LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM100319\
Data File : BM022928.D
Acq On : 03 Oct 2019 19:49
Operator : JU
Sample : K4983-17
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFDK7

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-BM092719MA.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	3.158	26	29	44	rVB	35043	63396	1.18%	0.093%
2	3.269	44	48	49	rBV	76951	86955	1.62%	0.128%
3	3.299	49	53	68	rVB	1809618	2149251	40.01%	3.153%
4	3.663	111	115	121	rVV	50140	67745	1.26%	0.099%
5	3.728	121	126	131	rVB	179831	225364	4.20%	0.331%
6	3.981	165	169	178	rVB2	56225	86535	1.61%	0.127%
7	4.263	212	217	225	rBV	392882	493864	9.19%	0.724%
8	4.334	225	229	237	rVB2	50513	74543	1.39%	0.109%
9	4.440	242	247	255	rBV	562835	721299	13.43%	1.058%
10	4.510	255	259	268	rVB	28137	44029	0.82%	0.065%
11	4.893	315	324	332	rBV	658957	871127	16.22%	1.278%
12	5.757	465	471	477	rBV	92582	131614	2.45%	0.193%
13	6.787	639	646	647	rBV	35167	50922	0.95%	0.075%
14	6.963	668	676	680	rBV	220344	436105	8.12%	0.640%
15	7.128	697	704	711	rBV	582055	924537	17.21%	1.356%
16	7.193	711	715	720	rVB	43822	68517	1.28%	0.101%
17	7.328	727	738	748	rBV	691679	1096970	20.42%	1.609%
18	7.445	752	758	769	rVB	127741	203940	3.80%	0.299%
19	7.793	808	817	826	rBV	944005	1516872	28.24%	2.225%
20	7.910	832	837	846	rVB2	68963	126173	2.35%	0.185%
21	8.169	873	881	888	rVB2	103148	181146	3.37%	0.266%
22	8.369	906	915	921	rVV5	17460	45036	0.84%	0.066%
23	8.440	921	927	931	rVV2	22888	41421	0.77%	0.061%
24	8.498	931	937	948	rVV	600307	960160	17.87%	1.409%
25	8.604	950	955	962	rVV2	20403	45547	0.85%	0.067%
26	8.898	1000	1005	1008	rBV	84593	134042	2.50%	0.197%
27	8.945	1008	1013	1025	rVV	735958	1304255	24.28%	1.913%
28	9.304	1068	1074	1080	rVB6	21714	43968	0.82%	0.064%
29	9.422	1085	1094	1099	rVB2	54291	98635	1.84%	0.145%
30	9.481	1099	1104	1115	rVB	64551	106745	1.99%	0.157%
31	9.669	1129	1136	1144	rBV	626637	1053589	19.61%	1.546%
32	9.839	1160	1165	1172	rVB	51775	88059	1.64%	0.129%
33	9.916	1172	1178	1184	rBV	66705	103502	1.93%	0.152%
34	9.992	1184	1191	1199	rVB3	130692	261665	4.87%	0.384%

LSC Area Percent Report

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Acq On : 03 Oct 2019 19:49
Operator : JU
Sample : K4983-17
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFDK7

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-BM092719MA.M
Title : SVOA CALIBRATION

35	10.134	1209	1215	1219	rVV4	27138	56658	1.05%	0.083%
36	10.210	1219	1228	1235	rVB	1019958	1716341	31.95%	2.518%
37	10.586	1283	1292	1297	rBV	1390122	2325878	43.29%	3.412%
38	10.634	1297	1300	1307	rVB	103795	174680	3.25%	0.256%
39	10.716	1307	1314	1326	rBV	748736	1336891	24.89%	1.961%
40	11.298	1408	1413	1418	rBV	36445	53109	0.99%	0.078%
41	11.610	1461	1466	1476	rVV10	16059	51273	0.95%	0.075%
42	11.780	1486	1495	1502	rVB7	31600	62687	1.17%	0.092%
43	11.928	1512	1520	1525	rBV	127346	211546	3.94%	0.310%
44	12.139	1547	1556	1558	rBV6	19494	45559	0.85%	0.067%
45	12.245	1569	1574	1580	rVB	195258	312018	5.81%	0.458%
46	12.357	1588	1593	1602	rBV5	33365	68486	1.27%	0.100%
47	12.469	1605	1612	1618	rVV	292426	467680	8.71%	0.686%
48	12.533	1619	1623	1632	rVB	24564	49878	0.93%	0.073%
49	12.627	1634	1639	1645	rBV5	31289	59817	1.11%	0.088%
50	12.716	1645	1654	1660	rBV5	49768	105200	1.96%	0.154%
51	13.139	1721	1726	1736	rBV7	21481	62147	1.16%	0.091%
52	13.222	1736	1740	1745	rBV6	23604	51146	0.95%	0.075%
53	13.275	1745	1749	1754	rVB	49873	74825	1.39%	0.110%
54	13.339	1754	1760	1766	rBV2	163044	263528	4.91%	0.387%
55	13.480	1777	1784	1791	rVB3	50114	130867	2.44%	0.192%
56	13.610	1800	1806	1810	rVB5	63490	130575	2.43%	0.192%
57	13.657	1810	1814	1818	rBV3	36245	61216	1.14%	0.090%
58	13.751	1825	1830	1835	rBV	115081	220320	4.10%	0.323%
59	13.804	1835	1839	1840	rVV	83807	113389	2.11%	0.166%
60	13.839	1840	1845	1849	rVV	1975542	2699880	50.26%	3.961%
61	13.886	1849	1853	1865	rVB	772060	1076286	20.03%	1.579%
62	13.986	1865	1870	1873	rBV	54110	84497	1.57%	0.124%
63	14.022	1873	1876	1883	rVB5	40558	80225	1.49%	0.118%
64	14.122	1884	1893	1904	rVB	2190525	3286134	61.17%	4.821%
65	14.351	1926	1932	1937	rVB8	32115	57342	1.07%	0.084%
66	14.433	1937	1946	1952	rBV	2111143	3310790	61.63%	4.857%
67	14.492	1952	1956	1963	rVB	200513	319741	5.95%	0.469%
68	14.633	1975	1980	1986	rBV	207774	321161	5.98%	0.471%
69	14.833	2009	2014	2015	rBV2	44036	71226	1.33%	0.104%
70	14.857	2015	2018	2022	rVV	60416	80658	1.50%	0.118%
71	14.904	2022	2026	2031	rVB3	45963	69711	1.30%	0.102%

LSC Area Percent Report

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Acq On : 03 Oct 2019 19:49
Operator : JU
Sample : K4983-17
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFDK7

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-BM092719MA.M
Title : SVOA CALIBRATION

72	15.069	2046	2054	2058	rBV3	55910	111672	2.08%	0.164%
73	15.204	2073	2077	2079	rBV	53899	80978	1.51%	0.119%
74	15.421	2108	2114	2119	rBV	2869466	4071309	75.79%	5.972%
75	15.474	2119	2123	2128	rVB	111366	174200	3.24%	0.256%
76	15.533	2128	2133	2141	rBV2	885262	1391489	25.90%	2.041%
77	15.657	2150	2154	2158	rBV2	82738	121869	2.27%	0.179%
78	15.933	2196	2201	2209	rVB2	178129	321266	5.98%	0.471%
79	16.098	2225	2229	2234	rBV	92267	133925	2.49%	0.196%
80	16.439	2282	2287	2291	rBV2	92190	147654	2.75%	0.217%
81	17.057	2388	2392	2396	rVB	51776	71777	1.34%	0.105%
82	17.168	2405	2411	2416	rBV2	2367835	3433931	63.92%	5.037%
83	17.210	2416	2418	2422	rVV	180252	208817	3.89%	0.306%
84	17.268	2422	2428	2438	rVB	3233441	4516999	84.08%	6.626%
85	17.557	2470	2477	2483	rBV4	82614	221405	4.12%	0.325%
86	18.068	2560	2564	2573	rBV	585362	796244	14.82%	1.168%
87	18.927	2706	2710	2720	rBV	984013	1262294	23.50%	1.852%
88	19.004	2720	2723	2729	rVB4	42852	69531	1.29%	0.102%
89	19.227	2758	2761	2763	rBV	50761	65769	1.22%	0.096%
90	19.251	2763	2765	2768	rVV	128528	141115	2.63%	0.207%
91	19.286	2769	2771	2776	rVB	94065	102567	1.91%	0.150%
92	19.351	2778	2782	2789	rBV2	122863	201163	3.74%	0.295%
93	19.562	2812	2818	2823	rBV	3924612	5372172	100.00%	7.881%
94	19.604	2823	2825	2833	rVB	269144	315083	5.87%	0.462%
95	20.080	2902	2906	2909	rBV2	168851	197619	3.68%	0.290%
96	20.115	2909	2912	2916	rVB	149798	154958	2.88%	0.227%
97	21.062	3070	3073	3080	rVB2	661761	821912	15.30%	1.206%
98	21.350	3117	3122	3127	rBV	2369421	2927603	54.50%	4.295%
99	23.468	3475	3482	3497	rVB	2193419	4217870	78.51%	6.187%
100	23.609	3500	3506	3524	rVB2	1472413	2943672	54.79%	4.318%

Sum of corrected areas: 68167756

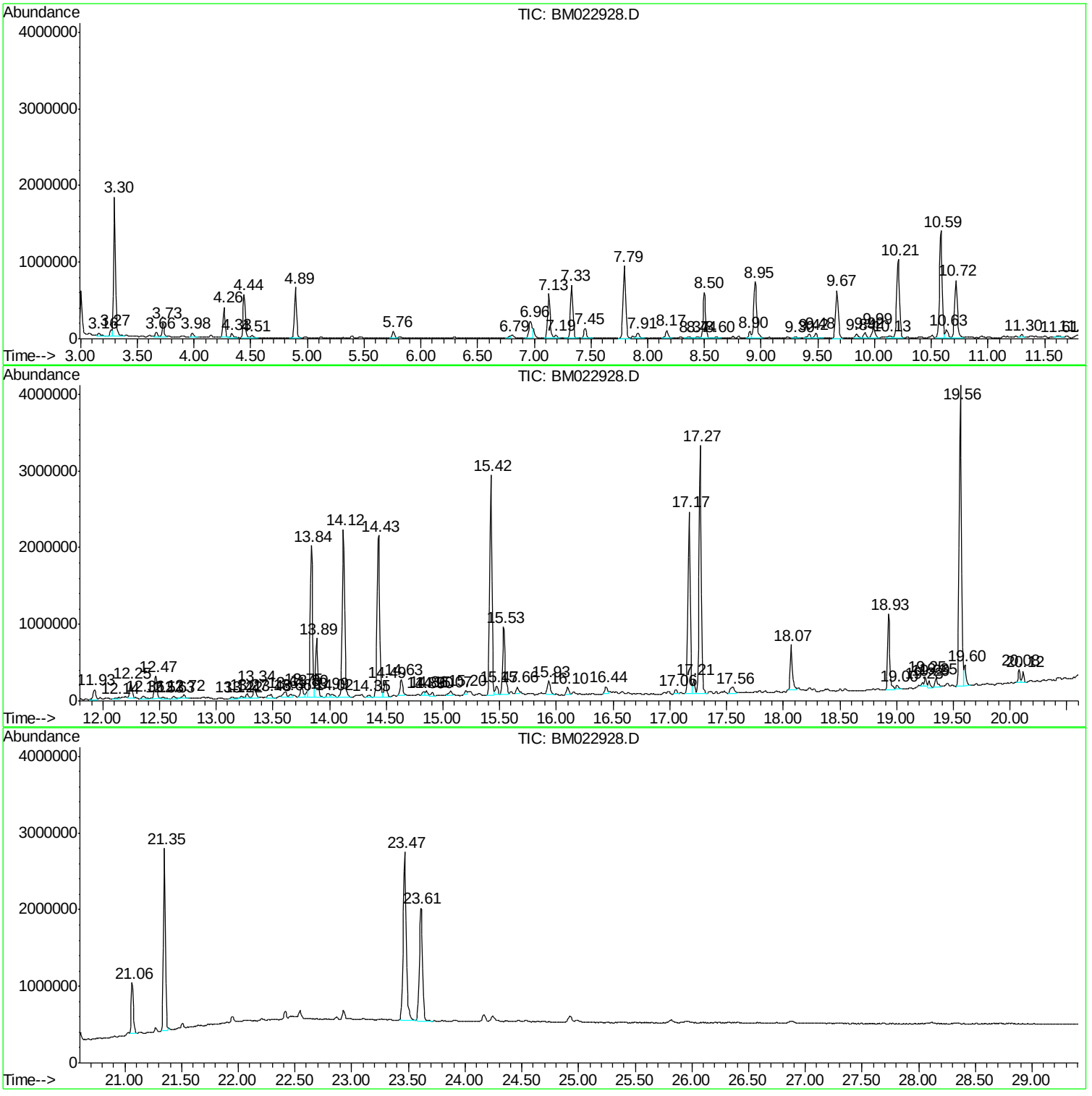
LSC Report - Integrated Chromatogram

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM100319\
Data File : BM022928.D
Acq On : 03 Oct 2019 19:49
Operator : JU
Sample : K4983-17
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFDK7

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM092719MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM100319\
Data File : BM022928.D
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Instrument : BNA_M
ClientSampleId : BFDK7

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM092719MA.M
Quant Title : SVOA CALIBRATION

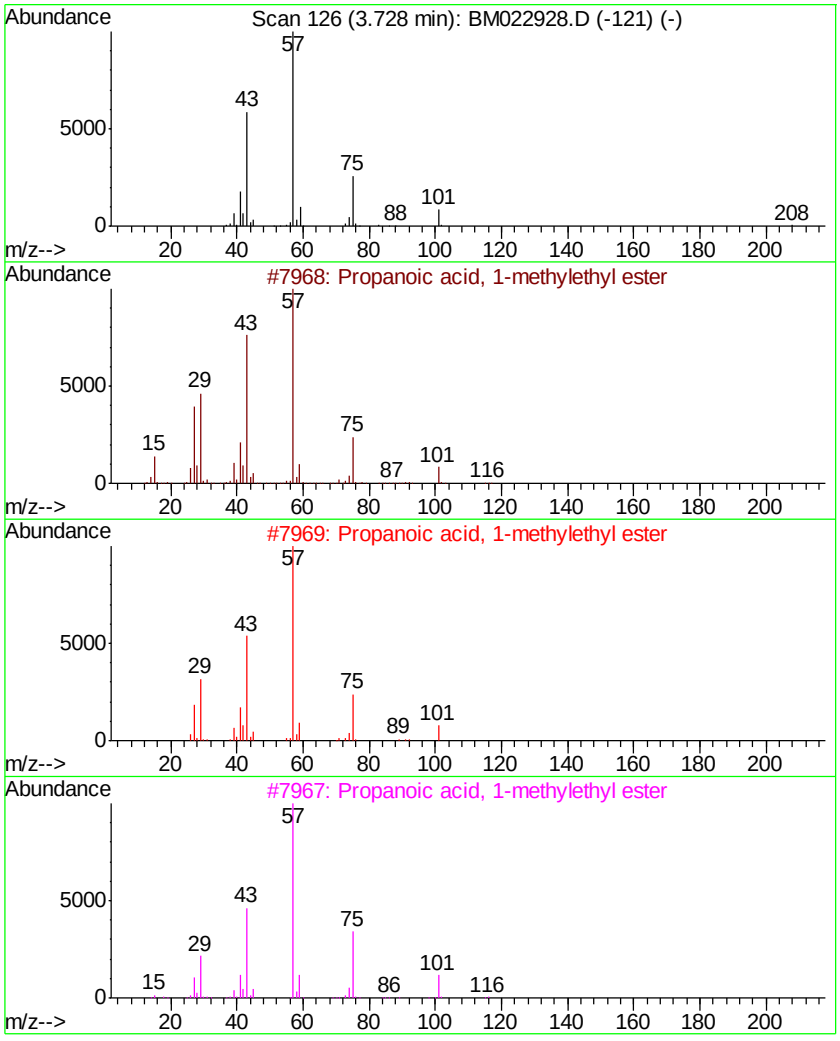
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

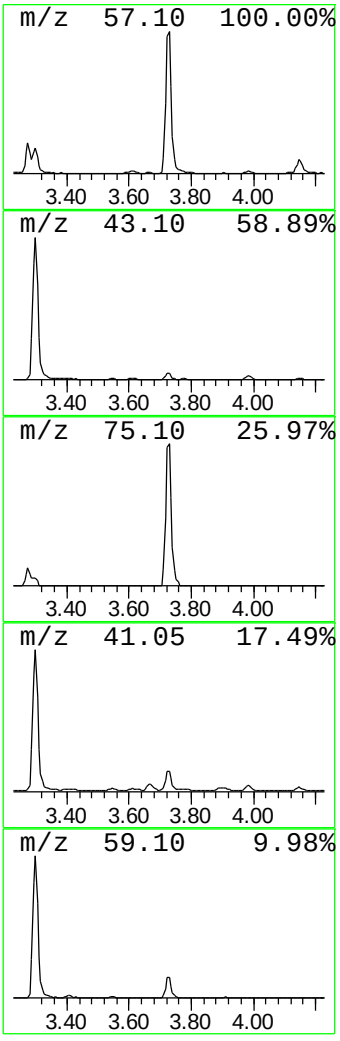
****

Peak Number 1 Propanoic acid, 1-methyleth... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.73	2.97 ng/ul	225364	1,4-Dichlorobenzene-d4	7.79

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propanoic acid, 1-methylethyl ester	116	C6H12O2	000637-78-5	78
2	Propanoic acid, 1-methylethyl ester	116	C6H12O2	000637-78-5	78
3	Propanoic acid, 1-methylethyl ester	116	C6H12O2	000637-78-5	64
4	Propanoic acid, 1-methylethyl ester	116	C6H12O2	000637-78-5	50
5	Propanoic acid, propyl ester	116	C6H12O2	000106-36-5	45





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM100319\
Data File : BM022928.D
Acq On : 03 Oct 2019 19:49
Operator : JU
Sample : K4983-17
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFDK7

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM092719MA.M
Quant Title : SVOA CALIBRATION

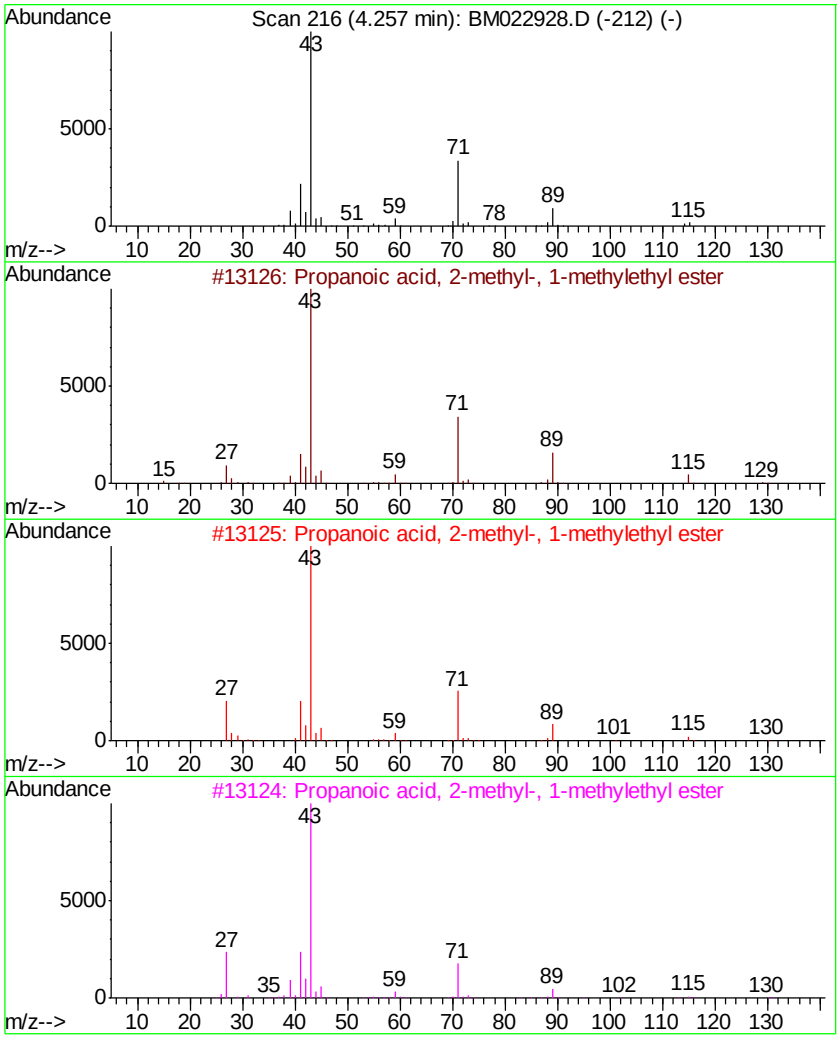
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

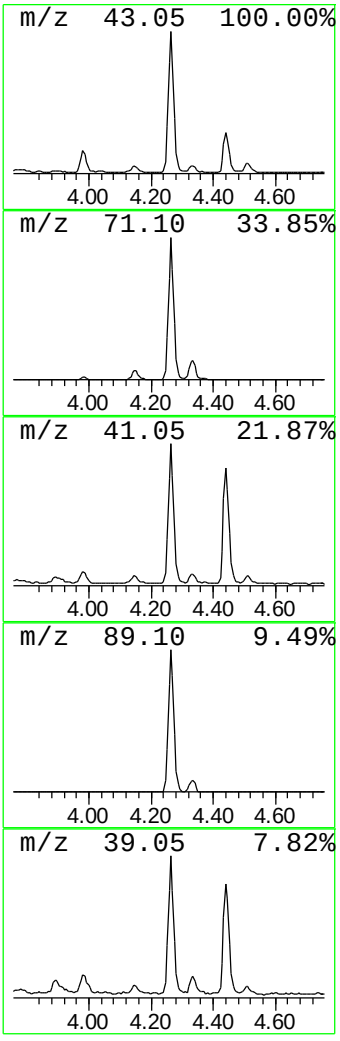
****

Peak Number 2 Propanoic acid, 2-methyl-, ... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.26	6.51 ng/ul	493864	1,4-Dichlorobenzene-d4	7.79

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propanoic acid, 2-methyl-, 1-met...	130	C7H14O2	000617-50-5	83
2	Propanoic acid, 2-methyl-, 1-met...	130	C7H14O2	000617-50-5	64
3	Propanoic acid, 2-methyl-, 1-met...	130	C7H14O2	000617-50-5	50
4	Propanoic acid, 2-methyl-, penty...	158	C9H18O2	002445-72-9	39
5	3,3-Dimethyl-2-pentanone	114	C7H14O	020669-04-9	33





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Instrument : BNA_M
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM092719MA.M
Quant Title : SVOA CALIBRATION

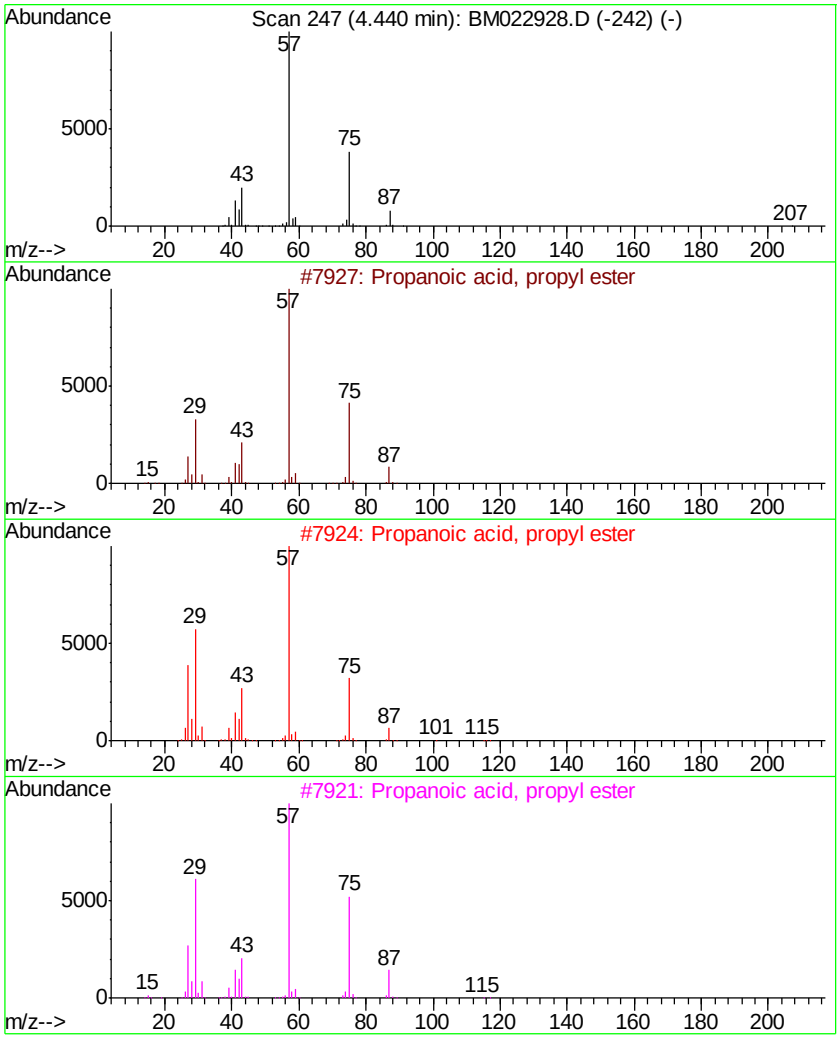
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

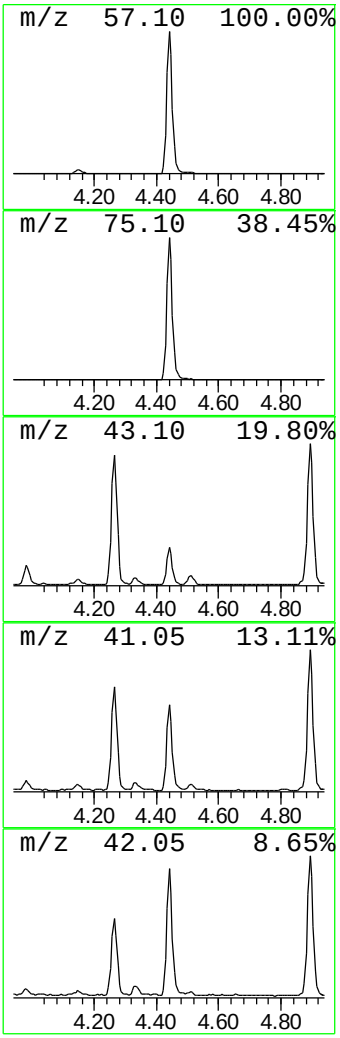
****

Peak Number 3 Propanoic acid, propyl ester Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.44	9.51 ng/ul	721299	1,4-Dichlorobenzene-d4	7.79

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propanoic acid, propyl ester	116	C6H12O2	000106-36-5	83
2	Propanoic acid, propyl ester	116	C6H12O2	000106-36-5	83
3	Propanoic acid, propyl ester	116	C6H12O2	000106-36-5	83
4	Propanoic acid, propyl ester	116	C6H12O2	000106-36-5	72
5	Propanoic acid, butyl ester	130	C7H14O2	000590-01-2	28





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM100319\
Data File : BM022928.D
Acq On : 03 Oct 2019 19:49
Operator : JU
Sample : K4983-17
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFDK7

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM092719MA.M
Quant Title : SVOA CALIBRATION

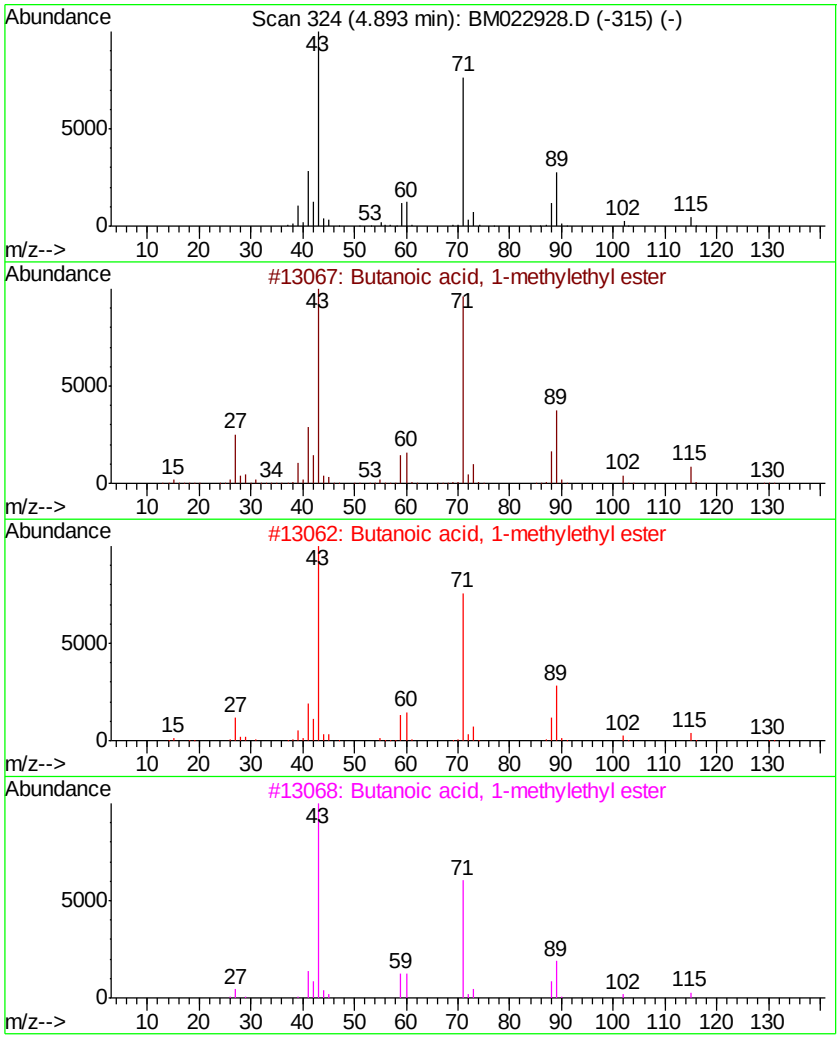
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

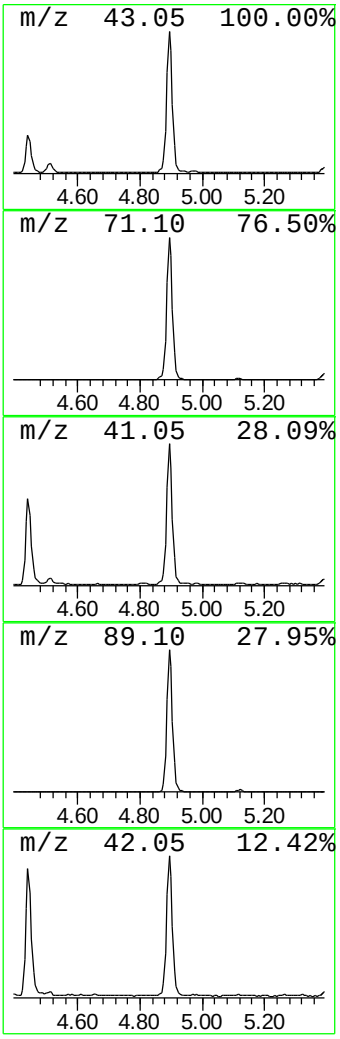
****

Peak Number 4 Butanoic acid, 1-methylethy... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.89	11.49 ng/ul	871127	1,4-Dichlorobenzene-d4	7.79

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butanoic acid, 1-methylethyl ester	130	C7H14O2	000638-11-9	90
2	Butanoic acid, 1-methylethyl ester	130	C7H14O2	000638-11-9	90
3	Butanoic acid, 1-methylethyl ester	130	C7H14O2	000638-11-9	83
4	Propanoic acid, 2-methyl-, 1-met...	130	C7H14O2	000617-50-5	78
5	Propanoic acid, 2-methyl-, propy...	130	C7H14O2	000644-49-5	56





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Acq On : 03 Oct 2019 19:49
Operator : JU
Sample : K4983-17
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFDK7

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM092719MA.M
Quant Title : SVOA CALIBRATION

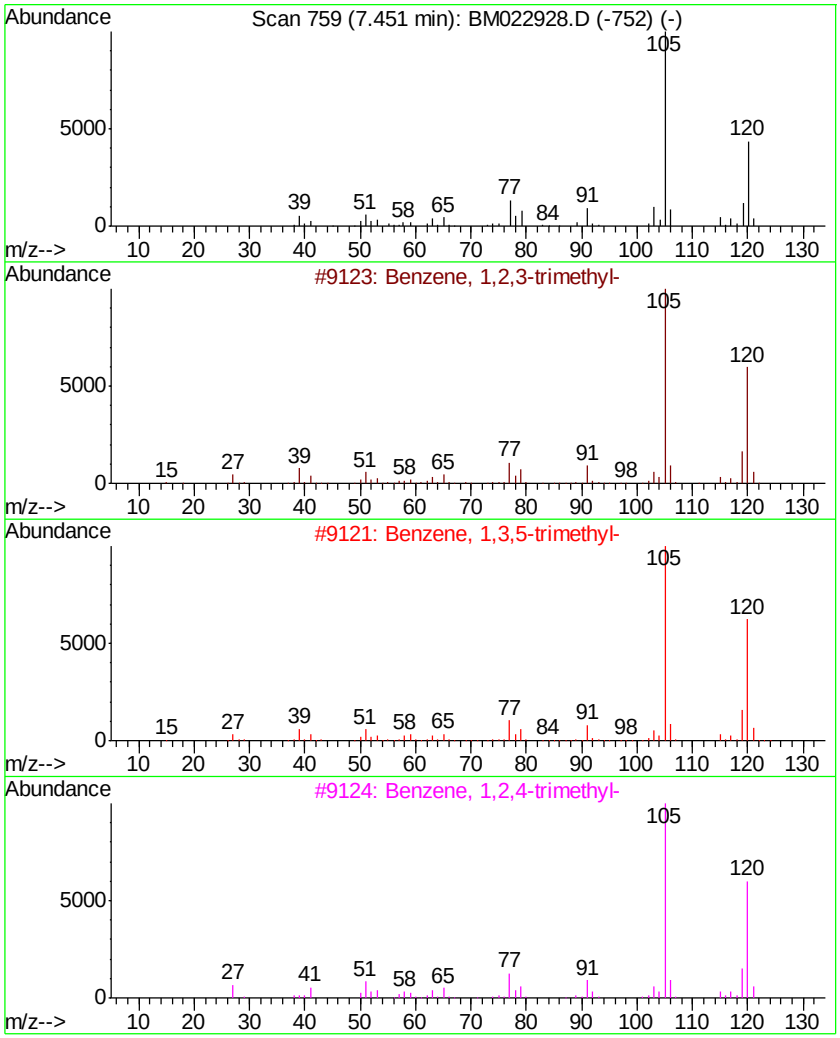
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

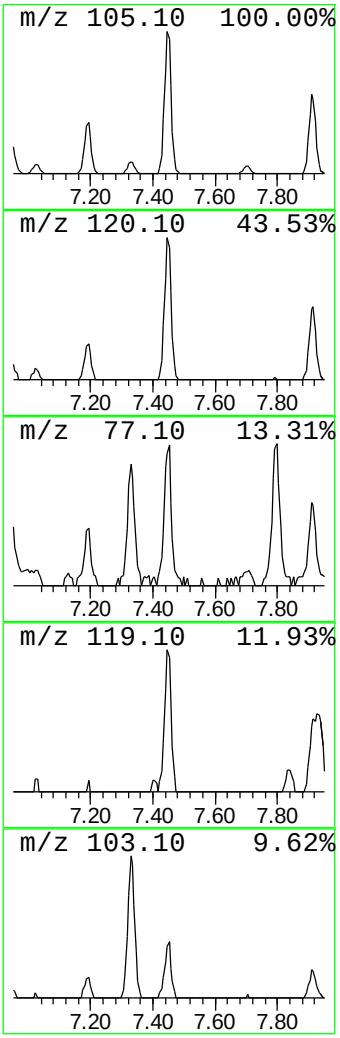
****

Peak Number 5 Benzene, 1,2,3-trimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.45	2.69 ng/ul	203940	1,4-Dichlorobenzene-d4	7.79

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
2	Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	95
3	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	95
4	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM100319\
Data File : BM022928.D
Acq On : 03 Oct 2019 19:49
Operator : JU
Sample : K4983-17
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFDK7

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM092719MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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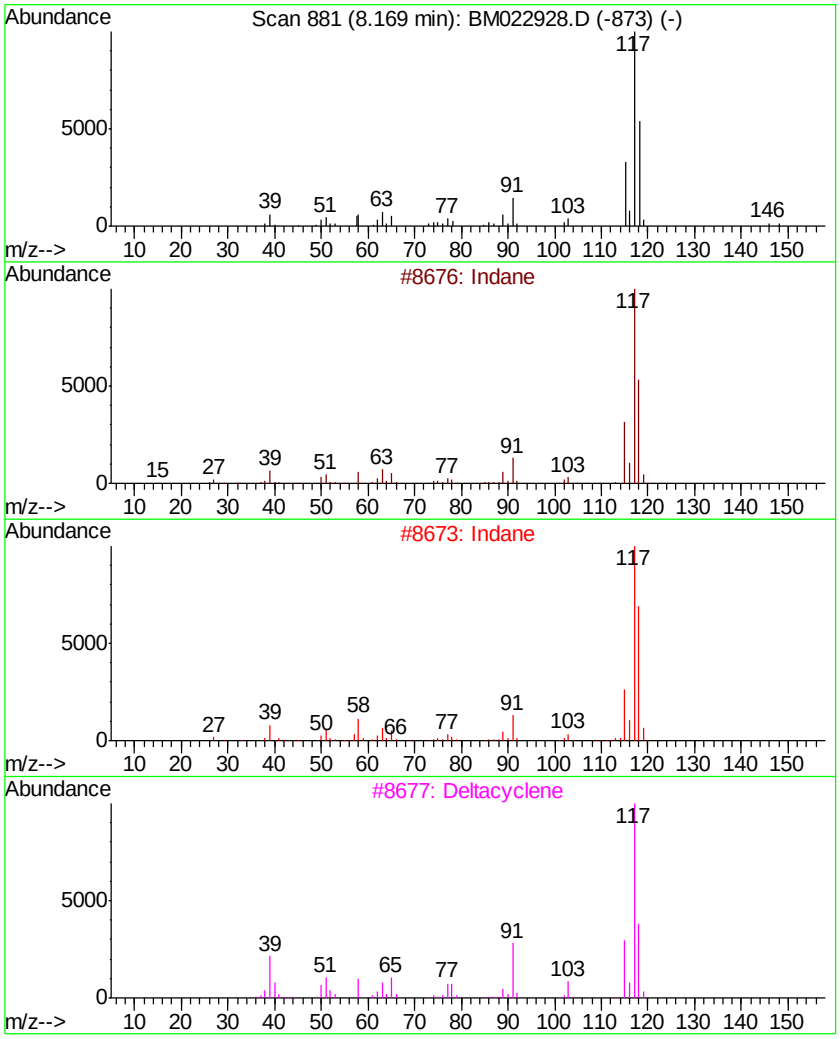
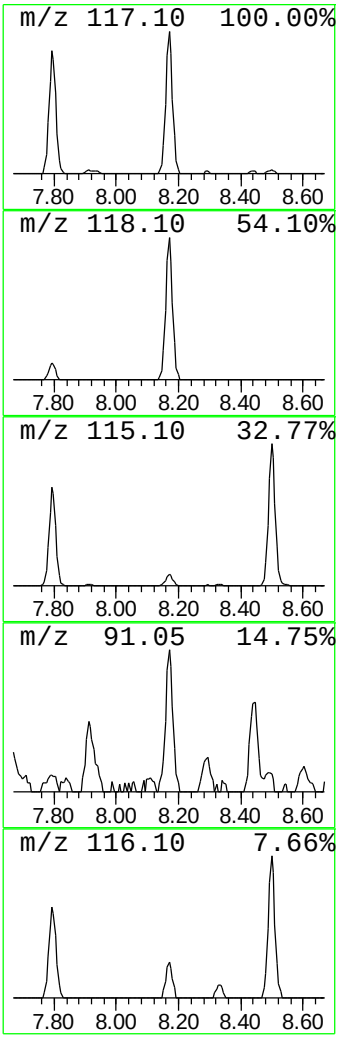
Peak Number 6 Indane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.17	2.39 ng/ul	181146	1,4-Dichlorobenzene-d4	7.79

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	93
2	Indane	118	C9H10	000496-11-7	83
3	Deltacyclene	118	C9H10	007785-10-6	74
4	Benzene, 1-propenyl-	118	C9H10	000637-50-3	74
5	Benzene, cyclopropyl-	118	C9H10	000873-49-4	72





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Instrument : BNA_M
ClientSampleId : BFDK7

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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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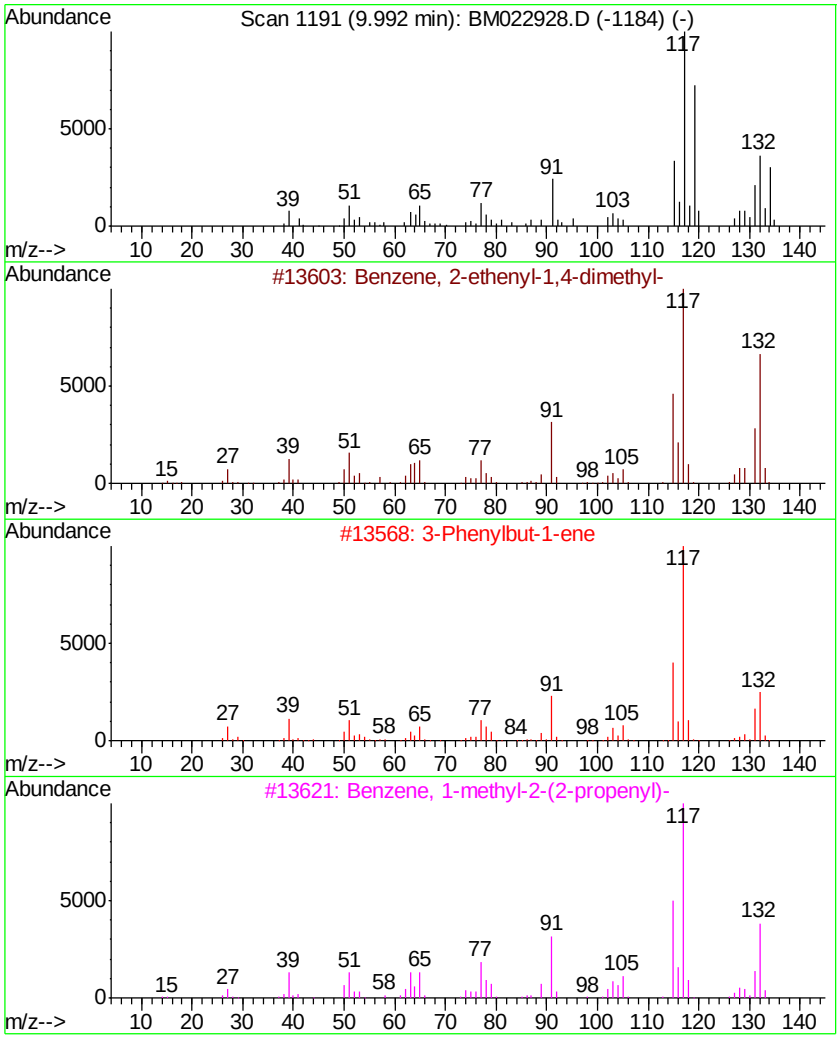
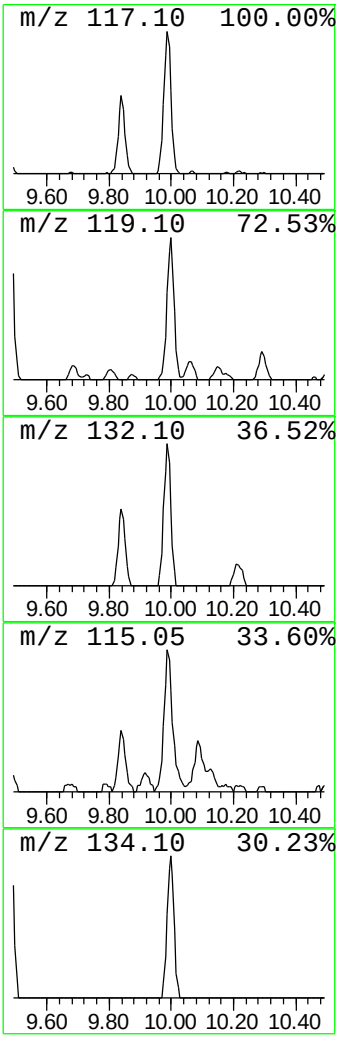
Peak Number 7 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.99	2.25 ng/ul	261665	Naphthalene-d8	10.59

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	90
2	3-Phenylbut-1-ene	132	C10H12	000934-10-1	70
3	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	64
4	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	64
5	1-Phenyl-1-butene	132	C10H12	000824-90-8	64





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Data File : BM022928.D
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Operator : JU
Sample : K4983-17
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFDK7

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM092719MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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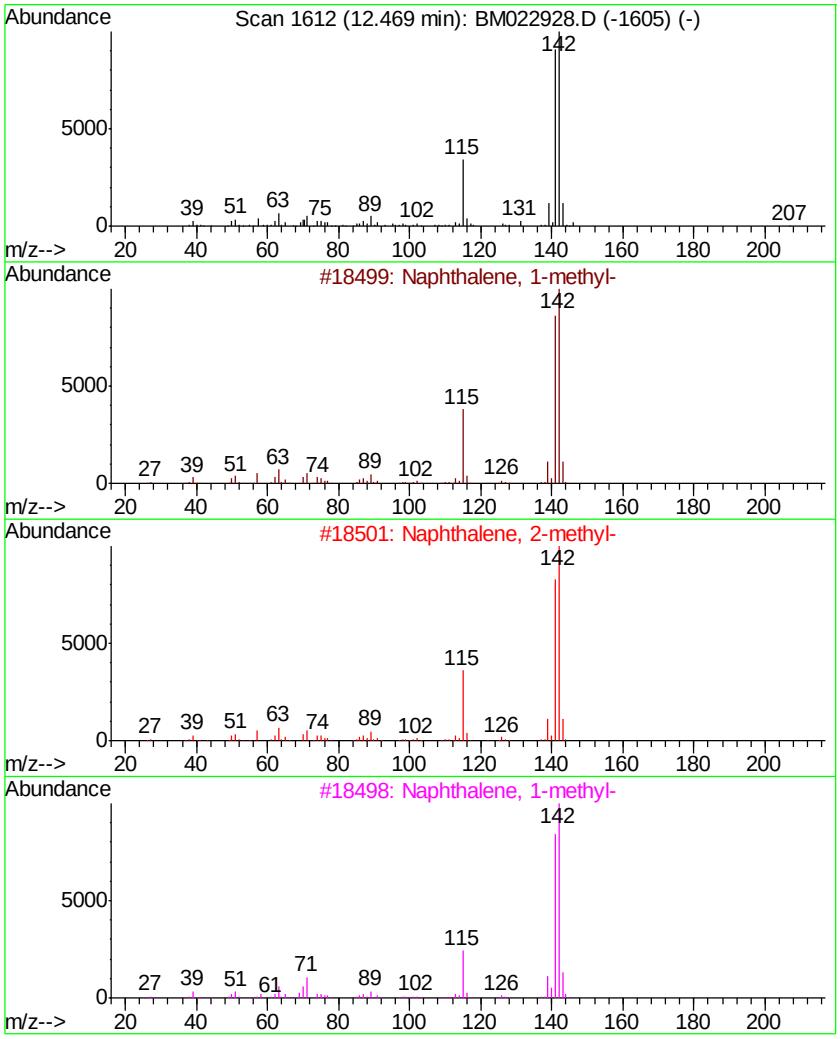
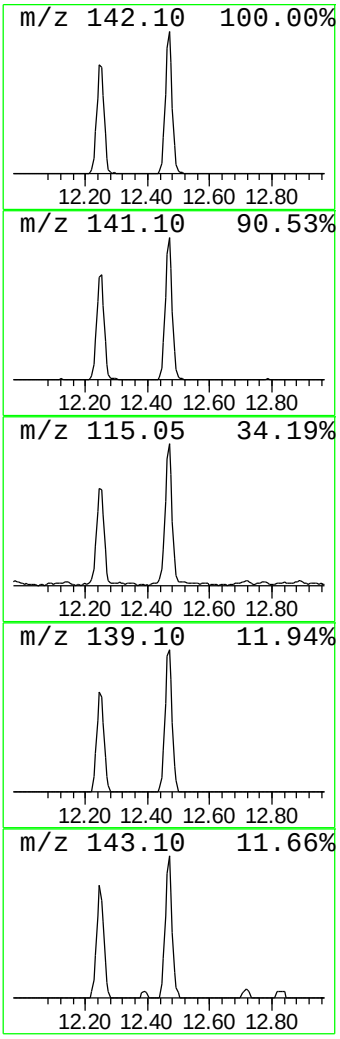
Peak Number 8 Naphthalene, 1-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.47	4.02 ng/ul	467680	Naphthalene-d8	10.59

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
2	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
3	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	94
4	1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	91
5	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	91





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ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFDK7

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Quant Title : SVOA CALIBRATION

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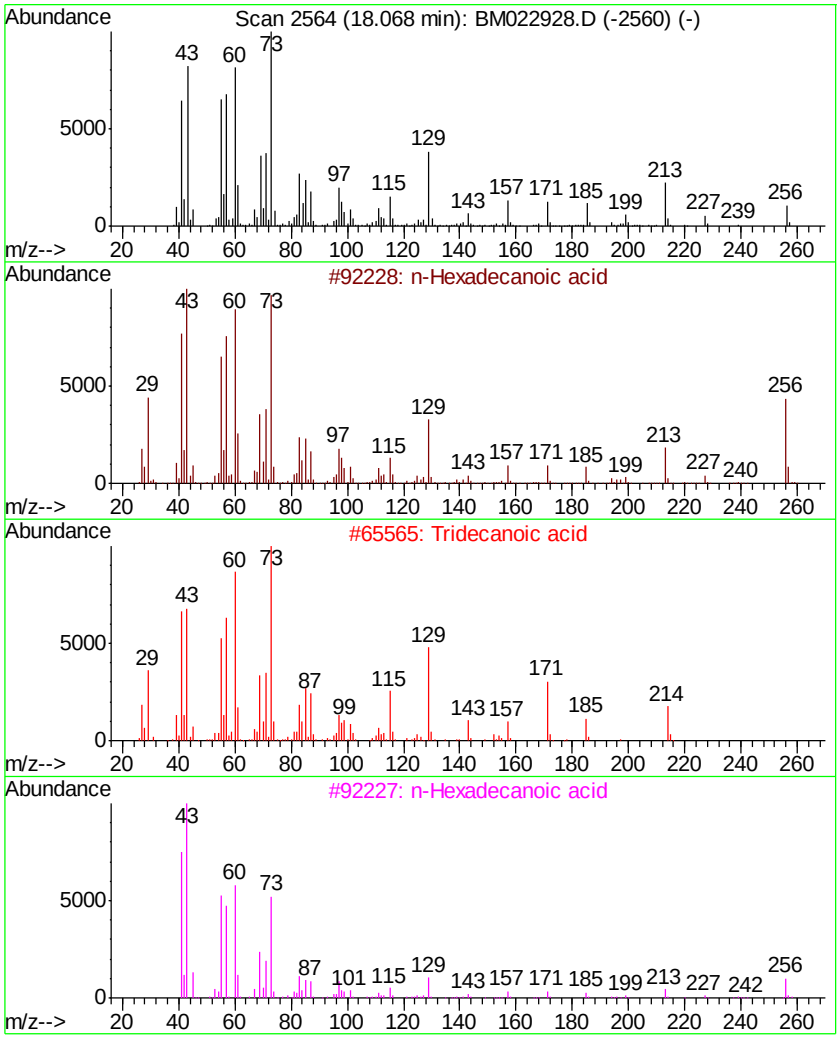
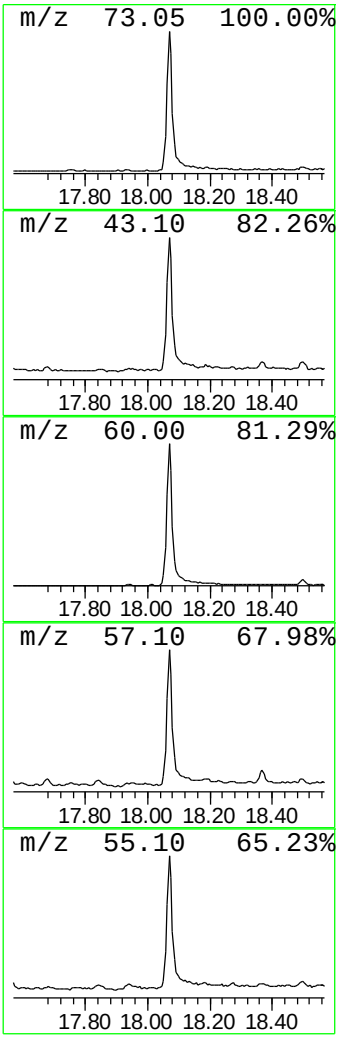
Peak Number 9 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.07	4.64 ng/ul	796244	Phenanthrene-d10	17.17

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2	Tridecanoic acid	214	C13H26O2	000638-53-9	95
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	94
4	Tridecanoic acid	214	C13H26O2	000638-53-9	93
5	Tetradecanoic acid	228	C14H28O2	000544-63-8	90





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Instrument : BNA_M
ClientSampleId : BFDK7

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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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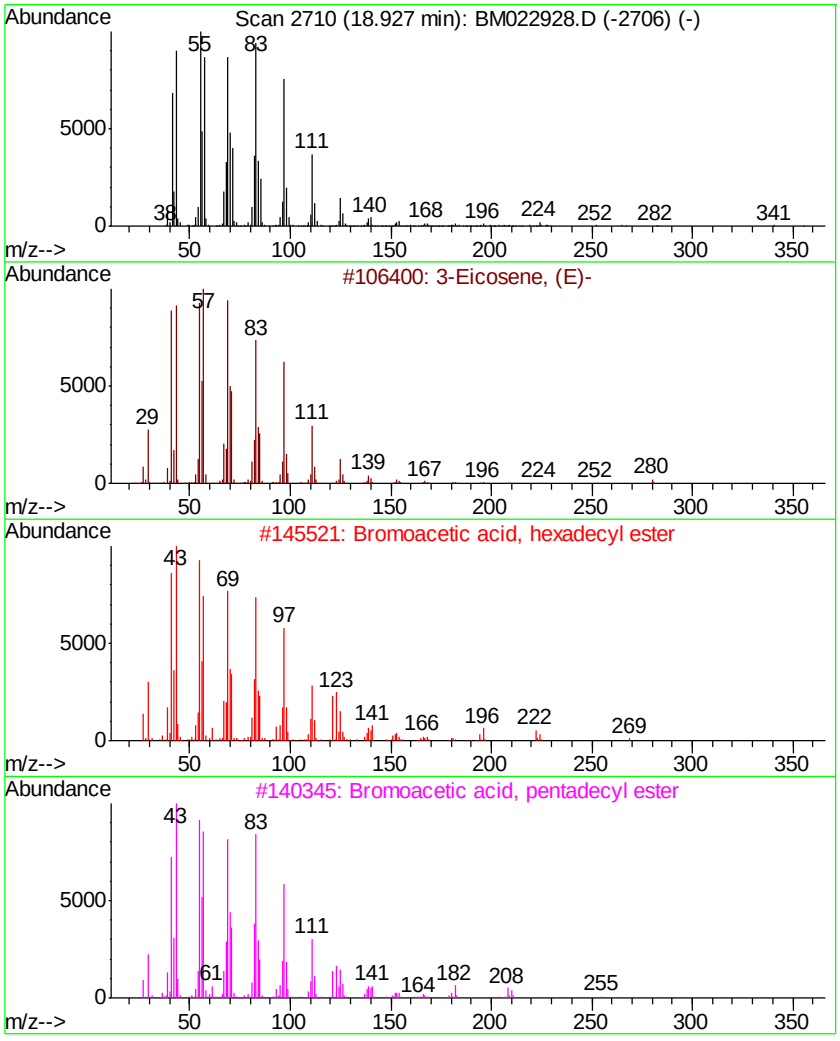
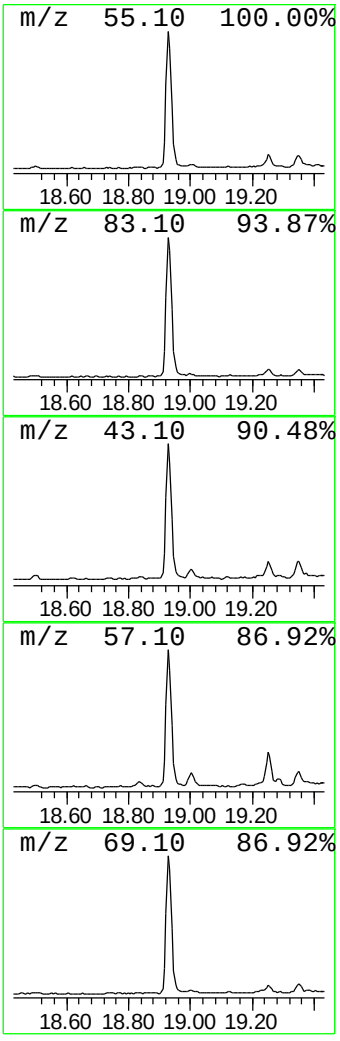
Peak Number 10 (DEL) Alkane: Cyclic18.93 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.93	7.35 ng/ul	1262290	Phenanthrene-d10	17.17

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Eicosene, (E)-	280	C20H40	074685-33-9	95
2	Bromoacetic acid, hexadecyl ester	362	C18H35BrO2	005454-48-8	94
3	Bromoacetic acid, pentadecyl ester	348	C17H33BrO2	131143-01-6	94
4	1-Octadecanol	270	C18H38O	000112-92-5	91
5	1-Hexadecanol	242	C16H34O	036653-82-4	91





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Instrument : BNA_M
ClientSampleId : BFDK7

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Quant Title : SVOA CALIBRATION

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TIC Integration Parameters: LSCINT.P

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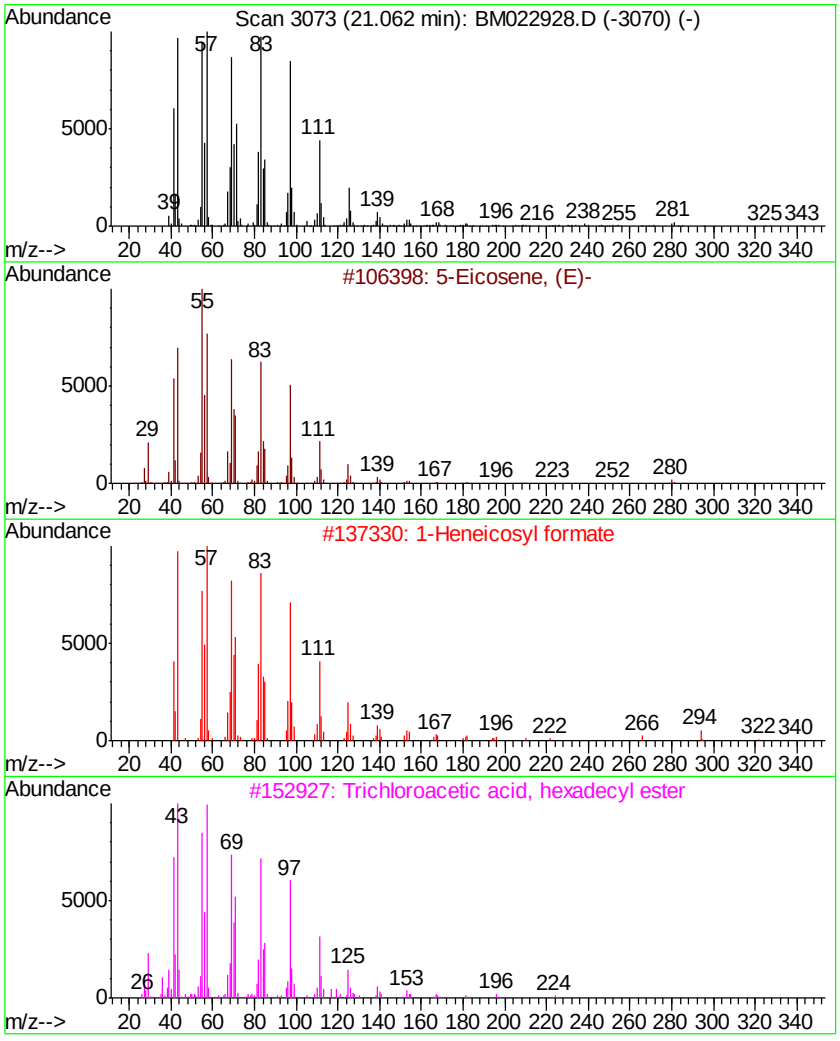
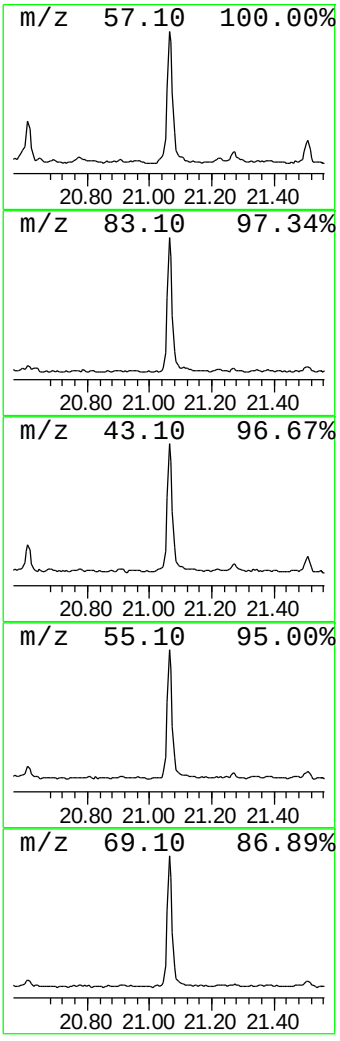
Peak Number 11 (DEL) Alkane: Cyclic21.06 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.06	5.61 ng/ul	821912	Chrysene-d12	21.35

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Eicosene, (E)-	280	C20H40	074685-30-6	99
2	1-Heneicosyl formate	340	C22H44O2	077899-03-7	94
3	Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	94
4	Cyclopentadecane	210	C15H30	000295-48-7	93
5	Heptafluorobutanoic acid, heptad...	452	C21H35F7O2	1000282-97-3	93





Tentatively Identified Compound (LSC) summary

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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Propanoic acid, 1...	3.73	3.0	ng/ul	225364	1	7.79	1516870	20.0
Propanoic acid, 2...	4.26	6.5	ng/ul	493864	1	7.79	1516870	20.0
Propanoic acid, p...	4.44	9.5	ng/ul	721299	1	7.79	1516870	20.0
Butanoic acid, 1-...	4.89	11.5	ng/ul	871127	1	7.79	1516870	20.0
Benzene, 1,2,3-tr...	7.45	2.7	ng/ul	203940	1	7.79	1516870	20.0
Indane	8.17	2.4	ng/ul	181146	1	7.79	1516870	20.0
Benzene, 2-etheny...	9.99	2.3	ng/ul	261665	2	10.59	2325880	20.0
Naphthalene, 1-me...	12.47	4.0	ng/ul	467680	2	10.59	2325880	20.0
n-Hexadecanoic acid	18.07	4.6	ng/ul	796244	4	17.17	3433930	20.0
(DEL) Alkane: Cyc...	18.93	7.3	ng/ul	1262290	4	17.17	3433930	20.0
(DEL) Alkane: Cyc...	21.06	5.6	ng/ul	821912	5	21.35	2927600	20.0