



Full wwPDB X-ray Structure Validation Report ⓘ

May 2, 2017 – 10:56 AM EDT

PDB ID : 5X5U
Title : Crystal structure of alpha-ketoglutarate-semialdehyde dehydrogenase (KGSADH) complexed with NAD
Authors : Son, H.-F.; Kim, K.-J.
Deposited on : 2017-02-17
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

<http://wwpdb.org/validation/2016/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 : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20029077
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20029077

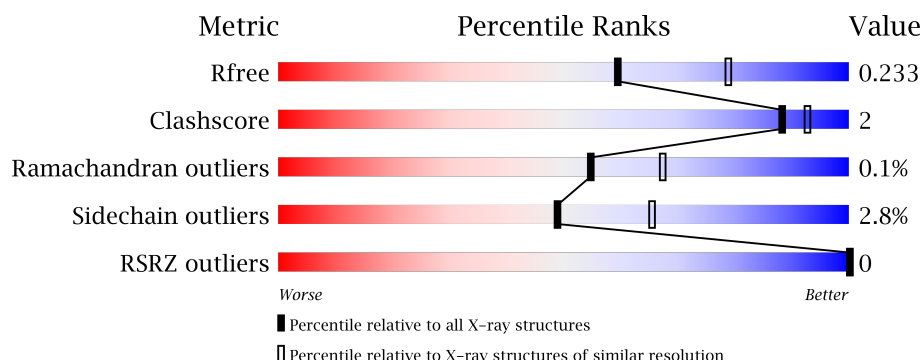
1 Overall quality at a glance

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}	100719	4130 (2.30-2.30)
Clashscore	112137	4751 (2.30-2.30)
Ramachandran outliers	110173	4705 (2.30-2.30)
Sidechain outliers	110143	4704 (2.30-2.30)
RSRZ outliers	101464	4156 (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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	505	 86% 8% • 6%
1	B	505	 87% 7% 6%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7748 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 Alpha-ketoglutaric semialdehyde dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	476	Total	C	N	O	S	0	0	0
			3548	2251	629	658	10			
1	B	476	Total	C	N	O	S	0	0	0
			3548	2251	629	658	10			

There are 50 discrepancies between the modelled and reference sequences:

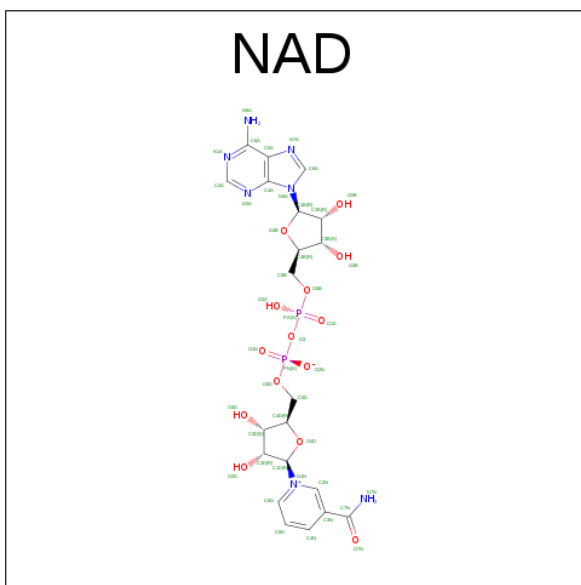
Chain	Residue	Modelled	Actual	Comment	Reference
A	-23	MET	-	expression tag	UNP Q1JUP4
A	-22	SER	-	expression tag	UNP Q1JUP4
A	-21	TYR	-	expression tag	UNP Q1JUP4
A	-20	TYR	-	expression tag	UNP Q1JUP4
A	-19	HIS	-	expression tag	UNP Q1JUP4
A	-18	HIS	-	expression tag	UNP Q1JUP4
A	-17	HIS	-	expression tag	UNP Q1JUP4
A	-16	HIS	-	expression tag	UNP Q1JUP4
A	-15	HIS	-	expression tag	UNP Q1JUP4
A	-14	HIS	-	expression tag	UNP Q1JUP4
A	-13	ASP	-	expression tag	UNP Q1JUP4
A	-12	TYR	-	expression tag	UNP Q1JUP4
A	-11	ASP	-	expression tag	UNP Q1JUP4
A	-10	ILE	-	expression tag	UNP Q1JUP4
A	-9	PRO	-	expression tag	UNP Q1JUP4
A	-8	THR	-	expression tag	UNP Q1JUP4
A	-7	THR	-	expression tag	UNP Q1JUP4
A	-6	GLU	-	expression tag	UNP Q1JUP4
A	-5	ASN	-	expression tag	UNP Q1JUP4
A	-4	ILE	-	expression tag	UNP Q1JUP4
A	-3	TYR	-	expression tag	UNP Q1JUP4
A	-2	PHE	-	expression tag	UNP Q1JUP4
A	-1	GLN	-	expression tag	UNP Q1JUP4
A	0	GLY	-	expression tag	UNP Q1JUP4
A	1	ALA	-	expression tag	UNP Q1JUP4

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-23	MET	-	expression tag	UNP Q1JUP4
B	-22	SER	-	expression tag	UNP Q1JUP4
B	-21	TYR	-	expression tag	UNP Q1JUP4
B	-20	TYR	-	expression tag	UNP Q1JUP4
B	-19	HIS	-	expression tag	UNP Q1JUP4
B	-18	HIS	-	expression tag	UNP Q1JUP4
B	-17	HIS	-	expression tag	UNP Q1JUP4
B	-16	HIS	-	expression tag	UNP Q1JUP4
B	-15	HIS	-	expression tag	UNP Q1JUP4
B	-14	HIS	-	expression tag	UNP Q1JUP4
B	-13	ASP	-	expression tag	UNP Q1JUP4
B	-12	TYR	-	expression tag	UNP Q1JUP4
B	-11	ASP	-	expression tag	UNP Q1JUP4
B	-10	ILE	-	expression tag	UNP Q1JUP4
B	-9	PRO	-	expression tag	UNP Q1JUP4
B	-8	THR	-	expression tag	UNP Q1JUP4
B	-7	THR	-	expression tag	UNP Q1JUP4
B	-6	GLU	-	expression tag	UNP Q1JUP4
B	-5	ASN	-	expression tag	UNP Q1JUP4
B	-4	ILE	-	expression tag	UNP Q1JUP4
B	-3	TYR	-	expression tag	UNP Q1JUP4
B	-2	PHE	-	expression tag	UNP Q1JUP4
B	-1	GLN	-	expression tag	UNP Q1JUP4
B	0	GLY	-	expression tag	UNP Q1JUP4
B	1	ALA	-	expression tag	UNP Q1JUP4

- Molecule 2 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (three-letter code: NAD) (formula: C₂₁H₂₇N₇O₁₄P₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	B	1	Total	C	N	O	P	0	0
			44	21	7	14	2		

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	3	3		

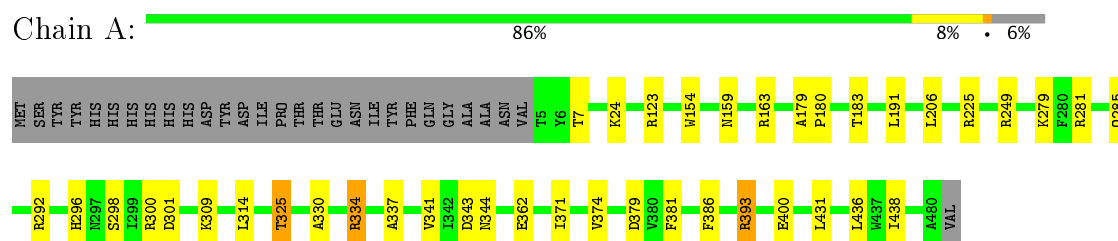
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	284	Total 284	O 284	0	0
4	B	274	Total 274	O 274	0	0

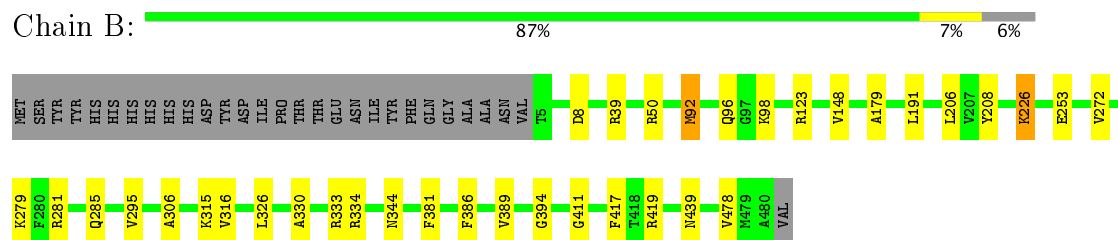
3 Residue-property plots [i](#)

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: Alpha-ketoglutaric semialdehyde dehydrogenase



- Molecule 1: Alpha-ketoglutaric semialdehyde dehydrogenase



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 2 2	Depositor
Cell constants a, b, c, α , β , γ	129.31Å 129.31Å 118.36Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	33.68 – 2.30 33.68 – 2.30	Depositor EDS
% Data completeness (in resolution range)	94.6 (33.68-2.30) 94.6 (33.68-2.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	19.55 (at 2.29Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.162 , 0.227 0.172 , 0.233	Depositor DCC
R_{free} test set	2184 reflections (5.39%)	DCC
Wilson B-factor (Å ²)	22.7	Xtriage
Anisotropy	0.005	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 26.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7748	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.88% 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: GOL, NAD

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	0.88	3/3625 (0.1%)	0.93	13/4938 (0.3%)
1	B	0.87	0/3625	0.92	9/4938 (0.2%)
All	All	0.87	3/7250 (0.0%)	0.93	22/9876 (0.2%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	400	GLU	CD-OE2	5.31	1.31	1.25
1	A	400	GLU	CG-CD	5.12	1.59	1.51
1	A	154	TRP	CB-CG	-5.01	1.41	1.50

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	225	ARG	NE-CZ-NH2	-10.35	115.13	120.30
1	B	92	MET	CG-SD-CE	-9.24	85.42	100.20
1	A	225	ARG	NE-CZ-NH1	9.16	124.88	120.30
1	B	50	ARG	NE-CZ-NH1	8.01	124.31	120.30
1	A	123	ARG	NE-CZ-NH2	-7.53	116.54	120.30
1	A	281	ARG	NE-CZ-NH1	7.02	123.81	120.30
1	A	123	ARG	NE-CZ-NH1	6.75	123.67	120.30
1	A	334	ARG	NE-CZ-NH1	6.20	123.40	120.30
1	A	163	ARG	NE-CZ-NH1	6.14	123.37	120.30
1	A	163	ARG	NE-CZ-NH2	-6.00	117.30	120.30
1	B	123	ARG	NE-CZ-NH1	5.64	123.12	120.30
1	B	281	ARG	NE-CZ-NH1	5.56	123.08	120.30
1	A	225	ARG	CG-CD-NE	-5.50	100.25	111.80
1	B	281	ARG	NE-CZ-NH2	-5.46	117.57	120.30
1	A	393	ARG	NE-CZ-NH2	-5.45	117.57	120.30
1	A	292	ARG	NE-CZ-NH1	5.35	122.97	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	50	ARG	NE-CZ-NH2	-5.25	117.67	120.30
1	A	249	ARG	NE-CZ-NH2	-5.24	117.68	120.30
1	B	419	ARG	NE-CZ-NH2	-5.22	117.69	120.30
1	B	419	ARG	NE-CZ-NH1	5.20	122.90	120.30
1	B	8	ASP	CB-CG-OD1	5.12	122.90	118.30
1	A	400	GLU	CA-CB-CG	5.08	124.58	113.40

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	3548	0	3571	16	0
1	B	3548	0	3571	16	0
2	A	44	0	26	0	0
2	B	44	0	26	3	0
3	A	6	0	8	0	0
4	A	284	0	0	1	0
4	B	274	0	0	1	2
All	All	7748	0	7202	33	2

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 (33) 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)
2:B:501:NAD:H51N	2:B:501:NAD:H6N	1.60	0.81
1:A:7:THR:HG23	4:A:771:HOH:O	1.84	0.77
1:B:333:ARG:NH1	2:B:501:NAD:O2N	2.19	0.74
1:B:92:MET:HE3	1:B:98:LYS:HE3	1.71	0.73
2:B:501:NAD:C5D	2:B:501:NAD:H6N	2.24	0.66
1:A:285:GLN:HE22	1:A:330:ALA:H	1.44	0.66
1:B:285:GLN:HE22	1:B:330:ALA:H	1.50	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:92:MET:HE3	1:B:98:LYS:CE	2.37	0.53
1:A:296:HIS:HD2	1:A:298:SER:OG	1.92	0.53
1:A:191:LEU:HD22	1:A:206:LEU:HB2	1.92	0.52
1:A:179:ALA:HB1	1:A:180:PRO:CD	2.43	0.49
1:A:344:ASN:ND2	1:A:379:ASP:H	2.11	0.48
1:A:371:ILE:HG22	1:A:374:VAL:HG22	1.96	0.47
1:B:333:ARG:HG3	4:B:789:HOH:O	2.14	0.47
1:B:148:VAL:HG22	1:B:226:LYS:HG3	1.96	0.46
1:A:179:ALA:HB1	1:A:180:PRO:HD2	1.98	0.45
1:A:180:PRO:HG2	1:A:183:THR:OG1	2.17	0.45
1:A:438:ILE:HD12	1:B:478:VAL:HG22	1.99	0.45
1:A:431:LEU:HD23	1:A:436:LEU:HD13	1.99	0.44
1:A:337:ALA:O	1:A:341:VAL:HG23	2.18	0.44
1:B:272:VAL:HG21	1:B:306:ALA:HB3	2.00	0.43
1:B:279:LYS:CE	1:B:389:VAL:O	2.66	0.43
1:B:179:ALA:HB3	1:B:208:TYR:CE1	2.54	0.42
1:A:314:LEU:HD22	1:A:325:THR:CG2	2.49	0.42
1:B:279:LYS:HE3	1:B:389:VAL:O	2.20	0.42
1:B:316:VAL:HG22	1:B:326:LEU:HD23	2.02	0.41
1:B:417:PHE:CD2	1:B:439:ASN:HA	2.55	0.41
1:B:295:VAL:O	1:B:394:GLY:HA2	2.20	0.41
1:A:300:ARG:NH1	1:A:301:ASP:OD1	2.54	0.41
1:A:344:ASN:HD21	1:A:379:ASP:H	1.67	0.41
1:B:279:LYS:O	1:B:279:LYS:HD2	2.21	0.41
1:A:191:LEU:CD2	1:A:206:LEU:HB2	2.51	0.40
1:B:191:LEU:HD22	1:B:206:LEU:HB2	2.01	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:B:864:HOH:O	4:B:864:HOH:O[5_455]	1.28	0.92
4:B:852:HOH:O	4:B:852:HOH:O[7_555]	1.45	0.75

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	474/505 (94%)	460 (97%)	14 (3%)	0	100	100
1	B	474/505 (94%)	460 (97%)	13 (3%)	1 (0%)	51	63
All	All	948/1010 (94%)	920 (97%)	27 (3%)	1 (0%)	55	67

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	411	GLY

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	359/385 (93%)	348 (97%)	11 (3%)	45	61
1	B	359/385 (93%)	350 (98%)	9 (2%)	53	70
All	All	718/770 (93%)	698 (97%)	20 (3%)	49	65

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	24	LYS
1	A	159	ASN
1	A	279	LYS
1	A	309	LYS
1	A	325	THR
1	A	334	ARG
1	A	343	ASP
1	A	362	GLU
1	A	381	PHE
1	A	386	PHE
1	A	393	ARG

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Mol	Chain	Res	Type
1	B	39	ARG
1	B	96	GLN
1	B	226	LYS
1	B	253	GLU
1	B	315	LYS
1	B	334	ARG
1	B	344	ASN
1	B	381	PHE
1	B	386	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	159	ASN
1	A	285	GLN
1	A	296	HIS
1	A	297	ASN
1	A	344	ASN
1	B	246	HIS
1	B	285	GLN
1	B	296	HIS
1	B	297	ASN
1	B	344	ASN

5.3.3 RNA [i](#)

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

3 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	NAD	A	501	-	41,48,48	1.34	6 (14%)	43,73,73	1.79	9 (20%)
3	GOL	A	502	-	5,5,5	0.47	0	5,5,5	0.76	0
2	NAD	B	501	-	41,48,48	1.03	3 (7%)	43,73,73	1.47	7 (16%)

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	NAD	A	501	-	-	0/22/62/62	0/5/5/5
3	GOL	A	502	-	-	0/4/4/4	0/0/0/0
2	NAD	B	501	-	-	0/22/62/62	0/5/5/5

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	501	NAD	O5D-C5D	-2.53	1.34	1.44
2	A	501	NAD	C2A-N3A	2.19	1.35	1.32
2	A	501	NAD	C3N-C7N	2.22	1.54	1.50
2	A	501	NAD	C2N-C3N	2.58	1.42	1.39
2	A	501	NAD	O4D-C1D	2.62	1.44	1.41
2	B	501	NAD	C5A-C4A	2.68	1.46	1.40
2	B	501	NAD	O4B-C1B	2.82	1.45	1.41
2	A	501	NAD	C5A-C4A	3.56	1.48	1.40
2	A	501	NAD	O4B-C1B	4.36	1.47	1.41

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	NAD	N3A-C2A-N1A	-6.56	123.15	128.86
2	B	501	NAD	N3A-C2A-N1A	-4.94	124.55	128.86
2	A	501	NAD	O7N-C7N-N7N	-4.29	116.48	122.58
2	B	501	NAD	C1B-N9A-C4A	-3.16	121.17	126.64

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	501	NAD	O2N-PN-O5D	-2.90	94.43	108.14
2	B	501	NAD	O4D-C4D-C5D	-2.49	100.99	109.40
2	A	501	NAD	C1B-N9A-C4A	-2.44	122.43	126.64
2	A	501	NAD	O4D-C4D-C3D	-2.40	100.40	105.17
2	B	501	NAD	C4A-C5A-N7A	-2.04	107.44	109.41
2	A	501	NAD	O2B-C2B-C3B	2.06	118.44	111.83
2	A	501	NAD	O2A-PA-O1A	2.25	123.94	112.28
2	B	501	NAD	O2N-PN-O1N	2.36	124.51	112.28
2	A	501	NAD	C3N-C7N-N7N	2.40	120.51	117.77
2	B	501	NAD	C3N-C2N-N1N	2.52	122.96	120.43
2	A	501	NAD	O7N-C7N-C3N	2.89	123.00	119.62
2	A	501	NAD	C4D-O4D-C1D	3.42	113.41	109.77

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	501	NAD	3	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	476/505 (94%)	-0.54	0 100 100	13, 19, 32, 53	0
1	B	476/505 (94%)	-0.50	0 100 100	14, 21, 34, 57	0
All	All	952/1010 (94%)	-0.52	0 100 100	13, 20, 34, 57	0

There are no RSRZ outliers to report.

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
3	GOL	A	502	6/6	0.90	0.17	1.18	32,36,38,39	0
2	NAD	A	501	44/44	0.90	0.15	0.91	26,45,49,51	0
2	NAD	B	501	44/44	0.91	0.14	0.41	24,37,49,56	0

6.5 Other polymers

There are no such residues in this entry.