



# Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 11:58 PM GMT

PDB ID : 3AX2  
Title : Crystal structure of rat TOM20-ALDH presequence complex: a disulfide-tethered complex with a non-optimized, long linker  
Authors : Saitoh, T.; Maita, Y.; Kohda, D.  
Deposited on : 2011-03-28  
Resolution : 1.90 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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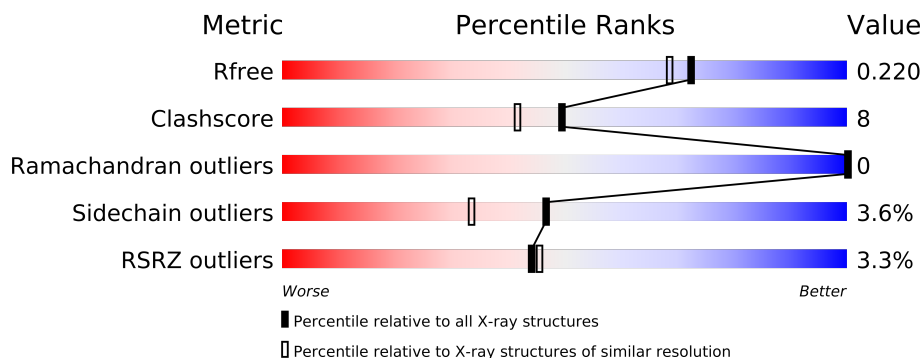
The following versions of software and data (see [references](#)) were used in the production of this report:

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 1.90 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	3684 (1.90-1.90)
Clashscore	79885	4465 (1.90-1.90)
Ramachandran outliers	78287	4413 (1.90-1.90)
Sidechain outliers	78261	4414 (1.90-1.90)
RSRZ outliers	66119	3686 (1.90-1.90)

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	73	
1	C	73	
1	E	73	
1	G	73	
2	B	16	
2	D	16	
2	F	16	
2	H	16	

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 2730 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 Mitochondrial import receptor subunit TOM20 homolog.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	70	Total	C	N	O	S	0	1	0
			551	354	88	107	2			
1	C	70	Total	C	N	O	S	0	1	0
			550	352	87	109	2			
1	E	65	Total	C	N	O	S	0	0	0
			510	328	81	99	2			
1	G	66	Total	C	N	O	S	0	0	0
			519	334	83	100	2			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	54	GLY	-	EXPRESSION TAG	UNP Q62760
A	55	PRO	-	EXPRESSION TAG	UNP Q62760
A	56	LEU	-	EXPRESSION TAG	UNP Q62760
A	57	GLY	-	EXPRESSION TAG	UNP Q62760
A	58	SER	-	EXPRESSION TAG	UNP Q62760
C	54	GLY	-	EXPRESSION TAG	UNP Q62760
C	55	PRO	-	EXPRESSION TAG	UNP Q62760
C	56	LEU	-	EXPRESSION TAG	UNP Q62760
C	57	GLY	-	EXPRESSION TAG	UNP Q62760
C	58	SER	-	EXPRESSION TAG	UNP Q62760
E	54	GLY	-	EXPRESSION TAG	UNP Q62760
E	55	PRO	-	EXPRESSION TAG	UNP Q62760
E	56	LEU	-	EXPRESSION TAG	UNP Q62760
E	57	GLY	-	EXPRESSION TAG	UNP Q62760
E	58	SER	-	EXPRESSION TAG	UNP Q62760
G	54	GLY	-	EXPRESSION TAG	UNP Q62760
G	55	PRO	-	EXPRESSION TAG	UNP Q62760
G	56	LEU	-	EXPRESSION TAG	UNP Q62760
G	57	GLY	-	EXPRESSION TAG	UNP Q62760
G	58	SER	-	EXPRESSION TAG	UNP Q62760

- Molecule 2 is a protein called Aldehyde dehydrogenase, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	16	Total	C	N	O	S	0	0	1
			107	65	22	19	1			
2	D	16	Total	C	N	O	S	0	2	1
			113	69	22	21	1			
2	F	16	Total	C	N	O	S	0	0	1
			107	65	22	19	1			
2	H	14	Total	C	N	O	S	0	0	1
			96	58	20	17	1			

There are 28 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	21	TYR	-	EXPRESSION TAG	UNP P11884
B	22	ALA	-	EXPRESSION TAG	UNP P11884
B	23	GLY	-	EXPRESSION TAG	UNP P11884
B	24	SER	-	EXPRESSION TAG	UNP P11884
B	25	GLY	-	EXPRESSION TAG	UNP P11884
B	26	CYS	-	EXPRESSION TAG	UNP P11884
B	27	NH2	-	AMIDATION	UNP P11884
D	21	TYR	-	EXPRESSION TAG	UNP P11884
D	22	ALA	-	EXPRESSION TAG	UNP P11884
D	23	GLY	-	EXPRESSION TAG	UNP P11884
D	24	SER	-	EXPRESSION TAG	UNP P11884
D	25	GLY	-	EXPRESSION TAG	UNP P11884
D	26	CYS	-	EXPRESSION TAG	UNP P11884
D	27	NH2	-	AMIDATION	UNP P11884
F	21	TYR	-	EXPRESSION TAG	UNP P11884
F	22	ALA	-	EXPRESSION TAG	UNP P11884
F	23	GLY	-	EXPRESSION TAG	UNP P11884
F	24	SER	-	EXPRESSION TAG	UNP P11884
F	25	GLY	-	EXPRESSION TAG	UNP P11884
F	26	CYS	-	EXPRESSION TAG	UNP P11884
F	27	NH2	-	AMIDATION	UNP P11884
H	21	TYR	-	EXPRESSION TAG	UNP P11884
H	22	ALA	-	EXPRESSION TAG	UNP P11884
H	23	GLY	-	EXPRESSION TAG	UNP P11884
H	24	SER	-	EXPRESSION TAG	UNP P11884
H	25	GLY	-	EXPRESSION TAG	UNP P11884
H	26	CYS	-	EXPRESSION TAG	UNP P11884
H	27	NH2	-	AMIDATION	UNP P11884

- Molecule 3 is PHOSPHATE ION (three-letter code: PO4) (formula: O<sub>4</sub>P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	O	P	0	0
			5	4	1		
3	D	1	Total	O	P	0	0
			5	4	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	37	Total	O	0	0
			37	37		
4	B	10	Total	O	0	0
			10	10		
4	C	49	Total	O	0	0
			49	49		
4	D	9	Total	O	0	0
			9	9		
4	E	30	Total	O	0	0
			30	30		
4	F	3	Total	O	0	0
			3	3		
4	G	21	Total	O	0	0
			21	21		
4	H	8	Total	O	0	0
			8	8		

### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: Mitochondrial import receptor subunit TOM20 homolog

Chain A: 



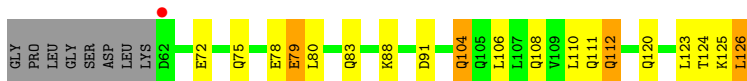
- Molecule 1: Mitochondrial import receptor subunit TOM20 homolog

Chain C: 



- Molecule 1: Mitochondrial import receptor subunit TOM20 homolog

Chain E: 



- Molecule 1: Mitochondrial import receptor subunit TOM20 homolog

Chain G: 



- Molecule 2: Aldehyde dehydrogenase, mitochondrial

Chain B: 



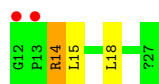
- Molecule 2: Aldehyde dehydrogenase, mitochondrial

Chain D: 



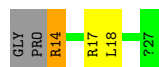
- Molecule 2: Aldehyde dehydrogenase, mitochondrial

Chain F: 



- Molecule 2: Aldehyde dehydrogenase, mitochondrial

Chain H: 



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 31	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	61.49Å 61.49Å 98.34Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	27.00 – 1.90 26.07 – 1.90	Depositor EDS
% Data completeness (in resolution range)	94.9 (27.00-1.90) 94.9 (26.07-1.90)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.32 (at 1.91Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, $R_{free}$	0.188 , 0.218 0.188 , 0.220	Depositor DCC
$R_{free}$ test set	1573 reflections (5.34%)	DCC
Wilson B-factor (Å <sup>2</sup> )	22.4	Xtriage
Anisotropy	0.062	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.38 , 37.7	EDS
Estimated twinning fraction	0.011 for -h,-k,l 0.032 for h,-h-k,-l 0.018 for -k,-h,-l	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.35$	Xtriage
Outliers	0 of 31049 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	2730	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.36% 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: PO4, NH2

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	1.34	4/562 (0.7%)	1.09	1/759 (0.1%)
1	C	1.34	2/561 (0.4%)	1.00	0/759
1	E	1.32	4/518 (0.8%)	1.07	2/702 (0.3%)
1	G	1.25	1/527 (0.2%)	0.99	0/713
2	B	1.40	0/107	1.23	1/142 (0.7%)
2	D	1.18	0/119	1.04	0/158
2	F	1.35	0/107	1.26	1/142 (0.7%)
2	H	1.35	0/95	1.17	0/125
All	All	1.32	11/2596 (0.4%)	1.06	5/3500 (0.1%)

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	78	GLU	CB-CG	-8.02	1.36	1.52
1	C	86	TYR	CD1-CE1	7.56	1.50	1.39
1	E	78	GLU	CB-CG	-6.71	1.39	1.52
1	E	104	GLN	CG-CD	5.82	1.64	1.51
1	A	64	GLU	CB-CG	5.67	1.62	1.52
1	E	79	GLU	CG-CD	5.66	1.60	1.51
1	A	70	PHE	CE2-CZ	5.51	1.47	1.37
1	E	79	GLU	CB-CG	5.46	1.62	1.52
1	C	72	GLU	CD-OE1	-5.19	1.20	1.25
1	A	73	GLU	CB-CG	5.15	1.61	1.52
1	G	82	ALA	CA-CB	5.02	1.62	1.52

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	14	ARG	NE-CZ-NH2	-7.54	116.53	120.30
1	A	110	LEU	CB-CG-CD2	-6.77	99.49	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	14	ARG	NE-CZ-NH1	-5.86	117.37	120.30
1	E	91	ASP	CB-CG-OD1	-5.31	113.52	118.30
1	E	78	GLU	OE1-CD-OE2	5.07	129.38	123.30

There are no chirality outliers.

There are no planarity outliers.

## 5.2 Close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	551	0	562	11	1
1	C	550	0	553	3	0
1	E	510	0	513	12	0
1	G	519	0	526	10	2
2	B	107	0	107	3	0
2	D	113	0	117	1	1
2	F	107	0	107	4	0
2	H	96	0	97	5	0
3	B	5	0	0	1	0
3	D	5	0	0	0	0
4	A	37	0	0	2	0
4	B	10	0	0	0	0
4	C	49	0	0	1	0
4	D	9	0	0	0	0
4	E	30	0	0	4	0
4	F	3	0	0	0	0
4	G	21	0	0	2	0
4	H	8	0	0	1	0
All	All	2730	0	2582	39	2

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 8.

All (39) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:79:GLU:OE2	2:F:14:ARG:NH1	1.95	0.99
1:C:67:GLN:HE22	2:D:18:LEU:HD23	1.33	0.92
1:E:75:GLN:OE1	4:E:157:HOH:O	2.01	0.78
1:G:61:LYS:HB2	4:G:135:HOH:O	1.83	0.77
1:A:74:ILE:HD12	2:F:15:LEU:HD23	1.70	0.72
4:G:137:HOH:O	2:H:17:ARG:HD2	1.93	0.68
1:E:111:GLN:HE21	1:E:112:GLN:HE22	1.40	0.68
1:E:126:LEU:HD13	4:E:130:HOH:O	1.93	0.67
2:B:12:GLY:N	3:B:2:PO4:O2	2.28	0.66
1:G:71:LEU:HD12	2:H:17:ARG:HG3	1.77	0.66
1:E:79:GLU:HG3	1:E:83:GLN:NE2	2.11	0.65
1:G:116:PRO:O	1:G:120:GLN:HG2	1.97	0.65
1:C:102:GLN:NE2	4:C:156:HOH:O	2.30	0.64
1:E:125:LYS:O	1:E:126:LEU:HB2	2.03	0.58
1:A:57:GLY:N	1:A:60:LEU:CD2	2.68	0.56
2:H:14:ARG:HB3	4:H:91:HOH:O	2.06	0.55
1:A:74:ILE:HD12	2:F:15:LEU:CD2	2.35	0.55
1:G:125:LYS:O	1:G:126:LEU:HB2	2.08	0.53
1:G:62:ASP:N	1:G:62:ASP:OD1	2.41	0.53
1:A:76:LEU:HD13	4:A:52:HOH:O	2.09	0.53
1:G:114:LEU:O	1:G:115:PRO:C	2.47	0.52
1:G:71:LEU:CD1	2:H:17:ARG:HG3	2.39	0.52
1:G:119:PHE:CE2	1:G:123:LEU:HD21	2.45	0.52
1:A:57:GLY:N	1:A:60:LEU:HD22	2.25	0.52
1:G:61:LYS:HG3	1:G:67:GLN:OE1	2.11	0.51
1:E:72:GLU:CD	4:E:156:HOH:O	2.49	0.49
1:E:126:LEU:HB2	4:E:130:HOH:O	2.12	0.49
1:E:80:LEU:HD13	1:E:88:LYS:HB3	1.94	0.48
1:A:125:LYS:O	1:A:126:LEU:HB2	2.13	0.47
2:B:15:LEU:HD21	1:E:106:LEU:HD11	1.97	0.46
1:G:67:GLN:HE22	2:H:18:LEU:HD23	1.81	0.46
2:B:15:LEU:HD22	1:E:110:LEU:HD21	1.98	0.46
1:A:72:GLU:O	1:A:76:LEU:HD13	2.18	0.44
1:E:120:GLN:O	1:E:124:THR:HG23	2.19	0.42
1:A:113:THR:HG21	2:F:18:LEU:HD12	2.02	0.42
1:C:122:LEU:C	1:C:122:LEU:HD23	2.40	0.41
1:A:76:LEU:CD1	4:A:52:HOH:O	2.65	0.40
1:A:100:CYS:SG	1:A:103:PRO:HA	2.61	0.40
1:E:104:GLN:O	1:E:108:GLN:HG3	2.22	0.40

All (2) 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)
1:A:84:GLY:O	1:G:86:TYR:OH[1_565]	1.98	0.22
2:D:18:LEU:CD1	1:G:113:THR:CG2[3_654]	2.06	0.14

## 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	69/73 (94%)	68 (99%)	1 (1%)	0	100	100
1	C	69/73 (94%)	69 (100%)	0	0	100	100
1	E	63/73 (86%)	62 (98%)	1 (2%)	0	100	100
1	G	64/73 (88%)	62 (97%)	2 (3%)	0	100	100
2	B	14/16 (88%)	14 (100%)	0	0	100	100
2	D	16/16 (100%)	16 (100%)	0	0	100	100
2	F	14/16 (88%)	13 (93%)	1 (7%)	0	100	100
2	H	12/16 (75%)	12 (100%)	0	0	100	100
All	All	321/356 (90%)	316 (98%)	5 (2%)	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	61/62 (98%)	59 (97%)	2 (3%)	50	37
1	C	61/62 (98%)	60 (98%)	1 (2%)	75	70
1	E	56/62 (90%)	53 (95%)	3 (5%)	31	17
1	G	57/62 (92%)	54 (95%)	3 (5%)	32	18

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	B	11/11 (100%)	11 (100%)	0	100	100
2	D	13/11 (118%)	13 (100%)	0	100	100
2	F	11/11 (100%)	11 (100%)	0	100	100
2	H	10/11 (91%)	9 (90%)	1 (10%)	11	4
All	All	280/292 (96%)	270 (96%)	10 (4%)	47	33

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	121	MET
1	A	126	LEU
1	C	125	LYS
1	E	112	GLN
1	E	123	LEU
1	E	126	LEU
1	G	61	LYS
1	G	123	LEU
1	G	126	LEU
2	H	14	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	67	GLN
1	A	75	GLN
1	C	67	GLN
1	E	67	GLN
1	E	75	GLN
1	E	83	GLN
1	E	102	GLN
1	E	108	GLN
1	E	111	GLN
1	E	120	GLN
1	G	67	GLN
1	G	83	GLN
1	G	104	GLN

### 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 ⓘ

2 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	$\# Z  > 2$
3	PO4	B	2	-	4,4,4	0.56	0	6,6,6	0.39	0
3	PO4	D	1	-	4,4,4	0.25	0	6,6,6	0.35	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
3	PO4	B	2	-	-	0/0/0/0	0/0/0/0
3	PO4	D	1	-	-	0/0/0/0	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	70/73 (95%)	-0.16	1 (1%) 72 74	11, 20, 35, 40	0
1	C	70/73 (95%)	-0.15	2 (2%) 49 50	11, 19, 36, 41	0
1	E	65/73 (89%)	-0.10	1 (1%) 70 72	12, 23, 39, 53	0
1	G	66/73 (90%)	0.17	4 (6%) 21 20	13, 28, 49, 52	0
2	B	16/16 (100%)	-0.17	0 100 100	16, 19, 33, 38	0
2	D	16/16 (100%)	-0.11	1 (6%) 19 19	16, 17, 42, 45	0
2	F	16/16 (100%)	-0.08	2 (12%) 5 4	18, 21, 44, 45	0
2	H	14/16 (87%)	-0.57	0 100 100	16, 17, 30, 37	0
All	All	333/356 (93%)	-0.09	11 (3%) 44 45	11, 21, 44, 53	0

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	62	ASP	4.2
1	G	61	LYS	3.6
1	G	86	TYR	2.6
1	C	58	SER	2.5
2	F	13	PRO	2.4
1	A	57	GLY	2.3
1	C	57	GLY	2.2
2	D	12	GLY	2.2
1	G	82	ALA	2.1
2	F	12	GLY	2.1
1	G	126	LEU	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
3	PO4	B	2	5/5	0.17	1.35	35,41,45,47	0
3	PO4	D	1	5/5	0.15	0.30	46,50,52,56	0

### 6.5 Other polymers

There are no such residues in this entry.