LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022721\
Data File : BM028950.D
Acq On : 27 Feb 2021 10:30
Operator : CG/JU
Sample : M1498-09
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB8G8

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\SOM-EPA-BM022721MA.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	3.122	19	23	27	rBV	2879	3450	0.92%	0.080%
2	3.158	27	29	35	rVB2	1551	1819	0.49%	0.042%
3	3.881	148	152	156	rVB2	1030	1212	0.32%	0.028%
4	4.352	229	232	235	rBB	1761	2390	0.64%	0.055%
5	5.040	345	349	356	rBB	22990	33447	8.93%	0.774%
6	6.957	670	675	684	rBB	18596	29760	7.95%	0.689%
7	7.110	695	701	708	rBB	53493	80901	21.61%	1.872%
8	7.304	728	734	741	rBV	66521	105547	28.20%	2.442%
9	7.769	807	813	824	rBB	66450	101263	27.05%	2.343%
10	8.134	872	875	880	rBB	1806	2482	0.66%	0.057%
11	8.504	932	938	947	rBV	52267	81900	21.88%	1.895%
12	8.946	1007	1013	1022	rBB	75506	117815	31.47%	2.726%
13	9.310	1073	1075	1078	rVB2	1278	1484	0.40%	0.034%
14	9.398	1086	1090	1093	rBB2	2764	3651	0.98%	0.084%
15	9.663	1130	1135	1142	rBB	65951	104317	27.87%	2.414%
16	10.204	1221	1227	1236	rBV	95860	155639	41.58%	3.601%
17	10.569	1283	1289	1298	rBV	87552	138279	36.94%	3.199%
18	10.728	1310	1316	1323	rBV2	11511	17827	4.76%	0.412%
19	11.928	1516	1520	1525	rVB2	3997	5696	1.52%	0.132%
20	12.139	1552	1556	1558	rVV	922	1292	0.35%	0.030%
21	12.169	1558	1561	1564	rVB2	1321	1453	0.39%	0.034%
22	12.216	1564	1569	1571	rBV3	834	1290	0.34%	0.030%
23	12.375	1590	1596	1598	rBV2	827	1390	0.37%	0.032%
24	12.463	1605	1611	1616	rBV2	2117	3185	0.85%	0.074%
25	12.528	1618	1622	1627	rVB3	776	1263	0.34%	0.029%
26	13.233	1736	1742	1746	rBV2	1323	1894	0.51%	0.044%
27	13.275	1746	1749	1753	rBV2	1130	1774	0.47%	0.041%
28	13.322	1753	1757	1760	rVB4	1521	2022	0.54%	0.047%
29	13.386	1765	1768	1773	rBV4	1287	2017	0.54%	0.047%
30	13.481	1782	1784	1789	rVB3	1107	1425	0.38%	0.033%
31	13.669	1811	1816	1819	rBV2	2032	3462	0.92%	0.080%
32	13.757	1827	1831	1835	rVV	1551	3202	0.86%	0.074%
33	13.851	1841	1847	1852	rVV	165396	219271	58.57%	5.073%
34	13.898	1852	1855	1861	rVB	5211	6268	1.67%	0.145%

LSC Area Percent Report

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Operator : CG/JU
Sample : M1498-09
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB8G8

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\SOM-EPA-BM022721MA.M
Title : SVOA CALIBRATION

35	13.998	1865	1872	1873	rBV3	1593	2377	0.63%	0.055%
36	14.022	1873	1876	1879	rVV3	1461	1874	0.50%	0.043%
37	14.075	1883	1885	1888	rVV2	3743	4442	1.19%	0.103%
38	14.128	1888	1894	1902	rVV	196968	280407	74.91%	6.488%
39	14.257	1911	1916	1923	rVV	4481	8106	2.17%	0.188%
40	14.363	1930	1934	1937	rVB2	1681	2019	0.54%	0.047%
41	14.433	1940	1946	1952	rBV	122271	177112	47.31%	4.098%
42	14.498	1952	1957	1962	rVB	19872	28496	7.61%	0.659%
43	14.580	1968	1971	1976	rVB3	1804	2662	0.71%	0.062%
44	14.686	1983	1989	1996	rBV	12463	24282	6.49%	0.562%
45	14.833	2010	2014	2018	rBV2	6075	7135	1.91%	0.165%
46	14.992	2038	2041	2046	rVB3	2351	3552	0.95%	0.082%
47	15.045	2046	2050	2051	rBV2	1448	1889	0.50%	0.044%
48	15.098	2057	2059	2062	rVB3	1185	1177	0.31%	0.027%
49	15.186	2070	2074	2079	rBV2	5366	8809	2.35%	0.204%
50	15.257	2085	2086	2092	rVB4	1813	2605	0.70%	0.060%
51	15.345	2097	2101	2102	rVV3	1576	1998	0.53%	0.046%
52	15.427	2109	2115	2121	rVV	236428	323941	86.54%	7.495%
53	15.480	2121	2124	2129	rVB3	9181	12733	3.40%	0.295%
54	15.575	2134	2140	2148	rVV	82217	115198	30.77%	2.665%
55	15.633	2148	2150	2152	rVV2	1125	1211	0.32%	0.028%
56	15.663	2152	2155	2158	rVV2	4186	4948	1.32%	0.114%
57	15.698	2158	2161	2168	rVV5	1458	2469	0.66%	0.057%
58	15.780	2172	2175	2179	rBV5	1377	2102	0.56%	0.049%
59	15.951	2198	2204	2213	rVB	40536	55839	14.92%	1.292%
60	16.092	2227	2228	2233	rVV5	1298	2362	0.63%	0.055%
61	16.145	2233	2237	2245	rVV2	11629	21225	5.67%	0.491%
62	16.204	2245	2247	2249	rVB3	1559	1333	0.36%	0.031%
63	16.286	2259	2261	2264	rBV3	1980	1907	0.51%	0.044%
64	16.369	2273	2275	2276	rBV2	1528	1175	0.31%	0.027%
65	16.386	2276	2278	2283	rVB3	2338	3426	0.92%	0.079%
66	16.463	2286	2291	2297	rBV	22022	30389	8.12%	0.703%
67	16.539	2303	2304	2310	rVB2	833	1154	0.31%	0.027%
68	16.722	2333	2335	2338	rVB3	1561	1558	0.42%	0.036%
69	16.957	2372	2375	2379	rBV3	1266	2003	0.54%	0.046%
70	16.992	2379	2381	2385	rVB3	2158	1922	0.51%	0.044%
71	17.033	2385	2388	2389	rBV3	1348	1477	0.39%	0.034%

LSC Area Percent Report

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Sample : M1498-09
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB8G8

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

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Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM022721MA.M
Title : SVOA CALIBRATION

72	17.092	2395	2398	2403	rVB5	1349	1894	0.51%	0.044%
73	17.174	2407	2412	2419	rVV	139612	185380	49.52%	4.289%
74	17.221	2419	2420	2424	rVV3	1644	1487	0.40%	0.034%
75	17.274	2424	2429	2436	rVB	252916	331297	88.50%	7.665%
76	17.404	2449	2451	2456	rVB6	1790	2005	0.54%	0.046%
77	17.457	2456	2460	2463	rBV6	2026	2956	0.79%	0.068%
78	17.598	2482	2484	2486	rBV3	1286	1205	0.32%	0.028%
79	17.874	2530	2531	2536	rVB4	1035	1457	0.39%	0.034%
80	17.939	2539	2542	2545	rBV4	2300	2804	0.75%	0.065%
81	18.068	2559	2564	2571	rVV	40681	53174	14.20%	1.230%
82	18.239	2590	2593	2598	rVB7	1953	2962	0.79%	0.069%
83	18.410	2620	2622	2626	rVB4	1435	1751	0.47%	0.041%
84	18.445	2626	2628	2629	rBV	2089	1231	0.33%	0.028%
85	18.551	2641	2646	2650	rBV2	3682	6906	1.84%	0.160%
86	18.804	2685	2689	2694	rVV	20499	26286	7.02%	0.608%
87	18.898	2701	2705	2709	rBV2	6335	9501	2.54%	0.220%
88	18.963	2713	2716	2724	rBV8	3065	6660	1.78%	0.154%
89	19.198	2753	2756	2760	rBV2	4714	5922	1.58%	0.137%
90	19.280	2767	2770	2772	rBV3	2155	2063	0.55%	0.048%
91	19.345	2777	2781	2785	rBV2	17629	22813	6.09%	0.528%
92	19.445	2795	2798	2801	rVB2	6427	7382	1.97%	0.171%
93	19.563	2812	2818	2823	rBV	282675	363783	97.18%	8.417%
94	19.615	2823	2827	2834	rVB	41154	50021	13.36%	1.157%
95	20.062	2900	2903	2907	rVB4	4007	4980	1.33%	0.115%
96	20.580	2989	2991	2994	rVB4	5725	4693	1.25%	0.109%
97	21.045	3066	3070	3075	rBV	48343	56064	14.98%	1.297%
98	21.339	3115	3120	3126	rBV	172379	206851	55.26%	4.786%
99	23.398	3464	3470	3481	rVB	208030	374344	100.00%	8.661%
100	23.539	3488	3494	3501	rVB2	112046	196759	52.56%	4.552%

Sum of corrected areas: 4322032

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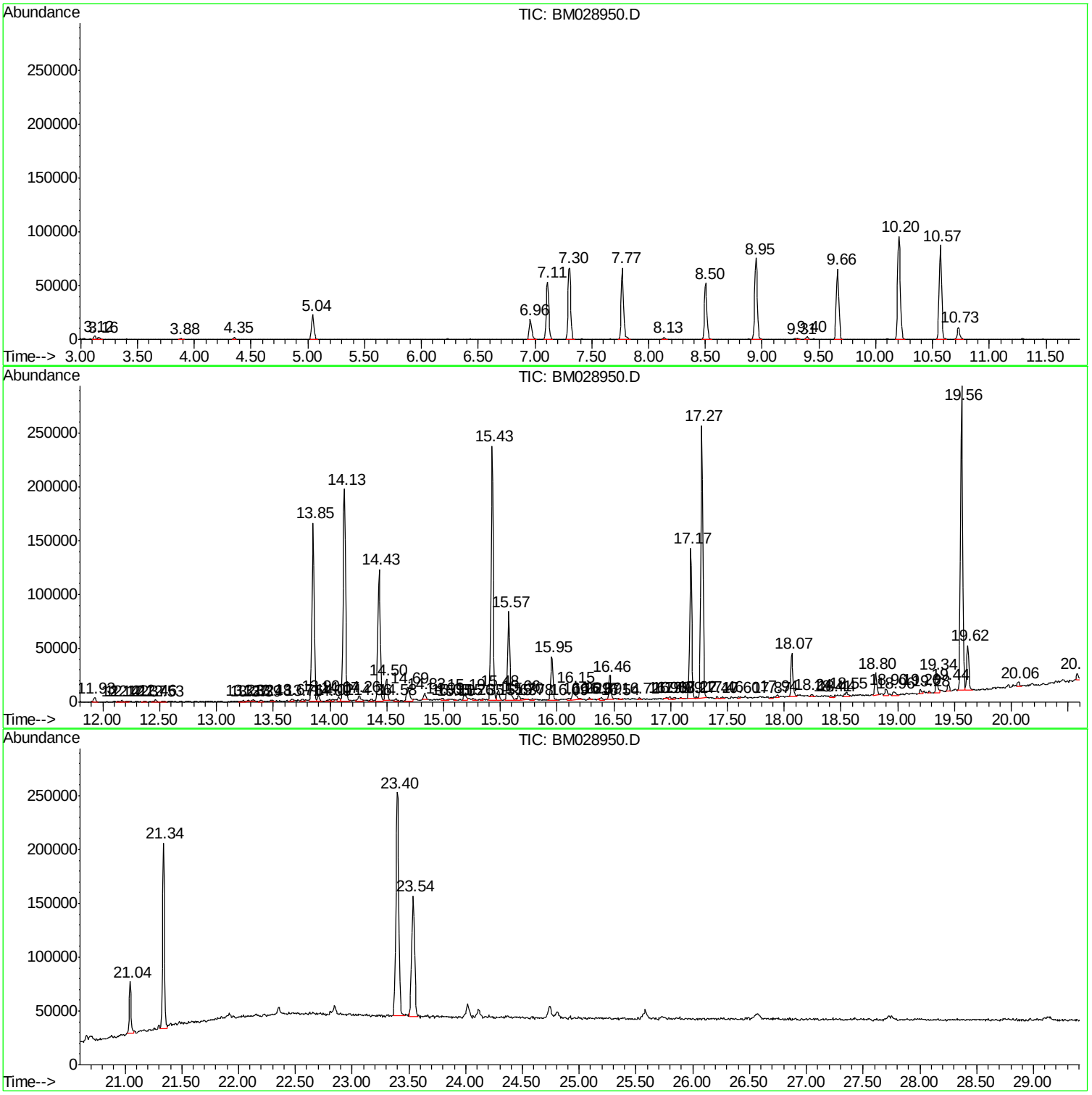
Instrument : BNA_M
ClientSampleId : CB8G8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM022721MA.M
Quant Title : SVOA CALIBRATION

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



Library Search Compound Report

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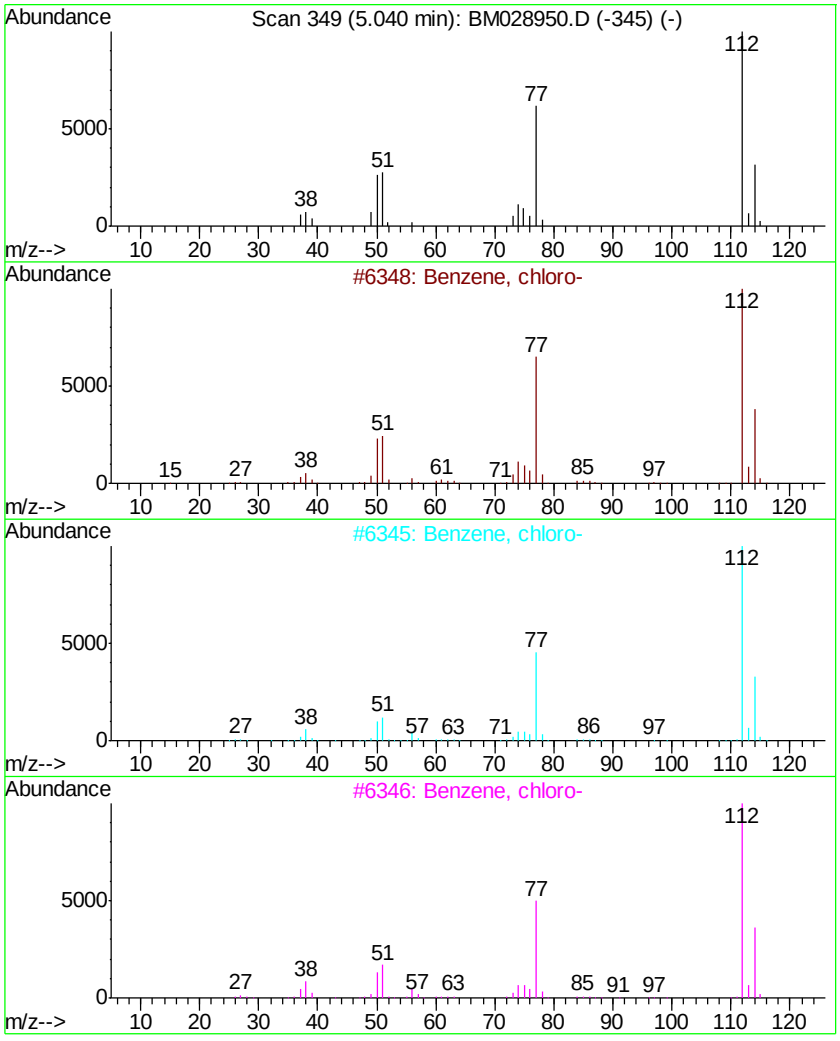
TIC Integration Parameters: LSCINT.P

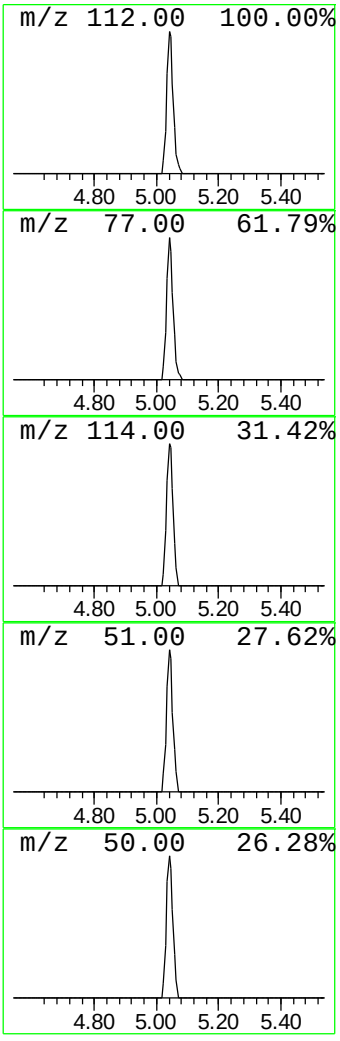
****

Peak Number 1 Benzene, chloro- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.04	6.61 ng/ul	33447	1,4-Dichlorobenzene-d4	7.77

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, chloro-	112	C6H5Cl	000108-90-7	91
2	Benzene, chloro-	112	C6H5Cl	000108-90-7	91
3	Benzene, chloro-	112	C6H5Cl	000108-90-7	91
4	Benzene, chloro-	112	C6H5Cl	000108-90-7	90
5	2-Furancarboxylic acid	112	C5H4O3	000088-14-2	16





Library Search Compound Report

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Data File : BM028950.D
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Sample : M1498-09
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB8G8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM022721MA.M
Quant Title : SVOA CALIBRATION

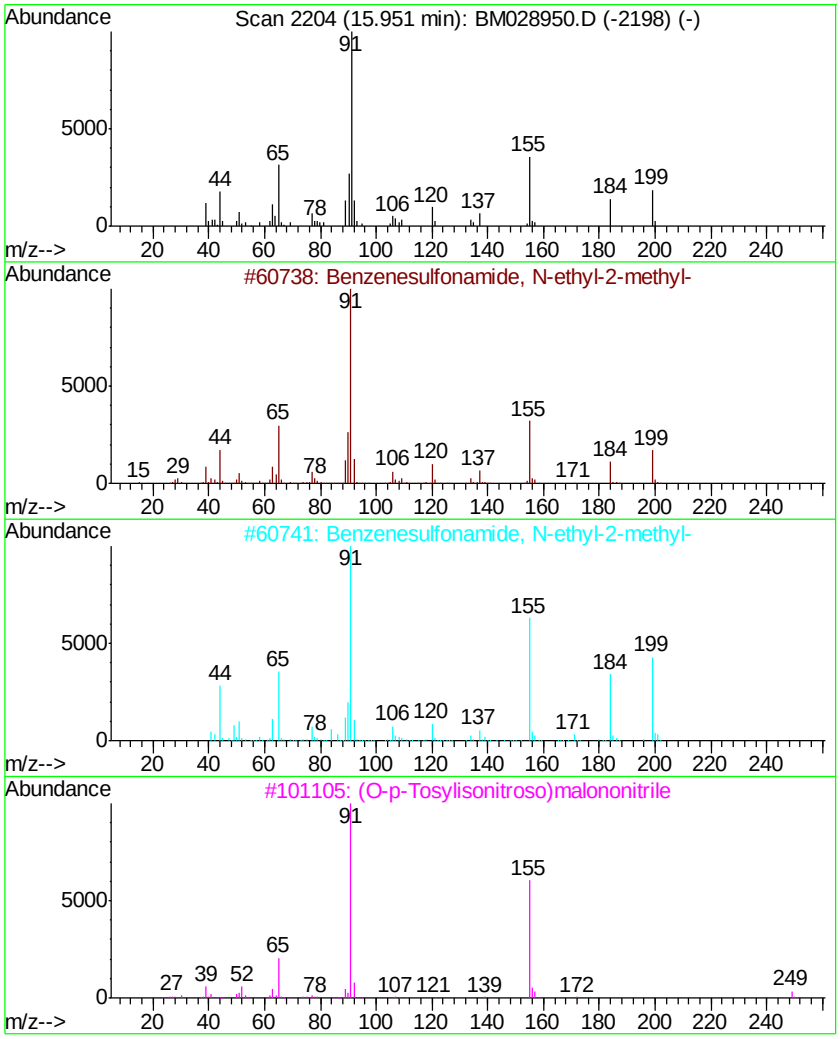
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

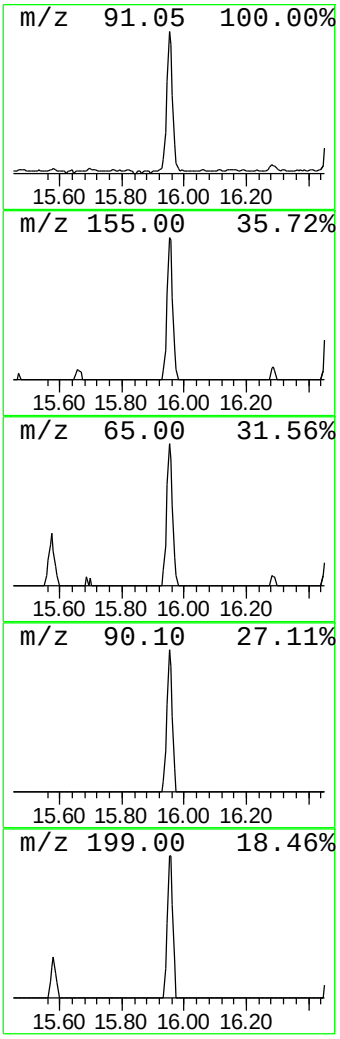
****

Peak Number 2 Benzenesulfonamide, N-ethyl... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.95	6.02 ng/ul	55839	Phenanthrene-d10	17.17

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenesulfonamide, N-ethyl-2-me...	199	C9H13NO2S	001077-56-1	98
2	Benzenesulfonamide, N-ethyl-2-me...	199	C9H13NO2S	001077-56-1	91
3	(O-p-Tosylisonitroso)malononitrile	249	C10H7N3O3S	020893-01-0	47
4	p-Toluenesulfonylacetonitrile	195	C9H9NO2S	005697-44-9	47
5	Ethyl(E/Z)-.alpha.-cyano-4-nitro...	416	C19H16N2O7S	1000287-26-8	47





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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM022721MA.M
Quant Title : SVOA CALIBRATION

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

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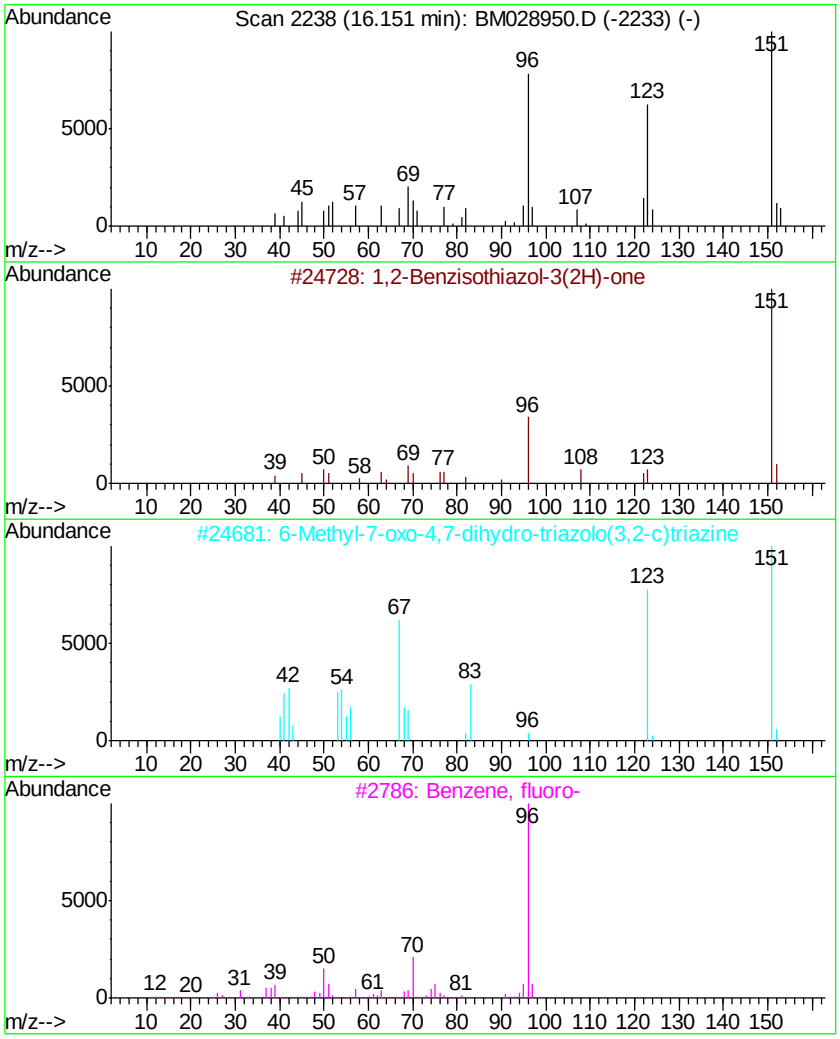
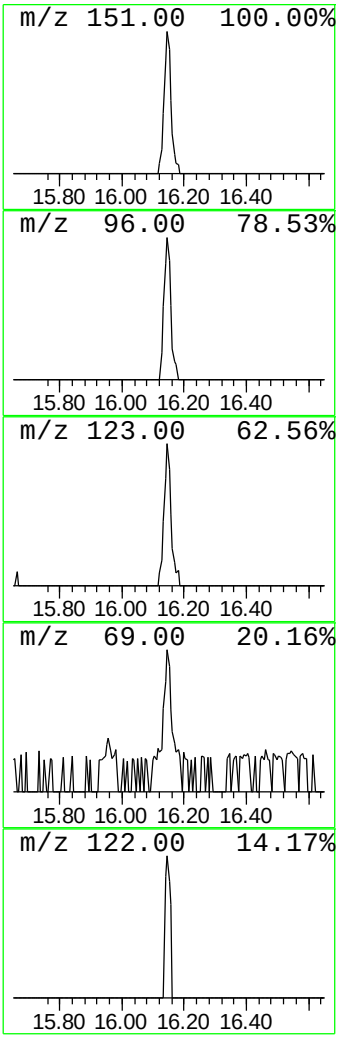
Peak Number 3 1,2-Benzisothiazol-3(2H)-one Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.15	2.29 ng/ul	21225	Phenanthrene-d10	17.17

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,2-Benzisothiazol-3(2H)-one	151	C7H5NOS	002634-33-5	64
2	6-Methyl-7-oxo-4,7-dihydro-triaz...	151	C5H5N5O	057250-39-2	32
3	Benzene, fluoro-	96	C6H5F	000462-06-6	30
4	RCH 108	151	C7H5NO3	002297-94-1	27
5	4-Methoxyformanilide	151	C8H9NO2	005470-34-8	16





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ClientSampleId : CB8G8

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Quant Title : SVOA CALIBRATION

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

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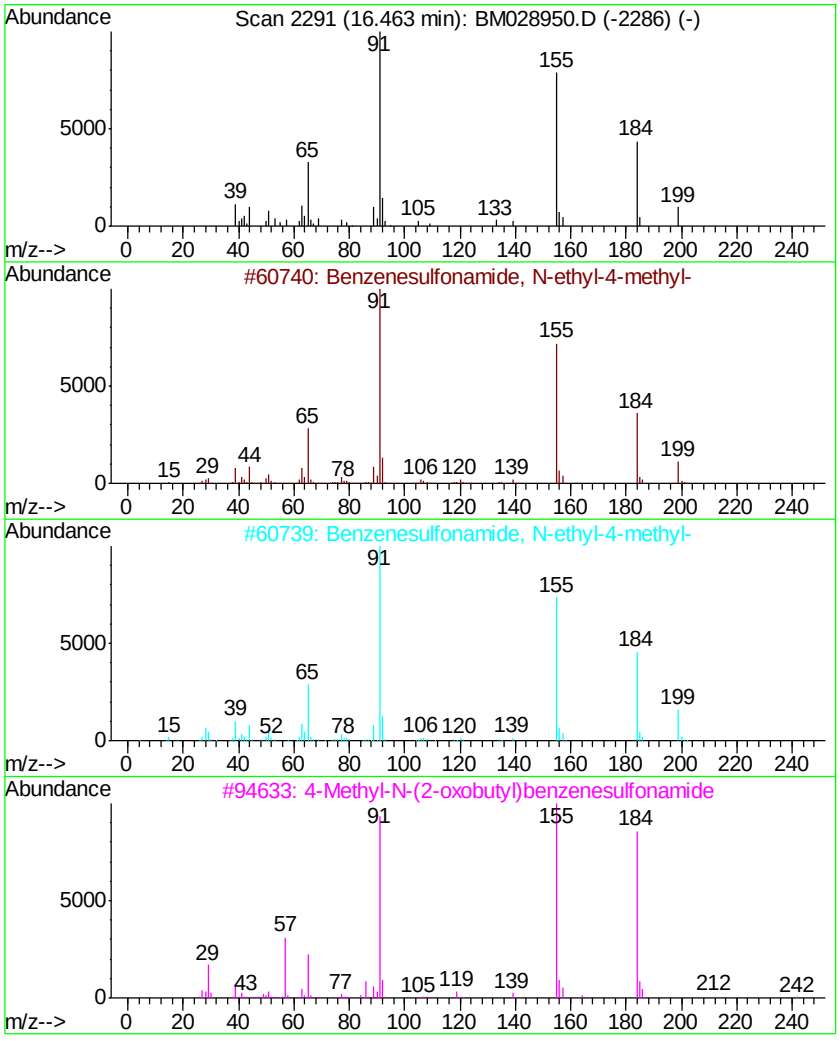
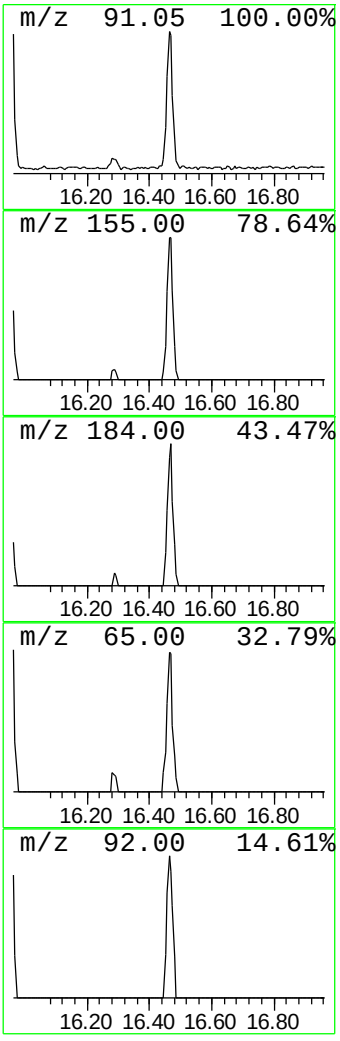
Peak Number 4 Benzenesulfonamide, N-ethyl... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.46	3.28 ng/ul	30389	Phenanthrene-d10	17.17

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	97
2	Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	90
3	4-Methyl-N-(2-oxobutyl)benzenesu...	241	C11H15NO3S	1000192-14-5	90
4	(Toluene-4-sulfonylamino)-acetic...	283	C13H17NO4S	1000190-48-0	59
5	Pyridine-2-carbonitrile, 3,5,6-t...	375	C13H8Cl3N3O2S	255713-77-0	53





Library Search Compound Report

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Data File : BM028950.D
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Operator : CG/JU
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Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB8G8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM022721MA.M
Quant Title : SVOA CALIBRATION

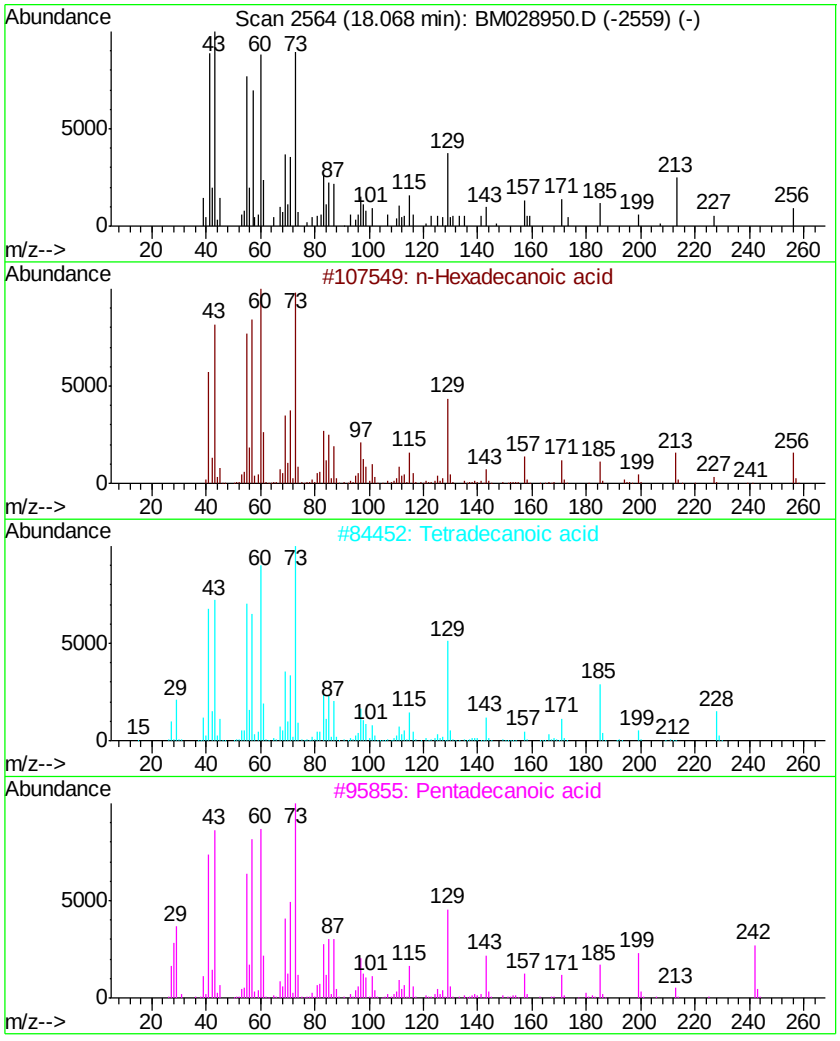
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

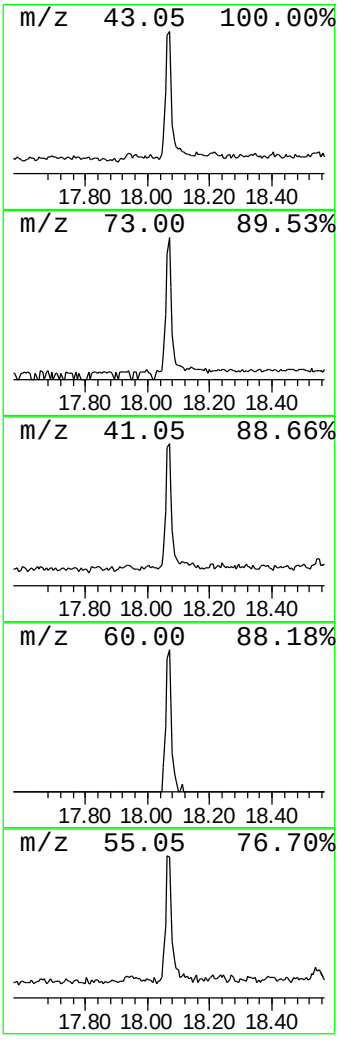
****

Peak Number 5 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.07	5.74 ng/ul	53174	Phenanthrene-d10	17.17

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
2	Tetradecanoic acid	228	C14H28O2	000544-63-8	83
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	81
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	80
5	Pentadecanoic acid	242	C15H30O2	001002-84-2	74





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Data File : BM028950.D
Acq On : 27 Feb 2021 10:30
Operator : CG/JU
Sample : M1498-09
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB8G8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM022721MA.M
Quant Title : SVOA CALIBRATION

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

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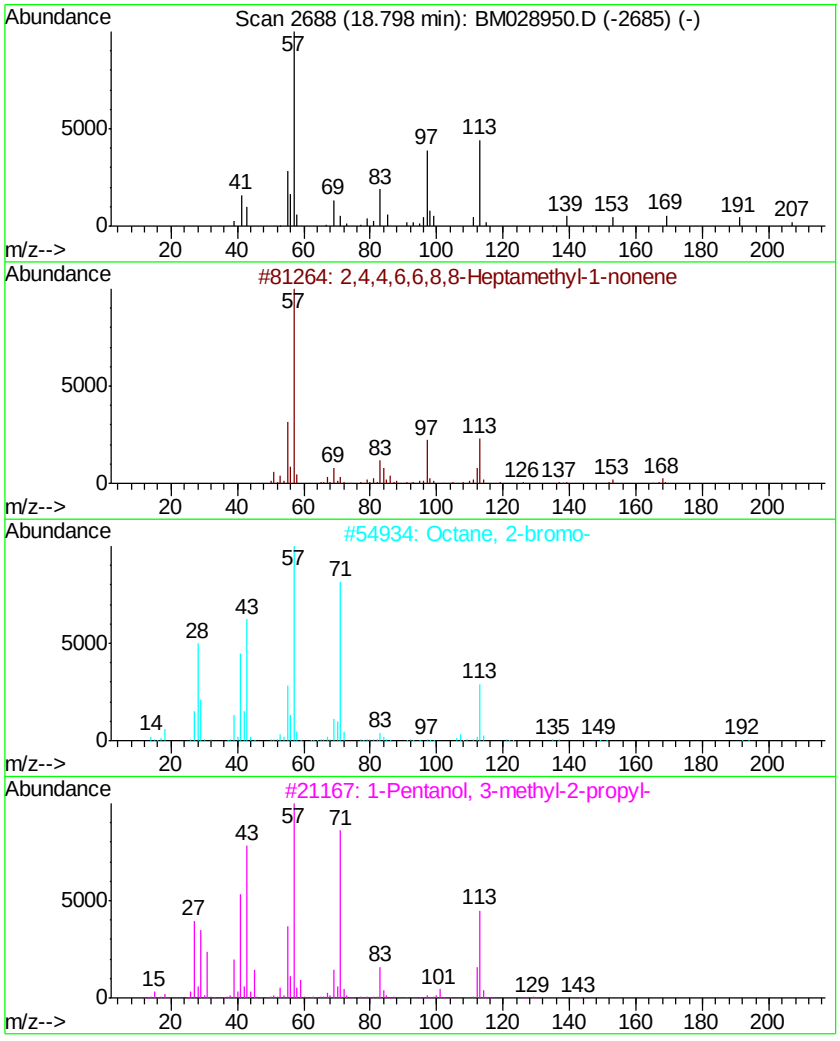
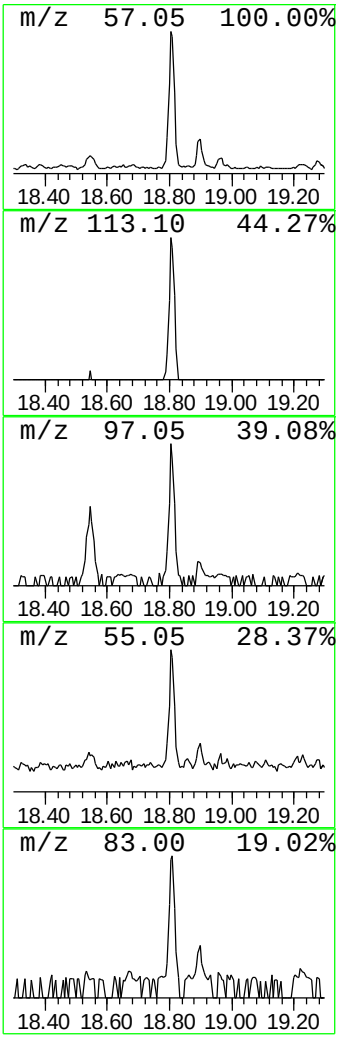
Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.80	2.84 ng/ul	26286	Phenanthrene-d10	17.17

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,4,6,6,8,8-Heptamethyl-1-nonene	224	C16H32	015796-04-0	64
2	Octane, 2-bromo-	192	C8H17Br	000557-35-7	38
3	1-Pentanol, 3-methyl-2-propyl-	144	C9H20O	054004-40-9	37
4	2,4,4-Trimethyl-1-pentanol, pent...	276	C11H17F5O2	1000365-51-0	35
5	Undecane, 6,6-dimethyl-	184	C13H28	017312-76-4	32





Library Search Compound Report

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Data File : BM028950.D
Acq On : 27 Feb 2021 10:30
Operator : CG/JU
Sample : M1498-09
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB8G8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM022721MA.M
Quant Title : SVOA CALIBRATION

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

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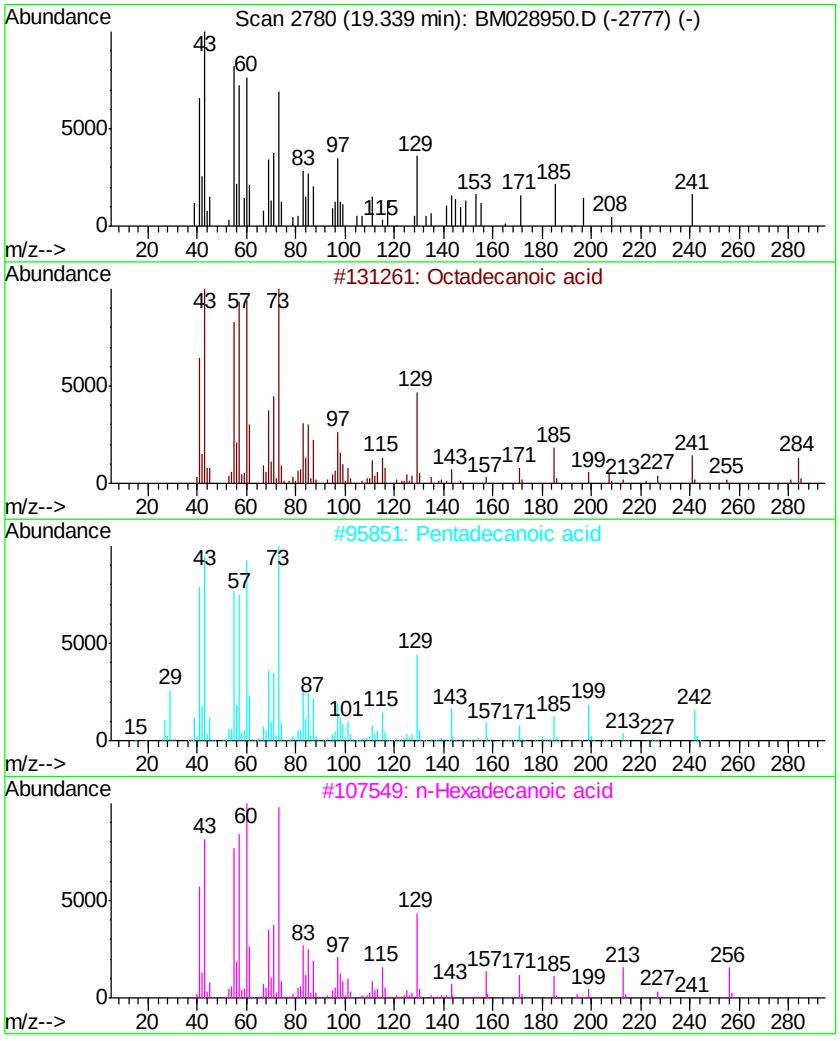
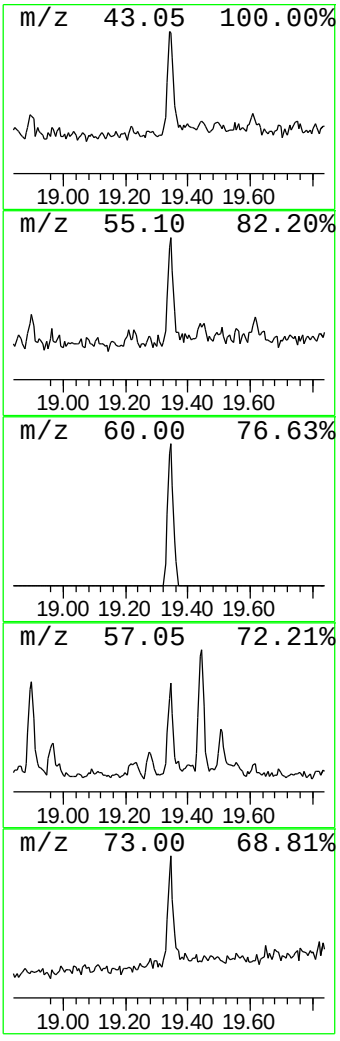
Peak Number 7 Octadecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.34	2.21 ng/ul	22813	Chrysene-d12	21.34

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	87
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	86
4	Pentadecanoic acid	242	C15H30O2	001002-84-2	80
5	Tetradecanoic acid	228	C14H28O2	000544-63-8	64





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Data File : BM028950.D
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Operator : CG/JU
Sample : M1498-09
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB8G8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM022721MA.M
Quant Title : SVOA CALIBRATION

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

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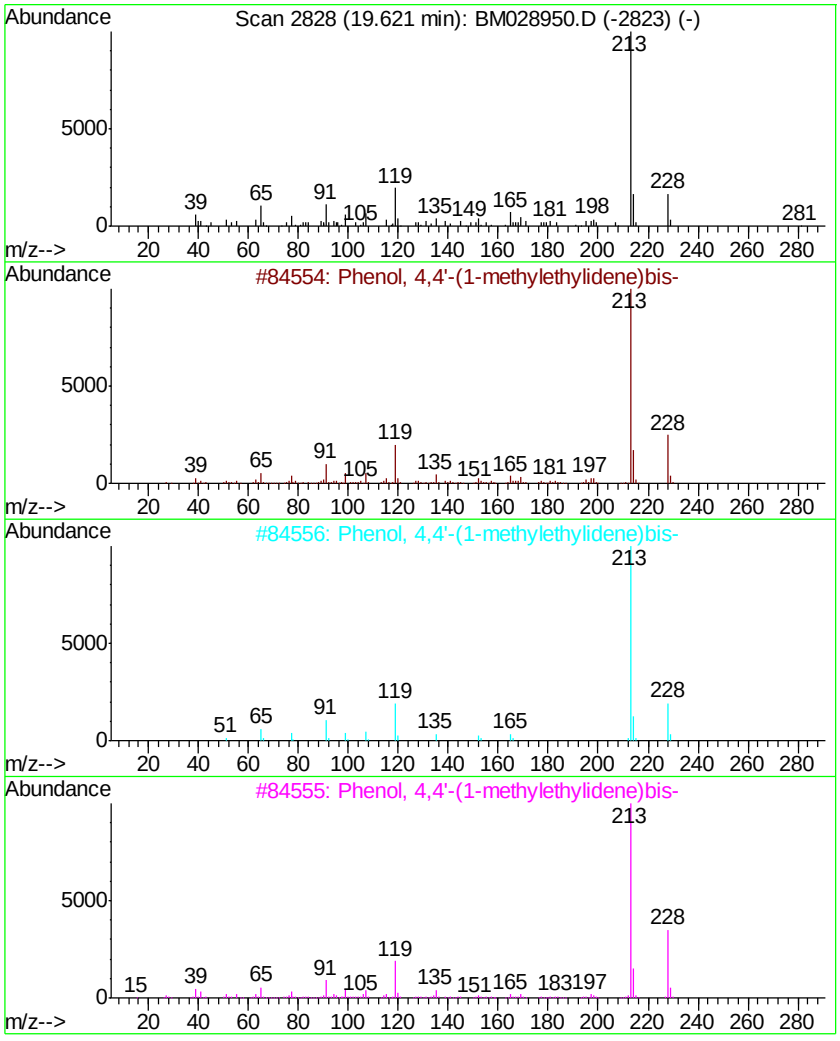
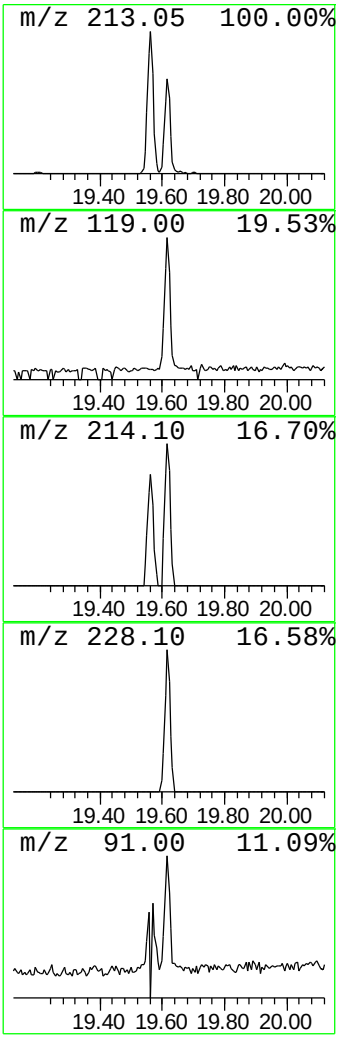
Peak Number 8 Phenol, 4,4'-(1-methylethyl... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.62	4.84 ng/ul	50021	Chrysene-d12	21.34

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	97
2	Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	94
3	Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	94
4	Phenol, 2,4'-isopropylidenedi-	228	C15H16O2	000837-08-1	72
5	2-Trimethylsilyl-3-trimethylsily...	228	C8H20N4Si2	1000373-05-5	58





Library Search Compound Report

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Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB8G8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM022721MA.M
Quant Title : SVOA CALIBRATION

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

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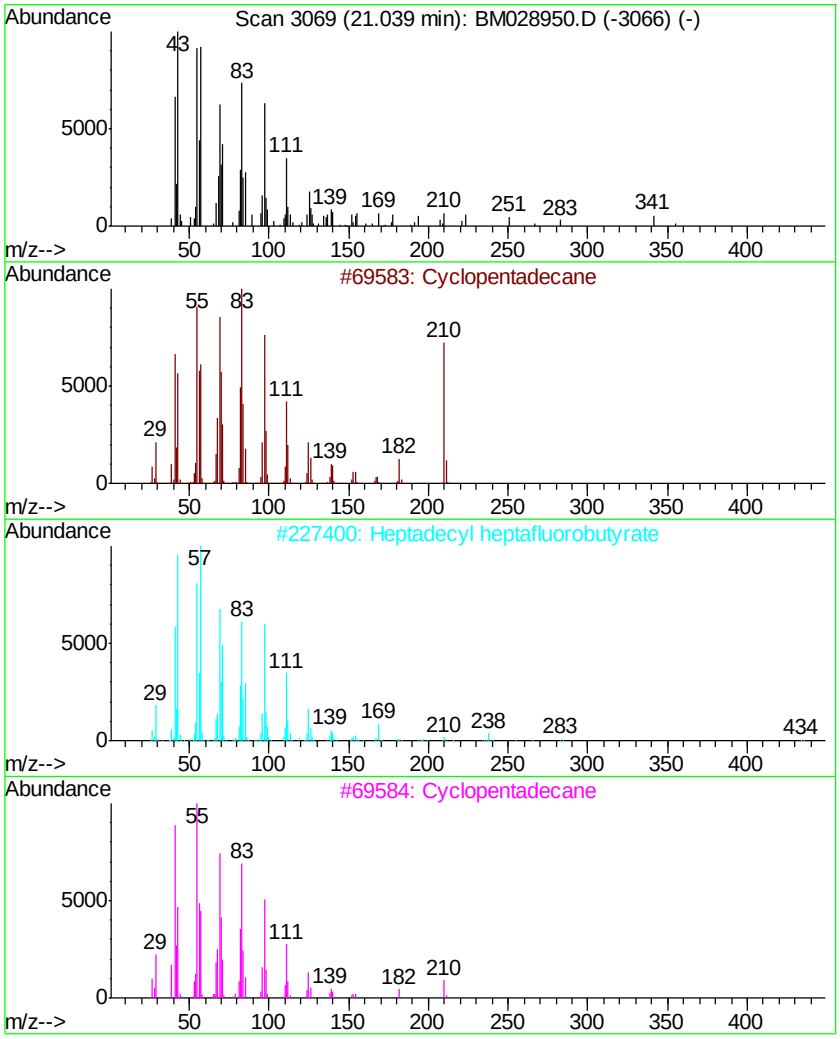
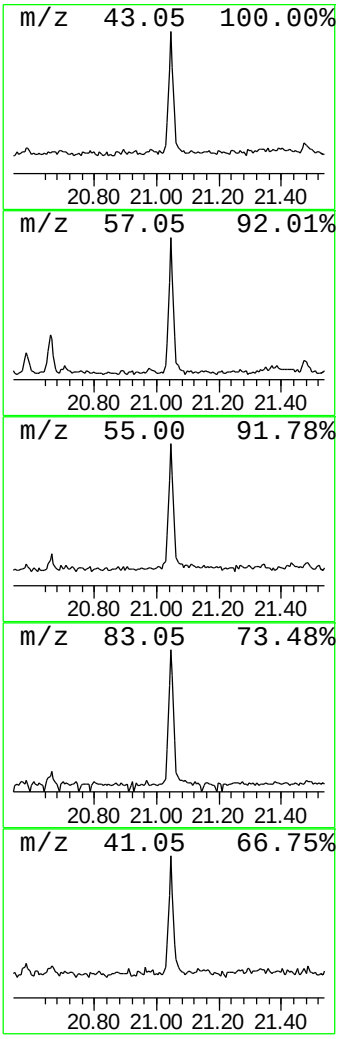
Peak Number 9 (DEL) Alkane: Cyclic21.04 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.04	5.42 ng/ul	56064	Chrysene-d12	21.34

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentadecane	210	C15H30	000295-48-7	96
2	Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	95
3	Cyclopentadecane	210	C15H30	000295-48-7	94
4	Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	93
5	Heptadecyl trifluoroacetate	352	C19H35F3O2	1000351-87-0	93





Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM022721\
Data File : BM028950.D
Acq On : 27 Feb 2021 10:30
Operator : CG/JU
Sample : M1498-09
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : CB8G8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM022721MA.M
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, chloro-	5.04	6.6	ng/ul	33447	1	7.77	101263	20.0
Benzenesulfonamid...	15.95	6.0	ng/ul	55839	4	17.17	185380	20.0
1,2-Benzisothiazo...	16.15	2.3	ng/ul	21225	4	17.17	185380	20.0
Benzenesulfonamid...	16.46	3.3	ng/ul	30389	4	17.17	185380	20.0
n-Hexadecanoic acid	18.07	5.7	ng/ul	53174	4	17.17	185380	20.0
(DEL) Alkane: Str...	18.80	2.8	ng/ul	26286	4	17.17	185380	20.0
Octadecanoic acid	19.34	2.2	ng/ul	22813	5	21.34	206851	20.0
Phenol, 4,4'-(1-m...	19.62	4.8	ng/ul	50021	5	21.34	206851	20.0
(DEL) Alkane: Cyc...	21.04	5.4	ng/ul	56064	5	21.34	206851	20.0