LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092520\
Data File : BM027537.D
Acq On : 25 Sep 2020 20:05
Operator : JU/CG
Sample : L4135-06
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** BG480

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-BM092420MA.M
Title : SVOA CALIBRATION

Signal : TIC: BM027537.D\data.ms

| peak # | R.T. min | first scan | max scan | last scan | PK TY | peak height | corr. area | corr. % max. | % of total |
|---|---|---|---|---|---|---|---|---|---|
| 1 | 3.063 | 9 | 13 | 20 | rVB | 144285 | 197427 | 5.27% | 0.442% |
| 2 | 3.311 | 51 | 55 | 59 | rBV | 23369 | 31399 | 0.84% | 0.070% |
| 3 | 5.152 | 361 | 368 | 380 | rVB | 1774066 | 2625327 | 70.12% | 5.875% |
| 4 | 6.310 | 558 | 565 | 574 | rBV | 119808 | 181870 | 4.86% | 0.407% |
| 5 | 6.799 | 641 | 648 | 657 | rBV3 | 17812 | 36622 | 0.98% | 0.082% |
| 6 | 6.969 | 670 | 677 | 685 | rBV | 95673 | 162947 | 4.35% | 0.365% |
| 7 | 7.140 | 699 | 706 | 713 | rBV | 272893 | 418205 | 11.17% | 0.936% |
| 8 | 7.234 | 718 | 722 | 727 | rVB | 33375 | 51050 | 1.36% | 0.114% |
| 9 | 7.340 | 734 | 740 | 748 | rBV | 360806 | 585249 | 15.63% | 1.310% |
| 10 | 7.416 | 748 | 753 | 759 | rVB | 37574 | 54892 | 1.47% | 0.123% |
| 11 | 7.516 | 764 | 770 | 776 | rVB | 62829 | 93466 | 2.50% | 0.209% |
| 12 | 7.810 | 808 | 820 | 831 | rBV | 392235 | 693851 | 18.53% | 1.553% |
| 13 | 8.010 | 850 | 854 | 863 | rVV | 59196 | 87909 | 2.35% | 0.197% |
| 14 | 8.181 | 875 | 883 | 890 | rVB2 | 31964 | 72116 | 1.93% | 0.161% |
| 15 | 8.504 | 932 | 938 | 948 | rVB | 297147 | 462619 | 12.36% | 1.035% |
| 16 | 8.798 | 980 | 988 | 997 | rVB4 | 15152 | 36589 | 0.98% | 0.082% |
| 17 | 8.910 | 1000 | 1007 | 1009 | rBV3 | 20254 | 33433 | 0.89% | 0.075% |
| 18 | 8.957 | 1009 | 1015 | 1024 | rVB | 376246 | 611799 | 16.34% | 1.369% |
| 19 | 9.034 | 1024 | 1028 | 1033 | rBV5 | 17341 | 29235 | 0.78% | 0.065% |
| 20 | 9.428 | 1091 | 1095 | 1101 | rVB | 21109 | 34395 | 0.92% | 0.077% |
| 21 | 9.675 | 1131 | 1137 | 1143 | rBV | 327230 | 517202 | 13.81% | 1.157% |
| 22 | 9.845 | 1161 | 1166 | 1172 | rBV5 | 20383 | 38941 | 1.04% | 0.087% |
| 23 | 10.010 | 1185 | 1194 | 1198 | rBV3 | 29460 | 60131 | 1.61% | 0.135% |
| 24 | 10.210 | 1221 | 1228 | 1236 | rVV | 655505 | 1063426 | 28.40% | 2.380% |
| 25 | 10.334 | 1236 | 1249 | 1257 | rVV7 | 35138 | 92109 | 2.46% | 0.206% |
| 26 | 10.410 | 1257 | 1262 | 1266 | rVV2 | 35753 | 54257 | 1.45% | 0.121% |
| 27 | 10.592 | 1286 | 1293 | 1299 | rVV | 540238 | 880223 | 23.51% | 1.970% |
| 28 | 10.722 | 1311 | 1315 | 1320 | rVV | 28846 | 52005 | 1.39% | 0.116% |
| 29 | 11.181 | 1388 | 1393 | 1402 | rVV5 | 27065 | 54319 | 1.45% | 0.122% |
| 30 | 11.604 | 1459 | 1465 | 1473 | rBB4 | 15540 | 31512 | 0.84% | 0.071% |
| 31 | 11.722 | 1475 | 1485 | 1491 | rBV5 | 19643 | 44188 | 1.18% | 0.099% |
| 32 | 11.933 | 1512 | 1521 | 1529 | rBV | 433167 | 722166 | 19.29% | 1.616% |
| 33 | 12.416 | 1596 | 1603 | 1610 | rBV9 | 19817 | 59642 | 1.59% | 0.133% |
| 34 | 12.692 | 1642 | 1650 | 1660 | rBV4 | 26699 | 58887 | 1.57% | 0.132% |
| 35 | 12.922 | 1683 | 1689 | 1693 | rVV3 | 25046 | 44799 | 1.20% | 0.100% |
| 36 | 12.969 | 1693 | 1697 | 1702 | rVV | 113369 | 155960 | 4.17% | 0.349% |

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092520\
Data File : BM027537.D
Acq On : 25 Sep 2020 20:05
Operator : JU/CG
Sample : L4135-06
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** BG480

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-BM092420MA.M
Title : SVOA CALIBRATION

| | | | | | | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 37 | 13.022 | 1702 | 1706 | 1710 | rVV6 | 18214 | 31794 | 0.85% | 0.071% |
| 38 | 13.086 | 1710 | 1717 | 1722 | rVV3 | 34450 | 57466 | 1.53% | 0.129% |
| 39 | 13.351 | 1754 | 1762 | 1768 | rBV3 | 12118 | 32001 | 0.85% | 0.072% |
| 40 | 13.492 | 1779 | 1786 | 1793 | rBV2 | 72841 | 127827 | 3.41% | 0.286% |
| 41 | 13.563 | 1793 | 1798 | 1803 | rBV6 | 19603 | 33386 | 0.89% | 0.075% |
| 42 | 13.733 | 1820 | 1827 | 1833 | rVB6 | 17493 | 37727 | 1.01% | 0.084% |
| 43 | 13.851 | 1840 | 1847 | 1853 | rBV | 1382551 | 1952762 | 52.15% | 4.370% |
| 44 | 13.980 | 1865 | 1869 | 1876 | rVV | 37323 | 53836 | 1.44% | 0.120% |
| 45 | 14.127 | 1887 | 1894 | 1903 | rVB | 1326795 | 2047718 | 54.69% | 4.582% |
| 46 | 14.216 | 1903 | 1909 | 1920 | rBV | 90336 | 147645 | 3.94% | 0.330% |
| 47 | 14.439 | 1939 | 1947 | 1951 | rBV | 1044210 | 1512977 | 40.41% | 3.386% |
| 48 | 14.474 | 1951 | 1953 | 1962 | rVB2 | 110400 | 159486 | 4.26% | 0.357% |
| 49 | 14.622 | 1973 | 1978 | 1988 | rBV | 100971 | 168856 | 4.51% | 0.378% |
| 50 | 14.980 | 2033 | 2039 | 2044 | rBV3 | 26772 | 47934 | 1.28% | 0.107% |
| 51 | 15.186 | 2069 | 2074 | 2083 | rBV | 73295 | 121466 | 3.24% | 0.272% |
| 52 | 15.339 | 2095 | 2100 | 2106 | rVB7 | 30479 | 48304 | 1.29% | 0.108% |
| 53 | 15.433 | 2109 | 2116 | 2123 | rVV | 1905861 | 2698378 | 72.07% | 6.038% |
| 54 | 15.539 | 2128 | 2134 | 2143 | rBV | 454624 | 670123 | 17.90% | 1.500% |
| 55 | 15.698 | 2151 | 2161 | 2171 | rVB | 2754265 | 3744180 | 100.00% | 8.379% |
| 56 | 15.939 | 2196 | 2202 | 2207 | rBV2 | 93894 | 164536 | 4.39% | 0.368% |
| 57 | 16.104 | 2224 | 2230 | 2235 | rBV | 161018 | 249139 | 6.65% | 0.558% |
| 58 | 16.157 | 2235 | 2239 | 2242 | rVB2 | 29123 | 40893 | 1.09% | 0.092% |
| 59 | 16.451 | 2284 | 2289 | 2292 | rBV2 | 60510 | 85107 | 2.27% | 0.190% |
| 60 | 16.486 | 2292 | 2295 | 2301 | rVB | 158622 | 205693 | 5.49% | 0.460% |
| 61 | 16.563 | 2301 | 2308 | 2311 | rBV2 | 32328 | 59342 | 1.58% | 0.133% |
| 62 | 16.651 | 2319 | 2323 | 2327 | rBV2 | 25381 | 34825 | 0.93% | 0.078% |
| 63 | 16.704 | 2329 | 2332 | 2339 | rVB7 | 19054 | 34456 | 0.92% | 0.077% |
| 64 | 17.139 | 2401 | 2406 | 2409 | rVV | 362878 | 474636 | 12.68% | 1.062% |
| 65 | 17.180 | 2409 | 2413 | 2421 | rVV | 1284028 | 1800586 | 48.09% | 4.029% |
| 66 | 17.280 | 2423 | 2430 | 2436 | rVV | 2146693 | 3065224 | 81.87% | 6.859% |
| 67 | 17.427 | 2451 | 2455 | 2463 | rVB6 | 22461 | 50918 | 1.36% | 0.114% |
| 68 | 17.674 | 2492 | 2497 | 2501 | rBV8 | 15208 | 30565 | 0.82% | 0.068% |
| 69 | 17.792 | 2513 | 2517 | 2521 | rVV2 | 43034 | 58498 | 1.56% | 0.131% |
| 70 | 17.845 | 2521 | 2526 | 2531 | rVV | 127022 | 190592 | 5.09% | 0.427% |
| 71 | 17.910 | 2533 | 2537 | 2546 | rVB | 47689 | 81161 | 2.17% | 0.182% |
| 72 | 18.086 | 2562 | 2567 | 2573 | rBV | 187391 | 263857 | 7.05% | 0.590% |
| 73 | 18.145 | 2573 | 2577 | 2583 | rVV | 137589 | 205039 | 5.48% | 0.459% |
| 74 | 18.215 | 2583 | 2589 | 2596 | rVB | 255350 | 358681 | 9.58% | 0.803% |
| 75 | 18.315 | 2601 | 2606 | 2609 | rVV | 213520 | 314769 | 8.41% | 0.704% |
| 76 | 18.368 | 2609 | 2615 | 2623 | rVB | 388629 | 673141 | 17.98% | 1.506% |

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092520\
Data File : BM027537.D
Acq On : 25 Sep 2020 20:05
Operator : JU/CG
Sample : L4135-06
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** BG480

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-BM092420MA.M
Title : SVOA CALIBRATION

| | | | | | | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 77 | 18.774 | 2677 | 2684 | 2690 | rBV5 | 34068 | 63245 | 1.69% | 0.142% |
| 78 | 18.904 | 2702 | 2706 | 2712 | rBV | 73946 | 105576 | 2.82% | 0.236% |
| 79 | 19.039 | 2725 | 2729 | 2732 | rBV | 38397 | 49099 | 1.31% | 0.110% |
| 80 | 19.074 | 2732 | 2735 | 2739 | rVB2 | 58019 | 63998 | 1.71% | 0.143% |
| 81 | 19.180 | 2749 | 2753 | 2763 | rVB3 | 106594 | 194363 | 5.19% | 0.435% |
| 82 | 19.292 | 2767 | 2772 | 2777 | rBV8 | 34801 | 70633 | 1.89% | 0.158% |
| 83 | 19.362 | 2781 | 2784 | 2795 | rVB2 | 71713 | 94353 | 2.52% | 0.211% |
| 84 | 19.457 | 2797 | 2800 | 2809 | rVB2 | 27100 | 53034 | 1.42% | 0.119% |
| 85 | 19.568 | 2813 | 2819 | 2824 | rBV | 2611439 | 3493849 | 93.31% | 7.819% |
| 86 | 19.615 | 2824 | 2827 | 2836 | rVB | 522604 | 615581 | 16.44% | 1.378% |
| 87 | 19.780 | 2852 | 2855 | 2859 | rVV2 | 54473 | 74831 | 2.00% | 0.167% |
| 88 | 19.821 | 2859 | 2862 | 2865 | rVV2 | 41728 | 51359 | 1.37% | 0.115% |
| 89 | 19.862 | 2865 | 2869 | 2874 | rVB2 | 46941 | 67386 | 1.80% | 0.151% |
| 90 | 19.974 | 2884 | 2888 | 2891 | rBV | 64931 | 84628 | 2.26% | 0.189% |
| 91 | 20.015 | 2891 | 2895 | 2901 | rVB | 105034 | 133288 | 3.56% | 0.298% |
| 92 | 20.068 | 2901 | 2904 | 2906 | rBV3 | 35521 | 42727 | 1.14% | 0.096% |
| 93 | 20.174 | 2919 | 2922 | 2926 | rVB3 | 28335 | 32911 | 0.88% | 0.074% |
| 94 | 20.298 | 2939 | 2943 | 2949 | rVB2 | 44573 | 59410 | 1.59% | 0.133% |
| 95 | 20.609 | 2993 | 2996 | 3002 | rVB7 | 28281 | 37042 | 0.99% | 0.083% |
| 96 | 21.080 | 3073 | 3076 | 3085 | rBV | 194967 | 270324 | 7.22% | 0.605% |
| 97 | 21.356 | 3118 | 3123 | 3130 | rBV | 1444303 | 1780141 | 47.54% | 3.984% |
| 98 | 21.480 | 3140 | 3144 | 3151 | rBV2 | 146327 | 252002 | 6.73% | 0.564% |
| 99 | 23.486 | 3478 | 3485 | 3493 | rBV | 1492199 | 2779063 | 74.22% | 6.219% |
| 100 | 23.627 | 3502 | 3509 | 3519 | rVB | 854442 | 1651819 | 44.12% | 3.696% |

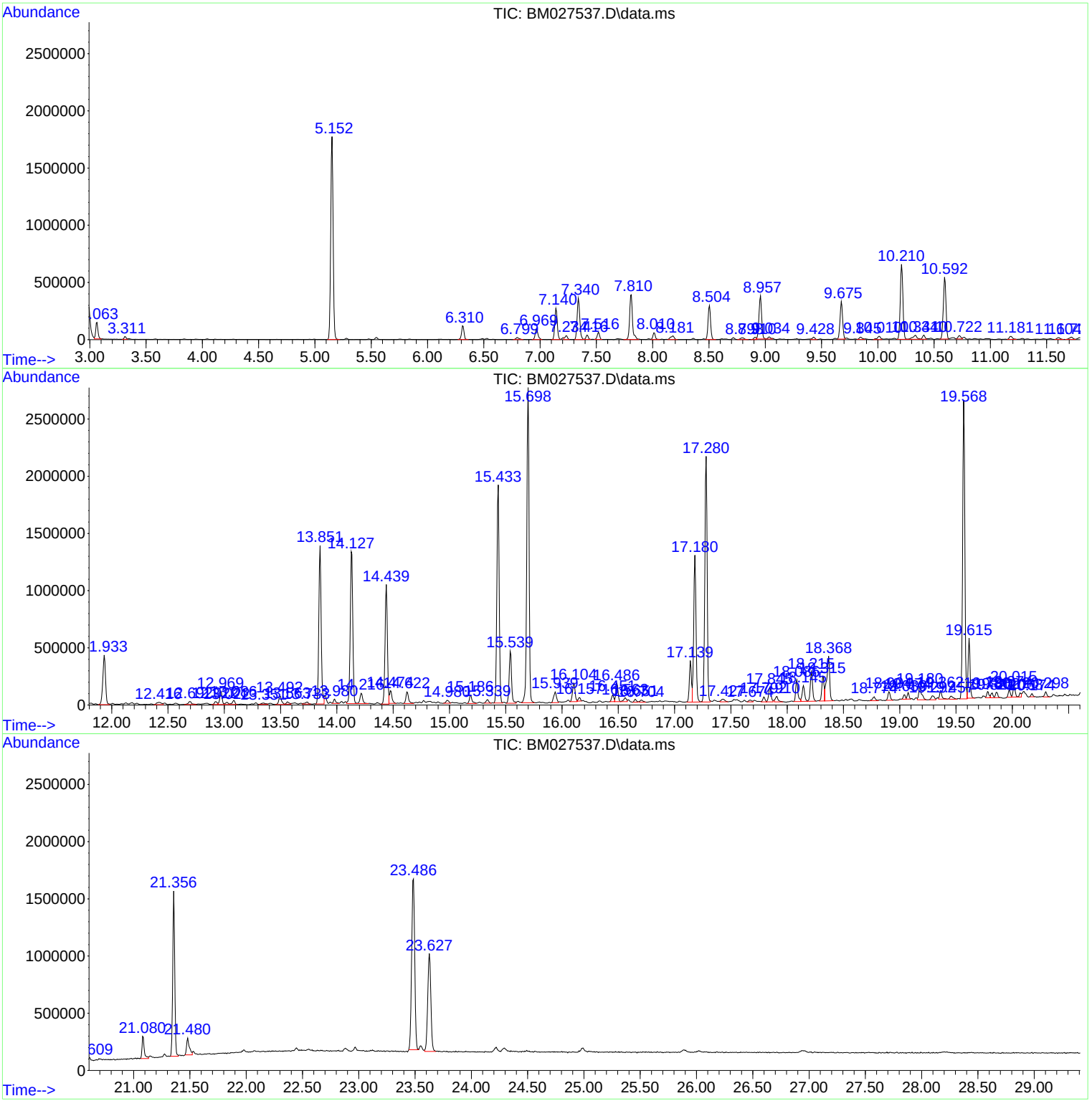
Sum of corrected areas: 44686353

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092520\
Data File : BM027537.D
Acq On : 25 Sep 2020 20:05
Operator : JU/CG
Sample : L4135-06
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :**
BNA_M
**ClientSampleId :**
BG480

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM092420MA.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092520\
Data File : BM027537.D
Acq On : 25 Sep 2020 20:05
Operator : JU/CG
Sample : L4135-06
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** BG480

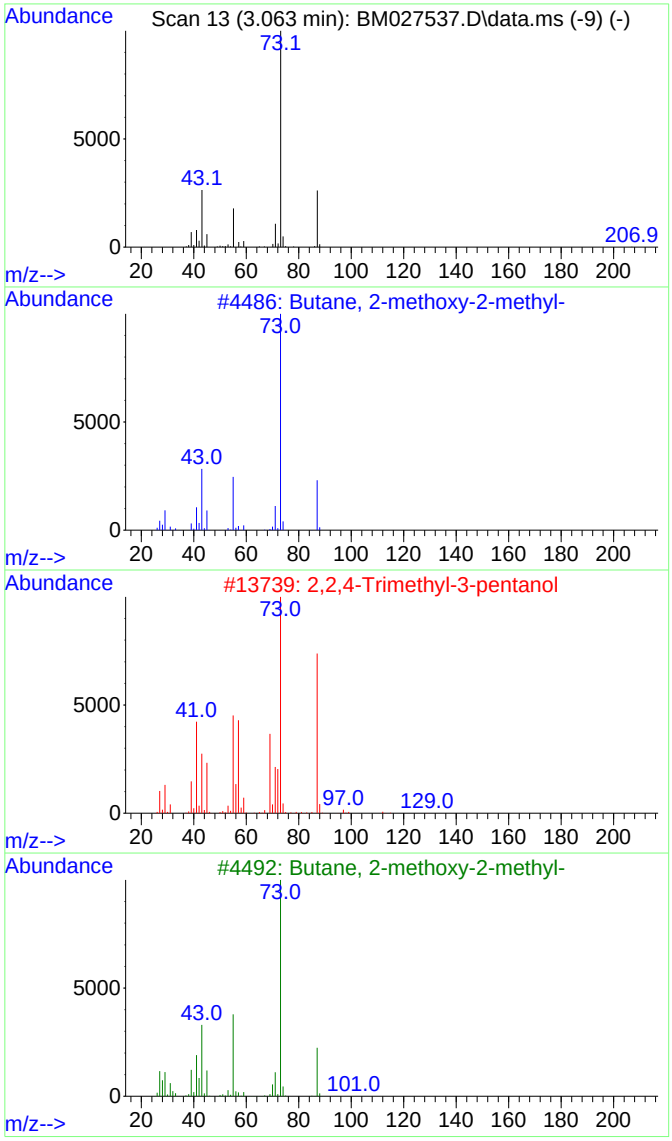
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM092420MA.M
Quant Title : SVOA CALIBRATION

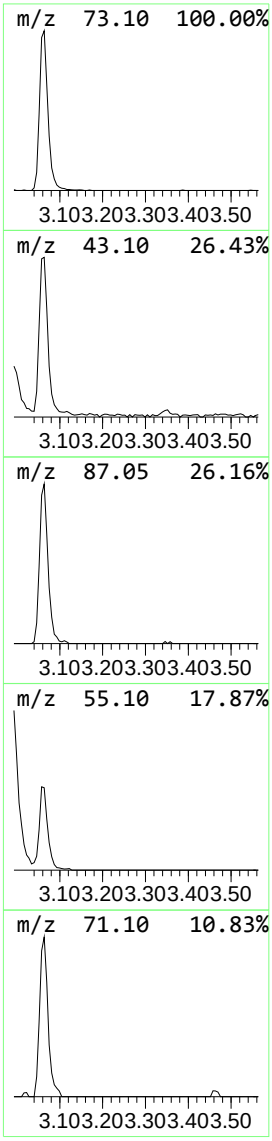
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 1 unknown-01 Concentration Rank 4

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 3.063 | 5.69 ng/ul | 197427 | 1,4-Dichlorobenzene-d4 | 7.810 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Butane, 2-methoxy-2-methyl- | 102 | C6H14O | 000994-05-8 | 47 |
| 2 | 2,2,4-Trimethyl-3-pentanol | 130 | C8H18O | 005162-48-1 | 40 |
| 3 | Butane, 2-methoxy-2-methyl- | 102 | C6H14O | 000994-05-8 | 39 |
| 4 | N-Ethylformamide | 73 | C3H7NO | 000627-45-2 | 9 |
| 5 | Silane, tetramethyl- | 88 | C4H12Si | 000075-76-3 | 9 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092520\
Data File : BM027537.D
Acq On : 25 Sep 2020 20:05
Operator : JU/CG
Sample : L4135-06
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** BG480

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM092420MA.M
Quant Title : SVOA CALIBRATION

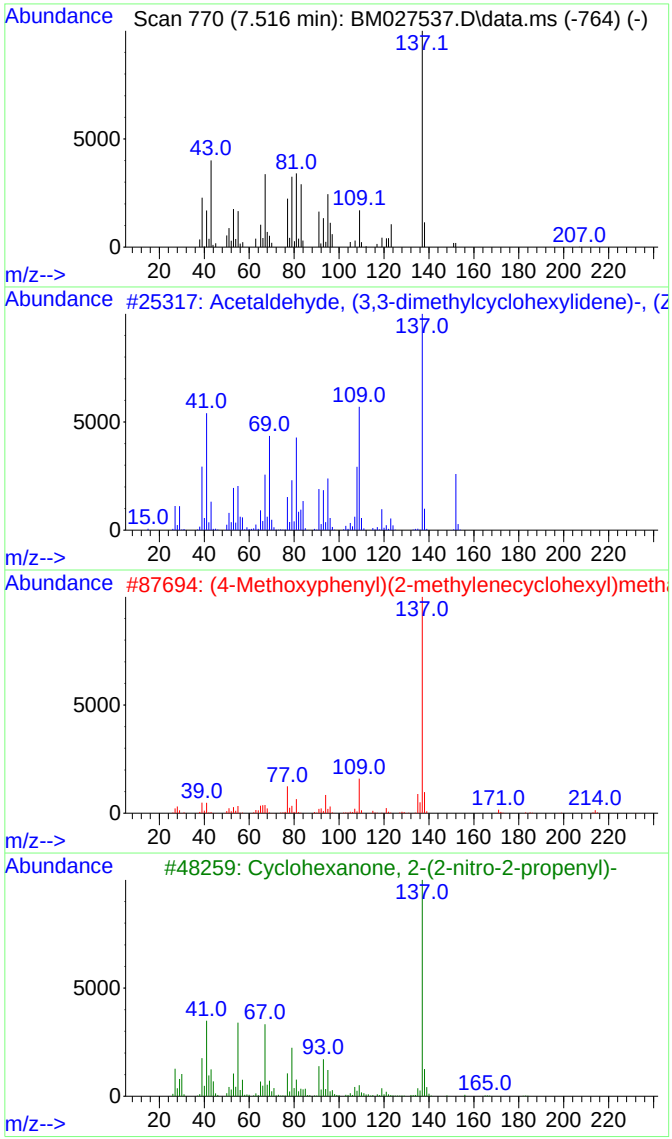
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

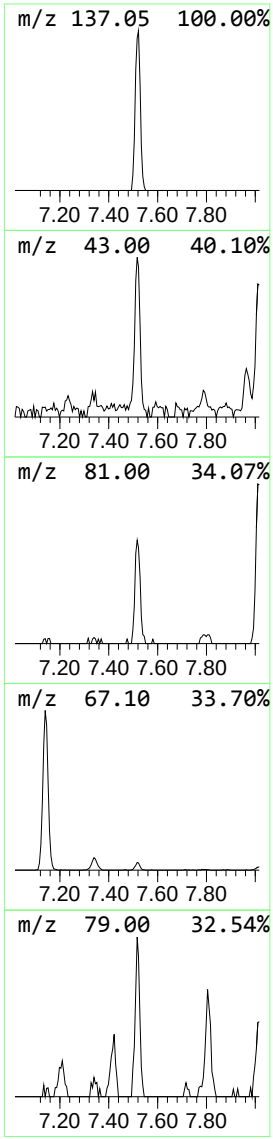
**********

Peak Number 4 unknown-02 Concentration Rank 12

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 7.516 | 2.69 ng/ul | 93466 | 1,4-Dichlorobenzene-d4 | 7.810 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Acetaldehyde, (3,3-dimethylcyclo... | 152 | C10H16O | 026532-24-1 | 45 |
| 2 | (4-Methoxyphenyl)(2-methylenecyc... | 232 | C15H20O2 | 1000191-91-2 | 43 |
| 3 | Cyclohexanone, 2-(2-nitro-2-prop... | 183 | C9H13NO3 | 078551-08-3 | 42 |
| 4 | Methanethioamide, N-phenyl- | 137 | C7H7NS | 000637-51-4 | 38 |
| 5 | 1,2,4-Triazolo[4,3-b][1,2,4,5]te... | 137 | C3H3N7 | 220696-99-1 | 37 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092520\
Data File : BM027537.D
Acq On : 25 Sep 2020 20:05
Operator : JU/CG
Sample : L4135-06
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** BG480

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM092420MA.M
Quant Title : SVOA CALIBRATION

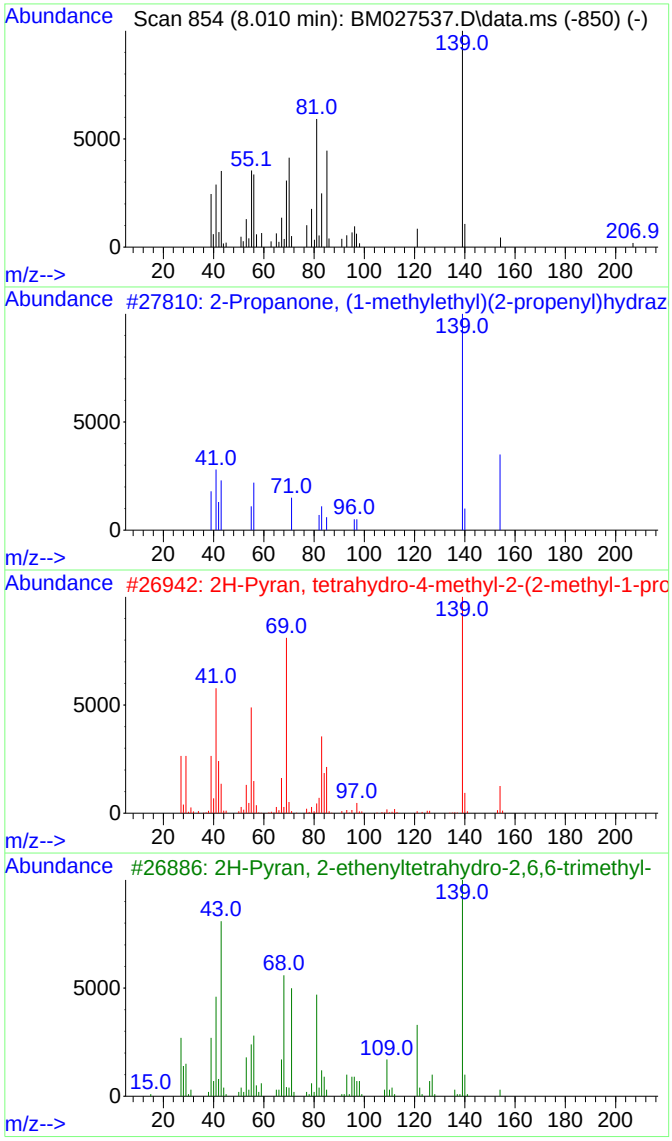
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

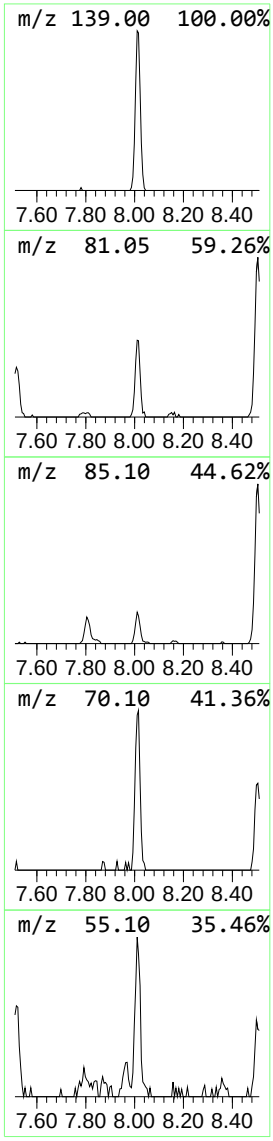
**********

Peak Number 5 unknown-03 Concentration Rank 13

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 8.010 | 2.53 ng/ul | 87909 | 1,4-Dichlorobenzene-d4 | 7.810 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2-Propanone, (1-methylethyl)(2-p... | 154 | C9H18N2 | 110213-09-7 | 38 |
| 2 | 2H-Pyran, tetrahydro-4-methyl-2-... | 154 | C10H18O | 016409-43-1 | 37 |
| 3 | 2H-Pyran, 2-ethenyltetrahydro-2,... | 154 | C10H18O | 007392-19-0 | 32 |
| 4 | 5-Dimethylamino-furan-2-carbalde... | 139 | C7H9NO2 | 1000317-69-8 | 27 |
| 5 | Pentan-2-yl 3-chlorobenzoate | 226 | C12H15ClO2 | 097989-35-0 | 25 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092520\
Data File : BM027537.D
Acq On : 25 Sep 2020 20:05
Operator : JU/CG
Sample : L4135-06
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :**
BNA_M
**ClientSampleId :**
BG480

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM092420MA.M
Quant Title : SVOA CALIBRATION

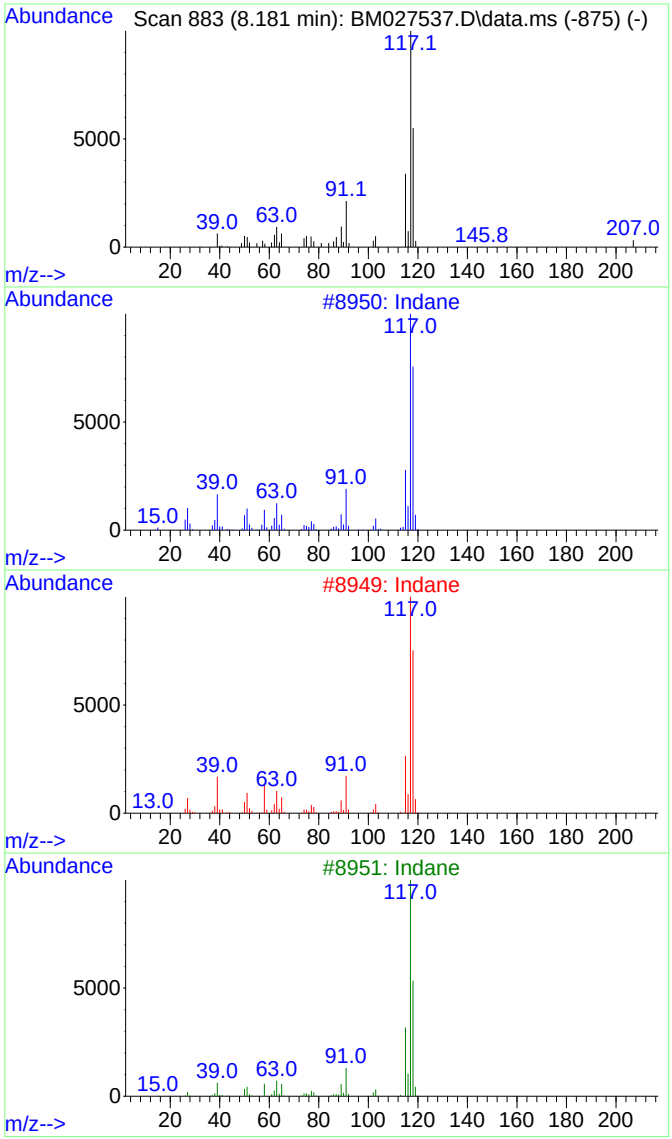
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

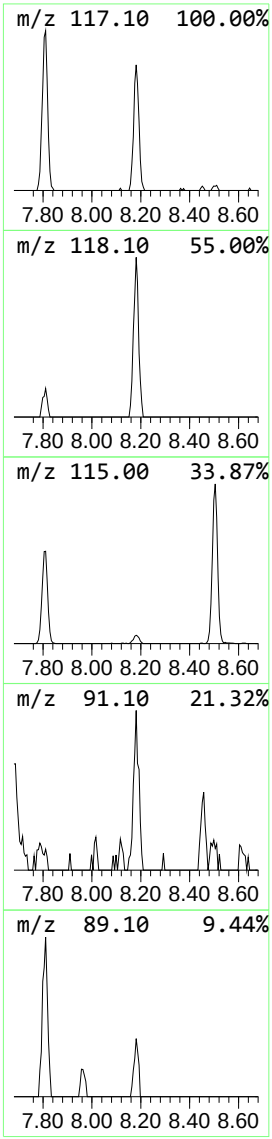
**********

Peak Number 6 Indane Concentration Rank 19

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 8.181 | 2.08 ng/ul | 72116 | 1,4-Dichlorobenzene-d4 | 7.810 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Indane | 118 | C9H10 | 000496-11-7 | 83 |
| 2 | Indane | 118 | C9H10 | 000496-11-7 | 74 |
| 3 | Indane | 118 | C9H10 | 000496-11-7 | 74 |
| 4 | Benzene, cyclopropyl- | 118 | C9H10 | 000873-49-4 | 64 |
| 5 | Tetracyclo[3.3.1.0(2,8).0(4,6)]-... | 118 | C9H10 | 1000191-13-7 | 59 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092520\
Data File : BM027537.D
Acq On : 25 Sep 2020 20:05
Operator : JU/CG
Sample : L4135-06
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :**
BNA_M
**ClientSampleId :**
BG480

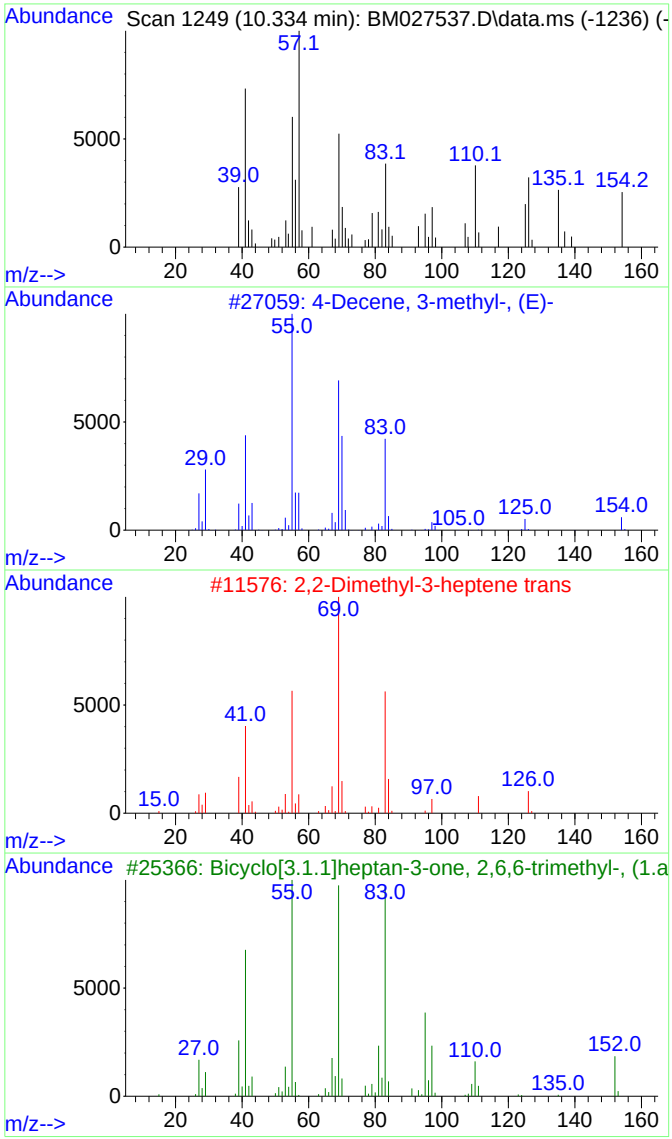
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM092420MA.M
Quant Title : SVOA CALIBRATION

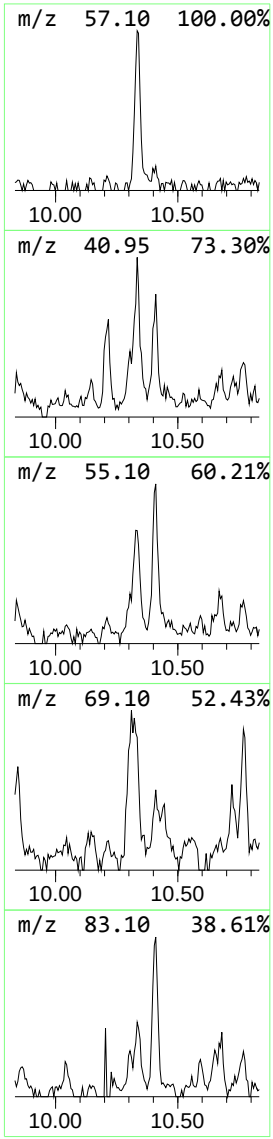
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 18

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 10.334 | 2.09 ng/ul | 92109 | Naphthalene-d8 | 10.592 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 4-Decene, 3-methyl-, (E)- | 154 | C11H22 | 062338-47-0 | 30 |
| 2 | 2,2-Dimethyl-3-heptene trans | 126 | C9H18 | 019550-75-5 | 25 |
| 3 | Bicyclo[3.1.1]heptan-3-one, 2,6,... | 152 | C10H16O | 015358-88-0 | 22 |
| 4 | Bicyclo[3.1.1]heptan-3-one, 2,6,... | 152 | C10H16O | 000547-60-4 | 22 |
| 5 | 3-Heptene, 2,6-dimethyl- | 126 | C9H18 | 002738-18-3 | 22 |



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Data File : BM027537.D
Acq On : 25 Sep 2020 20:05
Operator : JU/CG
Sample : L4135-06
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** BG480

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM092420MA.M
Quant Title : SVOA CALIBRATION

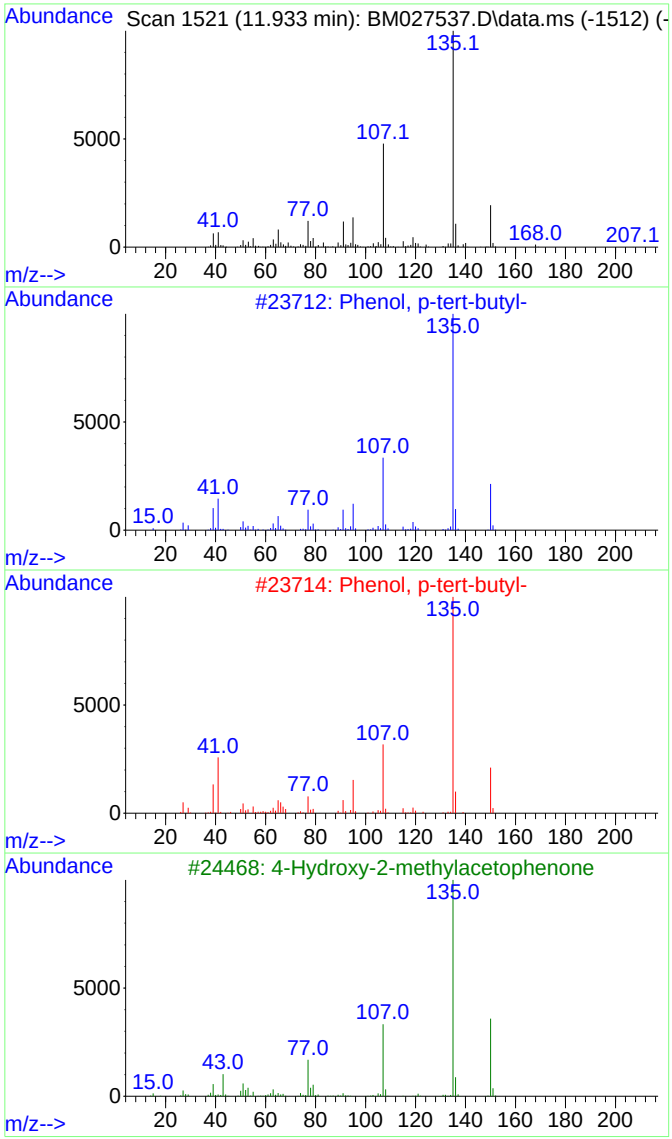
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

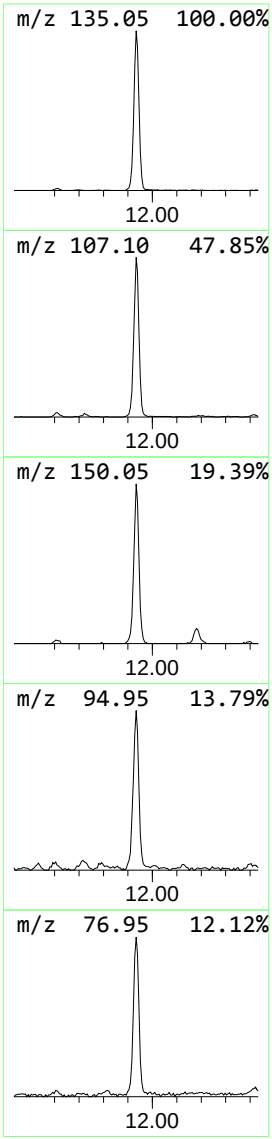
**********

Peak Number 8 Phenol, p-tert-butyl- Concentration Rank 2

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 11.934 | 16.41 ng/ul | 722166 | Naphthalene-d8 | 10.592 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Phenol, p-tert-butyl- | 150 | C10H14O | 000098-54-4 | 94 |
| 2 | Phenol, p-tert-butyl- | 150 | C10H14O | 000098-54-4 | 90 |
| 3 | 4-Hydroxy-2-methylacetophenone | 150 | C9H10O2 | 000875-59-2 | 87 |
| 4 | Phenol, p-tert-butyl- | 150 | C10H14O | 000098-54-4 | 87 |
| 5 | Phenol, m-tert-butyl- | 150 | C10H14O | 000585-34-2 | 81 |



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Data File : BM027537.D
Acq On : 25 Sep 2020 20:05
Operator : JU/CG
Sample : L4135-06
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** BG480

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM092420MA.M
Quant Title : SVOA CALIBRATION

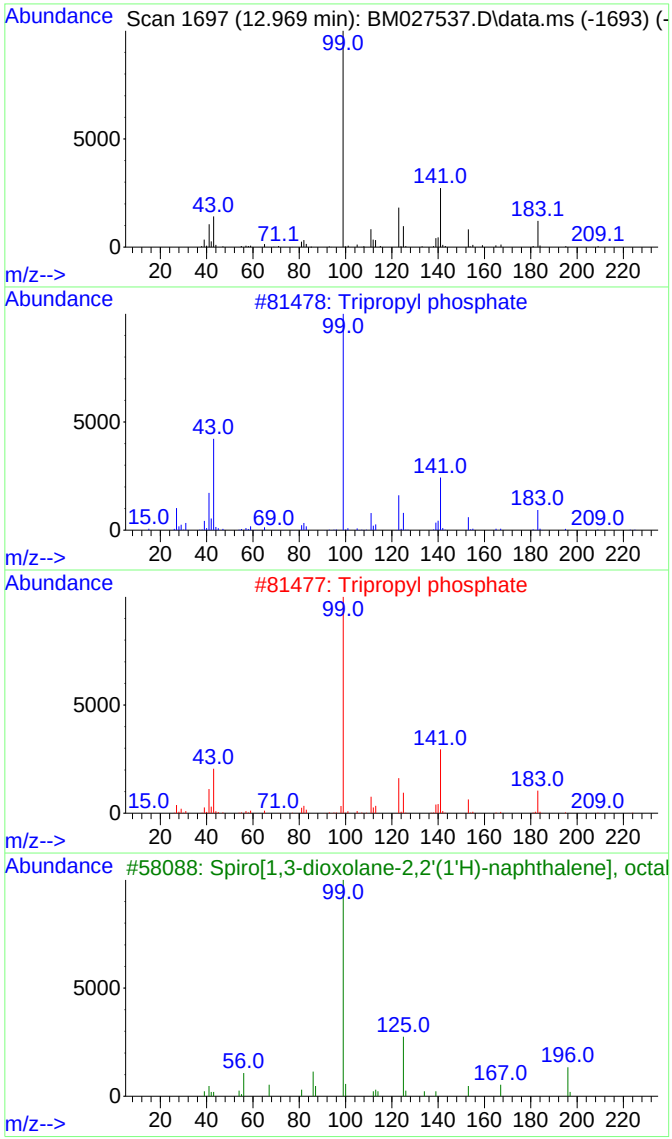
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

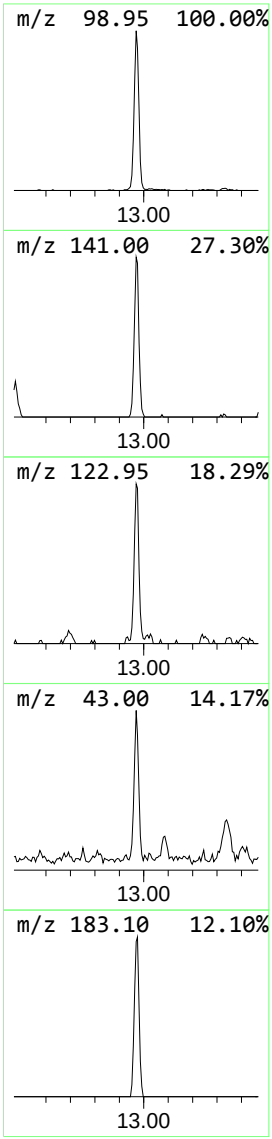
**********

Peak Number 9 Tripropyl phosphate Concentration Rank 20

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 12.969 | 2.06 ng/ul | 155960 | Acenaphthene-d10 | 14.439 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Tripropyl phosphate | 224 | C9H21O4P | 000513-08-6 | 90 |
| 2 | Tripropyl phosphate | 224 | C9H21O4P | 000513-08-6 | 90 |
| 3 | Spiro[1,3-dioxolane-2,2'(1'H)-na... | 196 | C12H20O2 | 006857-86-9 | 28 |
| 4 | n-Butyl methylphosphonofluoridate | 154 | C5H12FO2P | 000352-63-6 | 9 |
| 5 | 2-Methyl-2-propyl methylphosphon... | 154 | C5H12FO2P | 013273-12-6 | 9 |



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Data File : BM027537.D
Acq On : 25 Sep 2020 20:05
Operator : JU/CG
Sample : L4135-06
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** BG480

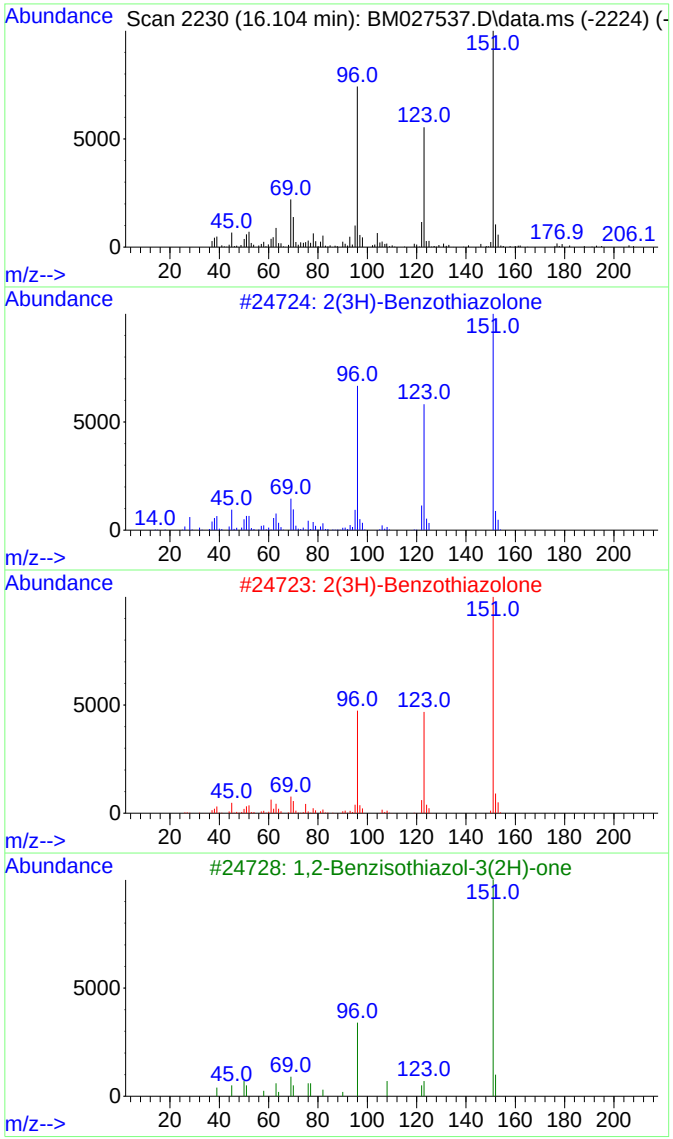
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM092420MA.M
Quant Title : SVOA CALIBRATION

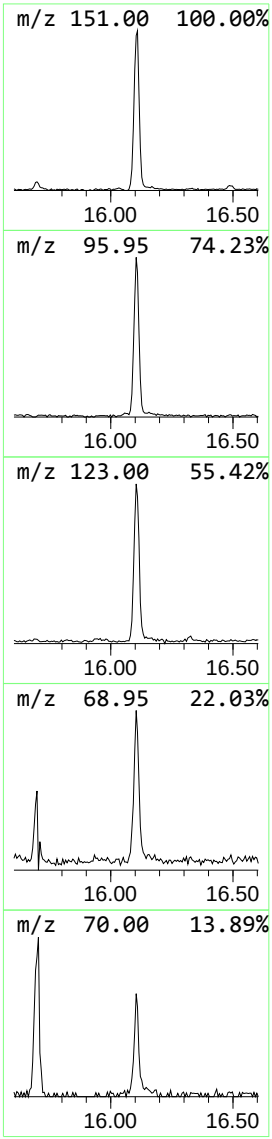
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 10 2(3H)-Benzothiazolone Concentration Rank 11

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 16.104 | 2.77 ng/ul | 249139 | Phenanthrene-d10 | 17.180 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2(3H)-Benzothiazolone | 151 | C7H5NOS | 000934-34-9 | 94 |
| 2 | 2(3H)-Benzothiazolone | 151 | C7H5NOS | 000934-34-9 | 94 |
| 3 | 1,2-Benzisothiazol-3(2H)-one | 151 | C7H5NOS | 002634-33-5 | 35 |
| 4 | Thieno[2,3-b]pyridine-N-oxide | 151 | C7H5NOS | 025557-50-0 | 32 |
| 5 | 4-Amino-2,1,3-benzothiadiazole | 151 | C6H5N3S | 000767-64-6 | 27 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092520\
Data File : BM027537.D
Acq On : 25 Sep 2020 20:05
Operator : JU/CG
Sample : L4135-06
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :**
BNA_M
**ClientSampleId :**
BG480

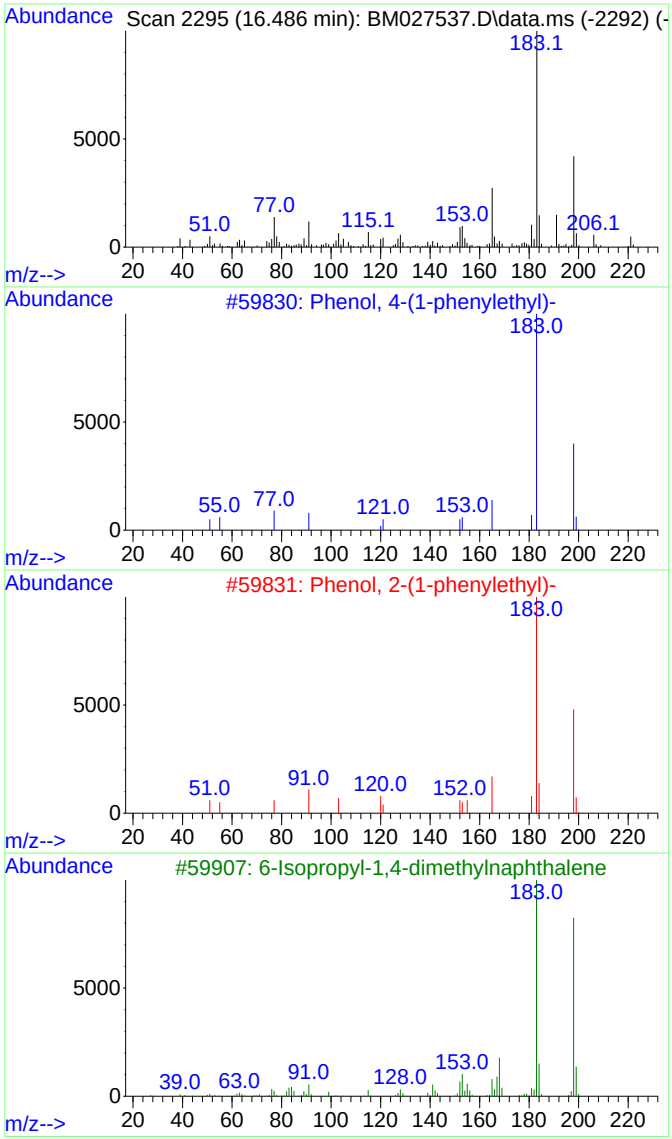
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM092420MA.M
Quant Title : SVOA CALIBRATION

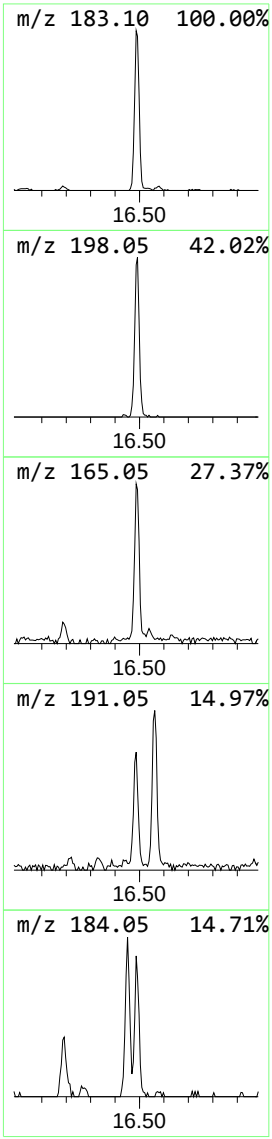
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 11 Phenol, 4-(1-phenylethyl)- Concentration Rank 14

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 16.486 | 2.28 ng/ul | 205693 | Phenanthrene-d10 | 17.180 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Phenol, 4-(1-phenylethyl)- | 198 | C14H14O | 001988-89-2 | 81 |
| 2 | Phenol, 2-(1-phenylethyl)- | 198 | C14H14O | 004237-44-9 | 81 |
| 3 | 6-Isopropyl-1,4-dimethylnaphthalene | 198 | C15H18 | 000489-77-0 | 74 |
| 4 | Azulene, 1,4-dimethyl-7-(1-methy... | 198 | C15H18 | 000489-84-9 | 72 |
| 5 | Azulene, 1,4-dimethyl-7-(1-methy... | 198 | C15H18 | 000489-84-9 | 64 |



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Acq On : 25 Sep 2020 20:05
Operator : JU/CG
Sample : L4135-06
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** BG480

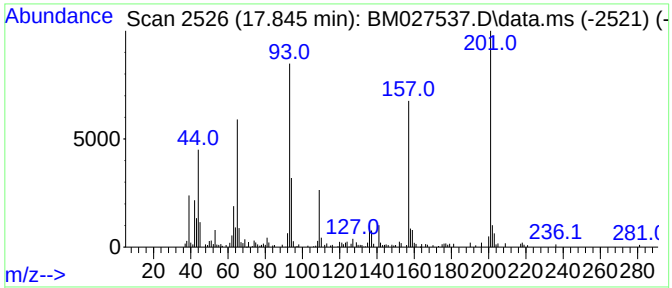
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM092420MA.M
Quant Title : SVOA CALIBRATION

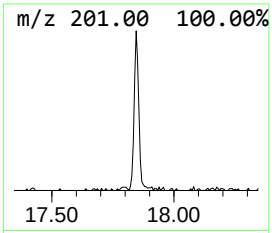
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

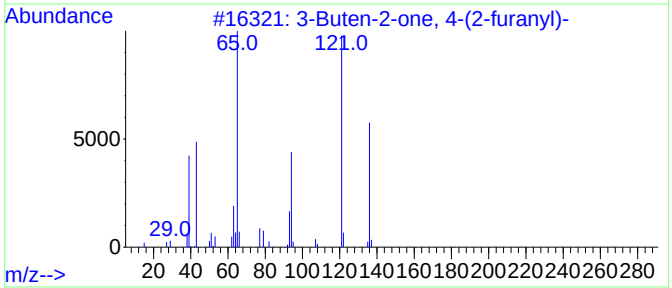
**********
Peak Number 12 unknown-04 Concentration Rank 17

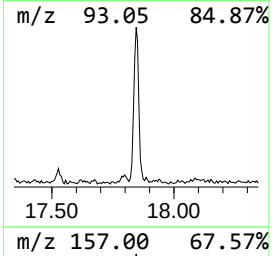
| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 17.845 | 2.12 ng/ul | 190592 | Phenanthrene-d10 | 17.180 |

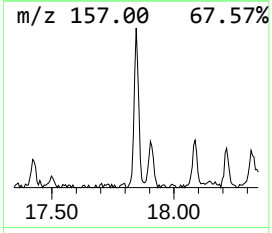
| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 3-Buten-2-one, 4-(2-furanyl)- | 136 | C8H8O2 | 000623-15-4 | 16 |
| 2 | 2-Pyridylacetamide | 136 | C7H8N2O | 003724-16-1 | 14 |
| 3 | 2,Beta-dinitrostyrene | 194 | C8H6N2O4 | 003156-39-6 | 12 |
| 4 | Benzene, 1-methyl-2-nitroso- | 121 | C7H7NO | 000611-23-4 | 12 |
| 5 | Acetamide, N-[4-(1H-imidazol-1-y... | 201 | C11H11N3O | 1000337-71-7 | 11 |

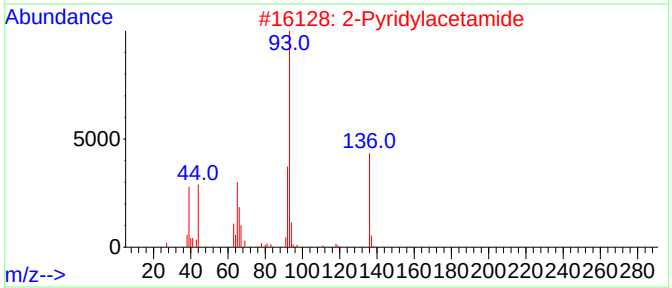


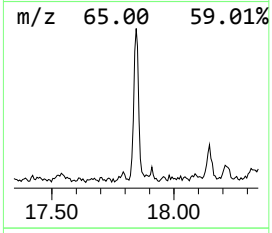



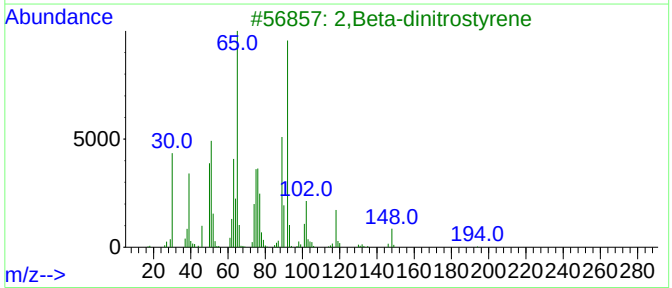



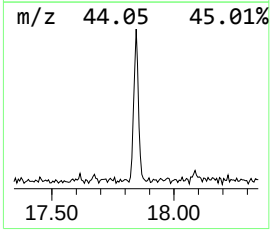



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Data File : BM027537.D
Acq On : 25 Sep 2020 20:05
Operator : JU/CG
Sample : L4135-06
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :**
BNA_M
**ClientSampleId :**
BG480

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM092420MA.M
Quant Title : SVOA CALIBRATION

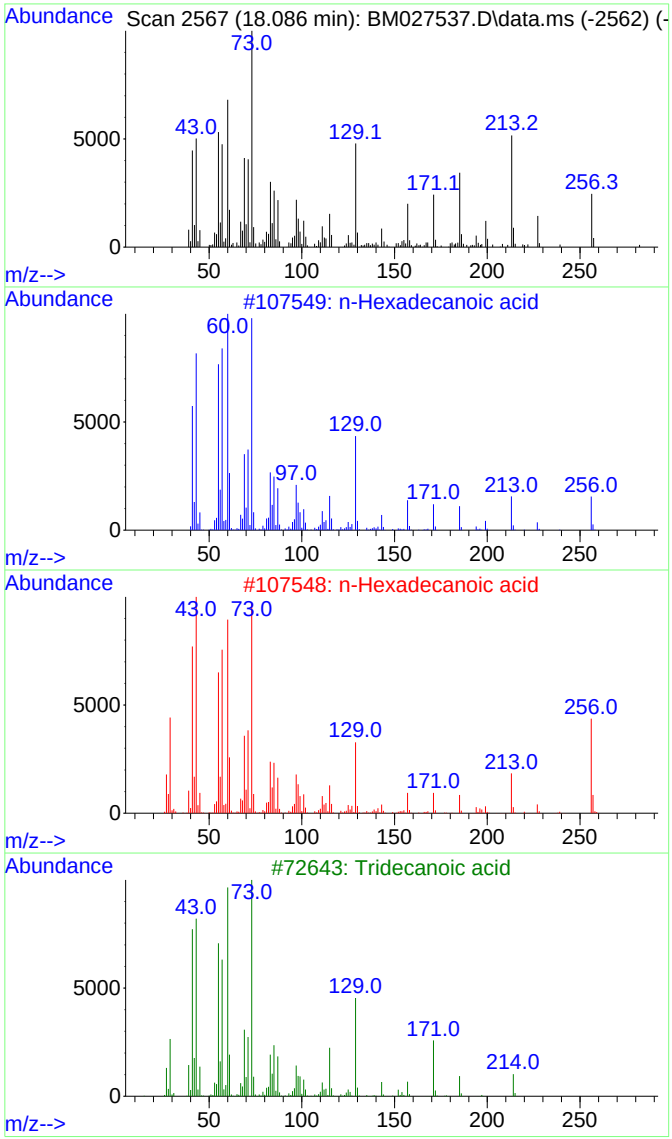
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

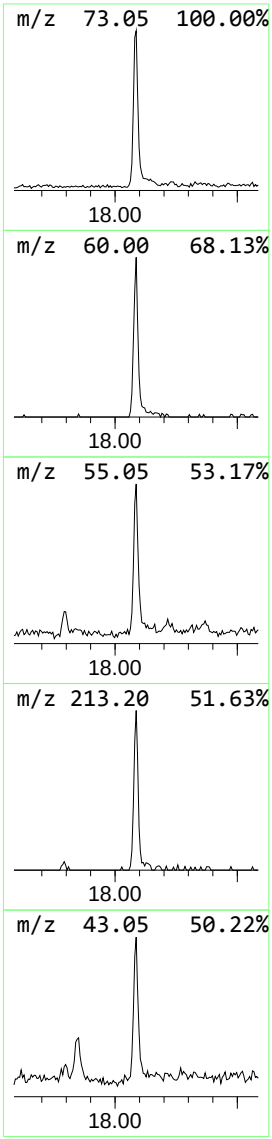
**********

Peak Number 13 n-Hexadecanoic acid Concentration Rank 9

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 18.086 | 2.93 ng/ul | 263857 | Phenanthrene-d10 | 17.180 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 96 |
| 2 | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 96 |
| 3 | Tridecanoic acid | 214 | C13H26O2 | 000638-53-9 | 83 |
| 4 | Dodecanoic acid | 200 | C12H24O2 | 000143-07-7 | 50 |
| 5 | Tridecanoic acid | 214 | C13H26O2 | 000638-53-9 | 47 |



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Acq On : 25 Sep 2020 20:05
Operator : JU/CG
Sample : L4135-06
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** BG480

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM092420MA.M
Quant Title : SVOA CALIBRATION

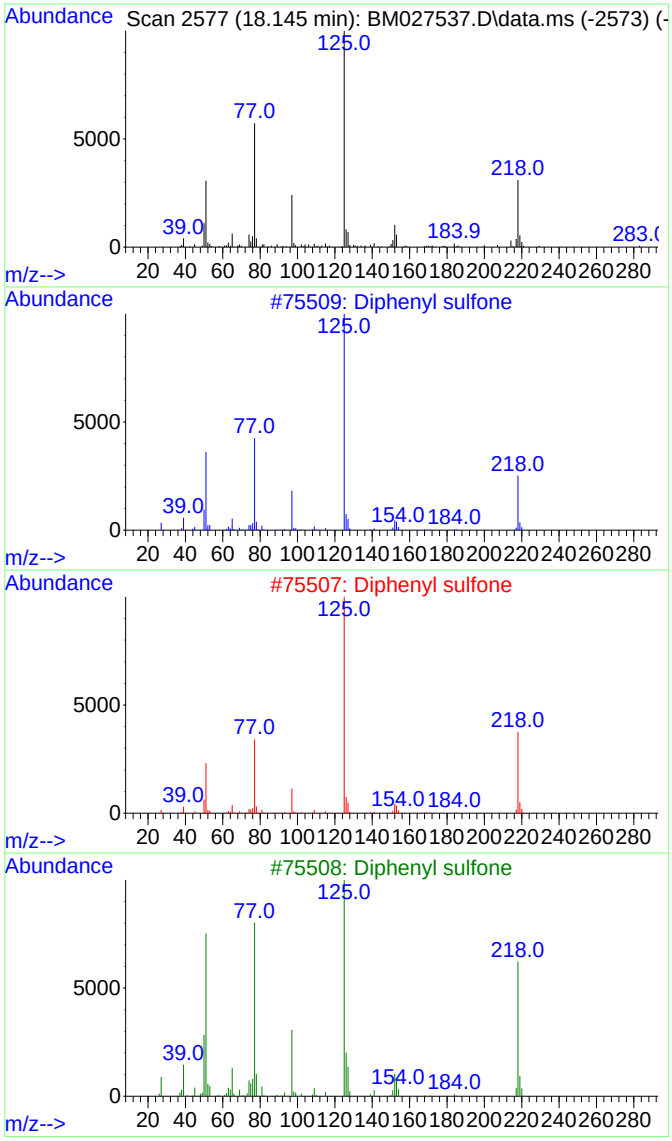
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

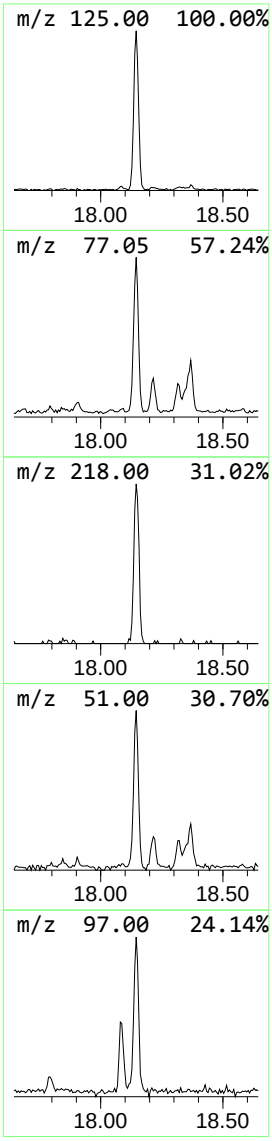
**********

Peak Number 14 Diphenyl sulfone Concentration Rank 15

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 18.145 | 2.28 ng/ul | 205039 | Phenanthrene-d10 | 17.180 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Diphenyl sulfone | 218 | C12H10O2S | 000127-63-9 | 91 |
| 2 | Diphenyl sulfone | 218 | C12H10O2S | 000127-63-9 | 74 |
| 3 | Diphenyl sulfone | 218 | C12H10O2S | 000127-63-9 | 60 |
| 4 | Benzene, (ethenylsulfonyl)- | 168 | C8H8O2S | 005535-48-8 | 53 |
| 5 | 2-Propenoic acid, 3-(phenylsulfo... | 226 | C10H10O4S | 063068-63-3 | 53 |



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Data File : BM027537.D
Acq On : 25 Sep 2020 20:05
Operator : JU/CG
Sample : L4135-06
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :**
BNA_M
**ClientSampleId :**
BG480

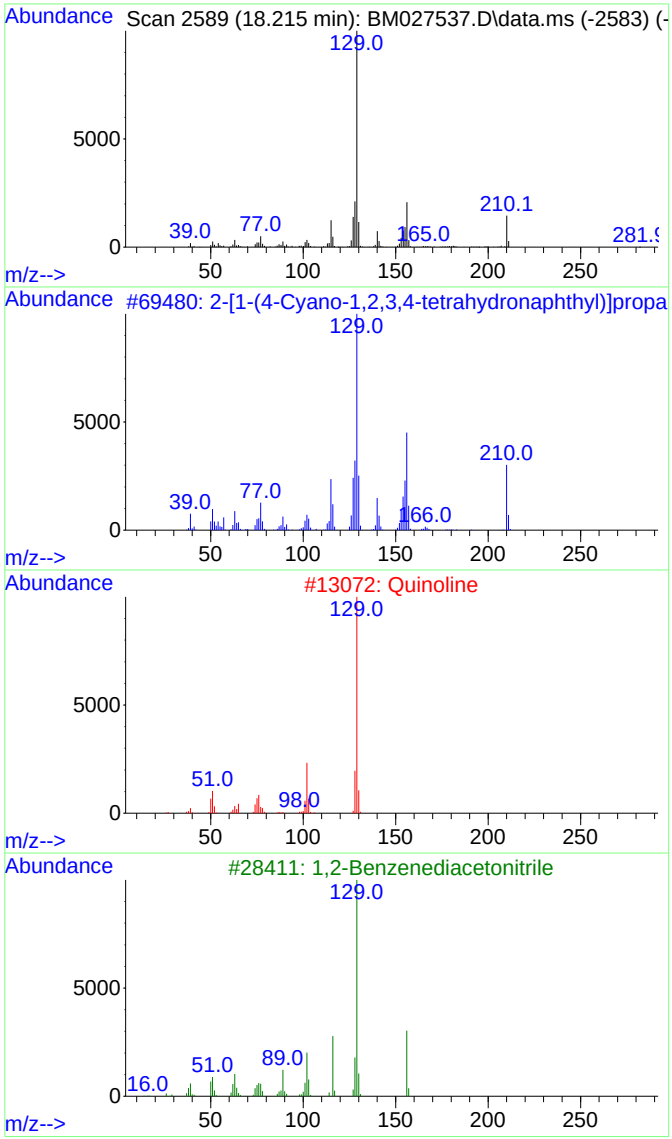
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM092420MA.M
Quant Title : SVOA CALIBRATION

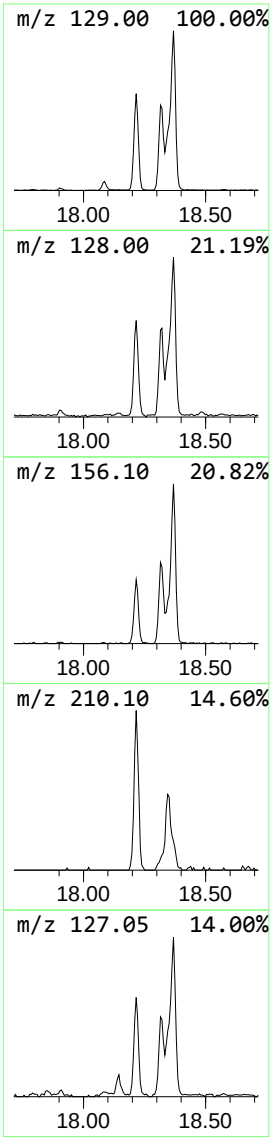
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 15 2-[1-(4-Cyano-1,2,3,4-tetra... Concentration Rank 6

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 18.215 | 3.98 ng/ul | 358681 | Phenanthrene-d10 | 17.180 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2-[1-(4-Cyano-1,2,3,4-tetrahydro... | 210 | C14H14N2 | 057964-39-3 | 91 |
| 2 | Quinoline | 129 | C9H7N | 000091-22-5 | 52 |
| 3 | 1,2-Benzenediacetonitrile | 156 | C10H8N2 | 000613-73-0 | 50 |
| 4 | 3-[1-(4-Cyano-1,2,3,4-tetrahydro... | 210 | C14H14N2 | 057964-40-6 | 49 |
| 5 | Quinolinium, 1-ethyl-, iodide | 285 | C11H12IN | 000634-35-5 | 45 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092520\
Data File : BM027537.D
Acq On : 25 Sep 2020 20:05
Operator : JU/CG
Sample : L4135-06
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :**
BNA_M
**ClientSampleId :**
BG480

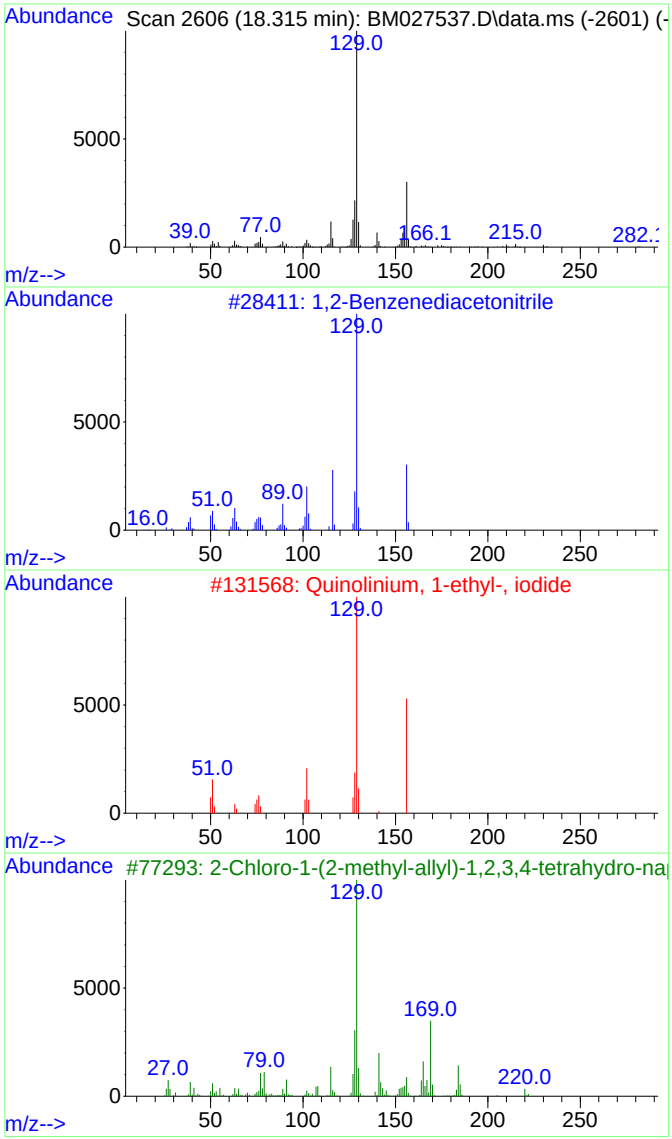
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM092420MA.M
Quant Title : SVOA CALIBRATION

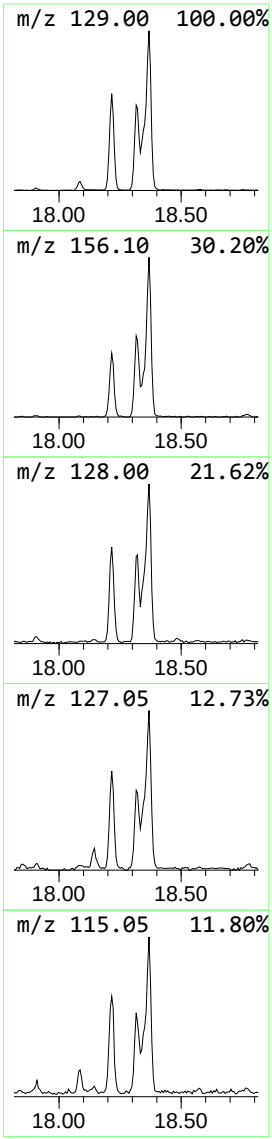
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 16 1,2-Benzenediacetonitrile Concentration Rank 7

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 18.315 | 3.50 ng/ul | 314769 | Phenanthrene-d10 | 17.180 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1,2-Benzenediacetonitrile | 156 | C10H8N2 | 000613-73-0 | 64 |
| 2 | Quinolinium, 1-ethyl-, iodide | 285 | C11H12IN | 000634-35-5 | 56 |
| 3 | 2-Chloro-1-(2-methyl-allyl)-1,2,... | 220 | C14H17Cl | 1000185-82-6 | 56 |
| 4 | Isoquinoline | 129 | C9H7N | 000119-65-3 | 42 |
| 5 | 3-[1-(4-Cyano-1,2,3,4-tetrahydro... | 210 | C14H14N2 | 057964-40-6 | 40 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092520\
Data File : BM027537.D
Acq On : 25 Sep 2020 20:05
Operator : JU/CG
Sample : L4135-06
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :**
BNA_M
**ClientSampleId :**
BG480

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM092420MA.M
Quant Title : SVOA CALIBRATION

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

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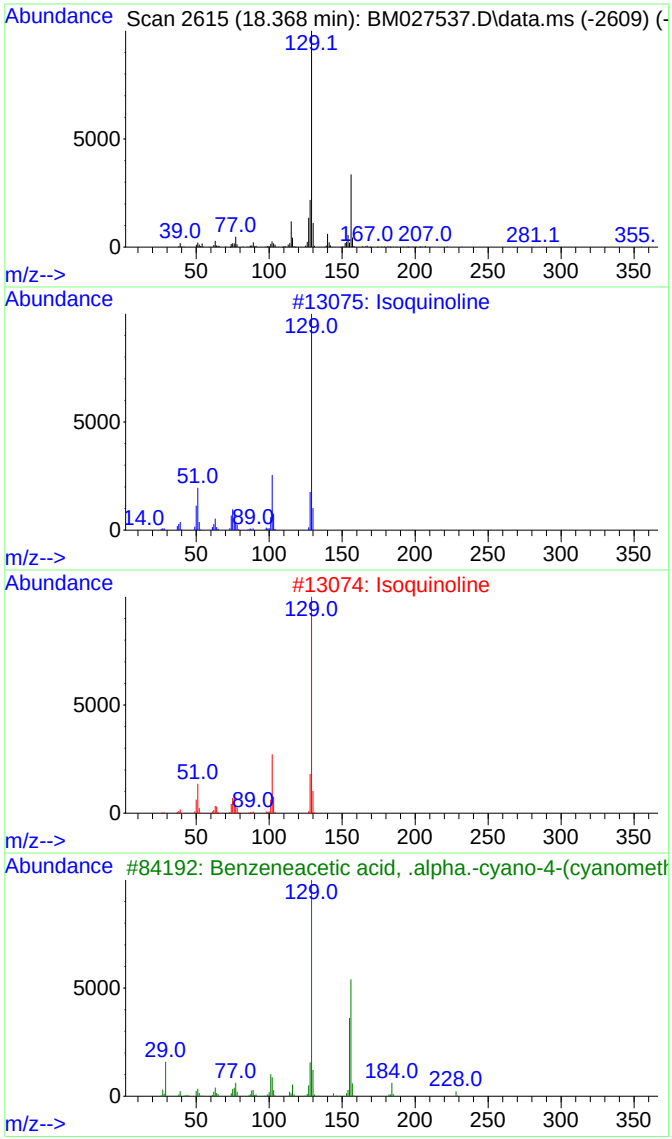
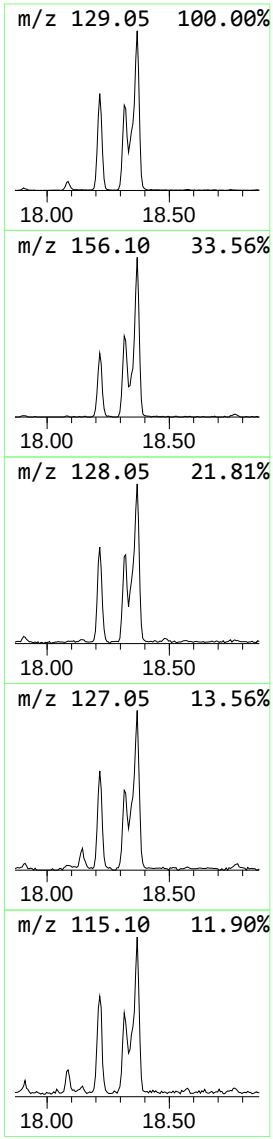
Peak Number 17 Isoquinoline Concentration Rank 3

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 18.368 | 7.48 ng/ul | 673141 | Phenanthrene-d10 | 17.180 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Isoquinoline | 129 | C9H7N | 000119-65-3 | 52 |
| 2 | Isoquinoline | 129 | C9H7N | 000119-65-3 | 50 |
| 3 | Benzeneacetic acid, .alpha.-cyan... | 228 | C13H12N2O2 | 134882-04-5 | 45 |
| 4 | Naphthalene, 1,2-dihydro-2-methyl- | 144 | C11H12 | 021564-79-4 | 43 |
| 5 | 2-Propenenitrile, 3-phenyl-, (E)- | 129 | C9H7N | 001885-38-7 | 40 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092520\
Data File : BM027537.D
Acq On : 25 Sep 2020 20:05
Operator : JU/CG
Sample : L4135-06
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** BG480

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM092420MA.M
Quant Title : SVOA CALIBRATION

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

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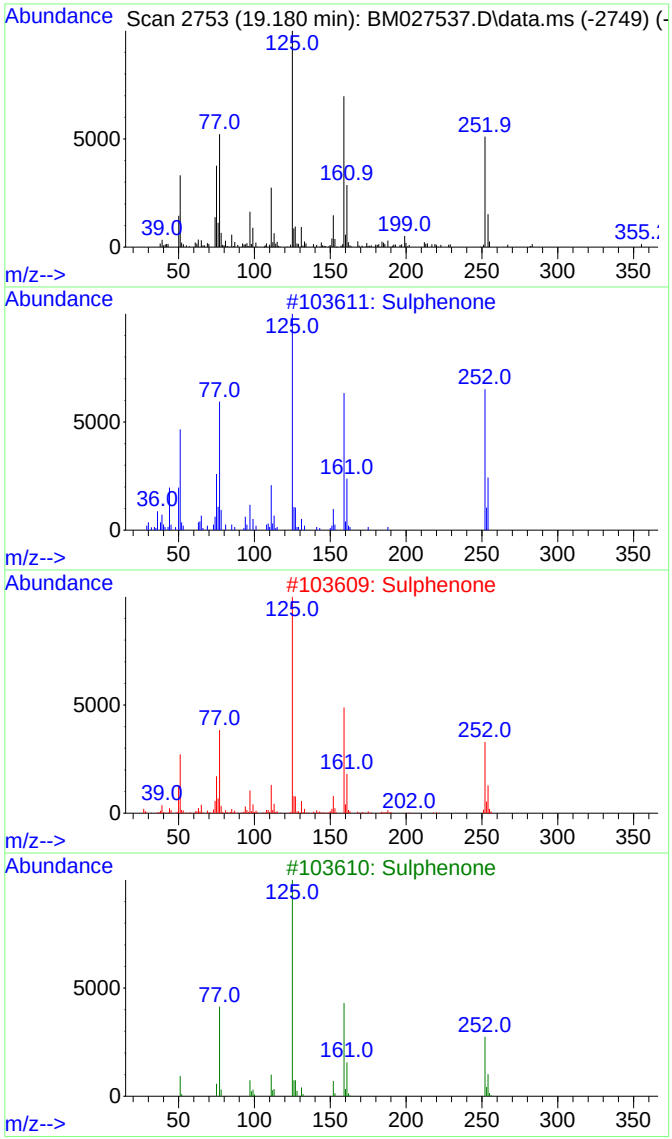
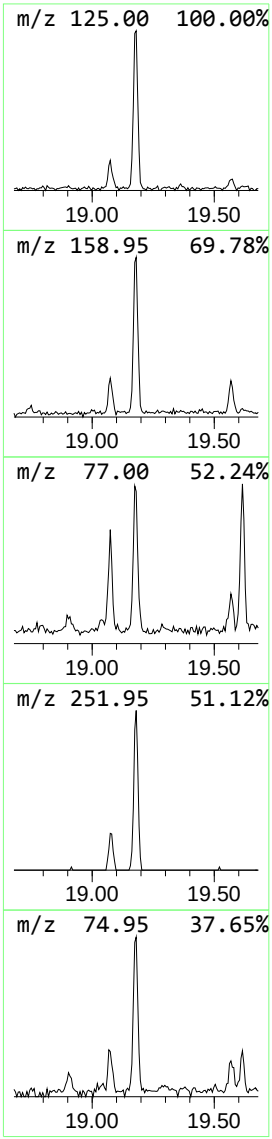
Peak Number 18 Sulphenone Concentration Rank 16

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 19.180 | 2.16 ng/ul | 194363 | Phenanthrene-d10 | 17.180 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Sulphenone | 252 | C12H9ClO2S | 000080-00-2 | 95 |
| 2 | Sulphenone | 252 | C12H9ClO2S | 000080-00-2 | 83 |
| 3 | Sulphenone | 252 | C12H9ClO2S | 000080-00-2 | 64 |
| 4 | Sulphenone | 252 | C12H9ClO2S | 000080-00-2 | 64 |
| 5 | Bis[phenylsulfonyl]-4-carboxyl p... | 416 | C20H16O6S2 | 081269-08-1 | 35 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092520\
Data File : BM027537.D
Acq On : 25 Sep 2020 20:05
Operator : JU/CG
Sample : L4135-06
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** BG480

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM092420MA.M
Quant Title : SVOA CALIBRATION

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

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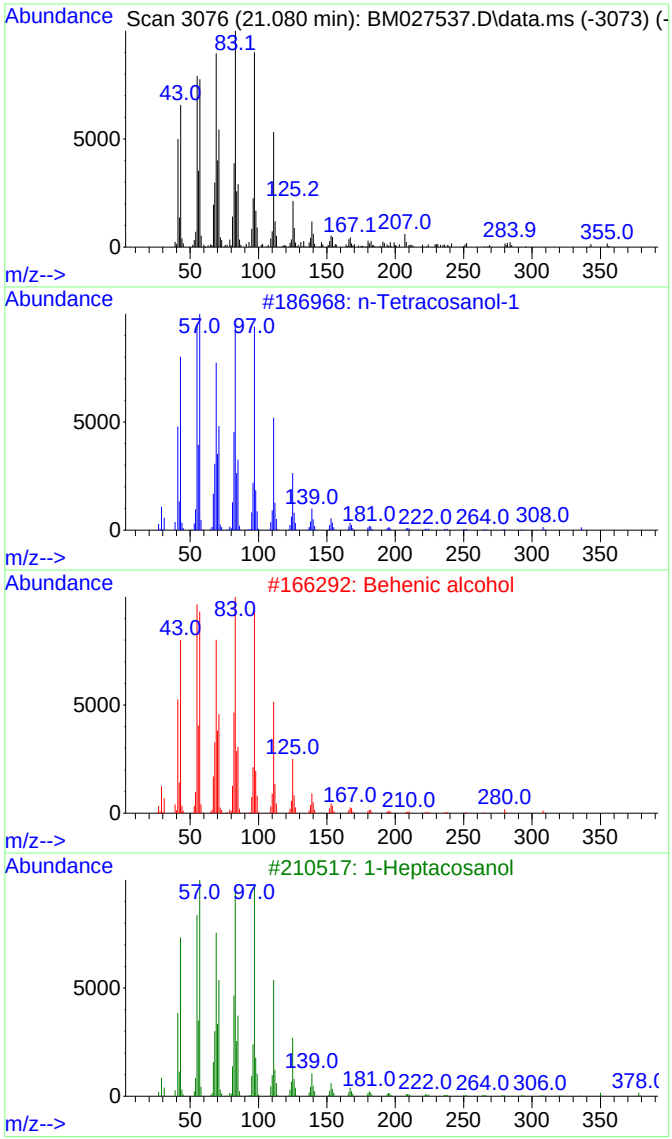
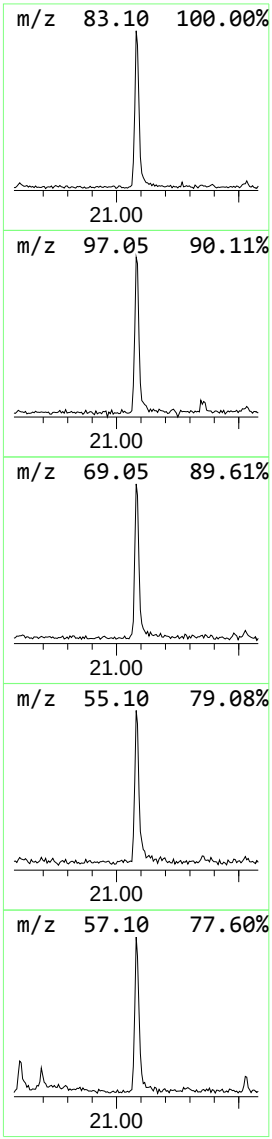
Peak Number 19 n-Tetracosanol-1 Concentration Rank 8

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 21.080 | 3.04 ng/ul | 270324 | Chrysene-d12 | 21.356 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | n-Tetracosanol-1 | 354 | C24H50O | 000506-51-4 | 95 |
| 2 | Behenic alcohol | 326 | C22H46O | 000661-19-8 | 95 |
| 3 | 1-Heptacosanol | 396 | C27H56O | 002004-39-9 | 93 |
| 4 | n-Nonadecanol-1 | 284 | C19H40O | 001454-84-8 | 93 |
| 5 | Octacosanol | 410 | C28H58O | 000557-61-9 | 90 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092520\
Data File : BM027537.D
Acq On : 25 Sep 2020 20:05
Operator : JU/CG
Sample : L4135-06
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :**
BNA_M
**ClientSampleId :**
BG480

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM092420MA.M
Quant Title : SVOA CALIBRATION

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

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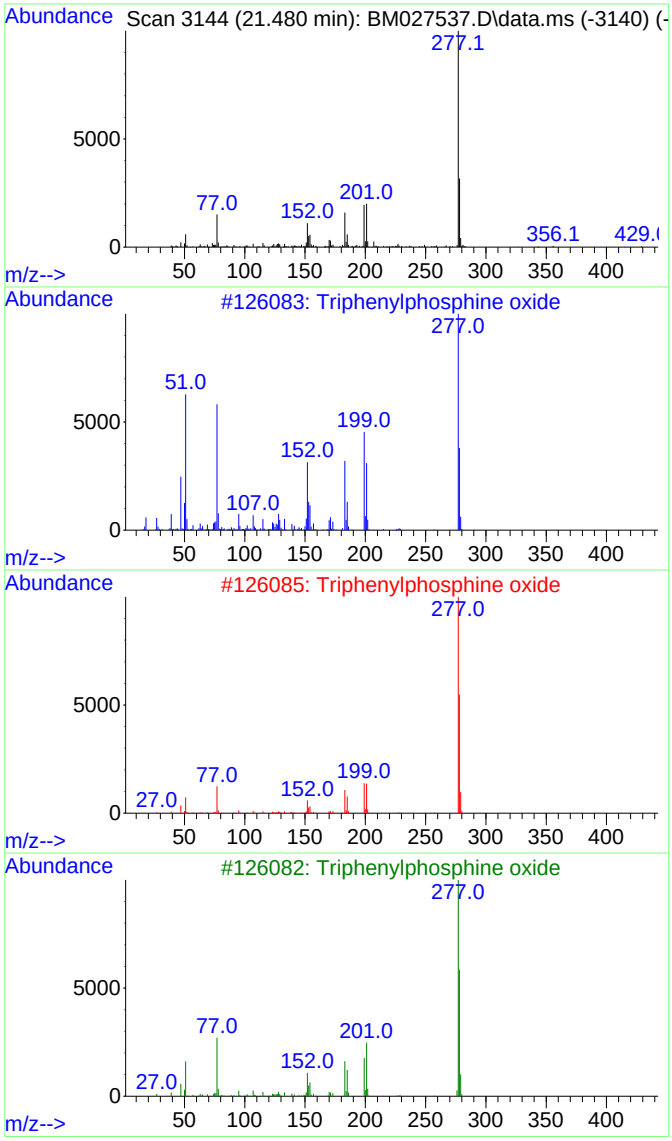
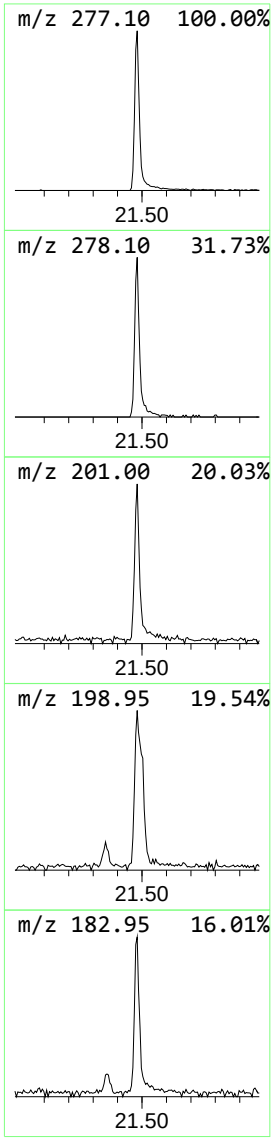
Peak Number 20 Triphenylphosphine oxide Concentration Rank 10

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 21.480 | 2.83 ng/ul | 252002 | Chrysene-d12 | 21.356 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Triphenylphosphine oxide | 278 | C18H15OP | 000791-28-6 | 91 |
| 2 | Triphenylphosphine oxide | 278 | C18H15OP | 000791-28-6 | 64 |
| 3 | Triphenylphosphine oxide | 278 | C18H15OP | 000791-28-6 | 46 |
| 4 | Triphenylphosphine oxide | 278 | C18H15OP | 000791-28-6 | 42 |
| 5 | Cyclotetrasilazane, 2,2,4,4,6,6,... | 292 | C8H28N4Si4 | 001020-84-4 | 28 |



Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092520\
Data File : BM027537.D
Acq On : 25 Sep 2020 20:05
Operator : JU/CG
Sample : L4135-06
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** BG480

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM092420MA.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 |
| unknown-01 | 3.063 | 5.7 | ng/ul | 197427 | 1 | 7.810 | 693851 | 20.0 |
| unknown-02 | 7.516 | 2.7 | ng/ul | 93466 | 1 | 7.810 | 693851 | 20.0 |
| unknown-03 | 8.010 | 2.5 | ng/ul | 87909 | 1 | 7.810 | 693851 | 20.0 |
| Indane | 8.181 | 2.1 | ng/ul | 72116 | 1 | 7.810 | 693851 | 20.0 |
| (DEL) Alkane: S... | 10.334 | 2.1 | ng/ul | 92109 | 2 | 10.592 | 880223 | 20.0 |
| Phenol, p-tert-... | 11.934 | 16.4 | ng/ul | 722166 | 2 | 10.592 | 880223 | 20.0 |
| Tripropyl phosp... | 12.969 | 2.1 | ng/ul | 155960 | 3 | 14.439 | 1512980 | 20.0 |
| 2(3H)-Benzothia... | 16.104 | 2.8 | ng/ul | 249139 | 4 | 17.180 | 1800590 | 20.0 |
| Phenol, 4-(1-ph... | 16.486 | 2.3 | ng/ul | 205693 | 4 | 17.180 | 1800590 | 20.0 |
| unknown-04 | 17.845 | 2.1 | ng/ul | 190592 | 4 | 17.180 | 1800590 | 20.0 |
| n-Hexadecanoic ... | 18.086 | 2.9 | ng/ul | 263857 | 4 | 17.180 | 1800590 | 20.0 |
| Diphenyl sulfone | 18.145 | 2.3 | ng/ul | 205039 | 4 | 17.180 | 1800590 | 20.0 |
| 2-[1-(4-Cyano-1... | 18.215 | 4.0 | ng/ul | 358681 | 4 | 17.180 | 1800590 | 20.0 |
| 1,2-Benzenediac... | 18.315 | 3.5 | ng/ul | 314769 | 4 | 17.180 | 1800590 | 20.0 |
| Isoquinoline | 18.368 | 7.5 | ng/ul | 673141 | 4 | 17.180 | 1800590 | 20.0 |
| Sulphenone | 19.180 | 2.2 | ng/ul | 194363 | 4 | 17.180 | 1800590 | 20.0 |
| n-Tetracosanol-1 | 21.080 | 3.0 | ng/ul | 270324 | 5 | 21.356 | 1780140 | 20.0 |
| Triphenylphosph... | 21.480 | 2.8 | ng/ul | 252002 | 5 | 21.356 | 1780140 | 20.0 |