LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP021721\
Data File : BP004771.D
Acq On : 17 Feb 2021 14:04
Operator : CG/JU
Sample : M1407-02
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DBGJ3

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-BP020821MA.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	7	11	15	rBV	10970	12804	0.98%	0.099%
2	6.934	628	637	646	rBV	46879	82150	6.28%	0.634%
3	7.099	658	665	673	rBV	131023	204999	15.67%	1.583%
4	7.299	692	699	707	rBV	170116	274058	20.94%	2.117%
5	7.763	771	778	785	rBV	192140	297150	22.71%	2.295%
6	8.469	891	898	906	rBV	121077	190214	14.54%	1.469%
7	8.604	915	921	932	rVB2	40515	68255	5.22%	0.527%
8	8.916	968	974	985	rVB	160649	274149	20.95%	2.117%
9	9.181	1012	1019	1025	rBV2	29007	46569	3.56%	0.360%
10	9.346	1042	1047	1051	rBV5	5727	9171	0.70%	0.071%
11	9.640	1089	1097	1104	rBV	151050	247998	18.95%	1.915%
12	9.763	1110	1118	1126	rVB4	21082	37978	2.90%	0.293%
13	9.957	1144	1151	1156	rBV8	5177	9305	0.71%	0.072%
14	10.175	1181	1188	1197	rVV	215242	361814	27.65%	2.794%
15	10.557	1246	1253	1260	rVV	221967	366461	28.00%	2.830%
16	10.704	1269	1278	1293	rBV3	41649	92576	7.07%	0.715%
17	10.916	1308	1314	1322	rBV7	17408	40629	3.10%	0.314%
18	11.146	1348	1353	1363	rVB4	11595	22284	1.70%	0.172%
19	11.587	1416	1428	1432	rBV	22955	38765	2.96%	0.299%
20	11.640	1432	1437	1441	rVV	15828	24680	1.89%	0.191%
21	11.687	1441	1445	1450	rVB4	7566	13547	1.04%	0.105%
22	11.769	1452	1459	1465	rVB9	3368	8852	0.68%	0.068%
23	11.916	1479	1484	1487	rBV4	5241	9398	0.72%	0.073%
24	12.134	1514	1521	1522	rVV5	5341	10938	0.84%	0.084%
25	12.163	1522	1526	1532	rVB6	9840	18760	1.43%	0.145%
26	12.263	1539	1543	1549	rVV5	6234	11962	0.91%	0.092%
27	12.322	1549	1553	1557	rVB5	5414	7974	0.61%	0.062%
28	12.445	1564	1574	1581	rVB4	19278	36026	2.75%	0.278%
29	12.557	1585	1593	1596	rBV	16353	26138	2.00%	0.202%
30	12.593	1596	1599	1606	rVV5	5933	11488	0.88%	0.089%
31	12.740	1619	1624	1629	rVB8	5129	7641	0.58%	0.059%
32	12.822	1632	1638	1643	rBV5	8847	15874	1.21%	0.123%
33	12.928	1647	1656	1664	rVB7	16899	49610	3.79%	0.383%
34	13.087	1677	1683	1685	rBV4	5767	10104	0.77%	0.078%

LSC Area Percent Report

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Operator : CG/JU
Sample : M1407-02
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DBGJ3

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-BP020821MA.M
Title : SVOA CALIBRATION

35	13.122	1685	1689	1693	rBV5	8970	15373	1.17%	0.119%
36	13.198	1698	1702	1706	rVV2	9631	17858	1.36%	0.138%
37	13.251	1706	1711	1719	rVV9	10085	23928	1.83%	0.185%
38	13.375	1728	1732	1736	rVV7	4288	7741	0.59%	0.060%
39	13.428	1736	1741	1743	rVV4	5000	8297	0.63%	0.064%
40	13.463	1743	1747	1754	rVB3	13502	21035	1.61%	0.162%
41	13.581	1762	1767	1776	rVB7	16799	38303	2.93%	0.296%
42	13.669	1776	1782	1787	rBV10	4012	8601	0.66%	0.066%
43	13.728	1787	1792	1797	rVV4	10081	17338	1.32%	0.134%
44	13.822	1802	1808	1814	rVV	386070	534152	40.82%	4.125%
45	13.904	1818	1822	1827	rVB2	15009	21702	1.66%	0.168%
46	14.104	1849	1856	1868	rVB	416325	617538	47.19%	4.769%
47	14.410	1902	1908	1914	rVV	317406	463776	35.44%	3.582%
48	14.475	1914	1919	1925	rVB	135262	196261	15.00%	1.516%
49	14.539	1925	1930	1935	rBV	18168	24789	1.89%	0.191%
50	14.616	1936	1943	1950	rBV	35961	57413	4.39%	0.443%
51	14.687	1950	1955	1959	rBV2	14265	21138	1.62%	0.163%
52	14.751	1963	1966	1974	rVB7	5616	8704	0.67%	0.067%
53	14.822	1974	1978	1984	rVB7	8174	15066	1.15%	0.116%
54	14.957	1995	2001	2006	rBV3	36303	56438	4.31%	0.436%
55	15.034	2011	2014	2020	rVB3	14864	22539	1.72%	0.174%
56	15.245	2046	2050	2053	rBV4	6366	9838	0.75%	0.076%
57	15.339	2057	2066	2072	rVV2	82186	141180	10.79%	1.090%
58	15.404	2072	2077	2083	rVV	515401	765709	58.52%	5.914%
59	15.463	2083	2087	2092	rVV2	58986	90544	6.92%	0.699%
60	15.522	2092	2097	2105	rVV2	212141	359167	27.45%	2.774%
61	15.592	2105	2109	2115	rVV8	11578	30432	2.33%	0.235%
62	15.651	2115	2119	2121	rVV5	11944	19953	1.52%	0.154%
63	15.686	2121	2125	2133	rVV7	16092	37686	2.88%	0.291%
64	15.751	2133	2136	2142	rVB5	6781	11737	0.90%	0.091%
65	15.834	2142	2150	2155	rBV6	12744	33435	2.56%	0.258%
66	15.904	2155	2162	2171	rVB9	21033	57231	4.37%	0.442%
67	16.122	2191	2199	2205	rBV9	10432	26008	1.99%	0.201%
68	16.428	2244	2251	2256	rBV	43568	68266	5.22%	0.527%
69	16.545	2267	2271	2274	rBV6	5045	7654	0.58%	0.059%
70	16.581	2274	2277	2284	rVB7	6201	8808	0.67%	0.068%
71	16.810	2310	2316	2327	rVB3	43090	71251	5.44%	0.550%

LSC Area Percent Report

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Operator : CG/JU
Sample : M1407-02
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DBGJ3

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-BP020821MA.M
Title : SVOA CALIBRATION

72	16.951	2335	2340	2351	rVV10	8900	23863	1.82%	0.184%
73	17.122	2365	2369	2371	rBV3	13031	16131	1.23%	0.125%
74	17.163	2371	2376	2381	rVV	380955	549832	42.02%	4.246%
75	17.204	2381	2383	2386	rVV	34583	46589	3.56%	0.360%
76	17.263	2388	2393	2404	rVB	647854	888671	67.91%	6.863%
77	17.539	2435	2440	2442	rBV3	22788	30749	2.35%	0.237%
78	17.569	2442	2445	2452	rVB2	26359	37067	2.83%	0.286%
79	17.798	2475	2484	2487	rBV2	6243	14814	1.13%	0.114%
80	17.910	2500	2503	2508	rBV7	9075	14690	1.12%	0.113%
81	17.998	2514	2518	2523	rVB2	17133	23483	1.79%	0.181%
82	18.057	2523	2528	2535	rBV2	90210	143869	10.99%	1.111%
83	18.151	2540	2544	2551	rVB	24058	37015	2.83%	0.286%
84	18.222	2552	2556	2560	rVB6	5350	7678	0.59%	0.059%
85	18.304	2565	2570	2573	rBV4	9004	15147	1.16%	0.117%
86	18.469	2593	2598	2602	rBV5	9296	15977	1.22%	0.123%
87	18.516	2603	2606	2611	rVV6	6707	10421	0.80%	0.080%
88	19.010	2685	2690	2695	rBV5	10562	19025	1.45%	0.147%
89	19.145	2709	2713	2716	rBV2	9421	14097	1.08%	0.109%
90	19.292	2734	2738	2745	rVB2	54916	69003	5.27%	0.533%
91	19.392	2752	2755	2759	rVB3	7936	9944	0.76%	0.077%
92	19.504	2768	2774	2787	rVB	862542	1124172	85.91%	8.682%
93	19.733	2810	2813	2818	rVB7	8228	9517	0.73%	0.074%
94	19.898	2838	2841	2844	rBV4	5565	7787	0.60%	0.060%
95	20.963	3018	3022	3028	rBV2	76874	100620	7.69%	0.777%
96	21.245	3065	3070	3076	rBV	570403	699985	53.49%	5.406%
97	22.845	3339	3342	3348	rVB4	24736	33135	2.53%	0.256%
98	23.404	3429	3437	3450	rBV	649132	1308568	100.00%	10.106%
99	23.551	3455	3462	3471	rVB2	366375	730053	55.79%	5.638%
100	24.127	3556	3560	3568	rVB4	37361	68822	5.26%	0.532%

Sum of corrected areas: 12947934

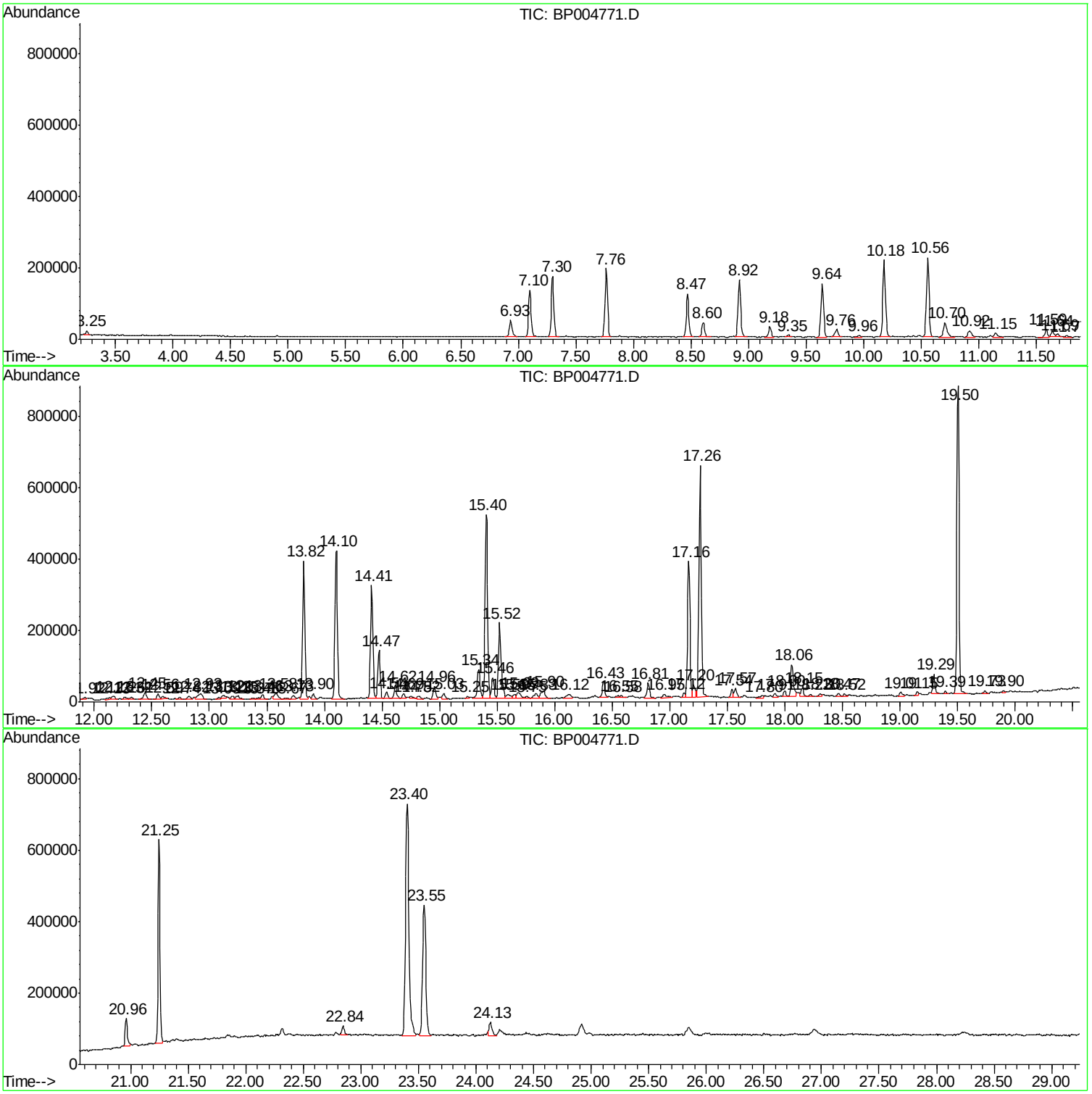
LSC Report - Integrated Chromatogram

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP021721\
Data File : BP004771.D
Acq On : 17 Feb 2021 14:04
Operator : CG/JU
Sample : M1407-02
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DBGJ3

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP020821MA.M
Quant Title : SVOA CALIBRATION

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



Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP021721\
Data File : BP004771.D
Acq On : 17 Feb 2021 14:04
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Instrument : BNA_P
ClientSampleId : DBGJ3

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP020821MA.M
Quant Title : SVOA CALIBRATION

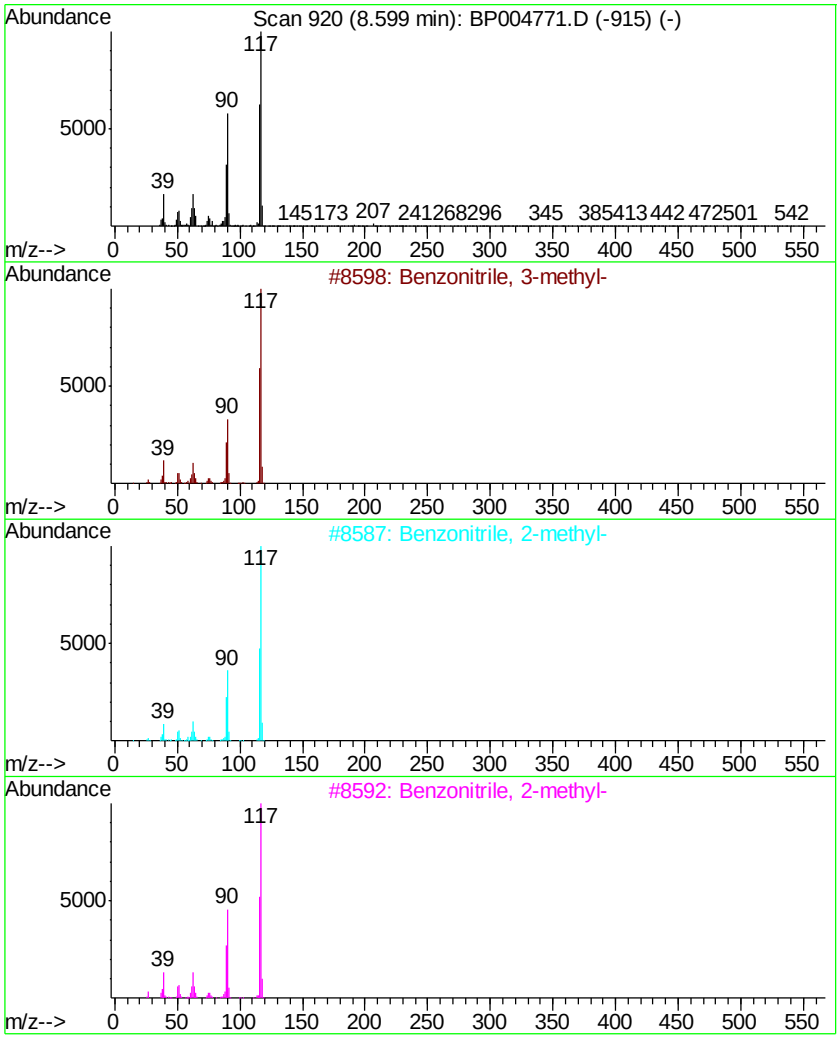
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

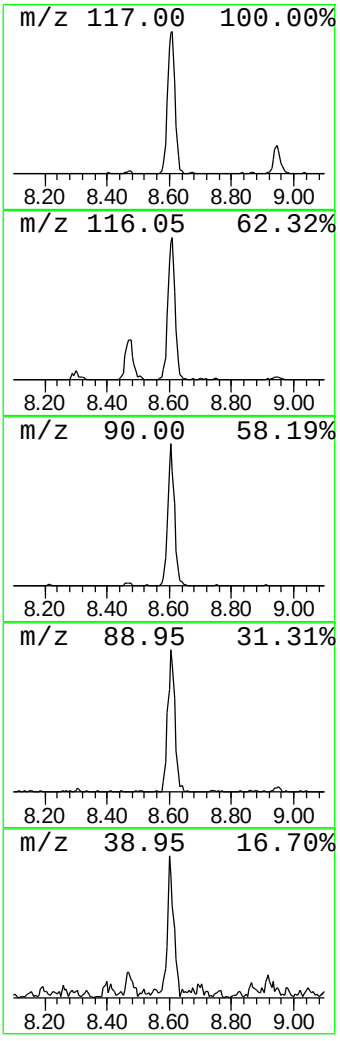
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.60	4.59 ng/ul	68255	1,4-Dichlorobenzene-d4	7.76

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzonitrile, 3-methyl-	117	C8H7N	000620-22-4	97
2	Benzonitrile, 2-methyl-	117	C8H7N	000529-19-1	97
3	Benzonitrile, 2-methyl-	117	C8H7N	000529-19-1	97
4	Benzonitrile, 2-methyl-	117	C8H7N	000529-19-1	96
5	Benzonitrile, 4-methyl-	117	C8H7N	000104-85-8	96





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP021721\
Data File : BP004771.D
Acq On : 17 Feb 2021 14:04
Operator : CG/JU
Sample : M1407-02
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DBGJ3

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP020821MA.M
Quant Title : SVOA CALIBRATION

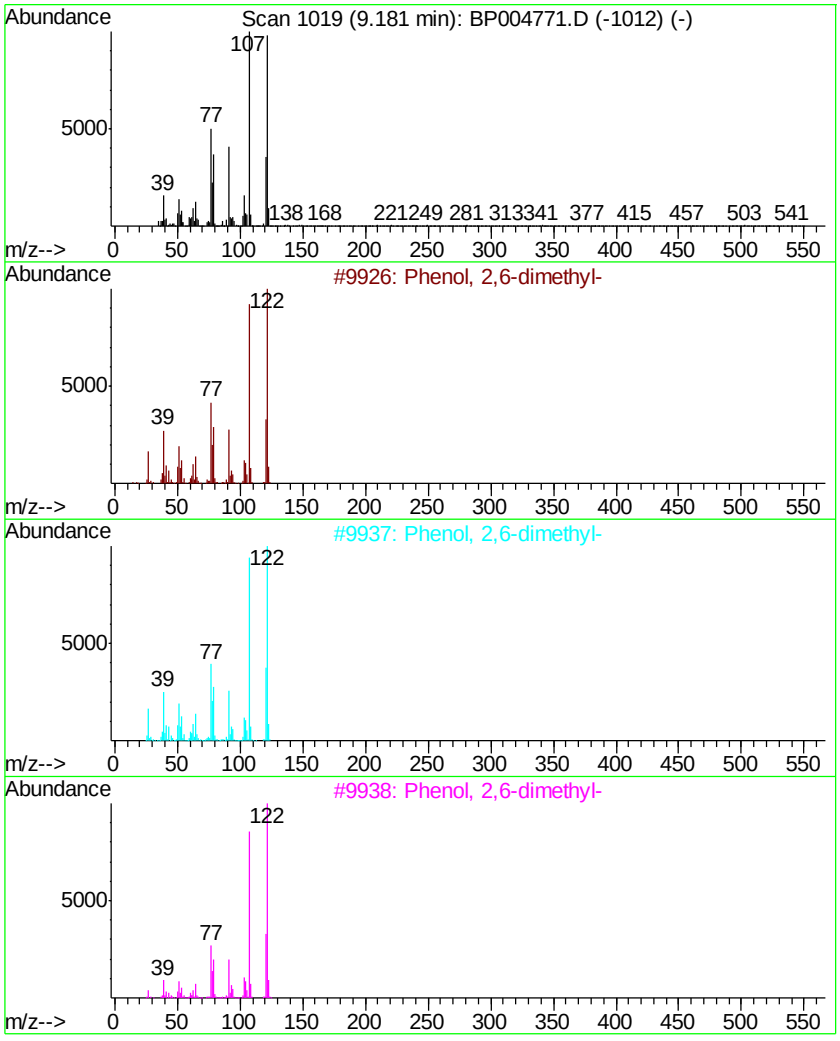
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

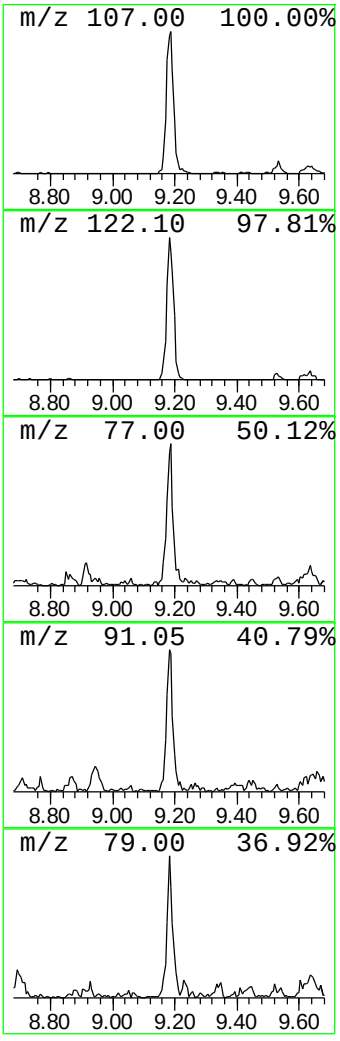
****

Peak Number 2 Phenol, 2,6-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.18	2.54 ng/ul	46569	Naphthalene-d8	10.56

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2,6-dimethyl-	122	C8H10O	000576-26-1	94
2	Phenol, 2,6-dimethyl-	122	C8H10O	000576-26-1	94
3	Phenol, 2,6-dimethyl-	122	C8H10O	000576-26-1	91
4	Phenol, 2,3-dimethyl-	122	C8H10O	000526-75-0	91
5	3-Methylbenzyl alcohol	122	C8H10O	000587-03-1	90





Library Search Compound Report

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Instrument : BNA_P
ClientSampleId : DBGJ3

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP020821MA.M
Quant Title : SVOA CALIBRATION

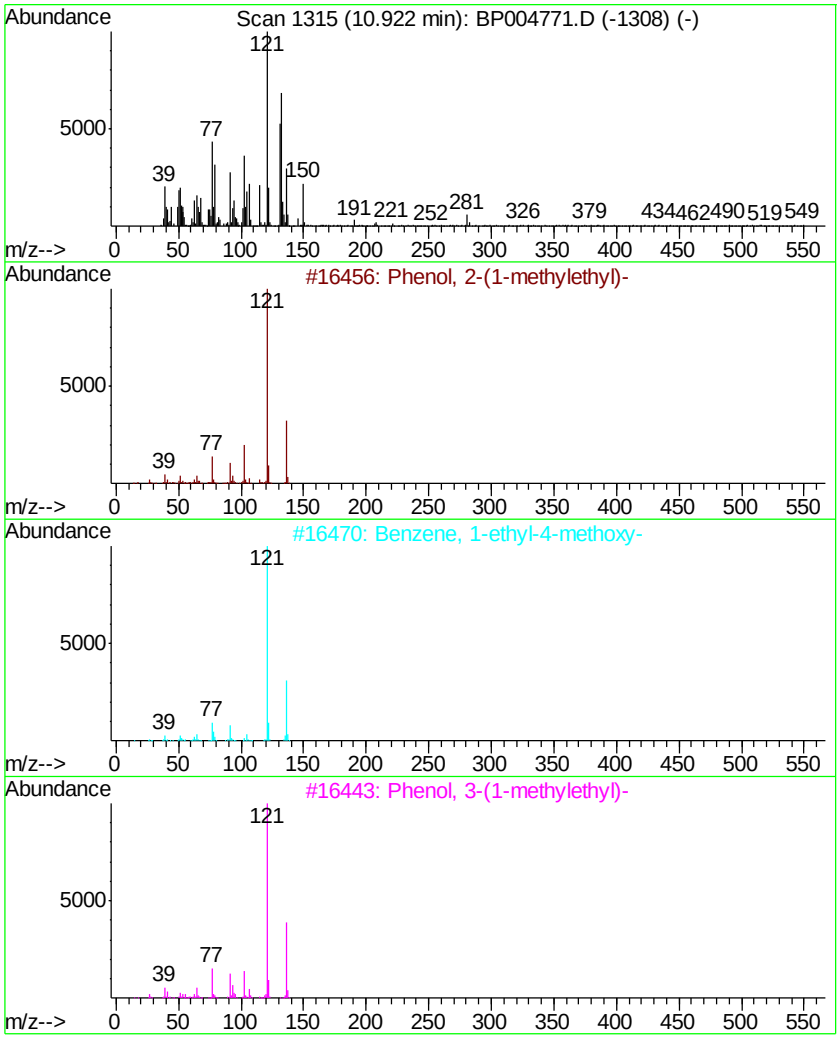
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

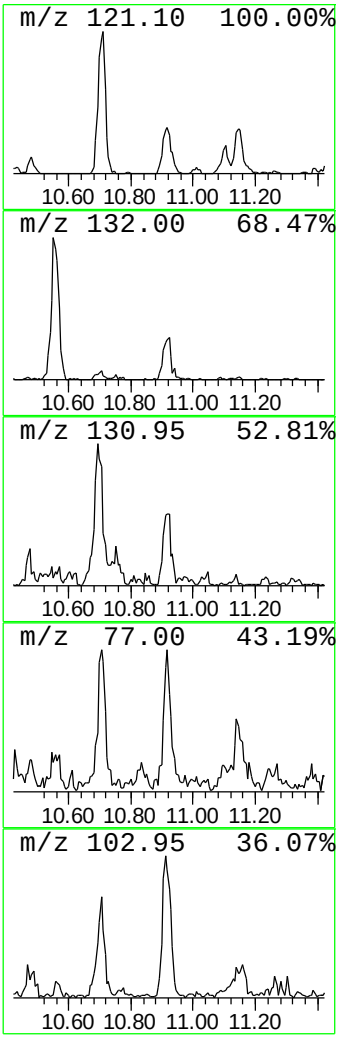
****

Peak Number 3 Phenol, 2-(1-methylethyl)- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.92	2.22 ng/ul	40629	Naphthalene-d8	10.56

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2-(1-methylethyl)-	136	C9H12O	000088-69-7	58
2	Benzene, 1-ethyl-4-methoxy-	136	C9H12O	001515-95-3	53
3	Phenol, 3-(1-methylethyl)-	136	C9H12O	000618-45-1	52
4	Benzene, 1-ethyl-4-methoxy-	136	C9H12O	001515-95-3	49
5	Phenol, 3-(1-methylethyl)-	136	C9H12O	000618-45-1	46





Library Search Compound Report

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Data File : BP004771.D
Acq On : 17 Feb 2021 14:04
Operator : CG/JU
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Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DBGJ3

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP020821MA.M
Quant Title : SVOA CALIBRATION

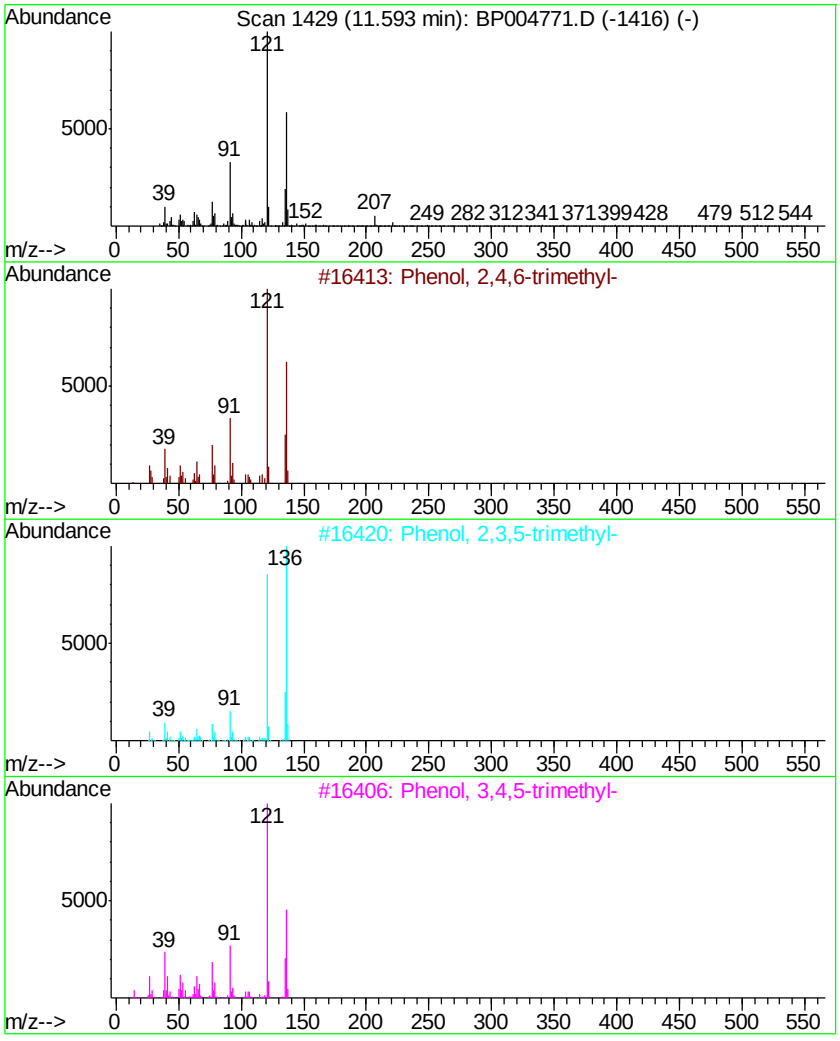
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

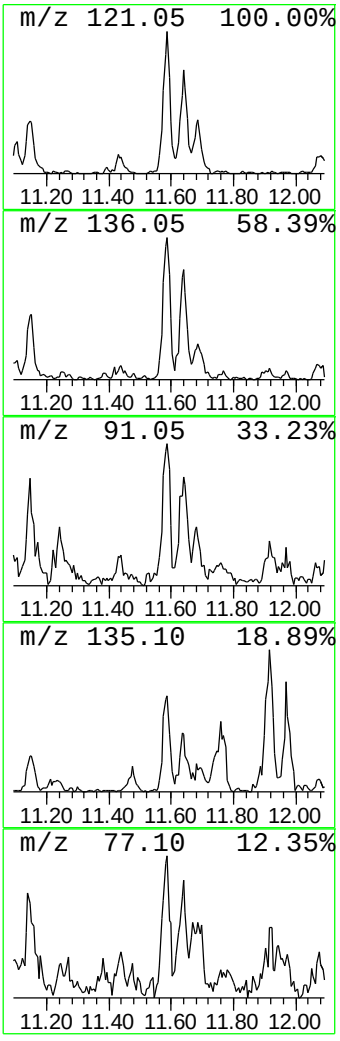
****

Peak Number 4 Phenol, 2,4,6-trimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.59	2.12 ng/ul	38765	Naphthalene-d8	10.56

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2,4,6-trimethyl-	136	C9H12O	000527-60-6	94
2	Phenol, 2,3,5-trimethyl-	136	C9H12O	000697-82-5	91
3	Phenol, 3,4,5-trimethyl-	136	C9H12O	000527-54-8	91
4	Phenol, 2,4,5-trimethyl-	136	C9H12O	000496-78-6	90
5	Phenol, 2,4,6-trimethyl-	136	C9H12O	000527-60-6	87





Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP021721\
Data File : BP004771.D
Acq On : 17 Feb 2021 14:04
Operator : CG/JU
Sample : M1407-02
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DBGJ3

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP020821MA.M
Quant Title : SVOA CALIBRATION

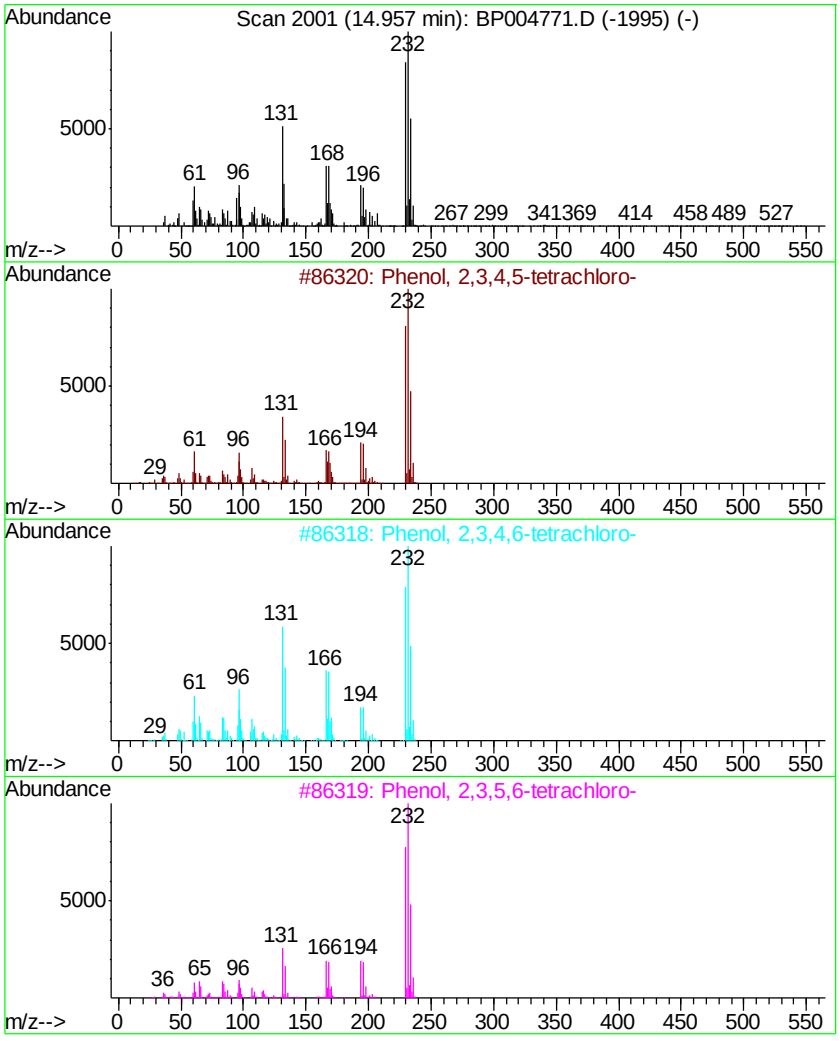
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

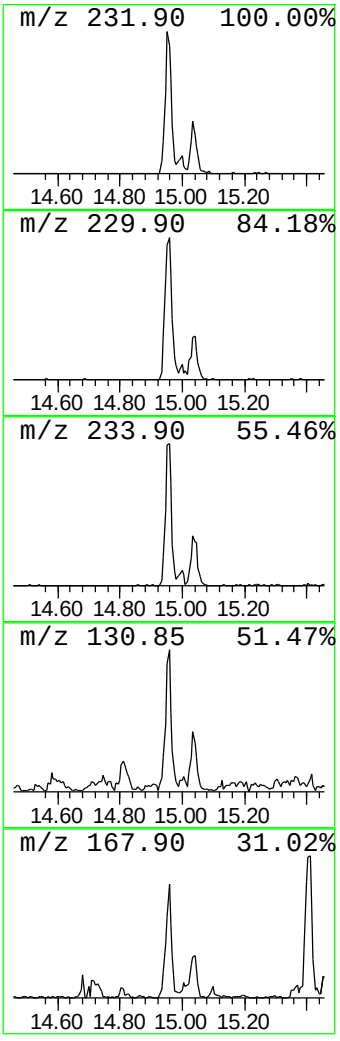
****

Peak Number 5 Phenol, 2,3,4,5-tetrachloro- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.96	2.43 ng/ul	56438	Acenaphthene-d10	14.41

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2,3,4,5-tetrachloro-	230	C6H2Cl4O	004901-51-3	98
2	Phenol, 2,3,4,6-tetrachloro-	230	C6H2Cl4O	000058-90-2	98
3	Phenol, 2,3,5,6-tetrachloro-	230	C6H2Cl4O	000935-95-5	96
4	Phenol, 2,3,4,6-tetrachloro-	230	C6H2Cl4O	000058-90-2	96
5	Phenol, 2,3,5,6-tetrachloro-	230	C6H2Cl4O	000935-95-5	94





Library Search Compound Report

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Data File : BP004771.D
Acq On : 17 Feb 2021 14:04
Operator : CG/JU
Sample : M1407-02
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DBGJ3

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP020821MA.M
Quant Title : SVOA CALIBRATION

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

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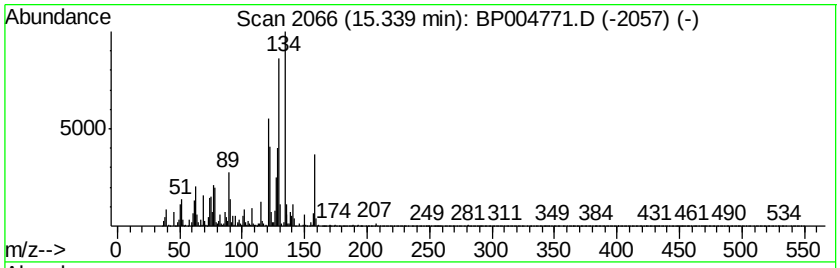
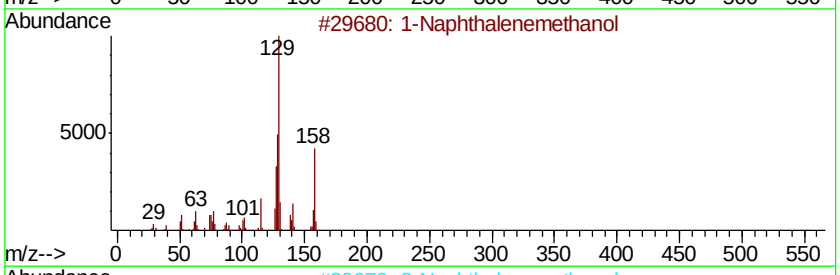
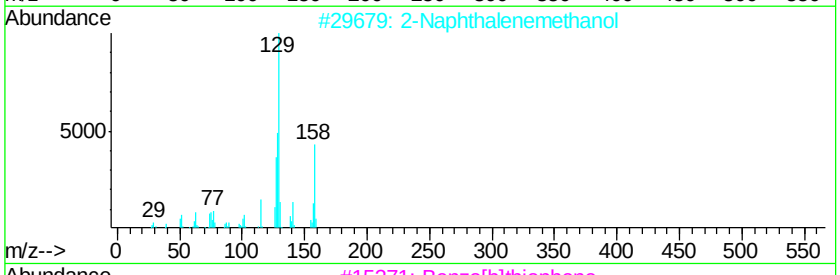
Peak Number 6 1-Naphthalenemethanol Concentration Rank 1

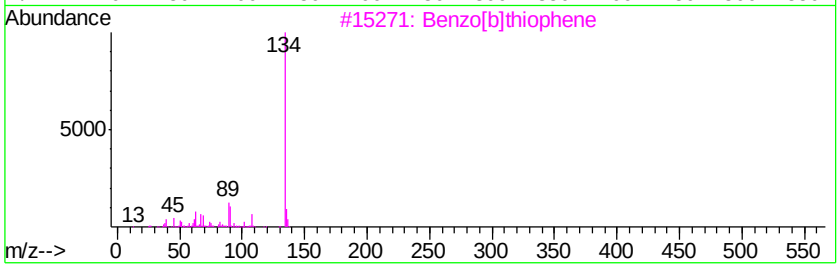
R.T.	EstConc	Area	Relative to ISTD	R.T.
15.34	6.09 ng/ul	141180	Acenaphthene-d10	14.41

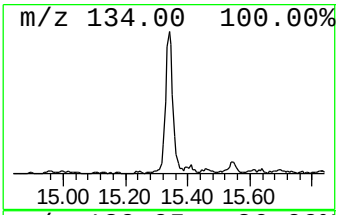
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Naphthalenemethanol	158	C11H10O	004780-79-4	84
2	2-Naphthalenemethanol	158	C11H10O	001592-38-7	66
3	Benzo[b]thiophene	134	C8H6S	000095-15-8	42
4	Benzene, 1,3-hexadienyl-	158	C12H14	041635-77-2	38
5	1-Naphthalenemethanol	158	C11H10O	004780-79-4	38

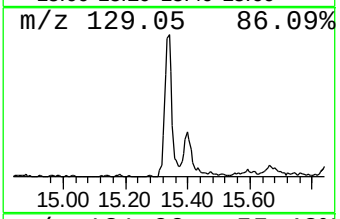


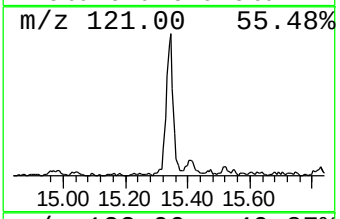


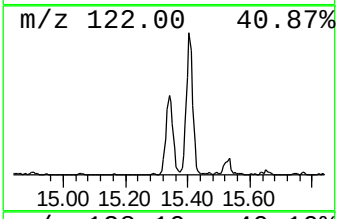


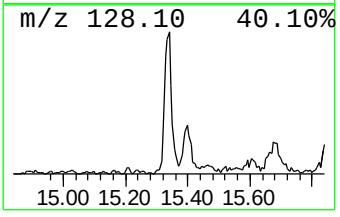












Library Search Compound Report

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Data File : BP004771.D
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Operator : CG/JU
Sample : M1407-02
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DBGJ3

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP020821MA.M
Quant Title : SVOA CALIBRATION

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

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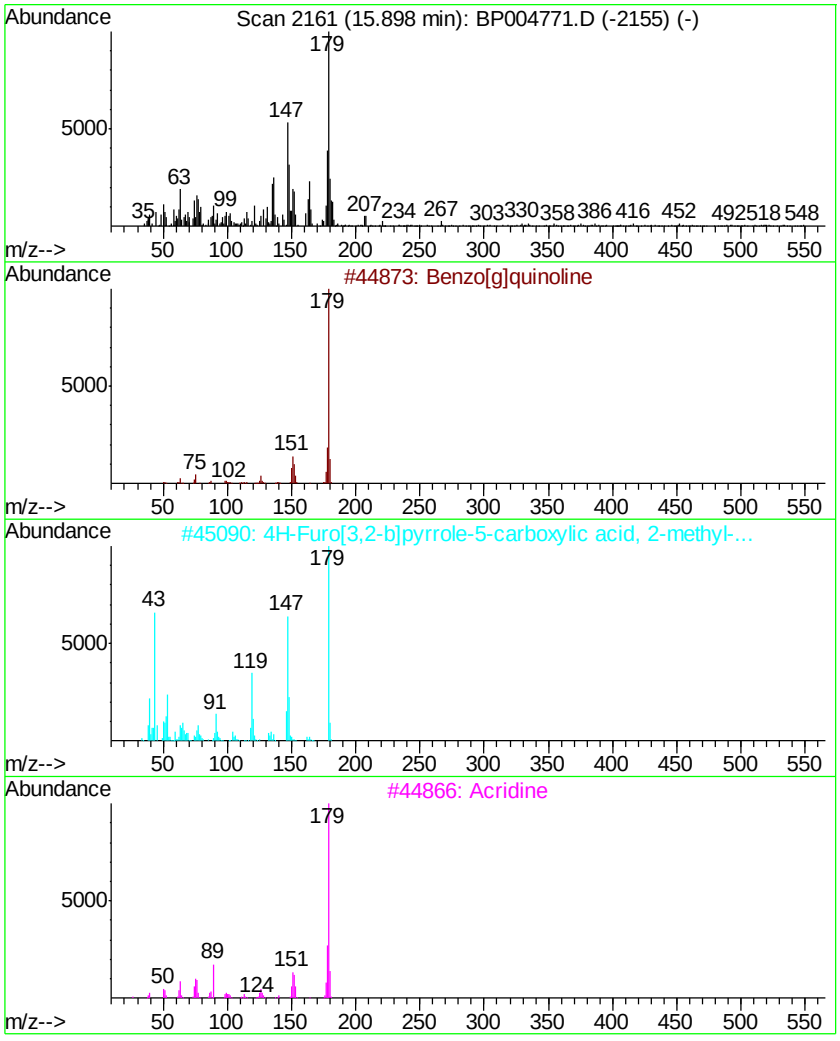
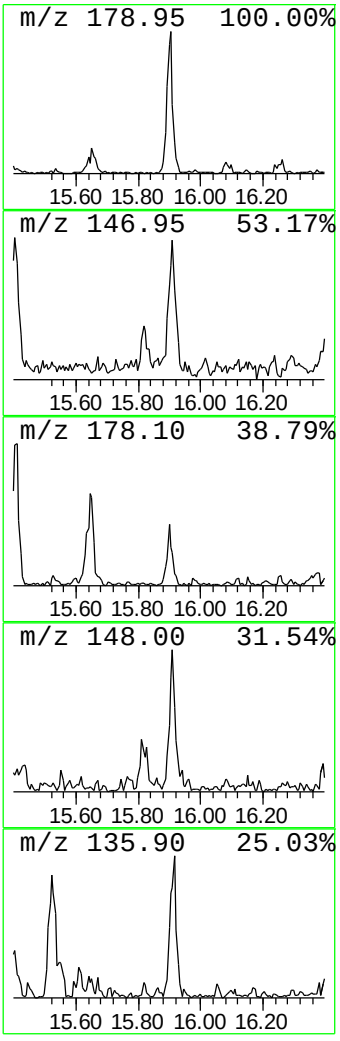
Peak Number 7 unknown-01 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.90	2.08 ng/ul	57231	Phenanthrene-d10	17.16

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[g]quinoline	179	C13H9N	000260-36-6	47
2	4H-Furo[3,2-b]pyrrole-5-carboxyl...	179	C9H9NO3	1000318-89-7	38
3	Acridine	179	C13H9N	000260-94-6	38
4	Cloxyquin	179	C9H6ClNO	000130-16-5	38
5	p-Phenylbenzonitrile	179	C13H9N	002920-38-9	38





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Instrument : BNA_P
ClientSampleId : DBGJ3

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Quant Title : SVOA CALIBRATION

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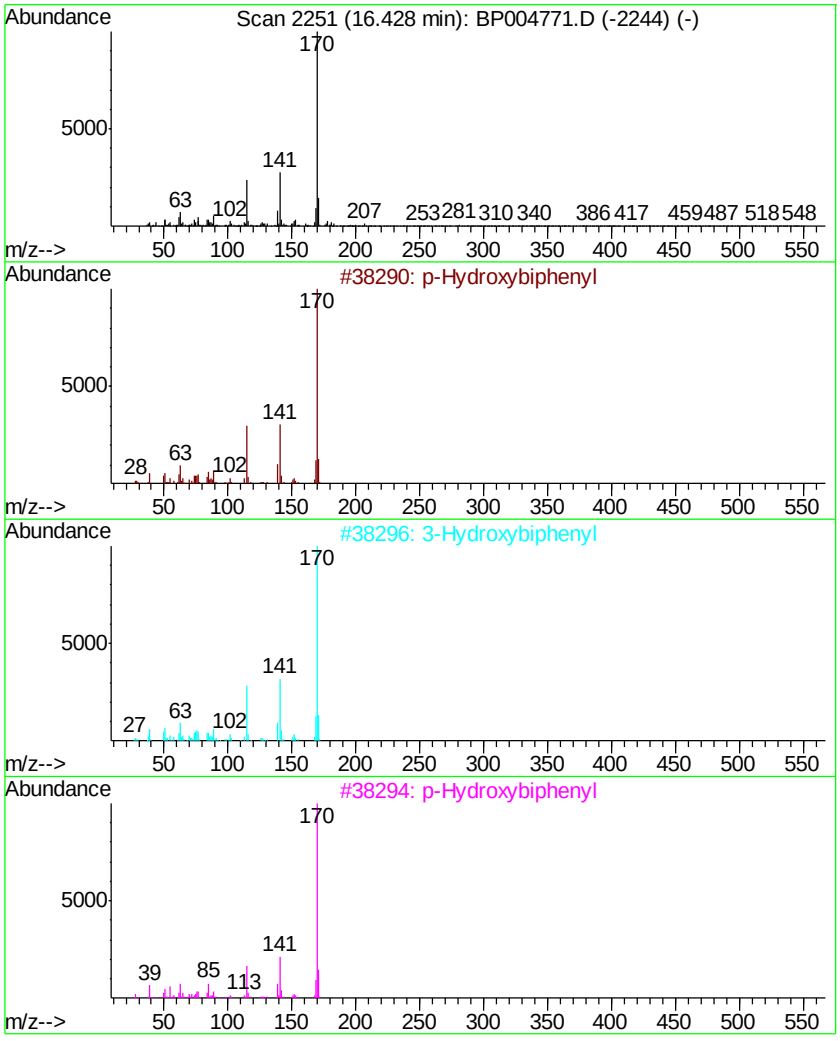
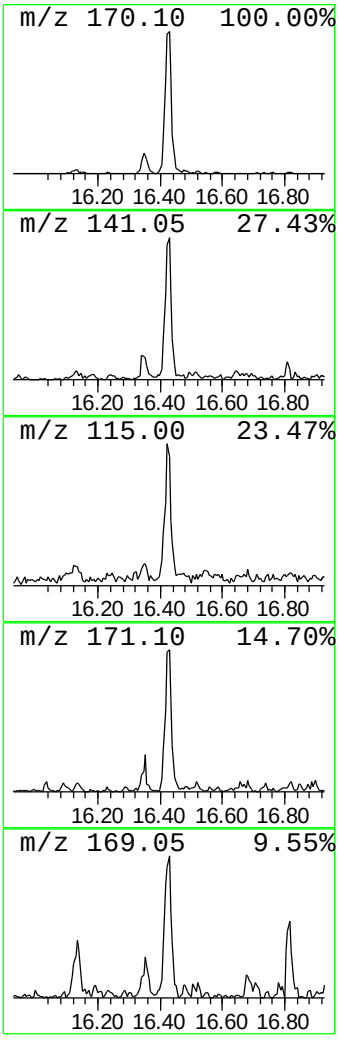
Peak Number 8 p-Hydroxybiphenyl Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.43	2.48 ng/ul	68266	Phenanthrene-d10	17.16

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	p-Hydroxybiphenyl	170	C12H10O	000092-69-3	96
2	3-Hydroxybiphenyl	170	C12H10O	000580-51-8	94
3	p-Hydroxybiphenyl	170	C12H10O	000092-69-3	93
4	o-Hydroxybiphenyl	170	C12H10O	000090-43-7	87
5	o-Hydroxybiphenyl	170	C12H10O	000090-43-7	80





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Data File : BP004771.D
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Operator : CG/JU
Sample : M1407-02
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : DBGJ3

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP020821MA.M
Quant Title : SVOA CALIBRATION

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

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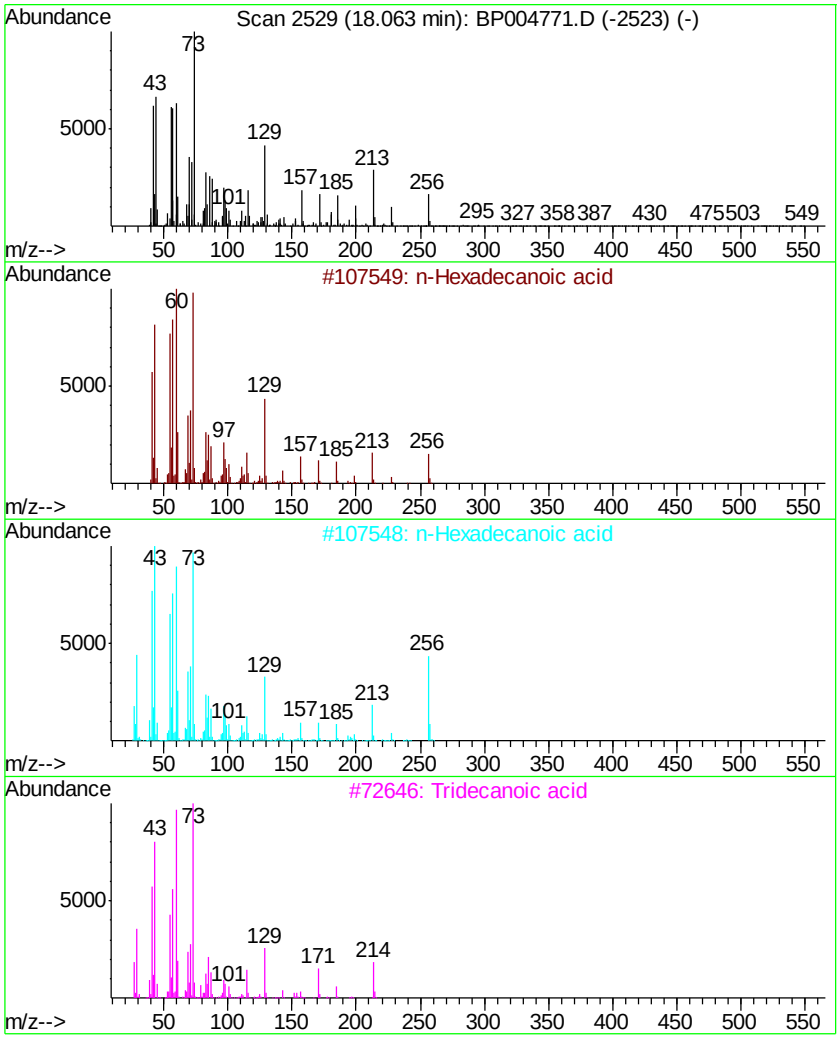
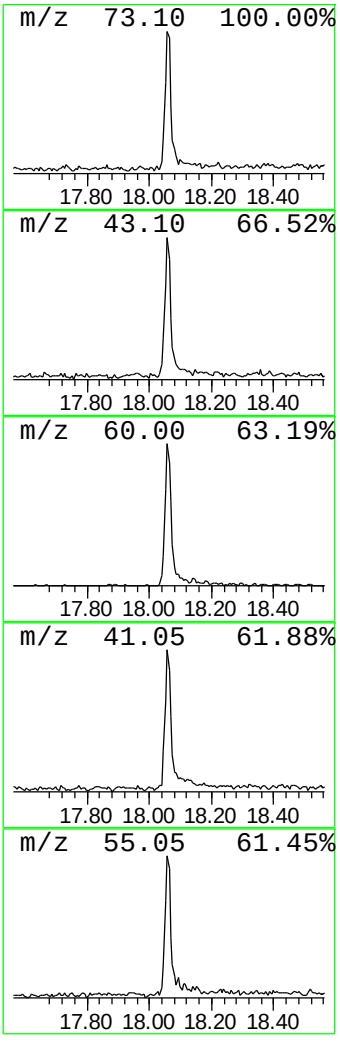
Peak Number 9 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.06	5.23 ng/ul	143869	Phenanthrene-d10	17.16

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93
3	Tridecanoic acid	214	C13H26O2	000638-53-9	89
4	Dodecanoic acid	200	C12H24O2	000143-07-7	70
5	Pentadecanoic acid	242	C15H30O2	001002-84-2	64





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Instrument : BNA_P
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Quant Title : SVOA CALIBRATION

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TIC Integration Parameters: LSCINT.P

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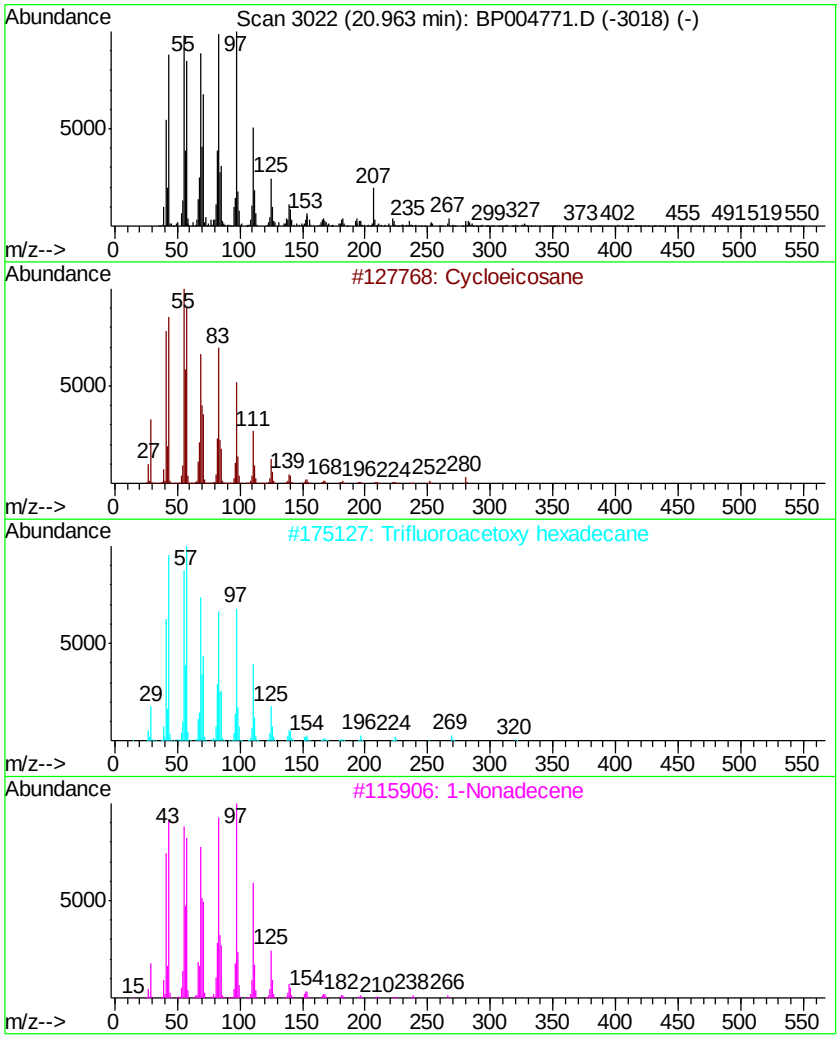
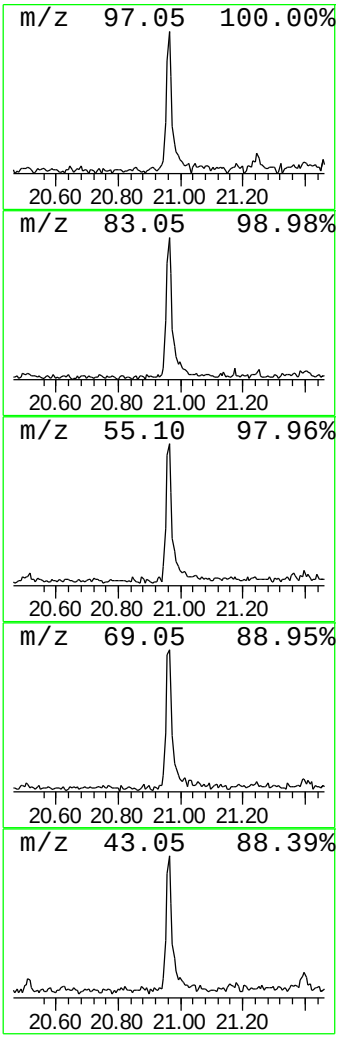
Peak Number 10 (DEL) Alkane: Cyclic20.96 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.96	2.87 ng/ul	100620	Chrysene-d12	21.25

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cycloeicosane	280	C20H40	000296-56-0	95
2	Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	91
3	1-Nonadecene	266	C19H38	018435-45-5	90
4	Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	90
5	Heptacosyl acetate	438	C29H58O2	1000351-78-2	89





Tentatively Identified Compound (LSC) summary

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Misc :
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Instrument : BNA_P
ClientSampleId : DBGJ3

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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
Benzonitrile, 3-m...	8.60	4.6	ng/ul	68255	1	7.76	297150	20.0
Phenol, 2,6-dimet...	9.18	2.5	ng/ul	46569	2	10.56	366461	20.0
Phenol, 2-(1-meth...	10.92	2.2	ng/ul	40629	2	10.56	366461	20.0
Phenol, 2,4,6-tri...	11.59	2.1	ng/ul	38765	2	10.56	366461	20.0
Phenol, 2,3,4,5-t...	14.96	2.4	ng/ul	56438	3	14.41	463776	20.0
1-Naphthalenemeth...	15.34	6.1	ng/ul	141180	3	14.41	463776	20.0
unknown-01	15.90	2.1	ng/ul	57231	4	17.16	549832	20.0
p-Hydroxybiphenyl	16.43	2.5	ng/ul	68266	4	17.16	549832	20.0
n-Hexadecanoic acid	18.06	5.2	ng/ul	143869	4	17.16	549832	20.0
(DEL) Alkane: Cyc...	20.96	2.9	ng/ul	100620	5	21.25	699985	20.0