



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 05:08 AM UTC

PDB ID : 2XHK / pdb_00002xhk
Title : Crystal structure of transcription factor NtcA from *Synechococcus elongatus* bound to 2-oxoglutarate
Authors : Llacer, J.L.; Castells, M.A.; Rubio, V.
Deposited on : 2010-06-17
Resolution : 2.30 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

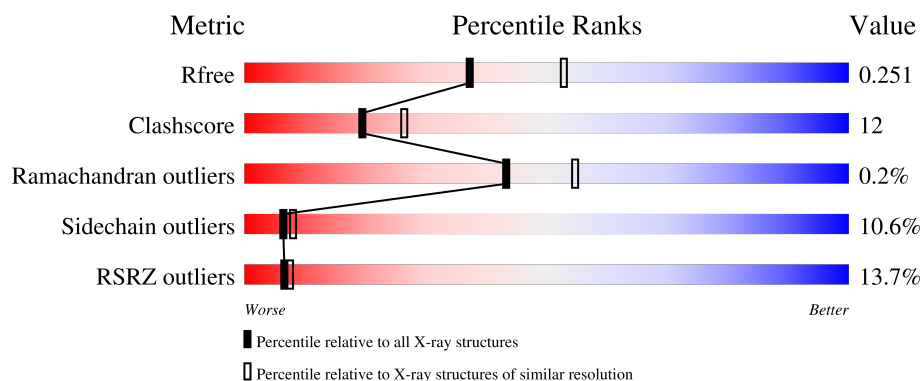
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	222	14% 71% 22% • •
1	B	222	13% 77% 15% 5% ••

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
2	AKG	A	519	-	X	X	-
2	AKG	B	519	-	X	-	-

2 Entry composition [i](#)

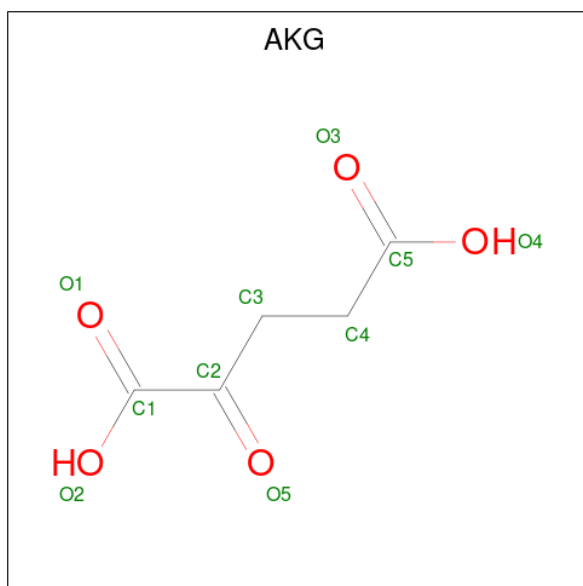
There are 3 unique types of molecules in this entry. The entry contains 3550 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 GLOBAL NITROGEN REGULATOR.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	215	Total	C	N	O	S	0	1	0
			1698	1092	293	305	8			
1	B	217	Total	C	N	O	S	0	2	0
			1716	1101	299	308	8			

- 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		

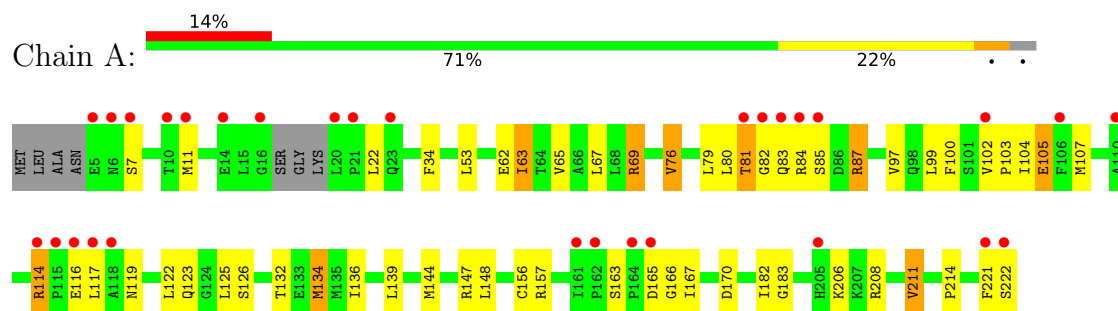
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	50	Total 50	O 50	0	0
3	B	66	Total 66	O 66	0	0

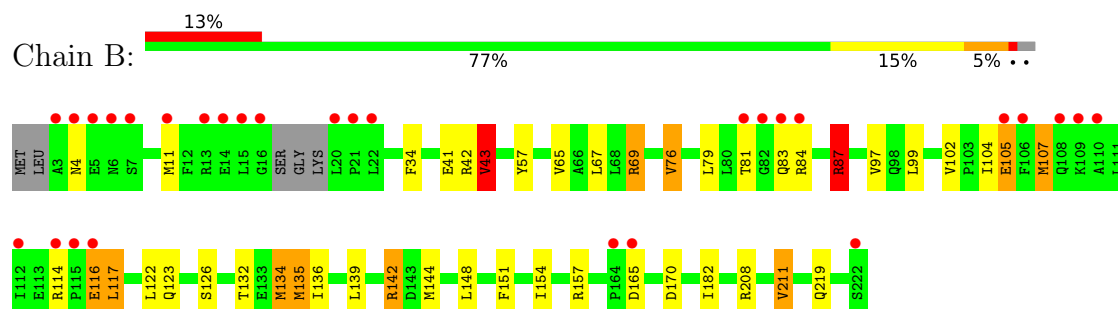
3 Residue-property plots [i](#)

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: GLOBAL NITROGEN REGULATOR



• Molecule 1: GLOBAL NITROGEN REGULATOR



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	68.22Å 70.81Å 149.96Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.30 50.00 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.8 (50.00-2.30) 99.8 (50.00-2.30)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.29 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.5.0072	Depositor
R, R_{free}	0.204 , 0.238 0.227 , 0.251	Depositor DCC
R_{free} test set	1673 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	41.8	Xtriage
Anisotropy	0.483	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 45.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.023 for k,h,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	3550	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.04% 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	0.84	0/1728	0.91	0/2331
1	B	0.90	1/1749 (0.1%)	0.99	4/2359 (0.2%)
All	All	0.87	1/3477 (0.0%)	0.95	4/4690 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	211	VAL	CA-CB	5.01	1.60	1.54

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	43	VAL	CB-CA-C	-6.91	100.88	111.31
1	B	135	MET	CG-SD-CE	-5.64	88.50	100.90
1	B	87	ARG	CB-CG-CD	5.42	123.76	111.30
1	B	87	ARG	CD-NE-CZ	5.14	131.60	124.40

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	1698	0	1771	47	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	1716	0	1786	41	0
2	A	10	0	4	4	0
2	B	10	0	4	1	0
3	A	50	0	0	4	0
3	B	66	0	0	3	0
All	All	3550	0	3565	81	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 12.

All (81) 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:B:11:MET:SD	1:B:117:LEU:HD11	1.66	1.35
1:A:170:ASP:OD1	1:A:208:ARG:NH2	1.79	1.15
1:B:11:MET:SD	1:B:117:LEU:CD1	2.37	1.12
1:A:81:THR:HG22	1:A:82:GLY:HA2	1.36	1.06
2:A:519:AKG:O2	3:A:2050:HOH:O	1.89	0.89
1:B:11:MET:CG	1:B:117:LEU:HD11	2.08	0.82
2:B:519:AKG:O2	3:B:2066:HOH:O	1.97	0.82
1:A:83:GLN:OE1	1:A:85:SER:OG	2.01	0.75
1:A:144:MET:HE1	1:A:182:ILE:HD12	1.71	0.73
1:B:34:PHE:CD1	1:B:87:ARG:HG2	2.26	0.70
1:A:79:LEU:HD23	1:A:104:ILE:HD11	1.71	0.70
1:B:144:MET:CE	1:B:182:ILE:HD12	2.22	0.69
1:A:34:PHE:CD1	1:A:87:ARG:HG2	2.29	0.68
1:B:11:MET:SD	1:B:117:LEU:HD12	2.34	0.67
1:A:122:LEU:HD23	1:B:122:LEU:HD23	1.80	0.64
1:A:114:ARG:HB3	1:A:117:LEU:HD23	1.80	0.63
1:A:222:SER:HB2	1:B:57:TYR:HD1	1.63	0.63
1:B:81:THR:O	1:B:84:ARG:HB3	1.98	0.63
1:A:81:THR:CG2	1:A:82:GLY:HA2	2.20	0.62
1:A:80:LEU:O	1:A:83:GLN:NE2	2.31	0.61
1:A:53:LEU:HD13	2:A:519:AKG:H41	1.83	0.61
1:B:69:ARG:HD3	3:B:2021:HOH:O	2.00	0.61
1:A:53:LEU:CD1	2:A:519:AKG:H41	2.30	0.61
1:B:41:GLU:HB3	1:B:83:GLN:HG2	1.82	0.61
1:B:105:GLU:H	1:B:105:GLU:CD	2.10	0.60
1:A:134[B]:MET:SD	1:A:157:ARG:NH2	2.75	0.59
1:A:99:LEU:C	1:A:99:LEU:HD12	2.29	0.58
1:A:136:ILE:HD11	1:B:132:THR:HG22	1.87	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:125:LEU:HD22	1:B:122:LEU:HD22	1.86	0.57
1:A:22:LEU:HD23	1:A:100:PHE:CD1	2.40	0.56
1:A:183:GLY:O	1:B:142:ARG:HD3	2.06	0.56
1:B:11:MET:CE	1:B:117:LEU:HD12	2.36	0.56
1:B:134[A]:MET:HE3	1:B:154:ILE:HD11	1.87	0.55
1:A:156:CYS:SG	1:A:211:VAL:HG13	2.46	0.55
1:B:41:GLU:CB	1:B:83:GLN:HG2	2.37	0.55
1:A:69:ARG:HD3	3:A:2013:HOH:O	2.07	0.54
1:B:4:ASN:CG	1:B:123:GLN:HE21	2.16	0.54
1:A:11:MET:HG3	1:A:117:LEU:HD12	1.91	0.53
1:B:135:MET:HE1	1:B:151:PHE:CD1	2.44	0.53
1:B:144:MET:HE2	1:B:182:ILE:HD12	1.91	0.52
1:A:63:ILE:CD1	3:A:2006:HOH:O	2.58	0.52
1:A:11:MET:SD	1:A:117:LEU:HD11	2.50	0.51
1:B:11:MET:HE2	1:B:107:MET:HE1	1.93	0.51
1:A:11:MET:HG3	1:A:117:LEU:CD1	2.41	0.51
1:A:119:ASN:HD21	1:A:123:GLN:HE21	1.60	0.50
1:A:167:ILE:HG12	1:A:214:PRO:HG3	1.94	0.50
1:B:11:MET:CE	1:B:117:LEU:CD1	2.90	0.49
1:A:144:MET:CE	1:A:182:ILE:HD12	2.42	0.49
1:B:134[B]:MET:SD	1:B:157:ARG:NH2	2.85	0.49
1:B:11:MET:CE	1:B:107:MET:HE1	2.42	0.49
1:A:166:GLY:C	1:A:167:ILE:HD13	2.38	0.48
1:A:144:MET:HE1	1:A:182:ILE:CD1	2.42	0.48
1:B:117:LEU:HD13	1:B:117:LEU:O	2.14	0.48
1:A:114:ARG:HB3	1:A:117:LEU:CD2	2.44	0.47
1:A:114:ARG:CB	1:A:117:LEU:HD23	2.44	0.47
1:B:41:GLU:HG3	1:B:42:ARG:HG3	1.97	0.47
1:A:126:SER:HB3	1:B:76:VAL:HG13	1.97	0.47
1:B:43:VAL:HG21	1:B:107:MET:HG3	1.97	0.46
1:A:63:ILE:HD11	3:A:2006:HOH:O	2.15	0.46
1:B:170:ASP:OD1	1:B:208:ARG:NH1	2.33	0.45
1:B:11:MET:HG3	1:B:117:LEU:HD11	1.94	0.44
1:A:63:ILE:HD12	1:A:63:ILE:HA	1.76	0.44
1:B:117:LEU:CD1	1:B:117:LEU:O	2.65	0.44
1:A:119:ASN:ND2	1:A:123:GLN:HE21	2.15	0.43
1:A:99:LEU:HD12	1:A:99:LEU:O	2.18	0.43
1:A:105:GLU:H	1:A:105:GLU:CD	2.27	0.43
1:A:132:THR:HG22	1:B:136:ILE:HD11	2.01	0.43
1:B:79:LEU:HD23	1:B:104:ILE:HD11	2.01	0.43
1:B:144:MET:HE1	1:B:182:ILE:HD12	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:62:GLU:O	1:A:63:ILE:HD13	2.19	0.42
1:A:166:GLY:O	1:A:167:ILE:HD13	2.20	0.42
1:B:99:LEU:HD12	1:B:99:LEU:C	2.44	0.42
1:A:53:LEU:HD11	2:A:519:AKG:H41	2.01	0.42
1:A:221:PHE:O	1:A:222:SER:C	2.62	0.42
1:A:76:VAL:HG13	1:B:126:SER:HB3	2.01	0.42
1:B:4:ASN:HB3	3:B:2001:HOH:O	2.19	0.42
1:A:132:THR:CG2	1:B:136:ILE:HD11	2.50	0.41
1:B:116:GLU:CD	1:B:116:GLU:C	2.89	0.41
1:A:102:VAL:HG12	1:A:103:PRO:O	2.21	0.41
1:A:84:ARG:CZ	1:A:84:ARG:HA	2.51	0.41
1:B:144:MET:HE3	1:B:144:MET:HA	2.03	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	212/222 (96%)	206 (97%)	5 (2%)	1 (0%)	24	31
1	B	215/222 (97%)	212 (99%)	3 (1%)	0	100	100
All	All	427/444 (96%)	418 (98%)	8 (2%)	1 (0%)	43	55

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	7	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	189/194 (97%)	168 (89%)	21 (11%)	6	7
1	B	190/194 (98%)	169 (89%)	21 (11%)	6	7
All	All	379/388 (98%)	337 (89%)	42 (11%)	6	7

All (42) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	63	ILE
1	A	65	VAL
1	A	67	LEU
1	A	69	ARG
1	A	76	VAL
1	A	81	THR
1	A	87	ARG
1	A	97	VAL
1	A	105	GLU
1	A	107	MET
1	A	114	ARG
1	A	116	GLU
1	A	134[A]	MET
1	A	134[B]	MET
1	A	139	LEU
1	A	147	ARG
1	A	148	LEU
1	A	163	SER
1	A	165	ASP
1	A	206	LYS
1	A	211	VAL
1	B	43	VAL
1	B	65	VAL
1	B	67	LEU
1	B	69	ARG
1	B	76	VAL
1	B	87	ARG

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Mol	Chain	Res	Type
1	B	97	VAL
1	B	102	VAL
1	B	105	GLU
1	B	107	MET
1	B	114	ARG
1	B	116	GLU
1	B	117	LEU
1	B	134[A]	MET
1	B	134[B]	MET
1	B	139	LEU
1	B	142	ARG
1	B	148	LEU
1	B	165	ASP
1	B	211	VAL
1	B	219	GLN

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	26	GLN
1	A	71	ASN
1	A	108	GLN
1	A	119	ASN
1	A	141	HIS
1	A	220	GLN
1	B	71	ASN
1	B	108	GLN
1	B	123	GLN
1	B	176	GLN
1	B	219	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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

2 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	A	519	-	9,9,9	3.17	4 (44%)	11,11,11	2.62	5 (45%)
2	AKG	B	519	-	9,9,9	3.62	4 (44%)	11,11,11	2.44	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
2	AKG	A	519	-	-	5/9/9/9	-
2	AKG	B	519	-	-	5/9/9/9	-

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	519	AKG	C2-C1	-8.59	1.40	1.53
2	A	519	AKG	C2-C1	-6.78	1.43	1.53
2	B	519	AKG	C3-C2	5.31	1.57	1.51
2	A	519	AKG	C3-C2	4.59	1.56	1.51
2	A	519	AKG	C4-C5	3.28	1.58	1.50
2	A	519	AKG	O2-C1	-2.68	1.23	1.30
2	B	519	AKG	O2-C1	-2.66	1.23	1.30
2	B	519	AKG	O4-C5	-2.15	1.23	1.30

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	519	AKG	O5-C2-C1	-4.99	112.53	119.67
2	A	519	AKG	C3-C2-C1	4.50	123.50	115.86
2	B	519	AKG	C4-C3-C2	-4.20	105.02	112.91
2	B	519	AKG	O5-C2-C1	-3.86	114.16	119.67
2	B	519	AKG	O1-C1-C2	-3.48	117.48	121.81
2	A	519	AKG	O1-C1-C2	-3.18	117.85	121.81
2	B	519	AKG	C3-C2-C1	2.82	120.64	115.86
2	A	519	AKG	C4-C3-C2	-2.73	107.78	112.91
2	A	519	AKG	O4-C5-C4	2.28	121.20	114.00
2	B	519	AKG	C3-C4-C5	-2.16	107.94	113.67

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	519	AKG	C1-C2-C3-C4
2	A	519	AKG	C2-C3-C4-C5
2	B	519	AKG	C1-C2-C3-C4
2	B	519	AKG	C2-C3-C4-C5
2	A	519	AKG	C3-C4-C5-O4
2	A	519	AKG	C3-C4-C5-O3
2	B	519	AKG	C3-C4-C5-O4
2	B	519	AKG	C3-C4-C5-O3
2	A	519	AKG	O5-C2-C3-C4
2	B	519	AKG	O5-C2-C3-C4

There are no ring outliers.

2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	519	AKG	4	0
2	B	519	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	215/222 (96%)	0.79	30 (13%) 6 7	20, 42, 84, 98	2 (0%)
1	B	217/222 (97%)	0.67	29 (13%) 7 8	21, 39, 87, 108	3 (1%)
All	All	432/444 (97%)	0.73	59 (13%) 6 7	20, 41, 85, 108	5 (1%)

All (59) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	20	LEU	8.5
1	B	3	ALA	6.3
1	A	222	SER	6.0
1	A	82	GLY	6.0
1	A	20	LEU	5.2
1	B	21	PRO	5.2
1	B	112	ILE	4.9
1	B	116	GLU	4.8
1	B	16	GLY	4.8
1	A	5	GLU	4.7
1	A	205	HIS	4.6
1	A	83	GLN	4.4
1	A	106	PHE	4.4
1	A	16	GLY	4.3
1	B	6	ASN	4.2
1	A	115	PRO	4.2
1	B	4	ASN	4.0
1	B	83	GLN	3.9
1	B	82	GLY	3.8
1	A	6	ASN	3.7
1	B	115	PRO	3.7
1	B	22	LEU	3.6
1	A	7	SER	3.5
1	A	84	ARG	3.5

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Mol	Chain	Res	Type	RSRZ
1	B	106	PHE	3.5
1	A	11	MET	3.4
1	A	117	LEU	3.3
1	A	164	PRO	3.3
1	A	114	ARG	3.2
1	B	5	GLU	3.2
1	B	81	THR	3.1
1	B	7	SER	3.0
1	B	11	MET	2.9
1	A	21	PRO	2.8
1	A	165	ASP	2.8
1	A	116	GLU	2.8
1	B	108	GLN	2.7
1	A	81	THR	2.7
1	B	109	LYS	2.7
1	A	14	GLU	2.6
1	A	161	ILE	2.6
1	B	164	PRO	2.6
1	B	105	GLU	2.5
1	B	15	LEU	2.5
1	A	85	SER	2.5
1	B	114	ARG	2.4
1	A	110	ALA	2.4
1	A	10	THR	2.4
1	B	165	ASP	2.3
1	B	13	ARG	2.3
1	A	118	ALA	2.3
1	B	84	ARG	2.2
1	B	14	GLU	2.2
1	A	23	GLN	2.2
1	A	221	PHE	2.1
1	A	162	PRO	2.0
1	A	102	VAL	2.0
1	B	222	SER	2.0
1	B	110	ALA	2.0

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

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	B	519	10/10	0.81	0.15	34,41,43,53	0
2	AKG	A	519	10/10	0.84	0.13	35,37,49,55	0

6.5 Other polymers [i](#)

There are no such residues in this entry.