245:VAL:HG11	1.98	0.46
1:B:372:SER:OG	1:B:373:GLU:N	2.49	0.45
1:B:88:ILE:HG21	1:B:95:TYR:HB3	1.97	0.45
1:A:482:ALA:O	1:A:486:ARG:HB2	2.16	0.45
1:B:19:THR:HA	1:B:181:HIS:HB3	1.98	0.45
1:A:278:LEU:HD23	1:A:305:PRO:HG3	1.99	0.45
1:A:287:LYS:HB3	1:A:287:LYS:HE3	1.62	0.44
1:A:309:ASN:HB2	1:A:325:LEU:HD11	1.99	0.44
1:B:241:GLU:OE2	1:B:244:ARG:NE	2.49	0.44
1:B:119:LEU:HD13	1:B:133:LEU:HD13	2.00	0.44
1:B:111:VAL:HG23	1:B:116:VAL:HG22	1.98	0.44
1:B:201:ASP:O	1:B:226:ALA:HB1	2.18	0.43
1:B:75:MET:HE3	1:B:75:MET:HB3	1.71	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:380:GLU:O	1:A:383:ALA:HB3	2.18	0.43
1:A:173:ARG:NH1	1:A:173:ARG:HG2	2.33	0.43
1:B:406:ASP:CG	1:B:421:ARG:HE	2.22	0.43
1:B:26:ARG:HB3	1:B:26:ARG:HE	1.61	0.43
1:B:347:ARG:NE	1:B:355:ASP:OD1	2.48	0.43
1:B:458:PHE:CE1	1:B:464:ARG:HG3	2.54	0.43
1:B:241:GLU:HA	1:B:244:ARG:HD3	2.01	0.43
1:B:298:LYS:HA	1:B:298:LYS:HD2	1.35	0.43
1:B:349:VAL:HA	1:B:354:ARG:O	2.19	0.42
1:A:378:ARG:HG3	1:A:381:LEU:HD12	2.01	0.42
1:A:302:LEU:O	1:A:324:ARG:NH1	2.53	0.42
1:B:373:GLU:C	1:B:400:ARG:HH22	2.23	0.42
1:A:99:ASP:OD2	1:A:101:ARG:HB2	2.20	0.42
1:B:56:LEU:HA	1:B:56:LEU:HD12	1.87	0.42
1:B:363:GLU:OE2	1:B:405:GLY:N	2.45	0.42
1:A:289:THR:HB	3:A:708:HOH:O	2.20	0.41
1:A:438:VAL:HG12	1:A:453:VAL:HG11	2.03	0.41
1:A:117:VAL:O	1:A:132:PRO:HG2	2.21	0.41
1:B:19:THR:HG23	1:B:181:HIS:ND1	2.36	0.41
1:B:518:ARG:HA	1:B:521:VAL:HG22	2.03	0.41
1:A:349:VAL:HG22	1:A:363:GLU:HB3	2.03	0.41
1:A:512:ILE:HA	1:A:512:ILE:HD12	1.90	0.40
1:B:526:ASP:HA	1:B:529:ARG:HD3	2.01	0.40
1:A:174:PRO:O	1:A:378:ARG:NH2	2.54	0.40
1:A:425:GLN:NE2	1:A:434:GLU:OE2	2.52	0.40
1:B:512:ILE:HG13	1:B:512:ILE:H	1.50	0.40
1:B:517:ASP:CG	1:B:520:SER:HB2	2.42	0.40
1:A:288:LEU:O	1:A:335:SER:OG	2.30	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:269:LEU:O	1:B:196:SER:OG[3_445]	2.18	0.02

## 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	464/529 (88%)	453 (98%)	11 (2%)	0	100	100
1	B	464/529 (88%)	450 (97%)	14 (3%)	0	100	100
All	All	928/1058 (88%)	903 (97%)	25 (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	Percentiles	
1	A	382/420 (91%)	334 (87%)	48 (13%)	4	3
1	B	379/420 (90%)	350 (92%)	29 (8%)	13	15
All	All	761/840 (91%)	684 (90%)	77 (10%)	7	7

All (77) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	23	LEU
1	A	26	ARG
1	A	42	GLU
1	A	56	LEU
1	A	59	LEU
1	A	70	ARG
1	A	74	ARG
1	A	133	LEU

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
1	A	137	GLU
1	A	142	LEU
1	A	145	ARG
1	A	160	ASP
1	A	173	ARG
1	A	175	LYS
1	A	194	LEU
1	A	202	ARG
1	A	213	PHE
1	A	244	ARG
1	A	247	ARG
1	A	286	GLU
1	A	299	ARG
1	A	303	ASP
1	A	307	LEU
1	A	313	ILE
1	A	324	ARG
1	A	335	SER
1	A	342	PRO
1	A	351	GLU
1	A	354	ARG
1	A	378	ARG
1	A	380	GLU
1	A	385	ARG
1	A	403	HIS
1	A	404	SER
1	A	409	SER
1	A	410	ARG
1	A	421	ARG
1	A	423	ASP
1	A	424	LEU
1	A	426	VAL
1	A	427	LYS
1	A	450	ASP
1	A	486	ARG
1	A	513	ASN
1	A	515[A]	LYS
1	A	515[B]	LYS
1	A	518	ARG
1	A	528	ARG
1	B	25	SER
1	B	26	ARG

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Mol	Chain	Res	Type
1	B	43	LYS
1	B	51	SER
1	B	56	LEU
1	B	85	MET
1	B	119	LEU
1	B	155	ARG
1	B	194	LEU
1	B	213	PHE
1	B	244	ARG
1	B	298	LYS
1	B	299	ARG
1	B	313	ILE
1	B	330	LEU
1	B	341	LEU
1	B	342	PRO
1	B	380	GLU
1	B	421	ARG
1	B	428	LEU
1	B	431	HIS
1	B	435	LEU
1	B	465	LEU
1	B	472	GLN
1	B	483	ARG
1	B	508	LEU
1	B	512	ILE
1	B	520	SER
1	B	529	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	403	HIS
1	A	456	HIS
1	B	38	HIS

### 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](#)

3 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	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
2	ACT	A	601	-	1,3,3	0.87	0	0,3,3	0.00	-
2	ACT	B	602	-	1,3,3	0.87	0	0,3,3	0.00	-
2	ACT	B	601	-	1,3,3	1.39	0	0,3,3	0.00	-

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<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	477/529 (90%)	-0.06	4 (0%) 86 86	11, 33, 58, 74	0
1	B	477/529 (90%)	-0.07	4 (0%) 86 86	16, 32, 60, 78	0
All	All	954/1058 (90%)	-0.06	8 (0%) 86 86	11, 33, 59, 78	0

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	389	GLY	4.4
1	B	460	PRO	3.4
1	B	430	GLY	2.8
1	B	412	ALA	2.4
1	A	129	TYR	2.3
1	A	124	GLY	2.1
1	A	426	VAL	2.0
1	A	424	LEU	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. 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	B-factors(Å <sup>2</sup> )	Q<0.9
2	ACT	B	602	4/4	0.83	0.29	47,50,51,52	0
2	ACT	A	601	4/4	0.94	0.18	48,51,51,53	0
2	ACT	B	601	4/4	0.95	0.14	29,30,32,37	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.