



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 27, 2014 – 04:51 PM GMT

PDB ID : 1GCO
Title : CRYSTAL STRUCTURE OF GLUCOSE DEHYDROGENASE COM-
PLEXED WITH NAD+
Authors : Yamamoto, K.; Kurisu, G.; Kusunoki, M.; Tabata, S.; Urabe, I.; Osaki, S.
Deposited on : 2000-08-07
Resolution : 1.70 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

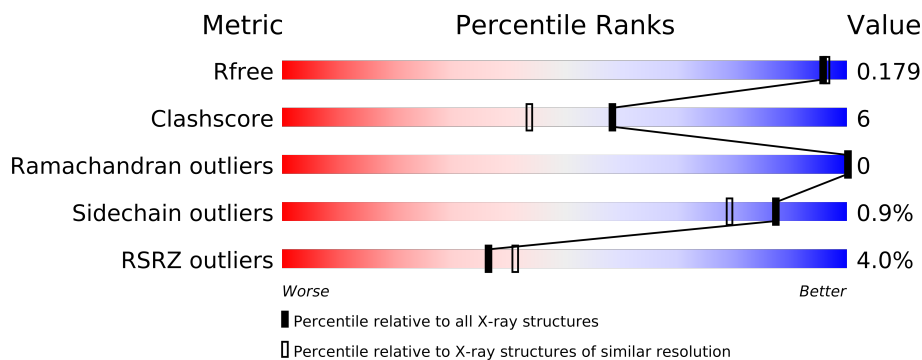
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 1.70 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}	66092	2456 (1.70-1.70)
Clashscore	79885	2929 (1.70-1.70)
Ramachandran outliers	78287	2878 (1.70-1.70)
Sidechain outliers	78261	2878 (1.70-1.70)
RSRZ outliers	66119	2456 (1.70-1.70)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	261	
1	B	261	
1	E	261	
1	F	261	

2 Entry composition i

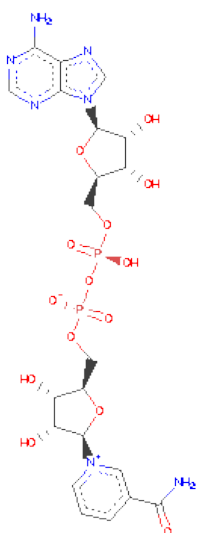
There are 3 unique types of molecules in this entry. The entry contains 8626 atoms, of which 0 are hydrogen and 0 are deuterium.

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 GLUCOSE DEHYDROGENASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	261	Total	C	N	O	S	0	0	0
			1972	1248	325	389	10			
1	B	261	Total	C	N	O	S	0	0	0
			1972	1248	325	389	10			
1	E	261	Total	C	N	O	S	0	0	0
			1972	1248	325	389	10			
1	F	261	Total	C	N	O	S	0	0	0
			1972	1248	325	389	10			

- Molecule 2 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (three-letter code: NAD) (formula: $C_{21}H_{27}N_7O_{14}P_2$).



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		

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

- Molecule 3 is water.

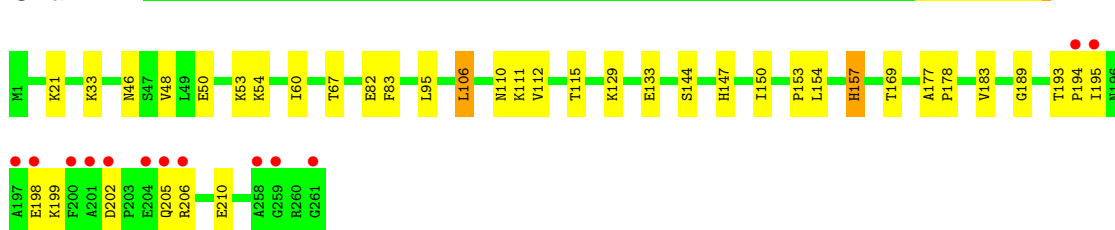
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	142	Total	O	0	0
			142	142		
3	B	135	Total	O	0	0
			135	135		
3	E	145	Total	O	0	0
			145	145		
3	F	140	Total	O	0	0
			140	140		

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 errors 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.

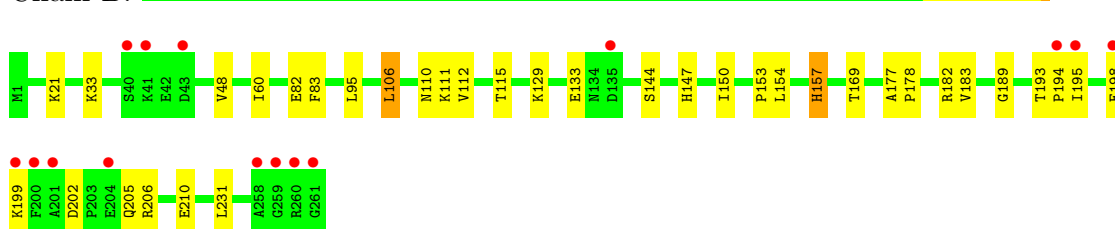
• Molecule 1: GLUCOSE DEHYDROGENASE

Chain A:



• Molecule 1: GLUCOSE DEHYDROGENASE

Chain B:



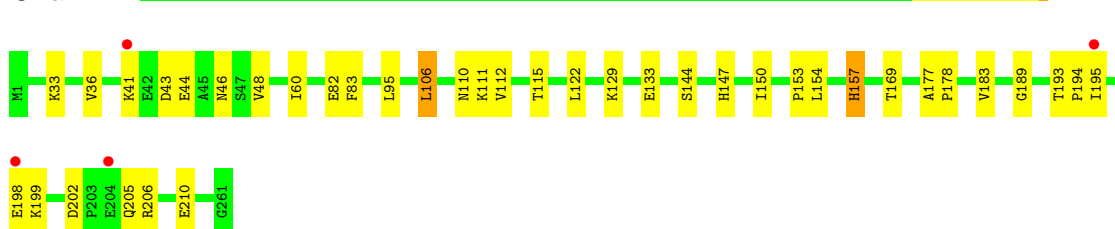
• Molecule 1: GLUCOSE DEHYDROGENASE

Chain E:



• Molecule 1: GLUCOSE DEHYDROGENASE

Chain F:



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	120.67Å 66.56Å 119.53Å 90.00° 93.21° 90.00°	Depositor
Resolution (Å)	34.39 – 1.70 55.03 – 1.70	Depositor EDS
% Data completeness (in resolution range)	90.7 (34.39-1.70) 91.1 (55.03-1.70)	Depositor EDS
R_{merge}	0.03	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	9.37 (at 1.70Å)	Xtriage
Refinement program	CNS 0.9	Depositor
R, R_{free}	0.178 , 0.192 0.180 , 0.179	Depositor DCC
R_{free} test set	4751 reflections (5.01%)	DCC
Wilson B-factor (Å ²)	17.6	Xtriage
Anisotropy	0.355	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 31.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 94820 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	8626	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.91% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 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	0.28	0/2006	0.56	0/2712
1	B	0.29	0/2006	0.57	0/2712
1	E	0.29	0/2006	0.57	0/2712
1	F	0.28	0/2006	0.56	0/2712
All	All	0.28	0/8024	0.56	0/10848

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1972	0	1974	26	0
1	B	1972	0	1974	23	0
1	E	1972	0	1974	25	0
1	F	1972	0	1974	28	0
2	A	44	0	26	2	0
2	B	44	0	26	2	0
2	E	44	0	26	1	0
2	F	44	0	26	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	142	0	0	3	0
3	B	135	0	0	0	0
3	E	145	0	0	1	0
3	F	140	0	0	1	0
All	All	8626	0	8000	100	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 6.

All (100) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:E:40:SER:HA	1:E:63:LYS:HE3	1.52	0.88
1:E:193:THR:HB	1:E:194:PRO:HD2	1.64	0.79
1:B:193:THR:HB	1:B:194:PRO:HD2	1.64	0.79
1:F:193:THR:HB	1:F:194:PRO:HD2	1.65	0.78
1:A:193:THR:HB	1:A:194:PRO:HD2	1.65	0.77
1:F:41:LYS:HE3	1:F:43:ASP:HB2	1.75	0.69
1:E:153:PRO:O	1:E:154:LEU:HB2	1.96	0.66
1:F:60:ILE:HD13	1:F:82:GLU:HG2	1.79	0.65
1:A:153:PRO:O	1:A:154:LEU:HB2	1.96	0.65
1:B:153:PRO:O	1:B:154:LEU:HB2	1.97	0.65
1:E:41:LYS:HE3	1:E:43:ASP:OD2	1.97	0.65
1:B:60:ILE:HD13	1:B:82:GLU:HG2	1.79	0.65
1:A:60:ILE:HD13	1:A:82:GLU:HG2	1.79	0.64
1:A:193:THR:OG1	1:A:195:ILE:HG22	1.98	0.64
1:F:153:PRO:O	1:F:154:LEU:HB2	1.97	0.64
1:E:60:ILE:HD13	1:E:82:GLU:HG2	1.79	0.64
1:F:193:THR:OG1	1:F:195:ILE:HG22	1.98	0.63
1:A:106:LEU:HD22	1:A:110:ASN:ND2	2.14	0.63
1:E:193:THR:OG1	1:E:195:ILE:HG22	1.98	0.63
1:B:193:THR:OG1	1:B:195:ILE:HG22	1.99	0.63
1:B:106:LEU:HD22	1:B:110:ASN:ND2	2.14	0.63
1:E:106:LEU:HD22	1:E:110:ASN:ND2	2.15	0.61
1:F:106:LEU:HD22	1:F:110:ASN:ND2	2.15	0.61
1:A:189:GLY:O	2:A:601:NAD:H4N	2.04	0.58
1:A:53:LYS:HE3	3:A:667:HOH:O	2.03	0.57
1:E:147:HIS:HD2	1:E:150:ILE:O	1.88	0.57
1:F:147:HIS:HD2	1:F:150:ILE:O	1.90	0.55
1:B:189:GLY:O	2:B:701:NAD:H4N	2.06	0.55
1:F:189:GLY:O	2:F:901:NAD:H4N	2.07	0.55
1:B:147:HIS:HD2	1:B:150:ILE:O	1.89	0.55
1:A:147:HIS:HD2	1:A:150:ILE:O	1.89	0.54

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:129:LYS:O	1:B:133:GLU:HG3	2.08	0.53
1:A:46:ASN:O	1:A:50:GLU:HG2	2.09	0.53
1:E:189:GLY:O	2:E:801:NAD:H4N	2.09	0.52
1:E:36:VAL:HG11	1:E:48:VAL:HG11	1.92	0.52
1:B:21:LYS:HG3	1:B:48:VAL:HG22	1.91	0.52
1:A:129:LYS:O	1:A:133:GLU:HG3	2.11	0.51
1:E:38:TYR:CZ	1:E:42:GLU:HA	2.46	0.51
1:F:129:LYS:O	1:F:133:GLU:HG3	2.11	0.51
1:A:21:LYS:HG3	1:A:48:VAL:HG22	1.93	0.51
1:E:129:LYS:O	1:E:133:GLU:HG3	2.11	0.50
1:F:111:LYS:NZ	3:F:934:HOH:O	2.44	0.50
1:F:95:LEU:HD11	1:F:111:LYS:HG2	1.94	0.49
1:A:95:LEU:HD11	1:A:111:LYS:HG2	1.95	0.49
1:E:95:LEU:HD11	1:E:111:LYS:HG2	1.95	0.49
1:B:95:LEU:HD11	1:B:111:LYS:HG2	1.95	0.48
1:E:43:ASP:HA	1:E:46:ASN:ND2	2.29	0.47
1:F:144:SER:O	2:F:901:NAD:H6N	2.15	0.46
1:A:111:LYS:NZ	3:A:695:HOH:O	2.48	0.46
1:B:169:THR:HG23	1:B:183:VAL:HG12	1.98	0.46
1:A:206:ARG:O	1:A:210:GLU:HG3	2.15	0.46
1:A:177:ALA:N	1:A:178:PRO:CD	2.79	0.46
1:E:33:LYS:HD3	1:E:83:PHE:CD2	2.51	0.46
1:E:206:ARG:O	1:E:210:GLU:HG3	2.16	0.45
1:B:202:ASP:OD2	1:B:205:GLN:HG3	2.17	0.45
1:B:206:ARG:O	1:B:210:GLU:HG3	2.16	0.45
1:B:33:LYS:HD3	1:B:83:PHE:CD2	2.51	0.45
1:F:36:VAL:HG11	1:F:48:VAL:HG11	1.98	0.45
1:E:177:ALA:N	1:E:178:PRO:CD	2.80	0.45
1:F:199:LYS:NZ	1:F:199:LYS:HB2	2.32	0.45
1:E:202:ASP:OD2	1:E:205:GLN:HG3	2.17	0.45
1:F:177:ALA:N	1:F:178:PRO:CD	2.80	0.45
1:A:33:LYS:HD3	1:A:83:PHE:CD2	2.51	0.45
1:A:67:THR:HA	3:A:679:HOH:O	2.16	0.45
1:A:199:LYS:HB2	1:A:199:LYS:NZ	2.32	0.44
1:F:202:ASP:OD2	1:F:205:GLN:HG3	2.17	0.44
1:B:199:LYS:NZ	1:B:199:LYS:HB2	2.32	0.44
1:F:43:ASP:HA	1:F:46:ASN:ND2	2.32	0.44
1:B:177:ALA:N	1:B:178:PRO:CD	2.80	0.44
1:F:33:LYS:HD3	1:F:83:PHE:CD2	2.52	0.44
1:E:199:LYS:HB2	1:E:199:LYS:NZ	2.32	0.44
1:A:106:LEU:HD22	1:A:110:ASN:HD21	1.82	0.44
1:B:144:SER:O	2:B:701:NAD:H6N	2.18	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:F:206:ARG:O	1:F:210:GLU:HG3	2.17	0.44
1:A:202:ASP:OD2	1:A:205:GLN:HG3	2.17	0.43
1:F:41:LYS:O	1:F:44:GLU:HG3	2.18	0.43
1:A:54:LYS:HE3	1:A:54:LYS:HB2	1.88	0.43
1:F:41:LYS:HG2	1:F:43:ASP:H	1.84	0.42
1:E:169:THR:HG23	1:E:183:VAL:HG12	2.00	0.42
1:F:193:THR:HB	1:F:194:PRO:CD	2.44	0.42
1:A:144:SER:O	2:A:601:NAD:H6N	2.19	0.42
1:F:169:THR:HG23	1:F:183:VAL:HG12	2.01	0.42
1:E:106:LEU:HG	1:F:122:LEU:HD21	2.01	0.42
1:E:67:THR:HA	3:E:899:HOH:O	2.19	0.42
1:E:182:ARG:NH2	1:E:231:LEU:O	2.50	0.42
1:B:193:THR:HB	1:B:194:PRO:CD	2.44	0.42
1:A:95:LEU:CD2	1:A:115:THR:HB	2.50	0.42
1:E:122:LEU:HD21	1:F:106:LEU:HG	2.01	0.41
1:E:95:LEU:CD2	1:E:115:THR:HB	2.50	0.41
1:F:95:LEU:CD2	1:F:115:THR:HB	2.50	0.41
1:B:95:LEU:CD2	1:B:115:THR:HB	2.51	0.41
1:B:193:THR:HG1	1:B:195:ILE:HG22	1.84	0.41
1:A:169:THR:HG23	1:A:183:VAL:HG12	2.02	0.41
1:F:112:VAL:HG21	1:F:157:HIS:CD2	2.56	0.41
1:F:198:GLU:HG2	1:F:199:LYS:N	2.36	0.41
1:A:198:GLU:HG2	1:A:199:LYS:N	2.36	0.40
1:B:198:GLU:HG2	1:B:199:LYS:N	2.36	0.40
1:A:112:VAL:HG21	1:A:157:HIS:CD2	2.56	0.40
1:B:182:ARG:NH2	1:B:231:LEU:O	2.50	0.40
1:B:112:VAL:HG21	1:B:157:HIS:CD2	2.57	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	259/261 (99%)	247 (95%)	12 (5%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	259/261 (99%)	247 (95%)	12 (5%)	0	100	100
1	E	259/261 (99%)	248 (96%)	11 (4%)	0	100	100
1	F	259/261 (99%)	246 (95%)	13 (5%)	0	100	100
All	All	1036/1044 (99%)	988 (95%)	48 (5%)	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	211/211 (100%)	209 (99%)	2 (1%)	87	79
1	B	211/211 (100%)	209 (99%)	2 (1%)	87	79
1	E	211/211 (100%)	209 (99%)	2 (1%)	87	79
1	F	211/211 (100%)	209 (99%)	2 (1%)	87	79
All	All	844/844 (100%)	836 (99%)	8 (1%)	87	79

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	106	LEU
1	A	157	HIS
1	B	106	LEU
1	B	157	HIS
1	E	106	LEU
1	E	157	HIS
1	F	106	LEU
1	F	157	HIS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	110	ASN

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Mol	Chain	Res	Type
1	A	147	HIS
1	B	147	HIS
1	E	46	ASN
1	E	147	HIS
1	F	46	ASN
1	F	147	HIS

5.3.3 RNA ⓘ

There are no RNA chains 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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

4 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	601	-	48,48,48	2.01	8 (16%)	73,73,73	1.48	9 (12%)
2	NAD	B	701	-	48,48,48	2.02	10 (20%)	73,73,73	1.47	9 (12%)
2	NAD	E	801	-	48,48,48	2.01	9 (18%)	73,73,73	1.46	9 (12%)
2	NAD	F	901	-	48,48,48	2.02	8 (16%)	73,73,73	1.48	9 (12%)

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	601	-	-	0/30/62/62	0/3/5/5
2	NAD	B	701	-	-	0/30/62/62	0/3/5/5
2	NAD	E	801	-	-	0/30/62/62	0/3/5/5
2	NAD	F	901	-	-	0/30/62/62	0/3/5/5

All (35) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	701	NAD	C2N-C3N	7.66	1.49	1.38
2	F	901	NAD	C2N-C3N	7.56	1.49	1.38
2	E	801	NAD	C2N-C3N	7.45	1.49	1.38
2	A	601	NAD	C2N-C3N	7.41	1.49	1.38
2	A	601	NAD	C4N-C3N	7.05	1.51	1.39
2	E	801	NAD	C4N-C3N	6.90	1.51	1.39
2	B	701	NAD	C4N-C3N	6.85	1.51	1.39
2	F	901	NAD	C4N-C3N	6.67	1.50	1.39
2	A	601	NAD	C2N-N1N	4.65	1.41	1.35
2	E	801	NAD	C2N-N1N	4.51	1.41	1.35
2	B	701	NAD	C2N-N1N	4.44	1.41	1.35
2	F	901	NAD	C2N-N1N	4.40	1.40	1.35
2	E	801	NAD	C5N-C4N	4.24	1.48	1.39
2	A	601	NAD	C5N-C4N	4.15	1.48	1.39
2	F	901	NAD	C5N-C4N	4.15	1.48	1.39
2	B	701	NAD	C5N-C4N	4.11	1.48	1.39
2	F	901	NAD	PA-O3	-3.30	1.53	1.59
2	F	901	NAD	C6N-C5N	-3.09	1.31	1.38
2	B	701	NAD	C6N-C5N	-2.98	1.31	1.38
2	A	601	NAD	PA-O3	-2.86	1.54	1.59
2	E	801	NAD	C6N-C5N	-2.86	1.31	1.38
2	E	801	NAD	PA-O3	-2.82	1.54	1.59
2	A	601	NAD	C6N-C5N	-2.77	1.32	1.38
2	B	701	NAD	PA-O3	-2.74	1.54	1.59
2	E	801	NAD	C2D-C1D	-2.58	1.49	1.53
2	F	901	NAD	C2D-C1D	-2.50	1.49	1.53
2	F	901	NAD	C6N-N1N	2.46	1.42	1.35
2	B	701	NAD	C2D-C1D	-2.42	1.50	1.53
2	E	801	NAD	C6N-N1N	2.41	1.42	1.35
2	A	601	NAD	C6N-N1N	2.40	1.42	1.35
2	B	701	NAD	C6N-N1N	2.38	1.42	1.35
2	A	601	NAD	C2D-C1D	-2.28	1.50	1.53
2	B	701	NAD	C2A-N1A	2.16	1.38	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	701	NAD	C3N-C7N	-2.03	1.47	1.50
2	E	801	NAD	C2A-N1A	2.00	1.37	1.33

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	601	NAD	C5N-C4N-C3N	-5.59	113.06	120.32
2	B	701	NAD	C5N-C4N-C3N	-5.54	113.12	120.32
2	F	901	NAD	C5N-C4N-C3N	-5.54	113.13	120.32
2	E	801	NAD	C5N-C4N-C3N	-5.45	113.24	120.32
2	E	801	NAD	C3N-C7N-N7N	5.05	123.51	117.77
2	F	901	NAD	C3N-C7N-N7N	5.02	123.48	117.77
2	A	601	NAD	C3N-C7N-N7N	5.00	123.46	117.77
2	B	701	NAD	C3N-C7N-N7N	4.89	123.34	117.77
2	F	901	NAD	C6N-C5N-C4N	3.95	125.71	119.44
2	B	701	NAD	C6N-C5N-C4N	3.91	125.65	119.44
2	A	601	NAD	C6N-C5N-C4N	3.91	125.64	119.44
2	E	801	NAD	C6N-C5N-C4N	3.78	125.44	119.44
2	F	901	NAD	C4B-O4B-C1B	-3.02	106.47	109.75
2	B	701	NAD	C4B-O4B-C1B	-3.00	106.49	109.75
2	A	601	NAD	C4B-O4B-C1B	-2.98	106.52	109.75
2	E	801	NAD	C4B-O4B-C1B	-2.96	106.54	109.75
2	F	901	NAD	O7N-C7N-C3N	-2.79	116.43	119.58
2	B	701	NAD	O7N-C7N-C3N	-2.70	116.54	119.58
2	A	601	NAD	C6N-N1N-C2N	-2.67	119.02	122.04
2	E	801	NAD	O7N-C7N-C3N	-2.66	116.58	119.58
2	A	601	NAD	O7N-C7N-C3N	-2.58	116.67	119.58
2	F	901	NAD	C6N-N1N-C2N	-2.57	119.13	122.04
2	B	701	NAD	O4B-C1B-C2B	-2.51	102.93	106.77
2	A	601	NAD	O4B-C1B-C2B	-2.49	102.95	106.77
2	B	701	NAD	C6N-N1N-C2N	-2.48	119.24	122.04
2	F	901	NAD	O4B-C1B-C2B	-2.46	103.00	106.77
2	E	801	NAD	C6N-N1N-C2N	-2.45	119.28	122.04
2	A	601	NAD	C5A-C6A-N6A	2.43	126.21	120.72
2	F	901	NAD	C5A-C6A-N6A	2.43	126.21	120.72
2	E	801	NAD	O4B-C1B-C2B	-2.43	103.05	106.77
2	B	701	NAD	C5A-C6A-N6A	2.39	126.12	120.72
2	E	801	NAD	C5A-C6A-N6A	2.38	126.10	120.72
2	B	701	NAD	O2N-PN-O5D	2.23	117.42	108.19
2	A	601	NAD	O2N-PN-O5D	2.14	117.02	108.19
2	E	801	NAD	O2N-PN-O5D	2.06	116.69	108.19
2	F	901	NAD	O2N-PN-O5D	2.02	116.52	108.19

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

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 ⓘ

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	261/261 (100%)	-0.06	13 (4%)	28 32	11, 18, 35, 42	0
1	B	261/261 (100%)	0.07	15 (5%)	23 27	11, 17, 37, 43	0
1	E	261/261 (100%)	0.04	10 (3%)	38 43	9, 17, 34, 42	0
1	F	261/261 (100%)	-0.01	4 (1%)	70 76	9, 17, 36, 41	0
All	All	1044/1044 (100%)	0.01	42 (4%)	36 41	9, 17, 37, 43	0

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	195	ILE	12.6
1	A	195	ILE	8.9
1	B	195	ILE	6.2
1	E	198	GLU	5.3
1	F	195	ILE	5.0
1	B	261	GLY	4.6
1	B	198	GLU	4.6
1	B	201	ALA	4.5
1	E	261	GLY	4.5
1	A	259	GLY	4.4
1	A	198	GLU	4.3
1	A	261	GLY	4.2
1	B	41	LYS	3.7
1	E	259	GLY	3.4
1	A	201	ALA	3.2
1	F	41	LYS	3.2
1	B	204	GLU	3.1
1	A	204	GLU	3.1
1	B	135	ASP	2.7
1	A	197	ALA	2.7
1	A	200	PHE	2.7

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Mol	Chain	Res	Type	RSRZ
1	E	258	ALA	2.6
1	A	202	ASP	2.6
1	E	201	ALA	2.6
1	B	43	ASP	2.6
1	E	197	ALA	2.6
1	B	40	SER	2.5
1	E	199	LYS	2.3
1	A	205	GLN	2.3
1	B	259	GLY	2.3
1	E	194	PRO	2.2
1	B	199	LYS	2.2
1	B	194	PRO	2.2
1	B	200	PHE	2.2
1	A	194	PRO	2.1
1	E	63	LYS	2.1
1	F	198	GLU	2.1
1	F	204	GLU	2.1
1	A	206	ARG	2.1
1	A	258	ALA	2.0
1	B	258	ALA	2.0
1	B	260	ARG	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 ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	NAD	B	701	44/44	0.09	-0.16	14,18,24,25	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	NAD	F	901	44/44	0.09	-0.20	12,18,23,24	0
2	NAD	A	601	44/44	0.08	-0.22	14,18,24,25	0
2	NAD	E	801	44/44	0.09	-0.26	14,18,24,25	0

6.5 Other polymers

There are no such residues in this entry.