LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121820\
Data File : BP004390.D
Acq On : 19 Dec 2020 12:48
Operator : CG/JU
Sample : L5128-03
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : CB8C2

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_P\METHODS\SOM-EPA-BP121820MA.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	4.481	216	220	230	rVB2	71977	118978	2.24%	0.192%
2	5.158	332	335	340	rVB	37826	54319	1.02%	0.088%
3	6.993	639	647	660	rVB2	591248	1117939	21.04%	1.802%
4	7.152	668	674	685	rBV	604794	974052	18.33%	1.570%
5	7.358	702	709	716	rBV	730530	1210690	22.78%	1.952%
6	7.822	782	788	795	rBV	766353	1226429	23.08%	1.977%
7	7.893	795	800	807	rVB6	19618	36845	0.69%	0.059%
8	8.193	847	851	858	rVB	45355	81259	1.53%	0.131%
9	8.358	877	879	888	rVB6	8844	18513	0.35%	0.030%
10	8.469	893	898	901	rVV3	15486	23874	0.45%	0.038%
11	8.528	901	908	918	rVV	933701	1512614	28.47%	2.439%
12	8.716	935	940	948	rVB3	13512	28265	0.53%	0.046%
13	8.799	948	954	965	rBV	198435	340921	6.42%	0.550%
14	8.922	970	975	978	rBV4	30957	52203	0.98%	0.084%
15	8.975	978	984	995	rVB	741239	1236191	23.26%	1.993%
16	9.163	1010	1016	1018	rBV6	7117	13175	0.25%	0.021%
17	9.334	1039	1045	1051	rBV2	34416	61086	1.15%	0.098%
18	9.416	1053	1059	1071	rVB4	61418	136681	2.57%	0.220%
19	9.563	1080	1084	1089	rVB5	13513	21257	0.40%	0.034%
20	9.699	1101	1107	1122	rVB	621740	1055654	19.87%	1.702%
21	9.828	1122	1129	1133	rBV9	14521	31288	0.59%	0.050%
22	9.910	1141	1143	1152	rVB9	10541	21920	0.41%	0.035%
23	10.028	1152	1163	1170	rBV9	17548	49280	0.93%	0.079%
24	10.105	1170	1176	1179	rBV3	11511	21069	0.40%	0.034%
25	10.152	1179	1184	1192	rVB6	22669	53443	1.01%	0.086%
26	10.240	1192	1199	1208	rBV	1082375	1835119	34.54%	2.959%
27	10.387	1217	1224	1226	rBV6	8221	14559	0.27%	0.023%
28	10.422	1226	1230	1237	rVB8	9552	19176	0.36%	0.031%
29	10.528	1242	1248	1255	rVV3	34219	69419	1.31%	0.112%
30	10.616	1255	1263	1270	rVV	1099669	1876555	35.32%	3.026%
31	10.669	1270	1272	1277	rVV3	13748	20963	0.39%	0.034%
32	10.752	1277	1286	1300	rVV	1038276	1781579	33.53%	2.872%
33	10.869	1302	1306	1314	rVV5	20668	45425	0.85%	0.073%
34	11.034	1328	1334	1341	rVB10	12108	25040	0.47%	0.040%

LSC Area Percent Report

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Operator : CG/JU
Sample : L5128-03
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : CB8C2

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_P\METHODS\SOM-EPA-BP121820MA.M
Title : SVOA CALIBRATION

35	11.181	1351	1359	1360	rBV7	10049	19513	0.37%	0.031%
36	11.199	1360	1362	1370	rVB8	10002	19201	0.36%	0.031%
37	11.328	1380	1384	1389	rVB3	12182	17339	0.33%	0.028%
38	11.457	1398	1406	1411	rVB8	16106	40906	0.77%	0.066%
39	11.581	1423	1427	1434	rVB8	12876	23951	0.45%	0.039%
40	11.675	1439	1443	1450	rVV10	7345	19568	0.37%	0.032%
41	11.822	1463	1468	1472	rBV4	14605	24745	0.47%	0.040%
42	11.963	1481	1492	1497	rBV	136542	241704	4.55%	0.390%
43	12.004	1497	1499	1503	rVB2	16069	21087	0.40%	0.034%
44	12.122	1513	1519	1525	rBV3	65327	122562	2.31%	0.198%
45	12.187	1525	1530	1532	rVV6	10371	18451	0.35%	0.030%
46	12.228	1533	1537	1545	rVB9	31673	64927	1.22%	0.105%
47	12.334	1550	1555	1560	rVB4	17708	32564	0.61%	0.053%
48	12.387	1560	1564	1576	rVB7	15271	34972	0.66%	0.056%
49	12.499	1577	1583	1587	rBV5	20307	30687	0.58%	0.049%
50	12.616	1599	1603	1608	rBV6	18808	33321	0.63%	0.054%
51	12.657	1608	1610	1614	rVV4	11472	13212	0.25%	0.021%
52	12.722	1616	1621	1625	rVB7	14939	25378	0.48%	0.041%
53	12.804	1630	1635	1640	rBV9	12575	29031	0.55%	0.047%
54	13.175	1691	1698	1702	rBV9	15744	35931	0.68%	0.058%
55	13.263	1707	1713	1716	rVV4	25341	51124	0.96%	0.082%
56	13.304	1716	1720	1728	rVV4	32581	75756	1.43%	0.122%
57	13.375	1728	1732	1735	rVV5	8416	16643	0.31%	0.027%
58	13.416	1737	1739	1746	rVB6	11595	21927	0.41%	0.035%
59	13.516	1751	1756	1763	rVV6	22160	46866	0.88%	0.076%
60	13.646	1773	1778	1782	rBV7	16003	27371	0.52%	0.044%
61	13.693	1782	1786	1791	rBV8	16686	34713	0.65%	0.056%
62	13.798	1797	1804	1805	rBV7	17302	28859	0.54%	0.047%
63	13.875	1811	1817	1822	rBV	1787143	2461546	46.33%	3.969%
64	13.922	1822	1825	1831	rVB	109312	138387	2.60%	0.223%
65	14.022	1837	1842	1848	rBV7	25598	59228	1.11%	0.095%
66	14.075	1848	1851	1856	rVB5	15204	24903	0.47%	0.040%
67	14.163	1856	1866	1879	rVB	2086715	3207440	60.36%	5.171%
68	14.275	1880	1885	1893	rBV2	55543	109356	2.06%	0.176%
69	14.393	1899	1905	1909	rBV4	67305	104833	1.97%	0.169%
70	14.469	1909	1918	1925	rBV	1718644	2638600	49.66%	4.254%
71	14.534	1925	1929	1936	rVB	45577	75302	1.42%	0.121%

LSC Area Percent Report

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Operator : CG/JU
Sample : L5128-03
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : CB8C2

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_P\METHODS\SOM-EPA-BP121820MA.M
Title : SVOA CALIBRATION

72	14.675	1948	1953	1960	rBV	554645	854311	16.08%	1.377%
73	15.216	2039	2045	2051	rBV	699127	990115	18.63%	1.596%
74	15.334	2062	2065	2067	rBV4	12683	15141	0.28%	0.024%
75	15.363	2068	2070	2077	rVB6	22123	36769	0.69%	0.059%
76	15.469	2081	2088	2094	rBV	2603481	3752253	70.62%	6.050%
77	15.592	2103	2109	2116	rBV	649353	951235	17.90%	1.534%
78	15.698	2122	2127	2141	rVB3	76399	200168	3.77%	0.323%
79	15.981	2169	2175	2182	rBV	782673	1083262	20.39%	1.747%
80	16.151	2197	2204	2210	rBV	263644	442725	8.33%	0.714%
81	16.316	2228	2232	2237	rBV2	38329	58602	1.10%	0.094%
82	16.498	2257	2263	2267	rBV	421818	589893	11.10%	0.951%
83	16.545	2267	2271	2276	rVB	84504	121655	2.29%	0.196%
84	16.634	2283	2286	2290	rVB5	17425	23916	0.45%	0.039%
85	16.763	2305	2308	2319	rVB2	49124	90157	1.70%	0.145%
86	17.028	2349	2353	2359	rVB3	72438	111749	2.10%	0.180%
87	17.151	2372	2374	2377	rVB4	16029	15585	0.29%	0.025%
88	17.228	2381	2387	2394	rVV	2090473	3099576	58.33%	4.997%
89	17.328	2398	2404	2414	rVB2	2840255	4122475	77.58%	6.647%
90	18.122	2535	2539	2545	rBV2	105387	176528	3.32%	0.285%
91	18.875	2664	2667	2670	rVB3	40395	47330	0.89%	0.076%
92	19.298	2736	2739	2742	rBV2	45339	49868	0.94%	0.080%
93	19.569	2780	2785	2790	rBV	3778968	5063115	95.29%	8.163%
94	19.610	2790	2792	2802	rVB	372349	498902	9.39%	0.804%
95	21.033	3031	3034	3042	rBV2	500655	754364	14.20%	1.216%
96	21.098	3042	3045	3053	rVB	137449	191786	3.61%	0.309%
97	21.322	3078	3083	3093	rBV	2949550	3666672	69.01%	5.912%
98	22.551	3288	3292	3301	rVB	541105	759678	14.30%	1.225%
99	23.527	3451	3458	3471	rVB2	2566842	5313542	100.00%	8.567%
100	23.680	3476	3484	3494	rVB	1845968	3763963	70.84%	6.069%

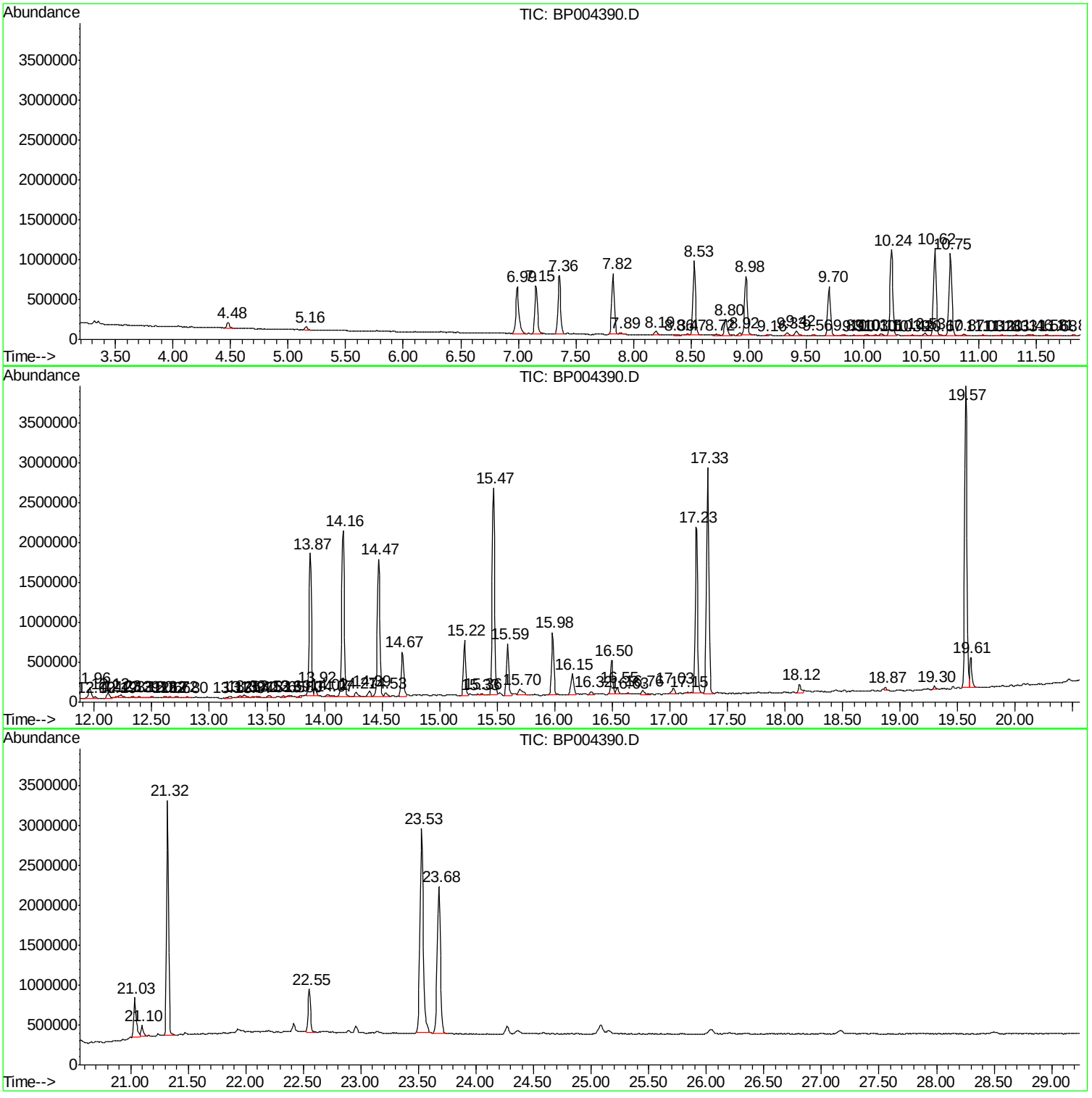
Sum of corrected areas: 62023019

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121820\
Data File : BP004390.D
Acq On : 19 Dec 2020 12:48
Operator : CG/JU
Sample : L5128-03
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB8C2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP121820MA.M
Quant Title : SVOA CALIBRATION

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



Library Search Compound Report

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ClientSampleId : CB8C2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP121820MA.M
Quant Title : SVOA CALIBRATION

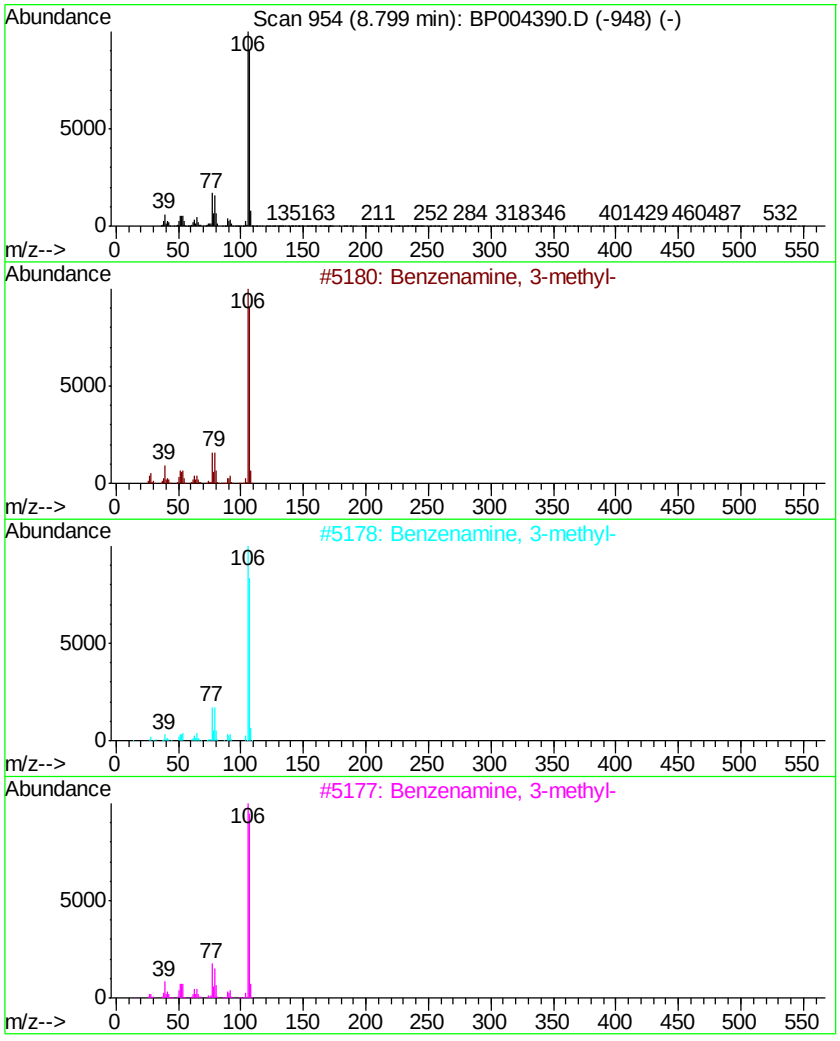
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

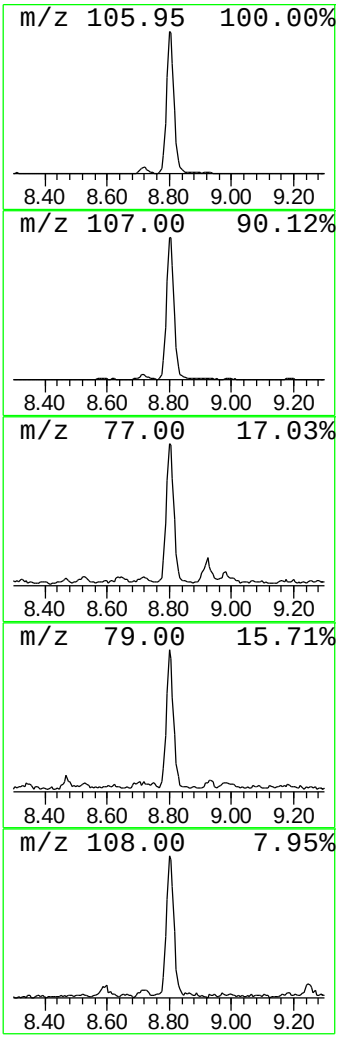
****

Peak Number 1 Benzenamine, 3-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.80	5.56 ng/ul	340921	1,4-Dichlorobenzene-d4	7.82

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenamine, 3-methyl-	107	C7H9N	000108-44-1	97
2	Benzenamine, 3-methyl-	107	C7H9N	000108-44-1	97
3	Benzenamine, 3-methyl-	107	C7H9N	000108-44-1	97
4	Benzenamine, 3-methyl-	107	C7H9N	000108-44-1	95
5	o-Toluidine	107	C7H9N	000095-53-4	94





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121820\
Data File : BP004390.D
Acq On : 19 Dec 2020 12:48
Operator : CG/JU
Sample : L5128-03
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : CB8C2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP121820MA.M
Quant Title : SVOA CALIBRATION

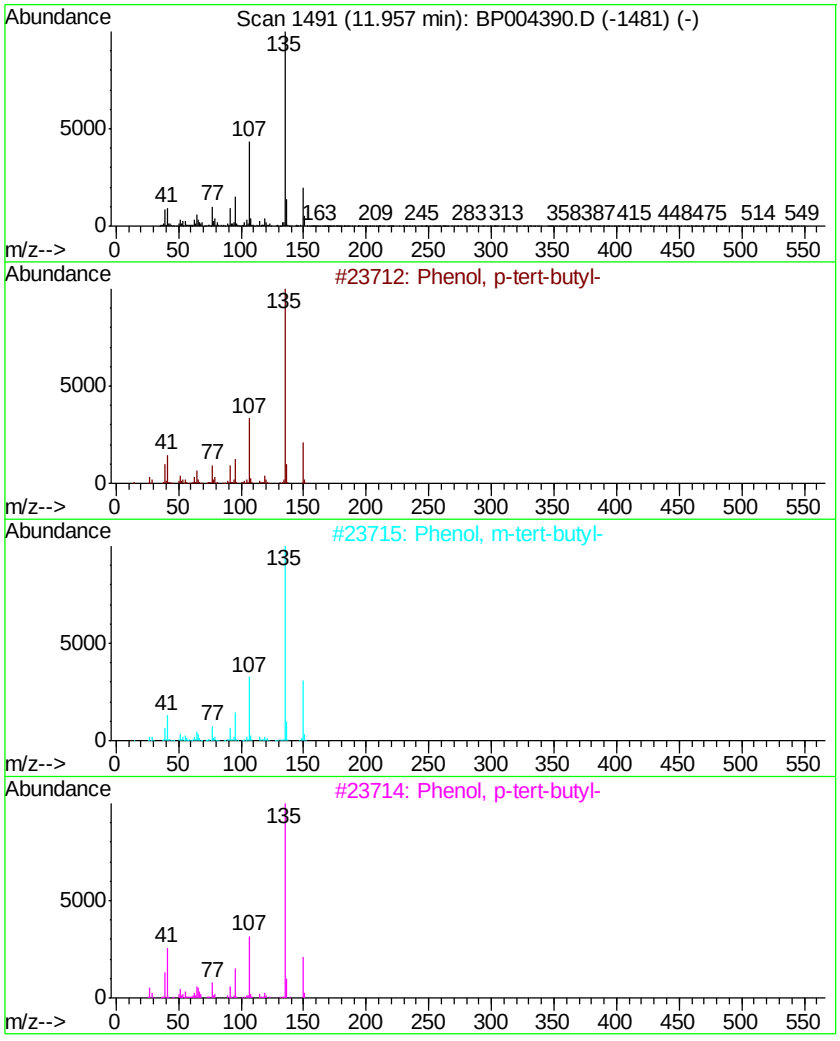
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

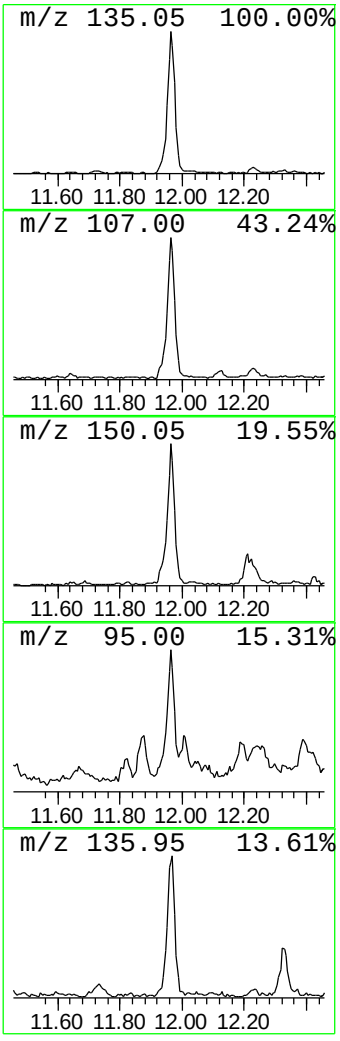
****

Peak Number 2 Phenol, p-tert-butyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.96	2.58 ng/ul	241704	Naphthalene-d8	10.62

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	97
2	Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	91
3	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	91
4	Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	86
5	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	81





Library Search Compound Report

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Data File : BP004390.D
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Sample : L5128-03
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : CB8C2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP121820MA.M
Quant Title : SVOA CALIBRATION

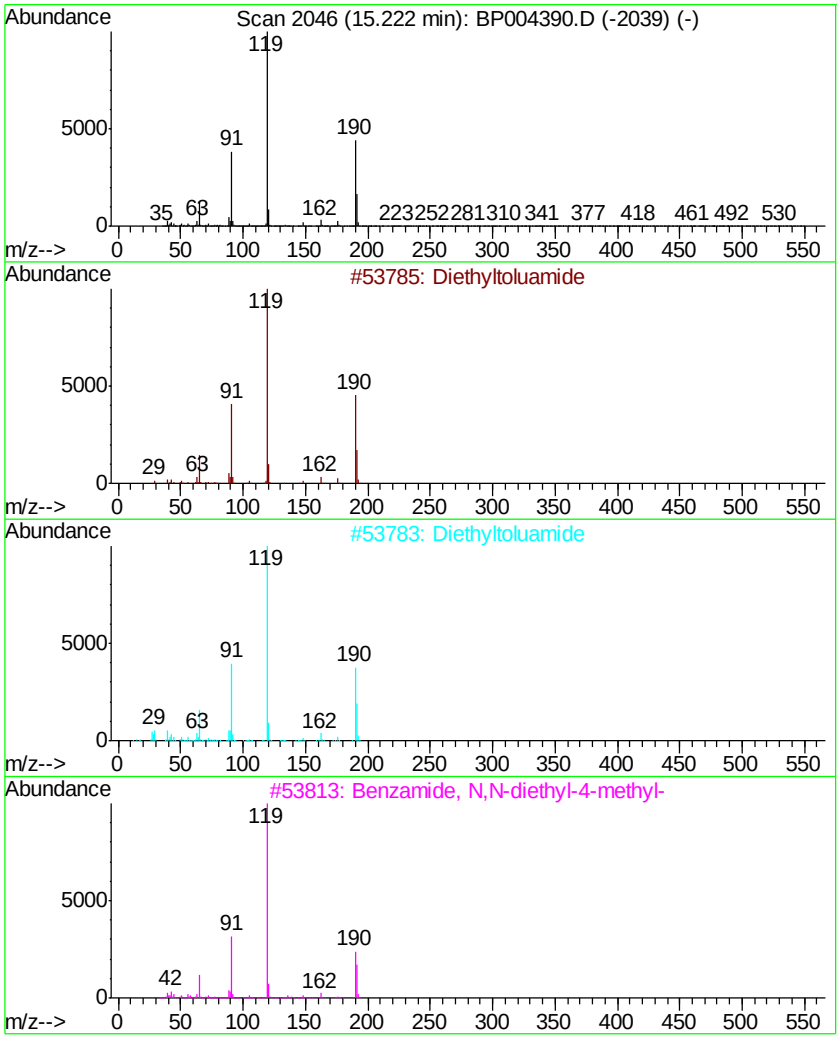
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

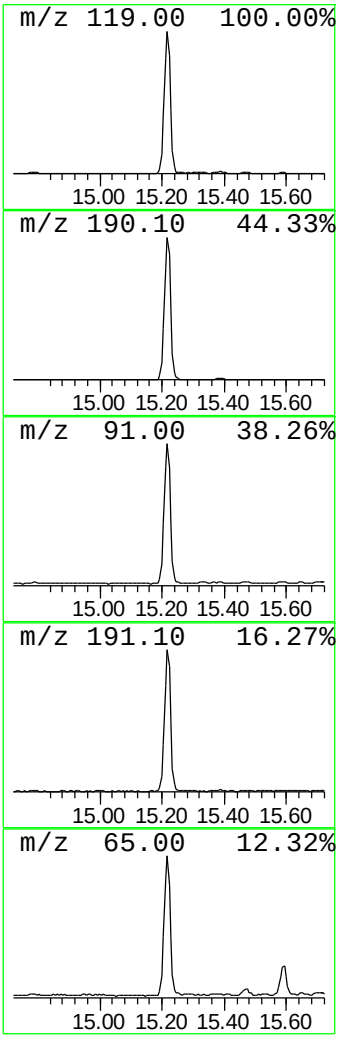
****

Peak Number 3 Diethyltoluamide Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.22	7.50 ng/ul	990115	Acenaphthene-d10	14.47

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Diethyltoluamide	191	C12H17NO	000134-62-3	96
2	Diethyltoluamide	191	C12H17NO	000134-62-3	94
3	Benzamide, N,N-diethyl-4-methyl-	191	C12H17NO	002728-05-4	93
4	Diethyltoluamide	191	C12H17NO	000134-62-3	81
5	Diethyltoluamide	191	C12H17NO	000134-62-3	60





Library Search Compound Report

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Data File : BP004390.D
Acq On : 19 Dec 2020 12:48
Operator : CG/JU
Sample : L5128-03
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : CB8C2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP121820MA.M
Quant Title : SVOA CALIBRATION

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

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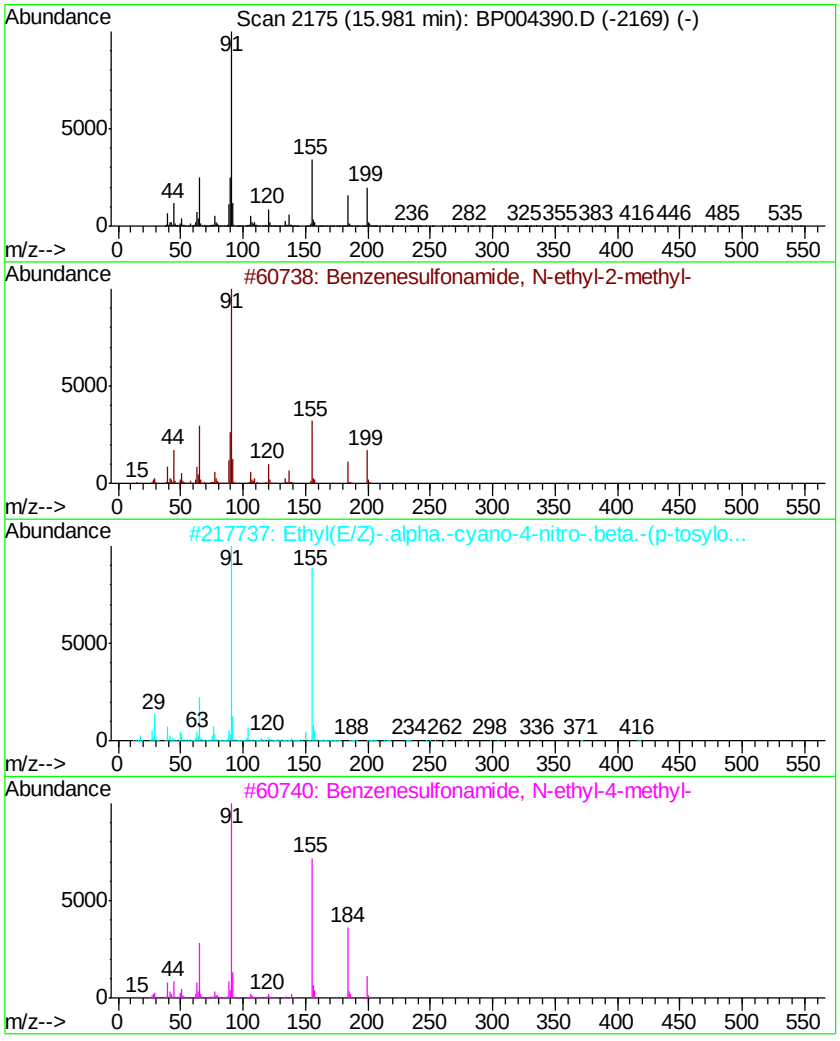
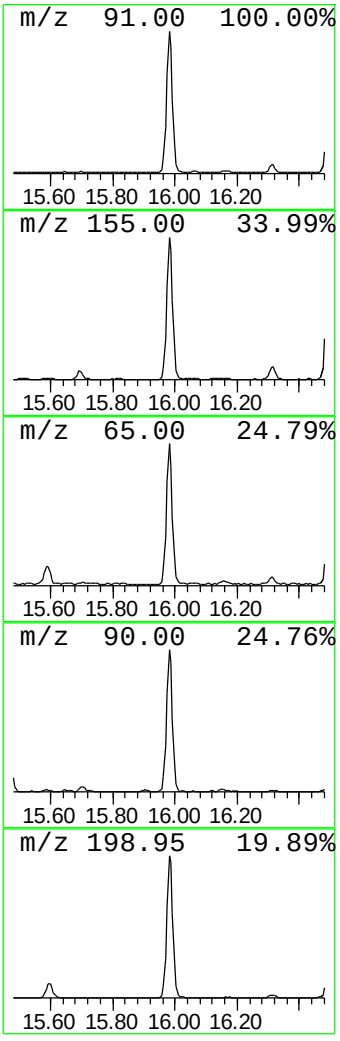
Peak Number 4 Benzenesulfonamide, N-ethyl... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.98	6.99 ng/ul	1083260	Phenanthrene-d10	17.23

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenesulfonamide, N-ethyl-2-me...	199	C9H13NO2S	001077-56-1	96
2	Ethyl(E/Z)-.alpha.-cyano-4-nitro...	416	C19H16N2O7S	1000287-26-8	55
3	Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	53
4	N3, N.alpha.-bis*ptoluenesulfony...	462	C20H22N4O5S2	1000130-21-9	53
5	Benzenesulfonamide, N-ethyl-2-me...	199	C9H13NO2S	001077-56-1	52





Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121820\
Data File : BP004390.D
Acq On : 19 Dec 2020 12:48
Operator : CG/JU
Sample : L5128-03
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : CB8C2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP121820MA.M
Quant Title : SVOA CALIBRATION

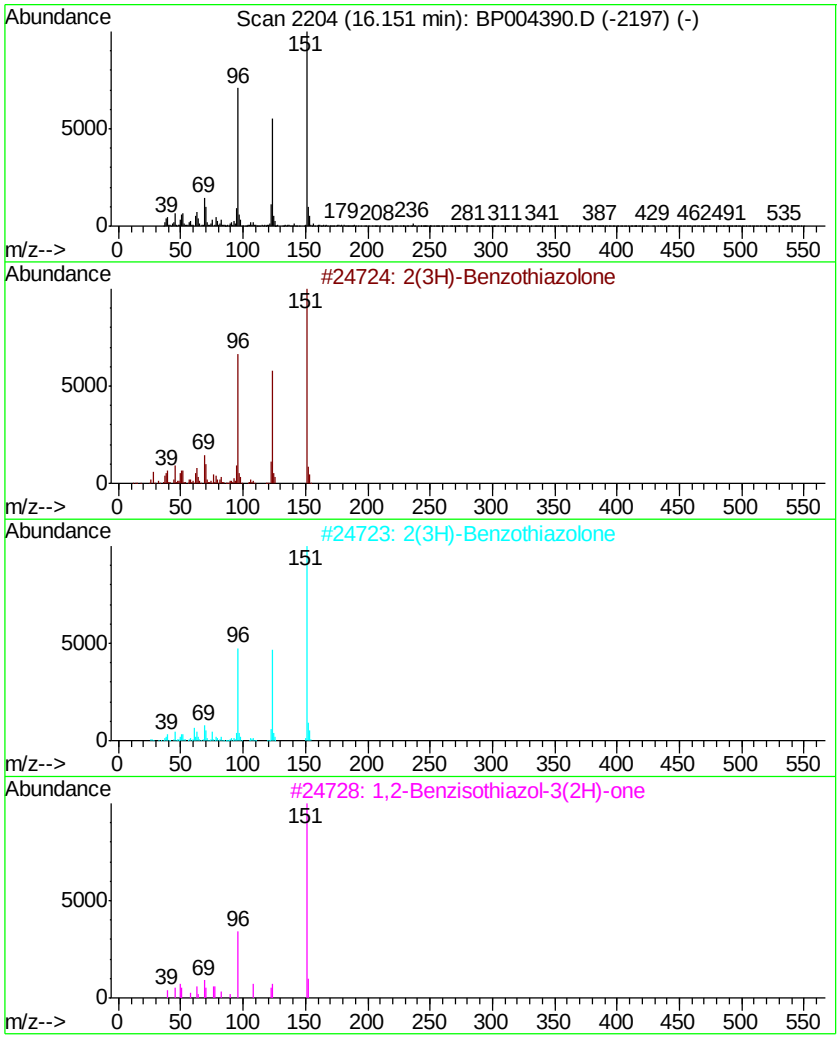
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

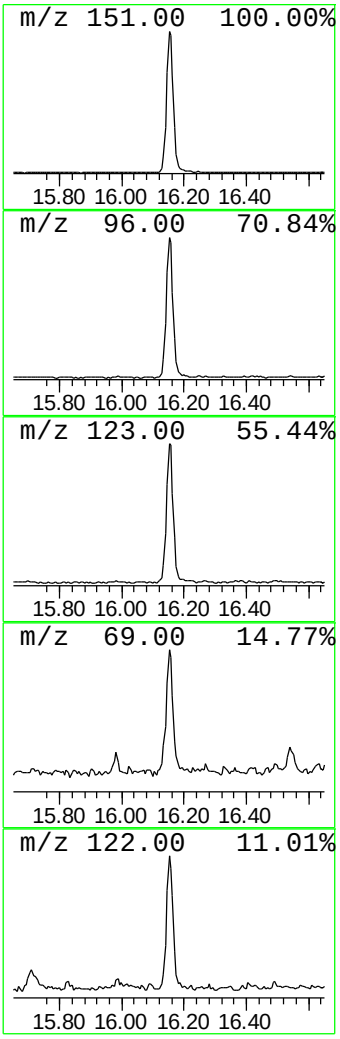
****

Peak Number 5 2(3H)-Benzothiazolone Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.15	2.86 ng/ul	442725	Phenanthrene-d10	17.23

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2(3H)-Benzothiazolone	151	C7H5NOS	000934-34-9	94
2	2(3H)-Benzothiazolone	151	C7H5NOS	000934-34-9	94
3	1,2-Benzisothiazol-3(2H)-one	151	C7H5NOS	002634-33-5	38
4	Thieno[2,3-b]pyridine-N-oxide	151	C7H5NOS	025557-50-0	38
5	2H-Pyran-2-carboxylic acid, 5-et...	196	C10H12O4	019776-81-9	37





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Acq On : 19 Dec 2020 12:48
Operator : CG/JU
Sample : L5128-03
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : CB8C2

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Quant Title : SVOA CALIBRATION

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

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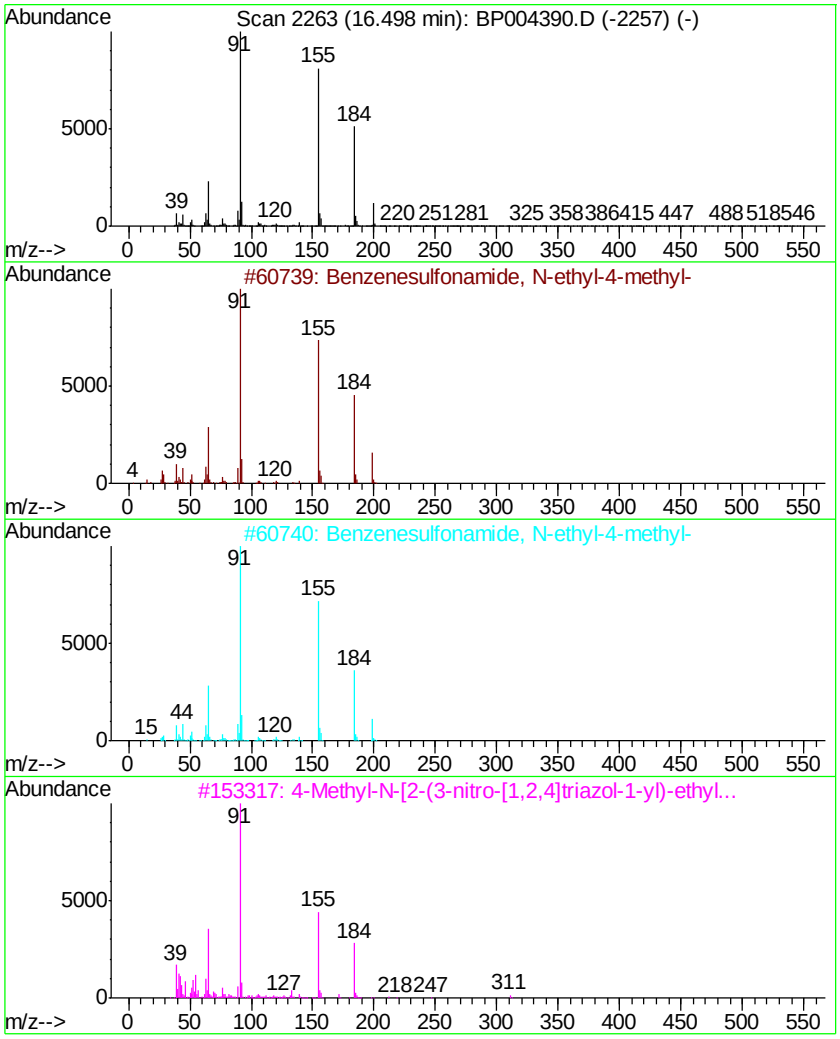
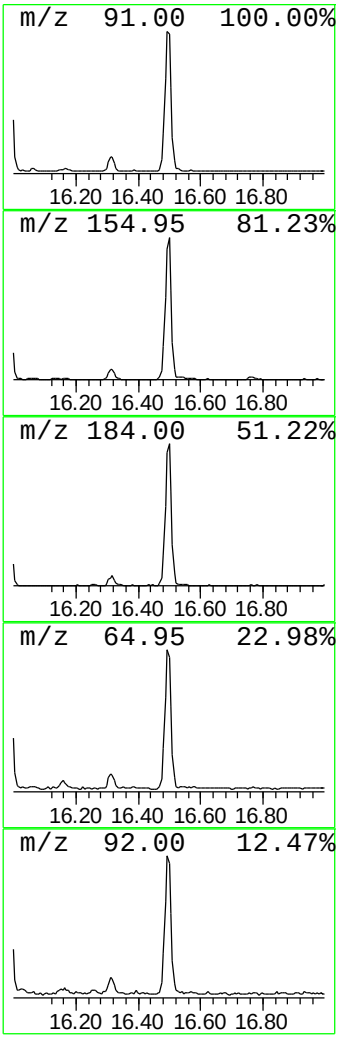
Peak Number 6 Benzenesulfonamide, N-ethyl... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.50	3.81 ng/ul	589893	Phenanthrene-d10	17.23

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	96
3	4-Methyl-N-[2-(3-nitro-[1,2,4]tr...	311	C11H13N5O4S	1000301-03-5	83
4	Furazan-3-carboxylic acid, 4-ami...	325	C12H15N5O4S	1000304-69-4	74
5	(Toluene-4-sulfonylamino)-acetic...	283	C13H17NO4S	1000190-48-0	72





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Data File : BP004390.D
Acq On : 19 Dec 2020 12:48
Operator : CG/JU
Sample : L5128-03
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : CB8C2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP121820MA.M
Quant Title : SVOA CALIBRATION

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

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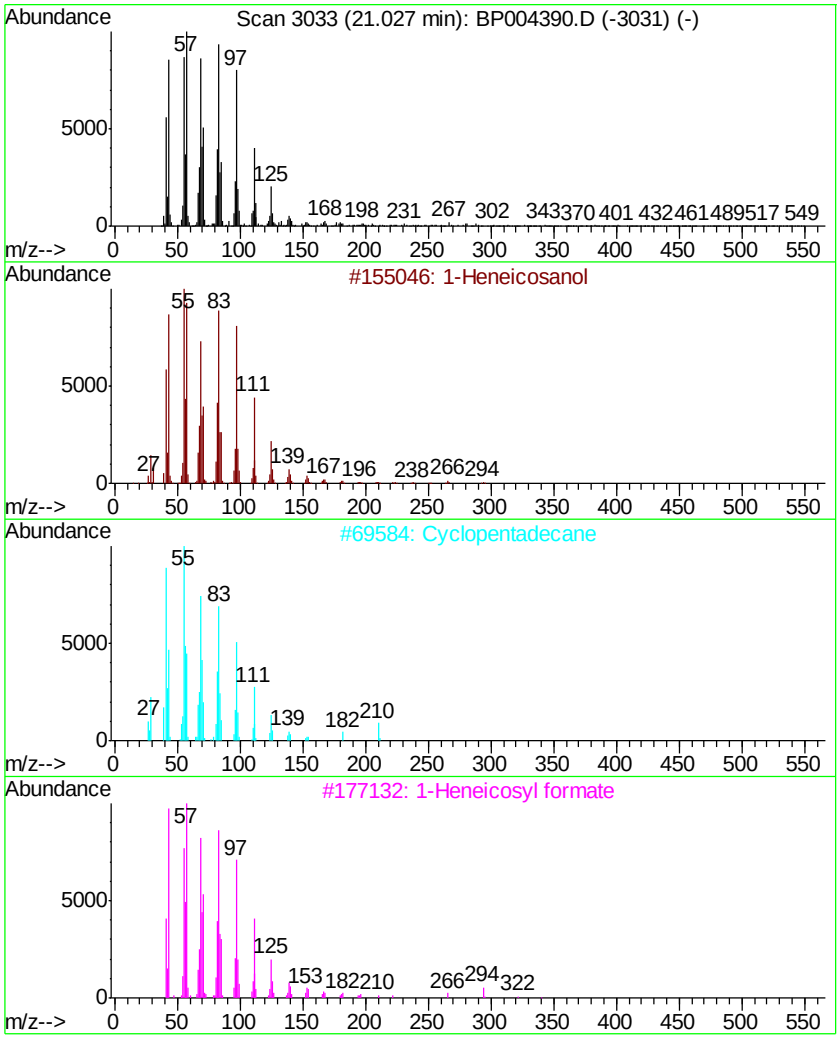
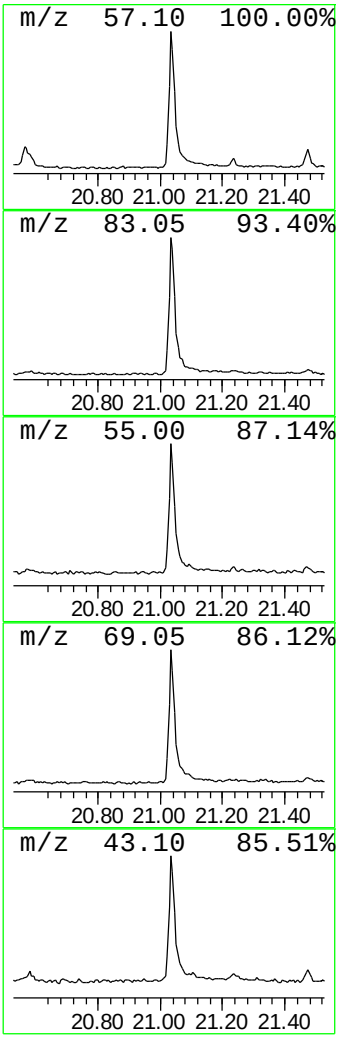
Peak Number 7 1-Heneicosanol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.03	4.11 ng/ul	754364	Chrysene-d12	21.32

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heneicosanol	312	C21H44O	015594-90-8	95
2	Cyclopentadecane	210	C15H30	000295-48-7	95
3	1-Heneicosyl formate	340	C22H44O2	077899-03-7	94
4	Behenic alcohol	326	C22H46O	000661-19-8	94
5	n-Nonadecanol-1	284	C19H40O	001454-84-8	94





Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121820\
Data File : BP004390.D
Acq On : 19 Dec 2020 12:48
Operator : CG/JU
Sample : L5128-03
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : CB8C2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP121820MA.M
Quant Title : SVOA CALIBRATION

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

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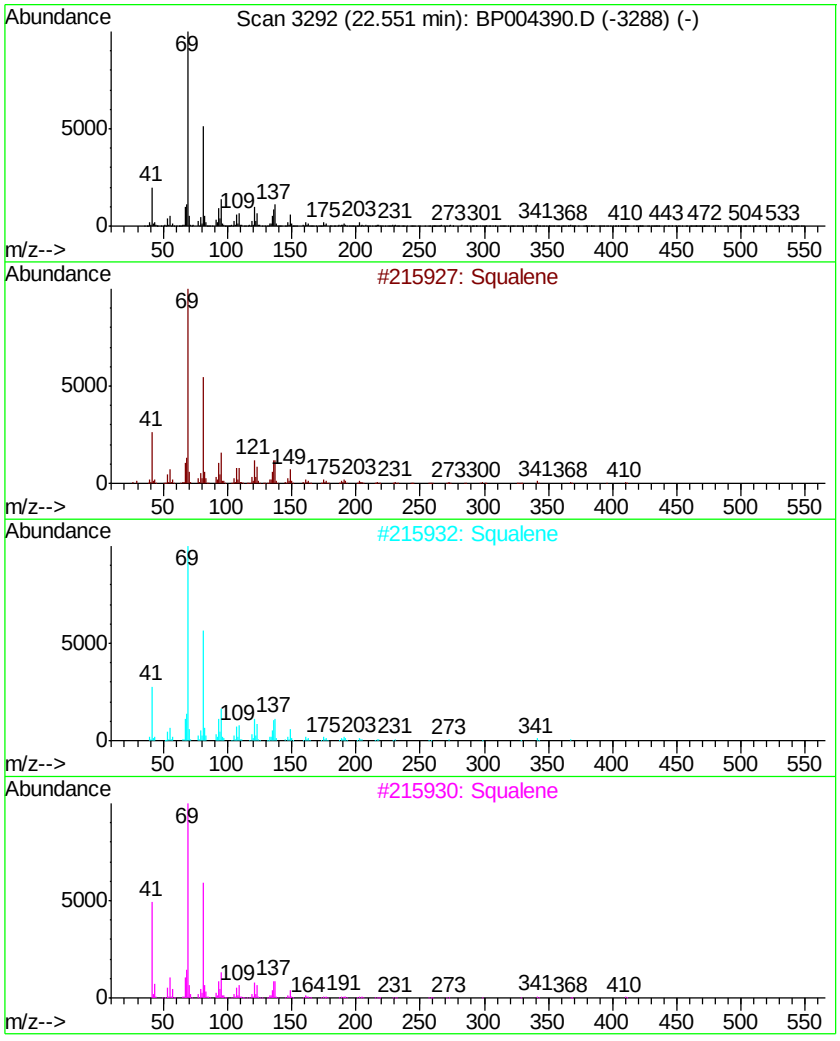
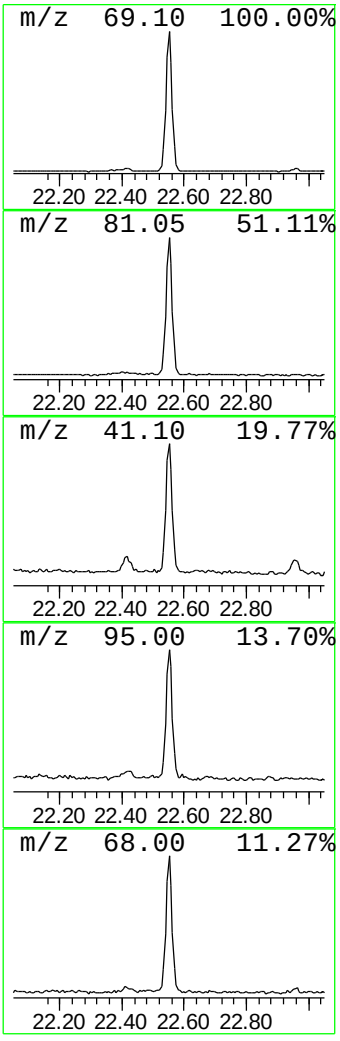
Peak Number 8 Squalene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.55	4.04 ng/ul	759678	Perylene-d12	23.68

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Squalene	410	C30H50	000111-02-4	96
2	Squalene	410	C30H50	000111-02-4	94
3	Squalene	410	C30H50	000111-02-4	93
4	Supraene	410	C30H50	007683-64-9	91
5	Docosa-2,6,10,14,18-pentaen-22-a...	384	C27H44O	1000163-04-7	90





Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP121820\
Data File : BP004390.D
Acq On : 19 Dec 2020 12:48
Operator : CG/JU
Sample : L5128-03
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : CB8C2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP121820MA.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
Benzenamine, 3-me...	8.80	5.6	ng/ul	340921	1	7.82	1226430	20.0
Phenol, p-tert-bu...	11.96	2.6	ng/ul	241704	2	10.62	1876560	20.0
Diethyltoluamide	15.22	7.5	ng/ul	990115	3	14.47	2638600	20.0
Benzenesulfonamid...	15.98	7.0	ng/ul	1083260	4	17.23	3099580	20.0
2(3H)-Benzothiazo...	16.15	2.9	ng/ul	442725	4	17.23	3099580	20.0
Benzenesulfonamid...	16.50	3.8	ng/ul	589893	4	17.23	3099580	20.0
1-Heneicosanol	21.03	4.1	ng/ul	754364	5	21.32	3666670	20.0
Squalene	22.55	4.0	ng/ul	759678	6	23.68	3763960	20.0