



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 07:44 PM GMT

PDB ID : 1GY8
Title : TRYPANOSOMA BRUCEI UDP-GALACTOSE 4' EPIMERASE
Authors : Shaw, M.P.; Bond, C.S.; Hunter, W.N.
Deposited on : 2002-04-21
Resolution : 2.00 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 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
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) : trunk26865

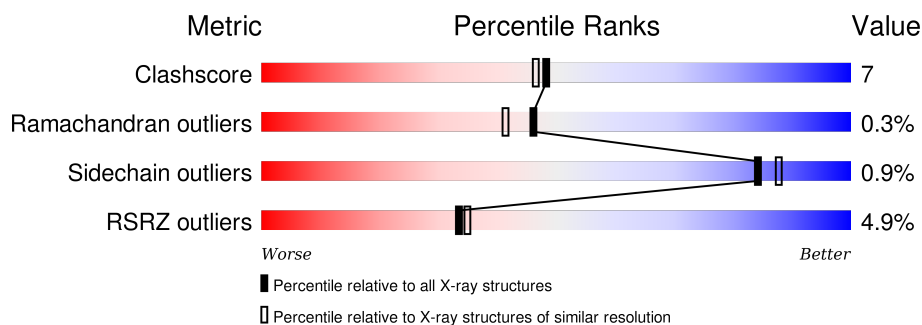
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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))
Clashscore	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)
RSRZ outliers	91569	6262 (2.00-2.00)

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	397	<div> <div>5%</div> <div>80% 12% • 8%</div> </div>
1	B	397	<div> <div>3%</div> <div>76% 15% • 8%</div> </div>
1	C	397	<div> <div>6%</div> <div>78% 14% • 7%</div> </div>
1	D	397	<div> <div>4%</div> <div>79% 12% • 8%</div> </div>

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	NAD	A	1382	-	-	-	X
2	NAD	B	1382	-	-	-	X
2	NAD	C	1382	-	-	-	X
2	NAD	D	1382	-	-	-	X

2 Entry composition [i](#)

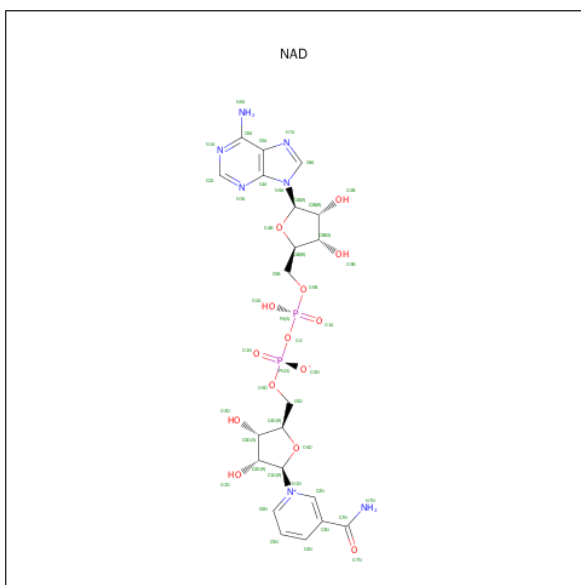
There are 4 unique types of molecules in this entry. The entry contains 12972 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 UDP-GALACTOSE 4-EPIMERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	367	Total	C	N	O	S	8	0	0
			2856	1798	509	532	17			
1	B	366	Total	C	N	O	S	13	0	0
			2848	1793	508	530	17			
1	C	370	Total	C	N	O	S	9	0	0
			2874	1808	512	537	17			
1	D	364	Total	C	N	O	S	12	0	0
			2833	1784	506	527	16			

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



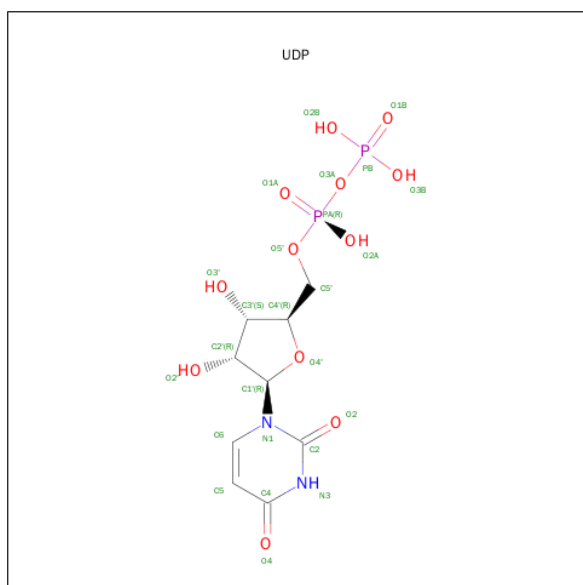
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	B	1	Total	C	N	O	P	0	0
			44	21	7	14	2		

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

- Molecule 3 is URIDINE-5'-DIPHOSPHATE (three-letter code: UDP) (formula: $C_9H_{14}N_2O_{12}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			25	9	2	12	2		
3	B	1	Total	C	N	O	P	0	0
			25	9	2	12	2		
3	C	1	Total	C	N	O	P	0	0
			25	9	2	12	2		
3	D	1	Total	C	N	O	P	0	0
			25	9	2	12	2		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	337	Total	O	0	0
			337	337		
4	B	350	Total	O	0	0
			350	350		
4	C	345	Total	O	0	0
			345	345		

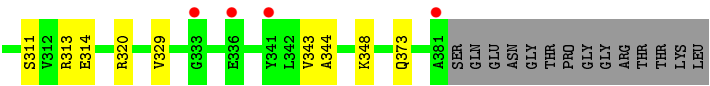
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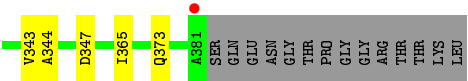
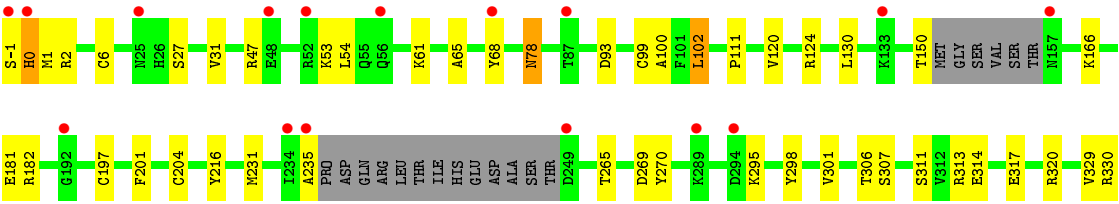
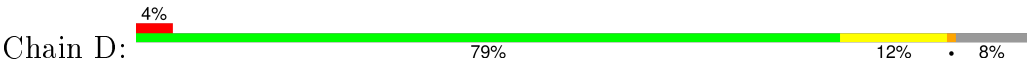
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	D	253	Total 253	O 253	0	0

• Molecule 1: UDP-GALACTOSE 4-EPIMERASE





● Molecule 1: UDP-GALACTOSE 4-EPIMERASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	102.19Å 112.53Å 160.59Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	100.00 – 2.00 20.04 – 2.00	Depositor EDS
% Data completeness (in resolution range)	89.5 (100.00-2.00) 85.3 (20.04-2.00)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.91 (at 2.01Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.216 , 0.245 0.229 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	22.1	Xtriage
Anisotropy	0.418	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 51.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	2 of 205073 reflections (0.001%)	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	12972	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 47.41 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 9.9652e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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: UDP, 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	0.32	0/2921	0.58	0/3956
1	B	0.32	0/2912	0.59	0/3943
1	C	0.33	0/2939	0.59	0/3981
1	D	0.32	0/2897	0.58	0/3923
All	All	0.32	0/11669	0.59	0/15803

There are no bond length outliers.

There are no bond angle outliers.

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	2856	0	2807	35	0
1	B	2848	0	2803	53	0
1	C	2874	0	2824	45	0
1	D	2833	0	2787	38	0
2	A	44	0	26	0	0
2	B	44	0	26	0	0
2	C	44	0	26	0	0
2	D	44	0	26	0	0
3	A	25	0	11	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	25	0	11	0	0
3	C	25	0	11	0	0
3	D	25	0	11	0	0
4	A	337	0	0	4	0
4	B	350	0	0	6	0
4	C	345	0	0	7	0
4	D	253	0	0	2	0
All	All	12972	0	11369	167	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 7.

All (167) 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:156:THR:HG23	1:B:157:ASN:H	1.35	0.91
1:B:308:ARG:HD2	1:B:360:ASP:OD1	1.79	0.82
1:B:159:GLU:HG3	1:B:160:PRO:HD2	1.65	0.78
1:B:56:GLN:HG2	4:B:2073:HOH:O	1.85	0.76
1:C:265:THR:HG23	1:C:313:ARG:NH1	2.02	0.74
1:B:54:LEU:HD21	1:B:61:LYS:HE2	1.68	0.74
1:D:150:THR:OG1	1:D:166:LYS:HD3	1.93	0.69
1:C:78:ASN:C	1:C:78:ASN:HD22	1.95	0.69
1:B:78:ASN:HD22	1:B:78:ASN:C	1.97	0.69
1:B:371:LYS:HA	1:B:374:ARG:NH1	2.08	0.68
1:B:6:CYS:HA	1:B:31:VAL:HG13	1.74	0.68
1:D:78:ASN:C	1:D:78:ASN:HD22	1.94	0.68
1:A:294:ASP:O	1:A:296:SER:N	2.27	0.67
1:B:265:THR:HG23	1:B:313:ARG:NH1	2.09	0.67
1:D:265:THR:HG23	1:D:313:ARG:NH1	2.10	0.67
1:A:204:CYS:SG	1:A:270:TYR:HB3	2.36	0.66
1:D:6:CYS:HA	1:D:31:VAL:HG13	1.78	0.65
1:C:49:ASN:ND2	1:C:52:ARG:HH22	1.94	0.65
1:B:216:TYR:H	1:B:373:GLN:HE22	1.43	0.65
1:A:216:TYR:H	1:A:373:GLN:HE22	1.44	0.65
1:D:216:TYR:H	1:D:373:GLN:HE22	1.43	0.65
1:B:265:THR:HG23	1:B:313:ARG:HH11	1.62	0.65
1:C:348:LYS:HE3	4:C:2301:HOH:O	1.96	0.64
1:D:204:CYS:SG	1:D:365:ILE:HG21	2.38	0.64
1:C:216:TYR:H	1:C:373:GLN:HE22	1.44	0.64
1:D:47:ARG:HD3	1:D:68:TYR:CE2	2.34	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:52:ARG:O	1:C:56:GLN:HG3	2.00	0.62
1:B:204:CYS:SG	1:B:365:ILE:HG21	2.40	0.62
1:B:156:THR:HG23	1:B:157:ASN:N	2.14	0.61
1:D:295:LYS:HA	1:D:298:TYR:CD1	2.36	0.61
1:A:78:ASN:C	1:A:78:ASN:HD22	2.04	0.60
1:B:204:CYS:SG	1:B:270:TYR:HB3	2.42	0.60
1:C:204:CYS:SG	1:C:270:TYR:HB3	2.41	0.60
1:A:54:LEU:HD22	1:A:65:ALA:HB1	1.84	0.60
1:D:166:LYS:HE2	4:D:2134:HOH:O	2.02	0.59
1:D:317:GLU:OE1	1:D:320:ARG:NH1	2.34	0.59
1:C:4:LEU:HD11	1:C:31:VAL:HG23	1.85	0.59
1:B:72:GLU:OE2	1:B:88:ARG:NH2	2.36	0.59
1:A:191:TYR:CZ	1:B:111:PRO:HD2	2.38	0.58
1:D:182:ARG:HH11	1:D:182:ARG:HG3	1.67	0.58
1:A:265:THR:HG23	1:A:313:ARG:NH1	2.18	0.58
1:B:347:ASP:O	1:B:351:GLU:HG3	2.05	0.57
1:D:265:THR:HG23	1:D:313:ARG:HH11	1.69	0.57
1:C:285:ASP:HB2	4:C:2246:HOH:O	2.03	0.57
1:A:251:ARG:HG2	1:A:328:PRO:HB2	1.86	0.57
1:D:54:LEU:HD11	1:D:61:LYS:HG2	1.87	0.56
1:C:99:CYS:O	1:C:100:ALA:HB2	2.04	0.56
1:A:132:HIS:HE1	4:A:2094:HOH:O	1.88	0.56
1:B:330:ARG:HG3	1:B:330:ARG:O	2.06	0.56
1:C:163:ILE:HD13	1:C:348:LYS:HG3	1.87	0.55
1:B:52:ARG:O	1:B:56:GLN:HG3	2.07	0.55
1:A:131:LEU:HD23	4:A:2094:HOH:O	2.06	0.55
1:D:269:ASP:HB2	1:D:343:VAL:HA	1.89	0.55
1:A:263:ASP:O	1:A:313:ARG:NH1	2.40	0.54
1:D:231:MET:O	1:D:235:ALA:N	2.39	0.54
1:C:132:HIS:HD2	4:C:2158:HOH:O	1.90	0.54
1:A:269:ASP:HB2	1:A:343:VAL:HA	1.89	0.54
1:C:269:ASP:HB2	1:C:343:VAL:HA	1.88	0.54
1:B:306:THR:O	1:B:307:SER:HB2	2.07	0.54
1:D:329:VAL:CG1	1:D:330:ARG:N	2.71	0.54
1:B:269:ASP:HB2	1:B:343:VAL:HA	1.90	0.53
1:C:311:SER:OG	1:C:314:GLU:HG3	2.08	0.53
1:B:329:VAL:CG1	1:B:330:ARG:N	2.71	0.53
1:A:150:THR:OG1	1:A:166:LYS:HD3	2.08	0.52
1:D:182:ARG:HG3	1:D:182:ARG:NH1	2.24	0.52
1:B:157:ASN:HA	4:B:2170:HOH:O	2.09	0.52
1:D:78:ASN:C	1:D:78:ASN:ND2	2.63	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:313:ARG:NH2	1:B:331:GLU:OE2	2.42	0.52
1:A:307:SER:HA	1:A:344:ALA:O	2.10	0.52
1:B:104:VAL:HG12	4:B:2300:HOH:O	2.09	0.52
1:D:1:MET:HG3	1:D:93:ASP:CB	2.41	0.51
1:D:204:CYS:SG	1:D:270:TYR:HB3	2.51	0.51
1:D:99:CYS:O	1:D:100:ALA:HB2	2.10	0.51
1:B:1:MET:HG3	1:B:93:ASP:CB	2.41	0.51
1:C:215:HIS:NE2	1:C:217:GLN:NE2	2.58	0.51
1:C:1:MET:HG3	1:C:93:ASP:CB	2.41	0.51
1:D:201:PHE:CZ	1:D:344:ALA:HB2	2.46	0.51
1:C:201:PHE:CE1	1:C:344:ALA:HB2	2.46	0.50
1:D:311:SER:OG	1:D:314:GLU:HG3	2.11	0.50
1:B:374:ARG:NH1	4:B:2330:HOH:O	2.44	0.50
1:B:201:PHE:CZ	1:B:344:ALA:HB2	2.47	0.50
1:A:1:MET:HG3	1:A:93:ASP:CB	2.41	0.50
1:B:99:CYS:O	1:B:100:ALA:HB2	2.11	0.50
1:B:294:ASP:HA	4:B:2261:HOH:O	2.11	0.50
1:A:201:PHE:CE1	1:A:344:ALA:HB2	2.47	0.49
1:A:311:SER:OG	1:A:314:GLU:HG3	2.12	0.49
1:B:311:SER:OG	1:B:314:GLU:HG3	2.12	0.49
1:B:201:PHE:CE1	1:B:344:ALA:HB2	2.47	0.49
1:A:201:PHE:CZ	1:A:344:ALA:HB2	2.47	0.49
1:B:231:MET:O	1:B:235:ALA:N	2.46	0.49
1:C:215:HIS:CD2	1:C:217:GLN:HB2	2.48	0.49
1:B:78:ASN:C	1:B:78:ASN:ND2	2.65	0.49
1:D:201:PHE:CE1	1:D:344:ALA:HB2	2.47	0.49
1:C:201:PHE:CZ	1:C:344:ALA:HB2	2.47	0.49
1:A:6:CYS:HA	1:A:31:VAL:HG13	1.94	0.49
1:A:49:ASN:HB3	4:A:2056:HOH:O	2.12	0.48
1:C:78:ASN:C	1:C:78:ASN:ND2	2.64	0.48
1:C:4:LEU:HD11	1:C:31:VAL:CG2	2.44	0.47
1:A:111:PRO:HD2	1:B:191:TYR:CZ	2.50	0.47
1:C:191:TYR:CZ	1:D:111:PRO:HD2	2.49	0.47
1:C:150:THR:OG1	1:C:166:LYS:HD3	2.15	0.47
1:D:54:LEU:HD21	1:D:61:LYS:HE2	1.96	0.47
1:A:55:GLN:CD	1:A:56:GLN:N	2.69	0.47
1:B:-1:SER:O	1:B:0:HIS:C	2.53	0.46
1:A:120:VAL:O	1:A:124:ARG:HG2	2.15	0.46
1:B:159:GLU:HG3	1:B:160:PRO:CD	2.40	0.46
1:A:306:THR:O	1:A:307:SER:HB2	2.16	0.46
1:C:72:GLU:OE2	1:C:88:ARG:NH2	2.49	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:62:PRO:HD2	1:B:65:ALA:HB2	1.98	0.46
1:D:120:VAL:O	1:D:124:ARG:HG2	2.16	0.46
1:B:54:LEU:HD11	1:B:61:LYS:HG2	1.98	0.45
1:C:120:VAL:O	1:C:124:ARG:HG2	2.16	0.45
1:A:99:CYS:O	1:A:100:ALA:HB2	2.17	0.45
1:A:252:MET:HA	1:A:253:PRO:HD3	1.82	0.45
1:B:197:CYS:HB2	1:B:301:VAL:HG12	1.99	0.45
1:D:197:CYS:HB2	1:D:301:VAL:HG12	1.99	0.45
1:C:307:SER:HA	1:C:344:ALA:O	2.17	0.44
1:C:295:LYS:HA	1:C:298:TYR:CD1	2.52	0.44
1:B:156:THR:O	1:B:157:ASN:HB3	2.17	0.44
1:B:147:GLY:HA3	1:B:168:SER:O	2.17	0.44
1:C:262:PRO:HD3	4:C:2109:HOH:O	2.16	0.44
1:B:72:GLU:CD	1:B:88:ARG:HH22	2.21	0.44
1:D:306:THR:O	1:D:307:SER:HB2	2.18	0.44
1:A:163:ILE:HD13	1:A:348:LYS:HG3	2.00	0.44
1:C:265:THR:HG23	1:C:313:ARG:HH12	1.81	0.43
1:C:263:ASP:O	1:C:313:ARG:NH1	2.52	0.43
1:B:102:LEU:HD23	1:B:102:LEU:N	2.33	0.43
1:C:57:SER:HB2	4:C:2068:HOH:O	2.17	0.43
1:B:120:VAL:O	1:B:124:ARG:HG2	2.18	0.43
1:C:47:ARG:HG2	1:C:48:GLU:OE2	2.18	0.43
1:A:197:CYS:HB2	1:A:301:VAL:HG12	2.00	0.43
1:A:295:LYS:HA	1:A:298:TYR:CD1	2.53	0.43
1:D:330:ARG:HG3	1:D:330:ARG:O	2.18	0.43
1:D:-1:SER:O	1:D:0:HIS:C	2.56	0.43
1:C:102:LEU:N	1:C:102:LEU:HD23	2.34	0.43
1:B:130:LEU:HA	1:B:130:LEU:HD12	1.92	0.43
1:B:1:MET:HG3	1:B:93:ASP:HB2	2.00	0.42
1:C:49:ASN:HD22	1:C:52:ARG:HH22	1.66	0.42
1:A:78:ASN:ND2	1:A:78:ASN:C	2.72	0.42
1:D:1:MET:HG3	1:D:93:ASP:HB2	1.99	0.42
1:C:197:CYS:HB2	1:C:301:VAL:HG12	2.01	0.42
1:C:114:TYR:O	1:C:118:ASN:HB2	2.20	0.42
1:C:215:HIS:HD2	1:C:217:GLN:HB2	1.84	0.42
1:A:1:MET:HG3	1:A:93:ASP:HB2	2.02	0.42
1:A:102:LEU:HD23	1:A:102:LEU:N	2.34	0.42
1:C:320:ARG:NH2	1:C:329:VAL:HG13	2.35	0.42
1:C:236:PRO:O	1:C:237:ASP:C	2.58	0.42
1:A:111:PRO:HD2	1:B:191:TYR:OH	2.20	0.42
1:B:307:SER:HA	1:B:344:ALA:O	2.19	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:54:LEU:HB2	4:A:2065:HOH:O	2.20	0.41
1:D:61:LYS:HB3	1:D:65:ALA:CB	2.50	0.41
1:C:181:GLU:HG2	1:C:197:CYS:SG	2.61	0.41
1:B:181:GLU:HG2	1:B:197:CYS:SG	2.61	0.41
1:C:50:VAL:O	1:C:54:LEU:HB2	2.21	0.41
1:D:53:LYS:NZ	4:D:2053:HOH:O	2.49	0.41
1:C:147:GLY:HA3	1:C:168:SER:O	2.20	0.41
1:D:102:LEU:N	1:D:102:LEU:HD23	2.34	0.41
1:C:1:MET:HG3	1:C:93:ASP:HB2	2.02	0.41
1:D:181:GLU:HG2	1:D:197:CYS:SG	2.60	0.41
1:D:130:LEU:HD12	1:D:130:LEU:HA	1.92	0.41
1:C:79:GLU:HB2	4:C:2093:HOH:O	2.21	0.41
1:A:130:LEU:HA	1:A:130:LEU:HD12	1.91	0.40
1:B:163:ILE:HG13	1:B:164:ASN:N	2.36	0.40
1:C:308:ARG:HD2	4:C:2264:HOH:O	2.21	0.40
1:B:53:LYS:HG3	4:B:2060:HOH:O	2.21	0.40
1:B:114:TYR:O	1:B:118:ASN:HB2	2.21	0.40
1:D:2:ARG:HD3	1:D:27:SER:HB2	2.04	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	361/397 (91%)	348 (96%)	12 (3%)	1 (0%)	46	41
1	B	360/397 (91%)	348 (97%)	10 (3%)	2 (1%)	30	22
1	C	364/397 (92%)	356 (98%)	7 (2%)	1 (0%)	46	41
1	D	358/397 (90%)	350 (98%)	7 (2%)	1 (0%)	46	41
All	All	1443/1588 (91%)	1402 (97%)	36 (2%)	5 (0%)	46	41

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	295	LYS
1	B	0	HIS
1	D	0	HIS
1	B	100	ALA
1	C	100	ALA

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	306/331 (92%)	303 (99%)	3 (1%)	82	85
1	B	305/331 (92%)	303 (99%)	2 (1%)	88	91
1	C	308/331 (93%)	305 (99%)	3 (1%)	82	85
1	D	303/331 (92%)	300 (99%)	3 (1%)	82	85
All	All	1222/1324 (92%)	1211 (99%)	11 (1%)	84	88

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	78	ASN
1	A	80	ASP
1	A	102	LEU
1	B	78	ASN
1	B	102	LEU
1	C	78	ASN
1	C	102	LEU
1	C	125	LEU
1	D	78	ASN
1	D	102	LEU
1	D	347	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	25	ASN

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Mol	Chain	Res	Type
1	A	49	ASN
1	A	55	GLN
1	A	78	ASN
1	A	83	ASN
1	A	132	HIS
1	A	373	GLN
1	B	49	ASN
1	B	78	ASN
1	B	83	ASN
1	B	132	HIS
1	B	373	GLN
1	C	49	ASN
1	C	78	ASN
1	C	83	ASN
1	C	132	HIS
1	C	217	GLN
1	C	373	GLN
1	D	49	ASN
1	D	56	GLN
1	D	78	ASN
1	D	83	ASN
1	D	132	HIS
1	D	373	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 ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

8 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	1382	-	38,48,48	1.53	6 (15%)	47,73,73	1.87	4 (8%)
3	UDP	A	1383	-	18,26,26	2.24	5 (27%)	26,40,40	1.70	4 (15%)
2	NAD	B	1382	-	38,48,48	1.60	7 (18%)	47,73,73	1.89	4 (8%)
3	UDP	B	1383	-	18,26,26	2.21	5 (27%)	26,40,40	1.68	4 (15%)
2	NAD	C	1382	-	38,48,48	1.56	6 (15%)	47,73,73	1.88	4 (8%)
3	UDP	C	1383	-	18,26,26	2.21	5 (27%)	26,40,40	1.73	4 (15%)
2	NAD	D	1382	-	38,48,48	1.63	7 (18%)	47,73,73	1.88	4 (8%)
3	UDP	D	1383	-	18,26,26	2.23	5 (27%)	26,40,40	1.68	4 (15%)

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	1382	-	-	0/22/62/62	0/5/5/5
3	UDP	A	1383	-	-	0/12/32/32	0/2/2/2
2	NAD	B	1382	-	-	0/22/62/62	0/5/5/5
3	UDP	B	1383	-	-	0/12/32/32	0/2/2/2
2	NAD	C	1382	-	-	0/22/62/62	0/5/5/5
3	UDP	C	1383	-	-	0/12/32/32	0/2/2/2
2	NAD	D	1382	-	-	0/22/62/62	0/5/5/5
3	UDP	D	1383	-	-	0/12/32/32	0/2/2/2

All (46) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1382	NAD	C2A-N3A	2.02	1.35	1.32
2	D	1382	NAD	O4B-C1B	2.02	1.43	1.41
2	B	1382	NAD	C4A-N3A	2.07	1.38	1.35
2	C	1382	NAD	C4A-N3A	2.08	1.38	1.35
2	A	1382	NAD	C4A-N3A	2.21	1.38	1.35
3	B	1383	UDP	PA-O1A	2.49	1.60	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	1383	UDP	PA-O1A	2.50	1.60	1.51
2	D	1382	NAD	C4A-N3A	2.64	1.39	1.35
3	D	1383	UDP	PA-O1A	2.66	1.60	1.51
3	A	1383	UDP	PA-O1A	2.68	1.61	1.51
2	A	1382	NAD	C6N-N1N	2.87	1.43	1.35
2	C	1382	NAD	C3N-C7N	2.90	1.55	1.50
2	A	1382	NAD	C5N-C4N	2.90	1.44	1.38
2	C	1382	NAD	C6N-N1N	2.95	1.43	1.35
2	B	1382	NAD	C5N-C4N	2.98	1.45	1.38
3	C	1383	UDP	PB-O1B	2.98	1.61	1.51
2	C	1382	NAD	C5N-C4N	3.01	1.45	1.38
2	A	1382	NAD	C4N-C3N	3.04	1.44	1.39
3	B	1383	UDP	PB-O1B	3.04	1.61	1.51
2	D	1382	NAD	C5N-C4N	3.05	1.45	1.38
3	D	1383	UDP	PB-O1B	3.05	1.61	1.51
3	A	1383	UDP	PB-O1B	3.06	1.61	1.51
2	C	1382	NAD	C4N-C3N	3.07	1.44	1.39
2	D	1382	NAD	C4N-C3N	3.08	1.44	1.39
2	D	1382	NAD	C6N-N1N	3.10	1.43	1.35
2	A	1382	NAD	C3N-C7N	3.10	1.55	1.50
2	B	1382	NAD	C6N-N1N	3.14	1.43	1.35
2	B	1382	NAD	C3N-C7N	3.20	1.55	1.50
2	B	1382	NAD	C4N-C3N	3.29	1.44	1.39
2	D	1382	NAD	C3N-C7N	3.33	1.55	1.50
3	B	1383	UDP	O4'-C1'	3.94	1.46	1.41
3	D	1383	UDP	O4'-C1'	3.99	1.46	1.41
3	A	1383	UDP	O4'-C1'	4.14	1.46	1.41
3	C	1383	UDP	C4-N3	4.27	1.41	1.33
3	C	1383	UDP	C6-N1	4.53	1.42	1.35
3	A	1383	UDP	C4-N3	4.54	1.41	1.33
2	A	1382	NAD	O4D-C1D	4.59	1.47	1.41
3	C	1383	UDP	O4'-C1'	4.60	1.47	1.41
3	D	1383	UDP	C6-N1	4.61	1.42	1.35
3	D	1383	UDP	C4-N3	4.62	1.41	1.33
3	B	1383	UDP	C6-N1	4.63	1.42	1.35
3	A	1383	UDP	C6-N1	4.67	1.42	1.35
3	B	1383	UDP	C4-N3	4.68	1.41	1.33
2	D	1382	NAD	O4D-C1D	4.69	1.47	1.41
2	B	1382	NAD	O4D-C1D	4.86	1.47	1.41
2	C	1382	NAD	O4D-C1D	5.01	1.47	1.41

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1382	NAD	N3A-C2A-N1A	-10.69	120.71	128.89
2	D	1382	NAD	N3A-C2A-N1A	-10.67	120.72	128.89
2	C	1382	NAD	N3A-C2A-N1A	-10.65	120.74	128.89
2	A	1382	NAD	N3A-C2A-N1A	-10.64	120.75	128.89
3	D	1383	UDP	PA-O3A-PB	-2.73	123.52	132.67
3	B	1383	UDP	PA-O3A-PB	-2.68	123.67	132.67
3	C	1383	UDP	PA-O3A-PB	-2.65	123.77	132.67
3	A	1383	UDP	PA-O3A-PB	-2.64	123.81	132.67
2	D	1382	NAD	C1B-N9A-C4A	-2.63	122.97	126.94
2	B	1382	NAD	C1B-N9A-C4A	-2.60	123.02	126.94
2	A	1382	NAD	C1B-N9A-C4A	-2.50	123.17	126.94
2	C	1382	NAD	C1B-N9A-C4A	-2.48	123.20	126.94
2	D	1382	NAD	C4A-C5A-N7A	2.20	111.50	109.48
2	C	1382	NAD	C4A-C5A-N7A	2.25	111.55	109.48
2	A	1382	NAD	C4A-C5A-N7A	2.29	111.59	109.48
2	B	1382	NAD	C4A-C5A-N7A	2.33	111.63	109.48
3	B	1383	UDP	O4'-C1'-N1	2.59	113.53	108.08
3	A	1383	UDP	O4'-C1'-N1	2.62	113.61	108.08
3	D	1383	UDP	O4'-C1'-N1	2.63	113.62	108.08
3	C	1383	UDP	O4'-C1'-N1	2.65	113.67	108.08
2	A	1382	NAD	O4D-C1D-N1N	2.97	111.40	108.13
3	D	1383	UDP	O3A-PA-O5'	3.11	111.18	102.94
3	B	1383	UDP	O3A-PA-O5'	3.14	111.27	102.94
2	D	1382	NAD	O4D-C1D-N1N	3.15	111.59	108.13
2	C	1382	NAD	O4D-C1D-N1N	3.18	111.62	108.13
2	B	1382	NAD	O4D-C1D-N1N	3.18	111.63	108.13
3	C	1383	UDP	O3A-PA-O5'	3.25	111.56	102.94
3	A	1383	UDP	O3A-PA-O5'	3.33	111.78	102.94
3	D	1383	UDP	C4-N3-C2	6.22	120.30	114.14
3	A	1383	UDP	C4-N3-C2	6.28	120.36	114.14
3	B	1383	UDP	C4-N3-C2	6.29	120.37	114.14
3	C	1383	UDP	C4-N3-C2	6.47	120.55	114.14

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	367/397 (92%)	0.15	20 (5%)	29 31	22, 30, 57, 85	2 (0%)
1	B	366/397 (92%)	0.10	13 (3%)	46 48	21, 31, 51, 82	3 (0%)
1	C	370/397 (93%)	0.24	22 (5%)	26 27	20, 30, 53, 76	3 (0%)
1	D	364/397 (91%)	0.24	17 (4%)	35 37	23, 33, 53, 73	3 (0%)
All	All	1467/1588 (92%)	0.18	72 (4%)	33 35	20, 31, 54, 85	11 (0%)

All (72) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	157	ASN	7.6
1	C	150	THR	7.3
1	D	157	ASN	7.0
1	C	237	ASP	7.0
1	B	381	ALA	6.9
1	A	56	GLN	6.4
1	A	237	ASP	6.2
1	A	150	THR	6.0
1	B	156	THR	5.7
1	B	0	HIS	5.4
1	A	151	MET	5.1
1	D	381	ALA	5.0
1	C	157	ASN	5.0
1	B	151	MET	4.9
1	C	151	MET	4.5
1	A	58	ASP	4.4
1	C	60	PRO	4.4
1	C	236	PRO	4.4
1	A	54	LEU	4.3
1	B	235	ALA	4.3
1	C	-1	SER	4.0

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Mol	Chain	Res	Type	RSRZ
1	A	157	ASN	4.0
1	D	235	ALA	3.8
1	A	60	PRO	3.8
1	A	250	LYS	3.8
1	C	248	THR	3.7
1	A	57	SER	3.6
1	C	246	ALA	3.6
1	C	333	GLY	3.5
1	C	59	GLY	3.5
1	A	249	ASP	3.5
1	C	247	SER	3.4
1	C	56	GLN	3.3
1	A	294	ASP	3.3
1	D	-1	SER	3.2
1	D	25	ASN	3.2
1	B	25	ASN	3.1
1	B	-1	SER	3.1
1	C	258	ASP	3.0
1	D	87	THR	3.0
1	D	249	ASP	3.0
1	B	234	ILE	2.9
1	A	328	PRO	2.9
1	A	52	ARG	2.9
1	D	133	LYS	2.8
1	D	294	ASP	2.8
1	D	48	GLU	2.7
1	A	381	ALA	2.7
1	C	158	ALA	2.7
1	B	330	ARG	2.7
1	B	52	ARG	2.6
1	C	381	ALA	2.5
1	C	336	GLU	2.5
1	C	293	ASN	2.5
1	C	341	TYR	2.5
1	D	56	GLN	2.4
1	A	59	GLY	2.4
1	A	230	VAL	2.4
1	D	192	GLY	2.4
1	D	52	ARG	2.3
1	D	0	HIS	2.3
1	A	-1	SER	2.2
1	A	55	GLN	2.2

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Mol	Chain	Res	Type	RSRZ
1	D	234	ILE	2.1
1	D	289	LYS	2.1
1	B	249	ASP	2.1
1	C	292	PRO	2.1
1	D	68	TYR	2.1
1	C	260	PRO	2.1
1	A	236	PRO	2.1
1	C	159	GLU	2.0
1	B	159	GLU	2.0

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
2	NAD	C	1382	44/44	0.89	0.23	8.24	21,23,25,25	18
2	NAD	B	1382	44/44	0.86	0.22	6.82	22,24,26,26	18
2	NAD	D	1382	44/44	0.82	0.23	6.53	24,27,28,29	18
2	NAD	A	1382	44/44	0.87	0.23	6.05	21,23,24,26	18
3	UDP	B	1383	25/25	0.98	0.11	0.15	23,27,29,30	0
3	UDP	C	1383	25/25	0.94	0.13	-0.06	28,29,31,34	0
3	UDP	D	1383	25/25	0.97	0.11	-0.18	24,27,29,31	0
3	UDP	A	1383	25/25	0.96	0.10	-0.78	24,28,30,31	0

6.5 Other polymers [i](#)

There are no such residues in this entry.