LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM060420\
Data File : BM026268.D
Acq On : 04 Jun 2020 21:40
Operator : JU/CG
Sample : L2844-01
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0AE5

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-BM051320MA.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	3.052	25	28	34	rVV	46749	59249	1.56%	0.124%
2	3.116	35	39	46	rVB	27492	44223	1.16%	0.092%
3	4.963	349	353	364	rVB	248775	350185	9.20%	0.732%
4	5.510	440	446	451	rBV	49008	76462	2.01%	0.160%
5	6.663	631	642	654	rBV2	177574	347435	9.13%	0.726%
6	6.816	662	668	675	rBV	530845	825379	21.68%	1.726%
7	7.004	693	700	709	rBV	597977	908892	23.87%	1.900%
8	7.134	717	722	727	rBV2	25920	40308	1.06%	0.084%
9	7.463	772	778	786	rVV	515194	802121	21.07%	1.677%
10	7.551	786	793	795	rVV6	16972	37696	0.99%	0.079%
11	7.575	795	797	805	rVV2	25827	40932	1.08%	0.086%
12	7.810	830	837	845	rVV2	30345	68602	1.80%	0.143%
13	8.181	894	900	908	rBV	462863	703093	18.47%	1.470%
14	8.463	941	948	956	rBV	53200	80381	2.11%	0.168%
15	8.604	966	972	982	rVV	654607	1049662	27.57%	2.195%
16	8.716	987	991	996	rVB	25666	40976	1.08%	0.086%
17	9.075	1045	1052	1058	rBV2	20113	31725	0.83%	0.066%
18	9.322	1085	1094	1108	rBV	568340	956194	25.11%	1.999%
19	9.486	1116	1122	1129	rVV	47664	82083	2.16%	0.172%
20	9.633	1140	1147	1152	rVV3	18417	35354	0.93%	0.074%
21	9.686	1152	1156	1163	rVB5	16630	28888	0.76%	0.060%
22	9.857	1178	1185	1196	rBV	777284	1294562	34.00%	2.707%
23	10.063	1215	1220	1225	rBV7	15518	27790	0.73%	0.058%
24	10.122	1225	1230	1236	rBV2	24184	41823	1.10%	0.087%
25	10.222	1240	1247	1253	rVV	696635	1087726	28.57%	2.274%
26	10.269	1253	1255	1262	rVB	28770	41378	1.09%	0.087%
27	10.381	1265	1274	1279	rBV2	32651	73056	1.92%	0.153%
28	10.639	1307	1318	1324	rBV8	15712	46057	1.21%	0.096%
29	10.833	1348	1351	1362	rVV8	14536	34570	0.91%	0.072%
30	11.210	1411	1415	1420	rBV7	18153	38557	1.01%	0.081%
31	11.322	1430	1434	1442	rBV5	14655	28424	0.75%	0.059%
32	11.398	1442	1447	1452	rBV	103519	158361	4.16%	0.331%
33	11.498	1458	1464	1470	rVB7	19246	38477	1.01%	0.080%
34	11.563	1470	1475	1486	rBV5	48353	104415	2.74%	0.218%

LSC Area Percent Report

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Operator : JU/CG
Sample : L2844-01
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0AE5

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-BM051320MA.M
Title : SVOA CALIBRATION

35	11.792	1503	1514	1516	rBV9	23450	59972	1.58%	0.125%
36	11.822	1516	1519	1524	rVV3	42124	69687	1.83%	0.146%
37	11.869	1524	1527	1531	rVV4	24549	39033	1.03%	0.082%
38	11.933	1533	1538	1545	rVB9	17476	34549	0.91%	0.072%
39	12.075	1555	1562	1568	rVB3	21364	52428	1.38%	0.110%
40	12.163	1573	1577	1580	rVV5	19368	37443	0.98%	0.078%
41	12.216	1580	1586	1589	rVV6	31706	74095	1.95%	0.155%
42	12.351	1605	1609	1614	rVB	36206	49726	1.31%	0.104%
43	12.592	1642	1650	1655	rBV2	134524	227200	5.97%	0.475%
44	12.663	1656	1662	1665	rVV3	28778	59375	1.56%	0.124%
45	12.739	1671	1675	1682	rVB9	16891	32381	0.85%	0.068%
46	12.933	1701	1708	1719	rBV5	33094	102990	2.71%	0.215%
47	13.163	1741	1747	1753	rBV	368715	519348	13.64%	1.086%
48	13.257	1757	1763	1766	rBV3	19673	37379	0.98%	0.078%
49	13.292	1766	1769	1775	rVB2	24762	35218	0.92%	0.074%
50	13.474	1795	1800	1803	rBV8	16218	30716	0.81%	0.064%
51	13.533	1804	1810	1815	rVV	1466803	1979247	51.98%	4.138%
52	13.580	1815	1818	1823	rVV	275340	353636	9.29%	0.739%
53	13.798	1843	1855	1865	rBV	1611056	2580857	67.79%	5.396%
54	14.110	1902	1908	1914	rVV	992470	1482039	38.93%	3.099%
55	14.163	1914	1917	1925	rVB3	42987	67225	1.77%	0.141%
56	14.345	1943	1948	1961	rVB2	105010	252733	6.64%	0.528%
57	14.480	1966	1971	1976	rBV2	61692	116864	3.07%	0.244%
58	14.680	2000	2005	2009	rVB4	30963	58722	1.54%	0.123%
59	14.733	2011	2014	2019	rVB7	22865	30480	0.80%	0.064%
60	14.786	2019	2023	2027	rBV3	37105	52597	1.38%	0.110%
61	14.963	2045	2053	2057	rBV7	28328	67105	1.76%	0.140%
62	15.110	2072	2078	2081	rBV	2017932	3050659	80.13%	6.378%
63	15.139	2081	2083	2091	rVV	1663732	1844065	48.43%	3.855%
64	15.245	2095	2101	2115	rVB	728193	1028223	27.01%	2.150%
65	15.386	2121	2125	2132	rVB2	255849	398838	10.48%	0.834%
66	15.651	2165	2170	2174	rVB	109338	147307	3.87%	0.308%
67	15.727	2180	2183	2186	rVV	31731	40525	1.06%	0.085%
68	15.786	2190	2193	2200	rVB6	32543	72534	1.91%	0.152%
69	15.974	2221	2225	2231	rBV	59390	101294	2.66%	0.212%
70	16.033	2231	2235	2242	rVV3	71283	154903	4.07%	0.324%
71	16.092	2242	2245	2249	rVB2	48925	63757	1.67%	0.133%

LSC Area Percent Report

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Sample : L2844-01
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0AE5

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
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Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Title : SVOA CALIBRATION

72	16.180	2255	2260	2264	rBV3	44036	72380	1.90%	0.151%
73	16.239	2266	2270	2274	rVB	160092	198760	5.22%	0.416%
74	16.286	2275	2278	2280	rBV3	31646	41559	1.09%	0.087%
75	16.357	2287	2290	2294	rVB	61528	72046	1.89%	0.151%
76	16.433	2299	2303	2307	rVB2	65417	84420	2.22%	0.176%
77	16.480	2308	2311	2315	rVB	32267	39845	1.05%	0.083%
78	16.680	2342	2345	2350	rVV	60379	79793	2.10%	0.167%
79	16.745	2352	2356	2363	rVB2	46143	66186	1.74%	0.138%
80	16.862	2369	2376	2384	rVB	1213896	1750130	45.97%	3.659%
81	16.957	2386	2392	2400	rBV	2228952	3213200	84.39%	6.718%
82	17.057	2404	2409	2417	rBV	2583861	3088271	81.11%	6.457%
83	17.245	2438	2441	2445	rVB4	20806	28641	0.75%	0.060%
84	17.792	2530	2534	2537	rBV5	44821	65012	1.71%	0.136%
85	17.921	2552	2556	2563	rVB3	81984	115512	3.03%	0.242%
86	18.327	2621	2625	2629	rBV2	43848	53026	1.39%	0.111%
87	18.609	2669	2673	2680	rBV3	68266	85142	2.24%	0.178%
88	18.780	2698	2702	2706	rBV3	30800	43210	1.13%	0.090%
89	18.992	2734	2738	2743	rBV	44946	56974	1.50%	0.119%
90	19.039	2743	2746	2750	rBV	43367	51780	1.36%	0.108%
91	19.262	2778	2784	2792	rBV	2710289	3807360	100.00%	7.960%
92	19.321	2792	2794	2799	rVB3	31848	38193	1.00%	0.080%
93	20.286	2953	2958	2961	rBV	345328	426880	11.21%	0.892%
94	20.809	3043	3047	3054	rBV	489944	624487	16.40%	1.306%
95	21.062	3085	3090	3098	rBV	1678942	2072856	54.44%	4.334%
96	21.192	3108	3112	3118	rBV	368948	428619	11.26%	0.896%
97	22.580	3345	3348	3354	rVB	161714	194708	5.11%	0.407%
98	23.044	3420	3427	3433	rBV	2047119	3601341	94.59%	7.529%
99	23.109	3435	3438	3442	rVV	195654	279142	7.33%	0.584%
100	23.174	3443	3449	3459	rVB	1178849	2121104	55.71%	4.435%

Sum of corrected areas: 47830763

LSC Report - Integrated Chromatogram

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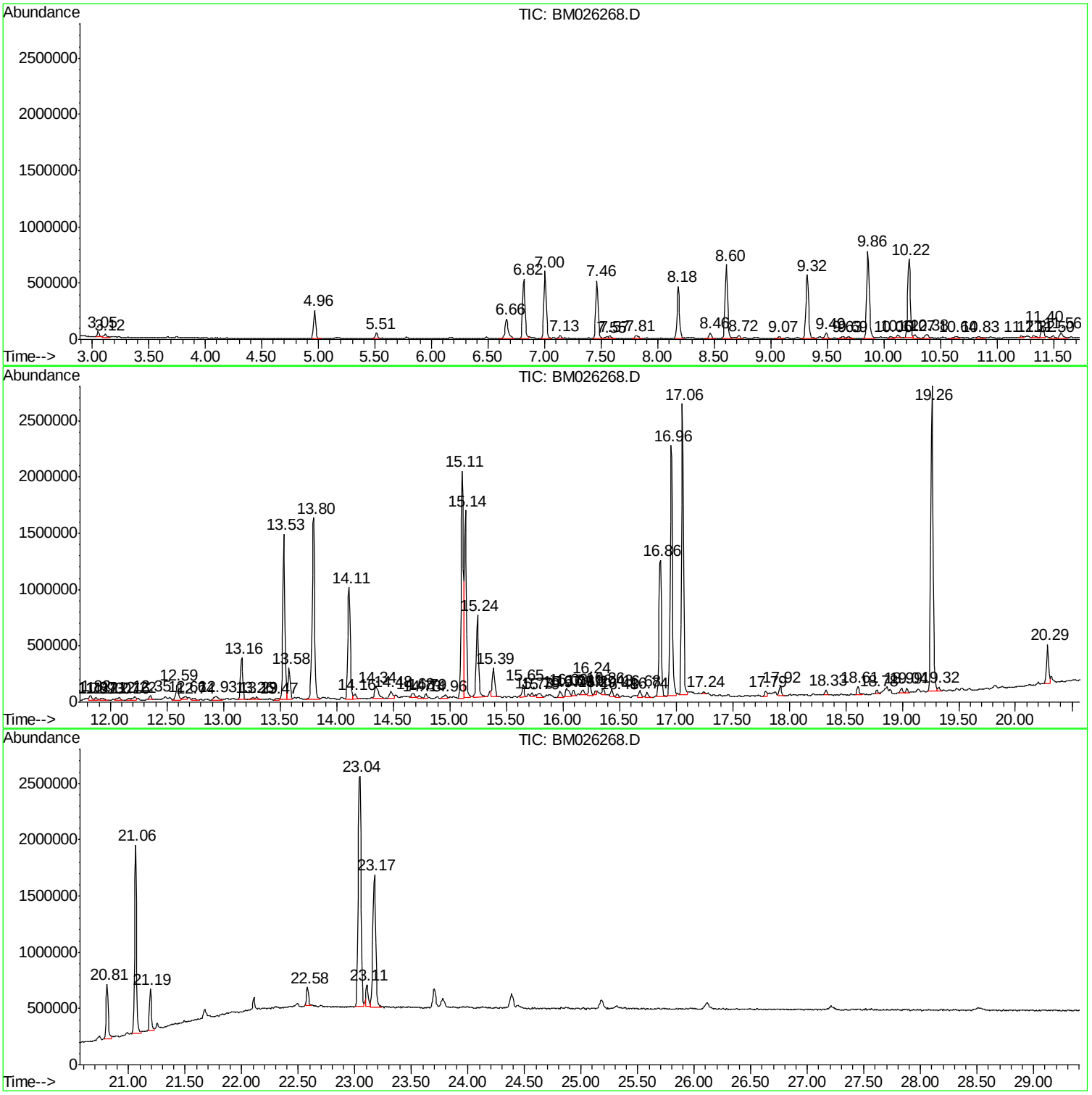
Data File : BM026268.D
Acq On : 04 Jun 2020 21:40
Operator : JU/CG
Sample : L2844-01
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0AE5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.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\BM060420\
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Instrument : BNA_M
ClientSampleId : C0AE5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

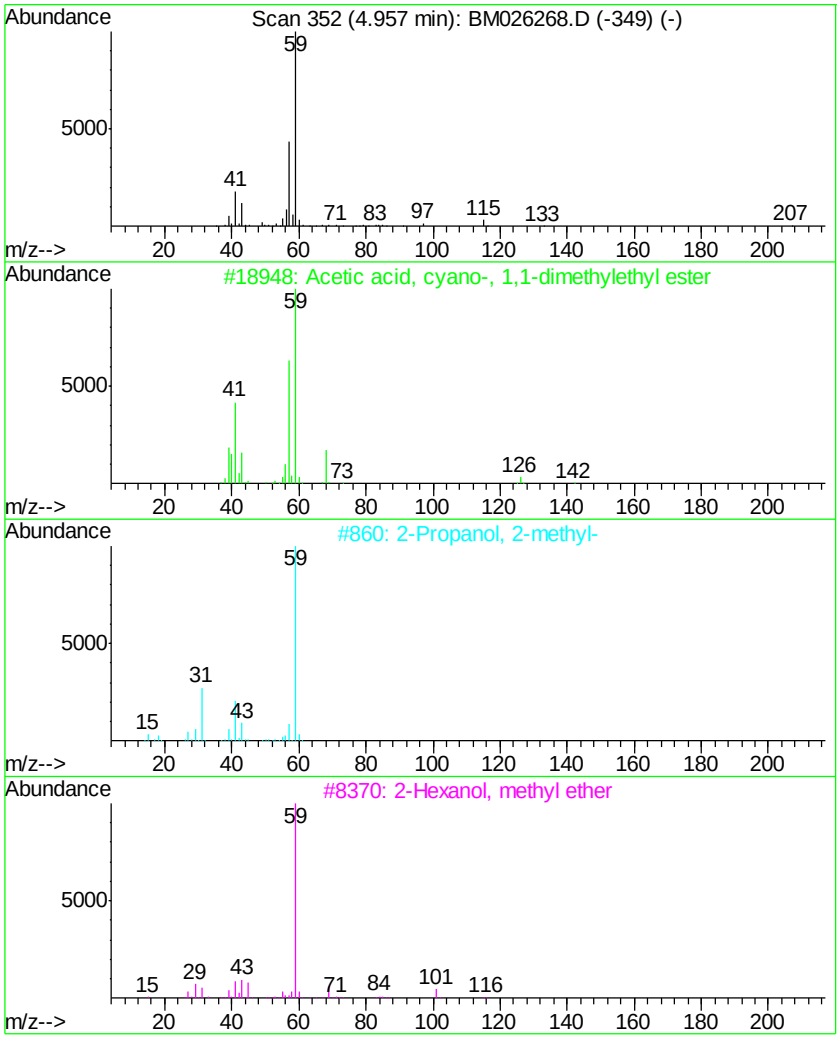
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

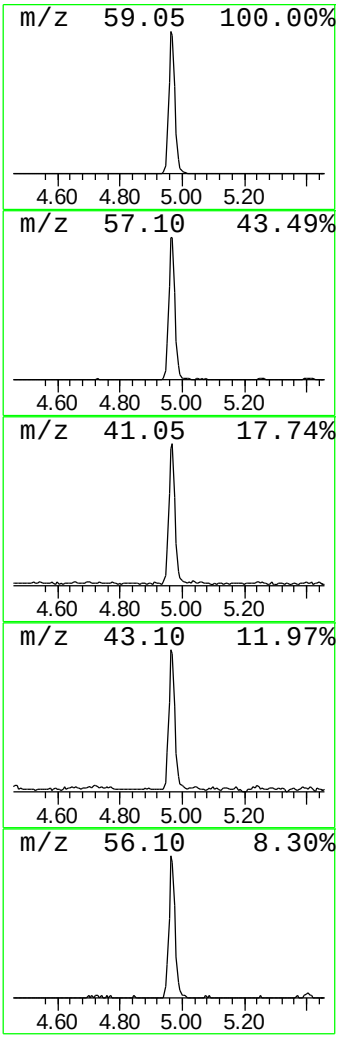
****

Peak Number 1 Acetic acid, cyano-, 1,1-di... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.96	8.73 ng/ul	350185	1,4-Dichlorobenzene-d4	7.46

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	64
2	2-Propanol, 2-methyl-	74	C4H10O	000075-65-0	38
3	2-Hexanol, methyl ether	116	C7H16O	1000333-92-0	9
4	Propane, 2-ethoxy-2-methyl-	102	C6H14O	000637-92-3	9
5	Propane, 2-ethoxy-2-methyl-	102	C6H14O	000637-92-3	9





Library Search Compound Report

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Acq On : 04 Jun 2020 21:40
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Sample : L2844-01
Misc :
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Instrument : BNA_M
ClientSampleId : C0AE5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

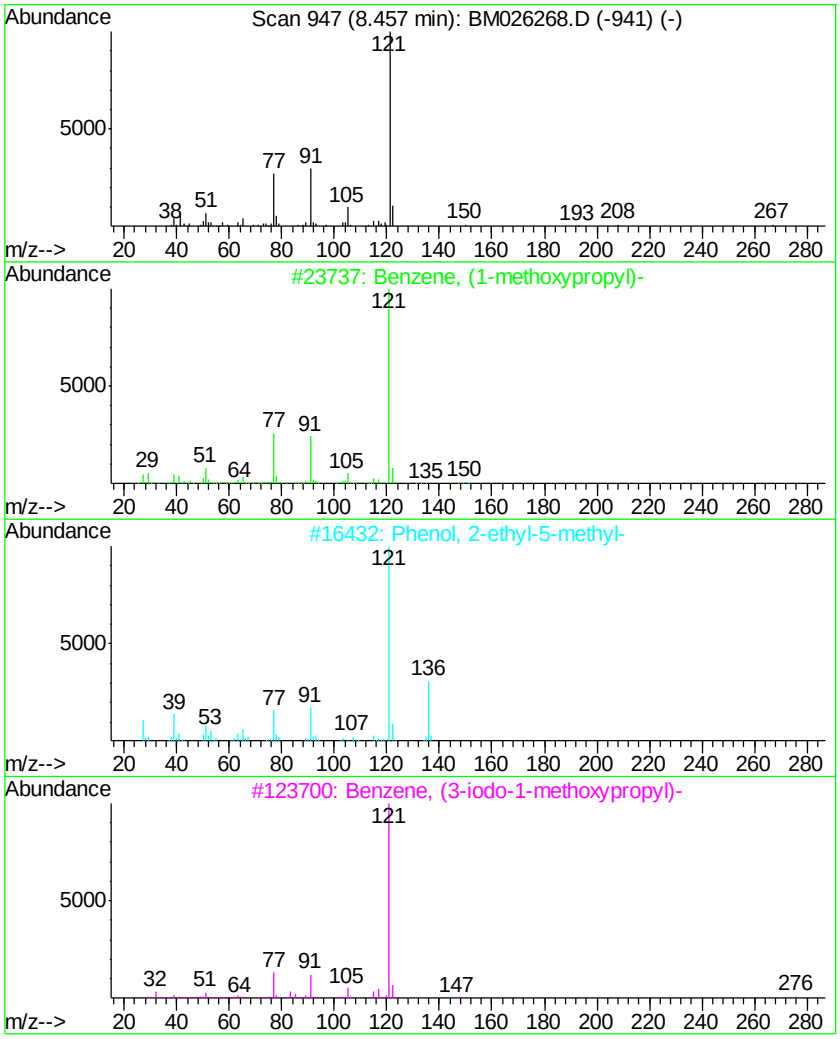
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

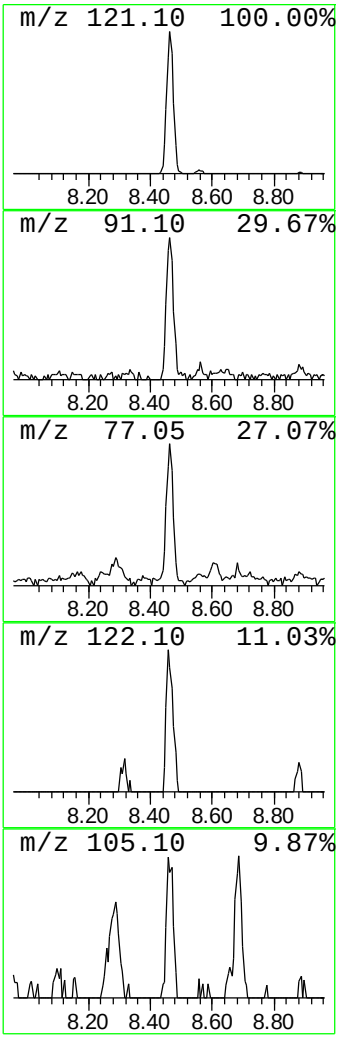
****

Peak Number 2 Benzene, (1-methoxypropyl)- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.46	2.00 ng/ul	80381	1,4-Dichlorobenzene-d4	7.46

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (1-methoxypropyl)-	150	C10H14O	059588-12-4	72
2	Phenol, 2-ethyl-5-methyl-	136	C9H12O	001687-61-2	72
3	Benzene, (3-iodo-1-methoxypropyl)-	276	C10H13IO	1000327-31-9	64
4	2-Methyl-6-propylphenol	150	C10H14O	003520-52-3	59
5	3-(4-Methoxyphenyl)propionic acid	180	C10H12O3	001929-29-9	59





Library Search Compound Report

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

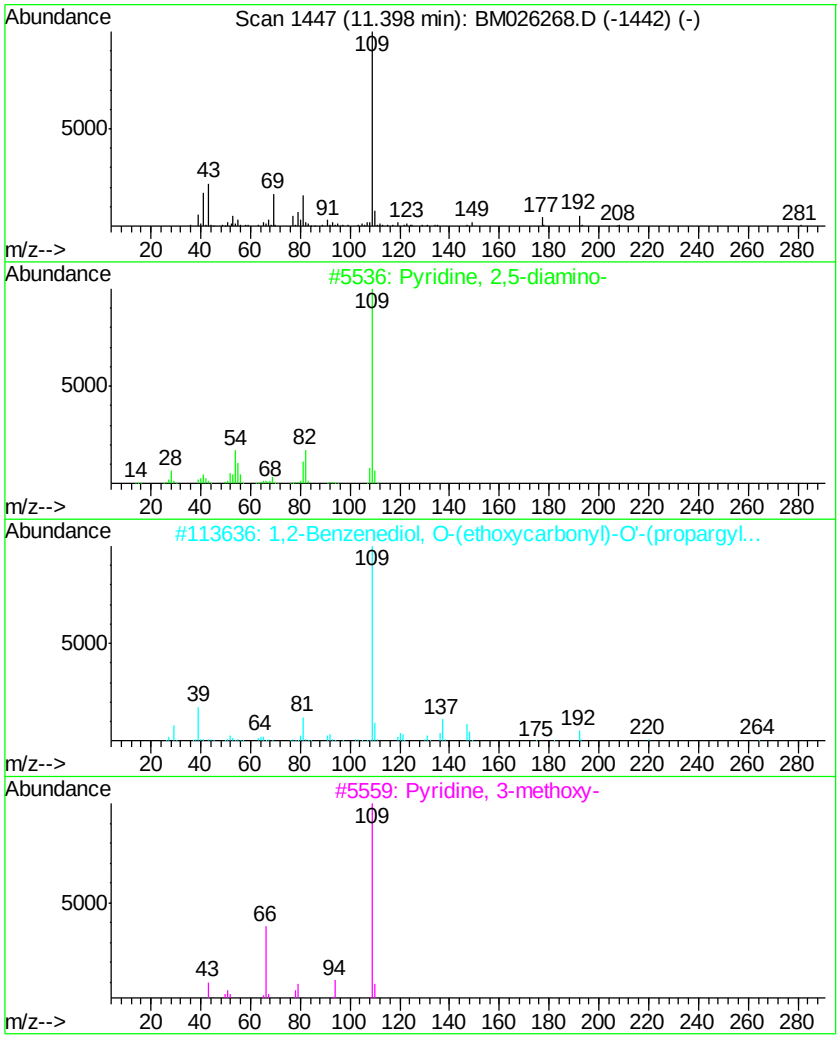
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

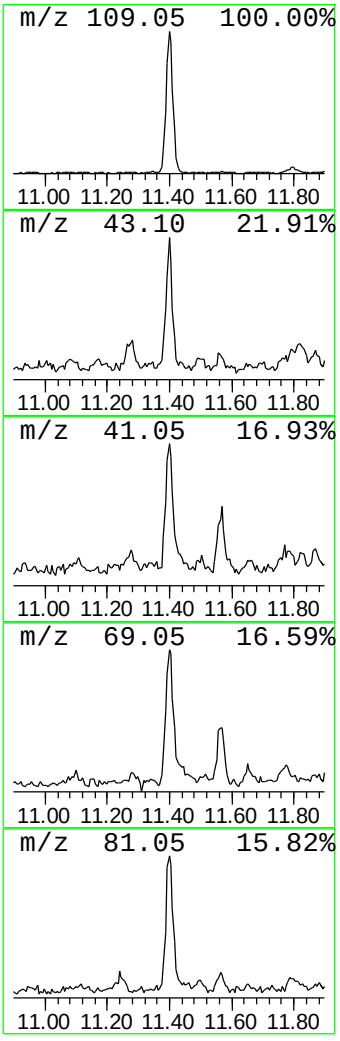
****

Peak Number 3 Pyridine, 2,5-diamino- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.40	2.91 ng/ul	158361	Naphthalene-d8	10.22

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyridine, 2,5-diamino-	109	C5H7N3	004318-76-7	59
2	1,2-Benzenediol, O-(ethoxycarbon...	264	C13H12O6	1000329-71-6	50
3	Pyridine, 3-methoxy-	109	C6H7NO	007295-76-3	50
4	Phenol, 3-amino-	109	C6H7NO	000591-27-5	45
5	7-(2,6-Dimethyl-hepta-1,5-dienyl...	272	C20H32	1000190-62-8	45





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Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0AE5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

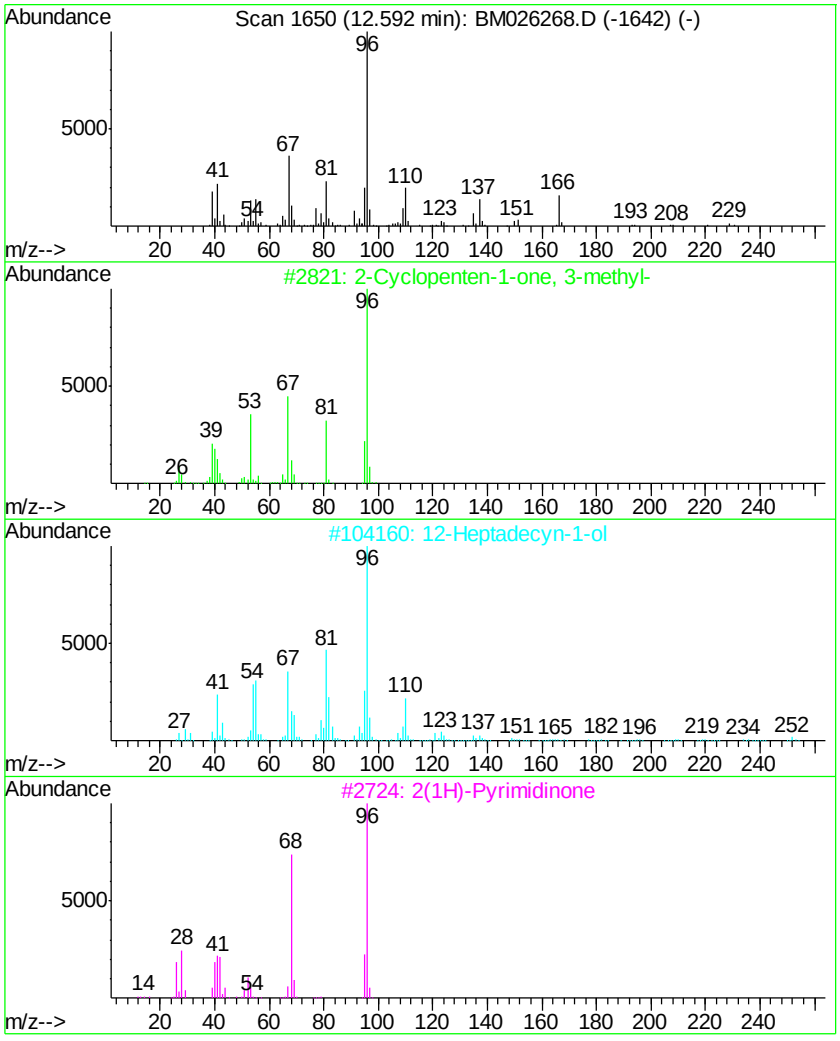
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

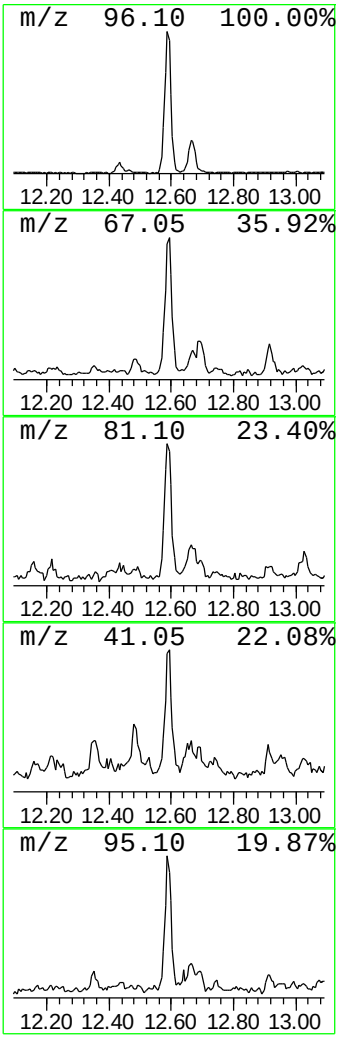
****

Peak Number 4 2-Cyclopenten-1-one, 3-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.59	3.07 ng/ul	227200	Acenaphthene-d10	14.11

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Cyclopenten-1-one, 3-methyl-	96	C6H8O	002758-18-1	58
2	12-Heptadecyn-1-ol	252	C17H32O	056554-76-8	45
3	2(1H)-Pyrimidinone	96	C4H4N2O	000557-01-7	43
4	Cyclohexane, 1-ethenyl-2-methyl-...	124	C9H16	034780-45-5	42
5	2-Cyclohexen-1-one, 4,4,6-trimet...	138	C9H14O	013395-73-8	40





Library Search Compound Report

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Data File : BM026268.D
Acq On : 04 Jun 2020 21:40
Operator : JU/CG
Sample : L2844-01
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0AE5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

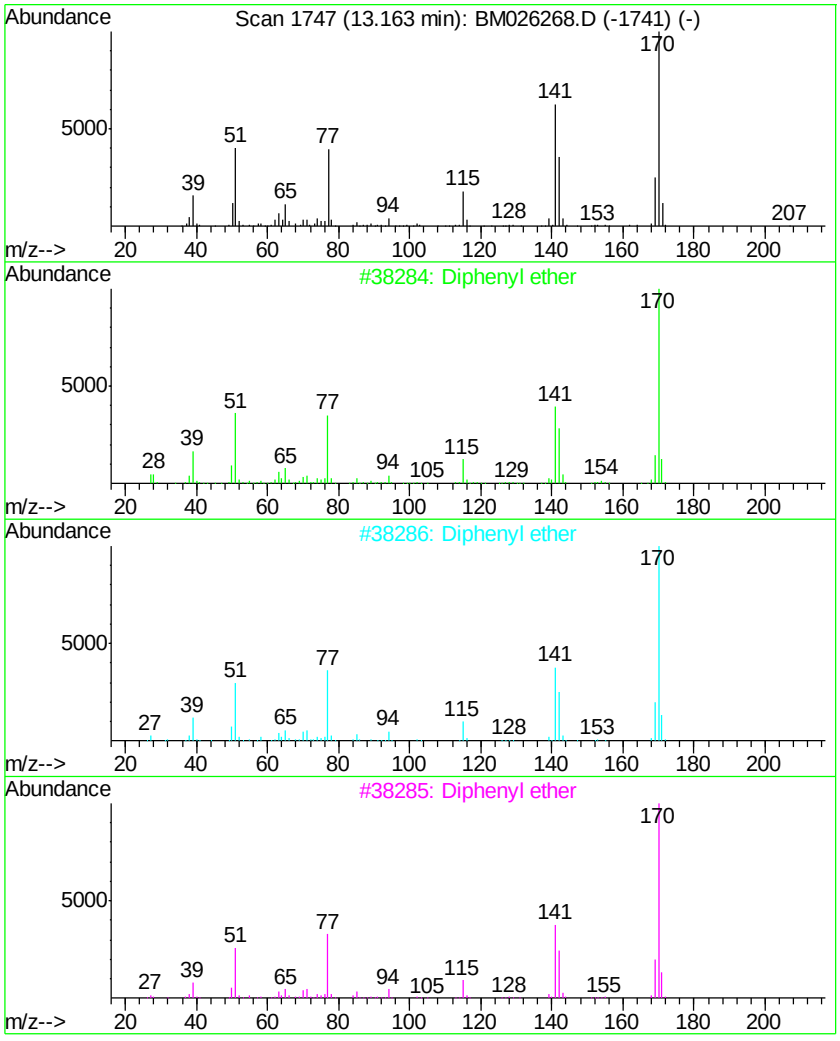
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

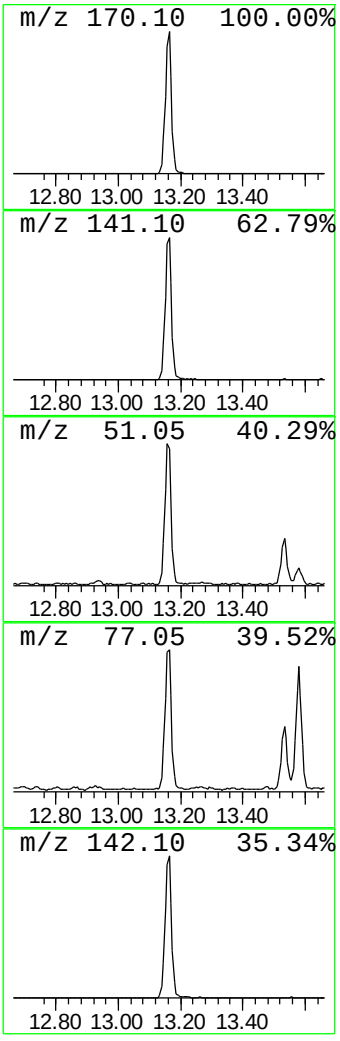
****

Peak Number 5 Diphenyl ether Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.16	7.01 ng/ul	519348	Acenaphthene-d10	14.11

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Diphenyl ether	170	C12H10O	000101-84-8	83
2	Diphenyl ether	170	C12H10O	000101-84-8	83
3	Diphenyl ether	170	C12H10O	000101-84-8	72
4	Diphenyl ether	170	C12H10O	000101-84-8	72
5	3-Hydroxybiphenyl	170	C12H10O	000580-51-8	43





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM060420\
Data File : BM026268.D
Acq On : 04 Jun 2020 21:40
Operator : JU/CG
Sample : L2844-01
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0AE5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

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

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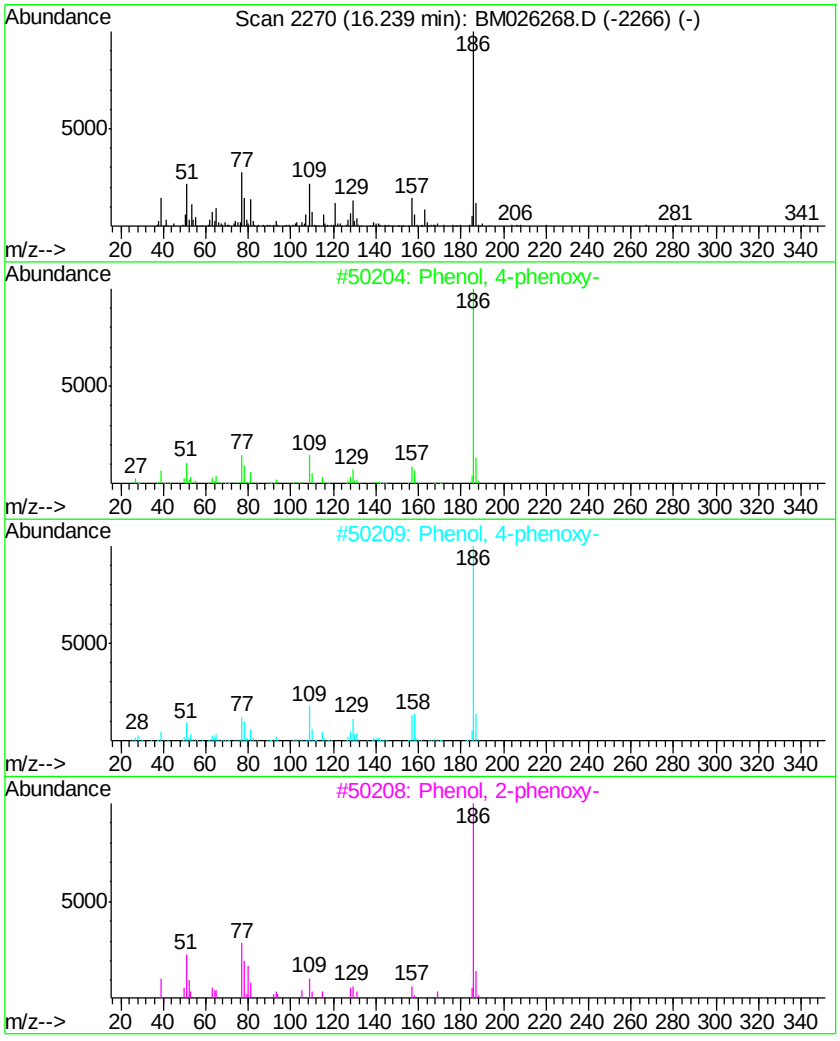
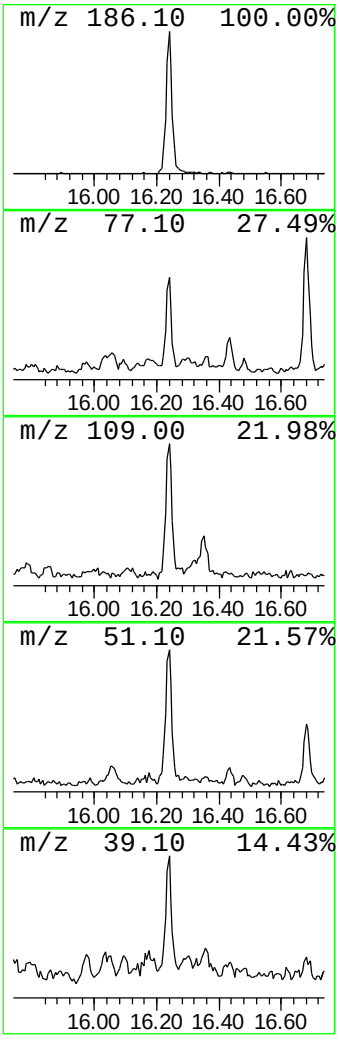
Peak Number 6 Phenol, 4-phenoxy- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.24	2.27 ng/ul	198760	Phenanthrene-d10	16.86

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 4-phenoxy-	186	C12H10O2	000831-82-3	93
2	Phenol, 4-phenoxy-	186	C12H10O2	000831-82-3	93
3	Phenol, 2-phenoxy-	186	C12H10O2	002417-10-9	64
4	Phenol, 2-phenoxy-	186	C12H10O2	002417-10-9	64
5	1,4-Naphthalenedione, 2,3-dimethyl-	186	C12H10O2	002197-57-1	49





Library Search Compound Report

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Data File : BM026268.D
Acq On : 04 Jun 2020 21:40
Operator : JU/CG
Sample : L2844-01
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0AE5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

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

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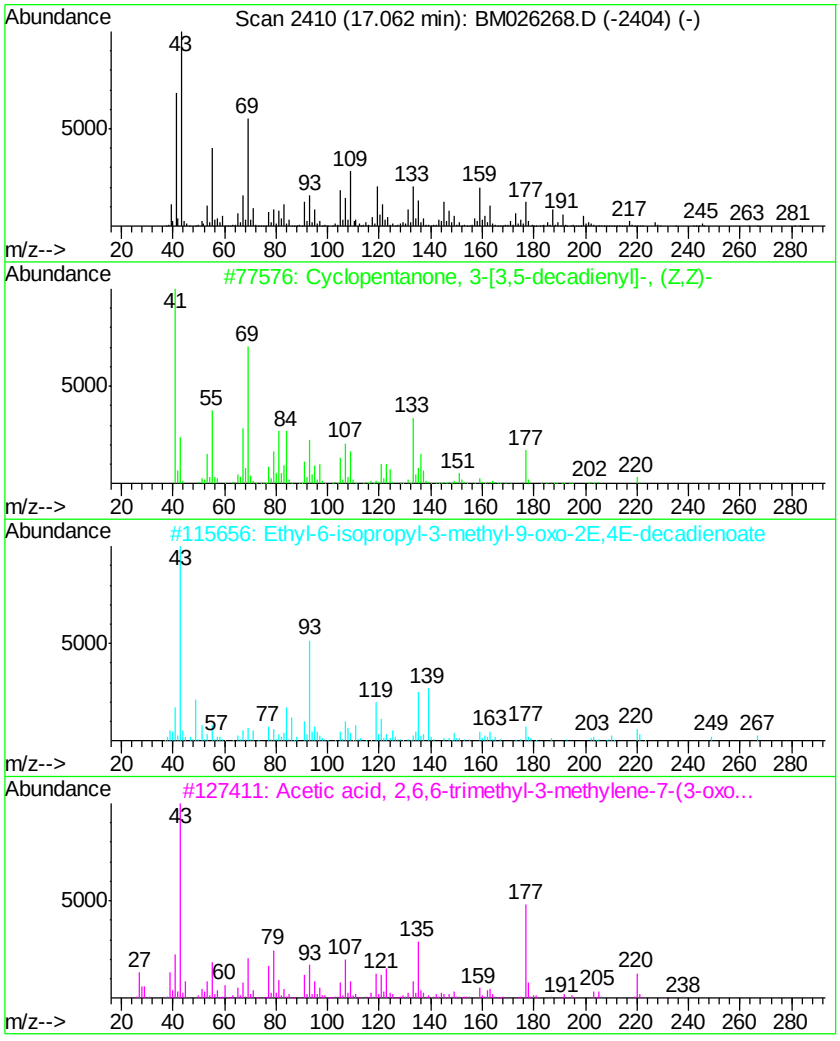
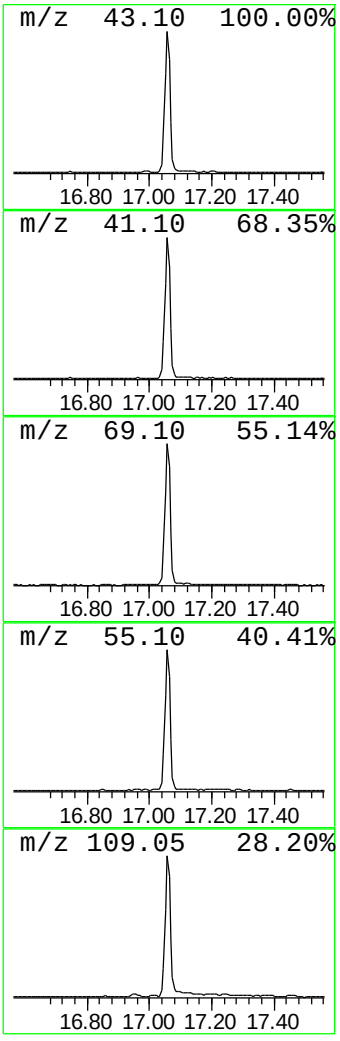
Peak Number 7 unknown-01 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.06	35.29 ng/ul	3088270	Phenanthrene-d10	16.86

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentanone, 3-[3,5-decadienyl...	220	C15H24O	1000131-99-8	35
2	Ethyl-6-isopropyl-3-methyl-9-oxo...	266	C16H26O3	063896-58-2	32
3	Acetic acid, 2,6,6-trimethyl-3-m...	280	C16H24O4	1000185-41-4	32
4	Tricyclo[5.2.2.0(1,6)]undecan-3-...	220	C15H24O	1000159-37-6	25
5	2H-1-Benzopyran, 3,5,8,8a-tetrah...	192	C13H20O	072468-40-7	25





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Operator : JU/CG
Sample : L2844-01
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0AE5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

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

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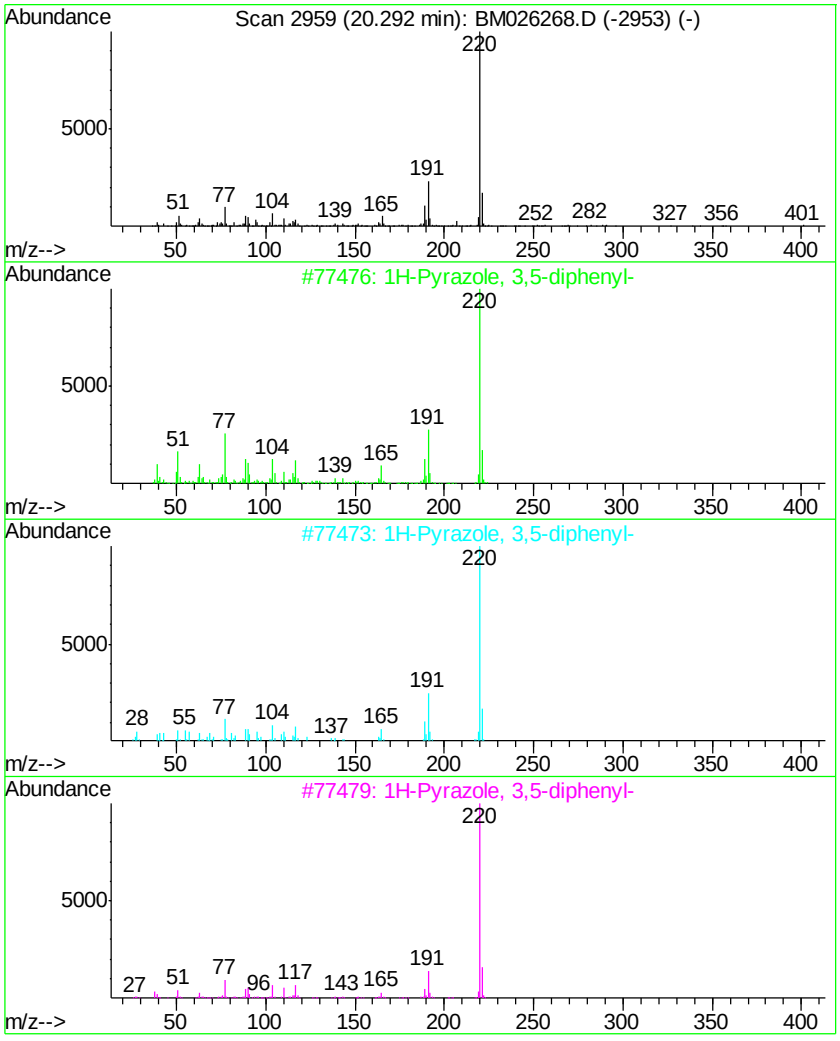
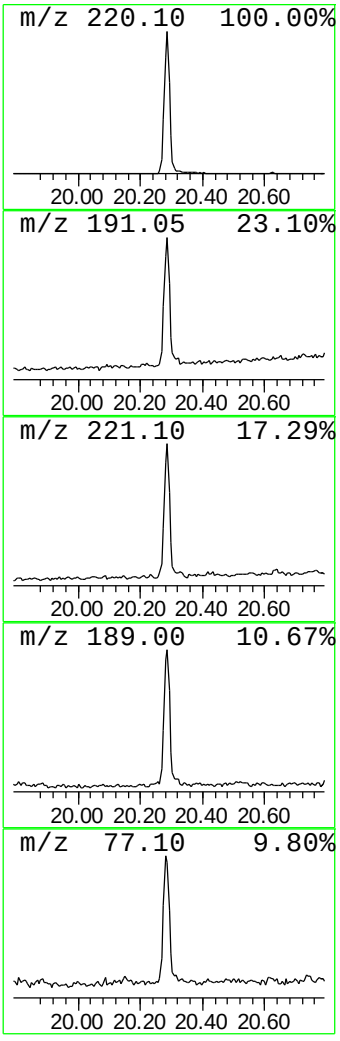
Peak Number 8 1H-Pyrazole, 3,5-diphenyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.29	4.12 ng/ul	426880	Chrysene-d12	21.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Pyrazole, 3,5-diphenyl-	220	C15H12N2	001145-01-3	97
2	1H-Pyrazole, 3,5-diphenyl-	220	C15H12N2	001145-01-3	93
3	1H-Pyrazole, 3,5-diphenyl-	220	C15H12N2	001145-01-3	93
4	1H-Pyrazole, 3,5-diphenyl-	220	C15H12N2	001145-01-3	93
5	6,7,8,9-Benzo[b]fluorene	220	C17H16	1000080-15-6	86





Library Search Compound Report

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Operator : JU/CG
Sample : L2844-01
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0AE5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

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

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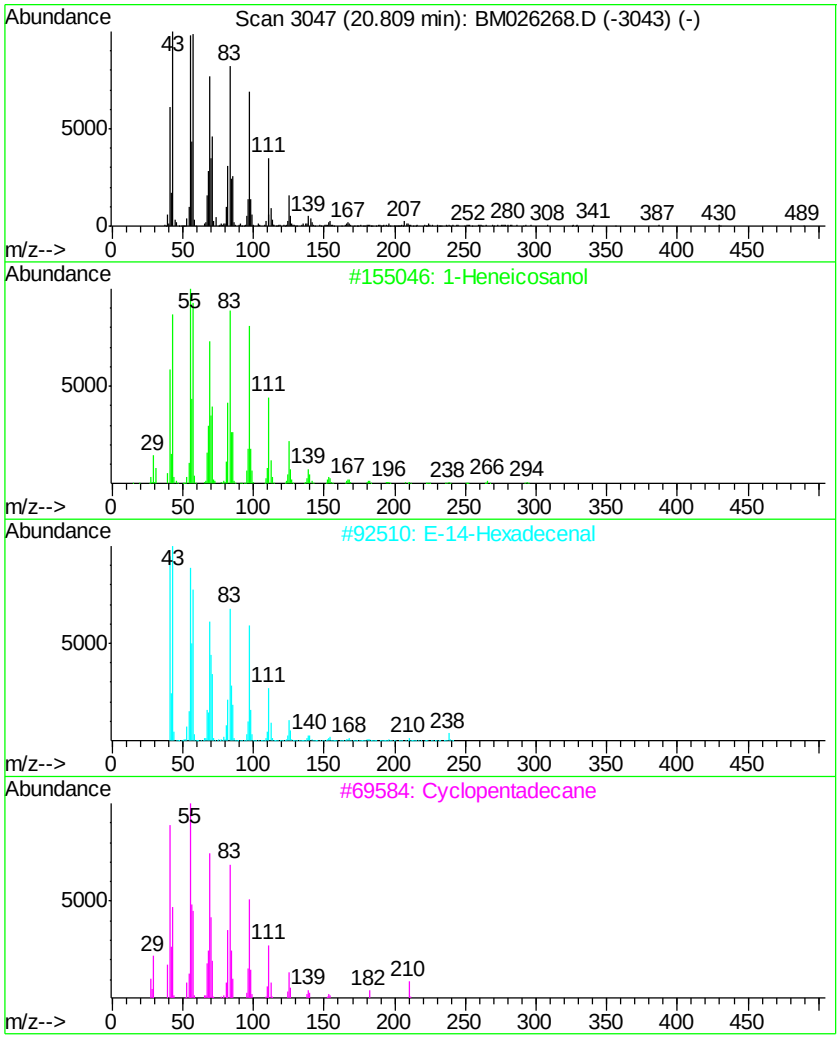
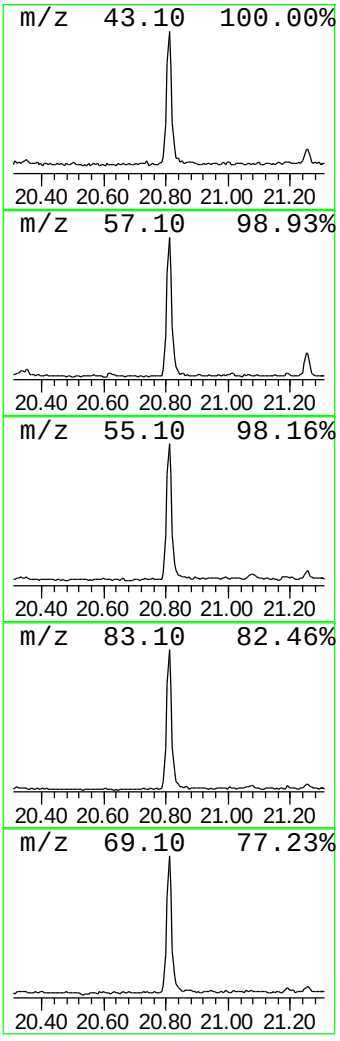
Peak Number 9 1-Heneicosanol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.81	6.03 ng/ul	624487	Chrysene-d12	21.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heneicosanol	312	C21H44O	015594-90-8	94
2	E-14-Hexadecenal	238	C16H30O	330207-53-9	93
3	Cyclopentadecane	210	C15H30	000295-48-7	93
4	Cyclotetracosane	336	C24H48	000297-03-0	93
5	1-Docosene	308	C22H44	001599-67-3	92





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Data File : BM026268.D
Acq On : 04 Jun 2020 21:40
Operator : JU/CG
Sample : L2844-01
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0AE5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

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

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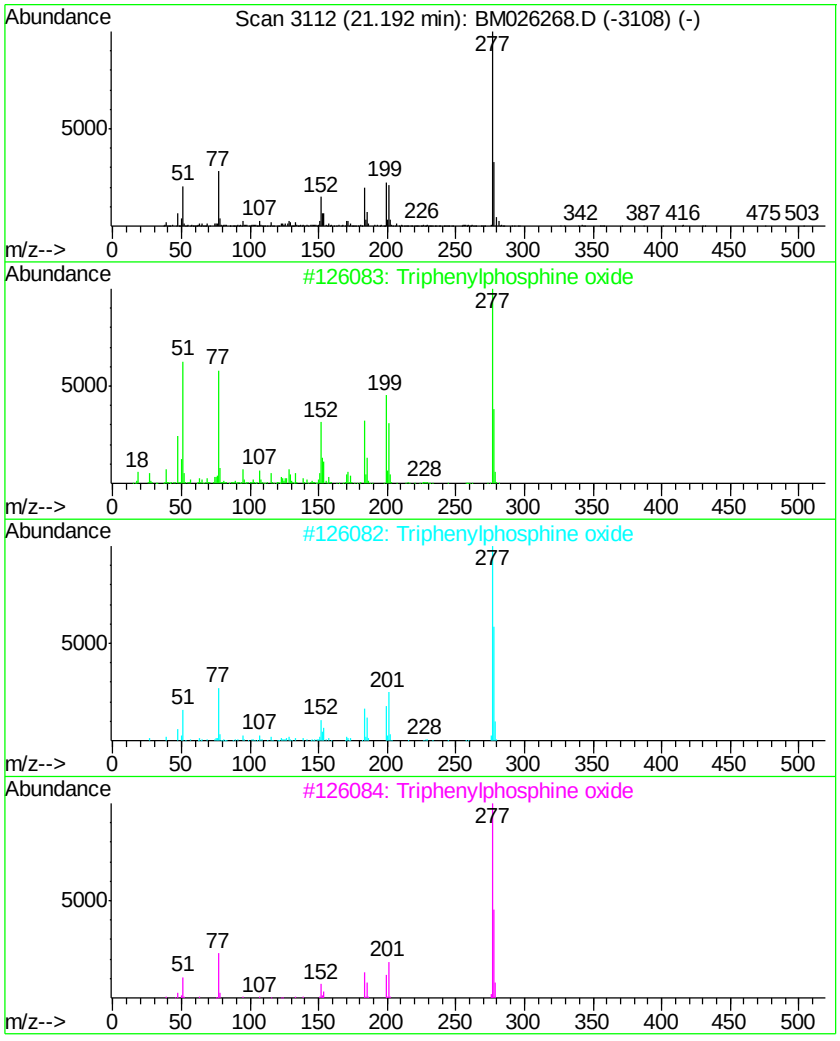
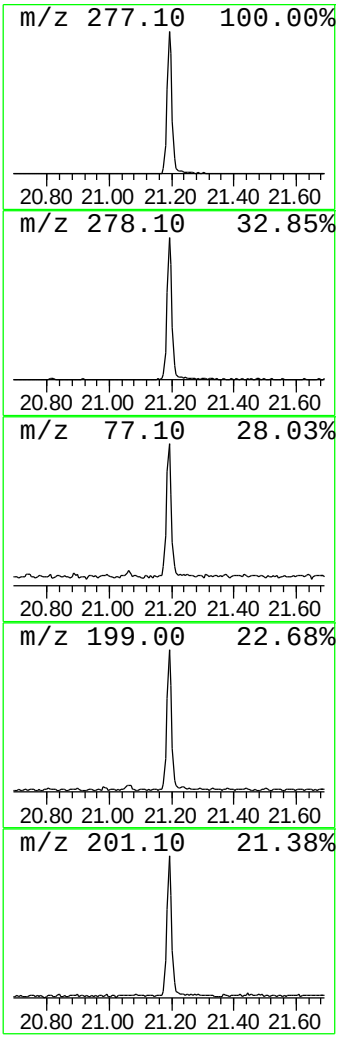
Peak Number 10 Triphenylphosphine oxide Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.19	4.14 ng/ul	428619	Chrysene-d12	21.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Triphenylphosphine oxide	278	C18H15OP	000791-28-6	96
2	Triphenylphosphine oxide	278	C18H15OP	000791-28-6	83
3	Triphenylphosphine oxide	278	C18H15OP	000791-28-6	50
4	Formaldehyde, (triphenylphosphor...	304	C19H17N2P	015990-54-2	50
5	Triphenylphosphine oxide	278	C18H15OP	000791-28-6	46





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Sample : L2844-01
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0AE5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

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

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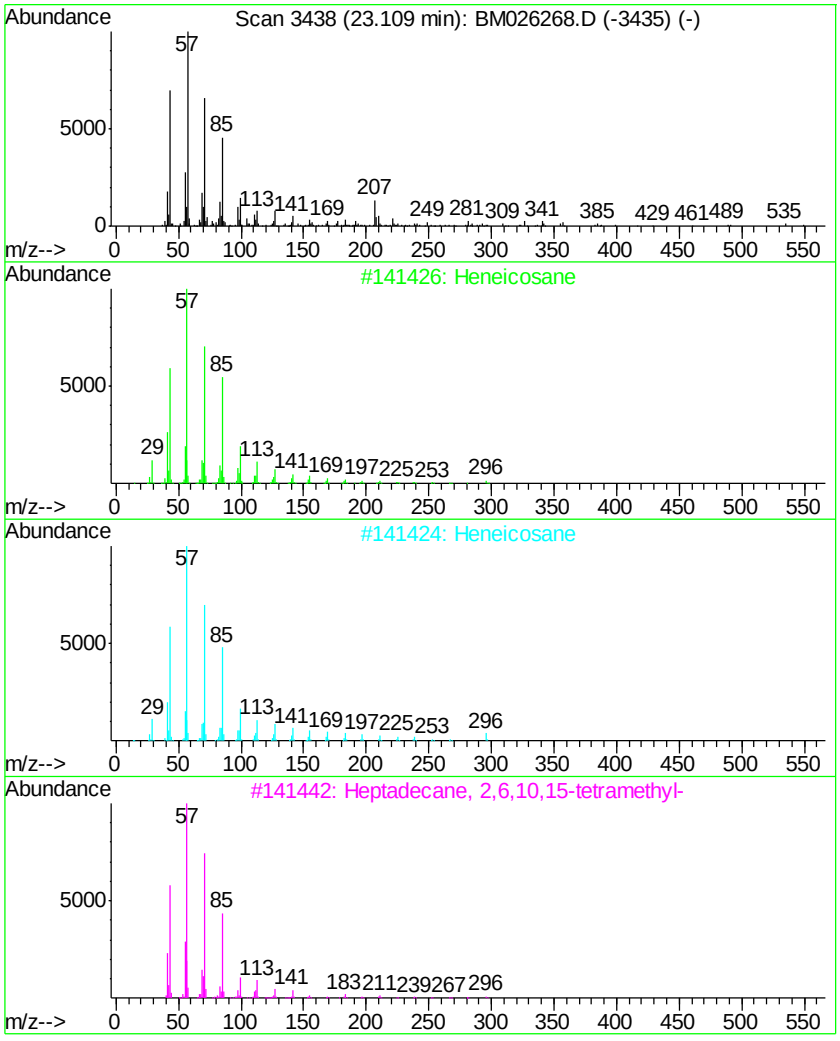
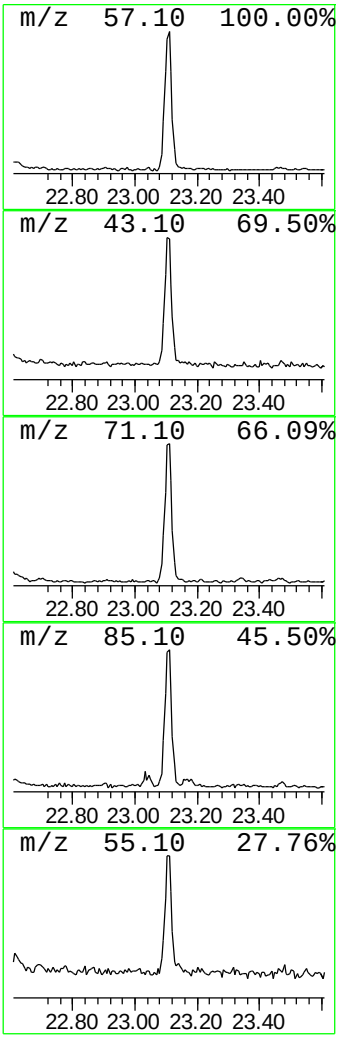
Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.11	2.63 ng/ul	279142	Perylene-d12	23.17

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	96
2	Heneicosane	296	C21H44	000629-94-7	95
3	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	93
4	Tricosane	324	C23H48	000638-67-5	89
5	Tricosane	324	C23H48	000638-67-5	89





Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM060420\
Data File : BM026268.D
Acq On : 04 Jun 2020 21:40
Operator : JU/CG
Sample : L2844-01
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0AE5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.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
Acetic acid, cyan...	4.96	8.7	ng/ul	350185	1	7.46	802121	20.0
Benzene, (1-metho...	8.46	2.0	ng/ul	80381	1	7.46	802121	20.0
Pyridine, 2,5-dia...	11.40	2.9	ng/ul	158361	2	10.22	1087730	20.0
2-Cyclopenten-1-o...	12.59	3.1	ng/ul	227200	3	14.11	1482040	20.0
Diphenyl ether	13.16	7.0	ng/ul	519348	3	14.11	1482040	20.0
Phenol, 4-phenoxy-	16.24	2.3	ng/ul	198760	4	16.86	1750130	20.0
unknown-01	17.06	35.3	ng/ul	3088270	4	16.86	1750130	20.0
1H-Pyrazole, 3,5-...	20.29	4.1	ng/ul	426880	5	21.06	2072860	20.0
1-Heneicosanol	20.81	6.0	ng/ul	624487	5	21.06	2072860	20.0
Triphenylphosphin...	21.19	4.1	ng/ul	428619	5	21.06	2072860	20.0
(DEL) Alkane: Str...	23.11	2.6	ng/ul	279142	6	23.17	2121100	20.0