



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 08:31 AM UTC

PDB ID : 5MLD / pdb_00005mld
Title : Crystal Structure of RosB with bound intermediate AFP (8-demethyl-8-aminoriboflavin-5'-phosphate)
Authors : Konjik, V.; Bruenle, S.; Demmer, U.; Vanselow, A.; Sandhoff, R.; Mack, M.; Ermler, U.
Deposited on : 2016-12-06
Resolution : 1.70 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

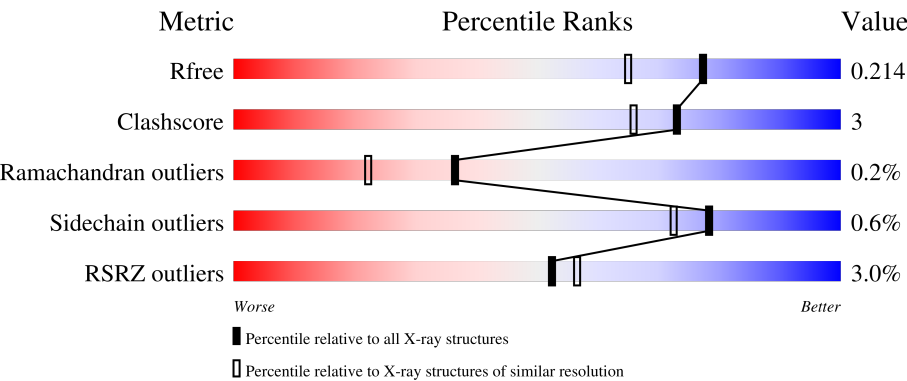
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.70 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	5551 (1.70-1.70)
Clashscore	190562	5924 (1.70-1.70)
Ramachandran outliers	187476	5846 (1.70-1.70)
Sidechain outliers	187428	5846 (1.70-1.70)
RSRZ outliers	180081	5554 (1.70-1.70)

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	257	5% 86% 11% .
1	B	257	5% 89% 7% ..
1	C	257	5% 93% 5% .
1	D	257	2% 92% 6% .

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Mol	Chain	Length	Quality of chain
1	E	257	2%88%10%•
1	F	257	2%89%8%••
1	G	257	%91%7%•
1	H	257	2%91%7%•

2 Entry composition

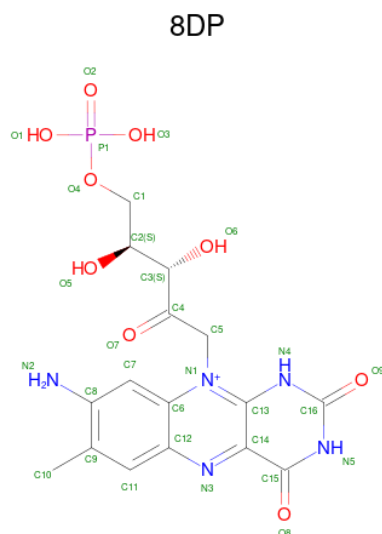
There are 3 unique types of molecules in this entry. The entry contains 18138 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 BRAMP domain protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	252	Total	C	N	O	S	0	8	0
			2054	1292	358	395	9			
1	B	252	Total	C	N	O	S	0	6	0
			2041	1285	355	392	9			
1	C	252	Total	C	N	O	S	0	1	0
			2000	1259	350	383	8			
1	D	252	Total	C	N	O	S	0	0	0
			1995	1256	350	381	8			
1	E	252	Total	C	N	O	S	0	2	0
			2008	1263	352	385	8			
1	F	252	Total	C	N	O	S	0	1	0
			2003	1260	351	384	8			
1	G	252	Total	C	N	O	S	0	1	0
			2000	1259	350	383	8			
1	H	252	Total	C	N	O	S	0	2	0
			2005	1262	351	384	8			

- Molecule 2 is 8-demethyl-8-aminoriboflavin-5'-phosphate (CCD ID: 8DP) (formula: C₁₆H₁₉N₅O₉P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total 31	C 16	N 5	O 9	P 1	0	0
2	B	1	Total 31	C 16	N 5	O 9	P 1	0	0
2	C	1	Total 31	C 16	N 5	O 9	P 1	0	0
2	D	1	Total 31	C 16	N 5	O 9	P 1	0	0
2	E	1	Total 31	C 16	N 5	O 9	P 1	0	0
2	F	1	Total 31	C 16	N 5	O 9	P 1	0	0
2	G	1	Total 31	C 16	N 5	O 9	P 1	0	0
2	H	1	Total 31	C 16	N 5	O 9	P 1	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	246	Total O 246 246	0	0
3	B	205	Total O 205 205	0	0
3	C	191	Total O 191 191	0	0
3	D	246	Total O 246 246	0	0

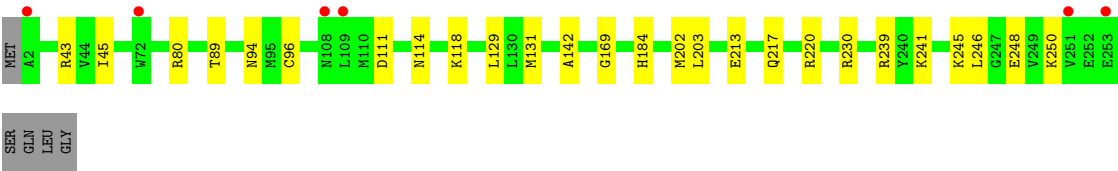
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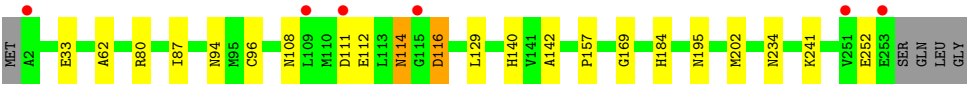
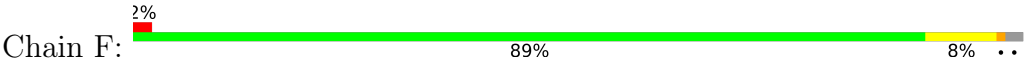
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	E	217	Total 217	O 217	0	0
3	F	221	Total 221	O 221	0	0
3	G	247	Total 247	O 247	0	0
3	H	211	Total 211	O 211	0	0

- Molecule 1: BRAMP domain protein

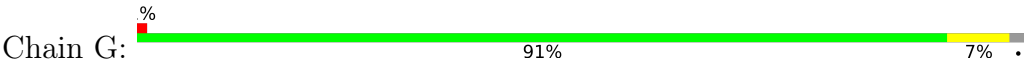




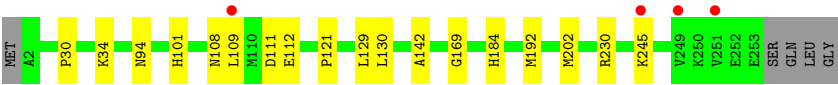
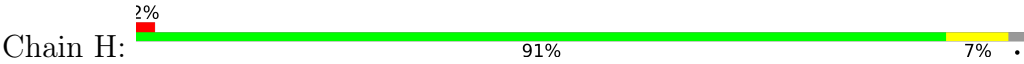
• Molecule 1: BRAMP domain protein



• Molecule 1: BRAMP domain protein



• Molecule 1: BRAMP domain protein



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	131.49Å 71.13Å 214.80Å 90.00° 92.08° 90.00°	Depositor
Resolution (Å)	47.54 – 1.70 47.54 – 1.70	Depositor EDS
% Data completeness (in resolution range)	92.9 (47.54-1.70) 92.9 (47.54-1.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.11	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.63 (at 1.70Å)	Xtriage
Refinement program	PHENIX (1.10.1_2155: ???)	Depositor
R, R_{free}	0.174 , 0.213 0.175 , 0.214	Depositor DCC
R_{free} test set	10149 reflections (4.66%)	wwPDB-VP
Wilson B-factor (Å ²)	23.6	Xtriage
Anisotropy	0.092	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 41.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.013 for -h,-k,l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	18138	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.60% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: 8DP

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	0.31	0/2101	0.56	0/2853
1	B	0.30	0/2085	0.59	2/2831 (0.1%)
1	C	0.30	0/2044	0.54	0/2776
1	D	0.30	0/2036	0.54	0/2765
1	E	0.34	1/2052 (0.0%)	0.53	0/2787
1	F	0.31	0/2044	0.53	0/2776
1	G	0.31	0/2044	0.54	0/2776
1	H	0.28	0/2052	0.53	0/2787
All	All	0.31	1/16458 (0.0%)	0.54	2/22351 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	241	LYS	CD-CE	-5.43	1.36	1.52

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	115	GLY	CA-C-N	7.17	134.60	121.70
1	B	115	GLY	C-N-CA	7.17	134.60	121.70

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	2054	0	2019	23	0
1	B	2041	0	2007	21	0
1	C	2000	0	1967	13	0
1	D	1995	0	1963	14	0
1	E	2008	0	1972	20	0
1	F	2003	0	1966	18	0
1	G	2000	0	1967	18	0
1	H	2005	0	1973	13	0
2	A	31	0	0	0	0
2	B	31	0	0	0	0
2	C	31	0	0	0	0
2	D	31	0	0	0	0
2	E	31	0	0	0	0
2	F	31	0	0	0	0
2	G	31	0	0	0	0
2	H	31	0	0	0	0
3	A	246	0	0	1	0
3	B	205	0	0	2	0
3	C	191	0	0	1	0
3	D	246	0	0	2	0
3	E	217	0	0	3	0
3	F	221	0	0	3	0
3	G	247	0	0	0	0
3	H	211	0	0	1	0
All	All	18138	0	15834	103	0

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 (103) 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:115:GLY:HA3	1:B:116:ASP:HB2	1.33	1.06
1:E:250:LYS:NZ	3:E:401:HOH:O	2.19	0.74
1:E:213:GLU:OE2	1:E:217:GLN:NE2	2.22	0.72
1:F:80:ARG:NH2	1:F:112:GLU:OE2	2.24	0.71
1:A:202:MET:HE2	1:B:202:MET:HE2	1.74	0.69
1:C:202:MET:HE2	1:D:202:MET:HE2	1.73	0.69
1:B:111[A]:ASP:OD2	1:D:241:LYS:NZ	2.26	0.69
1:D:220:ARG:NH2	3:D:401:HOH:O	2.23	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:111[A]:ASP:HB2	1:C:239:ARG:HH11	1.58	0.67
1:A:110[B]:MET:O	3:A:401:HOH:O	2.14	0.64
1:H:34:LYS:NZ	3:H:402:HOH:O	2.31	0.64
1:D:61:ASN:HB3	1:E:248:GLU:HG3	1.80	0.64
1:E:245:LYS:NZ	3:E:403:HOH:O	2.31	0.64
1:B:115:GLY:HA3	1:B:116:ASP:CB	2.20	0.63
1:F:202:MET:HE2	1:H:202:MET:HE2	1.82	0.62
1:C:80:ARG:HH22	1:C:112:GLU:CD	2.08	0.60
1:E:202:MET:HE2	1:G:202:MET:HE2	1.83	0.60
1:B:109[A]:LEU:HD12	1:D:239:ARG:NH1	2.15	0.60
1:E:184:HIS:CG	1:F:169:GLY:HA2	2.38	0.59
1:A:112[A]:GLU:HB2	1:A:121:PRO:HG3	1.85	0.58
1:G:80:ARG:NH2	1:G:112:GLU:OE1	2.28	0.56
1:A:230:ARG:HH12	1:C:253:GLU:HG2	1.71	0.55
1:B:115:GLY:CA	1:B:116:ASP:HB2	2.21	0.55
1:E:169:GLY:HA2	1:F:184:HIS:CG	2.42	0.54
1:E:239:ARG:NH1	1:F:111:ASP:OD1	2.40	0.53
1:H:30:PRO:O	1:H:34:LYS:HE2	2.09	0.53
1:H:130:LEU:HD21	1:H:192:MET:HE2	1.91	0.53
1:F:108:ASN:ND2	3:F:403:HOH:O	2.35	0.53
1:A:80:ARG:NH2	1:A:112[A]:GLU:OE1	2.43	0.52
1:H:112:GLU:O	1:H:121:PRO:HD3	2.10	0.52
1:D:18:ARG:HG2	3:D:557:HOH:O	2.09	0.52
1:B:221:GLU:OE1	3:B:401:HOH:O	2.19	0.52
1:D:248:GLU:OE1	1:D:250:LYS:NZ	2.33	0.52
1:C:239:ARG:NH2	3:C:404:HOH:O	2.44	0.51
1:B:184:HIS:CG	1:D:169:GLY:HA2	2.46	0.51
1:C:80:ARG:NH2	1:C:112:GLU:OE2	2.44	0.51
1:F:33:GLU:OE2	3:F:401:HOH:O	2.19	0.51
1:G:87:ILE:O	1:G:129:LEU:HD12	2.11	0.50
1:C:248:GLU:OE2	1:C:250:LYS:NZ	2.37	0.50
1:G:239:ARG:HD2	1:H:109:LEU:HD13	1.93	0.50
1:A:112[B]:GLU:O	1:A:121:PRO:HD3	2.11	0.50
1:G:253:GLU:C	1:H:230:ARG:HH22	2.20	0.50
1:F:80:ARG:HD2	3:F:410:HOH:O	2.11	0.49
1:A:239:ARG:NH1	1:C:111:ASP:HB2	2.27	0.49
1:E:111[B]:ASP:OD1	1:E:111[B]:ASP:N	2.41	0.49
1:G:253:GLU:O	1:H:230:ARG:NH2	2.41	0.49
1:F:94:ASN:OD1	1:F:140:HIS:ND1	2.32	0.48
1:B:105:GLU:O	1:B:109[A]:LEU:HD22	2.14	0.47
1:E:96:CYS:O	1:H:101:HIS:CE1	2.68	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:112[B]:GLU:O	1:A:120:PHE:HA	2.15	0.47
1:A:107:LEU:C	1:A:109[B]:LEU:H	2.23	0.46
1:B:252:GLU:OE2	1:D:250:LYS:C	2.59	0.45
1:A:169:GLY:HA2	1:C:184:HIS:CG	2.52	0.45
1:B:96:CYS:O	1:C:101:HIS:CE1	2.69	0.45
1:A:226:VAL:HG11	1:A:251:VAL:HG21	1.98	0.45
1:A:206:THR:OG1	1:B:202:MET:HE3	2.16	0.45
1:B:111[A]:ASP:HB2	1:D:239:ARG:HE	1.81	0.45
1:E:43:ARG:HG2	1:E:45:ILE:HG22	1.99	0.45
1:E:217:GLN:HG2	1:E:220:ARG:NH2	2.32	0.45
1:F:129:LEU:CD2	1:F:142:ALA:HA	2.47	0.45
1:B:80:ARG:NH2	1:B:112[A]:GLU:OE2	2.21	0.44
1:G:169:GLY:HA2	1:H:184:HIS:CG	2.53	0.44
1:A:101:HIS:CE1	1:D:96:CYS:O	2.70	0.44
1:B:178:LYS:HA	1:B:178:LYS:HD2	1.83	0.44
1:A:18:ARG:HD3	1:G:18:ARG:HG2	1.99	0.43
1:E:220:ARG:NH1	1:F:234:ASN:OD1	2.47	0.43
1:G:217:GLN:HG2	1:G:220:ARG:HH22	1.83	0.43
1:B:108:ASN:HB3	3:B:569:HOH:O	2.18	0.43
3:E:517:HOH:O	1:F:241:LYS:HE2	2.17	0.43
1:A:96:CYS:O	1:D:101:HIS:CE1	2.72	0.43
1:G:129:LEU:HG	1:G:131:MET:HG3	2.00	0.43
1:G:184:HIS:CG	1:H:169:GLY:HA2	2.53	0.43
1:A:129:LEU:CD2	1:A:142:ALA:HA	2.49	0.43
1:B:169:GLY:HA2	1:D:184:HIS:CG	2.54	0.43
1:G:57:TYR:HE1	1:H:245:LYS:HE3	1.83	0.43
1:E:129:LEU:CD2	1:E:142:ALA:HA	2.48	0.43
1:E:230:ARG:HD2	1:E:246:LEU:O	2.18	0.43
1:F:96:CYS:O	1:G:101:HIS:CE1	2.71	0.43
1:G:80:ARG:HH22	1:G:112:GLU:CD	2.23	0.43
1:A:217:GLN:HG2	1:A:220:ARG:HH22	1.84	0.42
1:G:217:GLN:HG2	1:G:220:ARG:NH2	2.33	0.42
1:C:87:ILE:O	1:C:129:LEU:HD12	2.20	0.42
1:H:129:LEU:CD2	1:H:142:ALA:HA	2.50	0.42
1:F:87:ILE:O	1:F:129:LEU:HD12	2.19	0.42
1:D:129:LEU:CD2	1:D:142:ALA:HA	2.50	0.41
1:A:153:GLY:C	1:B:158:PRO:HB3	2.45	0.41
1:E:203:LEU:HD23	1:G:202:MET:HE1	2.03	0.41
1:A:239:ARG:HH11	1:C:111:ASP:HB2	1.86	0.41
1:A:252:GLU:HB3	1:A:253:GLU:H	1.55	0.41
1:A:77:GLU:HG3	1:A:81:GLU:OE2	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:80:ARG:HH22	1:E:114:ASN:HD22	1.68	0.41
1:E:118:LYS:HA	1:E:118:LYS:HD3	1.92	0.41
1:F:62:ALA:HB2	1:G:16:PRO:HD2	2.03	0.41
1:F:114:ASN:O	1:F:116:ASP:N	2.54	0.41
1:G:129:LEU:CD2	1:G:142:ALA:HA	2.51	0.41
1:B:131:MET:O	1:B:163:TYR:HA	2.20	0.41
1:A:54:TRP:CD1	1:A:63:GLY:HA3	2.56	0.40
1:B:27:LYS:NZ	1:B:180[B]:ASP:OD1	2.39	0.40
1:B:129:LEU:CD2	1:B:142:ALA:HA	2.51	0.40
1:C:129:LEU:CD2	1:C:142:ALA:HA	2.51	0.40
1:E:89:THR:O	1:E:131:MET:HA	2.21	0.40
1:E:239:ARG:HH11	1:F:111:ASP:CG	2.29	0.40
1:F:157:PRO:HB2	1:F:195:ASN:HB3	2.04	0.40

There are no symmetry-related clashes.

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	258/257 (100%)	245 (95%)	12 (5%)	1 (0%)	30	16
1	B	256/257 (100%)	246 (96%)	8 (3%)	2 (1%)	16	5
1	C	251/257 (98%)	247 (98%)	4 (2%)	0	100	100
1	D	250/257 (97%)	247 (99%)	3 (1%)	0	100	100
1	E	252/257 (98%)	247 (98%)	5 (2%)	0	100	100
1	F	251/257 (98%)	244 (97%)	5 (2%)	2 (1%)	16	5
1	G	251/257 (98%)	246 (98%)	5 (2%)	0	100	100
1	H	252/257 (98%)	246 (98%)	6 (2%)	0	100	100
All	All	2021/2056 (98%)	1968 (97%)	48 (2%)	5 (0%)	36	28

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	252	GLU
1	B	116	ASP
1	B	115	GLY
1	F	116	ASP
1	F	114	ASN

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	219/215 (102%)	218 (100%)	1 (0%)	81	76
1	B	217/215 (101%)	215 (99%)	2 (1%)	70	62
1	C	212/215 (99%)	211 (100%)	1 (0%)	81	76
1	D	211/215 (98%)	210 (100%)	1 (0%)	81	76
1	E	213/215 (99%)	212 (100%)	1 (0%)	81	76
1	F	212/215 (99%)	211 (100%)	1 (0%)	81	76
1	G	212/215 (99%)	211 (100%)	1 (0%)	81	76
1	H	213/215 (99%)	210 (99%)	3 (1%)	59	46
All	All	1709/1720 (99%)	1698 (99%)	11 (1%)	78	72

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	94	ASN
1	B	109[A]	LEU
1	B	109[B]	LEU
1	C	108	ASN
1	D	94	ASN
1	E	94	ASN
1	F	252	GLU
1	G	94	ASN
1	H	94	ASN

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Mol	Chain	Res	Type
1	H	108	ASN
1	H	111	ASP

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	94	ASN
1	B	94	ASN
1	B	108	ASN
1	B	148	ASN
1	C	147	ASN
1	C	159	ASN
1	C	204	GLN
1	D	147	ASN
1	E	147	ASN
1	G	139	HIS
1	G	244	GLN
1	H	139	HIS
1	H	204	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

8 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	8DP	B	301	-	32,33,33	2.63	9 (28%)	36,50,50	1.73	8 (22%)
2	8DP	C	301	-	32,33,33	2.48	8 (25%)	36,50,50	2.10	8 (22%)
2	8DP	D	301	-	32,33,33	2.58	11 (34%)	36,50,50	1.85	8 (22%)
2	8DP	H	301	-	32,33,33	2.47	9 (28%)	36,50,50	1.97	9 (25%)
2	8DP	F	301	-	32,33,33	2.57	10 (31%)	36,50,50	1.88	8 (22%)
2	8DP	A	301	-	32,33,33	2.55	11 (34%)	36,50,50	1.72	8 (22%)
2	8DP	G	301	-	32,33,33	2.52	9 (28%)	36,50,50	2.05	9 (25%)
2	8DP	E	301	-	32,33,33	2.58	10 (31%)	36,50,50	1.69	8 (22%)

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	8DP	B	301	-	-	8/18/18/18	0/3/3/3
2	8DP	C	301	-	-	4/18/18/18	0/3/3/3
2	8DP	D	301	-	-	4/18/18/18	0/3/3/3
2	8DP	H	301	-	-	4/18/18/18	0/3/3/3
2	8DP	F	301	-	-	5/18/18/18	0/3/3/3
2	8DP	A	301	-	-	4/18/18/18	0/3/3/3
2	8DP	G	301	-	-	4/18/18/18	0/3/3/3
2	8DP	E	301	-	-	4/18/18/18	0/3/3/3

All (77) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	301	8DP	O9-C16	9.48	1.43	1.23
2	A	301	8DP	O9-C16	8.86	1.42	1.23
2	D	301	8DP	O9-C16	8.83	1.42	1.23
2	G	301	8DP	O9-C16	8.79	1.42	1.23
2	E	301	8DP	O9-C16	8.74	1.41	1.23
2	H	301	8DP	O9-C16	8.49	1.41	1.23
2	F	301	8DP	O9-C16	8.37	1.41	1.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	301	8DP	O8-C15	8.11	1.42	1.23
2	E	301	8DP	O8-C15	7.99	1.42	1.23
2	D	301	8DP	O8-C15	7.95	1.41	1.23
2	B	301	8DP	O8-C15	7.75	1.41	1.23
2	C	301	8DP	O9-C16	7.53	1.39	1.23
2	C	301	8DP	O8-C15	7.34	1.40	1.23
2	A	301	8DP	O8-C15	7.11	1.39	1.23
2	H	301	8DP	O8-C15	6.84	1.39	1.23
2	G	301	8DP	O8-C15	6.64	1.38	1.23
2	G	301	8DP	C6-N1	-4.23	1.35	1.40
2	H	301	8DP	C6-N1	-4.05	1.35	1.40
2	C	301	8DP	C6-N1	-3.95	1.35	1.40
2	F	301	8DP	C6-N1	-3.74	1.36	1.40
2	C	301	8DP	C14-C15	-3.69	1.40	1.47
2	C	301	8DP	C14-N3	3.66	1.37	1.32
2	H	301	8DP	C14-C15	-3.63	1.40	1.47
2	G	301	8DP	C14-N3	3.63	1.37	1.32
2	A	301	8DP	C6-N1	-3.61	1.36	1.40
2	B	301	8DP	C6-N1	-3.58	1.36	1.40
2	D	301	8DP	C6-N1	-3.53	1.36	1.40
2	G	301	8DP	C14-C15	-3.52	1.40	1.47
2	E	301	8DP	C6-N1	-3.52	1.36	1.40
2	F	301	8DP	C14-C15	-3.50	1.40	1.47
2	D	301	8DP	C14-C15	-3.49	1.40	1.47
2	A	301	8DP	C14-C15	-3.49	1.40	1.47
2	B	301	8DP	C14-C15	-3.46	1.40	1.47
2	E	301	8DP	C14-C15	-3.17	1.41	1.47
2	C	301	8DP	C15-N5	-2.99	1.33	1.37
2	A	301	8DP	C14-N3	2.99	1.36	1.32
2	F	301	8DP	C14-N3	2.89	1.36	1.32
2	H	301	8DP	C14-N3	2.88	1.36	1.32
2	E	301	8DP	C14-N3	2.79	1.36	1.32
2	G	301	8DP	C15-N5	-2.66	1.33	1.37
2	B	301	8DP	C14-N3	2.60	1.36	1.32
2	D	301	8DP	C16-N4	-2.47	1.32	1.36
2	C	301	8DP	P1-O2	2.46	1.58	1.50
2	A	301	8DP	C16-N4	-2.41	1.32	1.36
2	B	301	8DP	C1-C2	2.37	1.55	1.51
2	E	301	8DP	C1-C2	2.35	1.55	1.51
2	H	301	8DP	C16-N4	-2.31	1.32	1.36
2	A	301	8DP	C15-N5	-2.29	1.34	1.37
2	C	301	8DP	C8-N2	2.29	1.45	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	H	301	8DP	C15-N5	-2.27	1.34	1.37
2	F	301	8DP	C8-N2	2.26	1.45	1.38
2	D	301	8DP	C1-C2	2.25	1.54	1.51
2	G	301	8DP	C8-N2	2.24	1.45	1.38
2	F	301	8DP	C16-N4	-2.23	1.32	1.36
2	A	301	8DP	P1-O2	2.21	1.57	1.50
2	B	301	8DP	C16-N4	-2.20	1.32	1.36
2	F	301	8DP	C1-C2	2.19	1.54	1.51
2	B	301	8DP	C8-N2	2.18	1.45	1.38
2	D	301	8DP	C14-N3	2.17	1.35	1.32
2	F	301	8DP	C15-N5	-2.16	1.34	1.37
2	G	301	8DP	C16-N4	-2.14	1.32	1.36
2	H	301	8DP	C8-N2	2.14	1.45	1.38
2	D	301	8DP	C8-N2	2.13	1.45	1.38
2	D	301	8DP	C12-N3	-2.13	1.34	1.37
2	A	301	8DP	O5-C2	-2.13	1.38	1.43
2	F	301	8DP	P1-O2	2.11	1.57	1.50
2	G	301	8DP	O7-C4	2.10	1.25	1.21
2	D	301	8DP	C15-N5	-2.08	1.34	1.37
2	B	301	8DP	C15-N5	-2.07	1.34	1.37
2	E	301	8DP	C15-N5	-2.06	1.34	1.37
2	E	301	8DP	C8-N2	2.05	1.44	1.38
2	E	301	8DP	C5-N1	2.05	1.54	1.49
2	E	301	8DP	C16-N4	-2.04	1.32	1.36
2	D	301	8DP	O7-C4	2.04	1.25	1.21
2	A	301	8DP	O7-C4	2.03	1.25	1.21
2	A	301	8DP	C8-N2	2.02	1.44	1.38
2	H	301	8DP	C1-C2	2.00	1.54	1.51

All (66) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	301	8DP	C13-N4-C16	-6.48	117.98	128.11
2	F	301	8DP	C13-N4-C16	-5.90	118.89	128.11
2	G	301	8DP	O5-C2-C3	5.42	117.54	109.67
2	G	301	8DP	C13-N4-C16	-5.35	119.75	128.11
2	G	301	8DP	C12-C6-N1	5.30	121.50	117.75
2	D	301	8DP	C13-N4-C16	-5.29	119.83	128.11
2	B	301	8DP	C13-N4-C16	-5.01	120.28	128.11
2	H	301	8DP	C13-N4-C16	-4.96	120.35	128.11
2	A	301	8DP	C13-N4-C16	-4.90	120.45	128.11
2	C	301	8DP	N4-C16-N5	4.87	124.57	116.17

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	301	8DP	C13-N4-C16	-4.85	120.53	128.11
2	B	301	8DP	C12-C6-N1	4.78	121.13	117.75
2	E	301	8DP	C12-C6-N1	4.61	121.01	117.75
2	F	301	8DP	C12-C6-N1	4.35	120.83	117.75
2	H	301	8DP	C12-C6-N1	4.33	120.81	117.75
2	C	301	8DP	C12-C6-N1	4.32	120.80	117.75
2	D	301	8DP	C12-C6-N1	4.16	120.69	117.75
2	D	301	8DP	O5-C2-C3	4.06	115.58	109.67
2	H	301	8DP	N4-C16-N5	4.04	123.15	116.17
2	C	301	8DP	O9-C16-N5	-3.82	115.08	121.86
2	C	301	8DP	O5-C2-C3	3.74	115.11	109.67
2	F	301	8DP	N4-C16-N5	3.71	122.58	116.17
2	A	301	8DP	O5-C2-C3	3.61	114.92	109.67
2	D	301	8DP	N4-C16-N5	3.58	122.36	116.17
2	H	301	8DP	C14-C15-N5	3.48	120.71	114.07
2	H	301	8DP	C15-N5-C16	-3.46	118.72	126.47
2	G	301	8DP	N4-C16-N5	3.45	122.13	116.17
2	H	301	8DP	O5-C2-C3	3.45	114.69	109.67
2	F	301	8DP	O9-C16-N5	-3.40	115.82	121.86
2	A	301	8DP	C12-C6-N1	3.40	120.15	117.75
2	E	301	8DP	N4-C16-N5	3.32	121.91	116.17
2	B	301	8DP	O5-C2-C3	3.31	114.48	109.67
2	C	301	8DP	C15-N5-C16	-3.29	119.10	126.47
2	G	301	8DP	C14-C15-N5	3.26	120.30	114.07
2	A	301	8DP	N4-C16-N5	3.26	121.80	116.17
2	H	301	8DP	O9-C16-N5	-3.10	116.36	121.86
2	F	301	8DP	O5-C2-C3	3.00	114.03	109.67
2	C	301	8DP	C14-C15-N5	2.98	119.75	114.07
2	H	301	8DP	C4-C5-N1	-2.97	105.77	111.59
2	C	301	8DP	C4-C5-N1	-2.96	105.79	111.59
2	A	301	8DP	O9-C16-N5	-2.91	116.68	121.86
2	B	301	8DP	N4-C16-N5	2.90	121.17	116.17
2	A	301	8DP	C14-C15-N5	2.87	119.55	114.07
2	G	301	8DP	C15-N5-C16	-2.77	120.27	126.47
2	E	301	8DP	C4-C5-N1	-2.71	106.27	111.59
2	G	301	8DP	O9-C16-N5	-2.71	117.05	121.86
2	B	301	8DP	C4-C5-N1	-2.67	106.36	111.59
2	D	301	8DP	C4-C5-N1	-2.64	106.41	111.59
2	D	301	8DP	O9-C16-N5	-2.59	117.26	121.86
2	A	301	8DP	C4-C5-N1	-2.59	106.52	111.59
2	E	301	8DP	O9-C16-N5	-2.55	117.33	121.86
2	A	301	8DP	C15-N5-C16	-2.54	120.78	126.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	301	8DP	C15-N5-C16	-2.50	120.86	126.47
2	F	301	8DP	C4-C5-N1	-2.50	106.69	111.59
2	E	301	8DP	C15-N5-C16	-2.49	120.89	126.47
2	E	301	8DP	C14-C15-N5	2.49	118.82	114.07
2	H	301	8DP	O8-C15-C14	-2.42	119.52	124.32
2	B	301	8DP	C14-C15-N5	2.39	118.63	114.07
2	G	301	8DP	C4-C5-N1	-2.38	106.93	111.59
2	F	301	8DP	C14-C15-N5	2.33	118.51	114.07
2	F	301	8DP	C15-N5-C16	-2.32	121.27	126.47
2	B	301	8DP	O9-C16-N5	-2.31	117.75	121.86
2	D	301	8DP	C14-C15-N5	2.29	118.45	114.07
2	G	301	8DP	O8-C15-C14	-2.22	119.93	124.32
2	E	301	8DP	O5-C2-C3	2.14	112.79	109.67
2	B	301	8DP	C15-N5-C16	-2.06	121.85	126.47

There are no chirality outliers.

All (37) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	301	8DP	O7-C4-C5-N1
2	A	301	8DP	O4-C1-C2-C3
2	A	301	8DP	O4-C1-C2-O5
2	B	301	8DP	O7-C4-C5-N1
2	B	301	8DP	O4-C1-C2-C3
2	B	301	8DP	O4-C1-C2-O5
2	C	301	8DP	O7-C4-C5-N1
2	C	301	8DP	O4-C1-C2-C3
2	C	301	8DP	O4-C1-C2-O5
2	D	301	8DP	O7-C4-C5-N1
2	D	301	8DP	O4-C1-C2-C3
2	D	301	8DP	O4-C1-C2-O5
2	E	301	8DP	C3-C4-C5-N1
2	E	301	8DP	O7-C4-C5-N1
2	E	301	8DP	O4-C1-C2-C3
2	E	301	8DP	O4-C1-C2-O5
2	F	301	8DP	C3-C4-C5-N1
2	F	301	8DP	O7-C4-C5-N1
2	F	301	8DP	O4-C1-C2-C3
2	F	301	8DP	O4-C1-C2-O5
2	G	301	8DP	O7-C4-C5-N1
2	G	301	8DP	O4-C1-C2-C3
2	G	301	8DP	O4-C1-C2-O5

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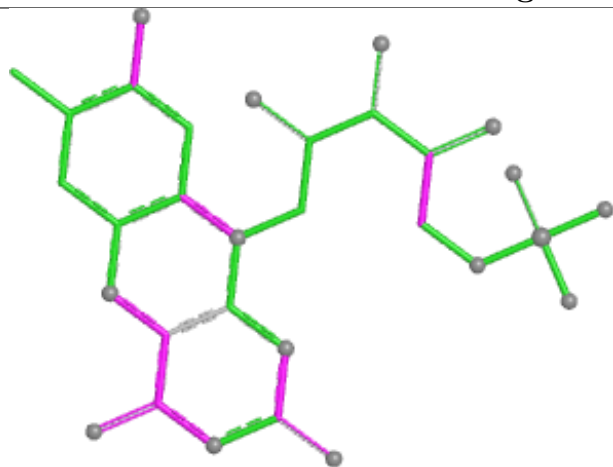
Mol	Chain	Res	Type	Atoms
2	H	301	8DP	O7-C4-C5-N1
2	H	301	8DP	O4-C1-C2-C3
2	H	301	8DP	O4-C1-C2-O5
2	B	301	8DP	O5-C2-C3-O6
2	B	301	8DP	C1-C2-C3-O6
2	B	301	8DP	C1-C2-C3-C4
2	B	301	8DP	O5-C2-C3-C4
2	A	301	8DP	C3-C4-C5-N1
2	B	301	8DP	C3-C4-C5-N1
2	C	301	8DP	C3-C4-C5-N1
2	D	301	8DP	C3-C4-C5-N1
2	G	301	8DP	C3-C4-C5-N1
2	H	301	8DP	C3-C4-C5-N1
2	F	301	8DP	O5-C2-C3-O6

There are no ring outliers.

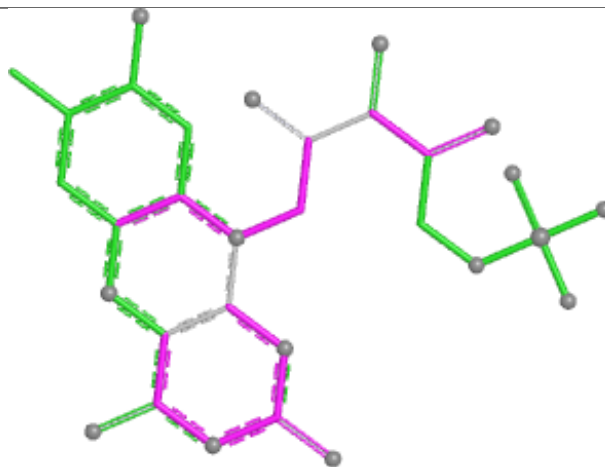
No monomer is involved in short contacts.

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.

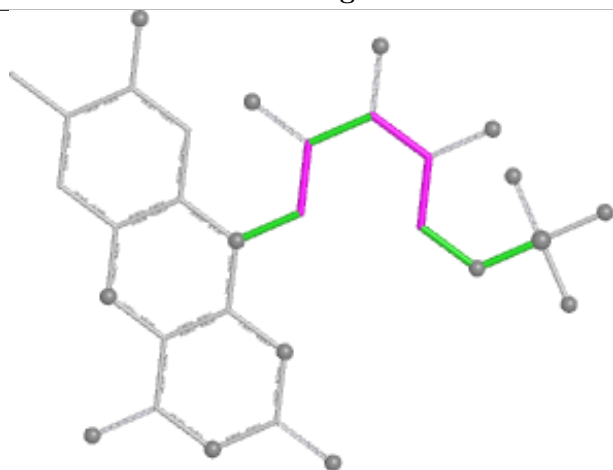
Ligand 8DP B 301



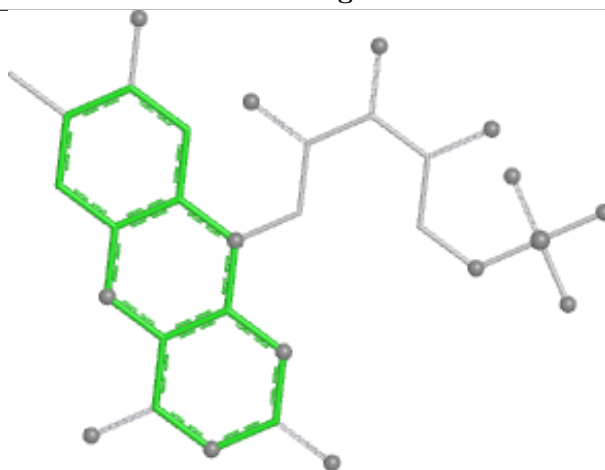
Bond lengths



Bond angles

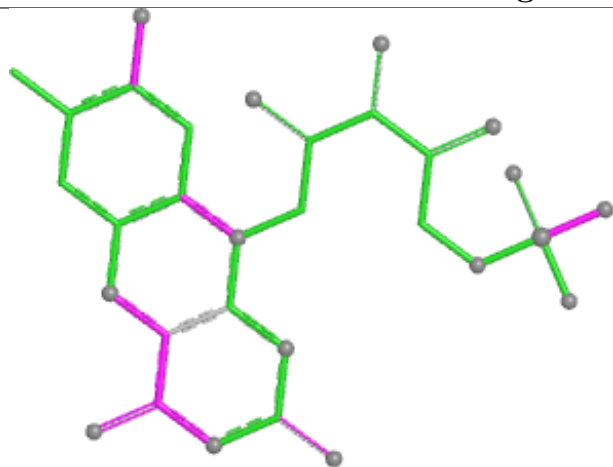


Torsions

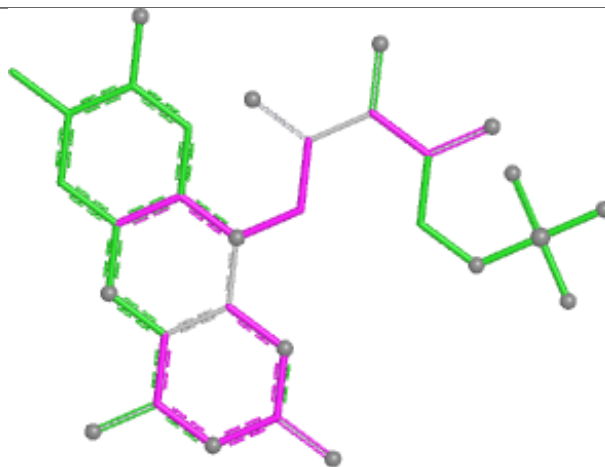


Rings

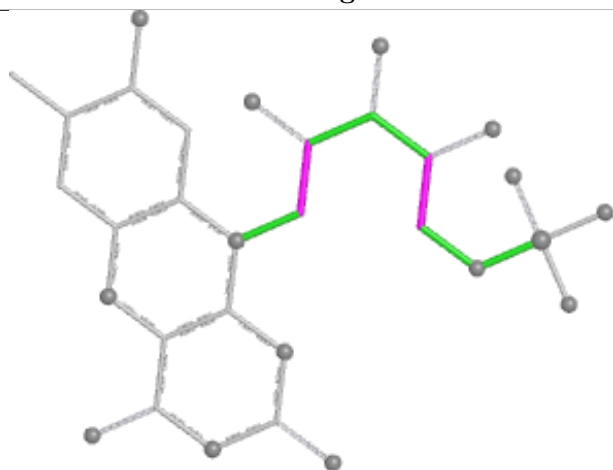
Ligand 8DP C 301



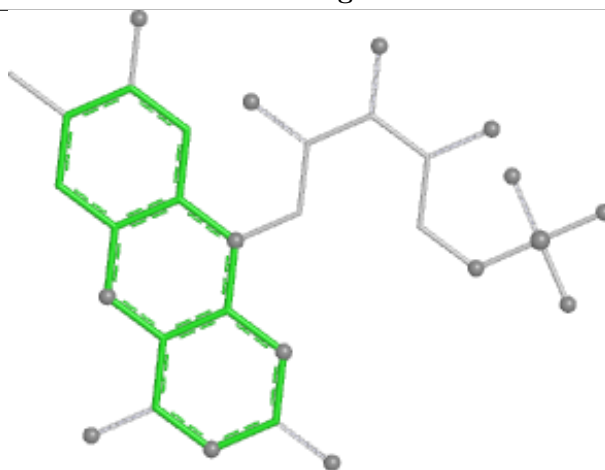
Bond lengths



Bond angles

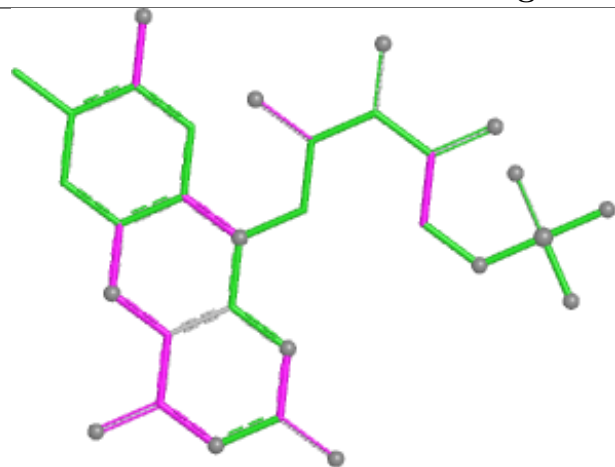


Torsions

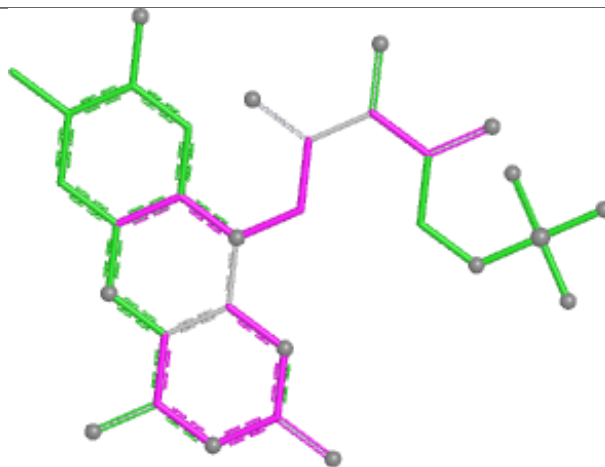


Rings

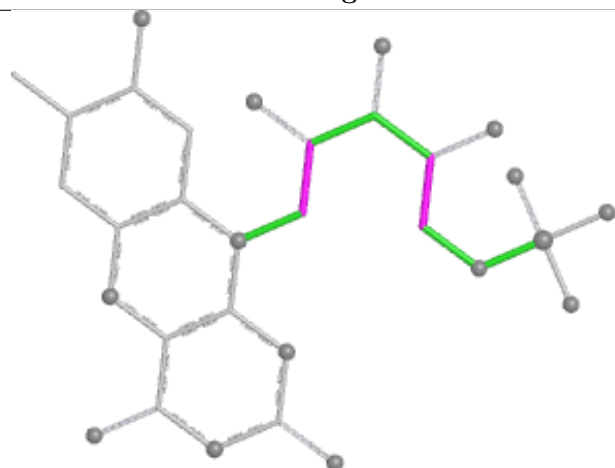
Ligand 8DP D 301



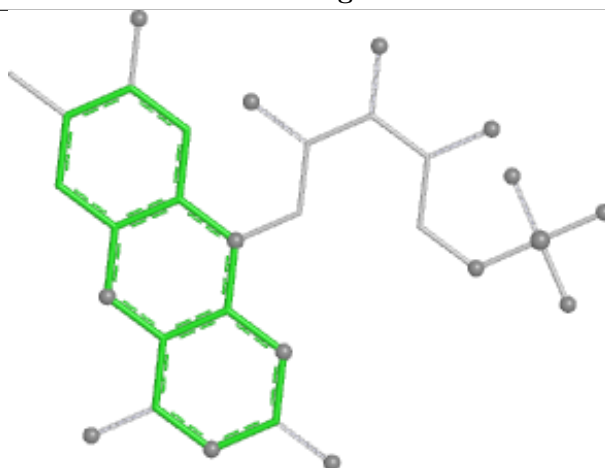
Bond lengths



Bond angles

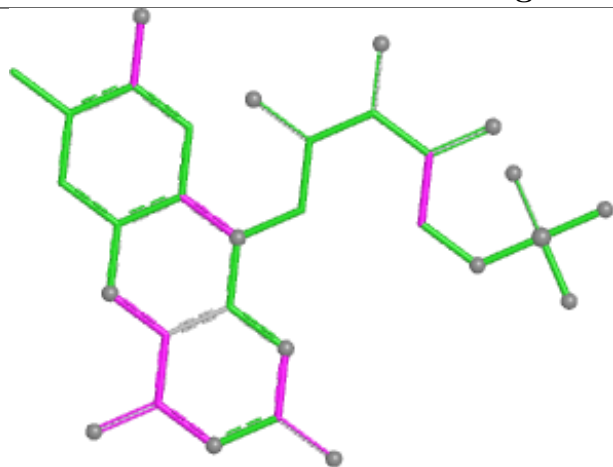


Torsions

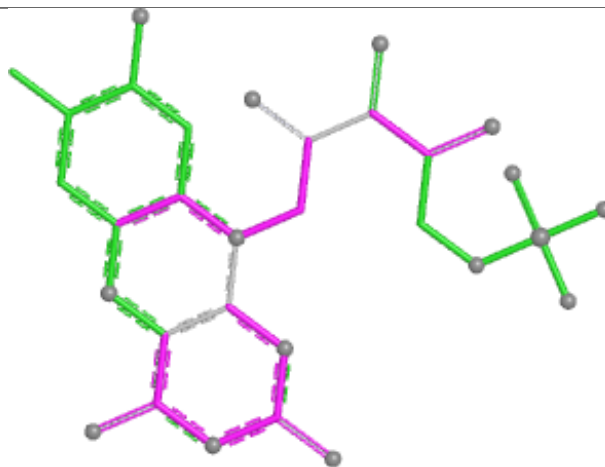


Rings

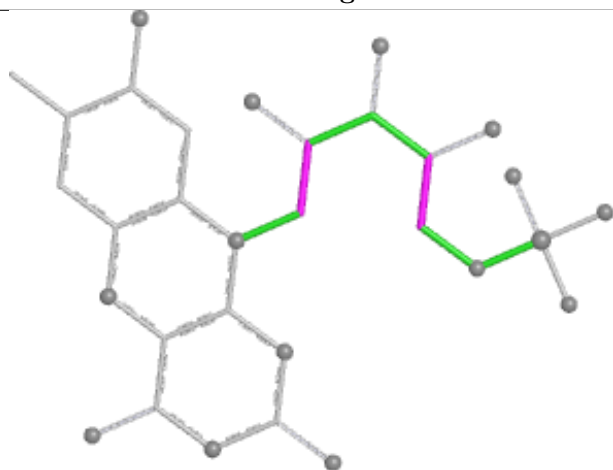
Ligand 8DP H 301



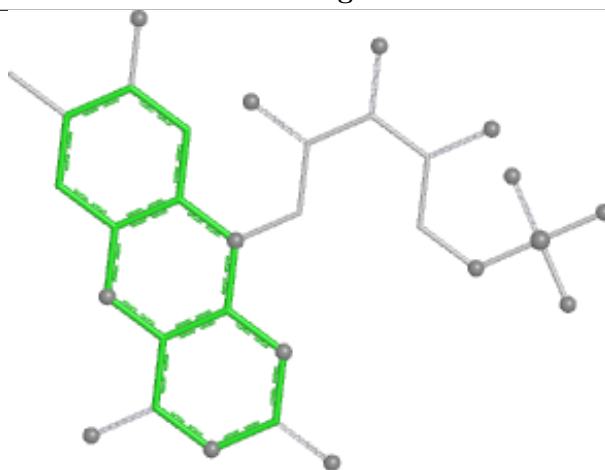
Bond lengths



Bond angles

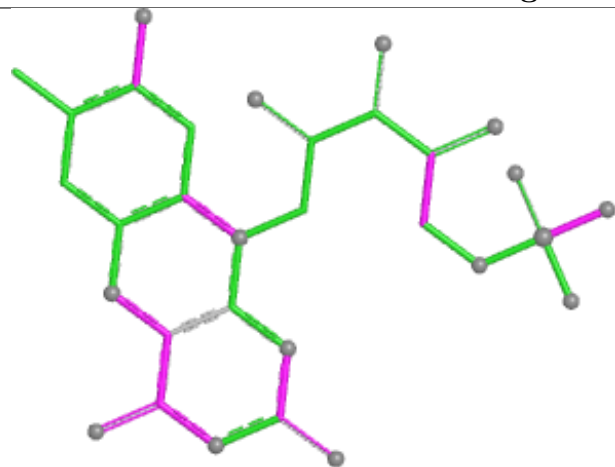


Torsions

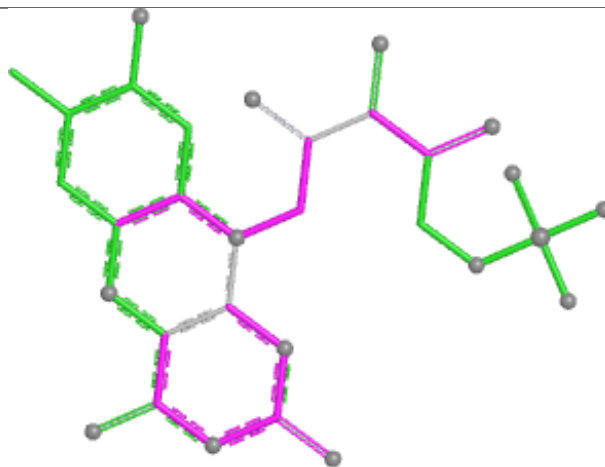


Rings

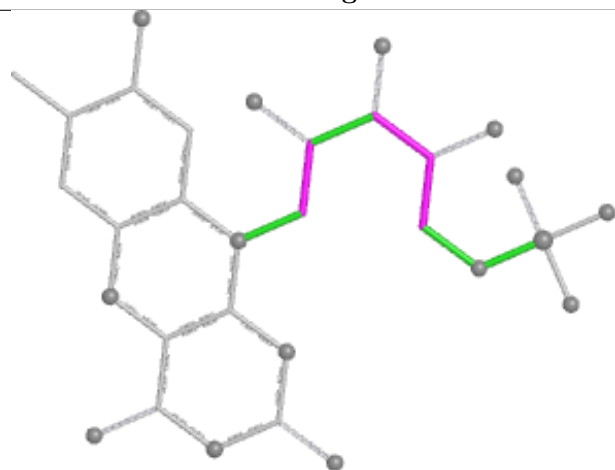
Ligand 8DP F 301



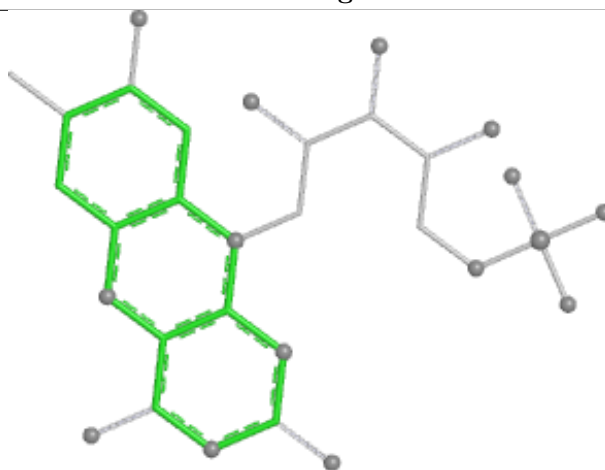
Bond lengths



Bond angles

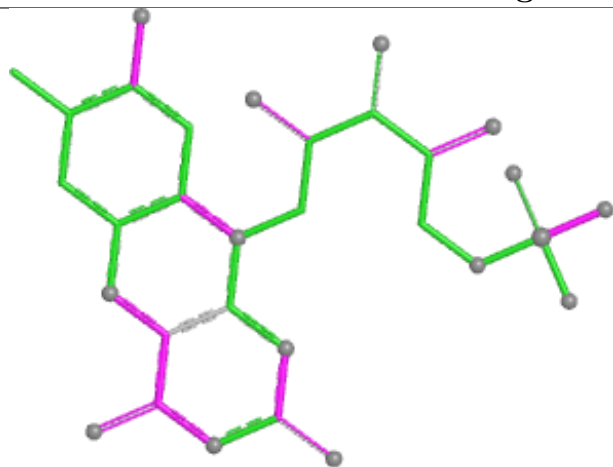


Torsions

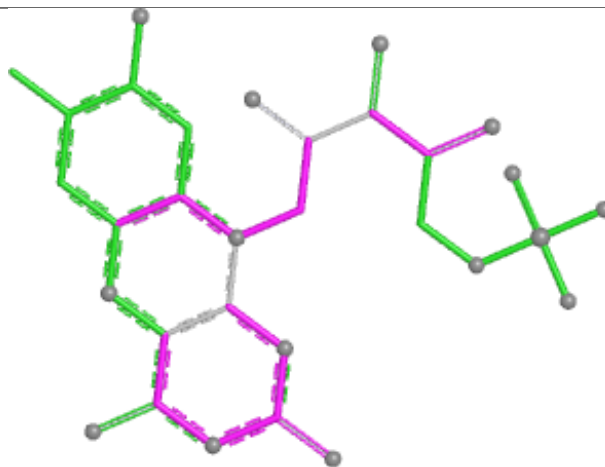


Rings

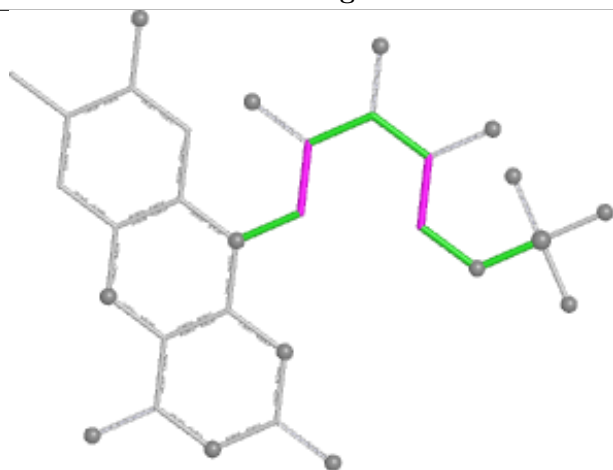
Ligand 8DP A 301



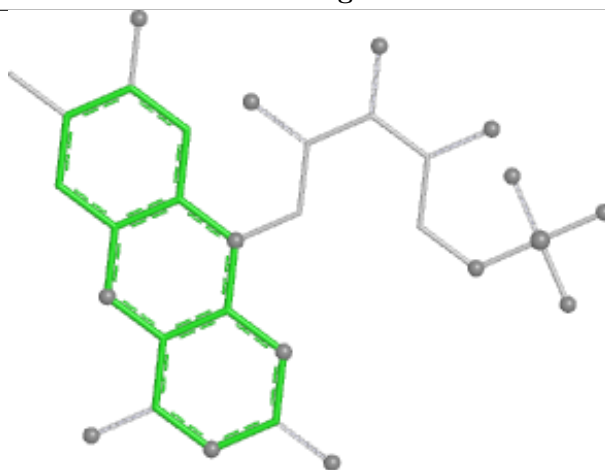
Bond lengths



Bond angles

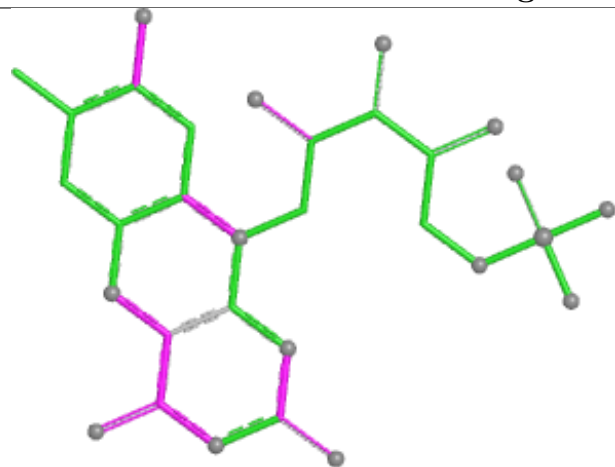


Torsions

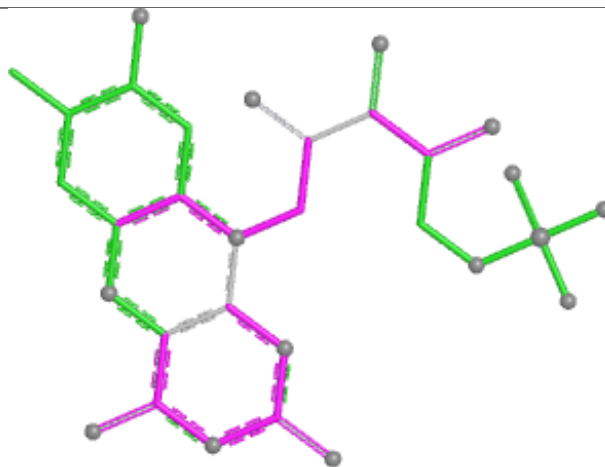


Rings

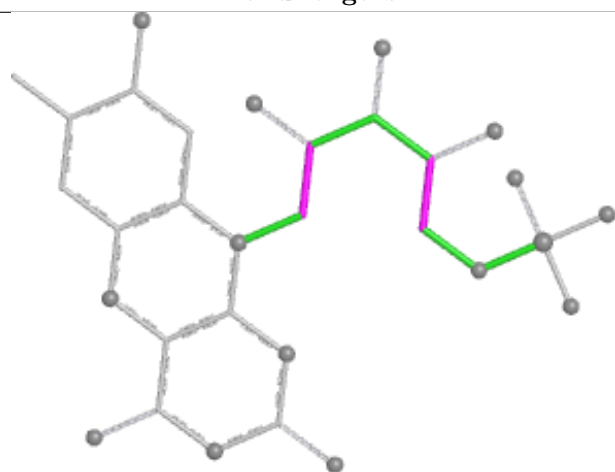
Ligand 8DP G 301



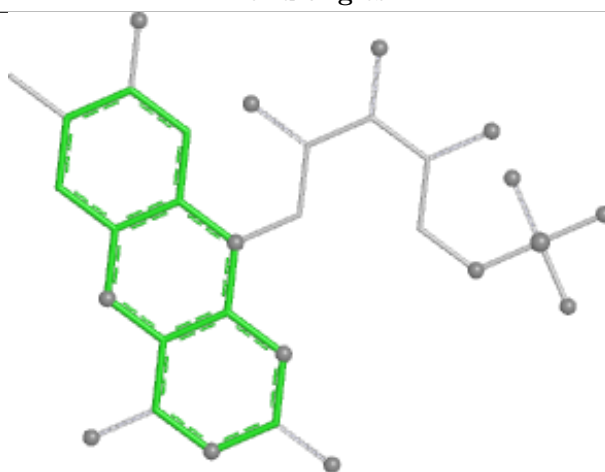
Bond lengths



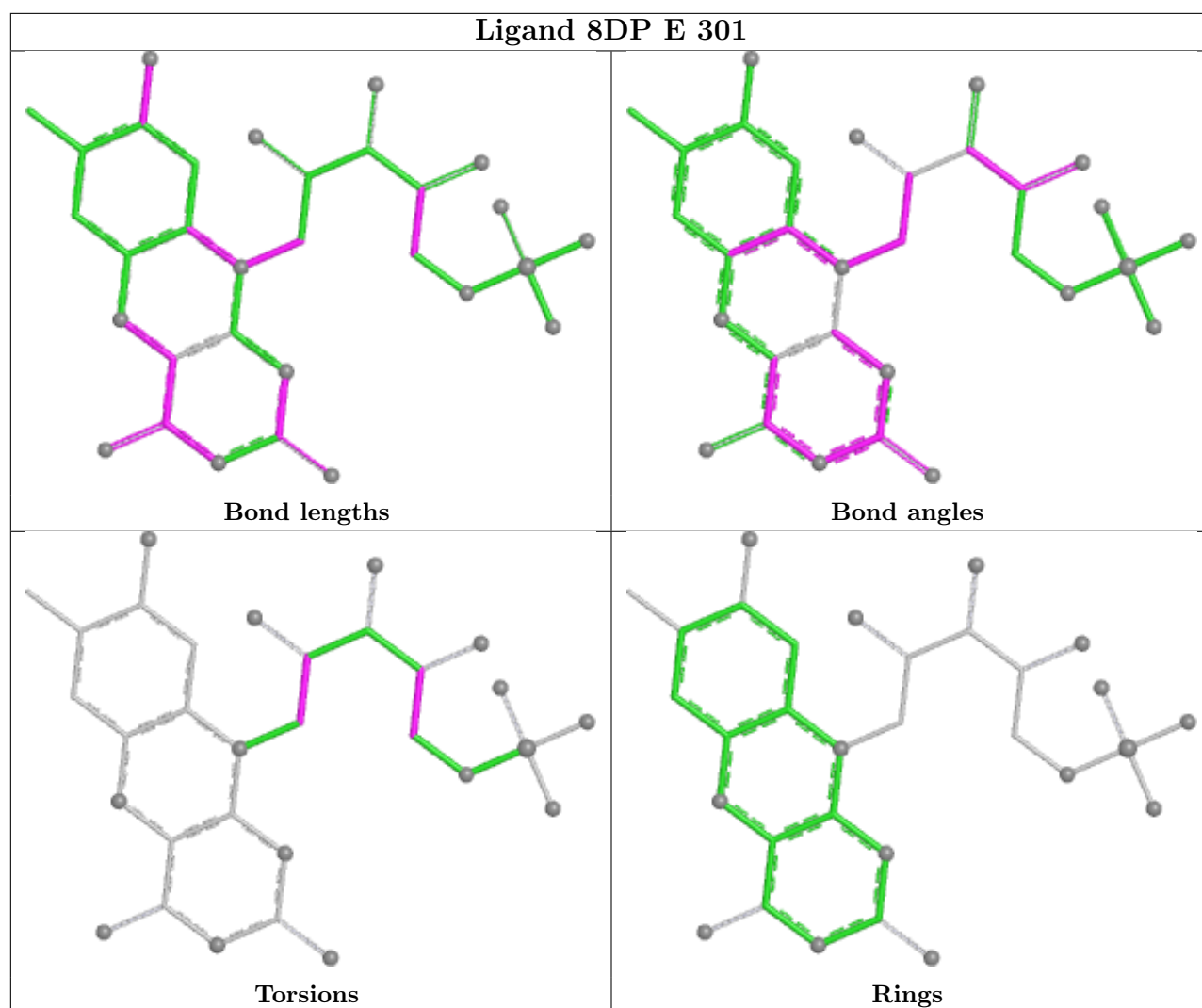
Bond angles



Torsions



Rings



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	252/257 (98%)	-0.07	13 (5%)	33 36	13, 21, 40, 95	8 (3%)
1	B	252/257 (98%)	0.11	12 (4%)	35 39	17, 25, 40, 88	6 (2%)
1	C	252/257 (98%)	0.17	13 (5%)	33 36	16, 28, 45, 93	1 (0%)
1	D	252/257 (98%)	-0.02	5 (1%)	65 69	17, 25, 39, 94	0
1	E	252/257 (98%)	0.07	6 (2%)	59 64	15, 26, 40, 82	2 (0%)
1	F	252/257 (98%)	-0.01	6 (2%)	59 64	11, 26, 43, 79	1 (0%)
1	G	252/257 (98%)	-0.16	2 (0%)	82 85	14, 22, 40, 99	1 (0%)
1	H	252/257 (98%)	0.18	4 (1%)	70 75	16, 29, 43, 102	2 (0%)
All	All	2016/2056 (98%)	0.03	61 (3%)	52 56	11, 25, 42, 102	21 (1%)

All (61) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	109[A]	LEU	4.8
1	C	251	VAL	4.7
1	A	109[A]	LEU	4.7
1	A	251	VAL	3.9
1	B	112[A]	GLU	3.8
1	A	117	THR	3.8
1	F	115	GLY	3.7
1	E	2	ALA	3.3
1	C	117	THR	3.3
1	A	112[A]	GLU	3.3
1	C	253	GLU	3.3
1	B	115	GLY	3.2
1	B	114	ASN	3.1
1	C	2	ALA	3.1
1	B	113[A]	LEU	3.0
1	E	253	GLU	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	114	ASN	3.0
1	D	2	ALA	3.0
1	F	2	ALA	2.9
1	A	111[A]	ASP	2.9
1	B	2	ALA	2.8
1	B	111[A]	ASP	2.8
1	F	251	VAL	2.8
1	C	250	LYS	2.8
1	D	253	GLU	2.8
1	A	2	ALA	2.7
1	H	251	VAL	2.7
1	C	113	LEU	2.7
1	C	111	ASP	2.6
1	B	117	THR	2.6
1	A	113[A]	LEU	2.5
1	F	253	GLU	2.4
1	D	56	ASP	2.4
1	B	253	GLU	2.4
1	C	116	ASP	2.4
1	A	250	LYS	2.4
1	E	251	VAL	2.4
1	G	251	VAL	2.4
1	C	115	GLY	2.3
1	E	109	LEU	2.3
1	F	111	ASP	2.3
1	B	252	GLU	2.3
1	D	251	VAL	2.3
1	H	109	LEU	2.3
1	D	250	LYS	2.3
1	A	115	GLY	2.2
1	C	114	ASN	2.2
1	C	252	GLU	2.2
1	C	249	VAL	2.2
1	E	108	ASN	2.2
1	G	250	LYS	2.2
1	H	249	VAL	2.1
1	E	72	TRP	2.1
1	C	245	LYS	2.1
1	A	247	GLY	2.1
1	H	245	LYS	2.1
1	B	251	VAL	2.1
1	A	253	GLU	2.0

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Mol	Chain	Res	Type	RSRZ
1	F	109	LEU	2.0
1	B	116	ASP	2.0
1	A	249	VAL	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](#)

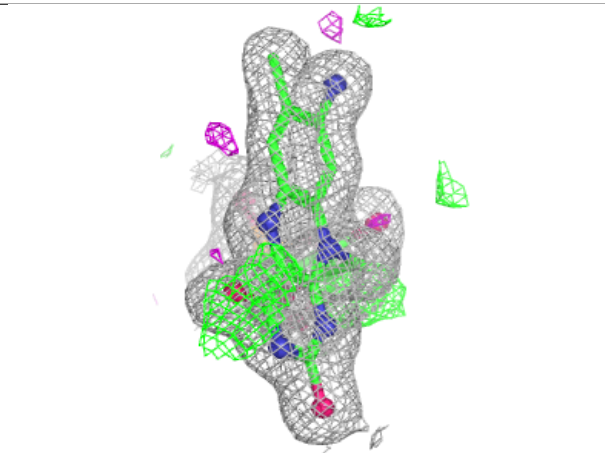
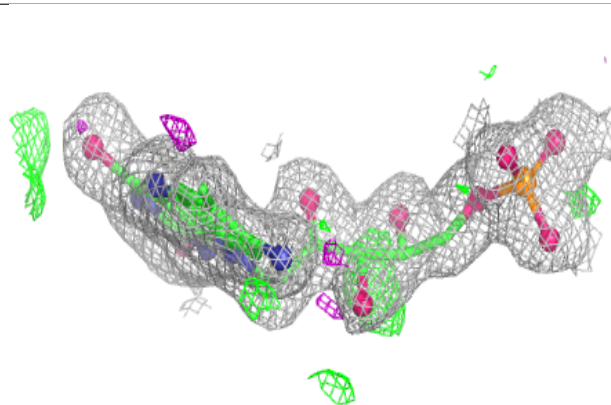
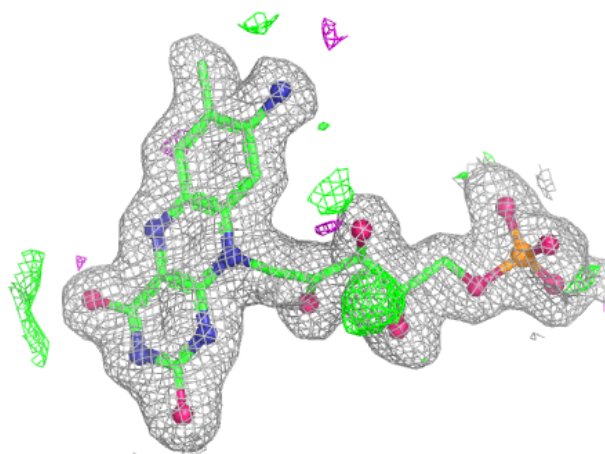
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	8DP	B	301	31/31	0.96	0.07	20,23,30,33	0
2	8DP	E	301	31/31	0.96	0.06	19,24,32,33	0
2	8DP	H	301	31/31	0.96	0.06	16,21,28,34	0
2	8DP	D	301	31/31	0.97	0.06	16,21,27,33	0
2	8DP	A	301	31/31	0.97	0.06	14,17,29,32	0
2	8DP	F	301	31/31	0.97	0.06	18,22,29,33	0
2	8DP	G	301	31/31	0.97	0.06	14,18,24,31	0
2	8DP	C	301	31/31	0.97	0.06	16,19,25,32	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.

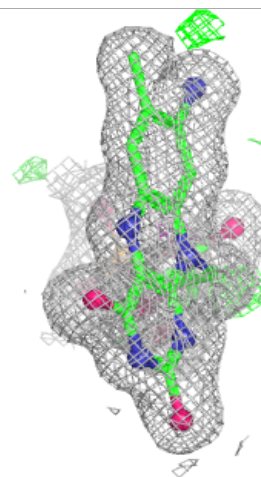
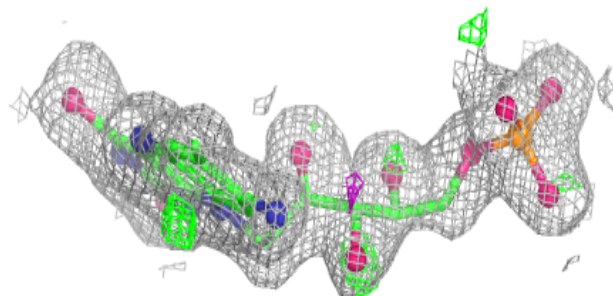
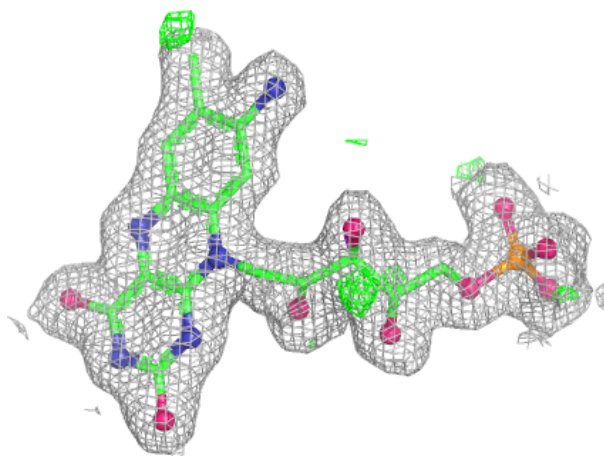
Electron density around 8DP B 301:

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



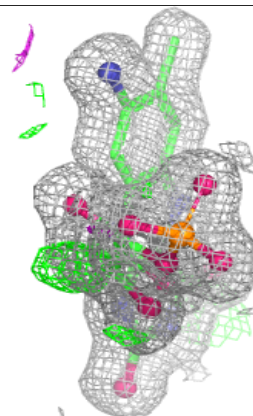
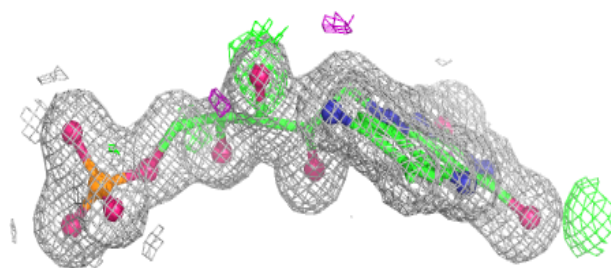
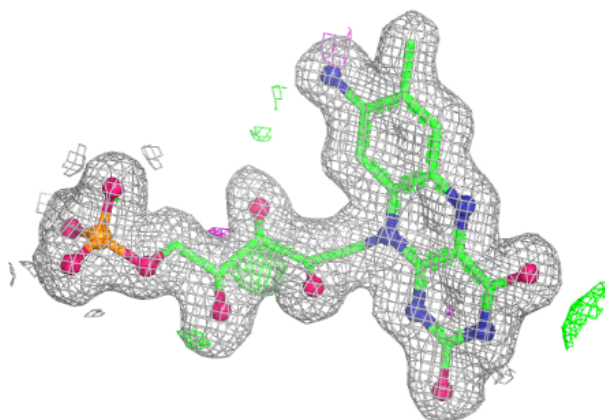
Electron density around 8DP E 301:

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



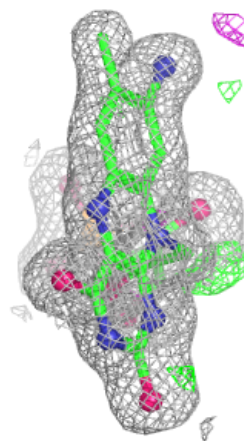
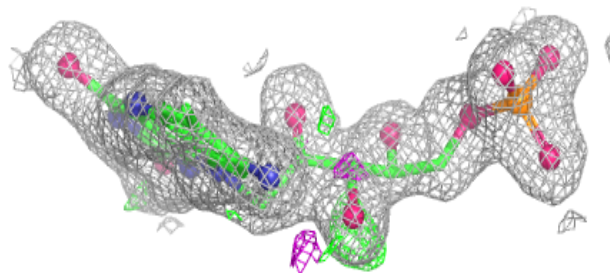
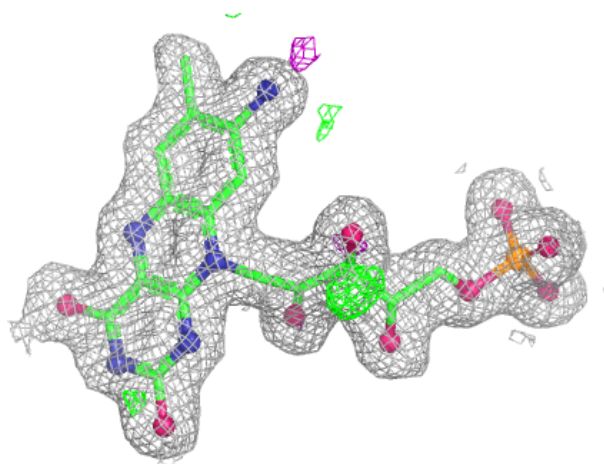
Electron density around 8DP H 301:

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



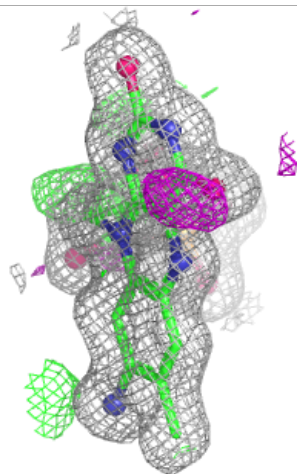
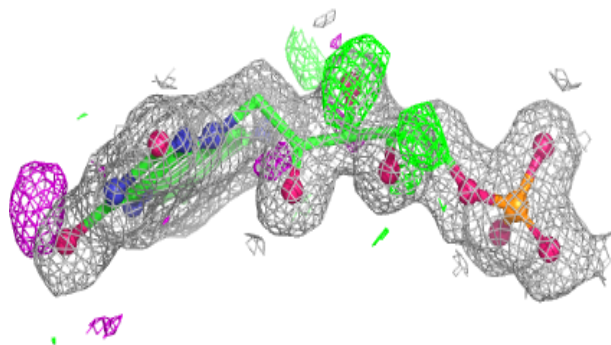
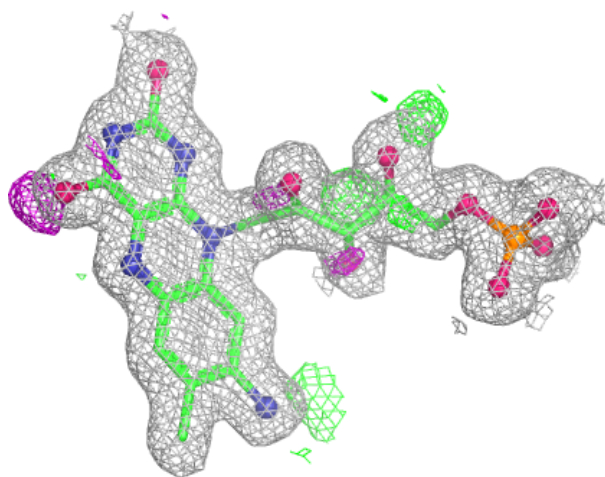
Electron density around 8DP D 301:

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



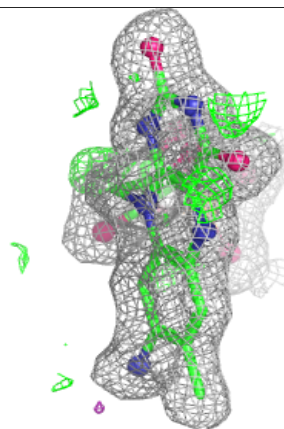
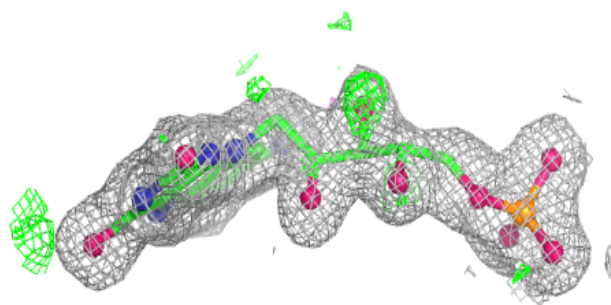
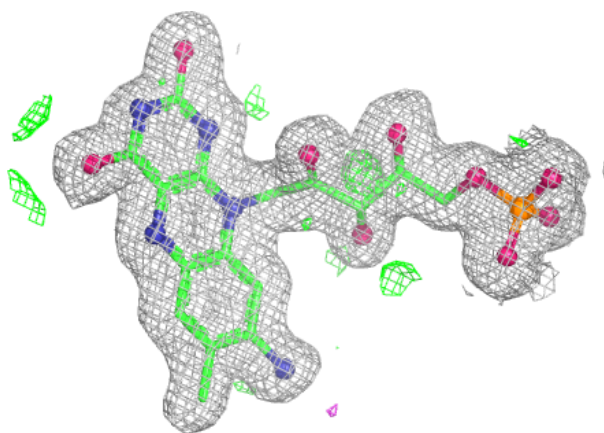
Electron density around 8DP A 301:

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



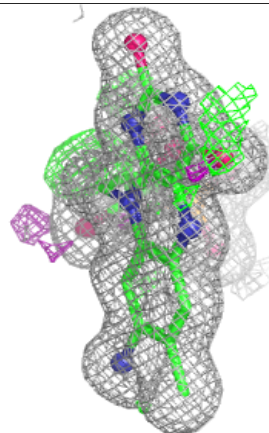
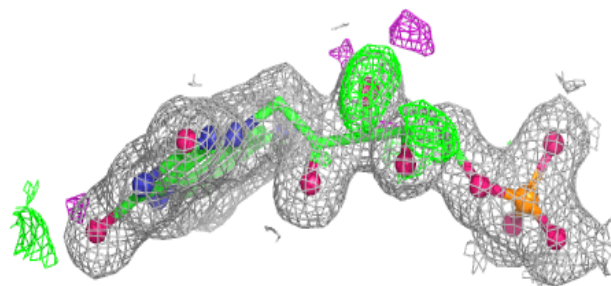
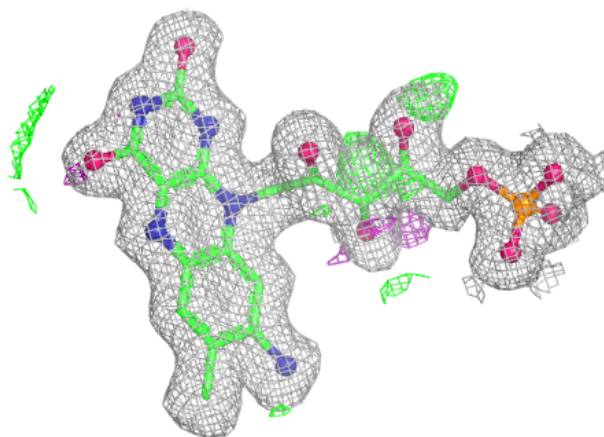
Electron density around 8DP F 301:

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



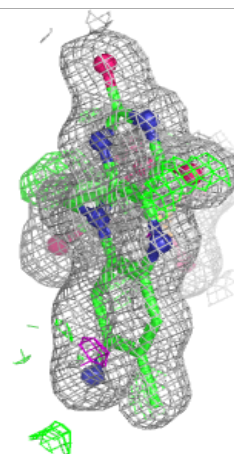
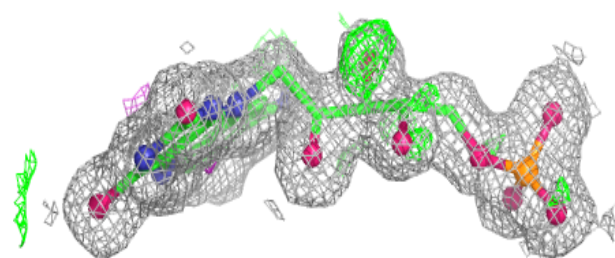
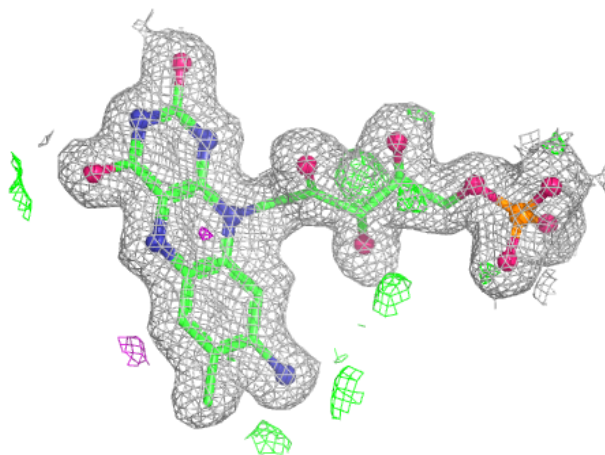
Electron density around 8DP G 301:

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



Electron density around 8DP C 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.