



Full wwPDB X-ray Structure Validation Report ⓘ

Jul 15, 2026 – 02:07 PM EDT

PDB ID : 9Y7T / pdb_00009y7t
Title : Crystal structure of Candida auris dihydrofolate reductase in complex with inhibitor 1172 and NADPH
Authors : Erlandsen, H.; Krucinska, J.; Wright, D.
Deposited on : 2025-09-11
Resolution : 1.98 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	:	9.0.015 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.50

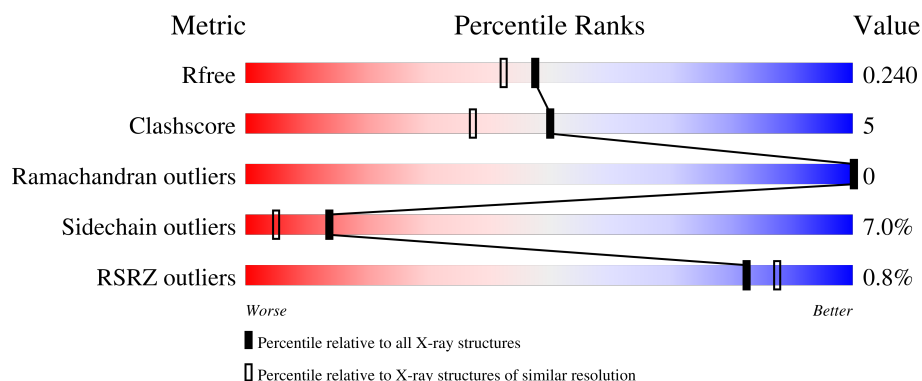
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.98 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	1506 (1.98-1.98)
Clashscore	190562	1534 (1.98-1.98)
Ramachandran outliers	187476	1518 (1.98-1.98)
Sidechain outliers	187428	1518 (1.98-1.98)
RSRZ outliers	180081	1506 (1.98-1.98)

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	204	 85% 11% ..
1	B	204	 69% 22% 7% ..

2 Entry composition [i](#)

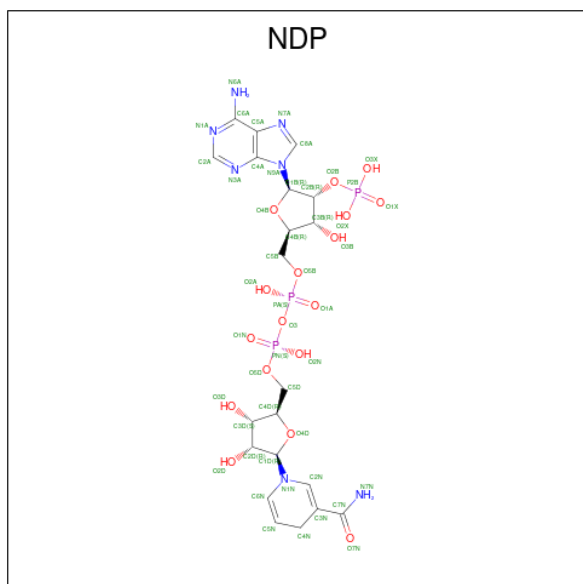
There are 6 unique types of molecules in this entry. The entry contains 3643 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Dihydrofolate reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	200	Total	C	N	O	S	0	0	0
			1632	1047	280	298	7			
1	B	200	Total	C	N	O	S	0	1	0
			1635	1049	280	299	7			

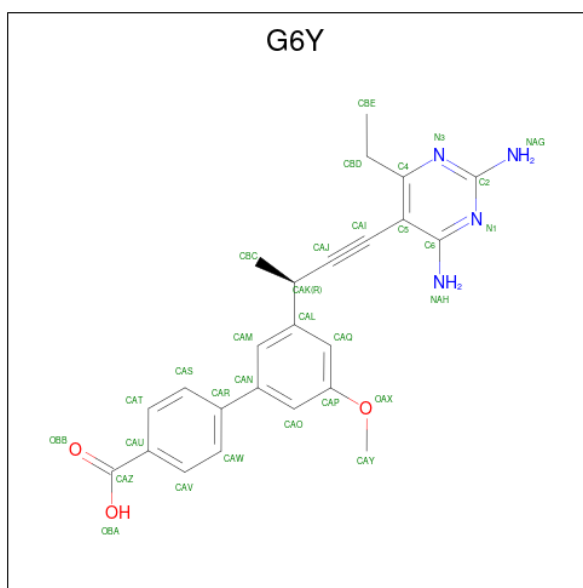
- Molecule 2 is NADPH DIHYDRO-NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (CCD ID: NDP) (formula: $C_{21}H_{30}N_7O_{17}P_3$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
2	B	1	Total	C	N	O	P	0	0
			48	21	7	17	3		

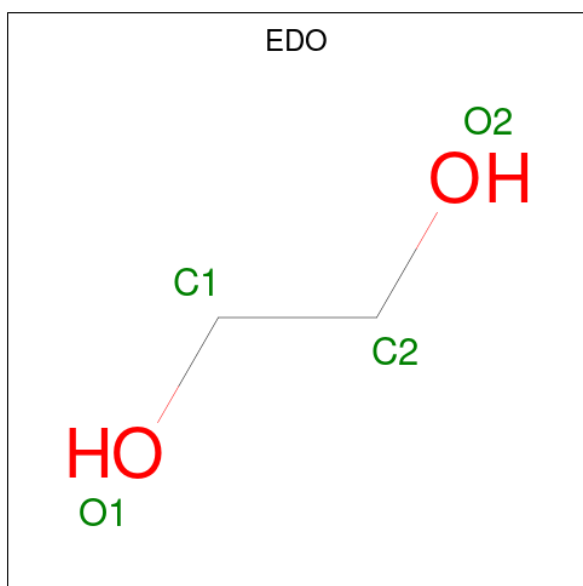
- Molecule 3 is 3'-[(2R)-4-(2,4-diamino-6-ethylpyrimidin-5-yl)but-3-yn-2-yl]-5'-methoxy[1,1'-biphenyl]-4-carboxylic acid (CCD ID: G6Y) (formula: $C_{24}H_{24}N_4O_3$) (labeled as "Ligand of

Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total 31	C 24	N 4	O 3	0	0
3	B	1	Total 31	C 24	N 4	O 3	0	0

- Molecule 4 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $\text{C}_2\text{H}_6\text{O}_2$).



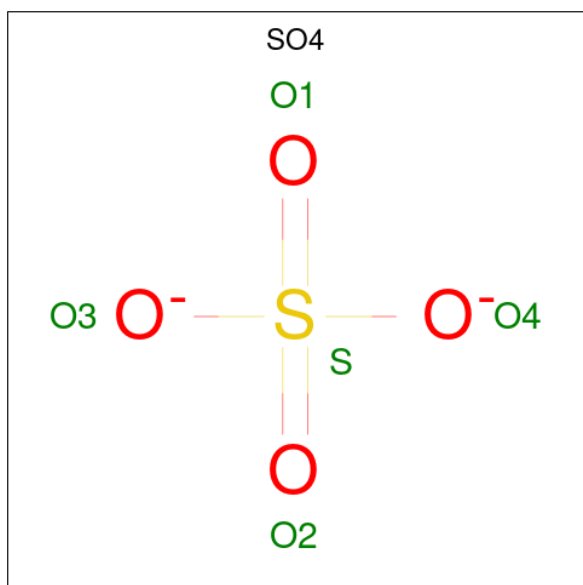
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			4	2	2		

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

- Molecule 5 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	O	S	0	0
			5	4	1		
5	A	1	Total	O	S	0	0
			5	4	1		
5	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	124	Total	O	0	0
			124	124		
6	B	67	Total	O	0	0
			67	67		

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.

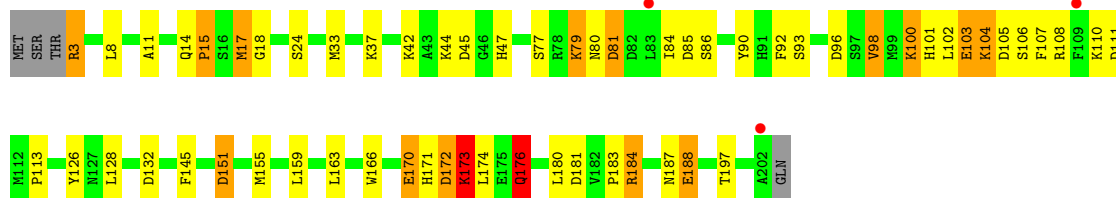
- Molecule 1: Dihydrofolate reductase

Chain A: 



- Molecule 1: Dihydrofolate reductase

Chain B: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	65.15Å 72.34Å 104.91Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.72 – 1.98 29.72 – 1.98	Depositor EDS
% Data completeness (in resolution range)	99.8 (29.72-1.98) 99.8 (29.72-1.98)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.78 (at 1.98Å)	Xtriage
Refinement program	REFMAC 5.8.0425	Depositor
R, R_{free}	0.168 , 0.233 0.180 , 0.240	Depositor DCC
R_{free} test set	1703 reflections (4.83%)	wwPDB-VP
Wilson B-factor (Å ²)	35.3	Xtriage
Anisotropy	0.084	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 38.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	3643	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.95% 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: EDO, G6Y, NDP, SO4

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	1.00	0/1670	1.46	11/2251 (0.5%)
1	B	0.93	0/1676	1.60	24/2259 (1.1%)
All	All	0.97	0/3346	1.53	35/4510 (0.8%)

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	5
1	B	0	3
All	All	0	8

There are no bond length outliers.

All (35) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	17	MET	CG-SD-CE	-11.50	75.59	100.90
1	A	202	ALA	CA-C-O	-10.87	102.33	120.80
1	A	22	LYS	CB-CA-C	-8.55	100.00	111.63
1	B	96	ASP	CA-CB-CG	8.18	120.78	112.60
1	B	188	GLU	CB-CA-C	-7.94	96.11	109.53
1	B	184	ARG	CB-CA-C	7.68	122.48	109.50
1	B	173	LYS	N-CA-CB	7.20	121.39	110.30
1	B	103	GLU	CA-C-N	7.16	130.83	120.38
1	B	103	GLU	C-N-CA	7.16	130.83	120.38
1	B	81	ASP	CB-CA-C	-7.10	101.65	111.73
1	B	197	THR	CA-CB-OG1	6.94	120.02	109.60
1	B	151	ASP	CA-CB-CG	6.78	119.38	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	176	GLN	N-CA-CB	-6.67	100.31	110.12
1	B	85	ASP	CA-CB-CG	6.60	119.20	112.60
1	B	100	LYS	N-CA-CB	6.50	119.41	109.98
1	B	15	PRO	N-CA-CB	-6.48	95.47	102.60
1	A	33	MET	CG-SD-CE	-6.38	86.87	100.90
1	B	176	GLN	CB-CA-C	6.20	121.08	110.79
1	A	23	GLY	CA-C-N	-5.71	113.34	122.53
1	A	23	GLY	C-N-CA	-5.71	113.34	122.53
1	B	170	GLU	CB-CG-CD	5.70	122.29	112.60
1	B	92	PHE	CA-CB-CG	-5.58	108.22	113.80
1	B	98	VAL	N-CA-CB	5.55	118.91	110.58
1	B	14	GLN	CB-CA-C	5.50	117.35	109.26
1	A	184	ARG	CB-CA-C	-5.50	100.72	109.80
1	A	66	ARG	CB-CA-C	-5.47	101.26	108.76
1	A	101	HIS	CA-CB-CG	-5.43	108.36	113.80
1	B	80	ASN	N-CA-C	5.41	118.34	111.69
1	B	96	ASP	CB-CA-C	-5.37	102.44	110.88
1	B	132	ASP	CA-CB-CG	5.34	117.94	112.60
1	A	14	GLN	CG-CD-NE2	-5.31	108.43	116.40
1	B	188	GLU	N-CA-CB	5.31	118.07	110.06
1	A	107	PHE	CA-CB-CG	-5.25	108.55	113.80
1	B	126	TYR	CB-CA-C	-5.02	102.46	110.79
1	A	168	ARG	CG-CD-NE	-5.00	100.99	112.00

There are no chirality outliers.

All (8) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	108	ARG	Sidechain
1	A	184	ARG	Sidechain
1	A	3	ARG	Sidechain
1	A	55	ARG	Sidechain
1	A	66	ARG	Sidechain
1	B	108	ARG	Sidechain
1	B	3	ARG	Sidechain
1	B	81	ASP	Peptide

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	1632	0	1631	6	0
1	B	1635	0	1636	30	0
2	A	48	0	26	2	0
2	B	48	0	26	2	0
3	A	31	0	0	0	0
3	B	31	0	0	1	0
4	A	8	0	12	0	0
4	B	4	0	6	0	0
5	A	10	0	0	0	0
5	B	5	0	0	0	0
6	A	124	0	0	1	1
6	B	67	0	0	4	0
All	All	3643	0	3337	36	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 5.

All (36) 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:166:TRP:O	6:B:401:HOH:O	1.94	0.85
1:B:47:HIS:ND1	1:B:113:PRO:HG2	1.94	0.82
1:A:11:ALA:H	2:A:301:NDP:H72N	1.29	0.77
1:B:11:ALA:H	2:B:301:NDP:H72N	1.43	0.66
1:B:183:PRO:HB2	1:B:187:ASN:HD21	1.60	0.66
1:B:176:GLN:O	1:B:176:GLN:NE2	2.32	0.63
1:B:172:ASP:O	1:B:176:GLN:HB2	2.02	0.59
1:B:47:HIS:ND1	1:B:113:PRO:CG	2.64	0.59
1:B:101:HIS:C	1:B:101:HIS:CD2	2.81	0.59
1:B:170:GLU:H	1:B:173:LYS:HE3	1.68	0.58
1:B:45:ASP:OD1	6:B:402:HOH:O	2.17	0.56
1:B:171:HIS:O	1:B:174:LEU:HB3	2.06	0.55
1:B:42:LYS:HG2	1:B:180:LEU:HD21	1.89	0.54
1:B:104:LYS:H	1:B:104:LYS:HE2	1.72	0.54
1:A:176:GLN:HE21	1:B:128:LEU:HD11	1.75	0.52
1:B:103:GLU:HG3	1:B:107:PHE:CE2	2.45	0.52
1:B:103:GLU:HG3	1:B:107:PHE:CZ	2.45	0.52
1:B:145:PHE:HZ	1:B:151:ASP:HB2	1.77	0.50
1:B:44:LYS:HB2	1:B:47:HIS:CD2	2.47	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:301:NDP:O3X	6:B:403:HOH:O	2.20	0.47
1:B:176:GLN:NE2	1:B:176:GLN:C	2.73	0.47
1:A:42:LYS:HE3	6:A:433:HOH:O	2.15	0.46
1:B:24:SER:HB2	6:B:438:HOH:O	2.16	0.46
1:A:81:ASP:HB2	1:A:83:LEU:HD23	1.99	0.45
1:B:37:LYS:HG3	3:B:302:G6Y:OBB	2.16	0.45
1:B:77:SER:O	1:B:93:SER:HA	2.16	0.45
1:B:110:LYS:O	1:B:111:ASP:C	2.60	0.45
1:B:98:VAL:O	1:B:102:LEU:HG	2.17	0.44
1:B:79:LYS:H	1:B:79:LYS:HG2	1.60	0.43
1:B:18:GLY:HA3	1:B:155:MET:HE3	2.01	0.42
1:B:84:ILE:HD13	1:B:90:TYR:CE1	2.55	0.42
1:B:163:LEU:HD23	1:B:163:LEU:HA	1.92	0.42
1:A:30:LYS:HB3	1:A:30:LYS:HE3	1.75	0.41
1:A:19:ILE:O	2:A:301:NDP:H2N	2.21	0.41
1:B:42:LYS:HG2	1:B:180:LEU:CD2	2.50	0.41
1:B:104:LYS:HE2	1:B:104:LYS:N	2.35	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)
6:A:510:HOH:O	6:A:521:HOH:O[4_445]	2.10	0.10

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	198/204 (97%)	194 (98%)	4 (2%)	0	100	100
1	B	199/204 (98%)	192 (96%)	7 (4%)	0	100	100
All	All	397/408 (97%)	386 (97%)	11 (3%)	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	179/183 (98%)	172 (96%)	7 (4%)	28	17
1	B	180/183 (98%)	161 (89%)	19 (11%)	6	1
All	All	359/366 (98%)	333 (93%)	26 (7%)	14	5

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	LYS
1	A	96	ASP
1	A	104	LYS
1	A	154	GLN
1	A	159	LEU
1	A	174	LEU
1	A	201	LYS
1	B	3	ARG
1	B	8	LEU
1	B	15	PRO
1	B	17	MET
1	B	33	MET
1	B	79	LYS
1	B	86[A]	SER
1	B	86[B]	SER
1	B	100	LYS
1	B	104	LYS
1	B	105	ASP
1	B	106	SER
1	B	159	LEU
1	B	172	ASP
1	B	173	LYS
1	B	176	GLN
1	B	181	ASP
1	B	184	ARG

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Mol	Chain	Res	Type
1	B	188	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	14	GLN
1	A	31	ASN
1	A	47	HIS
1	A	87	ASN
1	A	101	HIS
1	A	176	GLN
1	B	101	HIS
1	B	176	GLN
1	B	187	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](#)

10 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	NDP	A	301	-	51,52,52	1.13	2 (3%)	71,80,80	1.51	12 (16%)
4	EDO	A	303	-	3,3,3	0.35	0	2,2,2	0.60	0
3	G6Y	A	302	-	32,33,33	1.17	2 (6%)	43,46,46	1.26	3 (6%)
4	EDO	A	304	-	3,3,3	0.25	0	2,2,2	0.22	0
5	SO4	A	306	-	4,4,4	0.29	0	6,6,6	0.14	0
3	G6Y	B	302	-	32,33,33	1.44	2 (6%)	43,46,46	1.04	2 (4%)
5	SO4	B	304	-	4,4,4	0.29	0	6,6,6	0.26	0
5	SO4	A	305	-	4,4,4	0.24	0	6,6,6	0.04	0
2	NDP	B	301	-	51,52,52	1.09	3 (5%)	71,80,80	1.08	7 (9%)
4	EDO	B	303	-	3,3,3	0.46	0	2,2,2	0.21	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	NDP	A	301	-	-	2/34/77/77	0/5/5/5
4	EDO	A	303	-	-	0/1/1/1	-
3	G6Y	A	302	-	-	4/18/21/21	0/3/3/3
4	EDO	A	304	-	-	0/1/1/1	-
3	G6Y	B	302	-	-	4/18/21/21	0/3/3/3
2	NDP	B	301	-	-	1/34/77/77	0/5/5/5
4	EDO	B	303	-	-	1/1/1/1	-

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	302	G6Y	C5-C4	-6.21	1.33	1.40
2	A	301	NDP	P2B-O2B	5.23	1.68	1.59
2	B	301	NDP	P2B-O2B	4.34	1.67	1.59
3	A	302	G6Y	C5-C4	-4.20	1.35	1.40
2	B	301	NDP	P2B-O2X	-3.39	1.42	1.54
3	B	302	G6Y	OBA-CAZ	-3.38	1.20	1.30
2	B	301	NDP	PN-O3	-2.74	1.56	1.59
2	A	301	NDP	PN-O1N	-2.20	1.43	1.50
3	A	302	G6Y	OBA-CAZ	-2.17	1.24	1.30

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	302	G6Y	CBD-C4-C5	4.62	125.73	122.54
2	A	301	NDP	O3-PN-O1N	-4.50	97.17	110.70
2	A	301	NDP	C3N-C7N-N7N	-4.40	109.85	117.67
2	A	301	NDP	O2B-P2B-O1X	-3.26	97.71	109.33
2	A	301	NDP	O2N-PN-O1N	3.06	126.67	112.44
2	B	301	NDP	O2B-P2B-O1X	-2.94	98.84	109.33
2	A	301	NDP	C6N-N1N-C2N	2.91	122.43	119.32
2	A	301	NDP	C1D-N1N-C2N	-2.84	116.45	121.14
2	B	301	NDP	O3D-C3D-C2D	-2.71	103.12	111.82
3	A	302	G6Y	C2-N3-C4	2.65	118.19	116.26
3	B	302	G6Y	CBE-CBD-C4	-2.61	107.54	114.59
2	A	301	NDP	O3X-P2B-O2X	2.55	117.38	107.80
2	B	301	NDP	O3-PN-O1N	-2.50	103.19	110.70
2	A	301	NDP	O2X-P2B-O2B	-2.39	96.54	105.85
3	A	302	G6Y	CBC-CAK-CAJ	-2.35	106.72	110.09
2	A	301	NDP	O3X-P2B-O1X	2.34	119.96	110.83
2	A	301	NDP	P2B-O2B-C2B	-2.29	117.31	123.43
2	A	301	NDP	O7N-C7N-C3N	2.28	125.18	120.90
2	B	301	NDP	O3B-C3B-C4B	-2.15	104.92	111.08
2	B	301	NDP	O7N-C7N-C3N	2.11	124.88	120.90
2	A	301	NDP	O3-PA-O1A	-2.08	104.45	110.70
2	B	301	NDP	O3B-C3B-C2B	2.05	116.92	111.19
2	B	301	NDP	O3X-P2B-O2X	2.01	115.32	107.80
3	B	302	G6Y	OAX-CAP-CAO	-2.00	110.45	120.00

There are no chirality outliers.

All (12) torsion outliers are listed below:

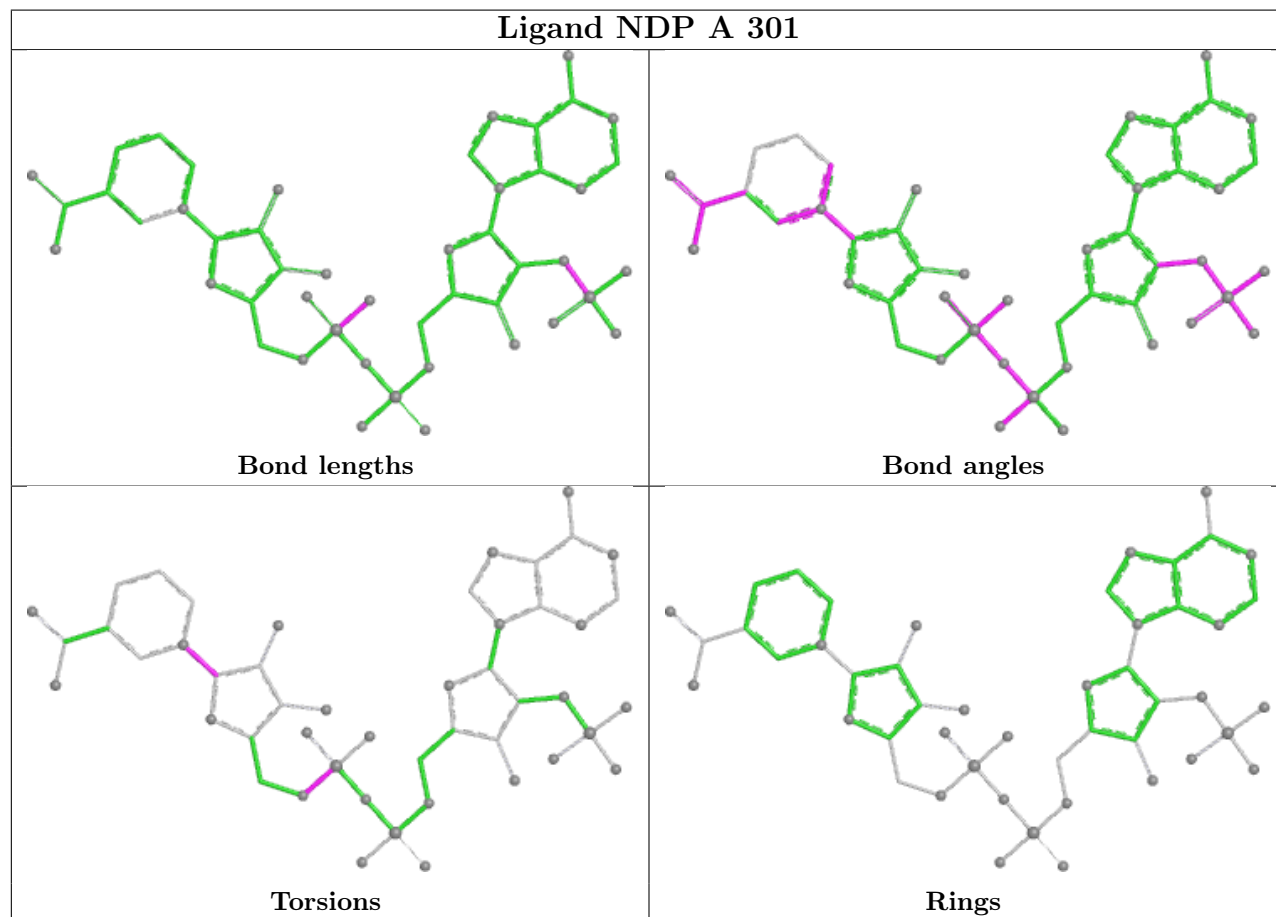
Mol	Chain	Res	Type	Atoms
2	A	301	NDP	C5D-O5D-PN-O1N
3	A	302	G6Y	C5-CAI-CAJ-CAK
3	B	302	G6Y	C5-CAI-CAJ-CAK
2	A	301	NDP	O4D-C1D-N1N-C2N
2	B	301	NDP	O4D-C1D-N1N-C2N
3	A	302	G6Y	C4-C5-CAI-CAJ
3	B	302	G6Y	C4-C5-CAI-CAJ
4	B	303	EDO	O1-C1-C2-O2
3	A	302	G6Y	CAJ-CAK-CAL-CAM
3	A	302	G6Y	CAJ-CAK-CAL-CAQ
3	B	302	G6Y	CAJ-CAK-CAL-CAM
3	B	302	G6Y	CAJ-CAK-CAL-CAQ

There are no ring outliers.

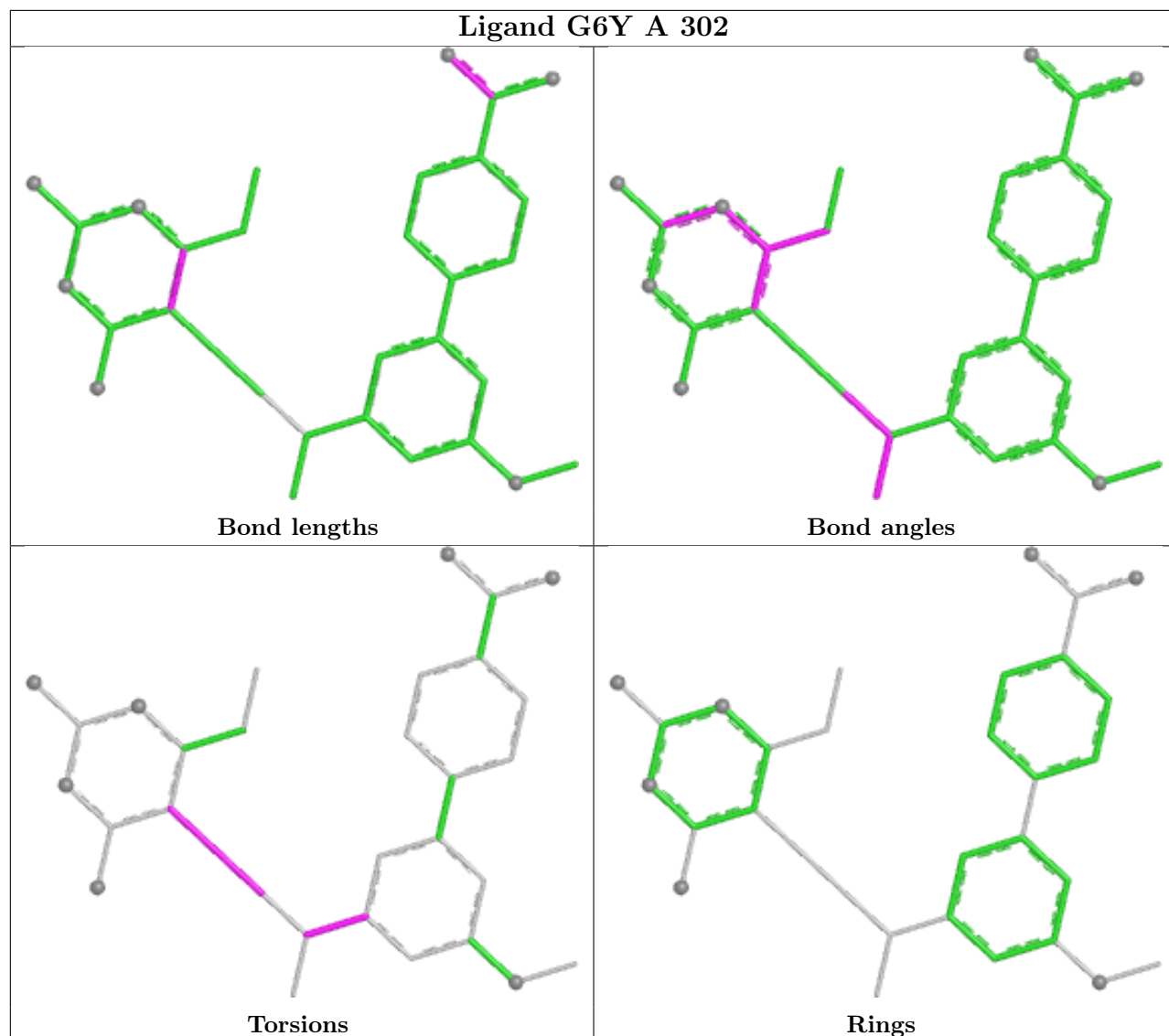
3 monomers are involved in 5 short contacts:

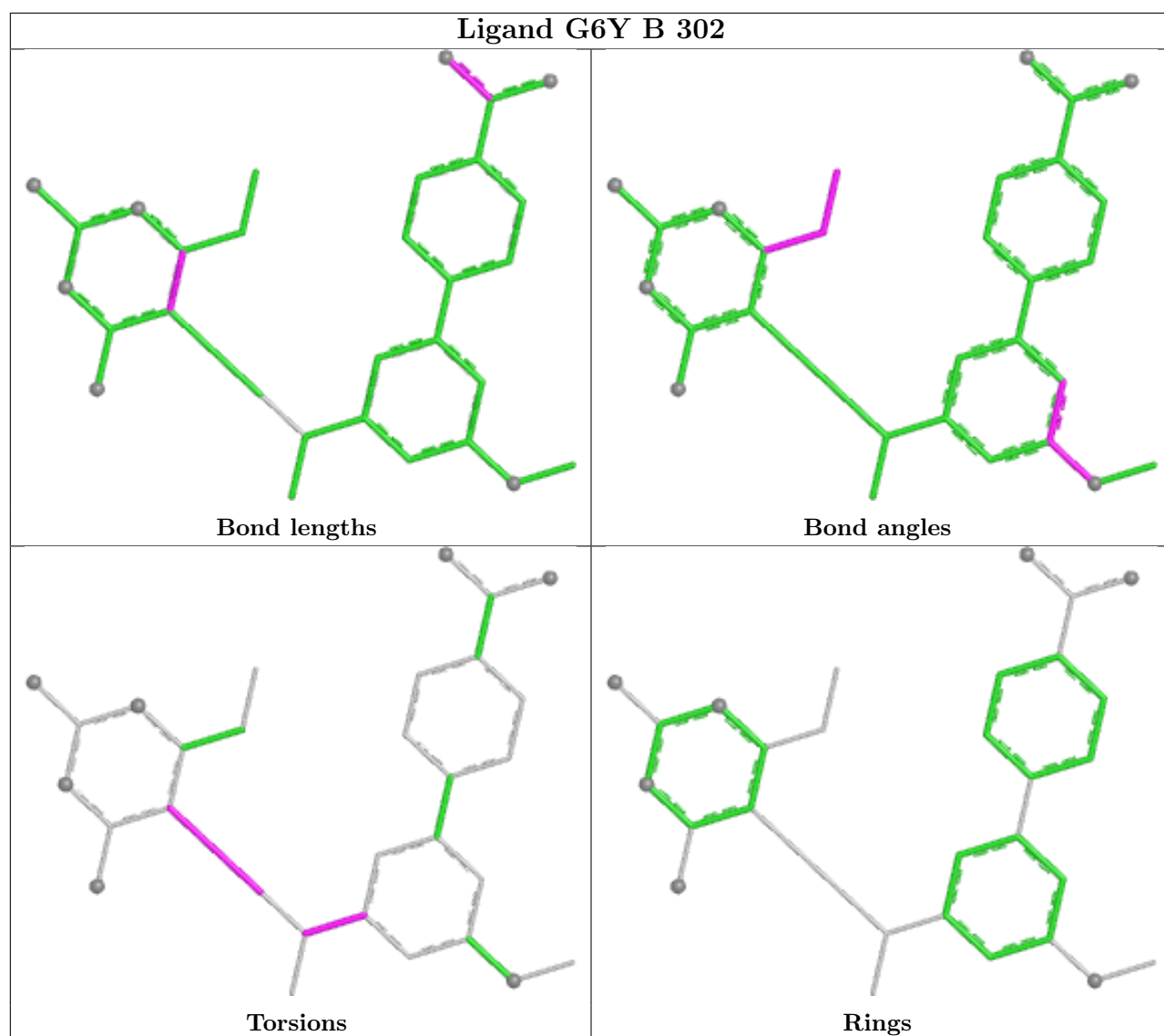
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	301	NDP	2	0
3	B	302	G6Y	1	0
2	B	301	NDP	2	0

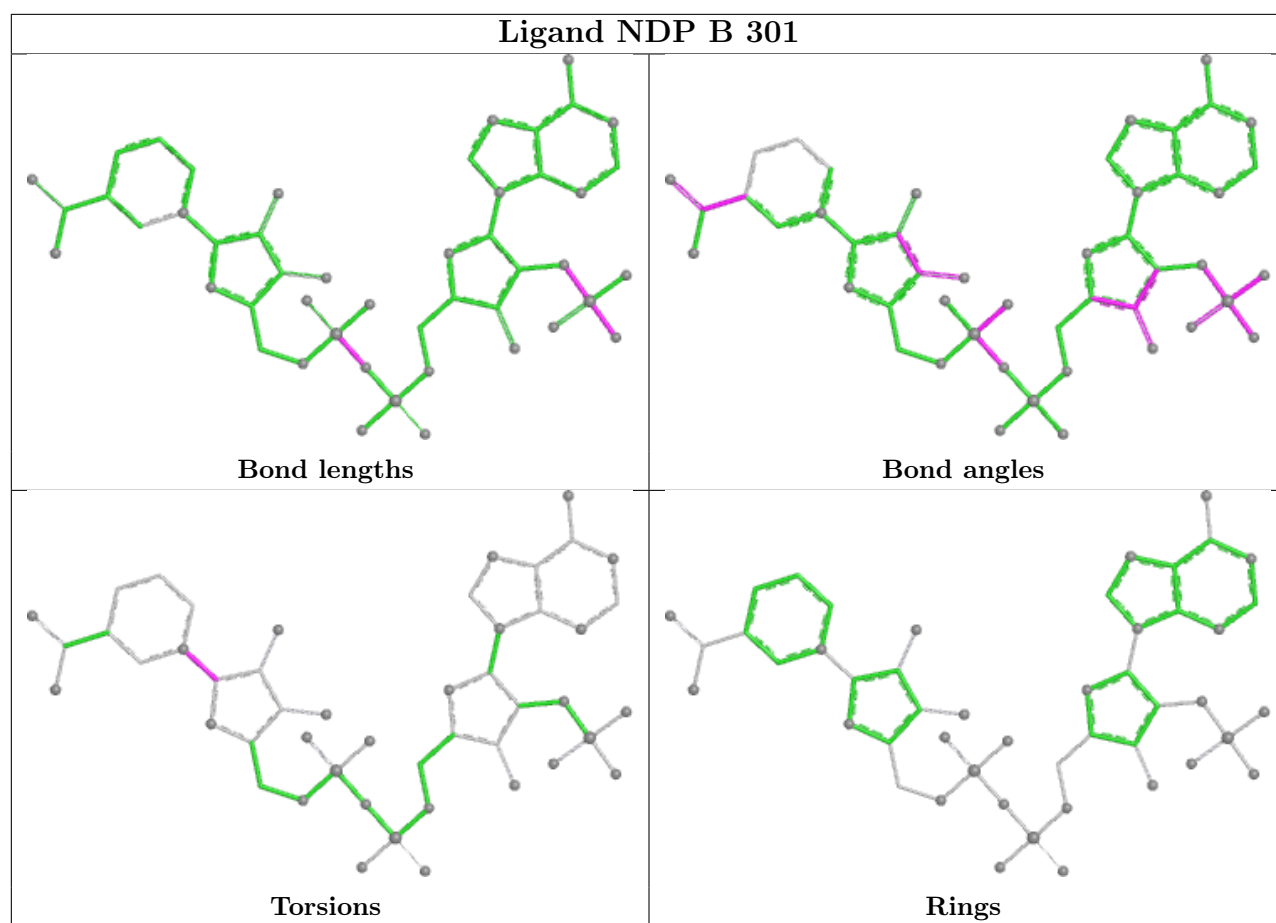
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.



Ligand G6Y A 302







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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	200/204 (98%)	-0.20	0	100 100	23, 34, 60, 73	0
1	B	200/204 (98%)	0.31	3 (1%)	72 80	29, 45, 76, 93	1 (0%)
All	All	400/408 (98%)	0.06	3 (0%)	82 87	23, 39, 71, 93	1 (0%)

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	202	ALA	2.2
1	B	83	LEU	2.1
1	B	109	PHE	2.0

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

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

6.3 Carbohydrates [i](#)

There are no 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
5	SO4	A	305	5/5	0.71	0.10	87,108,117,124	0
5	SO4	A	306	5/5	0.79	0.09	88,98,135,135	0

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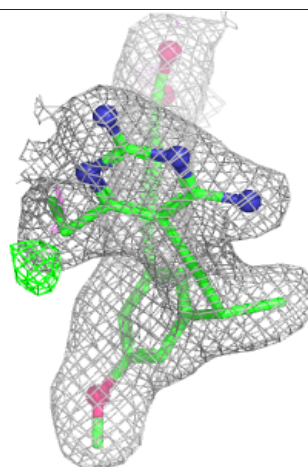
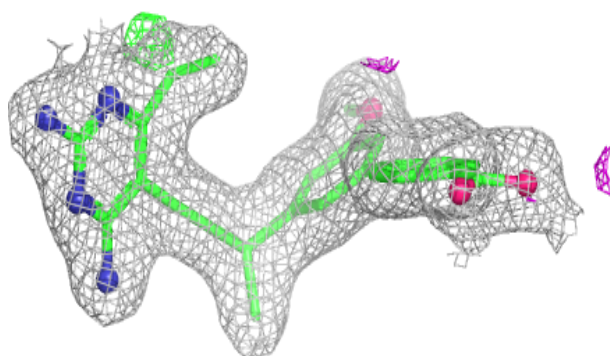
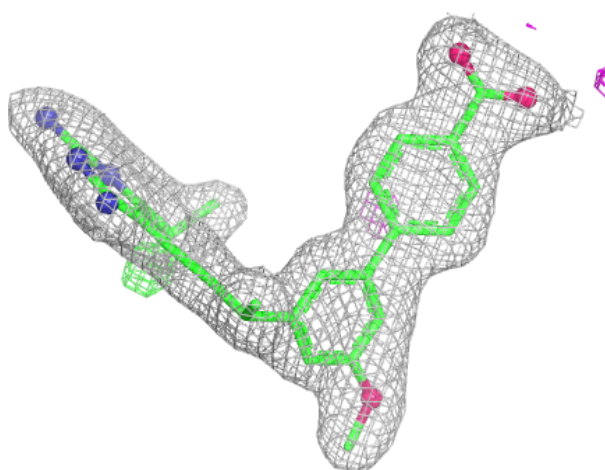
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	EDO	A	304	4/4	0.85	0.13	53,55,58,61	0
5	SO4	B	304	5/5	0.85	0.08	86,86,93,117	0
4	EDO	A	303	4/4	0.86	0.14	47,60,63,66	0
4	EDO	B	303	4/4	0.88	0.12	61,63,67,69	0
3	G6Y	B	302	31/31	0.95	0.07	25,34,41,44	0
3	G6Y	A	302	31/31	0.96	0.06	22,31,39,39	0
2	NDP	A	301	48/48	0.97	0.06	20,33,48,53	0
2	NDP	B	301	48/48	0.97	0.06	25,35,54,57	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.

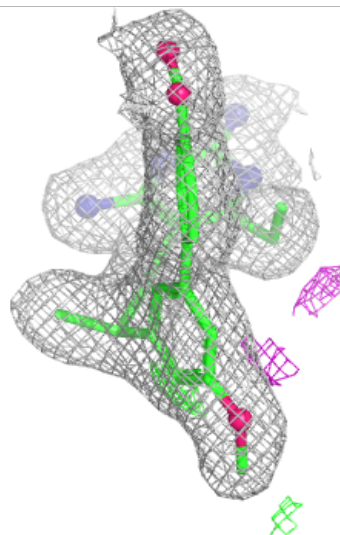
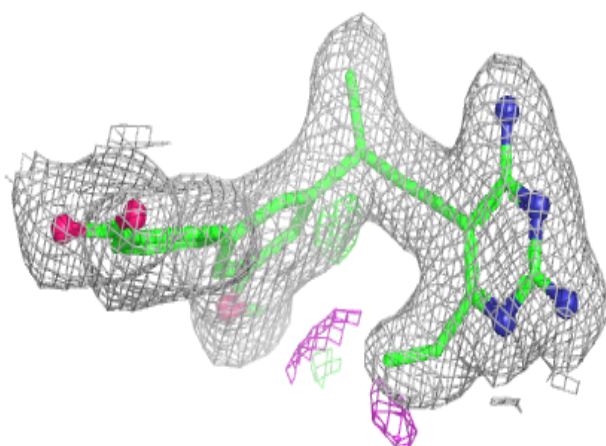
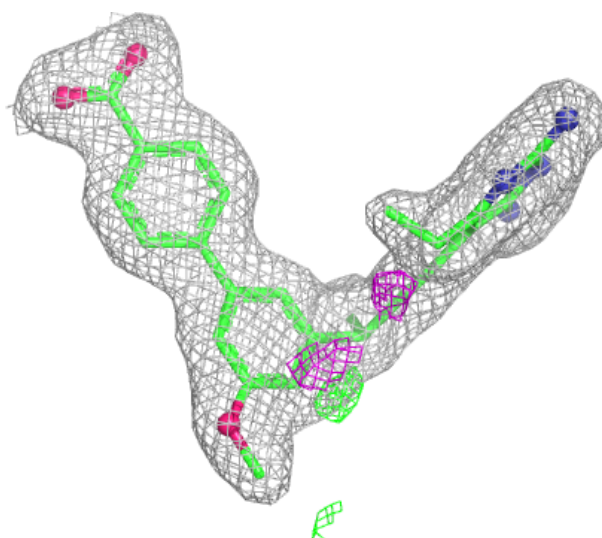
Electron density around G6Y B 302:

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



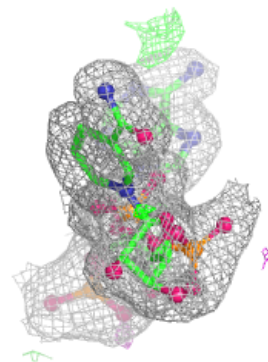
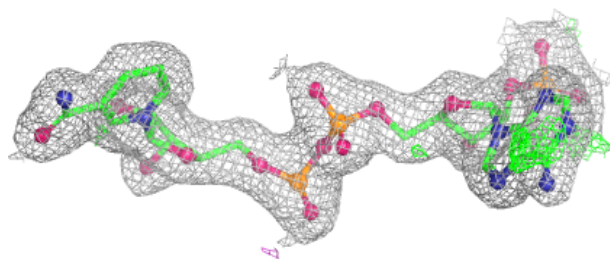
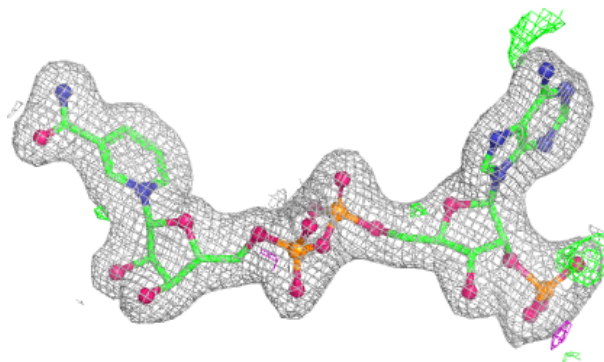
Electron density around G6Y A 302:

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

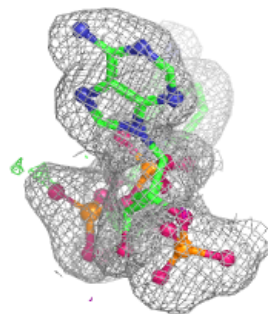
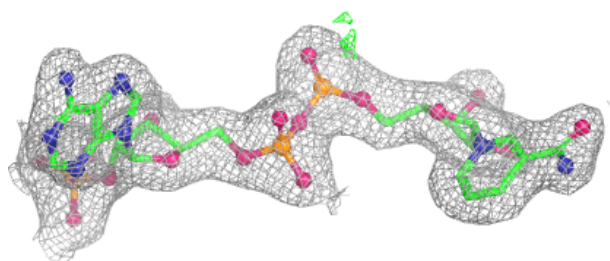
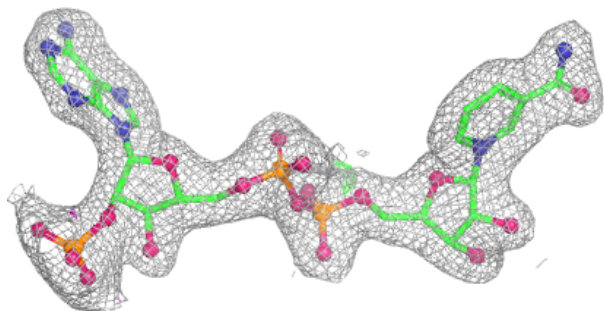


Electron density around NDP A 301:

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

**Electron density around NDP B 301:**

$2mF_o-DF_c$ (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.