



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 17, 2023 – 09:56 AM EDT

PDB ID : 4XC8
Title : Isobutyryl-CoA mutase fused with bound butyryl-CoA, GDP, and Mg and without cobalamin (apo-IcmF/GDP)
Authors : Jost, M.; Drennan, C.L.
Deposited on : 2014-12-17
Resolution : 3.25 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 : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtrriage (Phenix) : 1.13
EDS : 2.35.1
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) : 2.35.1

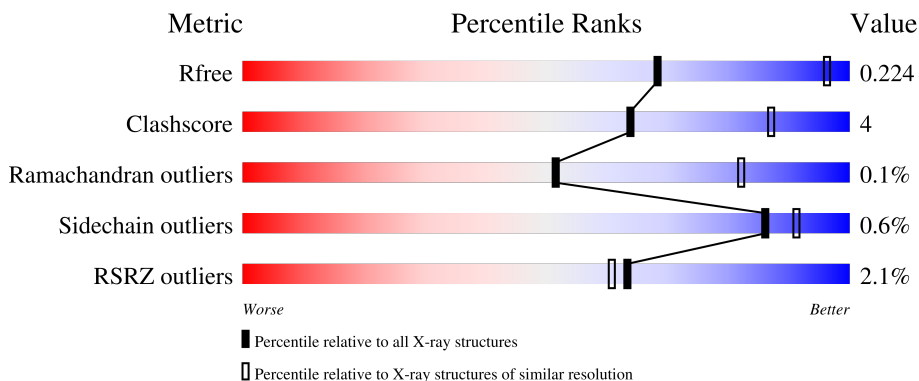
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

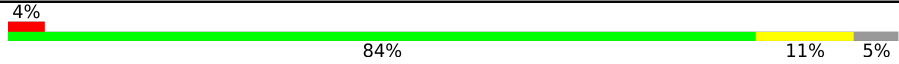

The reported resolution of this entry is 3.25 Å.

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	1191 (3.30-3.22)
Clashscore	141614	1251 (3.30-3.22)
Ramachandran outliers	138981	1229 (3.30-3.22)
Sidechain outliers	138945	1228 (3.30-3.22)
RSRZ outliers	127900	1154 (3.30-3.22)

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	1113	 4% 84% 11% 5%
1	B	1113	 84% 11% 5%

2 Entry composition i

There are 4 unique types of molecules in this entry. The entry contains 16245 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
			Total	C	N	O	S			
1	A	1058	Total	C	N	O	S	0	0	0
			7995	5004	1418	1537	36			
1	B	1060	Total	C	N	O	S	0	0	0
			8098	5058	1446	1558	36			

There are 40 discrepancies between the modelled and reference sequences:

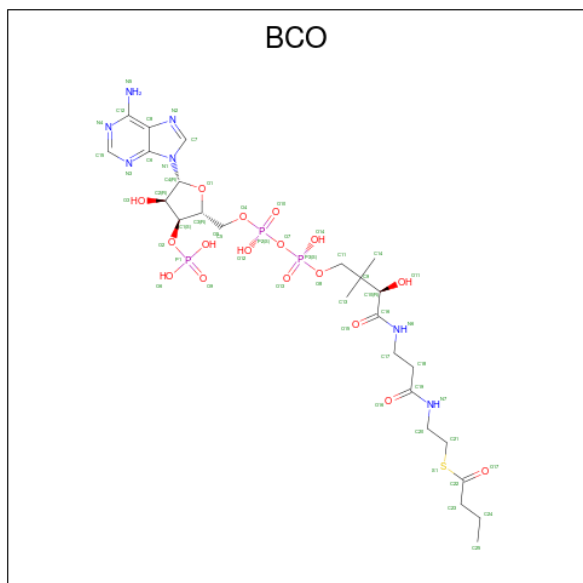
Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	initiating methionine	UNP Q1LRY0
A	-18	GLY	-	expression tag	UNP Q1LRY0
A	-17	SER	-	expression tag	UNP Q1LRY0
A	-16	SER	-	expression tag	UNP Q1LRY0
A	-15	HIS	-	expression tag	UNP Q1LRY0
A	-14	HIS	-	expression tag	UNP Q1LRY0
A	-13	HIS	-	expression 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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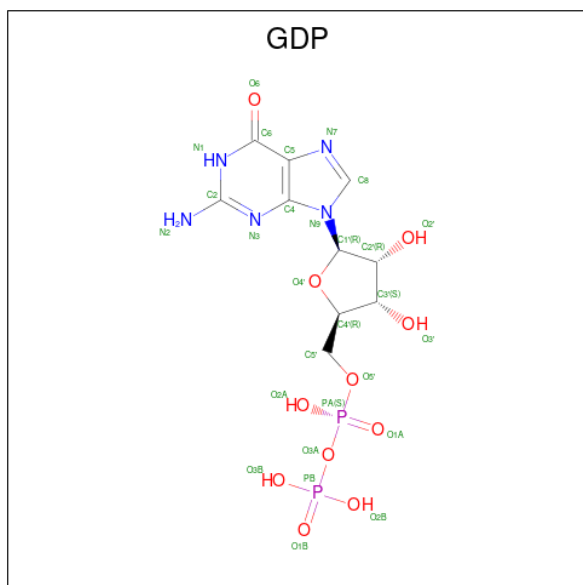
Chain	Residue	Modelled	Actual	Comment	Reference
B	-14	HIS	-	expression tag	UNP Q1LRY0
B	-13	HIS	-	expression tag	UNP Q1LRY0
B	-12	HIS	-	expression tag	UNP Q1LRY0
B	-11	HIS	-	expression tag	UNP Q1LRY0
B	-10	HIS	-	expression tag	UNP Q1LRY0
B	-9	SER	-	expression tag	UNP Q1LRY0
B	-8	SER	-	expression tag	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 Butyryl Coenzyme A (three-letter code: BCO) (formula: $C_{25}H_{42}N_7O_{17}P_3S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	
			Total	C	N	O	P			S
2	A	1	53	25	7	17	3	1	0	0
2	B	1	39	16	5	15	3		0	0

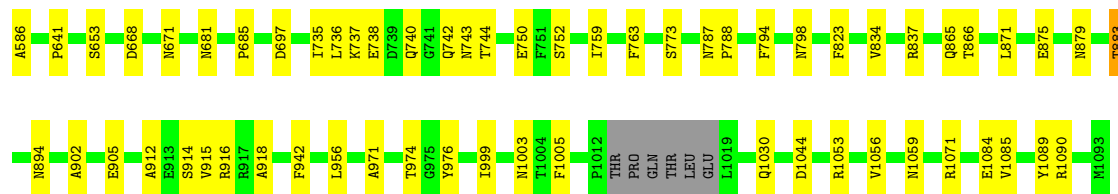
- Molecule 3 is GUANOSINE-5'-DIPHOSPHATE (three-letter code: GDP) (formula: $C_{10}H_{15}N_5O_{11}P_2$).



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

- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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



4 Data and refinement statistics

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, α , β , γ	316.51Å 316.51Å 342.27Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	99.16 – 3.25 99.16 – 3.25	Depositor EDS
% Data completeness (in resolution range)	99.3 (99.16-3.25) 99.3 (99.16-3.25)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.15	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.84 (at 3.26Å)	Xtrriage
Refinement program	PHENIX (phenix.refine: dev_1760)	Depositor
R, R_{free}	0.205 , 0.222 0.210 , 0.224	Depositor DCC
R_{free} test set	5148 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	84.0	Xtrriage
Anisotropy	0.483	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 67.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	16245	wwPDB-VP
Average B, all atoms (Å ²)	94.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.16% 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, BCO

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	0/8137	0.40	3/11034 (0.0%)
1	B	0.21	0/8239	0.40	3/11158 (0.0%)
All	All	0.21	0/16376	0.40	6/22192 (0.0%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	552	ARG	NE-CZ-NH2	-11.21	114.70	120.30
1	A	552	ARG	NE-CZ-NH1	-11.05	114.77	120.30
1	B	552	ARG	NE-CZ-NH1	11.05	125.83	120.30
1	A	552	ARG	NE-CZ-NH2	10.85	125.72	120.30
1	A	552	ARG	CD-NE-CZ	5.49	131.28	123.60
1	B	552	ARG	CD-NE-CZ	5.47	131.26	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	7995	0	7674	68	0
1	B	8098	0	7861	70	0
2	A	53	0	40	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	39	0	23	0	0
3	A	28	0	12	2	0
3	B	28	0	12	1	0
4	A	2	0	0	0	0
4	B	2	0	0	0	0
All	All	16245	0	15622	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:1092:ASN:OD1	3:A:1102:GDP:O3'	2.08	0.70
1:A:916:ARG:NH1	1:B:875:GLU:OE1	2.23	0.69
1:B:412:VAL:HG13	1:B:422:LEU:HD11	1.77	0.66
1:A:908:THR:OG1	1:A:909:THR:N	2.29	0.64
1:A:265:ARG:NH2	3:A:1102:GDP:O2A	2.34	0.60
1:A:735:ILE:HG13	1:A:752:SER:HB3	1.84	0.60
1:A:1082:LEU:O	1:A:1086:GLY:N	2.34	0.60
1:B:454:ALA:HA	1:B:956:LEU:HD22	1.84	0.60
1:B:265:ARG:NH2	3:B:1102:GDP:O2A	2.35	0.59
1:B:902:ALA:HB3	1:B:905:GLU:HG3	1.84	0.59
1:A:250:PRO:HD3	1:A:312:SER:HB2	1.85	0.59
1:B:735:ILE:HG13	1:B:752:SER:HB3	1.84	0.59
1:A:47:ARG:HH11	1:A:259:LEU:HD11	1.68	0.59
1:A:365:GLN:HG3	1:A:369:ARG:HH22	1.69	0.58
1:A:31:THR:HG22	1:A:82:ALA:HB3	1.86	0.57
1:A:91:VAL:HA	1:A:127:LEU:HD11	1.86	0.57
1:B:250:PRO:HD3	1:B:312:SER:HB2	1.87	0.57
1:B:91:VAL:HA	1:B:127:LEU:HD11	1.87	0.57
1:B:31:THR:HG22	1:B:82:ALA:HB3	1.86	0.56
1:A:640:ARG:NH1	1:A:642:ASP:OD2	2.39	0.55
1:A:529:VAL:HG12	1:A:538:ARG:HG2	1.87	0.55
1:A:35:LEU:HD22	1:A:642:ASP:HB3	1.88	0.55
1:B:685:PRO:HB3	1:B:759:ILE:HD11	1.89	0.55
1:A:356:ILE:HD12	1:A:368:TRP:HA	1.89	0.54
1:A:731:VAL:HG12	1:A:771:PHE:HE1	1.71	0.54
1:B:1053:ARG:HG2	1:B:1085:VAL:HB	1.90	0.54
1:A:30:VAL:HG23	1:A:78:VAL:HG11	1.90	0.54
1:B:572:GLU:HG2	1:B:581:THR:HG21	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:685:PRO:HB3	1:A:759:ILE:HD11	1.91	0.53
1:A:879:ASN:O	1:A:883:THR:OG1	2.28	0.52
1:B:742:GLN:HB3	1:B:744:THR:OG1	2.09	0.52
1:A:572:GLU:HG2	1:A:581:THR:HG21	1.92	0.52
1:B:200:LEU:HD11	1:B:298:ALA:HA	1.91	0.52
1:B:30:VAL:HG23	1:B:78:VAL:HG11	1.90	0.52
1:A:732:GLN:HA	1:A:775:SER:HB3	1.92	0.51
1:A:672:PRO:O	1:A:728:ARG:NH2	2.41	0.51
1:B:794:PHE:O	1:B:798:ASN:ND2	2.43	0.51
1:B:879:ASN:O	1:B:883:THR:OG1	2.28	0.51
1:A:454:ALA:HA	1:A:956:LEU:HD22	1.93	0.50
1:B:96:TYR:HB2	1:B:641:PRO:HB2	1.93	0.50
1:B:231:ARG:NH1	1:B:1084:GLU:HA	2.26	0.50
1:B:971:ALA:HB1	1:B:976:TYR:HB3	1.94	0.49
1:A:561:ASP:HB2	1:B:561:ASP:HB2	1.94	0.49
1:A:916:ARG:NH2	1:B:1005:PHE:O	2.46	0.49
1:B:228:GLU:OE1	1:B:232:ARG:NH1	2.46	0.49
1:B:248:ILE:HD13	1:B:289:SER:H	1.77	0.49
1:A:794:PHE:O	1:A:798:ASN:ND2	2.45	0.49
1:A:787:ASN:HB2	1:A:788:PRO:HD2	1.95	0.48
1:A:861:LYS:NZ	2:A:1101:BCO:O6	2.39	0.48
1:A:228:GLU:OE1	1:A:232:ARG:NH1	2.46	0.48
1:B:787:ASN:HB2	1:B:788:PRO:HD2	1.95	0.47
1:A:776:ILE:HD13	1:A:799:GLY:HA2	1.96	0.47
1:A:942:PHE:CZ	1:B:557:PRO:HB3	2.50	0.47
1:A:971:ALA:HB1	1:A:976:TYR:HB3	1.95	0.47
1:A:779:TYR:CD2	1:A:825:SER:HB2	2.50	0.47
1:A:485:GLN:HA	1:A:490:GLU:HG2	1.98	0.46
1:A:542:ILE:HD13	1:A:550:LYS:HD3	1.97	0.46
1:A:598:PHE:HB2	1:A:899:HIS:NE2	2.31	0.46
1:B:740:GLN:NE2	1:B:798:ASN:OD1	2.46	0.45
1:A:640:ARG:HA	1:A:641:PRO:HD3	1.86	0.45
1:A:777:SER:HA	1:A:823:PHE:HB3	1.98	0.45
1:A:599:ALA:HB3	1:A:611:PHE:CE1	2.52	0.44
1:B:267:ASN:ND2	1:B:1089:TYR:O	2.50	0.44
1:B:287:GLU:HG2	1:B:288:ILE:HG23	1.98	0.44
1:A:589:ARG:NE	1:A:592:GLU:HB2	2.33	0.44
1:B:1056:VAL:HB	1:B:1085:VAL:HG21	2.00	0.44
1:A:598:PHE:H	1:A:899:HIS:CD2	2.36	0.44
1:B:528:TYR:HB2	1:B:541:LEU:HD21	1.99	0.44
1:A:697:ASP:OD1	1:A:1071:ARG:NH1	2.51	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:60:LEU:HD21	1:B:263:ARG:HD3	1.99	0.44
1:B:271:HIS:CG	1:B:272:PRO:HD2	2.53	0.44
1:A:441:ILE:HG22	1:A:442:VAL:HG23	2.00	0.43
1:B:443:PRO:HA	1:B:444:PRO:HD3	1.85	0.43
1:B:26:LYS:HA	1:B:26:LYS:HD2	1.62	0.43
1:A:528:TYR:HB2	1:A:541:LEU:HD21	1.99	0.43
1:B:653:SER:O	1:B:1030:GLN:NE2	2.51	0.43
1:A:681:ASN:HA	1:A:733:ALA:HA	2.00	0.43
1:B:209:VAL:HA	1:B:210:PRO:HD3	1.87	0.43
1:A:871:LEU:HD22	1:A:879:ASN:HB3	2.01	0.43
1:A:210:PRO:HG2	1:A:420:MET:HB2	2.01	0.43
1:A:866:THR:HG21	1:A:883:THR:HG22	2.01	0.43
1:B:327:LEU:HD21	1:B:418:ARG:NH2	2.34	0.43
1:B:398:ARG:NH2	1:B:440:VAL:HA	2.34	0.43
1:B:668:ASP:HB3	1:B:671:ASN:HB2	2.01	0.43
1:A:327:LEU:HD21	1:A:418:ARG:NH2	2.34	0.43
1:A:653:SER:O	1:A:1030:GLN:NE2	2.52	0.43
1:A:45:ILE:O	1:A:49:ILE:HG12	2.18	0.42
1:A:271:HIS:CG	1:A:272:PRO:HD2	2.54	0.42
1:A:834:VAL:HG12	1:A:837:ARG:HD3	2.01	0.42
1:B:697:ASP:OD1	1:B:1071:ARG:NH1	2.52	0.42
1:B:871:LEU:HD22	1:B:879:ASN:HB3	2.01	0.42
1:B:210:PRO:HG2	1:B:420:MET:HB2	2.01	0.42
1:B:738:GLU:O	1:B:743:ASN:N	2.51	0.42
1:B:866:THR:HG21	1:B:883:THR:HG22	2.01	0.42
1:A:757:GLY:HA3	1:A:809:ARG:NH1	2.35	0.42
1:A:1044:ASP:OD1	1:A:1044:ASP:N	2.53	0.42
1:B:763:PHE:CE2	1:B:773:SER:HB3	2.55	0.42
1:B:912:ALA:O	1:B:916:ARG:HG3	2.19	0.42
1:A:265:ARG:O	1:A:1090:ARG:NH1	2.53	0.42
1:A:902:ALA:HB1	1:A:914:SER:OG	2.20	0.42
1:B:902:ALA:HB1	1:B:914:SER:OG	2.19	0.42
1:B:28:ARG:HG2	1:B:56:GLU:HB2	2.02	0.42
1:B:736:LEU:HB3	1:B:798:ASN:OD1	2.20	0.42
1:A:58:ILE:HG13	1:A:260:LEU:HD12	2.02	0.41
1:B:44:ASN:HA	1:B:47:ARG:HG2	2.01	0.41
1:A:200:LEU:HD11	1:A:298:ALA:HA	2.02	0.41
1:A:1059:ASN:O	1:A:1059:ASN:ND2	2.54	0.41
1:B:834:VAL:HG12	1:B:837:ARG:HD3	2.02	0.41
1:B:184:ILE:HG23	1:B:295:VAL:HG22	2.01	0.41
1:A:575:PRO:O	1:A:586:ALA:HA	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:681:ASN:HD21	1:B:737:LYS:HE2	1.85	0.41
1:B:32:ALA:HB2	1:B:60:LEU:HB2	2.02	0.41
1:B:41:ALA:O	1:B:45:ILE:HG12	2.21	0.41
1:B:823:PHE:HZ	1:B:865:GLN:HE21	1.69	0.41
1:B:912:ALA:O	1:B:915:VAL:HG12	2.21	0.41
1:A:291:ALA:O	1:A:295:VAL:HG23	2.20	0.41
1:A:505:LEU:O	1:A:510:ARG:NH1	2.54	0.41
1:B:267:ASN:HD21	1:B:1090:ARG:HA	1.85	0.41
1:B:505:LEU:O	1:B:510:ARG:NH1	2.54	0.41
1:B:1044:ASP:N	1:B:1044:ASP:OD1	2.54	0.41
1:A:598:PHE:HD1	1:A:598:PHE:HA	1.72	0.41
1:B:172:VAL:HG21	1:B:183:LEU:HD22	2.03	0.41
1:B:398:ARG:HD3	1:B:750:GLU:OE2	2.20	0.41
1:B:999:ILE:N	1:B:1003:ASN:OD1	2.52	0.41
1:B:58:ILE:HG13	1:B:260:LEU:HD12	2.03	0.41
1:A:60:LEU:HD21	1:A:263:ARG:HD3	2.03	0.40
1:B:575:PRO:O	1:B:586:ALA:HA	2.20	0.40
1:A:557:PRO:HB3	1:B:942:PHE:CZ	2.56	0.40
1:B:375:VAL:HG12	1:B:389:MET:HE2	2.03	0.40
1:A:146:LEU:O	1:A:150:ILE:HG12	2.22	0.40
1:A:596:ARG:HD3	1:A:677:SER:HB3	2.02	0.40
1:B:44:ASN:O	1:B:48:ARG:HG3	2.21	0.40
1:B:902:ALA:HA	1:B:918:ALA:HB2	2.03	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	1052/1113 (94%)	1001 (95%)	50 (5%)	1 (0%)	51 82
1	B	1054/1113 (95%)	1015 (96%)	38 (4%)	1 (0%)	51 82

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	2106/2226 (95%)	2016 (96%)	88 (4%)	2 (0%)	51 82

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	115	GLY
1	A	285	GLY

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	802/906 (88%)	798 (100%)	4 (0%)	88 93
1	B	828/906 (91%)	822 (99%)	6 (1%)	84 90
All	All	1630/1812 (90%)	1620 (99%)	10 (1%)	86 91

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	349	ASP
1	A	598	PHE
1	A	883	THR
1	A	894	ASN
1	B	346	ASP
1	B	349	ASP
1	B	883	THR
1	B	894	ASN
1	B	974	THR
1	B	1059	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	865	GLN

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Mol	Chain	Res	Type
1	A	894	ASN
1	A	1030	GLN
1	B	439	ASN
1	B	894	ASN
1	B	1030	GLN
1	B	1059	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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 4 are monoatomic - leaving 4 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$
3	GDP	B	1102	-	24,30,30	2.73	11 (45%)	30,47,47	1.31	5 (16%)
2	BCO	A	1101	-	47,55,55	2.13	10 (21%)	58,81,81	1.84	13 (22%)
2	BCO	B	1101	-	33,41,55	1.79	4 (12%)	42,64,81	1.49	5 (11%)
3	GDP	A	1102	4	24,30,30	2.73	10 (41%)	30,47,47	1.35	5 (16%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the

Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GDP	B	1102	-	-	0/12/32/32	0/3/3/3
2	BCO	A	1101	-	-	18/50/70/70	0/3/3/3
2	BCO	B	1101	-	-	9/30/52/70	0/3/3/3
3	GDP	A	1102	4	-	2/12/32/32	0/3/3/3

All (35) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	1102	GDP	O6-C6	6.85	1.37	1.23
3	A	1102	GDP	O6-C6	6.73	1.37	1.23
2	B	1101	BCO	C2-C4	-6.29	1.44	1.53
2	A	1101	BCO	C2-C4	-6.10	1.44	1.53
2	A	1101	BCO	C19-N7	6.00	1.47	1.33
3	B	1102	GDP	C2-N2	5.96	1.48	1.34
3	A	1102	GDP	C2-N2	5.96	1.48	1.34
2	A	1101	BCO	C16-N6	5.87	1.46	1.33
2	A	1101	BCO	C2-C1	-5.02	1.41	1.52
2	B	1101	BCO	C2-C1	-4.79	1.42	1.52
2	A	1101	BCO	C23-C22	4.06	1.55	1.50
3	A	1102	GDP	O5'-C5'	-3.92	1.29	1.44
3	A	1102	GDP	C5-C4	-3.84	1.33	1.43
3	B	1102	GDP	O5'-C5'	-3.83	1.30	1.44
3	A	1102	GDP	C2-N1	-3.82	1.28	1.37
2	B	1101	BCO	C12-N5	3.78	1.47	1.34
3	B	1102	GDP	C5-C4	-3.77	1.33	1.43
2	A	1101	BCO	C12-N5	3.76	1.47	1.34
3	B	1102	GDP	C2-N1	-3.75	1.28	1.37
3	A	1102	GDP	C2'-C1'	-3.39	1.48	1.53
3	B	1102	GDP	C2'-C1'	-3.14	1.49	1.53
2	A	1101	BCO	O17-C22	3.05	1.26	1.21
2	A	1101	BCO	C22-S1	2.91	1.83	1.76
3	B	1102	GDP	C2'-C3'	-2.77	1.45	1.53
3	A	1102	GDP	C2-N3	2.65	1.39	1.33
3	A	1102	GDP	PA-O2A	-2.64	1.42	1.55
3	A	1102	GDP	C2'-C3'	-2.62	1.46	1.53
3	B	1102	GDP	PA-O2A	-2.62	1.43	1.55
3	B	1102	GDP	C2-N3	2.50	1.39	1.33
2	A	1101	BCO	O1-C3	-2.35	1.39	1.45
2	B	1101	BCO	O1-C3	-2.34	1.39	1.45
2	A	1101	BCO	C1-C3	-2.23	1.46	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1102	GDP	C6-N1	-2.16	1.34	1.37
3	B	1102	GDP	PB-O3B	-2.13	1.46	1.54
3	B	1102	GDP	C6-N1	-2.05	1.34	1.37

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1101	BCO	C23-C22-S1	5.88	120.31	113.46
2	B	1101	BCO	N3-C15-N4	-5.21	120.53	128.68
2	A	1101	BCO	N3-C15-N4	-5.13	120.66	128.68
2	B	1101	BCO	O8-C11-C9	4.59	117.92	110.55
2	A	1101	BCO	O8-C11-C9	4.21	117.32	110.55
2	A	1101	BCO	C18-C17-N6	3.98	119.94	111.90
2	A	1101	BCO	O17-C22-C23	-3.55	119.80	123.99
2	A	1101	BCO	P3-O7-P2	-3.43	121.07	132.83
3	A	1102	GDP	C5-C6-N1	3.31	119.79	113.95
3	B	1102	GDP	C5-C6-N1	3.10	119.43	113.95
2	A	1101	BCO	C21-C20-N7	3.04	118.81	112.42
3	A	1102	GDP	PA-O3A-PB	-3.02	122.48	132.83
2	B	1101	BCO	P3-O7-P2	-3.00	122.54	132.83
3	B	1102	GDP	PA-O3A-PB	-2.98	122.60	132.83
3	A	1102	GDP	C2-N1-C6	-2.80	119.94	125.10
2	B	1101	BCO	O4-C5-C3	2.75	118.47	108.99
2	A	1101	BCO	C17-C18-C19	2.71	116.87	112.36
3	A	1102	GDP	C8-N7-C5	2.69	108.11	102.99
3	B	1102	GDP	C2-N1-C6	-2.68	120.16	125.10
3	B	1102	GDP	C8-N7-C5	2.64	108.01	102.99
3	A	1102	GDP	O6-C6-C5	-2.47	119.54	124.37
2	A	1101	BCO	C21-S1-C22	2.47	109.55	101.87
2	B	1101	BCO	C1-C2-C4	2.38	105.17	99.89
2	A	1101	BCO	O4-C5-C3	2.37	117.15	108.99
2	A	1101	BCO	C1-C2-C4	2.37	105.13	99.89
2	A	1101	BCO	O17-C22-S1	-2.20	119.75	122.61
3	B	1102	GDP	O6-C6-C5	-2.15	120.17	124.37
2	A	1101	BCO	C24-C23-C22	2.08	117.88	112.72

There are no chirality outliers.

All (29) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1101	BCO	C5-O4-P2-O10
2	A	1101	BCO	C5-O4-P2-O7

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Mol	Chain	Res	Type	Atoms
2	A	1101	BCO	C11-O8-P3-O13
2	A	1101	BCO	N6-C17-C18-C19
2	A	1101	BCO	N7-C20-C21-S1
2	A	1101	BCO	C20-C21-S1-C22
2	B	1101	BCO	C11-O8-P3-O7
2	B	1101	BCO	C11-O8-P3-O14
2	B	1101	BCO	C11-O8-P3-O13
2	B	1101	BCO	O11-C10-C9-C11
2	A	1101	BCO	O1-C3-C5-O4
2	B	1101	BCO	O11-C10-C9-C13
2	B	1101	BCO	O11-C10-C9-C14
2	B	1101	BCO	C16-C10-C9-C14
2	A	1101	BCO	C23-C22-S1-C21
2	A	1101	BCO	C11-O8-P3-O7
2	A	1101	BCO	C11-O8-P3-O14
2	A	1101	BCO	C2-C1-O2-P1
2	A	1101	BCO	C3-C1-O2-P1
2	A	1101	BCO	C1-C3-C5-O4
2	B	1101	BCO	C16-C10-C9-C13
3	A	1102	GDP	PB-O3A-PA-O2A
2	A	1101	BCO	O17-C22-S1-C21
2	B	1101	BCO	P3-O7-P2-O4
2	A	1101	BCO	C1-O2-P1-O5
2	A	1101	BCO	P3-O7-P2-O12
2	A	1101	BCO	P3-O7-P2-O10
3	A	1102	GDP	C5'-O5'-PA-O1A
2	A	1101	BCO	C17-C18-C19-O16

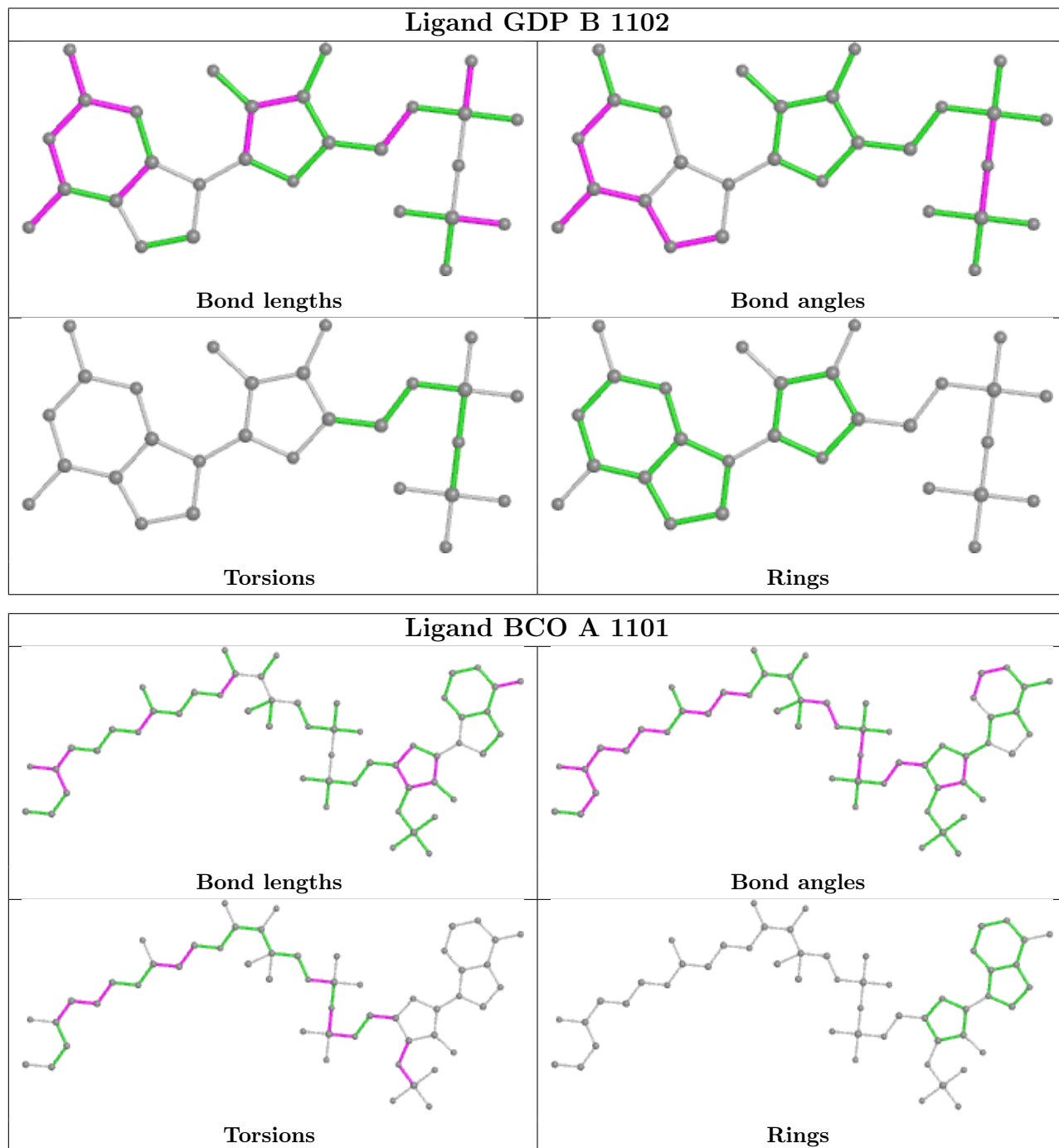
There are no ring outliers.

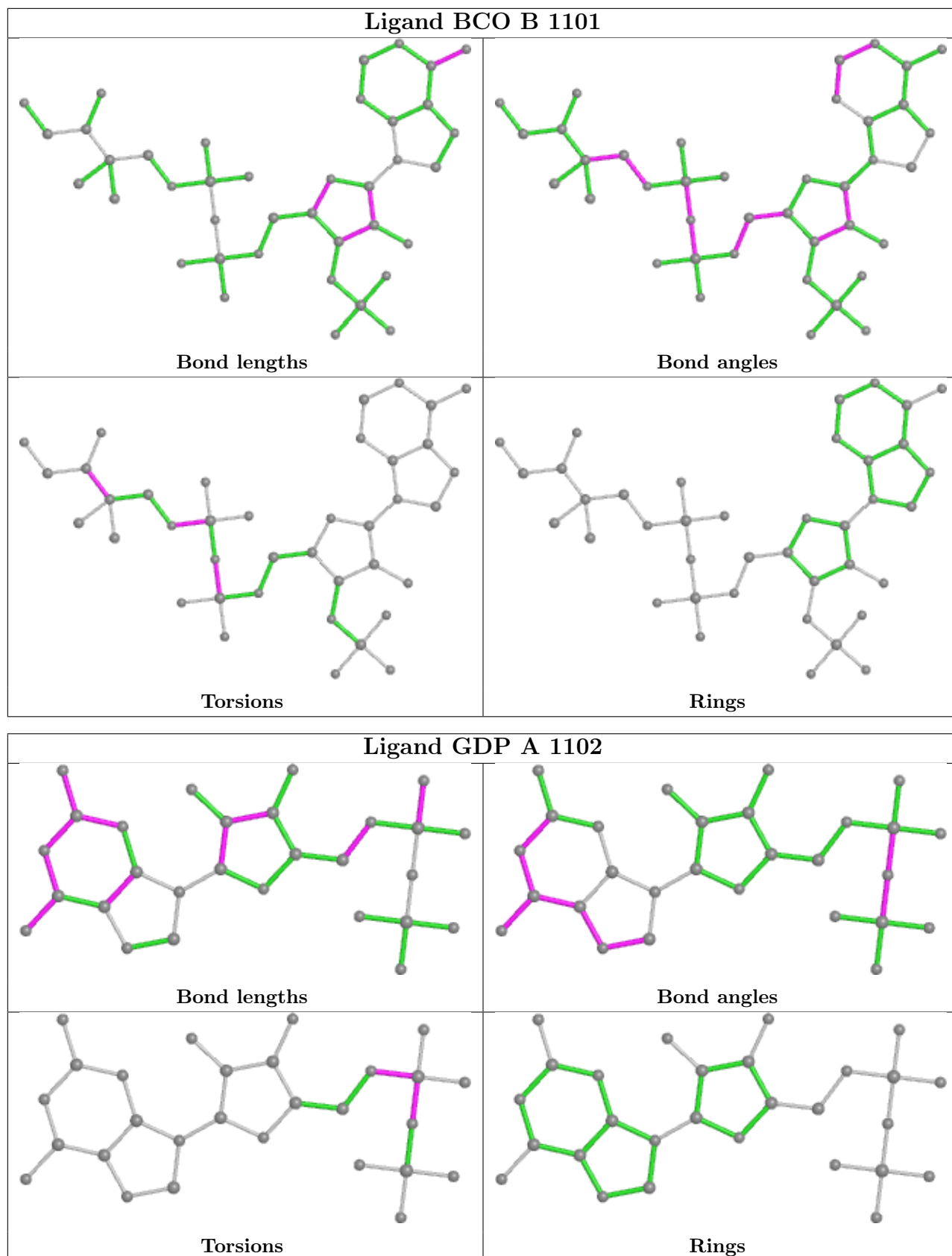
3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	1102	GDP	1	0
2	A	1101	BCO	1	0
3	A	1102	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	1058/1113 (95%)	0.50	40 (3%) 40 37	63, 102, 148, 200	0
1	B	1060/1113 (95%)	0.22	4 (0%) 92 92	53, 79, 129, 182	0
All	All	2118/2226 (95%)	0.36	44 (2%) 63 61	53, 91, 143, 200	0

All (44) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	620	ALA	5.8
1	A	494	LEU	5.3
1	A	280	ALA	3.5
1	B	489	HIS	3.5
1	A	570	MET	3.2
1	A	246	ILE	3.0
1	A	279	LEU	3.0
1	A	157	CYS	3.0
1	A	148	GLY	3.0
1	A	200	LEU	3.0
1	A	85	SER	2.9
1	A	119	ILE	2.7
1	A	57	VAL	2.7
1	B	488	GLY	2.6
1	A	108	HIS	2.6
1	A	805	ALA	2.6
1	A	32	ALA	2.6
1	A	43	ILE	2.6
1	A	184	ILE	2.5
1	A	266	MET	2.5
1	A	491	SER	2.5
1	A	187	LEU	2.4
1	A	816	PHE	2.4
1	A	258	ALA	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	113	GLY	2.4
1	B	477	LEU	2.4
1	A	49	ILE	2.3
1	A	171	THR	2.3
1	A	66	VAL	2.3
1	A	31	THR	2.2
1	A	47	ARG	2.2
1	A	118	VAL	2.2
1	A	711	PRO	2.2
1	A	275	PHE	2.1
1	B	491	SER	2.1
1	A	311	THR	2.1
1	A	336	PHE	2.1
1	A	33	ALA	2.1
1	A	196	LEU	2.1
1	A	489	HIS	2.1
1	A	58	ILE	2.1
1	A	27	VAL	2.0
1	A	60	LEU	2.0
1	A	83	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 monosaccharides 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.

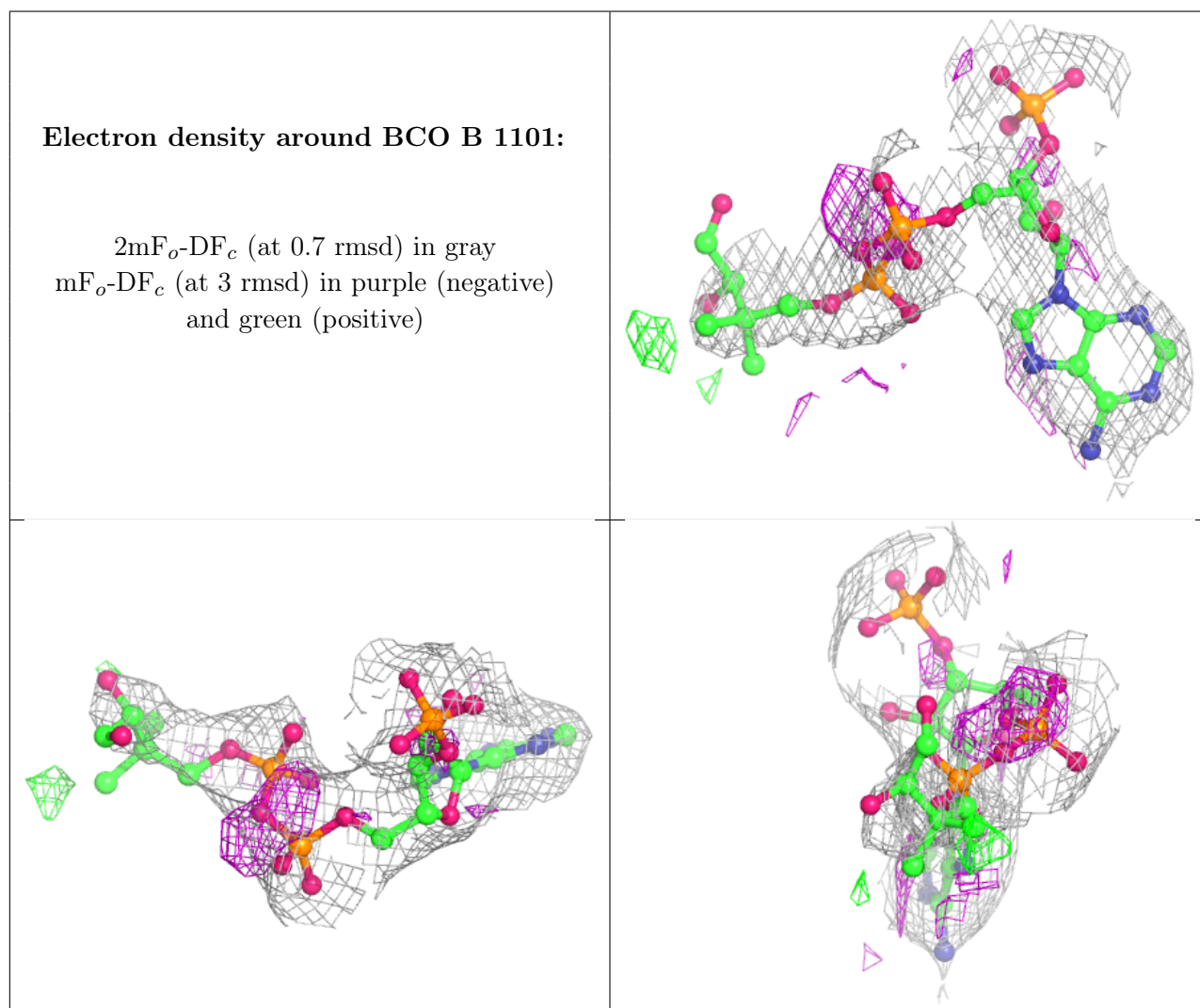
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	BCO	B	1101	39/53	0.86	0.21	108,139,221,267	0
4	MG	A	1104	1/1	0.89	0.10	66,66,66,66	0
4	MG	B	1104	1/1	0.93	0.19	61,61,61,61	0

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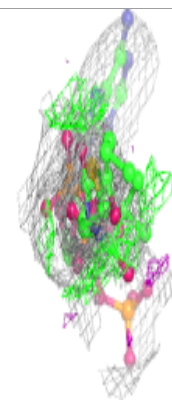
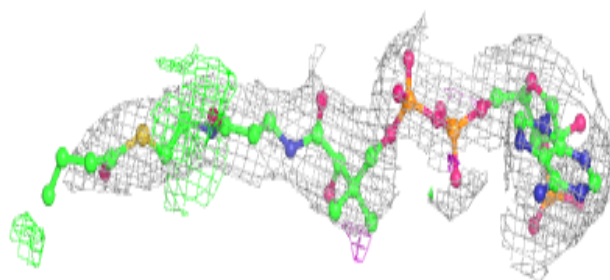
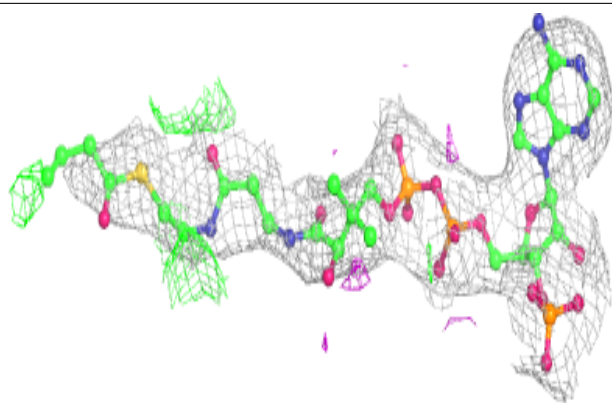
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	BCO	A	1101	53/53	0.94	0.24	65,105,132,135	0
3	GDP	A	1102	28/28	0.95	0.20	73,86,96,99	0
4	MG	B	1103	1/1	0.96	0.12	82,82,82,82	0
4	MG	A	1103	1/1	0.97	0.13	77,77,77,77	0
3	GDP	B	1102	28/28	0.97	0.21	52,59,74,76	0

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.

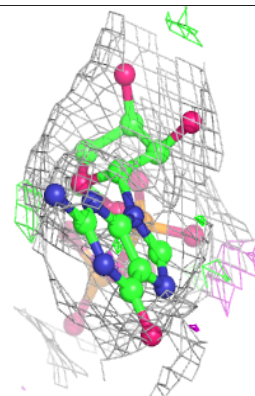
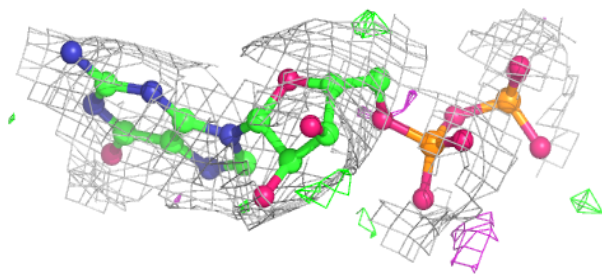
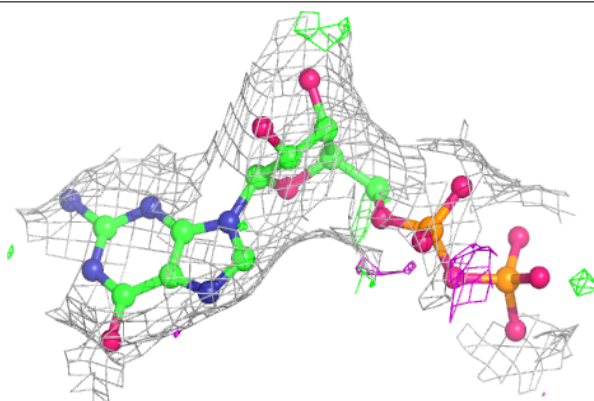


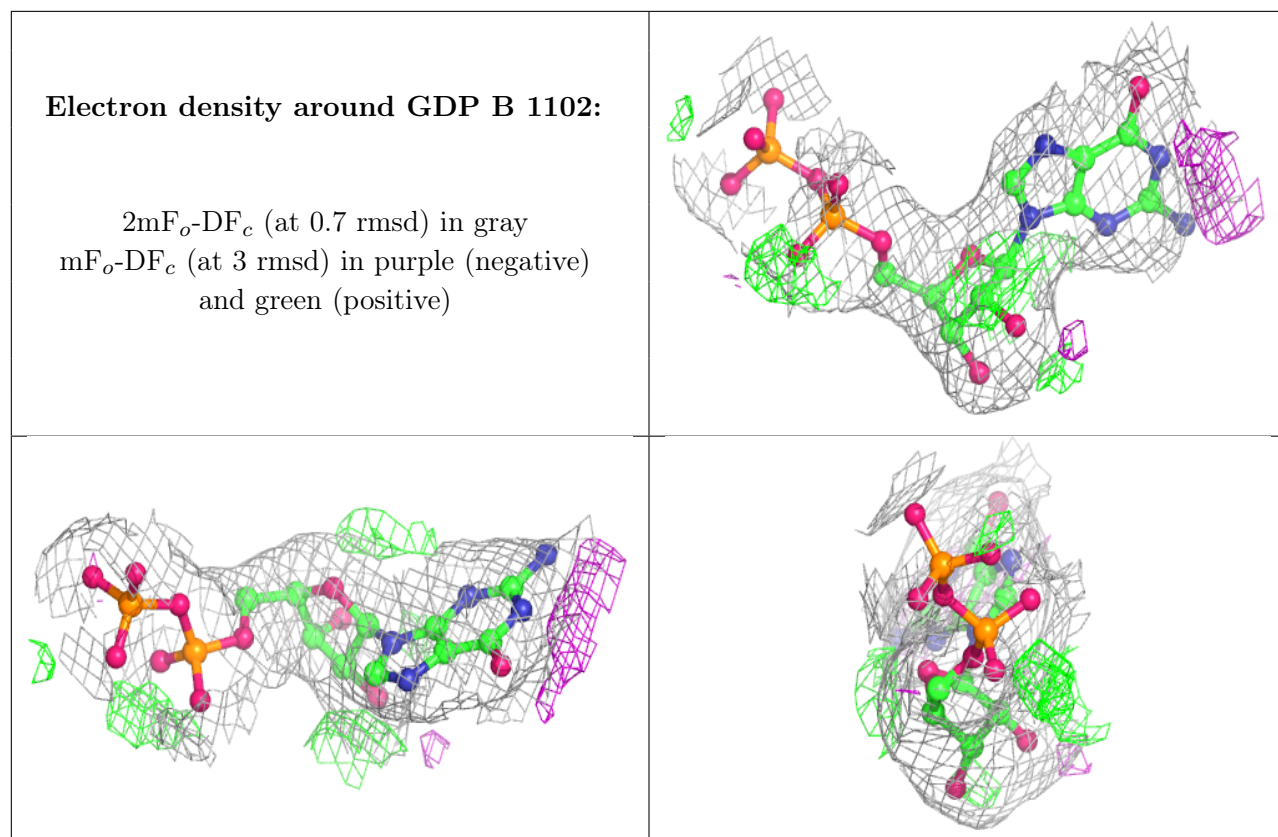
Electron density around BCO A 1101:

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

**Electron density around GDP A 1102:**

$2mF_o-DF_c$ (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.