LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG112120\
Data File : BG047410.D
Acq On : 22 Nov 2020 1:14
Operator : CG/JU
Sample : L4711-04
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0AZ8

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_G\METHODS\SOM-EPA-BG111920MA.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	3.616	44	47	52	rVB2	8040	10333	0.87%	0.073%
2	3.898	91	95	99	rVV2	8345	13557	1.15%	0.095%
3	3.951	103	104	108	rVB2	3153	3085	0.26%	0.022%
4	3.992	108	111	113	rVB3	2726	3208	0.27%	0.023%
5	4.015	113	115	119	rBV2	3445	4137	0.35%	0.029%
6	4.292	160	162	165	rBV3	2737	3063	0.26%	0.022%
7	4.386	176	178	184	rVB2	2518	4053	0.34%	0.029%
8	4.603	212	215	219	rVB4	5435	8272	0.70%	0.058%
9	4.715	231	234	238	rBV3	2474	3711	0.31%	0.026%
10	5.232	319	322	324	rVV2	2637	2772	0.23%	0.020%
11	5.337	338	340	345	rVB2	2829	3386	0.29%	0.024%
12	5.907	436	437	442	rBV2	2169	2757	0.23%	0.019%
13	6.248	492	495	497	rBV2	2802	3129	0.26%	0.022%
14	6.771	580	584	587	rBV	1963	2971	0.25%	0.021%
15	7.223	657	661	665	rBV3	7830	10634	0.90%	0.075%
16	7.411	685	693	702	rBV	211161	390723	33.06%	2.751%
17	7.588	716	723	729	rBV	122866	205205	17.36%	1.445%
18	7.799	752	759	765	rVV2	203441	367338	31.08%	2.586%
19	7.852	765	768	775	rVB2	23711	37441	3.17%	0.264%
20	8.175	820	823	826	rBV	1867	2852	0.24%	0.020%
21	8.275	831	840	847	rBV	162906	283874	24.02%	1.999%
22	8.968	951	958	965	rBV	259661	456711	38.65%	3.216%
23	9.262	1002	1008	1009	rBV	2039	2731	0.23%	0.019%
24	9.286	1009	1012	1013	rBV	3792	3720	0.31%	0.026%
25	9.444	1030	1039	1048	rBV	194904	367907	31.13%	2.590%
26	9.856	1105	1109	1112	rVB	5369	8214	0.70%	0.058%
27	10.173	1155	1163	1172	rVB2	196294	352872	29.86%	2.485%
28	10.367	1192	1196	1197	rBV	3164	3613	0.31%	0.025%
29	10.713	1247	1255	1261	rVV	279554	531801	45.00%	3.744%
30	10.778	1264	1266	1269	rVV	3242	3110	0.26%	0.022%
31	10.819	1269	1273	1277	rVV	5942	8019	0.68%	0.056%
32	11.101	1314	1321	1330	rBV	205484	381002	32.24%	2.683%
33	11.225	1333	1342	1350	rBV2	229491	445956	37.74%	3.140%
34	11.565	1398	1400	1405	rVV2	2912	3436	0.29%	0.024%

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Misc :
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Instrument : BNA_G
ClientSampleId : C0AZ8

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_G\METHODS\SOM-EPA-BG111920MA.M
Title : SVOA CALIBRATION

35	11.612	1405	1408	1413	rVV	2308	3807	0.32%	0.027%
36	12.018	1475	1477	1481	rVB2	2670	2808	0.24%	0.020%
37	12.094	1485	1490	1495	rVB4	36499	61958	5.24%	0.436%
38	12.394	1537	1541	1542	rBV2	3808	4603	0.39%	0.032%
39	12.411	1542	1544	1545	rVV2	3670	3336	0.28%	0.023%
40	12.429	1545	1547	1553	rVB3	3832	4501	0.38%	0.032%
41	12.664	1584	1587	1592	rVB4	4543	5548	0.47%	0.039%
42	12.958	1635	1637	1641	rVB2	2196	2684	0.23%	0.019%
43	13.199	1674	1678	1680	rBV	2273	3540	0.30%	0.025%
44	13.252	1685	1687	1694	rVB	4944	5821	0.49%	0.041%
45	13.334	1698	1701	1706	rVB3	2012	3100	0.26%	0.022%
46	13.722	1761	1767	1770	rVB2	4578	7812	0.66%	0.055%
47	13.769	1770	1775	1781	rBV2	56190	85274	7.22%	0.600%
48	13.992	1810	1813	1818	rBV2	3009	4209	0.36%	0.030%
49	14.062	1820	1825	1828	rBV2	2044	3240	0.27%	0.023%
50	14.280	1855	1862	1866	rBV	486351	706287	59.76%	4.973%
51	14.327	1866	1870	1881	rVB	194549	286334	24.23%	2.016%
52	14.503	1898	1900	1901	rBV	3399	3237	0.27%	0.023%
53	14.515	1901	1902	1906	rVV	4822	4710	0.40%	0.033%
54	14.585	1906	1914	1922	rVV	501128	792783	67.08%	5.582%
55	14.762	1942	1944	1946	rBV2	4115	2934	0.25%	0.021%
56	14.832	1950	1956	1958	rBV	2127	2782	0.24%	0.020%
57	14.885	1958	1965	1971	rBV2	322395	519519	43.96%	3.658%
58	14.938	1971	1974	1980	rVB3	11579	17375	1.47%	0.122%
59	15.050	1986	1993	2001	rBV	220271	386120	32.67%	2.719%
60	15.226	2015	2023	2024	rBV6	19718	35341	2.99%	0.249%
61	15.255	2024	2028	2033	rVV2	95939	148243	12.54%	1.044%
62	15.414	2052	2055	2057	rVB2	3804	2776	0.23%	0.020%
63	15.484	2065	2067	2069	rBV	3025	2887	0.24%	0.020%
64	15.655	2092	2096	2100	rBV4	5824	7690	0.65%	0.054%
65	15.807	2117	2122	2126	rBV4	22687	35951	3.04%	0.253%
66	15.872	2126	2133	2141	rVB	613437	947527	80.18%	6.671%
67	15.966	2145	2149	2157	rVB	125305	203406	17.21%	1.432%
68	16.107	2170	2173	2182	rVB3	10439	17509	1.48%	0.123%
69	16.266	2197	2200	2202	rBV2	3360	4722	0.40%	0.033%
70	16.360	2214	2216	2217	rBV2	5593	3530	0.30%	0.025%
71	16.436	2224	2229	2234	rBV5	15295	25119	2.13%	0.177%

LSC Area Percent Report

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Sample : L4711-04
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0AZ8

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
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Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG111920MA.M
Title : SVOA CALIBRATION

72	16.595	2254	2256	2260	rBV4	3348	3249	0.27%	0.023%
73	16.648	2263	2265	2271	rVB3	7698	8628	0.73%	0.061%
74	16.689	2271	2272	2276	rBV3	2597	3409	0.29%	0.024%
75	16.783	2286	2288	2290	rBV3	4545	2850	0.24%	0.020%
76	16.812	2290	2293	2295	rVV3	5143	4850	0.41%	0.034%
77	16.847	2296	2299	2305	rVB4	9237	14586	1.23%	0.103%
78	16.983	2318	2322	2323	rBV2	5494	6088	0.52%	0.043%
79	17.077	2335	2338	2342	rBV6	4763	8414	0.71%	0.059%
80	17.270	2369	2371	2374	rBV3	4487	4680	0.40%	0.033%
81	17.353	2381	2385	2391	rVB8	5167	8664	0.73%	0.061%
82	17.623	2424	2431	2436	rBV	385793	622792	52.70%	4.385%
83	17.658	2436	2437	2441	rVV	32684	39177	3.32%	0.276%
84	17.717	2441	2447	2454	rVB	607485	952074	80.56%	6.703%
85	17.817	2462	2464	2466	rVB3	4166	3270	0.28%	0.023%
86	18.193	2525	2528	2532	rBV5	6386	8411	0.71%	0.059%
87	18.410	2561	2565	2571	rVB4	52470	68069	5.76%	0.479%
88	18.481	2573	2577	2584	rBV3	85681	135159	11.44%	0.952%
89	18.540	2584	2587	2592	rVB7	11717	20356	1.72%	0.143%
90	18.681	2608	2611	2615	rVB6	9156	14515	1.23%	0.102%
91	18.945	2654	2656	2659	rBV4	7534	9346	0.79%	0.066%
92	19.092	2679	2681	2685	rBV5	7190	8158	0.69%	0.057%
93	19.251	2706	2708	2712	rBV2	26091	31889	2.70%	0.225%
94	19.650	2772	2776	2782	rVB	85713	121343	10.27%	0.854%
95	19.985	2827	2833	2843	rVB	726892	1181800	100.00%	8.321%
96	21.513	3088	3093	3099	rBV2	241542	363652	30.77%	2.560%
97	21.900	3153	3159	3165	rBV2	325044	563357	47.67%	3.967%
98	24.403	3580	3585	3595	rVB2	221229	484437	40.99%	3.411%
99	25.073	3692	3699	3708	rVB	296610	809513	68.50%	5.700%
100	25.308	3733	3739	3748	rVB3	157785	405341	34.30%	2.854%

Sum of corrected areas: 14202693

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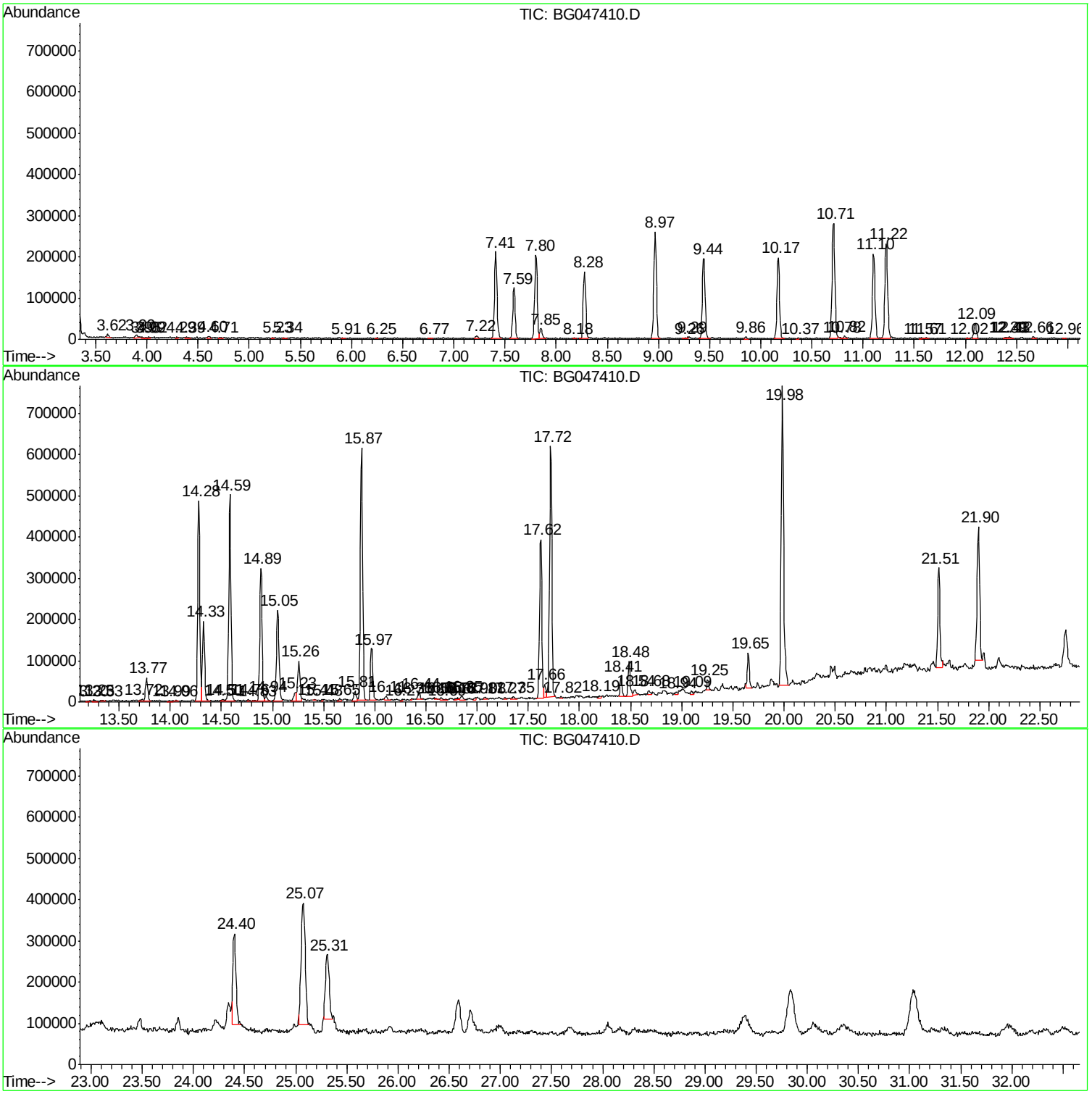
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AZ8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG111920MA.M
Quant Title : SVOA CALIBRATION

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



Library Search Compound Report

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Instrument : BNA_G
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Quant Title : SVOA CALIBRATION

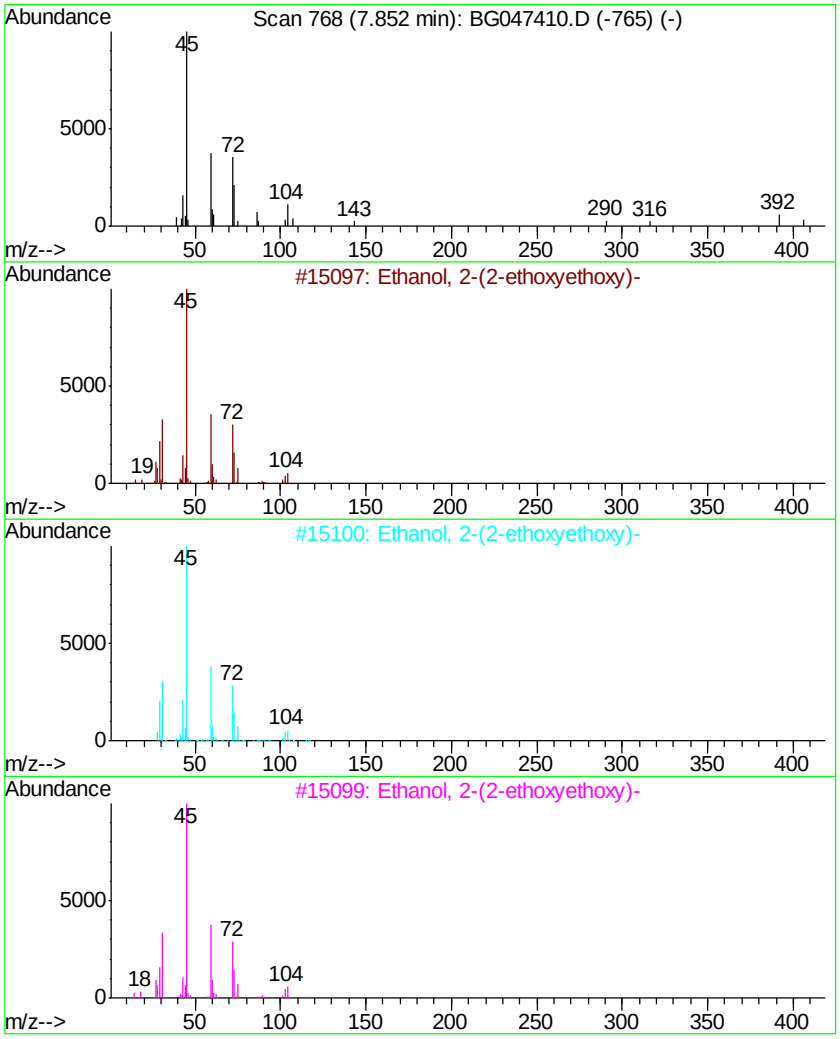
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

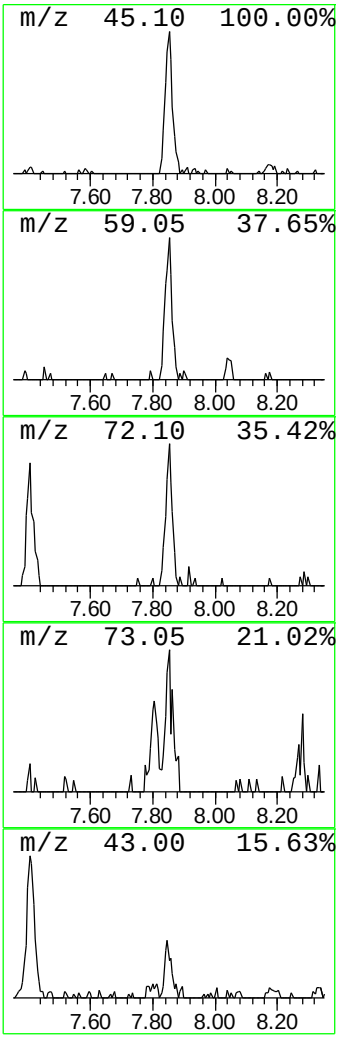
****

Peak Number 1 Ethanol, 2-(2-ethoxyethoxy)- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.85	2.64 ng/ul	37441	1,4-Dichlorobenzene-d4	8.28

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	64
2	Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	56
3	Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	50
4	Diethyl carbitol	162	C8H18O3	000112-36-7	50
5	Ethyl 2-(2-(2-ethoxyethoxy)ethox...	220	C10H20O5	1000366-77-8	45





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG112120\
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Sample : L4711-04
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0AZ8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG111920MA.M
Quant Title : SVOA CALIBRATION

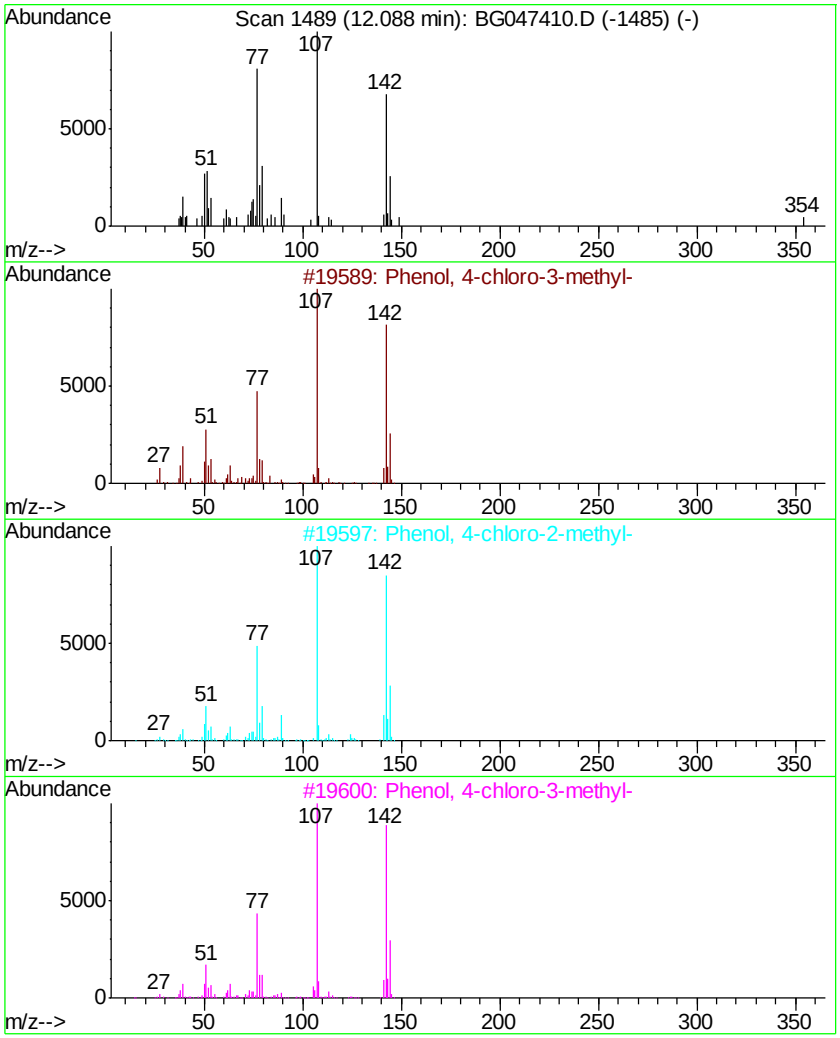
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

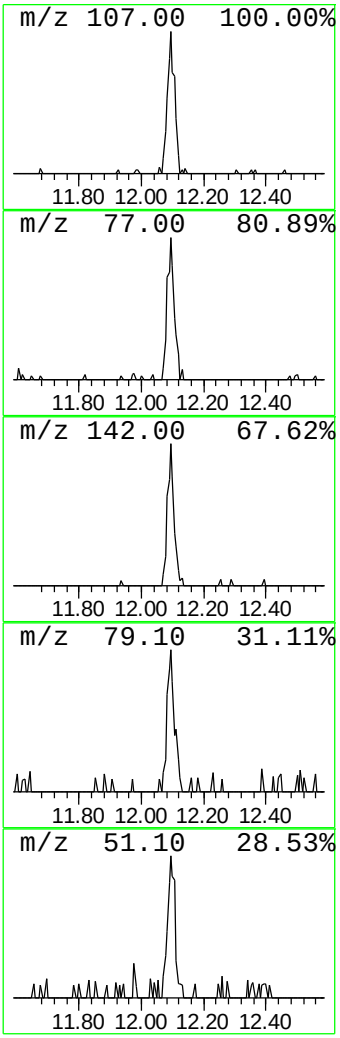
****

Peak Number 2 Phenol, 4-chloro-3-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.09	3.25 ng/ul	61958	Naphthalene-d8	11.10

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 4-chloro-3-methyl-	142	C7H7ClO	000059-50-7	93
2	Phenol, 4-chloro-2-methyl-	142	C7H7ClO	001570-64-5	86
3	Phenol, 4-chloro-3-methyl-	142	C7H7ClO	000059-50-7	76
4	Phenol, 4-chloro-3-methyl-	142	C7H7ClO	000059-50-7	76
5	3-Chlorobenzyl alcohol	142	C7H7ClO	000873-63-2	64





Library Search Compound Report

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG111920MA.M
Quant Title : SVOA CALIBRATION

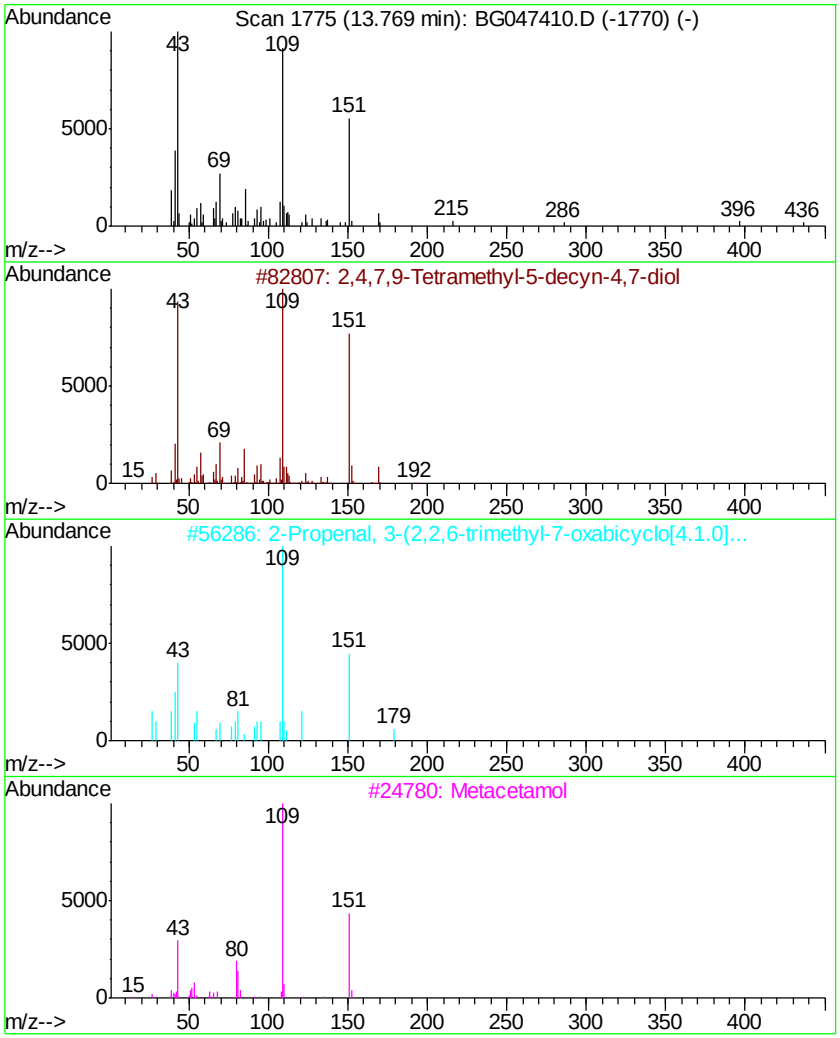
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

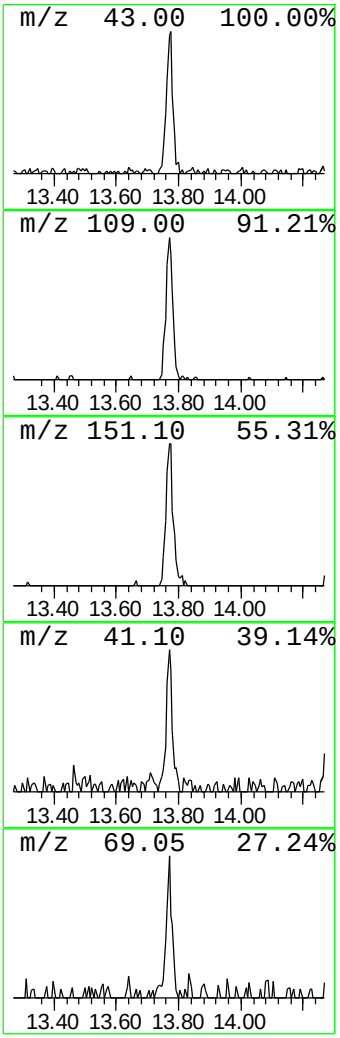
****

Peak Number 3 2,4,7,9-Tetramethyl-5-decyn... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.77	3.28 ng/ul	85274	Acenaphthene-d10	14.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2	000126-86-3	91
2	2-Propenal, 3-(2,2,6-trimethyl-7...	194	C12H18O2	055759-91-6	78
3	Metacetamol	151	C8H9NO2	000621-42-1	45
4	m-Isopropoxyaniline	151	C9H13NO	041406-00-2	45
5	Acetaminophen	151	C8H9NO2	000103-90-2	45





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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG111920MA.M
Quant Title : SVOA CALIBRATION

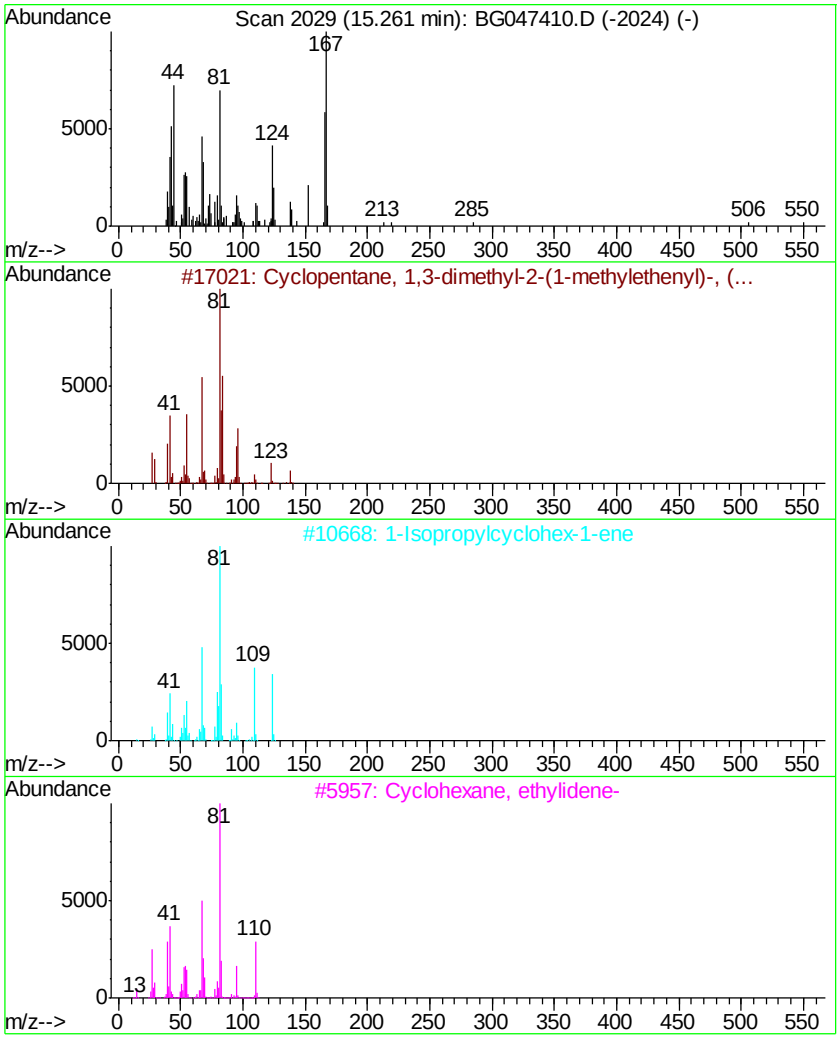
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

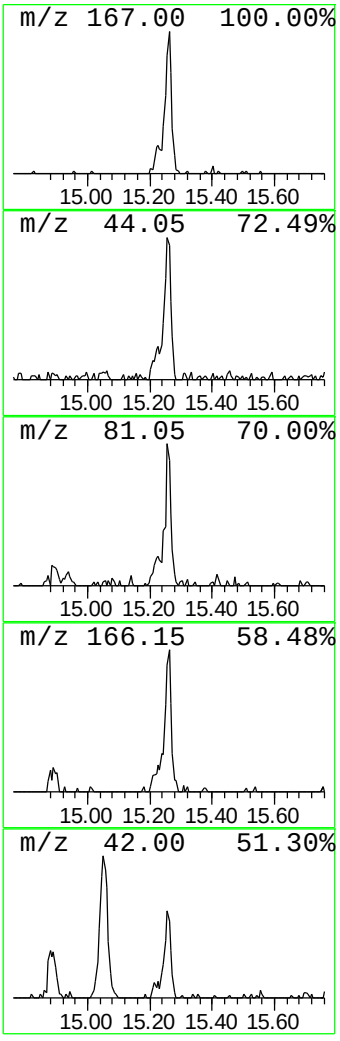
****

Peak Number 4 unknown-01 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.26	5.71 ng/ul	148243	Acenaphthene-d10	14.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, 1,3-dimethyl-2-(1-...	138	C10H18	061142-31-2	18
2	1-Isopropylcyclohex-1-ene	124	C9H16	004292-04-0	14
3	Cyclohexane, ethylidene-	110	C8H14	001003-64-1	14
4	Cyclohexane, ethylidene-	110	C8H14	001003-64-1	14
5	1-Cyano-4-methylnaphthalene	167	C12H9N	036062-93-8	11





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ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0AZ8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG111920MA.M
Quant Title : SVOA CALIBRATION

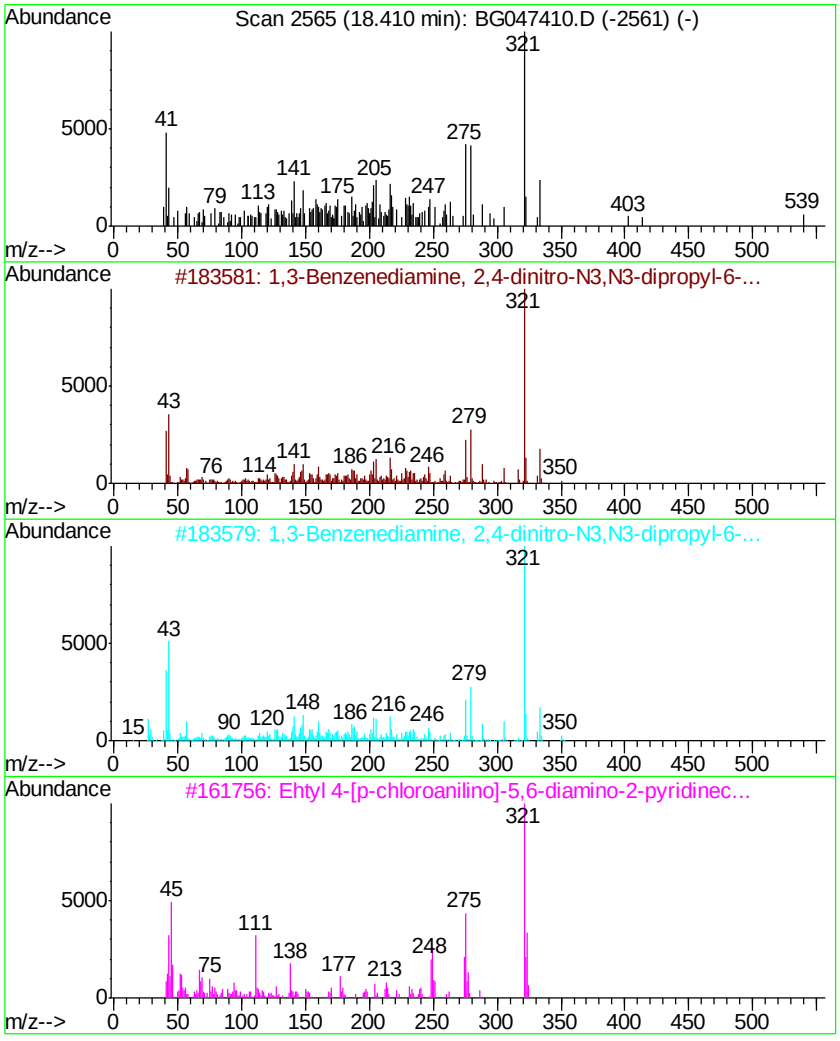
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

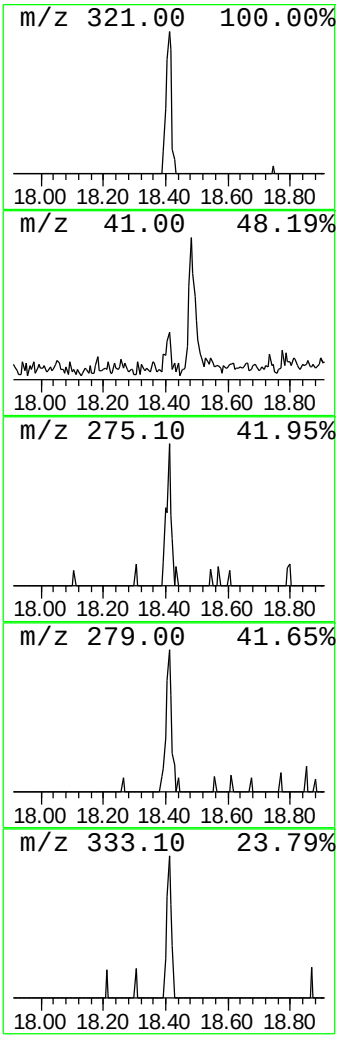
****

Peak Number 5 1,3-Benzenediamine, 2,4-din... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.41	2.19 ng/ul	68069	Phenanthrene-d10	17.62

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Benzenediamine, 2,4-dinitro-...	350	C13H17F3N4O4	029091-21-2	86
2	1,3-Benzenediamine, 2,4-dinitro-...	350	C13H17F3N4O4	029091-21-2	74
3	Ehtyl 4-[p-chloroanilino]-5,6-di...	321	C14H16ClN5O2	019270-38-3	49
4	3-(2-Hydroxy-5-nitrophenyl)-2-(6...	322	C17H14N4O3	1000311-98-2	38
5	5-Allyl-2-(4-iodo-phenyl)-pyridine	321	C14H12IN	302602-55-7	30





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG112120\
Data File : BG047410.D
Acq On : 22 Nov 2020 1:14
Operator : CG/JU
Sample : L4711-04
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0AZ8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG111920MA.M
Quant Title : SVOA CALIBRATION

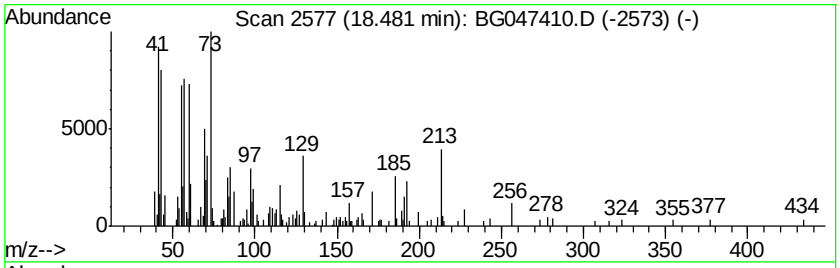
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

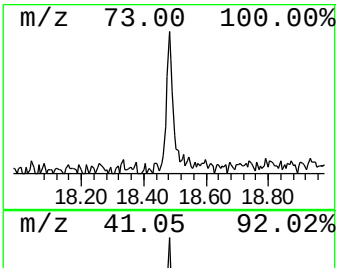
****

Peak Number 6 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.48	4.34 ng/ul	135159	Phenanthrene-d10	17.62

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	87
3	Tridecanoic acid	214	C13H26O2	000638-53-9	70
4	Tridecanoic acid	214	C13H26O2	000638-53-9	70
5	Octadecanoic acid	284	C18H36O2	000057-11-4	64





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Data File : BG047410.D
Acq On : 22 Nov 2020 1:14
Operator : CG/JU
Sample : L4711-04
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0AZ8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG111920MA.M
Quant Title : SVOA CALIBRATION

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

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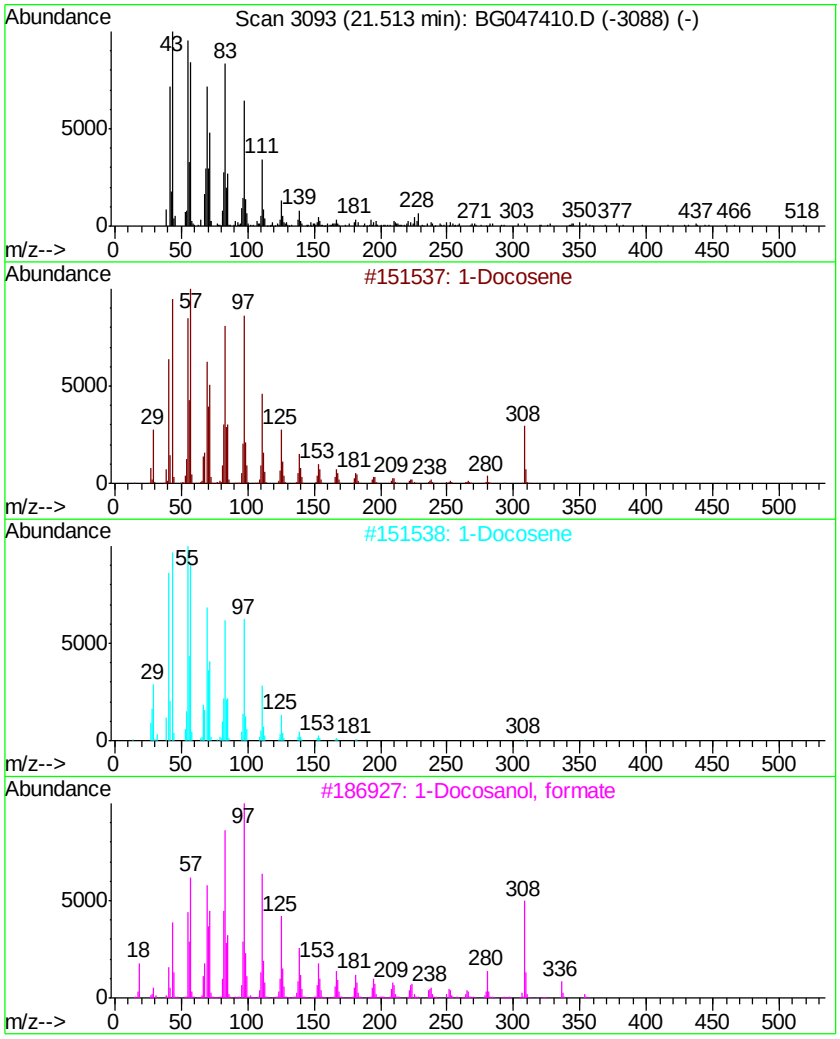
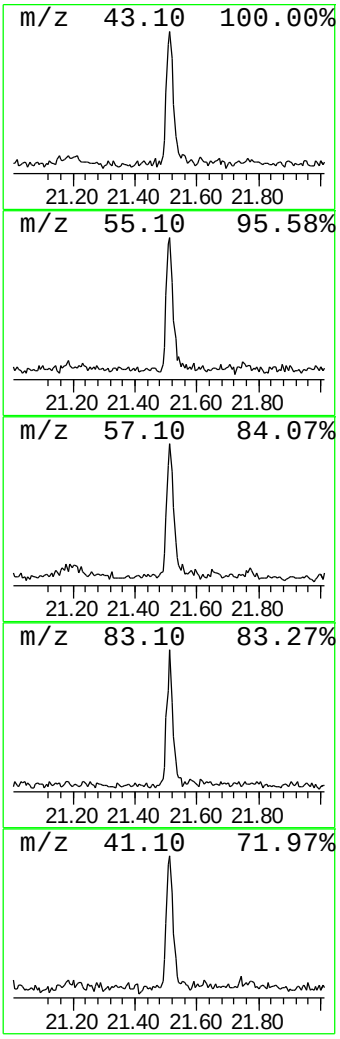
Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.51	12.91 ng/ul	363652	Chrysene-d12	21.90

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Docosene	308	C22H44	001599-67-3	97
2	1-Docosene	308	C22H44	001599-67-3	96
3	1-Docosanol, formate	354	C23H46O2	015155-62-1	94
4	Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	93
5	1-Heptacosanol	396	C27H56O	002004-39-9	91





Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG112120\
Data File : BG047410.D
Acq On : 22 Nov 2020 1:14
Operator : CG/JU
Sample : L4711-04
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0AZ8

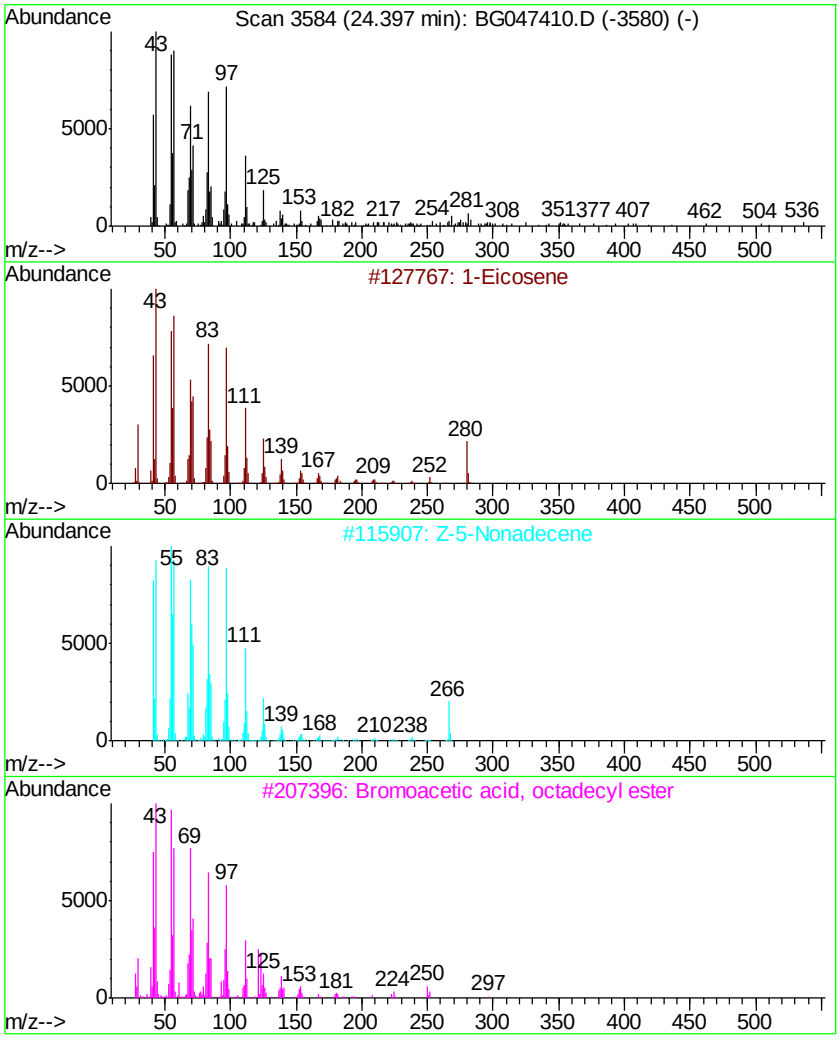
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG111920MA.M
Quant Title : SVOA CALIBRATION

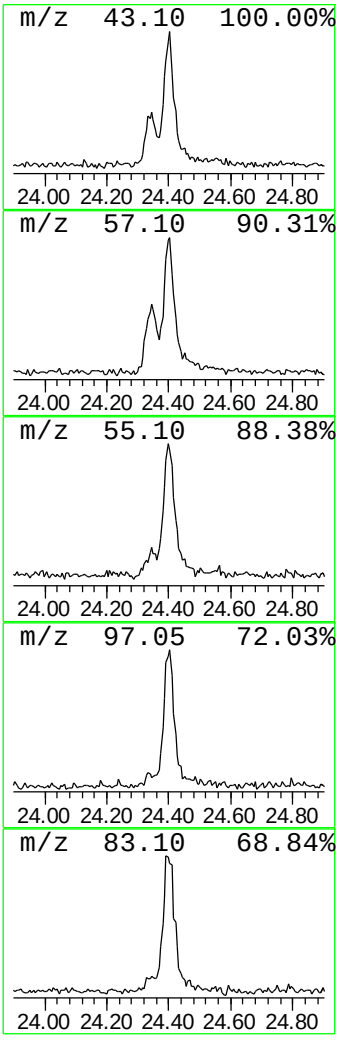
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.40	23.90 ng/ul	484437	Perylene-d12	25.31

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Eicosene	280	C20H40	003452-07-1	97
2	Z-5-Nonadecene	266	C19H38	1000131-11-8	92
3	Bromoacetic acid, octadecyl ester	390	C20H39BrO2	018992-03-5	90
4	1-Eicosene	280	C20H40	003452-07-1	90
5	13-Tetradecen-1-ol acetate	254	C16H30O2	056221-91-1	90





Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG112120\
Data File : BG047410.D
Acq On : 22 Nov 2020 1:14
Operator : CG/JU
Sample : L4711-04
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : C0AZ8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG111920MA.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
Ethanol, 2-(2-eth...	7.85	2.6	ng/ul	37441	1	8.28	283874	20.0
Phenol, 4-chloro-...	12.09	3.3	ng/ul	61958	2	11.10	381002	20.0
2,4,7,9-Tetrameth...	13.77	3.3	ng/ul	85274	3	14.89	519519	20.0
unknown-01	15.26	5.7	ng/ul	148243	3	14.89	519519	20.0
1,3-Benzenediamin...	18.41	2.2	ng/ul	68069	4	17.62	622792	20.0
n-Hexadecanoic acid	18.48	4.3	ng/ul	135159	4	17.62	622792	20.0
(DEL) Alkane: Str...	21.51	12.9	ng/ul	363652	5	21.90	563357	20.0
(DEL) Alkane: Str...	24.40	23.9	ng/ul	484437	6	25.31	405341	20.0