Full wwPDB X-ray Structure Validation Report ⓘ

May 29, 2020 – 04:36 pm BST

PDB ID	:	2WIB
Title	:	Crystal Structures of the N-terminal Intracellular Domain of FeoB from Klebsiella Pneumoniae in GDP binding state
Authors	:	Hung, K.-W.; Chang, Y.-W.; Chen, J.-H.; Chen, Y.-C.; Sun, Y.-J.; Hsiao, C.-D.; Huang, T.-H.
Deposited on	:	2009-05-09
Resolution	:	2.56 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:

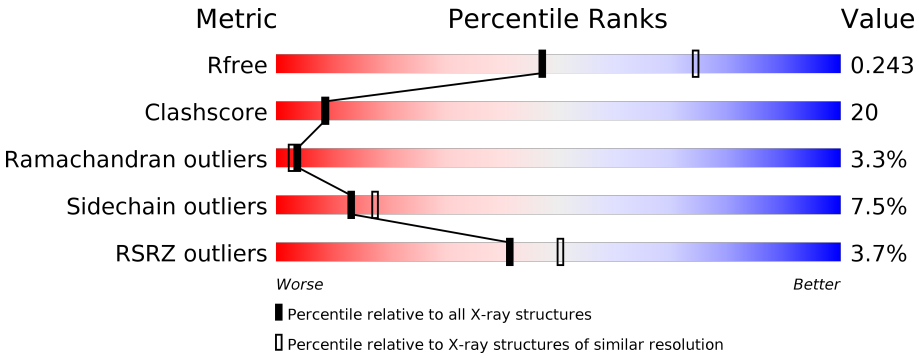
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.11
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber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

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.56 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	1279 (2.58-2.54)
Clashscore	141614	1327 (2.58-2.54)
Ramachandran outliers	138981	1312 (2.58-2.54)
Sidechain outliers	138945	1312 (2.58-2.54)
RSRZ outliers	127900	1269 (2.58-2.54)

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 <=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	267	2% 61% 30% 6%
1	B	267	5% 54% 30% 10%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3839 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.

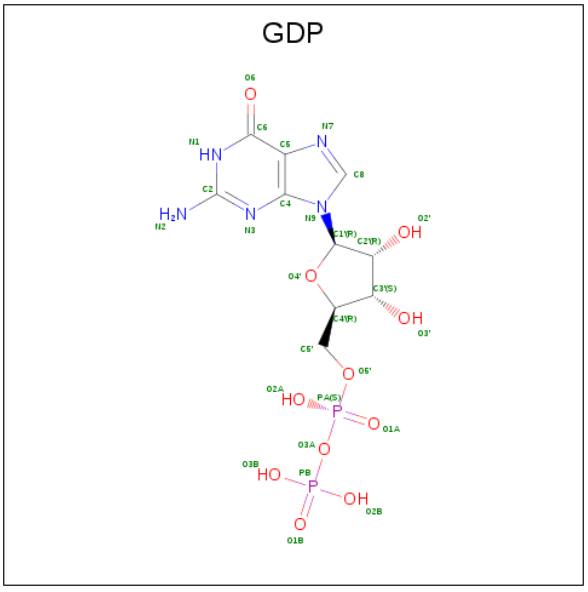
- Molecule 1 is a protein called FERROUS IRON TRANSPORT PROTEIN B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	261	Total	C	N	O	S	0	0	0
			1874	1175	328	362	9			
1	B	257	Total	C	N	O	S	0	0	0
			1792	1126	312	345	9			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	129	GLN	LYS	conflict	UNP A6TF32
B	129	GLN	LYS	conflict	UNP A6TF32

- Molecule 2 is GUANOSINE-5'-DIPHOSPHATE (three-letter code: GDP) (formula: $C_{10}H_{15}N_5O_{11}P_2$).



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

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

- Molecule 3 is water.

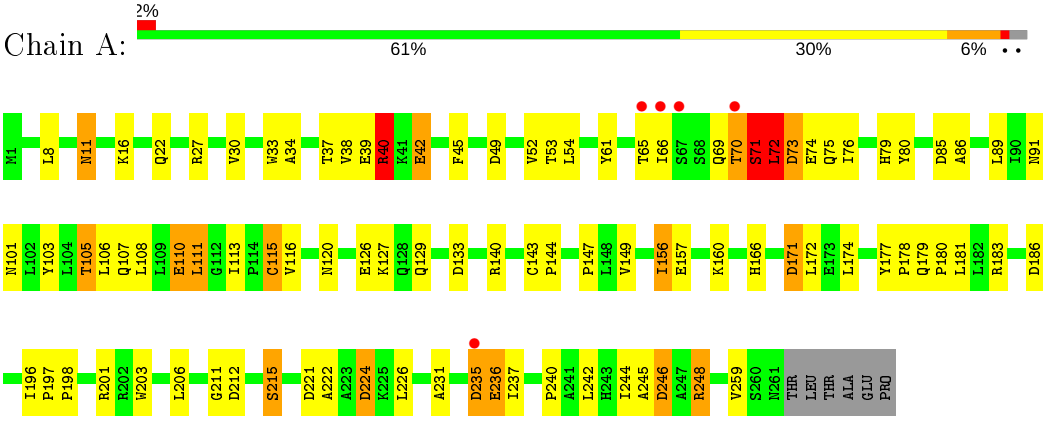
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	87	Total	O	0	0
			87	87		
3	B	30	Total	O	0	0
			30	30		

3 Residue-property plots (i)

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.

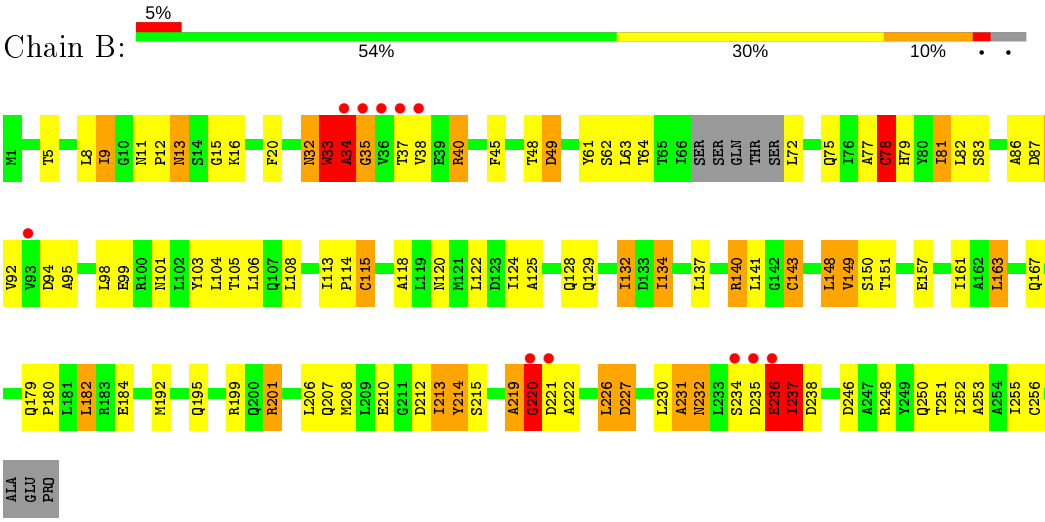
- Molecule 1: FERROUS IRON TRANSPORT PROTEIN B

Chain A:



- Molecule 1: FERROUS IRON TRANSPORT PROTEIN B

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α, β, γ	86.55Å 184.69Å 38.17Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.53 – 2.56 19.53 – 2.54	Depositor EDS
% Data completeness (in resolution range)	82.4 (19.53-2.56) 90.1 (19.53-2.54)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	13.23 (at 2.53Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.222 , 0.236 0.231 , 0.243	Depositor DCC
R_{free} test set	949 reflections (4.84%)	wwPDB-VP
Wilson B-factor (Å^2)	44.2	Xtriage
Anisotropy	0.474	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.27 , 44.4	EDS
L-test for twinning[2]	$< \|L\| > = 0.46$, $< L^2 > = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.91	EDS
Total number of atoms	3839	wwPDB-VP
Average B, all atoms (Å^2)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 15.41% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

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

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	#$\|Z\|$ >5	RMSZ	#$\|Z\|$ >5
1	A	1.52	18/1899 (0.9%)	1.46	21/2599 (0.8%)
1	B	1.45	11/1814 (0.6%)	1.55	31/2484 (1.2%)
All	All	1.49	29/3713 (0.8%)	1.51	52/5083 (1.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand.A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	3
1	B	1	13
All	All	1	16

All (29) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	237	ILE	N-CA	11.27	1.68	1.46
1	A	236	GLU	N-CA	9.96	1.66	1.46
1	A	177	TYR	CE2-CZ	8.28	1.49	1.38
1	A	42	GLU	CB-CG	-8.06	1.36	1.52
1	A	115	CYS	CB-SG	-7.74	1.69	1.82
1	A	49	ASP	CB-CG	7.58	1.67	1.51
1	B	78	CYS	CB-SG	-7.30	1.69	1.82
1	B	49	ASP	CB-CG	7.02	1.66	1.51
1	A	126	GLU	CA-CB	6.99	1.69	1.53
1	A	38	VAL	CB-CG2	6.75	1.67	1.52
1	A	110	GLU	CD-OE1	6.45	1.32	1.25
1	B	253	ALA	CA-CB	-6.27	1.39	1.52
1	B	143	CYS	CB-SG	6.14	1.92	1.82
1	B	220	GLY	CA-C	-6.04	1.42	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	151	THR	C-O	-5.93	1.12	1.23
1	B	45	PHE	CE2-CZ	5.84	1.48	1.37
1	A	53	THR	CB-CG2	5.83	1.71	1.52
1	A	157	GLU	CB-CG	5.80	1.63	1.52
1	A	157	GLU	CG-CD	5.61	1.60	1.51
1	A	30	VAL	CB-CG1	-5.59	1.41	1.52
1	A	183	ARG	CG-CD	5.57	1.65	1.51
1	A	61	TYR	CE1-CZ	5.52	1.45	1.38
1	B	149	VAL	CB-CG1	5.49	1.64	1.52
1	A	39	GLU	CD-OE2	5.46	1.31	1.25
1	B	157	GLU	CB-CG	5.36	1.62	1.52
1	A	231	ALA	CA-CB	5.32	1.63	1.52
1	A	224	ASP	CA-CB	5.31	1.65	1.53
1	B	13	ASN	CB-CG	5.14	1.62	1.51
1	A	11	ASN	N-CA	-5.11	1.36	1.46

All (52) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	173	GLU	O-C-N	-13.84	100.55	122.70
1	A	72	LEU	O-C-N	-9.74	107.11	122.70
1	B	173	GLU	CB-CA-C	9.30	129.01	110.40
1	B	237	ILE	O-C-N	-9.27	107.86	122.70
1	B	32	ASN	C-N-CA	-8.79	99.73	121.70
1	B	201	ARG	NE-CZ-NH2	-8.66	115.97	120.30
1	B	220	GLY	N-CA-C	8.37	134.03	113.10
1	B	237	ILE	N-CA-C	-8.23	88.78	111.00
1	A	111	LEU	CB-CG-CD1	-8.05	97.32	111.00
1	A	221	ASP	O-C-N	-8.02	109.88	122.70
1	A	85	ASP	CB-CG-OD1	8.01	125.51	118.30
1	A	40	ARG	NE-CZ-NH2	-7.91	116.34	120.30
1	B	173	GLU	CA-C-N	7.89	134.56	117.20
1	A	129	GLN	O-C-N	-7.88	110.09	122.70
1	A	133	ASP	CB-CG-OD1	7.34	124.90	118.30
1	B	172	LEU	O-C-N	-7.23	111.14	122.70
1	B	220	GLY	O-C-N	-7.19	111.19	122.70
1	B	82	LEU	CB-CG-CD1	-7.02	99.06	111.00
1	B	38	VAL	O-C-N	-6.73	111.93	122.70
1	B	221	ASP	O-C-N	-6.64	112.08	122.70
1	B	182	LEU	CB-CG-CD2	-6.52	99.91	111.00
1	B	33	TRP	C-N-CA	6.38	137.64	121.70
1	A	105	THR	CA-CB-CG2	-6.36	103.49	112.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	236	GLU	N-CA-C	-6.27	94.06	111.00
1	B	149	VAL	CB-CA-C	6.17	123.13	111.40
1	B	134	ILE	CG1-CB-CG2	-6.10	97.98	111.40
1	B	172	LEU	C-N-CA	-6.09	106.48	121.70
1	B	140	ARG	NE-CZ-NH2	-6.03	117.28	120.30
1	B	34	ALA	N-CA-CB	5.92	118.39	110.10
1	B	113	ILE	CB-CA-C	-5.92	99.76	111.60
1	A	248	ARG	NE-CZ-NH1	-5.87	117.37	120.30
1	B	232	ASN	N-CA-C	5.82	126.70	111.00
1	B	49	ASP	CB-CG-OD2	5.76	123.48	118.30
1	A	235	ASP	C-N-CA	-5.74	107.34	121.70
1	A	242	LEU	CB-CG-CD1	5.73	120.74	111.00
1	A	140	ARG	NE-CZ-NH1	-5.70	117.45	120.30
1	B	220	GLY	CA-C-O	5.59	130.67	120.60
1	A	129	GLN	CA-C-O	5.53	131.72	120.10
1	A	246	ASP	CB-CG-OD1	-5.52	113.33	118.30
1	A	221	ASP	CA-C-O	5.45	131.55	120.10
1	B	38	VAL	CA-C-N	5.43	129.15	117.20
1	A	52	VAL	CB-CA-C	-5.42	101.10	111.40
1	B	237	ILE	CA-C-O	5.36	131.36	120.10
1	A	100	ARG	CB-CA-C	5.33	121.07	110.40
1	B	173	GLU	N-CA-CB	5.30	120.14	110.60
1	A	66	ILE	O-C-N	-5.20	114.37	122.70
1	B	143	CYS	N-CA-C	-5.13	97.14	111.00
1	B	163	LEU	CB-CG-CD1	-5.08	102.36	111.00
1	A	171	ASP	CB-CG-OD1	-5.08	113.73	118.30
1	B	236	GLU	C-N-CA	-5.02	109.14	121.70
1	A	171	ASP	CB-CG-OD2	5.01	122.81	118.30
1	B	235	ASP	O-C-N	5.01	130.72	122.70

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	B	173	GLU	CA

All (16) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	236	GLU	Peptide
1	A	71	SER	Mainchain
1	A	72	LEU	Mainchain
1	B	171	ASP	Mainchain

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Mol	Chain	Res	Type	Group
1	B	219	ALA	Mainchain,Peptide
1	B	220	GLY	Mainchain,Peptide
1	B	226	LEU	Mainchain
1	B	231	ALA	Mainchain
1	B	236	GLU	Mainchain
1	B	237	ILE	Mainchain
1	B	32	ASN	Mainchain
1	B	34	ALA	Peptide
1	B	35	GLY	Mainchain
1	B	37	THR	Mainchain

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	1874	0	1800	71	0
1	B	1792	0	1688	79	0
2	A	28	0	12	0	0
2	B	28	0	12	1	0
3	A	87	0	0	5	0
3	B	30	0	0	2	0
All	All	3839	0	3512	148	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 20.

All (148) 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:237:ILE:N	1:B:237:ILE:CA	1.68	1.53
1:B:220:GLY:HA2	1:B:222:ALA:H	1.34	0.93
1:B:237:ILE:N	1:B:237:ILE:C	2.25	0.90
1:B:236:GLU:C	1:B:237:ILE:CA	2.41	0.89
1:A:22:GLN:HB3	1:A:156:ILE:HD12	1.54	0.87
1:B:137:LEU:HG	1:B:141:LEU:CD1	2.05	0.85
1:A:178:PRO:HD2	1:A:181:LEU:HD12	1.61	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:192:MET:CE	1:B:201:ARG:HB3	2.10	0.81
1:B:120:ASN:HD21	1:B:150:SER:H	1.28	0.80
1:B:137:LEU:HG	1:B:141:LEU:HD11	1.62	0.78
1:B:167:GLN:HA	1:B:167:GLN:OE1	1.86	0.75
1:A:240:PRO:O	1:A:244:ILE:HG13	1.87	0.75
1:A:201:ARG:HG3	1:A:201:ARG:HH11	1.51	0.75
1:B:192:MET:CE	1:B:201:ARG:CB	2.65	0.74
1:A:65:THR:HG21	1:A:70:THR:HG21	1.69	0.74
1:A:127:LYS:NZ	1:B:94:ASP:OD2	2.21	0.74
1:B:128:GLN:O	3:B:2016:HOH:O	2.04	0.73
1:B:210:GLU:HG3	1:B:210:GLU:O	1.90	0.71
1:A:22:GLN:HB3	1:A:156:ILE:CD1	2.21	0.70
1:B:115:CYS:O	1:B:143:CYS:HB2	1.91	0.69
1:B:81:ILE:HG21	1:B:89:LEU:HD11	1.74	0.69
1:A:45:PHE:CD1	1:A:160:LYS:HE2	2.28	0.68
1:B:11:ASN:HD21	1:B:101:ASN:HA	1.58	0.67
1:A:192:MET:HB2	1:A:201:ARG:HD2	1.77	0.67
1:B:137:LEU:HG	1:B:141:LEU:HD12	1.77	0.66
1:A:65:THR:N	1:A:212:ASP:OD2	2.27	0.65
1:B:192:MET:HE1	1:B:201:ARG:HB3	1.78	0.64
1:B:11:ASN:ND2	1:B:101:ASN:HA	2.12	0.64
1:B:192:MET:CE	1:B:201:ARG:HA	2.28	0.64
1:A:22:GLN:CB	1:A:156:ILE:HD12	2.28	0.63
1:B:192:MET:HE3	1:B:201:ARG:HA	1.80	0.63
1:A:65:THR:O	1:A:212:ASP:HA	1.99	0.62
1:B:234:SER:C	1:B:236:GLU:H	2.01	0.62
1:A:201:ARG:NH1	1:A:201:ARG:HG3	2.12	0.62
1:A:22:GLN:CB	1:A:156:ILE:CD1	2.78	0.61
1:B:213:ILE:O	1:B:214:TYR:C	2.38	0.61
1:B:137:LEU:CG	1:B:141:LEU:HD11	2.31	0.61
1:B:237:ILE:N	1:B:237:ILE:CB	2.57	0.61
1:A:111:LEU:HD13	1:A:113:ILE:HD11	1.82	0.61
1:B:48:THR:HG22	1:B:49:ASP:OD1	2.01	0.60
1:B:95:ALA:CB	1:B:122:LEU:HD23	2.31	0.60
1:A:11:ASN:HD21	1:A:101:ASN:HA	1.65	0.59
1:B:210:GLU:CG	1:B:210:GLU:O	2.49	0.59
1:B:33:TRP:CD1	1:B:40:ARG:HB3	2.37	0.59
1:A:27:ARG:HD2	3:A:2022:HOH:O	2.02	0.59
1:B:192:MET:CE	1:B:201:ARG:CA	2.81	0.59
1:B:192:MET:HE2	1:B:201:ARG:HB3	1.84	0.58
1:B:5:THR:HG22	1:B:86:ALA:HA	1.84	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:186:ASP:HB3	3:A:2071:HOH:O	2.04	0.58
1:B:192:MET:HE1	1:B:201:ARG:CA	2.34	0.57
1:A:40:ARG:NH1	1:A:42:GLU:OE2	2.37	0.57
1:A:144:PRO:HD3	1:A:166:HIS:CE1	2.39	0.57
1:A:72:LEU:O	1:A:73:ASP:C	2.42	0.57
1:A:212:ASP:OD1	1:A:215:SER:HB2	2.06	0.56
1:B:118:ALA:HB1	1:B:148:LEU:HD13	1.89	0.55
1:A:116:VAL:HG13	1:A:144:PRO:HB2	1.88	0.55
1:A:89:LEU:CD1	1:A:108:LEU:HD22	2.36	0.54
1:A:65:THR:CG2	1:A:70:THR:HG21	2.38	0.54
1:A:115:CYS:O	1:A:143:CYS:HB2	2.07	0.54
1:A:69:GLN:CB	3:A:2030:HOH:O	2.55	0.54
1:A:111:LEU:HD21	1:A:203:TRP:CE3	2.43	0.54
1:A:89:LEU:HD12	1:A:108:LEU:HD22	1.89	0.54
1:B:192:MET:HE1	1:B:201:ARG:CB	2.37	0.54
1:B:64:THR:HA	1:B:212:ASP:OD2	2.09	0.53
1:B:88:MET:HE3	3:B:2022:HOH:O	2.08	0.53
1:B:120:ASN:HD21	1:B:150:SER:N	2.03	0.53
1:A:11:ASN:ND2	1:A:101:ASN:HA	2.24	0.53
1:B:87:ASP:O	1:B:114:PRO:HG2	2.09	0.53
1:A:33:TRP:CE2	1:A:40:ARG:HG3	2.44	0.52
1:B:140:ARG:HB3	1:B:255:ILE:HD12	1.93	0.51
1:B:213:ILE:O	1:B:215:SER:N	2.44	0.51
1:A:8:LEU:HG	1:A:16:LYS:HD2	1.93	0.50
1:B:72:LEU:HA	1:B:75:GLN:CB	2.42	0.50
1:B:180:PRO:O	1:B:184:GLU:HG2	2.11	0.50
1:B:63:LEU:HD11	1:B:108:LEU:HD23	1.94	0.50
1:B:124:ILE:O	1:B:125:ALA:C	2.50	0.49
1:A:106:LEU:O	1:A:110:GLU:HG3	2.12	0.49
1:A:203:TRP:O	1:A:206:LEU:HB2	2.14	0.48
1:B:237:ILE:O	1:B:238:ASP:C	2.48	0.48
1:B:120:ASN:ND2	1:B:150:SER:H	2.05	0.48
1:B:99:GLU:HB2	1:B:256:CYS:SG	2.53	0.48
1:B:220:GLY:HA2	1:B:222:ALA:N	2.16	0.48
1:A:22:GLN:CB	1:A:156:ILE:HD11	2.44	0.47
1:A:222:ALA:C	1:A:224:ASP:N	2.64	0.47
1:A:42:GLU:HA	1:A:54:LEU:O	2.14	0.47
1:B:163:LEU:HD23	1:B:163:LEU:HA	1.53	0.47
1:A:149:VAL:O	1:A:149:VAL:HG22	2.14	0.47
1:B:227:ASP:HA	1:B:230:LEU:CB	2.44	0.47
1:B:78:CYS:O	1:B:79:HIS:C	2.52	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:246:ASP:O	1:B:250:GLN:HB2	2.15	0.47
1:B:8:LEU:HG	1:B:16:LYS:HG3	1.97	0.47
1:A:111:LEU:HD13	1:A:113:ILE:CD1	2.45	0.46
1:A:143:CYS:O	1:A:144:PRO:C	2.52	0.46
1:A:171:ASP:C	1:A:171:ASP:OD1	2.53	0.46
1:B:16:LYS:O	1:B:20:PHE:N	2.43	0.46
1:B:179:GLN:NE2	1:B:182:LEU:HD12	2.29	0.46
1:A:186:ASP:HA	1:A:189:ALA:HB3	1.98	0.46
1:B:81:ILE:CG2	1:B:89:LEU:HD11	2.43	0.46
1:A:45:PHE:HD1	1:A:160:LYS:HE2	1.80	0.46
1:A:147:PRO:HG2	3:A:2057:HOH:O	2.16	0.46
1:A:127:LYS:NZ	1:B:13:ASN:O	2.49	0.46
1:B:230:LEU:C	1:B:232:ASN:N	2.69	0.46
1:A:226:LEU:HA	1:A:226:LEU:HD13	1.84	0.45
1:B:184:GLU:HG3	1:B:232:ASN:ND2	2.32	0.45
1:B:98:LEU:HD22	1:B:132:ILE:HD13	1.99	0.45
1:A:98:LEU:O	1:A:99:GLU:C	2.55	0.44
1:B:64:THR:OG1	1:B:207:GLN:NE2	2.51	0.44
1:B:98:LEU:HD22	1:B:132:ILE:CD1	2.46	0.44
1:B:8:LEU:HD11	1:B:92:VAL:CG2	2.47	0.44
1:A:178:PRO:CD	1:A:181:LEU:HD12	2.42	0.44
1:B:103:TYR:HA	1:B:252:ILE:HD13	1.99	0.44
1:B:179:GLN:N	1:B:180:PRO:HD2	2.33	0.44
1:A:206:LEU:HA	1:A:206:LEU:HD23	1.63	0.43
1:A:103:TYR:O	1:A:107:GLN:HG3	2.18	0.43
1:A:80:TYR:HE2	1:A:86:ALA:HB2	1.83	0.43
1:A:33:TRP:HH2	1:A:80:TYR:CD1	2.36	0.43
1:B:251:THR:O	1:B:255:ILE:HG12	2.19	0.43
1:A:172:LEU:HA	1:A:172:LEU:HD23	1.74	0.43
1:B:182:LEU:HD23	1:B:182:LEU:HA	1.75	0.42
1:A:79:HIS:HB2	3:A:2075:HOH:O	2.20	0.42
1:A:197:PRO:HA	1:A:198:PRO:HD3	1.83	0.42
1:A:80:TYR:CE2	1:A:86:ALA:HB2	2.54	0.42
1:B:12:PRO:HG2	1:B:61:TYR:CE2	2.55	0.42
1:B:220:GLY:CA	1:B:222:ALA:H	2.17	0.42
1:A:211:GLY:O	1:A:212:ASP:C	2.58	0.42
1:B:120:ASN:HA	1:B:148:LEU:O	2.19	0.42
1:A:179:GLN:CB	1:A:180:PRO:HD3	2.49	0.41
1:B:104:LEU:O	1:B:105:THR:C	2.57	0.41
1:A:120:ASN:HD22	1:A:120:ASN:HA	1.59	0.41
1:B:161:ILE:HD13	1:B:161:ILE:N	2.34	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:9:ILE:O	1:B:91:ASN:HA	2.21	0.41
1:A:196:ILE:HG22	1:A:201:ARG:HG2	2.03	0.41
1:A:70:THR:HG23	1:A:71:SER:H	1.85	0.41
1:B:92:VAL:HA	1:B:118:ALA:HB3	2.02	0.41
1:B:103:TYR:HA	1:B:252:ILE:CD1	2.51	0.41
1:A:22:GLN:HB2	1:A:156:ILE:CD1	2.49	0.41
1:A:222:ALA:C	1:A:224:ASP:H	2.24	0.41
1:A:22:GLN:HB2	1:A:156:ILE:HD11	2.03	0.41
1:B:15:GLY:HA2	2:B:1263:GDP:PA	2.61	0.41
1:A:22:GLN:HA	1:A:22:GLN:HE21	1.86	0.41
1:A:33:TRP:CH2	1:A:80:TYR:CD1	3.09	0.41
1:A:111:LEU:HB3	1:A:113:ILE:HD12	2.03	0.40
1:A:245:ALA:O	1:A:246:ASP:C	2.58	0.40
1:A:91:ASN:ND2	1:A:105:THR:HG23	2.35	0.40
1:B:192:MET:SD	1:B:219:ALA:HB1	2.60	0.40
1:A:248:ARG:HD2	1:A:248:ARG:HH11	1.61	0.40
1:A:73:ASP:O	1:A:75:GLN:N	2.55	0.40
1:B:206:LEU:HA	1:B:206:LEU:HD23	1.60	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	259/267 (97%)	233 (90%)	19 (7%)	7 (3%)	5 4
1	B	253/267 (95%)	219 (87%)	24 (10%)	10 (4%)	3 2
All	All	512/534 (96%)	452 (88%)	43 (8%)	17 (3%)	4 3

All (17) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	214	TYR
1	A	73	ASP
1	A	74	GLU
1	A	237	ILE
1	B	35	GLY
1	B	213	ILE
1	B	226	LEU
1	B	231	ALA
1	A	71	SER
1	B	34	ALA
1	B	129	GLN
1	B	237	ILE
1	A	34	ALA
1	A	70	THR
1	A	235	ASP
1	B	77	ALA
1	B	227	ASP

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	182/221 (82%)	175 (96%)	7 (4%)	33 44
1	B	166/221 (75%)	147 (89%)	19 (11%)	5 6
All	All	348/442 (79%)	322 (92%)	26 (8%)	13 17

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

Mol	Chain	Res	Type
1	A	37	THR
1	A	40	ARG
1	A	76	ILE
1	A	156	ILE
1	A	174	LEU
1	A	215	SER
1	A	259	VAL

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Mol	Chain	Res	Type
1	B	9	ILE
1	B	33	TRP
1	B	40	ARG
1	B	62	SER
1	B	78	CYS
1	B	81	ILE
1	B	83	SER
1	B	88	MET
1	B	106	LEU
1	B	115	CYS
1	B	132	ILE
1	B	134	ILE
1	B	148	LEU
1	B	149	VAL
1	B	172	LEU
1	B	195	GLN
1	B	199	ARG
1	B	208	MET
1	B	248	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	22	GLN
1	A	51	GLN
1	A	91	ASN
1	A	120	ASN
1	A	195	GLN
1	B	11	ASN
1	B	22	GLN
1	B	51	GLN
1	B	91	ASN
1	B	120	ASN
1	B	129	GLN
1	B	179	GLN
1	B	207	GLN
1	B	232	ASN

5.3.3 RNA

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 carbohydrates in this entry.

5.6 Ligand geometry (i)

2 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	GDP	B	1263	-	24,30,30	3.33	6 (25%)	31,47,47	2.78	12 (38%)
2	GDP	A	1262	-	24,30,30	2.42	8 (33%)	31,47,47	3.02	14 (45%)

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	GDP	B	1263	-	-	0/12/32/32	0/3/3/3
2	GDP	A	1262	-	-	3/12/32/32	0/3/3/3

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1263	GDP	C6-N1	8.84	1.48	1.33
2	B	1263	GDP	C2-N1	8.25	1.50	1.35
2	A	1262	GDP	C6-N1	7.17	1.45	1.33
2	B	1263	GDP	PB-O3B	6.90	1.81	1.54
2	B	1263	GDP	O4'-C1'	5.25	1.48	1.41
2	A	1262	GDP	C8-N7	4.37	1.42	1.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1262	GDP	PB-O2B	-4.03	1.39	1.54
2	B	1263	GDP	C2'-C1'	-3.81	1.48	1.53
2	A	1262	GDP	C2-N1	3.71	1.42	1.35
2	A	1262	GDP	C2'-C1'	-3.27	1.48	1.53
2	B	1263	GDP	O4'-C4'	3.08	1.51	1.45
2	A	1262	GDP	PB-O3B	2.57	1.64	1.54
2	A	1262	GDP	C5'-C4'	2.46	1.59	1.51
2	A	1262	GDP	C2'-C3'	-2.09	1.47	1.53

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	1262	GDP	C2-N3-C4	8.10	124.61	115.36
2	B	1263	GDP	N3-C2-N1	-7.54	117.16	127.22
2	B	1263	GDP	C2-N3-C4	7.00	123.35	115.36
2	A	1262	GDP	N3-C2-N1	-6.73	118.24	127.22
2	B	1263	GDP	N2-C2-N1	6.40	127.20	117.25
2	A	1262	GDP	C5-C6-N1	-5.82	115.47	123.43
2	A	1262	GDP	C4-C5-N7	-5.48	103.69	109.40
2	A	1262	GDP	PA-O3A-PB	-4.43	117.64	132.83
2	A	1262	GDP	N2-C2-N1	3.69	123.00	117.25
2	A	1262	GDP	O4'-C4'-C5'	-3.38	98.27	109.37
2	B	1263	GDP	C5'-C4'-C3'	-3.25	103.01	115.18
2	B	1263	GDP	C4-C5-N7	-3.24	106.03	109.40
2	A	1262	GDP	O2B-PB-O1B	3.24	123.35	110.68
2	A	1262	GDP	C6-N1-C2	3.13	120.90	115.93
2	B	1263	GDP	O2B-PB-O1B	2.97	122.30	110.68
2	B	1263	GDP	O3B-PB-O1B	-2.94	99.17	110.68
2	B	1263	GDP	C5-C6-N1	-2.89	119.47	123.43
2	A	1262	GDP	O3B-PB-O3A	2.81	114.05	104.64
2	B	1263	GDP	C3'-C2'-C1'	2.77	105.14	100.98
2	B	1263	GDP	O3'-C3'-C4'	-2.73	103.16	111.05
2	A	1262	GDP	C3'-C2'-C1'	2.52	104.77	100.98
2	B	1263	GDP	O3B-PB-O3A	2.48	112.95	104.64
2	A	1262	GDP	C5'-C4'-C3'	-2.30	106.55	115.18
2	A	1262	GDP	O2B-PB-O3A	-2.20	97.26	104.64
2	A	1262	GDP	O3B-PB-O2B	-2.00	99.99	107.64
2	B	1263	GDP	O4'-C1'-C2'	-2.00	104.00	106.93

There are no chirality outliers.

All (3) torsion outliers are listed below:

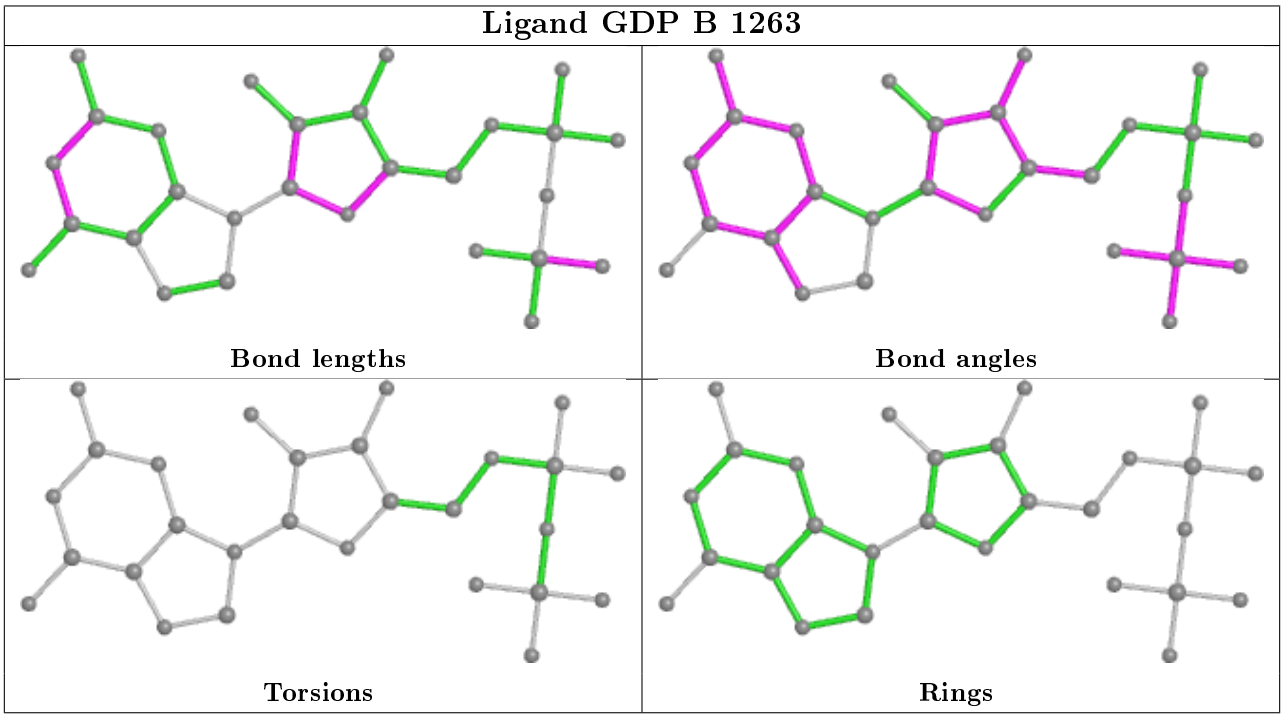
Mol	Chain	Res	Type	Atoms
2	A	1262	GDP	C5'-O5'-PA-O3A
2	A	1262	GDP	C3'-C4'-C5'-O5'
2	A	1262	GDP	C5'-O5'-PA-O2A

There are no ring outliers.

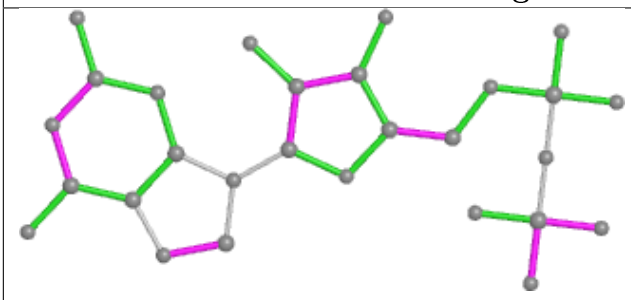
1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	1263	GDP	1	0

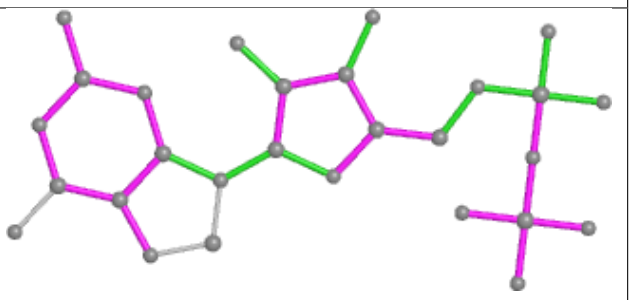
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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



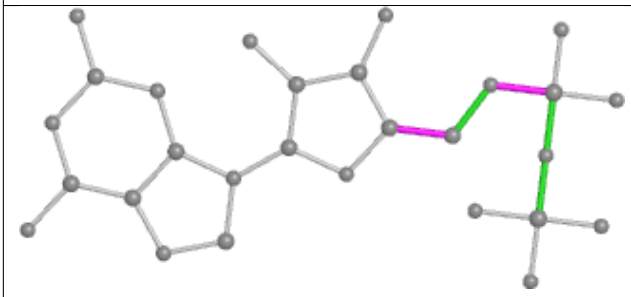
Ligand GDP A 1262



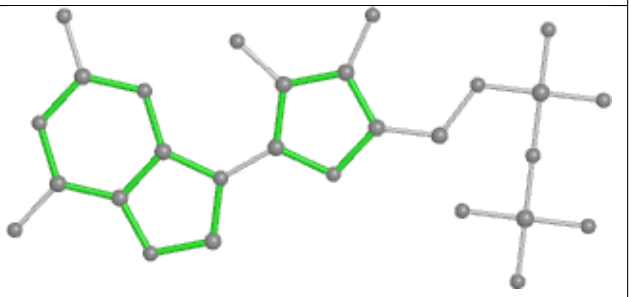
Bond lengths



Bond angles



Torsions



Rings

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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, 95^{th} 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	261/267 (97%)	-0.31	5 (1%) 66 74	27, 44, 75, 88	0
1	B	257/267 (96%)	-0.02	14 (5%) 25 32	33, 53, 93, 101	0
All	All	518/534 (97%)	-0.17	19 (3%) 41 50	27, 49, 88, 101	0

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	36	VAL	6.0
1	B	236	GLU	4.9
1	B	37	THR	4.7
1	B	220	GLY	4.3
1	A	66	ILE	3.9
1	B	234	SER	3.9
1	B	35	GLY	3.8
1	B	34	ALA	3.5
1	B	172	LEU	3.3
1	A	235	ASP	3.2
1	A	70	THR	3.0
1	A	65	THR	2.8
1	B	38	VAL	2.8
1	B	221	ASP	2.7
1	B	262	THR	2.5
1	B	93	VAL	2.5
1	A	67	SER	2.3
1	B	235	ASP	2.2
1	B	171	ASP	2.1

6.2 Non-standard residues in protein, DNA, RNA chains

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)

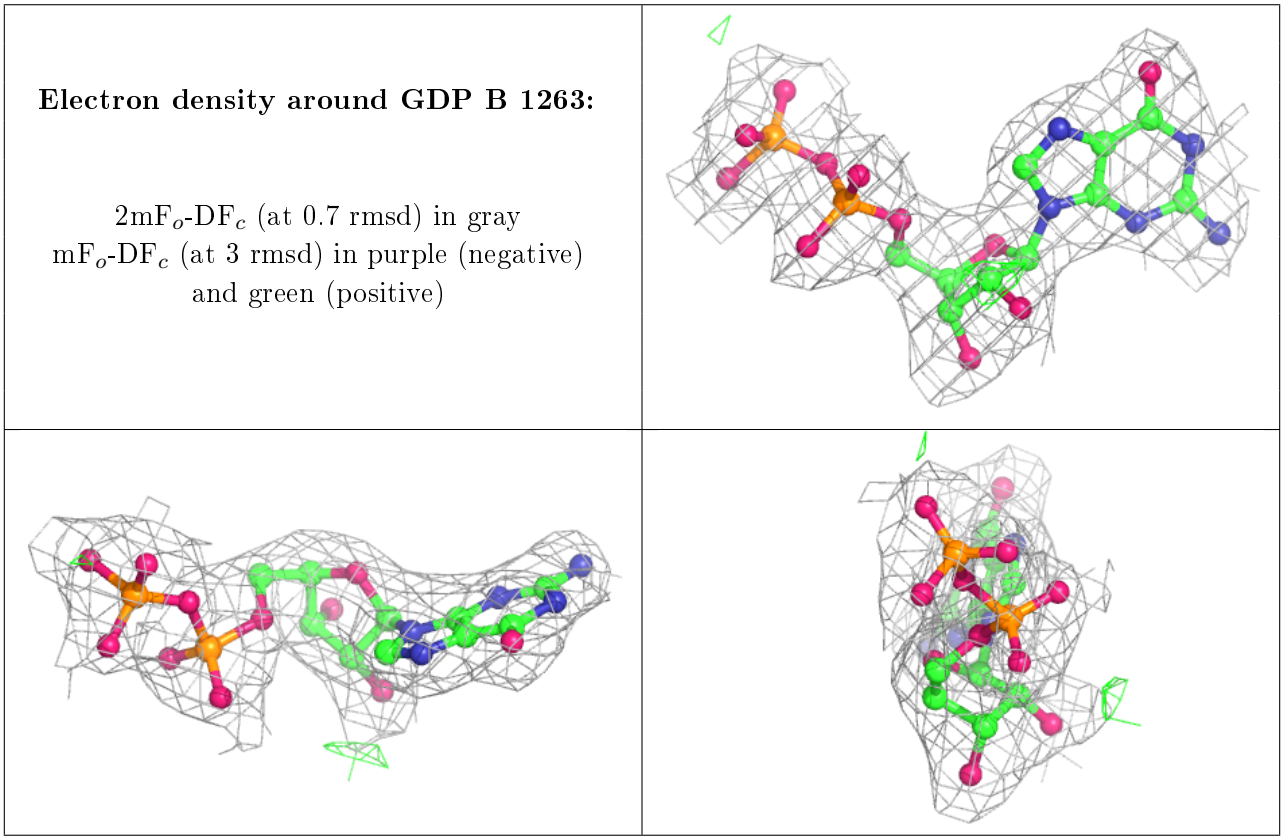
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' 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	GDP	B	1263	28/28	0.97	0.10	33,35,38,40	0
2	GDP	A	1262	28/28	0.97	0.10	28,32,35,37	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.

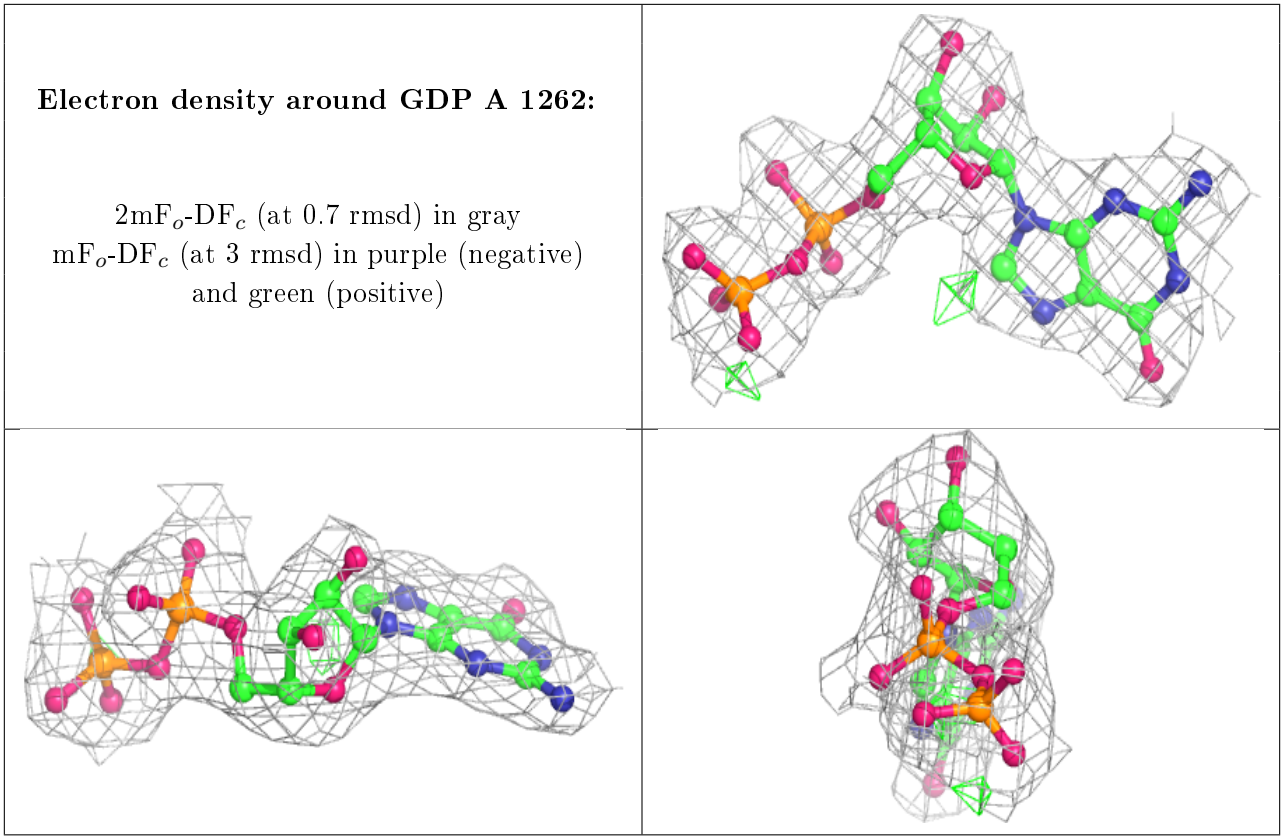
Electron density around GDP B 1263:

$2mF_o$-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 1262:

$2mF_o$-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.