



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 26, 2018 – 01:54 PM EST

PDB ID : 2C59
Title : gdp-mannose-3', 5' -epimerase (arabidopsis thaliana), with gdp-alpha-d-mannose and gdp-beta-l-galactose bound in the active site.
Authors : Major, L.L.; Wolucka, B.A.; Naismith, J.H.
Deposited on : 2005-10-26
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 : 1.7.2 (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

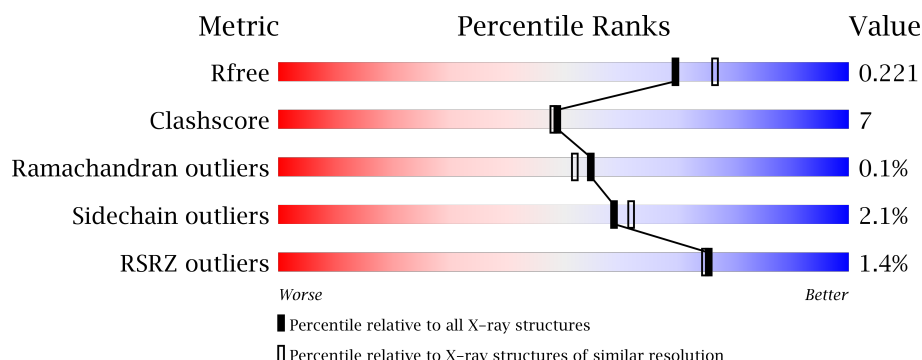
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 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))
R_{free}	100719	6609 (2.00-2.00)
Clashscore	112137	7775 (2.00-2.00)
Ramachandran outliers	110173	7679 (2.00-2.00)
Sidechain outliers	110143	7678 (2.00-2.00)
RSRZ outliers	101464	6696 (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	379	<div> <div>2%</div> <div> <div></div> <div>82%</div> <div>13%</div> <div></div> </div> <div></div> </div>
1	B	379	<div> <div>%</div> <div> <div></div> <div>83%</div> <div>12%</div> <div></div> </div> <div></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
5	FMT	A	1382	-	-	-	X
6	BTB	B	1376	-	-	-	X

2 Entry composition [i](#)

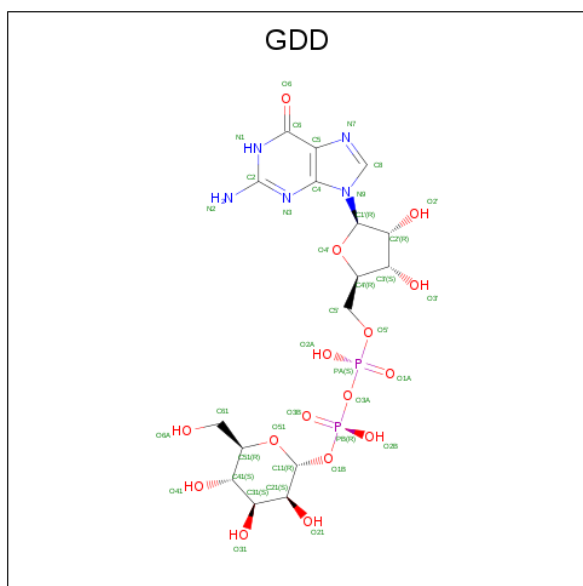
There are 7 unique types of molecules in this entry. The entry contains 7048 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 GDP-MANNOSE-3', 5'-EPIMERASE.

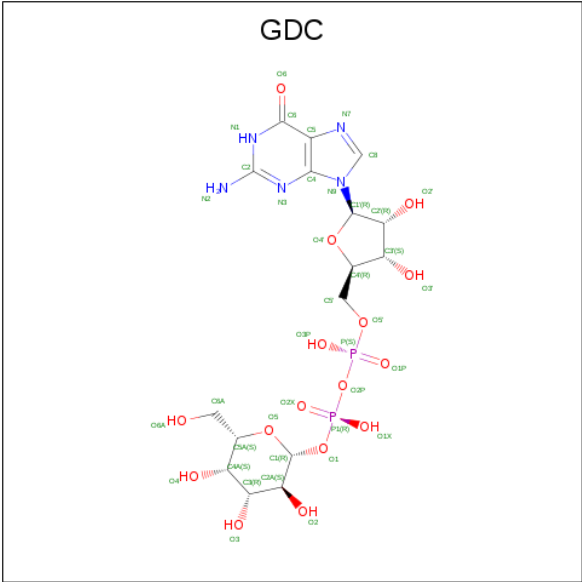
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	364	Total	C	N	O	S	0	7	0
			2944	1864	504	554	22			
1	B	362	Total	C	N	O	S	0	16	0
			2975	1888	511	553	23			

- Molecule 2 is GUANOSINE-5'-DIPHOSPHATE-ALPHA-D-MANNOSE (three-letter code: GDD) (formula: $C_{16}H_{25}N_5O_{16}P_2$).



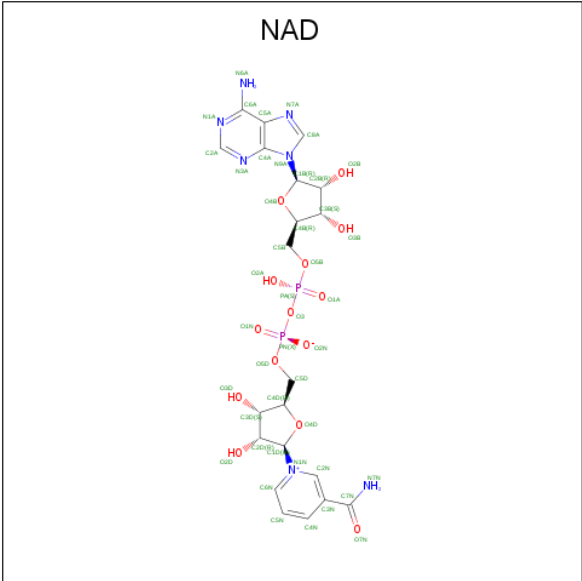
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	1
			39	16	5	16	2		
2	B	1	Total	C	N	O	P	0	1
			39	16	5	16	2		

- Molecule 3 is GUANOSINE-5'-DIPHOSPHATE-BETA-L-GALACTOSE (three-letter code: GDC) (formula: $C_{16}H_{25}N_5O_{16}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	1
			39	16	5	16	2		
3	B	1	Total	C	N	O	P	0	1
			39	16	5	16	2		

- Molecule 4 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (three-letter code: NAD) (formula: C₂₁H₂₇N₇O₁₄P₂).



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

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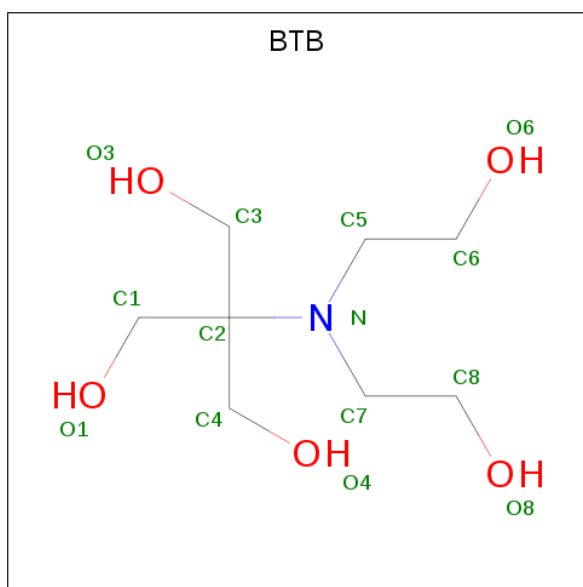
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	B	1	Total	C	N	O	P	0	0
			44	21	7	14	2		

- Molecule 5 is FORMIC ACID (three-letter code: FMT) (formula: CH₂O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			3	1	2		
5	A	1	Total	C	O	0	0
			3	1	2		
5	A	1	Total	C	O	0	0
			3	1	2		
5	B	1	Total	C	O	0	0
			3	1	2		
5	B	1	Total	C	O	0	0
			3	1	2		
5	B	1	Total	C	O	0	0
			3	1	2		
5	B	1	Total	C	O	0	0
			3	1	2		

- Molecule 6 is 2-[BIS-(2-HYDROXY-ETHYL)-AMINO]-2-HYDROXYMETHYL-PROPAN E-1,3-DIOL (three-letter code: BTB) (formula: C₈H₁₉NO₅).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	B	1	Total	C	N	O	0	0
			14	8	1	5		

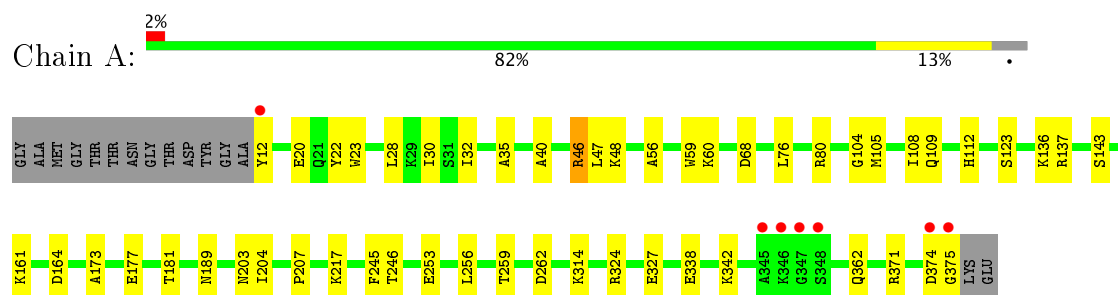
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	391	Total	O	0	0
			391	391		
7	B	456	Total	O	0	0
			456	456		

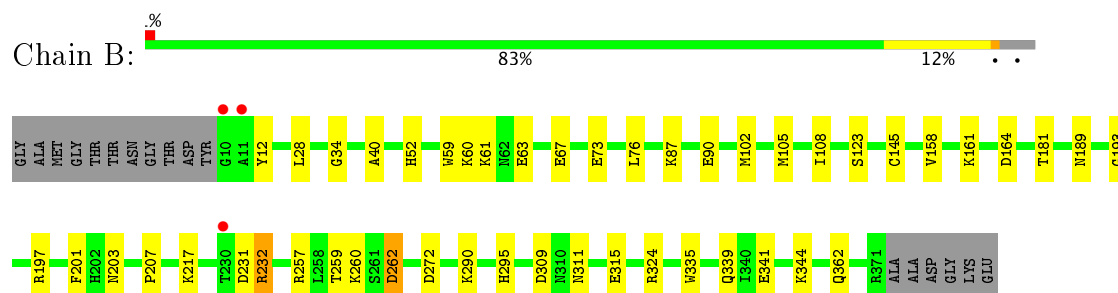
3 Residue-property plots

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: GDP-MANNOSE-3', 5'-EPIMERASE



- Molecule 1: GDP-MANNOSE-3', 5'-EPIMERASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	62.81Å 83.81Å 66.29Å 90.00° 98.91° 90.00°	Depositor
Resolution (Å)	29.88 – 2.00 29.64 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.6 (29.88-2.00) 99.6 (29.64-2.00)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.10 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.147 , 0.223 0.145 , 0.221	Depositor DCC
R_{free} test set	2324 reflections (5.35%)	DCC
Wilson B-factor (Å ²)	23.9	Xtriage
Anisotropy	0.052	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 68.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	7048	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.02% 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: FMT, GDC, GDD, NAD, BTB

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	0.94	2/3037 (0.1%)	0.82	1/4086 (0.0%)
1	B	0.93	1/3102 (0.0%)	0.82	2/4169 (0.0%)
All	All	0.93	3/6139 (0.0%)	0.82	3/8255 (0.0%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	201	PHE	CE1-CZ	5.66	1.48	1.37
1	A	253	GLU	CB-CG	5.23	1.62	1.52
1	A	245	PHE	CE1-CZ	5.17	1.47	1.37

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	272	ASP	CB-CG-OD1	5.95	123.65	118.30
1	B	309	ASP	CB-CG-OD1	5.56	123.30	118.30
1	A	80	ARG	NE-CZ-NH2	-5.29	117.65	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	2944	0	2864	41	0
1	B	2975	0	2917	37	0
2	A	39	0	23	1	0
2	B	39	0	23	3	0
3	A	39	0	23	1	0
3	B	39	0	23	0	0
4	A	44	0	26	0	0
4	B	44	0	26	1	0
5	A	9	0	3	0	0
5	B	15	0	5	1	0
6	B	14	0	19	4	0
7	A	391	0	0	10	1
7	B	456	0	0	15	1
All	All	7048	0	5952	83	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 7.

All (83) 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:262[B]:ASP:OD2	7:B:2312:HOH:O	1.62	1.15
1:B:87:LYS:HD2	7:B:2140:HOH:O	1.50	1.11
1:A:374:ASP:HB2	1:A:375:GLY:HA3	1.28	1.08
1:B:262[A]:ASP:OD2	7:B:2313:HOH:O	1.73	1.07
1:A:177[A]:GLU:OE2	7:A:2181:HOH:O	1.76	1.02
1:A:262[B]:ASP:OD1	7:A:2241:HOH:O	1.80	0.97
1:A:105:MET:CE	1:A:108[B]:ILE:HD11	1.96	0.96
1:B:28:LEU:H	1:B:52:HIS:HD2	1.11	0.90
1:B:231:ASP:HB3	1:B:232:ARG:HH21	1.34	0.89
1:A:105:MET:HE3	1:A:108[B]:ILE:HD11	1.58	0.86
1:A:327[B]:GLU:OE2	7:A:2330:HOH:O	1.93	0.84
1:B:262[A]:ASP:OD1	7:B:2311:HOH:O	1.94	0.84
1:B:145:CYS:SG	2:B:1377[A]:GDD:O41	2.20	0.80
1:B:73:GLU:OE1	6:B:1376:BTB:H71	1.83	0.78
1:A:177[A]:GLU:OE2	7:A:2183:HOH:O	2.06	0.74
1:A:108[B]:ILE:HD13	1:A:109:GLN:N	2.02	0.73
1:A:105:MET:HA	1:A:108[B]:ILE:HD12	1.71	0.71
1:A:324:ARG:HB2	1:A:327[B]:GLU:HG3	1.72	0.70
1:A:136:LYS:NZ	7:A:2134:HOH:O	2.25	0.69
1:B:28:LEU:H	1:B:52:HIS:CD2	2.03	0.68
1:B:60[A]:LYS:HE2	1:B:63:GLU:OE2	1.94	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:362:GLN:HG2	7:B:2194:HOH:O	1.93	0.67
1:A:108[B]:ILE:HD13	1:A:109:GLN:H	1.58	0.67
1:A:374:ASP:HB2	1:A:375:GLY:CA	2.16	0.63
1:B:257:ARG:HA	1:B:260:LYS:HE2	1.82	0.61
1:A:374:ASP:CB	1:A:375:GLY:HA3	2.15	0.60
6:B:1376:BTB:H81	7:B:2141:HOH:O	2.00	0.60
1:B:231:ASP:HB3	1:B:232:ARG:NH2	2.14	0.59
1:B:145:CYS:HG	2:B:1377[A]:GDD:C41	2.13	0.58
1:A:60:LYS:HG2	1:A:371:ARG:HB2	1.86	0.58
1:A:12:TYR:OH	1:A:338:GLU:OE1	2.22	0.57
1:A:177[A]:GLU:HG2	7:A:2181:HOH:O	2.05	0.57
1:A:22:TYR:HB2	1:A:256:LEU:HD13	1.87	0.56
6:B:1376:BTB:H31	7:B:2448:HOH:O	2.06	0.56
1:A:48:LYS:HE2	1:A:68:ASP:O	2.05	0.56
1:B:232:ARG:HD3	1:B:295:HIS:CE1	2.41	0.55
1:A:342:LYS:NZ	7:A:2336:HOH:O	2.40	0.55
1:A:105:MET:HA	1:A:108[B]:ILE:CD1	2.36	0.54
1:B:311[B]:ASN:ND2	7:B:2365:HOH:O	2.40	0.54
1:B:108:ILE:HD13	2:B:1377[A]:GDD:O31	2.07	0.54
1:A:23:TRP:CD1	1:A:28:LEU:HD11	2.43	0.53
6:B:1376:BTB:H41	6:B:1376:BTB:C8	2.39	0.52
1:A:105:MET:HE1	1:A:108[B]:ILE:HD11	1.87	0.51
1:A:177[A]:GLU:CG	7:A:2181:HOH:O	2.58	0.51
1:A:204:ILE:HD13	1:A:246:THR:HB	1.92	0.50
1:B:60[A]:LYS:HE2	7:B:2092:HOH:O	2.11	0.50
1:B:61[A]:LYS:NZ	1:B:67:GLU:OE2	2.36	0.50
1:A:137:ARG:HD3	1:A:259:THR:O	2.13	0.49
1:B:28:LEU:N	1:B:52:HIS:HD2	1.94	0.49
1:A:20:GLU:O	1:A:46:ARG:NH2	2.45	0.49
1:A:143:SER:OG	2:A:1377[A]:GDD:O41	2.31	0.48
1:A:327[B]:GLU:CD	7:A:2329:HOH:O	2.51	0.48
1:B:123:SER:HB3	1:B:181:THR:HG21	1.95	0.48
1:A:161:LYS:O	1:A:164:ASP:HB2	2.14	0.47
1:B:102[B]:MET:HG2	4:B:1380:NAD:O3	2.13	0.47
1:B:52:HIS:HE1	7:B:2077:HOH:O	1.97	0.47
1:A:203:ASN:N	3:A:1378[B]:GDC:O6A	2.40	0.47
1:A:112:HIS:HD2	1:A:173:ALA:H	1.63	0.46
1:B:231:ASP:CB	1:B:232:ARG:HH21	2.17	0.46
1:B:197:ARG:HG3	1:B:259:THR:HA	1.98	0.46
1:B:12:TYR:HB3	1:B:207:PRO:HG2	1.98	0.46
1:B:90[A]:GLU:HG3	7:B:2143:HOH:O	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:35:ALA:HB3	1:A:56:ALA:HB1	1.97	0.45
1:B:108:ILE:HD12	7:B:2244:HOH:O	2.16	0.45
1:A:104:GLY:O	1:A:108[B]:ILE:HG23	2.17	0.45
1:A:12:TYR:HB3	1:A:207:PRO:HG2	2.00	0.43
1:B:335:TRP:O	1:B:339:GLN:HG2	2.18	0.43
1:B:105[A]:MET:HE1	1:B:108:ILE:HD11	1.99	0.43
1:B:324[B]:ARG:NH1	7:B:2381:HOH:O	2.52	0.42
1:B:61[A]:LYS:HE2	1:B:76:LEU:HD22	2.00	0.42
1:A:204:ILE:HA	1:A:246:THR:O	2.20	0.42
1:B:161:LYS:O	1:B:164:ASP:HB2	2.20	0.42
1:A:362:GLN:HB2	1:A:362:GLN:HE21	1.67	0.42
1:B:193:GLY:HA2	7:B:2270:HOH:O	2.18	0.42
1:A:123:SER:HB3	1:A:181:THR:HG21	2.01	0.41
1:A:30:ILE:HD13	1:A:47:LEU:HD13	2.03	0.41
1:B:290:LYS:O	7:B:2345:HOH:O	2.22	0.41
1:A:32:ILE:HG22	1:A:40:ALA:HB1	2.03	0.41
1:B:34:GLY:O	1:B:40:ALA:HB3	2.22	0.40

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:2288:HOH:O	7:B:2088:HOH:O[2_656]	2.05	0.15

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	369/379 (97%)	357 (97%)	12 (3%)	0	100	100
1	B	376/379 (99%)	368 (98%)	7 (2%)	1 (0%)	44	40
All	All	745/758 (98%)	725 (97%)	19 (3%)	1 (0%)	55	52

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	203	ASN

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	317/319 (99%)	312 (98%)	5 (2%)	68	72
1	B	325/319 (102%)	315 (97%)	10 (3%)	45	44
All	All	642/638 (101%)	627 (98%)	15 (2%)	59	58

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	46	ARG
1	A	59	TRP
1	A	76	LEU
1	A	189	ASN
1	A	217	LYS
1	B	59	TRP
1	B	158	VAL
1	B	189	ASN
1	B	217	LYS
1	B	232	ARG
1	B	262[A]	ASP
1	B	262[B]	ASP
1	B	341	GLU
1	B	344[A]	LYS
1	B	344[B]	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	42	HIS

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Mol	Chain	Res	Type
1	A	112	HIS
1	A	362	GLN
1	B	52	HIS
1	B	362	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

15 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	GDD	A	1377[A]	-	35,42,42	1.05	4 (11%)	44,65,65	1.82	9 (20%)
3	GDC	A	1378[B]	-	35,42,42	1.11	4 (11%)	44,65,65	2.35	11 (25%)
4	NAD	A	1379	-	41,48,48	1.60	3 (7%)	43,73,73	2.26	7 (16%)
5	FMT	A	1380	-	0,2,2	0.00	-	0,1,1	0.00	-
5	FMT	A	1381	-	0,2,2	0.00	-	0,1,1	0.00	-
5	FMT	A	1382	-	0,2,2	0.00	-	0,1,1	0.00	-
5	FMT	B	1371	-	0,2,2	0.00	-	0,1,1	0.00	-
5	FMT	B	1372	-	0,2,2	0.00	-	0,1,1	0.00	-
5	FMT	B	1373	-	0,2,2	0.00	-	0,1,1	0.00	-

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	FMT	B	1374	-	0,2,2	0.00	-	0,1,1	0.00	-
5	FMT	B	1375	-	0,2,2	0.00	-	0,1,1	0.00	-
6	BTB	B	1376	-	13,13,13	1.21	1 (7%)	9,16,16	1.70	3 (33%)
2	GDD	B	1377[A]	-	35,42,42	1.12	3 (8%)	44,65,65	1.85	9 (20%)
3	GDC	B	1378[B]	-	35,42,42	1.21	3 (8%)	44,65,65	1.93	8 (18%)
4	NAD	B	1380	-	41,48,48	1.46	3 (7%)	43,73,73	1.99	5 (11%)

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	GDD	A	1377[A]	-	-	0/19/59/59	0/4/4/4
3	GDC	A	1378[B]	-	-	0/19/59/59	0/4/4/4
4	NAD	A	1379	-	-	0/22/62/62	0/5/5/5
5	FMT	A	1380	-	-	0/0/0/0	0/0/0/0
5	FMT	A	1381	-	-	0/0/0/0	0/0/0/0
5	FMT	A	1382	-	-	0/0/0/0	0/0/0/0
5	FMT	B	1371	-	-	0/0/0/0	0/0/0/0
5	FMT	B	1372	-	-	0/0/0/0	0/0/0/0
5	FMT	B	1373	-	-	0/0/0/0	0/0/0/0
5	FMT	B	1374	-	-	0/0/0/0	0/0/0/0
5	FMT	B	1375	-	-	0/0/0/0	0/0/0/0
6	BTB	B	1376	-	-	0/21/21/21	0/0/0/0
2	GDD	B	1377[A]	-	-	0/19/59/59	0/4/4/4
3	GDC	B	1378[B]	-	-	0/19/59/59	0/4/4/4
4	NAD	B	1380	-	-	0/22/62/62	0/5/5/5

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	1380	NAD	O4B-C4B	-2.15	1.40	1.45
3	A	1378[B]	GDC	C6-N1	2.08	1.36	1.33
2	A	1377[A]	GDD	C6-N1	2.08	1.36	1.33
3	B	1378[B]	GDC	C6-N1	2.17	1.37	1.33
2	B	1377[A]	GDD	C6-N1	2.17	1.37	1.33
3	A	1378[B]	GDC	O4'-C1'	2.22	1.44	1.41
2	A	1377[A]	GDD	O4'-C1'	2.22	1.44	1.41
3	A	1378[B]	GDC	C5-C4	2.45	1.46	1.40
2	A	1377[A]	GDD	C5-C4	2.45	1.46	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1378[B]	GDC	C6-C5	2.68	1.46	1.41
2	A	1377[A]	GDD	C6-C5	2.68	1.46	1.41
3	B	1378[B]	GDC	C5-C4	2.74	1.46	1.40
2	B	1377[A]	GDD	C5-C4	2.74	1.46	1.40
6	B	1376	BTB	C5-N	2.95	1.52	1.48
4	A	1379	NAD	C2A-N1A	3.17	1.39	1.33
4	B	1380	NAD	C2A-N3A	3.20	1.37	1.32
3	B	1378[B]	GDC	C6-C5	3.72	1.48	1.41
2	B	1377[A]	GDD	C6-C5	3.72	1.48	1.41
4	A	1379	NAD	C2A-N3A	4.22	1.39	1.32
4	B	1380	NAD	O7N-C7N	7.02	1.38	1.24
4	A	1379	NAD	O7N-C7N	7.06	1.38	1.24

All (52) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1379	NAD	N3A-C2A-N1A	-10.24	119.94	128.86
4	B	1380	NAD	N3A-C2A-N1A	-9.88	120.25	128.86
4	A	1379	NAD	O7N-C7N-C3N	-5.42	113.28	119.62
3	B	1378[B]	GDC	C6-C5-C4	-4.79	116.08	120.84
2	B	1377[A]	GDD	C6-C5-C4	-4.79	116.08	120.84
3	A	1378[B]	GDC	C6-C5-C4	-4.04	116.83	120.84
2	A	1377[A]	GDD	C6-C5-C4	-4.04	116.83	120.84
3	A	1378[B]	GDC	N3-C2-N1	-3.93	121.72	127.46
2	A	1377[A]	GDD	N3-C2-N1	-3.93	121.72	127.46
3	B	1378[B]	GDC	N3-C2-N1	-3.79	121.93	127.46
2	B	1377[A]	GDD	N3-C2-N1	-3.79	121.93	127.46
2	B	1377[A]	GDD	O51-C11-O1B	-3.58	106.69	111.36
3	B	1378[B]	GDC	C5-C6-N1	-3.51	118.48	123.48
2	B	1377[A]	GDD	C5-C6-N1	-3.51	118.48	123.48
6	B	1376	BTB	C5-N-C2	-3.29	104.00	113.70
3	A	1378[B]	GDC	C5-C6-N1	-3.17	118.98	123.48
2	A	1377[A]	GDD	C5-C6-N1	-3.17	118.98	123.48
3	A	1378[B]	GDC	C4'-O4'-C1'	-3.15	106.42	109.77
2	A	1377[A]	GDD	C4'-O4'-C1'	-3.15	106.42	109.77
4	B	1380	NAD	O7N-C7N-C3N	-3.14	115.95	119.62
4	A	1379	NAD	C4A-C5A-N7A	-3.03	106.48	109.41
3	A	1378[B]	GDC	O5-C1-C2A	-2.95	104.60	110.30
2	B	1377[A]	GDD	O21-C21-C31	-2.67	104.54	110.36
2	A	1377[A]	GDD	PB-O1B-C11	-2.48	109.91	119.74
2	B	1377[A]	GDD	O41-C41-C31	-2.39	105.16	110.36
4	A	1379	NAD	C4B-O4B-C1B	-2.24	107.38	109.77

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	1380	NAD	C4B-O4B-C1B	-2.23	107.40	109.77
2	A	1377[A]	GDD	O51-C11-O1B	-2.20	108.49	111.36
3	B	1378[B]	GDC	O3-C3-C4A	-2.02	105.97	110.36
4	A	1379	NAD	C1B-N9A-C4A	-2.01	123.16	126.64
4	B	1380	NAD	O2D-C2D-C1D	2.11	118.23	111.61
6	B	1376	BTB	C6-C5-N	2.14	119.87	111.34
4	A	1379	NAD	C2A-N1A-C6A	2.27	122.74	118.77
3	B	1378[B]	GDC	N2-C2-N1	2.35	120.99	117.24
2	B	1377[A]	GDD	N2-C2-N1	2.35	120.99	117.24
6	B	1376	BTB	O8-C8-C7	2.40	121.81	111.27
3	A	1378[B]	GDC	O4-C4A-C3	2.70	116.23	110.36
2	A	1377[A]	GDD	O1B-C11-C21	3.20	114.25	108.38
3	A	1378[B]	GDC	O2P-P1-O1	3.56	108.42	102.05
3	A	1378[B]	GDC	C1-O5-C5A	4.11	121.47	113.72
3	A	1378[B]	GDC	C6-N1-C2	4.15	122.02	116.06
2	A	1377[A]	GDD	C6-N1-C2	4.15	122.02	116.06
3	B	1378[B]	GDC	C6-N1-C2	4.44	122.44	116.06
2	B	1377[A]	GDD	C6-N1-C2	4.44	122.44	116.06
3	B	1378[B]	GDC	C2-N3-C4	4.82	120.78	115.16
2	B	1377[A]	GDD	C2-N3-C4	4.82	120.78	115.16
3	B	1378[B]	GDC	O1-C1-C2A	4.84	117.25	108.38
4	B	1380	NAD	C3N-C7N-N7N	5.20	123.71	117.77
3	A	1378[B]	GDC	C2-N3-C4	5.23	121.27	115.16
2	A	1377[A]	GDD	C2-N3-C4	5.23	121.27	115.16
4	A	1379	NAD	C3N-C7N-N7N	5.87	124.48	117.77
3	A	1378[B]	GDC	O1-C1-C2A	8.07	123.16	108.38

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

6 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1377[A]	GDD	1	0
3	A	1378[B]	GDC	1	0
5	B	1373	FMT	1	0
6	B	1376	BTB	4	0
2	B	1377[A]	GDD	3	0
4	B	1380	NAD	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 ⓘ

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	364/379 (96%)	-0.38	7 (1%) 67 66	14, 20, 40, 51	6 (1%)
1	B	362/379 (95%)	-0.47	3 (0%) 86 85	13, 19, 32, 46	6 (1%)
All	All	726/758 (95%)	-0.42	10 (1%) 75 75	13, 19, 36, 51	12 (1%)

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	347	GLY	5.0
1	A	374	ASP	3.8
1	A	345	ALA	3.2
1	A	375	GLY	3.1
1	B	10	GLY	3.0
1	A	12	TYR	3.0
1	A	346	LYS	2.9
1	B	11	ALA	2.8
1	B	230[A]	THR	2.4
1	A	348	SER	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	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
5	FMT	A	1382	3/3	0.81	0.25	4.96	38,38,40,40	0
6	BTB	B	1376	14/14	0.77	0.21	2.93	23,28,31,34	14
5	FMT	B	1373	3/3	0.93	0.21	0.62	46,46,46,47	0
3	GDC	B	1378[B]	39/39	0.98	0.13	0.35	3,14,18,19	39
3	GDC	A	1378[B]	39/39	0.98	0.11	0.22	2,15,19,20	39
2	GDD	B	1377[A]	39/39	0.98	0.12	0.18	10,16,18,19	39
2	GDD	A	1377[A]	39/39	0.98	0.11	0.03	14,18,26,29	39
4	NAD	B	1380	44/44	0.97	0.10	-0.58	14,18,21,24	0
4	NAD	A	1379	44/44	0.97	0.08	-0.70	12,19,23,25	0
5	FMT	A	1380	3/3	0.97	0.07	-0.74	34,34,36,36	0
5	FMT	B	1374	3/3	0.99	0.05	-1.38	20,20,22,24	0
5	FMT	B	1372	3/3	0.96	0.17	-	35,35,36,36	0
5	FMT	A	1381	3/3	0.78	0.30	-	58,58,58,59	0
5	FMT	B	1371	3/3	0.85	0.35	-	71,71,71,71	0
5	FMT	B	1375	3/3	0.89	0.41	-	42,42,42,43	0

6.5 Other polymers

There are no such residues in this entry.