



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 5, 2026 – 06:14 PM UTC

PDB ID : 2QFY / pdb_00002qfy
Title : Crystal structure of *Saccharomyces cerevesiae* mitochondrial NADP(+)-dependent isocitrate dehydrogenase in complex with α -ketoglutarate
Authors : Peng, Y.J.; Ding, J.P.
Deposited on : 2007-06-28
Resolution : 2.10 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 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

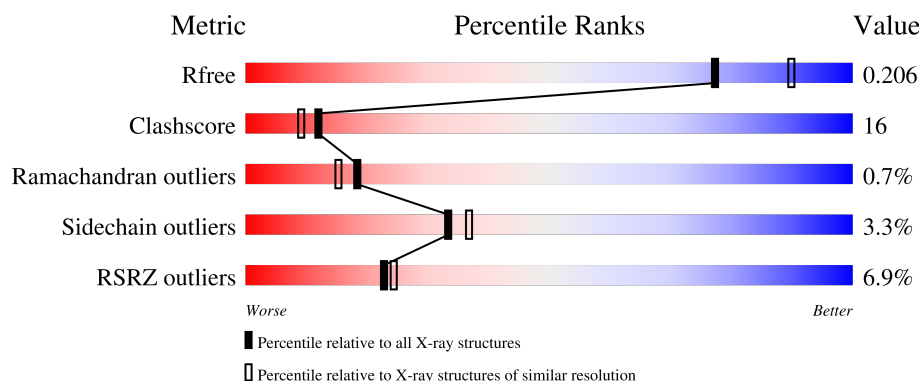
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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	427	2% 72% 22% • •
1	B	427	4% 72% 22% • •
1	C	427	10% 64% 30% • •
1	D	427	9% 65% 29% • •
1	E	427	9% 67% 28% • •

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Mol	Chain	Length	Quality of chain
1	F	427	6% 65% 29%

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 20960 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 Isocitrate dehydrogenase [NADP].

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	411	Total	C	N	O	S	0	0	0
			3256	2076	551	617	12			
1	B	411	Total	C	N	O	S	0	0	0
			3253	2073	551	617	12			
1	C	411	Total	C	N	O	S	0	0	0
			3253	2073	551	617	12			
1	D	411	Total	C	N	O	S	0	0	0
			3256	2076	551	617	12			
1	E	411	Total	C	N	O	S	0	0	0
			3253	2073	551	617	12			
1	F	411	Total	C	N	O	S	0	0	0
			3256	2076	551	617	12			

There are 84 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-13	MET	-	expression tag	UNP P21954
A	-12	HIS	-	expression tag	UNP P21954
A	-11	HIS	-	expression tag	UNP P21954
A	-10	HIS	-	expression tag	UNP P21954
A	-9	HIS	-	expression tag	UNP P21954
A	-8	HIS	-	expression tag	UNP P21954
A	-7	HIS	-	expression tag	UNP P21954
A	-6	ALA	-	expression tag	UNP P21954
A	-5	MET	-	expression tag	UNP P21954
A	-4	GLY	-	expression tag	UNP P21954
A	-3	ILE	-	expression tag	UNP P21954
A	-2	PRO	-	expression tag	UNP P21954
A	-1	GLY	-	expression tag	UNP P21954
A	0	HIS	-	expression tag	UNP P21954
B	-13	MET	-	expression tag	UNP P21954
B	-12	HIS	-	expression tag	UNP P21954
B	-11	HIS	-	expression tag	UNP P21954

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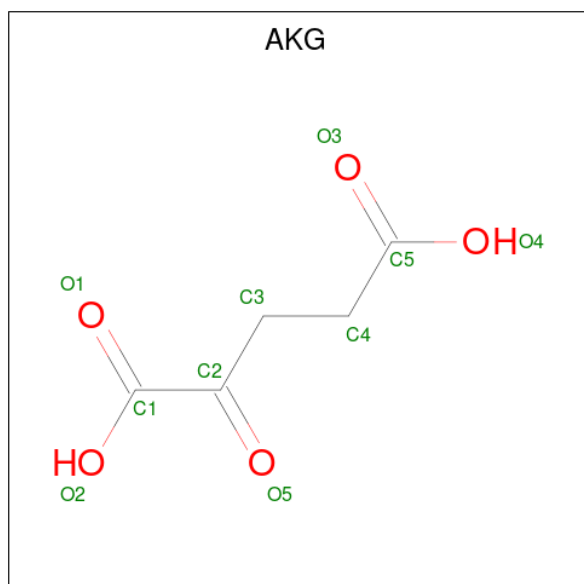
Chain	Residue	Modelled	Actual	Comment	Reference
B	-10	HIS	-	expression tag	UNP P21954
B	-9	HIS	-	expression tag	UNP P21954
B	-8	HIS	-	expression tag	UNP P21954
B	-7	HIS	-	expression tag	UNP P21954
B	-6	ALA	-	expression tag	UNP P21954
B	-5	MET	-	expression tag	UNP P21954
B	-4	GLY	-	expression tag	UNP P21954
B	-3	ILE	-	expression tag	UNP P21954
B	-2	PRO	-	expression tag	UNP P21954
B	-1	GLY	-	expression tag	UNP P21954
B	0	HIS	-	expression tag	UNP P21954
C	-13	MET	-	expression tag	UNP P21954
C	-12	HIS	-	expression tag	UNP P21954
C	-11	HIS	-	expression tag	UNP P21954
C	-10	HIS	-	expression tag	UNP P21954
C	-9	HIS	-	expression tag	UNP P21954
C	-8	HIS	-	expression tag	UNP P21954
C	-7	HIS	-	expression tag	UNP P21954
C	-6	ALA	-	expression tag	UNP P21954
C	-5	MET	-	expression tag	UNP P21954
C	-4	GLY	-	expression tag	UNP P21954
C	-3	ILE	-	expression tag	UNP P21954
C	-2	PRO	-	expression tag	UNP P21954
C	-1	GLY	-	expression tag	UNP P21954
C	0	HIS	-	expression tag	UNP P21954
D	-13	MET	-	expression tag	UNP P21954
D	-12	HIS	-	expression tag	UNP P21954
D	-11	HIS	-	expression tag	UNP P21954
D	-10	HIS	-	expression tag	UNP P21954
D	-9	HIS	-	expression tag	UNP P21954
D	-8	HIS	-	expression tag	UNP P21954
D	-7	HIS	-	expression tag	UNP P21954
D	-6	ALA	-	expression tag	UNP P21954
D	-5	MET	-	expression tag	UNP P21954
D	-4	GLY	-	expression tag	UNP P21954
D	-3	ILE	-	expression tag	UNP P21954
D	-2	PRO	-	expression tag	UNP P21954
D	-1	GLY	-	expression tag	UNP P21954
D	0	HIS	-	expression tag	UNP P21954
E	-13	MET	-	expression tag	UNP P21954
E	-12	HIS	-	expression tag	UNP P21954
E	-11	HIS	-	expression tag	UNP P21954

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Chain	Residue	Modelled	Actual	Comment	Reference
E	-10	HIS	-	expression tag	UNP P21954
E	-9	HIS	-	expression tag	UNP P21954
E	-8	HIS	-	expression tag	UNP P21954
E	-7	HIS	-	expression tag	UNP P21954
E	-6	ALA	-	expression tag	UNP P21954
E	-5	MET	-	expression tag	UNP P21954
E	-4	GLY	-	expression tag	UNP P21954
E	-3	ILE	-	expression tag	UNP P21954
E	-2	PRO	-	expression tag	UNP P21954
E	-1	GLY	-	expression tag	UNP P21954
E	0	HIS	-	expression tag	UNP P21954
F	-13	MET	-	expression tag	UNP P21954
F	-12	HIS	-	expression tag	UNP P21954
F	-11	HIS	-	expression tag	UNP P21954
F	-10	HIS	-	expression tag	UNP P21954
F	-9	HIS	-	expression tag	UNP P21954
F	-8	HIS	-	expression tag	UNP P21954
F	-7	HIS	-	expression tag	UNP P21954
F	-6	ALA	-	expression tag	UNP P21954
F	-5	MET	-	expression tag	UNP P21954
F	-4	GLY	-	expression tag	UNP P21954
F	-3	ILE	-	expression tag	UNP P21954
F	-2	PRO	-	expression tag	UNP P21954
F	-1	GLY	-	expression tag	UNP P21954
F	0	HIS	-	expression tag	UNP P21954

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



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

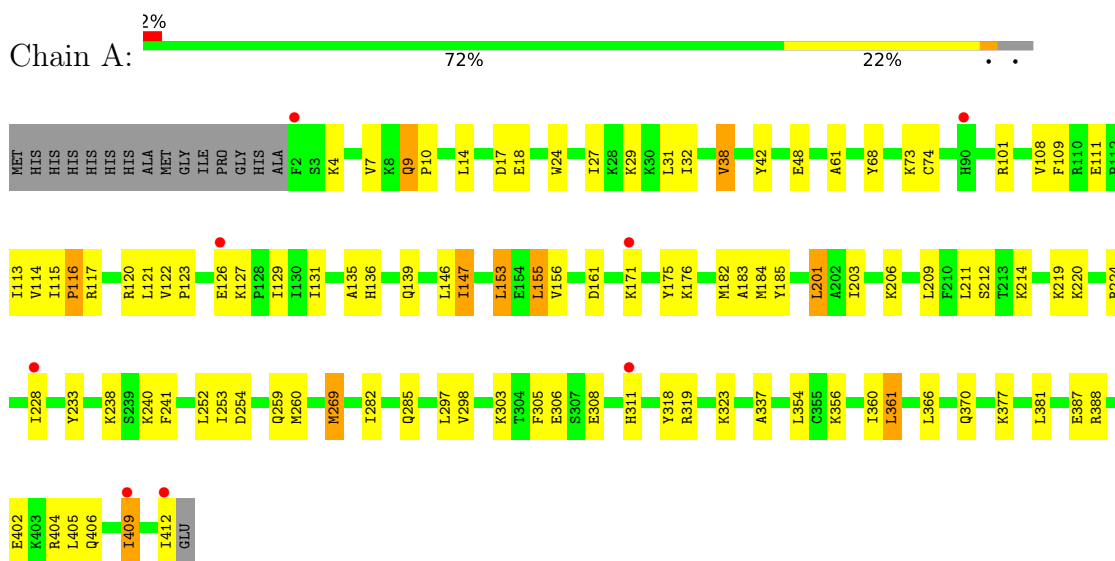
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	392	Total	O	0	0
			392	392		
3	B	328	Total	O	0	0
			328	328		
3	C	167	Total	O	0	0
			167	167		
3	D	151	Total	O	0	0
			151	151		
3	E	130	Total	O	0	0
			130	130		
3	F	205	Total	O	0	0
			205	205		

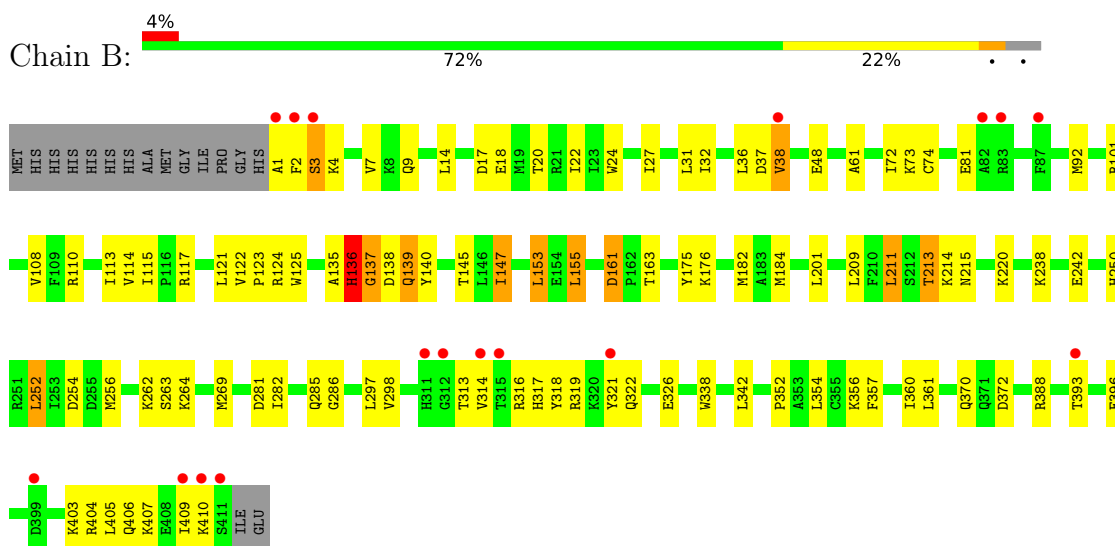
3 Residue-property plots

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: Isocitrate dehydrogenase [NADP]

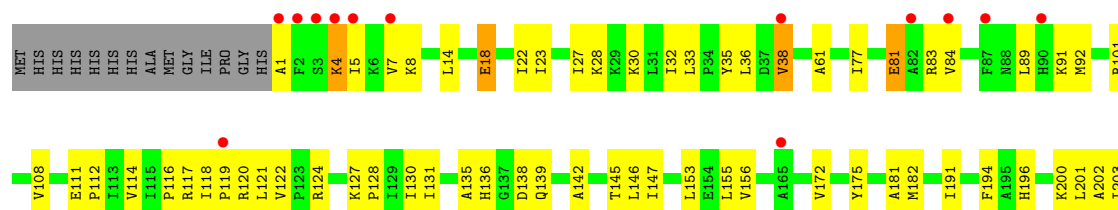


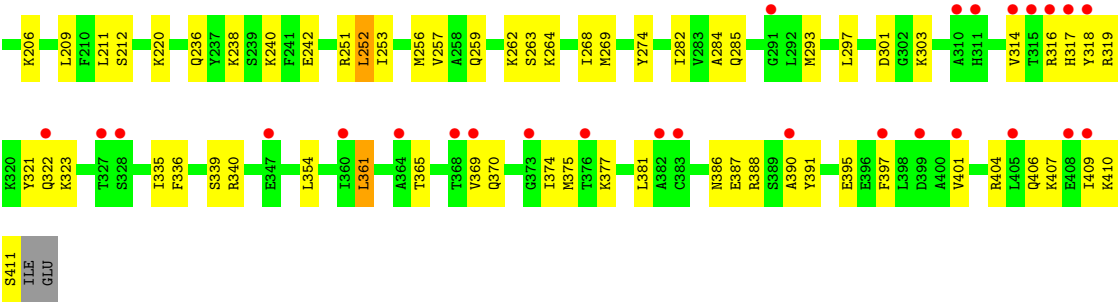
• Molecule 1: Isocitrate dehydrogenase [NADP]



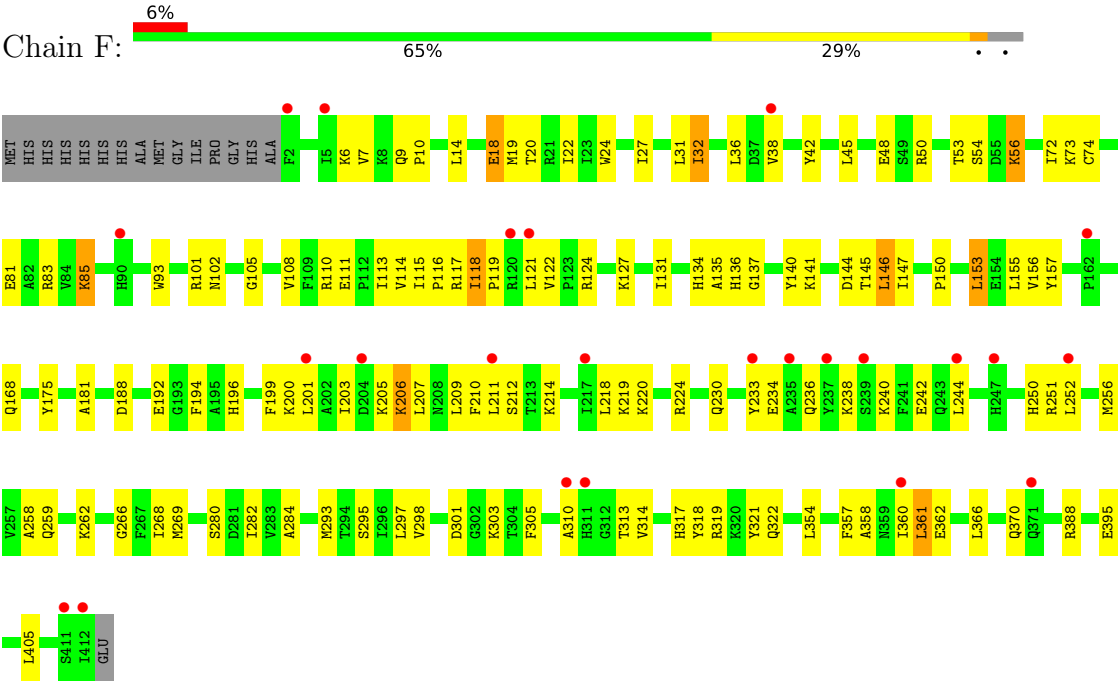
• Molecule 1: Isocitrate dehydrogenase [NADP]







● Molecule 1: Isocitrate dehydrogenase [NADP]



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	74.90Å 98.48Å 190.60Å 90.00° 98.40° 90.00°	Depositor
Resolution (Å)	43.65 – 2.10 43.65 – 2.10	Depositor EDS
% Data completeness (in resolution range)	99.2 (43.65-2.10) 99.2 (43.65-2.10)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.72 (at 2.08Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.206 , 0.247 0.208 , 0.206	Depositor DCC
R_{free} test set	8055 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å ²)	38.4	Xtriage
Anisotropy	0.262	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 48.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.013 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	20960	wwPDB-VP
Average B, all atoms (Å ²)	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.93% 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

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	0.43	1/3319 (0.0%)	0.94	15/4477 (0.3%)
1	B	0.41	0/3316	1.03	22/4473 (0.5%)
1	C	0.35	0/3316	0.90	7/4473 (0.2%)
1	D	0.34	0/3319	0.87	11/4477 (0.2%)
1	E	0.33	0/3316	0.88	9/4473 (0.2%)
1	F	0.35	0/3319	0.90	13/4477 (0.3%)
All	All	0.37	1/19905 (0.0%)	0.92	77/26850 (0.3%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	269	MET	SD-CE	-5.30	1.66	1.79

The worst 5 of 77 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	3	SER	N-CA-C	-26.45	68.87	109.86
1	B	3	SER	CB-CA-C	7.64	131.33	113.15
1	B	115	ILE	N-CA-C	-7.57	100.66	108.15
1	E	147	ILE	CA-C-N	7.49	127.03	119.24
1	E	147	ILE	C-N-CA	7.49	127.03	119.24

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	3256	0	3302	76	0
1	B	3253	0	3299	97	0
1	C	3253	0	3299	131	0
1	D	3256	0	3302	127	0
1	E	3253	0	3299	119	0
1	F	3256	0	3302	108	0
2	A	10	0	4	1	0
2	B	10	0	4	3	0
2	C	10	0	4	1	0
2	D	10	0	4	3	0
2	E	10	0	4	1	0
2	F	10	0	4	2	0
3	A	392	0	0	13	0
3	B	328	0	0	10	0
3	C	167	0	0	9	0
3	D	151	0	0	4	0
3	E	130	0	0	1	0
3	F	205	0	0	8	0
All	All	20960	0	19827	617	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 16.

The worst 5 of 617 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:182:MET:HE1	1:D:218:LEU:HD13	1.42	1.00
1:D:117:ARG:H	1:D:370:GLN:HE22	1.00	0.98
1:A:7:VAL:HG21	1:A:38:VAL:HG13	1.46	0.96
1:D:79:PRO:HG3	1:D:93:TRP:HB2	1.51	0.92
1:C:117:ARG:HD2	1:C:370:GLN:HG2	1.51	0.92

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	409/427 (96%)	395 (97%)	14 (3%)	0	100	100
1	B	409/427 (96%)	392 (96%)	15 (4%)	2 (0%)	24	22
1	C	409/427 (96%)	378 (92%)	25 (6%)	6 (2%)	8	4
1	D	409/427 (96%)	384 (94%)	23 (6%)	2 (0%)	24	22
1	E	409/427 (96%)	382 (93%)	23 (6%)	4 (1%)	12	9
1	F	409/427 (96%)	394 (96%)	13 (3%)	2 (0%)	24	22
All	All	2454/2562 (96%)	2325 (95%)	113 (5%)	16 (1%)	18	15

5 of 16 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	137	GLY
1	E	4	LYS
1	C	311	HIS
1	D	18	GLU
1	D	311	HIS

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	355/367 (97%)	339 (96%)	16 (4%)	24	25
1	B	354/367 (96%)	341 (96%)	13 (4%)	30	33
1	C	354/367 (96%)	343 (97%)	11 (3%)	35	39

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	D	355/367 (97%)	344 (97%)	11 (3%)	35	39
1	E	354/367 (96%)	347 (98%)	7 (2%)	48	56
1	F	355/367 (97%)	343 (97%)	12 (3%)	32	35
All	All	2127/2202 (97%)	2057 (97%)	70 (3%)	33	37

5 of 70 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	361	LEU
1	F	56	LYS
1	F	206	LYS
1	B	213	THR
1	B	211	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 49 such sidechains are listed below:

Mol	Chain	Res	Type
1	D	371	GLN
1	E	367	ASN
1	E	166	GLN
1	E	259	GLN
1	F	88	ASN

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 oligosaccharides in this entry.

5.6 Ligand geometry

6 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
2	AKG	B	2002	-	9,9,9	1.28	1 (11%)	11,11,11	1.50	2 (18%)
2	AKG	E	2005	-	9,9,9	1.26	1 (11%)	11,11,11	1.51	2 (18%)
2	AKG	C	2003	-	9,9,9	1.28	1 (11%)	11,11,11	1.48	2 (18%)
2	AKG	D	2004	-	9,9,9	1.27	1 (11%)	11,11,11	1.50	2 (18%)
2	AKG	F	2006	-	9,9,9	1.27	1 (11%)	11,11,11	1.51	2 (18%)
2	AKG	A	2001	-	9,9,9	1.29	1 (11%)	11,11,11	1.50	2 (18%)

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	AKG	B	2002	-	-	4/9/9/9	-
2	AKG	E	2005	-	-	1/9/9/9	-
2	AKG	C	2003	-	-	2/9/9/9	-
2	AKG	D	2004	-	-	1/9/9/9	-
2	AKG	F	2006	-	-	3/9/9/9	-
2	AKG	A	2001	-	-	3/9/9/9	-

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

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	2001	AKG	C3-C2	2.18	1.53	1.51
2	C	2003	AKG	C3-C2	2.15	1.53	1.51
2	D	2004	AKG	C3-C2	2.09	1.53	1.51
2	B	2002	AKG	C3-C2	2.08	1.53	1.51
2	F	2006	AKG	C3-C2	2.07	1.53	1.51

The worst 5 of 12 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	2006	AKG	O1-C1-C2	-3.08	117.98	121.81
2	D	2004	AKG	O1-C1-C2	-3.06	118.01	121.81
2	A	2001	AKG	O1-C1-C2	-3.04	118.03	121.81
2	C	2003	AKG	O1-C1-C2	-3.04	118.03	121.81
2	E	2005	AKG	O1-C1-C2	-3.03	118.05	121.81

There are no chirality outliers.

5 of 14 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	2001	AKG	C1-C2-C3-C4
2	B	2002	AKG	C1-C2-C3-C4
2	C	2003	AKG	C1-C2-C3-C4
2	D	2004	AKG	C1-C2-C3-C4
2	E	2005	AKG	C1-C2-C3-C4

There are no ring outliers.

6 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	2002	AKG	3	0
2	E	2005	AKG	1	0
2	C	2003	AKG	1	0
2	D	2004	AKG	3	0
2	F	2006	AKG	2	0
2	A	2001	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 [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	411/427 (96%)	0.07	8 (1%) 66 69	23, 34, 55, 89	0
1	B	411/427 (96%)	0.24	17 (4%) 41 43	21, 36, 65, 103	0
1	C	411/427 (96%)	0.78	43 (10%) 11 12	31, 48, 86, 103	0
1	D	411/427 (96%)	0.75	37 (9%) 15 16	34, 49, 89, 103	0
1	E	411/427 (96%)	0.93	40 (9%) 13 14	36, 53, 101, 128	0
1	F	411/427 (96%)	0.59	24 (5%) 29 30	31, 46, 73, 108	0
All	All	2466/2562 (96%)	0.56	169 (6%) 23 24	21, 45, 83, 128	0

The worst 5 of 169 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	2	PHE	5.7
1	C	291	GLY	5.5
1	F	412	ILE	5.3
1	E	1	ALA	5.3
1	C	313	THR	4.5

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
2	AKG	C	2003	10/10	0.65	0.23	75,77,80,80	10
2	AKG	D	2004	10/10	0.66	0.18	87,88,91,91	0
2	AKG	B	2002	10/10	0.68	0.30	76,77,80,80	10
2	AKG	E	2005	10/10	0.68	0.23	83,84,87,87	10
2	AKG	F	2006	10/10	0.75	0.25	80,82,85,87	9
2	AKG	A	2001	10/10	0.76	0.20	72,74,76,77	0

6.5 Other polymers [i](#)

There are no such residues in this entry.