Full wwPDB X-ray Structure Validation Report (i)

Apr 7, 2022 – 01:14 PM EDT

PDB ID	:	1FTN
Title	:	CRYSTAL STRUCTURE OF THE HUMAN RHOA/GDP COMPLEX
Authors	:	Wei, Y.; Zhang, Y.; Derewenda, U.; Liu, X.; Minor, W.; Nakamoto, R.K.; Somlyo, A.V.; Somlyo, A.P.; Derewenda, Z.S.
Deposited on	:	1997-03-13
Resolution	:	2.10 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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

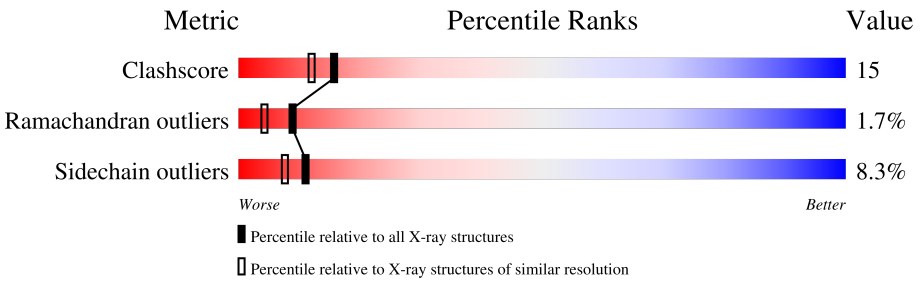
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.27

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 2.10 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))
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)

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 <=5%

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	193	

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 1653 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 TRANSFORMING PROTEIN RHOA (H12).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	177	1405	886	239	270	10	26	0	0

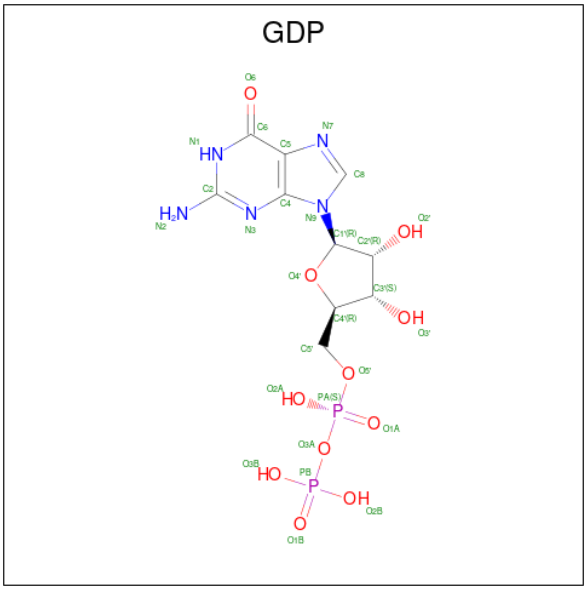
There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	25	ASN	PHE	engineered mutation	UNP P61586

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Mg		
2	A	1	1	1	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

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	A	219	219	219	0	0

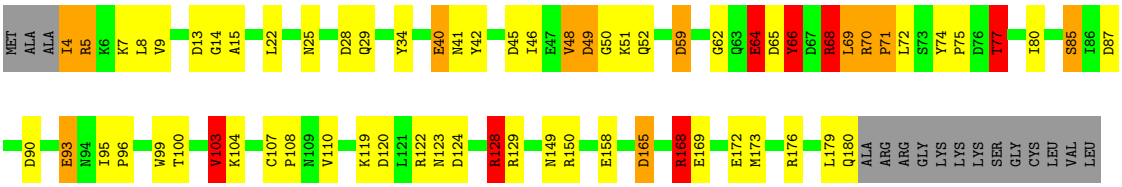
3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. 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. 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.

Note EDS was not executed.

• Molecule 1: TRANSFORMING PROTEIN RHOA (H12)

Chain A:



4 Data and refinement statistics (i)

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	31.92Å 65.96Å 83.37Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	8.00 – 2.10	Depositor
% Data completeness (in resolution range)	(Not available) (8.00-2.10)	Depositor
R_{merge}	0.04	Depositor
R_{sym}	0.04	Depositor
Refinement program	SHELX-90	Depositor
R, R_{free}	0.186 , 0.304	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	1653	wwPDB-VP
Average B, all atoms (Å^2)	31.0	wwPDB-VP

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: MG, 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	1.00	2/1432 (0.1%)	2.12	48/1937 (2.5%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	5

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	93	GLU	CD-OE2	5.72	1.31	1.25
1	A	93	GLU	CG-CD	5.37	1.60	1.51

All (48) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	122	ARG	NE-CZ-NH2	-16.29	112.16	120.30
1	A	176	ARG	NE-CZ-NH2	-15.44	112.58	120.30
1	A	150	ARG	NE-CZ-NH2	-15.29	112.65	120.30
1	A	168	ARG	CD-NE-CZ	15.06	144.68	123.60
1	A	5	ARG	NE-CZ-NH2	-14.70	112.95	120.30
1	A	5	ARG	CD-NE-CZ	14.56	143.99	123.60
1	A	5	ARG	NE-CZ-NH1	14.11	127.36	120.30
1	A	176	ARG	NE-CZ-NH1	13.16	126.88	120.30
1	A	68	ARG	NE-CZ-NH1	13.15	126.87	120.30
1	A	87	ASP	CB-CG-OD2	11.15	128.34	118.30
1	A	122	ARG	NH1-CZ-NH2	9.72	130.09	119.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	34	TYR	CB-CG-CD1	-8.99	115.61	121.00
1	A	124	ASP	CB-CG-OD2	8.78	126.21	118.30
1	A	40	GLU	OE1-CD-OE2	8.63	133.66	123.30
1	A	28	ASP	CB-CG-OD1	7.80	125.32	118.30
1	A	165	ASP	CB-CG-OD2	7.69	125.22	118.30
1	A	129	ARG	NE-CZ-NH2	-7.29	116.66	120.30
1	A	34	TYR	CB-CG-CD2	7.00	125.20	121.00
1	A	59	ASP	CB-CG-OD1	6.96	124.57	118.30
1	A	68	ARG	NH1-CZ-NH2	-6.77	111.96	119.40
1	A	129	ARG	CD-NE-CZ	6.60	132.84	123.60
1	A	45	ASP	CB-CG-OD1	6.43	124.09	118.30
1	A	176	ARG	CD-NE-CZ	6.43	132.60	123.60
1	A	42	TYR	CB-CG-CD2	-6.42	117.15	121.00
1	A	103	VAL	CA-CB-CG2	6.33	120.39	110.90
1	A	59	ASP	CB-CG-OD2	-6.12	112.79	118.30
1	A	150	ARG	NE-CZ-NH1	6.11	123.35	120.30
1	A	158	GLU	OE1-CD-OE2	-5.95	116.16	123.30
1	A	4	ILE	N-CA-CB	5.83	124.21	110.80
1	A	90	ASP	N-CA-CB	5.82	121.08	110.60
1	A	13	ASP	CB-CG-OD1	5.64	123.38	118.30
1	A	93	GLU	OE1-CD-OE2	-5.58	116.60	123.30
1	A	66	TYR	CB-CG-CD1	5.57	124.34	121.00
1	A	49	ASP	CB-CG-OD2	-5.54	113.31	118.30
1	A	65	ASP	CA-C-N	5.50	129.30	117.20
1	A	128	ARG	CD-NE-CZ	5.50	131.29	123.60
1	A	65	ASP	CA-C-O	-5.45	108.65	120.10
1	A	46	ILE	CB-CG1-CD1	5.45	129.15	113.90
1	A	77	THR	CA-CB-OG1	5.43	120.40	109.00
1	A	70	ARG	NE-CZ-NH2	5.35	122.97	120.30
1	A	122	ARG	NE-CZ-NH1	-5.21	117.69	120.30
1	A	180	GLN	CB-CA-C	-5.19	100.03	110.40
1	A	103	VAL	CA-CB-CG1	5.15	118.63	110.90
1	A	85	SER	CB-CA-C	-5.12	100.37	110.10
1	A	120	ASP	CB-CG-OD2	5.06	122.85	118.30
1	A	65	ASP	C-N-CA	5.06	134.35	121.70
1	A	180	GLN	N-CA-C	5.03	124.58	111.00
1	A	128	ARG	CB-CG-CD	5.01	124.64	111.60

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	48	VAL	Peptide
1	A	49	ASP	Peptide
1	A	50	GLY	Peptide
1	A	62	GLY	Mainchain
1	A	64	GLU	Peptide

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	1405	0	1396	41	0
2	A	1	0	0	0	0
3	A	28	0	12	0	0
4	A	219	0	0	16	1
All	All	1653	0	1408	41	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 15.

All (41) 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:107:CYS:HB3	1:A:110:VAL:HG13	1.56	0.87
1:A:40:GLU:HG2	4:A:398:HOH:O	1.86	0.75
1:A:128:ARG:HH11	1:A:128:ARG:HG2	1.54	0.71
1:A:149:ASN:ND2	4:A:392:HOH:O	2.23	0.71
1:A:66:TYR:HA	4:A:370:HOH:O	1.90	0.70
1:A:68:ARG:NH2	4:A:463:HOH:O	2.26	0.68
1:A:66:TYR:OH	4:A:522:HOH:O	2.09	0.66
1:A:4:ILE:HG22	1:A:52:GLN:O	1.96	0.65
1:A:68:ARG:HD3	4:A:464:HOH:O	1.96	0.65
1:A:107:CYS:HB3	1:A:110:VAL:CG1	2.27	0.63
1:A:128:ARG:HG2	1:A:128:ARG:NH1	2.13	0.62
1:A:95:ILE:HB	1:A:96:PRO:HD3	1.81	0.61
1:A:168:ARG:HG2	4:A:403:HOH:O	2.00	0.60
1:A:25:ASN:ND2	4:A:491:HOH:O	2.33	0.60
1:A:7:LYS:HE2	1:A:77:THR:HG22	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:72:LEU:HD22	1:A:72:LEU:H	1.72	0.54
1:A:74:TYR:N	1:A:75:PRO:CD	2.72	0.53
1:A:128:ARG:HG3	4:A:443:HOH:O	2.08	0.52
1:A:119:LYS:NZ	1:A:165:ASP:OD2	2.26	0.51
1:A:100:THR:O	1:A:104:LYS:HG2	2.13	0.48
1:A:168:ARG:HE	1:A:168:ARG:CA	2.27	0.48
1:A:169:GLU:OE1	1:A:169:GLU:N	2.45	0.47
1:A:5:ARG:HD2	4:A:509:HOH:O	2.14	0.47
1:A:93:GLU:O	1:A:96:PRO:HD2	2.14	0.47
1:A:9:VAL:HB	1:A:80:ILE:HD13	1.97	0.46
1:A:99:TRP:O	1:A:103:VAL:HG13	2.16	0.46
1:A:22:LEU:CD1	1:A:59:ASP:HB2	2.46	0.44
1:A:40:GLU:HA	4:A:386:HOH:O	2.17	0.43
1:A:74:TYR:O	1:A:77:THR:HG23	2.19	0.43
1:A:172:GLU:HB3	4:A:485:HOH:O	2.19	0.43
1:A:95:ILE:HB	1:A:96:PRO:CD	2.49	0.42
1:A:70:ARG:HB3	1:A:71:PRO:HD3	2.01	0.42
1:A:108:PRO:HG3	4:A:503:HOH:O	2.19	0.42
1:A:8:LEU:HD23	1:A:8:LEU:C	2.41	0.41
1:A:69:LEU:HA	1:A:72:LEU:HD23	2.01	0.41
1:A:14:GLY:O	1:A:15:ALA:HB3	2.21	0.41
1:A:173:MET:HE1	4:A:459:HOH:O	2.21	0.41
1:A:66:TYR:CA	4:A:370:HOH:O	2.59	0.41
1:A:68:ARG:HB2	4:A:525:HOH:O	2.21	0.40
1:A:51:LYS:CG	1:A:52:GLN:H	2.33	0.40
1:A:9:VAL:HG21	1:A:77:THR:HG21	2.02	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:368:HOH:O	4:A:422:HOH:O[4_455]	2.06	0.14

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	175/193 (91%)	164 (94%)	8 (5%)	3 (2%)	9 4

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	41	ASN
1	A	66	TYR
1	A	64	GLU

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	156/167 (93%)	143 (92%)	13 (8%)	11 7

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	29	GLN
1	A	48	VAL
1	A	64	GLU
1	A	68	ARG
1	A	69	LEU
1	A	71	PRO
1	A	77	THR
1	A	85	SER
1	A	103	VAL
1	A	123	ASN
1	A	128	ARG
1	A	168	ARG
1	A	179	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 2 ligands modelled in this entry, 1 is monoatomic - leaving 1 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	A	200	2	24,30,30	1.64	5 (20%)	31,47,47	2.34	9 (29%)

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	A	200	2	-	0/12/32/32	0/3/3/3

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	200	GDP	C5-C4	-4.10	1.30	1.40
3	A	200	GDP	C4-N3	-2.91	1.31	1.35
3	A	200	GDP	PB-O1B	2.64	1.59	1.50
3	A	200	GDP	C2-N1	2.64	1.40	1.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	200	GDP	C6-N1	2.34	1.37	1.33

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	A	200	GDP	C5-C6-N1	-7.26	113.50	123.43
3	A	200	GDP	N3-C2-N1	-4.48	121.24	127.22
3	A	200	GDP	C2-N1-C6	3.87	122.08	115.93
3	A	200	GDP	C4-C5-N7	3.76	113.31	109.40
3	A	200	GDP	O2'-C2'-C1'	-3.24	98.87	110.85
3	A	200	GDP	O4'-C1'-C2'	-3.23	102.21	106.93
3	A	200	GDP	O3B-PB-O1B	2.87	121.93	110.68
3	A	200	GDP	N2-C2-N3	2.85	122.44	117.79
3	A	200	GDP	O2B-PB-O3A	-2.54	96.10	104.64

There are no chirality outliers.

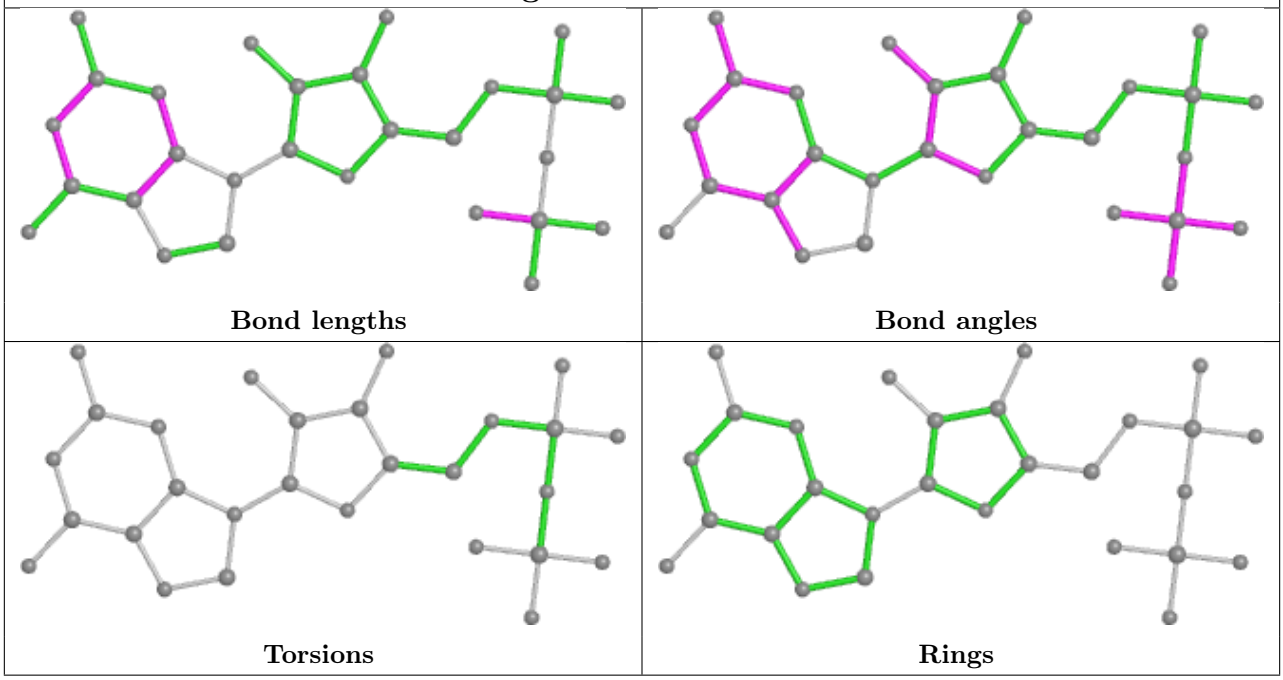
There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less 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 200



5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data (i)

6.1 Protein, DNA and RNA chains (i)

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates (i)

EDS was not executed - this section is therefore empty.

6.4 Ligands (i)

EDS was not executed - this section is therefore empty.

6.5 Other polymers (i)

EDS was not executed - this section is therefore empty.