

Data Path : Z:\HPCHEM1\BNA M\DATA\BM031317\
 Data File : BM009247.D
 Acq On : 14 Mar 2017 09:58
 Operator : SJ/MA
 Sample : I2181-17
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0BD4

Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 1 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-2016\SOM-EPA-BM031317MA.M
 Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	7.034	663	671	678	rBV	147960	250686	3.22%	0.453%
2	7.210	694	701	709	rBV	541450	867189	11.12%	1.566%
3	7.404	728	734	742	rVB	534192	873594	11.20%	1.577%
4	7.875	808	814	820	rBV	450615	743753	9.54%	1.343%
5	8.569	926	932	941	rVB2	468230	732416	9.39%	1.323%
6	8.969	992	1000	1006	rBV2	123380	201528	2.58%	0.364%
7	9.040	1006	1012	1018	rVV	789207	1325731	17.00%	2.394%
8	9.763	1126	1135	1144	rVB	641975	1078677	13.83%	1.948%
9	10.292	1218	1225	1233	rBV	901619	1498300	19.22%	2.706%
10	10.581	1270	1274	1283	rVB2	72413	121519	1.56%	0.219%
11	10.675	1283	1290	1296	rVV	688355	1161549	14.90%	2.097%
12	13.922	1835	1842	1854	rBV	2231073	3077133	39.46%	5.556%
13	14.210	1884	1891	1898	rBV	2220904	3226496	41.38%	5.826%
14	14.516	1937	1943	1952	rBV2	1193752	1903224	24.41%	3.437%
15	14.710	1969	1976	1986	rBV	197795	328955	4.22%	0.594%
16	15.510	2104	2112	2120	rBV	3198762	4449501	57.06%	8.035%
17	15.627	2125	2132	2143	rBV	1226151	1640151	21.03%	2.962%
18	17.263	2404	2410	2420	rBV2	1777225	2622674	33.64%	4.736%
19	17.363	2420	2427	2435	rBV2	3771371	5356975	68.70%	9.673%
20	17.921	2516	2522	2527	rBV2	933235	1141434	14.64%	2.061%
21	19.656	2811	2817	2827	rBV	5158008	7011639	89.92%	12.661%
22	21.439	3115	3120	3129	rBV	3160022	3952065	50.69%	7.136%
23	23.627	3485	3492	3501	rVB2	3961656	7797303	100.00%	14.080%
24	23.780	3511	3518	3525	rVB2	2071623	4017026	51.52%	7.254%

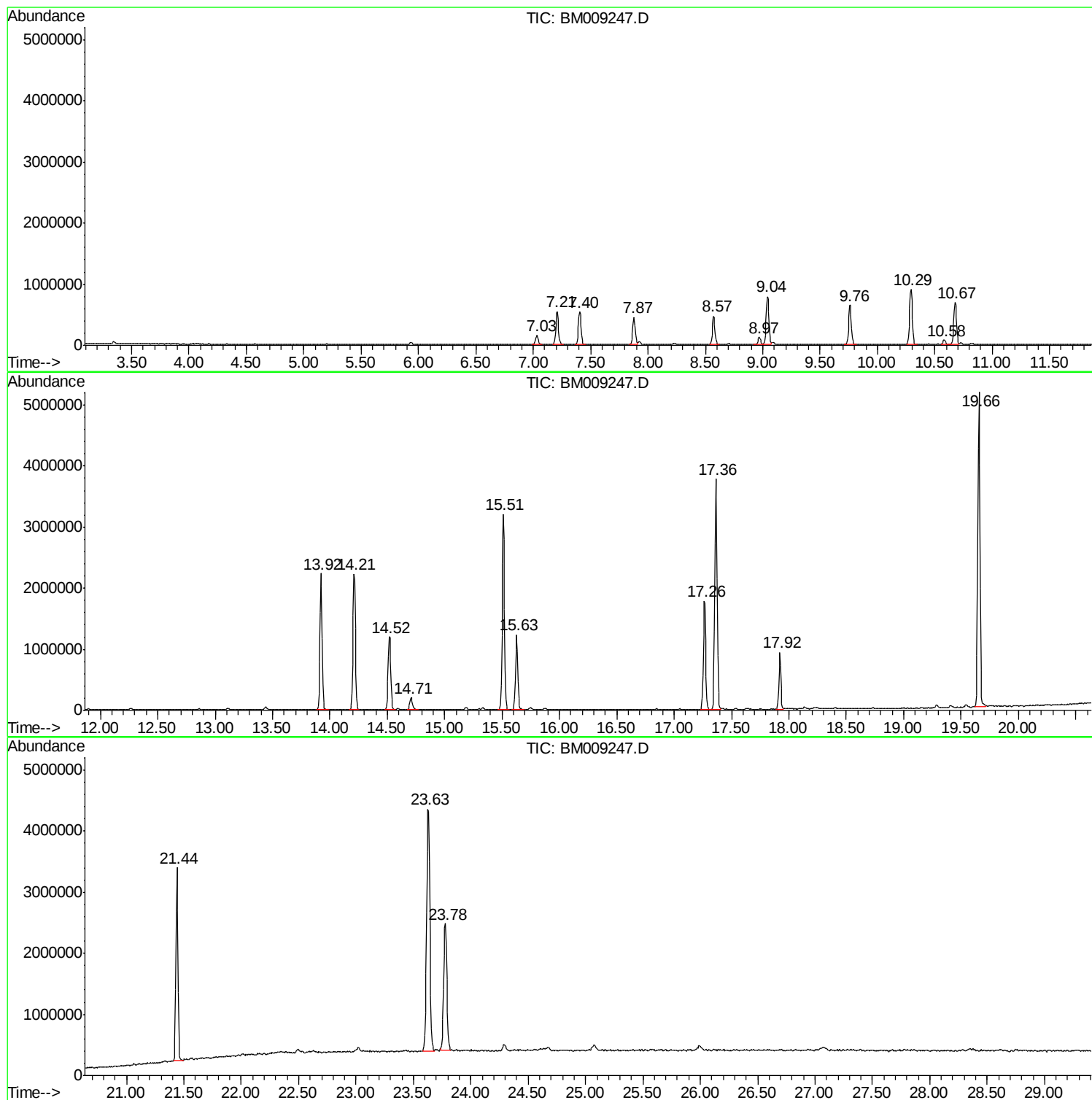
Sum of corrected areas: 55379518

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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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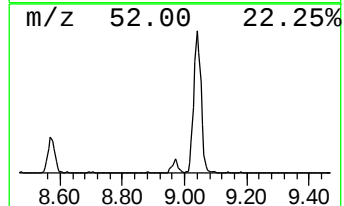
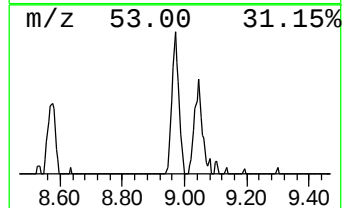
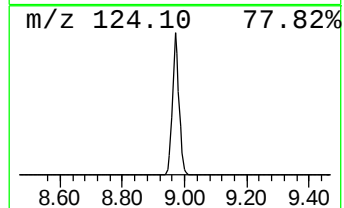
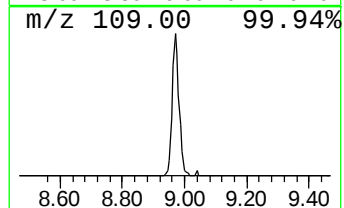
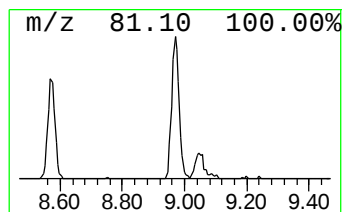
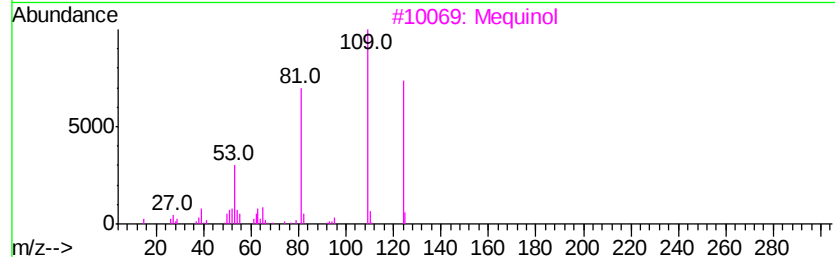
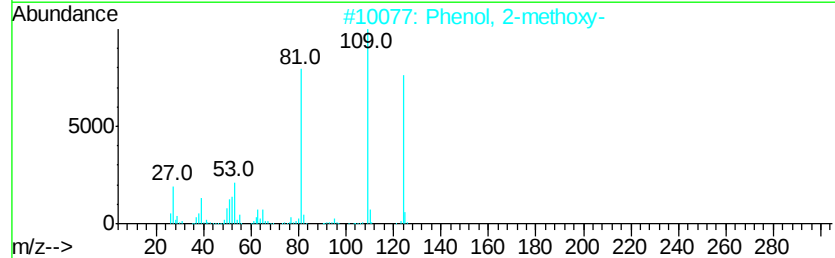
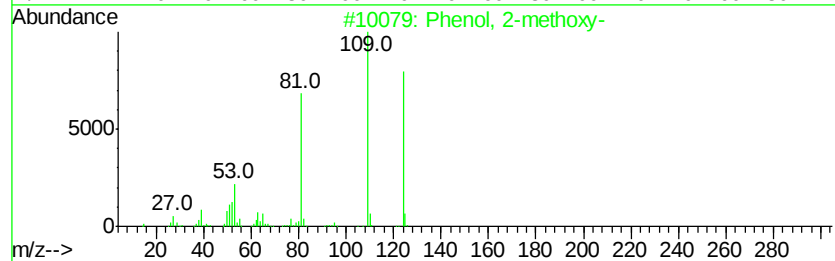
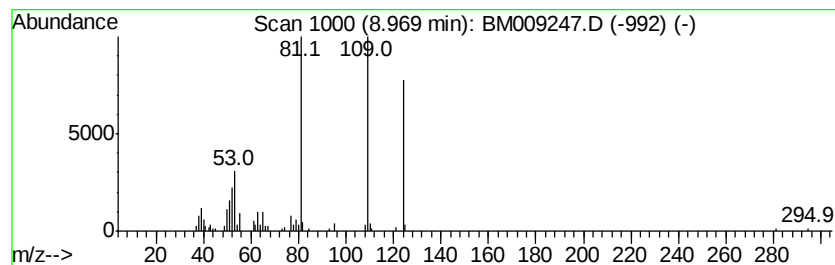
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-2016\SOM-EPA-BM031317MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Phenol, 2-methoxy- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.97	5.42 ng/ul	201528	1,4-Dichlorobenzene-d4	7.87

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenol, 2-methoxy-	124	C7H8O2	000090-05-1	76
2		Phenol, 2-methoxy-	124	C7H8O2	000090-05-1	76
3		Mequinol	124	C7H8O2	000150-76-5	68
4		Mequinol	124	C7H8O2	000150-76-5	49
5		Phenol, 2-methoxy-	124	C7H8O2	000090-05-1	46



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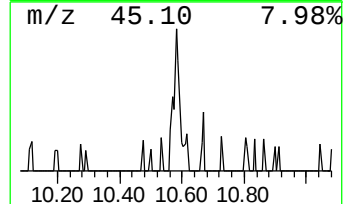
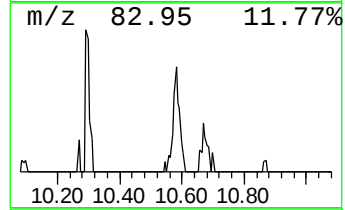
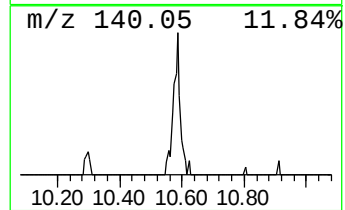
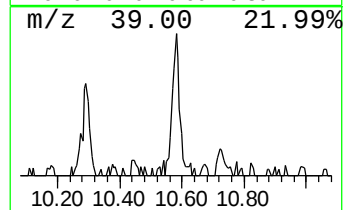
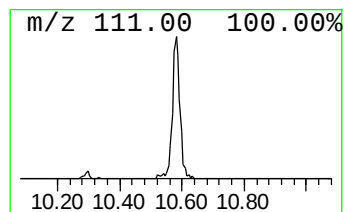
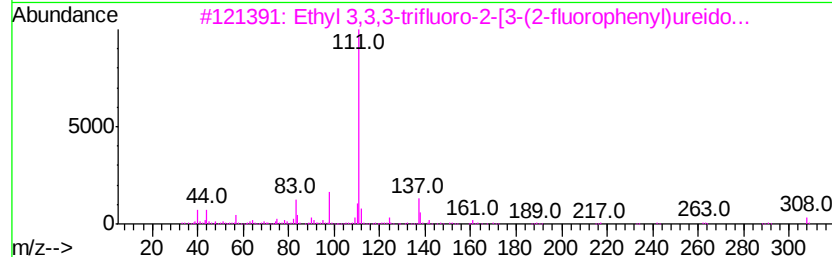
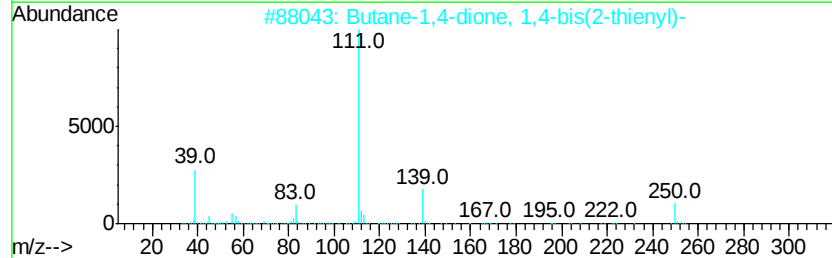
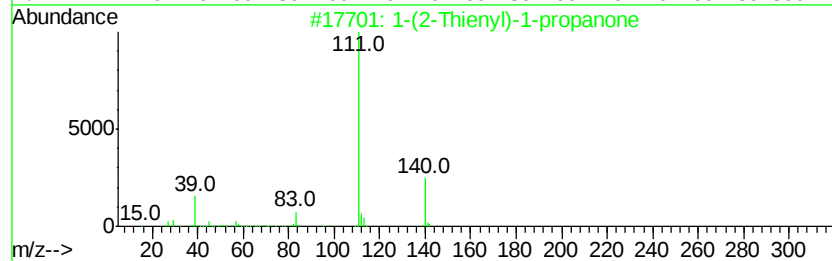
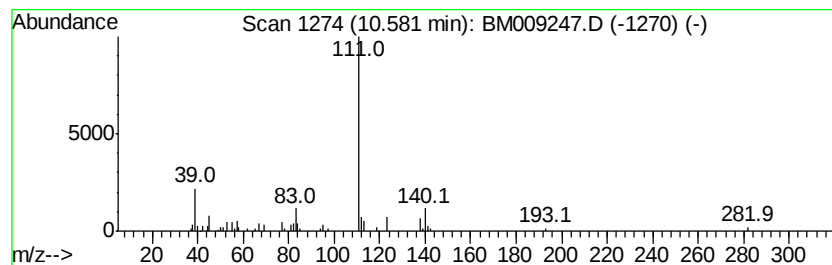
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Peak Number 2 1-(2-Thienyl)-1-propanone Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.58	2.09 ng/ul	121519	Napthalene-d8	10.67
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		1-(2-Thienyl)-1-propanone	140 C7H8OS	013679-75-9 72
2		Butane-1,4-dione, 1,4-bis(2-thie...	250 C12H10O2S2	013669-05-1 64
3		Ethyl 3,3,3-trifluoro-2-[3-(2-fl...	308 C12H12F4N2O3	1000225-32-1 64
4		Thiophene-2-carboxamide, N-(2-ch...	282 C11H7ClN2O3S	312592-64-6 59
5		4-Fluorophenyl 2-thiophenecarbox...	222 C11H7F02S	1000298-21-9 59



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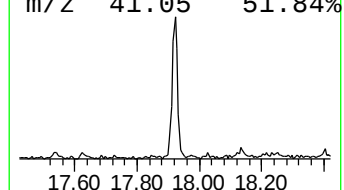
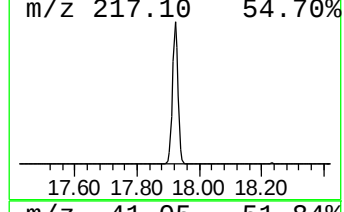
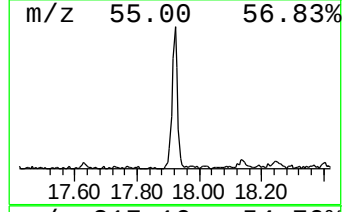
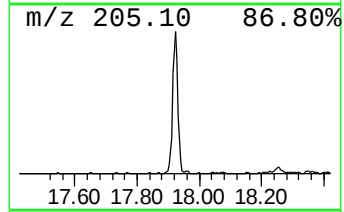
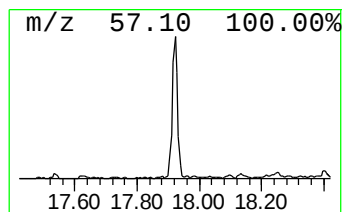
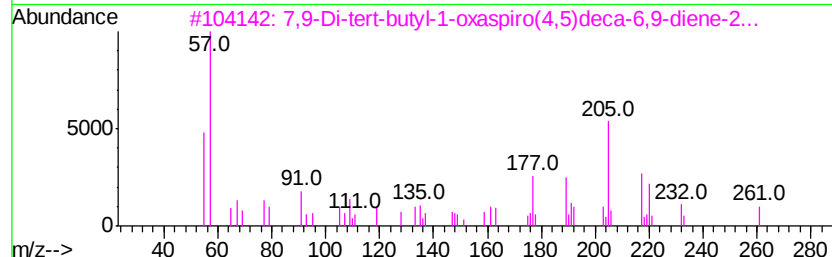
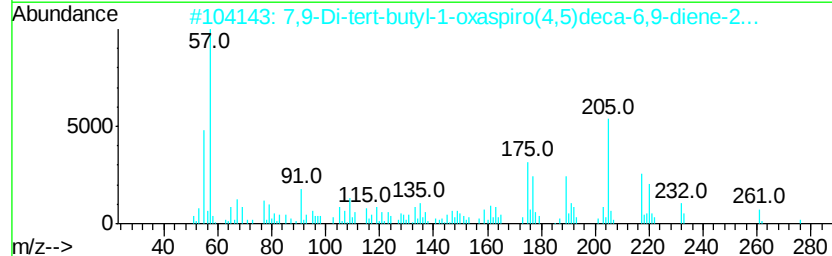
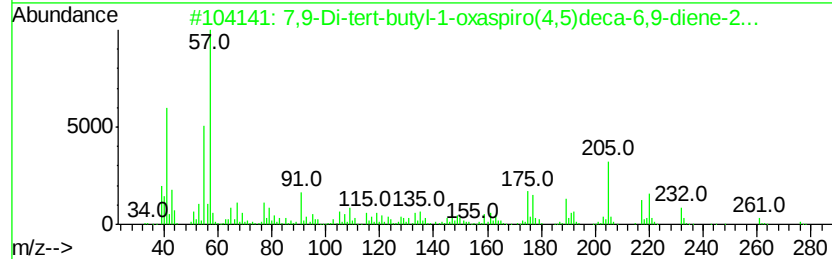
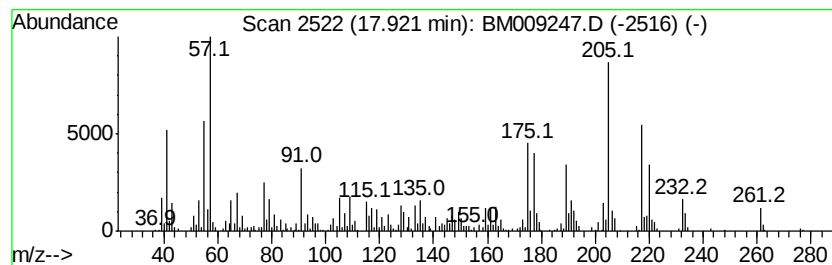
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Peak Number 3 7,9-Di-tert-butyl-1-oxaspiro... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.92	8.70 ng/ul	1141430	Phenanthrene-d10	17.27

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			7,9-Di-tert-butyl-1-oxaspiro(4,5)deca-6,9-diene-2-one	276	C17H24O3	082304-66-3	87
2			7,9-Di-tert-butyl-1-oxaspiro(4,5)deca-6,9-diene-2-one	276	C17H24O3	082304-66-3	60
3			7,9-Di-tert-butyl-1-oxaspiro(4,5)deca-6,9-diene-2-one	276	C17H24O3	082304-66-3	55
4			2,5-di-tert-Butyl-1,4-benzoquinone	220	C14H2O2	002460-77-7	47
5			Ethanone, 1-(9-anthracenyl)-	220	C16H12O	000784-04-3	46



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Phenol, 2-methoxy-	8.97	5.4	ng/ul	201528	1	7.87	743753	20.0
1-(2-Thienyl)-1-p...	10.58	2.1	ng/ul	121519	2	10.67	1161550	20.0
7,9-Di-tert-butyl...	17.92	8.7	ng/ul	1141430	4	17.27	2622670	20.0