



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 5, 2019 – 07:38 PM EDT

PDB ID : 1A2N
Title : STRUCTURE OF THE C115A MUTANT OF MURA COMPLEXED WITH
THE FLUORINATED ANALOG OF THE REACTION TETRAHEDRAL
INTERMEDIATE
Authors : Skarzynski, T.
Deposited on : 1998-01-06
Resolution : 2.80 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.8.0 (224370), CSD as540be (2019)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20171227.v01 (using entries in the PDB archive December 27th 2017)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.4

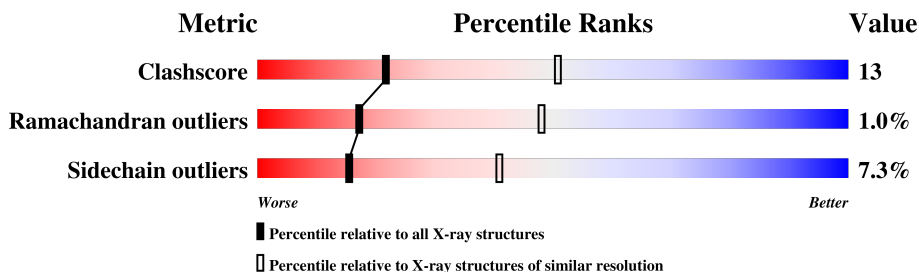
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	122126	3209 (2.80-2.80)
Ramachandran outliers	120053	3158 (2.80-2.80)
Sidechain outliers	120020	3160 (2.80-2.80)

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 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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	419	 63% 30% 5%

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 3449 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.

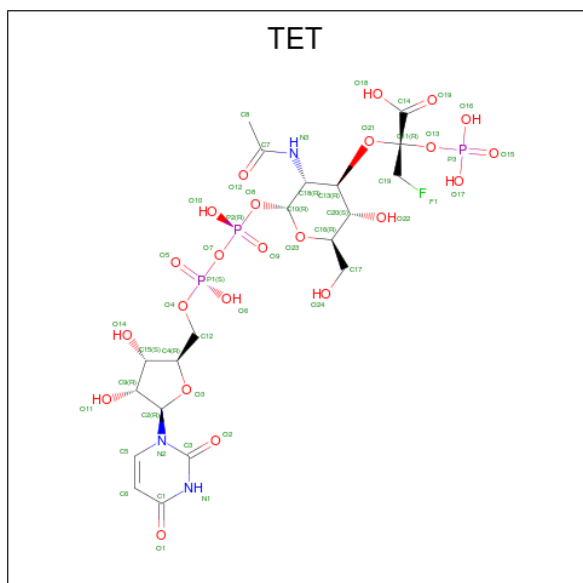
- Molecule 1 is a protein called UDP-N-ACETYLGLUCOSAMINE ENOLPYRUVYL TRANSFERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	418	3128	1965	557	591	15	0	0	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	115	ALA	CYS	ENGINEERED	UNP P0A749

- Molecule 2 is URIDINE-DIPHOSPHATE-2(N-ACETYLGLUCOSAMINYL-3-FLUORO-2-PHOSPHONOOXY)PROPIONIC ACID (three-letter code: TET) (formula: $C_{20}H_{31}FN_3O_{23}P_3$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
			Total	C	F	N	O	P		
2	A	1	50	20	1	3	23	3	0	0

- Molecule 3 is water.

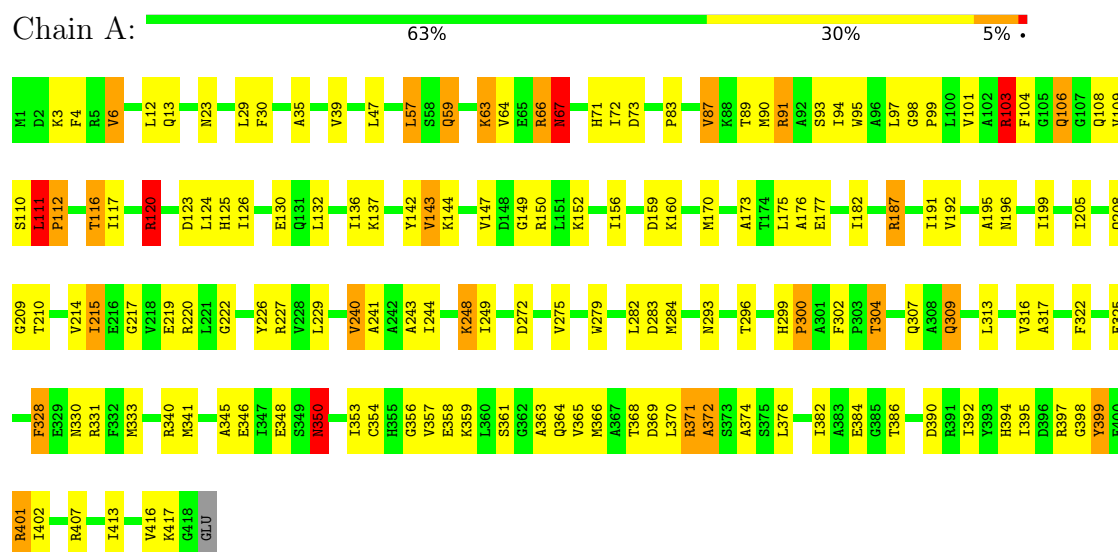
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	271	Total 271	O 271	0	0

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. 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	111.15Å 111.15Å 67.51Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	10.00 – 2.80	Depositor
% Data completeness (in resolution range)	99.2 (10.00-2.80)	Depositor
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	PROLSQ, X-PLOR 3.1	Depositor
R, R_{free}	0.175 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3449	wwPDB-VP
Average B, all atoms (Å ²)	23.0	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: TET

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	1.01	6/3172 (0.2%)	1.79	55/4297 (1.3%)

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	1	2

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	346	GLU	CG-CD	9.86	1.66	1.51
1	A	346	GLU	CB-CG	7.58	1.66	1.52
1	A	279	TRP	CD2-CE2	7.25	1.50	1.41
1	A	346	GLU	CA-CB	7.04	1.69	1.53
1	A	95	TRP	CD2-CE2	5.83	1.48	1.41
1	A	227	ARG	NE-CZ	5.07	1.39	1.33

All (55) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	120	ARG	NE-CZ-NH1	12.94	126.77	120.30
1	A	67	ASN	N-CA-CB	12.18	132.52	110.60
1	A	227	ARG	NE-CZ-NH1	11.63	126.12	120.30
1	A	187	ARG	NE-CZ-NH1	10.71	125.65	120.30
1	A	346	GLU	OE1-CD-OE2	-10.56	110.62	123.30
1	A	187	ARG	NE-CZ-NH2	-9.65	115.48	120.30
1	A	401	ARG	NE-CZ-NH1	8.98	124.79	120.30

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	407	ARG	NE-CZ-NH2	-8.84	115.88	120.30
1	A	91	ARG	CD-NE-CZ	-8.48	111.72	123.60
1	A	220	ARG	NE-CZ-NH2	-8.43	116.09	120.30
1	A	120	ARG	CD-NE-CZ	7.92	134.68	123.60
1	A	191	ILE	CA-CB-CG1	-7.78	96.22	111.00
1	A	346	GLU	CA-CB-CG	7.72	130.38	113.40
1	A	150	ARG	NE-CZ-NH2	-7.54	116.53	120.30
1	A	66	ARG	CB-CA-C	-7.36	95.69	110.40
1	A	366	MET	CA-CB-CG	-7.30	100.89	113.30
1	A	226	TYR	CB-CG-CD1	-7.16	116.71	121.00
1	A	39	VAL	CG1-CB-CG2	-6.92	99.82	110.90
1	A	132	LEU	CB-CG-CD1	-6.79	99.45	111.00
1	A	91	ARG	N-CA-CB	-6.68	98.58	110.60
1	A	346	GLU	CG-CD-OE1	6.50	131.29	118.30
1	A	354	CYS	CB-CA-C	-6.37	97.66	110.40
1	A	397	ARG	NE-CZ-NH2	-6.34	117.13	120.30
1	A	150	ARG	NE-CZ-NH1	6.29	123.45	120.30
1	A	331	ARG	NE-CZ-NH1	6.28	123.44	120.30
1	A	103	ARG	NE-CZ-NH2	-6.23	117.19	120.30
1	A	340	ARG	CD-NE-CZ	-6.12	115.03	123.60
1	A	346	GLU	CB-CG-CD	5.97	130.32	114.20
1	A	372	ALA	CB-CA-C	-5.79	101.41	110.10
1	A	143	VAL	CB-CA-C	-5.78	100.41	111.40
1	A	173	ALA	CB-CA-C	-5.72	101.52	110.10
1	A	91	ARG	CA-CB-CG	5.64	125.81	113.40
1	A	87	VAL	CA-CB-CG2	-5.58	102.53	110.90
1	A	120	ARG	CB-CG-CD	5.50	125.91	111.60
1	A	220	ARG	NE-CZ-NH1	5.42	123.01	120.30
1	A	227	ARG	NE-CZ-NH2	-5.41	117.60	120.30
1	A	111	LEU	N-CA-CB	-5.40	99.60	110.40
1	A	159	ASP	N-CA-CB	-5.40	100.88	110.60
1	A	350	ASN	N-CA-C	5.39	125.57	111.00
1	A	300	PRO	N-CA-CB	-5.36	96.70	102.60
1	A	348	GLU	CB-CA-C	-5.32	99.77	110.40
1	A	328	PHE	CB-CA-C	-5.27	99.86	110.40
1	A	4	PHE	CB-CG-CD2	-5.24	117.14	120.80
1	A	215	ILE	CB-CA-C	-5.21	101.17	111.60
1	A	364	GLN	CA-C-N	-5.21	105.74	117.20
1	A	142	TYR	CB-CG-CD1	-5.18	117.89	121.00
1	A	399	TYR	CB-CG-CD2	-5.14	117.91	121.00
1	A	371	ARG	CD-NE-CZ	-5.09	116.47	123.60
1	A	112	PRO	C-N-CA	-5.08	111.63	122.30

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	325	GLU	C-N-CA	-5.08	109.01	121.70
1	A	240	VAL	CA-CB-CG1	5.07	118.51	110.90
1	A	364	GLN	CB-CA-C	5.06	120.53	110.40
1	A	120	ARG	CG-CD-NE	5.04	122.37	111.80
1	A	386	THR	CA-CB-CG2	5.03	119.44	112.40
1	A	106	GLN	CA-CB-CG	-5.02	102.36	113.40

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	67	ASN	CA

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	120	ARG	Sidechain
1	A	187	ARG	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	3128	0	3209	79	0
2	A	50	0	26	5	0
3	A	271	0	0	7	1
All	All	3449	0	3235	82	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 13.

All (82) 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:111:LEU:HD13	1:A:143:VAL:HG22	1.62	0.80
1:A:64:VAL:HG22	1:A:72:ILE:HG12	1.65	0.78
1:A:116:THR:HG22	1:A:333:MET:HE1	1.65	0.77

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:116:THR:HG22	1:A:333:MET:CE	2.17	0.74
1:A:370:LEU:HG	1:A:371:ARG:HD3	1.69	0.74
1:A:243:ALA:HA	1:A:284:MET:HG3	1.73	0.71
2:A:420:TET:O17	2:A:420:TET:H192	1.93	0.66
2:A:420:TET:O17	2:A:420:TET:C19	2.46	0.63
1:A:30:PHE:HZ	1:A:90:MET:HE1	1.66	0.60
1:A:12:LEU:HD13	1:A:241:ALA:HB1	1.83	0.60
1:A:94:ILE:HA	1:A:109:VAL:HG11	1.84	0.59
1:A:152:LYS:HD2	1:A:177:GLU:HB3	1.84	0.57
2:A:420:TET:H9	3:A:555:HOH:O	2.04	0.57
1:A:35:ALA:HB1	1:A:222:GLY:O	2.06	0.55
1:A:47:LEU:HD13	1:A:398:GLY:HA2	1.89	0.55
1:A:87:VAL:HG13	1:A:93:SER:OG	2.07	0.54
1:A:176:ALA:O	1:A:217:GLY:HA3	2.07	0.54
1:A:316:VAL:HA	1:A:357:VAL:O	2.08	0.54
1:A:152:LYS:HD2	1:A:177:GLU:O	2.08	0.54
1:A:215:ILE:HD12	1:A:215:ILE:N	2.24	0.53
1:A:195:ALA:HB3	1:A:208:GLN:HG3	1.90	0.52
1:A:365:VAL:HB	1:A:376:LEU:HD13	1.92	0.52
1:A:120:ARG:HA	2:A:420:TET:O11	2.10	0.52
1:A:341:MET:HG2	1:A:363:ALA:HB3	1.92	0.52
1:A:97:LEU:O	1:A:101:VAL:HG23	2.10	0.51
1:A:358:GLU:HA	3:A:640:HOH:O	2.10	0.51
1:A:120:ARG:HD2	1:A:328:PHE:CE1	2.46	0.51
1:A:293:ASN:ND2	1:A:322:PHE:H	2.09	0.51
1:A:98:GLY:HA3	3:A:632:HOH:O	2.11	0.50
1:A:120:ARG:HG2	3:A:555:HOH:O	2.11	0.50
1:A:350:ASN:N	1:A:350:ASN:HD22	2.11	0.49
1:A:374:ALA:HA	1:A:395:ILE:HD11	1.94	0.49
1:A:111:LEU:HD13	1:A:143:VAL:CG2	2.39	0.48
1:A:272:ASP:HB3	1:A:283:ASP:HB3	1.94	0.48
1:A:66:ARG:O	1:A:67:ASN:HB2	2.14	0.48
1:A:130:GLU:HG3	1:A:136:ILE:HD12	1.96	0.48
1:A:103:ARG:HG2	1:A:104:PHE:CE2	2.48	0.47
1:A:196:ASN:HA	1:A:199:ILE:HD12	1.96	0.47
1:A:57:LEU:HD23	1:A:72:ILE:HD13	1.95	0.47
1:A:192:VAL:HG22	1:A:209:GLY:N	2.30	0.47
1:A:63:LYS:HB2	1:A:73:ASP:HB3	1.96	0.47
1:A:244:ILE:HG22	1:A:313:LEU:HA	1.96	0.47
1:A:240:VAL:O	1:A:244:ILE:HG12	2.15	0.46
1:A:345:ALA:HA	1:A:353:ILE:O	2.15	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:3:LYS:HE3	1:A:416:VAL:HG23	1.97	0.46
1:A:126:ILE:HG23	1:A:136:ILE:HD13	1.97	0.46
1:A:299:HIS:CG	1:A:300:PRO:HA	2.51	0.46
1:A:244:ILE:HD12	1:A:382:ILE:HD13	1.96	0.46
1:A:248:LYS:HE3	1:A:248:LYS:HB2	1.87	0.46
1:A:170:MET:SD	1:A:195:ALA:HB2	2.57	0.45
1:A:296:THR:HG21	1:A:304:THR:HA	1.97	0.45
1:A:99:PRO:HD3	3:A:632:HOH:O	2.16	0.45
1:A:3:LYS:HG2	1:A:390:ASP:HA	1.99	0.45
1:A:117:ILE:HA	1:A:330:ASN:O	2.16	0.45
1:A:6:VAL:HG13	1:A:413:ILE:HG12	1.99	0.44
1:A:64:VAL:HA	1:A:71:HIS:O	2.18	0.44
1:A:175:LEU:HD11	3:A:454:HOH:O	2.16	0.44
1:A:392:ILE:HA	1:A:392:ILE:HD12	1.64	0.44
1:A:399:TYR:HB3	1:A:402:ILE:HB	1.99	0.44
1:A:413:ILE:HD13	1:A:413:ILE:HG21	1.72	0.43
1:A:309:GLN:NE2	1:A:309:GLN:H	2.16	0.43
1:A:359:LYS:HE2	1:A:384:GLU:HB2	2.00	0.43
1:A:369:ASP:HA	1:A:394:HIS:CD2	2.53	0.43
1:A:23:ASN:OD1	2:A:420:TET:H83	2.19	0.42
1:A:317:ALA:O	1:A:356:GLY:HA3	2.19	0.42
1:A:137:LYS:HG3	1:A:144:LYS:HB2	2.01	0.42
1:A:116:THR:HG21	1:A:368:THR:O	2.19	0.42
1:A:175:LEU:HD23	1:A:175:LEU:HA	1.86	0.42
1:A:147:VAL:HG13	1:A:149:GLY:O	2.20	0.42
1:A:111:LEU:HA	1:A:112:PRO:HD3	1.91	0.42
1:A:106:GLN:NE2	1:A:108:GLN:HE22	2.18	0.42
1:A:243:ALA:HA	1:A:284:MET:CG	2.48	0.41
1:A:125:HIS:H	1:A:125:HIS:CD2	2.38	0.41
1:A:156:ILE:O	1:A:182:ILE:HA	2.20	0.41
1:A:372:ALA:HB1	3:A:654:HOH:O	2.20	0.41
1:A:59:GLN:HG2	1:A:83:PRO:HG3	2.01	0.41
1:A:205:ILE:HA	1:A:214:VAL:O	2.19	0.41
1:A:350:ASN:HD22	1:A:350:ASN:H	1.68	0.41
1:A:244:ILE:CG2	1:A:313:LEU:HA	2.51	0.40
1:A:249:ILE:HG21	1:A:249:ILE:HD13	1.82	0.40
1:A:124:LEU:HD11	1:A:160:LYS:HG3	2.03	0.40
1:A:296:THR:HG22	1:A:302:PHE:HD1	1.86	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)
3:A:467:HOH:O	3:A:643:HOH:O[1_554]	1.61	0.59

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%)	389 (94%)	23 (6%)	4 (1%)	17 48

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	67	ASN
1	A	123	ASP
1	A	417	LYS
1	A	304	THR

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	329/331 (99%)	305 (93%)	24 (7%)	15 41

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	VAL
1	A	13	GLN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	29	LEU
1	A	57	LEU
1	A	59	GLN
1	A	63	LYS
1	A	67	ASN
1	A	89	THR
1	A	91	ARG
1	A	103	ARG
1	A	110	SER
1	A	111	LEU
1	A	116	THR
1	A	210	THR
1	A	219	GLU
1	A	229	LEU
1	A	248	LYS
1	A	275	VAL
1	A	282	LEU
1	A	307	GLN
1	A	309	GLN
1	A	350	ASN
1	A	361	SER
1	A	401	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	7	GLN
1	A	13	GLN
1	A	42	GLN
1	A	59	GLN
1	A	67	ASN
1	A	106	GLN
1	A	184	ASN
1	A	253	ASN
1	A	293	ASN
1	A	309	GLN
1	A	350	ASN
1	A	394	HIS

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 ⓘ

1 ligand is 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	TET	A	420	-	39,52,52	1.80	12 (30%)	48,80,80	2.29	17 (35%)

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	TET	A	420	-	-	9/27/83/83	0/3/3/3

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	420	TET	C3-N1	-4.14	1.30	1.38
2	A	420	TET	P3-O13	-3.34	1.53	1.59
2	A	420	TET	O11-C9	3.34	1.50	1.43

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	420	TET	O14-C15	3.03	1.50	1.43
2	A	420	TET	O21-C13	2.86	1.45	1.42
2	A	420	TET	P3-O17	-2.66	1.44	1.54
2	A	420	TET	P2-O9	-2.43	1.42	1.50
2	A	420	TET	O24-C17	2.33	1.52	1.42
2	A	420	TET	P3-O15	-2.32	1.43	1.50
2	A	420	TET	O22-C20	2.30	1.48	1.43
2	A	420	TET	O3-C4	-2.16	1.40	1.45
2	A	420	TET	C15-C4	2.07	1.58	1.53

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	420	TET	O22-C20-C13	5.98	126.07	109.94
2	A	420	TET	O7-P2-O8	5.61	113.80	102.48
2	A	420	TET	O11-C9-C15	4.91	127.73	111.80
2	A	420	TET	O13-P3-O15	-4.34	92.64	109.39
2	A	420	TET	C6-C1-N1	-3.97	114.38	123.28
2	A	420	TET	P2-O7-P1	-3.53	121.35	132.57
2	A	420	TET	C13-C18-N3	-3.40	105.03	110.90
2	A	420	TET	O14-C15-C4	3.05	119.88	111.07
2	A	420	TET	O14-C15-C9	-2.93	102.30	111.80
2	A	420	TET	O3-C2-N2	-2.85	102.49	108.06
2	A	420	TET	O12-C7-C8	2.55	126.60	122.07
2	A	420	TET	O8-P2-O9	-2.54	99.94	109.47
2	A	420	TET	O22-C20-C16	-2.41	103.27	109.29
2	A	420	TET	O24-C17-C16	2.24	119.06	111.29
2	A	420	TET	O23-C16-C17	2.19	111.89	106.43
2	A	420	TET	O17-P3-O16	2.15	115.94	107.57
2	A	420	TET	O17-P3-O15	-2.11	102.23	110.53

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	420	TET	C12-O4-P1-O5
2	A	420	TET	C12-O4-P1-O6
2	A	420	TET	O23-C10-O8-P2
2	A	420	TET	C11-O13-P3-O15
2	A	420	TET	P2-O7-P1-O5
2	A	420	TET	C13-C18-N3-C7
2	A	420	TET	C12-O4-P1-O7

Continued on next page...

Continued from previous page...

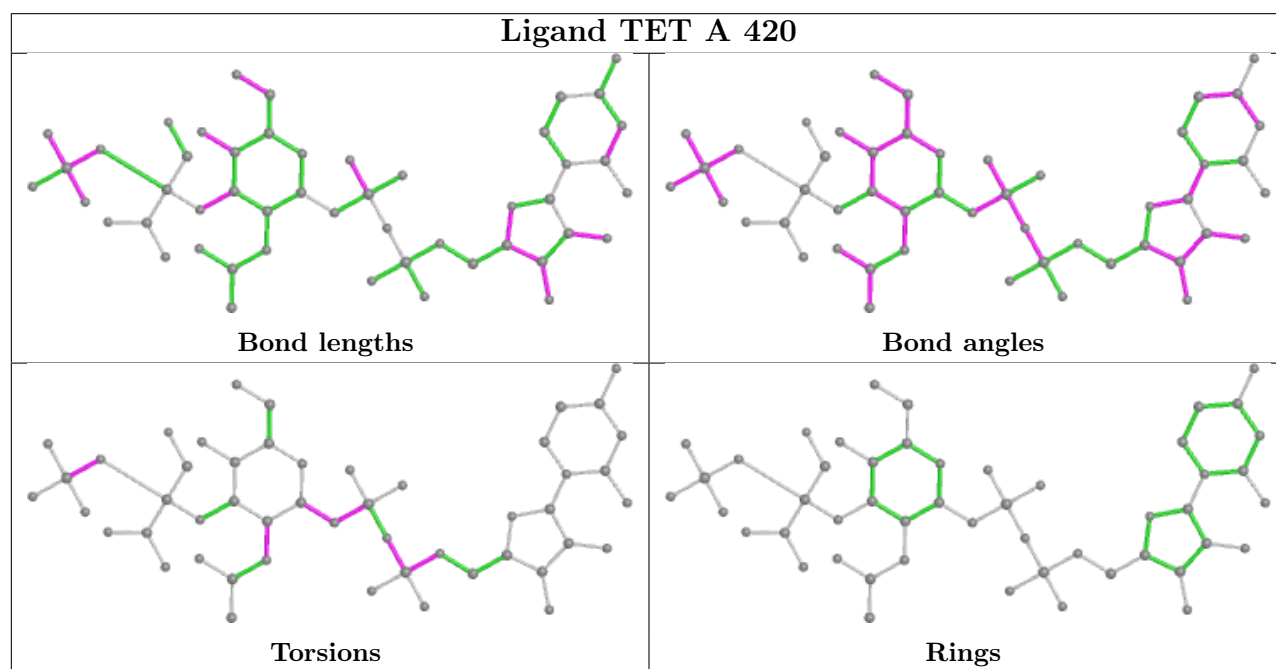
Mol	Chain	Res	Type	Atoms
2	A	420	TET	C10-O8-P2-O7
2	A	420	TET	C10-O8-P2-O9

There are no ring outliers.

1 monomer is involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	420	TET	5	0

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.