LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102519\
Data File : BM023405.D
Acq On : 25 Oct 2019 14:09
Operator : JU
Sample : K5344-05
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BF6L2

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-BM102319MA.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	38	rVB	107719	171108	1.88%	0.135%
2	3.875	147	151	161	rVB	481341	689661	7.57%	0.544%
3	5.175	367	372	379	rVB	423486	623623	6.85%	0.492%
4	5.305	388	394	405	rVB	1420300	2713031	29.79%	2.140%
5	5.669	450	456	462	rBV	799721	1185232	13.02%	0.935%
6	5.763	469	472	481	rVB2	24211	45698	0.50%	0.036%
7	6.146	532	537	546	rVB3	27522	52953	0.58%	0.042%
8	6.322	563	567	573	rVB4	25191	48950	0.54%	0.039%
9	6.628	614	619	622	rBV	71435	111835	1.23%	0.088%
10	6.663	622	625	631	rVB3	48458	77292	0.85%	0.061%
11	6.734	632	637	642	rVB	287359	424125	4.66%	0.335%
12	6.810	642	650	653	rBV2	544976	1053426	11.57%	0.831%
13	6.869	658	660	667	rVB	209143	265817	2.92%	0.210%
14	6.975	672	678	684	rBV	1352855	2118370	23.26%	1.671%
15	7.034	684	688	693	rVB	160675	239867	2.63%	0.189%
16	7.169	704	711	720	rVB	1589943	2509464	27.56%	1.979%
17	7.293	725	732	739	rVB	717315	1114216	12.24%	0.879%
18	7.522	766	771	779	rVB9	32079	78841	0.87%	0.062%
19	7.634	783	790	800	rBV	2066265	3392262	37.25%	2.676%
20	7.716	800	804	805	rVV	86669	118196	1.30%	0.093%
21	7.751	805	810	818	rVB	279235	500436	5.50%	0.395%
22	8.004	844	853	861	rVB2	160873	432201	4.75%	0.341%
23	8.128	871	874	878	rBV4	44238	72419	0.80%	0.057%
24	8.187	878	884	888	rVV2	66084	113743	1.25%	0.090%
25	8.281	894	900	904	rVV	133101	219551	2.41%	0.173%
26	8.340	904	910	922	rVV	1341968	2236447	24.56%	1.764%
27	8.451	922	929	936	rVB	611768	1014812	11.14%	0.800%
28	8.587	948	952	957	rVV	73946	118185	1.30%	0.093%
29	8.640	957	961	968	rVB	81364	132838	1.46%	0.105%
30	8.740	968	978	980	rBV	244029	405231	4.45%	0.320%
31	8.781	980	985	998	rVV	1777450	3101662	34.06%	2.446%
32	8.875	998	1001	1005	rVV5	31834	49752	0.55%	0.039%
33	8.945	1005	1013	1026	rVV	338238	687139	7.55%	0.542%
34	9.063	1026	1033	1040	rVV4	64549	157759	1.73%	0.124%

LSC Area Percent Report

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Acq On : 25 Oct 2019 14:09
Operator : JU
Sample : K5344-05
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BF6L2

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-BM102319MA.M
Title : SVOA CALIBRATION

35	9.257	1058	1066	1071	rVV4	138240	251529	2.76%	0.198%
36	9.322	1071	1077	1085	rVV	168337	291452	3.20%	0.230%
37	9.498	1100	1107	1115	rBV	1413516	2383328	26.17%	1.880%
38	9.604	1120	1125	1129	rVV2	96170	182952	2.01%	0.144%
39	9.675	1133	1137	1147	rVB	124851	233780	2.57%	0.184%
40	9.769	1147	1153	1157	rBV7	55197	141604	1.56%	0.112%
41	9.822	1157	1162	1171	rVB3	207327	424957	4.67%	0.335%
42	9.904	1171	1176	1183	rBV3	109506	187677	2.06%	0.148%
43	10.040	1189	1199	1210	rVB	2436329	4085141	44.86%	3.222%
44	10.128	1210	1214	1221	rVB3	28205	56464	0.62%	0.045%
45	10.222	1221	1230	1236	rBV8	28257	82187	0.90%	0.065%
46	10.281	1236	1240	1246	rBV4	53555	98107	1.08%	0.077%
47	10.345	1246	1251	1256	rBV	65984	117669	1.29%	0.093%
48	10.416	1256	1263	1267	rBV	2827089	4762364	52.30%	3.756%
49	10.463	1267	1271	1279	rVB	781620	1295562	14.23%	1.022%
50	10.792	1321	1327	1330	rBV7	33399	66122	0.73%	0.052%
51	10.839	1330	1335	1338	rVV6	33766	58372	0.64%	0.046%
52	10.969	1352	1357	1366	rVB8	47772	115767	1.27%	0.091%
53	11.057	1366	1372	1373	rBV5	34276	52350	0.57%	0.041%
54	11.234	1396	1402	1403	rBV6	28222	46137	0.51%	0.036%
55	11.328	1414	1418	1425	rVB3	41584	81751	0.90%	0.064%
56	11.522	1444	1451	1459	rVB8	42399	125224	1.38%	0.099%
57	11.616	1461	1467	1471	rBV8	29511	52552	0.58%	0.041%
58	11.769	1483	1493	1497	rBV3	67124	173574	1.91%	0.137%
59	11.857	1504	1508	1513	rBV8	20765	42492	0.47%	0.034%
60	11.975	1523	1528	1531	rVV3	38231	83547	0.92%	0.066%
61	12.010	1531	1534	1538	rVV2	41973	65454	0.72%	0.052%
62	12.086	1538	1547	1553	rVB	499911	827024	9.08%	0.652%
63	12.304	1575	1584	1590	rVB	333612	560353	6.15%	0.442%
64	12.386	1590	1598	1604	rBV5	64274	125941	1.38%	0.099%
65	12.481	1609	1614	1618	rVB8	28887	47022	0.52%	0.037%
66	13.128	1718	1724	1731	rVV5	125101	245380	2.69%	0.194%
67	13.198	1731	1736	1747	rVV6	50092	145064	1.59%	0.114%
68	13.328	1748	1758	1765	rVB6	60391	196288	2.16%	0.155%
69	13.451	1773	1779	1785	rBV5	69868	181390	1.99%	0.143%
70	13.604	1799	1805	1810	rBV2	127084	244050	2.68%	0.192%
71	13.657	1810	1814	1815	rVV	89468	108971	1.20%	0.086%

LSC Area Percent Report

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Operator : JU
Sample : K5344-05
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BF6L2

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-BM102319MA.M
Title : SVOA CALIBRATION

72	13.698	1815	1821	1825	rVV	4516593	6244748	68.58%	4.925%
73	13.745	1825	1829	1836	rVB	2409419	3148668	34.58%	2.483%
74	13.839	1841	1845	1848	rBV2	90932	133637	1.47%	0.105%
75	13.869	1848	1850	1860	rVB6	45174	93749	1.03%	0.074%
76	13.969	1861	1867	1874	rBV	4864952	7537375	82.77%	5.945%
77	14.198	1902	1906	1912	rVB7	79456	158275	1.74%	0.125%
78	14.280	1913	1920	1926	rBV2	4302974	6476920	71.13%	5.109%
79	14.333	1926	1929	1935	rVB2	203728	364222	4.00%	0.287%
80	14.486	1950	1955	1958	rBV	440161	647869	7.11%	0.511%
81	14.516	1958	1960	1964	rVB	225320	276076	3.03%	0.218%
82	14.680	1985	1988	1993	rBV4	32750	56110	0.62%	0.044%
83	14.727	1993	1996	2000	rVV4	61870	81805	0.90%	0.065%
84	14.904	2023	2026	2032	rBV6	48370	85868	0.94%	0.068%
85	15.280	2083	2090	2096	rBV	6339399	9105951	100.00%	7.182%
86	15.333	2096	2099	2104	rVB2	112295	157973	1.73%	0.125%
87	15.404	2106	2111	2117	rVV	1075945	1481060	16.26%	1.168%
88	15.457	2117	2120	2125	rBV6	62913	100336	1.10%	0.079%
89	17.027	2382	2387	2392	rBV2	4512739	6234035	68.46%	4.917%
90	17.068	2392	2394	2398	rVV	205255	248555	2.73%	0.196%
91	17.127	2398	2404	2415	rVB	6187141	8570013	94.11%	6.759%
92	17.410	2449	2452	2456	rVB	129357	154326	1.69%	0.122%
93	17.951	2540	2544	2553	rBV	1368405	2062485	22.65%	1.627%
94	19.239	2759	2763	2769	rBV2	440083	656533	7.21%	0.518%
95	19.427	2790	2795	2806	rVB	6052922	8249831	90.60%	6.507%
96	20.968	3053	3057	3061	rBV	1086434	1270214	13.95%	1.002%
97	21.009	3062	3064	3068	rVB	320420	353206	3.88%	0.279%
98	21.221	3096	3100	3106	rBV	4055690	5153405	56.59%	4.065%
99	23.286	3445	3451	3458	rVB	4299055	7564680	83.07%	5.966%
100	23.421	3469	3474	3486	rVB	2910512	5602684	61.53%	4.419%

Sum of corrected areas: 126786349

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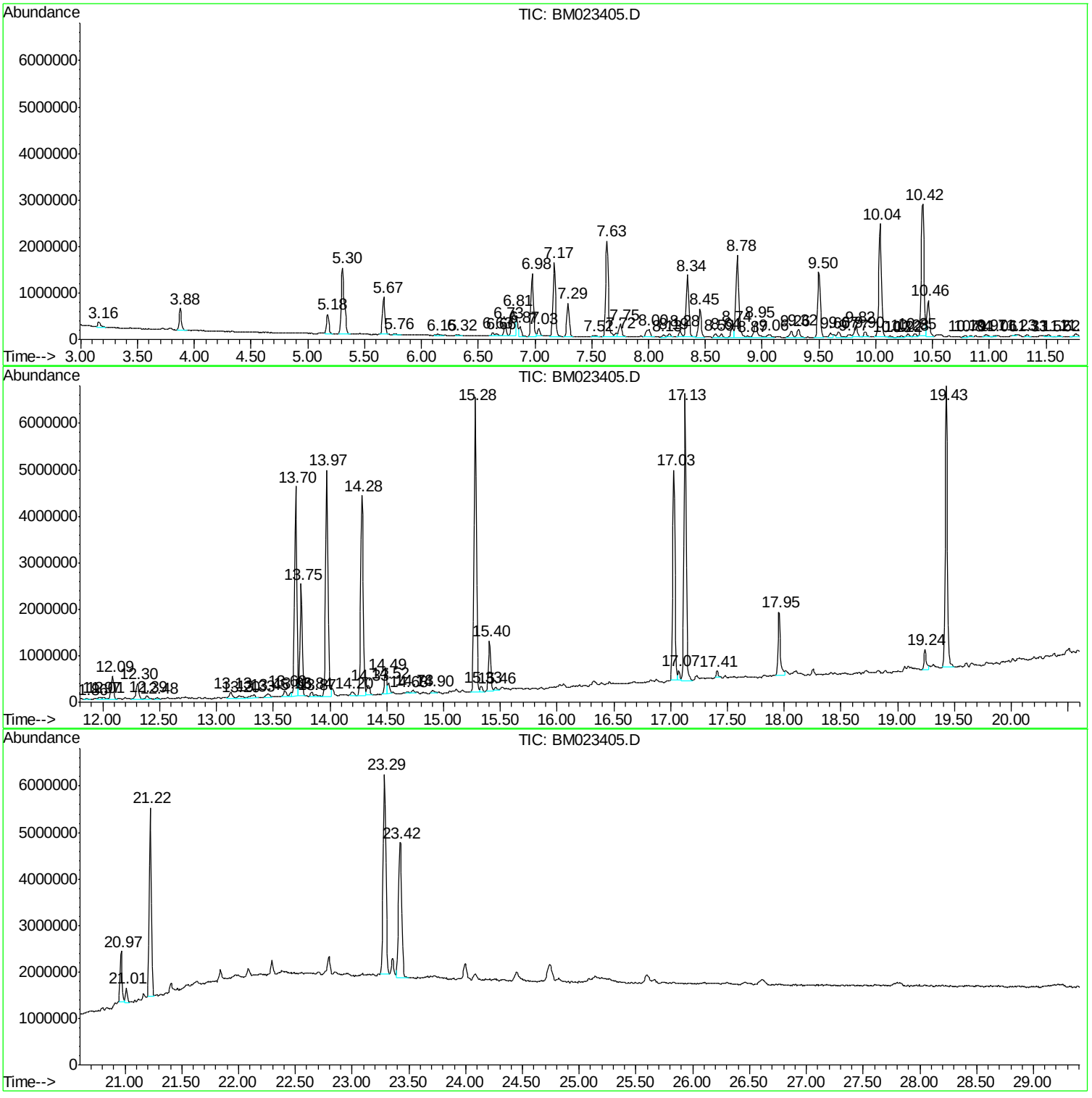
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BF6L2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.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_M
ClientSampleId : BF6L2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
Quant Title : SVOA CALIBRATION

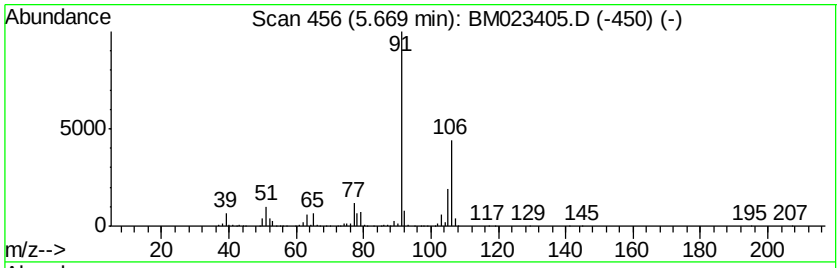
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

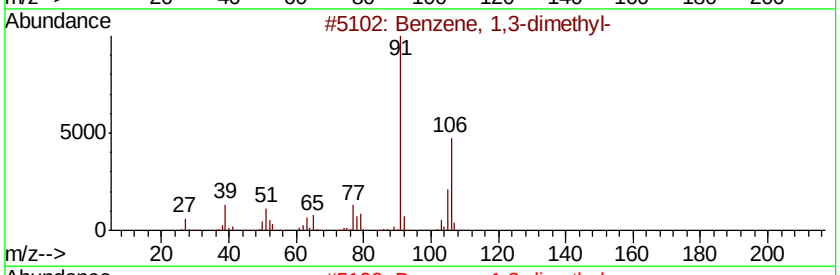
****

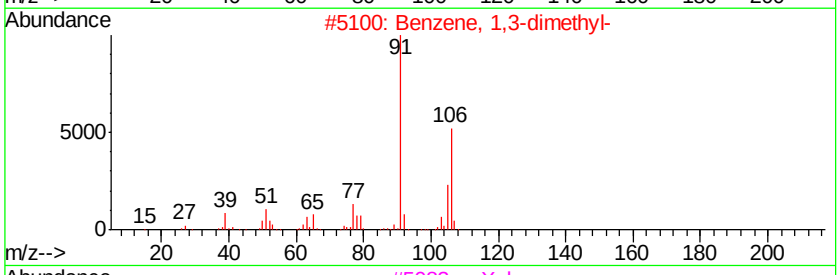
Peak Number 4 Benzene, 1,3-dimethyl- Concentration Rank 2

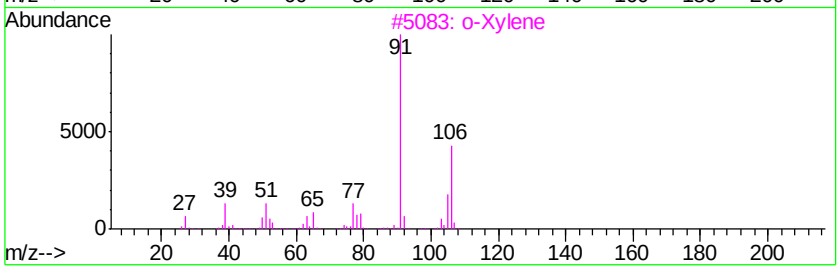
R.T.	EstConc	Area	Relative to ISTD	R.T.
5.67	6.99 ng/ul	1185230	1,4-Dichlorobenzene-d4	7.63

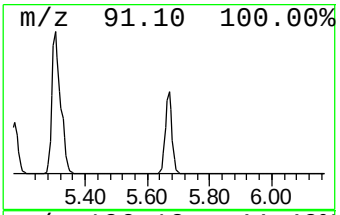
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	97
2	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	97
3	o-Xylene	106	C8H10	000095-47-6	97
4	p-Xylene	106	C8H10	000106-42-3	97
5	o-Xylene	106	C8H10	000095-47-6	97

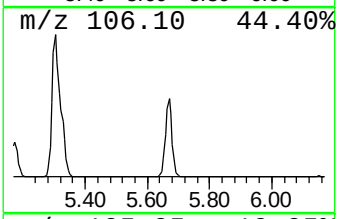


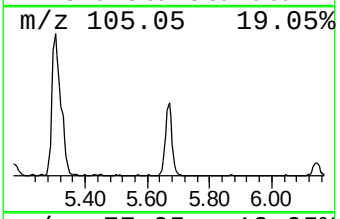


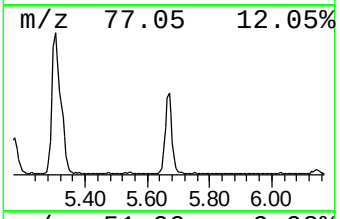


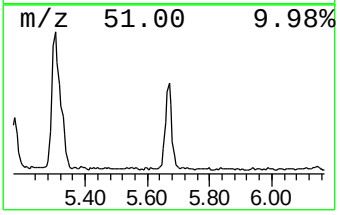












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Instrument : BNA_M
ClientSampleId : BF6L2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
Quant Title : SVOA CALIBRATION

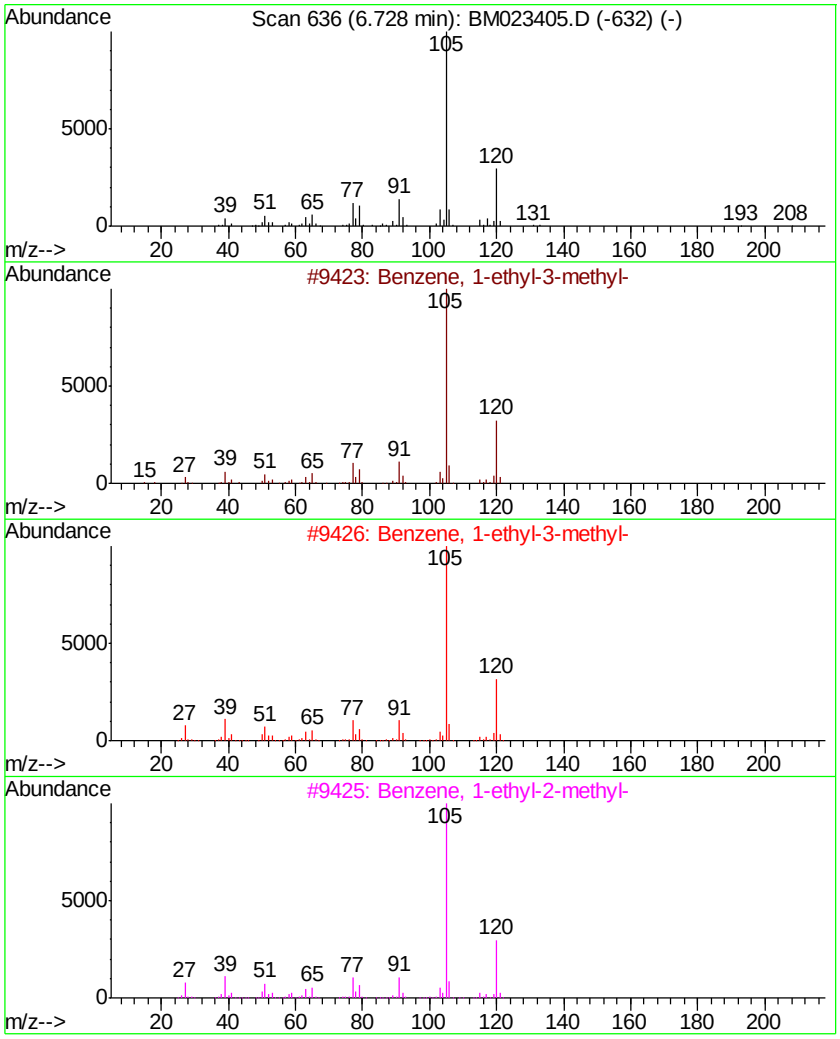
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

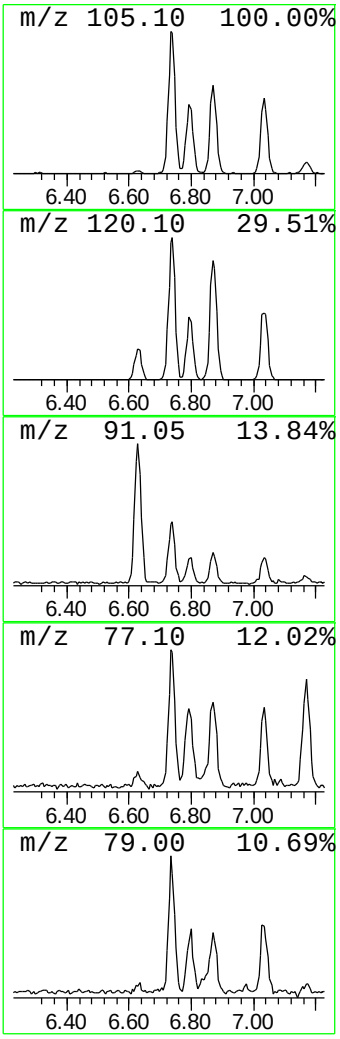
****

Peak Number 5 Benzene, 1-ethyl-3-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.73	2.50 ng/ul	424125	1,4-Dichlorobenzene-d4	7.63

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	97
2	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
3	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
4	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	93





Library Search Compound Report

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Data File : BM023405.D
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Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BF6L2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
Quant Title : SVOA CALIBRATION

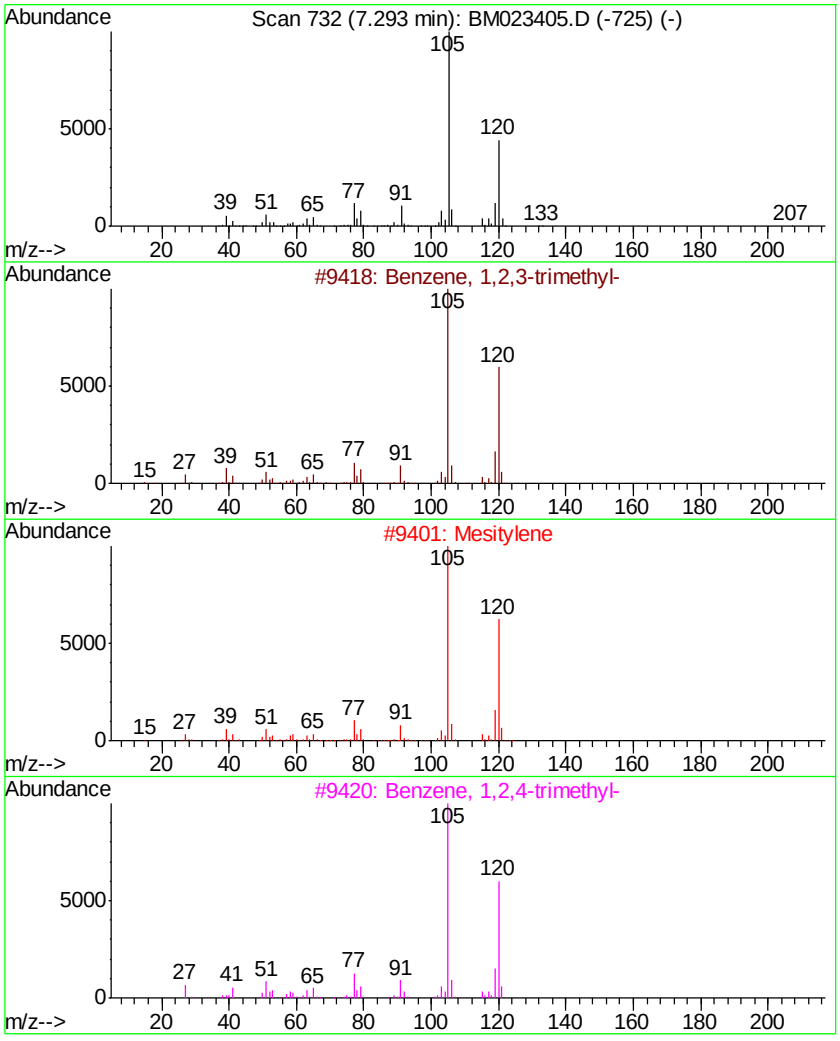
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

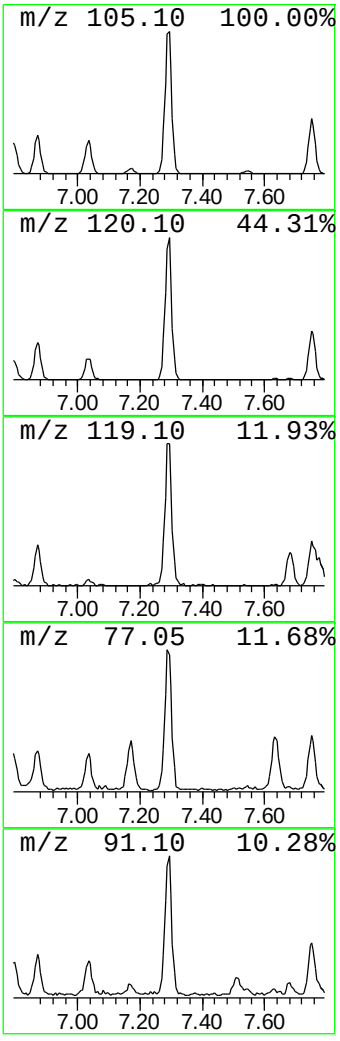
****

Peak Number 6 Benzene, 1,2,3-trimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.29	6.57 ng/ul	1114220	1,4-Dichlorobenzene-d4	7.63

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





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102519\
Data File : BM023405.D
Acq On : 25 Oct 2019 14:09
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Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BF6L2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
Quant Title : SVOA CALIBRATION

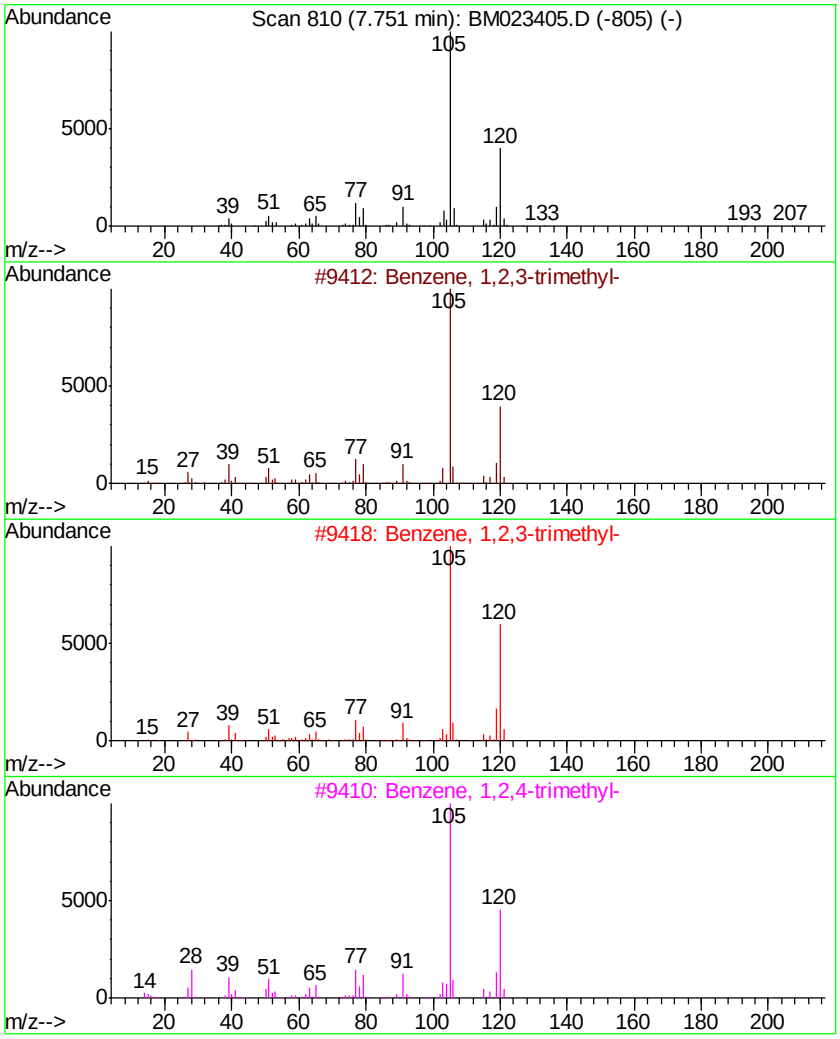
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

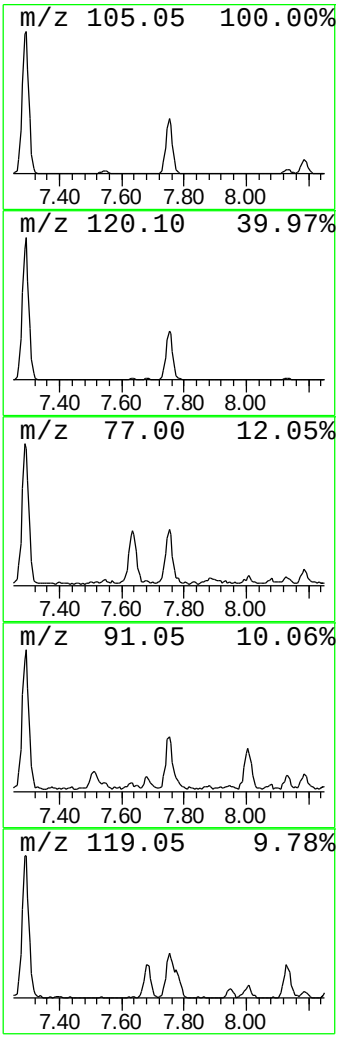
****

Peak Number 7 Benzene, 1,2,4-trimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.75	2.95 ng/ul	500436	1,4-Dichlorobenzene-d4	7.63

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
2	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
3	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
4	Mesitylene	120	C9H12	000108-67-8	91
5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102519\
Data File : BM023405.D
Acq On : 25 Oct 2019 14:09
Operator : JU
Sample : K5344-05
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BF6L2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
Quant Title : SVOA CALIBRATION

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

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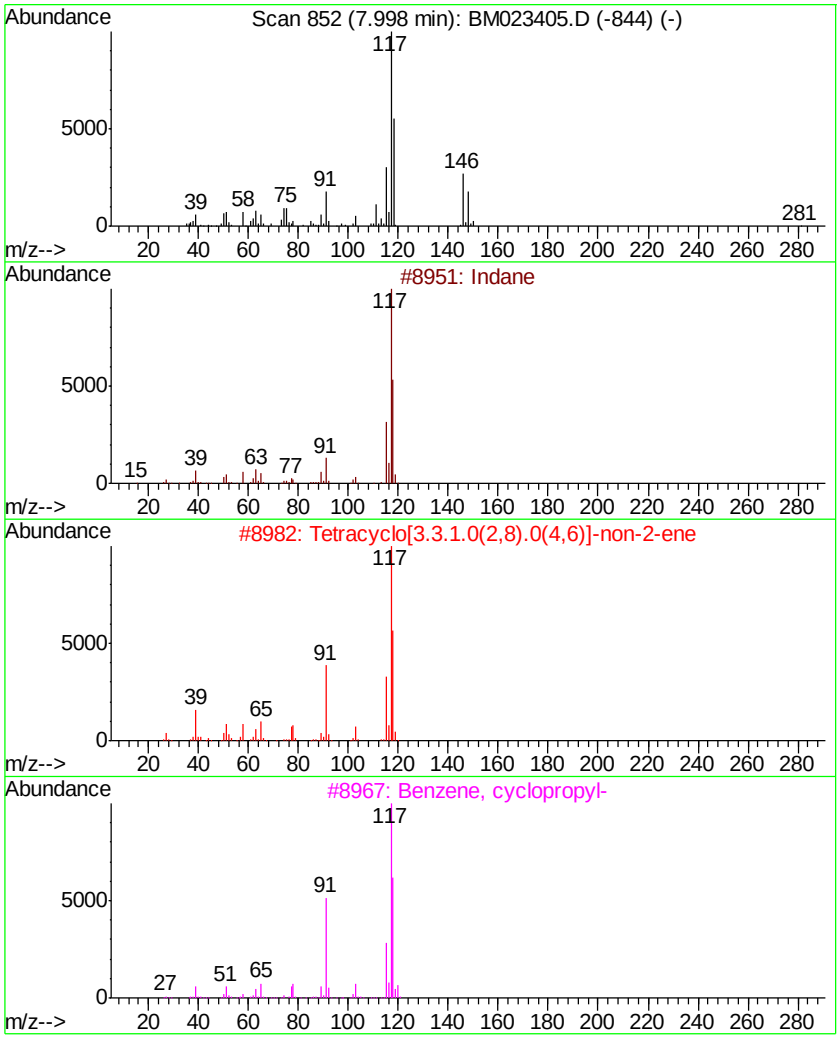
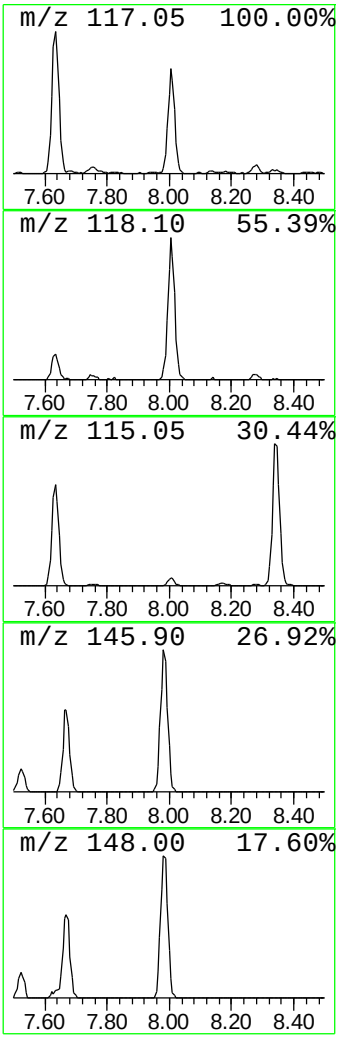
Peak Number 8 Indane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.00	2.55 ng/ul	432201	1,4-Dichlorobenzene-d4	7.63

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	81
2	Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	72
3	Benzene, cyclopropyl-	118	C9H10	000873-49-4	72
4	Indane	118	C9H10	000496-11-7	55
5	Indane	118	C9H10	000496-11-7	55





Library Search Compound Report

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Data File : BM023405.D
Acq On : 25 Oct 2019 14:09
Operator : JU
Sample : K5344-05
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BF6L2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
Quant Title : SVOA CALIBRATION

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

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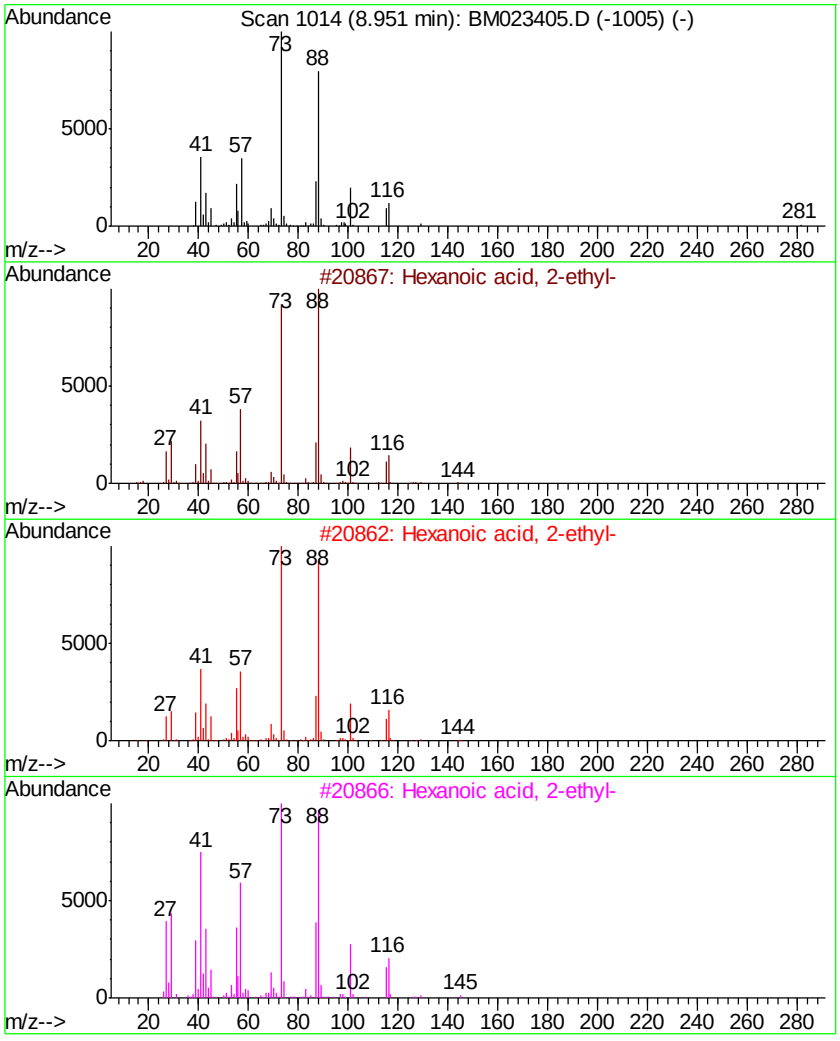
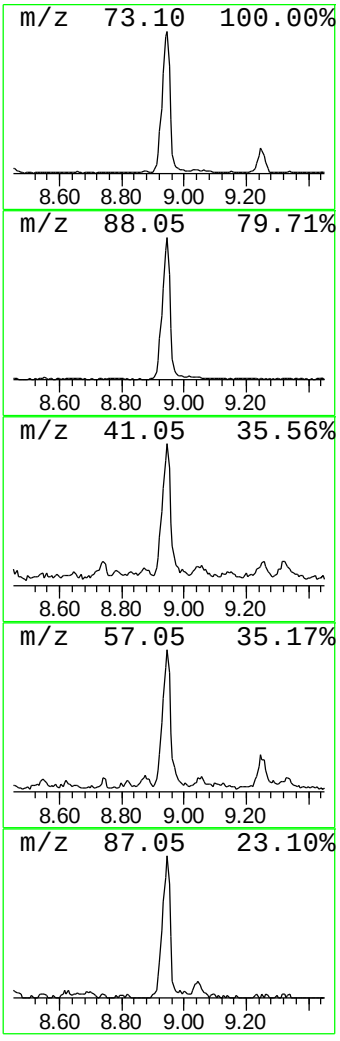
Peak Number 9 Hexanoic acid, 2-ethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.95	4.05 ng/ul	687139	1,4-Dichlorobenzene-d4	7.63

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	86
2	Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	72
3	Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	47
4	Heptanoic acid, 2-ethyl-	158	C9H18O2	003274-29-1	40
5	Methyl 2,3,4-tri-O-methyl-6-deox...	220	C10H20O5	072983-16-5	40





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Data File : BM023405.D
Acq On : 25 Oct 2019 14:09
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Sample : K5344-05
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BF6L2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
Quant Title : SVOA CALIBRATION

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

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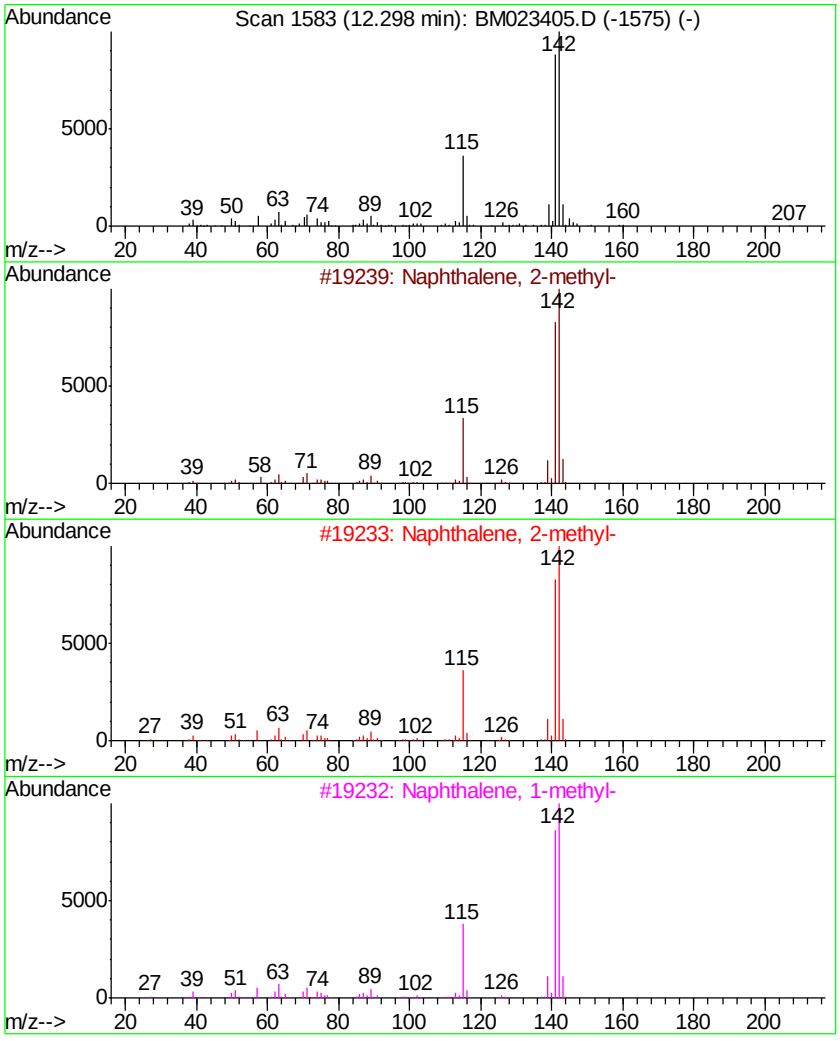
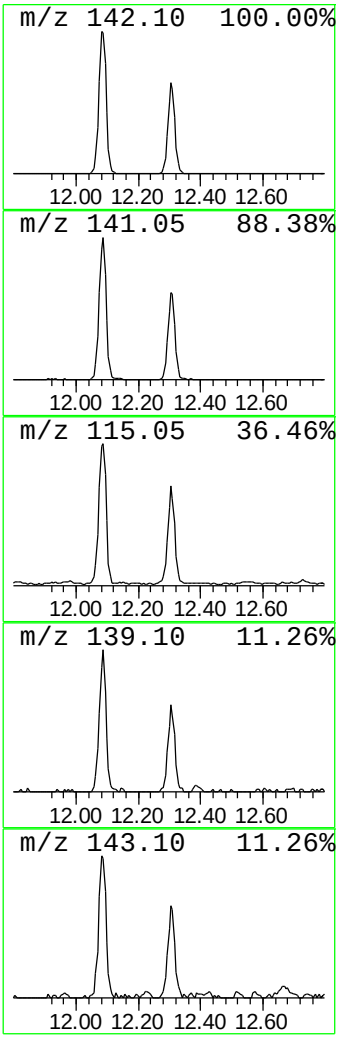
Peak Number 10 Naphthalene, 1-methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.30	2.35 ng/ul	560353	Naphthalene-d8	10.42

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





Library Search Compound Report

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Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BF6L2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
Quant Title : SVOA CALIBRATION

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

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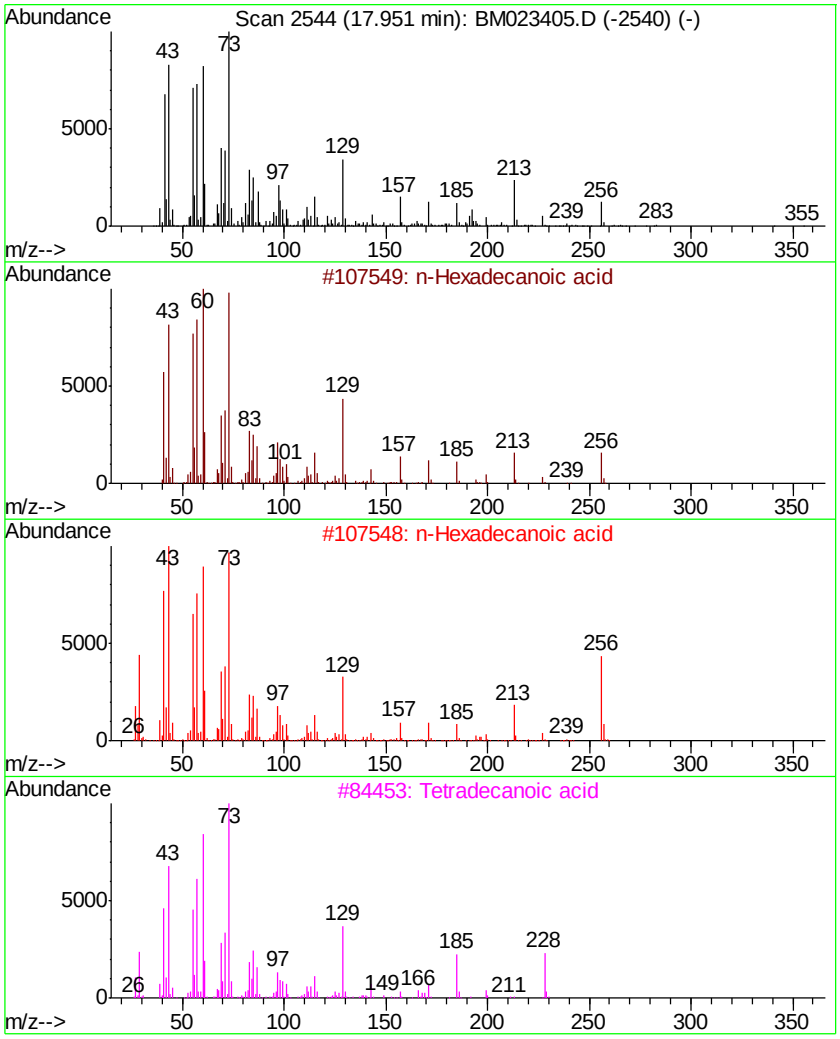
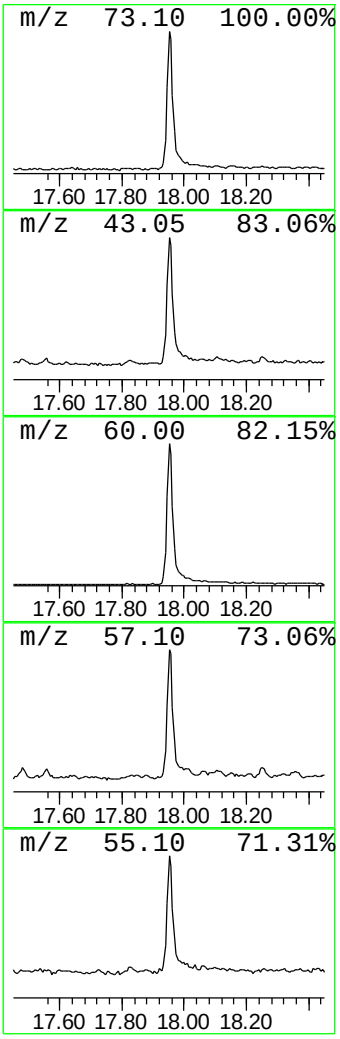
Peak Number 11 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.95	6.62 ng/ul	2062490	Phenanthrene-d10	17.03

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	98
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	94
4	Tridecanoic acid	214	C13H26O2	000638-53-9	93
5	Tridecanoic acid	214	C13H26O2	000638-53-9	93





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102519\
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Acq On : 25 Oct 2019 14:09
Operator : JU
Sample : K5344-05
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BF6L2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
Quant Title : SVOA CALIBRATION

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

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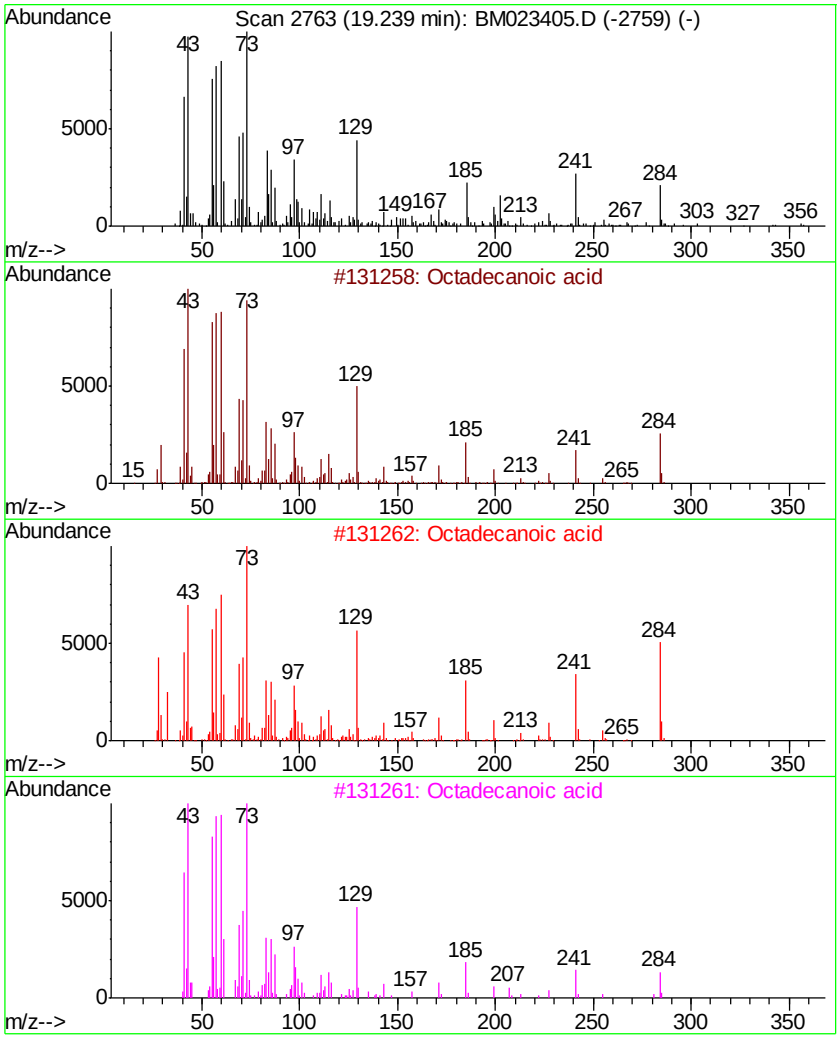
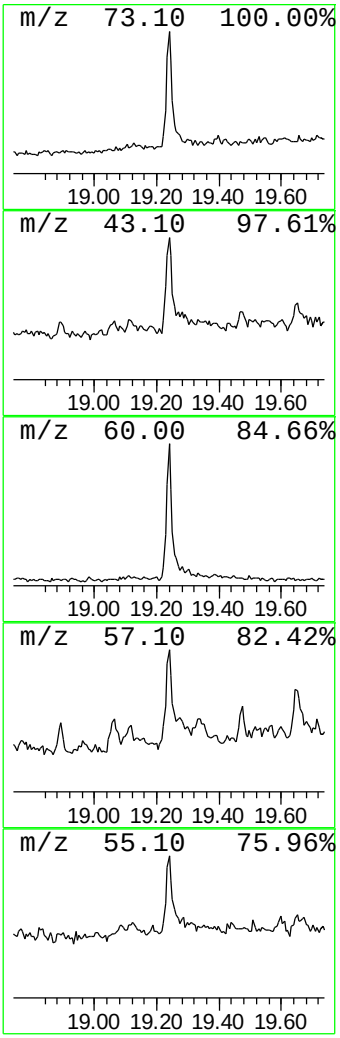
Peak Number 12 Octadecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.24	2.55 ng/ul	656533	Chrysene-d12	21.22

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2	Octadecanoic acid	284	C18H36O2	000057-11-4	99
3	Octadecanoic acid	284	C18H36O2	000057-11-4	97
4	Octadecanoic acid	284	C18H36O2	000057-11-4	95
5	Octadecanoic acid	284	C18H36O2	000057-11-4	91





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Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BF6L2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
Quant Title : SVOA CALIBRATION

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

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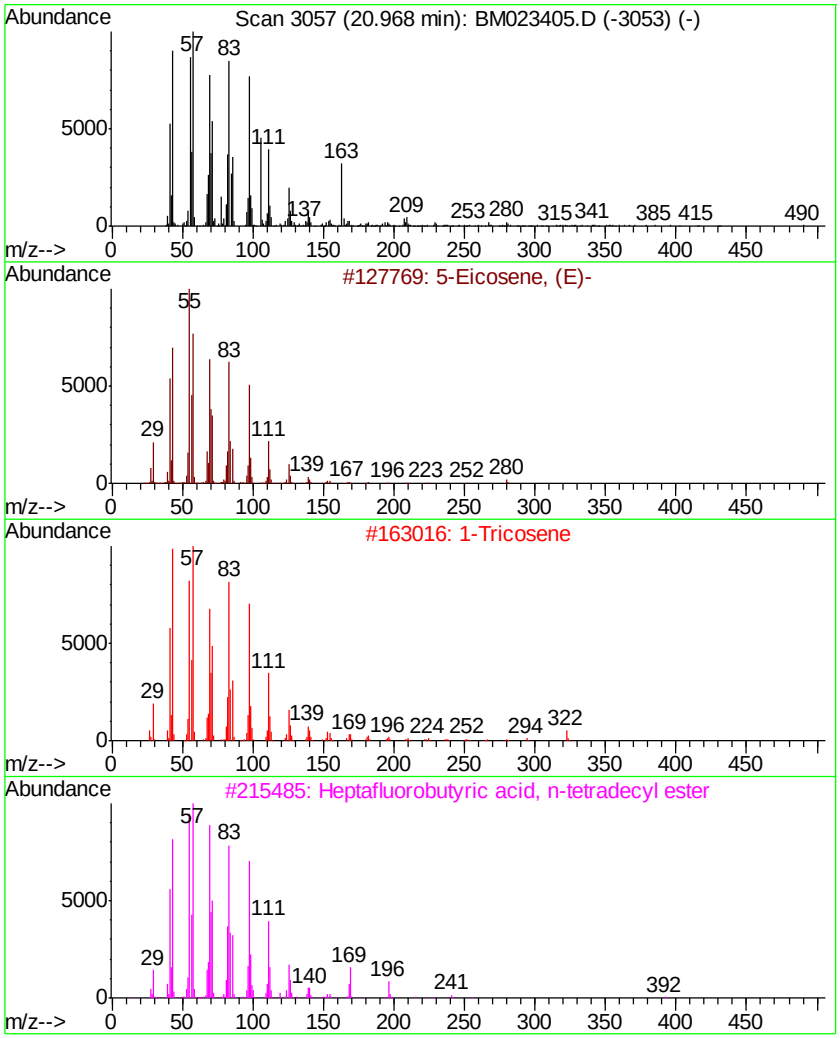
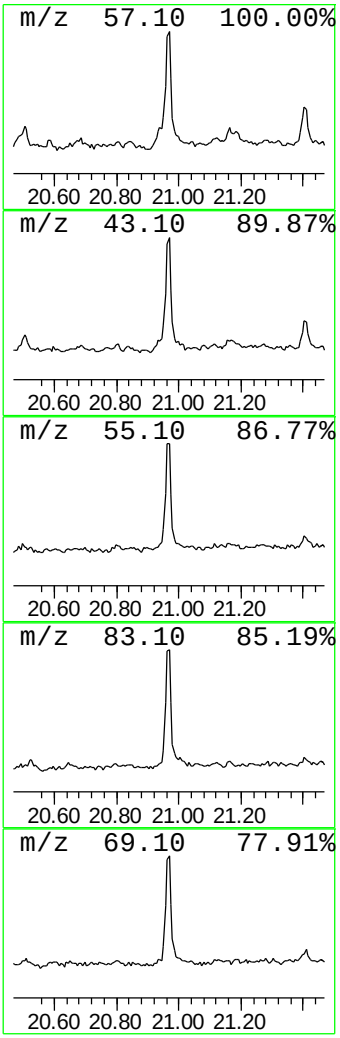
Peak Number 13 (DEL) Alkane: Cyclic20.97 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.97	4.93 ng/ul	1270210	Chrysene-d12	21.22

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Eicosene, (E)-	280	C20H40	074685-30-6	98
2	1-Tricosene	322	C23H46	018835-32-0	98
3	Heptafluorobutyric acid, n-tetra...	410	C18H29F7O2	007365-36-8	94
4	Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	94
5	Behenic alcohol	326	C22H46O	000661-19-8	93





Tentatively Identified Compound (LSC) summary

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Data File : BM023405.D
Acq On : 25 Oct 2019 14:09
Operator : JU
Sample : K5344-05
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BF6L2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.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, 1,3-dime...	5.67	7.0	ng/ul	1185230	1	7.63	3392260	20.0
Benzene, 1-ethyl-...	6.73	2.5	ng/ul	424125	1	7.63	3392260	20.0
Benzene, 1,2,3-tr...	7.29	6.6	ng/ul	1114220	1	7.63	3392260	20.0
Benzene, 1,2,4-tr...	7.75	3.0	ng/ul	500436	1	7.63	3392260	20.0
Indane	8.00	2.5	ng/ul	432201	1	7.63	3392260	20.0
Hexanoic acid, 2-...	8.95	4.0	ng/ul	687139	1	7.63	3392260	20.0
Naphthalene, 1-me...	12.30	2.4	ng/ul	560353	2	10.42	4762360	20.0
n-Hexadecanoic acid	17.95	6.6	ng/ul	2062490	4	17.03	6234040	20.0
Octadecanoic acid	19.24	2.5	ng/ul	656533	5	21.22	5153410	20.0
(DEL) Alkane: Cyc...	20.97	4.9	ng/ul	1270210	5	21.22	5153410	20.0