



# wwPDB X-ray Structure Validation Summary Report

Feb 27, 2014 – 10:31 AM GMT

PDB ID : 1A2V  
Title : COPPER AMINE OXIDASE FROM HANSENULA POLYMORPHA  
Authors : Li, R.; Mathews, F.S.  
Deposited on : 1998-01-12  
Resolution : 2.40 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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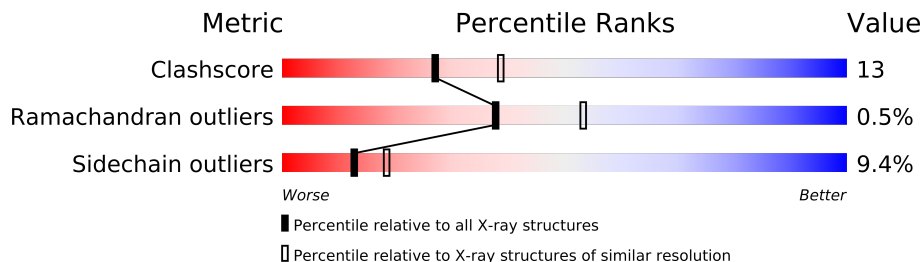
The following versions of software and data (see [references](#)) were used in the production of this report:

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 2.40 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	2789 (2.40-2.40)
Ramachandran outliers	78287	2736 (2.40-2.40)
Sidechain outliers	78261	2737 (2.40-2.40)

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	655	
1	B	655	
1	C	655	
1	D	655	
1	E	655	
1	F	655	

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 33726 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 METHYLAMINE OXIDASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	655	Total	C	N	O	S	0	0	0
			5194	3305	892	974	23			
1	B	655	Total	C	N	O	S	0	0	0
			5194	3305	892	974	23			
1	C	655	Total	C	N	O	S	0	0	0
			5194	3305	892	974	23			
1	D	655	Total	C	N	O	S	0	0	0
			5194	3305	892	974	23			
1	E	655	Total	C	N	O	S	0	0	0
			5194	3305	892	974	23			
1	F	655	Total	C	N	O	S	0	0	0
			5194	3305	892	974	23			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	405	TPQ	TYR	MODIFIED RESIDUE	UNP P12807
B	405	TPQ	TYR	MODIFIED RESIDUE	UNP P12807
C	405	TPQ	TYR	MODIFIED RESIDUE	UNP P12807
D	405	TPQ	TYR	MODIFIED RESIDUE	UNP P12807
E	405	TPQ	TYR	MODIFIED RESIDUE	UNP P12807
F	405	TPQ	TYR	MODIFIED RESIDUE	UNP P12807

- Molecule 2 is COPPER (II) ION (three-letter code: CU) (formula: Cu).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	D	1	Total	Cu	0	0
			1	1		
2	E	1	Total	Cu	0	0
			1	1		
2	B	1	Total	Cu	0	0
			1	1		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	C	1	Total 1	Cu 1	0	0
2	A	1	Total 1	Cu 1	0	0
2	F	1	Total 1	Cu 1	0	0

- Molecule 3 is water.

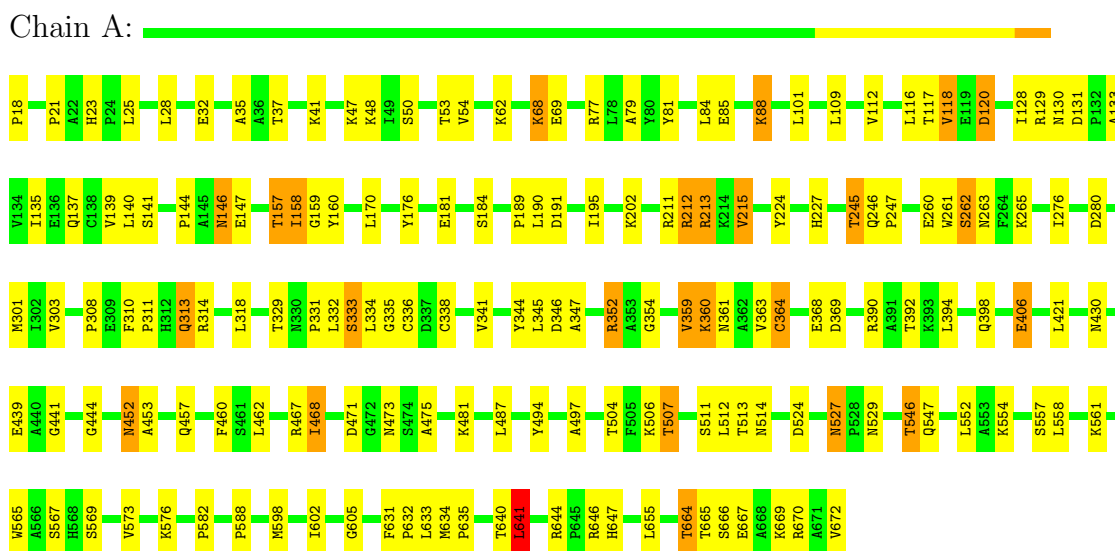
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	499	Total 499	O 499	0	0
3	B	453	Total 453	O 453	0	0
3	C	474	Total 474	O 474	0	0
3	D	415	Total 415	O 415	0	0
3	E	335	Total 335	O 335	0	0
3	F	380	Total 380	O 380	0	0

### 3 Residue-property plots

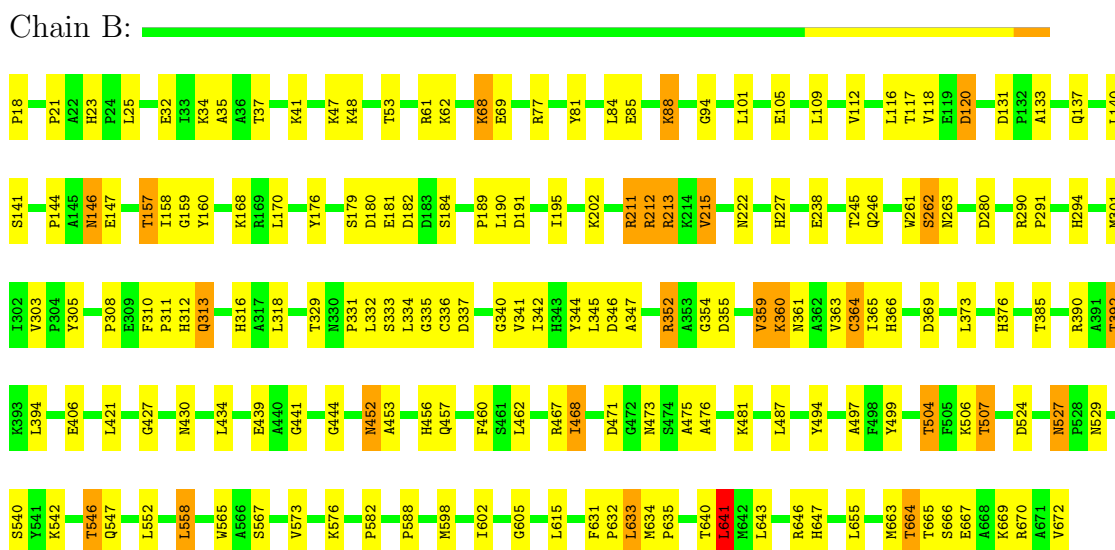
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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.

Note EDS was not executed.

#### • Molecule 1: METHYLAMINE OXIDASE



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Age Group	Percentage
18-24	35%
25-34	25%
35-44	15%
45-54	10%
55-64	5%
65-74	3%
75-84	2%
85+	1%

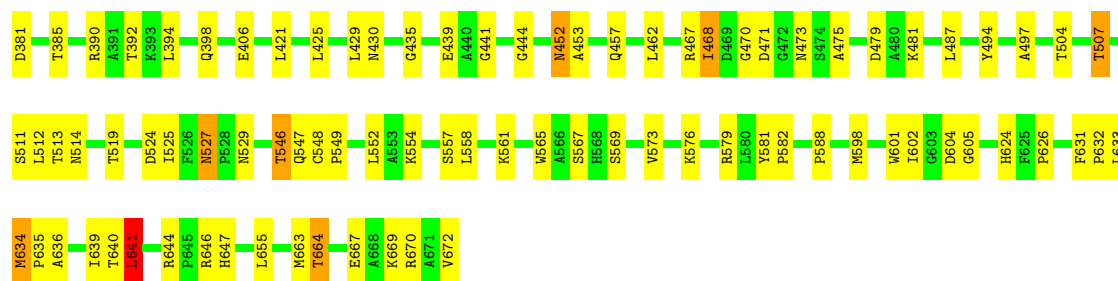


Age Group	Percentage
18-24	45%
25-34	35%
35-44	15%
45-54	5%
55-64	2%
65-74	1%
75-84	1%
85+	1%



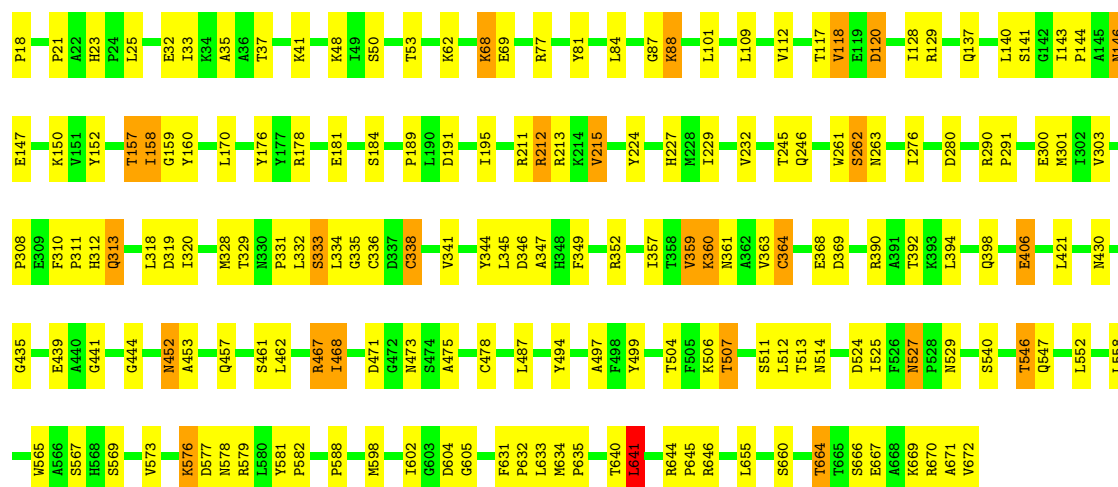
Frequency	Percentage
Daily	75%
Weekly	20%
Monthly	5%





• Molecule 1: METHYLAMINE OXIDASE

Chain F: 



## 4 Data and refinement statistics

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

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	138.77Å 148.22Å 234.01Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	100.00 – 2.40	Depositor
% Data completeness (in resolution range)	83.6 (100.00-2.40)	Depositor
$R_{merge}$	0.05	Depositor
$R_{sym}$	0.05	Depositor
Refinement program	X-PLOR 3.843	Depositor
R, $R_{free}$	0.184 , 0.224	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	33726	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	24.0	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: TPQ, CU

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	0.49	0/5329	0.74	2/7251 (0.0%)
1	B	0.53	1/5329 (0.0%)	0.77	2/7251 (0.0%)
1	C	0.53	0/5329	0.78	4/7251 (0.1%)
1	D	0.52	0/5329	0.77	2/7251 (0.0%)
1	E	0.50	0/5329	0.76	3/7251 (0.0%)
1	F	0.52	0/5329	0.77	3/7251 (0.0%)
All	All	0.52	1/31974 (0.0%)	0.76	16/43506 (0.0%)

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	D	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	364	CYS	CB-SG	-5.73	1.72	1.81

The worst 5 of 16 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	641	LEU	CA-CB-CG	7.25	131.96	115.30
1	A	641	LEU	CA-CB-CG	7.24	131.96	115.30
1	D	641	LEU	CA-CB-CG	7.03	131.47	115.30
1	E	641	LEU	CA-CB-CG	6.98	131.36	115.30
1	B	641	LEU	CA-CB-CG	6.80	130.95	115.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	305	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	5194	0	5033	129	0
1	B	5194	0	5033	141	0
1	C	5194	0	5033	129	0
1	D	5194	0	5033	150	0
1	E	5194	0	5033	150	0
1	F	5194	0	5033	140	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
2	E	1	0	0	0	0
2	F	1	0	0	0	0
3	A	499	0	0	7	0
3	B	453	0	0	6	0
3	C	474	0	0	5	0
3	D	415	0	0	9	0
3	E	335	0	0	5	0
3	F	380	0	0	7	0
All	All	33726	0	30198	791	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 13.

The worst 5 of 791 close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:546:THR:CG2	1:D:546:THR:HG21	1.57	1.32
1:A:546:THR:HG21	1:B:546:THR:CG2	1.64	1.27

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:E:546:THR:CG2	1:F:546:THR:HG21	1.64	1.26
1:A:546:THR:CG2	1:B:546:THR:HG21	1.67	1.24
1:C:546:THR:HG21	1:D:546:THR:CG2	1.69	1.22

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	652/655 (100%)	616 (94%)	33 (5%)	3 (0%)	38	53
1	B	652/655 (100%)	616 (94%)	34 (5%)	2 (0%)	50	68
1	C	652/655 (100%)	616 (94%)	32 (5%)	4 (1%)	33	47
1	D	652/655 (100%)	617 (95%)	32 (5%)	3 (0%)	38	53
1	E	652/655 (100%)	619 (95%)	30 (5%)	3 (0%)	38	53
1	F	652/655 (100%)	613 (94%)	36 (6%)	3 (0%)	38	53
All	All	3912/3930 (100%)	3697 (94%)	197 (5%)	18 (0%)	38	53

5 of 18 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	512	LEU
1	E	468	ILE
1	E	512	LEU
1	A	512	LEU
1	D	333	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	565/565 (100%)	512 (91%)	53 (9%)	13	18
1	B	565/565 (100%)	511 (90%)	54 (10%)	12	17
1	C	565/565 (100%)	511 (90%)	54 (10%)	12	17
1	D	565/565 (100%)	511 (90%)	54 (10%)	12	17
1	E	565/565 (100%)	511 (90%)	54 (10%)	12	17
1	F	565/565 (100%)	515 (91%)	50 (9%)	14	21
All	All	3390/3390 (100%)	3071 (91%)	319 (9%)	13	18

5 of 319 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	421	LEU
1	D	213	ARG
1	F	352	ARG
1	C	481	LYS
1	D	50	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 107 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	527	ASN
1	D	313	GLN
1	F	313	GLN
1	C	547	GLN
1	D	246	GLN

### 5.3.3 RNA ⓘ

There are no RNA chains in this entry.

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

6 non-standard protein/DNA/RNA residues 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	$\# Z  > 2$
1	TPQ	A	405	1	14,14,15	5.12	6 (42%)	17,19,21	1.05	1 (5%)
1	TPQ	B	405	1	14,14,15	4.97	6 (42%)	17,19,21	1.09	1 (5%)
1	TPQ	C	405	1	14,14,15	5.13	7 (50%)	17,19,21	1.13	1 (5%)
1	TPQ	D	405	1	14,14,15	5.63	5 (35%)	17,19,21	1.08	1 (5%)
1	TPQ	E	405	1	14,14,15	5.30	5 (35%)	17,19,21	1.12	1 (5%)
1	TPQ	F	405	1	14,14,15	5.05	5 (35%)	17,19,21	1.19	1 (5%)

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
1	TPQ	A	405	1	-	0/4/22/24	0/1/1/1
1	TPQ	B	405	1	-	0/4/22/24	0/1/1/1
1	TPQ	C	405	1	-	0/4/22/24	0/1/1/1
1	TPQ	D	405	1	-	0/4/22/24	0/1/1/1
1	TPQ	E	405	1	-	0/4/22/24	0/1/1/1
1	TPQ	F	405	1	-	0/4/22/24	0/1/1/1

The worst 5 of 34 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	405	TPQ	O-C	18.98	1.24	1.11
1	E	405	TPQ	O-C	18.07	1.23	1.11
1	A	405	TPQ	O-C	17.18	1.23	1.11
1	C	405	TPQ	O-C	17.16	1.23	1.11
1	F	405	TPQ	O-C	16.69	1.22	1.11

The worst 5 of 6 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	405	TPQ	C-CA-N	3.08	116.91	113.83
1	F	405	TPQ	C-CA-N	3.05	116.88	113.83
1	C	405	TPQ	C-CA-N	2.85	116.68	113.83
1	B	405	TPQ	C-CA-N	2.72	116.54	113.83
1	A	405	TPQ	C-CA-N	2.60	116.43	113.83

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

Of 6 ligands modelled in this entry, 6 are monoatomic - leaving 0 for Mogul analysis.

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 ⓘ

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.