wwPDB X-ray Structure Validation Summary Report (i)

May 16, 2020 – 10:58 pm BST

PDB ID	:	1DFO
Title	:	CRYSTAL STRUCTURE AT 2.4 ANGSTROM RESOLUTION OF E. COLI SERINE HYDROXYMETHYLTRANSFERASE IN COMPLEX WITH GLYCINE AND 5-FORMYL TETRAHYDROFOLATE
Authors	:	Scarsdale, J.N.; Radaev, S.; Kazanina, G.; Schirch, V.; Wright, H.T.
Deposited on	:	1999-11-20
Resolution	:	2.40 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 (i)) were used in the production of this report:

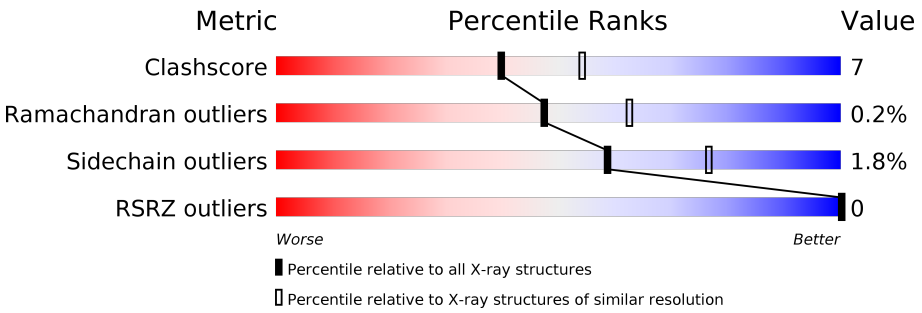
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.11
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (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 2.40 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)
RSRZ outliers	127900	3811 (2.40-2.40)

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	417	84% 14% •
1	B	417	84% 15% •
1	C	417	83% 16% •
1	D	417	84% 14% •

2 Entry composition (i)

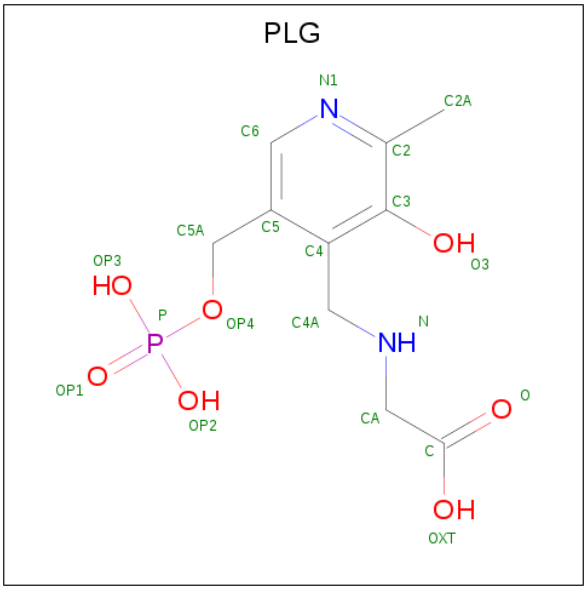
There are 4 unique types of molecules in this entry. The entry contains 13204 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 SERINE HYDROXYMETHYLTRANSFERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	416	3148	1995	541	598	14	0	0	0
1	B	417	3167	2006	544	602	15	0	0	0
1	C	417	3140	1989	539	597	15	0	0	0
1	D	416	3145	1993	539	599	14	0	0	0

- Molecule 2 is N-GLYCINE-[3-HYDROXY-2-METHYL-5-PHOSPHONOOXYMETHYL-PYRIDIN-4-YL-METHANE] (three-letter code: PLG) (formula: $C_{10}H_{15}N_2O_7P$).



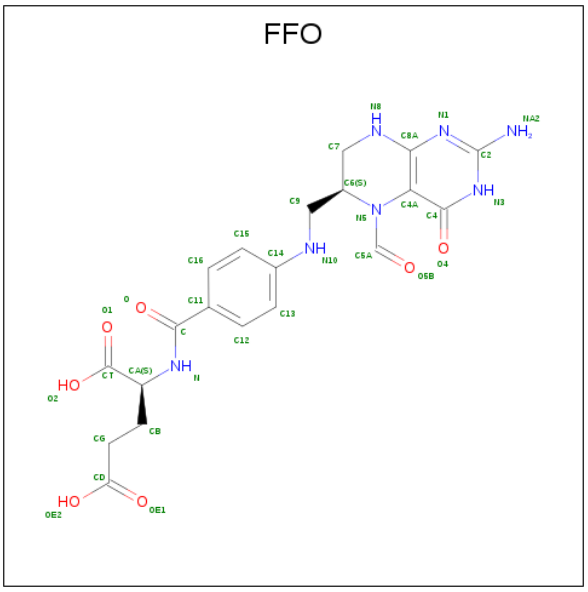
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
2	A	1	20	10	2	7	1	0	0
2	B	1	20	10	2	7	1	0	0

Continued on next page...

Continued from previous page...

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
2	C	1	20	10	2	7	1	0	0
			Total	C	N	O	P		
2	D	1	20	10	2	7	1	0	0

- Molecule 3 is N-[4-({[(6S)-2-amino-5-formyl-4-oxo-3,4,5,6,7,8-hexahydropteridin-6-yl]methyl }amino)benzoyl]-L-glutamic acid (three-letter code: FFO) (formula: $C_{20}H_{23}N_7O_7$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
3	A	1	34	20	7	7	0	0
			Total	C	N	O		
3	B	1	34	20	7	7	0	0
			Total	C	N	O		
3	C	1	34	20	7	7	0	0
			Total	C	N	O		
3	D	1	34	20	7	7	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	A	93	93	93	0	0
			Total	O		
4	B	94	94	94	0	0
			Total	O		
4	C	85	85	85	0	0

Continued on next page...

Continued from previous page...

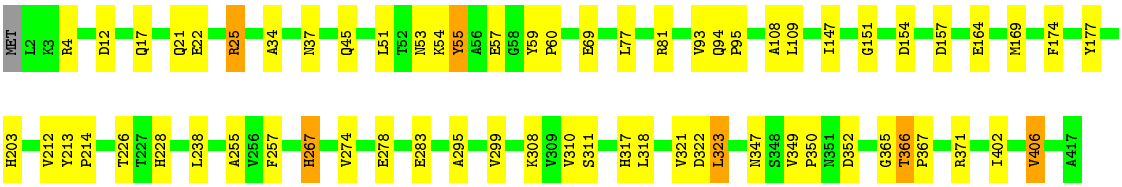
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	D	116	116	116	0	0

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.

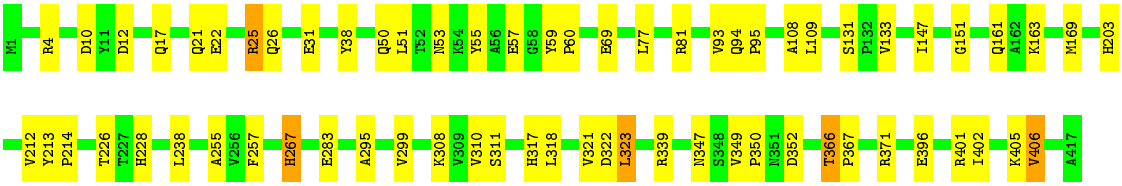
- Molecule 1: SERINE HYDROXYMETHYLTRANSFERASE

Chain A:



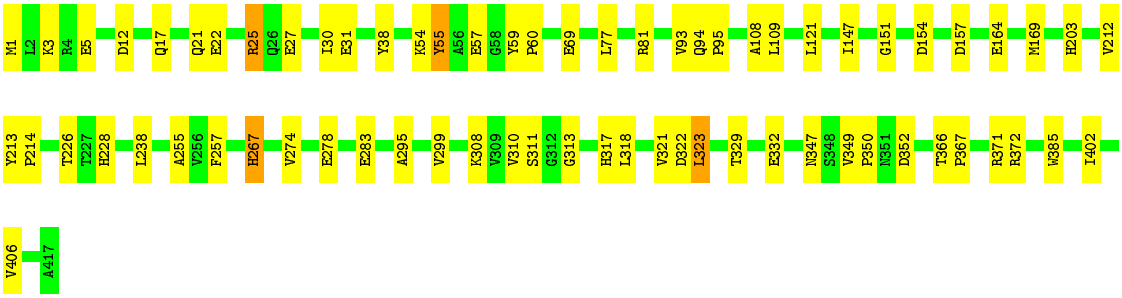
- Molecule 1: SERINE HYDROXYMETHYLTRANSFERASE

Chain B:



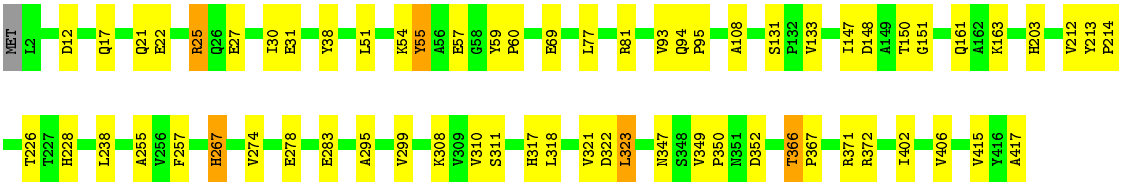
- Molecule 1: SERINE HYDROXYMETHYLTRANSFERASE

Chain C:



- Molecule 1: SERINE HYDROXYMETHYLTRANSFERASE

Chain D:


MET
L2
D12
Q17
Q21
E22
R25
Q26
E27
I30
E31
Y38
L51
K54
Y55
A56
E57
G58
Y59
P60
E69
L77
R81
V93
Q94
P95
A108
S131
P132
V133
I147
D148
A149
T150
G151
Q161
A162
K163
H203
V212
Y213
P214
T226
T227
H228
L238
A255
V256
F257
H267
V274
E278
E283
A295
V299
K308
V309
V310
S311
H317
L318
V321
D322
L323
N347
S348
V349
P350
N351
D352
T366
P367
R371
R372
I402
V406
V415
Y416
A417

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	77.46Å 172.24Å 95.05Å 90.00° 104.12° 90.00°	Depositor
Resolution (Å)	20.00 – 2.40 92.18 – 2.40	Depositor EDS
% Data completeness (in resolution range)	88.0 (20.00-2.40) 99.8 (92.18-2.40)	Depositor EDS
R_{merge}	0.23	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.54 (at 2.40Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.174 , 0.196 0.180 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å^2)	36.7	Xtriage
Anisotropy	0.331	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.31 , 41.1	EDS
L-test for twinning[2]	$< \|L\| > = 0.51$, $< L^2 > = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	13204	wwPDB-VP
Average B, all atoms (Å^2)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.73% 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 (i)

5.1 Standard geometry (i)

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

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	0.32	0/3216	0.59	0/4363
1	B	0.32	0/3235	0.58	0/4386
1	C	0.32	0/3208	0.58	0/4356
1	D	0.33	0/3213	0.59	0/4360
All	All	0.32	0/12872	0.59	0/17465

There are no bond length outliers.

There are no bond angle outliers.

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	3148	0	3081	46	0
1	B	3167	0	3108	48	0
1	C	3140	0	3056	55	0
1	D	3145	0	3070	47	0
2	A	20	0	11	1	0
2	B	20	0	11	1	0
2	C	20	0	11	0	0
2	D	20	0	11	0	0
3	A	34	0	21	0	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	34	0	21	0	0
3	C	34	0	21	0	0
3	D	34	0	21	0	0
4	A	93	0	0	2	0
4	B	94	0	0	6	0
4	C	85	0	0	4	0
4	D	116	0	0	4	0
All	All	13204	0	12443	185	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 7.

The worst 5 of 185 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:163:LYS:HD3	1:C:164:GLU:OE2	1.77	0.84
1:C:77:LEU:O	1:C:81:ARG:HG3	1.80	0.82
1:D:77:LEU:O	1:D:81:ARG:HG3	1.80	0.81
1:B:77:LEU:O	1:B:81:ARG:HG3	1.82	0.80
1:A:164:GLU:OE2	1:D:163:LYS:HD3	1.81	0.79

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	414/417 (99%)	402 (97%)	11 (3%)	1 (0%)	47	62
1	B	415/417 (100%)	403 (97%)	11 (3%)	1 (0%)	47	62
1	C	415/417 (100%)	401 (97%)	14 (3%)	0	100	100
1	D	414/417 (99%)	401 (97%)	12 (3%)	1 (0%)	47	62

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	1658/1668 (99%)	1607 (97%)	48 (3%)	3 (0%)	47	62

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	366	THR
1	A	366	THR
1	D	366	THR

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	320/329 (97%)	314 (98%)	6 (2%)	57	75
1	B	323/329 (98%)	316 (98%)	7 (2%)	52	71
1	C	317/329 (96%)	312 (98%)	5 (2%)	62	79
1	D	319/329 (97%)	314 (98%)	5 (2%)	62	79
All	All	1279/1316 (97%)	1256 (98%)	23 (2%)	59	76

5 of 23 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	347	ASN
1	B	406	VAL
1	D	323	LEU
1	B	396	GLU
1	C	25	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 21 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	394	ASN
1	C	203	HIS

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	D	267	HIS
1	B	290	GLN
1	D	289	GLN

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 carbohydrates in this entry.

5.6 Ligand geometry (i)

8 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	PLG	C	3001	-	17,20,20	1.87	5 (29%)	23,28,28	2.68	10 (43%)
3	FFO	D	4002	-	28,36,36	2.48	13 (46%)	28,50,50	3.87	15 (53%)
3	FFO	B	2002	-	28,36,36	2.48	13 (46%)	28,50,50	3.90	15 (53%)
3	FFO	A	1002	-	28,36,36	2.39	13 (46%)	28,50,50	3.89	15 (53%)
2	PLG	D	4001	-	17,20,20	1.85	5 (29%)	23,28,28	2.74	11 (47%)
2	PLG	B	2001	-	17,20,20	1.91	6 (35%)	23,28,28	2.70	11 (47%)
2	PLG	A	1001	-	17,20,20	2.00	5 (29%)	23,28,28	2.68	11 (47%)
3	FFO	C	3002	-	28,36,36	2.40	12 (42%)	28,50,50	3.88	15 (53%)

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	PLG	C	3001	-	-	3/10/12/12	0/1/1/1
3	FFO	D	4002	-	-	3/18/37/37	0/2/3/3
3	FFO	B	2002	-	-	3/18/37/37	0/2/3/3
3	FFO	A	1002	-	-	3/18/37/37	0/2/3/3
2	PLG	D	4001	-	-	2/10/12/12	0/1/1/1
2	PLG	B	2001	-	-	3/10/12/12	0/1/1/1
2	PLG	A	1001	-	-	3/10/12/12	0/1/1/1
3	FFO	C	3002	-	-	3/18/37/37	0/2/3/3

The worst 5 of 72 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	2002	FFO	C4-C4A	5.49	1.49	1.41
3	D	4002	FFO	C4-C4A	5.37	1.48	1.41
3	A	1002	FFO	C4-C4A	5.18	1.48	1.41
3	C	3002	FFO	C4-C4A	5.16	1.48	1.41
2	A	1001	PLG	C4A-C4	-4.89	1.45	1.51

The worst 5 of 103 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	A	1002	FFO	NA2-C2-N1	9.22	131.59	117.25
3	B	2002	FFO	NA2-C2-N1	9.15	131.48	117.25
3	C	3002	FFO	NA2-C2-N1	9.11	131.43	117.25
3	D	4002	FFO	NA2-C2-N1	9.11	131.41	117.25
3	A	1002	FFO	C4-N3-C2	7.18	127.34	115.93

There are no chirality outliers.

5 of 23 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	3001	PLG	C5-C4-C4A-N
2	D	4001	PLG	C5-C4-C4A-N
2	B	2001	PLG	C5-C4-C4A-N
2	A	1001	PLG	C5-C4-C4A-N
2	C	3001	PLG	C3-C4-C4A-N

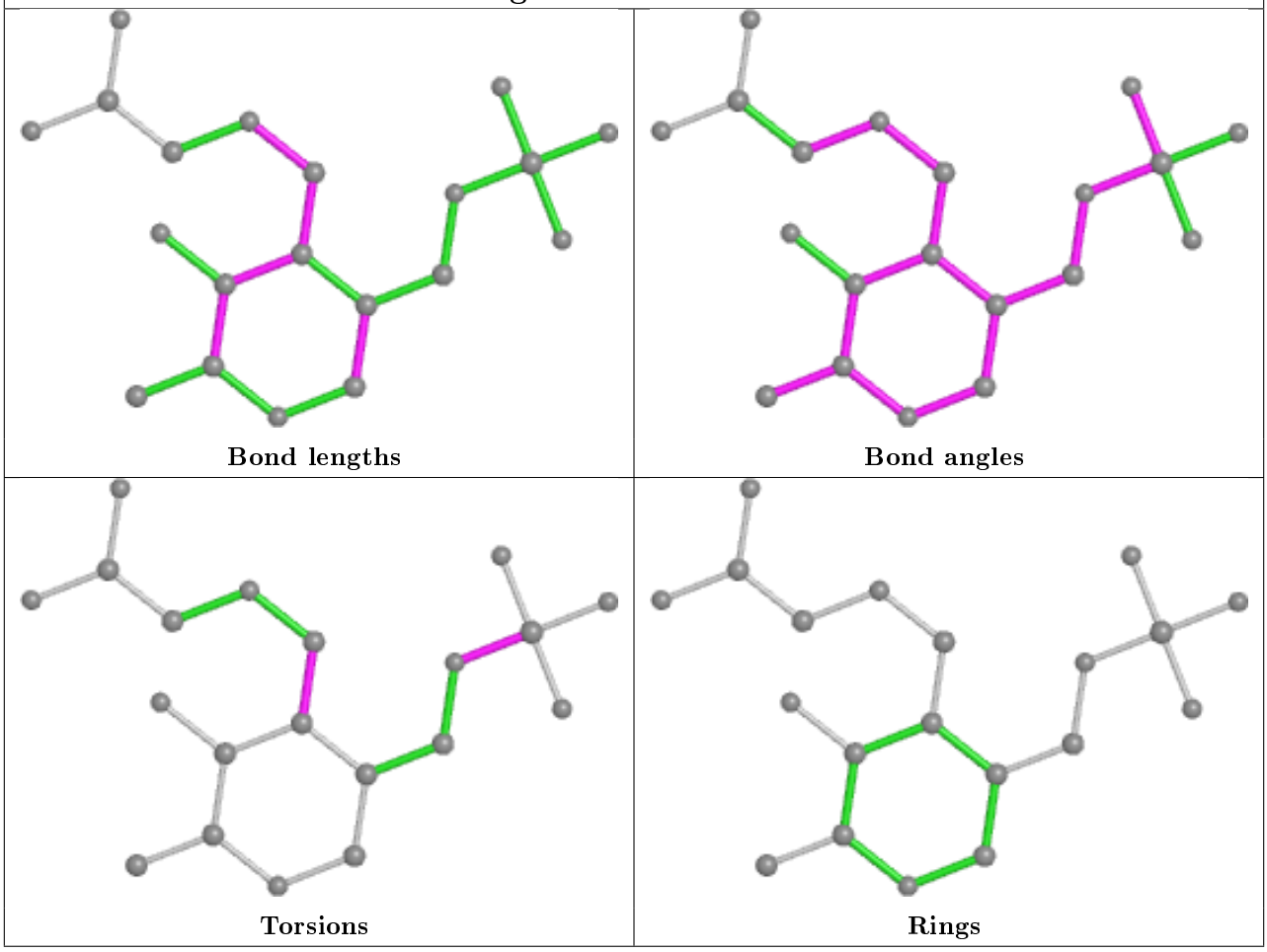
There are no ring outliers.

2 monomers are involved in 2 short contacts:

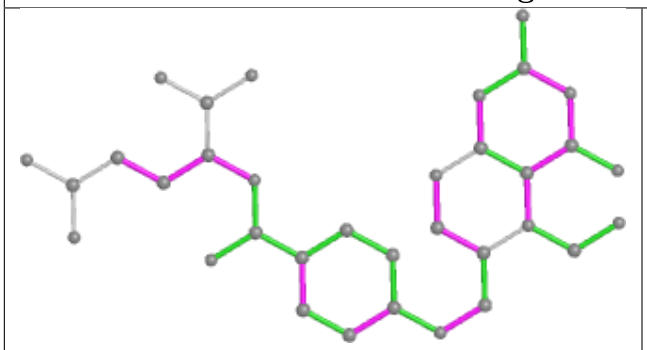
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	2001	PLG	1	0
2	A	1001	PLG	1	0

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.

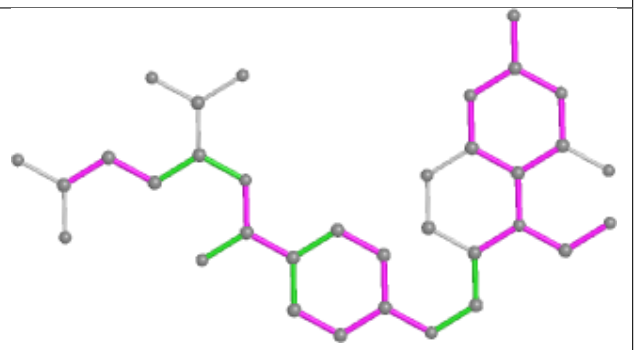
Ligand PLG C 3001



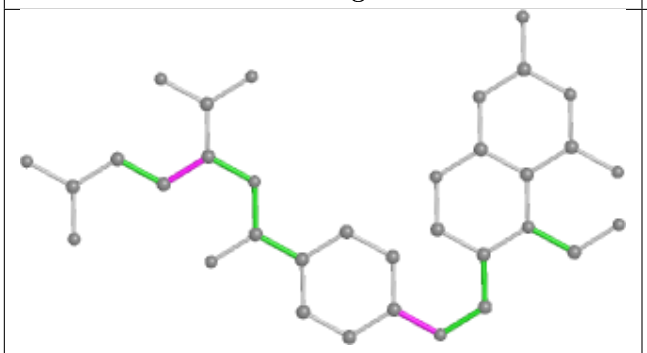
Ligand FFO D 4002



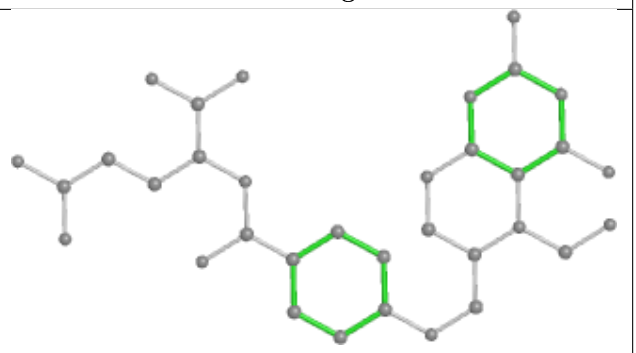
Bond lengths



Bond angles

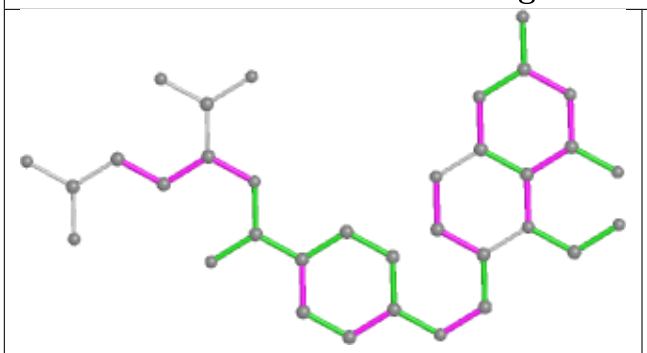


Torsions

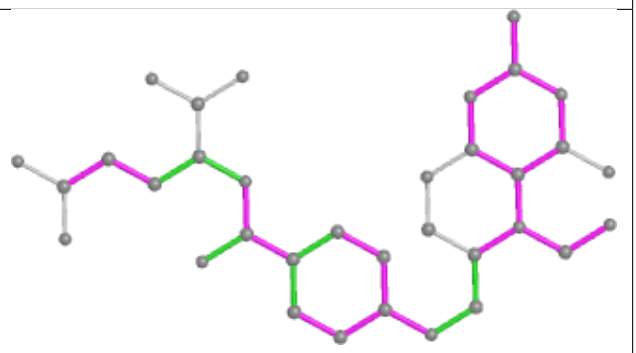


Rings

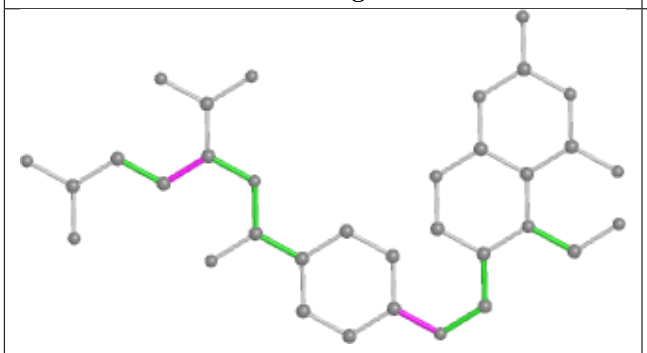
Ligand FFO B 2002



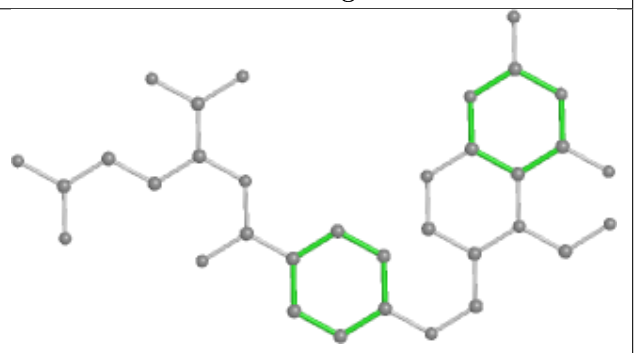
Bond lengths



Bond angles

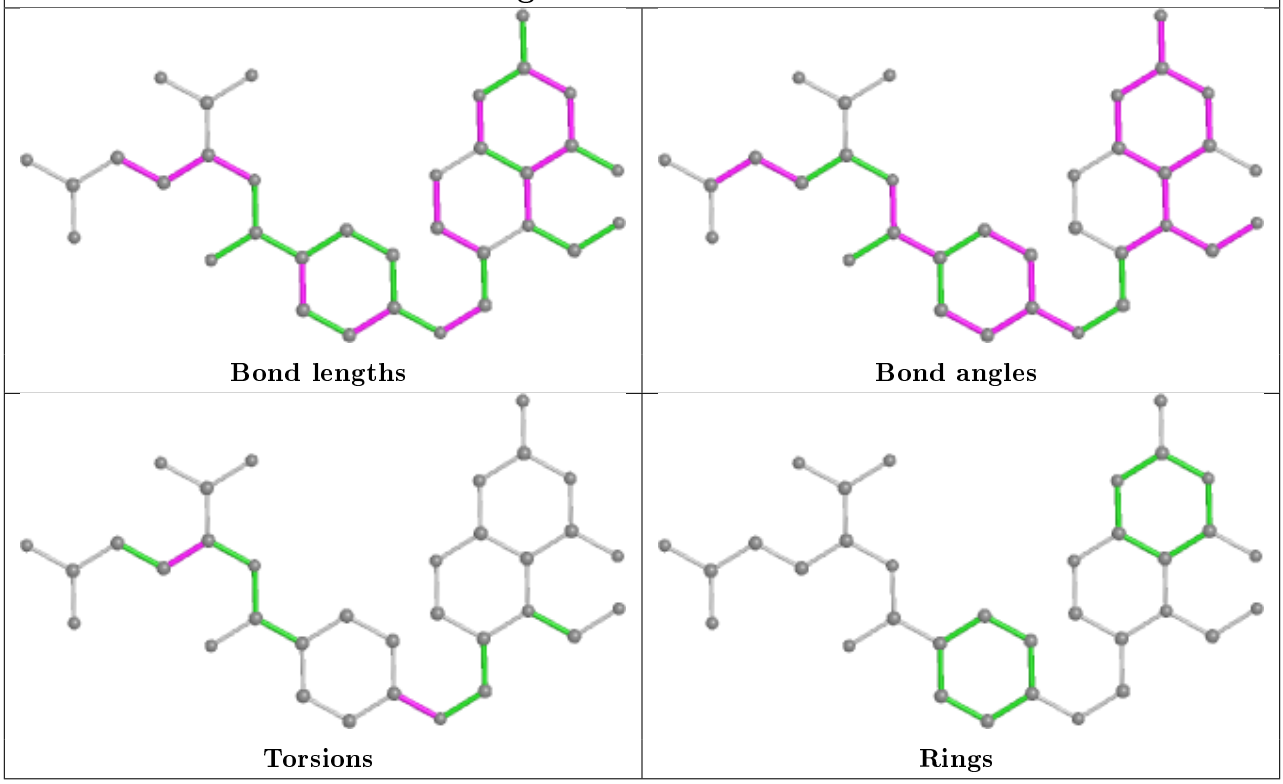


Torsions

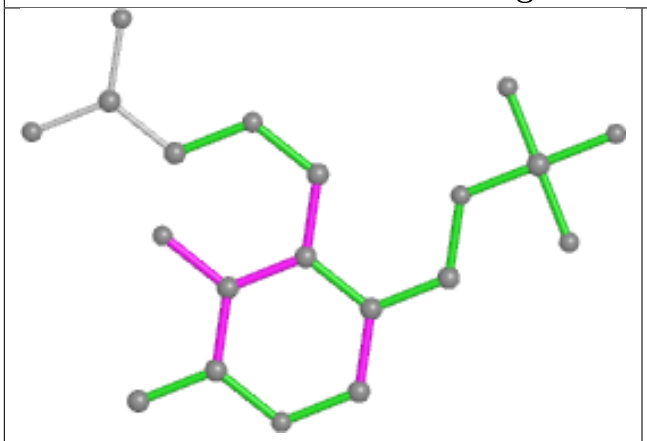


Rings

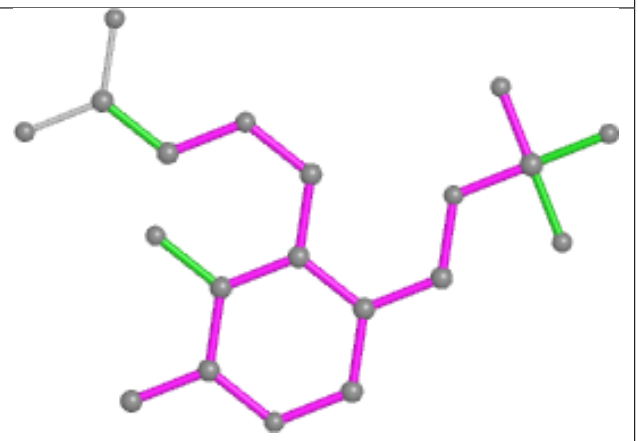
Ligand FFO A 1002



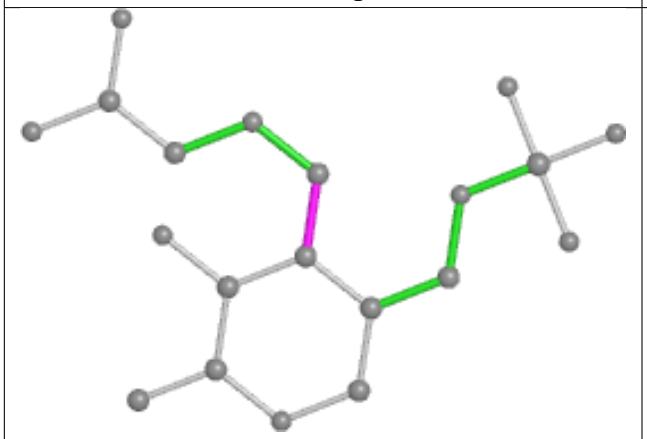
Ligand PLG D 4001



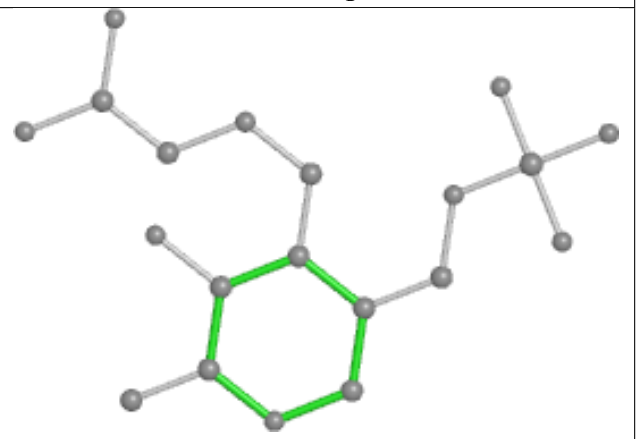
Bond lengths



Bond angles

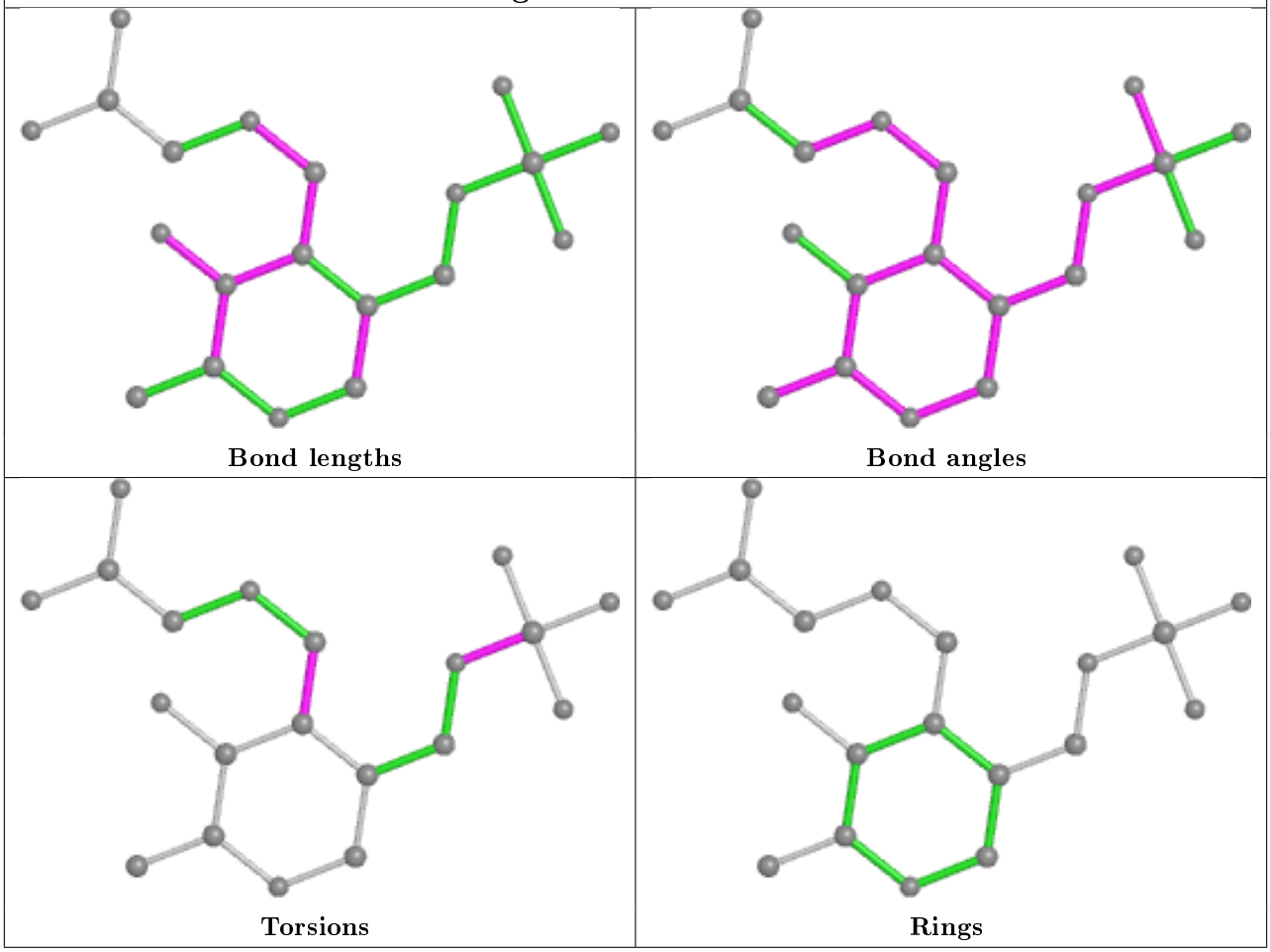


Torsions

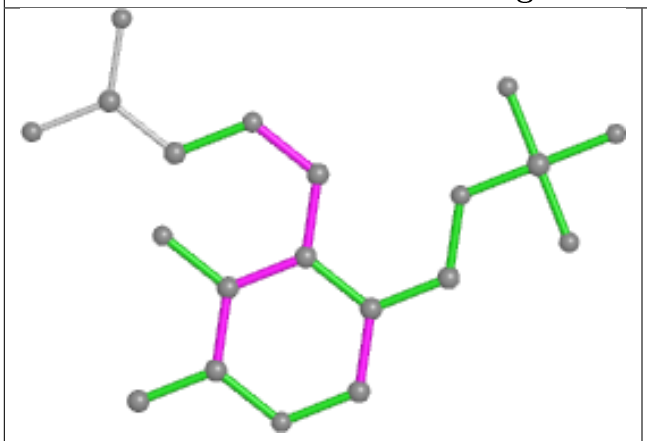


Rings

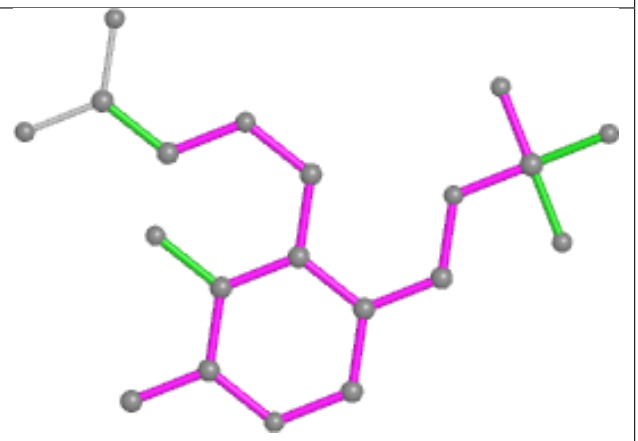
Ligand PLG B 2001



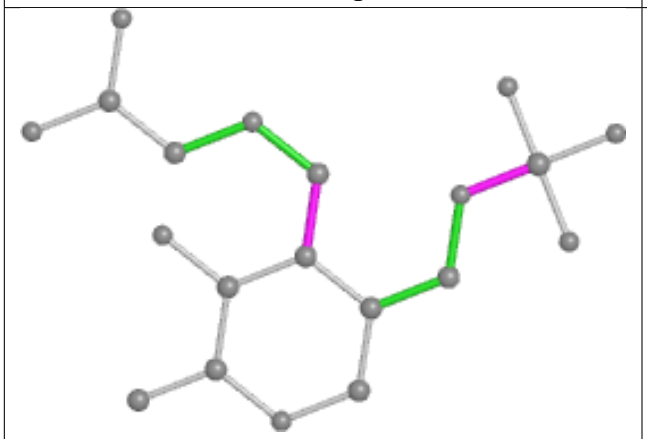
Ligand PLG A 1001



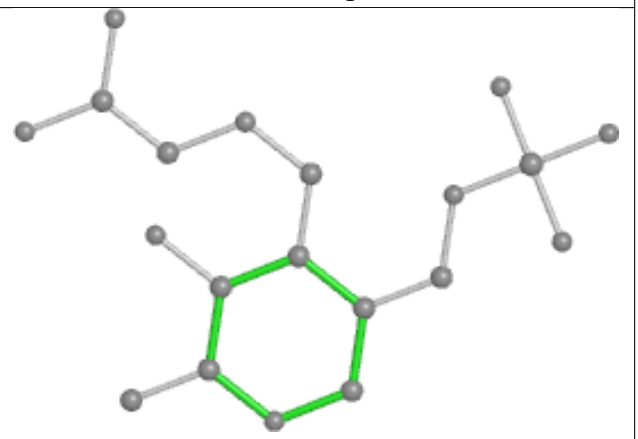
Bond lengths



Bond angles

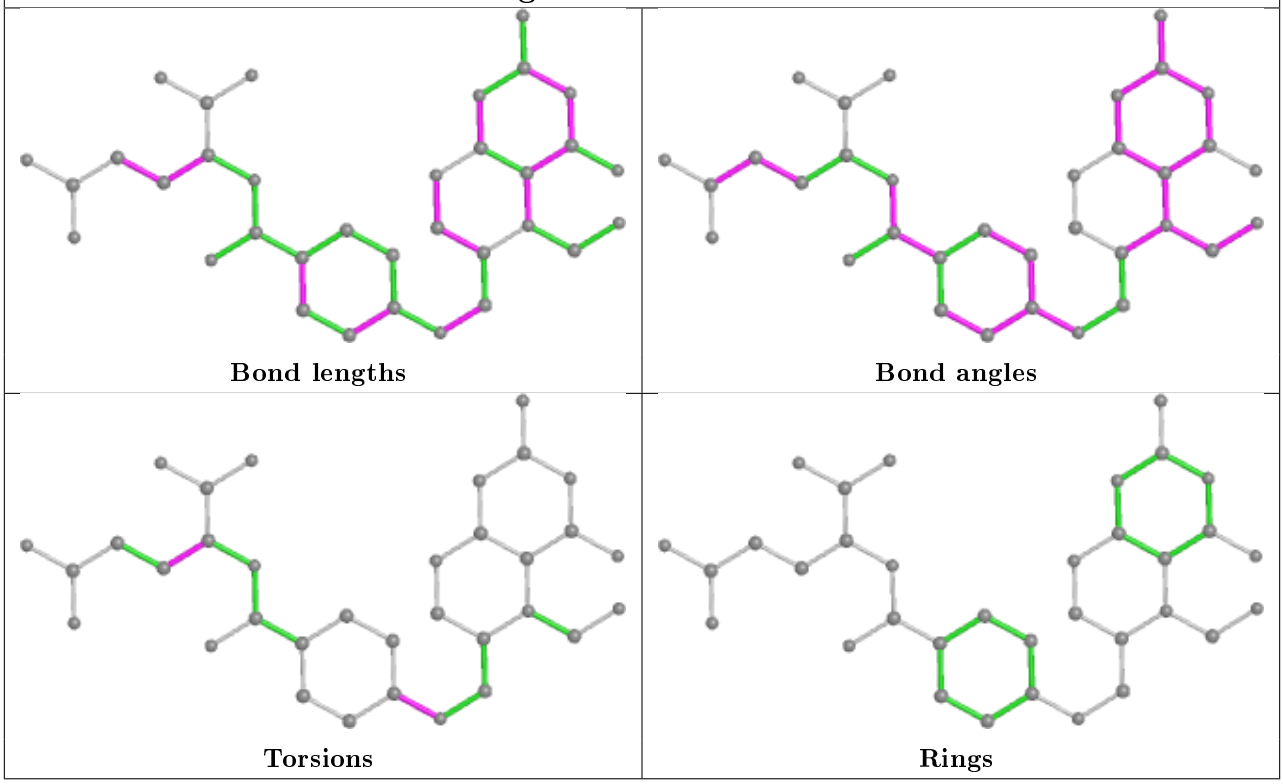


Torsions



Rings

Ligand FFO C 3002



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)

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	416/417 (99%)	-0.04	0	100	100	22, 33, 52, 69	0
1	B	417/417 (100%)	-0.03	0	100	100	21, 33, 54, 66	0
1	C	417/417 (100%)	-0.07	0	100	100	21, 36, 55, 70	0
1	D	416/417 (99%)	-0.02	0	100	100	22, 33, 53, 67	0
All	All	1666/1668 (99%)	-0.04	0	100	100	21, 34, 54, 70	0

There are no RSRZ outliers to report.

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

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

6.3 Carbohydrates (i)

There are no carbohydrates in this entry.

6.4 Ligands (i)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	FFO	B	2002	34/34	0.93	0.18	37,41,58,63	0
3	FFO	D	4002	34/34	0.94	0.19	35,41,60,64	0
3	FFO	C	3002	34/34	0.94	0.19	29,41,52,57	0
3	FFO	A	1002	34/34	0.95	0.18	28,39,49,54	0

Continued on next page...

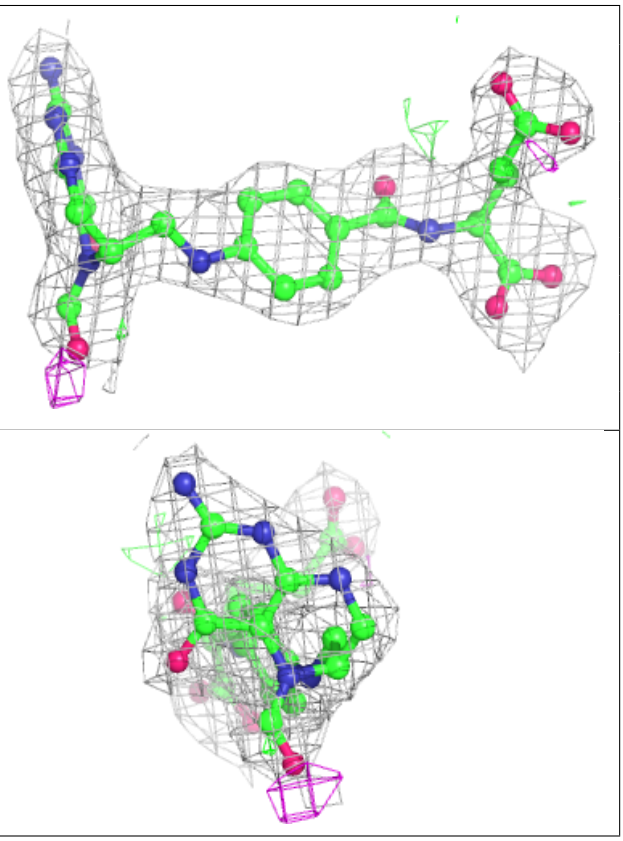
Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	PLG	B	2001	20/20	0.98	0.19	27,30,36,36	0
2	PLG	A	1001	20/20	0.98	0.17	25,28,31,32	0
2	PLG	D	4001	20/20	0.98	0.18	26,30,34,37	0
2	PLG	C	3001	20/20	0.99	0.17	25,29,34,36	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.

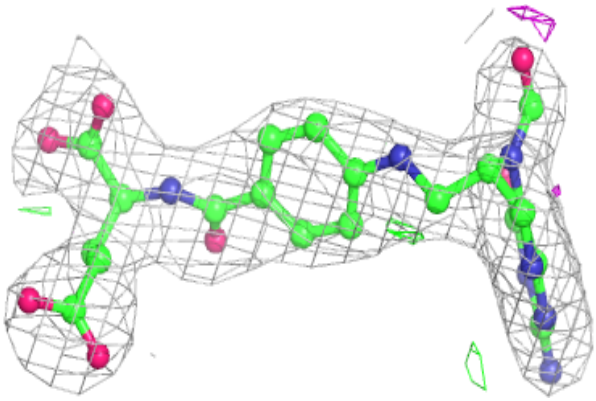
Electron density around FFO B 2002:

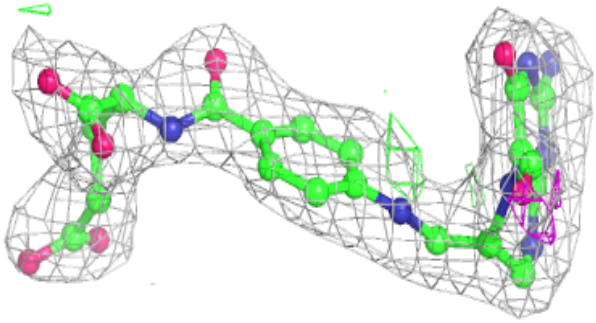
2mF$_o$-DF$_c$ (at 0.7 rmsd) in gray
mF$_o$-DF$_c$ (at 3 rmsd) in purple (negative)
and green (positive)

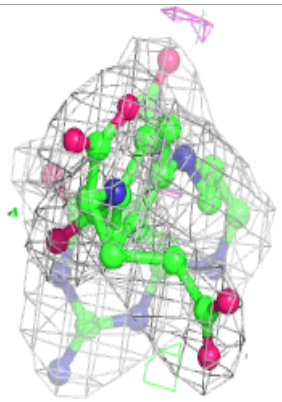


Electron density around FFO D 4002:

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

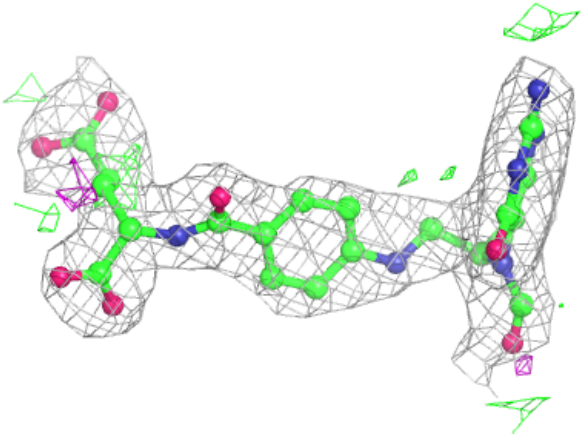


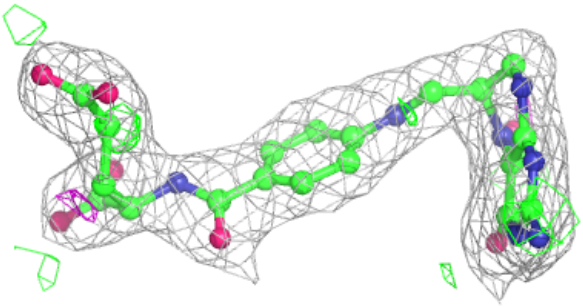


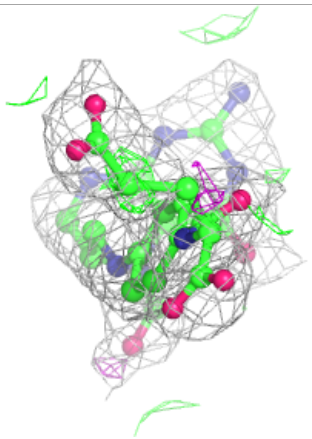


Electron density around FFO C 3002:

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

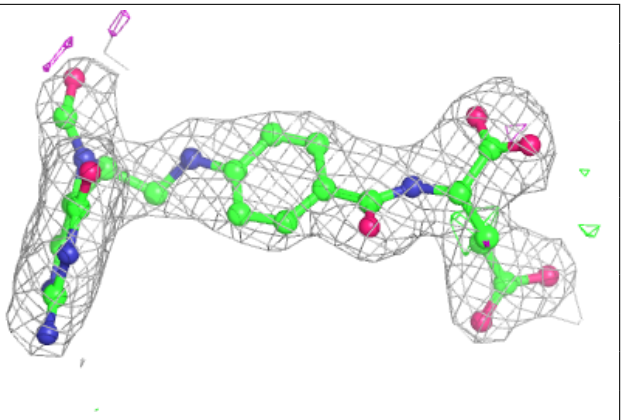


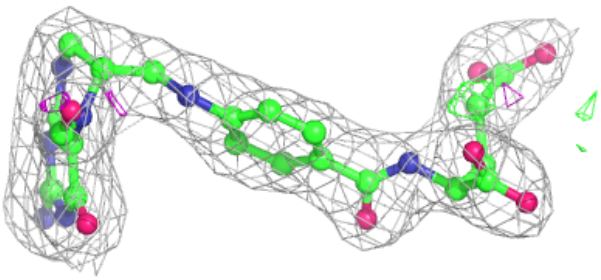


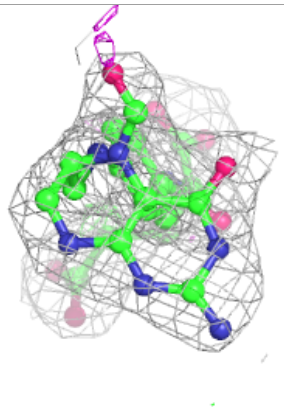


Electron density around FFO A 1002:

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

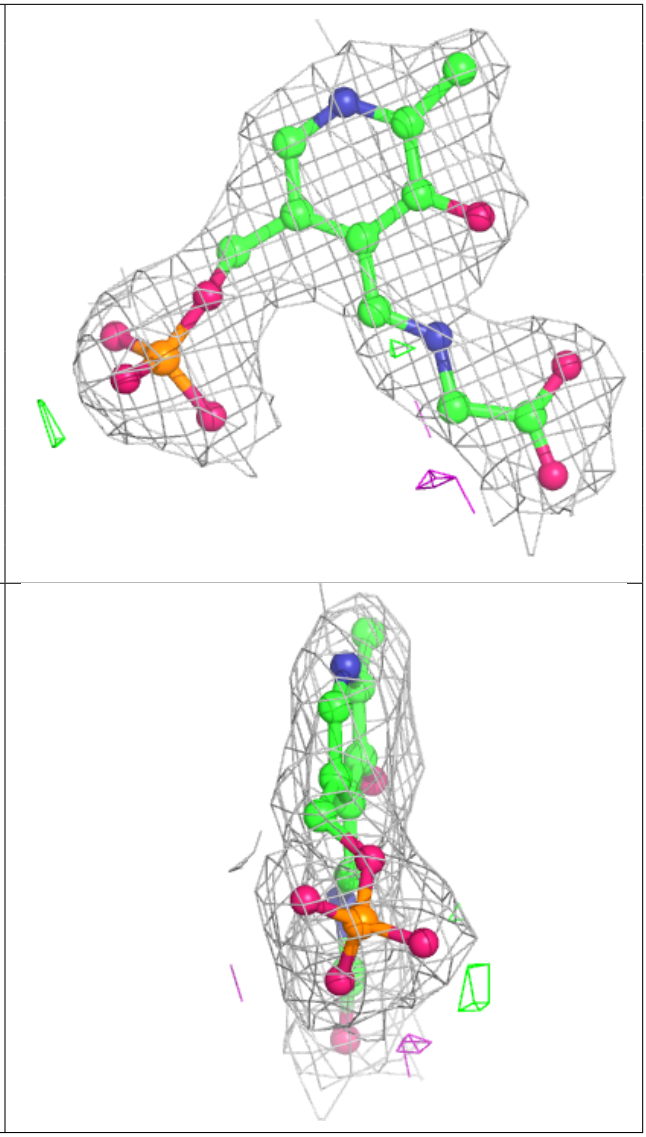


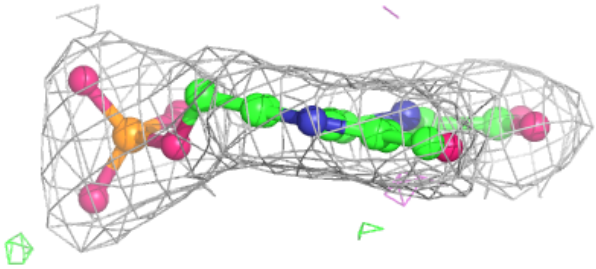




Electron density around PLG B 2001:

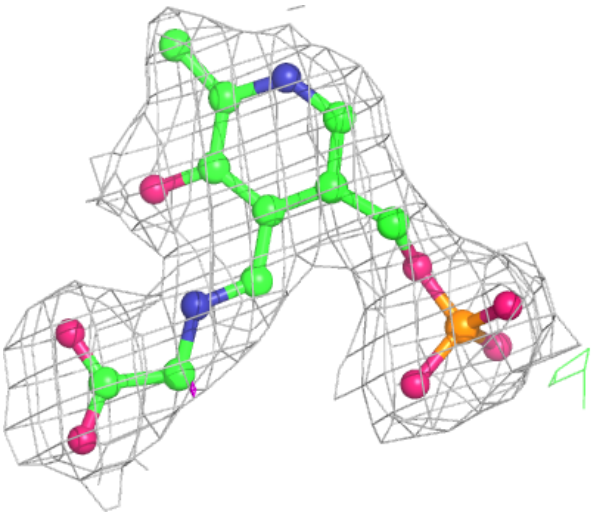
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

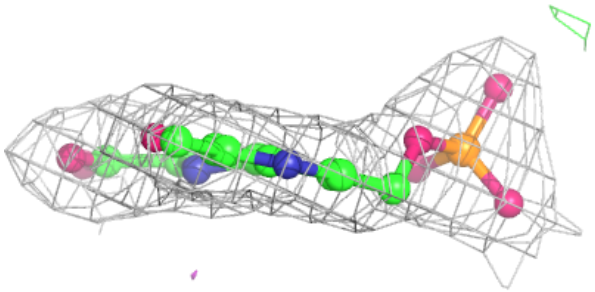


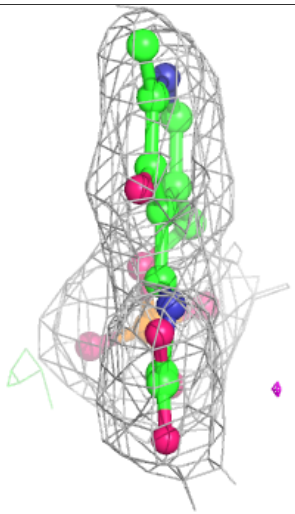


Electron density around PLG A 1001:

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

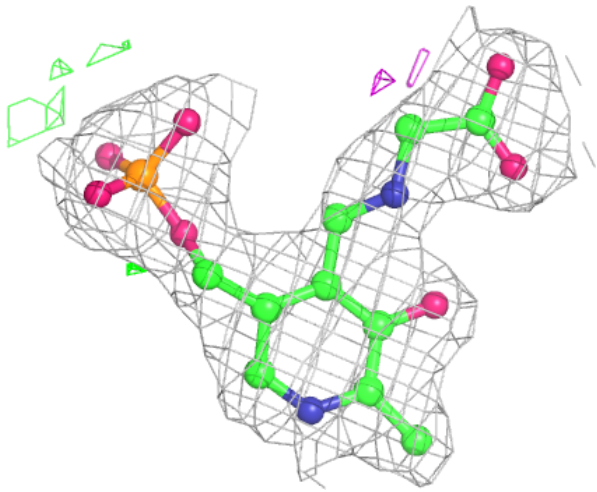


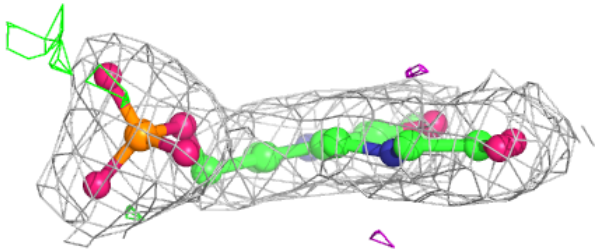


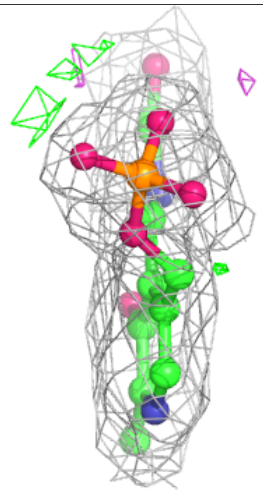


Electron density around PLG D 4001:

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

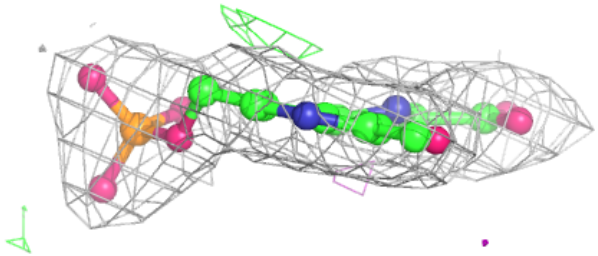


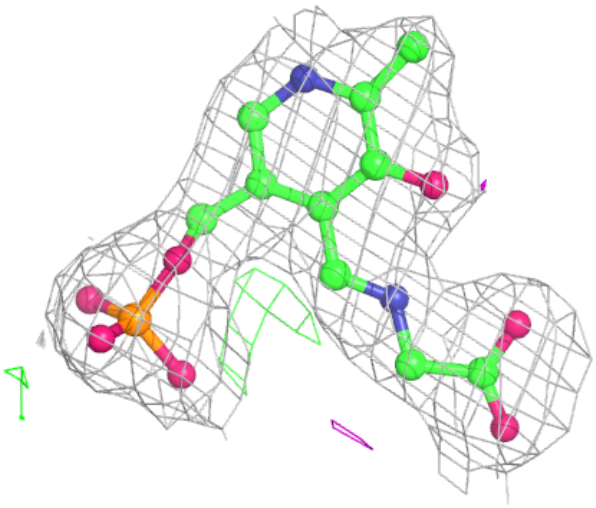


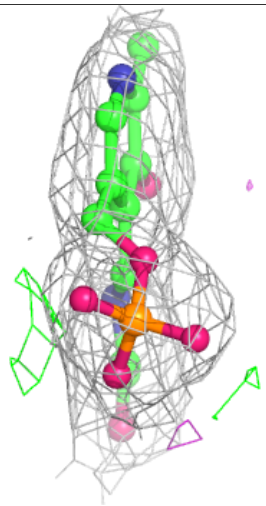


Electron density around PLG C 3001:

$2mF_o$-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.