



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 25, 2018 – 11:30 AM EST

PDB ID : 1OC2
Title : The structure of NADH in the dTDP-D-glucose dehydratase (RmlB) enzyme
Authors : Beis, K.; Naismith, J.H.
Deposited on : 2003-02-04
Resolution : 1.50 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-20030736
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-20030736

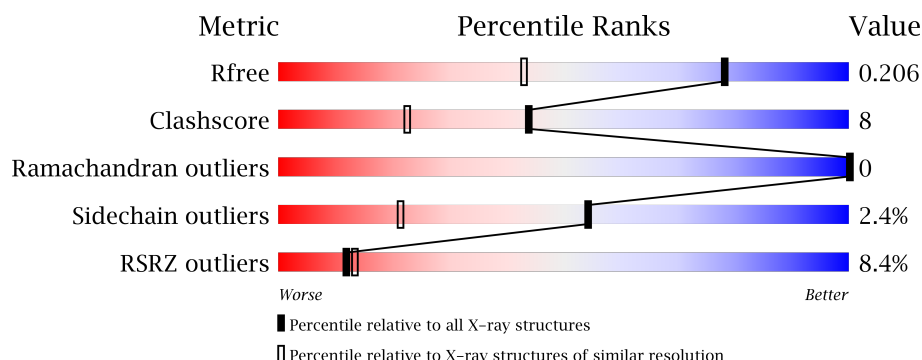
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.50 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	2279 (1.50-1.50)
Clashscore	112137	2503 (1.50-1.50)
Ramachandran outliers	110173	2445 (1.50-1.50)
Sidechain outliers	110143	2443 (1.50-1.50)
RSRZ outliers	101464	2305 (1.50-1.50)

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	348	
1	B	348	

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	TDX	A	1349	-	-	-	X
4	SO4	A	1352	-	-	X	X

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 6465 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.

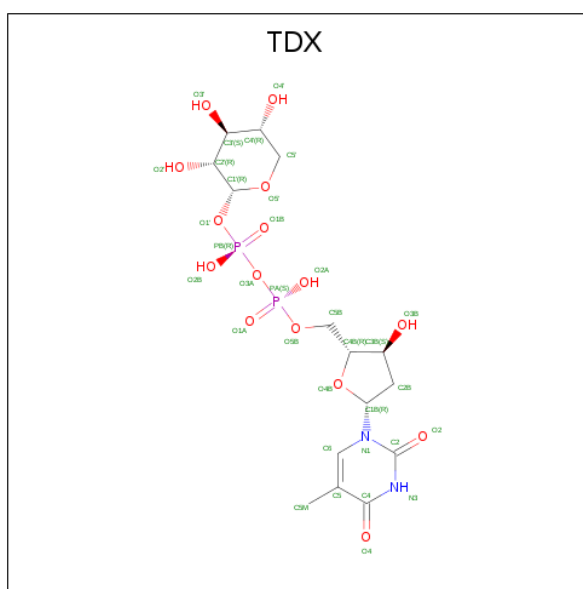
- Molecule 1 is a protein called DTDP-GLUCOSE 4,6-DEHYDRATASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	346	Total	C	N	O	S	0	13	0
			2787	1760	474	550	3			
1	B	347	Total	C	N	O	S	0	4	1
			2750	1737	471	539	3			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	313	GLU	GLN	conflict	UNP Q8GIP9
B	313	GLU	GLN	conflict	UNP Q8GIP9

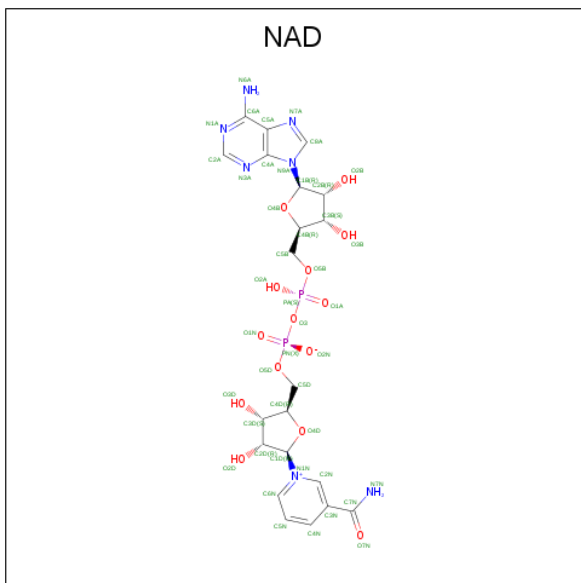
- Molecule 2 is THYMIDINE-5'-DIPHOSPHO-BETA-D-XYLOSE (three-letter code: TDX) (formula: C₁₅H₂₄N₂O₁₅P₂).



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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	B	1	Total	C	N	O	P	0	0
			34	15	2	15	2		

- Molecule 3 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (three-letter code: NAD) (formula: $C_{21}H_{27}N_7O_{14}P_2$).



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

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



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

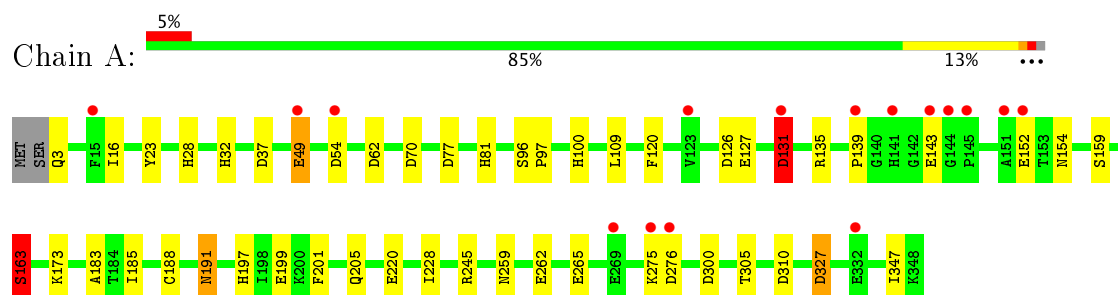
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	404	Total	O	0	0
			404	404		
5	B	353	Total	O	0	0
			353	353		

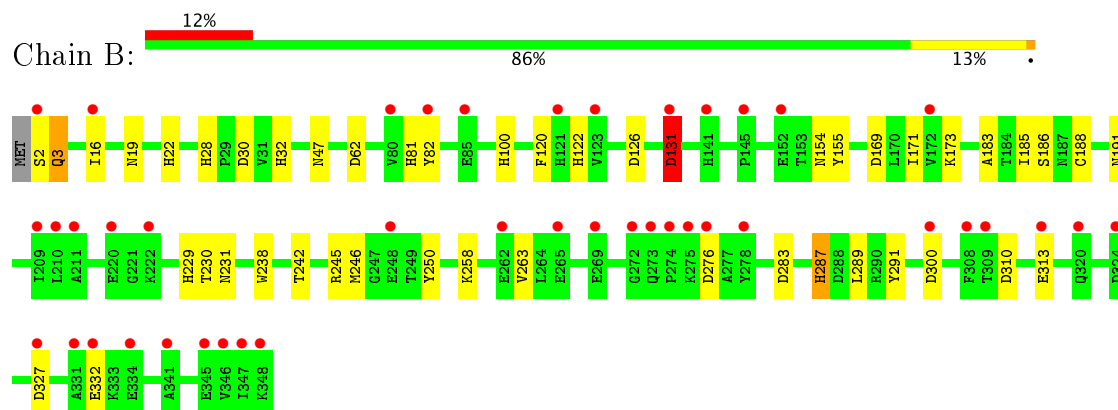
3 Residue-property plots [i](#)

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: DTDP-GLUCOSE 4,6-DEHYDRATASE



• Molecule 1: DTDP-GLUCOSE 4,6-DEHYDRATASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	61.41Å 94.83Å 183.54Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	91.29 – 1.50 34.74 – 1.50	Depositor EDS
% Data completeness (in resolution range)	98.3 (91.29-1.50) 98.3 (34.74-1.50)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.50 (at 1.50Å)	Xtriage
Refinement program	REFMAC 5.1.19	Depositor
R, R_{free}	0.170 , 0.200 0.183 , 0.206	Depositor DCC
R_{free} test set	16689 reflections (10.97%)	DCC
Wilson B-factor (Å ²)	15.1	Xtriage
Anisotropy	0.075	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 54.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6465	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.93% 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: TDX, SO4, 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	1.06	9/2924 (0.3%)	1.01	17/3973 (0.4%)
1	B	0.90	2/2838 (0.1%)	0.90	9/3860 (0.2%)
All	All	0.98	11/5762 (0.2%)	0.96	26/7833 (0.3%)

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	49[A]	GLU	CD-OE2	16.95	1.44	1.25
1	A	49[B]	GLU	CD-OE2	16.95	1.44	1.25
1	B	131[A]	ASP	CB-CG	-10.24	1.30	1.51
1	B	131[B]	ASP	CB-CG	-10.24	1.30	1.51
1	A	163[A]	SER	CB-OG	-9.14	1.30	1.42
1	A	163[B]	SER	CB-OG	-9.14	1.30	1.42
1	A	131[A]	ASP	CB-CG	-7.90	1.35	1.51
1	A	131[B]	ASP	CB-CG	-7.90	1.35	1.51
1	A	131[A]	ASP	CA-CB	6.10	1.67	1.53
1	A	131[B]	ASP	CA-CB	6.10	1.67	1.53
1	A	220	GLU	CD-OE2	5.04	1.31	1.25

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	49[A]	GLU	OE1-CD-OE2	-10.90	110.22	123.30
1	A	49[B]	GLU	OE1-CD-OE2	-10.90	110.22	123.30
1	A	276	ASP	CB-CG-OD2	9.26	126.64	118.30
1	B	276	ASP	CB-CG-OD2	8.50	125.95	118.30
1	A	54	ASP	CB-CG-OD2	8.04	125.54	118.30
1	A	300	ASP	CB-CG-OD2	7.97	125.48	118.30
1	A	126	ASP	CB-CG-OD1	7.32	124.89	118.30
1	A	3	GLN	CA-CB-CG	7.24	129.33	113.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	310	ASP	CB-CG-OD2	7.18	124.76	118.30
1	A	310	ASP	CB-CG-OD2	6.96	124.56	118.30
1	B	245	ARG	NE-CZ-NH1	6.91	123.75	120.30
1	B	300	ASP	CB-CG-OD2	6.83	124.45	118.30
1	B	327	ASP	CB-CG-OD2	6.38	124.04	118.30
1	A	131[A]	ASP	CB-CA-C	-6.03	98.34	110.40
1	A	131[B]	ASP	CB-CA-C	-6.03	98.34	110.40
1	A	131[A]	ASP	CB-CG-OD2	5.87	123.58	118.30
1	A	131[B]	ASP	CB-CG-OD2	5.87	123.58	118.30
1	B	30	ASP	CB-CG-OD1	5.81	123.53	118.30
1	B	291	TYR	CB-CG-CD2	-5.62	117.63	121.00
1	A	77	ASP	CB-CG-OD2	5.59	123.33	118.30
1	A	70	ASP	CB-CG-OD2	5.53	123.28	118.30
1	A	135	ARG	NE-CZ-NH1	5.46	123.03	120.30
1	A	327	ASP	CB-CG-OD2	5.38	123.15	118.30
1	B	126	ASP	CB-CG-OD1	5.21	122.99	118.30
1	B	245	ARG	NE-CZ-NH2	-5.16	117.72	120.30
1	A	37	ASP	CB-CG-OD2	5.14	122.93	118.30

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	2787	0	2655	32	0
1	B	2750	0	2627	44	0
2	A	34	0	22	3	0
2	B	34	0	22	2	0
3	A	44	0	26	4	0
3	B	44	0	26	4	0
4	A	15	0	0	4	0
5	A	404	0	0	16	0
5	B	353	0	0	19	0
All	All	6465	0	5378	84	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 8.

All (84) 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:19[B]:ASN:ND2	1:B:230:THR:HG23	1.58	1.18
1:A:265[A]:GLU:OE2	5:A:2278:HOH:O	1.60	1.18
1:A:265[A]:GLU:CD	5:A:2278:HOH:O	1.82	1.16
4:A:1352:SO4:O1	5:A:2398:HOH:O	1.71	1.07
1:A:131[B]:ASP:OD1	5:A:2169:HOH:O	1.71	1.07
4:A:1352:SO4:S	5:A:2398:HOH:O	2.17	1.02
1:B:154[B]:ASN:OD1	5:B:2196:HOH:O	1.80	0.97
1:B:19[B]:ASN:HD22	1:B:230:THR:HG23	1.14	0.97
4:A:1352:SO4:O1	5:A:2401:HOH:O	1.82	0.96
1:A:131[B]:ASP:OD2	5:A:2170:HOH:O	1.84	0.95
1:B:154[B]:ASN:OD1	5:B:2194:HOH:O	1.89	0.90
4:A:1352:SO4:S	5:A:2401:HOH:O	2.32	0.88
1:B:19[B]:ASN:HD22	1:B:230:THR:CG2	1.87	0.85
1:B:2:SER:OG	5:B:2001:HOH:O	1.95	0.84
1:B:19[B]:ASN:ND2	1:B:230:THR:CG2	2.42	0.82
1:B:19[B]:ASN:HD21	1:B:231:ASN:HD22	1.28	0.81
1:B:19[B]:ASN:CG	5:B:2015:HOH:O	2.20	0.81
1:B:19[A]:ASN:HB3	5:B:2015:HOH:O	1.82	0.79
1:B:229:HIS:HD2	1:B:231:ASN:H	1.31	0.78
1:B:22:HIS:HE1	1:B:47:ASN:O	1.68	0.77
1:B:2:SER:HA	5:B:2002:HOH:O	1.85	0.76
1:B:19[B]:ASN:ND2	5:B:2015:HOH:O	2.21	0.73
1:B:131[B]:ASP:OD1	5:B:2165:HOH:O	2.07	0.73
1:B:19[B]:ASN:HD21	1:B:231:ASN:ND2	1.88	0.72
1:A:131[B]:ASP:CG	5:A:2169:HOH:O	2.24	0.69
1:A:327:ASP:HB2	5:A:2352:HOH:O	1.98	0.64
1:B:3:GLN:NE2	1:B:242:THR:O	2.31	0.63
2:A:1349:TDX:H5D	5:A:2211:HOH:O	1.98	0.61
1:B:2:SER:CA	5:B:2002:HOH:O	2.45	0.60
3:B:1348:NAD:H4N	5:B:2218:HOH:O	2.02	0.59
1:A:81[A]:HIS:CE1	1:A:109:LEU:HD12	2.38	0.58
1:A:154:ASN:HA	1:B:154[B]:ASN:OD1	2.03	0.58
1:A:188:CYS:HB3	3:A:1350:NAD:H5N	1.84	0.57
1:B:283:ASP:OD1	1:B:287:HIS:HD2	1.87	0.57
1:B:131[B]:ASP:CG	5:B:2165:HOH:O	2.42	0.57
1:A:197:HIS:HD2	1:A:199:GLU:H	1.54	0.56
1:B:188:CYS:HB3	3:B:1348:NAD:H5N	1.88	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:16:ILE:HD11	3:A:1350:NAD:C7N	2.37	0.55
1:B:313:GLU:HG3	5:B:2312:HOH:O	2.06	0.54
1:A:173:LYS:HE2	1:A:185:ILE:HD12	1.89	0.54
1:B:131[A]:ASP:CG	1:B:289:LEU:H	2.12	0.53
1:B:173:LYS:HE2	1:B:185:ILE:HD12	1.92	0.51
1:A:305:THR:HG23	5:A:2329:HOH:O	2.10	0.51
1:A:81[A]:HIS:ND1	1:A:109:LEU:CD1	2.74	0.51
1:A:81[A]:HIS:CE1	1:A:109:LEU:CD1	2.94	0.50
3:B:1348:NAD:C3N	2:B:1349:TDX:H4'	2.42	0.50
1:B:155:TYR:OH	5:B:2199:HOH:O	2.18	0.50
1:B:122:HIS:HE1	1:B:169:ASP:OD1	1.95	0.50
1:A:127:GLU:OE1	2:A:1349:TDX:H5B	2.11	0.49
1:B:62:ASP:OD2	1:B:100:HIS:HE1	1.96	0.49
1:A:23:TYR:OH	1:A:28:HIS:HE1	1.95	0.49
1:B:131[A]:ASP:OD1	1:B:289:LEU:N	2.47	0.48
1:A:265[B]:GLU:HG2	5:A:2278:HOH:O	2.14	0.47
1:A:259:ASN:O	1:A:262[A]:GLU:HG2	2.15	0.47
1:B:122:HIS:CE1	1:B:169:ASP:OD1	2.67	0.47
1:A:275:LYS:O	5:A:2278:HOH:O	2.21	0.47
1:A:62:ASP:OD2	1:A:100:HIS:HE1	1.98	0.46
1:A:347:ILE:O	5:A:2385:HOH:O	2.21	0.46
1:B:173:LYS:CE	1:B:185:ILE:HD12	2.46	0.46
1:B:120:PHE:O	1:B:183:ALA:HA	2.17	0.45
1:A:32:HIS:HE1	5:A:2068:HOH:O	1.99	0.45
1:A:197:HIS:CD2	1:A:199:GLU:H	2.35	0.45
2:B:1349:TDX:H5D	5:B:2218:HOH:O	2.18	0.44
1:A:159:SER:O	1:A:163[A]:SER:HB2	2.18	0.44
1:B:19[B]:ASN:ND2	1:B:231:ASN:HD22	2.06	0.44
1:A:120:PHE:O	1:A:183:ALA:HA	2.17	0.44
1:B:16:ILE:HD11	3:B:1348:NAD:C7N	2.48	0.43
2:A:1349:TDX:H4'	3:A:1350:NAD:C3N	2.48	0.43
1:A:16:ILE:HD11	3:A:1350:NAD:N7N	2.34	0.43
1:A:173:LYS:CE	1:A:185:ILE:HD12	2.49	0.43
1:A:96[B]:SER:OG	1:A:97:PRO:HD3	2.19	0.42
1:B:32:HIS:HE1	5:B:2073:HOH:O	2.02	0.42
1:A:201:PHE:O	1:A:205:GLN:HG2	2.20	0.41
1:B:81:HIS:HD2	1:B:122:HIS:HD2	1.68	0.41
1:B:258:LYS:HD3	5:B:2269:HOH:O	2.20	0.41
1:B:28:HIS:CE1	1:B:238:TRP:CE2	3.08	0.41
1:B:28:HIS:HD2	5:B:2001:HOH:O	2.03	0.41
1:B:186:SER:HA	1:B:250:TYR:O	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:16:ILE:HG22	1:B:82:TYR:CE2	2.56	0.40
1:B:263:VAL:HA	5:B:2269:HOH:O	2.21	0.40
1:A:191:ASN:HA	1:A:228:ILE:O	2.21	0.40
1:A:163[A]:SER:HB3	1:B:171:ILE:HG22	2.02	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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	357/348 (103%)	348 (98%)	9 (2%)	0	100	100
1	B	349/348 (100%)	341 (98%)	8 (2%)	0	100	100
All	All	706/696 (101%)	689 (98%)	17 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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	302/291 (104%)	290 (96%)	12 (4%)	36	7
1	B	293/291 (101%)	286 (98%)	7 (2%)	54	21
All	All	595/582 (102%)	576 (97%)	19 (3%)	54	12

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	49[A]	GLU
1	A	49[B]	GLU
1	A	131[A]	ASP
1	A	131[B]	ASP
1	A	139	PRO
1	A	143	GLU
1	A	152[A]	GLU
1	A	152[B]	GLU
1	A	163[A]	SER
1	A	163[B]	SER
1	A	191	ASN
1	A	245	ARG
1	B	3	GLN
1	B	131[A]	ASP
1	B	131[B]	ASP
1	B	191	ASN
1	B	246	MET
1	B	287	HIS
1	B	332	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (28) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	28	HIS
1	A	32	HIS
1	A	100	HIS
1	A	122	HIS
1	A	141	HIS
1	A	196	GLN
1	A	197	HIS
1	A	231	ASN
1	A	260	ASN
1	A	280	HIS
1	A	307	GLN
1	A	339	ASN
1	B	22	HIS
1	B	28	HIS
1	B	32	HIS
1	B	87	HIS
1	B	100	HIS
1	B	122	HIS

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Mol	Chain	Res	Type
1	B	191	ASN
1	B	196	GLN
1	B	197	HIS
1	B	229	HIS
1	B	233	HIS
1	B	260	ASN
1	B	273	GLN
1	B	280	HIS
1	B	287	HIS
1	B	339	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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

7 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	TDX	A	1349	-	28,36,36	0.98	2 (7%)	39,55,55	2.23	6 (15%)
3	NAD	A	1350	-	41,48,48	2.03	6 (14%)	43,73,73	3.01	11 (25%)
4	SO4	A	1351	-	4,4,4	0.45	0	6,6,6	0.68	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	SO4	A	1352	-	4,4,4	0.25	0	6,6,6	0.34	0
4	SO4	A	1353	-	4,4,4	0.54	0	6,6,6	0.67	0
3	NAD	B	1348	-	41,48,48	2.34	11 (26%)	43,73,73	3.10	8 (18%)
2	TDX	B	1349	-	28,36,36	0.74	0	39,55,55	1.85	4 (10%)

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	TDX	A	1349	-	-	0/17/50/50	0/3/3/3
3	NAD	A	1350	-	-	0/22/62/62	0/5/5/5
4	SO4	A	1351	-	-	0/0/0/0	0/0/0/0
4	SO4	A	1352	-	-	0/0/0/0	0/0/0/0
4	SO4	A	1353	-	-	0/0/0/0	0/0/0/0
3	NAD	B	1348	-	-	0/22/62/62	0/5/5/5
2	TDX	B	1349	-	-	0/17/50/50	0/3/3/3

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	1348	NAD	C2B-C1B	-2.71	1.49	1.53
3	B	1348	NAD	O4D-C4D	-2.67	1.39	1.45
2	A	1349	TDX	C2-N3	-2.63	1.33	1.38
3	B	1348	NAD	C6N-C5N	-2.61	1.32	1.38
2	A	1349	TDX	C4'-C3'	-2.42	1.49	1.52
3	A	1350	NAD	C6N-C5N	-2.22	1.33	1.38
3	A	1350	NAD	C2A-N3A	2.00	1.35	1.32
3	B	1348	NAD	O4B-C1B	2.04	1.44	1.41
3	B	1348	NAD	C5A-C4A	2.09	1.45	1.40
3	B	1348	NAD	O3D-C3D	2.50	1.48	1.43
3	B	1348	NAD	C4A-N3A	2.69	1.39	1.35
3	B	1348	NAD	C7N-N7N	2.70	1.38	1.33
3	A	1350	NAD	C4A-N3A	3.35	1.40	1.35
3	A	1350	NAD	C6N-N1N	3.53	1.44	1.35
3	B	1348	NAD	C6N-N1N	4.03	1.45	1.35
3	A	1350	NAD	C5N-C4N	4.96	1.48	1.38
3	B	1348	NAD	C5N-C4N	5.93	1.50	1.38
3	A	1350	NAD	C4N-C3N	9.35	1.54	1.39
3	B	1348	NAD	C4N-C3N	9.94	1.55	1.39

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1348	NAD	C5N-C4N-C3N	-13.90	104.00	120.35
3	A	1350	NAD	C5N-C4N-C3N	-11.35	107.00	120.35
3	A	1350	NAD	O7N-C7N-C3N	-7.87	110.42	119.62
3	B	1348	NAD	O7N-C7N-C3N	-7.56	110.78	119.62
3	B	1348	NAD	N3A-C2A-N1A	-7.11	122.67	128.86
3	A	1350	NAD	N3A-C2A-N1A	-6.98	122.78	128.86
2	A	1349	TDX	C5-C4-N3	-6.79	117.76	125.24
2	B	1349	TDX	C5-C4-N3	-5.01	119.71	125.24
3	A	1350	NAD	C3N-C2N-N1N	-3.96	116.44	120.43
2	A	1349	TDX	C4'-C3'-C2'	-2.89	105.84	110.88
3	A	1350	NAD	C4B-O4B-C1B	-2.62	106.98	109.77
2	A	1349	TDX	O4'-C4'-C5'	-2.57	103.95	109.18
2	B	1349	TDX	O4B-C1B-N1	-2.41	103.72	107.78
3	B	1348	NAD	C4A-C5A-N7A	-2.33	107.16	109.41
3	B	1348	NAD	C5A-C6A-N6A	-2.07	116.25	120.47
2	A	1349	TDX	O4B-C1B-N1	-2.04	104.34	107.78
3	B	1348	NAD	C6N-C5N-C4N	2.13	122.66	119.44
3	B	1348	NAD	C2N-C3N-C4N	2.14	120.70	118.26
3	A	1350	NAD	C6N-C5N-C4N	2.28	122.88	119.44
2	A	1349	TDX	C2B-C1B-N1	2.33	119.74	114.23
2	B	1349	TDX	C5M-C5-C6	2.56	123.77	118.67
3	A	1350	NAD	O7N-C7N-N7N	2.62	126.31	122.58
3	A	1350	NAD	N6A-C6A-N1A	3.10	124.91	118.77
3	A	1350	NAD	C2A-N1A-C6A	3.59	125.05	118.77
3	A	1350	NAD	C3N-C7N-N7N	4.42	122.82	117.77
3	A	1350	NAD	C2N-C3N-C4N	6.45	125.62	118.26
3	B	1348	NAD	C3N-C7N-N7N	8.04	126.96	117.77
2	B	1349	TDX	C4-N3-C2	8.13	122.27	115.16
2	A	1349	TDX	C4-N3-C2	9.65	123.60	115.16

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5 monomers are involved in 15 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1349	TDX	3	0
3	A	1350	NAD	4	0
4	A	1352	SO4	4	0
3	B	1348	NAD	4	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	1349	TDX	2	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	346/348 (99%)	0.42	16 (4%) 33 37	10, 14, 23, 35	5 (1%)
1	B	347/348 (99%)	0.82	42 (12%) 5 5	12, 18, 28, 39	9 (2%)
All	All	693/696 (99%)	0.62	58 (8%) 12 13	10, 16, 27, 39	14 (2%)

All (58) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	121	HIS	7.7
1	B	272	GLY	6.9
1	B	85	GLU	5.4
1	A	145	PRO	5.3
1	B	141	HIS	5.2
1	B	275	LYS	5.1
1	A	141	HIS	4.2
1	B	334	GLU	4.2
1	A	139	PRO	3.9
1	B	2	SER	3.9
1	B	347	ILE	3.7
1	B	278	TYR	3.6
1	B	276	ASP	3.4
1	A	131[A]	ASP	3.4
1	B	123	VAL	3.4
1	A	275	LYS	3.2
1	B	332	GLU	3.1
1	A	332	GLU	3.0
1	B	262	GLU	3.0
1	B	331	ALA	3.0
1	B	327	ASP	2.9
1	B	16	ILE	2.9
1	B	274	PRO	2.8
1	B	265	GLU	2.8

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Mol	Chain	Res	Type	RSRZ
1	B	80	VAL	2.7
1	B	313	GLU	2.6
1	B	82	TYR	2.6
1	B	145	PRO	2.6
1	B	341	ALA	2.6
1	A	15	PHE	2.6
1	A	152[A]	GLU	2.6
1	B	273	GLN	2.5
1	B	348	LYS	2.5
1	A	276	ASP	2.5
1	B	308	PHE	2.4
1	B	269	GLU	2.4
1	A	123	VAL	2.4
1	B	209	ILE	2.4
1	A	143	GLU	2.3
1	B	222	LYS	2.3
1	B	248	GLU	2.3
1	A	151	ALA	2.3
1	B	210	LEU	2.2
1	A	144	GLY	2.2
1	B	220	GLU	2.2
1	B	345	GLU	2.2
1	B	172	VAL	2.2
1	B	152	GLU	2.2
1	B	320	GLN	2.1
1	A	54	ASP	2.1
1	A	49[A]	GLU	2.1
1	B	300	ASP	2.1
1	B	309	THR	2.1
1	B	211	ALA	2.1
1	A	269	GLU	2.0
1	B	324	ASP	2.0
1	B	346	VAL	2.0
1	B	131[A]	ASP	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 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(\AA^2)	Q<0.9
4	SO4	A	1352	5/5	0.75	0.79	12.57	90,90,90,90	0
2	TDX	A	1349	34/34	0.96	0.15	2.33	12,14,17,19	0
4	SO4	A	1353	5/5	0.98	0.18	1.42	29,31,33,37	0
3	NAD	A	1350	44/44	0.96	0.15	1.42	12,14,25,26	0
2	TDX	B	1349	34/34	0.95	0.12	0.73	18,20,22,25	0
3	NAD	B	1348	44/44	0.96	0.13	0.01	16,18,31,32	0
4	SO4	A	1351	5/5	0.97	0.15	-	26,26,30,30	0

6.5 Other polymers [i](#)

There are no such residues in this entry.