LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062320\
Data File : BM026507.D
Acq On : 23 Jun 2020 19:38
Operator : JU/CG
Sample : L3027-07
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BG2L0

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-BM061220MA.M
Title : SVOA CALIBRATION

Signal : TIC: BM026510.D

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.969	11	14	24	rVB	27458	42837	1.82%	0.117%
2	3.499	101	104	111	rVB4	10232	14449	0.61%	0.040%
3	3.581	113	118	128	rBV3	16930	36421	1.55%	0.100%
4	3.663	129	132	137	rVB3	9725	12774	0.54%	0.035%
5	4.557	280	284	287	rBV	14606	19536	0.83%	0.053%
6	6.587	622	629	642	rBV	436599	728176	30.97%	1.994%
7	6.740	647	655	667	rBV	336024	536305	22.81%	1.468%
8	6.922	679	686	694	rBV	457282	708199	30.12%	1.939%
9	7.022	700	703	710	rVB8	4009	7583	0.32%	0.021%
10	7.381	758	764	774	rBV	474072	717572	30.52%	1.965%
11	7.675	809	814	823	rBV3	12754	38960	1.66%	0.107%
12	8.104	880	887	902	rBV	565141	904264	38.45%	2.476%
13	8.204	902	904	912	rVB5	8950	15024	0.64%	0.041%
14	8.528	952	959	972	rBV	467807	752684	32.01%	2.061%
15	8.869	1011	1017	1023	rBV	27858	48392	2.06%	0.133%
16	8.981	1032	1036	1041	rVB3	4829	7606	0.32%	0.021%
17	9.239	1069	1080	1090	rBV	397696	641550	27.28%	1.757%
18	9.534	1124	1130	1139	rBV2	20496	43268	1.84%	0.118%
19	9.775	1165	1171	1178	rBV	520562	896674	38.13%	2.455%
20	9.839	1178	1182	1193	rVB	44851	94527	4.02%	0.259%
21	10.139	1224	1233	1242	rBV	683732	1100719	46.81%	3.014%
22	10.292	1252	1259	1277	rBV	427451	793032	33.72%	2.171%
23	10.675	1318	1324	1351	rBV	485674	985093	41.89%	2.697%
24	11.733	1500	1504	1511	rVB2	6643	12458	0.53%	0.034%
25	11.927	1529	1537	1539	rBV3	7495	16144	0.69%	0.044%
26	12.345	1602	1608	1612	rBV4	9318	17598	0.75%	0.048%
27	12.957	1707	1712	1720	rBV5	9701	17888	0.76%	0.049%
28	13.045	1720	1727	1733	rVV5	7147	15941	0.68%	0.044%
29	13.463	1792	1798	1803	rBV	1117607	1497464	63.68%	4.100%
30	13.510	1803	1806	1814	rVB	151481	202688	8.62%	0.555%
31	13.721	1836	1842	1857	rBV	1077870	1535754	65.31%	4.205%
32	14.039	1889	1896	1904	rBV	1077747	1566267	66.61%	4.289%
33	14.280	1931	1937	1956	rBV	334449	708296	30.12%	1.939%
34	14.516	1969	1977	1985	rBV	85924	131242	5.58%	0.359%

LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062320\
Data File : BM026507.D
Acq On : 23 Jun 2020 19:38
Operator : JU/CG
Sample : L3027-07
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BG2L0

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-BM061220MA.M
Title : SVOA CALIBRATION

35	14.621	1991	1995	2002	rVB	34322	52694	2.24%	0.144%
36	15.039	2060	2066	2069	rBV	1323269	1867021	79.40%	5.112%
37	15.068	2069	2071	2081	rVB	608295	707277	30.08%	1.937%
38	15.186	2085	2091	2104	rBV	527176	771075	32.79%	2.111%
39	15.280	2104	2107	2112	rVB3	15776	23659	1.01%	0.065%
40	15.427	2129	2132	2139	rVB6	3644	7328	0.31%	0.020%
41	15.515	2139	2147	2150	rBV2	12585	22257	0.95%	0.061%
42	15.545	2150	2152	2155	rVV2	12304	14013	0.60%	0.038%
43	15.580	2155	2158	2164	rVB	26890	33443	1.42%	0.092%
44	15.768	2183	2190	2194	rBV3	15149	26790	1.14%	0.073%
45	15.815	2194	2198	2205	rVV	99261	132577	5.64%	0.363%
46	15.904	2208	2213	2218	rBV	235692	309701	13.17%	0.848%
47	15.968	2219	2224	2229	rVV3	283902	527214	22.42%	1.444%
48	16.027	2229	2234	2238	rVV2	250865	404559	17.20%	1.108%
49	16.068	2238	2241	2246	rVV	161165	209334	8.90%	0.573%
50	16.127	2246	2251	2252	rVV	92581	125780	5.35%	0.344%
51	16.168	2256	2258	2262	rVV	61850	81860	3.48%	0.224%
52	16.221	2262	2267	2275	rVV4	160962	337976	14.37%	0.925%
53	16.292	2275	2279	2283	rVV	251659	360272	15.32%	0.986%
54	16.333	2283	2286	2288	rVV2	101126	147905	6.29%	0.405%
55	16.362	2288	2291	2299	rVV	168935	229887	9.78%	0.629%
56	16.510	2312	2316	2319	rVV2	10829	14451	0.61%	0.040%
57	16.539	2319	2321	2325	rVV6	7286	12193	0.52%	0.033%
58	16.580	2325	2328	2331	rVV3	10254	17260	0.73%	0.047%
59	16.627	2333	2336	2339	rVV2	14027	20625	0.88%	0.056%
60	16.662	2340	2342	2345	rVV3	7309	10046	0.43%	0.028%
61	16.704	2346	2349	2351	rVV4	10535	11093	0.47%	0.030%
62	16.733	2352	2354	2358	rVB2	6382	8354	0.36%	0.023%
63	16.792	2358	2364	2375	rBV	1479113	2007654	85.38%	5.497%
64	16.892	2375	2381	2397	rVB2	1547058	2111548	89.79%	5.782%
65	17.027	2400	2404	2406	rVV4	9975	11986	0.51%	0.033%
66	17.057	2406	2409	2414	rVB3	12199	16678	0.71%	0.046%
67	17.180	2425	2430	2433	rBV5	7639	11742	0.50%	0.032%
68	17.674	2508	2514	2518	rVB5	3442	6912	0.29%	0.019%
69	17.733	2518	2524	2531	rBV3	57057	107256	4.56%	0.294%
70	17.839	2539	2542	2547	rVB5	4792	7277	0.31%	0.020%
71	17.892	2547	2551	2552	rBV4	9002	11463	0.49%	0.031%

LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062320\
Data File : BM026507.D
Acq On : 23 Jun 2020 19:38
Operator : JU/CG
Sample : L3027-07
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BG2L0

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-BM061220MA.M
Title : SVOA CALIBRATION

72	17.968	2561	2564	2567	rBV5	6740	8349	0.36%	0.023%
73	18.256	2609	2613	2616	rBV5	6142	10150	0.43%	0.028%
74	18.356	2625	2630	2635	rBV9	5994	9985	0.42%	0.027%
75	18.427	2639	2642	2646	rBV4	11444	13850	0.59%	0.038%
76	18.833	2708	2711	2717	rVB	16357	21260	0.90%	0.058%
77	18.933	2723	2728	2733	rVB3	27027	34360	1.46%	0.094%
78	19.027	2740	2744	2750	rBV3	31060	46728	1.99%	0.128%
79	19.086	2750	2754	2758	rVV	17042	24613	1.05%	0.067%
80	19.198	2767	2773	2780	rBV2	1744821	2344963	99.72%	6.421%
81	19.592	2836	2840	2843	rBV	29088	35097	1.49%	0.096%
82	19.627	2843	2846	2850	rVB	52127	56924	2.42%	0.156%
83	20.174	2935	2939	2942	rBV5	10796	18169	0.77%	0.050%
84	20.298	2957	2960	2963	rBV2	26663	27569	1.17%	0.075%
85	20.568	3003	3006	3009	rBV	64637	70730	3.01%	0.194%
86	20.603	3009	3012	3015	rVB	149778	142259	6.05%	0.390%
87	20.762	3034	3039	3052	rBV	202874	347443	14.78%	0.951%
88	21.003	3075	3080	3088	rBV	1803621	2351524	100.00%	6.439%
89	21.074	3089	3092	3097	rVB3	40651	44740	1.90%	0.123%
90	21.197	3110	3113	3116	rBV2	70309	71934	3.06%	0.197%
91	21.433	3150	3153	3155	rBV	64488	59901	2.55%	0.164%
92	21.462	3155	3158	3162	rVB	124857	133616	5.68%	0.366%
93	21.615	3181	3184	3192	rBV2	85663	103101	4.38%	0.282%
94	22.044	3254	3257	3262	rVB	118776	136957	5.82%	0.375%
95	22.333	3303	3306	3310	rVB	125530	146604	6.23%	0.401%
96	22.509	3333	3336	3341	rVB	137720	180894	7.69%	0.495%
97	22.962	3407	3413	3420	rBV	1146377	1846178	78.51%	5.055%
98	23.027	3420	3424	3429	rVV	175986	247194	10.51%	0.677%
99	23.091	3429	3435	3449	rVB	1214978	2112945	89.85%	5.786%
100	23.609	3519	3523	3531	rVB	133621	226580	9.64%	0.620%

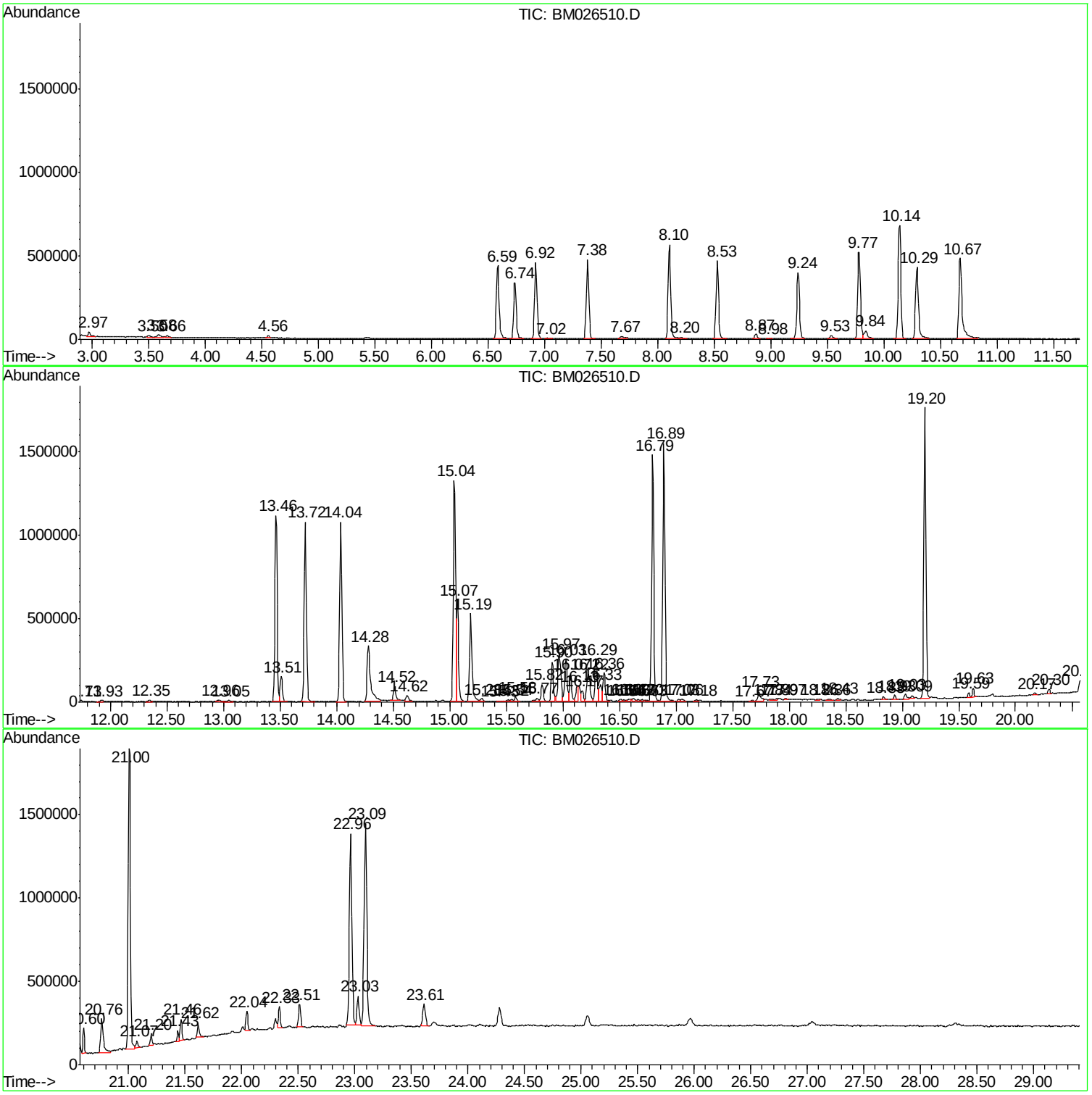
Sum of corrected areas: 36521162

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062320\
Data File : BM026507.D
Acq On : 23 Jun 2020 19:38
Operator : JU/CG
Sample : L3027-07
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BG2L0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.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 M\DATA\BM062320\
Data File : BM026507.D
Acq On : 23 Jun 2020 19:38
Operator : JU/CG
Sample : L3027-07
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BG2L0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.M
Quant Title : SVOA CALIBRATION

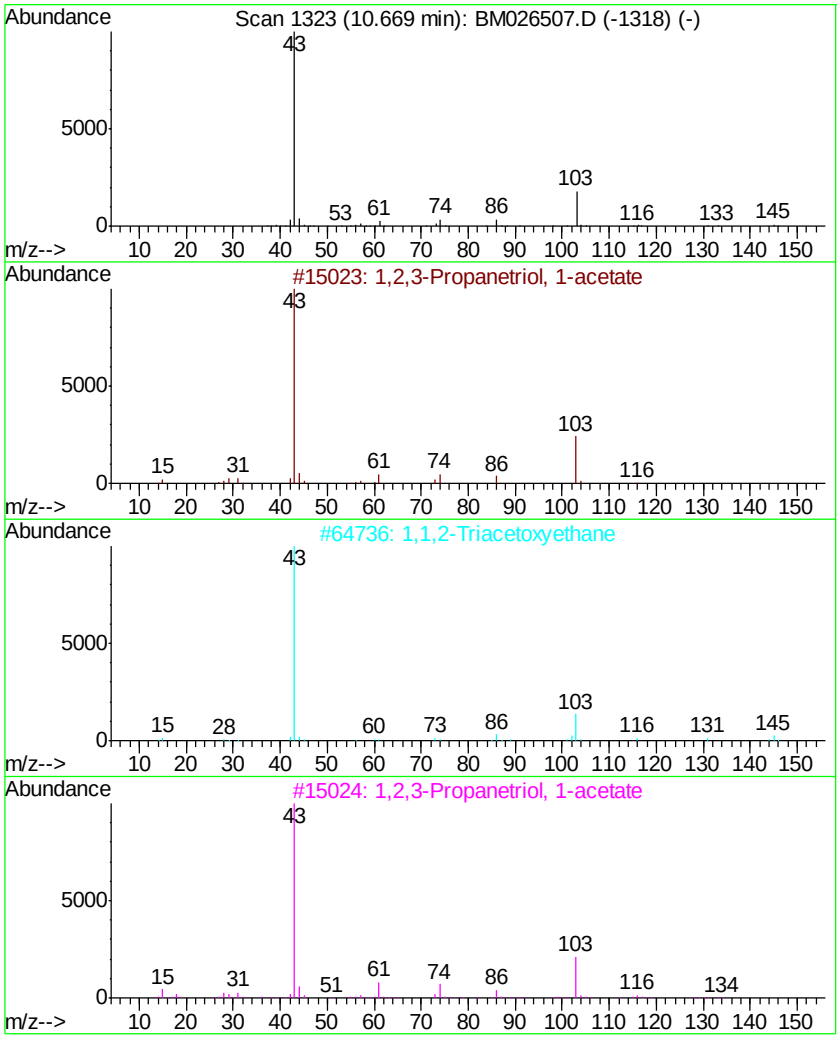
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

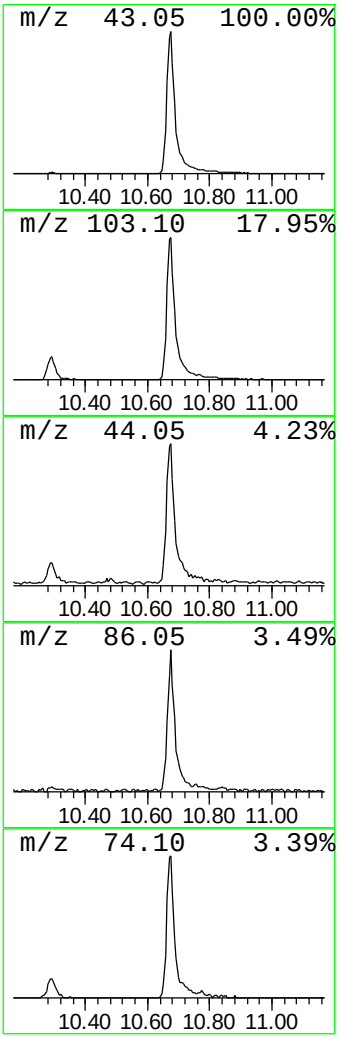
****

Peak Number 1 1,2,3-Propanetriol, 1-acetate Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.67	17.90 ng/ul	985093	Naphthalene-d8	10.14

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,2,3-Propanetriol, 1-acetate	134	C5H10O4	000106-61-6	53
2	1,1,2-Triacetoxyethane	204	C8H12O6	002983-35-9	38
3	1,2,3-Propanetriol, 1-acetate	134	C5H10O4	000106-61-6	36
4	Glycerol 1,2-diacetate	176	C7H12O5	000102-62-5	9
5	1,2,3-Propanetriol, 1-acetate	134	C5H10O4	000106-61-6	9





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062320\
Data File : BM026507.D
Acq On : 23 Jun 2020 19:38
Operator : JU/CG
Sample : L3027-07
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BG2L0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.M
Quant Title : SVOA CALIBRATION

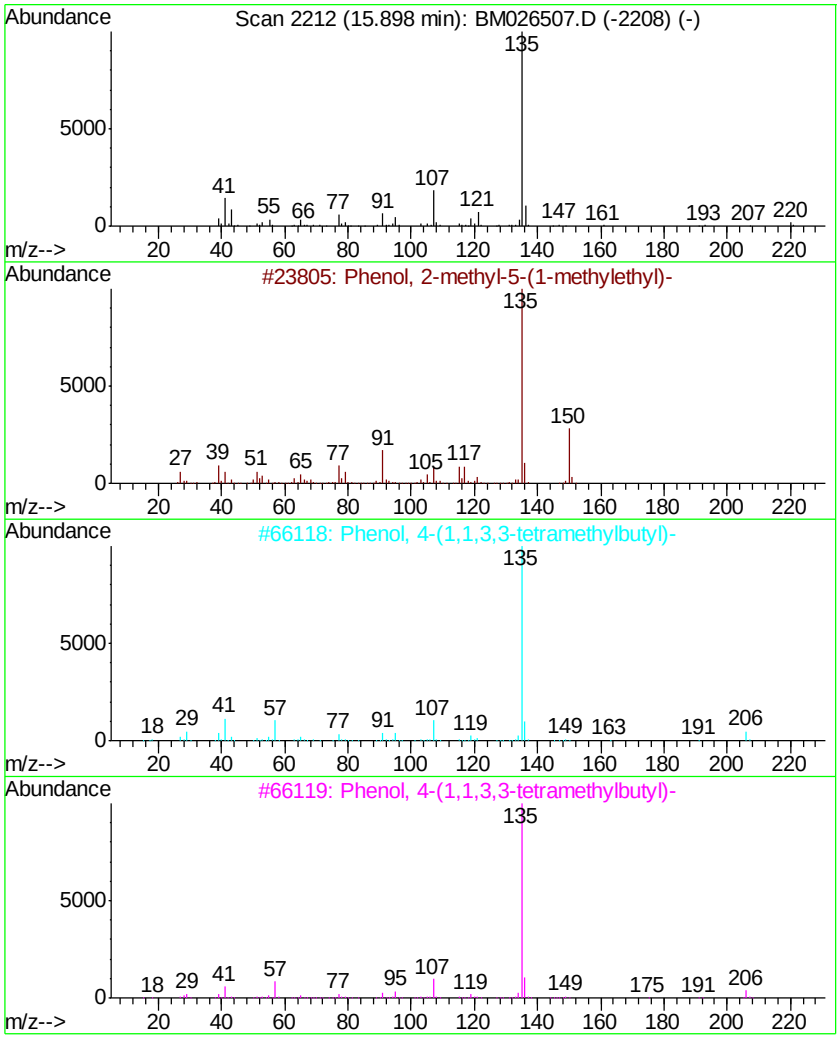
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

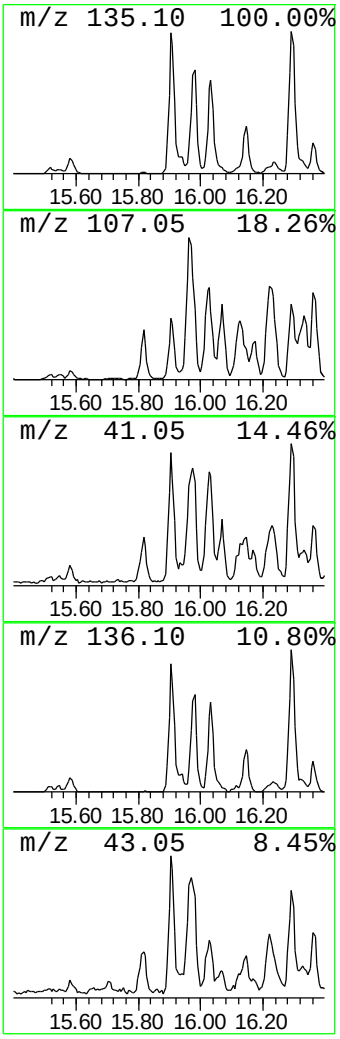
****

Peak Number 2 Phenol, 2-methyl-5-(1-methy... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.90	3.09 ng/ul	309701	Phenanthrene-d10	16.79

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2-methyl-5-(1-methylethyl)-	150	C10H14O	000499-75-2	78
2	Phenol, 4-(1,1,3,3-tetramethylbu...	206	C14H22O	000140-66-9	78
3	Phenol, 4-(1,1,3,3-tetramethylbu...	206	C14H22O	000140-66-9	78
4	4-(1,1-Dimethylpropyl)phenyl ace...	206	C13H18O2	1000373-45-3	78
5	4-tert-Butylphenyl acetate	192	C12H16O2	003056-64-2	64





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062320\
Data File : BM026507.D
Acq On : 23 Jun 2020 19:38
Operator : JU/CG
Sample : L3027-07
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BG2L0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.M
Quant Title : SVOA CALIBRATION

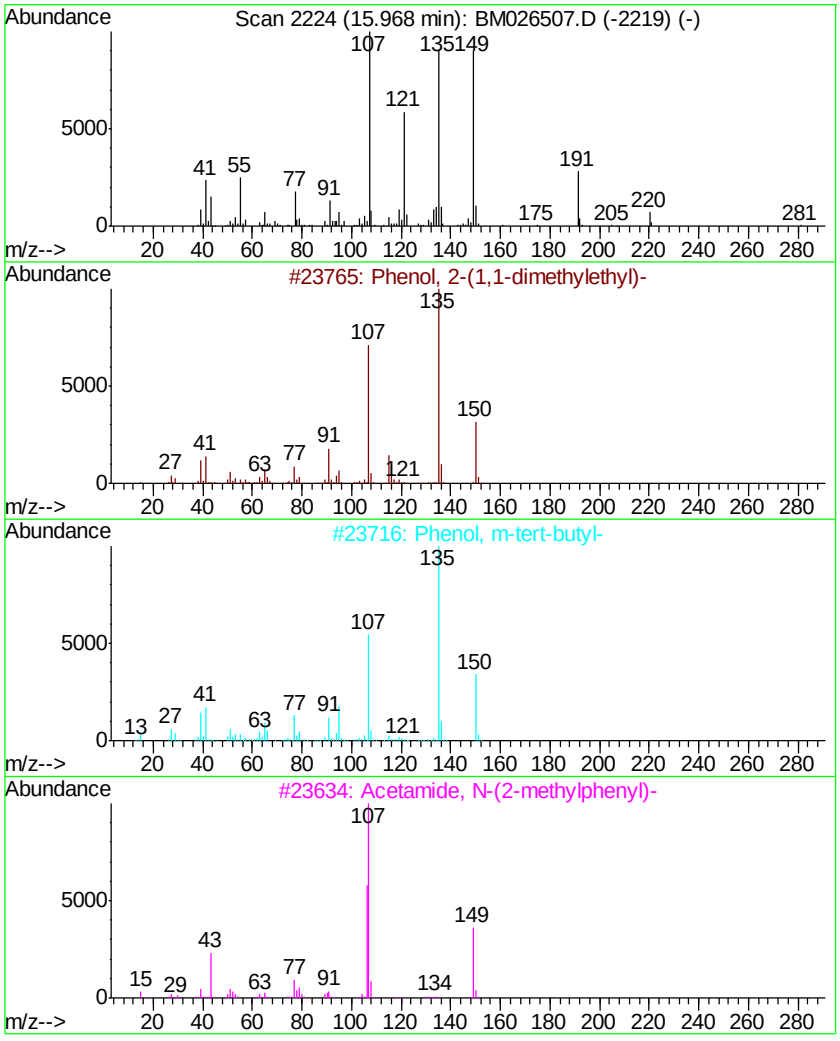
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

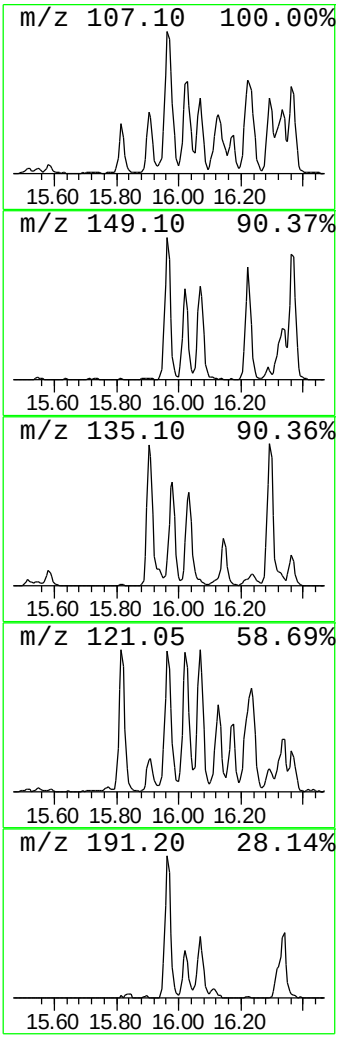
****

Peak Number 3 unknown-01 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.97	5.25 ng/ul	527214	Phenanthrene-d10	16.79

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2-(1,1-dimethylethyl)-	150	C10H14O	000088-18-6	46
2	Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	43
3	Acetamide, N-(2-methylphenyl)-	149	C9H11NO	000120-66-1	43
4	Acetamide, N-methyl-N-phenyl-	149	C9H11NO	000579-10-2	30
5	Acetamide, N-(4-methylphenyl)-	149	C9H11NO	000103-89-9	25





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062320\
Data File : BM026507.D
Acq On : 23 Jun 2020 19:38
Operator : JU/CG
Sample : L3027-07
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BG2L0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.M
Quant Title : SVOA CALIBRATION

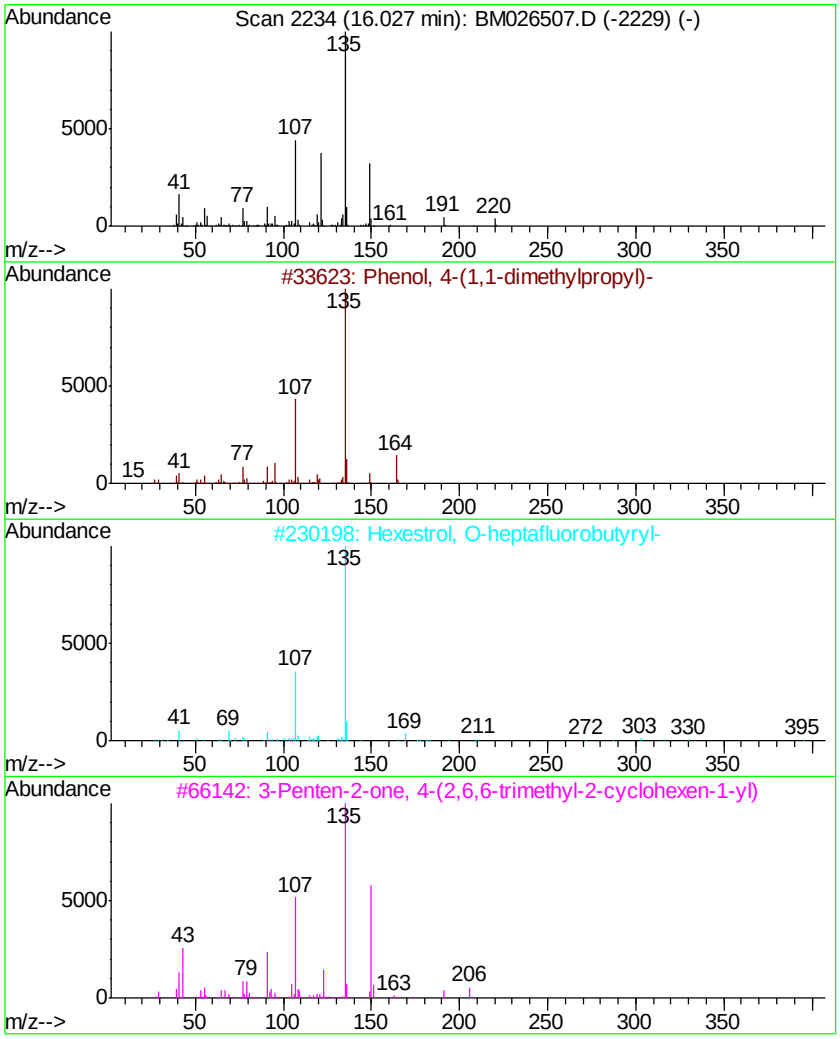
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

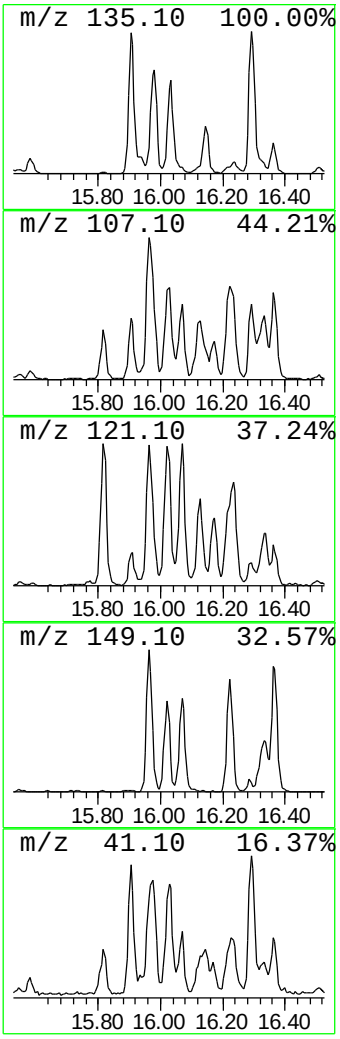
****

Peak Number 4 Phenol, 4-(1,1-dimethylprop... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.03	4.03 ng/ul	404559	Phenanthrene-d10	16.79

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	59
2	Hexestrol, O-heptafluorobutyryl-	466	C22H21F7O3	1000365-31-9	59
3	3-Penten-2-one, 4-(2,6,6-trimeth...	206	C14H22O	114933-28-7	59
4	3,3'-Dimethoxybenzil	270	C16H14O4	040101-17-5	53
5	Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	53





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062320\
Data File : BM026507.D
Acq On : 23 Jun 2020 19:38
Operator : JU/CG
Sample : L3027-07
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BG2L0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.M
Quant Title : SVOA CALIBRATION

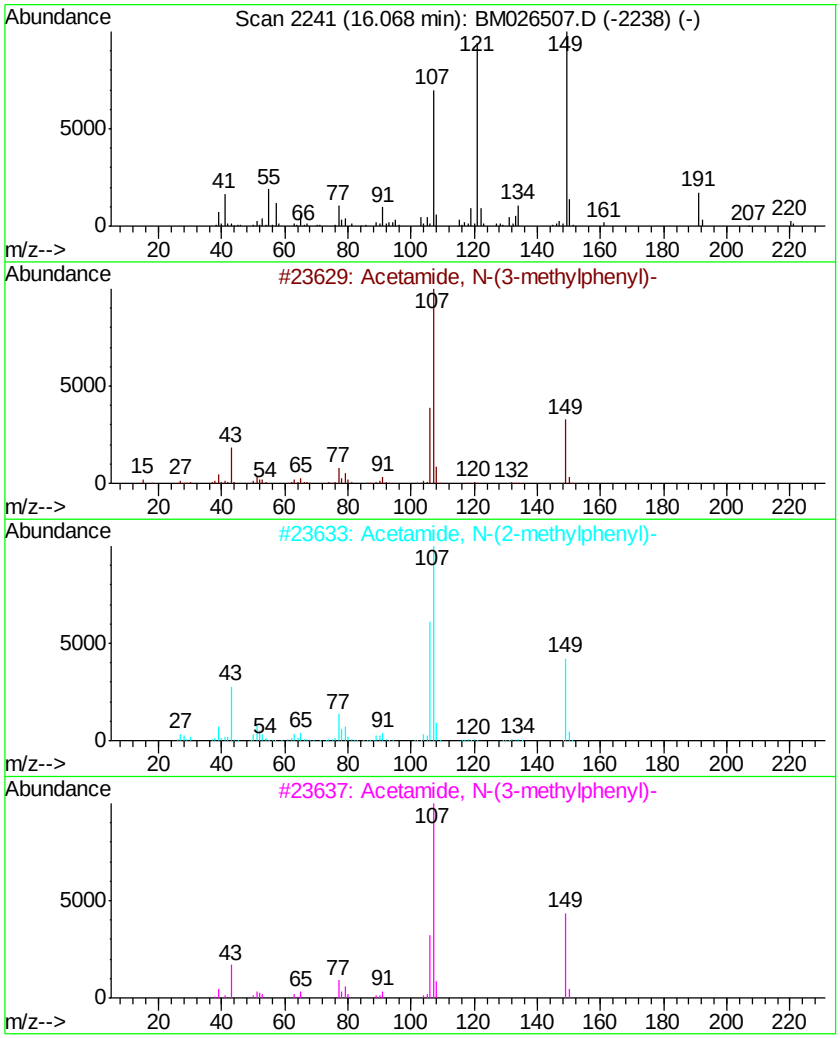
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

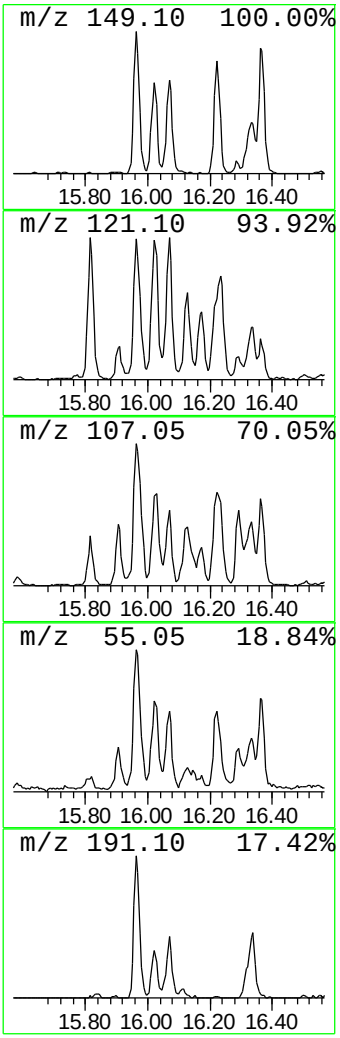
****

Peak Number 5 unknown-02 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.07	2.09 ng/ul	209334	Phenanthrene-d10	16.79

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetamide, N-(3-methylphenyl)-	149	C9H11NO	000537-92-8	45
2	Acetamide, N-(2-methylphenyl)-	149	C9H11NO	000120-66-1	45
3	Acetamide, N-(3-methylphenyl)-	149	C9H11NO	000537-92-8	43
4	Acetamide, N-(2-methylphenyl)-	149	C9H11NO	000120-66-1	38
5	Isoxazole, 4,5-dihydro-5-[(4-met...	191	C11H13NO2	1000362-14-0	38





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Data File : BM026507.D
Acq On : 23 Jun 2020 19:38
Operator : JU/CG
Sample : L3027-07
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BG2L0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.M
Quant Title : SVOA CALIBRATION

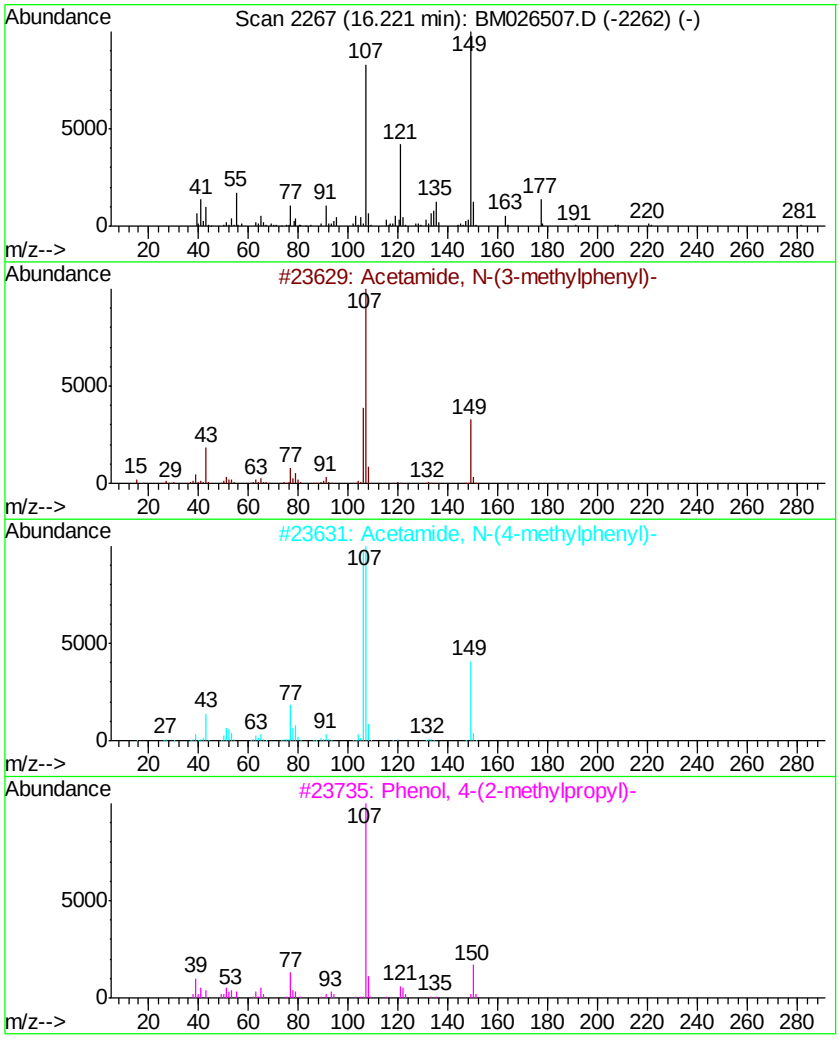
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

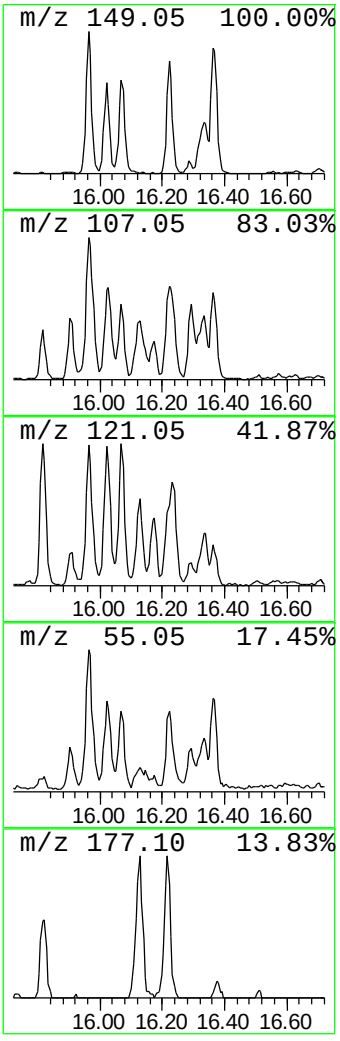
****

Peak Number 6 unknown-03 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.22	3.37 ng/ul	337976	Phenanthrene-d10	16.79

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetamide, N-(3-methylphenyl)-	149	C9H11NO	000537-92-8	64
2	Acetamide, N-(4-methylphenyl)-	149	C9H11NO	000103-89-9	43
3	Phenol, 4-(2-methylpropyl)-	150	C10H14O	004167-74-2	41
4	Phenol, 2-methyl-4-(1,1,3,3-tetr...	220	C15H24O	002219-84-3	38
5	2-Methyl-2-(4'-methoxyphenyl)pen...	206	C13H18O2	185700-49-6	38





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Data File : BM026507.D
Acq On : 23 Jun 2020 19:38
Operator : JU/CG
Sample : L3027-07
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BG2L0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.M
Quant Title : SVOA CALIBRATION

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

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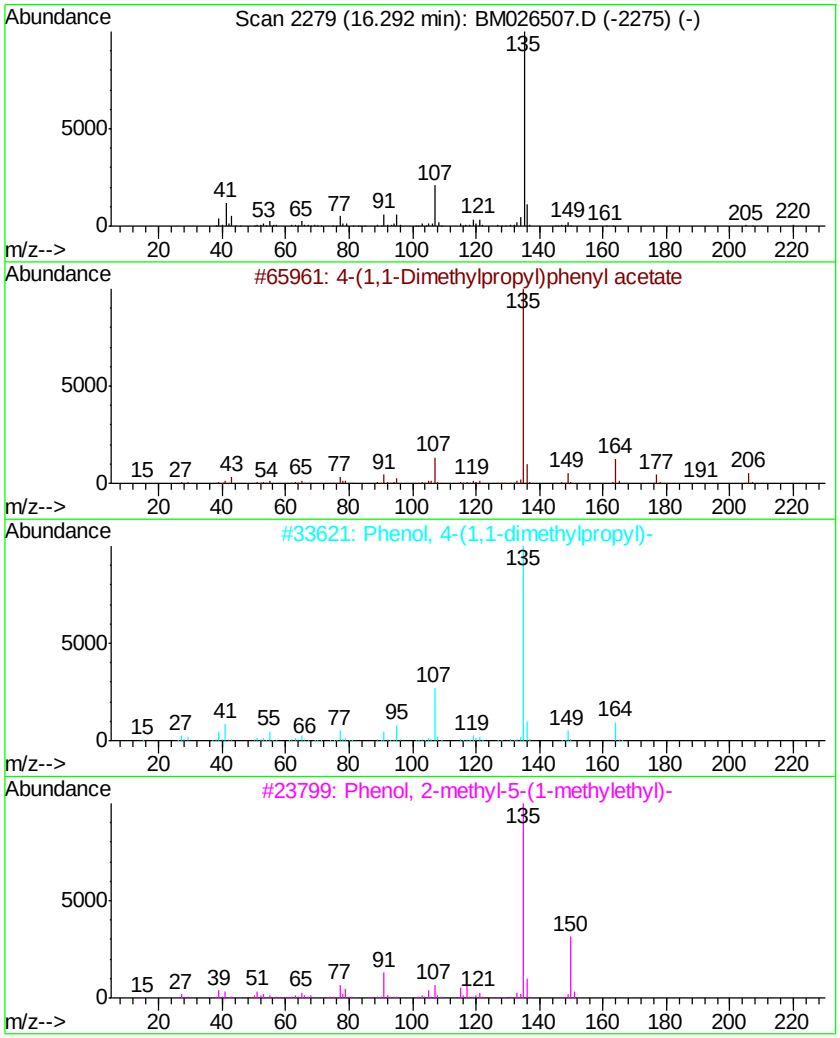
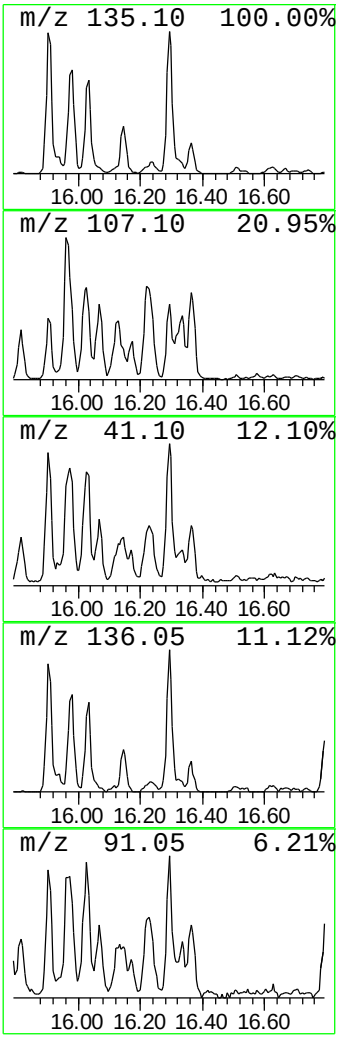
Peak Number 7 4-(1,1-Dimethylpropyl)pheny... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.29	3.59 ng/ul	360272	Phenanthrene-d10	16.79

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-(1,1-Dimethylpropyl)phenyl ace...	206	C13H18O2	1000373-45-3	83
2	Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	78
3	Phenol, 2-methyl-5-(1-methylethyl)-	150	C10H14O	000499-75-2	72
4	((1,2-Diethylethylene)bis(p-phen...	354	C22H26O4	004547-76-6	72
5	Hexestrol, O-heptafluorobutyryl-	466	C22H21F7O3	1000365-31-9	72





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Data File : BM026507.D
Acq On : 23 Jun 2020 19:38
Operator : JU/CG
Sample : L3027-07
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BG2L0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.M
Quant Title : SVOA CALIBRATION

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

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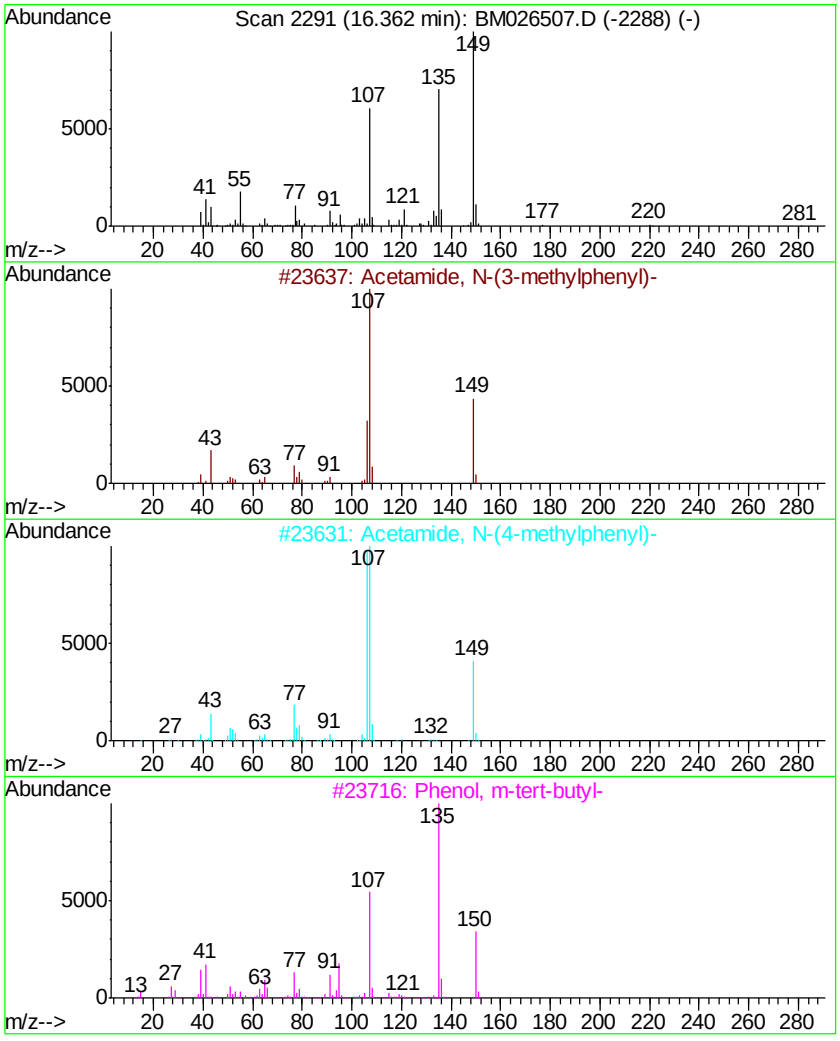
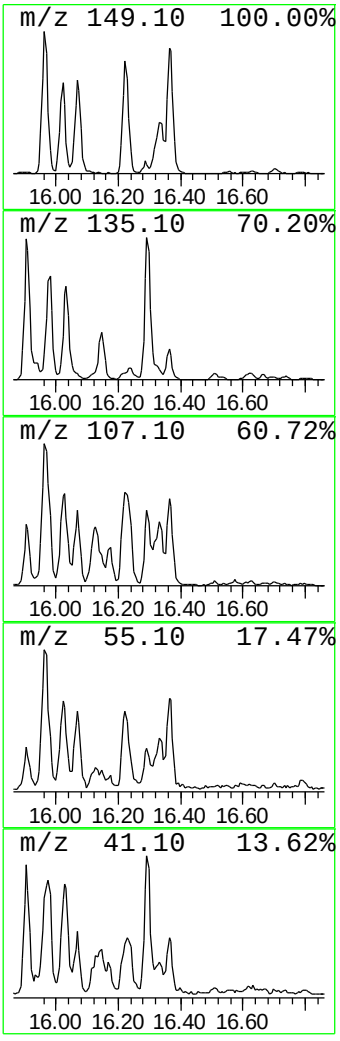
Peak Number 8 Acetamide, N-(3-methylphenyl)- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.36	2.29 ng/ul	229887	Phenanthrene-d10	16.79

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetamide, N-(3-methylphenyl)-	149	C9H11NO	000537-92-8	52
2	Acetamide, N-(4-methylphenyl)-	149	C9H11NO	000103-89-9	43
3	Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	38
4	Phenol, 2-(1,1-dimethylethyl)-	150	C10H14O	000088-18-6	38
5	5H-Pyrrolo(3,2-d)pyrimidine-2,4-...	149	C6H7N5	1000244-21-4	38





Library Search Compound Report

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Data File : BM026507.D
Acq On : 23 Jun 2020 19:38
Operator : JU/CG
Sample : L3027-07
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BG2L0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.M
Quant Title : SVOA CALIBRATION

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

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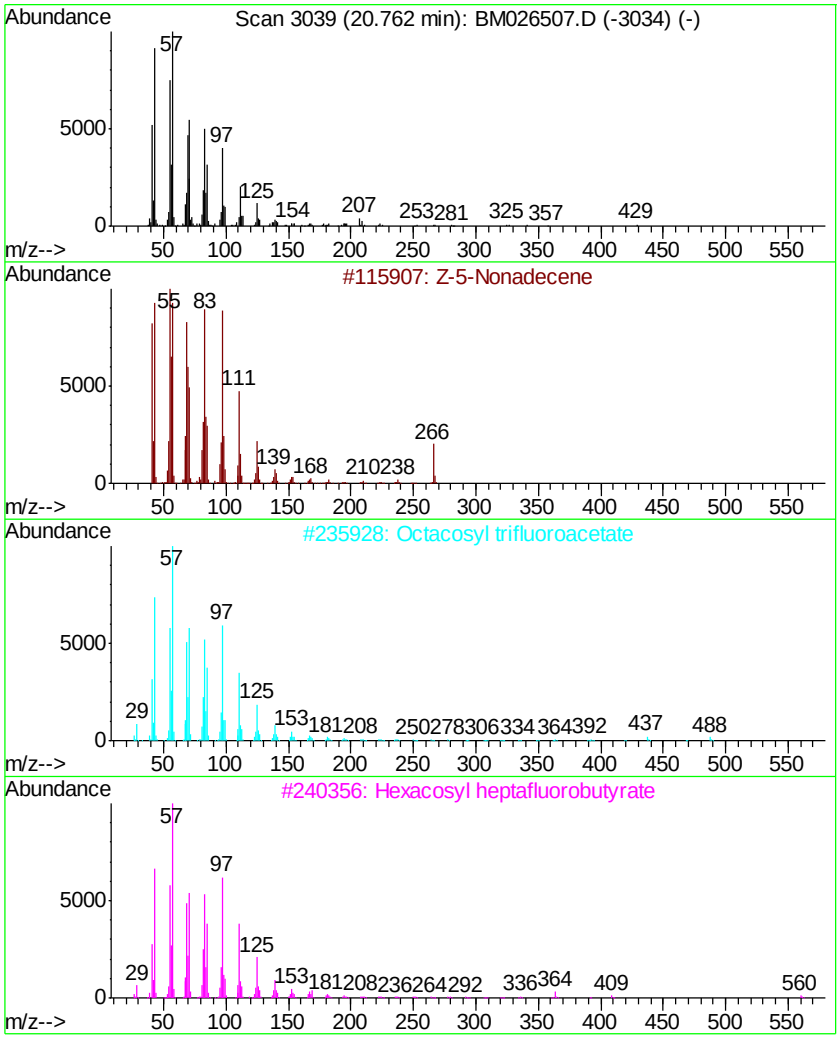
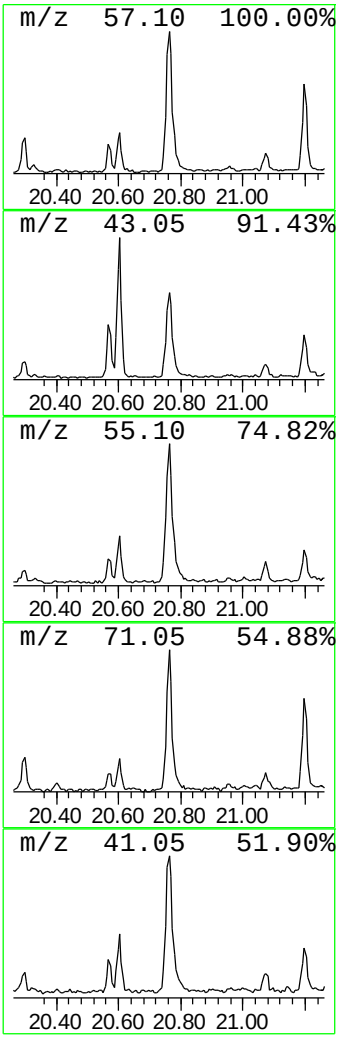
Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.76	2.96 ng/ul	347443	Chrysene-d12	21.00

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Z-5-Nonadecene	266	C19H38	1000131-11-8	92
2	Octacosyl trifluoroacetate	506	C30H57F3O2	1000351-74-9	91
3	Hexacosyl heptafluorobutyrate	578	C30H53F7O2	1000351-83-3	91
4	Oxalic acid, isobutyl octadecyl ...	398	C24H46O4	1000309-38-3	91
5	Oxalic acid, isobutyl hexadecyl ...	370	C22H42O4	1000309-38-1	91





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM062320\
Data File : BM026507.D
Acq On : 23 Jun 2020 19:38
Operator : JU/CG
Sample : L3027-07
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BG2L0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.M
Quant Title : SVOA CALIBRATION

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

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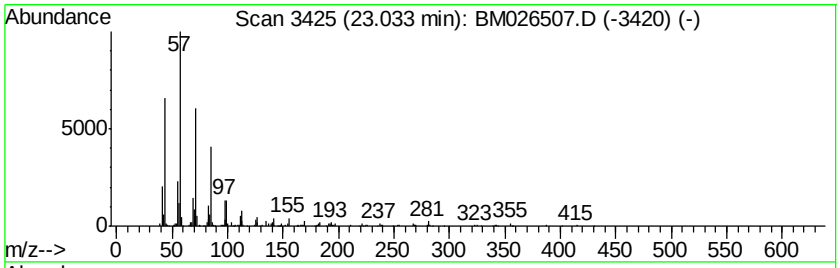
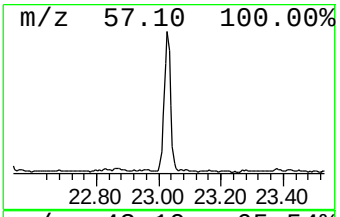
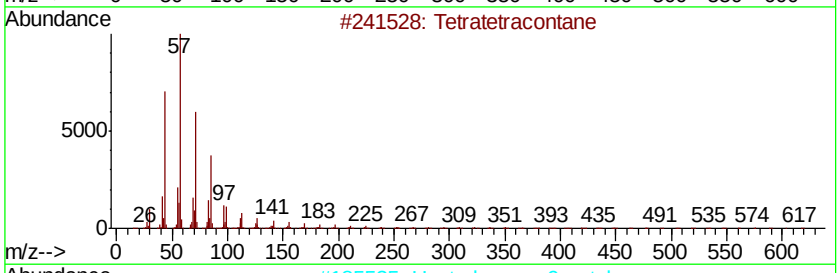
Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 8

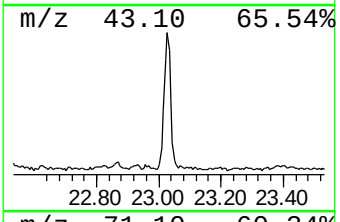
R.T.	EstConc	Area	Relative to ISTD	R.T.
23.03	2.34 ng/ul	247194	Perylene-d12	23.09

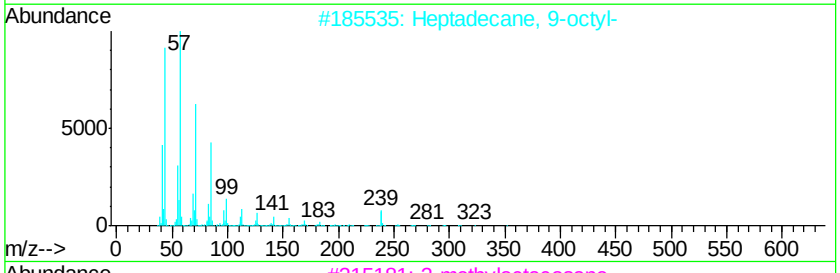
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetratetracontane	619	C44H90	007098-22-8	90
2	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	90
3	2-methyloctacosane	408	C29H60	1000376-72-8	90
4	Octacosane	394	C28H58	000630-02-4	90
5	Docosane, 11-butyl-	366	C26H54	013475-76-8	90

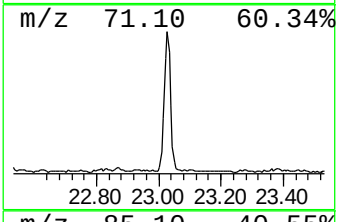


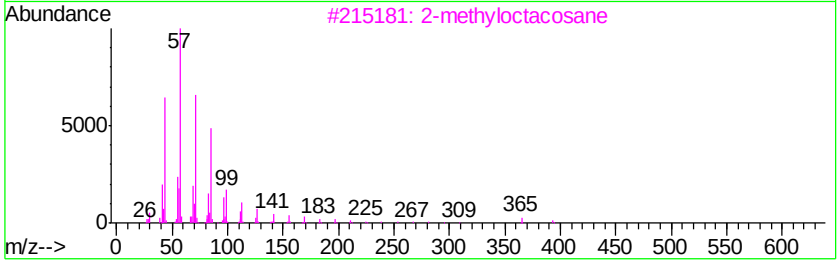


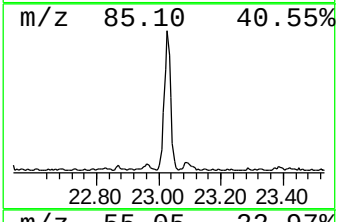


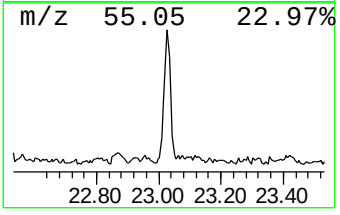












Library Search Compound Report

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Data File : BM026507.D
Acq On : 23 Jun 2020 19:38
Operator : JU/CG
Sample : L3027-07
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BG2L0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.M
Quant Title : SVOA CALIBRATION

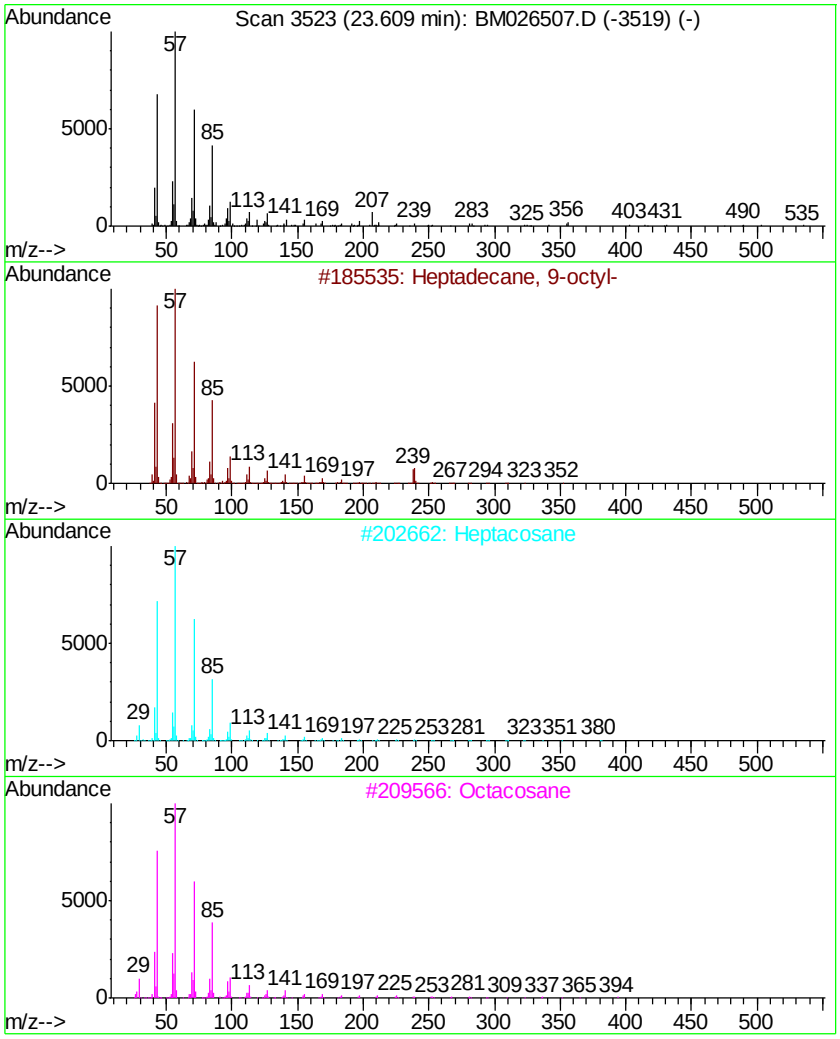
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

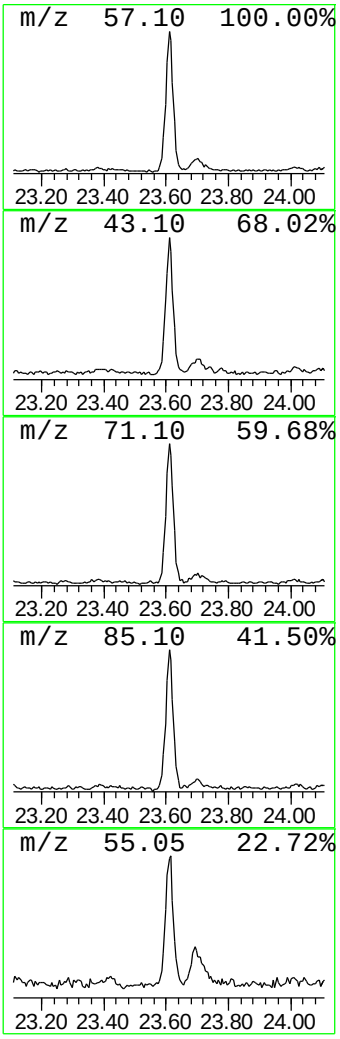
****

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.61	2.14 ng/ul	226580	Perylene-d12	23.09

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	90
2	Heptacosane	380	C27H56	000593-49-7	90
3	Octacosane	394	C28H58	000630-02-4	90
4	Eicosane, 9-octyl-	394	C28H58	013475-77-9	87
5	Tetracosane	338	C24H50	000646-31-1	87





Tentatively Identified Compound (LSC) summary

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Data File : BM026507.D
Acq On : 23 Jun 2020 19:38
Operator : JU/CG
Sample : L3027-07
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BG2L0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM061220MA.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
1,2,3-Propanetrio...	10.67	17.9	ng/ul	985093	2	10.14	1100720	20.0
Phenol, 2-methyl-...	15.90	3.1	ng/ul	309701	4	16.79	2007650	20.0
unknown-01	15.97	5.3	ng/ul	527214	4	16.79	2007650	20.0
Phenol, 4-(1,1-di...	16.03	4.0	ng/ul	404559	4	16.79	2007650	20.0
unknown-02	16.07	2.1	ng/ul	209334	4	16.79	2007650	20.0
unknown-03	16.22	3.4	ng/ul	337976	4	16.79	2007650	20.0
4-(1,1-Dimethylpr...	16.29	3.6	ng/ul	360272	4	16.79	2007650	20.0
Acetamide, N-(3-m...	16.36	2.3	ng/ul	229887	4	16.79	2007650	20.0
(DEL) Alkane: Str...	20.76	3.0	ng/ul	347443	5	21.00	2351520	20.0
(DEL) Alkane: Str...	23.03	2.3	ng/ul	247194	6	23.09	2112950	20.0
(DEL) Alkane: Str...	23.61	2.1	ng/ul	226580	6	23.09	2112950	20.0