



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 13, 2026 – 06:38 AM UTC

PDB ID : 9C3Y / pdb_00009c3y
Title : Crystal structure of biphenyl synthase from *Malus domestica* complexed with tetraketide-CoA mimetic
Authors : Re, R.N.; Noel, J.P.; Burkart, M.D.
Deposited on : 2024-06-02
Resolution : 1.19 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)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

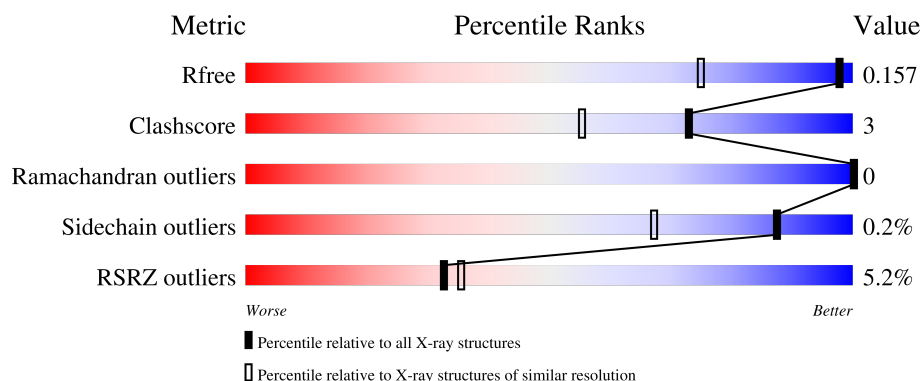
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.19 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}	180053	1216 (1.20-1.20)
Clashscore	190562	1265 (1.20-1.20)
Ramachandran outliers	187476	1226 (1.20-1.20)
Sidechain outliers	187428	1226 (1.20-1.20)
RSRZ outliers	180081	1214 (1.20-1.20)

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 $\leq 5\%$. 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	390	3% 92% 5% .
1	B	390	7% 91% 5% . .

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 13343 atoms, of which 6215 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.

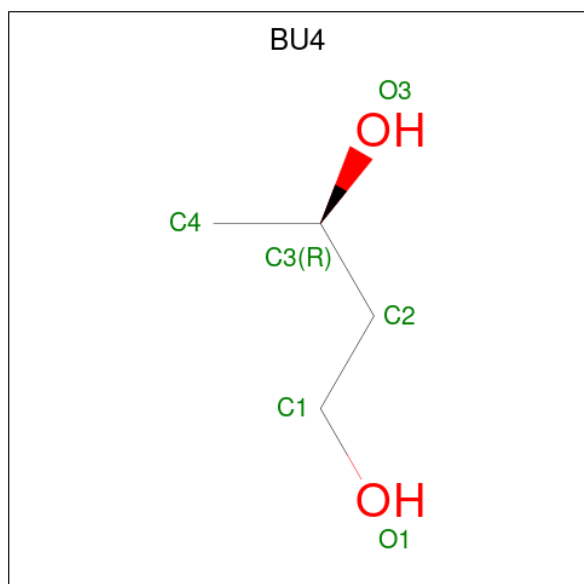
- Molecule 1 is a protein called BIS3 biphenyl synthase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	379	Total	C	H	N	O	S	0	15	0
			6107	1937	3070	522	560	18			
1	B	378	Total	C	H	N	O	S	0	14	0
			6057	1923	3041	517	559	17			

There are 4 discrepancies between the modelled and reference sequences:

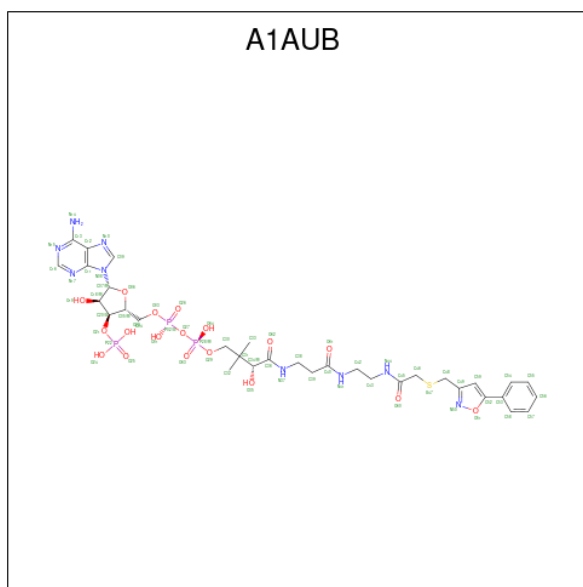
Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	expression tag	UNP K9MST3
A	0	SER	-	expression tag	UNP K9MST3
B	-1	GLY	-	expression tag	UNP K9MST3
B	0	SER	-	expression tag	UNP K9MST3

- Molecule 2 is (3R)-butane-1,3-diol (CCD ID: BU4) (formula: C₄H₁₀O₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	H	O	0	0
			16	4	10	2		
2	B	1	Total	C	H	O	0	0
			16	4	10	2		

- Molecule 3 is [(2R,3S,4R,5R)-5-(6-amino-9H-purin-9-yl)-4-hydroxy-3-(phosphonoxy)oxolan-2-yl]methyl (14R)-14-hydroxy-15,15-dimethyl-4,9,13-trioxo-1-(5-phenyl-1,2-oxazol-3-yl)-2-thia-5,8,12-triazahexadecan-16-yl dihydrogen diphosphate (CCD ID: A1AUB) (formula: $C_{33}H_{46}N_9O_{18}P_3S$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms							ZeroOcc	AltConf
3	A	1	Total	C	H	N	O	P	S	0	0
			106	33	42	9	18	3	1		
3	B	1	Total	C	H	N	O	P	S	0	0
			106	33	42	9	18	3	1		

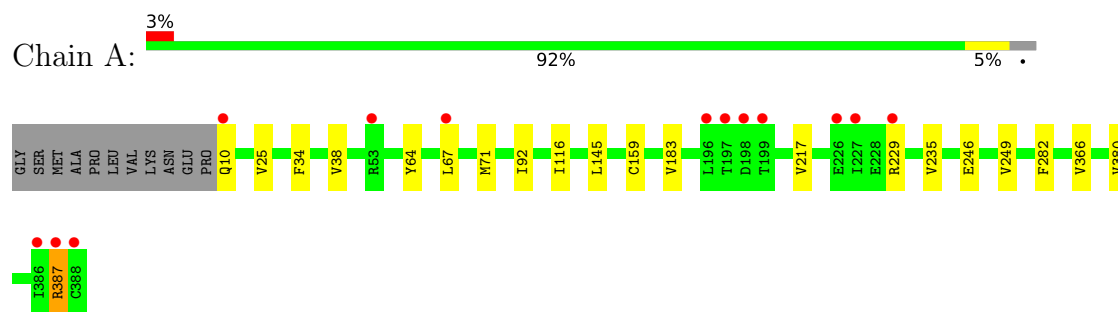
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	479	Total	O	0	0
			479	479		
4	B	456	Total	O	0	0
			456	456		

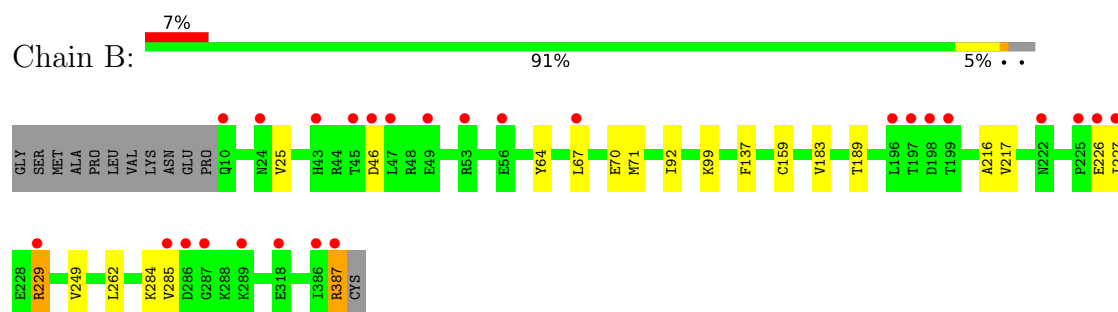
3 Residue-property plots [i](#)

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 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: BIS3 biphenyl synthase



- Molecule 1: BIS3 biphenyl synthase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	55.31Å 112.92Å 62.66Å 90.00° 93.19° 90.00°	Depositor
Resolution (Å)	54.72 – 1.19 54.72 – 1.19	Depositor EDS
% Data completeness (in resolution range)	97.8 (54.72-1.19) 92.1 (54.72-1.19)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.60 (at 1.19Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.151 , 0.163 (Not available) , 0.157	Depositor DCC
R_{free} test set	11989 reflections (4.90%)	wwPDB-VP
Wilson B-factor (Å ²)	8.8	Xtriage
Anisotropy	0.278	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 36.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	13343	wwPDB-VP
Average B, all atoms (Å ²)	15.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.52% 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: CSD, BU4, A1AUB

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	0.38	0/3126	0.61	0/4229
1	B	0.48	2/3102 (0.1%)	0.68	4/4198 (0.1%)
All	All	0.43	2/6228 (0.0%)	0.65	4/8427 (0.0%)

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	B	0	2

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	285[A]	VAL	C-O	5.34	1.29	1.23
1	B	285[B]	VAL	C-O	5.34	1.29	1.23

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	284[A]	LYS	CA-C-N	-5.31	117.72	123.08
1	B	284[A]	LYS	C-N-CA	-5.31	117.72	123.08
1	B	284[B]	LYS	CA-C-N	-5.31	117.72	123.08
1	B	284[B]	LYS	C-N-CA	-5.31	117.72	123.08

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	229	ARG	Sidechain
1	B	387	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	3037	3070	3089	18	0
1	B	3016	3041	3063	16	0
2	A	6	10	10	0	0
2	B	6	10	10	1	0
3	A	64	42	0	5	0
3	B	64	42	0	4	0
4	A	479	0	0	3	0
4	B	456	0	0	3	1
All	All	7128	6215	6172	39	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 3.

All (39) 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)
3:A:401:A1AUB:C05	3:A:401:A1AUB:O06	1.66	1.30
3:B:401:A1AUB:O06	3:B:401:A1AUB:C05	1.67	1.28
1:A:183:VAL:HB	1:A:217[B]:VAL:HG22	1.57	0.85
1:B:183:VAL:HB	1:B:217[B]:VAL:HG22	1.61	0.83
1:A:71[B]:MET:HE3	1:A:92:ILE:HD13	1.69	0.72
1:A:235[B]:VAL:HG22	1:A:380:VAL:O	1.90	0.72
1:A:10:GLN:N	4:A:501:HOH:O	2.26	0.68
1:B:25[B]:VAL:HG22	1:B:64:TYR:CE2	2.30	0.66
1:B:70:GLU:OE1	4:B:501:HOH:O	2.14	0.64
1:B:71[A]:MET:HE3	1:B:92:ILE:HD13	1.79	0.64
1:B:229:ARG:NH2	4:B:506:HOH:O	2.34	0.59
1:A:67:LEU:HD12	1:A:71[B]:MET:HE2	1.85	0.57
1:B:262:LEU:HD21	3:B:401:A1AUB:S47	2.44	0.57
3:A:401:A1AUB:O63	3:A:401:A1AUB:O35	2.23	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:249:VAL:HG13	3:B:401:A1AUB:S47	2.46	0.55
1:A:25[B]:VAL:HG22	1:A:64:TYR:CE2	2.43	0.54
1:B:67:LEU:HD12	1:B:71[A]:MET:HE2	1.90	0.52
1:A:249:VAL:HG13	3:A:401:A1AUB:S47	2.52	0.49
1:B:226:GLU:C	1:B:227:ILE:HD12	2.38	0.48
1:A:235[B]:VAL:CG2	1:A:380:VAL:HG12	2.44	0.47
3:A:401:A1AUB:O06	3:A:401:A1AUB:C04	2.53	0.47
1:B:46:ASP:N	1:B:46:ASP:OD1	2.47	0.47
1:B:99[B]:LYS:HG3	4:B:720:HOH:O	2.13	0.47
3:B:401:A1AUB:O06	3:B:401:A1AUB:C04	2.52	0.46
1:A:116:ILE:HG23	1:A:145:LEU:HD23	1.97	0.46
1:A:246[B]:GLU:HG2	1:B:137:PHE:HZ	1.81	0.46
3:A:401:A1AUB:O35	3:A:401:A1AUB:P28	2.74	0.46
1:A:235[B]:VAL:HG21	1:A:380:VAL:HG12	2.00	0.44
1:A:246[B]:GLU:HG3	4:A:842:HOH:O	2.17	0.44
1:A:282:PHE:HE1	1:A:366[B]:VAL:HG11	1.83	0.43
1:B:387:ARG:NH1	1:B:387:ARG:CG	2.81	0.43
1:A:235[B]:VAL:HG22	1:A:380:VAL:C	2.44	0.43
1:A:387:ARG:NH2	4:A:517:HOH:O	2.53	0.42
1:A:34:PHE:O	1:A:38:VAL:HG22	2.19	0.42
1:A:229:ARG:O	1:A:229:ARG:HD2	2.20	0.42
1:B:387:ARG:CG	1:B:387:ARG:HH11	2.33	0.41
1:B:189:THR:HB	2:B:400:BU4:H4B	2.03	0.41
1:B:183:VAL:O	1:B:216:ALA:HA	2.20	0.41
1:A:67:LEU:CD1	1:A:71[B]:MET:HE2	2.50	0.41

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)
4:B:916:HOH:O	4:B:931:HOH:O[1_554]	1.99	0.21

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	390/390 (100%)	382 (98%)	8 (2%)	0	100	100
1	B	388/390 (100%)	380 (98%)	8 (2%)	0	100	100
All	All	778/780 (100%)	762 (98%)	16 (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	173/328 (53%)	172 (99%)	1 (1%)	78	52
1	B	264/328 (80%)	264 (100%)	0	100	100
All	All	437/656 (67%)	436 (100%)	1 (0%)	87	67

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	387	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	252	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

2 non-standard protein/DNA/RNA residues 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
1	CSD	A	159[A]	1	4,7,8	1.39	1 (25%)	1,8,10	1.20	0
1	CSD	B	159[A]	1	4,7,8	1.93	1 (25%)	1,8,10	0.71	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
1	CSD	A	159[A]	1	-	0/2/6/8	-
1	CSD	B	159[A]	1	-	0/2/6/8	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	159[A]	CSD	OD1-SG	-3.28	1.44	1.47
1	A	159[A]	CSD	OD1-SG	-2.06	1.45	1.47

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

4 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$
3	A1AUB	B	401	-	66,68,68	3.06	24 (36%)	91,99,99	2.97	22 (24%)
3	A1AUB	A	401	-	66,68,68	3.05	24 (36%)	91,99,99	2.90	26 (28%)
2	BU4	A	400	-	5,5,5	0.28	0	5,5,5	0.82	0
2	BU4	B	400	-	5,5,5	0.30	0	5,5,5	0.13	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
3	A1AUB	B	401	-	-	11/59/77/77	0/5/5/5
3	A1AUB	A	401	-	-	17/59/77/77	0/5/5/5
2	BU4	A	400	-	-	1/3/3/3	-
2	BU4	B	400	-	-	0/3/3/3	-

All (48) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	401	A1AUB	O06-C05	10.01	1.67	1.45
3	A	401	A1AUB	O06-C05	9.56	1.66	1.45
3	A	401	A1AUB	C36-N37	7.50	1.51	1.33
3	A	401	A1AUB	C20-C05	-7.28	1.34	1.52
3	B	401	A1AUB	C20-C05	-7.24	1.34	1.52
3	B	401	A1AUB	C36-N37	7.12	1.50	1.33
3	B	401	A1AUB	O51-C52	-7.05	1.29	1.36
3	A	401	A1AUB	O51-C52	-6.97	1.29	1.36
3	A	401	A1AUB	P02-O27	6.52	1.66	1.59
3	B	401	A1AUB	P02-O27	6.51	1.66	1.59
3	A	401	A1AUB	C40-N41	5.82	1.47	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	401	A1AUB	O06-C07	-5.62	1.29	1.42
3	B	401	A1AUB	C40-N41	5.46	1.46	1.33
3	B	401	A1AUB	C45-N44	5.45	1.46	1.33
3	A	401	A1AUB	C45-N44	5.41	1.46	1.33
3	B	401	A1AUB	P28-O27	5.19	1.65	1.59
3	A	401	A1AUB	P28-O27	5.13	1.65	1.59
3	B	401	A1AUB	O06-C07	-5.10	1.30	1.42
3	A	401	A1AUB	C13-N14	5.08	1.47	1.34
3	B	401	A1AUB	C13-N14	5.05	1.47	1.34
3	B	401	A1AUB	O51-N50	-5.03	1.33	1.42
3	A	401	A1AUB	O51-N50	-5.00	1.33	1.42
3	A	401	A1AUB	P22-O21	4.73	1.67	1.59
3	B	401	A1AUB	P22-O21	4.71	1.67	1.59
3	B	401	A1AUB	O19-C18	-4.23	1.32	1.43
3	B	401	A1AUB	C46-C45	4.10	1.56	1.51
3	A	401	A1AUB	O19-C18	-3.80	1.33	1.43
3	A	401	A1AUB	C46-C45	3.73	1.56	1.51
3	B	401	A1AUB	C59-C52	3.50	1.40	1.35
3	A	401	A1AUB	C59-C52	3.45	1.40	1.35
3	B	401	A1AUB	O62-C36	-3.30	1.17	1.23
3	A	401	A1AUB	O62-C36	-3.05	1.17	1.23
3	A	401	A1AUB	C30-C31	3.00	1.57	1.52
3	B	401	A1AUB	C53-C52	2.97	1.54	1.48
3	A	401	A1AUB	C53-C52	2.97	1.54	1.48
3	B	401	A1AUB	C18-C07	2.94	1.62	1.53
3	B	401	A1AUB	C09-N08	-2.76	1.32	1.37
3	A	401	A1AUB	C18-C07	2.71	1.62	1.53
3	A	401	A1AUB	C09-N08	-2.69	1.33	1.37
3	B	401	A1AUB	O61-C40	-2.61	1.18	1.23
3	B	401	A1AUB	C30-C31	2.54	1.56	1.52
3	A	401	A1AUB	O61-C40	-2.31	1.18	1.23
3	A	401	A1AUB	O60-C45	-2.28	1.18	1.23
3	B	401	A1AUB	C12-N10	-2.23	1.35	1.39
3	A	401	A1AUB	P28-O29	2.14	1.67	1.59
3	A	401	A1AUB	C12-N10	-2.13	1.35	1.39
3	B	401	A1AUB	P28-O29	2.06	1.67	1.59
3	B	401	A1AUB	O60-C45	-2.04	1.19	1.23

All (48) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	401	A1AUB	O51-C52-C53	18.36	132.29	116.74

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	401	A1AUB	O51-C52-C53	18.16	132.12	116.74
3	B	401	A1AUB	C49-C48-S47	-11.33	101.91	113.60
3	A	401	A1AUB	C49-C48-S47	-8.89	104.43	113.60
3	A	401	A1AUB	C53-C52-C59	-6.49	118.47	133.12
3	B	401	A1AUB	C53-C52-C59	-6.38	118.71	133.12
3	A	401	A1AUB	C12-C11-N17	-6.16	118.23	126.72
3	B	401	A1AUB	C12-C11-N17	-5.99	118.47	126.72
3	B	401	A1AUB	C48-S47-C46	5.50	113.37	100.69
3	A	401	A1AUB	C48-S47-C46	4.89	111.96	100.69
3	B	401	A1AUB	C59-C49-N50	-4.72	106.88	111.41
3	A	401	A1AUB	C59-C49-N50	-4.68	106.92	111.41
3	A	401	A1AUB	N17-C11-N08	4.41	134.66	127.17
3	A	401	A1AUB	C58-C53-C52	-4.35	113.85	120.77
3	B	401	A1AUB	C58-C53-C52	-4.31	113.92	120.77
3	B	401	A1AUB	N17-C16-N15	-4.23	122.18	128.58
3	B	401	A1AUB	N17-C11-N08	4.20	134.31	127.17
3	A	401	A1AUB	N17-C16-N15	-4.18	122.25	128.58
3	A	401	A1AUB	C16-N17-C11	3.86	121.26	111.83
3	B	401	A1AUB	C16-N17-C11	3.86	121.26	111.83
3	B	401	A1AUB	C42-N41-C40	-3.34	116.60	122.82
3	B	401	A1AUB	C11-C12-N10	-3.16	106.97	110.58
3	A	401	A1AUB	O51-N50-C49	3.16	109.76	105.74
3	B	401	A1AUB	N08-C09-N10	-3.10	109.54	113.94
3	B	401	A1AUB	O51-N50-C49	3.08	109.66	105.74
3	A	401	A1AUB	C11-C12-N10	-3.07	107.07	110.58
3	B	401	A1AUB	C12-N10-C09	3.03	108.22	103.45
3	A	401	A1AUB	N08-C09-N10	-3.02	109.66	113.94
3	A	401	A1AUB	C12-N10-C09	2.98	108.13	103.45
3	A	401	A1AUB	C20-C18-C07	2.85	106.15	99.89
3	B	401	A1AUB	O06-C07-N08	2.76	113.38	108.09
3	B	401	A1AUB	C52-O51-N50	2.74	110.11	108.73
3	B	401	A1AUB	C54-C53-C52	2.69	125.04	120.77
3	A	401	A1AUB	C18-C07-N08	-2.53	107.02	113.30
3	B	401	A1AUB	C39-C38-N37	-2.51	106.65	112.00
3	A	401	A1AUB	C54-C53-C52	2.50	124.75	120.77
3	A	401	A1AUB	C38-C39-C40	2.27	116.17	112.39
3	B	401	A1AUB	C11-N08-C09	2.21	108.06	105.74
3	A	401	A1AUB	C11-N08-C09	2.19	108.03	105.74
3	A	401	A1AUB	C58-C53-C54	2.17	121.33	118.57
3	A	401	A1AUB	O06-C05-C20	2.15	109.45	104.92
3	A	401	A1AUB	C18-C20-C05	2.15	107.00	103.24
3	A	401	A1AUB	O06-C07-N08	2.14	112.20	108.09

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	401	A1AUB	C52-O51-N50	2.14	109.81	108.73
3	A	401	A1AUB	C43-N44-C45	-2.12	118.89	122.82
3	A	401	A1AUB	P22-O21-C20	-2.06	117.92	123.43
3	B	401	A1AUB	C48-C49-C59	-2.05	123.20	128.94
3	B	401	A1AUB	C55-C54-C53	-2.03	118.36	120.36

There are no chirality outliers.

All (29) torsion outliers are listed below:

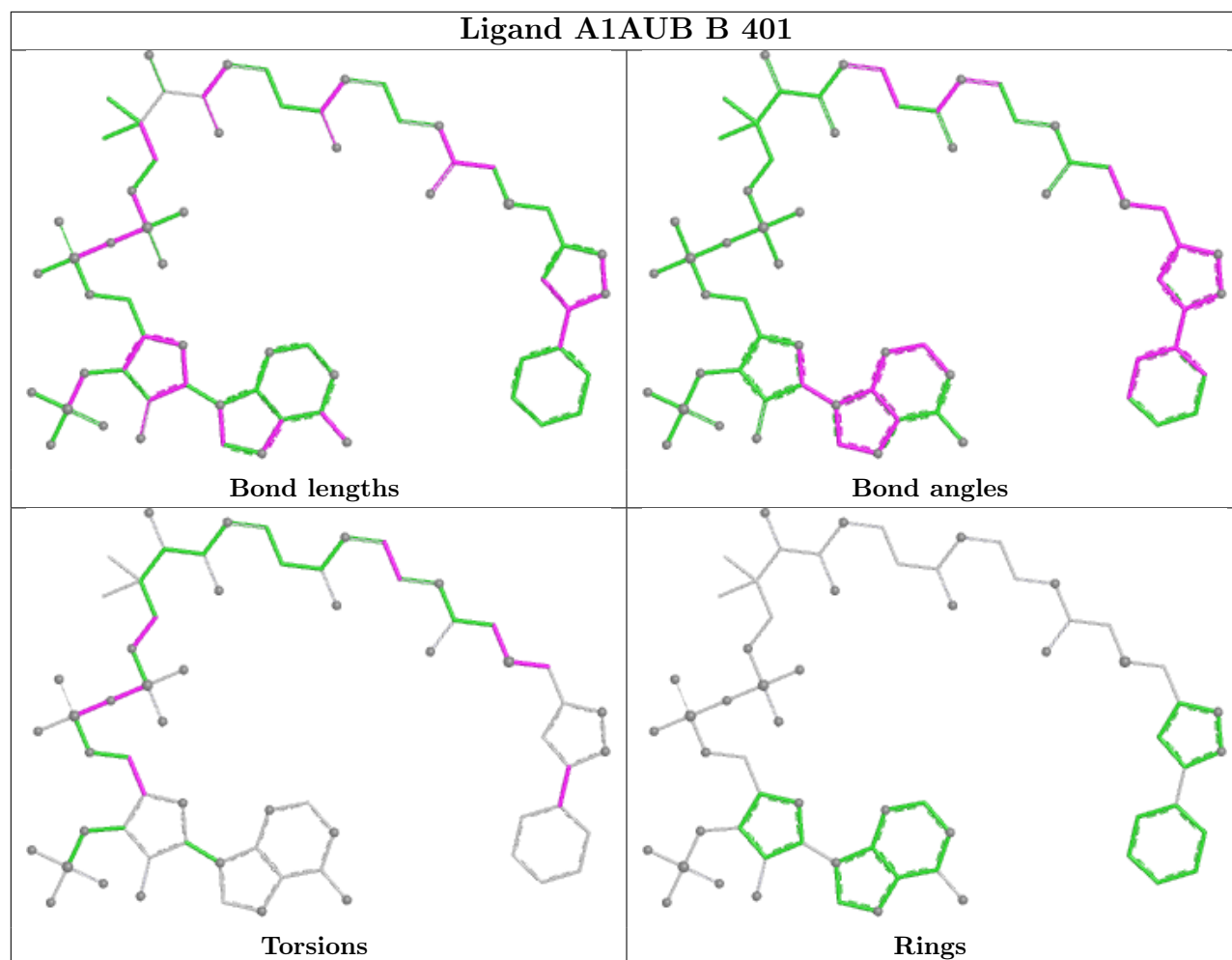
Mol	Chain	Res	Type	Atoms
3	A	401	A1AUB	O29-C30-C31-C32
3	A	401	A1AUB	O29-C30-C31-C33
3	A	401	A1AUB	O29-C30-C31-C34
3	A	401	A1AUB	C31-C30-O29-P28
3	A	401	A1AUB	C49-C48-S47-C46
3	A	401	A1AUB	C30-O29-P28-O27
3	A	401	A1AUB	C30-O29-P28-O63
3	A	401	A1AUB	C30-O29-P28-O64
3	B	401	A1AUB	C49-C48-S47-C46
3	A	401	A1AUB	C59-C52-C53-C54
3	A	401	A1AUB	O51-C52-C53-C54
3	A	401	A1AUB	C59-C52-C53-C58
3	A	401	A1AUB	O51-C52-C53-C58
3	A	401	A1AUB	N41-C42-C43-N44
3	B	401	A1AUB	C59-C52-C53-C54
3	B	401	A1AUB	N41-C42-C43-N44
3	A	401	A1AUB	C45-C46-S47-C48
3	B	401	A1AUB	C45-C46-S47-C48
3	B	401	A1AUB	C59-C52-C53-C58
3	B	401	A1AUB	P28-O27-P02-O26
3	B	401	A1AUB	O03-C04-C05-O06
2	A	400	BU4	O1-C1-C2-C3
3	A	401	A1AUB	C20-O21-P22-O24
3	B	401	A1AUB	P02-O27-P28-O63
3	B	401	A1AUB	C31-C30-O29-P28
3	B	401	A1AUB	O51-C52-C53-C54
3	A	401	A1AUB	O35-C34-C36-N37
3	A	401	A1AUB	P02-O27-P28-O64
3	B	401	A1AUB	P28-O27-P02-O01

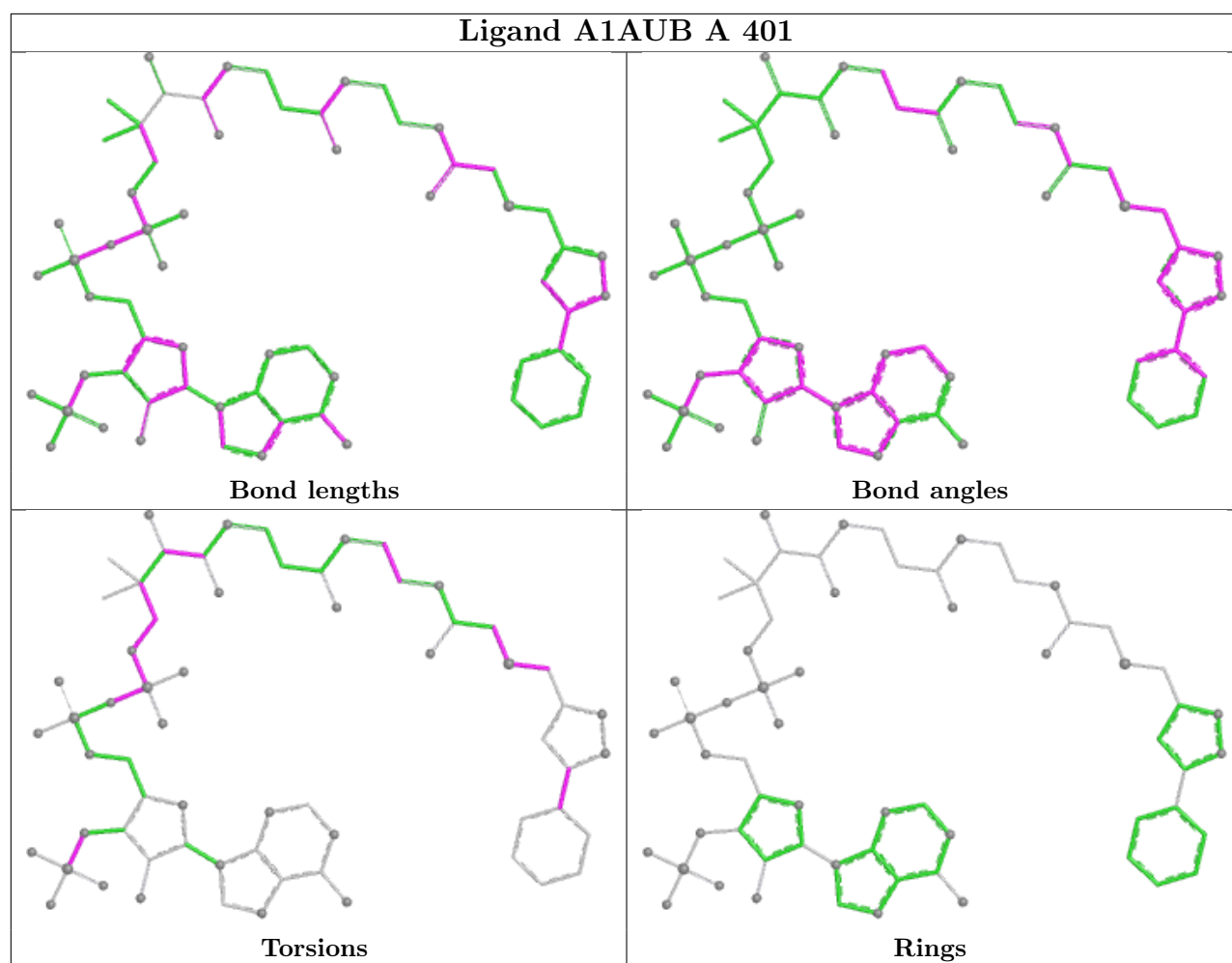
There are no ring outliers.

3 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	401	A1AUB	4	0
3	A	401	A1AUB	5	0
2	B	400	BU4	1	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 [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2		OWAB(Å ²)	Q<0.9
1	A	378/390 (96%)	0.21	13 (3%)	48 48	4, 11, 26, 70	14 (3%)
1	B	377/390 (96%)	0.30	26 (6%)	23 24	5, 11, 32, 60	13 (3%)
All	All	755/780 (96%)	0.26	39 (5%)	33 35	4, 11, 30, 70	27 (3%)

All (39) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	388	CYS	9.9
1	B	229	ARG	5.0
1	A	387	ARG	4.9
1	B	285[A]	VAL	4.8
1	B	386	ILE	4.7
1	B	225	PRO	4.3
1	B	387	ARG	4.3
1	B	227	ILE	3.9
1	B	45	THR	3.5
1	B	286	ASP	3.5
1	A	386	ILE	3.4
1	A	227	ILE	3.4
1	B	10	GLN	3.3
1	B	196	LEU	3.2
1	A	229	ARG	3.0
1	B	226	GLU	3.0
1	A	196	LEU	3.0
1	B	46	ASP	2.9
1	A	10	GLN	2.9
1	A	197	THR	2.8
1	B	222	ASN	2.8
1	B	289	LYS	2.7
1	B	49	GLU	2.6
1	A	226	GLU	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	67	LEU	2.5
1	B	53	ARG	2.5
1	B	287	GLY	2.5
1	B	24[A]	ASN	2.4
1	B	47	LEU	2.4
1	B	43	HIS	2.4
1	A	53[A]	ARG	2.4
1	B	197	THR	2.3
1	B	199	THR	2.3
1	A	199	THR	2.3
1	B	318	GLU	2.2
1	B	198	ASP	2.2
1	A	67	LEU	2.1
1	B	56	GLU	2.1
1	A	198	ASP	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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. 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	B-factors(Å ²)	Q<0.9
1	CSD	A	159[A]	8/9	0.96	0.08	8,12,15,25	0
1	CSD	B	159[A]	8/9	0.97	0.08	8,11,14,24	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides 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. 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	B-factors(Å ²)	Q<0.9
3	A1AUB	A	401	64/64	0.56	0.21	20,52,98,115	0

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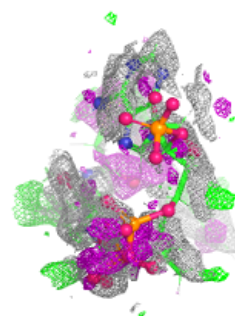
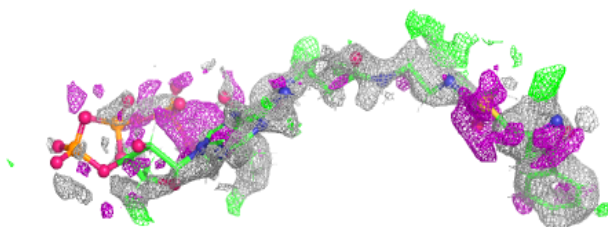
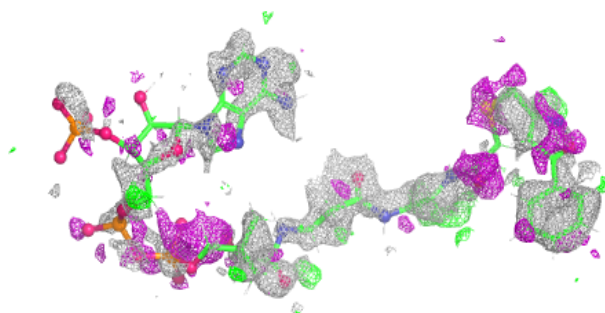
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	A1AUB	B	401	64/64	0.58	0.20	20,51,91,110	0
2	BU4	B	400	6/6	0.90	0.12	14,24,29,29	0
2	BU4	A	400	6/6	0.92	0.10	15,19,23,23	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

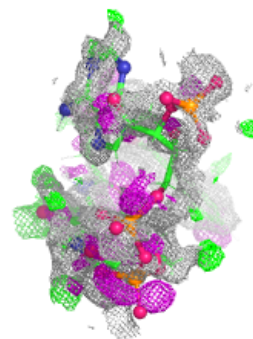
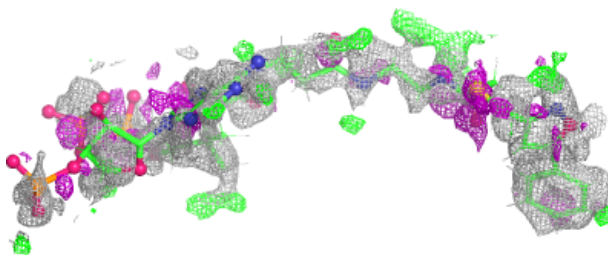
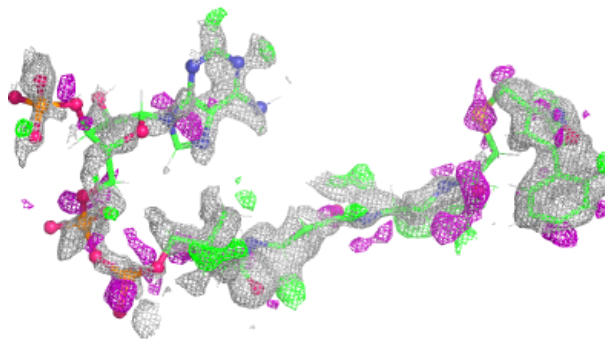
Electron density around A1AUB A 401:

2mF_o-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around A1AUB B 401:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.