Full wwPDB X-ray Structure Validation Report (i)

Jun 12, 2024 – 05:38 PM EDT

PDB ID	:	1EU1
Title	:	THE CRYSTAL STRUCTURE OF RHODOBACTER SPHAEROIDES DIMETHYLSULFOXIDE REDUCTASE REVEALS TWO DISTINCT MOLYBDENUM COORDINATION ENVIRONMENTS.
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Deposited on	:	2000-04-13
Resolution	:	1.30 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 types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

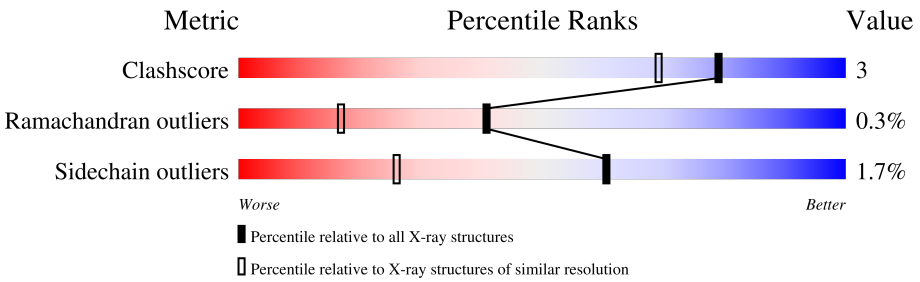
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 1.30 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	1101 (1.30-1.30)
Ramachandran outliers	138981	1058 (1.30-1.30)
Sidechain outliers	138945	1058 (1.30-1.30)

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 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	780	91% 7% • •

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	GLC	A	2002	X	-	-	-
2	GLC	A	2003	X	-	-	-
2	GLC	A	2004	X	-	-	-
7	O	A	3002	-	-	X	-

2 Entry composition (i)

There are 9 unique types of molecules in this entry. The entry contains 7307 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.

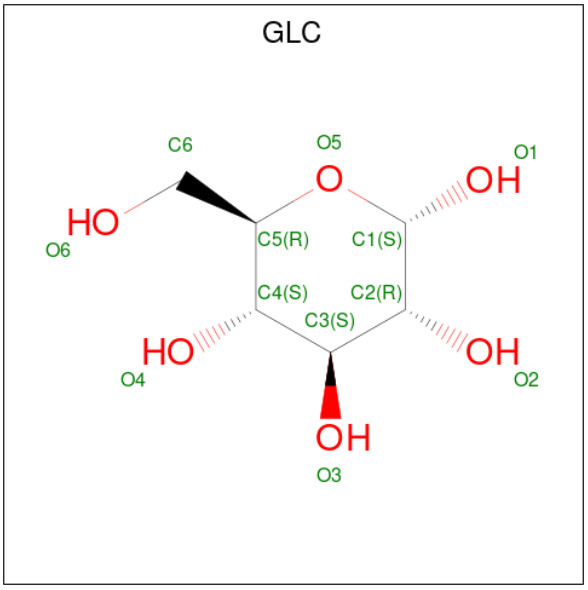
- Molecule 1 is a protein called DIMETHYL SULFOXIDE REDUCTASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	767	6048	3829	1034	1150	35	0	34	0

There is a discrepancy between the modelled and reference sequences:

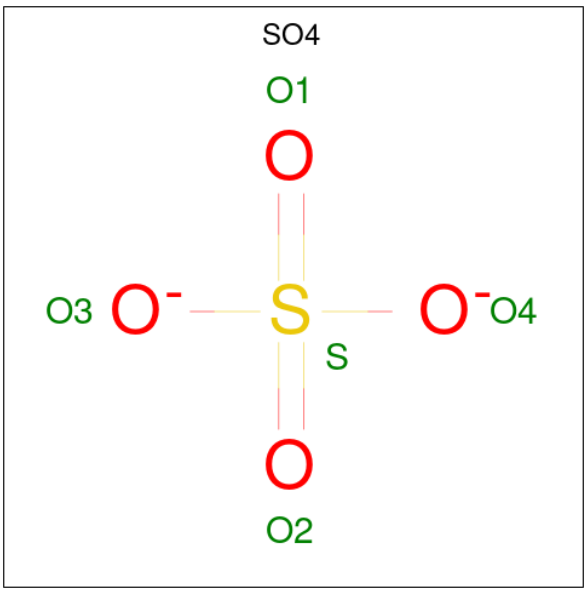
Chain	Residue	Modelled	Actual	Comment	Reference
A	219	CSD	CYS	modified residue	UNP Q57366

- Molecule 2 is alpha-D-glucopyranose (three-letter code: GLC) (formula: $C_6H_{12}O_6$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
2	A	1	12	6	6	0	0
2	A	1	12	6	6	0	0
2	A	1	12	6	6	0	0

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O_4S).

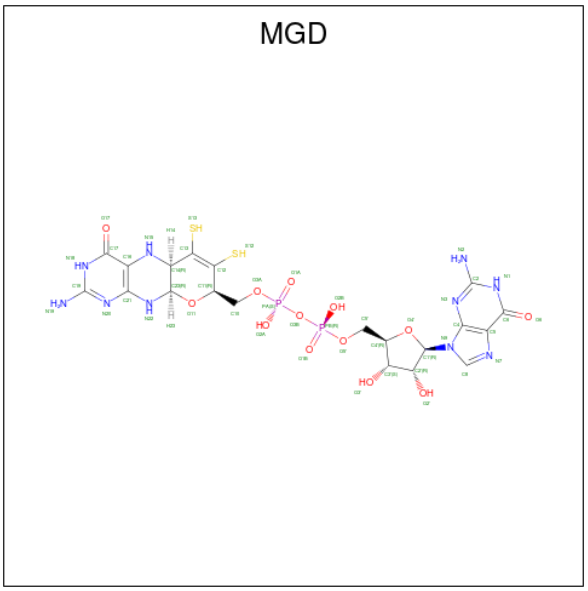


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total O S 5 4 1	0	0

- Molecule 4 is CADMIUM ION (three-letter code: CD) (formula: Cd).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total Cd 1 1	0	0

- Molecule 5 is 2-AMINO-5,6-DIMERCAPTO-7-METHYL-3,7,8A,9-TETRAHYDRO-8-OXA-1,3,9,10-TETRAAZA-ANTHRACEN-4-ONE GUANOSINE DINUCLEOTIDE (three-letter code: MGD) (formula: $C_{20}H_{26}N_{10}O_{13}P_2S_2$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
			Total	C	N	O	P	S		
5	A	1	75	30	15	22	4	4	0	1
			Total	C	N	O	P	S		
5	A	1	83	35	15	25	4	4	0	1

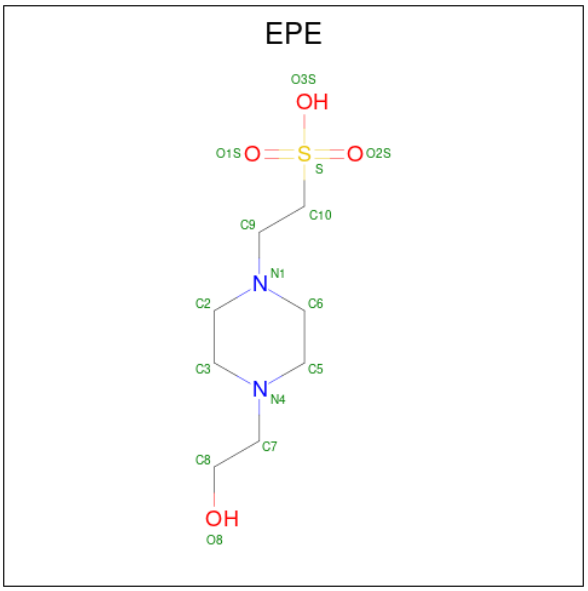
- Molecule 6 is MOLYBDENUM(VI) ION (three-letter code: 6MO) (formula: Mo).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Mo		
6	A	1	2	2	0	1

- Molecule 7 is OXYGEN ATOM (three-letter code: O) (formula: O).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
7	A	2	3	3	0	1

- Molecule 8 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (three-letter code: EPE) (formula: $C_8H_{18}N_2O_4S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	S		
8	A	1	15	8	2	4	1	0	0

- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
9	A	1039	1039	1039	0	0

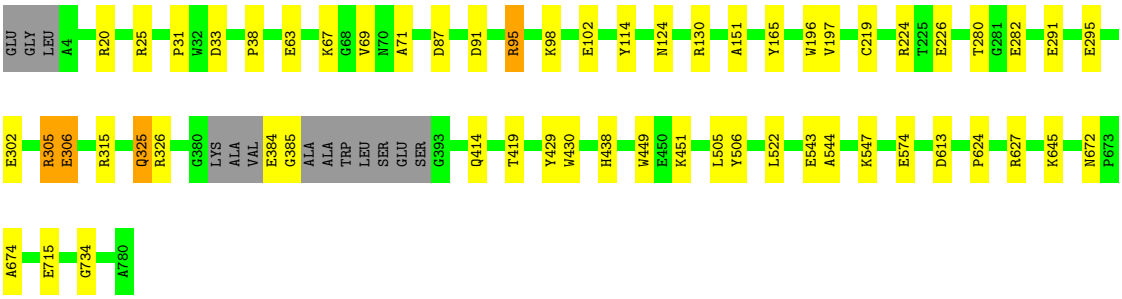
3 Residue-property plots (i)

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: DIMETHYL SULFOXIDE REDUCTASE

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 2	Depositor
Cell constants a, b, c, α, β, γ	102.18Å 141.72Å 59.81Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 1.30	Depositor
% Data completeness (in resolution range)	89.8 (10.00-1.30)	Depositor
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	REFMAC	Depositor
R, R_{free}	0.121 , 0.145	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	7307	wwPDB-VP
Average B, all atoms (Å^2)	15.0	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: O, 6MO, EPE, SO4, CD, GLC, MGD, CSD

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	0.52	0/6328	1.11	24/8580 (0.3%)

There are no bond length outliers.

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	25	ARG	CD-NE-CZ	14.89	144.44	123.60
1	A	130[A]	ARG	NE-CZ-NH1	-11.92	114.34	120.30
1	A	130[B]	ARG	NE-CZ-NH1	-11.92	114.34	120.30
1	A	305[A]	ARG	NE-CZ-NH2	-11.49	114.56	120.30
1	A	305[B]	ARG	NE-CZ-NH2	-11.49	114.56	120.30
1	A	305[A]	ARG	NE-CZ-NH1	10.55	125.57	120.30
1	A	305[B]	ARG	NE-CZ-NH1	10.55	125.57	120.30
1	A	87	ASP	CB-CG-OD2	9.58	126.93	118.30
1	A	25	ARG	NE-CZ-NH1	8.31	124.45	120.30
1	A	91	ASP	CB-CG-OD1	7.88	125.39	118.30
1	A	130[A]	ARG	CD-NE-CZ	7.84	134.58	123.60
1	A	130[B]	ARG	CD-NE-CZ	7.84	134.58	123.60
1	A	130[A]	ARG	NE-CZ-NH2	7.51	124.05	120.30
1	A	130[B]	ARG	NE-CZ-NH2	7.51	124.05	120.30
1	A	95[A]	ARG	NE-CZ-NH1	7.38	123.99	120.30
1	A	95[B]	ARG	NE-CZ-NH1	7.38	123.99	120.30
1	A	87	ASP	CB-CG-OD1	-7.32	111.71	118.30
1	A	219	CSD	O-C-N	-7.21	111.16	122.70
1	A	95[A]	ARG	CD-NE-CZ	6.59	132.83	123.60
1	A	95[B]	ARG	CD-NE-CZ	6.59	132.83	123.60
1	A	613	ASP	CB-CG-OD1	5.81	123.53	118.30
1	A	326	ARG	NE-CZ-NH1	5.79	123.20	120.30
1	A	315	ARG	NE-CZ-NH1	5.54	123.07	120.30
1	A	224	ARG	CD-NE-CZ	5.51	131.32	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	6048	0	5816	32	1
2	A	36	0	36	0	1
3	A	5	0	0	0	0
4	A	1	0	0	0	0
5	A	158	0	62	0	0
6	A	2	0	0	1	0
7	A	3	0	0	3	0
8	A	15	0	18	2	0
9	A	1039	0	0	6	0
All	All	7307	0	5932	34	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 3.

All (34) 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:114[B]:TYR:OH	7:A:3002:O:O	1.59	1.19
1:A:124:ASN:HD21	1:A:384:GLU:HA	1.28	0.94
1:A:544:ALA:HA	1:A:547[A]:LYS:HE2	1.53	0.90
1:A:196:TRP:HE1	1:A:385:GLY:HA3	1.38	0.88
1:A:151:ALA:H	1:A:325:GLN:HE21	1.35	0.71
1:A:124:ASN:ND2	1:A:384:GLU:HA	2.04	0.71
1:A:291:GLU:OE2	1:A:305[A]:ARG:NH2	2.23	0.70
1:A:124:ASN:HD21	1:A:384:GLU:CA	2.06	0.61
1:A:196:TRP:NE1	1:A:385:GLY:HA3	2.13	0.61
6:A:3000[A]:6MO:MO	7:A:3002:O:O	1.71	0.61
1:A:672:ASN:HD22	1:A:674:ALA:H	1.50	0.59
1:A:98:LYS:HE3	1:A:522:LEU:HD23	1.88	0.55
1:A:196:TRP:HZ2	1:A:384:GLU:N	2.05	0.54
1:A:98:LYS:CE	1:A:522:LEU:HD23	2.38	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:102:GLU:HG3	9:A:3588:HOH:O	2.09	0.51
1:A:505[B]:LEU:HD13	1:A:506:TYR:CE1	2.48	0.49
1:A:63:GLU:HG3	1:A:71:ALA:HB1	1.96	0.47
1:A:414[A]:GLN:NE2	1:A:419:THR:OG1	2.34	0.45
1:A:280:THR:OG1	1:A:282[B]:GLU:HG3	2.15	0.45
1:A:672:ASN:ND2	1:A:674:ALA:H	2.13	0.45
1:A:645[B]:LYS:HD2	9:A:3745:HOH:O	2.17	0.45
1:A:196:TRP:HE1	1:A:385:GLY:CA	2.20	0.44
1:A:31:PRO:HB3	1:A:38:PRO:HB3	2.00	0.44
1:A:624:PRO:HG2	1:A:627:ARG:HB3	1.99	0.43
1:A:69:VAL:HG12	9:A:3989:HOH:O	2.18	0.43
1:A:543:GLU:O	1:A:547[A]:LYS:HG3	2.18	0.42
1:A:95[B]:ARG:HH11	1:A:95[B]:ARG:HG2	1.83	0.42
7:A:3001[B]:O:O	8:A:2001:EPE:H82	2.20	0.41
1:A:165:TYR:CZ	8:A:2001:EPE:H31	2.56	0.41
1:A:451[B]:LYS:HE2	9:A:3034:HOH:O	2.20	0.41
1:A:226[B]:GLU:HG2	9:A:3226:HOH:O	2.20	0.41
1:A:547[A]:LYS:HE3	9:A:3434:HOH:O	2.21	0.40
1:A:302:GLU:O	1:A:306[B]:GLU:HG3	2.22	0.40
1:A:20[B]:ARG:HH11	1:A:20[B]:ARG:HD2	1.77	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)
1:A:67[A]:LYS:NZ	2:A:2003:GLC:C5[1_554]	2.07	0.13

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	794/780 (102%)	774 (98%)	18 (2%)	2 (0%)	41 17

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	197	VAL
1	A	734	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	638/614 (104%)	627 (98%)	11 (2%)	60 26

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

Mol	Chain	Res	Type
1	A	33	ASP
1	A	295	GLU
1	A	306[A]	GLU
1	A	306[B]	GLU
1	A	325	GLN
1	A	429	TYR
1	A	430	TRP
1	A	438	HIS
1	A	449	TRP
1	A	574	GLU
1	A	715	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	241	GLN
1	A	325	GLN
1	A	458	GLN
1	A	606	ASN
1	A	672	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)

1 non-standard protein/DNA/RNA residue is 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	$\#\|Z\| > 2$
1	CSD	A	219	1	3,7,8	1.68	1 (33%)	1,8,10	2.07	1 (100%)

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
1	CSD	A	219	1	-	2/2/6/8	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	219	CSD	CB-SG	2.33	1.92	1.79

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	219	CSD	OD1-SG-CB	2.07	109.47	105.54

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	219	CSD	N-CA-CB-SG

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Mol	Chain	Res	Type	Atoms
1	A	219	CSD	CA-CB-SG-OD1

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 15 ligands modelled in this entry, 6 are monoatomic - leaving 9 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	$\#\|Z\| > 2$
2	GLC	A	2003	-	12,12,12	1.05	0	17,17,17	1.42	3 (17%)
5	MGD	A	1002[B]	6	41,52,52	1.94	15 (36%)	40,81,81	1.16	3 (7%)
5	MGD	A	1002[A]	-	41,52,52	1.97	14 (34%)	40,81,81	1.19	2 (5%)
3	SO4	A	2005	-	4,4,4	0.61	0	6,6,6	0.27	0
2	GLC	A	2004	-	12,12,12	1.11	1 (8%)	17,17,17	1.02	0
8	EPE	A	2001	-	15,15,15	1.98	3 (20%)	18,20,20	1.54	3 (16%)
2	GLC	A	2002	-	12,12,12	1.20	1 (8%)	17,17,17	0.95	1 (5%)
5	MGD	A	1001[B]	6	41,52,52	1.81	10 (24%)	40,81,81	1.08	3 (7%)
5	MGD	A	1001[A]	6	41,52,52	1.80	10 (24%)	40,81,81	1.07	3 (7%)

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	GLC	A	2003	-	1/1/5/5	0/2/22/22	0/1/1/1
5	MGD	A	1002[B]	6	-	0/18/66/66	0/6/6/6
5	MGD	A	1002[A]	-	-	4/18/66/66	0/6/6/6
2	GLC	A	2004	-	1/1/5/5	0/2/22/22	0/1/1/1
8	EPE	A	2001	-	-	1/9/19/19	0/1/1/1
2	GLC	A	2002	-	1/1/5/5	0/2/22/22	0/1/1/1
5	MGD	A	1001[B]	6	-	5/18/66/66	0/6/6/6
5	MGD	A	1001[A]	6	-	7/18/66/66	0/6/6/6

All (54) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	A	2001	EPE	O3S-S	5.62	1.67	1.47
5	A	1001[B]	MGD	C14-N15	5.44	1.52	1.46
5	A	1001[A]	MGD	C14-N15	5.29	1.52	1.46
5	A	1002[A]	MGD	C10-C11	4.81	1.58	1.52
5	A	1002[B]	MGD	C14-N15	4.35	1.51	1.46
5	A	1002[B]	MGD	C10-C11	4.26	1.57	1.52
5	A	1001[A]	MGD	C21-N22	4.01	1.39	1.35
5	A	1001[B]	MGD	C21-N22	4.00	1.39	1.35
5	A	1002[A]	MGD	C14-N15	3.95	1.51	1.46
5	A	1002[B]	MGD	O4'-C1'	3.95	1.46	1.41
5	A	1002[A]	MGD	C23-C14	3.81	1.56	1.53
5	A	1002[A]	MGD	O4'-C1'	3.76	1.46	1.41
8	A	2001	EPE	C10-S	3.72	1.82	1.77
5	A	1001[A]	MGD	O11-C11	3.64	1.48	1.43
5	A	1001[B]	MGD	O11-C11	3.60	1.48	1.43
5	A	1002[B]	MGD	C23-C14	3.38	1.56	1.53
5	A	1001[A]	MGD	O4'-C1'	3.16	1.45	1.41
5	A	1001[B]	MGD	O4'-C1'	3.16	1.45	1.41
5	A	1002[A]	MGD	C6-N1	2.93	1.42	1.37
5	A	1002[B]	MGD	C6-N1	2.93	1.42	1.37
5	A	1002[B]	MGD	O17-C17	2.87	1.29	1.23
5	A	1002[A]	MGD	O17-C17	2.85	1.29	1.23
5	A	1002[A]	MGD	C21-N22	2.77	1.38	1.35
5	A	1002[A]	MGD	C21-N20	2.73	1.40	1.36
5	A	1002[B]	MGD	C21-N20	2.71	1.40	1.36
5	A	1001[A]	MGD	C21-N20	2.59	1.40	1.36
5	A	1001[B]	MGD	C21-N20	2.59	1.40	1.36
5	A	1002[A]	MGD	C19-N19	2.58	1.40	1.34
8	A	2001	EPE	C7-N4	2.56	1.53	1.47
5	A	1002[B]	MGD	C19-N19	2.49	1.40	1.34
5	A	1002[B]	MGD	C21-N22	2.41	1.38	1.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	1002[B]	MGD	O11-C11	2.41	1.47	1.43
5	A	1002[A]	MGD	O11-C23	2.32	1.47	1.43
5	A	1002[A]	MGD	C2'-C3'	2.31	1.59	1.53
5	A	1001[B]	MGD	O17-C17	2.30	1.28	1.23
5	A	1001[A]	MGD	O17-C17	2.29	1.27	1.23
5	A	1002[A]	MGD	C2-N3	2.28	1.38	1.33
5	A	1002[B]	MGD	C2-N3	2.28	1.38	1.33
5	A	1002[A]	MGD	O11-C11	2.28	1.46	1.43
5	A	1002[B]	MGD	C2'-C3'	2.26	1.59	1.53
5	A	1001[A]	MGD	O6-C6	2.21	1.27	1.23
5	A	1001[B]	MGD	O6-C6	2.21	1.27	1.23
5	A	1002[B]	MGD	O11-C23	2.17	1.46	1.43
5	A	1001[A]	MGD	C6-N1	2.15	1.41	1.37
5	A	1001[B]	MGD	C6-N1	2.15	1.41	1.37
5	A	1001[A]	MGD	C3'-C4'	2.09	1.58	1.53
5	A	1001[B]	MGD	C3'-C4'	2.09	1.58	1.53
5	A	1002[B]	MGD	C17-N18	2.07	1.42	1.38
5	A	1001[A]	MGD	C17-N18	2.06	1.42	1.38
2	A	2004	GLC	O1-C1	2.04	1.46	1.39
2	A	2002	GLC	O1-C1	2.02	1.46	1.39
5	A	1001[B]	MGD	C17-N18	2.01	1.42	1.38
5	A	1002[A]	MGD	C2-N2	2.00	1.39	1.34
5	A	1002[B]	MGD	C2-N2	2.00	1.39	1.34

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
8	A	2001	EPE	O3S-S-O1S	-3.93	101.68	111.27
2	A	2003	GLC	C1-C2-C3	-3.16	103.75	110.31
5	A	1001[A]	MGD	O4'-C1'-C2'	-3.04	102.48	106.93
5	A	1001[B]	MGD	O4'-C1'-C2'	-3.04	102.48	106.93
5	A	1002[B]	MGD	O4'-C1'-C2'	-3.03	102.50	106.93
5	A	1002[A]	MGD	O11-C23-N22	-2.90	105.58	108.57
5	A	1001[B]	MGD	O11-C23-C14	2.78	110.82	108.96
8	A	2001	EPE	O1S-S-C10	2.73	110.21	106.92
5	A	1002[B]	MGD	O11-C23-N22	-2.66	105.83	108.57
2	A	2003	GLC	O1-C1-O5	-2.55	102.72	110.38
5	A	1002[A]	MGD	C19-N18-C17	-2.41	120.70	125.10
5	A	1002[B]	MGD	C19-N18-C17	-2.40	120.72	125.10
5	A	1001[A]	MGD	O11-C23-C14	2.16	110.41	108.96
2	A	2002	GLC	O1-C1-O5	-2.15	103.92	110.38
2	A	2003	GLC	O1-C1-C2	2.12	115.01	109.03

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
5	A	1001[A]	MGD	O6-C6-C5	2.02	128.32	124.37
5	A	1001[B]	MGD	O6-C6-C5	2.02	128.32	124.37
8	A	2001	EPE	O2S-S-O1S	2.02	120.95	113.95

All (3) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	A	2002	GLC	C1
2	A	2003	GLC	C1
2	A	2004	GLC	C1

All (17) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	1001[A]	MGD	PA-O3B-PB-O5'
5	A	1001[A]	MGD	C5'-O5'-PB-O1B
5	A	1001[A]	MGD	C5'-O5'-PB-O2B
5	A	1001[B]	MGD	PA-O3B-PB-O5'
5	A	1001[B]	MGD	C5'-O5'-PB-O1B
5	A	1001[B]	MGD	C5'-O5'-PB-O2B
5	A	1002[A]	MGD	O3A-C10-C11-O11
5	A	1002[A]	MGD	O3A-C10-C11-C12
8	A	2001	EPE	C10-C9-N1-C2
5	A	1001[B]	MGD	C5'-O5'-PB-O3B
5	A	1002[A]	MGD	C5'-O5'-PB-O3B
5	A	1001[A]	MGD	O3A-C10-C11-C12
5	A	1001[A]	MGD	O3A-C10-C11-O11
5	A	1001[A]	MGD	O4'-C4'-C5'-O5'
5	A	1001[A]	MGD	C5'-O5'-PB-O3B
5	A	1001[B]	MGD	O4'-C4'-C5'-O5'
5	A	1002[A]	MGD	C5'-O5'-PB-O1B

There are no ring outliers.

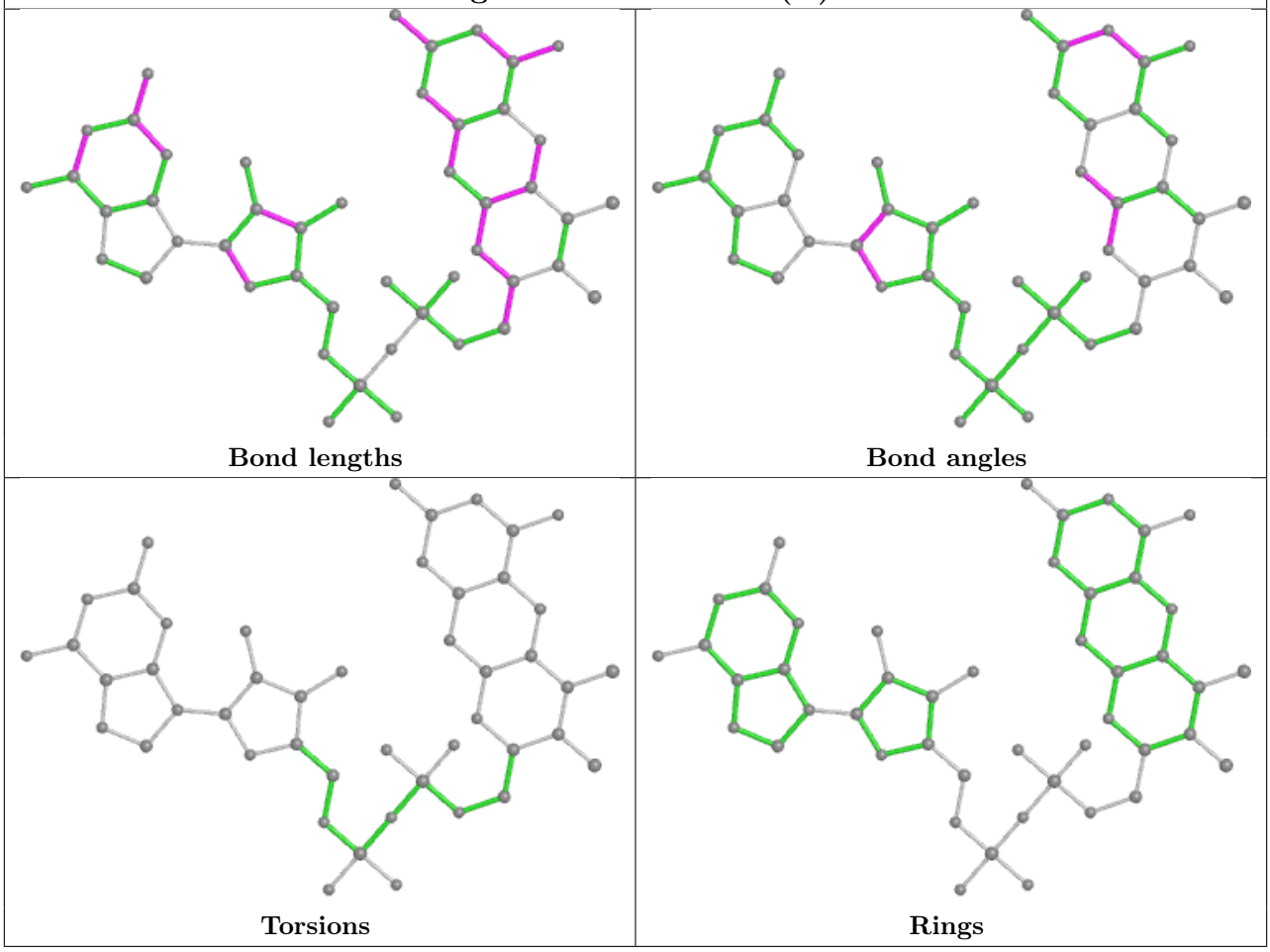
2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	2003	GLC	0	1
8	A	2001	EPE	2	0

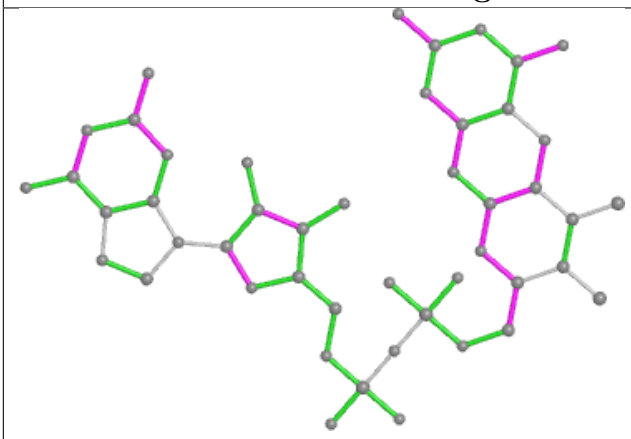
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will

also be included. For torsion angles, if less 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.

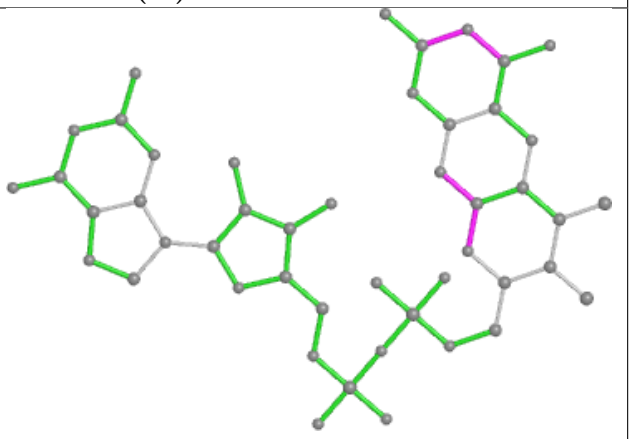
Ligand MGD A 1002 (B)



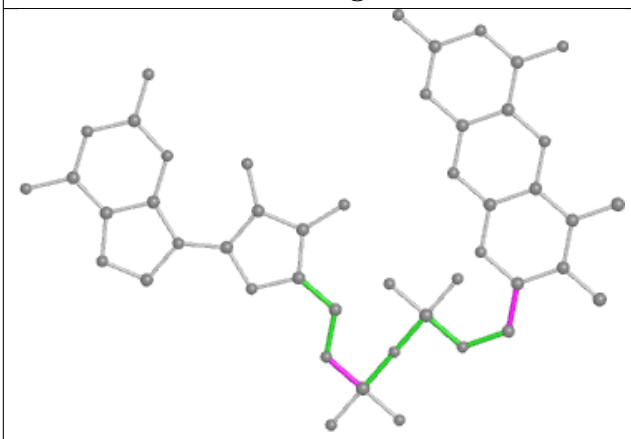
Ligand MGD A 1002 (A)



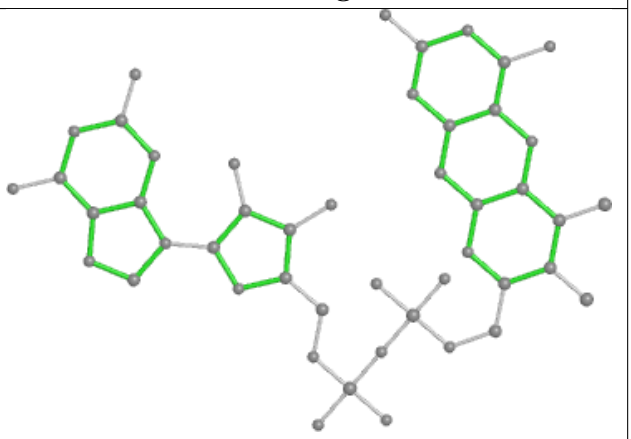
Bond lengths



Bond angles

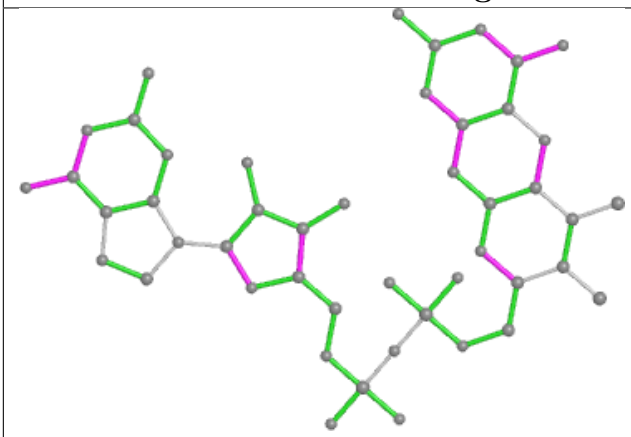


Torsions

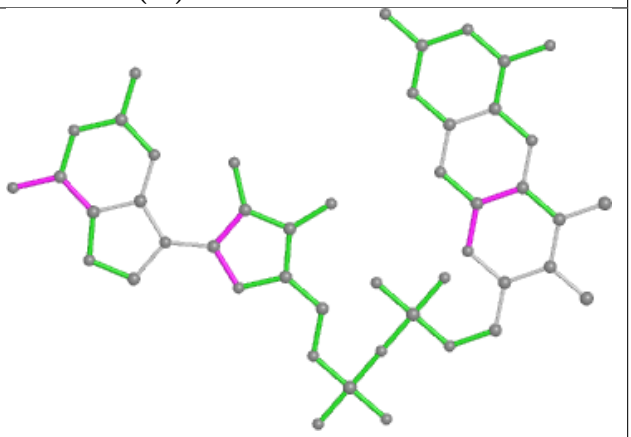


Rings

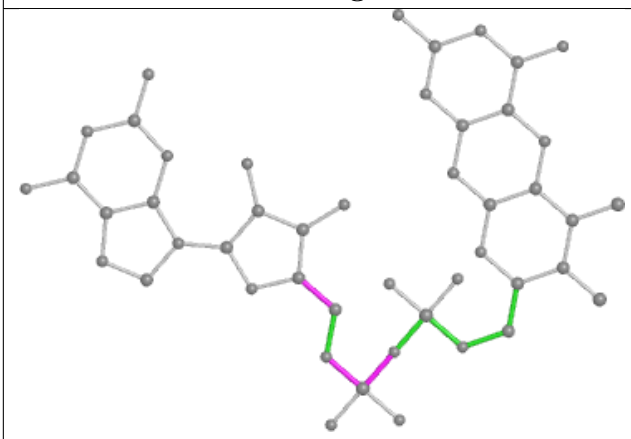
Ligand MGD A 1001 (B)



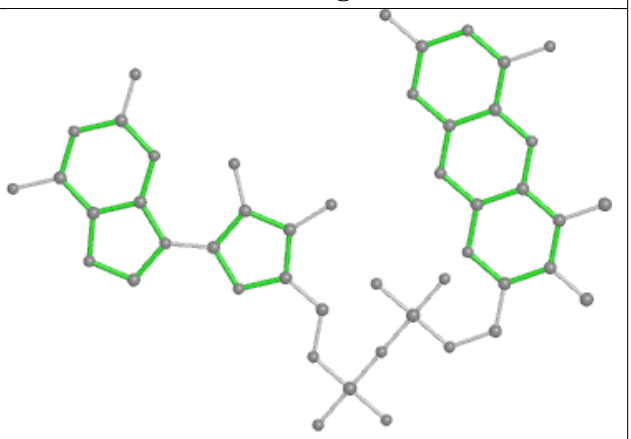
Bond lengths



Bond angles

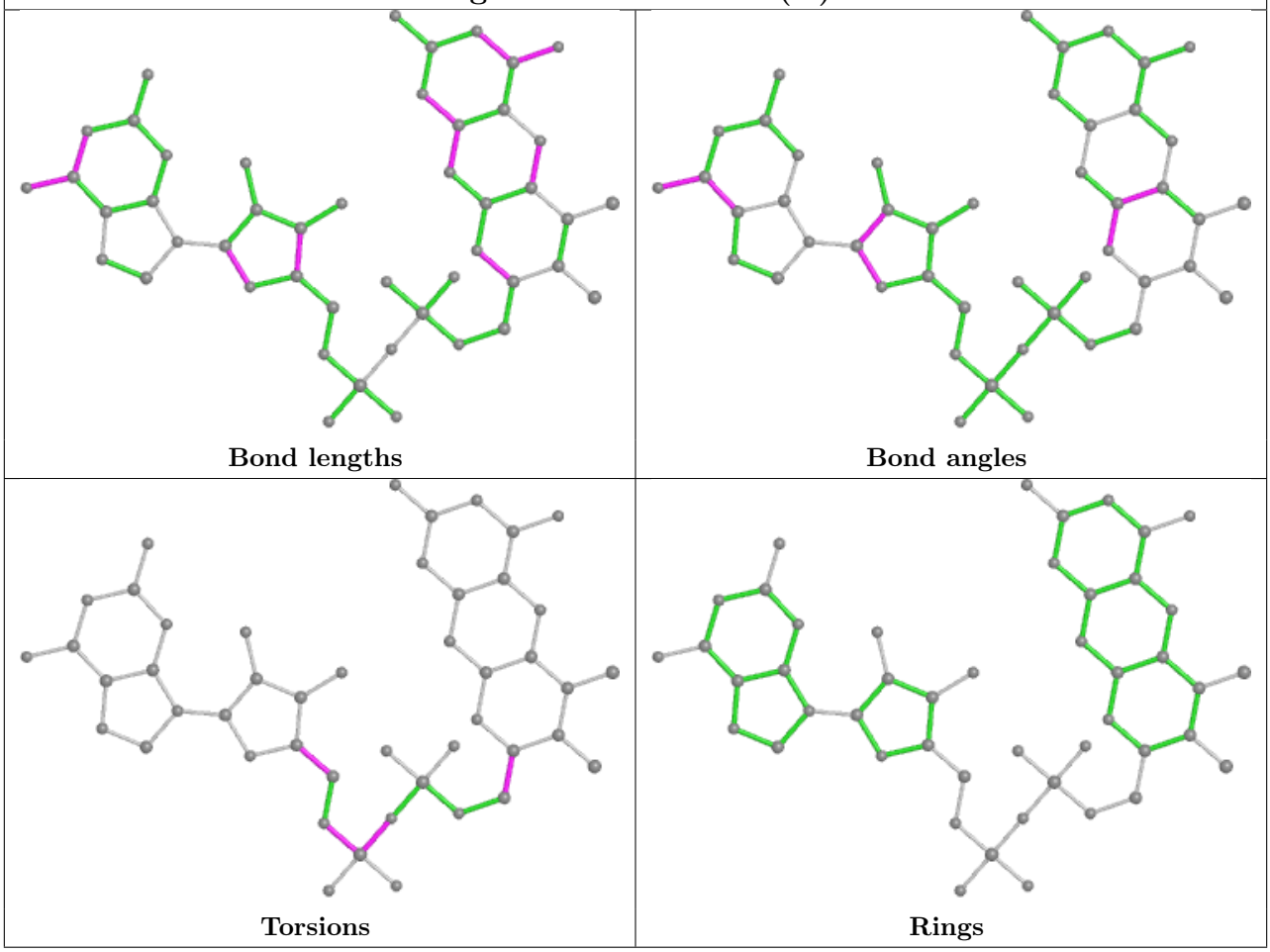


Torsions



Rings

Ligand MGD A 1001 (A)



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.