



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 27, 2017 – 09:29 PM EDT

PDB ID : 5I95
Title : Crystal Structure of Human Mitochondrial Isocitrate Dehydrogenase R140Q
Mutant Homodimer bound to NADPH and alpha-Ketoglutaric acid
Authors : Zhang, B.; Jin, L.; Wu, W.; Jiang, F.; DeLaBarre, B.; Travins, J.A.; Padyana,
A.K.
Deposited on : 2016-02-19
Resolution : 1.54 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

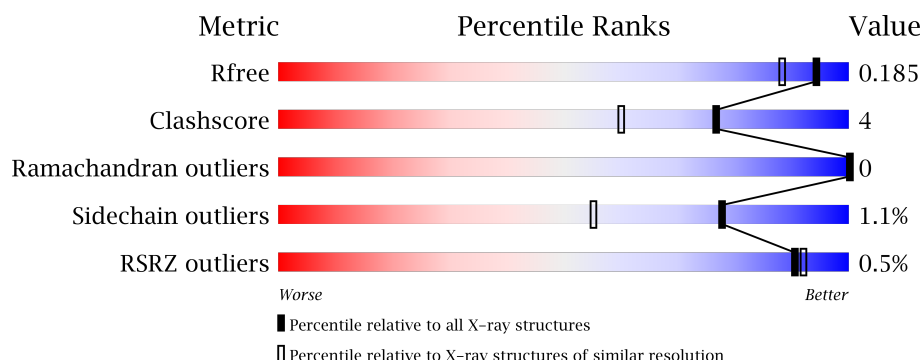
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION


The reported resolution of this entry is 1.54 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	1773 (1.56-1.52)
Clashscore	112137	1845 (1.56-1.52)
Ramachandran outliers	110173	1810 (1.56-1.52)
Sidechain outliers	110143	1808 (1.56-1.52)
RSRZ outliers	101464	1774 (1.56-1.52)

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	424	

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
4	GOL	A	506	-	-	-	X
5	ACT	A	507	-	-	-	X

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 4134 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.

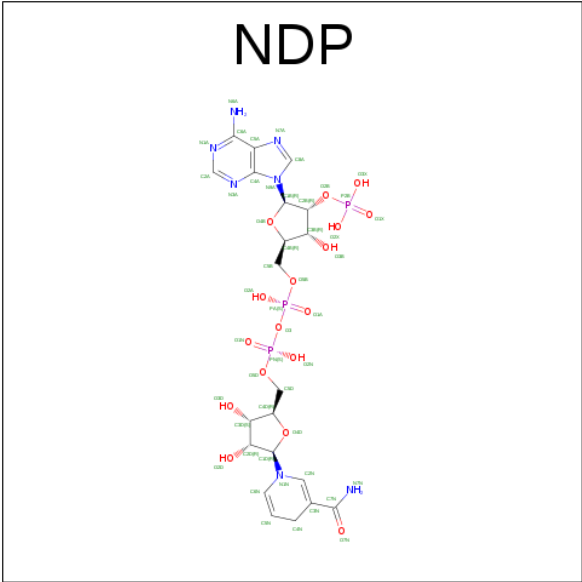
- Molecule 1 is a protein called Isocitrate dehydrogenase [NADP], mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	413	Total	C	N	O	S	0	28	0
			3455	2204	591	636	24			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	39	MET	-	initiating methionine	UNP P48735
A	140	GLN	ARG	engineered mutation	UNP P48735
A	453	LEU	-	expression tag	UNP P48735
A	454	GLU	-	expression tag	UNP P48735
A	455	HIS	-	expression tag	UNP P48735
A	456	HIS	-	expression tag	UNP P48735
A	457	HIS	-	expression tag	UNP P48735
A	458	HIS	-	expression tag	UNP P48735
A	459	HIS	-	expression tag	UNP P48735
A	460	HIS	-	expression tag	UNP P48735
A	461	HIS	-	expression tag	UNP P48735
A	462	HIS	-	expression tag	UNP P48735

- Molecule 2 is NADPH DIHYDRO-NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (three-letter code: NDP) (formula: C₂₁H₃₀N₇O₁₇P₃).

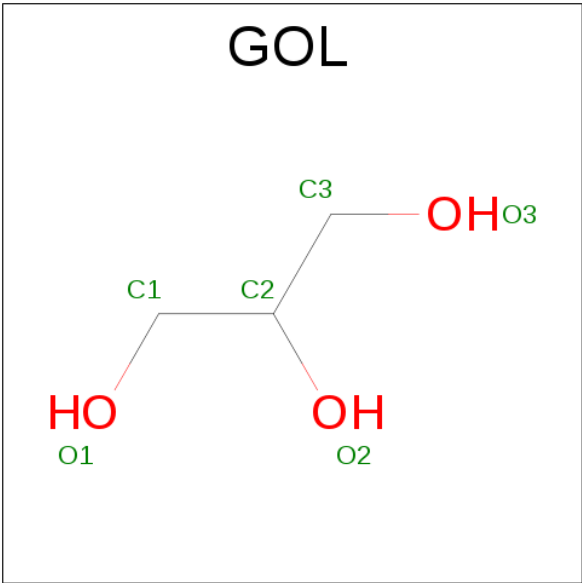


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0
			48	21	7	17	3	

- Molecule 3 is CALCIUM ION (three-letter code: CA) (formula: Ca).

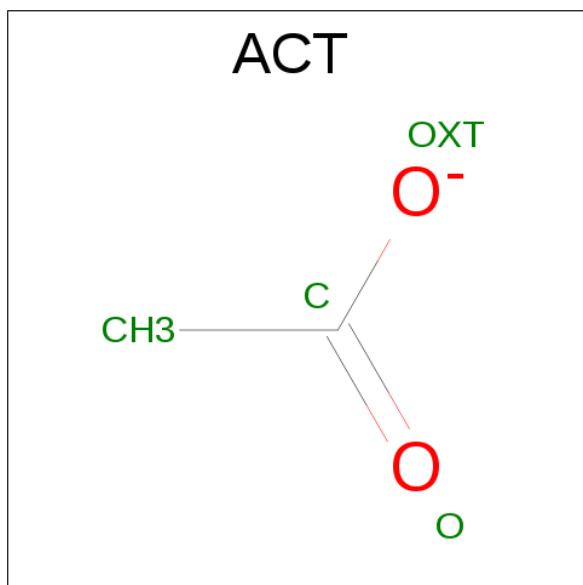
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Ca	0	1
			2	2		

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



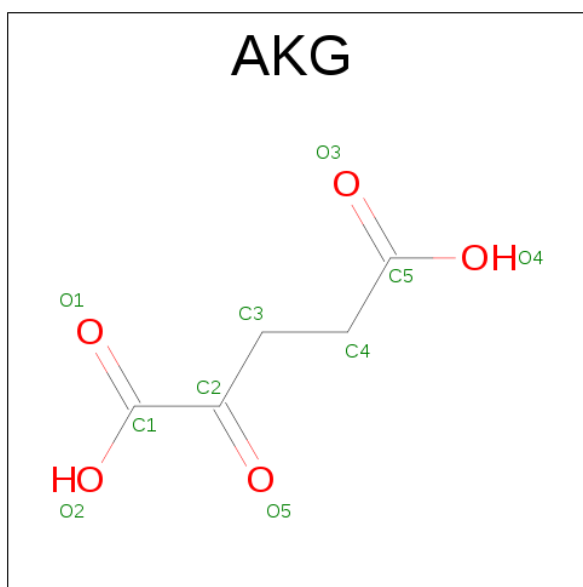
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		

- Molecule 5 is ACETATE ION (three-letter code: ACT) (formula: $C_2H_3O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			4	2	2		

- Molecule 6 is 2-OXOGLUTARIC ACID (three-letter code: AKG) (formula: $C_5H_6O_5$).



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

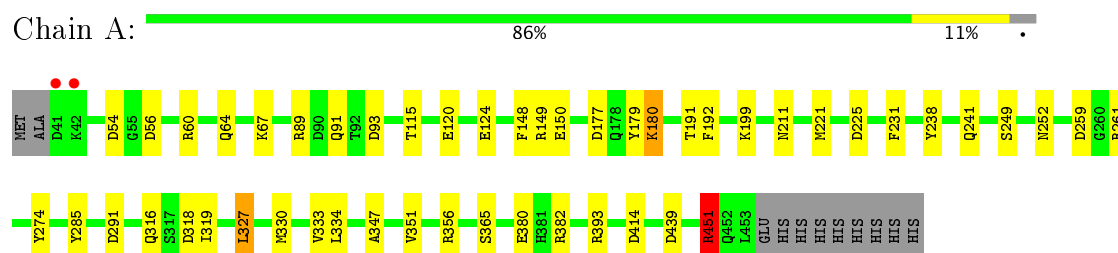
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	589	Total	O	0	3
			591	591		

3 Residue-property plots

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: Isocitrate dehydrogenase [NADP], mitochondrial



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	67.18 Å 154.88 Å 93.57 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.83 – 1.54 29.27 – 1.54	Depositor EDS
% Data completeness (in resolution range)	99.4 (29.83-1.54) 99.4 (29.27-1.54)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.42 (at 1.54 Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.142 , 0.174 0.156 , 0.185	Depositor DCC
R_{free} test set	3624 reflections (5.31%)	DCC
Wilson B-factor (Å ²)	13.7	Xtriage
Anisotropy	0.032	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 54.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	4134	wwPDB-VP
Average B, all atoms (Å ²)	18.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.67% 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: GOL, CA, AKG, NDP, ACT

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	1.28	9/3595 (0.3%)	1.29	31/4847 (0.6%)

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	249	SER	CA-CB	6.61	1.62	1.52
1	A	451[A]	ARG	CB-CG	-6.54	1.34	1.52
1	A	451[B]	ARG	CB-CG	-6.54	1.34	1.52
1	A	180	LYS	CB-CG	-5.85	1.36	1.52
1	A	179	TYR	N-CA	-5.70	1.34	1.46
1	A	150	GLU	CB-CG	-5.61	1.41	1.52
1	A	365	SER	CA-CB	-5.27	1.45	1.52
1	A	120	GLU	CD-OE1	5.12	1.31	1.25
1	A	150	GLU	CG-CD	5.11	1.59	1.51

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	451[A]	ARG	CB-CG-CD	10.81	139.71	111.60
1	A	451[B]	ARG	CB-CG-CD	10.81	139.71	111.60
1	A	60	ARG	NE-CZ-NH2	9.78	125.19	120.30
1	A	54	ASP	CB-CG-OD1	9.22	126.60	118.30
1	A	177	ASP	CB-CG-OD1	9.05	126.45	118.30
1	A	180	LYS	CD-CE-NZ	-7.65	94.11	111.70
1	A	382	ARG	NE-CZ-NH1	7.34	123.97	120.30
1	A	356	ARG	NE-CZ-NH1	7.22	123.91	120.30
1	A	231	PHE	CB-CG-CD1	-7.12	115.81	120.80
1	A	259	ASP	CB-CG-OD2	7.12	124.70	118.30
1	A	56	ASP	CB-CG-OD2	6.88	124.49	118.30
1	A	382	ARG	NE-CZ-NH2	-6.76	116.92	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	93	ASP	CB-CG-OD1	6.62	124.26	118.30
1	A	451[A]	ARG	NE-CZ-NH2	6.42	123.51	120.30
1	A	451[B]	ARG	NE-CZ-NH2	6.42	123.51	120.30
1	A	327[A]	LEU	CB-CG-CD1	-6.30	100.30	111.00
1	A	327[B]	LEU	CB-CG-CD1	-6.30	100.30	111.00
1	A	238	TYR	CB-CG-CD1	6.29	124.77	121.00
1	A	414	ASP	CB-CG-OD2	6.12	123.81	118.30
1	A	192	PHE	CB-CG-CD1	6.05	125.04	120.80
1	A	274	TYR	CG-CD1-CE1	5.93	126.05	121.30
1	A	149	ARG	NE-CZ-NH1	-5.66	117.47	120.30
1	A	89	ARG	NE-CZ-NH2	-5.65	117.47	120.30
1	A	285	TYR	CB-CG-CD1	-5.57	117.66	121.00
1	A	148	PHE	CB-CG-CD2	5.56	124.69	120.80
1	A	225	ASP	CB-CG-OD1	5.47	123.22	118.30
1	A	149	ARG	NE-CZ-NH2	5.42	123.01	120.30
1	A	439	ASP	CB-CG-OD1	5.35	123.11	118.30
1	A	285	TYR	CB-CG-CD2	5.21	124.12	121.00
1	A	393	ARG	NE-CZ-NH1	5.18	122.89	120.30
1	A	261	ARG	NE-CZ-NH1	5.03	122.81	120.30

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	3455	0	3475	23	0
2	A	48	0	26	2	0
3	A	2	0	0	1	0
4	A	24	0	32	0	0
5	A	4	0	3	0	0
6	A	10	0	4	1	0
7	A	591	0	0	17	1
All	All	4134	0	3540	25	1

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 4.

All (25) 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)
3:A:502[B]:CA:CA	7:A:619:HOH:O	1.66	0.70
1:A:333[A]:VAL:CG2	7:A:977:HOH:O	2.42	0.67
1:A:380[B]:GLU:OE2	7:A:603[B]:HOH:O	2.14	0.66
1:A:380[A]:GLU:OE2	7:A:603[A]:HOH:O	2.17	0.60
1:A:241[B]:GLN:OE1	7:A:602:HOH:O	2.16	0.59
1:A:221[B]:MET:HE1	7:A:745:HOH:O	2.05	0.56
1:A:221[B]:MET:CE	7:A:745:HOH:O	2.54	0.55
2:A:501:NDP:H42N	6:A:508:AKG:C2	2.38	0.53
1:A:199[B]:LYS:NZ	7:A:606:HOH:O	2.38	0.52
1:A:347:ALA:HB3	7:A:672:HOH:O	2.11	0.51
1:A:316:GLN:OE1	1:A:319[A]:ILE:HD12	2.11	0.50
1:A:333[A]:VAL:HG23	7:A:977:HOH:O	2.09	0.49
1:A:64:GLN:OE1	7:A:604:HOH:O	2.20	0.49
1:A:115:THR:O	2:A:501:NDP:H2N	2.13	0.48
1:A:333[A]:VAL:HG21	7:A:977:HOH:O	2.10	0.48
1:A:330[B]:MET:HE3	1:A:330[B]:MET:HB2	1.83	0.47
1:A:451[B]:ARG:NH2	7:A:617:HOH:O	2.47	0.47
1:A:67:LYS:NZ	7:A:618:HOH:O	2.47	0.47
1:A:180:LYS:NZ	7:A:601:HOH:O	2.00	0.45
1:A:180:LYS:CE	7:A:601:HOH:O	2.60	0.45
1:A:124[A]:GLU:HG2	7:A:667:HOH:O	2.19	0.43
1:A:291:ASP:OD1	1:A:291:ASP:N	2.49	0.43
1:A:318[B]:ASP:OD1	1:A:327[B]:LEU:HD11	2.18	0.43
1:A:191[B]:THR:HG22	1:A:211[B]:ASN:OD1	2.19	0.41

All (1) 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)
7:A:855:HOH:O	7:A:947:HOH:O[3_454]	1.97	0.23

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	440/424 (104%)	425 (97%)	15 (3%)	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	382/364 (105%)	377 (99%)	5 (1%)	73	46

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	91	GLN
1	A	252	ASN
1	A	351	VAL
1	A	451[A]	ARG
1	A	451[B]	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	91	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 9 ligands modelled in this entry, 2 are monoatomic - leaving 7 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	$\# Z > 2$
2	NDP	A	501	-	43,52,52	1.78	8 (18%)	49,80,80	1.63	11 (22%)
4	GOL	A	503	-	5,5,5	0.55	0	5,5,5	0.74	0
4	GOL	A	504	-	5,5,5	0.72	0	5,5,5	1.17	1 (20%)
4	GOL	A	505	-	5,5,5	0.94	0	5,5,5	1.25	0
4	GOL	A	506	-	5,5,5	0.73	0	5,5,5	0.85	0
5	ACT	A	507	-	1,3,3	0.59	0	0,3,3	0.00	-
6	AKG	A	508	3	3,9,9	1.66	1 (33%)	4,11,11	2.37	3 (75%)

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	NDP	A	501	-	-	0/30/77/77	0/5/5/5
4	GOL	A	503	-	-	0/4/4/4	0/0/0/0
4	GOL	A	504	-	-	0/4/4/4	0/0/0/0
4	GOL	A	505	-	-	0/4/4/4	0/0/0/0
4	GOL	A	506	-	-	0/4/4/4	0/0/0/0
5	ACT	A	507	-	-	0/0/0/0	0/0/0/0
6	AKG	A	508	3	-	0/3/9/9	0/0/0/0

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	NDP	C4N-C5N	-4.43	1.39	1.49
2	A	501	NDP	O4B-C4B	-2.79	1.38	1.45
2	A	501	NDP	C2A-N1A	-2.59	1.29	1.33
2	A	501	NDP	O4B-C1B	2.39	1.44	1.41
2	A	501	NDP	O3D-C3D	2.56	1.48	1.43
2	A	501	NDP	O2D-C2D	2.69	1.49	1.43
6	A	508	AKG	O5-C2	2.83	1.27	1.22
2	A	501	NDP	C6N-C5N	5.16	1.42	1.33
2	A	501	NDP	C2A-N3A	5.24	1.40	1.32

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	NDP	N3A-C2A-N1A	-3.97	125.40	128.86
2	A	501	NDP	O7N-C7N-N7N	-3.18	115.20	122.92
6	A	508	AKG	C4-C3-C2	-3.07	105.95	113.04
2	A	501	NDP	O4D-C1D-C2D	-2.70	100.66	106.64
6	A	508	AKG	C3-C2-C1	-2.66	115.57	121.60
6	A	508	AKG	C3-C4-C5	-2.27	108.78	112.66
2	A	501	NDP	C1D-N1N-C2N	-2.25	117.27	121.09
2	A	501	NDP	C2B-C3B-C4B	-2.14	97.08	101.95
2	A	501	NDP	C5A-C6A-N1A	-2.09	113.38	119.70
2	A	501	NDP	O3D-C3D-C2D	-2.06	105.23	111.83
4	A	504	GOL	O1-C1-C2	2.25	121.41	110.07
2	A	501	NDP	O3B-C3B-C2B	2.38	117.95	111.18
2	A	501	NDP	N6A-C6A-N1A	3.39	125.48	118.77
2	A	501	NDP	C4A-C5A-N7A	3.59	112.88	109.41
2	A	501	NDP	C2A-N1A-C6A	4.13	125.99	118.77

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	501	NDP	2	0
6	A	508	AKG	1	0

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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	413/424 (97%)	-0.30	2 (0%) 90 92	7, 14, 28, 50	1 (0%)

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	41	ASP	2.7
1	A	42	LYS	2.2

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
4	GOL	A	506	6/6	0.92	0.20	5.67	23,28,37,46	0
5	ACT	A	507	4/4	0.89	0.19	3.87	35,38,39,43	0
6	AKG	A	508	10/10	0.93	0.09	1.53	17,21,24,28	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
4	GOL	A	503	6/6	0.95	0.10	1.44	23,25,27,28	0
4	GOL	A	504	6/6	0.87	0.11	0.82	25,40,47,52	0
4	GOL	A	505	6/6	0.93	0.10	0.48	21,33,38,39	0
2	NDP	A	501	48/48	0.99	0.04	-1.67	8,9,12,20	0
3	CA	A	502[B]	1/1	0.98	0.04	-2.05	13,13,13,13	1
3	CA	A	502[A]	1/1	0.98	0.04	-2.52	9,9,9,9	1

6.5 Other polymers [i](#)

There are no such residues in this entry.