LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM111717\
Data File : BM013029.D
Acq On : 18 Nov 2017 08:39
Operator : SJ/JU
Sample : I6405-04
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BE2S6

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

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

Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM111617.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	3.252	25	28	36	rVB	45911	61066	1.30%	0.113%
2	3.969	145	150	155	rBV	100128	125952	2.67%	0.234%
3	4.481	232	237	243	rVB3	44227	67589	1.43%	0.126%
4	5.069	331	337	345	rVB	354960	508508	10.78%	0.945%
5	5.257	363	369	378	rVB	673657	945699	20.06%	1.757%
6	5.410	385	395	401	rVB	232968	443229	9.40%	0.824%
7	5.752	445	453	458	rBV	165040	253303	5.37%	0.471%
8	6.222	529	533	538	rVB2	63986	94063	1.99%	0.175%
9	6.704	609	615	623	rBV2	60616	94946	2.01%	0.176%
10	6.887	638	646	654	rBV3	187761	401807	8.52%	0.747%
11	7.057	667	675	680	rBV	560526	828020	17.56%	1.539%
12	7.110	680	684	689	rVB	77171	110919	2.35%	0.206%
13	7.251	702	708	715	rVB	655994	1007293	21.36%	1.872%
14	7.369	722	728	733	rBV	318440	486014	10.31%	0.903%
15	7.716	777	787	798	rBV	610077	1042674	22.11%	1.937%
16	7.828	801	806	815	rVB2	100515	178965	3.80%	0.333%
17	8.087	841	850	854	rBV3	108704	221080	4.69%	0.411%
18	8.145	854	860	865	rVV	334300	514776	10.92%	0.957%
19	8.204	865	870	875	rVV3	55010	88179	1.87%	0.164%
20	8.263	875	880	885	rVV3	76851	124601	2.64%	0.232%
21	8.340	885	893	900	rVV	434730	763422	16.19%	1.419%
22	8.416	900	906	915	rVB	491360	784126	16.63%	1.457%
23	8.816	968	974	977	rBV3	64158	98346	2.09%	0.183%
24	8.869	977	983	989	rVB	719916	1160791	24.62%	2.157%
25	9.334	1058	1062	1067	rBV2	56903	84247	1.79%	0.157%
26	9.392	1067	1072	1080	rVB2	50606	88044	1.87%	0.164%
27	9.581	1096	1104	1111	rBV	578480	979629	20.78%	1.820%
28	9.869	1148	1153	1156	rBV2	45154	78560	1.67%	0.146%
29	9.910	1156	1160	1167	rVB7	61716	130756	2.77%	0.243%
30	10.116	1187	1195	1202	rBV	938160	1522549	32.29%	2.829%
31	10.498	1253	1260	1265	rBV	763830	1271751	26.97%	2.363%
32	10.545	1265	1268	1274	rVB	67556	110467	2.34%	0.205%
33	10.857	1315	1321	1327	rBV2	135740	219657	4.66%	0.408%
34	11.769	1472	1476	1482	rVB3	58144	83583	1.77%	0.155%

LSC Area Percent Report

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Operator : SJ/JU
Sample : I6405-04
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BE2S6

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

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

Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM111617.M
Title : SVOA CALIBRATION

35	11.845	1482	1489	1493	rBV4	28683	49526	1.05%	0.092%
36	12.269	1555	1561	1568	rBV3	55941	103694	2.20%	0.193%
37	12.463	1590	1594	1600	rVB2	66950	94630	2.01%	0.176%
38	12.698	1629	1634	1640	rBV	43198	72139	1.53%	0.134%
39	12.939	1668	1675	1679	rBV3	52275	82681	1.75%	0.154%
40	13.257	1723	1729	1739	rBV5	56662	105624	2.24%	0.196%
41	13.627	1787	1792	1799	rBV	80235	115263	2.44%	0.214%
42	13.769	1810	1816	1822	rBV	1732004	2382055	50.52%	4.426%
43	13.845	1822	1829	1837	rBV	687744	1102383	23.38%	2.048%
44	14.045	1857	1863	1871	rVB	1823155	2674096	56.71%	4.969%
45	14.357	1909	1916	1924	rBV2	1128280	1728898	36.67%	3.212%
46	14.469	1930	1935	1944	rVB2	89344	130484	2.77%	0.242%
47	14.551	1944	1949	1957	rBV	186700	275649	5.85%	0.512%
48	14.910	2005	2010	2016	rBV	165632	389363	8.26%	0.723%
49	14.998	2022	2025	2033	rVB9	39552	70700	1.50%	0.131%
50	15.263	2063	2070	2074	rBV6	43449	72621	1.54%	0.135%
51	15.351	2079	2085	2093	rVB2	2550500	4324565	91.72%	8.035%
52	15.463	2098	2104	2110	rBV	712057	1011362	21.45%	1.879%
53	15.616	2127	2130	2135	rVB	91371	122309	2.59%	0.227%
54	15.863	2165	2172	2179	rBV6	103080	213705	4.53%	0.397%
55	15.980	2185	2192	2195	rBV3	73731	138514	2.94%	0.257%
56	16.016	2195	2198	2204	rVV3	92790	168315	3.57%	0.313%
57	16.104	2208	2213	2221	rVB3	56292	128435	2.72%	0.239%
58	16.198	2223	2229	2233	rBV	91636	137374	2.91%	0.255%
59	16.257	2235	2239	2245	rVB2	116529	192915	4.09%	0.358%
60	16.316	2245	2249	2253	rBV2	92054	127164	2.70%	0.236%
61	16.363	2253	2257	2261	rVB2	53809	72328	1.53%	0.134%
62	16.416	2261	2266	2268	rBV4	31018	52599	1.12%	0.098%
63	16.521	2280	2284	2289	rVB4	73853	119964	2.54%	0.223%
64	16.580	2289	2294	2297	rBV2	143353	213958	4.54%	0.398%
65	16.616	2297	2300	2304	rVV3	107250	155696	3.30%	0.289%
66	16.651	2304	2306	2311	rVB2	45210	59405	1.26%	0.110%
67	16.986	2357	2363	2368	rBV7	54934	106958	2.27%	0.199%
68	17.104	2375	2383	2389	rBV	1482697	2164992	45.92%	4.023%
69	17.204	2392	2400	2409	rBV	2620194	3805596	80.71%	7.071%
70	17.274	2409	2412	2418	rVV6	28986	59051	1.25%	0.110%
71	17.345	2421	2424	2429	rVB5	32113	48150	1.02%	0.089%

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Misc :
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Instrument : BNA_M
ClientSampleId : BE2S6

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
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Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM111617.M
Title : SVOA CALIBRATION

72	17.468	2439	2445	2449	rBV2	65180	115040	2.44%	0.214%
73	18.010	2533	2537	2541	rBV5	35440	50519	1.07%	0.094%
74	18.233	2571	2575	2580	rBV2	73329	102021	2.16%	0.190%
75	18.421	2603	2607	2614	rBV2	37222	61003	1.29%	0.113%
76	18.580	2630	2634	2639	rVB6	64789	97018	2.06%	0.180%
77	18.927	2690	2693	2698	rBV2	70453	94639	2.01%	0.176%
78	19.139	2725	2729	2734	rBV2	41922	71309	1.51%	0.132%
79	19.227	2740	2744	2746	rBV5	60584	78025	1.65%	0.145%
80	19.292	2752	2755	2758	rBV4	67573	74641	1.58%	0.139%
81	19.380	2767	2770	2774	rVB3	64293	88076	1.87%	0.164%
82	19.498	2784	2790	2795	rVV	3416014	4539903	96.29%	8.436%
83	19.551	2795	2799	2806	rVB2	91268	149486	3.17%	0.278%
84	19.686	2818	2822	2826	rVB5	96090	124406	2.64%	0.231%
85	20.227	2910	2914	2919	rVB	218556	273545	5.80%	0.508%
86	21.286	3089	3094	3100	rBV	2225365	2722553	57.74%	5.059%
87	23.380	3442	3450	3457	rBV2	2631324	4714981	100.00%	8.761%
88	23.521	3467	3474	3481	rVB	1401953	2613663	55.43%	4.856%

Sum of corrected areas: 53818600

LSC Report - Integrated Chromatogram

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Data File : BM013029.D
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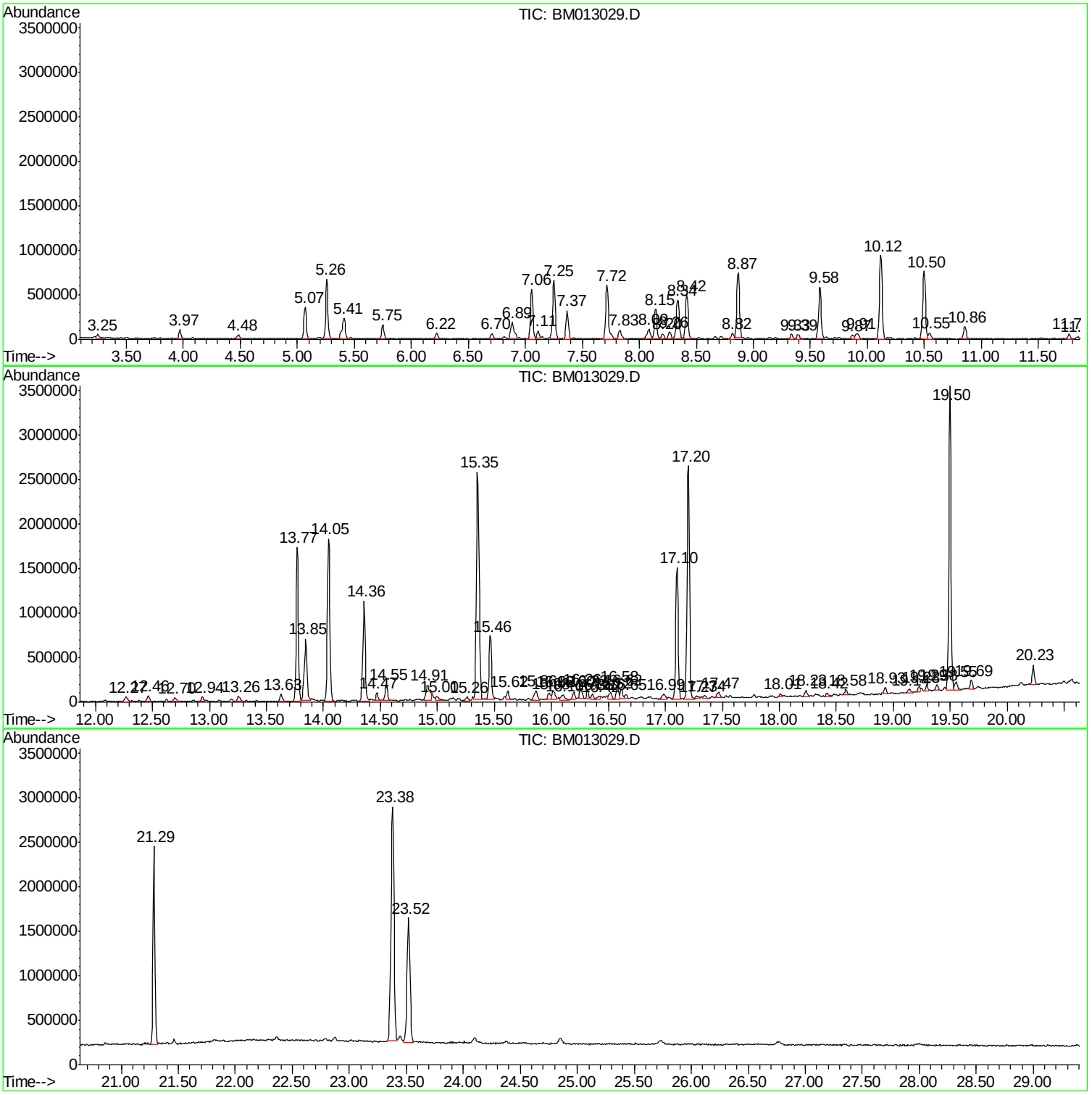
Operator : SJ/JU
Sample : I6405-04
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BE2S6

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM111617.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\HPCHEM1\BNA M\DATA\BM111717\
Data File : BM013029.D
Acq On : 18 Nov 2017 08:39
Operator : SJ/JU
Sample : I6405-04
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BE2S6

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM111617.M
Quant Title : SVOA CALIBRATION

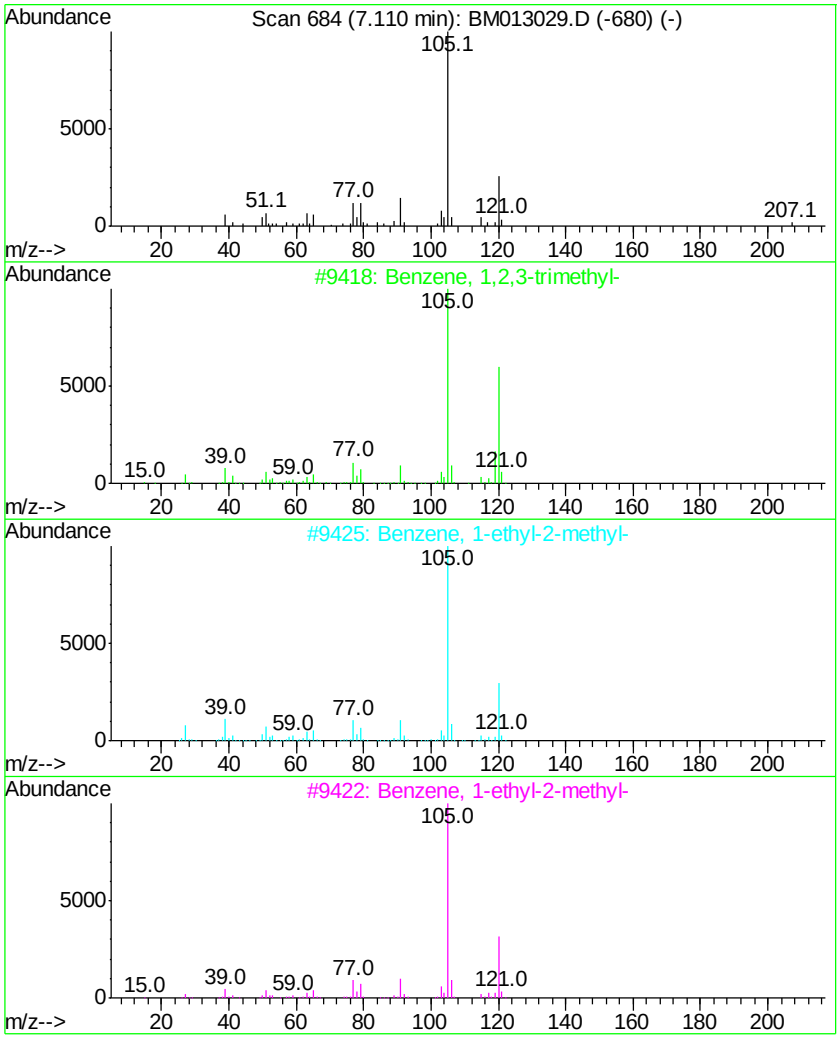
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

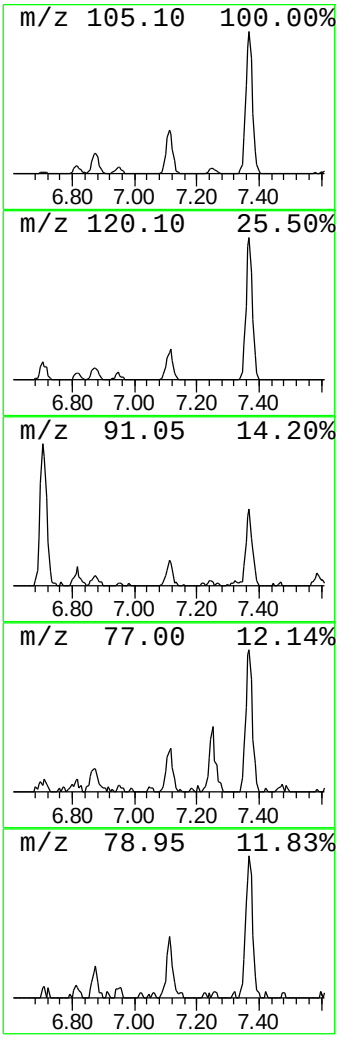
****

Peak Number 6 Benzene, 1,2,3-trimethyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.11	2.13 ng/ul	110919	1,4-Dichlorobenzene-d4	7.72

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	90
2	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	90
3	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	90
4	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	87
5	Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	87





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM111717\
Data File : BM013029.D
Acq On : 18 Nov 2017 08:39
Operator : SJ/JU
Sample : I6405-04
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BE2S6

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM111617.M
Quant Title : SVOA CALIBRATION

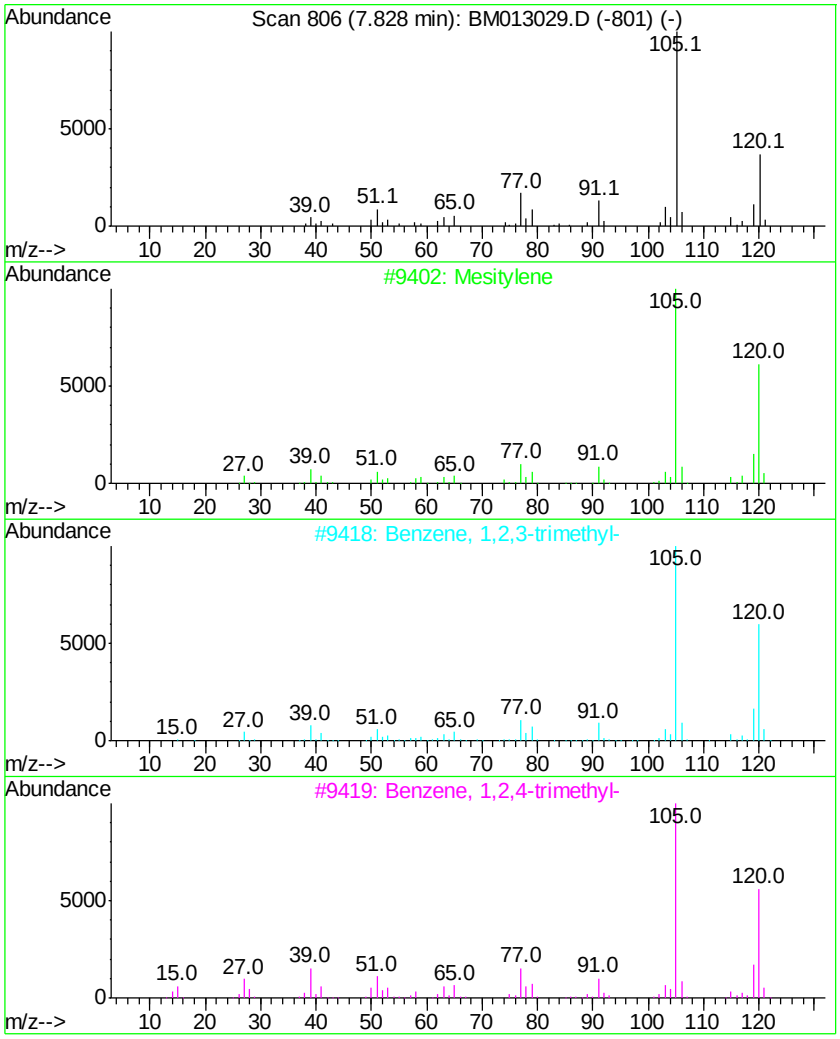
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

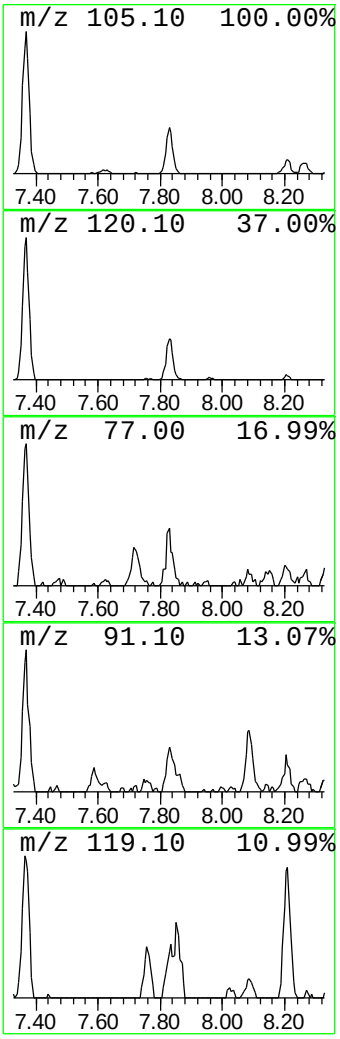
****

Peak Number 8 Mesitylene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.83	3.43 ng/ul	178965	1,4-Dichlorobenzene-d4	7.72

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Mesitylene	120	C9H12	000108-67-8	94
2	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
3	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
4	Mesitylene	120	C9H12	000108-67-8	91
5	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91





Data Path : Z:\HPCHEM1\BNA M\DATA\BM111717\
Data File : BM013029.D
Acq On : 18 Nov 2017 08:39
Operator : SJ/JU
Sample : I6405-04
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BE2S6

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM111617.M
Quant Title : SVOA CALIBRATION

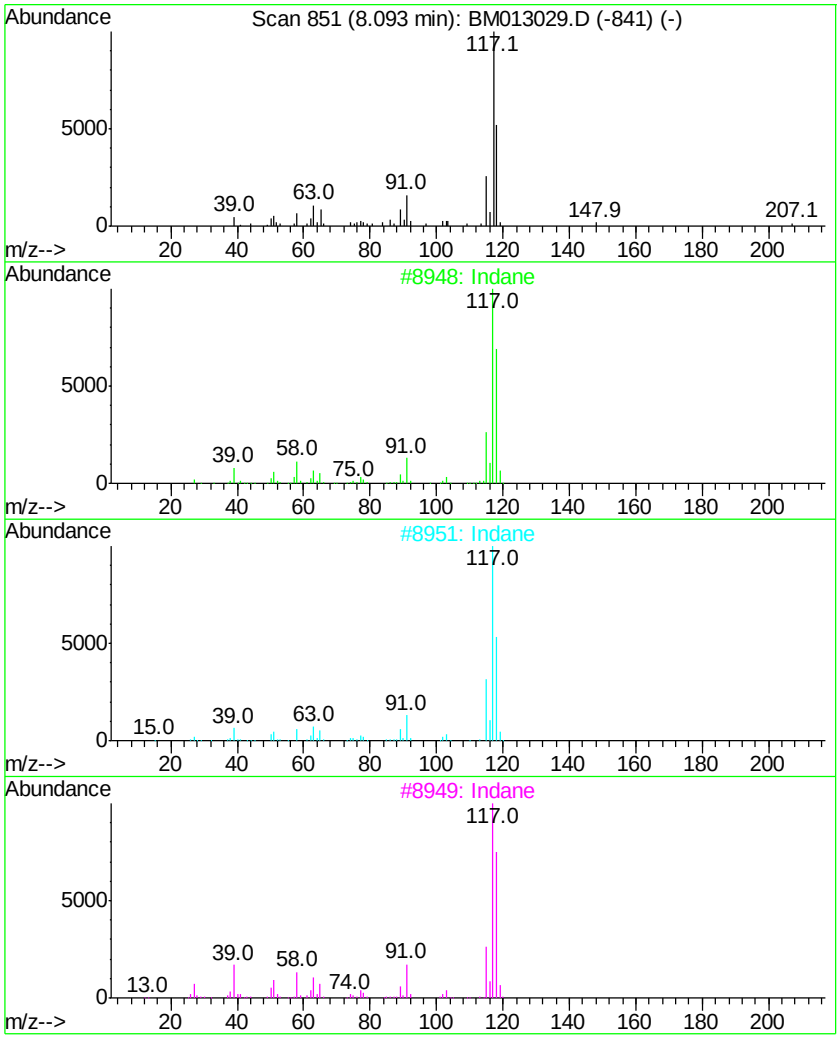
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

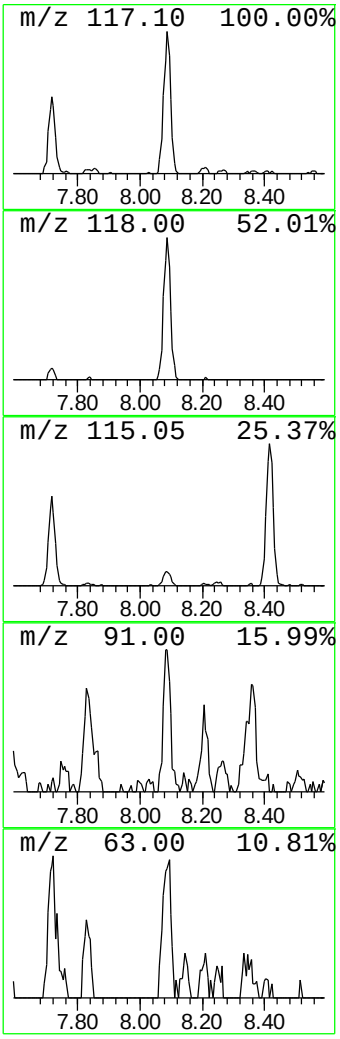
****

Peak Number 9 Indane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.09	4.24 ng/ul	221080	1,4-Dichlorobenzene-d4	7.72

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	87
2	Indane	118	C9H10	000496-11-7	87
3	Indane	118	C9H10	000496-11-7	87
4	Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	81
5	Indane	118	C9H10	000496-11-7	81





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Acq On : 18 Nov 2017 08:39
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ALS Vial : 39 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BE2S6

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM111617.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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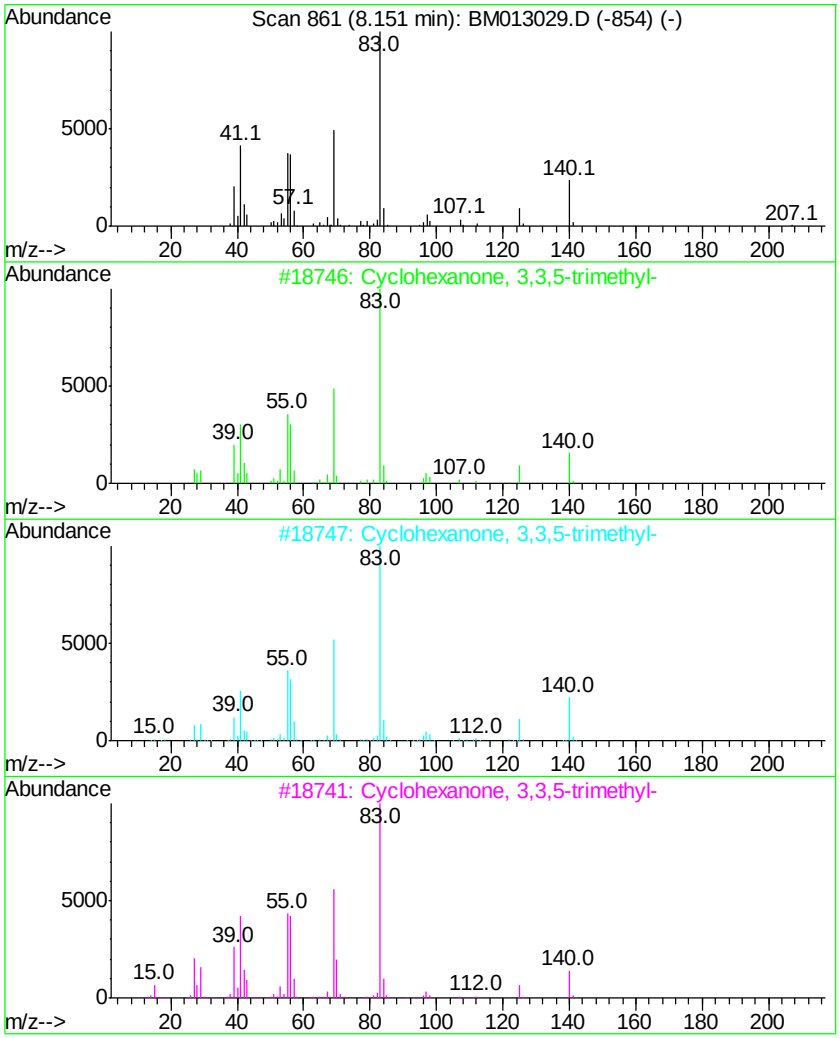
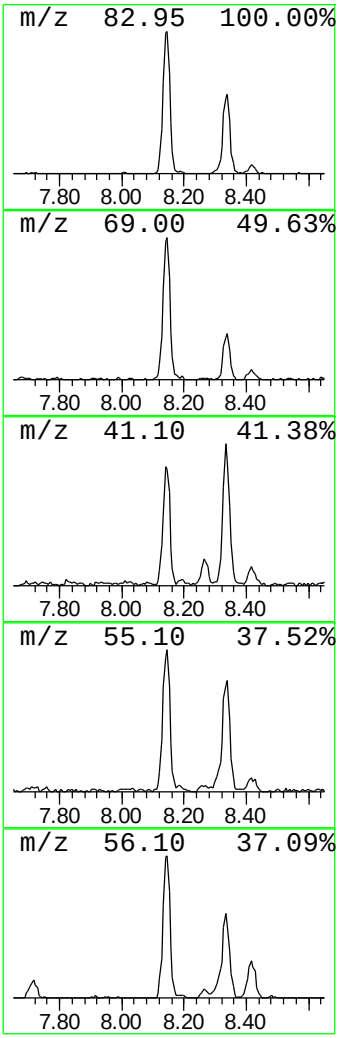
Peak Number 10 Cyclohexanone, 3,3,5-trimet... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.15	9.87 ng/ul	514776	1,4-Dichlorobenzene-d4	7.72

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	95
2	Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	94
3	Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	91
4	Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	78
5	Cyclohexane, butyl-	140	C10H20	001678-93-9	64





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM111717\
Data File : BM013029.D
Acq On : 18 Nov 2017 08:39
Operator : SJ/JU
Sample : I6405-04
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BE2S6

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM111617.M
Quant Title : SVOA CALIBRATION

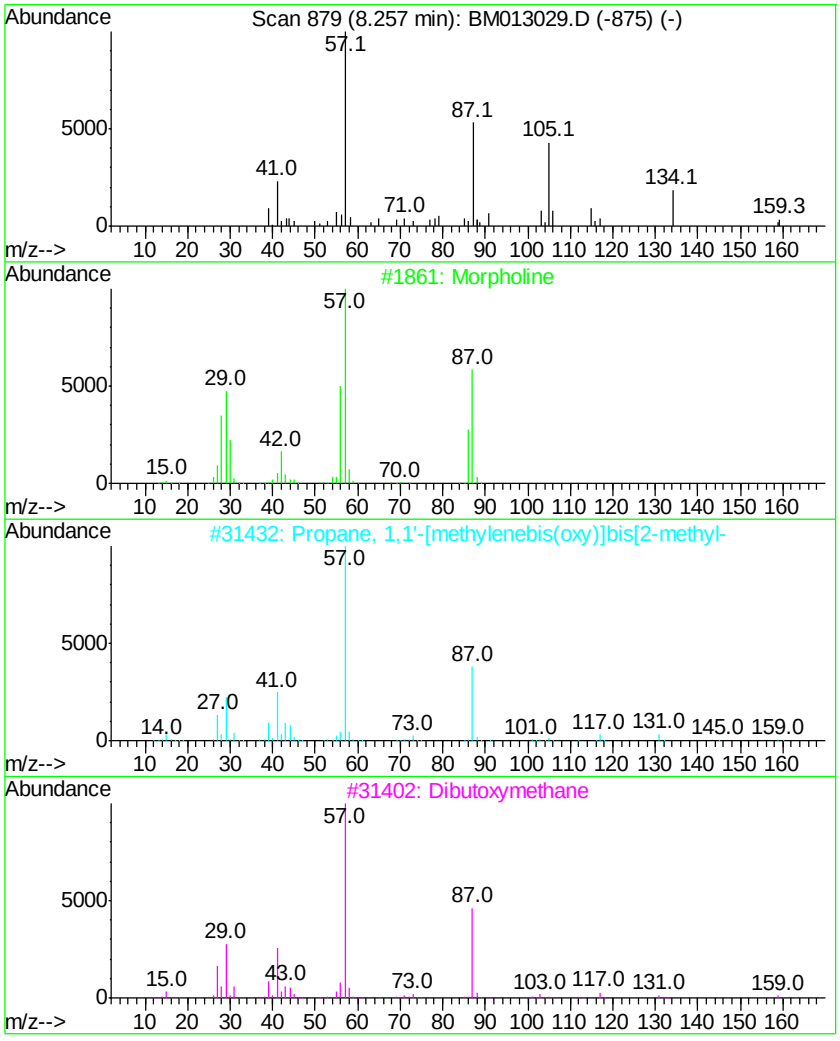
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

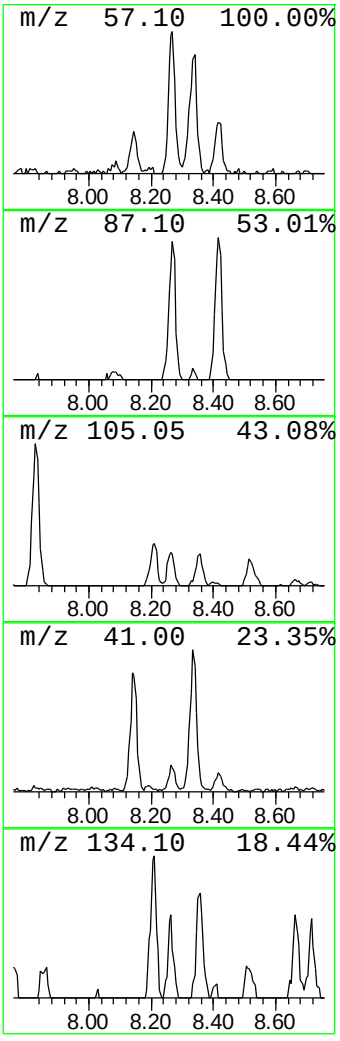
****

Peak Number 11 unknown-01 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.26	2.39 ng/ul	124601	1,4-Dichlorobenzene-d4	7.72

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Morpholine	87	C4H9NO	000110-91-8	43
2	Propane, 1,1'-[methylenebis(oxy)...	160	C9H20O2	002568-91-4	37
3	Dibutoxymethane	160	C9H20O2	002568-90-3	37
4	Morpholine	87	C4H9NO	000110-91-8	32
5	Morpholine	87	C4H9NO	000110-91-8	32





Data Path : Z:\HPCHEM1\BNA M\DATA\BM111717\
Data File : BM013029.D
Acq On : 18 Nov 2017 08:39
Operator : SJ/JU
Sample : I6405-04
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BE2S6

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM111617.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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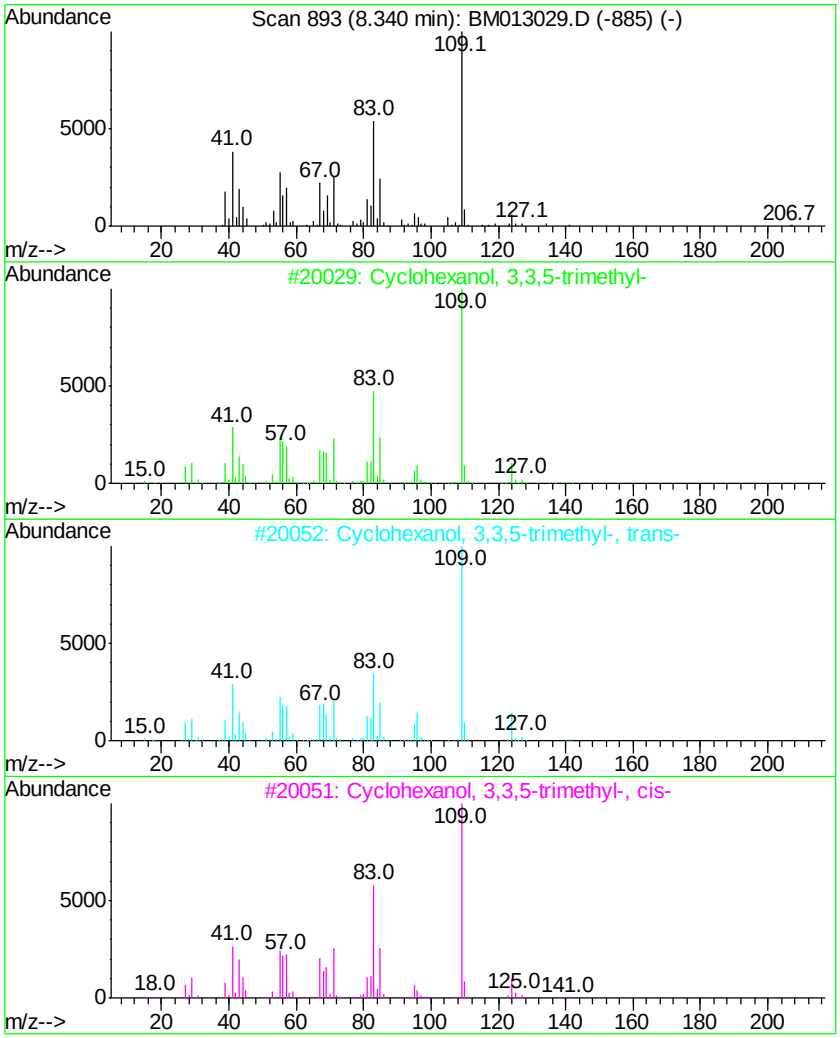
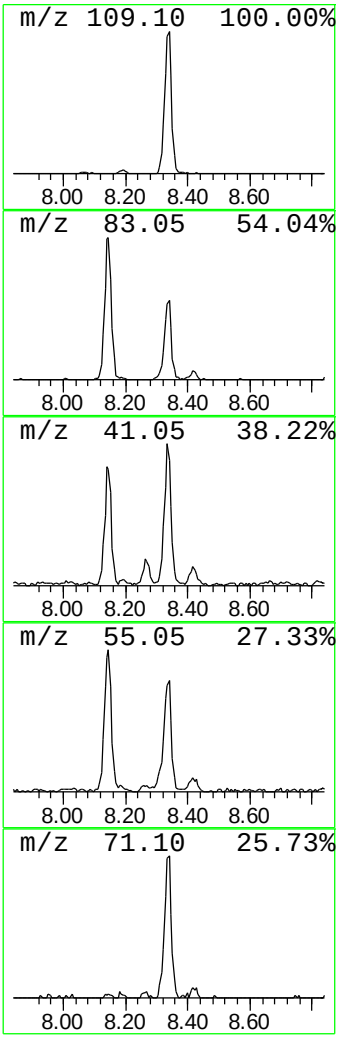
Peak Number 12 Cyclohexanol, 3,3,5-trimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.34	14.64 ng/ul	763422	1,4-Dichlorobenzene-d4	7.72

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexanol, 3,3,5-trimethyl-	142	C9H18O	000116-02-9	72
2	Cyclohexanol, 3,3,5-trimethyl-, ...	142	C9H18O	000767-54-4	64
3	Cyclohexanol, 3,3,5-trimethyl-, ...	142	C9H18O	000933-48-2	53
4	Cyclohexanol, 3,3,5-trimethyl-, ...	184	C11H20O2	024691-16-5	43
5	2-(1-Cyclopent-1-enyl-1-methylet...	192	C13H20O	1000190-57-4	43





Data Path : Z:\HPCHEM1\BNA M\DATA\BM111717\
Data File : BM013029.D
Acq On : 18 Nov 2017 08:39
Operator : SJ/JU
Sample : I6405-04
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BE2S6

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM111617.M
Quant Title : SVOA CALIBRATION

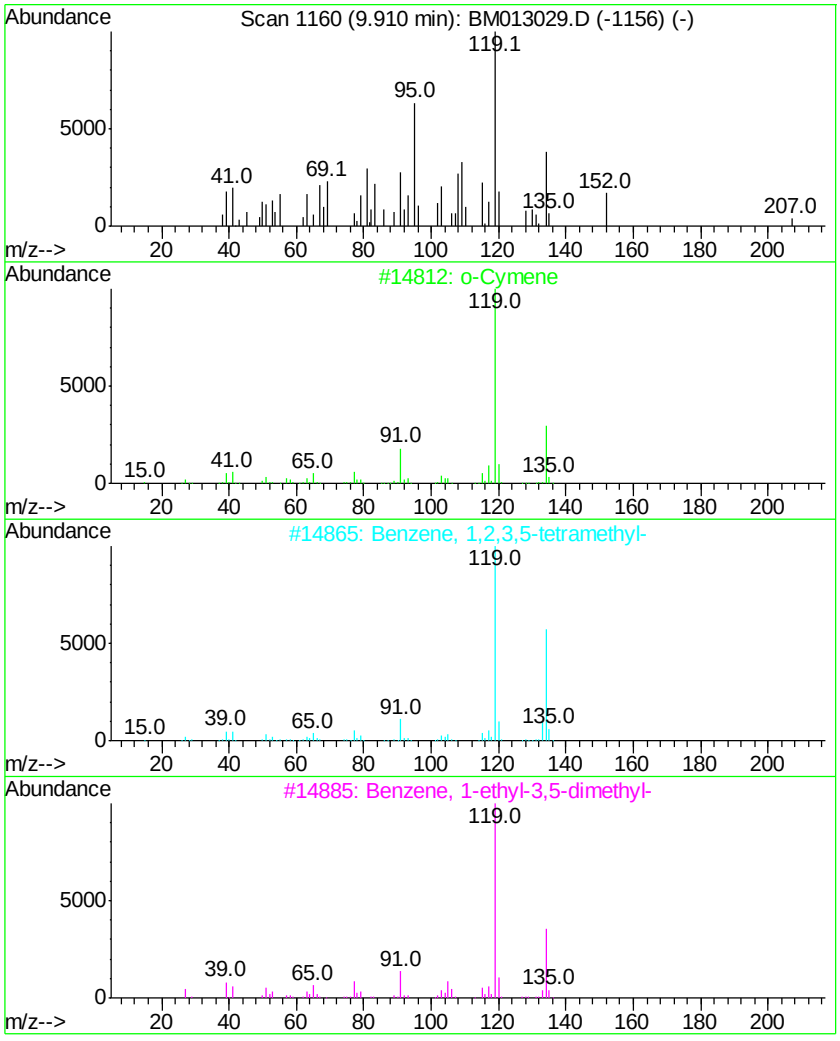
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

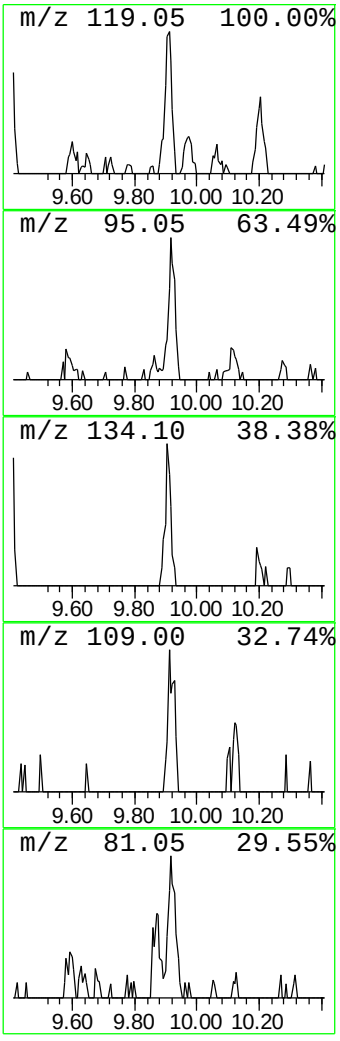
****

Peak Number 13 o-Cymene Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.91	2.06 ng/ul	130756	Naphthalene-d8	10.50

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Cymene	134	C10H14	000527-84-4	46
2	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	46
3	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	43
4	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	43
5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	43





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM111717\
Data File : BM013029.D
Acq On : 18 Nov 2017 08:39
Operator : SJ/JU
Sample : I6405-04
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BE2S6

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM111617.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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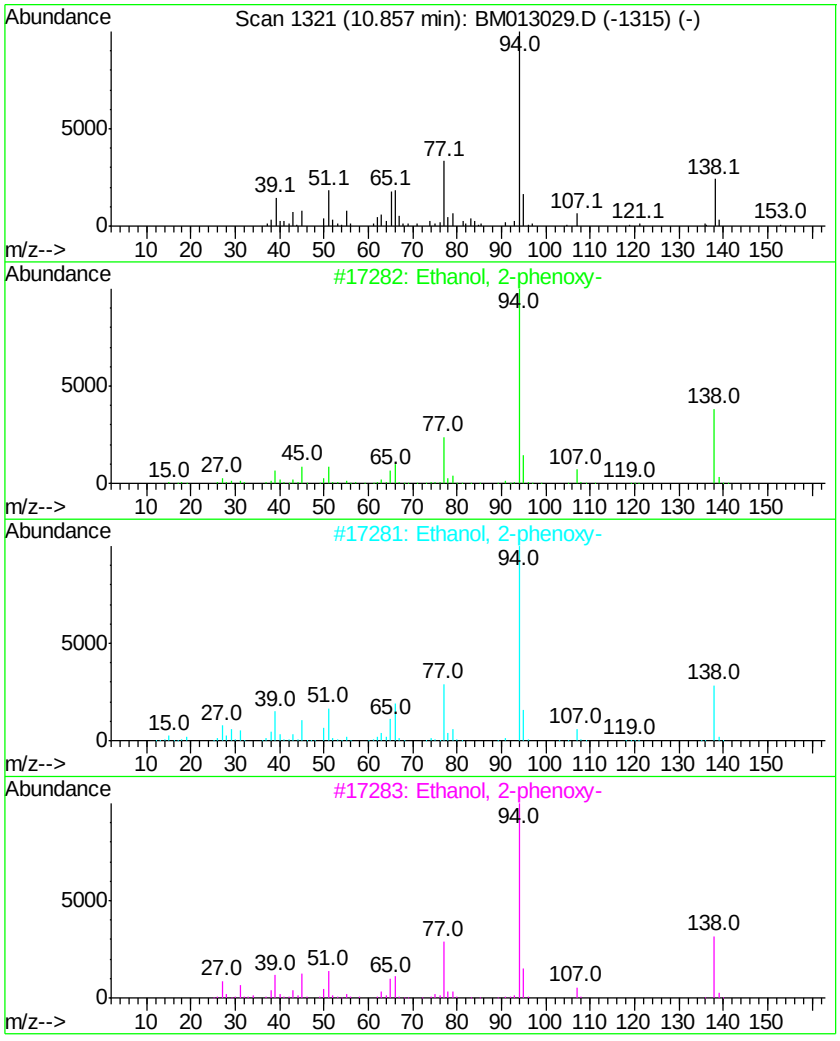
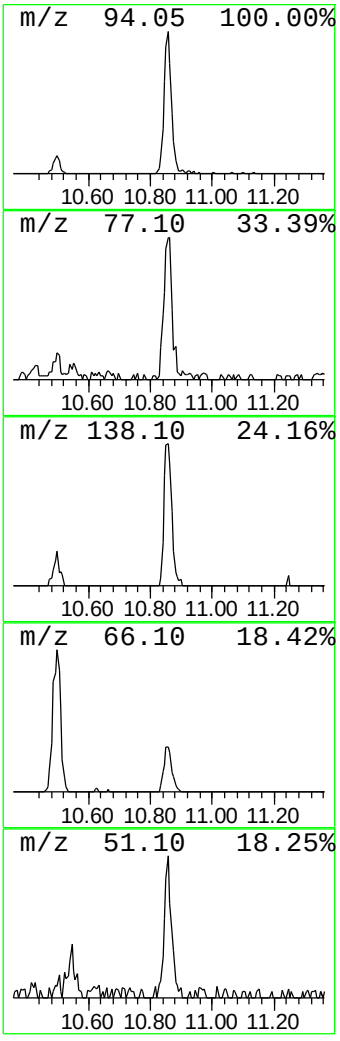
Peak Number 14 Ethanol, 2-phenoxy- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.86	3.45 ng/ul	219657	Naphthalene-d8	10.50

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-phenoxy-	138	C8H10O2	000122-99-6	91
2	Ethanol, 2-phenoxy-	138	C8H10O2	000122-99-6	91
3	Ethanol, 2-phenoxy-	138	C8H10O2	000122-99-6	91
4	2-Pyridinecarboxylic acid, 3-amino-	138	C6H6N2O2	001462-86-8	53
5	Benzene, (1,1-dimethylethoxy)-	150	C10H14O	006669-13-2	53





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Acq On : 18 Nov 2017 08:39
Operator : SJ/JU
Sample : I6405-04
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BE2S6

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM111617.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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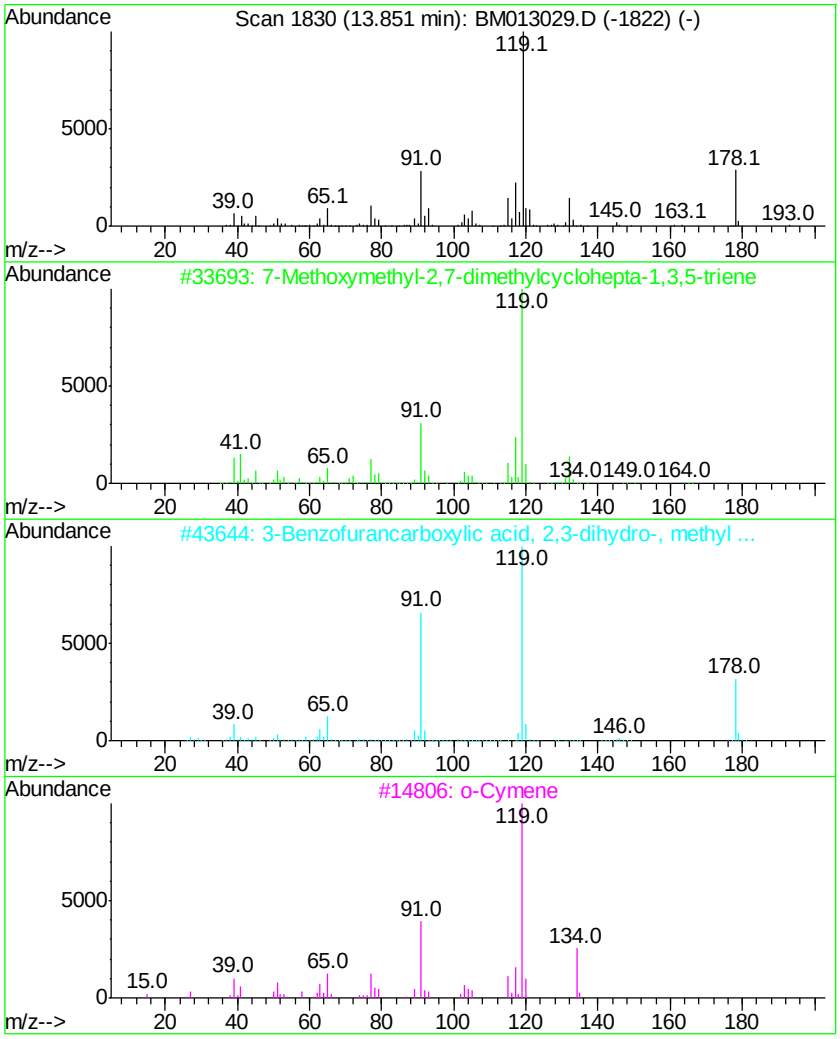
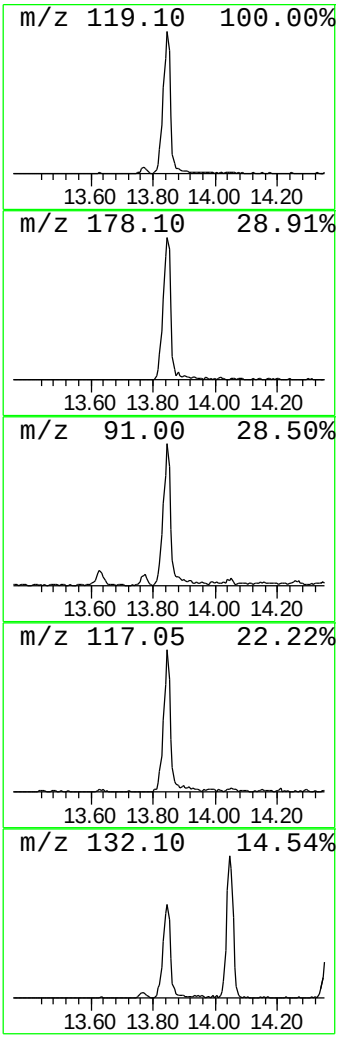
Peak Number 15 7-Methoxymethyl-2,7-dimethy... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.85	12.75 ng/ul	1102380	Acenaphthene-d10	14.36

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	7-Methoxymethyl-2,7-dimethylcycl...	164	C11H16O	073992-48-0	64
2	3-Benzofurancarboxylic acid, 2,3...	178	C10H10O3	039891-56-0	64
3	o-Cymene	134	C10H14	000527-84-4	53
4	1-(2-Methoxy-1-methylethyl)-2-me...	164	C11H16O	1000210-02-1	53
5	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	53





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Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BE2S6

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM111617.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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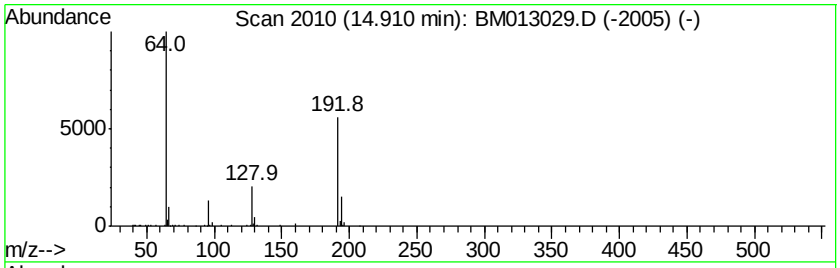
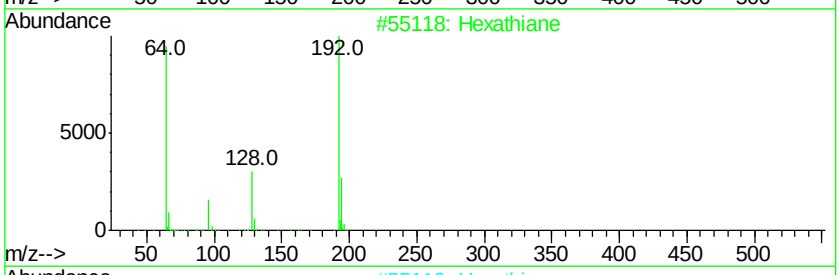
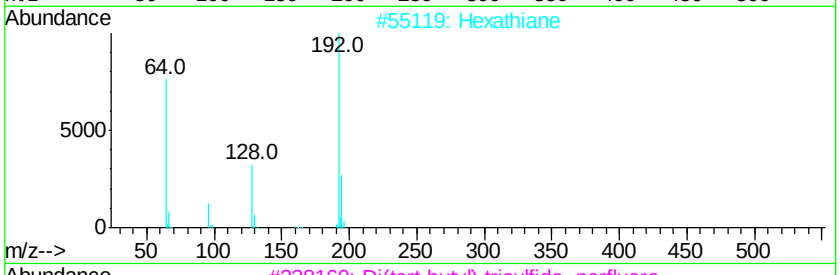
Peak Number 16 Hexathiane Concentration Rank 9

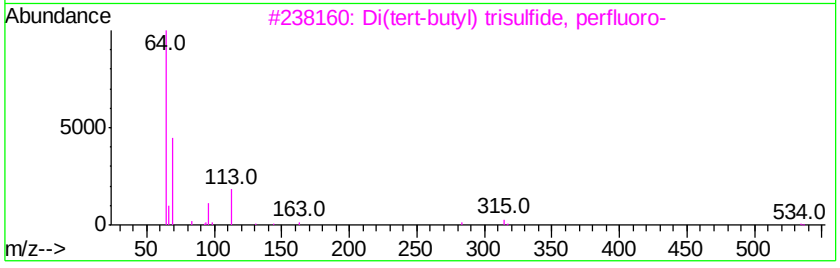
R.T.	EstConc	Area	Relative to ISTD	R.T.
14.91	4.50 ng/ul	389363	Acenaphthene-d10	14.36

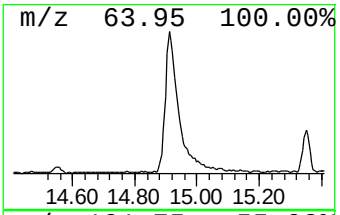
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexathiane	192	S6	013798-23-7	94
2	Hexathiane	192	S6	013798-23-7	91
3	Di(tert-butyl) trisulfide, perfl...	534	C8F18S3	016005-64-4	39
4	1,2-Phenylene diisothiocyanate	192	C8H4N2S2	071105-17-4	9
5	Sulfur dioxide	64	O2S	007446-09-5	7

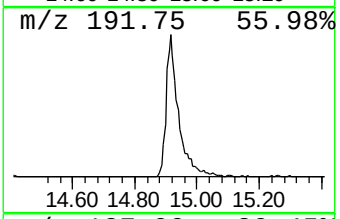


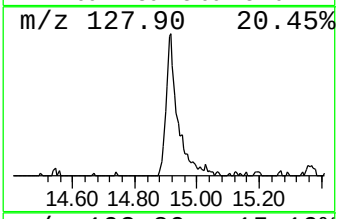


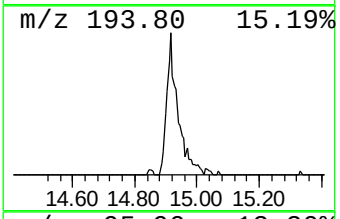


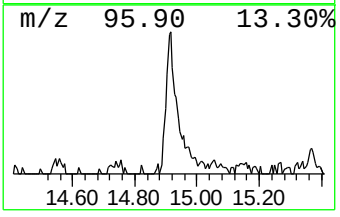












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Instrument : BNA_M
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Quant Title : SVOA CALIBRATION

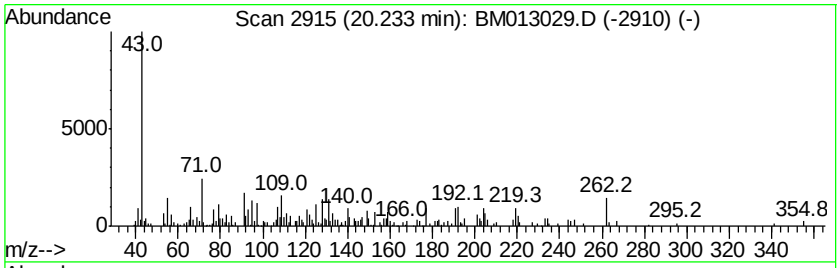
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

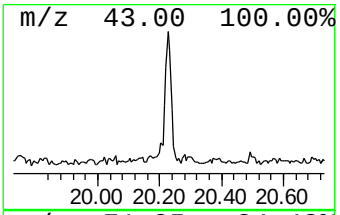
****

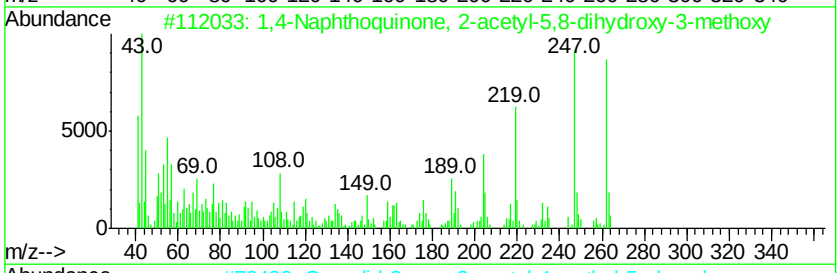
Peak Number 17 unknown-02 Concentration Rank 17

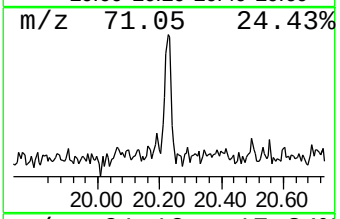
R.T.	EstConc	Area	Relative to ISTD	R.T.
20.23	2.01 ng/ul	273545	Chrysene-d12	21.29

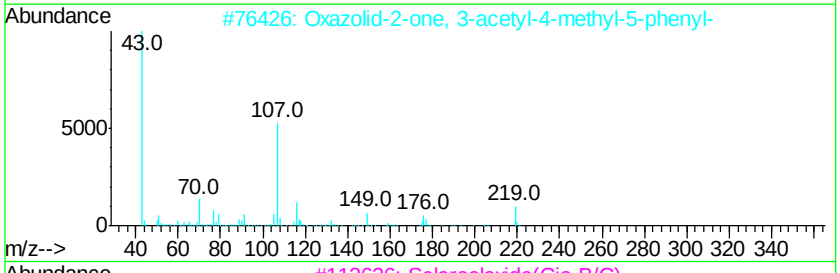
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,4-Naphthoquinone, 2-acetyl-5,8...	262	C13H10O6	014090-55-2	35
2	Oxazolid-2-one, 3-acetyl-4-methy...	219	C12H13NO3	063204-87-5	22
3	Sclareoloxide(Cis-B/C)	262	C18H30O	1000215-78-4	20
4	3,4-Heptadien-2-one, 3-cyclopent...	192	C13H20O	063922-50-9	18
5	2,4-Pentadienoic acid, 5-(8-hydr...	280	C15H20O5	024394-14-7	16

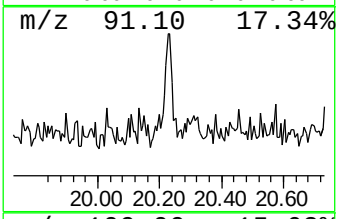


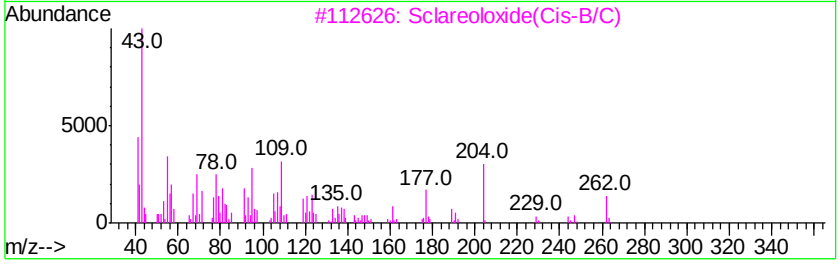


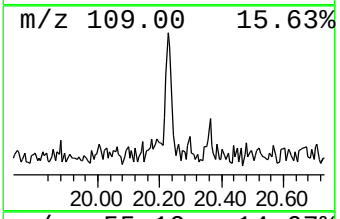


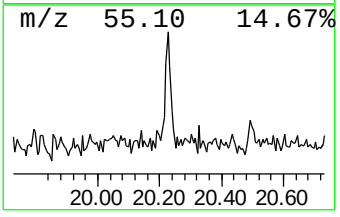












Tentatively Identified Compound (LSC) summary

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Misc :
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Instrument : BNA_M
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Benzene, 1,2,3-tr...	7.11	2.1	ng/ul	110919	1	7.72	1042670	20.0
Mesitylene	7.83	3.4	ng/ul	178965	1	7.72	1042670	20.0
Indane	8.09	4.2	ng/ul	221080	1	7.72	1042670	20.0
Cyclohexanone, 3,...	8.15	9.9	ng/ul	514776	1	7.72	1042670	20.0
unknown-01	8.26	2.4	ng/ul	124601	1	7.72	1042670	20.0
Cyclohexanol, 3,3...	8.34	14.6	ng/ul	763422	1	7.72	1042670	20.0
o-Cymene	9.91	2.1	ng/ul	130756	2	10.50	1271750	20.0
Ethanol, 2-phenoxy-	10.86	3.5	ng/ul	219657	2	10.50	1271750	20.0
7-Methoxymethyl-2...	13.85	12.8	ng/ul	1102380	3	14.36	1728900	20.0
Hexathiane	14.91	4.5	ng/ul	389363	3	14.36	1728900	20.0
unknown-02	20.23	2.0	ng/ul	273545	5	21.29	2722550	20.0