LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111723\
Data File : BM042905.D
Acq On : 19 Nov 2023 14:24
Operator : MA/JU
Sample : 05370-16
Misc :
ALS Vial : 63 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : GCDR0

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % 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:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM111323.MA.M
Title : SVOA CALIBRATION

Signal : TIC: BM042905.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.169	28	31	40	rVB4	9515	14703	1.37%	0.137%
2	3.616	104	107	123	rBV	37558	78884	7.34%	0.737%
3	4.922	327	329	332	rBV4	1411	1643	0.15%	0.015%
4	5.110	359	361	365	rVB3	1447	1533	0.14%	0.014%
5	5.669	455	456	462	rBV4	1407	1784	0.17%	0.017%
6	5.869	485	490	500	rVB6	7067	16297	1.52%	0.152%
7	6.187	538	544	548	rBV2	9138	13813	1.28%	0.129%
8	6.992	676	681	690	rBV5	15551	40737	3.79%	0.380%
9	7.163	704	710	719	rBV	116841	210813	19.61%	1.968%
10	7.334	733	739	752	rBV2	120509	223794	20.81%	2.090%
11	7.569	778	779	784	rVB4	1111	1515	0.14%	0.014%
12	7.810	814	820	832	rBV	125392	229459	21.34%	2.143%
13	8.539	938	944	954	rBV	58925	127249	11.83%	1.188%
14	9.016	1018	1025	1042	rBV	151366	302488	28.13%	2.825%
15	9.728	1140	1146	1161	rBV	120348	236773	22.02%	2.211%
16	10.251	1229	1235	1256	rBV	129450	297021	27.62%	2.773%
17	10.404	1260	1261	1263	rVV	2543	2102	0.20%	0.020%
18	10.433	1263	1266	1272	rVV5	3933	7594	0.71%	0.071%
19	10.516	1277	1280	1285	rVB3	3276	6205	0.58%	0.058%
20	10.622	1291	1298	1309	rBV	176781	299584	27.86%	2.797%
21	10.804	1324	1329	1351	rVV	88348	261663	24.34%	2.443%
22	10.992	1355	1361	1372	rVV9	9800	25411	2.36%	0.237%
23	11.104	1375	1380	1382	rBV6	1514	2125	0.20%	0.020%
24	11.139	1382	1386	1387	rBV3	3441	3684	0.34%	0.034%
25	11.157	1387	1389	1395	rVB5	3309	5169	0.48%	0.048%
26	11.245	1403	1404	1408	rBV3	1392	1554	0.14%	0.015%
27	11.333	1415	1419	1423	rVV3	3143	3061	0.28%	0.029%
28	11.416	1423	1433	1440	rVV6	10440	32747	3.05%	0.306%
29	11.492	1442	1446	1451	rVV5	5215	10919	1.02%	0.102%
30	11.545	1451	1455	1465	rVV6	8716	19900	1.85%	0.186%
31	11.686	1475	1479	1480	rVV3	1820	2181	0.20%	0.020%
32	11.745	1483	1489	1492	rBV6	995	2046	0.19%	0.019%
33	11.845	1501	1506	1510	rVB3	1444	2219	0.21%	0.021%
34	12.233	1568	1572	1576	rBV5	1441	2282	0.21%	0.021%
35	12.322	1582	1587	1595	rBV2	10652	18682	1.74%	0.174%
36	12.392	1595	1599	1600	rBV3	1668	1905	0.18%	0.018%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111723\
Data File : BM042905.D
Acq On : 19 Nov 2023 14:24
Operator : MA/JU
Sample : 05370-16
Misc :
ALS Vial : 63 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GCDR0

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % 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:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM111323.MA.M
Title : SVOA CALIBRATION

37	12.439	1605	1607	1613	rVB5	1684	3702	0.34%	0.035%
38	12.545	1619	1625	1631	rBV5	2701	4502	0.42%	0.042%
39	13.145	1721	1727	1732	rBV7	2467	4968	0.46%	0.046%
40	13.316	1748	1756	1762	rBV3	10543	18258	1.70%	0.170%
41	13.410	1767	1772	1784	rVB8	6369	14682	1.37%	0.137%
42	13.780	1830	1835	1844	rVV	15061	26710	2.48%	0.249%
43	13.892	1848	1854	1866	rVV	356307	500211	46.52%	4.671%
44	14.133	1889	1895	1896	rBV4	9933	13719	1.28%	0.128%
45	14.168	1896	1901	1918	rVB	363320	550720	51.22%	5.142%
46	14.468	1947	1952	1963	rBV	253090	369411	34.36%	3.449%
47	14.704	1987	1992	1993	rBV5	1525	2246	0.21%	0.021%
48	14.792	2004	2007	2009	rBV2	2898	3526	0.33%	0.033%
49	14.968	2034	2037	2044	rVB6	4516	6188	0.58%	0.058%
50	15.051	2048	2051	2058	rVB5	1525	1928	0.18%	0.018%
51	15.204	2074	2077	2082	rVB5	2202	2663	0.25%	0.025%
52	15.251	2082	2085	2087	rBV3	1438	2129	0.20%	0.020%
53	15.280	2087	2090	2096	rVV5	5494	9973	0.93%	0.093%
54	15.321	2096	2097	2103	rVV4	2518	3871	0.36%	0.036%
55	15.462	2115	2121	2132	rBV	459373	669474	62.26%	6.251%
56	15.615	2141	2147	2159	rBV	117394	207251	19.27%	1.935%
57	15.727	2164	2166	2172	rVB5	3451	4868	0.45%	0.045%
58	15.874	2187	2191	2195	rVV2	14076	19169	1.78%	0.179%
59	15.915	2195	2198	2201	rVV3	7479	9755	0.91%	0.091%
60	15.951	2201	2204	2206	rVV2	2751	3314	0.31%	0.031%
61	16.086	2225	2227	2233	rVB5	1251	1968	0.18%	0.018%
62	16.174	2236	2242	2251	rVB	36644	48158	4.48%	0.450%
63	16.268	2251	2258	2262	rBV	114376	156379	14.54%	1.460%
64	16.321	2262	2267	2274	rVV3	130128	256592	23.86%	2.396%
65	16.386	2274	2278	2282	rVV2	113421	185064	17.21%	1.728%
66	16.427	2282	2285	2290	rVV	64671	93864	8.73%	0.876%
67	16.486	2290	2295	2297	rVV	33065	54338	5.05%	0.507%
68	16.504	2297	2298	2301	rVV	31243	37011	3.44%	0.346%
69	16.533	2301	2303	2307	rVV	30586	37664	3.50%	0.352%
70	16.586	2307	2312	2319	rVV4	82494	150227	13.97%	1.403%
71	16.657	2319	2324	2329	rVV	89797	136313	12.68%	1.273%
72	16.727	2333	2336	2343	rVB	42674	61534	5.72%	0.575%
73	16.798	2346	2348	2353	rVB5	2062	2605	0.24%	0.024%
74	16.845	2353	2356	2357	rBV3	1883	2112	0.20%	0.020%
75	16.862	2357	2359	2363	rBV4	1645	2204	0.20%	0.021%
76	16.945	2369	2373	2374	rBV3	2005	2309	0.21%	0.022%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111723\
Data File : BM042905.D
Acq On : 19 Nov 2023 14:24
Operator : MA/JU
Sample : 05370-16
Misc :
ALS Vial : 63 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : GCDR0

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % 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:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM111323.MA.M
Title : SVOA CALIBRATION

77	16.980	2376	2379	2384	rVB6	3323	5012	0.47%	0.047%
78	17.056	2390	2392	2397	rVB3	3352	3825	0.36%	0.036%
79	17.221	2414	2420	2431	rBV	307991	441178	41.03%	4.120%
80	17.321	2431	2437	2451	rBV2	510927	764300	71.08%	7.137%
81	17.480	2461	2464	2467	rVB5	2653	3380	0.31%	0.032%
82	17.851	2523	2527	2533	rVB6	3383	4196	0.39%	0.039%
83	18.033	2554	2558	2562	rBV6	2776	6172	0.57%	0.058%
84	18.080	2562	2566	2570	rVV4	7481	11752	1.09%	0.110%
85	18.121	2570	2573	2579	rVB7	2778	4914	0.46%	0.046%
86	18.245	2591	2594	2595	rBV2	2439	2332	0.22%	0.022%
87	18.292	2598	2602	2609	rBV3	12638	19939	1.85%	0.186%
88	18.833	2689	2694	2699	rBV2	6629	12992	1.21%	0.121%
89	18.898	2699	2705	2713	rVB10	6893	16646	1.55%	0.155%
90	18.974	2713	2718	2722	rBV4	4328	7413	0.69%	0.069%
91	19.086	2735	2737	2741	rBV5	1715	2086	0.19%	0.019%
92	19.180	2750	2753	2756	rBV3	4860	5959	0.55%	0.056%
93	19.268	2763	2768	2773	rBV2	14749	30766	2.86%	0.287%
94	19.327	2776	2778	2779	rVV2	3155	2935	0.27%	0.027%
95	19.362	2782	2784	2790	rVB7	5130	8683	0.81%	0.081%
96	19.621	2822	2828	2837	rBV	730611	1006943	93.65%	9.402%
97	19.845	2864	2866	2867	rBV2	2303	1536	0.14%	0.014%
98	21.409	3127	3132	3144	rBV	338438	514090	47.81%	4.800%
99	23.609	3499	3506	3519	rBV	551875	1075246	100.00%	10.040%
100	23.762	3526	3532	3546	rVB	264360	566000	52.64%	5.285%

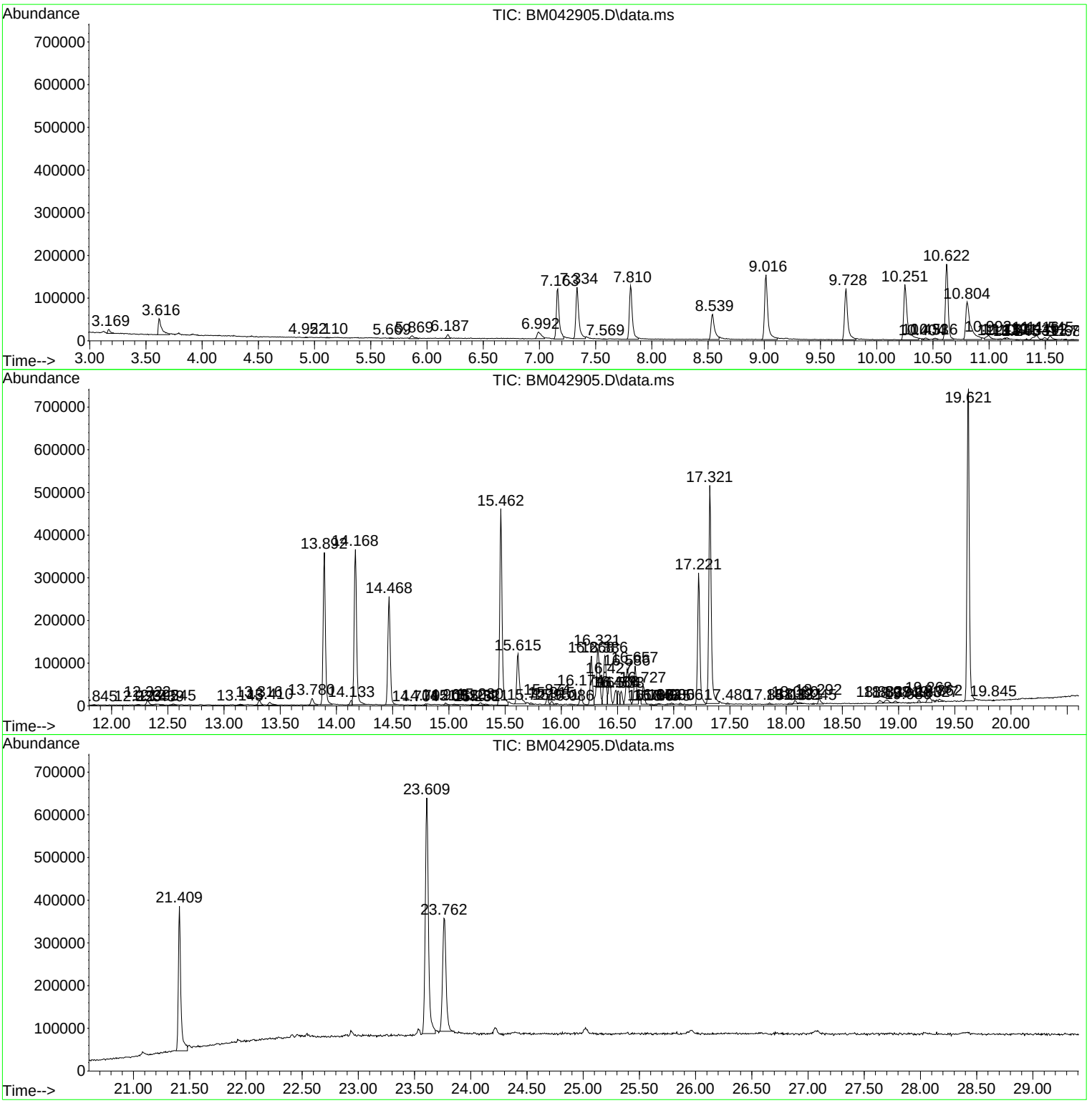
Sum of corrected areas: 10709414

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111723\
Data File : BM042905.D
Acq On : 19 Nov 2023 14:24
Operator : MA/JU
Sample : 05370-16
Misc :
ALS Vial : 63 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GCDR0

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM111323.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111723\
Data File : BM042905.D
Acq On : 19 Nov 2023 14:24
Operator : MA/JU
Sample : 05370-16
Misc :
ALS Vial : 63 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GCDR0

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM111323.MA.M
Quant Title : SVOA CALIBRATION

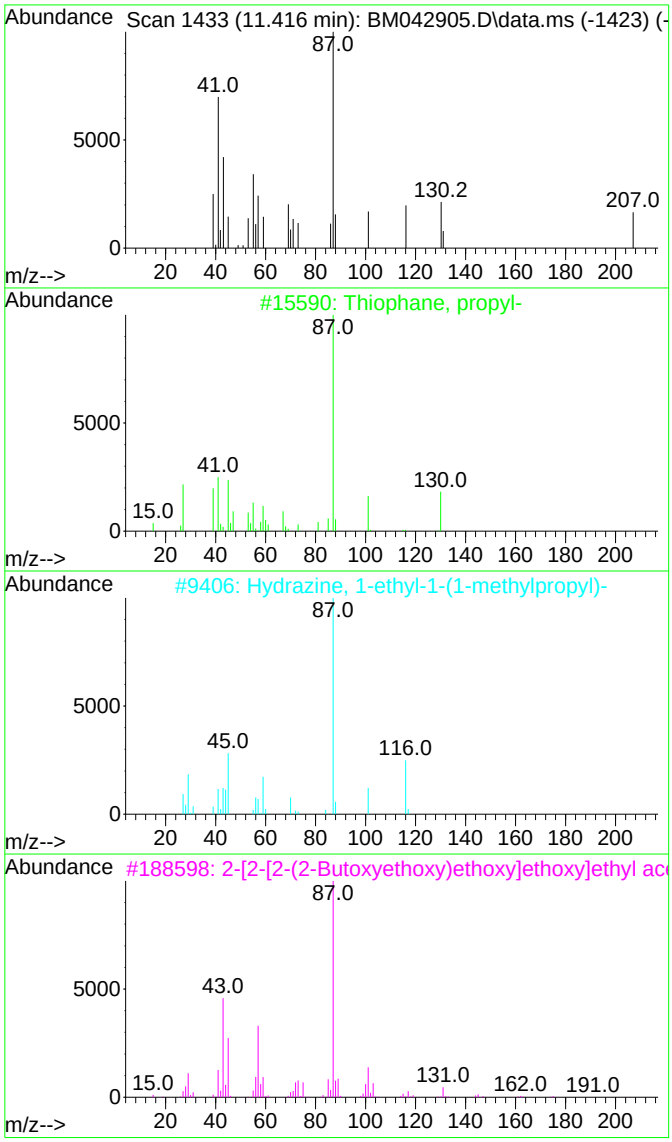
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

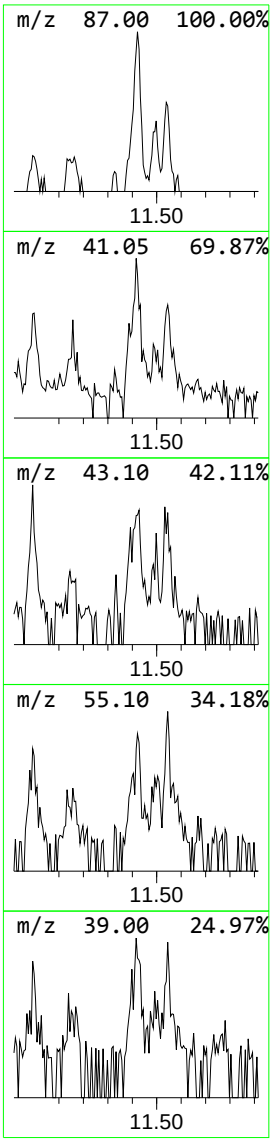
****

Peak Number 1 Thiophane, propyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.416	2.19 ng/ul	32747	Naphthalene-d8	10.622

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Thiophane, propyl-	130	C7H14S	001551-34-4	50
2	Hydrazine, 1-ethyl-1-(1-methylpr...	116	C6H16N2	020325-97-7	50
3	2-[2-[2-(2-Butoxyethoxy)ethoxy]e...	292	C14H28O6	1000351-93-9	47
4	.alpha.-D-Xylofuranoside, methyl...	178	C7H14O5	034338-86-8	47
5	Neodecanoic acid	172	C10H20O2	026896-20-8	45





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111723\
Data File : BM042905.D
Acq On : 19 Nov 2023 14:24
Operator : MA/JU
Sample : 05370-16
Misc :
ALS Vial : 63 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GCDR0

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM111323.MA.M
Quant Title : SVOA CALIBRATION

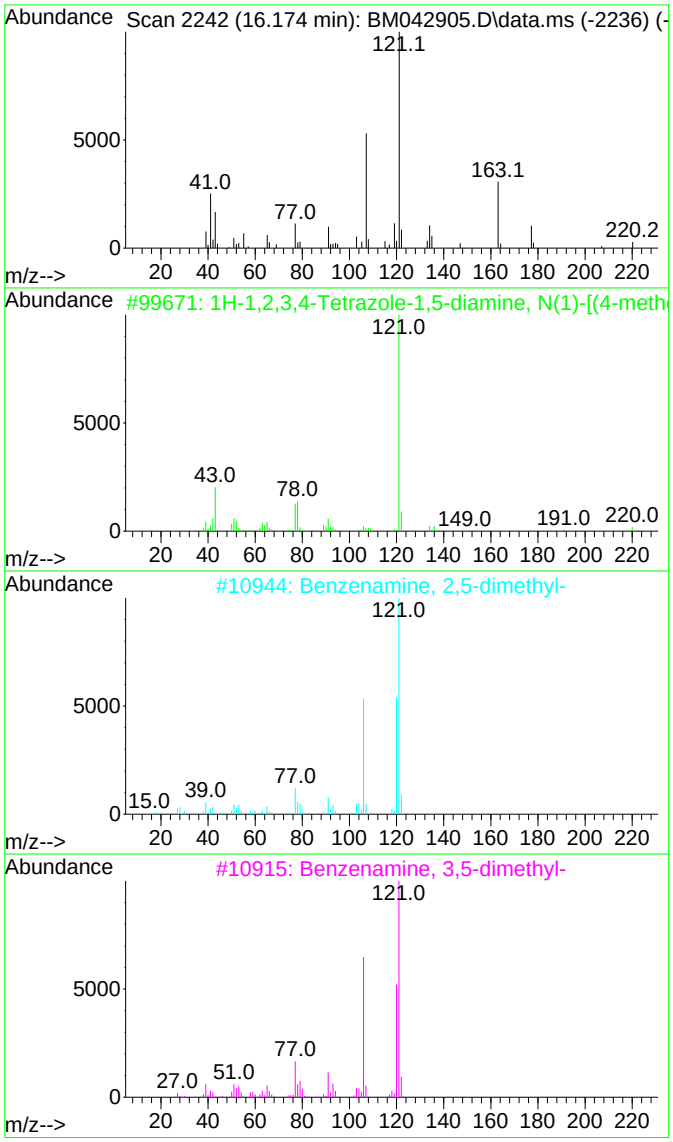
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

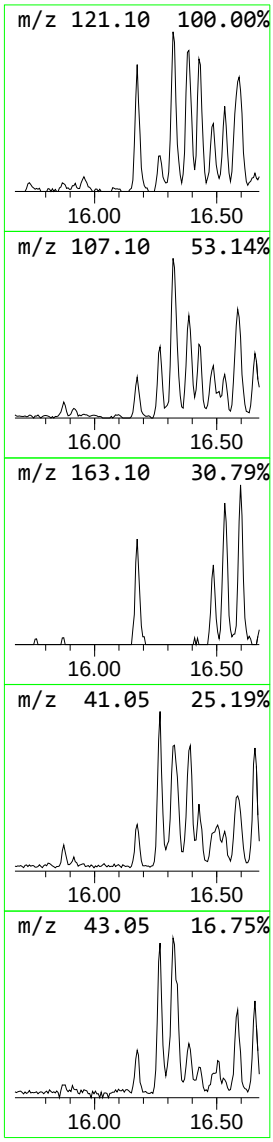
****

Peak Number 2 unknown-01 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.174	2.18 ng/ul	48158	Phenanthrene-d10	17.221

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-1,2,3,4-Tetrazole-1,5-diamine...	220	C9H12N6O	1000349-95-8	43
2	Benzenamine, 2,5-dimethyl-	121	C8H11N	000095-78-3	43
3	Benzenamine, 3,5-dimethyl-	121	C8H11N	000108-69-0	43
4	Acetamide, N-(2,4-dimethylphenyl)-	163	C10H13NO	002050-43-3	43
5	Carbonic acid, ethyl 4-isopropyl...	208	C12H16O3	1000331-34-2	38





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111723\
Data File : BM042905.D
Acq On : 19 Nov 2023 14:24
Operator : MA/JU
Sample : 05370-16
Misc :
ALS Vial : 63 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GCDR0

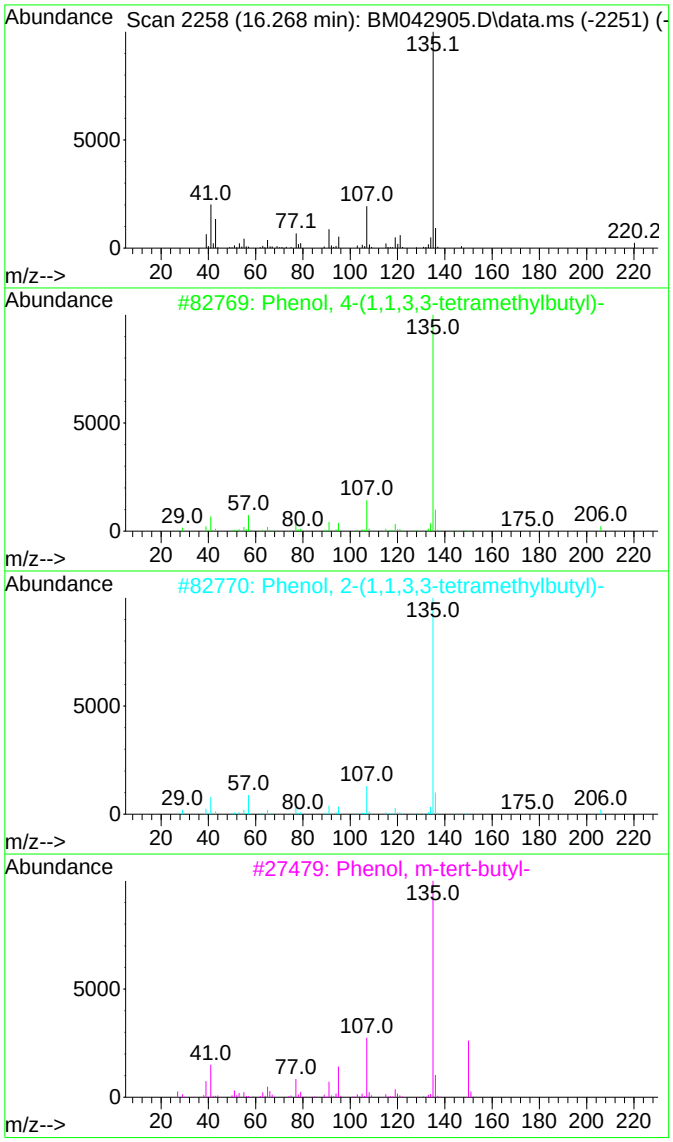
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM111323.MA.M
Quant Title : SVOA CALIBRATION

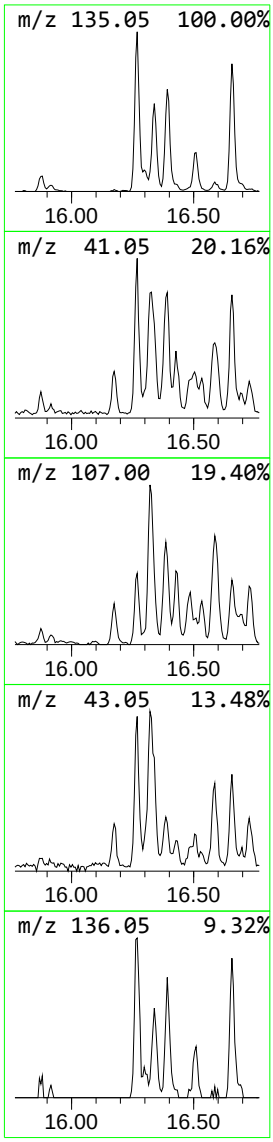
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Phenol, 4-(1,1,3,3-tetramet... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.268	7.09 ng/ul	156379	Phenanthrene-d10	17.221

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 4-(1,1,3,3-tetramethylbu...	206	C14H22O	000140-66-9	86
2	Phenol, 2-(1,1,3,3-tetramethylbu...	206	C14H22O	003884-95-5	78
3	Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	78
4	Phenol, 4-(1,1,3,3-tetramethylbu...	206	C14H22O	000140-66-9	72
5	((1,2-Diethylethylene)bis(p-phen...	354	C22H26O4	004547-76-6	72





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111723\
Data File : BM042905.D
Acq On : 19 Nov 2023 14:24
Operator : MA/JU
Sample : 05370-16
Misc :
ALS Vial : 63 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GCDR0

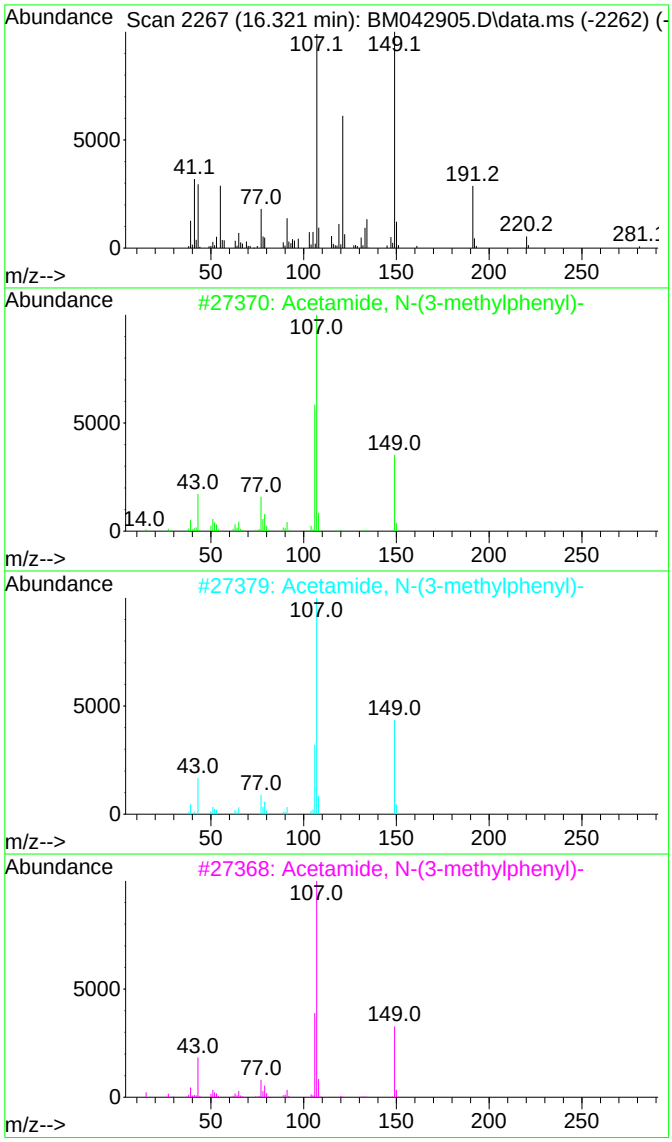
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM111323.MA.M
Quant Title : SVOA CALIBRATION

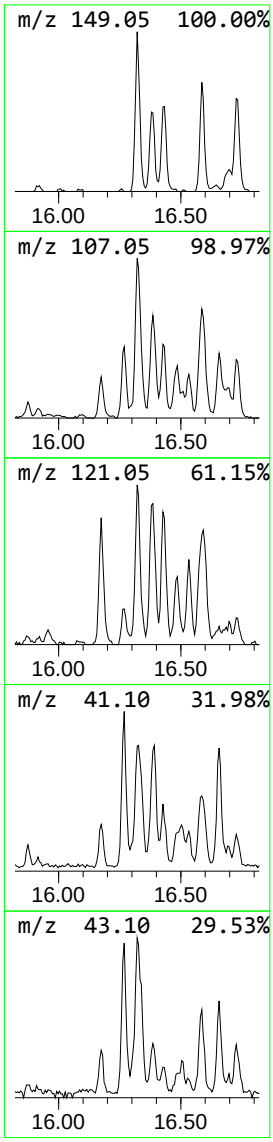
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Acetamide, N-(3-methylphenyl)- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.321	11.63 ng/ul	256592	Phenanthrene-d10	17.221

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetamide, N-(3-methylphenyl)-	149	C9H11NO	000537-92-8	58
2	Acetamide, N-(3-methylphenyl)-	149	C9H11NO	000537-92-8	58
3	Acetamide, N-(3-methylphenyl)-	149	C9H11NO	000537-92-8	58
4	Acetamide, N-(3-methylphenyl)-	149	C9H11NO	000537-92-8	58
5	Acetamide, N-(2-methylphenyl)-	149	C9H11NO	000120-66-1	46





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111723\
Data File : BM042905.D
Acq On : 19 Nov 2023 14:24
Operator : MA/JU
Sample : 05370-16
Misc :
ALS Vial : 63 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GCDR0

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM111323.MA.M
Quant Title : SVOA CALIBRATION

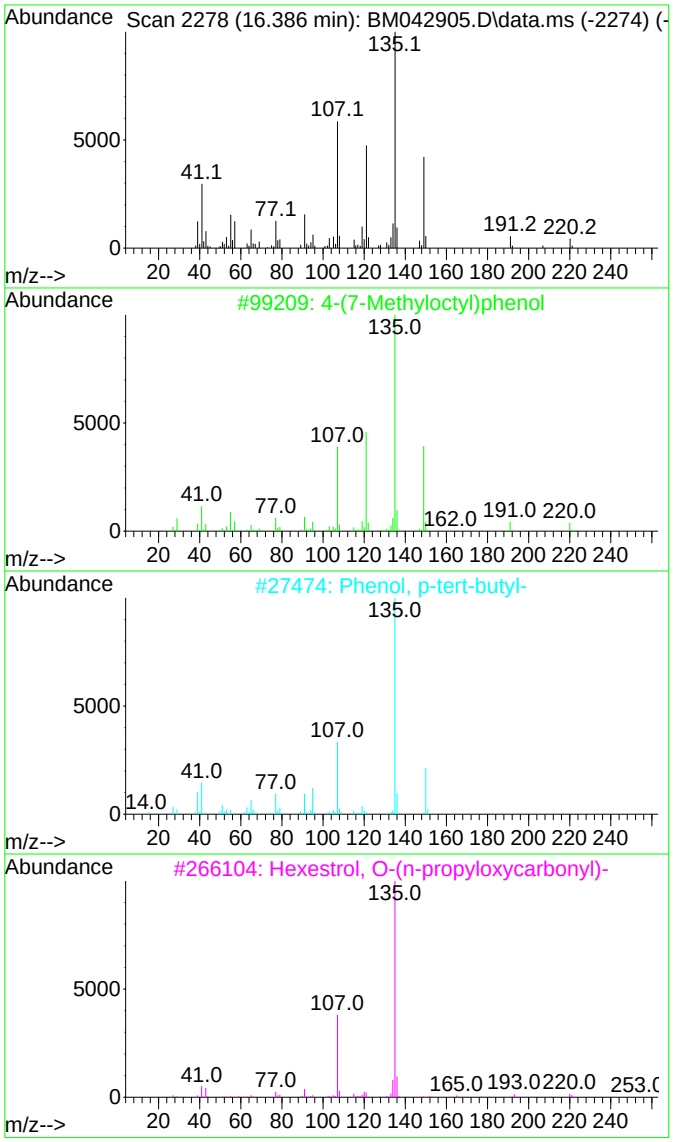
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

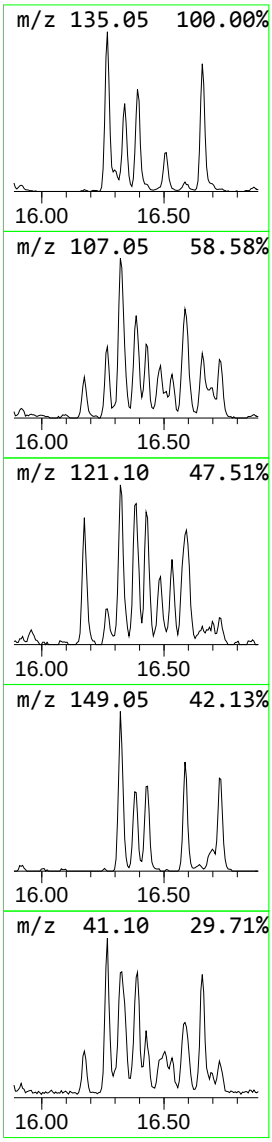
****

Peak Number 5 4-(7-Methyloctyl)phenol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.386	8.39 ng/ul	185064	Phenanthrene-d10	17.221

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-(7-Methyloctyl)phenol	220	C15H24O	024518-48-7	96
2	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	58
3	Hexestrol, O-(n-propyloxycarbonyl)-	356	C22H28O4	1000487-65-1	53
4	Hexestrol, O-methoxycarbonyl-	328	C20H24O4	1000487-66-3	53
5	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	53





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Data File : BM042905.D
Acq On : 19 Nov 2023 14:24
Operator : MA/JU
Sample : 05370-16
Misc :
ALS Vial : 63 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GCDR0

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM111323.MA.M
Quant Title : SVOA CALIBRATION

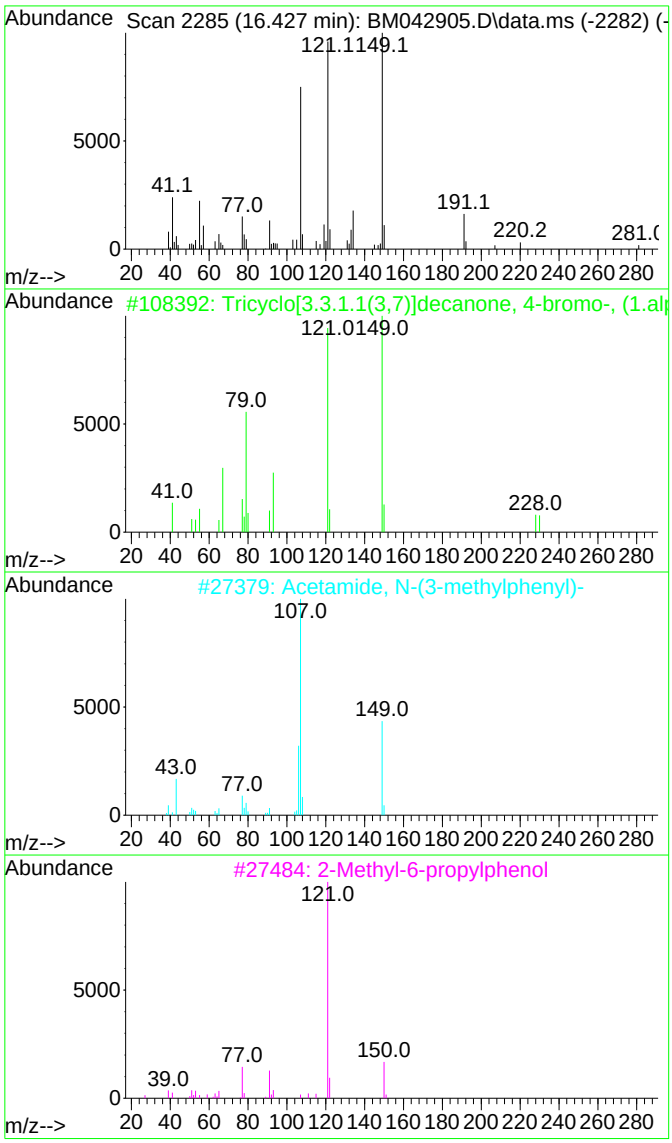
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

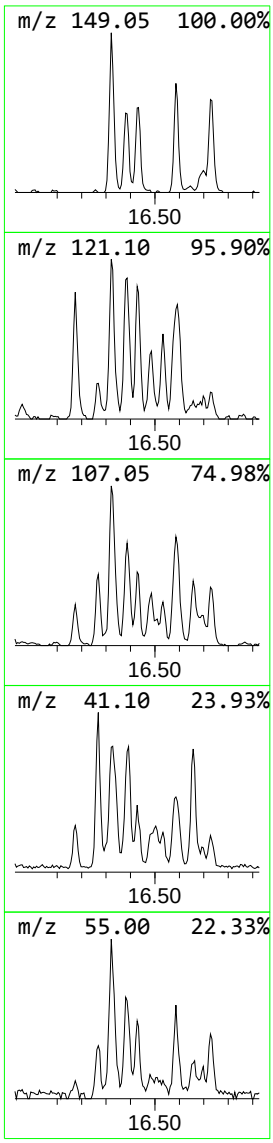
****

Peak Number 6 unknown-02 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.427	4.26 ng/ul	93864	Phenanthrene-d10	17.221

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tricyclo[3.3.1.1(3,7)]decanone, ...	228	C10H13BrO	032456-48-7	47
2	Acetamide, N-(3-methylphenyl)-	149	C9H11NO	000537-92-8	43
3	2-Methyl-6-propylphenol	150	C10H14O	003520-52-3	25
4	Phenol, 4-(1-methylpropyl)-	150	C10H14O	000099-71-8	25
5	3',4'-Formoxylidide	149	C9H11NO	006639-60-7	25





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111723\
Data File : BM042905.D
Acq On : 19 Nov 2023 14:24
Operator : MA/JU
Sample : 05370-16
Misc :
ALS Vial : 63 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GCDR0

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM111323.MA.M
Quant Title : SVOA CALIBRATION

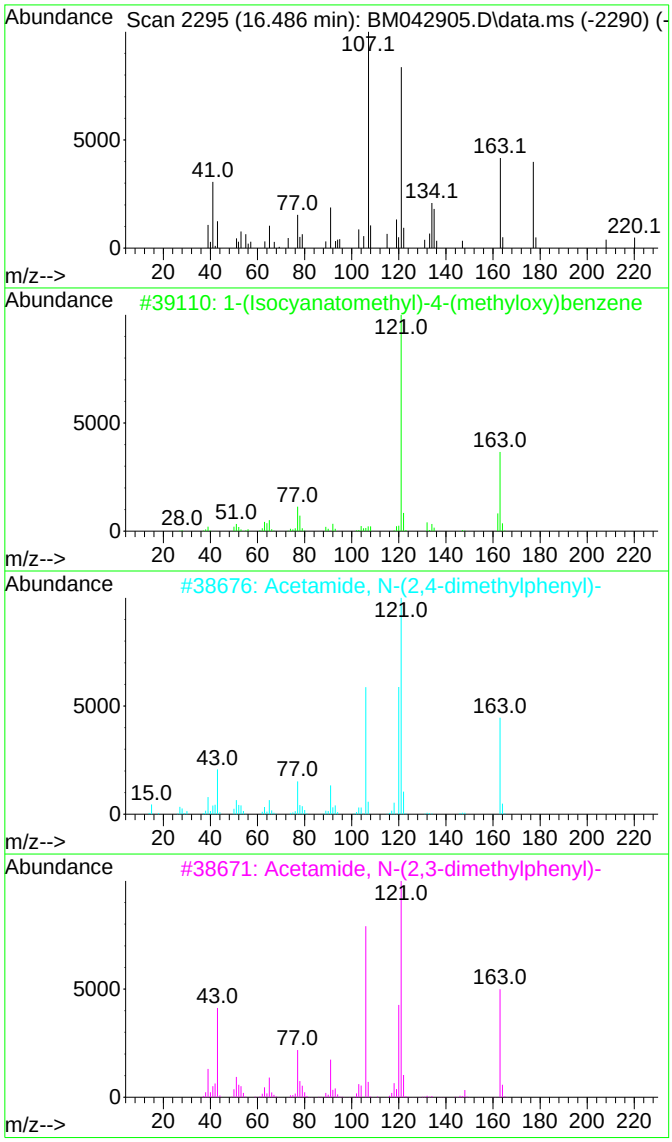
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

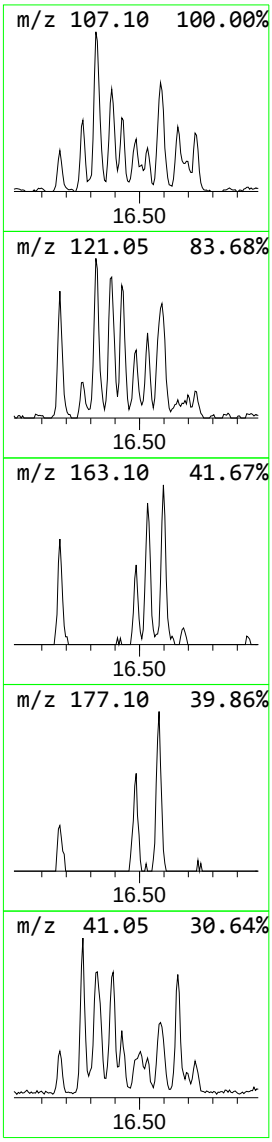
****

Peak Number 7 unknown-03 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.486	2.46 ng/ul	54338	Phenanthrene-d10	17.221

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-(Isocyanatomethyl)-4-(methylox...	163	C9H9NO2	056651-60-6	38
2	Acetamide, N-(2,4-dimethylphenyl)-	163	C10H13NO	002050-43-3	38
3	Acetamide, N-(2,3-dimethylphenyl)-	163	C10H13NO	000134-98-5	38
4	Acetamide, N-(2,6-dimethylphenyl)-	163	C10H13NO	002198-53-0	35
5	Propanamide, N-(3-methylphenyl)-...	177	C11H15NO	1000307-32-0	27





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Data File : BM042905.D
Acq On : 19 Nov 2023 14:24
Operator : MA/JU
Sample : 05370-16
Misc :
ALS Vial : 63 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GCDR0

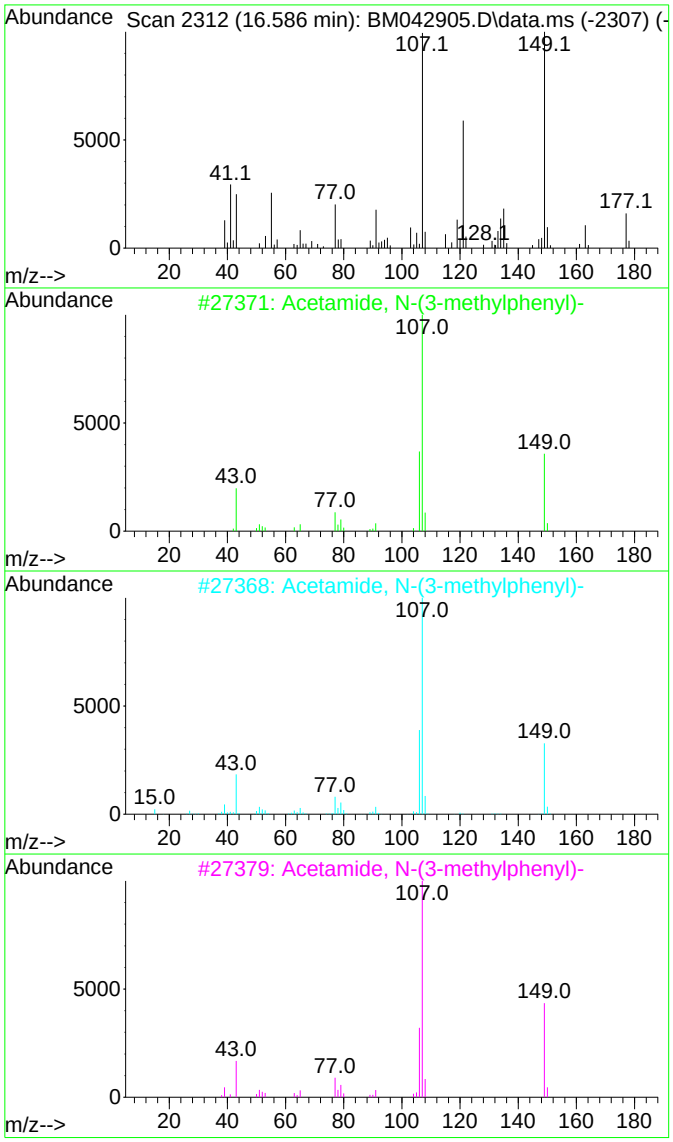
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM111323.MA.M
Quant Title : SVOA CALIBRATION

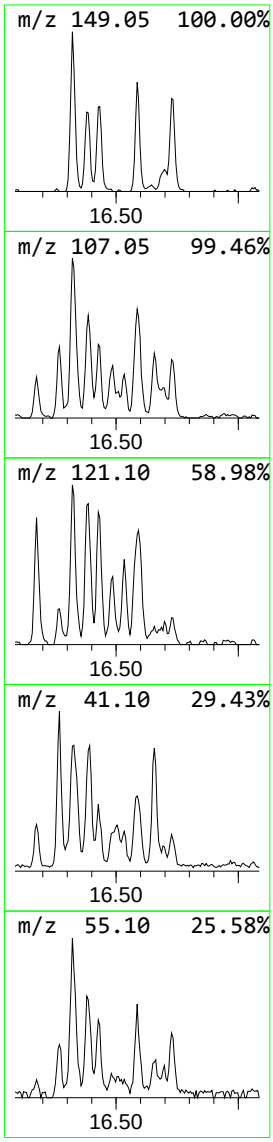
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 2-Methyl-4-hydroxybenzoxazole Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.586	6.81 ng/ul	150227	Phenanthrene-d10	17.221

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetamide, N-(3-methylphenyl)-	149	C9H11NO	000537-92-8	58
2	Acetamide, N-(3-methylphenyl)-	149	C9H11NO	000537-92-8	58
3	Acetamide, N-(3-methylphenyl)-	149	C9H11NO	000537-92-8	58
4	2-Methyl-4-hydroxybenzoxazole	149	C8H7NO2	051110-60-2	53
5	3,5-Diethyl-2-n-propylpyridine	177	C12H19N	004808-75-7	46





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111723\
Data File : BM042905.D
Acq On : 19 Nov 2023 14:24
Operator : MA/JU
Sample : 05370-16
Misc :
ALS Vial : 63 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GCDR0

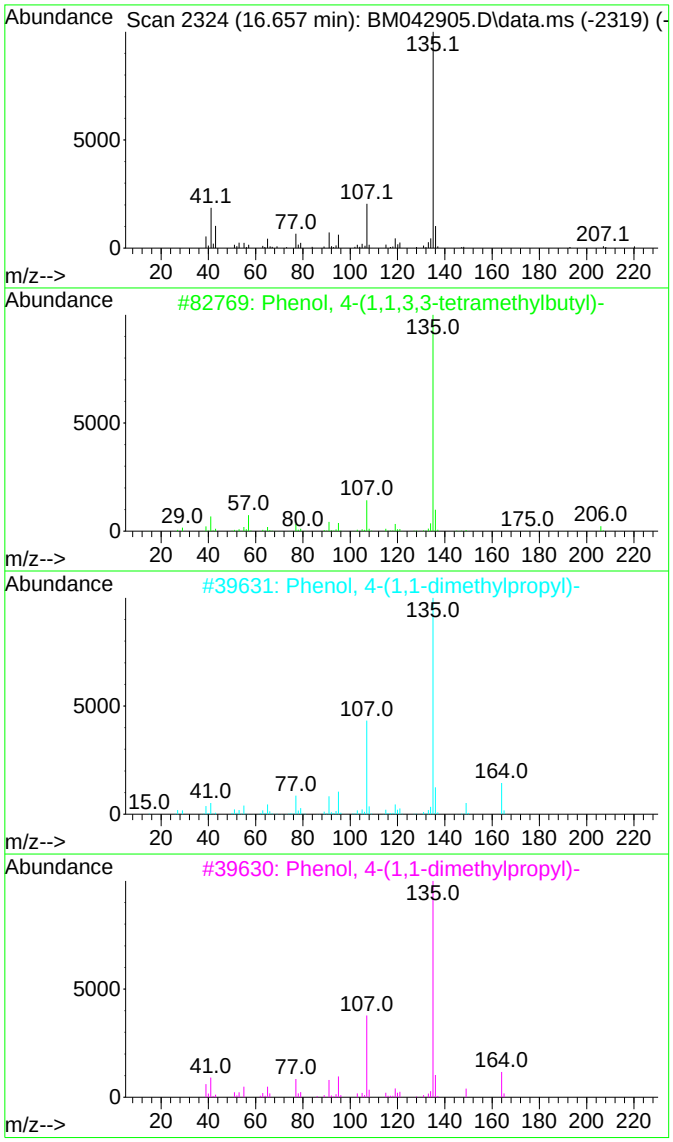
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM111323.MA.M
Quant Title : SVOA CALIBRATION

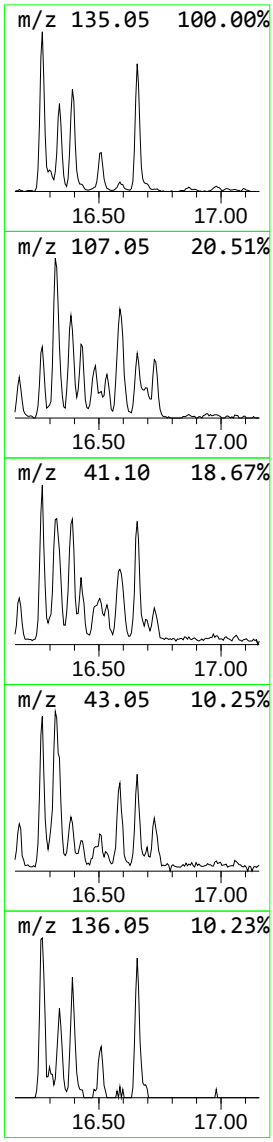
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Phenol, 4-(1,1-dimethylprop... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.657	6.18 ng/ul	136313	Phenanthrene-d10	17.221

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 4-(1,1,3,3-tetramethylbu...	206	C14H22O	000140-66-9	83
2	Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	78
3	Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	78
4	Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	78
5	Hexestrol, O-heptafluorobutyryl-	466	C22H21F7O3	1000365-31-9	72





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111723\
Data File : BM042905.D
Acq On : 19 Nov 2023 14:24
Operator : MA/JU
Sample : 05370-16
Misc :
ALS Vial : 63 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GCDR0

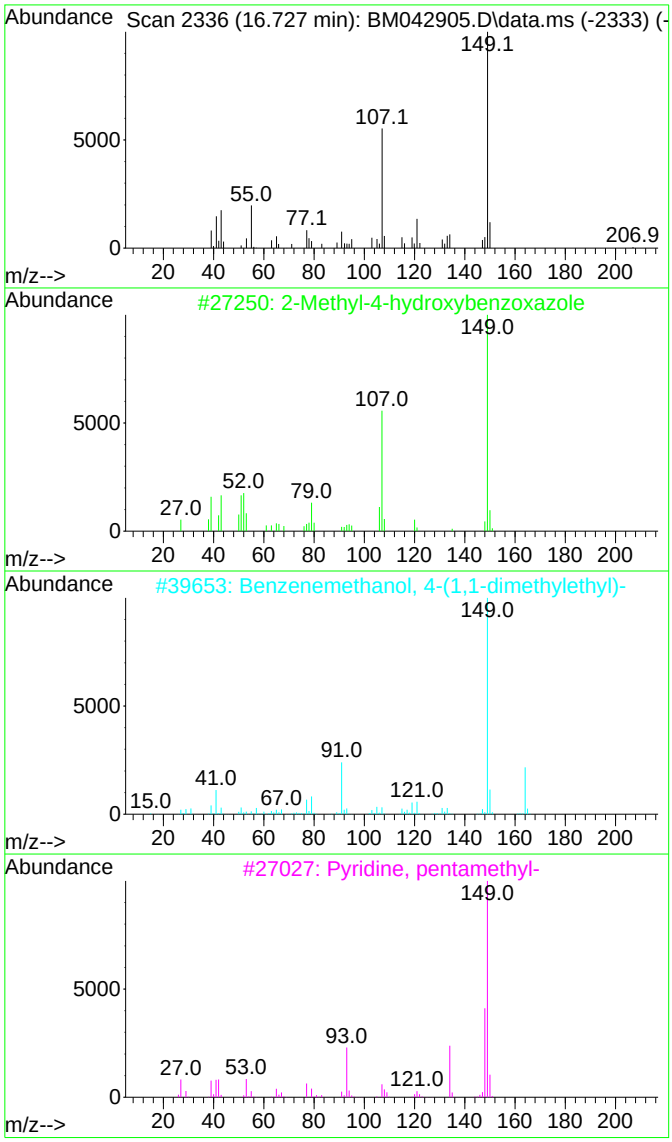
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM111323.MA.M
Quant Title : SVOA CALIBRATION

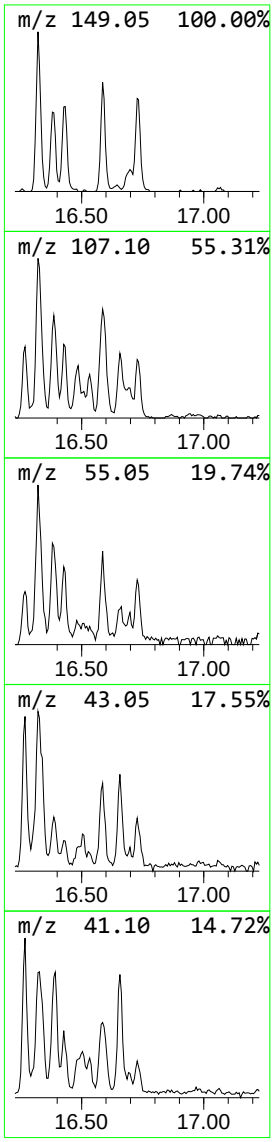
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Benzenemethanol, 4-(1,1-dim... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.727	2.79 ng/ul	61534	Phenanthrene-d10	17.221

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Methyl-4-hydroxybenzoxazole	149	C8H7NO2	051110-60-2	64
2	Benzenemethanol, 4-(1,1-dimethyl...	164	C11H16O	000877-65-6	53
3	Pyridine, pentamethyl-	149	C10H15N	003748-83-2	47
4	Benzofuran-2-one, 3-methyl-3-aza...	149	C8H7NO2	1010289-12-2	43
5	3,4-Methylenedioxypropiophenone	178	C10H10O3	028281-49-4	43





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111723\
Data File : BM042905.D
Acq On : 19 Nov 2023 14:24
Operator : MA/JU
Sample : 05370-16
Misc :
ALS Vial : 63 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GCDR0

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM111323.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Thiophane, propyl-	11.416	2.2	ng/ul	32747	2	10.622	299584	20.0
unknown-01	16.174	2.2	ng/ul	48158	4	17.221	441178	20.0
Phenol, 4-(1,1,...	16.268	7.1	ng/ul	156379	4	17.221	441178	20.0
Acetamide, N-(3...	16.321	11.6	ng/ul	256592	4	17.221	441178	20.0
4-(7-Methylocty...	16.386	8.4	ng/ul	185064	4	17.221	441178	20.0
unknown-02	16.427	4.3	ng/ul	93864	4	17.221	441178	20.0
unknown-03	16.486	2.5	ng/ul	54338	4	17.221	441178	20.0
2-Methyl-4-hydr...	16.586	6.8	ng/ul	150227	4	17.221	441178	20.0
Phenol, 4-(1,1-...	16.657	6.2	ng/ul	136313	4	17.221	441178	20.0
Benzenemethanol...	16.727	2.8	ng/ul	61534	4	17.221	441178	20.0