



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 04:49 AM UTC

PDB ID : 1UAE / pdb_00001uae
Title : STRUCTURE OF UDP-N-ACETYLGLUCOSAMINE ENOLPYRUVYL
TRANSFERASE
Authors : Skarzynski, T.
Deposited on : 1996-09-30
Resolution : 1.80 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

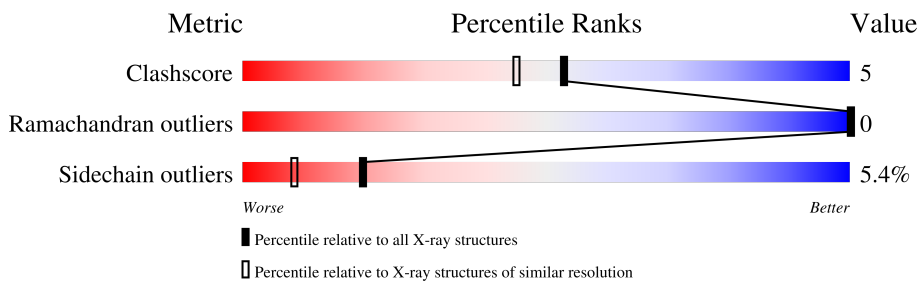
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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	190562	8479 (1.80-1.80)
Ramachandran outliers	187476	8391 (1.80-1.80)
Sidechain outliers	187428	8390 (1.80-1.80)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	419	

2 Entry composition [i](#)

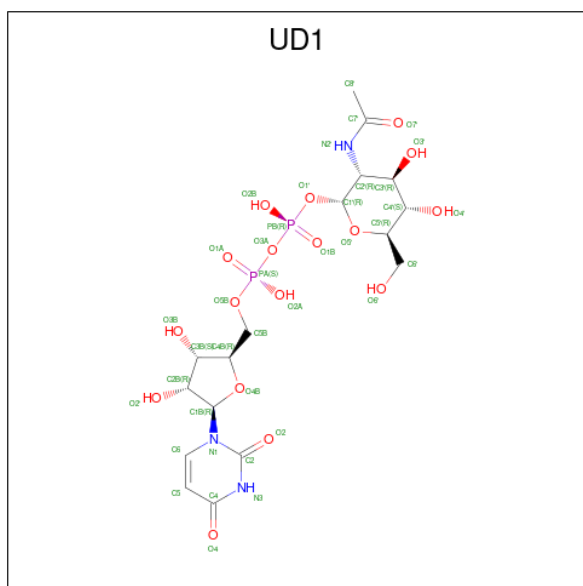
There are 4 unique types of molecules in this entry. The entry contains 3602 atoms, of which 0 are hydrogens and 0 are deuteriums.

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-N-ACETYLGLUCOSAMINE ENOLPYRUVYL TRANSFERASE.

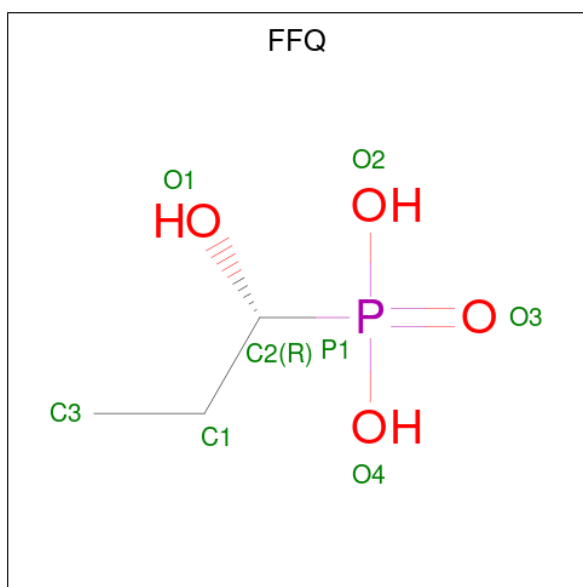
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	418	3133	1967	557	593	16	0	0	0

- Molecule 2 is URIDINE-DIPHOSPHATE-N-ACETYLGLUCOSAMINE (CCD ID: UD1) (formula: C₁₇H₂₇N₃O₁₇P₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
2	A	1	39	17	3	17	2	0	0

- Molecule 3 is [(1R)-1-hydroxypropyl]phosphonic acid (CCD ID: FFQ) (formula: C₃H₉O₄P).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	O	P	0	0
			8	3	4	1		

- Molecule 4 is water.

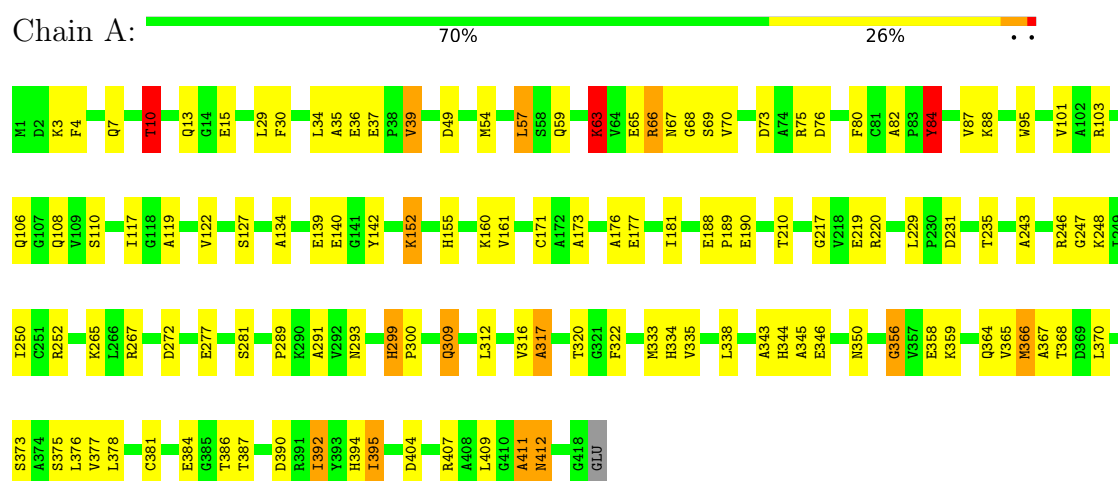
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	422	Total	O	0	0
			422	422		

3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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. 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. 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.

Note EDS was not executed.

• Molecule 1: UDP-N-ACETYLGLUCOSAMINE ENOLPYRUVYL TRANSFERASE



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 3 2 1	Depositor
Cell constants a, b, c, α , β , γ	110.66 Å 110.66 Å 67.55 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	6.00 – 1.80	Depositor
% Data completeness (in resolution range)	(Not available) (6.00-1.80)	Depositor
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.185 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3602	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: UD1, FFQ

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	1.31	9/3177 (0.3%)	2.01	87/4303 (2.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	344	HIS	CE1-NE2	5.58	1.38	1.32
1	A	299	HIS	CE1-NE2	5.42	1.38	1.32
1	A	335	VAL	CA-C	5.38	1.57	1.52
1	A	394	HIS	ND1-CE1	5.31	1.37	1.32
1	A	356	GLY	CA-C	5.26	1.57	1.52
1	A	344	HIS	CG-CD2	5.17	1.41	1.35
1	A	95	TRP	CD2-CE2	5.16	1.50	1.41
1	A	155	HIS	ND1-CE1	5.03	1.37	1.32
1	A	344	HIS	ND1-CE1	5.03	1.37	1.32

All (87) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	377	VAL	N-CA-C	10.00	120.02	110.42
1	A	68	GLY	CA-C-O	9.56	130.32	120.09
1	A	152	LYS	CA-C-N	-8.12	112.63	120.34
1	A	152	LYS	C-N-CA	-8.12	112.63	120.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	87	VAL	N-CA-C	7.61	118.39	110.62
1	A	404	ASP	N-CA-C	7.46	119.06	111.07
1	A	346	GLU	CA-C-N	-7.43	113.57	122.93
1	A	346	GLU	C-N-CA	-7.43	113.57	122.93
1	A	84	TYR	N-CA-C	7.27	120.13	111.33
1	A	140	GLU	CA-C-N	-7.05	110.18	122.09
1	A	140	GLU	C-N-CA	-7.05	110.18	122.09
1	A	231	ASP	N-CA-C	6.98	119.20	109.15
1	A	160	LYS	CB-CA-C	-6.84	95.03	110.07
1	A	101	VAL	N-CA-C	6.78	116.93	110.42
1	A	49	ASP	N-CA-C	6.77	118.66	111.28
1	A	365	VAL	CA-C-N	-6.67	112.47	122.74
1	A	365	VAL	C-N-CA	-6.67	112.47	122.74
1	A	358	GLU	CB-CG-CD	-6.63	101.33	112.60
1	A	117	ILE	N-CA-C	6.45	118.35	112.29
1	A	370	LEU	N-CA-C	6.40	118.83	111.02
1	A	122	VAL	CA-C-O	6.38	127.69	120.42
1	A	139	GLU	CB-CG-CD	-6.35	101.81	112.60
1	A	65	GLU	CA-C-N	-6.34	111.94	122.21
1	A	65	GLU	C-N-CA	-6.34	111.94	122.21
1	A	161	VAL	N-CA-C	6.15	116.99	109.30
1	A	30	PHE	N-CA-C	6.13	118.04	111.36
1	A	39	VAL	CA-CB-CG1	6.12	120.80	110.40
1	A	243	ALA	N-CA-C	6.11	117.61	111.07
1	A	267	ARG	N-CA-C	6.10	118.71	111.33
1	A	381	CYS	N-CA-C	6.09	118.00	111.36
1	A	39	VAL	N-CA-CB	-6.02	101.57	111.44
1	A	333	MET	N-CA-C	5.97	119.66	112.38
1	A	82	ALA	N-CA-C	5.96	117.39	109.83
1	A	368	THR	CA-C-O	5.94	125.68	119.14
1	A	127	SER	N-CA-C	5.94	118.23	111.11
1	A	76	ASP	CA-CB-CG	-5.91	106.69	112.60
1	A	320	THR	CA-C-N	-5.91	117.06	121.61
1	A	320	THR	C-N-CA	-5.91	117.06	121.61
1	A	316	VAL	CA-C-O	5.90	125.34	119.97
1	A	411	ALA	N-CA-C	5.84	118.74	110.50
1	A	247	GLY	N-CA-C	5.83	120.95	112.55
1	A	10	THR	CA-C-N	-5.82	113.48	122.09
1	A	10	THR	C-N-CA	-5.82	113.48	122.09
1	A	76	ASP	N-CA-C	5.82	119.94	112.26
1	A	75	ARG	CA-C-N	-5.78	114.01	122.86
1	A	75	ARG	C-N-CA	-5.78	114.01	122.86

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	272	ASP	CA-C-N	-5.75	115.68	123.10
1	A	272	ASP	C-N-CA	-5.75	115.68	123.10
1	A	190	GLU	N-CA-C	5.73	118.47	111.82
1	A	35	ALA	CA-C-N	-5.71	113.53	122.73
1	A	35	ALA	C-N-CA	-5.71	113.53	122.73
1	A	312	LEU	N-CA-C	5.67	117.13	111.07
1	A	409	LEU	CA-C-N	-5.62	113.76	122.69
1	A	409	LEU	C-N-CA	-5.62	113.76	122.69
1	A	407	ARG	CD-NE-CZ	-5.49	116.72	124.40
1	A	63	LYS	CG-CD-CE	5.45	123.84	111.30
1	A	65	GLU	N-CA-CB	-5.43	101.43	111.52
1	A	235	THR	N-CA-C	5.37	116.82	110.97
1	A	59	GLN	N-CA-C	5.35	118.91	112.38
1	A	345	ALA	CA-C-N	-5.35	114.50	122.74
1	A	345	ALA	C-N-CA	-5.35	114.50	122.74
1	A	317	ALA	CA-C-N	-5.32	115.35	122.42
1	A	317	ALA	C-N-CA	-5.32	115.35	122.42
1	A	291	ALA	CA-C-N	-5.31	114.30	122.68
1	A	291	ALA	C-N-CA	-5.31	114.30	122.68
1	A	117	ILE	CA-C-O	5.30	125.67	119.89
1	A	37	GLU	CA-C-N	-5.26	114.36	119.78
1	A	37	GLU	C-N-CA	-5.26	114.36	119.78
1	A	387	THR	CA-C-N	-5.25	115.97	123.11
1	A	387	THR	C-N-CA	-5.25	115.97	123.11
1	A	119	ALA	CA-C-O	5.24	126.02	120.10
1	A	376	LEU	CA-C-N	-5.22	113.98	120.56
1	A	376	LEU	C-N-CA	-5.22	113.98	120.56
1	A	387	THR	CA-CB-CG2	5.22	119.37	110.50
1	A	80	PHE	CA-CB-CG	-5.21	108.58	113.80
1	A	343	ALA	N-CA-C	5.20	117.43	110.35
1	A	57	LEU	N-CA-C	5.20	117.35	111.11
1	A	267	ARG	CD-NE-CZ	-5.19	117.13	124.40
1	A	375	SER	N-CA-C	5.15	116.89	111.28
1	A	338	LEU	N-CA-C	5.14	116.89	111.28
1	A	392	ILE	N-CA-C	5.11	119.64	113.00
1	A	134	ALA	N-CA-C	5.08	117.67	110.50
1	A	173	ALA	N-CA-C	5.08	118.58	112.38
1	A	210	THR	N-CA-CB	5.04	118.20	110.70
1	A	34	LEU	N-CA-C	5.01	117.63	111.82
1	A	252	ARG	NE-CZ-NH1	-5.00	116.50	121.50
1	A	366	MET	CA-CB-CG	-5.00	104.09	114.10

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	103	ARG	Sidechain
1	A	84	TYR	Sidechain

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	3133	0	3212	32	1
2	A	39	0	25	0	0
3	A	8	0	6	0	0
4	A	422	0	0	3	4
All	All	3602	0	3243	32	4

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 5.

All (32) 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:A:54:MET:SD	1:A:66:ARG:HG2	2.09	0.92
1:A:309:GLN:HE21	1:A:309:GLN:H	1.24	0.84
1:A:13:GLN:HE22	1:A:248:LYS:HG2	1.49	0.77
1:A:359:LYS:HE2	1:A:384:GLU:HB2	1.68	0.74
1:A:366:MET:HG3	1:A:390:ASP:HB3	1.71	0.73
1:A:66:ARG:HB3	1:A:70:VAL:HG22	1.73	0.69
1:A:10:THR:HG21	1:A:411:ALA:HA	1.77	0.66
1:A:250:ILE:HG12	1:A:281:SER:HB3	1.83	0.61
1:A:317:ALA:O	1:A:356:GLY:HA3	2.02	0.60
1:A:106:GLN:HE21	1:A:108:GLN:HE22	1.51	0.59
1:A:367:ALA:HB1	1:A:373:SER:HB3	1.85	0.59
1:A:7:GLN:HE22	1:A:386:THR:HG23	1.69	0.58
1:A:176:ALA:O	1:A:217:GLY:HA3	2.04	0.57
1:A:293:ASN:HD22	1:A:322:PHE:H	1.54	0.55
1:A:289:PRO:HD3	4:A:807:HOH:O	2.07	0.54
1:A:63:LYS:HB2	1:A:73:ASP:HB3	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:293:ASN:ND2	1:A:322:PHE:H	2.09	0.51
1:A:84:TYR:O	1:A:88:LYS:HG2	2.11	0.50
1:A:412:ASN:HD22	1:A:412:ASN:C	2.21	0.48
1:A:67:ASN:N	1:A:69:SER:O	2.39	0.47
1:A:4:PHE:CD2	1:A:392:ILE:HG12	2.50	0.46
1:A:334:HIS:H	1:A:334:HIS:CD2	2.34	0.46
1:A:152:LYS:HD2	1:A:177:GLU:HB3	1.98	0.46
1:A:36:GLU:CD	1:A:220:ARG:HG3	2.41	0.46
1:A:13:GLN:NE2	1:A:248:LYS:HG2	2.25	0.44
1:A:110:SER:HA	1:A:142:TYR:HA	2.00	0.44
1:A:181:ILE:HD13	1:A:181:ILE:HG21	1.81	0.43
1:A:171:CYS:HB2	4:A:740:HOH:O	2.19	0.42
1:A:188:GLU:HA	1:A:189:PRO:HD3	1.88	0.42
1:A:246:ARG:HG2	4:A:807:HOH:O	2.19	0.42
1:A:299:HIS:CG	1:A:300:PRO:HA	2.55	0.42
1:A:392:ILE:O	1:A:395:ILE:HG22	2.20	0.41

All (4) 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)
1:A:67:ASN:ND2	4:A:709:HOH:O[4_556]	1.80	0.40
4:A:567:HOH:O	4:A:782:HOH:O[1_556]	1.98	0.22
4:A:642:HOH:O	4:A:692:HOH:O[2_655]	2.06	0.14
4:A:536:HOH:O	4:A:624:HOH:O[3_665]	2.19	0.01

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	416/419 (99%)	407 (98%)	9 (2%)	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	331/332 (100%)	313 (95%)	18 (5%)	20 8

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	LYS
1	A	10	THR
1	A	15	GLU
1	A	29	LEU
1	A	39	VAL
1	A	57	LEU
1	A	63	LYS
1	A	66	ARG
1	A	219	GLU
1	A	229	LEU
1	A	265	LYS
1	A	277	GLU
1	A	309	GLN
1	A	350	ASN
1	A	364	GLN
1	A	378	LEU
1	A	395	ILE
1	A	412	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	7	GLN
1	A	13	GLN
1	A	59	GLN
1	A	108	GLN

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Mol	Chain	Res	Type
1	A	131	GLN
1	A	184	ASN
1	A	253	ASN
1	A	293	ASN
1	A	309	GLN
1	A	334	HIS
1	A	350	ASN
1	A	364	GLN
1	A	412	ASN

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

2 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	UD1	A	420	-	40,41,41	2.12	6 (15%)	59,62,62	1.14	6 (10%)
3	FFQ	A	421	-	5,7,7	3.10	2 (40%)	5,10,10	0.88	0

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	UD1	A	420	-	-	6/26/63/63	0/3/3/3
3	FFQ	A	421	-	-	0/8/8/8	-

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	420	UD1	PA-O3A	-10.10	1.48	1.59
3	A	421	FFQ	C1-C2	6.31	1.57	1.52
2	A	420	UD1	C6-N1	4.15	1.48	1.38
2	A	420	UD1	C4-N3	-3.81	1.32	1.38
2	A	420	UD1	PB-O3A	-3.69	1.55	1.59
2	A	420	UD1	C2-N3	-2.92	1.32	1.38
3	A	421	FFQ	P1-O3	-2.38	1.45	1.49
2	A	420	UD1	PA-O2A	-2.04	1.45	1.55

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	420	UD1	C5-C4-N3	2.58	118.41	114.80
2	A	420	UD1	O2-C2-N3	-2.49	116.90	121.49
2	A	420	UD1	N3-C2-N1	2.40	118.01	114.89
2	A	420	UD1	O5'-C1'-O1'	-2.39	108.24	111.36
2	A	420	UD1	O4-C4-C5	-2.19	121.39	125.16
2	A	420	UD1	C5-C6-N1	-2.05	118.51	121.84

There are no chirality outliers.

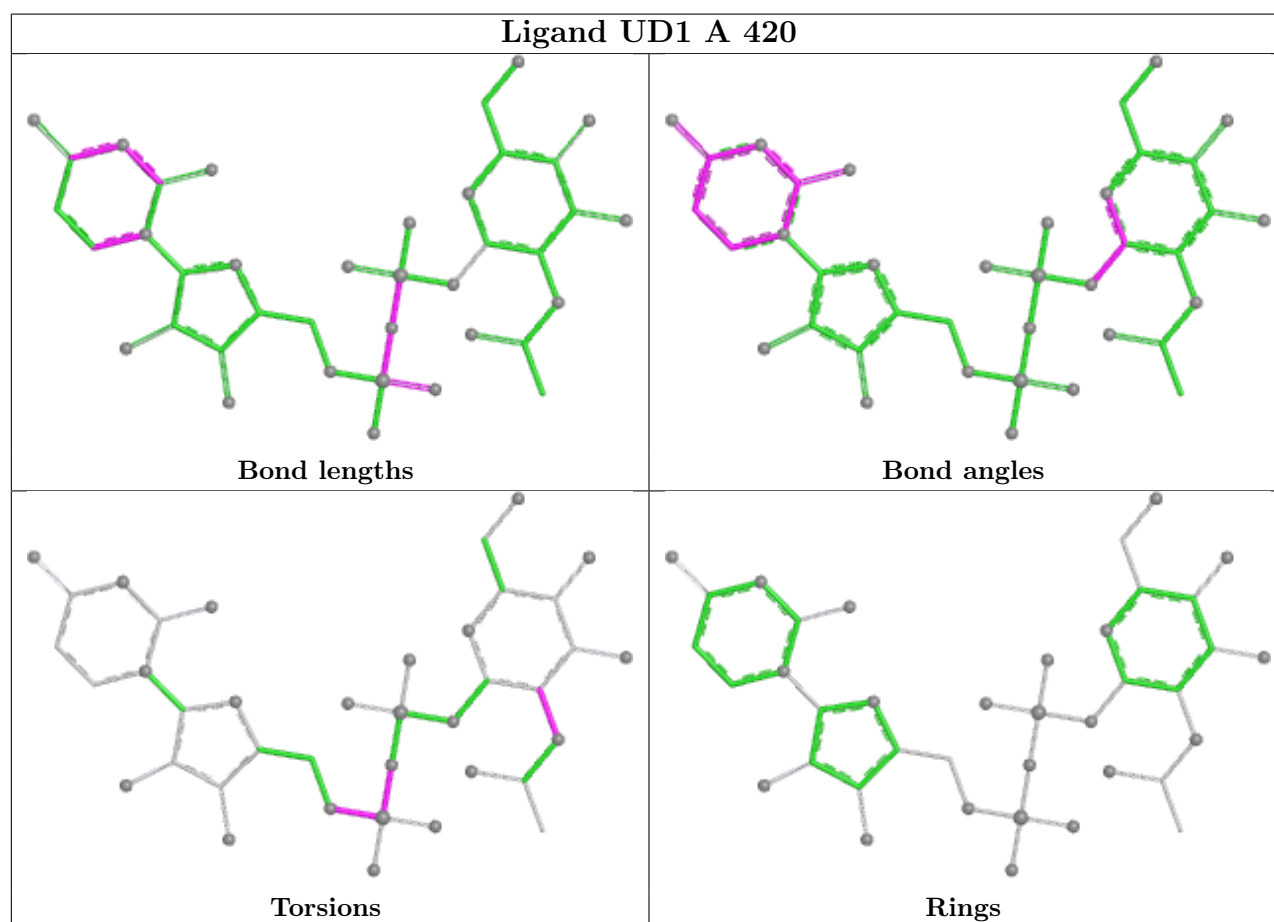
All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	420	UD1	C5B-O5B-PA-O1A
2	A	420	UD1	C5B-O5B-PA-O2A
2	A	420	UD1	C5B-O5B-PA-O3A
2	A	420	UD1	C3'-C2'-N2'-C7'
2	A	420	UD1	PB-O3A-PA-O1A
2	A	420	UD1	PB-O3A-PA-O2A

There are no ring outliers.

No monomer is involved in short contacts.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



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

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.