



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 07:30 AM UTC

PDB ID : 6F2B / pdb_00006f2b
Title : Crystal structure of the complex Fe(II)/alpha-ketoglutarate dependent dioxxygenase KDO1 with Fe(II)/alpha-ketoglutarate
Authors : Isabet, T.; Stura, E.A.; Legrand, P.; Zaparucha, A.; Bastard, K.
Deposited on : 2017-11-24
Resolution : 2.00 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 ⓘ 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 ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

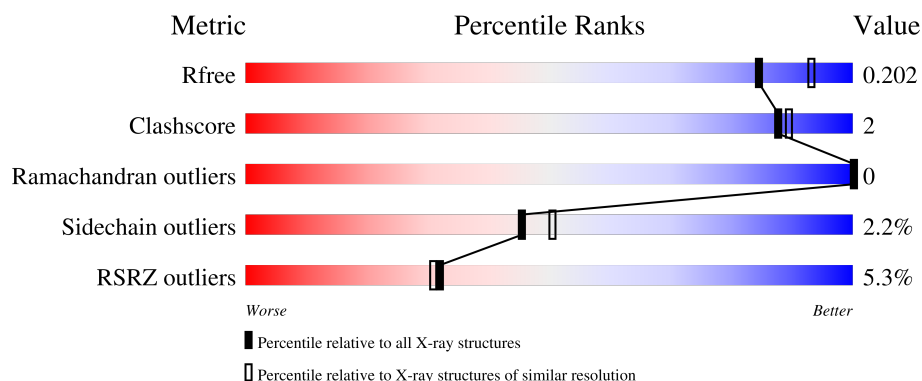
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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}	180053	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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 $\leq 5\%$. 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	358	5% 82% 8% 9%
1	B	358	5% 83% • 12%
1	C	358	6% 84% 7% 9%
1	D	358	3% 81% 6% • 13%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	AKG	B	402	-	X	-	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 11275 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.

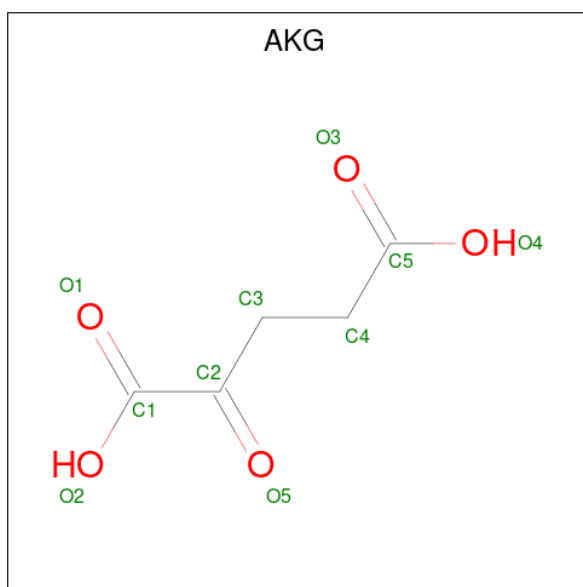
- Molecule 1 is a protein called L-lysine 3-hydroxylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	325	Total	C	N	O	S	0	5	0
			2541	1594	456	475	16			
1	B	316	Total	C	N	O	S	0	5	0
			2481	1558	446	464	13			
1	C	326	Total	C	N	O	S	0	10	0
			2594	1626	470	482	16			
1	D	312	Total	C	N	O	S	0	4	0
			2452	1541	439	459	13			

- Molecule 2 is FE (II) ION (CCD ID: FE2) (formula: Fe).

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

- Molecule 3 is 2-OXOGLUTARIC ACID (CCD ID: AKG) (formula: C₅H₆O₅).



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

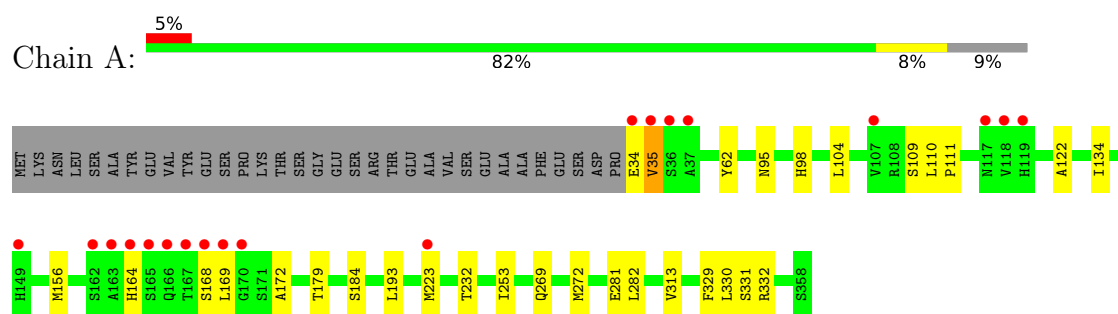
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	343	Total	O	0	0
			343	343		
4	B	235	Total	O	0	1
			235	235		
4	C	328	Total	O	0	3
			328	328		
4	D	257	Total	O	0	0
			257	257		

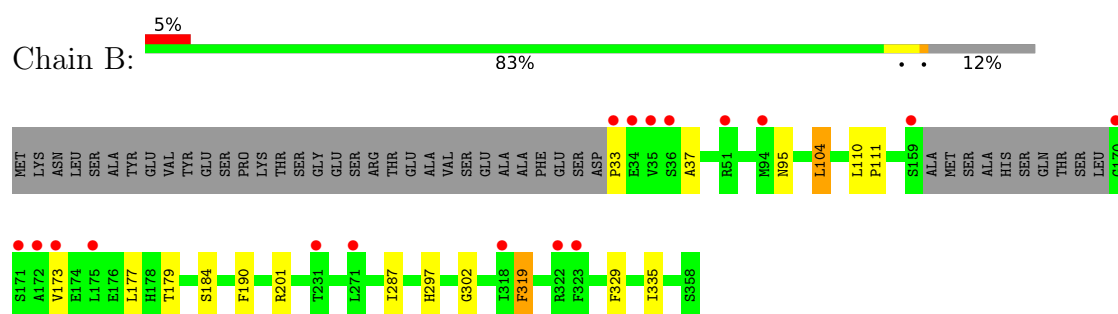
3 Residue-property plots [i](#)

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 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.

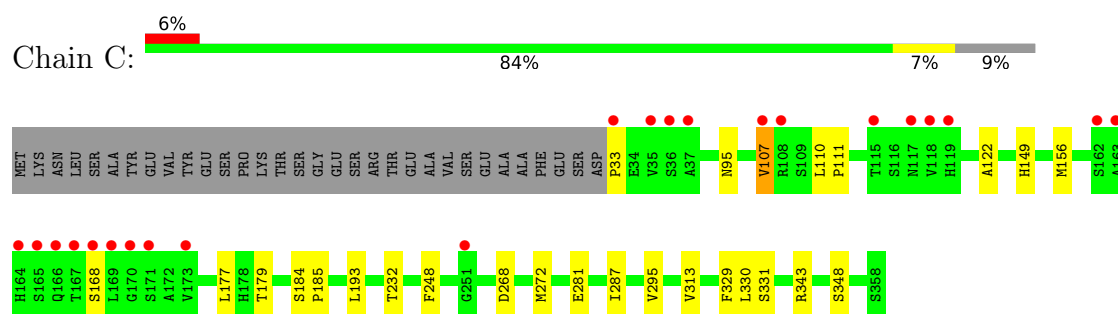
- Molecule 1: L-lysine 3-hydroxylase



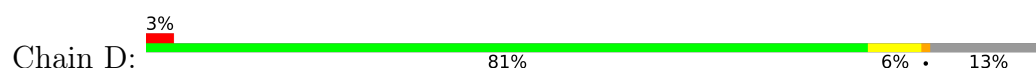
- Molecule 1: L-lysine 3-hydroxylase

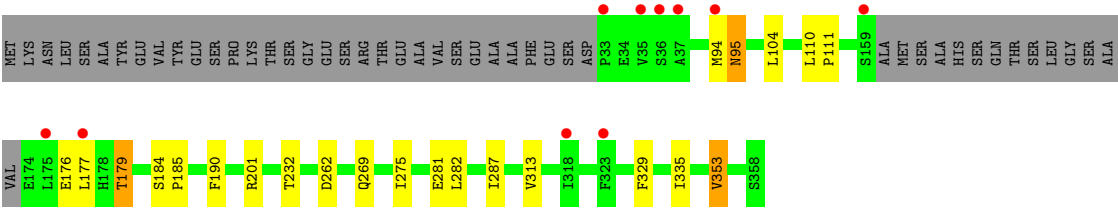


- Molecule 1: L-lysine 3-hydroxylase



- Molecule 1: L-lysine 3-hydroxylase





4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	56.82Å 67.97Å 110.43Å 107.78° 102.79° 93.45°	Depositor
Resolution (Å)	40.00 – 2.00 40.00 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.1 (40.00-2.00) 99.1 (40.00-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.28 (at 2.00Å)	Xtriage
Refinement program	BUSTER 2.10.3	Depositor
R, R_{free}	0.171 , 0.201 0.172 , 0.202	Depositor DCC
R_{free} test set	5033 reflections (4.94%)	wwPDB-VP
Wilson B-factor (Å ²)	34.5	Xtriage
Anisotropy	0.377	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 51.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.006 for -h,-k,h+k+l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	11275	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.65% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: FE2, AKG

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	0.89	0/2596	1.15	5/3524 (0.1%)
1	B	0.82	0/2534	1.14	7/3439 (0.2%)
1	C	0.84	0/2651	1.14	3/3598 (0.1%)
1	D	0.82	0/2505	1.15	6/3400 (0.2%)
All	All	0.84	0/10286	1.14	21/13961 (0.2%)

There are no bond length outliers.

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	37	ALA	N-CA-C	-8.99	101.28	112.88
1	B	95	ASN	CA-CB-CG	8.12	120.72	112.60
1	B	329	PHE	CA-CB-CG	7.86	121.66	113.80
1	D	329	PHE	CA-CB-CG	7.82	121.62	113.80
1	D	95	ASN	CA-CB-CG	7.77	120.37	112.60
1	A	329	PHE	CA-CB-CG	7.74	121.54	113.80
1	C	329	PHE	CA-CB-CG	7.72	121.52	113.80
1	C	95	ASN	CA-CB-CG	7.67	120.27	112.60
1	A	34	GLU	CA-C-N	6.46	128.70	120.56
1	A	34	GLU	C-N-CA	6.46	128.70	120.56
1	B	319	PHE	CA-CB-CG	6.43	120.23	113.80
1	D	353	VAL	N-CA-CB	6.20	117.89	110.95
1	C	184	SER	N-CA-C	5.76	117.97	109.62
1	D	184	SER	N-CA-C	5.67	117.84	109.62
1	B	184	SER	N-CA-C	5.62	117.76	109.62
1	A	184	SER	N-CA-C	5.62	117.76	109.62
1	D	95	ASN	N-CA-C	-5.48	103.16	110.55
1	D	179	THR	N-CA-C	-5.36	101.54	110.17
1	B	179	THR	N-CA-C	-5.19	101.24	109.96
1	B	95	ASN	N-CA-C	-5.13	103.62	110.55

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	95	ASN	CA-CB-CG	5.04	117.64	112.60

There are no chirality outliers.

There are no planarity outliers.

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	2541	0	2491	12	0
1	B	2481	0	2433	6	0
1	C	2594	0	2548	15	0
1	D	2452	0	2405	11	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
3	A	10	0	4	1	0
3	B	10	0	4	0	0
3	C	10	0	4	2	0
3	D	10	0	4	0	0
4	A	343	0	0	0	0
4	B	235	0	0	0	0
4	C	328	0	0	1	0
4	D	257	0	0	1	0
All	All	11275	0	9893	39	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 2.

All (39) 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:343:ARG:HH21	1:D:94:MET:HE1	1.53	0.74
1:D:179:THR:HG23	1:D:313:VAL:HG12	1.75	0.68
1:A:168:SER:HB2	1:A:172:ALA:HB3	1.79	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:179[A]:THR:HG23	1:C:313:VAL:HG12	1.79	0.63
1:A:179:THR:HG23	1:A:313:VAL:HG12	1.83	0.61
1:D:185:PRO:HD2	4:D:502:HOH:O	2.03	0.59
1:C:185:PRO:HD2	4:C:551:HOH:O	2.03	0.58
1:C:343:ARG:HH21	1:D:94:MET:CE	2.18	0.56
1:D:94:MET:HE1	1:D:262:ASP:HB2	1.90	0.53
1:C:268:ASP:O	1:C:272[B]:MET:HG2	2.09	0.53
1:A:164:HIS:HA	1:D:275:ILE:O	2.08	0.53
1:B:33:PRO:HD3	1:B:297:HIS:HB2	1.92	0.52
1:A:156[B]:MET:HB2	1:A:330:LEU:HB2	1.92	0.50
1:C:193:LEU:HD21	3:C:402:AKG:O1	2.12	0.49
1:A:110:LEU:HD12	1:A:111:PRO:HD2	1.96	0.48
1:D:177:LEU:HD23	1:D:269:GLN:HE21	1.78	0.48
1:A:193:LEU:HD21	3:A:402:AKG:O2	2.14	0.47
1:C:107:VAL:HG21	1:C:110:LEU:HD13	1.96	0.47
1:B:104:LEU:HB2	1:B:302:GLY:O	2.15	0.47
1:D:110:LEU:HD12	1:D:111:PRO:HD2	1.97	0.47
1:A:104:LEU:HD21	1:A:134:ILE:HD11	1.96	0.47
1:A:62:TYR:CD1	1:C:248:PHE:HB2	2.51	0.46
1:B:110:LEU:HD12	1:B:111:PRO:HD2	1.98	0.46
1:D:177:LEU:HD21	1:D:287:ILE:HG12	1.97	0.46
1:A:35:VAL:HG13	1:A:98:HIS:NE2	2.31	0.46
1:B:201:ARG:HG3	1:B:319:PHE:HB2	1.98	0.45
1:C:110:LEU:HD12	1:C:111:PRO:HD2	1.98	0.45
1:C:348:SER:OG	1:D:94:MET:HE3	2.16	0.45
1:C:177:LEU:HD21	1:C:287:ILE:HG12	1.97	0.45
1:C:156[B]:MET:HB3	1:C:330:LEU:HB2	1.99	0.43
1:B:177:LEU:HD21	1:B:287:ILE:HG12	2.01	0.43
1:B:190:PHE:HB2	1:B:335:ILE:HB	2.02	0.42
1:A:156[B]:MET:HG3	1:A:332:ARG:HB2	2.02	0.42
1:C:122:ALA:HB1	1:C:331:SER:HB2	2.03	0.41
1:D:190:PHE:HB2	1:D:335:ILE:HB	2.03	0.41
1:C:33:PRO:HB2	1:C:295:VAL:HG12	2.03	0.41
1:A:253:ILE:HD11	1:A:272:MET:HE3	2.03	0.40
1:A:122:ALA:HB1	1:A:331:SER:HB2	2.03	0.40
1:C:156[B]:MET:SD	3:C:402:AKG:H32	2.62	0.40

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	328/358 (92%)	320 (98%)	8 (2%)	0	100	100
1	B	317/358 (88%)	312 (98%)	5 (2%)	0	100	100
1	C	334/358 (93%)	330 (99%)	4 (1%)	0	100	100
1	D	312/358 (87%)	309 (99%)	3 (1%)	0	100	100
All	All	1291/1432 (90%)	1271 (98%)	20 (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	269/293 (92%)	261 (97%)	8 (3%)	36	38
1	B	262/293 (89%)	260 (99%)	2 (1%)	73	80
1	C	275/293 (94%)	269 (98%)	6 (2%)	45	50
1	D	260/293 (89%)	252 (97%)	8 (3%)	35	37
All	All	1066/1172 (91%)	1042 (98%)	24 (2%)	45	49

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	35	VAL
1	A	109	SER

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Mol	Chain	Res	Type
1	A	169	LEU
1	A	223	MET
1	A	232	THR
1	A	269	GLN
1	A	281	GLU
1	A	282	LEU
1	B	104	LEU
1	B	173	VAL
1	C	107	VAL
1	C	149[A]	HIS
1	C	149[B]	HIS
1	C	168	SER
1	C	232	THR
1	C	281	GLU
1	D	95	ASN
1	D	104	LEU
1	D	176	GLU
1	D	201	ARG
1	D	232	THR
1	D	281	GLU
1	D	282	LEU
1	D	353	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	131	GLN
1	A	269	GLN
1	B	106	GLN
1	B	131	GLN
1	B	314	HIS
1	C	131	GLN
1	C	269	GLN
1	C	283	GLN
1	D	106	GLN
1	D	131	GLN
1	D	269	GLN

5.3.3 RNA ⓘ

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 4 are monoatomic - leaving 4 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	$\# Z > 2$
3	AKG	D	402	2	9,9,9	0.21	0	11,11,11	0.71	0
3	AKG	C	402	2	9,9,9	0.31	0	11,11,11	1.08	2 (18%)
3	AKG	B	402	2	9,9,9	1.91	4 (44%)	11,11,11	3.13	5 (45%)
3	AKG	A	402	2	9,9,9	2.12	4 (44%)	11,11,11	2.77	5 (45%)

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	D	402	2	-	1/9/9/9	-
3	AKG	C	402	2	-	3/9/9/9	-
3	AKG	B	402	2	-	6/9/9/9	-
3	AKG	A	402	2	-	1/9/9/9	-

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	402	AKG	O1-C1	3.51	1.31	1.22

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	402	AKG	O2-C1	-3.24	1.21	1.30
3	A	402	AKG	O3-C5	3.15	1.32	1.22
3	B	402	AKG	O2-C1	-3.11	1.22	1.30
3	B	402	AKG	O1-C1	2.90	1.29	1.22
3	B	402	AKG	O3-C5	2.80	1.31	1.22
3	A	402	AKG	O4-C5	-2.70	1.21	1.30
3	B	402	AKG	O4-C5	-2.61	1.22	1.30

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	402	AKG	O1-C1-C2	-8.04	111.81	121.81
3	A	402	AKG	O1-C1-C2	-6.95	113.17	121.81
3	B	402	AKG	O2-C1-C2	4.32	125.57	113.59
3	A	402	AKG	O2-C1-C2	3.94	124.52	113.59
3	B	402	AKG	O3-C5-C4	-3.20	112.95	123.09
3	B	402	AKG	O4-C5-C4	3.03	123.57	114.00
3	C	402	AKG	C3-C2-C1	2.72	120.48	115.86
3	A	402	AKG	C3-C2-C1	2.63	120.33	115.86
3	A	402	AKG	O4-C5-C4	2.55	122.06	114.00
3	A	402	AKG	O3-C5-C4	-2.48	115.22	123.09
3	B	402	AKG	C3-C2-C1	2.17	119.55	115.86
3	C	402	AKG	O1-C1-C2	2.01	124.31	121.81

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	402	AKG	C1-C2-C3-C4
3	B	402	AKG	O1-C1-C2-O5
3	A	402	AKG	C1-C2-C3-C4
3	C	402	AKG	C1-C2-C3-C4
3	B	402	AKG	O2-C1-C2-O5
3	B	402	AKG	O2-C1-C2-C3
3	B	402	AKG	O1-C1-C2-C3
3	D	402	AKG	C1-C2-C3-C4
3	C	402	AKG	C3-C4-C5-O4
3	B	402	AKG	O5-C2-C3-C4
3	C	402	AKG	C3-C4-C5-O3

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	402	AKG	2	0
3	A	402	AKG	1	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

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, 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	325/358 (90%)	-0.00	19 (5%)	29 27	16, 36, 66, 106	5 (1%)
1	B	316/358 (88%)	0.25	17 (5%)	31 30	18, 44, 78, 110	5 (1%)
1	C	326/358 (91%)	0.09	22 (6%)	24 22	14, 38, 70, 119	10 (3%)
1	D	312/358 (87%)	0.11	10 (3%)	50 49	17, 43, 72, 101	4 (1%)
All	All	1279/1432 (89%)	0.11	68 (5%)	32 31	14, 40, 72, 119	24 (1%)

All (68) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	169	LEU	6.0
1	A	35	VAL	5.3
1	B	172	ALA	5.3
1	A	163	ALA	5.3
1	B	35	VAL	5.0
1	C	119[A]	HIS	4.9
1	B	173	VAL	4.8
1	C	170	GLY	4.7
1	C	118[A]	VAL	4.7
1	A	36	SER	4.6
1	A	169	LEU	4.4
1	C	163	ALA	4.3
1	B	33	PRO	4.3
1	C	164	HIS	4.3
1	C	107	VAL	4.2
1	C	165	SER	4.2
1	D	318	ILE	4.2
1	A	164	HIS	4.2
1	A	118	VAL	4.0
1	A	107	VAL	4.0
1	A	170	GLY	3.9

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Mol	Chain	Res	Type	RSRZ
1	A	168	SER	3.8
1	C	35	VAL	3.7
1	C	33	PRO	3.7
1	B	323	PHE	3.7
1	B	175	LEU	3.6
1	A	149[A]	HIS	3.6
1	B	170	GLY	3.5
1	B	318	ILE	3.4
1	C	36	SER	3.4
1	C	162	SER	3.3
1	D	33	PRO	3.2
1	B	271	LEU	3.2
1	A	162	SER	3.2
1	C	173	VAL	3.0
1	B	36	SER	3.0
1	D	35	VAL	3.0
1	C	37	ALA	2.9
1	D	159	SER	2.9
1	A	37	ALA	2.9
1	B	51[A]	ARG	2.7
1	A	165	SER	2.7
1	D	323	PHE	2.7
1	A	117	ASN	2.7
1	A	119	HIS	2.7
1	C	167	THR	2.6
1	C	166	GLN	2.6
1	D	36	SER	2.6
1	C	108[A]	ARG	2.6
1	B	171	SER	2.5
1	D	175	LEU	2.5
1	B	159	SER	2.4
1	A	167	THR	2.4
1	C	115	THR	2.4
1	D	94	MET	2.4
1	A	34	GLU	2.3
1	C	117	ASN	2.3
1	B	322	ARG	2.2
1	B	94	MET	2.2
1	B	231	THR	2.2
1	A	166	GLN	2.2
1	D	177	LEU	2.2
1	C	168	SER	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	171	SER	2.1
1	D	37	ALA	2.1
1	B	34	GLU	2.1
1	C	251	GLY	2.1
1	A	223	MET	2.0

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 oligosaccharides 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($\text{\AA}^2$)	Q<0.9
3	AKG	D	402	10/10	0.84	0.15	59,74,82,82	0
3	AKG	B	402	10/10	0.85	0.13	55,68,73,76	0
3	AKG	C	402	10/10	0.89	0.13	44,63,68,71	0
3	AKG	A	402	10/10	0.89	0.13	41,61,68,71	0
2	FE2	B	401	1/1	0.98	0.05	69,69,69,69	0
2	FE2	D	401	1/1	0.99	0.06	67,67,67,67	0
2	FE2	C	401	1/1	1.00	0.03	47,47,47,47	1
2	FE2	A	401	1/1	1.00	0.03	45,45,45,45	1

6.5 Other polymers [i](#)

There are no such residues in this entry.