

# Full wwPDB X-ray Structure Validation Report (i)

#### Aug 22, 2023 - 03:41 AM EDT

PDB ID	:	2Q47
Title	:	Ensemble refinement of the protein crystal structure of a putative phospho-
		protein phosphatase from Arabidopsis thaliana gene At1g05000
Authors	:	Levin, E.J.; Kondrashov, D.A.; Wesenberg, G.E.; Phillips Jr., G.N.; Center for
		Eukaryotic Structural Genomics (CESG)
Deposited on	:	2007-05-31
Resolution	:	3.30 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* A user guide is available at https://www.wwpdb.org/validation/2017/XrayValidationReportHelp with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at http://www.wwpdb.org/validation/2017/FAQs#types.

The following versions of software and data (see references (1)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
$\mathrm{EDS}$	:	2.35
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)	:	2.35

# 1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:  $X\text{-}RAY \, DIFFRACTION$ 

The reported resolution of this entry is 3.30 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.



Matria	Whole archive	Similar resolution
wietric	$(\# { m Entries})$	$(\# { m Entries},  { m resolution}  { m range}({ m \AA}))$
R <sub>free</sub>	130704	1149 (3.34-3.26)
Ramachandran outliers	138981	1183 (3.34-3.26)
Sidechain outliers	138945	1182 (3.34-3.26)

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 of the lower bar indicate the fraction of residues that contain outliers for >=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 <=5%

Mol	Chain	Length	Quality of chain	
1	1-A	151	91%	8% •
1	1-B	151	91%	9%
1	10-A	151	94%	6%
1	10-B	151	87%	13% •
1	11-A	151	94%	6%
1	11-B	151	92%	7% •
1	12-A	151	93%	7%
1	12-B	151	87%	12% •



Mol	Chain	Length	Quality of chain	
1	13-A	151	89%	11% •
1	13-B	151	79%	20% •
1	14-A	151	89%	11%
1	14-B	151	84%	15% •
1	15-A	151	81%	18% •
1	15-B	151	89%	11%
1	16-A	151	82%	17% ·
1	16-B	151	84%	16%
1	2-A	151	97%	•
1	2-B	151	93%	7%
1	3-A	151	95%	5%
1	3-B	151	86%	14%
1	4-A	151	91%	9%
1	4-B	151	94%	6%
1	5-A	151	95%	5%•
1	5-B	151	89%	11%
1	6-A	151	91%	9%
1	6-B	151	93%	7% •
1	7-A	151	92%	7% •
1	7-B	151	97%	•
1	8-A	151	97%	•
1	8-B	151	93%	7%
1	9-A	151	95%	5%
1	9-B	151	94%	6%



# 2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 40448 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.

Mol	Chain	Residues		Atoms					ZeroOcc	AltConf	Trace
1	1-A	151	Total 1224	C 789	N 216	0 211	${ m S}{ m 5}$	Se 3	0	0	0
1	2-A	151	Total 1224	C 789	N 216	0 211	$\frac{S}{5}$	${ m Se} \ 3$	0	0	0
1	3-A	151	Total 1224	C 789	N 216	0 211	$\frac{S}{5}$	${ m Se} \ 3$	0	0	0
1	4-A	151	Total 1224	C 789	N 216	0 211	${S \atop 5}$	${ m Se} \ 3$	0	0	0
1	5-A	151	Total 1224	C 789	N 216	0 211	${S \atop 5}$	${ m Se} \ 3$	0	0	0
1	6-A	151	Total 1224	C 789	N 216	0 211	${ m S}{ m 5}$	${ m Se} 3$	0	0	0
1	7-A	151	Total 1224	C 789	N 216	0 211	${S \atop 5}$	${ m Se} \ 3$	0	0	0
1	8-A	151	Total 1224	C 789	N 216	0 211	${ m S}{ m 5}$	${ m Se} \ 3$	0	0	0
1	9-A	151	Total 1224	C 789	N 216	0 211	$\frac{S}{5}$	${ m Se} \ 3$	0	0	0
1	10-A	151	Total 1224	C 789	N 216	0 211	${ m S}{ m 5}$	${ m Se} \ 3$	0	0	0
1	11-A	151	Total 1224	C 789	N 216	0 211	${S \atop 5}$	${ m Se} \ 3$	0	0	0
1	12-A	151	Total 1224	C 789	N 216	0 211	${ m S}{ m 5}$	${ m Se} 3$	0	0	0
1	13-A	151	Total 1224	C 789	N 216	0 211	${ m S}{ m 5}$	${ m Se} { m 3}$	0	0	0
1	14-A	151	Total 1224	C 789	N 216	O 211	${ m S}{ m 5}$	${ m Se} 3$	0	0	0
1	15-A	151	Total 1224	С 789	N 216	0 211	$\frac{S}{5}$	Se 3	0	0	0
1	16-A	151	Total 1224	C 789	N 216	0 211	$\frac{S}{5}$	Se 3	0	0	0

• Molecule 1 is a protein called Probable tyrosine-protein phosphatase At1g05000.



Mol	Chain	Residues	5		Atom	s			ZeroOcc	AltConf	Trace
1	1 D	1 5 1	Total	С	Ν	0	S	Se	0	0	0
	1-D	101	1224	789	216	211	5	3	0	0	0
1	٩D	151	Total	С	Ν	0	S	Se	0	0	0
	2-В	101	1224	789	216	211	5	3	0	0	0
1	лD	151	Total	С	Ν	0	S	Se	0	0	0
1	9-D	101	1224	789	216	211	5	3	0	0	0
1	4 D	151	Total	С	Ν	0	S	Se	0	0	0
	4-D	101	1224	789	216	211	5	3	0	0	0
1	5 P	151	Total	С	Ν	0	S	Se	0	0	0
1	0-D	101	1224	789	216	211	5	3	0	0	0
1	6 B	151	Total	С	Ν	0	S	Se	0	0	0
1	0-D	101	1224	789	216	211	5	3	0	0	0
1	7 B	151	Total	С	Ν	0	S	Se	0	Ο	0
1	1-D	101	1224	789	216	211	5	3	0	0	0
1	8-B	151	Total	С	Ν	Ο	$\mathbf{S}$	Se	0	0	Ο
	0-D	101	1224	789	216	211	5	3		0	0
1	9-B	151	Total	С	Ν	Ο	$\mathbf{S}$	Se	0	0	0
	5.0	101	1224	789	216	211	5	3	0	0	0
1	10-B	151	Total	С	Ν	Ο	$\mathbf{S}$	Se	0	0	0
	10 D	101	1224	789	216	211	5	3	0	0	0
1	11-B	151	Total	С	Ν	Ο	$\mathbf{S}$	Se	0	0	0
		101	1224	789	216	211	5	3			
1	12-B	151	Total	С	Ν	Ο	$\mathbf{S}$	Se	0	0	0
			1224	789	216	211	5	3	, , , , , , , , , , , , , , , , , , ,		<u> </u>
1	13-B	151	Total	С	Ν	0	S	Se	0	0	0
			1224	789	216	211	5	3			
1	14-B	151	Total	C	N	0	S	Se	0	0	0
			1224	789	216	211	5	3			
1	15-B	151	Total	C	N	0	S	Se	0	0	0
	_	-	1224	789	216	211	5	3	-	-	-
1	16-B	151	'Total	C	N	0	S	Se	0	0	0
			1224	789	216	211	5	3	0	Ŭ	Ŭ

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There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
А	61	MSE	MET	modified residue	UNP Q9ZVN4
А	133	MSE	MET	modified residue	UNP Q9ZVN4
А	195	MSE	MET	modified residue	UNP Q9ZVN4
В	61	MSE	MET	modified residue	UNP Q9ZVN4
В	133	MSE	MET	modified residue	UNP Q9ZVN4
В	195	MSE	MET	modified residue	UNP Q9ZVN4





Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	1-A	1	Total O S	0	0
			5 $4$ $1$	_	
2	2-A	1	Total O S	0	0
			5 $4$ $1$		
2	3-A	1	Total O S	0	0
			5 $4$ $1$		
2	4-A	1	Total O S	0	0
			5 $4$ $1$		
2	5-A	1	Total O S	0	0
			5 $4$ $1$		
2	6-A	1	Total O S	0	0
			5 $4$ $1$		
2	7-A	1	Total O S	0	0
			5 $4$ $1$		
2	8-A	1	Total O S	0	0
			5 $4$ $1$		
2	9-A	1	Total O S	0	0
			5 4 1		
2	10-A	1	Total O S	0	0
			5 4 1		
2	11-A	1	Total O S	0	0
			5 $4$ $1$	-	-
2	12-A	1	Total O S	0	0
		-	5 $4$ $1$		, in the second
2	13-A	1	Total O S	0	0
_	2 10-A		5 4 1	0	U



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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	14-A	1	Total O S	0	0
		-	5 $4$ $1$		0
2	15-A	1	Total O S	0	0
			$\begin{array}{ccc} 5 & 4 & 1 \\ \hline Total & O & S \end{array}$		
2	16-A	1	5 4 1	0	0
			Total O S	0	0
2	1-A	1	5 4 1	0	0
2	2 4	1	Total O S	0	0
	2-11	T	5 4 1	0	0
2	3-A	1	Total O S	0	0
			5 4 1		
2	4-A	1	$\begin{array}{ccc} 10tal & 0 & 5 \\ 5 & 4 & 1 \end{array}$	0	0
	- •		Total O S		
2	5-A	1	5 4 1	0	0
2	6 1	1	Total O S	0	0
	0-A	1	5 4 1	0	0
2	7-A	1	Total O S	0	0
		_	5 $4$ $1$		
2	8-A	1	Total O S	0	0
			Total O S		
2	9-A	1	5 4 1	0	0
0	10 4	1	Total O S	0	0
2	10-A	1	$5 \ 4 \ 1$	0	0
2	11-A	1	Total O S	0	0
		-	5 $4$ $1$		
2	12-A	1	Total O S	0	0
			Total O S		
2	13-A	1	5 4 1	0	0
0	14.4	1	Total O S	0	0
	14-A	1	5 4 1	0	0
2	15-A	1	Total O S	0	0
	10 11	1	5 $4$ $1$		
2	16-A	1	Total O S	0	0
			$\begin{array}{ccc} 0 & 4 & 1 \\ \hline Total & O & S \end{array}$		
2	1-B	1	5 4 1	0	0
	0.D		Total O S	0	
2	2-B		5 4 1	0	0



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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	3-B	1	Total O S $5 4 1$	0	0
2	4-B	1	$\begin{array}{ccc} 5 & 4 & 1 \\ \hline \text{Total} & \text{O} & \text{S} \\ 5 & 4 & 1 \end{array}$	0	0
2	5-B	1	$\begin{array}{ccc}  & & 1 \\  & & 1 \\  & & Total & O & S \\  & & 5 & 4 & 1 \end{array}$	0	0
2	6-B	1	$\begin{array}{ccc} \text{Total} & \text{O} & \text{S} \\ 5 & 4 & 1 \end{array}$	0	0
2	7-B	1	$\begin{array}{ccc} \text{Total} & \text{O} & \text{S} \\ 5 & 4 & 1 \end{array}$	0	0
2	8-B	1	$\begin{array}{ccc} \text{Total} & \text{O} & \text{S} \\ 5 & 4 & 1 \end{array}$	0	0
2	9-B	1	$\begin{array}{ccc} \text{Total} & \text{O} & \text{S} \\ 5 & 4 & 1 \end{array}$	0	0
2	10-B	1	$\begin{array}{ccc} \text{Total} & \text{O} & \text{S} \\ 5 & 4 & 1 \end{array}$	0	0
2	11-B	1	$\begin{array}{ccc} \text{Total} & \text{O} & \text{S} \\ 5 & 4 & 1 \end{array}$	0	0
2	12-B	1	$\begin{array}{ccc} \text{Total} & \text{O} & \text{S} \\ 5 & 4 & 1 \end{array}$	0	0
2	13-B	1	$\begin{array}{ccc} \text{Total} & \text{O} & \text{S} \\ 5 & 4 & 1 \end{array}$	0	0
2	14-B	1	$\begin{array}{ccc} \text{Total} & \text{O} & \text{S} \\ 5 & 4 & 1 \end{array}$	0	0
2	15-B	1	$\begin{array}{ccc} \text{Total} & \text{O} & \text{S} \\ 5 & 4 & 1 \end{array}$	0	0
2	16-B	1	$\begin{array}{ccc} \text{Total} & \text{O} & \text{S} \\ 5 & 4 & 1 \end{array}$	0	0
2	1-B	1	$\begin{array}{ccc} \text{Total} & \text{O} & \text{S} \\ 5 & 4 & 1 \end{array}$	0	0
2	2-B	1	$\begin{array}{ccc} \text{Total} & \text{O} & \text{S} \\ 5 & 4 & 1 \end{array}$	0	0
2	3-B	1	$\begin{array}{ccc} \text{Total} & \text{O} & \text{S} \\ 5 & 4 & 1 \end{array}$	0	0
2	4-B	1	$\begin{array}{ccc} \text{Total} & \text{O} & \text{S} \\ 5 & 4 & 1 \end{array}$	0	0
2	5-B	1	$\begin{array}{ccc} \text{Total} & \text{O} & \text{S} \\ 5 & 4 & 1 \end{array}$	0	0
2	6-B	1	$\begin{array}{c cc} Total & O & S \\ 5 & 4 & 1 \end{array}$	0	0
2	7-B	1	$\begin{array}{ccc} \text{Total} & \text{O} & \text{S} \\ 5 & 4 & 1 \end{array}$	0	0



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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	8-B	1	$\begin{array}{ccc} \text{Total} & \text{O} & \text{S} \\ 5 & 4 & 1 \end{array}$	0	0
2	9-B	1	$\begin{array}{ccc} \text{Total} & \text{O} & \text{S} \\ 5 & 4 & 1 \end{array}$	0	0
2	10-B	1	$\begin{array}{ccc} \text{Total} & \text{O} & \text{S} \\ 5 & 4 & 1 \end{array}$	0	0
2	11-B	1	$\begin{array}{ccc} \text{Total} & \text{O} & \text{S} \\ 5 & 4 & 1 \end{array}$	0	0
2	12-B	1	$\begin{array}{ccc} \text{Total} & \text{O} & \text{S} \\ 5 & 4 & 1 \end{array}$	0	0
2	13-B	1	$\begin{array}{ccc} \text{Total} & \text{O} & \text{S} \\ 5 & 4 & 1 \end{array}$	0	0
2	14-B	1	$\begin{array}{ccc} \text{Total} & \text{O} & \text{S} \\ 5 & 4 & 1 \end{array}$	0	0
2	15-B	1	$\begin{array}{ccc} \text{Total} & \text{O} & \text{S} \\ 5 & 4 & 1 \end{array}$	0	0
2	16-B	1	$\begin{array}{ccc} \text{Total} & \text{O} & \text{S} \\ 5 & 4 & 1 \end{array}$	0	0

• Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	1-A	31	Total O 31 31	0	0
3	2-A	31	Total         O           31         31	0	0
3	3-A	30	Total O 30 30	0	0
3	4-A	30	Total O 30 30	0	0
3	5-A	29	Total         O           29         29	0	0
3	6-A	28	Total         O           28         28	0	0
3	7-A	29	Total         O           29         29	0	0
3	8-A	29	Total         O           29         29	0	0
3	9-A	29	Total         O           29         29	0	0
3	10-A	28	Total         O           28         28	0	0



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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	11-A	30	Total O 30 30	0	0
3	12-A	31	Total O 31 31	0	0
3	13-A	29	Total O 29 29	0	0
3	14-A	31	Total O 31 31	0	0
3	15-A	31	Total         O           31         31	0	0
3	16-A	31	Total         O           31         31	0	0
3	1-B	29	TotalO2929	0	0
3	2-B	29	Total         O           29         29	0	0
3	3-B	30	Total         O           30         30	0	0
3	4-B	30	Total O 30 30	0	0
3	5-B	31	Total O 31 31	0	0
3	6-B	32	$\begin{array}{cc} \text{Total} & \text{O} \\ 32 & 32 \end{array}$	0	0
3	7-B	31	Total O 31 31	0	0
3	8-B	31	Total O 31 31	0	0
3	9-B	31	$\begin{array}{cc} \text{Total} & \text{O} \\ 31 & 31 \end{array}$	0	0
3	10-B	32	Total O 32 32	0	0
3	11-B	30	Total         O           30         30	0	0
3	12-B	29	Total O 29 29	0	0
3	13-B	31	Total O 31 31	0	0
3	14-B	29	TotalO2929	0	0
3	15-B	29	TotalO2929	0	0



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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	16-B	29	Total         O           29         29	0	0



# 3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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. 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. 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: Probable tyrosine-protein phosphatase At1g05000

Chain 1-A:	91%	8% •
H52 V62 D63 D63 C70 C70 C70 C70 C70 D73 D73 D73 D73 D73 D73 D73 D73 D73 D73	H149 8174 A183 A184 A184 A186 S202 S202	
• Molecule 1: Probable ty	rosine-protein phosphatase At1g0500	0
Chain 1-B:	91%	9%
HB2 F11 F11 Y96 N100 N100 N148 H149 H149 C150 C150 Q168 M169 M170	R181 A185 F198 S202	
• Molecule 1: Probable ty	rosine-protein phosphatase At1g0500	0
Chain 2-A:	97%	· ·
H5 2 569 57 4 77 4 196 196 196 8174 8174 8174		
• Molecule 1: Probable ty	rosine-protein phosphatase At1g0500	0
Chain 2-B:	93%	7%
H62 F77 F77 F96 H129 C160 C160 C160 C160 C166 C166 C166 C166	A 185 198 8 202	
• Molecule 1: Probable ty	rosine-protein phosphatase At1g0500	0
Chain 3-A:	95%	5%
HE2 N64 N64 F71 F71 F71 F71 F71 F71 F71 F71 F71 F71		
• Molecule 1: Probable ty	rosine-protein phosphatase At1g0500	0
Chain 3-B:	86%	14%

• Molecule 1: Probable tyrosine-protein phosphatase At1g05000





• Molecule 1: Probable tyrosine-protein phosphatase At1g05000

Chain 7-B:



• Molecule 1: Probable tyrosine-protein phosphatase At1g05000

Chain 8-A:	97%	•
H52 869 869 8117 1167 8202 8202		
• Molecule 1: Probable tyrosine-protein p	phosphatase At1g05000	

97%

Chain 8-B: 93% 7%

• Molecule 1: Probable tyrosine-protein phosphatase At1g05000

Chain 9-A: 95% 5%

• Molecule 1: Probable tyrosine-protein phosphatase At1g05000

Chain 9-B:

94%

6%



• Molecule 1: Probable tyrosine-protein phosphatase At1g05000

Chain 10-A: 94% 6%

• Molecule 1: Probable tyrosine-protein phosphatase At1g05000

Chain 10-B: 87% 13% •



# 

• Molecule 1: Probable tyrosine-protein phosphatase At1g05000

Chain 11-A: 94% 6%

 $\bullet$  Molecule 1: Probable tyrosine-protein phosphatase At1g05000

Chain 11-B:	92%	7% •
H52 N68 F71 F71 F71 F71 F117 N126 K130 V146 V146	H149 K165 F198 S202	
• Molecule 1: Probable t	yrosine-protein phosphatase $At1g05000$	
Chain 12-A:	93%	7%
H52 157 158 158 158 158 158 159 1104 1117 1117 1117 1117	8201 8201 8202	
• Molecule 1: Probable t	yrosine-protein phosphatase $At1g05000$	
Chain 12-B:	87%	12% •
H52 P56 F71 790 796 F103 F106 F103 F112	E117 1135 1135 1146 1147 1148 1148 1148 1148 1148 1197 1197 1197 1197	
• Molecule 1: Probable t	yrosine-protein phosphatase At1g05000	
Chain 13-A:	89%	11% •
H52 D63 S74 S74 K86 Y90 Y90 Y90 Y90 Y90 Y125 K142	M43 H149 C159 T173 S174 Q192 Q192 Q192	
• Molecule 1: Probable t	yrosine-protein phosphatase $At1g05000$	
Chain 13-B:	79%	20% •
H52 L57 L57 N64 R68 R68 R68 873 S74 S74 S74 S78 S74 C83 S78 C83 C83 C83 C83 C83 C83 C83 C83 C83 C8	R86 191 191 191 191 191 193 193 193 193 111 111	R165 R165 W170 W170 C171 A185 S201 S202
• Molecule 1: Probable t	yrosine-protein phosphatase $At1g05000$	
Chain 14-A:	89%	11%



#### H52 F59 G70 G70 G70 B12 F101 F101 F101 F101 F101 H149 H149 H149 H149 H149 G158 G158 G158 G158 G158 F198 F198 F198

• Molecule 1: Probable tyrosine-protein phosphatase At1g05000

Chain 14-B:	84%	15% •
H52 N64 D7 3 S74 F77 T82	Y96 F112 F112 1131 1131 1148 1148 1148 1148 1148 1148 1172 1172 1172 1168 1148 1148 1172 1172 1172 1185 1197 1172 1197 1172 1197 1	
• Molecule 1:	Probable tyrosine-protein phosphatase At1g05000	
Chain 15-A:	81%	18% •
H52 L57 N58 N64 D73 D73	E99 E99 E99 E99 E99 E117 E117 E117 E117	<b>5</b> 202
• Molecule 1:	Probable tyrosine-protein phosphatase At1g05000	
Chain 15-B:	89%	11%
H52 V62 G65 F77 Y96	1126 M133 V146 V146 C150 C150 C171 C171 C171 C171 C172 C174 C173 C174 C173 C174 C173 C174 C173 C174 C173 C176 C176 C176 C176 C176 C176 C176 C176	
• Molecule 1:	Probable tyrosine-protein phosphatase At1g05000	
Chain 16-A:	82%	17% •
H52 F71 S74 L80 L80 V00	190 190 190 190 100 100 110 113 113 113 113 113 113 11	
• Molecule 1:	Probable tyrosine-protein phosphatase At1g05000	
Chain 16-B:	84%	16%
H52 L53 P56 M61 V62 V62 D63 D63	3/4           3/4           3/4           878           878           878           878           878           878           878           878           878           878           878           878           878           878           933           933           933           933           933           933           933           933           933           933           933           933           933           934           934           934           934           934           934           934           934           934           935           935           936           936           937           936           937           936           937           936           937           936	



# 4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 3	Depositor
Cell constants	124.48Å 124.48Å 124.48Å	Deperitor
a, b, c, $\alpha$ , $\beta$ , $\gamma$	$90.00^{\circ}$ $90.00^{\circ}$ $90.00^{\circ}$	Depositor
$\mathbf{P}_{\text{assolution}}(\hat{\mathbf{A}})$	41.50 - 3.30	Depositor
Resolution (A)	44.01 - 3.30	EDS
% Data completeness	99.1 (41.50-3.30)	Depositor
(in resolution range)	99.1 (44.01-3.30)	EDS
R <sub>merge</sub>	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$< I/\sigma(I) > 1$	10.21 (at 3.32Å)	Xtriage
Refinement program	CNS 1.1	Depositor
D D.	0.160 , $0.234$	Depositor
$n, n_{free}$	0.204 , $0.253$	DCC
$R_{free}$ test set	1736 reflections $(9.21%)$	wwPDB-VP
Wilson B-factor $(Å^2)$	33.9	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent $k_{sol}(e/Å^3), B_{sol}(Å^2)$	0.32, 105.6	EDS
L-test for twinning <sup>2</sup>	$< L >=0.50, < L^2>=0.33$	Xtriage
Estimated twinning fraction	0.035 for l,-k,h	Xtriage
$F_o, F_c$ correlation	0.88	EDS
Total number of atoms	40448	wwPDB-VP
Average B, all atoms $(Å^2)$	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 3.45% of the height of the origin peak. No significant pseudotranslation is detected.

<sup>&</sup>lt;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.



<sup>&</sup>lt;sup>1</sup>Intensities estimated from amplitudes.

# 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: SO4

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).

Mal	Chain	Bond lengths		Bond angles	
	Ullaill	RMSZ	# Z  > 5	RMSZ	# Z  > 5
1	1-A	0.36	0/1252	0.58	0/1680
1	1-B	0.34	0/1252	0.55	0/1680
1	2-A	0.36	0/1252	0.57	0/1680
1	2-B	0.35	0/1252	0.56	0/1680
1	3-A	0.35	0/1252	0.56	0/1680
1	3-B	0.35	0/1252	0.57	0/1680
1	4-A	0.37	0/1252	0.58	0/1680
1	4-B	0.36	0/1252	0.56	0/1680
1	5-A	0.36	0/1252	0.57	0/1680
1	5-B	0.34	0/1252	0.55	0/1680
1	6-A	0.36	0/1252	0.58	0/1680
1	6-B	0.34	0/1252	0.56	0/1680
1	7-A	0.36	0/1252	0.56	0/1680
1	7-B	0.35	0/1252	0.54	0/1680
1	8-A	0.35	0/1252	0.56	0/1680
1	8-B	0.34	0/1252	0.55	0/1680
1	9-A	0.35	0/1252	0.57	0/1680
1	9-B	0.34	0/1252	0.55	0/1680
1	10-A	0.35	0/1252	0.55	0/1680
1	10-B	0.35	0/1252	0.55	0/1680
1	11-A	0.36	0/1252	0.57	0/1680
1	11-B	0.34	0/1252	0.55	0/1680
1	12-A	0.35	0/1252	0.57	0/1680
1	12-B	0.34	0/1252	0.57	0/1680
1	13-A	0.40	0/1252	0.62	0/1680
1	13-B	0.39	0/1252	0.62	0/1680
1	14-A	0.38	0/1252	0.65	0/1680
1	14-B	0.38	0/1252	0.62	0/1680
1	15-A	0.39	0/1252	0.66	0/1680
1	15-B	0.38	0/1252	0.63	0/1680
1	16-A	0.40	0/1252	0.65	0/1680
1	16-B	0.40	0/1252	0.61	0/1680



Mol	Chain	Bond	lengths	Bond angles		
		RMSZ	# Z  > 5	RMSZ	# Z  > 5	
All	All	0.36	0/40064	0.58	0/53760	

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 (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	1-A	1224	0	1233	0	0
1	1-B	1224	0	1233	0	0
1	2-A	1224	0	1233	0	0
1	2-B	1224	0	1233	0	0
1	3-A	1224	0	1233	0	0
1	3-B	1224	0	1233	0	0
1	4-A	1224	0	1233	0	0
1	4-B	1224	0	1233	0	0
1	5-A	1224	0	1233	0	0
1	5-B	1224	0	1233	0	0
1	6-A	1224	0	1233	0	0
1	6-B	1224	0	1233	0	0
1	7-A	1224	0	1233	0	0
1	7-B	1224	0	1233	0	0
1	8-A	1224	0	1233	0	0
1	8-B	1224	0	1233	0	0
1	9-A	1224	0	1233	0	0
1	9-B	1224	0	1233	0	0
1	10-A	1224	0	1233	0	0
1	10-B	1224	0	1233	0	0
1	11-A	1224	0	1233	0	0
1	11-B	1224	0	1233	0	0
1	12-A	1224	0	1233	0	0
1	12-B	1224	0	1233	0	0
1	13-A	1224	0	1233	0	0



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 Mol
 Chain
 Non-H
 H(model)
 H(added)
 Clashes
 Symm-Clashes

1	13-B	1224	0	1233	0	0
1	14-A	1224	0	1233	0	0
1	14-B	1224	0	1233	0	0
1	15-A	1224	0	1233	0	0
1	15-B	1224	0	1233	0	0
1	16-A	1224	0	1233	0	0
1	16-B	1224	0	1233	0	0
2	1-A	10	0	0	0	0
2	1-B	10	0	0	0	0
2	2-A	10	0	0	0	0
2	2-B	10	0	0	0	0
2	3-A	10	0	0	0	0
2	3-B	10	0	0	0	0
2	4-A	10	0	0	0	0
2	4-B	10	0	0	0	0
2	5-A	10	0	0	0	0
2	5-B	10	0	0	0	0
2	6-A	10	0	0	0	0
2	6-B	10	0	0	0	0
2	7-A	10	0	0	0	0
2	7-B	10	0	0	0	0
2	8-A	10	0	0	0	0
2	8-B	10	0	0	0	0
2	9-A	10	0	0	0	0
2	9-B	10	0	0	0	0
2	10-A	10	0	0	0	0
2	10-B	10	0	0	0	0
2	11-A	10	0	0	0	0
2	11-B	10	0	0	0	0
2	12-A	10	0	0	0	0
2	12-B	10	0	0	0	0
2	13-A	10	0	0	0	0
2	13-B	10	0	0	0	0
2	14-A	10	0	0	0	0
2	14-B	10	0	0	0	0
2	15-A	10	0	0	0	0
2	15-B	10	0	0	0	0
2	16-A	10	0	0	0	0
2	16-B	10	0	0	0	0
3	1-A	31	0	0	0	0
3	1-B	29	0	0	0	0
3	2-A	31	0	0	0	0



Continue contract c	Continued from previous page					
Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	2-B	29	0	0	0	0
3	3-A	30	0	0	0	0
3	3-B	30	0	0	0	0
3	4-A	30	0	0	0	0
3	4-B	30	0	0	0	0
3	5-A	29	0	0	0	0
3	5-B	31	0	0	0	0
3	6-A	28	0	0	0	0
3	6-B	32	0	0	0	0
3	7-A	29	0	0	0	0
3	7-B	31	0	0	0	0
3	8-A	29	0	0	0	0
3	8-B	31	0	0	0	0
3	9-A	29	0	0	0	0
3	9-B	31	0	0	0	0
3	10-A	28	0	0	0	0
3	10-B	32	0	0	0	0
3	11-A	30	0	0	0	0
3	11-B	30	0	0	0	0
3	12-A	31	0	0	0	0
3	12-B	29	0	0	0	0
3	13-A	29	0	0	0	0
3	13-B	31	0	0	0	0
3	14-A	31	0	0	0	0
3	14-B	29	0	0	0	0
3	15-A	31	0	0	0	0
3	15-B	29	0	0	0	0
3	16-A	31	0	0	0	0
3	16-B	29	0	0	0	0
All	All	40448	0	39456	0	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). Clashscore could not be calculated for this entry.

There are no clashes within the asymmetric unit.

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	Perce	entiles
1	1-A	149/151~(99%)	116 (78%)	26 (17%)	7 (5%)	2	14
1	1-B	149/151~(99%)	118 (79%)	24 (16%)	7 (5%)	2	14
1	2-A	149/151~(99%)	127 (85%)	21 (14%)	1 (1%)	22	54
1	2-B	149/151~(99%)	122 (82%)	22~(15%)	5 (3%)	3	22
1	3-A	149/151~(99%)	127 (85%)	17 (11%)	5 (3%)	3	22
1	3-B	149/151~(99%)	112 (75%)	29 (20%)	8 (5%)	2	12
1	4-A	149/151~(99%)	114 (76%)	30 (20%)	5 (3%)	3	22
1	4-B	149/151 (99%)	122 (82%)	22 (15%)	5 (3%)	3	22
1	5-A	149/151~(99%)	116 (78%)	28 (19%)	5 (3%)	3	22
1	5-B	149/151~(99%)	116 (78%)	20 (13%)	13 (9%)	1	5
1	6-A	149/151~(99%)	120 (80%)	20 (13%)	9 (6%)	1	10
1	6-B	149/151~(99%)	122 (82%)	23 (15%)	4 (3%)	5	26
1	7-A	149/151~(99%)	130 (87%)	13 (9%)	6 (4%)	3	18
1	7-B	149/151 (99%)	119 (80%)	28 (19%)	2 (1%)	12	40
1	8-A	149/151~(99%)	126 (85%)	21 (14%)	2 (1%)	12	40
1	8-B	149/151 (99%)	119 (80%)	27 (18%)	3 (2%)	7	32
1	9-A	149/151~(99%)	122 (82%)	22 (15%)	5 (3%)	3	22
1	9-B	149/151 (99%)	121 (81%)	24 (16%)	4 (3%)	5	26
1	10-A	149/151~(99%)	118 (79%)	27 (18%)	4 (3%)	5	26
1	10-B	149/151~(99%)	109 (73%)	29 (20%)	11 (7%)	1	7
1	11-A	149/151~(99%)	119 (80%)	26 (17%)	4 (3%)	5	26
1	11-B	149/151~(99%)	113 (76%)	32 (22%)	4 (3%)	5	26
1	12-A	149/151 (99%)	111 (74%)	31 (21%)	7 (5%)	2	14
1	12-B	149/151~(99%)	104 (70%)	32 (22%)	13 (9%)	1	5
1	13-A	149/151~(99%)	115 (77%)	26 (17%)	8 (5%)	2	12



Mol	Chain	Analysed	Favoured	Allowed	Outliers	Perc	entiles
1	13-B	149/151~(99%)	100 (67%)	32 (22%)	17 (11%)	0	2
1	14-A	149/151~(99%)	109 (73%)	30 (20%)	10 (7%)	1	8
1	14-B	149/151~(99%)	103 (69%)	34 (23%)	12 (8%)	1	6
1	15-A	149/151~(99%)	108 (72%)	26 (17%)	15 (10%)	0	3
1	15-B	149/151~(99%)	115 (77%)	26~(17%)	8 (5%)	2	12
1	16-A	149/151~(99%)	106 (71%)	26 (17%)	17 (11%)	0	2
1	16-B	149/151~(99%)	105 (70%)	36 (24%)	8 (5%)	2	12
All	All	4768/4832 (99%)	3704 (78%)	830 (17%)	234 (5%)	2	14

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All (234) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	1-A	62	VAL
1	2-B	170	TRP
1	3-A	64	ASN
1	3-A	140	ASP
1	3-B	71	PHE
1	5-B	74	SER
1	5-B	151	LYS
1	5-B	183	ALA
1	6-A	123	PHE
1	6-A	140	ASP
1	7-A	74	SER
1	8-A	201	SER
1	10-B	87	SER
1	10-B	95	PRO
1	10-B	183	ALA
1	10-B	198	PHE
1	12-A	64	ASN
1	12-A	170	TRP
1	13-A	142	LYS
1	13-A	143	ASN
1	13-B	68	ARG
1	13-B	71	PHE
1	13-B	78	SER
1	13-B	83	LEU
1	13-B	85	LEU
1	13-B	111	LEU
1	14-B	197	ILE
1	15-A	57	LEU



Mol	Chain	Res	Type
1	15-A	64	ASN
1	15-A	98	GLU
1	15-A	106	SER
1	15-A	151	LYS
1	15-A	185	ALA
1	15-A	197	ILE
1	15-B	62	VAL
1	16-A	98	GLU
1	16-A	106	SER
1	16-A	151	LYS
1	16-A	170	TRP
1	16-A	185	ALA
1	16-A	188	ARG
1	16-A	197	ILE
1	16-A	201	SER
1	16-B	186	LYS
1	1-A	70	GLY
1	1-A	106	SER
1	1-A	125	ASN
1	1-A	185	ALA
1	1-B	147	LEU
1	1-B	181	ARG
1	1-B	185	ALA
1	2-B	185	ALA
1	2-B	198	PHE
1	3-A	69	SER
1	3-B	198	PHE
1	4-A	186	LYS
1	4-B	63	ASP
1	4-B	151	LYS
1	4-B	152	ARG
1	5-A	70	GLY
1	5-A	119	ASN
1	5-A	201	SER
1	5-B	120	LYS
1	5-B	185	ALA
1	5-B	198	PHE
1	6-A	64	ASN
1	6-A	120	LYS
1	6-A	201	SER
1	6-B	185	ALA
1	6-B	201	SER



Mol	Chain	Res	Type
1	8-B	198	PHE
1	9-A	64	ASN
1	9-A	201	SER
1	9-B	85	LEU
1	10-A	201	SER
1	10-B	62	VAL
1	10-B	168	GLN
1	10-B	185	ALA
1	11-A	70	GLY
1	11-A	125	ASN
1	11-B	198	PHE
1	12-A	57	LEU
1	12-A	58	ASN
1	12-A	125	ASN
1	12-A	201	SER
1	12-B	87	SER
1	12-B	103	PHE
1	12-B	198	PHE
1	13-A	63	ASP
1	13-A	74	SER
1	13-A	125	ASN
1	13-A	170	TRP
1	13 <b>-</b> B	73	ASP
1	13-B	74	SER
1	13-B	86	ARG
1	13-B	108	GLY
1	13-B	150	CYS
1	13-B	154	LYS
1	13-B	185	ALA
1	14-A	62	VAL
1	14-A	80	LEU
1	14-A	168	GLN
1	14-B	74	SER
1	14-B	77	PHE
1	14-B	172	LEU
1	14-B	185	ALA
1	14-B	188	ARG
1	14-B	198	PHE
1	14-B	199	ASP
1	15-A	58	ASN
1	15-A	103	PHE
1	15-A	125	ASN



Mol	Chain	Res	Type
1	15-A	186	LYS
1	15-B	77	PHE
1	15-B	168	GLN
1	15-B	197	ILE
1	15-B	198	PHE
1	16-A	74	SER
1	16-B	63	ASP
1	16-B	185	ALA
1	1-A	183	ALA
1	1-B	168	GLN
1	2-A	69	SER
1	2-B	77	PHE
1	2-B	168	GLN
1	3-A	106	SER
1	3-A	141	GLU
1	3-B	185	ALA
1	4-A	64	ASN
1	4-B	154	LYS
1	4-B	198	PHE
1	5-A	97	PRO
1	5-B	115	GLY
1	5-B	117	GLU
1	5-B	125	ASN
1	5-B	168	GLN
1	5-B	188	ARG
1	6-B	188	ARG
1	7-A	69	SER
1	7-A	97	PRO
1	7-A	186	LYS
1	8-A	69	SER
1	9-B	117	GLU
1	10-A	64	ASN
1	10-A	125	ASN
1	10-A	150	CYS
1	10-B	106	SER
1	10-B	140	ASP
1	11-A	64	ASN
1	11-A	186	LYS
1	11-B	130	LYS
1	11-B	186	LYS
1	12-B	117	GLU
1	12-B	146	VAL



Mol	Chain	Res	Type
1	12-B	147	LEU
1	12-B	185	ALA
1	13-B	57	LEU
1	13-B	130	LYS
1	14-A	74	SER
1	14-A	185	ALA
1	14-B	125	ASN
1	15-B	65	GLY
1	15-B	172	LEU
1	15-B	183	ALA
1	16-A	80	LEU
1	16-A	97	PRO
1	16-A	107	ASN
1	16-B	74	SER
1	1-A	64	ASN
1	1-B	170	TRP
1	1-B	198	PHE
1	3-B	82	THR
1	3-B	130	LYS
1	4-A	97	PRO
1	5-A	186	LYS
1	6-A	122	PRO
1	6-A	141	GLU
1	7-A	63	ASP
1	7-A	200	VAL
1	9-A	62	VAL
1	10-B	143	ASN
1	10-B	154	LYS
1	11-B	125	ASN
1	12-B	90	TYR
1	12-B	125	ASN
1	12-B	197	ILE
1	13-A	117	GLU
1	13-A	186	LYS
1	13-B	93	PRO
1	14-A	70	GLY
1	14-A	98	GLU
1	14-A	198	PHE
1	14-B	82	THR
1	14-B	168	GLN
1	15-A	156	ARG
1	15-A	198	PHE



Mol	Chain	Res	Type
1	16-A	108	GLY
1	16-A	133	MSE
1	16-A	134	ALA
1	16-B	56	PRO
1	16-B	111	LEU
1	3-B	63	ASP
1	3-B	187	ALA
1	4-A	173	THR
1	5-B	150	CYS
1	7-B	168	GLN
1	8-B	87	SER
1	9-A	125	ASN
1	9-B	86	ARG
1	9-B	191	ASP
1	12-A	169	LYS
1	12-B	106	SER
1	13-B	201	SER
1	14-B	131	ILE
1	15-A	97	PRO
1	15-A	158	GLY
1	4-A	65	GLY
1	6-B	173	THR
1	7-B	143	ASN
1	12-B	148	ILE
1	1-B	148	ILE
1	5-B	122	PRO
1	14-A	158	GLY
1	16-B	93	PRO
1	6-A	62	VAL
1	12-B	56	PRO
1	14-A	108	GLY
1	16-B	145	PRO
1	6-A	108	GLY
1	3-B	197	ILE
1	9-A	124	VAL
1	16-A	200	VAL
1	8-B	62	VAL
1	16-A	93	PRO



#### 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 side chain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers Percent		ntiles
1	1-A	136/133~(102%)	129~(95%)	7~(5%)	24	54
1	1-B	136/133~(102%)	129~(95%)	7~(5%)	24	54
1	2-A	136/133~(102%)	132 (97%)	4 (3%)	42	69
1	2-B	136/133~(102%)	130 (96%)	6 (4%)	28	59
1	3-A	136/133~(102%)	133 (98%)	3 (2%)	52	74
1	3-B	136/133~(102%)	123 (90%)	13 (10%)	8	29
1	4-A	136/133~(102%)	128 (94%)	8 (6%)	19	49
1	4-B	136/133~(102%)	132 (97%)	4 (3%)	42	69
1	5-A	136/133~(102%)	132 (97%)	4 (3%)	42	69
1	5-B	136/133~(102%)	133 (98%)	3 (2%)	52	74
1	6-A	136/133~(102%)	132 (97%)	4 (3%)	42	69
1	6-B	136/133~(102%)	128 (94%)	8 (6%)	19	49
1	7-A	136/133~(102%)	128 (94%)	8 (6%)	19	49
1	7-B	136/133~(102%)	134 (98%)	2(2%)	65	81
1	8-A	136/133~(102%)	133 (98%)	3 (2%)	52	74
1	8-B	136/133~(102%)	129 (95%)	7 (5%)	24	54
1	9-A	136/133~(102%)	133 (98%)	3 (2%)	52	74
1	9-B	136/133~(102%)	131 (96%)	5 (4%)	34	63
1	10-A	136/133~(102%)	131 (96%)	5 (4%)	34	63
1	10-B	136/133~(102%)	126 (93%)	10 (7%)	13	40
1	11-A	136/133~(102%)	131 (96%)	5 (4%)	34	63
1	11-B	136/133~(102%)	127 (93%)	9~(7%)	16	46
1	12-A	136/133~(102%)	132 (97%)	4 (3%)	42	69
1	12-B	136/133~(102%)	127 (93%)	9~(7%)	16	46
1	13-A	136/133~(102%)	126 (93%)	10 (7%)	13	40
1	13-B	136/133~(102%)	119 (88%)	17 (12%)	4	19



Mol	Chain	Analysed	Rotameric	Outliers	Pe	erce	entiles
1	14-A	136/133~(102%)	130~(96%)	6 (4%)		28	59
1	14-B	136/133~(102%)	123 (90%)	13 (10%)		8	29
1	15-A	136/133~(102%)	120 (88%)	16 (12%)		5	21
1	15-B	136/133~(102%)	128 (94%)	8 (6%)		19	49
1	16-A	136/133~(102%)	125~(92%)	11 (8%)		11	36
1	16-B	136/133~(102%)	120 (88%)	16 (12%)		5	21
All	All	4352/4256~(102%)	4114 (94%)	238 (6%)		21	52

Continued from previous page...

All (238) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	1-A	64	ASN
1	1-A	71	PHE
1	1-A	73	ASP
1	1-A	96	TYR
1	1-A	100	ASN
1	1-A	149	HIS
1	1-A	174	SER
1	1-B	71	PHE
1	1-B	96	TYR
1	1-B	100	ASN
1	1-B	146	VAL
1	1-B	149	HIS
1	1-B	150	CYS
1	1-B	174	SER
1	2-A	74	SER
1	2-A	96	TYR
1	2-A	146	VAL
1	2-A	174	SER
1	2-B	96	TYR
1	2-B	129	HIS
1	2-B	146	VAL
1	2-B	150	CYS
1	2-B	157	THR
1	2-B	174	SER
1	3-A	71	PHE
1	3-A	96	TYR
1	3-A	104	LEU
1	3-B	52	HIS
1	3-B	61	MSE



Mol	Chain	Res	Type
1	3-B	78	SER
1	3-B	90	TYR
1	3-B	96	TYR
1	3-B	117	GLU
1	3-B	119	ASN
1	3-B	149	HIS
1	3-B	150	CYS
1	3-B	165	ARG
1	3-B	173	THR
1	3-B	174	SER
1	3-B	188	ARG
1	4-A	53	LEU
1	4-A	73	ASP
1	4-A	96	TYR
1	4-A	146	VAL
1	4-A	149	HIS
1	4-A	171	CYS
1	4-A	174	SER
1	4-A	181	ARG
1	4-B	58	ASN
1	4-B	96	TYR
1	4-B	149	HIS
1	4-B	169	LYS
1	5-A	96	TYR
1	5-A	119	ASN
1	5-A	136	LYS
1	5-A	149	HIS
1	5-B	71	PHE
1	5-B	96	TYR
1	5-B	146	VAL
1	6-A	69	SER
1	6-A	90	TYR
1	6-A	146	VAL
1	6-A	149	HIS
1	6-B	71	PHE
1	6-B	96	TYR
1	6-B	130	LYS
1	6-B	149	HIS
	6-B	167	LEU
1	6-B	171	CYS
1	6-B	173	THR
1	6-B	174	SER



Mol	Chain	Res	Type
1	7-A	63	ASP
1	7-A	71	PHE
1	7-A	74	SER
1	7-A	96	TYR
1	7-A	117	GLU
1	7-A	149	HIS
1	7-A	167	LEU
1	7-A	174	SER
1	7-B	96	TYR
1	7-B	146	VAL
1	8-A	96	TYR
1	8-A	117	GLU
1	8-A	167	LEU
1	8-B	68	ARG
1	8-B	96	TYR
1	8-B	101	LEU
1	8-B	112	PHE
1	8-B	149	HIS
1	8-B	150	CYS
1	8-B	157	THR
1	9-A	96	TYR
1	9-A	146	VAL
1	9-A	149	HIS
1	9-B	96	TYR
1	9-B	146	VAL
1	9-B	150	CYS
1	9-B	157	THR
1	9-B	173	THR
1	10-A	90	TYR
1	10-A	96	TYR
1	10-A	117	GLU
1	10-A	152	ARG
1	10-A	169	LYS
1	10-B	58	ASN
1	10-B	63	ASP
1	10-B	71	PHE
1	10-B	96	TYR
1	10-B	100	ASN
1	10-B	141	GLU
1	10-B	143	ASN
1	10-B	146	VAL
1	10-B	149	HIS



Mol	Chain	Res	Type
1	10-B	191	ASP
1	11-A	71	PHE
1	11-A	96	TYR
1	11-A	146	VAL
1	11-A	151	LYS
1	11-A	174	SER
1	11-B	58	ASN
1	11-B	71	PHE
1	11-B	96	TYR
1	11-B	101	LEU
1	11-B	117	GLU
1	11 <b>-</b> B	130	LYS
1	11-B	146	VAL
1	11-B	149	HIS
1	11-B	165	ARG
1	12-A	73	ASP
1	12-A	96	TYR
1	12-A	104	LEU
1	12-A	117	GLU
1	12-B	52	HIS
1	12-B	71	PHE
1	12-B	96	TYR
1	12-B	103	PHE
1	12-B	112	PHE
1	12-B	135	LEU
1	12-B	146	VAL
1	12-B	149	HIS
1	12-B	150	CYS
1	13-A	85	LEU
1	13-A	86	ARG
1	13-A	90	TYR
1	13-A	96	TYR
1	13-A	117	GLU
1	13-A	149	HIS
1	13-A	159	CYS
1	13-A	173	THR
1	13-A	174	SER
1	13-A	192	GLN
1	13-B	52	HIS
1	13-B	64	ASN
1	13-B	68	ARG
1	13-B	85	LEU



Mol	Chain	Res	Type
1	13-B	90	TYR
1	13-B	91	LEU
1	13-B	92	CYS
1	13-B	96	TYR
1	13-B	102	GLN
1	13-B	110	ARG
1	13-B	117	GLU
1	13-B	120	LYS
1	13-B	146	VAL
1	13-B	157	THR
1	13-B	165	ARG
1	13-B	169	LYS
1	13-B	171	CYS
1	14-A	59	PHE
1	14-A	96	TYR
1	14-A	101	LEU
1	14-A	117	GLU
1	14-A	146	VAL
1	14-A	149	HIS
1	14-B	64	ASN
1	14-B	73	ASP
1	14-B	74	SER
1	14-B	96	TYR
1	14-B	112	PHE
1	14 <b>-</b> B	146	VAL
1	14 <b>-</b> B	147	LEU
1	14 <b>-</b> B	149	HIS
1	14 <b>-</b> B	150	CYS
1	14-B	157	THR
1	14-B	165	ARG
1	14-B	171	CYS
1	14-B	174	SER
1	15-A	73	ASP
1	15-A	96	TYR
1	15-A	100	ASN
1	15-A	104	LEU
1	15-A	117	GLU
1	15-A	131	ILE
1	15-A	149	HIS
1	15-A	151	LYS
1	15-A	152	ARG
1	15-A	156	ARG



Mol	Chain	Res	Type
1	15-A	157	THR
1	15-A	169	LYS
1	15-A	174	SER
1	15-A	176	PHE
1	15-A	180	GLN
1	15-A	192	GLN
1	15-B	96	TYR
1	15-B	126	ILE
1	15-B	133	MSE
1	15-B	146	VAL
1	15-B	150	CYS
1	15-B	156	ARG
1	15-B	171	CYS
1	15-B	174	SER
1	16-A	71	PHE
1	16-A	85	LEU
1	16-A	90	TYR
1	16-A	96	TYR
1	16-A	107	ASN
1	16-A	110	ARG
1	16-A	131	ILE
1	16-A	136	LYS
1	16-A	146	VAL
1	16-A	149	HIS
1	16-A	157	THR
1	16-B	53	LEU
1	16-B	61	MSE
1	16-B	73	ASP
1	16-B	76	ASN
1	16-B	78	SER
1	16-B	96	TYR
1	16-B	110	ARG
1	16-B	113	GLN
1	16-B	117	GLU
1	16-B	126	ILE
1	16-B	131	ILE
1	16-B	135	LEU
1	16-B	150	CYS
1	16-B	156	ARG
1	16-B	165	ARG
1	16-B	174	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (125)



such sidechains are listed below:

$\mathbf{Mol}$	Chain	Res	Type
1	1-A	64	ASN
1	1-A	81	GLN
1	1-A	129	HIS
1	1-A	143	ASN
1	1-B	100	ASN
1	1-B	129	HIS
1	1-B	192	GLN
1	2-A	102	GLN
1	2-A	192	GLN
1	2-B	76	ASN
1	2-B	143	ASN
1	2-B	192	GLN
1	3-A	100	ASN
1	3-A	129	HIS
1	3-A	168	GLN
1	3-A	192	GLN
1	3-B	52	HIS
1	3-B	64	ASN
1	3-B	107	ASN
1	3-B	143	ASN
1	4-A	76	ASN
1	4-A	102	GLN
1	4-A	129	HIS
1	4-A	143	ASN
1	4-A	144	HIS
1	4-A	192	GLN
1	4-B	168	GLN
1	4-B	192	GLN
1	5-A	119	ASN
1	5-A	129	HIS
1	5-B	113	GLN
1	5-B	168	GLN
1	5-B	180	GLN
1	5-B	192	GLN
1	6-A	64	ASN
1	6-A	144	HIS
1	6-A	149	HIS
1	6-B	113	GLN
1	6-B	180	GLN
1	6-B	192	GLN
1	7-A	129	HIS
1	7-A	143	ASN



Mol	Chain	Res	Type
1	7-A	192	GLN
1	7-B	81	GLN
1	7-B	113	GLN
1	8-A	102	GLN
1	8-A	129	HIS
1	8-B	64	ASN
1	8-B	81	GLN
1	8-B	107	ASN
1	8-B	168	GLN
1	8-B	180	GLN
1	8-B	192	GLN
1	9-A	64	ASN
1	9-A	76	ASN
1	9-A	81	GLN
1	9-A	107	ASN
1	9-A	129	HIS
1	9-A	143	ASN
1	9-B	143	ASN
1	9-B	192	GLN
1	10-A	102	GLN
1	10-A	113	GLN
1	10-A	129	HIS
1	10-A	143	ASN
1	10-A	144	HIS
1	10-A	155	HIS
1	10-B	100	ASN
1	11-A	129	HIS
1	11-A	143	ASN
1	11-A	168	GLN
1	11-A	192	GLN
1	11-B	129	HIS
1	11-B	192	GLN
1	12-A	58	ASN
1	12-A	107	ASN
1	12-A	143	ASN
1	12-A	192	GLN
1	12-B	52	HIS
1	12-B	107	ASN
1	12-B	192	GLN
1	13-A	64	ASN
1	13-A	76	ASN
1	13-A	102	GLN



Mol	Chain	$\frac{1}{\mathbf{n} \cdot \mathbf{Res} \cdot \mathbf{Tvpe}}$			
1	13-A	129	HIS		
1	13-A	143	ASN		
1	13-A	168	GLN		
1	13-A	192	GLN		
1	13-B	58	ASN		
1	13-B	64	ASN		
1	13-B	102	GLN		
1	13-B	113	GLN		
1	13-B	143	ASN		
1	13-B	192	GLN		
1	14-A	102	GLN		
1	14-A	107	ASN		
1	14-A	129	HIS		
1	14-A	168	GLN		
1	14-A	192	GLN		
1	14-B	64	ASN		
1	14-B	81	GLN		
1	14-B	107	ASN		
1	14-B	143	ASN		
1	14-B	168	GLN		
1	14-B	192	GLN		
1	15-A	76	ASN		
1	15-A	100	ASN		
1	15-A	107	ASN		
1	15-A	129	HIS		
1	15-A	155	HIS		
1	15-A	180	GLN		
1	15-A	192	GLN		
1	15-B	125	ASN		
1	15-B	129	HIS		
1	15-B	143	ASN		
1	15-B	144	HIS		
1	15-B	168	GLN		
1	15-B	180	GLN		
1	15-B	192	GLN		
1	16-A	81	GLN		
1	16-A	107	ASN		
1	16-A	192	GLN		
1	16-B	76	ASN		
1	16-B	113	GLN		
1	16-B	144	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 monosaccharides in this entry.

#### 5.6 Ligand geometry (i)

64 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).

Mal	Type	Chain	Dog	Link	B	ond leng	$\operatorname{gths}$	Bond angles		
	Type	Ullalli	nes		Counts	RMSZ	# Z >2	Counts	RMSZ	# Z  > 2
2	SO4	3-A	204	-	4,4,4	0.26	0	$6,\!6,\!6$	0.11	0
2	SO4	6-A	204	-	4,4,4	0.26	0	$6,\!6,\!6$	0.11	0
2	SO4	8-B	203	-	4,4,4	0.24	0	$6,\!6,\!6$	0.05	0
2	SO4	1-A	203	-	4,4,4	0.23	0	$6,\!6,\!6$	0.07	0
2	SO4	6-B	203	-	4,4,4	0.23	0	$6,\!6,\!6$	0.05	0
2	SO4	12-A	203	-	4,4,4	0.25	0	$6,\!6,\!6$	0.06	0
2	SO4	5-B	203	-	4,4,4	0.22	0	$6,\!6,\!6$	0.05	0
2	SO4	11-B	204	-	4,4,4	0.24	0	$6,\!6,\!6$	0.10	0
2	SO4	5-B	204	-	4,4,4	0.21	0	$6,\!6,\!6$	0.12	0
2	SO4	9-B	203	-	4,4,4	0.24	0	$6,\!6,\!6$	0.05	0
2	SO4	7-A	203	-	4,4,4	0.27	0	$6,\!6,\!6$	0.05	0
2	SO4	15-B	204	-	4,4,4	0.27	0	$6,\!6,\!6$	0.10	0
2	SO4	7-A	204	-	4,4,4	0.26	0	$6,\!6,\!6$	0.11	0
2	SO4	3-B	203	-	4,4,4	0.24	0	$6,\!6,\!6$	0.06	0
2	SO4	10-A	203	-	4,4,4	0.21	0	$6,\!6,\!6$	0.07	0
2	SO4	4-A	204	-	4,4,4	0.26	0	6,6,6	0.11	0
2	SO4	9-B	204	-	4,4,4	0.28	0	$6,\!6,\!6$	0.07	0
2	SO4	1-B	204	-	4,4,4	0.26	0	6,6,6	0.08	0



	т	<u> </u>	Ъ	τ· 1	В	ond leng	gths	В	ond ang	gles
NIOI	Type	Chain	Res	Link	Counts	RMSZ	# Z  > 2	Counts	RMSZ	$\int \# Z  > 2$
2	SO4	4-A	203	-	4,4,4	0.24	0	$6,\!6,\!6$	0.06	0
2	SO4	5-A	203	-	4,4,4	0.24	0	$6,\!6,\!6$	0.06	0
2	SO4	8-A	203	-	4,4,4	0.25	0	$6,\!6,\!6$	0.08	0
2	SO4	7-B	204	-	4,4,4	0.28	0	$6,\!6,\!6$	0.07	0
2	SO4	7-B	203	-	$4,\!4,\!4$	0.24	0	$6,\!6,\!6$	0.05	0
2	SO4	16-B	203	-	$4,\!4,\!4$	0.24	0	$6,\!6,\!6$	0.05	0
2	SO4	12-B	203	-	$4,\!4,\!4$	0.23	0	$6,\!6,\!6$	0.05	0
2	SO4	11-A	204	-	$4,\!4,\!4$	0.27	0	$6,\!6,\!6$	0.11	0
2	SO4	11-B	203	-	$4,\!4,\!4$	0.23	0	$6,\!6,\!6$	0.06	0
2	SO4	13-B	203	-	$4,\!4,\!4$	0.25	0	$6,\!6,\!6$	0.09	0
2	SO4	16-A	203	-	$4,\!4,\!4$	0.24	0	$6,\!6,\!6$	0.06	0
2	SO4	15-B	203	-	$4,\!4,\!4$	0.21	0	$6,\!6,\!6$	0.06	0
2	SO4	4-B	204	-	$4,\!4,\!4$	0.27	0	$6,\!6,\!6$	0.06	0
2	SO4	12-B	204	-	$4,\!4,\!4$	0.27	0	$6,\!6,\!6$	0.08	0
2	SO4	3-A	203	-	$4,\!4,\!4$	0.23	0	$6,\!6,\!6$	0.06	0
2	SO4	12-A	204	-	$4,\!4,\!4$	0.26	0	$6,\!6,\!6$	0.11	0
2	SO4	16-B	204	-	$4,\!4,\!4$	0.17	0	$6,\!6,\!6$	0.13	0
2	SO4	1-A	204	-	$4,\!4,\!4$	0.25	0	$6,\!6,\!6$	0.08	0
2	SO4	14-A	204	-	$4,\!4,\!4$	0.27	0	$6,\!6,\!6$	0.13	0
2	SO4	10-B	203	-	$4,\!4,\!4$	0.26	0	$6,\!6,\!6$	0.09	0
2	SO4	14-A	203	-	$4,\!4,\!4$	0.22	0	$6,\!6,\!6$	0.07	0
2	SO4	2-A	204	-	$4,\!4,\!4$	0.26	0	$6,\!6,\!6$	0.10	0
2	SO4	16-A	204	-	$4,\!4,\!4$	0.25	0	$6,\!6,\!6$	0.13	0
2	SO4	8-B	204	-	$4,\!4,\!4$	0.28	0	$6,\!6,\!6$	0.07	0
2	SO4	15-A	203	-	$4,\!4,\!4$	0.27	0	$6,\!6,\!6$	0.11	0
2	SO4	15-A	204	-	4,4,4	0.25	0	$6,\!6,\!6$	0.08	0
2	SO4	6-B	204	-	$4,\!4,\!4$	0.27	0	$6,\!6,\!6$	0.08	0
2	SO4	10-A	204	-	4,4,4	0.26	0	$6,\!6,\!6$	0.06	0
2	SO4	9-A	204	-	$4,\!4,\!4$	0.26	0	$6,\!6,\!6$	0.11	0
2	SO4	2-B	203	-	4,4,4	0.24	0	$6,\!6,\!6$	0.05	0
2	SO4	9-A	203	-	4,4,4	0.23	0	$6,\!6,\!6$	0.05	0
2	SO4	11-A	203	-	4,4,4	0.24	0	$6,\!6,\!6$	0.08	0
2	SO4	13-A	204	-	4,4,4	0.27	0	$6,\!6,\!6$	0.13	0
2	SO4	13-A	203	-	4,4,4	0.24	0	$6,\!6,\!6$	0.07	0
2	SO4	1-B	203	-	4,4,4	0.24	0	$6,\!6,\!6$	0.05	0
2	SO4	4-B	203	-	4,4,4	0.25	0	$6,\!6,\!6$	0.08	0
2	SO4	14-B	203	-	4,4,4	0.25	0	$6,\!6,\!6$	0.06	0
2	SO4	2-A	203	-	4,4,4	0.26	0	$6,\!6,\!6$	0.06	0
2	SO4	8-A	204	-	4,4,4	0.26	0	$6,\!6,\!6$	0.11	0
2	SO4	10-B	204	-	4,4,4	0.27	0	$6,\!6,\!6$	0.07	0
2	SO4	6-A	203	-	4,4,4	0.26	0	$6,\!6,\!6$	0.05	0
2	SO4	14-B	204	-	4,4,4	0.27	0	$6,\!6,\!6$	0.10	0
2	SO4	5-A	204	-	4,4,4	0.26	0	$6,\!6,\!6$	0.11	0



Mal	Turne	no Chain Bog Link		Bond lengths			Bond angles			
INIOI	туре	Unam	nes	LIIIK	Counts	RMSZ	# Z >2	Counts	RMSZ	# Z >2
2	SO4	2-B	204	-	4,4,4	0.26	0	$6,\!6,\!6$	0.09	0
2	SO4	13-B	204	-	4,4,4	0.27	0	6,6,6	0.08	0
2	SO4	3-B	204	-	4,4,4	0.28	0	$6,\!6,\!6$	0.07	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.

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)

Unable to reproduce the depositors R factor - this section is therefore empty.

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

Unable to reproduce the depositors R factor - this section is therefore empty.

#### 6.3 Carbohydrates (i)

Unable to reproduce the depositors R factor - this section is therefore empty.

#### 6.4 Ligands (i)

Unable to reproduce the depositors R factor - this section is therefore empty.

#### 6.5 Other polymers (i)

Unable to reproduce the depositors R factor - this section is therefore empty.

