



Full wwPDB X-ray Structure Validation Report ⓘ

May 15, 2020 – 06:10 am BST

PDB ID : 5CJW
Title : Isobutyryl-CoA mutase fused with bound adenosylcobalamin, GDP, Mg (holo-IcmF/GDP), and substrate pivalyl-coenzyme A
Authors : Jost, M.; Drennan, C.L.
Deposited on : 2015-07-15
Resolution : 3.40 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

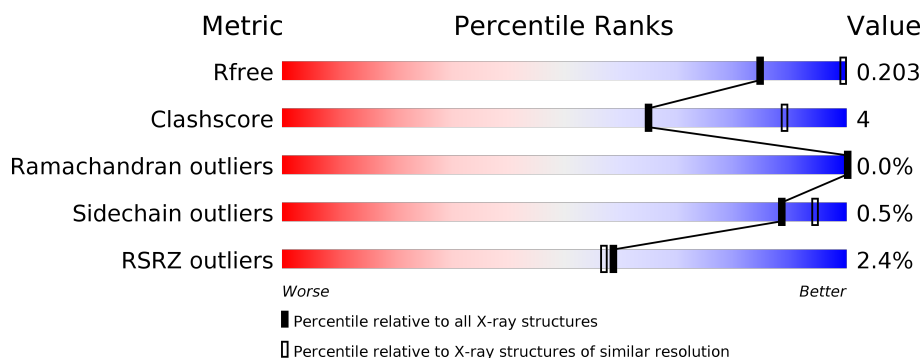
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.40 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}	130704	1026 (3.48-3.32)
Clashscore	141614	1055 (3.48-3.32)
Ramachandran outliers	138981	1038 (3.48-3.32)
Sidechain outliers	138945	1038 (3.48-3.32)
RSRZ outliers	127900	2173 (3.50-3.30)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria 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	1113	<div> <div>4%</div> <div> <div></div> <div>85%</div> <div>10%</div> <div>.</div> </div> </div>
1	B	1113	<div> <div>%</div> <div> <div></div> <div>86%</div> <div>9%</div> <div>5%</div> </div> </div>

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 16502 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 Isobutyryl-CoA mutase fused.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	1063	Total	C	N	O	S	0	0	0
			8098	5060	1453	1548	37			
1	B	1060	Total	C	N	O	S	0	0	0
			8045	5022	1447	1539	37			

There are 40 discrepancies between the modelled and reference sequences:

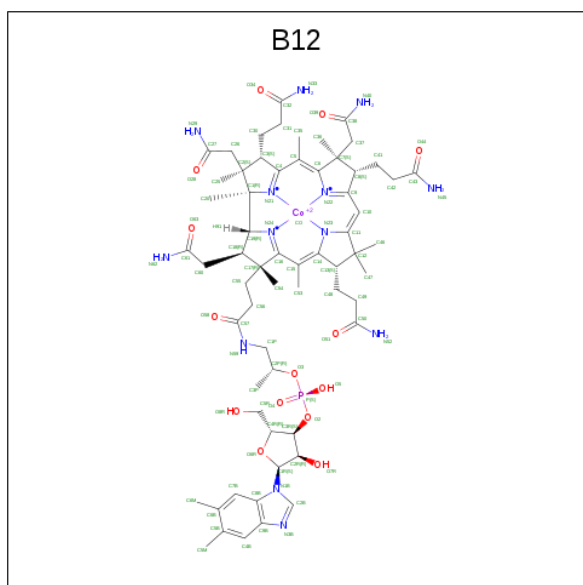
Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	initiating methionine	UNP Q1LRY0
A	-18	GLY	-	expression tag	UNP Q1LRY0
A	-17	SER	-	expression tag	UNP Q1LRY0
A	-16	SER	-	expression tag	UNP Q1LRY0
A	-15	HIS	-	expression tag	UNP Q1LRY0
A	-14	HIS	-	expression tag	UNP Q1LRY0
A	-13	HIS	-	expression tag	UNP Q1LRY0
A	-12	HIS	-	expression tag	UNP Q1LRY0
A	-11	HIS	-	expression tag	UNP Q1LRY0
A	-10	HIS	-	expression tag	UNP Q1LRY0
A	-9	SER	-	expression tag	UNP Q1LRY0
A	-8	SER	-	expression tag	UNP Q1LRY0
A	-7	GLY	-	expression tag	UNP Q1LRY0
A	-6	LEU	-	expression tag	UNP Q1LRY0
A	-5	VAL	-	expression tag	UNP Q1LRY0
A	-4	PRO	-	expression tag	UNP Q1LRY0
A	-3	ARG	-	expression tag	UNP Q1LRY0
A	-2	GLY	-	expression tag	UNP Q1LRY0
A	-1	SER	-	expression tag	UNP Q1LRY0
A	0	HIS	-	expression tag	UNP Q1LRY0
B	-19	MET	-	initiating methionine	UNP Q1LRY0
B	-18	GLY	-	expression tag	UNP Q1LRY0
B	-17	SER	-	expression tag	UNP Q1LRY0
B	-16	SER	-	expression tag	UNP Q1LRY0
B	-15	HIS	-	expression tag	UNP Q1LRY0

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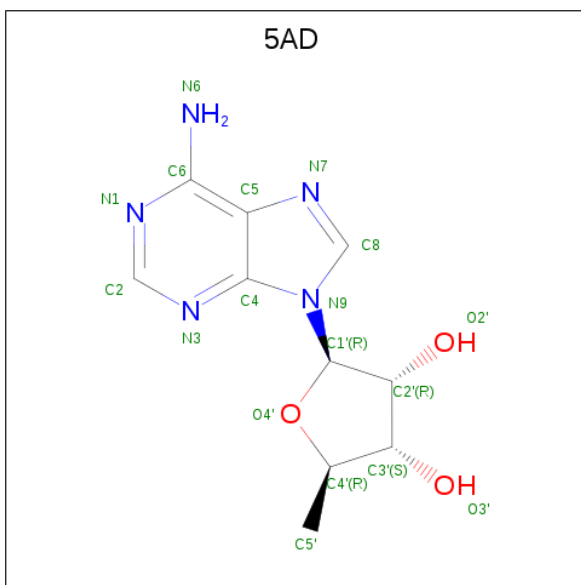
Chain	Residue	Modelled	Actual	Comment	Reference
B	-14	HIS	-	expression tag	UNP Q1LRY0
B	-13	HIS	-	expression tag	UNP Q1LRY0
B	-12	HIS	-	expression tag	UNP Q1LRY0
B	-11	HIS	-	expression tag	UNP Q1LRY0
B	-10	HIS	-	expression tag	UNP Q1LRY0
B	-9	SER	-	expression tag	UNP Q1LRY0
B	-8	SER	-	expression tag	UNP Q1LRY0
B	-7	GLY	-	expression tag	UNP Q1LRY0
B	-6	LEU	-	expression tag	UNP Q1LRY0
B	-5	VAL	-	expression tag	UNP Q1LRY0
B	-4	PRO	-	expression tag	UNP Q1LRY0
B	-3	ARG	-	expression tag	UNP Q1LRY0
B	-2	GLY	-	expression tag	UNP Q1LRY0
B	-1	SER	-	expression tag	UNP Q1LRY0
B	0	HIS	-	expression tag	UNP Q1LRY0

- Molecule 2 is COBALAMIN (three-letter code: B12) (formula: $C_{62}H_{89}CoN_{13}O_{14}P$).



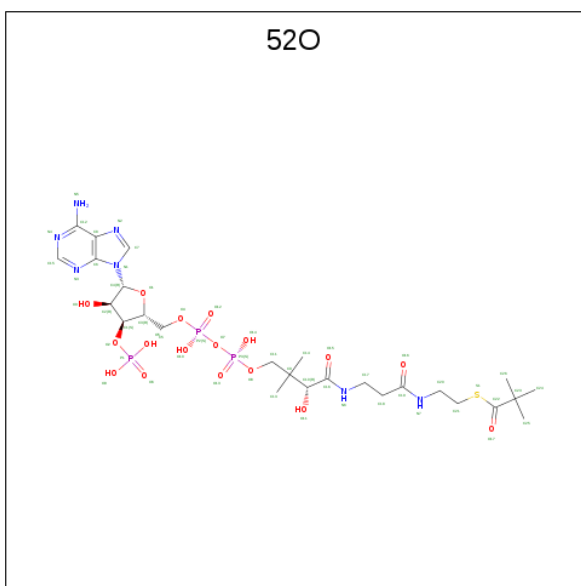
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total	C	Co	N	O	P	0	0
			91	62	1	13	14	1		
2	B	1	Total	C	Co	N	O	P	0	0
			91	62	1	13	14	1		

- Molecule 3 is 5'-DEOXYADENOSINE (three-letter code: 5AD) (formula: $C_{10}H_{13}N_5O_3$).



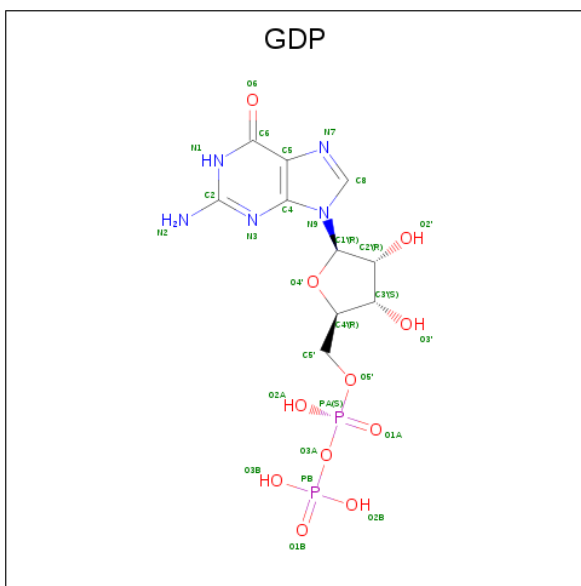
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	1
			36	20	10	6		

- Molecule 4 is pivalyl-coenzyme A (three-letter code: 52O) (formula: $C_{26}H_{44}N_7O_{17}P_3S$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	S	0	0
			54	26	7	17	3	1		
4	B	1	Total	C	N	O	P		0	0
			27	10	5	10	2			

- Molecule 5 is GUANOSINE-5'-DIPHOSPHATE (three-letter code: GDP) (formula:

$$\text{C}_{10}\text{H}_{15}\text{N}_5\text{O}_{11}\text{P}_2).$$


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
5	A	1	Total 28	C 10	N 5	O 11	P 2	0	0
5	B	1	Total 28	C 10	N 5	O 11	P 2	0	0

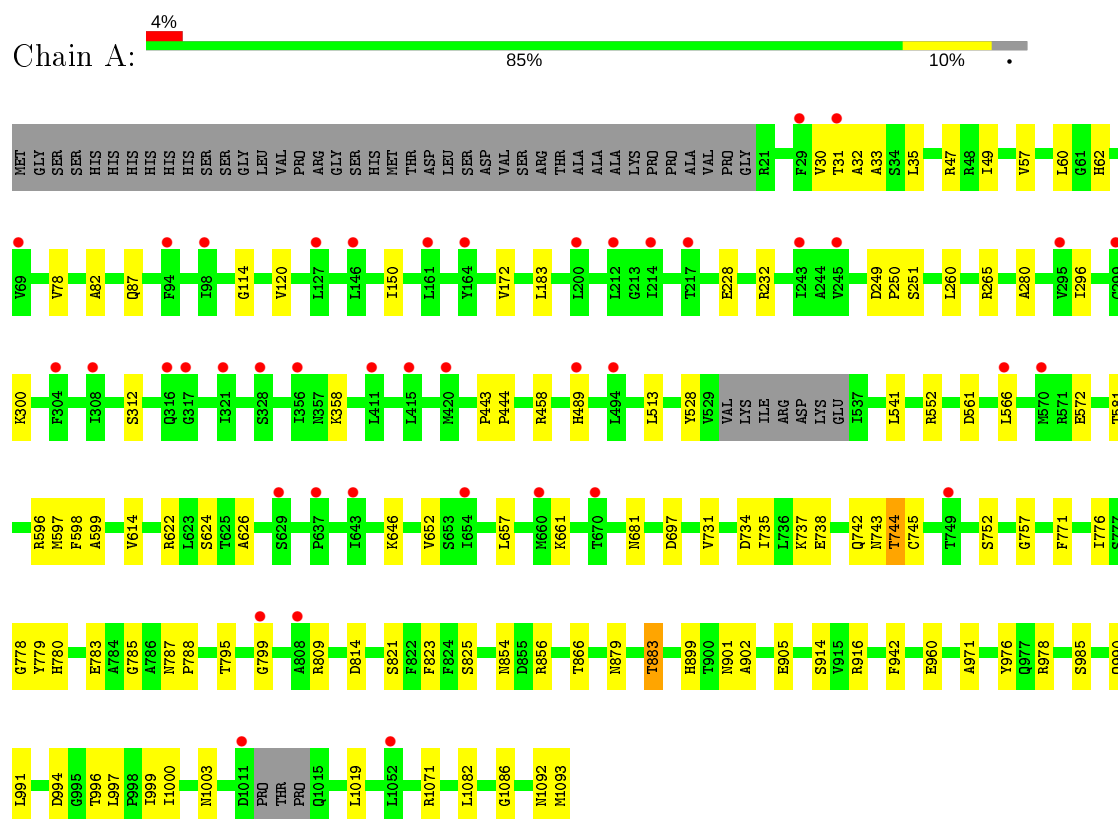
- Molecule 6 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	B	2	Total Mg 2 2	0	0
6	A	2	Total Mg 2 2	0	0

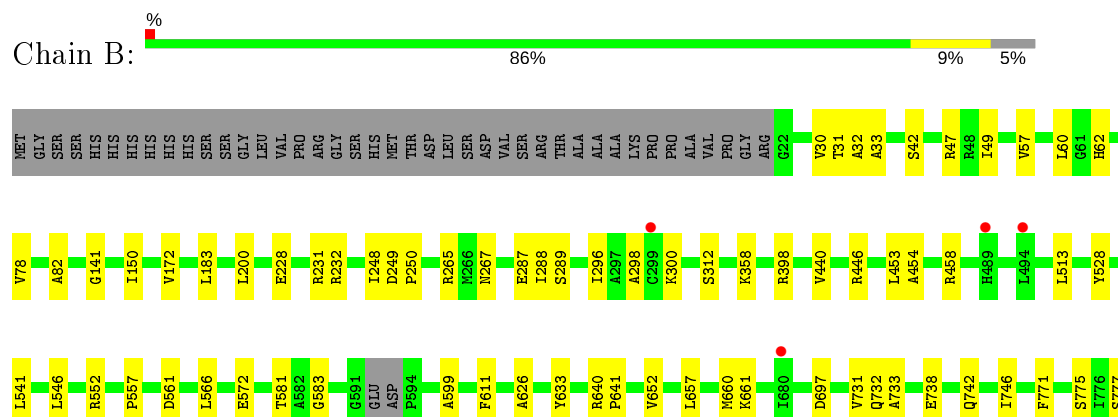
3 Residue-property plots

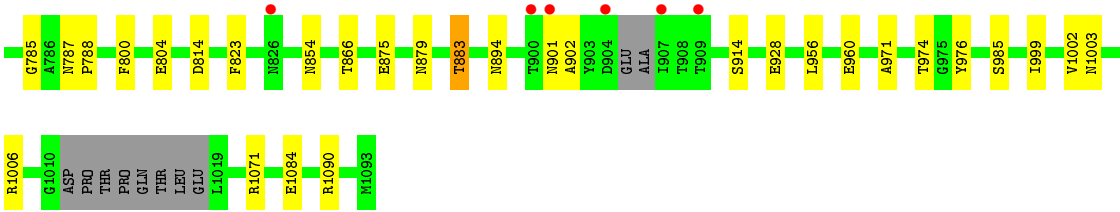
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are 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: Isobutyryl-CoA mutase fused



- Molecule 1: Isobutyryl-CoA mutase fused





4 Data and refinement statistics

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, α , β , γ	317.52Å 317.52Å 343.04Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	100.00 – 3.40 127.62 – 3.40	Depositor EDS
% Data completeness (in resolution range)	99.2 (100.00-3.40) 99.3 (127.62-3.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.14	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.04 (at 3.41Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.183 , 0.200 0.186 , 0.203	Depositor DCC
R_{free} test set	4549 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	104.9	Xtriage
Anisotropy	0.418	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 77.0	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	0.95	EDS
Total number of atoms	16502	wwPDB-VP
Average B, all atoms (Å ²)	108.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.99% 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: GDP, MG, 52O, B12, 5AD

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	0.21	0/8240	0.39	2/11165 (0.0%)
1	B	0.22	0/8183	0.39	3/11083 (0.0%)
All	All	0.21	0/16423	0.39	5/22248 (0.0%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	552	ARG	NE-CZ-NH2	-10.35	115.12	120.30
1	A	552	ARG	NE-CZ-NH1	10.05	125.33	120.30
1	B	552	ARG	NE-CZ-NH2	8.98	124.79	120.30
1	B	552	ARG	NE-CZ-NH1	-8.41	116.09	120.30
1	B	552	ARG	CD-NE-CZ	5.08	130.71	123.60

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	8098	0	7834	72	0
1	B	8045	0	7756	54	0
2	A	91	0	88	14	0
2	B	91	0	88	6	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	36	0	26	4	0
4	A	54	0	41	7	0
4	B	27	0	12	0	0
5	A	28	0	12	2	0
5	B	28	0	12	2	0
6	A	2	0	0	0	0
6	B	2	0	0	0	0
All	All	16502	0	15869	133	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 4.

All (133) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:87:GLN:HE22	2:A:1101:B12:H421	1.46	0.80
1:A:783:GLU:HB3	2:A:1101:B12:H532	1.68	0.76
1:B:454:ALA:HA	1:B:956:LEU:HD22	1.79	0.64
3:A:1102[A]:5AD:H5'1	4:A:1103:52O:H42	1.80	0.63
1:A:31:THR:HG22	1:A:82:ALA:HB3	1.82	0.61
1:A:905:GLU:O	2:A:1101:B12:N29	2.25	0.61
1:A:681:ASN:ND2	1:A:742:GLN:OE1	2.35	0.60
1:A:622:ARG:NH1	4:A:1103:52O:O13	2.21	0.59
1:A:734:ASP:HB3	1:A:737:LYS:HZ2	1.68	0.59
1:A:742:GLN:HE21	2:A:1101:B12:H411	1.67	0.58
1:A:990:GLN:O	1:A:994:ASP:HB2	2.04	0.58
1:B:31:THR:HG22	1:B:82:ALA:HB3	1.84	0.57
1:A:1082:LEU:O	1:A:1086:GLY:N	2.37	0.56
1:A:821:SER:OG	4:A:1103:52O:O11	2.19	0.56
1:A:856:ARG:NH1	4:A:1103:52O:O9	2.39	0.56
1:A:780:HIS:HA	2:A:1101:B12:H461	1.88	0.55
2:A:1101:B12:H261	3:A:1102[A]:5AD:H3'	1.89	0.55
1:A:35:LEU:HB3	1:A:646:LYS:HE3	1.89	0.55
1:A:33:ALA:HB3	1:A:62:HIS:HA	1.89	0.55
1:B:738:GLU:OE2	1:B:742:GLN:NE2	2.39	0.54
1:A:596:ARG:HG2	1:A:624:SER:HB2	1.89	0.54
1:B:1002:VAL:O	1:B:1006:ARG:NH1	2.36	0.54
1:A:561:ASP:HB2	1:B:561:ASP:HB2	1.90	0.54
1:B:33:ALA:HB3	1:B:62:HIS:HA	1.90	0.53
1:B:267:ASN:HD21	1:B:1090:ARG:HA	1.73	0.53
1:A:458:ARG:NH1	1:A:960:GLU:OE2	2.39	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:942:PHE:CZ	1:B:557:PRO:HB3	2.44	0.53
1:A:30:VAL:HG23	1:A:78:VAL:HG11	1.91	0.53
1:B:583:GLY:HA3	1:B:894:ASN:HD22	1.74	0.53
1:B:30:VAL:HG23	1:B:78:VAL:HG11	1.90	0.53
1:A:528:TYR:HB2	1:A:541:LEU:HD21	1.90	0.52
1:B:528:TYR:HB2	1:B:541:LEU:HD21	1.90	0.52
1:B:513:LEU:HD21	1:B:566:LEU:HB2	1.91	0.52
1:A:513:LEU:HD21	1:A:566:LEU:HB2	1.91	0.51
1:A:785:GLY:HA3	1:A:985:SER:HB2	1.92	0.51
1:A:735:ILE:HG13	1:A:752:SER:HB3	1.92	0.51
1:A:991:LEU:HD22	1:A:996:THR:HB	1.92	0.51
1:A:32:ALA:HB2	1:A:60:LEU:HB2	1.94	0.50
1:B:228:GLU:OE1	1:B:232:ARG:NH1	2.45	0.50
1:A:228:GLU:OE1	1:A:232:ARG:NH1	2.45	0.49
1:B:785:GLY:HA3	1:B:985:SER:HB2	1.94	0.49
1:B:32:ALA:HB2	1:B:60:LEU:HB2	1.95	0.49
1:A:916:ARG:NH1	1:B:875:GLU:OE2	2.45	0.49
1:A:598:PHE:CE2	4:A:1103:52O:H40	2.48	0.49
1:B:787:ASN:HB2	1:B:788:PRO:HD2	1.95	0.49
1:B:777:SER:HA	1:B:823:PHE:HB3	1.95	0.49
1:A:787:ASN:HB2	1:A:788:PRO:HD2	1.94	0.49
1:A:62:HIS:CE1	1:A:743:ASN:HB3	2.47	0.48
1:B:296:ILE:O	1:B:300:LYS:HG3	2.13	0.48
1:A:814:ASP:CG	1:A:854:ASN:HD22	2.17	0.48
1:A:1000:ILE:N	1:B:928:GLU:OE2	2.36	0.48
1:A:250:PRO:HD3	1:A:312:SER:HB2	1.96	0.48
1:B:572:GLU:HG2	1:B:581:THR:HG21	1.96	0.48
1:B:814:ASP:CG	1:B:854:ASN:HD22	2.17	0.48
1:A:731:VAL:HG12	1:A:771:PHE:HE1	1.79	0.47
1:B:879:ASN:O	1:B:883:THR:OG1	2.29	0.47
1:A:296:ILE:O	1:A:300:LYS:HG3	2.14	0.47
1:A:572:GLU:HG2	1:A:581:THR:HG21	1.97	0.47
1:A:697:ASP:OD1	1:A:1071:ARG:NH1	2.48	0.47
1:A:779:TYR:OH	3:A:1102[B]:5AD:O3'	2.19	0.47
1:B:231:ARG:NH1	1:B:1084:GLU:HA	2.30	0.47
1:A:978:ARG:NE	1:A:1093:MET:OXT	2.46	0.47
1:B:697:ASP:OD1	1:B:1071:ARG:NH1	2.48	0.46
1:A:734:ASP:O	1:A:737:LYS:HG2	2.16	0.46
1:B:458:ARG:NH1	1:B:960:GLU:OE2	2.43	0.46
1:B:657:LEU:HG	1:B:661:LYS:HE3	1.98	0.46
1:B:200:LEU:HD11	1:B:298:ALA:HA	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:757:GLY:HA3	1:A:809:ARG:NH2	2.31	0.45
1:A:779:TYR:CD2	1:A:825:SER:HB2	2.51	0.45
2:B:1101:B12:H353	2:B:1101:B12:H311	1.98	0.45
2:B:1101:B12:H473	2:B:1101:B12:H481	1.77	0.45
1:A:172:VAL:HG21	1:A:183:LEU:HD22	1.97	0.45
2:B:1101:B12:H531	2:B:1101:B12:H552	1.98	0.45
1:B:172:VAL:HG21	1:B:183:LEU:HD22	1.98	0.45
1:B:731:VAL:O	1:B:733:ALA:N	2.50	0.45
1:A:999:ILE:N	1:A:1003:ASN:OD1	2.46	0.45
1:A:776:ILE:HD13	1:A:799:GLY:HA2	1.98	0.45
1:A:598:PHE:H	1:A:899:HIS:CE1	2.34	0.45
1:B:250:PRO:HD3	1:B:312:SER:HB2	1.98	0.45
1:A:657:LEU:HG	1:A:661:LYS:HE3	1.98	0.45
1:B:971:ALA:HB1	1:B:976:TYR:HB3	1.99	0.45
1:B:248:ILE:HD13	1:B:289:SER:H	1.82	0.45
1:B:902:ALA:HB1	1:B:914:SER:OG	2.17	0.45
1:A:823:PHE:CE2	4:A:1103:52O:H34	2.52	0.45
1:B:446:ARG:NH2	1:B:804:GLU:OE2	2.49	0.45
1:A:902:ALA:HB1	1:A:914:SER:OG	2.17	0.45
2:A:1101:B12:H261	3:A:1102[B]:5AD:H5'2	2.00	0.44
1:A:49:ILE:HB	1:A:150:ILE:HG13	1.99	0.44
1:A:742:GLN:HG3	2:A:1101:B12:H472	1.99	0.43
1:B:866:THR:O	1:B:901:ASN:ND2	2.49	0.43
1:A:443:PRO:HA	1:A:444:PRO:HD3	1.88	0.43
1:A:597:MET:HB2	1:A:614:VAL:HG11	2.01	0.43
1:A:626:ALA:O	1:A:652:VAL:HG12	2.18	0.43
1:A:971:ALA:HB1	1:A:976:TYR:HB3	2.00	0.43
1:A:598:PHE:HA	1:A:624:SER:HB3	2.00	0.43
1:A:120:VAL:HG11	1:A:1019:LEU:HD11	2.00	0.43
1:A:879:ASN:O	1:A:883:THR:OG1	2.30	0.43
1:A:265:ARG:NH1	5:A:1104:GDP:O1A	2.52	0.43
1:B:453:LEU:HD13	1:B:800:PHE:CD2	2.54	0.43
1:B:47:ARG:HB2	1:B:57:VAL:HG11	2.01	0.42
1:B:599:ALA:HB3	1:B:611:PHE:CE1	2.54	0.42
2:A:1101:B12:H362	2:A:1101:B12:H351	2.00	0.42
1:B:287:GLU:HG2	1:B:288:ILE:HG23	2.01	0.42
1:A:738:GLU:HB3	1:A:745:CYS:HB2	2.01	0.42
1:B:141:GLY:HA3	2:B:1101:B12:H5R2	2.02	0.42
1:A:866:THR:O	1:A:901:ASN:ND2	2.52	0.42
1:A:779:TYR:CE1	1:A:783:GLU:HG3	2.54	0.42
1:B:49:ILE:HB	1:B:150:ILE:HG13	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:47:ARG:HB2	1:A:57:VAL:HG11	2.02	0.42
2:A:1101:B12:H252	2:A:1101:B12:H601	2.02	0.41
1:B:358:LYS:HG2	5:B:1103:GDP:C6	2.55	0.41
1:A:358:LYS:HG2	5:A:1104:GDP:C6	2.55	0.41
1:A:778:GLY:HA3	1:A:795:THR:OG1	2.21	0.41
1:B:626:ALA:O	1:B:652:VAL:HG12	2.20	0.41
1:B:265:ARG:NH1	5:B:1103:GDP:O2A	2.53	0.41
1:B:999:ILE:N	1:B:1003:ASN:OD1	2.47	0.41
1:B:640:ARG:HA	1:B:641:PRO:HD3	1.91	0.41
1:A:251:SER:HB2	1:A:280:ALA:HB1	2.03	0.41
1:A:114:GLY:HA3	2:A:1101:B12:C9B	2.50	0.41
1:A:681:ASN:HB2	1:A:744:THR:HB	2.02	0.41
1:B:660:MET:HB2	1:B:660:MET:HE2	1.93	0.41
2:A:1101:B12:H8	4:A:1103:52O:H36	2.03	0.41
2:B:1101:B12:H91	2:B:1101:B12:H262	1.88	0.41
1:B:732:GLN:HA	1:B:775:SER:HB3	2.03	0.41
2:A:1101:B12:H531	2:A:1101:B12:H552	2.03	0.40
1:B:42:SER:OG	2:B:1101:B12:N3B	2.46	0.40
1:A:735:ILE:HA	1:A:738:GLU:HG2	2.02	0.40
1:B:398:ARG:NH1	1:B:440:VAL:HA	2.36	0.40
1:A:997:LEU:HD22	1:B:546:LEU:HD11	2.04	0.40
1:B:633:TYR:HD2	1:B:746:ILE:HD13	1.86	0.40
1:A:599:ALA:O	1:A:626:ALA:N	2.46	0.40
1:B:731:VAL:HG23	1:B:771:PHE:CE1	2.57	0.40
2:A:1101:B12:H262	2:A:1101:B12:H91	1.93	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	1057/1113 (95%)	1009 (96%)	47 (4%)	1 (0%)	51 82

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	1052/1113 (94%)	1016 (97%)	36 (3%)	0	100	100
All	All	2109/2226 (95%)	2025 (96%)	83 (4%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	260	LEU

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	821/906 (91%)	816 (99%)	5 (1%)	86	94
1	B	810/906 (89%)	807 (100%)	3 (0%)	91	95
All	All	1631/1812 (90%)	1623 (100%)	8 (0%)	88	94

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	249	ASP
1	A	489	HIS
1	A	744	THR
1	A	883	THR
1	A	1092	ASN
1	B	249	ASP
1	B	883	THR
1	B	974	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	267	ASN
1	B	899	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 12 ligands modelled in this entry, 4 are monoatomic - leaving 8 for Mogul analysis.

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
4	52O	A	1103	-	48,56,56	0.80	1 (2%)	61,85,85	0.67	1 (1%)
3	5AD	A	1102[A]	-	17,20,20	0.64	0	15,30,30	0.79	1 (6%)
4	52O	B	1102	-	26,29,56	0.95	1 (3%)	31,45,85	0.86	2 (6%)
2	B12	A	1101	1,3	80,101,101	0.67	2 (2%)	101,166,166	1.14	11 (10%)
5	GDP	B	1103	6	24,30,30	2.62	10 (41%)	31,47,47	1.66	6 (19%)
2	B12	B	1101	1	80,101,101	0.65	2 (2%)	101,166,166	1.08	7 (6%)
3	5AD	A	1102[B]	2	17,20,20	0.66	0	15,30,30	0.76	1 (6%)
5	GDP	A	1104	6	24,30,30	2.59	10 (41%)	31,47,47	1.66	6 (19%)

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
4	52O	A	1103	-	-	16/53/73/73	0/3/3/3
3	5AD	A	1102[A]	-	-	0/0/20/20	0/3/3/3
4	52O	B	1102	-	-	4/11/31/73	0/3/3/3
2	B12	A	1101	1,3	-	14/51/223/223	0/3/11/11
5	GDP	B	1103	6	-	0/12/32/32	0/3/3/3
2	B12	B	1101	1	-	13/51/223/223	0/3/11/11
3	5AD	A	1102[B]	2	-	0/0/20/20	0/3/3/3
5	GDP	A	1104	6	-	0/12/32/32	0/3/3/3

All (26) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	1103	GDP	C2-N2	7.21	1.48	1.33
5	A	1104	GDP	C2-N2	7.16	1.48	1.33
5	B	1103	GDP	O6-C6	5.00	1.37	1.24
5	A	1104	GDP	O6-C6	4.95	1.37	1.24
5	A	1104	GDP	C2-N1	-4.01	1.28	1.35
5	B	1103	GDP	C2-N1	-3.95	1.28	1.35
5	A	1104	GDP	O5'-C5'	-3.93	1.29	1.44
5	B	1103	GDP	O5'-C5'	-3.90	1.29	1.44
4	A	1103	52O	C23-C22	-3.57	1.50	1.53
5	B	1103	GDP	C2'-C1'	-3.37	1.48	1.53
2	B	1101	B12	C16-C15	3.29	1.53	1.41
5	A	1104	GDP	C2'-C1'	-3.24	1.48	1.53
2	A	1101	B12	C16-C15	2.91	1.52	1.41
4	B	1102	52O	P2-O12	2.91	1.59	1.50
5	B	1103	GDP	C5-C4	-2.90	1.33	1.40
5	A	1104	GDP	C5-C4	-2.90	1.33	1.40
5	B	1103	GDP	C2'-C3'	-2.72	1.45	1.53
5	B	1103	GDP	PA-O2A	-2.62	1.43	1.55
5	A	1104	GDP	C2'-C3'	-2.60	1.46	1.53
2	A	1101	B12	C1P-C2P	2.42	1.57	1.51
2	B	1101	B12	C1P-C2P	2.30	1.57	1.51
5	A	1104	GDP	PA-O1A	-2.24	1.43	1.50
5	B	1103	GDP	PB-O3B	-2.15	1.46	1.54
5	A	1104	GDP	PB-O3B	-2.12	1.46	1.54
5	B	1103	GDP	C6-C5	2.12	1.45	1.41
5	A	1104	GDP	C6-C5	2.05	1.44	1.41

All (35) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	1104	GDP	C2-N3-C4	4.27	120.23	115.36
5	B	1103	GDP	C2-N3-C4	4.19	120.15	115.36
5	A	1104	GDP	N3-C2-N1	-4.06	121.80	127.22
5	B	1103	GDP	N3-C2-N1	-4.02	121.87	127.22
2	A	1101	B12	C19-C1-N21	3.41	105.66	102.16
2	A	1101	B12	C30-C3-C2	-3.38	111.97	119.13
5	B	1103	GDP	PA-O3A-PB	-3.30	121.49	132.83
5	A	1104	GDP	PA-O3A-PB	-3.28	121.58	132.83
2	A	1101	B12	C55-C17-C16	3.18	120.51	109.92
2	B	1101	B12	C55-C17-C16	3.01	119.96	109.92
2	B	1101	B12	C30-C3-C2	-3.01	112.75	119.13
5	B	1103	GDP	C5-C6-N1	-2.83	119.56	123.43
5	A	1104	GDP	C5-C6-N1	-2.73	119.69	123.43
5	B	1103	GDP	C6-N1-C2	2.66	120.16	115.93
5	A	1104	GDP	C6-C5-C4	-2.66	118.26	120.80
5	A	1104	GDP	C6-N1-C2	2.64	120.13	115.93
5	B	1103	GDP	C6-C5-C4	-2.57	118.35	120.80
2	A	1101	B12	C13-C14-C15	-2.54	122.45	131.68
2	A	1101	B12	C3-C4-C5	-2.49	122.67	131.68
2	B	1101	B12	C13-C14-C15	-2.48	122.67	131.68
2	B	1101	B12	C16-C15-C14	-2.46	120.43	124.27
4	B	1102	52O	O7-P2-O4	2.45	113.25	106.73
2	A	1101	B12	C16-C15-C14	-2.40	120.53	124.27
2	B	1101	B12	C3-C4-C5	-2.40	122.99	131.68
2	A	1101	B12	C20-C1-C19	-2.33	107.11	109.36
4	B	1102	52O	C8-C12-N5	2.30	123.85	120.35
2	B	1101	B12	C55-C17-C18	-2.30	106.70	111.14
3	A	1102[A]	5AD	C5-C6-N6	2.28	123.82	120.35
4	A	1103	52O	C8-C12-N5	2.27	123.80	120.35
3	A	1102[B]	5AD	C5-C6-N6	2.23	123.74	120.35
2	A	1101	B12	C15-C14-N23	2.20	128.36	124.64
2	B	1101	B12	C20-C1-C19	-2.17	107.26	109.36
2	A	1101	B12	C8-C9-N22	2.16	113.82	111.12
2	A	1101	B12	C2-C1-C19	-2.08	115.31	118.60
2	A	1101	B12	C5-C6-N22	-2.08	121.01	124.81

There are no chirality outliers.

All (47) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	1103	52O	C5-O4-P2-O10
4	A	1103	52O	O8-C11-C9-C10
4	A	1103	52O	C11-O8-P3-O14

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Mol	Chain	Res	Type	Atoms
4	B	1102	52O	C5-O4-P2-O7
4	B	1102	52O	C5-O4-P2-O12
4	B	1102	52O	C5-O4-P2-O10
4	B	1102	52O	C1-O2-P1-O6
2	A	1101	B12	C1P-C2P-O3-P
2	A	1101	B12	C3P-C2P-O3-P
2	B	1101	B12	C2-C3-C30-C31
2	B	1101	B12	C18-C60-C61-N62
4	A	1103	52O	C2-C1-O2-P1
2	A	1101	B12	C2-C3-C30-C31
4	A	1103	52O	O8-C11-C9-C13
4	A	1103	52O	O8-C11-C9-C14
2	B	1101	B12	C12-C13-C48-C49
2	B	1101	B12	C18-C60-C61-O63
2	B	1101	B12	C3R-O2-P-O3
2	B	1101	B12	C41-C42-C43-O44
2	A	1101	B12	C18-C60-C61-O63
2	A	1101	B12	C18-C60-C61-N62
2	B	1101	B12	C41-C42-C43-N45
2	A	1101	B12	C55-C56-C57-O58
2	B	1101	B12	C38-C37-C7-C36
4	A	1103	52O	C5-O4-P2-O7
4	A	1103	52O	C11-O8-P3-O7
4	A	1103	52O	S1-C22-C23-C24
4	A	1103	52O	C5-O4-P2-O12
4	A	1103	52O	C11-O8-P3-O13
2	A	1101	B12	C42-C41-C8-C9
2	B	1101	B12	C14-C13-C48-C49
2	A	1101	B12	C30-C31-C32-O34
2	A	1101	B12	C55-C56-C57-N59
2	A	1101	B12	C30-C31-C32-N33
2	A	1101	B12	C3R-O2-P-O3
2	A	1101	B12	C12-C13-C48-C49
2	B	1101	B12	C38-C37-C7-C8
4	A	1103	52O	C20-C21-S1-C22
2	A	1101	B12	N59-C1P-C2P-O3
2	B	1101	B12	C55-C56-C57-O58
4	A	1103	52O	C1-O2-P1-O9
4	A	1103	52O	S1-C22-C23-C25
4	A	1103	52O	O17-C22-C23-C24
4	A	1103	52O	P3-O7-P2-O10
2	B	1101	B12	C55-C56-C57-N59

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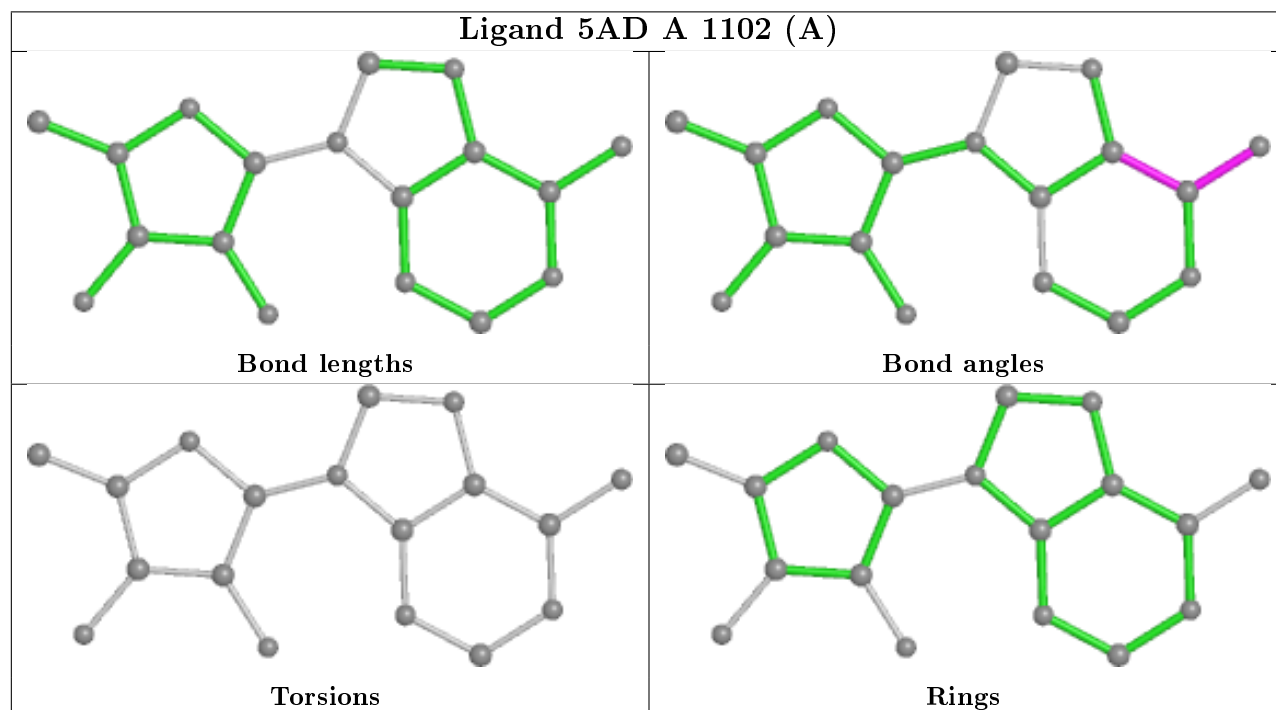
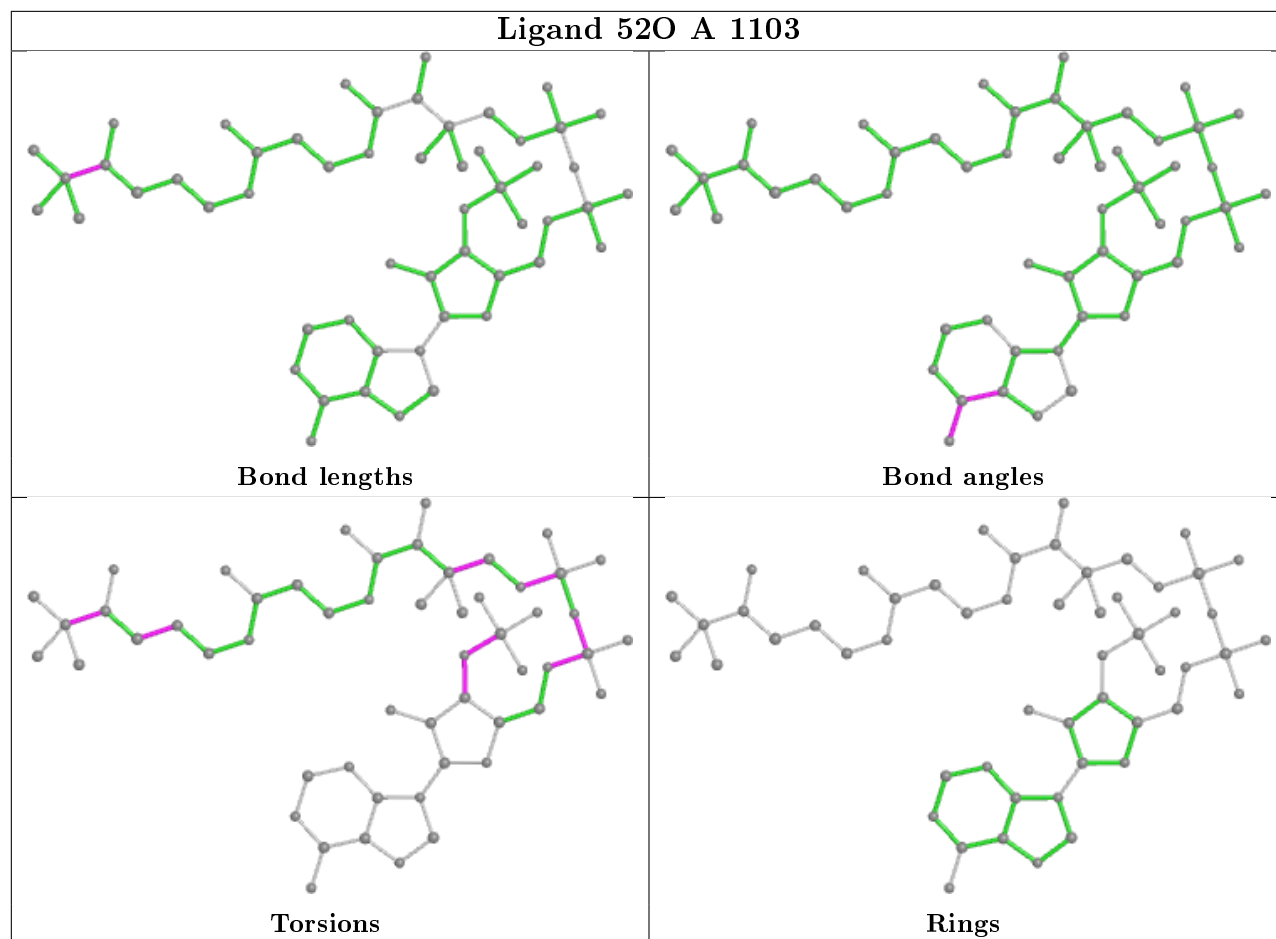
Mol	Chain	Res	Type	Atoms
2	A	1101	B12	C4-C3-C30-C31
2	B	1101	B12	C4-C3-C30-C31

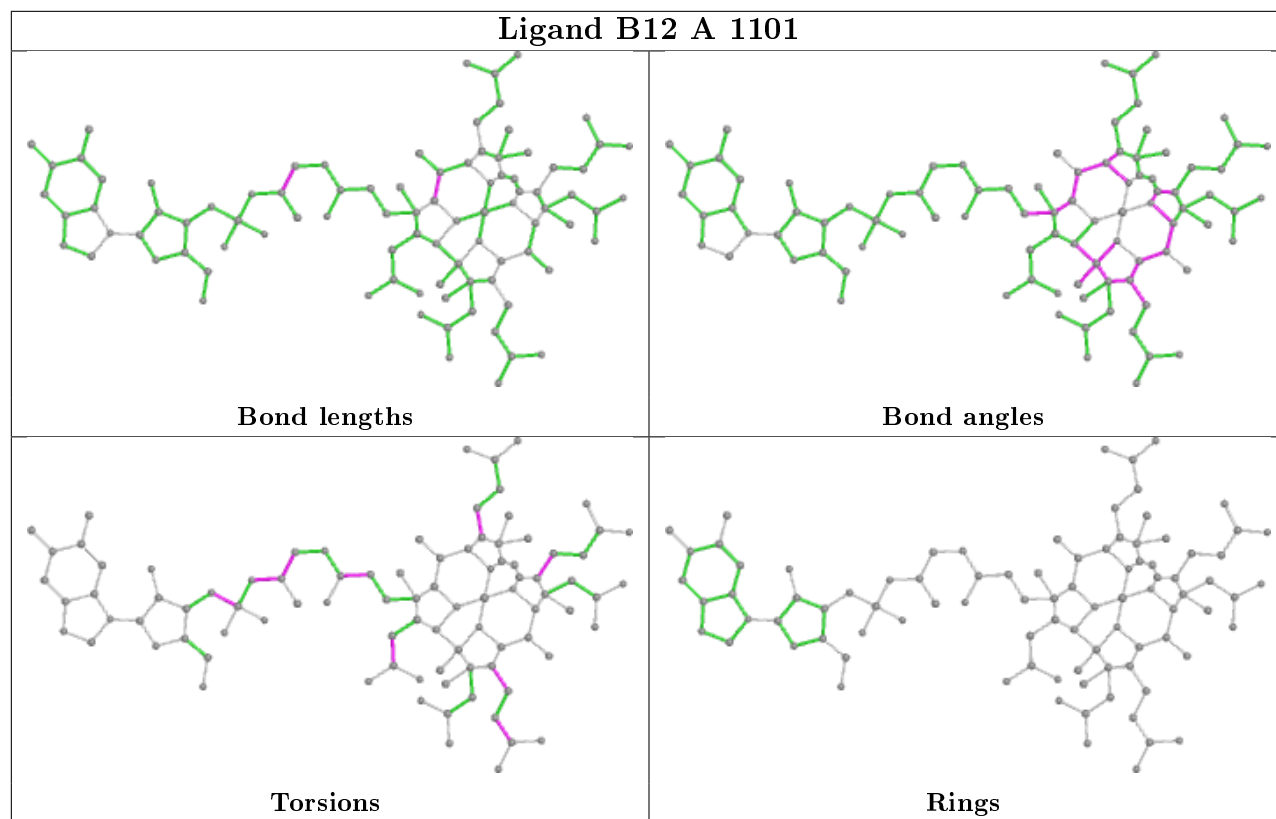
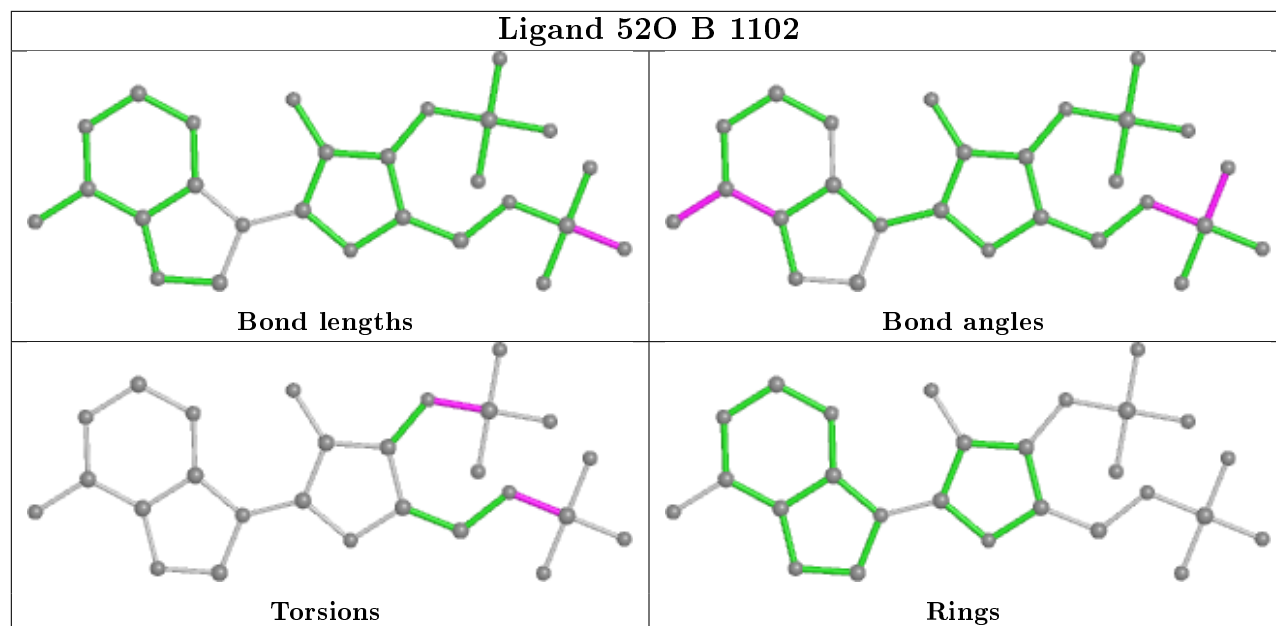
There are no ring outliers.

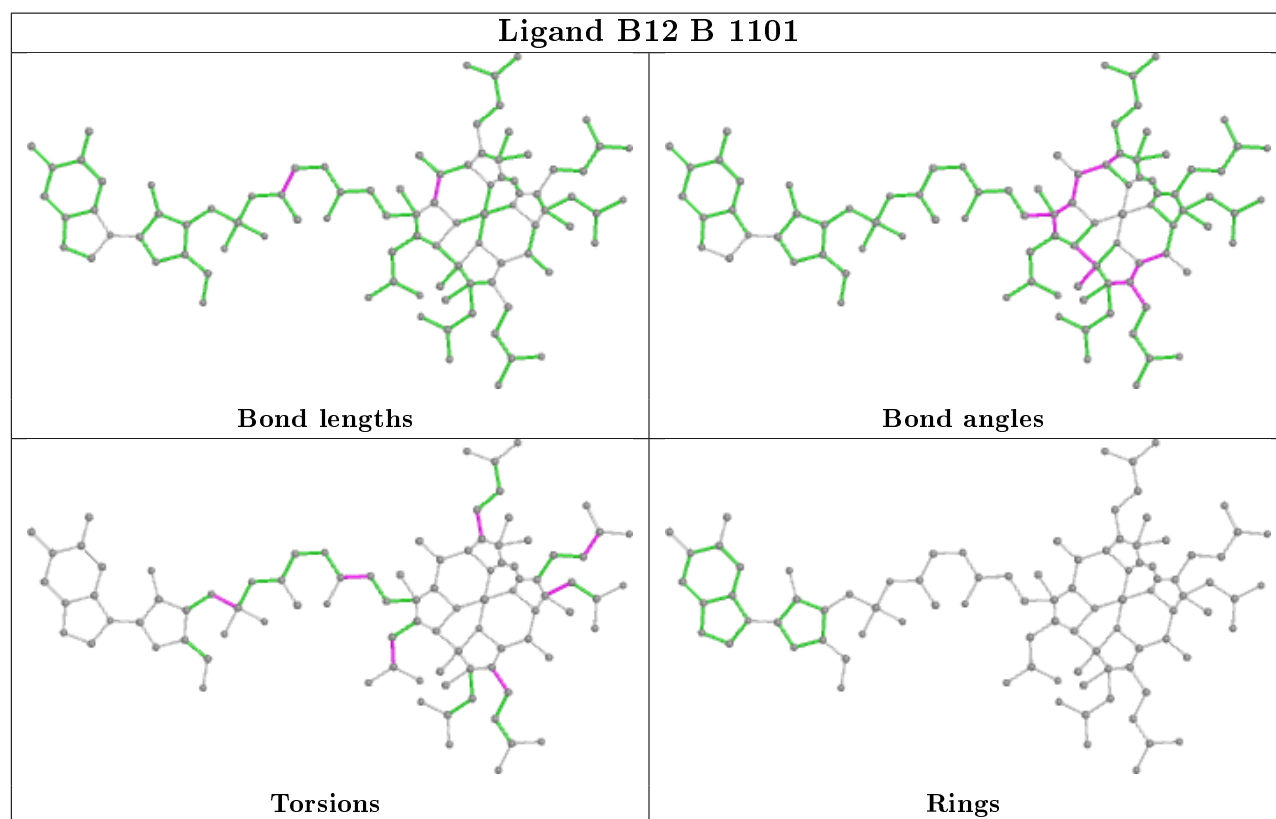
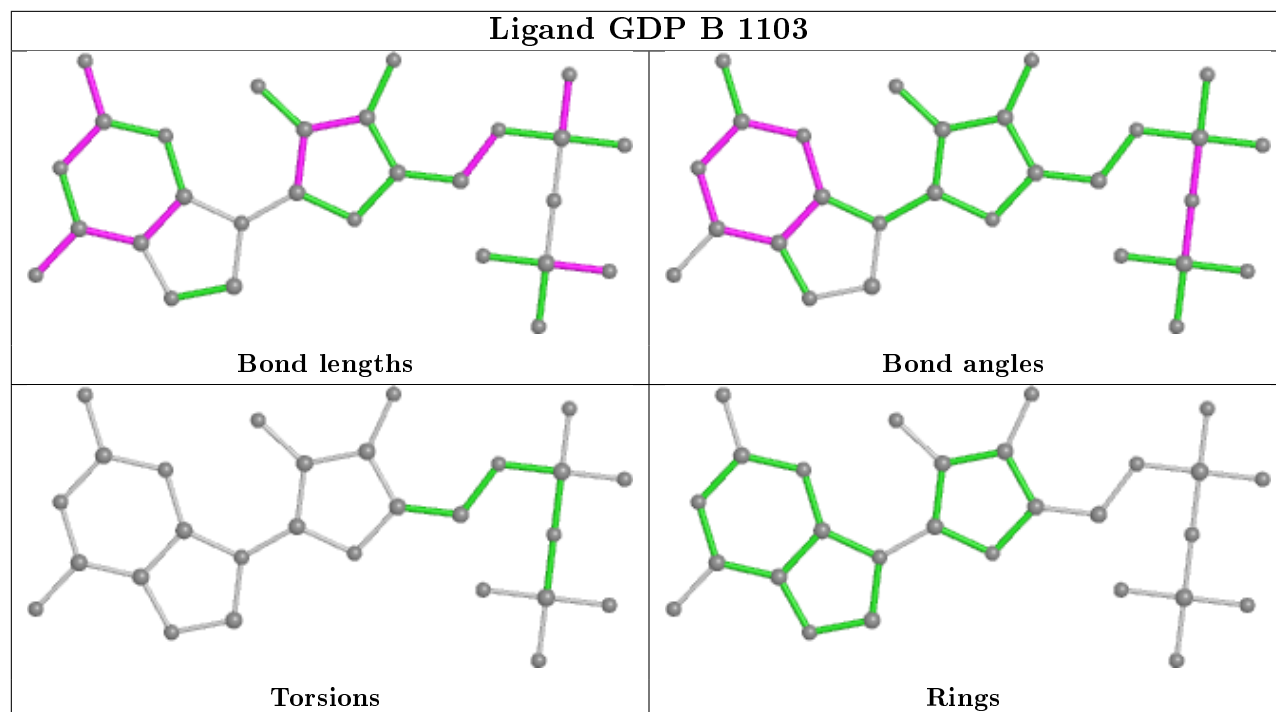
7 monomers are involved in 31 short contacts:

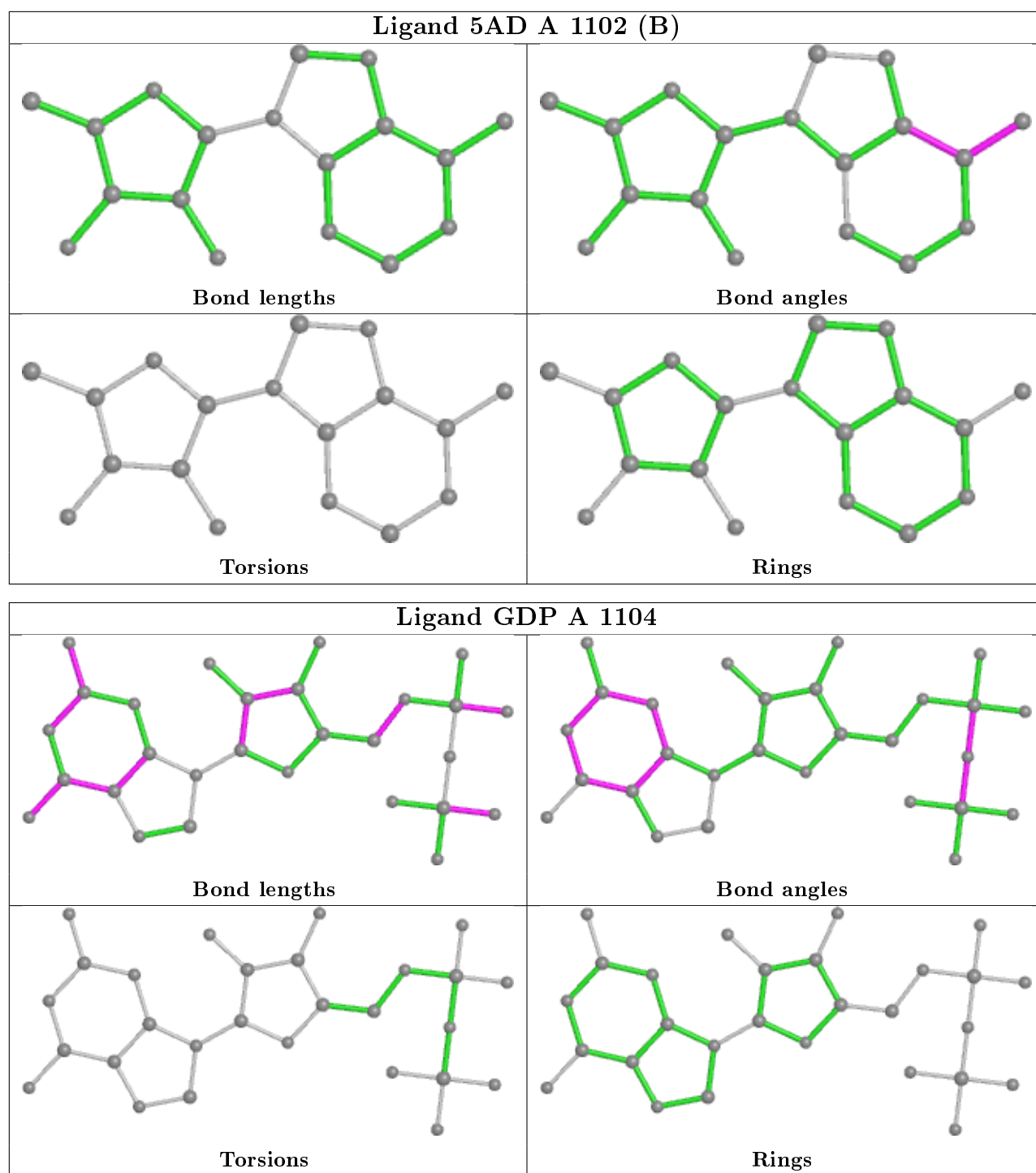
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1103	52O	7	0
3	A	1102[A]	5AD	2	0
2	A	1101	B12	14	0
5	B	1103	GDP	2	0
2	B	1101	B12	6	0
3	A	1102[B]	5AD	2	0
5	A	1104	GDP	2	0

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.









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	1063/1113 (95%)	0.48	42 (3%) 38 37	75, 108, 157, 214	0
1	B	1060/1113 (95%)	0.32	10 (0%) 84 83	71, 100, 157, 234	0
All	All	2123/2226 (95%)	0.40	52 (2%) 59 57	71, 104, 157, 234	0

All (52) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	420	MET	3.9
1	A	494	LEU	3.5
1	B	494	LEU	3.5
1	A	200	LEU	3.3
1	A	295	VAL	3.3
1	A	415	LEU	2.9
1	A	1011	ASP	2.8
1	A	98	ILE	2.8
1	A	146	LEU	2.7
1	A	94	PHE	2.7
1	A	316	GLN	2.7
1	B	680	ILE	2.6
1	A	127	LEU	2.6
1	A	808	ALA	2.6
1	A	660	MET	2.5
1	A	31	THR	2.5
1	A	321	ILE	2.5
1	B	826	ASN	2.5
1	A	654	ILE	2.4
1	A	299	CYS	2.4
1	A	217	THR	2.4
1	B	909	THR	2.4
1	A	489	HIS	2.4
1	B	299	CYS	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	214	ILE	2.4
1	A	749	THR	2.3
1	A	161	LEU	2.3
1	A	69	VAL	2.3
1	A	243	ILE	2.3
1	A	317	GLY	2.3
1	B	900	THR	2.3
1	A	629	SER	2.2
1	A	356	ILE	2.2
1	A	164	TYR	2.2
1	A	799	GLY	2.2
1	A	304	PHE	2.2
1	B	904	ASP	2.2
1	B	901	ASN	2.2
1	A	1052	LEU	2.2
1	A	411	LEU	2.1
1	A	328	SER	2.1
1	A	637	PRO	2.1
1	A	566	LEU	2.1
1	B	489	HIS	2.1
1	A	29	PHE	2.1
1	A	643	ILE	2.1
1	A	570	MET	2.1
1	A	212	LEU	2.0
1	A	670	THR	2.0
1	A	245	VAL	2.0
1	A	308	ILE	2.0
1	B	907	ILE	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 carbohydrates 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,

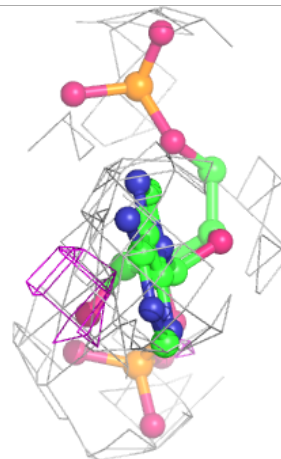
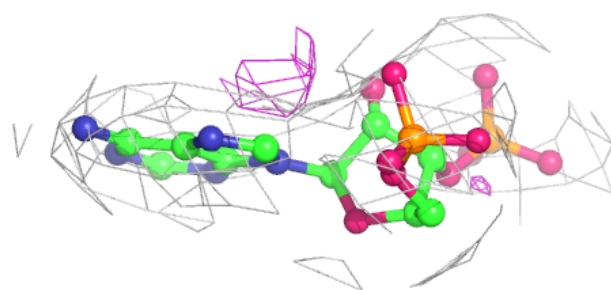
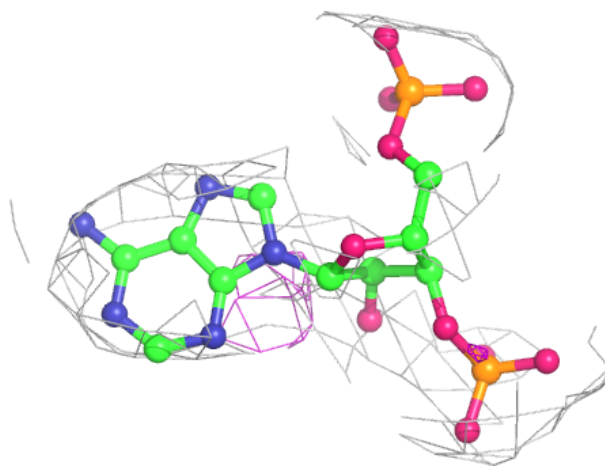
median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	52O	B	1102	27/54	0.81	0.18	131,172,224,225	0
3	5AD	A	1102[A]	18/18	0.91	0.82	171,173,177,178	18
3	5AD	A	1102[B]	18/18	0.91	0.82	170,174,177,179	18
2	B12	B	1101	91/91	0.92	0.30	106,130,151,162	0
4	52O	A	1103	54/54	0.92	0.26	75,113,160,165	0
6	MG	A	1106	1/1	0.94	0.10	107,107,107,107	0
6	MG	A	1105	1/1	0.95	0.15	78,78,78,78	0
6	MG	B	1105	1/1	0.96	0.22	70,70,70,70	0
2	B12	A	1101	91/91	0.96	0.35	65,83,104,110	0
5	GDP	B	1103	28/28	0.97	0.23	72,81,92,133	0
5	GDP	A	1104	28/28	0.97	0.17	95,113,118,122	0
6	MG	B	1104	1/1	0.99	0.23	70,70,70,70	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.

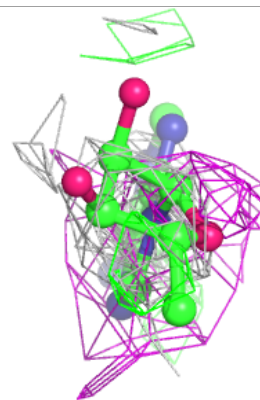
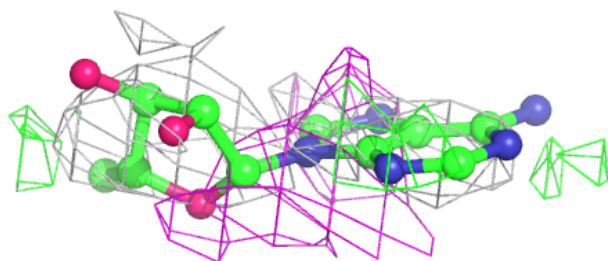
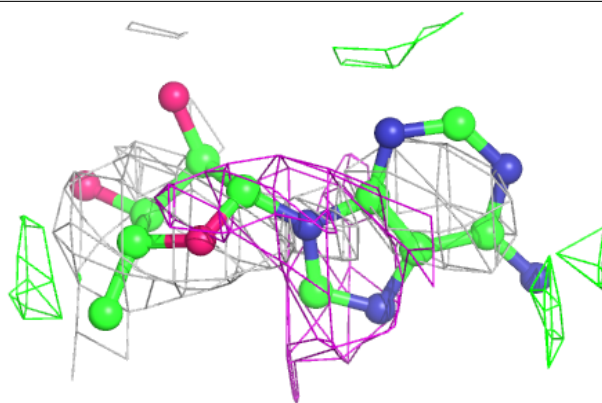
Electron density around 52O B 1102:

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

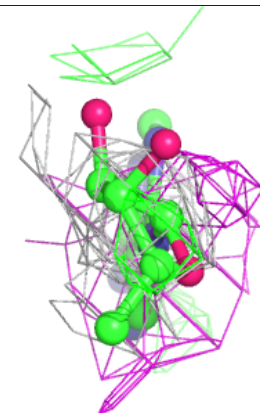
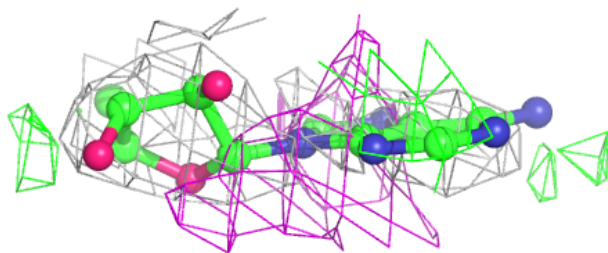
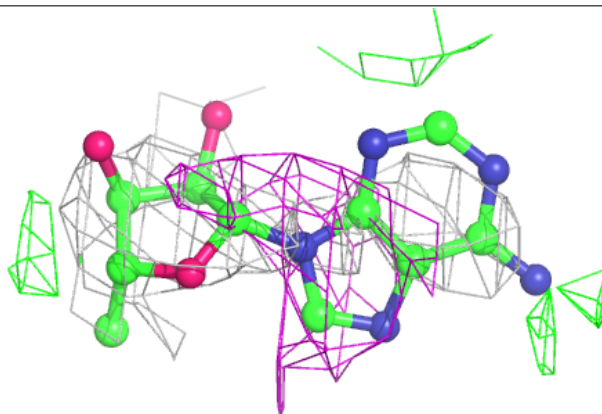


Electron density around 5AD A 1102 (A):

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

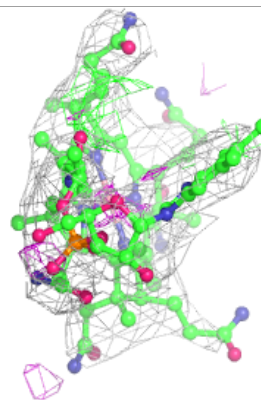
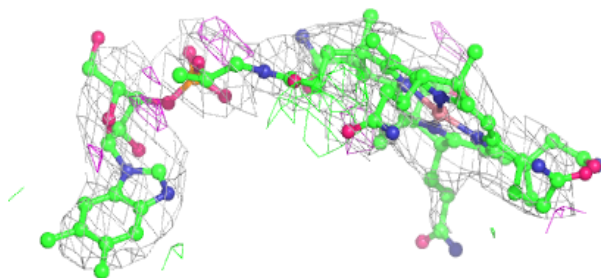
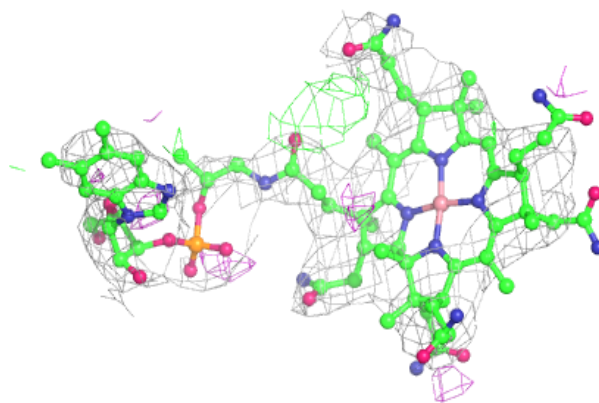
**Electron density around 5AD A 1102 (B):**

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

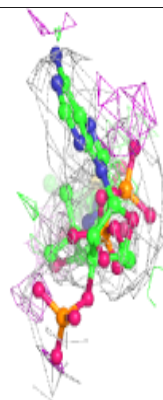
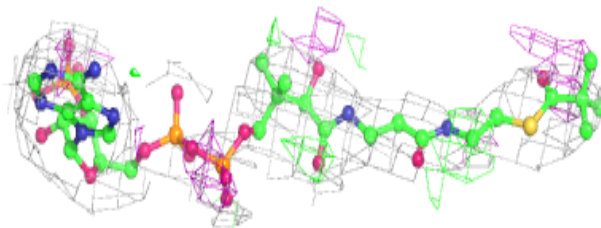
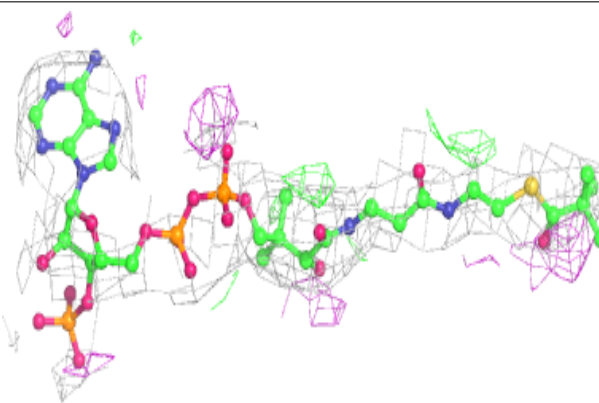


Electron density around B12 B 1101:

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

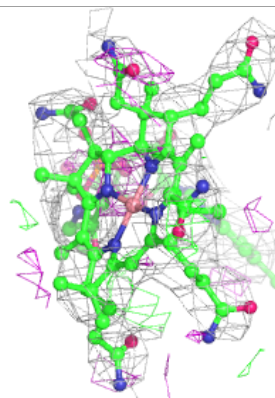
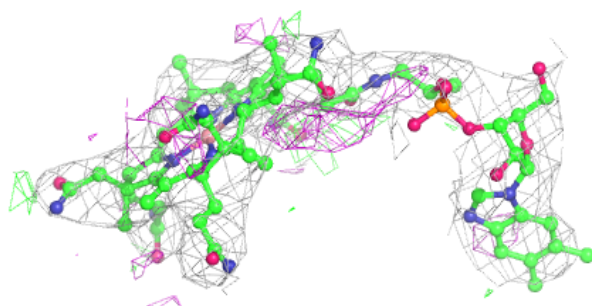
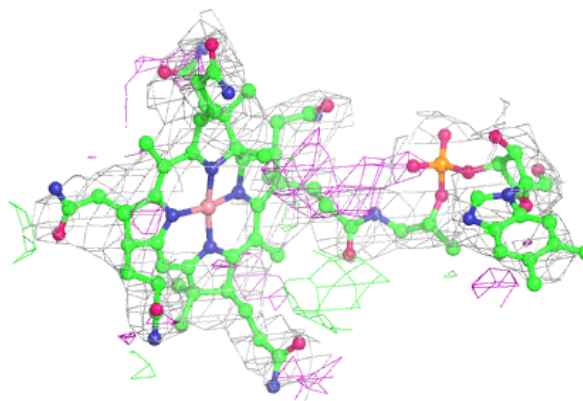
**Electron density around 52O A 1103:**

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

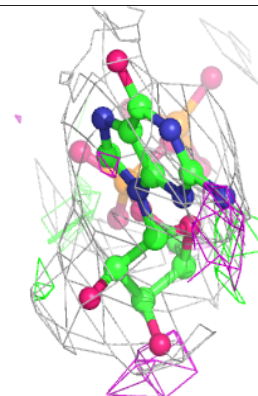
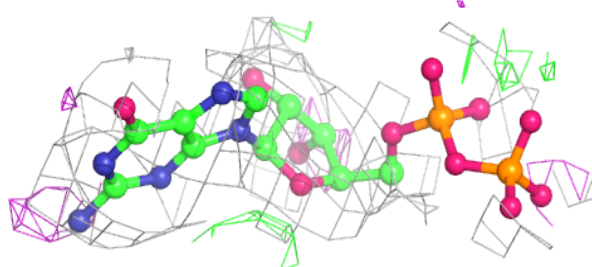
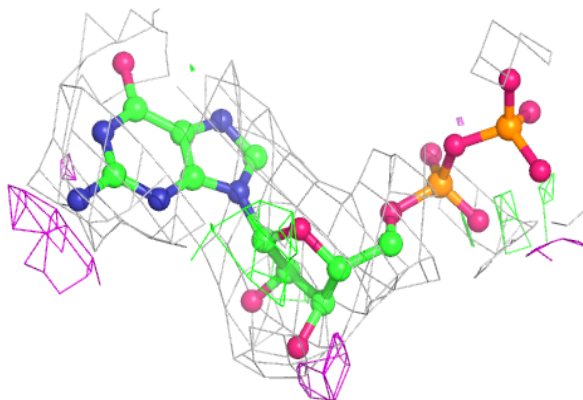


Electron density around B12 A 1101:

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

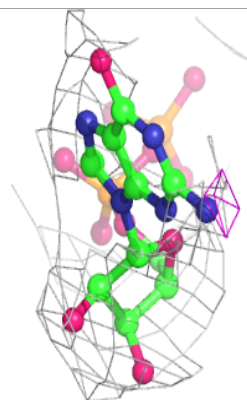
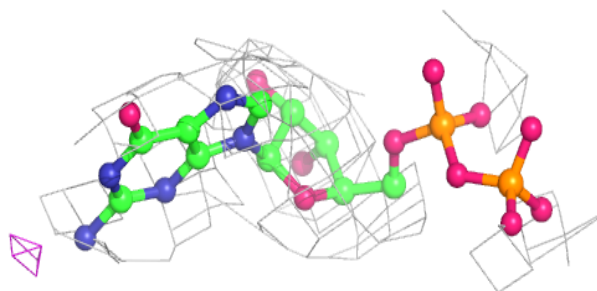
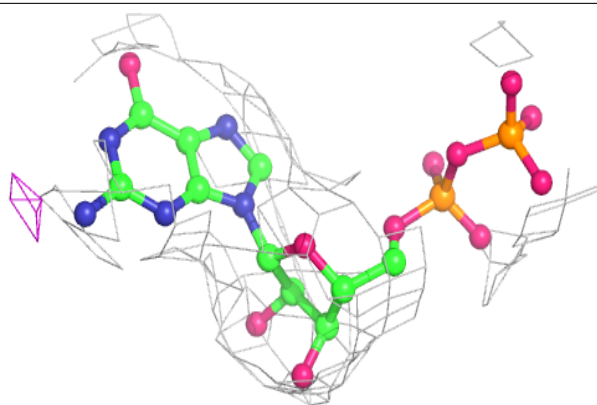
**Electron density around GDP B 1103:**

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



Electron density around GDP A 1104:

$2mF_o - DF_c$ (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.