



# wwPDB X-ray Structure Validation Summary Report

Feb 28, 2014 – 06:23 PM GMT

PDB ID : 4KBP  
Title : KIDNEY BEAN PURPLE ACID PHOSPHATASE  
Authors : Klabunde, T.; Strater, N.; Krebs, B.  
Deposited on : 1995-10-02  
Resolution : 2.70 Å(reported)

This is a wwPDB validation summary 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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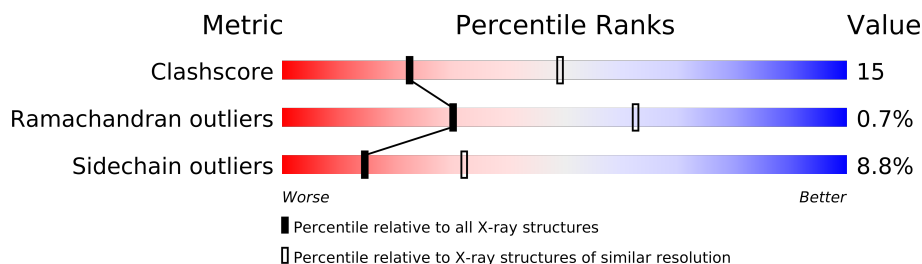
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : **NOT EXECUTED**  
EDS : **NOT EXECUTED**  
Percentile statistics : 21963  
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 2.70 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))
Clashscore	79885	1939 (2.70-2.70)
Ramachandran outliers	78287	1905 (2.70-2.70)
Sidechain outliers	78261	1905 (2.70-2.70)

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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	432	
1	B	432	
1	C	432	
1	D	432	

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 14340 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 PURPLE ACID PHOSPHATASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	424	Total	C	N	O	S	0	0	0
			3494	2243	605	636	10			
1	B	424	Total	C	N	O	S	0	0	0
			3494	2243	605	636	10			
1	C	424	Total	C	N	O	S	0	0	0
			3494	2243	605	636	10			
1	D	424	Total	C	N	O	S	0	0	0
			3494	2243	605	636	10			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	253	TYR	HIS	CONFLICT	UNP P80366
A	254	SER	ILE	CONFLICT	UNP P80366
B	253	TYR	HIS	CONFLICT	UNP P80366
B	254	SER	ILE	CONFLICT	UNP P80366
C	253	TYR	HIS	CONFLICT	UNP P80366
C	254	SER	ILE	CONFLICT	UNP P80366
D	253	TYR	HIS	CONFLICT	UNP P80366
D	254	SER	ILE	CONFLICT	UNP P80366

- Molecule 2 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula: C<sub>8</sub>H<sub>15</sub>NO<sub>6</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	B	1	Total	C	N	O	0	0
			14	8	1	5		
2	B	1	Total	C	N	O	0	0
			14	8	1	5		
2	B	1	Total	C	N	O	0	0
			14	8	1	5		
2	B	1	Total	C	N	O	0	0
			14	8	1	5		
2	C	1	Total	C	N	O	0	0
			14	8	1	5		
2	C	1	Total	C	N	O	0	0
			14	8	1	5		
2	C	1	Total	C	N	O	0	0
			14	8	1	5		
2	C	1	Total	C	N	O	0	0
			14	8	1	5		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	C	1	Total	C	N	O	0	0
			14	8	1	5		
2	D	1	Total	C	N	O	0	0
			14	8	1	5		
2	D	1	Total	C	N	O	0	0
			14	8	1	5		
2	D	1	Total	C	N	O	0	0
			14	8	1	5		
2	D	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 3 is FE (III) ION (three-letter code: FE) (formula: Fe).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Fe	0	0
			1	1		
3	A	1	Total	Fe	0	0
			1	1		
3	D	1	Total	Fe	0	0
			1	1		
3	C	1	Total	Fe	0	0
			1	1		

- Molecule 4 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Zn	0	0
			1	1		
4	A	1	Total	Zn	0	0
			1	1		
4	D	1	Total	Zn	0	0
			1	1		
4	C	1	Total	Zn	0	0
			1	1		

- Molecule 5 is PHOSPHATE ION (three-letter code: PO4) (formula: O<sub>4</sub>P).

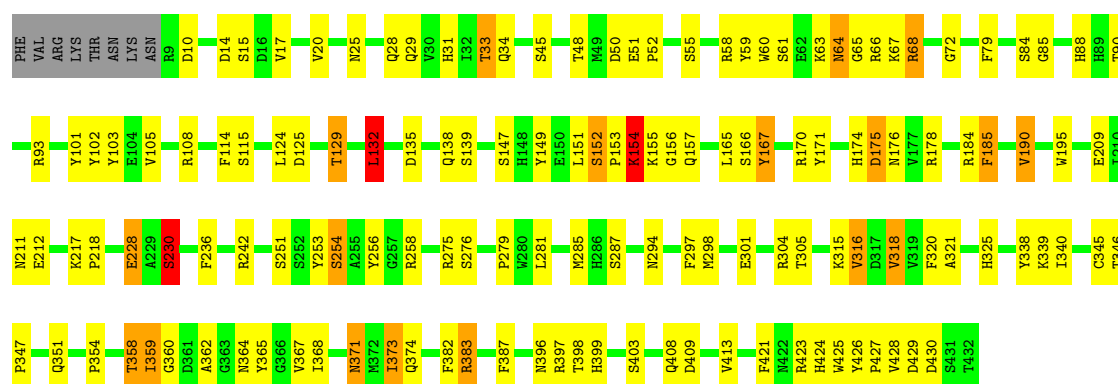


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	O	P	0	0
			5	4	1		
5	B	1	Total	O	P	0	0
			5	4	1		
5	C	1	Total	O	P	0	0
			5	4	1		
5	D	1	Total	O	P	0	0
			5	4	1		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	16	Total	O	0	0
			16	16		
6	B	13	Total	O	0	0
			13	13		
6	C	14	Total	O	0	0
			14	14		
6	D	13	Total	O	0	0
			13	13		







## 4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	132.70Å 347.30Å 128.70Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 2.70	Depositor
% Data completeness (in resolution range)	72.4 (10.00-2.70)	Depositor
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
Refinement program	X-PLOR 3.1	Depositor
R, $R_{free}$	0.192 , 0.228	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	14340	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	38.0	wwPDB-VP

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: PO4, FE, ZN, NAG

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	0.95	1/3613 (0.0%)	0.99	9/4912 (0.2%)
1	B	0.87	2/3613 (0.1%)	0.95	6/4912 (0.1%)
1	C	0.92	1/3613 (0.0%)	0.98	8/4912 (0.2%)
1	D	0.92	1/3613 (0.0%)	0.97	7/4912 (0.1%)
All	All	0.92	5/14452 (0.0%)	0.97	30/19648 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	2
1	D	0	2
All	All	0	4

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	345	CYS	CB-SG	-6.62	1.71	1.82
1	A	238	TYR	CD1-CE1	6.09	1.48	1.39
1	B	425	TRP	CB-CG	-5.96	1.39	1.50
1	C	345	CYS	CB-SG	-5.75	1.72	1.81
1	B	345	CYS	CB-SG	-5.10	1.73	1.81

The worst 5 of 30 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	135	ASP	CB-CG-OD1	-13.04	106.56	118.30
1	A	135	ASP	CB-CG-OD1	-12.71	106.87	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	135	ASP	CB-CG-OD1	-12.65	106.92	118.30
1	C	135	ASP	CB-CG-OD1	-12.01	107.49	118.30
1	C	135	ASP	CB-CG-OD2	8.18	125.66	118.30

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	223	TYR	Sidechain
1	B	426	TYR	Sidechain
1	D	167	TYR	Sidechain
1	D	253	TYR	Sidechain

## 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	3494	0	3306	94	0
1	B	3494	0	3305	113	0
1	C	3494	0	3305	98	0
1	D	3494	0	3306	106	0
2	A	70	0	65	2	0
2	B	70	0	65	1	0
2	C	70	0	65	1	0
2	D	70	0	65	1	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
5	A	5	0	0	0	0
5	B	5	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	C	5	0	0	0	0
5	D	5	0	0	0	0
6	A	16	0	0	0	0
6	B	13	0	0	4	0
6	C	14	0	0	2	0
6	D	13	0	0	2	0
All	All	14340	0	13482	410	0

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 15.

The worst 5 of 410 close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:398:THR:HG22	1:D:426:TYR:HB2	1.59	0.82
1:D:368:ILE:HG12	1:D:387:PHE:CZ	2.19	0.78
1:C:33:THR:HG23	1:C:195:TRP:O	1.87	0.75
1:B:426:TYR:N	1:B:427:PRO:HD3	2.01	0.75
1:B:275:ARG:NH2	1:B:315:LYS:O	2.21	0.73

There are no symmetry-related clashes.

## 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	422/432 (98%)	386 (92%)	34 (8%)	2 (0%)	38	70
1	B	422/432 (98%)	383 (91%)	34 (8%)	5 (1%)	19	45
1	C	422/432 (98%)	385 (91%)	34 (8%)	3 (1%)	30	62
1	D	422/432 (98%)	384 (91%)	36 (8%)	2 (0%)	38	70
All	All	1688/1728 (98%)	1538 (91%)	138 (8%)	12 (1%)	30	62

5 of 12 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	175	ASP
1	B	175	ASP
1	C	175	ASP
1	D	175	ASP
1	C	431	SER

### 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	373/381 (98%)	341 (91%)	32 (9%)	15	33
1	B	373/381 (98%)	340 (91%)	33 (9%)	14	32
1	C	373/381 (98%)	339 (91%)	34 (9%)	14	30
1	D	373/381 (98%)	341 (91%)	32 (9%)	15	33
All	All	1492/1524 (98%)	1361 (91%)	131 (9%)	14	32

5 of 131 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	359	ILE
1	C	129	THR
1	D	287	SER
1	B	378	GLU
1	C	33	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 53 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	364	ASN
1	C	34	GLN
1	D	364	ASN
1	B	399	HIS
1	C	25	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 32 ligands modelled in this entry, 8 are monoatomic - leaving 24 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z  > 2$	Counts	RMSZ	# $ Z  > 2$
2	NAG	A	433(A)	1	12,14,15	0.97	0	15,19,21	1.25	1 (6%)
2	NAG	A	434(A)	1	12,14,15	1.02	0	15,19,21	1.45	2 (13%)
2	NAG	A	435(A)	1	12,14,15	0.67	0	15,19,21	0.84	1 (6%)
2	NAG	A	436(A)	1	12,14,15	1.06	1 (8%)	15,19,21	1.33	2 (13%)
2	NAG	A	437(A)	1	12,14,15	0.81	0	15,19,21	1.12	1 (6%)
5	PO4	A	440	3,4	4,4,4	2.44	2 (50%)	6,6,6	0.32	0
2	NAG	B	433(A)	1	12,14,15	0.89	1 (8%)	15,19,21	1.04	0
2	NAG	B	434(A)	1	12,14,15	0.84	0	15,19,21	1.26	2 (13%)
2	NAG	B	435(A)	1	12,14,15	0.44	0	15,19,21	0.72	1 (6%)
2	NAG	B	436(A)	1	12,14,15	1.14	1 (8%)	15,19,21	1.25	3 (20%)
2	NAG	B	437(A)	1	12,14,15	0.75	0	15,19,21	1.01	0
5	PO4	B	440	3,4	4,4,4	1.51	1 (25%)	6,6,6	0.30	0
2	NAG	C	433(A)	1	12,14,15	0.99	1 (8%)	15,19,21	1.09	1 (6%)
2	NAG	C	434(A)	1	12,14,15	1.10	1 (8%)	15,19,21	1.45	3 (20%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	NAG	C	435(A)	1	12,14,15	0.63	0	15,19,21	0.82	1 (6%)
2	NAG	C	436(A)	1	12,14,15	1.07	1 (8%)	15,19,21	1.22	2 (13%)
2	NAG	C	437(A)	1	12,14,15	0.78	0	15,19,21	1.03	1 (6%)
5	PO4	C	440	3,4	4,4,4	1.84	2 (50%)	6,6,6	0.30	0
2	NAG	D	433(A)	1	12,14,15	0.79	0	15,19,21	1.26	3 (20%)
2	NAG	D	434(A)	1	12,14,15	1.09	1 (8%)	15,19,21	1.35	2 (13%)
2	NAG	D	435(A)	1	12,14,15	0.59	0	15,19,21	0.91	1 (6%)
2	NAG	D	436(A)	1	12,14,15	1.12	0	15,19,21	1.24	1 (6%)
2	NAG	D	437(A)	1	12,14,15	1.16	1 (8%)	15,19,21	1.23	1 (6%)
5	PO4	D	440	3,4	4,4,4	1.74	2 (50%)	6,6,6	0.31	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
2	NAG	A	433(A)	1	-	0/6/23/26	0/1/1/1
2	NAG	A	434(A)	1	-	0/6/23/26	0/1/1/1
2	NAG	A	435(A)	1	-	0/6/23/26	0/1/1/1
2	NAG	A	436(A)	1	-	0/6/23/26	0/1/1/1
2	NAG	A	437(A)	1	-	0/6/23/26	0/1/1/1
5	PO4	A	440	3,4	-	0/0/0/0	0/0/0/0
2	NAG	B	433(A)	1	-	0/6/23/26	0/1/1/1
2	NAG	B	434(A)	1	-	0/6/23/26	0/1/1/1
2	NAG	B	435(A)	1	-	0/6/23/26	0/1/1/1
2	NAG	B	436(A)	1	-	0/6/23/26	0/1/1/1
2	NAG	B	437(A)	1	-	0/6/23/26	0/1/1/1
5	PO4	B	440	3,4	-	0/0/0/0	0/0/0/0
2	NAG	C	433(A)	1	-	0/6/23/26	0/1/1/1
2	NAG	C	434(A)	1	-	0/6/23/26	0/1/1/1
2	NAG	C	435(A)	1	-	0/6/23/26	0/1/1/1
2	NAG	C	436(A)	1	-	0/6/23/26	0/1/1/1
2	NAG	C	437(A)	1	-	0/6/23/26	0/1/1/1
5	PO4	C	440	3,4	-	0/0/0/0	0/0/0/0
2	NAG	D	433(A)	1	-	0/6/23/26	0/1/1/1
2	NAG	D	434(A)	1	-	0/6/23/26	0/1/1/1
2	NAG	D	435(A)	1	-	0/6/23/26	0/1/1/1
2	NAG	D	436(A)	1	-	0/6/23/26	0/1/1/1
2	NAG	D	437(A)	1	-	0/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	PO4	D	440	3,4	-	0/0/0/0	0/0/0/0

The worst 5 of 15 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	440	PO4	P-O4	3.31	1.66	1.52
5	A	440	PO4	P-O2	-2.99	1.40	1.52
5	C	440	PO4	P-O4	2.44	1.62	1.52
2	C	433(A)	NAG	C3-C2	2.40	1.57	1.52
5	D	440	PO4	P-O4	2.40	1.62	1.52

The worst 5 of 29 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	434(A)	NAG	C4-C3-C2	-3.27	103.31	111.32
2	C	434(A)	NAG	C4-C3-C2	-3.27	103.31	111.32
2	D	434(A)	NAG	C4-C3-C2	-3.15	103.61	111.32
2	B	434(A)	NAG	C4-C3-C2	-3.09	103.75	111.32
2	A	434(A)	NAG	O3-C3-C2	2.48	114.29	109.09

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 ⓘ

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

### 6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

### 6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

### 6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.