



# Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 03:27 AM GMT

PDB ID : 2R5O  
Title : Crystal structure of the C-terminal domain of wzt  
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Deposited on : 2007-09-04  
Resolution : 1.30 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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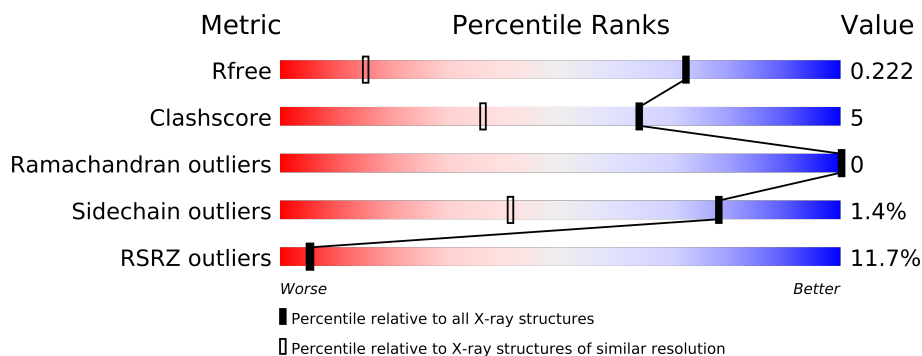
The following versions of software and data (see [references](#)) were used in the production of this report:

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 1.30 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	1025 (1.34-1.26)
Clashscore	79885	1140 (1.34-1.26)
Ramachandran outliers	78287	1093 (1.34-1.26)
Sidechain outliers	78261	1092 (1.34-1.26)
RSRZ outliers	66119	1025 (1.34-1.26)

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	188	
1	B	188	

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

Mol	Type	Chain	Res	Geometry	Electron density
4	PG4	A	432	-	X

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 3158 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 Putative ATP binding component of ABC-transporter.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	167	Total	C	N	O	S	0	13	0
			1390	897	217	270	6			
1	B	166	Total	C	N	O	S	0	15	0
			1383	894	218	265	6			

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	244	MET	-	EXPRESSION TAG	UNP Q47591
A	245	HIS	-	EXPRESSION TAG	UNP Q47591
A	246	HIS	-	EXPRESSION TAG	UNP Q47591
A	247	HIS	-	EXPRESSION TAG	UNP Q47591
A	248	HIS	-	EXPRESSION TAG	UNP Q47591
A	249	HIS	-	EXPRESSION TAG	UNP Q47591
A	250	HIS	-	EXPRESSION TAG	UNP Q47591
A	251	HIS	-	EXPRESSION TAG	UNP Q47591
A	252	HIS	-	EXPRESSION TAG	UNP Q47591
A	253	HIS	-	EXPRESSION TAG	UNP Q47591
A	254	HIS	-	EXPRESSION TAG	UNP Q47591
B	244	MET	-	EXPRESSION TAG	UNP Q47591
B	245	HIS	-	EXPRESSION TAG	UNP Q47591
B	246	HIS	-	EXPRESSION TAG	UNP Q47591
B	247	HIS	-	EXPRESSION TAG	UNP Q47591
B	248	HIS	-	EXPRESSION TAG	UNP Q47591
B	249	HIS	-	EXPRESSION TAG	UNP Q47591
B	250	HIS	-	EXPRESSION TAG	UNP Q47591
B	251	HIS	-	EXPRESSION TAG	UNP Q47591
B	252	HIS	-	EXPRESSION TAG	UNP Q47591
B	253	HIS	-	EXPRESSION TAG	UNP Q47591
B	254	HIS	-	EXPRESSION TAG	UNP Q47591

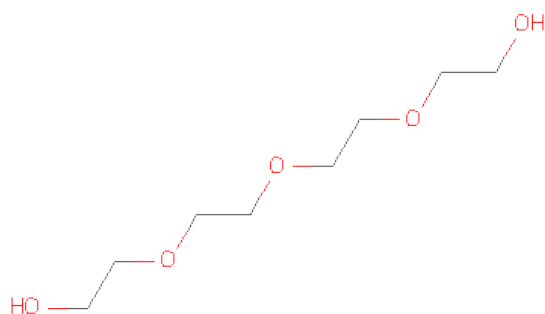
- Molecule 2 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	3	Total	Cl	0	0
			3	3		

- Molecule 3 is SODIUM ION (three-letter code: NA) (formula: Na).

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

- Molecule 4 is TETRAETHYLENE GLYCOL (three-letter code: PG4) (formula: C<sub>8</sub>H<sub>18</sub>O<sub>5</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			13	8	5		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	183	Total	O	0	0
			183	183		
5	B	181	Total	O	0	0
			181	181		



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 4 <sub>2</sub> 2 <sub>1</sub> 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	106.08Å 106.08Å 70.35Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 1.30 19.95 – 1.30	Depositor EDS
% Data completeness (in resolution range)	98.5 (20.00-1.30) 98.5 (19.95-1.30)	Depositor EDS
$R_{merge}$	0.05	Depositor
$R_{sym}$	0.06	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.53 (at 1.30Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, $R_{free}$	0.166 , 0.182 0.209 , 0.222	Depositor DCC
$R_{free}$ test set	4824 reflections (4.98%)	DCC
Wilson B-factor (Å <sup>2</sup> )	15.0	Xtriage
Anisotropy	0.006	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.42 , 46.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.35$	Xtriage
Outliers	0 of 96876 reflections	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	3158	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	18.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.08% 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: NA, PG4, CL

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	0.59	1/1474 (0.1%)	0.78	2/1998 (0.1%)
1	B	0.60	0/1478	0.75	1/2003 (0.0%)
All	All	0.60	1/2952 (0.0%)	0.76	3/4001 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	267	ASP	CG-OD1	5.24	1.37	1.25

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	326	ASP	CB-CG-OD1	5.35	123.12	118.30
1	A	402	ARG	NE-CZ-NH2	-5.17	117.72	120.30
1	A	343	PHE	CB-CG-CD1	-5.12	117.22	120.80

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	1390	0	1324	10	0
1	B	1383	0	1331	20	0
2	A	3	0	0	0	0
3	A	2	0	0	0	0
3	B	3	0	0	0	0
4	A	13	0	18	0	0
5	A	183	0	0	0	1
5	B	181	0	0	2	1
All	All	3158	0	2673	30	1

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 5.

All (30) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:335[B]:LEU:HD22	1:B:343[B]:PHE:CZ	1.63	1.33
1:B:424:LYS:NZ	5:B:583:HOH:O	1.71	1.18
1:B:335[B]:LEU:CD2	1:B:343[B]:PHE:HZ	1.67	1.06
1:B:335[B]:LEU:HD22	1:B:343[B]:PHE:HZ	0.79	0.95
1:B:312:ASN:HD21	1:B:415:MET:HG2	1.36	0.91
1:B:312:ASN:ND2	1:B:415:MET:HG2	1.90	0.86
1:B:329[B]:CYS:SG	1:B:358:ASN:ND2	2.55	0.80
1:B:265:SER:OG	1:B:267:ASP:OD1	2.07	0.73
1:A:393:PRO:HA	1:A:397[A]:MET:CG	2.22	0.69
1:B:335[B]:LEU:CD2	1:B:343[B]:PHE:CZ	2.56	0.66
1:A:413:LYS:HE2	1:A:417[B]:PHE:CE2	2.31	0.65
1:B:312:ASN:HD21	1:B:415:MET:CG	2.08	0.65
1:B:415:MET:HB3	5:B:529:HOH:O	1.98	0.64
1:B:312:ASN:ND2	1:B:415:MET:CG	2.63	0.62
1:A:393:PRO:HA	1:A:397[A]:MET:SD	2.41	0.61
1:A:393:PRO:HA	1:A:397[A]:MET:HG3	1.84	0.60
1:B:286:TRP:CZ2	1:B:402:ARG:HD2	2.41	0.56
1:B:329[B]:CYS:SG	1:B:358:ASN:CG	2.87	0.54
1:A:389:SER:HB3	1:A:397[B]:MET:HG2	1.94	0.49
1:B:334[A]:ILE:HG22	1:B:386:PHE:CD1	2.48	0.49
1:B:390:ALA:HB3	1:B:400:LEU:HD11	1.93	0.49
1:A:417[B]:PHE:CZ	1:A:422:ASP:HB2	2.51	0.46
1:A:397[A]:MET:HE3	1:A:397[A]:MET:HB3	1.90	0.43
1:B:337[B]:LYS:HE2	1:B:341:GLY:HA2	2.01	0.42
1:B:283[B]:GLU:OE1	1:B:403:ARG:HB3	2.20	0.41
1:B:294:ILE:CD1	1:B:320:LYS:HE3	2.51	0.40
1:B:337[A]:LYS:HE2	1:B:343[A]:PHE:CE1	2.56	0.40
1:A:393:PRO:CA	1:A:397[A]:MET:SD	3.08	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	Distance(Å)	Clash(Å)
5:A:473:HOH:O	5:B:482:HOH:O[6_555]	1.99	0.21

## 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	177/188 (94%)	169 (96%)	8 (4%)	0	100	100
1	B	179/188 (95%)	174 (97%)	5 (3%)	0	100	100
All	All	356/376 (95%)	343 (96%)	13 (4%)	0	100	100

There are no Ramachandran outliers to report.

### 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	161/167 (96%)	156 (97%)	5 (3%)	52	10
1	B	162/167 (97%)	161 (99%)	1 (1%)	92	72
All	All	323/334 (97%)	317 (98%)	6 (2%)	78	26

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	324[A]	GLU
1	A	324[B]	GLU
1	A	393	PRO
1	A	428[A]	THR

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Mol	Chain	Res	Type
1	A	428[B]	THR
1	B	363	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	312	ASN
1	B	358	ASN

### 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 9 ligands modelled in this entry, 8 are monoatomic - leaving 1 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	$\# Z  > 2$
4	PG4	A	432	-	12,12,12	0.54	0	11,11,11	0.55	0

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
4	PG4	A	432	-	-	0/10/10/10	0/0/0/0

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	167/188 (88%)	0.94	22 (13%) 4 4	13, 16, 23, 35	0
1	B	166/188 (88%)	0.74	18 (10%) 6 6	12, 15, 24, 29	0
All	All	333/376 (88%)	0.84	40 (12%) 5 5	12, 16, 24, 35	0

All (40) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	418	TRP	12.2
1	A	266	LEU	11.8
1	A	267	ASP	8.1
1	B	267	ASP	6.6
1	A	265	SER	5.9
1	B	343[A]	PHE	5.9
1	A	415	MET	5.4
1	B	354	GLU	5.0
1	B	266	LEU	4.9
1	B	265	SER	4.6
1	A	325	TYR	4.4
1	A	268	GLU	4.4
1	B	340	ASP	4.3
1	B	268	GLU	4.2
1	A	430[A]	TYR	4.1
1	B	355	GLY	3.9
1	A	269	ILE	3.9
1	A	416	ASP	3.9
1	A	311	GLY	3.8
1	B	404[A]	TYR	3.7
1	A	363	ARG	3.6
1	A	264	MET	3.5
1	B	269	ILE	3.3
1	B	339[A]	LEU	3.3

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Mol	Chain	Res	Type	RSRZ
1	A	340[A]	ASP	3.2
1	A	414	SER	3.0
1	B	353	SER	3.0
1	B	342	LEU	3.0
1	A	419	GLY	2.6
1	B	363	ARG	2.6
1	A	429	SER	2.6
1	A	394	GLN	2.6
1	A	417[A]	PHE	2.5
1	B	334[A]	ILE	2.5
1	A	293	ILE	2.4
1	B	264	MET	2.2
1	B	350	PHE	2.1
1	A	377	LEU	2.1
1	B	429	SER	2.0
1	A	407	ILE	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
4	PG4	A	432	13/13	0.15	2.42	38,39,42,43	0
2	CL	A	8	1/1	0.09	0.79	19,19,19,19	0
3	NA	B	12	1/1	0.11	-0.19	23,23,23,23	0
3	NA	B	10	1/1	0.08	-0.65	23,23,23,23	0
3	NA	A	11	1/1	0.06	-1.99	21,21,21,21	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
3	NA	B	9	1/1	0.04	-2.22	18,18,18,18	0
2	CL	A	1	1/1	0.04	-2.95	15,15,15,15	0
3	NA	A	13	1/1	0.02	-2.96	17,17,17,17	0
2	CL	A	2	1/1	0.03	-6.97	14,14,14,14	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.