



wwPDB X-ray Structure Validation Summary Report ⓘ

May 18, 2020 – 09:47 pm BST

PDB ID : 3AOW
Title : Crystal structure of Pyrococcus horikoshii kynurenine aminotransferase in complex with AKG
Authors : Okada, K.; Angkawidjaja, C.; Koga, Y.; Takano, K.; Kanaya, S.
Deposited on : 2010-10-07
Resolution : 1.56 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.11
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)	:	2.11

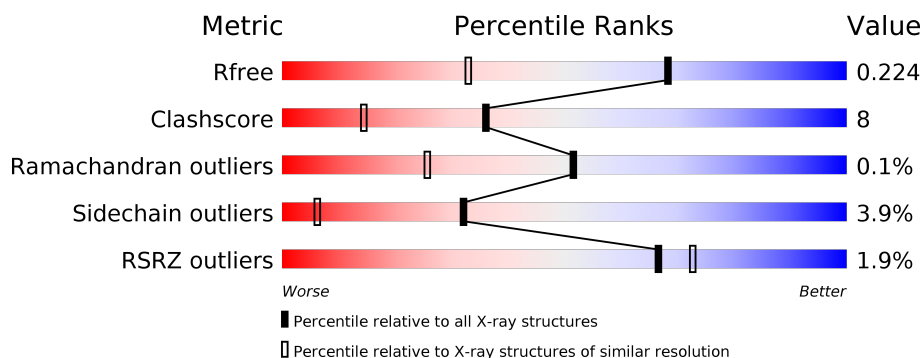
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.56 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}	130704	1483 (1.56-1.56)
Clashscore	141614	1529 (1.56-1.56)
Ramachandran outliers	138981	1498 (1.56-1.56)
Sidechain outliers	138945	1495 (1.56-1.56)
RSRZ outliers	127900	1465 (1.56-1.56)

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 ≥ 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 $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	448	<div> <div>2%</div> <div> <div></div> <div>73%</div> <div>15%</div> <div>••</div> <div>10%</div> </div> </div>
1	C	448	<div> <div>2%</div> <div> <div></div> <div>75%</div> <div>14%</div> <div>•</div> <div>10%</div> </div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7382 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.

- Molecule 1 is a protein called Putative uncharacterized protein PH0207.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	404	Total	C	N	O	S	0	0	0
			3247	2093	536	602	16			
1	C	404	Total	C	N	O	S	0	0	0
			3247	2093	536	602	16			

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	EXPRESSION TAG	UNP O57946
A	-18	GLY	-	EXPRESSION TAG	UNP O57946
A	-17	SER	-	EXPRESSION TAG	UNP O57946
A	-16	SER	-	EXPRESSION TAG	UNP O57946
A	-15	HIS	-	EXPRESSION TAG	UNP O57946
A	-14	HIS	-	EXPRESSION TAG	UNP O57946
A	-13	HIS	-	EXPRESSION TAG	UNP O57946
A	-12	HIS	-	EXPRESSION TAG	UNP O57946
A	-11	HIS	-	EXPRESSION TAG	UNP O57946
A	-10	HIS	-	EXPRESSION TAG	UNP O57946
A	-9	SER	-	EXPRESSION TAG	UNP O57946
A	-8	SER	-	EXPRESSION TAG	UNP O57946
A	-7	GLY	-	EXPRESSION TAG	UNP O57946
A	-6	LEU	-	EXPRESSION TAG	UNP O57946
A	-5	VAL	-	EXPRESSION TAG	UNP O57946
A	-4	PRO	-	EXPRESSION TAG	UNP O57946
A	-3	ARG	-	EXPRESSION TAG	UNP O57946
A	-2	GLY	-	EXPRESSION TAG	UNP O57946
A	-1	SER	-	EXPRESSION TAG	UNP O57946
A	0	HIS	-	EXPRESSION TAG	UNP O57946
C	-19	MET	-	EXPRESSION TAG	UNP O57946
C	-18	GLY	-	EXPRESSION TAG	UNP O57946
C	-17	SER	-	EXPRESSION TAG	UNP O57946
C	-16	SER	-	EXPRESSION TAG	UNP O57946
C	-15	HIS	-	EXPRESSION TAG	UNP O57946

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-14	HIS	-	EXPRESSION TAG	UNP O57946
C	-13	HIS	-	EXPRESSION TAG	UNP O57946
C	-12	HIS	-	EXPRESSION TAG	UNP O57946
C	-11	HIS	-	EXPRESSION TAG	UNP O57946
C	-10	HIS	-	EXPRESSION TAG	UNP O57946
C	-9	SER	-	EXPRESSION TAG	UNP O57946
C	-8	SER	-	EXPRESSION TAG	UNP O57946
C	-7	GLY	-	EXPRESSION TAG	UNP O57946
C	-6	LEU	-	EXPRESSION TAG	UNP O57946
C	-5	VAL	-	EXPRESSION TAG	UNP O57946
C	-4	PRO	-	EXPRESSION TAG	UNP O57946
C	-3	ARG	-	EXPRESSION TAG	UNP O57946
C	-2	GLY	-	EXPRESSION TAG	UNP O57946
C	-1	SER	-	EXPRESSION TAG	UNP O57946
C	0	HIS	-	EXPRESSION TAG	UNP O57946

- PLP
-
- The diagram shows the chemical structure of PLP (Pyridoxal Phosphate). It consists of a central pyridine ring. At the 2-position of the ring, there is a hydroxyl group (-OH) labeled O3. At the 3-position, there is a carboxylate group (-COO-) labeled O4A and O4P. At the 4-position, there is a phosphate group (-PO3H2) labeled O1P, O2P, O3P, and P. At the 5-position, there is a hydroxyl group (-OH) labeled O4P. At the 6-position, there is a hydroxyl group (-OH) labeled O3A. The nitrogen atom in the ring is labeled N1. The carbon atoms are labeled C2, C3, C4, C5, and C6. The phosphate group is shown with a central phosphorus atom (P) bonded to four oxygen atoms (O1P, O2P, O3P, O4P). The carboxylate group is shown with a central carbon atom (C4A) bonded to two oxygen atoms (O4A, O4P) and a hydrogen atom (H4A). The hydroxyl groups are shown with a central oxygen atom (O3, O3A, O4P) bonded to a hydrogen atom (H3, H3A, H4P).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total 15	C 8	N 1	O 5	P 1	0	0
2	C	1	Total 15	C 8	N 1	O 5	P 1	0	0

- 



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			10	5	5		
3	C	1	Total	C	O	0	0
			10	5	5		

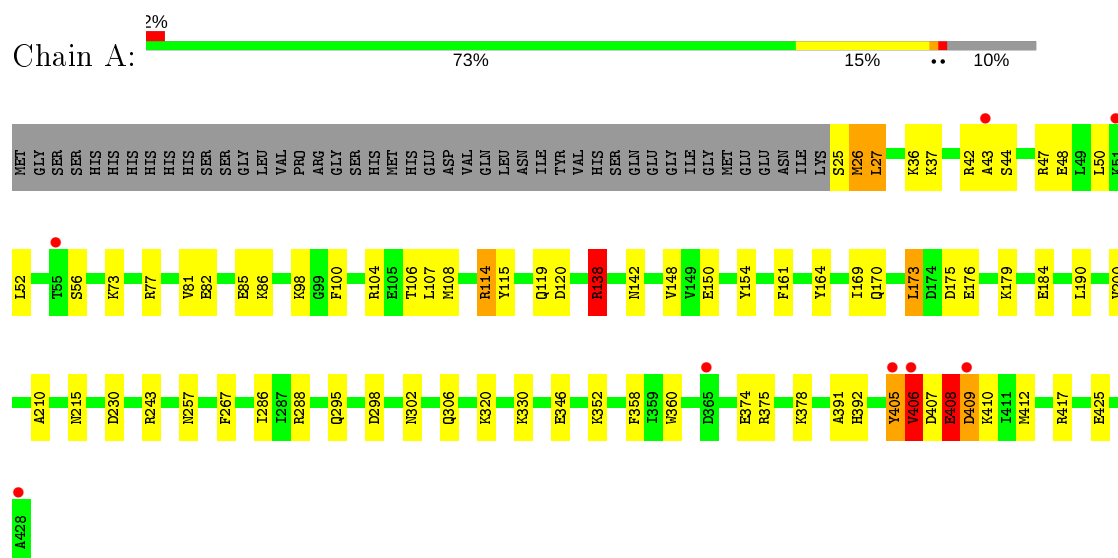
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	422	Total	O	0	0
			422	422		
4	C	416	Total	O	0	0
			416	416		

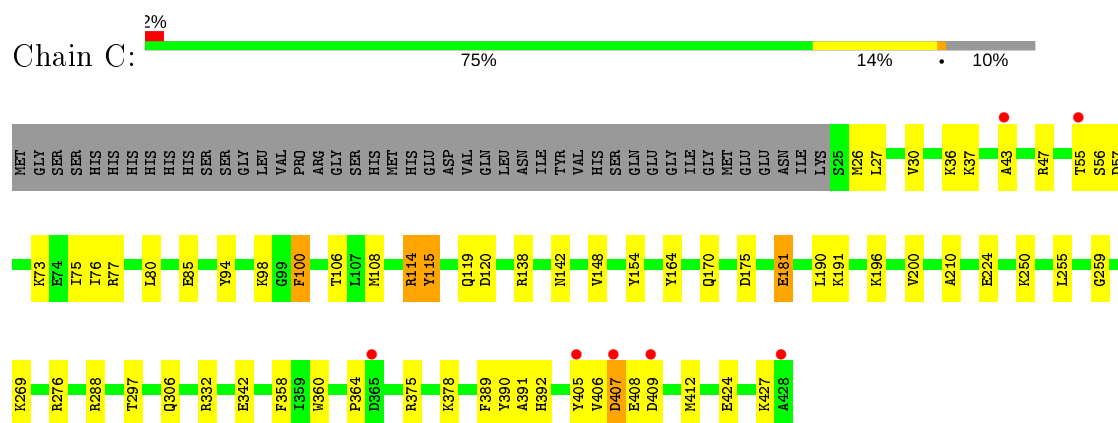
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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 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: Putative uncharacterized protein PH0207



- Molecule 1: Putative uncharacterized protein PH0207



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	85.77Å 70.98Å 136.72Å 90.00° 89.99° 90.00°	Depositor
Resolution (Å)	26.81 – 1.56 26.81 – 1.56	Depositor EDS
% Data completeness (in resolution range)	99.5 (26.81-1.56) 99.5 (26.81-1.56)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	0.04	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.84 (at 1.57Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.177 , 0.224 0.179 , 0.224	Depositor DCC
R_{free} test set	5780 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	20.6	Xtriage
Anisotropy	0.092	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 40.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.477 for -h,-k,l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	7382	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.71% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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.

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: AKG, PLP

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	1.25	3/3313 (0.1%)	1.18	14/4463 (0.3%)
1	C	1.25	3/3313 (0.1%)	1.16	6/4463 (0.1%)
All	All	1.25	6/6626 (0.1%)	1.17	20/8926 (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	A	0	3

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

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	115	TYR	CE2-CZ	-5.87	1.30	1.38
1	C	390	TYR	CG-CD2	5.70	1.46	1.39
1	A	330	LYS	CE-NZ	5.66	1.63	1.49
1	A	358	PHE	CD2-CE2	5.57	1.50	1.39
1	A	184	GLU	CB-CG	5.12	1.61	1.52

The worst 5 of 20 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	417	ARG	NE-CZ-NH2	-5.99	117.30	120.30
1	A	173	LEU	CB-CG-CD2	-5.98	100.84	111.00
1	A	409	ASP	CB-CG-OD2	5.96	123.66	118.30
1	C	98	LYS	CD-CE-NZ	-5.88	98.19	111.70
1	A	267	PHE	CB-CG-CD1	-5.87	116.69	120.80

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	27	LEU	Peptide
1	A	405	TYR	Peptide
1	A	406	VAL	Peptide

5.2 Too-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 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	A	3247	0	3306	60	0
1	C	3247	0	3306	55	0
2	A	15	0	6	1	0
2	C	15	0	6	2	0
3	A	10	0	4	1	0
3	C	10	0	4	2	0
4	A	422	0	0	19	0
4	C	416	0	0	18	0
All	All	7382	0	6632	105	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 8.

The worst 5 of 105 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:85:GLU:HG2	4:C:780:HOH:O	1.46	1.12
1:A:320:LYS:HE2	4:A:782:HOH:O	1.62	0.97
1:A:142:ASN:HD21	1:C:36:LYS:H	1.08	0.97
1:A:169:ILE:CD1	1:A:190:LEU:HD11	1.95	0.96
1:C:170:GLN:NE2	1:C:392:HIS:H	1.67	0.93

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	Percentiles	
1	A	402/448 (90%)	387 (96%)	14 (4%)	1 (0%)	47	23
1	C	402/448 (90%)	388 (96%)	14 (4%)	0	100	100
All	All	804/896 (90%)	775 (96%)	28 (4%)	1 (0%)	51	26

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	408	GLU

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 sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	349/388 (90%)	334 (96%)	15 (4%)	29	5
1	C	349/388 (90%)	337 (97%)	12 (3%)	37	9
All	All	698/776 (90%)	671 (96%)	27 (4%)	32	6

5 of 27 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	406	VAL
1	C	27	LEU
1	C	405	TYR
1	A	408	GLU
1	A	50	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 8 such sidechains are listed below:

Mol	Chain	Res	Type
1	A	306	GLN
1	C	306	GLN
1	C	142	ASN
1	A	170	GLN
1	C	119	GLN

5.3.3 RNA ⓘ

There are no RNA molecules 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 ⓘ

4 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	AKG	A	430	-	3,9,9	2.01	1 (33%)	4,11,11	4.32	3 (75%)
2	PLP	A	429	1	15,15,16	1.42	1 (6%)	20,22,23	1.48	4 (20%)
3	AKG	C	430	-	3,9,9	0.92	0	4,11,11	3.11	2 (50%)
2	PLP	C	429	1	15,15,16	1.48	3 (20%)	20,22,23	1.66	5 (25%)

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	AKG	A	430	-	-	1/3/9/9	-
2	PLP	A	429	1	-	0/6/6/8	0/1/1/1
3	AKG	C	430	-	-	0/3/9/9	-
2	PLP	C	429	1	-	0/6/6/8	0/1/1/1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	429	PLP	C3-C2	3.61	1.44	1.40
2	C	429	PLP	C2A-C2	2.88	1.55	1.50
3	A	430	AKG	O5-C2	2.86	1.26	1.22
2	C	429	PLP	C6-N1	2.52	1.39	1.34
2	C	429	PLP	C3-C2	2.11	1.43	1.40

The worst 5 of 14 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	430	AKG	C3-C4-C5	-6.41	101.92	112.67
3	C	430	AKG	C3-C4-C5	-5.34	103.72	112.67
3	A	430	AKG	C4-C3-C2	-4.33	103.81	113.14
2	C	429	PLP	C2A-C2-C3	3.38	125.07	120.89
3	A	430	AKG	O5-C2-C3	-3.34	114.68	120.38

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	430	AKG	C1-C2-C3-C4

There are no ring outliers.

4 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	430	AKG	1	0
2	A	429	PLP	1	0
3	C	430	AKG	2	0
2	C	429	PLP	2	0

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](#)

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, 95th 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 ²)	Q<0.9
1	A	404/448 (90%)	-0.32	8 (1%) 65 71	14, 21, 44, 60	0
1	C	404/448 (90%)	-0.31	7 (1%) 70 75	14, 21, 43, 58	0
All	All	808/896 (90%)	-0.31	15 (1%) 66 73	14, 21, 44, 60	0

The worst 5 of 15 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	55	THR	4.8
1	A	428	ALA	4.2
1	C	55	THR	3.9
1	A	409	ASP	3.6
1	A	405	TYR	3.4

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

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

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. The B-factors column lists the minimum, median, 95th 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	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	AKG	A	430	10/10	0.71	0.36	37,41,45,47	0
3	AKG	C	430	10/10	0.71	0.33	35,38,46,47	0
2	PLP	A	429	15/16	0.98	0.07	13,15,17,18	0
2	PLP	C	429	15/16	0.98	0.07	14,15,17,18	0

6.5 Other polymers [i](#)

There are no such residues in this entry.