LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN071620\
Data File : BN011190.D
Acq On : 16 Jul 2020 15:23
Operator : CG/JU
Sample : L3261-07
Misc :
ALS Vial : 6 Sample Multiplier: 1

**Instrument :** BNA_N
**ClientSampleId :** BE506

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_N\METHODS\SOM-EPA-BN071520MA.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 | 3.828 | 137 | 143 | 159 | rVB | 12924938 | 19072496 | 100.00% | 14.726% |
| 2 | 4.887 | 317 | 323 | 336 | rVB | 1830571 | 2761619 | 14.48% | 2.132% |
| 3 | 5.075 | 349 | 355 | 367 | rVV | 2334751 | 3298446 | 17.29% | 2.547% |
| 4 | 5.205 | 370 | 377 | 387 | rVV | 4958893 | 8801191 | 46.15% | 6.796% |
| 5 | 5.558 | 430 | 437 | 448 | rVB | 1768485 | 2579392 | 13.52% | 1.992% |
| 6 | 6.022 | 505 | 516 | 525 | rBV | 413420 | 683615 | 3.58% | 0.528% |
| 7 | 6.199 | 540 | 546 | 553 | rVB3 | 114319 | 190198 | 1.00% | 0.147% |
| 8 | 6.493 | 588 | 596 | 610 | rBV | 192802 | 358328 | 1.88% | 0.277% |
| 9 | 6.605 | 610 | 615 | 619 | rVV | 272255 | 404828 | 2.12% | 0.313% |
| 10 | 6.663 | 619 | 625 | 626 | rVV2 | 183715 | 313153 | 1.64% | 0.242% |
| 11 | 6.681 | 626 | 628 | 632 | rVV | 226552 | 400740 | 2.10% | 0.309% |
| 12 | 6.734 | 634 | 637 | 642 | rVV | 231941 | 350433 | 1.84% | 0.271% |
| 13 | 6.840 | 650 | 655 | 660 | rVV | 594099 | 913704 | 4.79% | 0.705% |
| 14 | 6.893 | 660 | 664 | 668 | rVV | 253419 | 396729 | 2.08% | 0.306% |
| 15 | 6.934 | 668 | 671 | 674 | rVV | 132543 | 191559 | 1.00% | 0.148% |
| 16 | 6.969 | 674 | 677 | 682 | rVV2 | 109826 | 176146 | 0.92% | 0.136% |
| 17 | 7.034 | 682 | 688 | 696 | rVV2 | 730348 | 1263738 | 6.63% | 0.976% |
| 18 | 7.146 | 701 | 707 | 714 | rVV | 883916 | 1400583 | 7.34% | 1.081% |
| 19 | 7.216 | 714 | 719 | 723 | rVV | 125134 | 196833 | 1.03% | 0.152% |
| 20 | 7.487 | 758 | 765 | 769 | rBV | 618656 | 1003793 | 5.26% | 0.775% |
| 21 | 7.522 | 769 | 771 | 779 | rVV2 | 157814 | 209599 | 1.10% | 0.162% |
| 22 | 7.605 | 779 | 785 | 796 | rVB | 296344 | 502561 | 2.64% | 0.388% |
| 23 | 7.699 | 796 | 801 | 806 | rBV | 152284 | 224343 | 1.18% | 0.173% |
| 24 | 7.852 | 814 | 827 | 832 | rBV2 | 417197 | 922699 | 4.84% | 0.712% |
| 25 | 7.910 | 832 | 837 | 845 | rVV2 | 577074 | 1136687 | 5.96% | 0.878% |
| 26 | 8.110 | 864 | 871 | 879 | rBV | 762877 | 1365434 | 7.16% | 1.054% |
| 27 | 8.193 | 879 | 885 | 891 | rVV | 504379 | 823629 | 4.32% | 0.636% |
| 28 | 8.252 | 891 | 895 | 912 | rVB2 | 251766 | 547051 | 2.87% | 0.422% |
| 29 | 8.581 | 942 | 951 | 954 | rBV | 147723 | 246489 | 1.29% | 0.190% |
| 30 | 8.622 | 954 | 958 | 964 | rVB | 646180 | 979678 | 5.14% | 0.756% |
| 31 | 8.704 | 969 | 972 | 980 | rVV5 | 113578 | 216015 | 1.13% | 0.167% |
| 32 | 8.799 | 980 | 988 | 999 | rVB | 278614 | 479550 | 2.51% | 0.370% |
| 33 | 9.269 | 1059 | 1068 | 1073 | rBV3 | 91399 | 230695 | 1.21% | 0.178% |
| 34 | 9.334 | 1073 | 1079 | 1086 | rVV | 651907 | 1096706 | 5.75% | 0.847% |

LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN071620\
Data File : BN011190.D
Acq On : 16 Jul 2020 15:23
Operator : CG/JU
Sample : L3261-07
Misc :
ALS Vial : 6 Sample Multiplier: 1

**Instrument :** BNA_N
**ClientSampleId :** BE506

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_N\METHODS\SOM-EPA-BN071520MA.M
Title : SVOA CALIBRATION

| | | | | | | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 35 | 9.663 | 1129 | 1135 | 1144 | rVV7 | 223412 | 486626 | 2.55% | 0.376% |
| 36 | 9.869 | 1160 | 1170 | 1180 | rBV | 1123747 | 1904722 | 9.99% | 1.471% |
| 37 | 10.234 | 1226 | 1232 | 1237 | rBV | 868630 | 1425009 | 7.47% | 1.100% |
| 38 | 10.287 | 1237 | 1241 | 1251 | rVB | 620159 | 987341 | 5.18% | 0.762% |
| 39 | 10.387 | 1253 | 1258 | 1267 | rBV5 | 114266 | 273411 | 1.43% | 0.211% |
| 40 | 10.651 | 1289 | 1303 | 1313 | rBV4 | 248053 | 631752 | 3.31% | 0.488% |
| 41 | 10.840 | 1329 | 1335 | 1343 | rVB2 | 426549 | 722907 | 3.79% | 0.558% |
| 42 | 11.246 | 1397 | 1404 | 1412 | rBV9 | 66996 | 228846 | 1.20% | 0.177% |
| 43 | 11.328 | 1412 | 1418 | 1424 | rVV | 253979 | 511414 | 2.68% | 0.395% |
| 44 | 11.598 | 1455 | 1464 | 1471 | rVV | 644229 | 1248911 | 6.55% | 0.964% |
| 45 | 11.904 | 1508 | 1516 | 1521 | rVB | 146200 | 250007 | 1.31% | 0.193% |
| 46 | 12.081 | 1540 | 1546 | 1549 | rVV5 | 83159 | 184759 | 0.97% | 0.143% |
| 47 | 12.128 | 1549 | 1554 | 1560 | rVB | 222664 | 376934 | 1.98% | 0.291% |
| 48 | 12.598 | 1628 | 1634 | 1642 | rBV3 | 107082 | 186007 | 0.98% | 0.144% |
| 49 | 12.675 | 1642 | 1647 | 1654 | rBV3 | 181402 | 312851 | 1.64% | 0.242% |
| 50 | 12.769 | 1659 | 1663 | 1669 | rVB | 161183 | 226333 | 1.19% | 0.175% |
| 51 | 13.040 | 1703 | 1709 | 1719 | rBV | 870213 | 1411868 | 7.40% | 1.090% |
| 52 | 13.539 | 1787 | 1794 | 1799 | rBV | 2229089 | 3136787 | 16.45% | 2.422% |
| 53 | 13.587 | 1799 | 1802 | 1805 | rVV | 295364 | 446407 | 2.34% | 0.345% |
| 54 | 13.628 | 1805 | 1809 | 1815 | rVV | 831772 | 1308882 | 6.86% | 1.011% |
| 55 | 13.804 | 1830 | 1839 | 1849 | rVB | 2321782 | 3511198 | 18.41% | 2.711% |
| 56 | 13.904 | 1849 | 1856 | 1859 | rBV | 131393 | 188822 | 0.99% | 0.146% |
| 57 | 14.116 | 1886 | 1892 | 1898 | rBV2 | 1421267 | 2171553 | 11.39% | 1.677% |
| 58 | 14.175 | 1898 | 1902 | 1908 | rVB2 | 464601 | 714350 | 3.75% | 0.552% |
| 59 | 14.351 | 1923 | 1932 | 1941 | rBV2 | 197219 | 433848 | 2.27% | 0.335% |
| 60 | 15.122 | 2056 | 2063 | 2066 | rBV | 3258133 | 4773883 | 25.03% | 3.686% |
| 61 | 15.151 | 2066 | 2068 | 2077 | rVB | 1346906 | 1536859 | 8.06% | 1.187% |
| 62 | 15.251 | 2077 | 2085 | 2090 | rBV | 1136838 | 1621879 | 8.50% | 1.252% |
| 63 | 15.392 | 2105 | 2109 | 2114 | rVV2 | 490876 | 678779 | 3.56% | 0.524% |
| 64 | 15.645 | 2147 | 2152 | 2163 | rBV3 | 313651 | 700547 | 3.67% | 0.541% |
| 65 | 15.816 | 2176 | 2181 | 2186 | rVV2 | 354655 | 584216 | 3.06% | 0.451% |
| 66 | 15.875 | 2188 | 2191 | 2200 | rVB8 | 96066 | 205483 | 1.08% | 0.159% |
| 67 | 16.192 | 2241 | 2245 | 2255 | rVB | 464502 | 603658 | 3.17% | 0.466% |
| 68 | 16.869 | 2352 | 2360 | 2366 | rBV2 | 1833210 | 3583361 | 18.79% | 2.767% |
| 69 | 16.969 | 2372 | 2377 | 2384 | rBV2 | 3762001 | 5355450 | 28.08% | 4.135% |
| 70 | 17.263 | 2423 | 2427 | 2432 | rVB | 174866 | 235571 | 1.24% | 0.182% |
| 71 | 17.439 | 2452 | 2457 | 2462 | rVV | 250506 | 413145 | 2.17% | 0.319% |

LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN071620\
Data File : BN011190.D
Acq On : 16 Jul 2020 15:23
Operator : CG/JU
Sample : L3261-07
Misc :
ALS Vial : 6 Sample Multiplier: 1

**Instrument :** BNA_N
**ClientSampleId :** BE506

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_N\METHODS\SOM-EPA-BN071520MA.M
Title : SVOA CALIBRATION

| | | | | | | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 72 | 17.563 | 2475 | 2478 | 2484 | rVV2 | 170771 | 263851 | 1.38% | 0.204% |
| 73 | 17.804 | 2515 | 2519 | 2523 | rBV | 378817 | 506159 | 2.65% | 0.391% |
| 74 | 17.922 | 2534 | 2539 | 2547 | rBV | 421107 | 614193 | 3.22% | 0.474% |
| 75 | 18.022 | 2552 | 2556 | 2559 | rVV2 | 426944 | 593878 | 3.11% | 0.459% |
| 76 | 18.075 | 2559 | 2565 | 2573 | rVB | 641732 | 1144390 | 6.00% | 0.884% |
| 77 | 18.369 | 2607 | 2615 | 2620 | rBV6 | 84206 | 198223 | 1.04% | 0.153% |
| 78 | 18.622 | 2654 | 2658 | 2663 | rBV | 139471 | 189879 | 1.00% | 0.147% |
| 79 | 18.792 | 2684 | 2687 | 2691 | rVB3 | 190339 | 248543 | 1.30% | 0.192% |
| 80 | 19.010 | 2719 | 2724 | 2733 | rBV5 | 187301 | 470215 | 2.47% | 0.363% |
| 81 | 19.104 | 2736 | 2740 | 2745 | rBV2 | 195226 | 287954 | 1.51% | 0.222% |
| 82 | 19.280 | 2764 | 2770 | 2776 | rVB | 4947627 | 6536664 | 34.27% | 5.047% |
| 83 | 19.351 | 2776 | 2782 | 2793 | rBV | 1110272 | 1690527 | 8.86% | 1.305% |
| 84 | 19.539 | 2812 | 2814 | 2819 | rVB2 | 179414 | 224395 | 1.18% | 0.173% |
| 85 | 19.592 | 2819 | 2823 | 2828 | rBV | 165993 | 193509 | 1.01% | 0.149% |
| 86 | 19.710 | 2836 | 2843 | 2847 | rBV3 | 154784 | 282741 | 1.48% | 0.218% |
| 87 | 19.798 | 2855 | 2858 | 2861 | rBV3 | 172532 | 238524 | 1.25% | 0.184% |
| 88 | 20.033 | 2894 | 2898 | 2904 | rVB | 362034 | 450560 | 2.36% | 0.348% |
| 89 | 20.357 | 2950 | 2953 | 2957 | rBV2 | 179962 | 224069 | 1.17% | 0.173% |
| 90 | 20.457 | 2966 | 2970 | 2975 | rBV3 | 160297 | 319631 | 1.68% | 0.247% |
| 91 | 20.851 | 3033 | 3037 | 3049 | rBV5 | 274767 | 539983 | 2.83% | 0.417% |
| 92 | 21.016 | 3062 | 3065 | 3070 | rVB2 | 212902 | 227235 | 1.19% | 0.175% |
| 93 | 21.086 | 3072 | 3077 | 3083 | rBV | 2686593 | 3570627 | 18.72% | 2.757% |
| 94 | 21.221 | 3096 | 3100 | 3107 | rBV | 1046075 | 1353486 | 7.10% | 1.045% |
| 95 | 22.151 | 3255 | 3258 | 3261 | rVB2 | 165023 | 194609 | 1.02% | 0.150% |
| 96 | 22.627 | 3336 | 3339 | 3348 | rVB2 | 225316 | 318699 | 1.67% | 0.246% |
| 97 | 23.092 | 3411 | 3418 | 3426 | rBV | 4163605 | 7105926 | 37.26% | 5.487% |
| 98 | 23.163 | 3426 | 3430 | 3434 | rVV2 | 287544 | 432150 | 2.27% | 0.334% |
| 99 | 23.221 | 3434 | 3440 | 3450 | rVB2 | 1917903 | 3470287 | 18.20% | 2.680% |
| 100 | 23.762 | 3528 | 3532 | 3537 | rVB2 | 200712 | 307847 | 1.61% | 0.238% |

Sum of corrected areas: 129511790

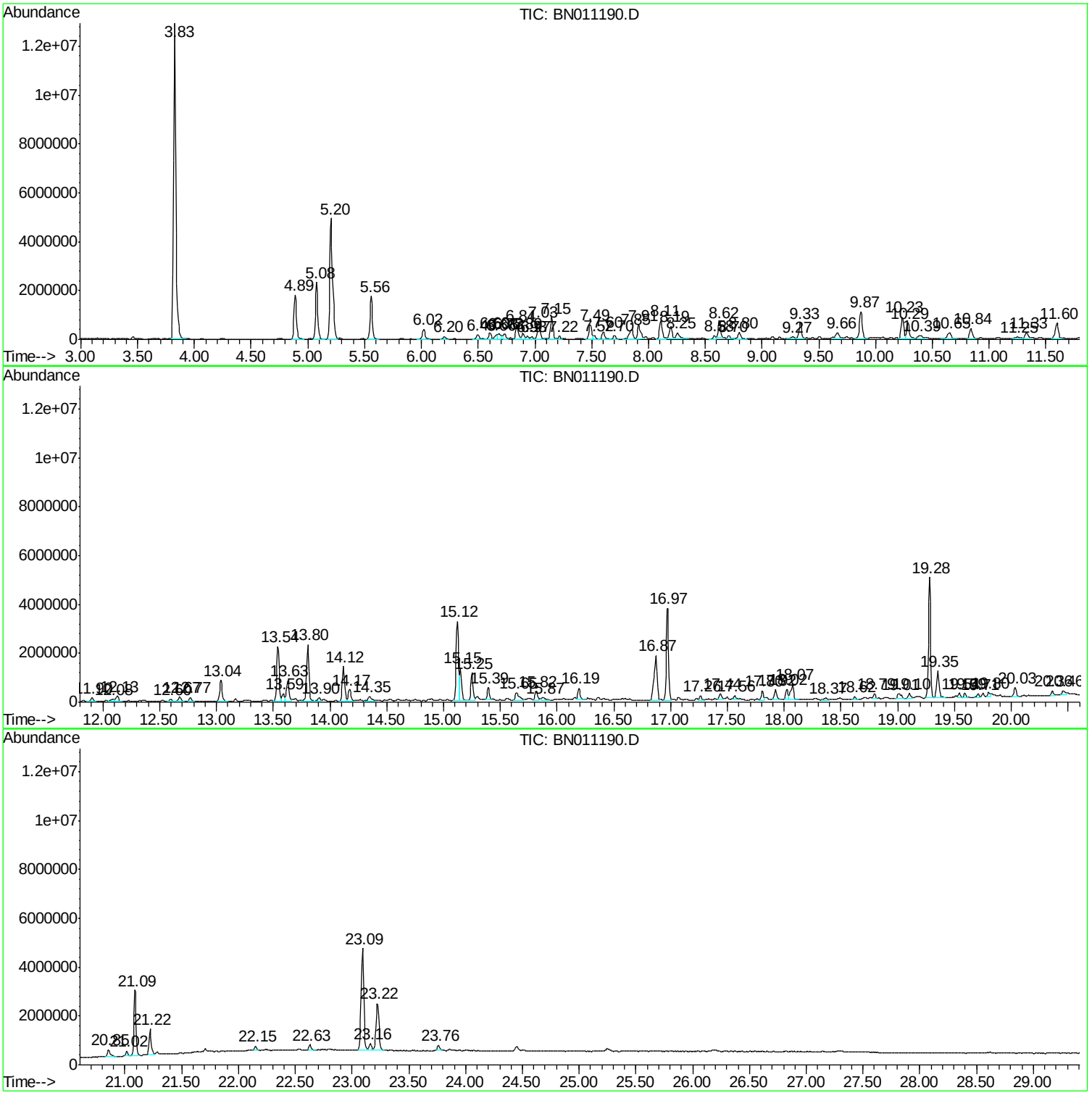
LSC Report - Integrated Chromatogram

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN071620\
Data File : BN011190.D
Acq On : 16 Jul 2020 15:23
Operator : CG/JU
Sample : L3261-07
Misc :
ALS Vial : 6 Sample Multiplier: 1

**Instrument :** BNA_N
**ClientSampleId :** BE506

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN071520MA.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 N\DATA\BN071620\
Data File : BN011190.D
Acq On : 16 Jul 2020 15:23
Operator : CG/JU
Sample : L3261-07
Misc :
ALS Vial : 6 Sample Multiplier: 1

**Instrument :** BNA_N
**ClientSampleId :** BE506

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN071520MA.M
Quant Title : SVOA CALIBRATION

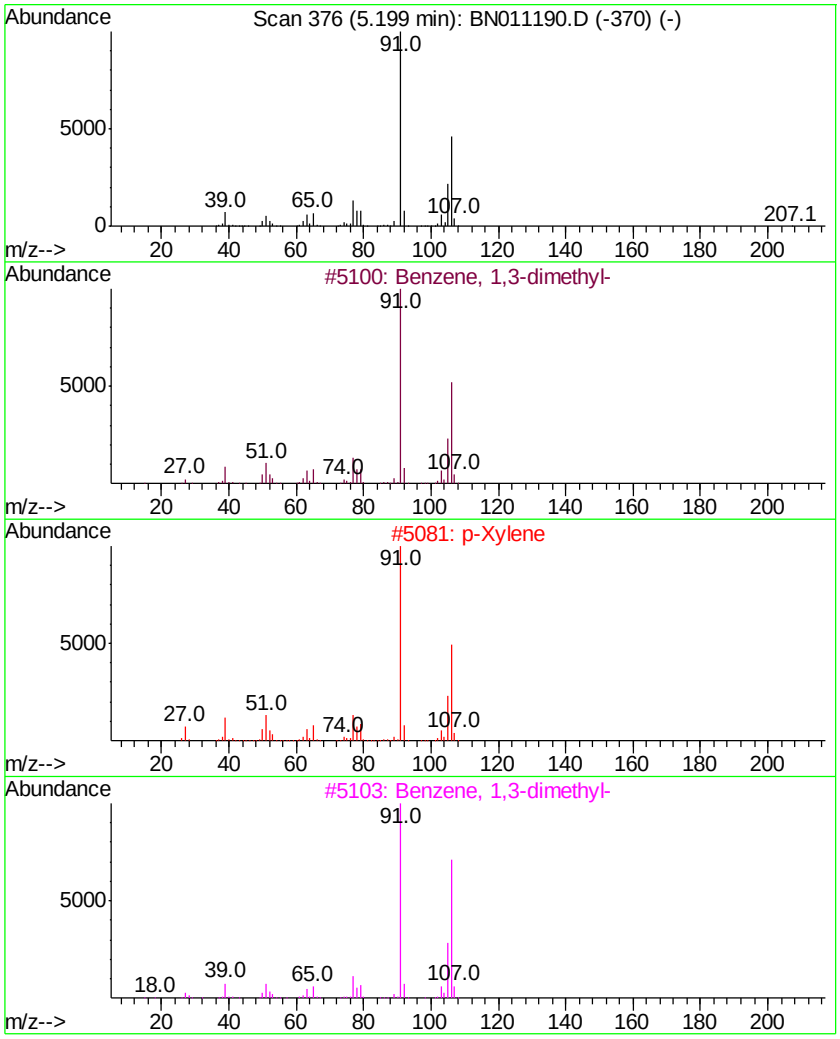
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

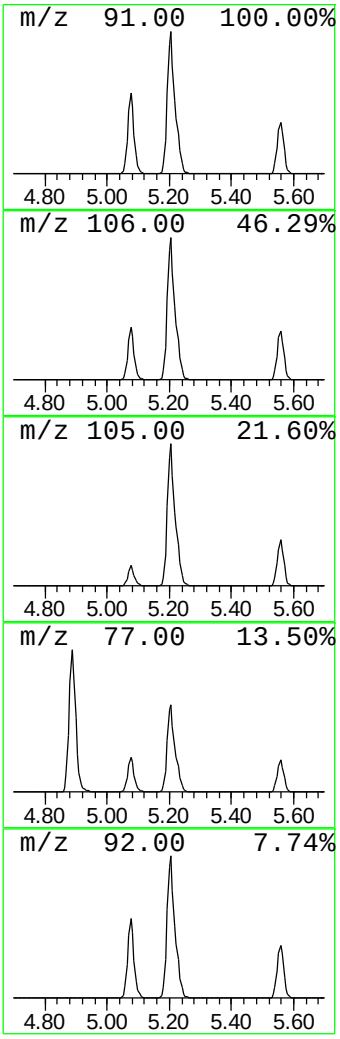
**********

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

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 5.20 | 175.36 ng/ul | 8801190 | 1,4-Dichlorobenzene-d4 | 7.49 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzene, 1,3-dimethyl- | 106 | C8H10 | 000108-38-3 | 95 |
| 2 | p-Xylene | 106 | C8H10 | 000106-42-3 | 95 |
| 3 | Benzene, 1,3-dimethyl- | 106 | C8H10 | 000108-38-3 | 95 |
| 4 | p-Xylene | 106 | C8H10 | 000106-42-3 | 95 |
| 5 | o-Xylene | 106 | C8H10 | 000095-47-6 | 95 |



# Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN071620\
Data File : BN011190.D
Acq On : 16 Jul 2020 15:23
Operator : CG/JU
Sample : L3261-07
Misc :
ALS Vial : 6 Sample Multiplier: 1

**Instrument :** BNA_N
**ClientSampleId :** BE506

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN071520MA.M
Quant Title : SVOA CALIBRATION

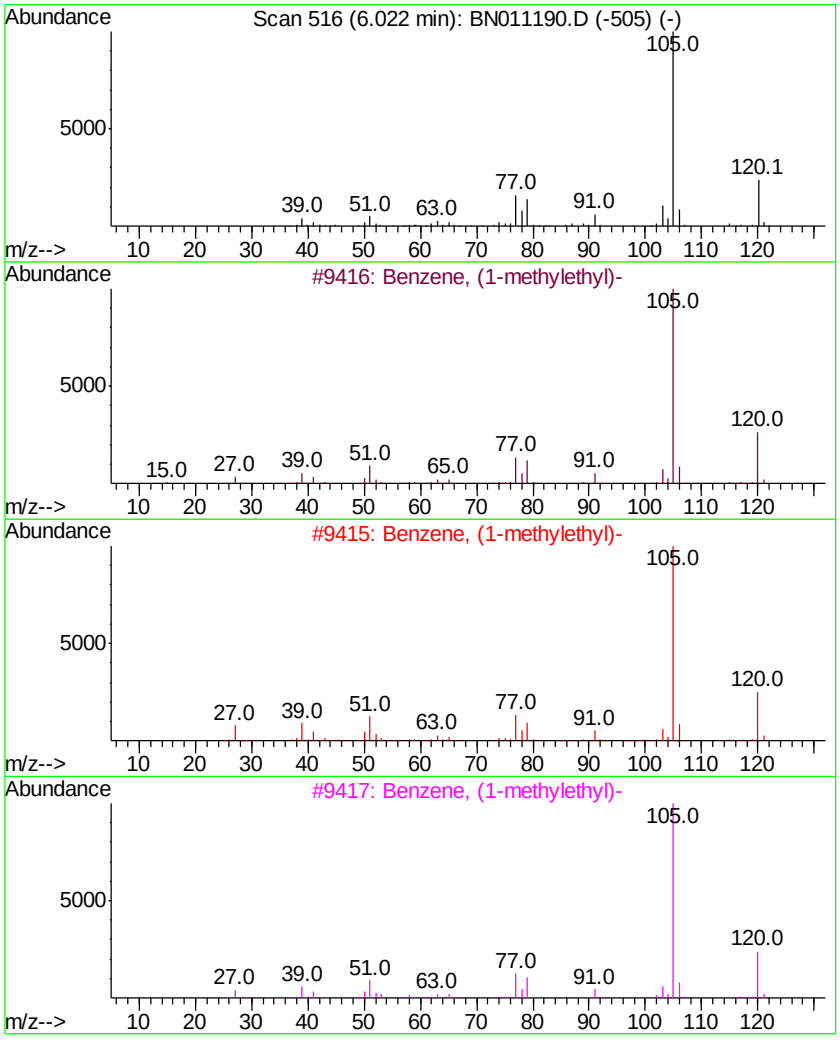
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

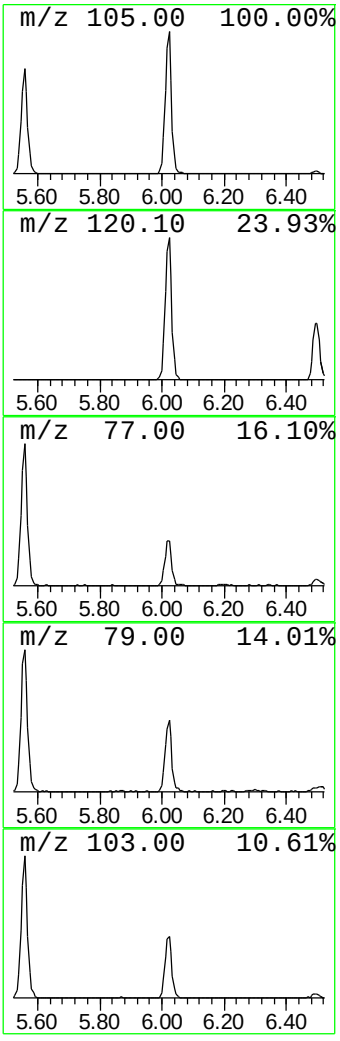
**********

Peak Number 6 Benzene, (1-methylethyl)- Concentration Rank 10

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 6.02 | 13.62 ng/ul | 683615 | 1,4-Dichlorobenzene-d4 | 7.49 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzene, (1-methylethyl)- | 120 | C9H12 | 000098-82-8 | 91 |
| 2 | Benzene, (1-methylethyl)- | 120 | C9H12 | 000098-82-8 | 91 |
| 3 | Benzene, (1-methylethyl)- | 120 | C9H12 | 000098-82-8 | 91 |
| 4 | Benzene, (1-methylethyl)- | 120 | C9H12 | 000098-82-8 | 87 |
| 5 | Benzene, 1-ethyl-3-methyl- | 120 | C9H12 | 000620-14-4 | 80 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN071620\
Data File : BN011190.D
Acq On : 16 Jul 2020 15:23
Operator : CG/JU
Sample : L3261-07
Misc :
ALS Vial : 6 Sample Multiplier: 1

**Instrument :** BNA_N
**ClientSampleId :** BE506

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN071520MA.M
Quant Title : SVOA CALIBRATION

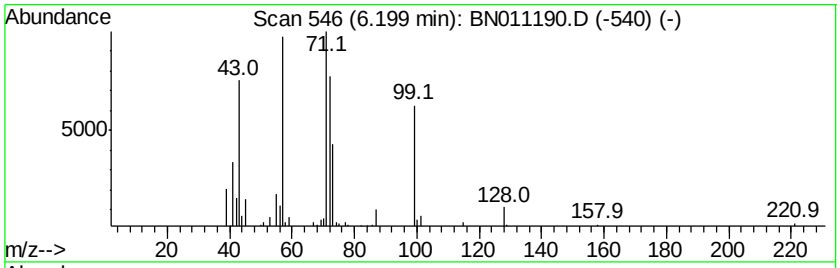
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

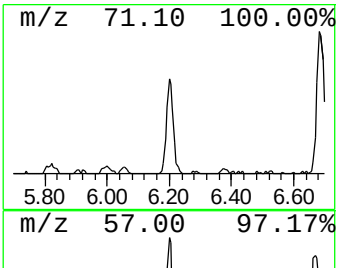
**********

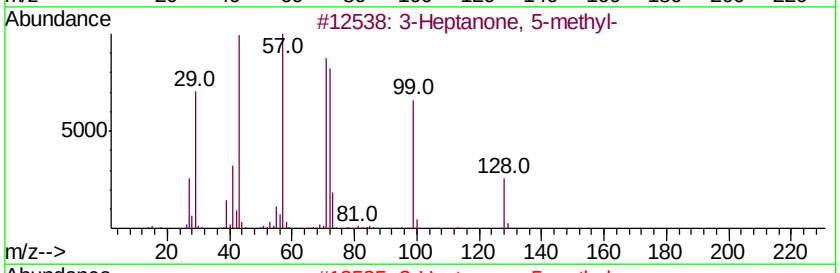
Peak Number 7 3-Heptanone, 5-methyl- Concentration Rank 28

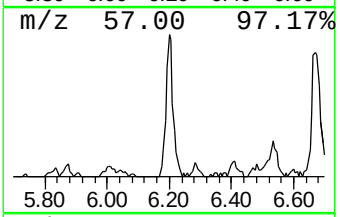
| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 6.20 | 3.79 ng/ul | 190198 | 1,4-Dichlorobenzene-d4 | 7.49 |

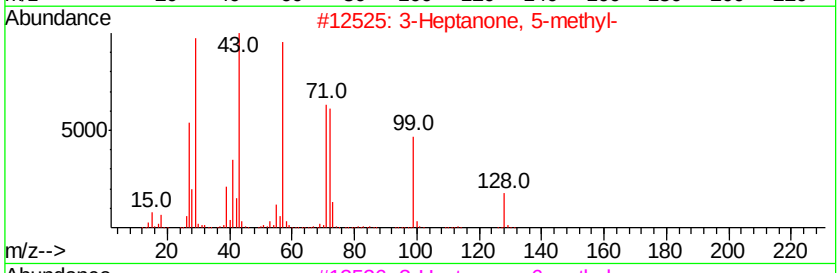
| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 3-Heptanone, 5-methyl- | 128 | C8H16O | 000541-85-5 | 87 |
| 2 | 3-Heptanone, 5-methyl- | 128 | C8H16O | 000541-85-5 | 87 |
| 3 | 3-Heptanone, 6-methyl- | 128 | C8H16O | 000624-42-0 | 86 |
| 4 | 3-Octanone | 128 | C8H16O | 000106-68-3 | 80 |
| 5 | 3-Octanone | 128 | C8H16O | 000106-68-3 | 64 |

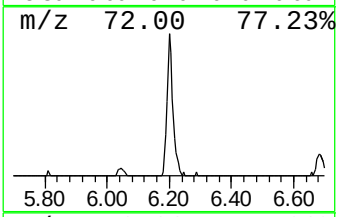


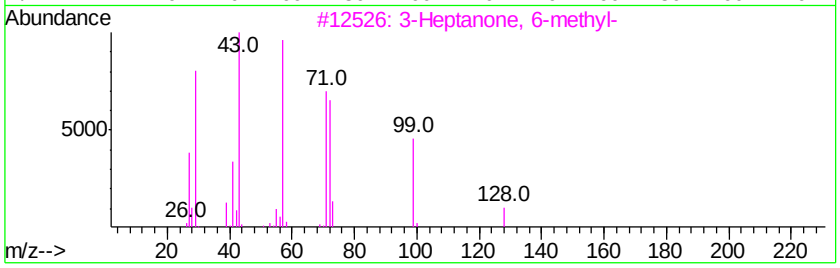



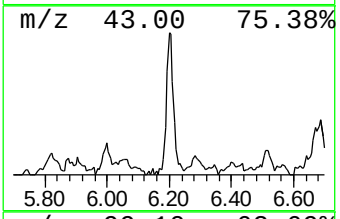



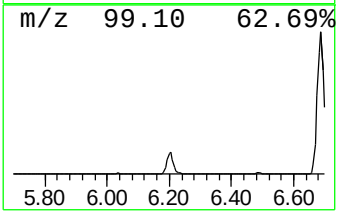



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Data File : BN011190.D
Acq On : 16 Jul 2020 15:23
Operator : CG/JU
Sample : L3261-07
Misc :
ALS Vial : 6 Sample Multiplier: 1

**Instrument :** BNA_N
**ClientSampleId :** BE506

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN071520MA.M
Quant Title : SVOA CALIBRATION

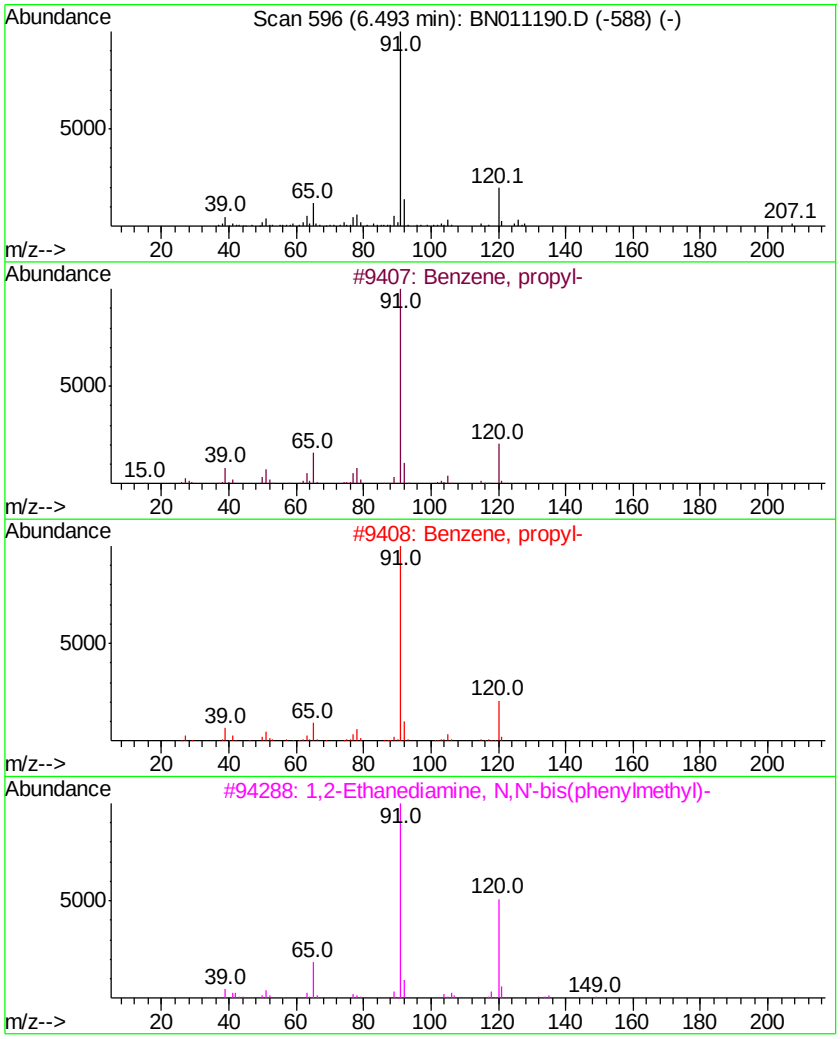
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

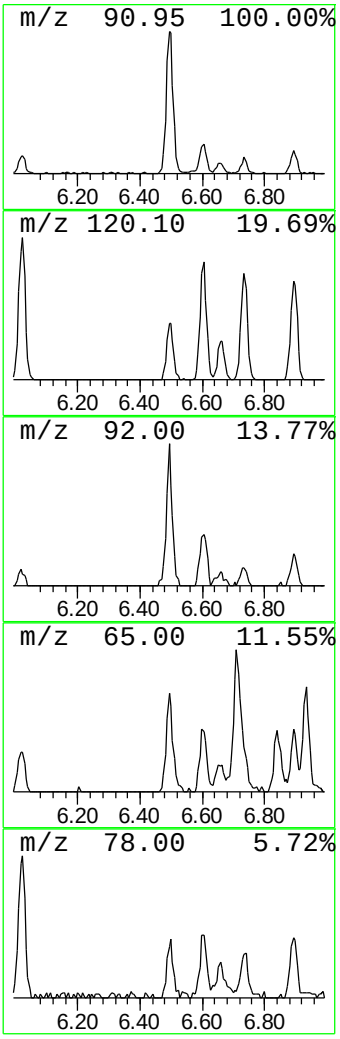
**********

Peak Number 8 Benzene, propyl- Concentration Rank 20

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 6.49 | 7.14 ng/ul | 358328 | 1,4-Dichlorobenzene-d4 | 7.49 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzene, propyl- | 120 | C9H12 | 000103-65-1 | 87 |
| 2 | Benzene, propyl- | 120 | C9H12 | 000103-65-1 | 81 |
| 3 | 1,2-Ethanediamine, N,N'-bis(phen... | 240 | C16H20N2 | 000140-28-3 | 72 |
| 4 | Ethanol, 2-[(phenylmethyl)amino]- | 151 | C9H13NO | 000104-63-2 | 72 |
| 5 | Benzene, propyl- | 120 | C9H12 | 000103-65-1 | 64 |



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Data File : BN011190.D
Acq On : 16 Jul 2020 15:23
Operator : CG/JU
Sample : L3261-07
Misc :
ALS Vial : 6 Sample Multiplier: 1

**Instrument :** BNA_N
**ClientSampleId :** BE506

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN071520MA.M
Quant Title : SVOA CALIBRATION

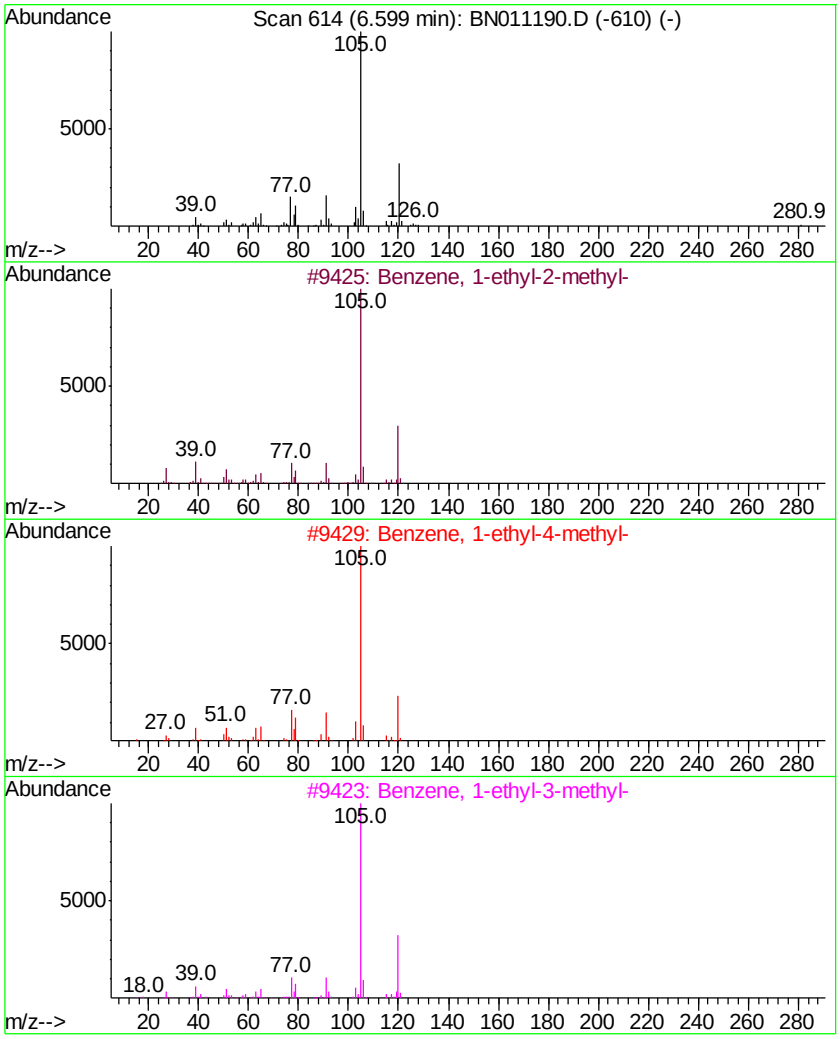
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

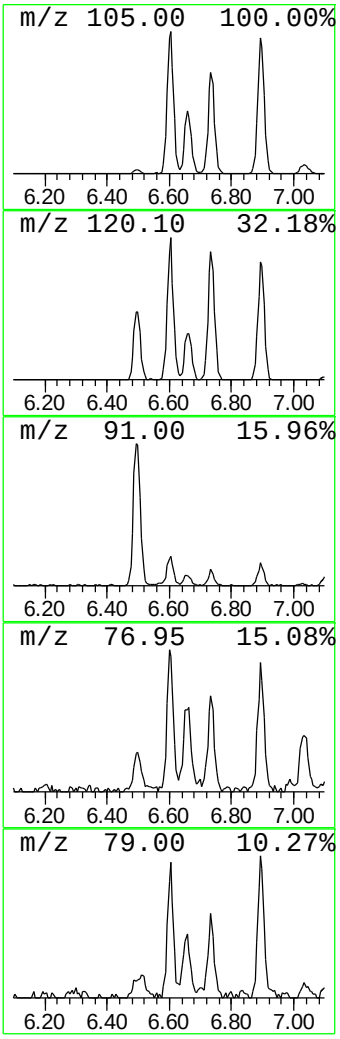
**********

Peak Number 9 Benzene, 1-ethyl-2-methyl- Concentration Rank 17

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 6.60 | 8.07 ng/ul | 404828 | 1,4-Dichlorobenzene-d4 | 7.49 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzene, 1-ethyl-2-methyl- | 120 | C9H12 | 000611-14-3 | 94 |
| 2 | Benzene, 1-ethyl-4-methyl- | 120 | C9H12 | 000622-96-8 | 93 |
| 3 | Benzene, 1-ethyl-3-methyl- | 120 | C9H12 | 000620-14-4 | 91 |
| 4 | Mesitylene | 120 | C9H12 | 000108-67-8 | 91 |
| 5 | Benzene, 1-ethyl-3-methyl- | 120 | C9H12 | 000620-14-4 | 91 |



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Data File : BN011190.D
Acq On : 16 Jul 2020 15:23
Operator : CG/JU
Sample : L3261-07
Misc :
ALS Vial : 6 Sample Multiplier: 1

**Instrument :** BNA_N
**ClientSampleId :** BE506

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN071520MA.M
Quant Title : SVOA CALIBRATION

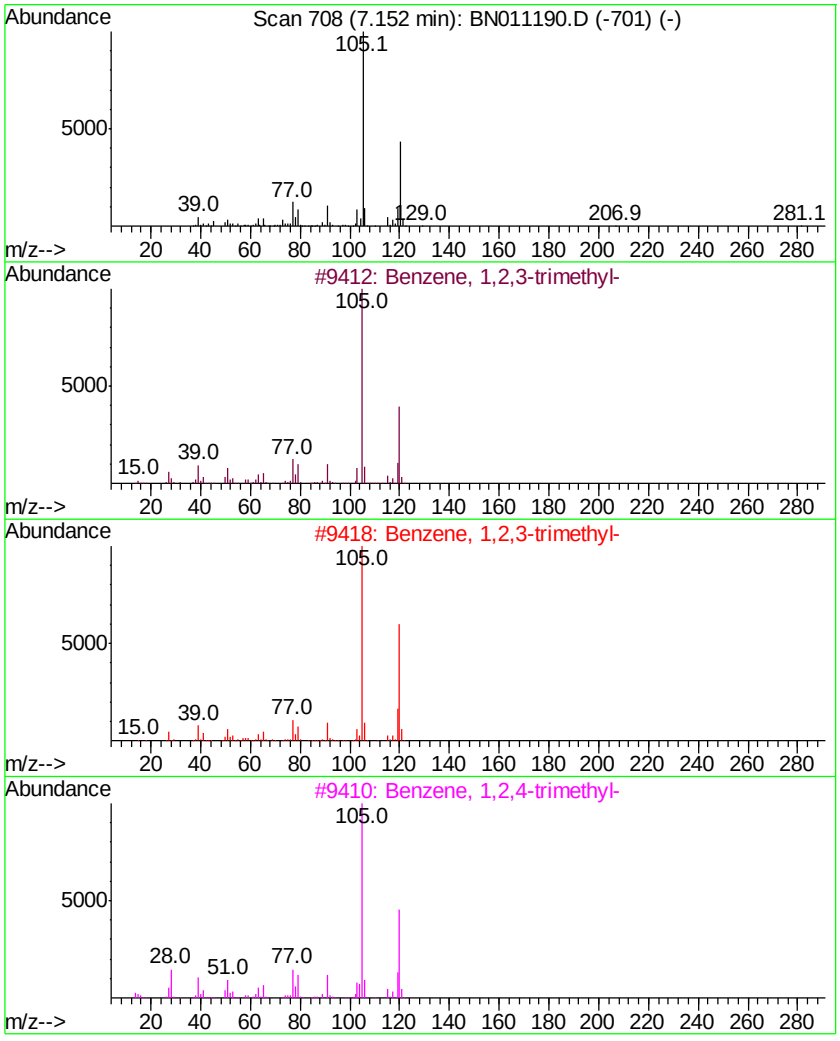
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

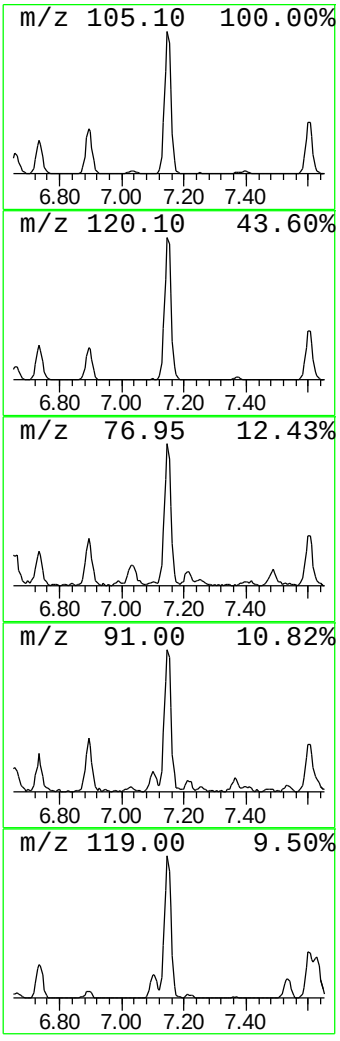
**********

Peak Number 10 Benzene, 1,2,3-trimethyl- Concentration Rank 6

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 7.15 | 27.91 ng/ul | 1400580 | 1,4-Dichlorobenzene-d4 | 7.49 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzene, 1,2,3-trimethyl- | 120 | C9H12 | 000526-73-8 | 97 |
| 2 | Benzene, 1,2,3-trimethyl- | 120 | C9H12 | 000526-73-8 | 95 |
| 3 | Benzene, 1,2,4-trimethyl- | 120 | C9H12 | 000095-63-6 | 94 |
| 4 | Mesitylene | 120 | C9H12 | 000108-67-8 | 93 |
| 5 | Benzene, 1-ethyl-4-methyl- | 120 | C9H12 | 000622-96-8 | 91 |



# Library Search Compound Report

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Data File : BN011190.D
Acq On : 16 Jul 2020 15:23
Operator : CG/JU
Sample : L3261-07
Misc :
ALS Vial : 6 Sample Multiplier: 1

**Instrument :** BNA_N
**ClientSampleId :** BE506

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN071520MA.M
Quant Title : SVOA CALIBRATION

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

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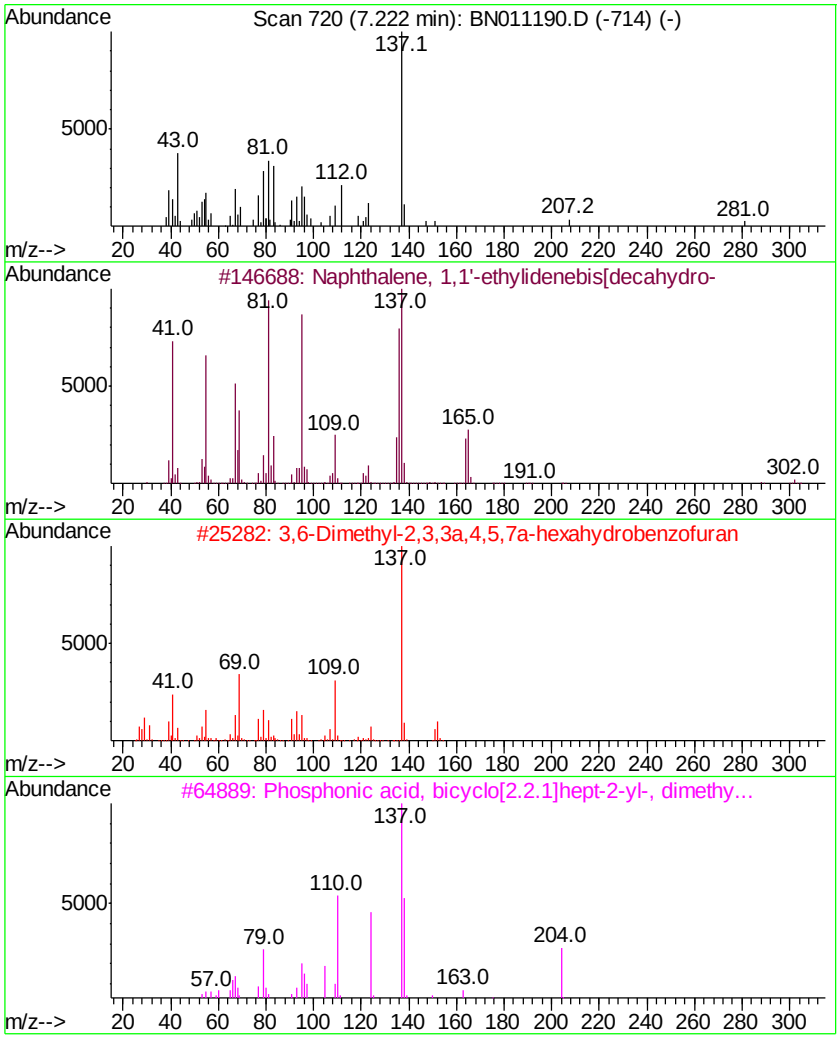
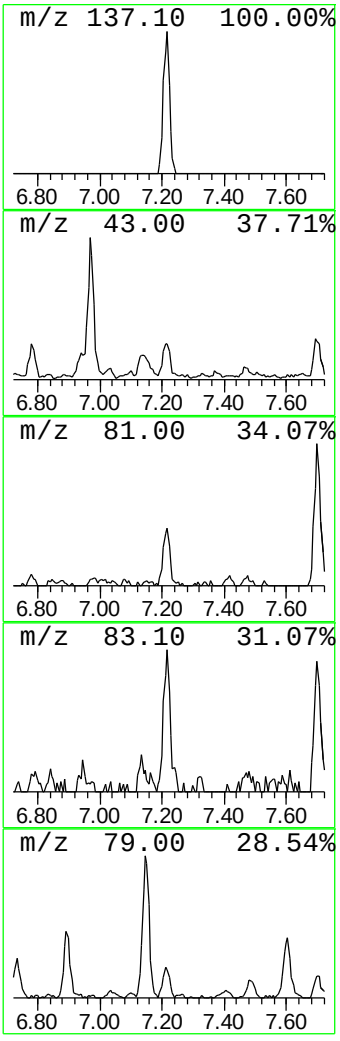
Peak Number 11 Naphthalene, 1,1'-ethyliden... Concentration Rank 26

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 7.22 | 3.92 ng/ul | 196833 | 1,4-Dichlorobenzene-d4 | 7.49 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Naphthalene, 1,1'-ethylidenebis[... | 302 | C22H38 | 054934-70-2 | 50 |
| 2 | 3,6-Dimethyl-2,3,3a,4,5,7a-hexah... | 152 | C10H16O | 070786-44-6 | 40 |
| 3 | Phosphonic acid, bicyclo[2.2.1]h... | 204 | C9H17O3P | 050457-69-7 | 40 |
| 4 | 3,5-Octadiene, 2,2,4,5,7,7-hexam... | 194 | C14H26 | 055712-52-2 | 38 |
| 5 | 3,6-Dimethyl-2,3,3a,4,5,7a-hexah... | 152 | C10H16O | 070786-44-6 | 38 |



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Data File : BN011190.D
Acq On : 16 Jul 2020 15:23
Operator : CG/JU
Sample : L3261-07
Misc :
ALS Vial : 6 Sample Multiplier: 1

**Instrument :** BNA_N
**ClientSampleId :** BE506

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN071520MA.M
Quant Title : SVOA CALIBRATION

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

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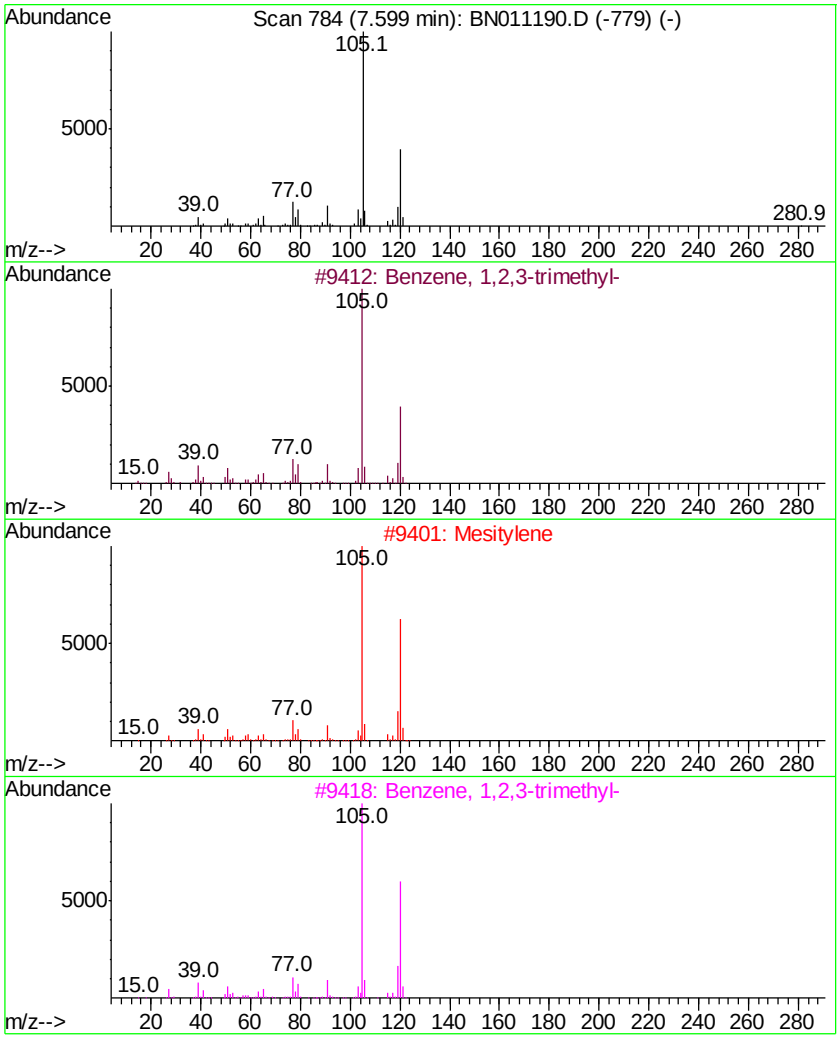
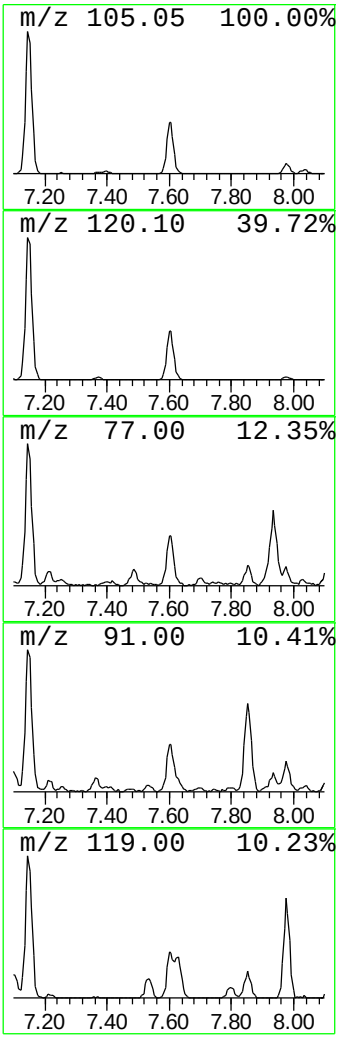
Peak Number 12 Mesitylene Concentration Rank 13

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 7.60 | 10.01 ng/ul | 502561 | 1,4-Dichlorobenzene-d4 | 7.49 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzene, 1,2,3-trimethyl- | 120 | C9H12 | 000526-73-8 | 97 |
| 2 | Mesitylene | 120 | C9H12 | 000108-67-8 | 94 |
| 3 | Benzene, 1,2,3-trimethyl- | 120 | C9H12 | 000526-73-8 | 94 |
| 4 | Benzene, 1,2,4-trimethyl- | 120 | C9H12 | 000095-63-6 | 94 |
| 5 | Benzene, 1-ethyl-2-methyl- | 120 | C9H12 | 000611-14-3 | 93 |



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Data File : BN011190.D
Acq On : 16 Jul 2020 15:23
Operator : CG/JU
Sample : L3261-07
Misc :
ALS Vial : 6 Sample Multiplier: 1

**Instrument :** BNA_N
**ClientSampleId :** BE506

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN071520MA.M
Quant Title : SVOA CALIBRATION

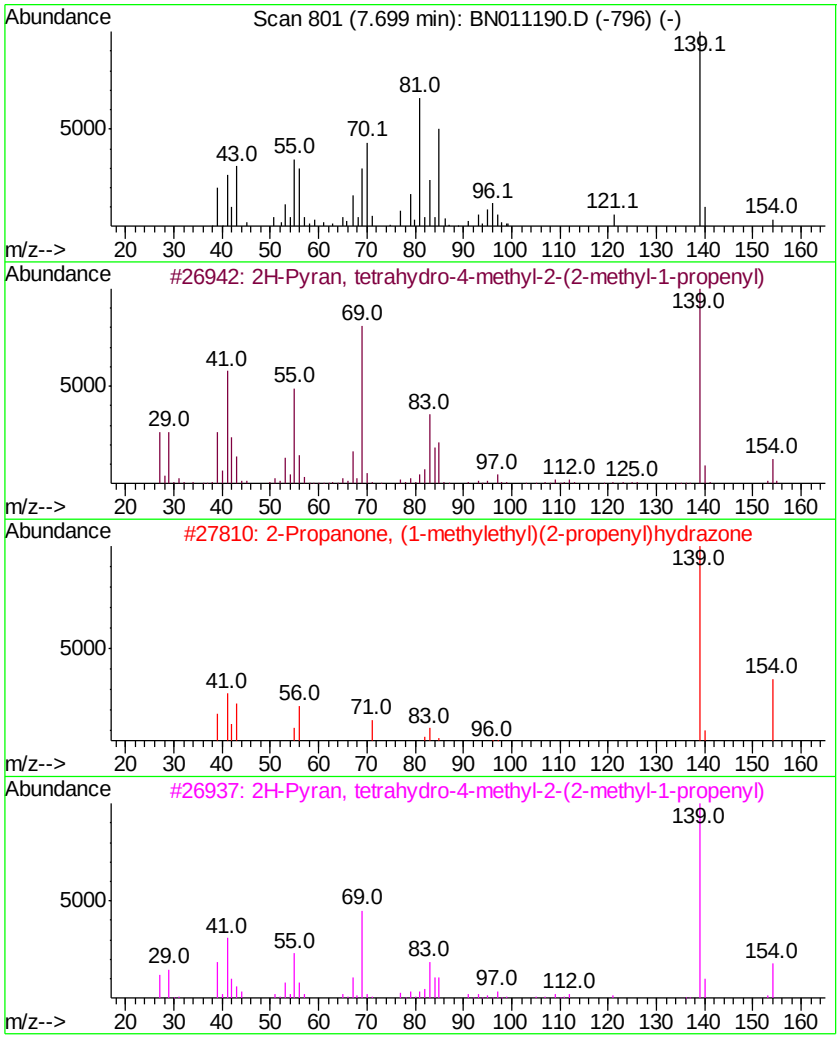
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

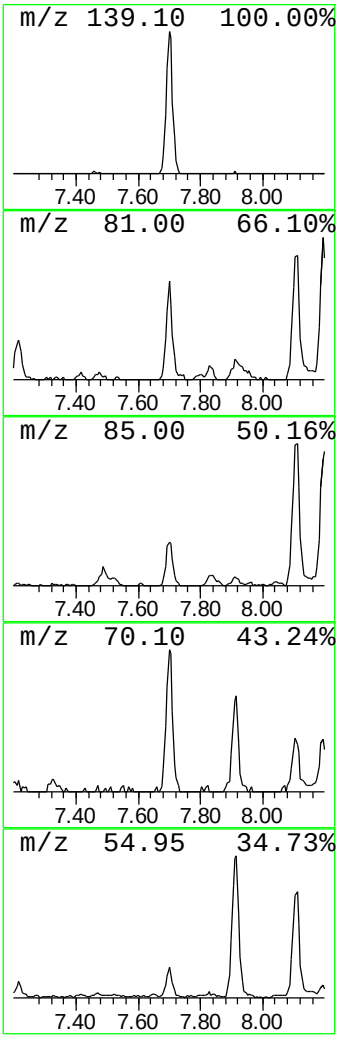
**********

Peak Number 13 unknown-01 Concentration Rank 24

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 7.70 | 4.47 ng/ul | 224343 | 1,4-Dichlorobenzene-d4 | 7.49 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2H-Pyran, tetrahydro-4-methyl-2-... | 154 | C10H18O | 016409-43-1 | 37 |
| 2 | 2-Propanone, (1-methylethyl)(2-p... | 154 | C9H18N2 | 110213-09-7 | 37 |
| 3 | 2H-Pyran, tetrahydro-4-methyl-2-... | 154 | C10H18O | 016409-43-1 | 37 |
| 4 | 2-Amino-4-hydroxy-3,6-dimethylpy... | 139 | C6H9N3O | 033796-41-7 | 35 |
| 5 | Glutaric acid, di(1-(4-fluorophe... | 376 | C21H22F2O4 | 1000377-11-0 | 35 |



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Data File : BN011190.D
Acq On : 16 Jul 2020 15:23
Operator : CG/JU
Sample : L3261-07
Misc :
ALS Vial : 6 Sample Multiplier: 1

**Instrument :** BNA_N
**ClientSampleId :** BE506

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN071520MA.M
Quant Title : SVOA CALIBRATION

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

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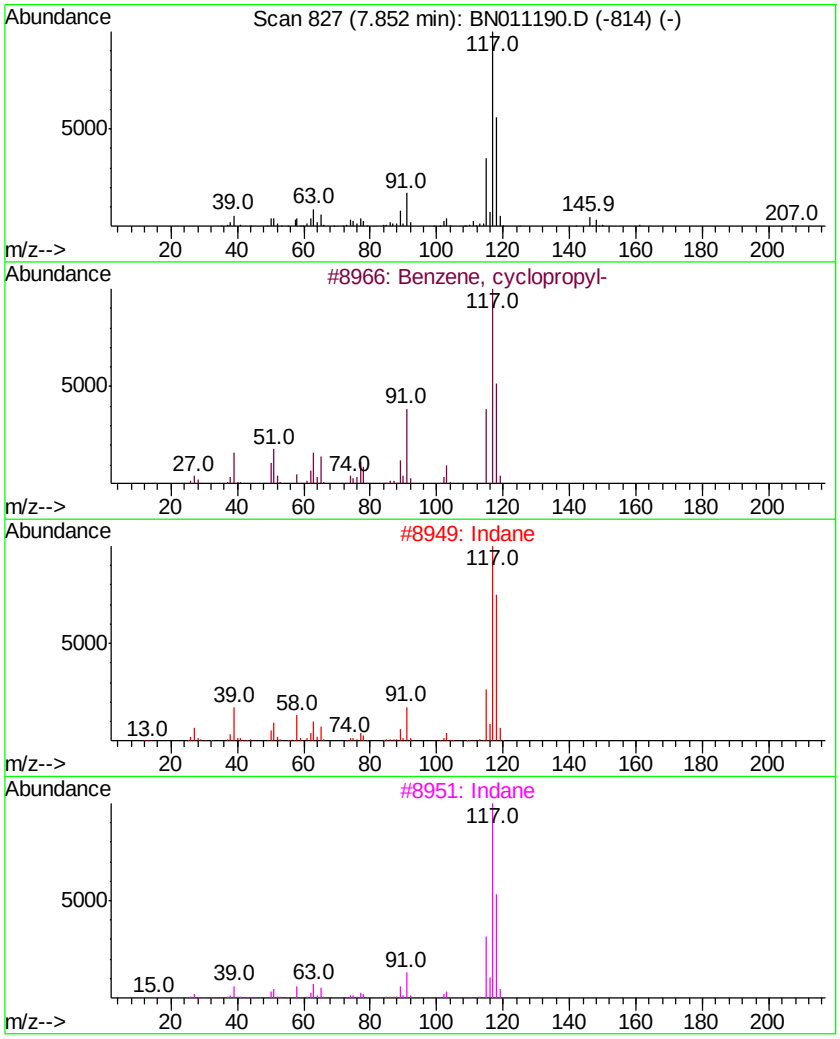
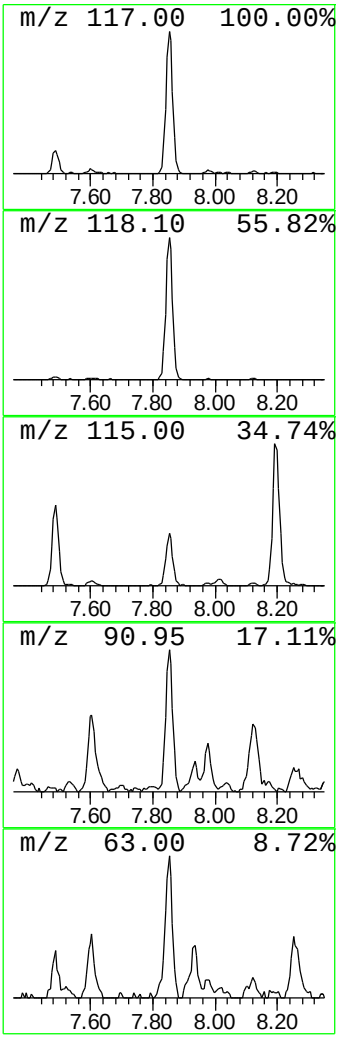
Peak Number 14 Benzene, cyclopropyl- Concentration Rank 8

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 7.85 | 18.38 ng/ul | 922699 | 1,4-Dichlorobenzene-d4 | 7.49 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzene, cyclopropyl- | 118 | C9H10 | 000873-49-4 | 94 |
| 2 | Indane | 118 | C9H10 | 000496-11-7 | 90 |
| 3 | Indane | 118 | C9H10 | 000496-11-7 | 76 |
| 4 | Benzene, cyclopropyl- | 118 | C9H10 | 000873-49-4 | 74 |
| 5 | (Z)-1-Phenylpropene | 118 | C9H10 | 000766-90-5 | 68 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN071620\
Data File : BN011190.D
Acq On : 16 Jul 2020 15:23
Operator : CG/JU
Sample : L3261-07
Misc :
ALS Vial : 6 Sample Multiplier: 1

**Instrument :** BNA_N
**ClientSampleId :** BE506

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN071520MA.M
Quant Title : SVOA CALIBRATION

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

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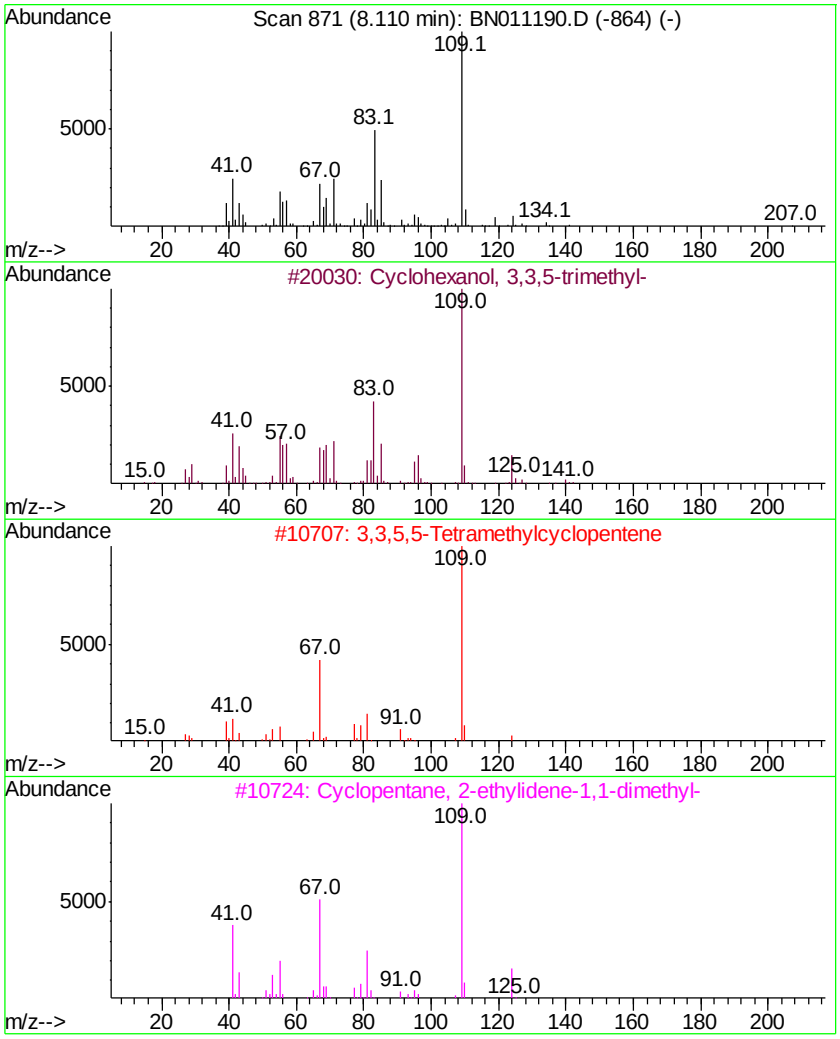
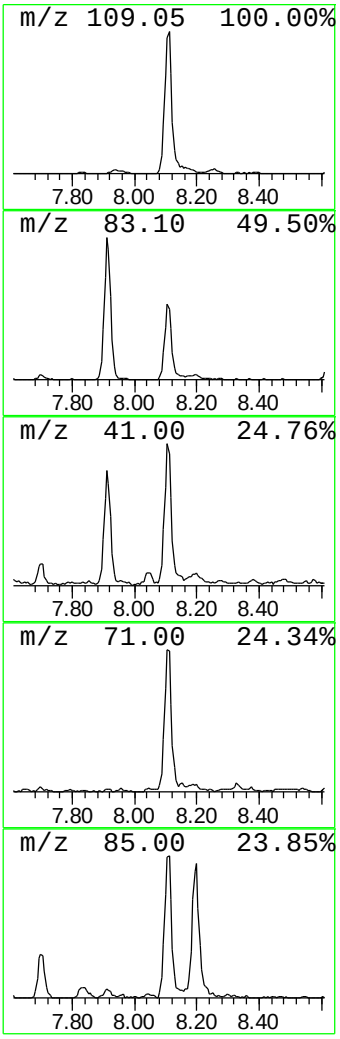
Peak Number 15 Cyclohexanol, 3,3,5-trimethyl- Concentration Rank 7

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 8.11 | 27.21 ng/ul | 1365430 | 1,4-Dichlorobenzene-d4 | 7.49 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Cyclohexanol, 3,3,5-trimethyl- | 142 | C9H18O | 000116-02-9 | 64 |
| 2 | 3,3,5,5-Tetramethylcyclopentene | 124 | C9H16 | 038667-10-6 | 43 |
| 3 | Cyclopentane, 2-ethylidene-1,1-d... | 124 | C9H16 | 056324-66-4 | 43 |
| 4 | 1-Buten-3-yne,4-trimethylsilyl | 124 | C7H12Si | 002696-32-4 | 38 |
| 5 | Cyclohexanol, 3,3,5-trimethyl-, ... | 142 | C9H18O | 000933-48-2 | 38 |



# Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN071620\
Data File : BN011190.D
Acq On : 16 Jul 2020 15:23
Operator : CG/JU
Sample : L3261-07
Misc :
ALS Vial : 6 Sample Multiplier: 1

**Instrument :** BNA_N
**ClientSampleId :** BE506

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN071520MA.M
Quant Title : SVOA CALIBRATION

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

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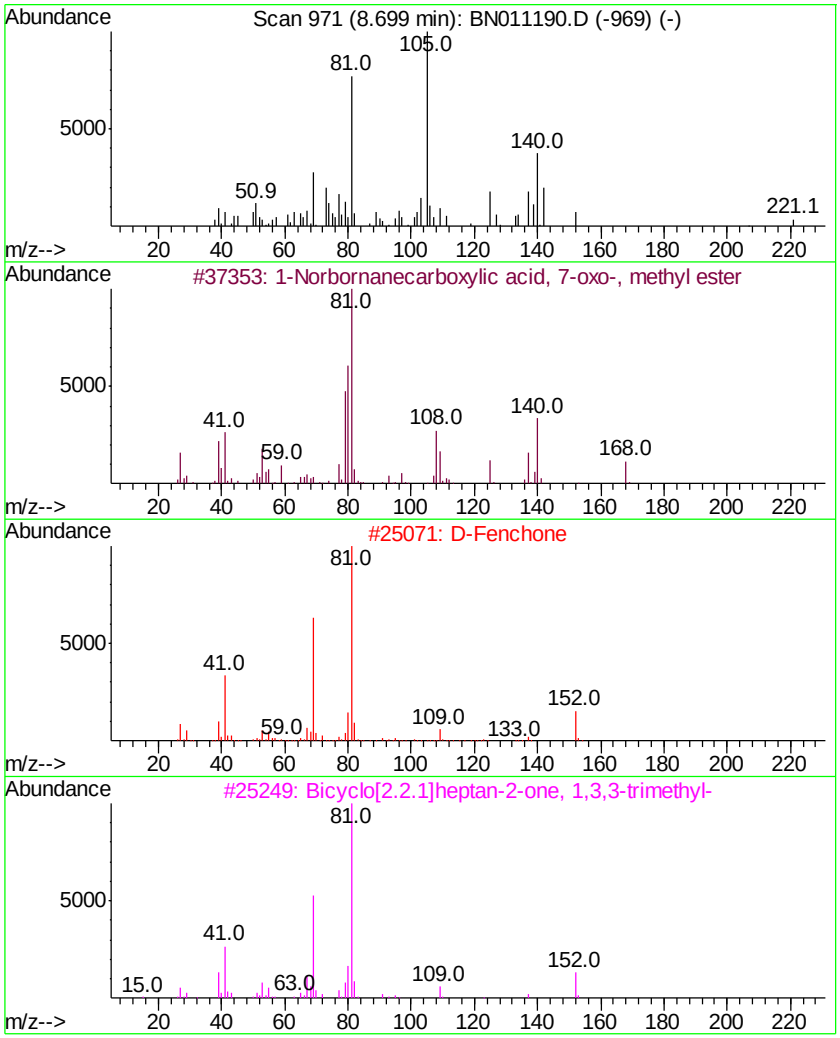
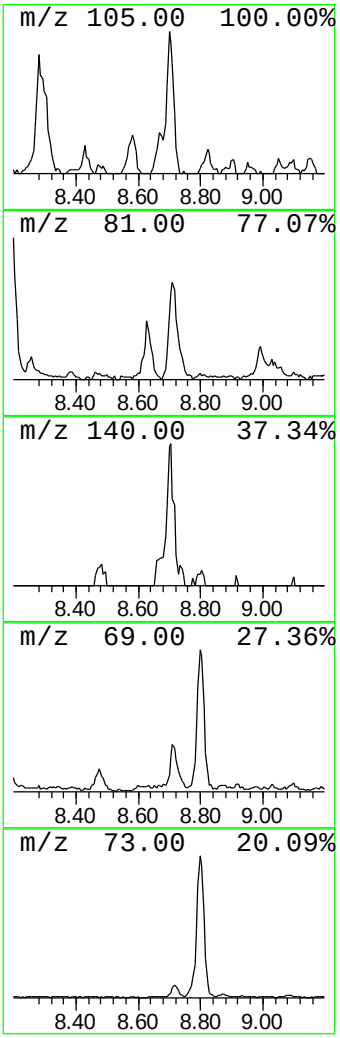
Peak Number 16 unknown-02 Concentration Rank 25

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 8.70 | 4.30 ng/ul | 216015 | 1,4-Dichlorobenzene-d4 | 7.49 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1-Norbornanecarboxylic acid, 7-o... | 168 | C9H12O3 | 090673-79-3 | 37 |
| 2 | D-Fenchone | 152 | C10H16O | 004695-62-9 | 35 |
| 3 | Bicyclo[2.2.1]heptan-2-one, 1,3,... | 152 | C10H16O | 001195-79-5 | 35 |
| 4 | Bicyclo[2.2.1]heptan-2-one, 1,3,... | 152 | C10H16O | 001195-79-5 | 35 |
| 5 | Cyclopropane, 1-methyl-1-isoprop... | 96 | C7H12 | 003422-07-9 | 30 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN071620\
Data File : BN011190.D
Acq On : 16 Jul 2020 15:23
Operator : CG/JU
Sample : L3261-07
Misc :
ALS Vial : 6 Sample Multiplier: 1

**Instrument :** BNA_N
**ClientSampleId :** BE506

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN071520MA.M
Quant Title : SVOA CALIBRATION

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

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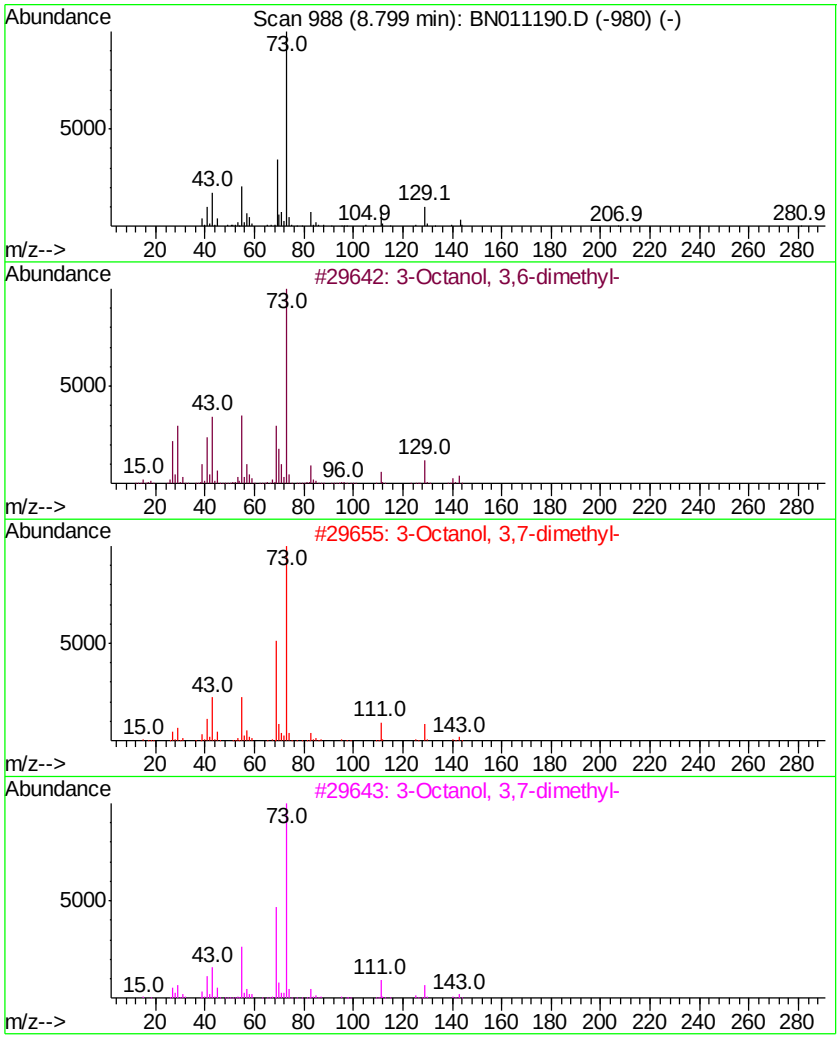
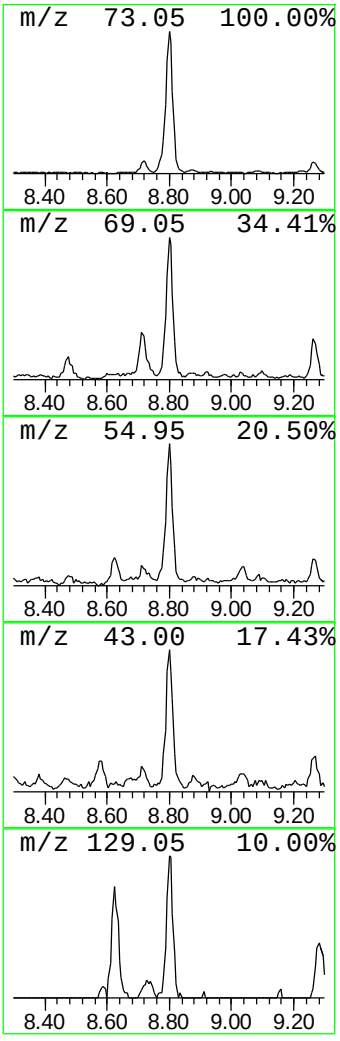
Peak Number 17 3-Octanol, 3,6-dimethyl- Concentration Rank 14

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 8.80 | 9.55 ng/ul | 479550 | 1,4-Dichlorobenzene-d4 | 7.49 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 3-Octanol, 3,6-dimethyl- | 158 | C10H22O | 000151-19-9 | 64 |
| 2 | 3-Octanol, 3,7-dimethyl- | 158 | C10H22O | 000078-69-3 | 59 |
| 3 | 3-Octanol, 3,7-dimethyl- | 158 | C10H22O | 000078-69-3 | 50 |
| 4 | 1,3-Dioxolane, 2-butyl- | 130 | C7H14O2 | 004360-76-3 | 37 |
| 5 | 3-Dodecanol, 3,7,11-trimethyl- | 228 | C15H32O | 007278-65-1 | 27 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN071620\
Data File : BN011190.D
Acq On : 16 Jul 2020 15:23
Operator : CG/JU
Sample : L3261-07
Misc :
ALS Vial : 6 Sample Multiplier: 1

**Instrument :** BNA_N
**ClientSampleId :** BE506

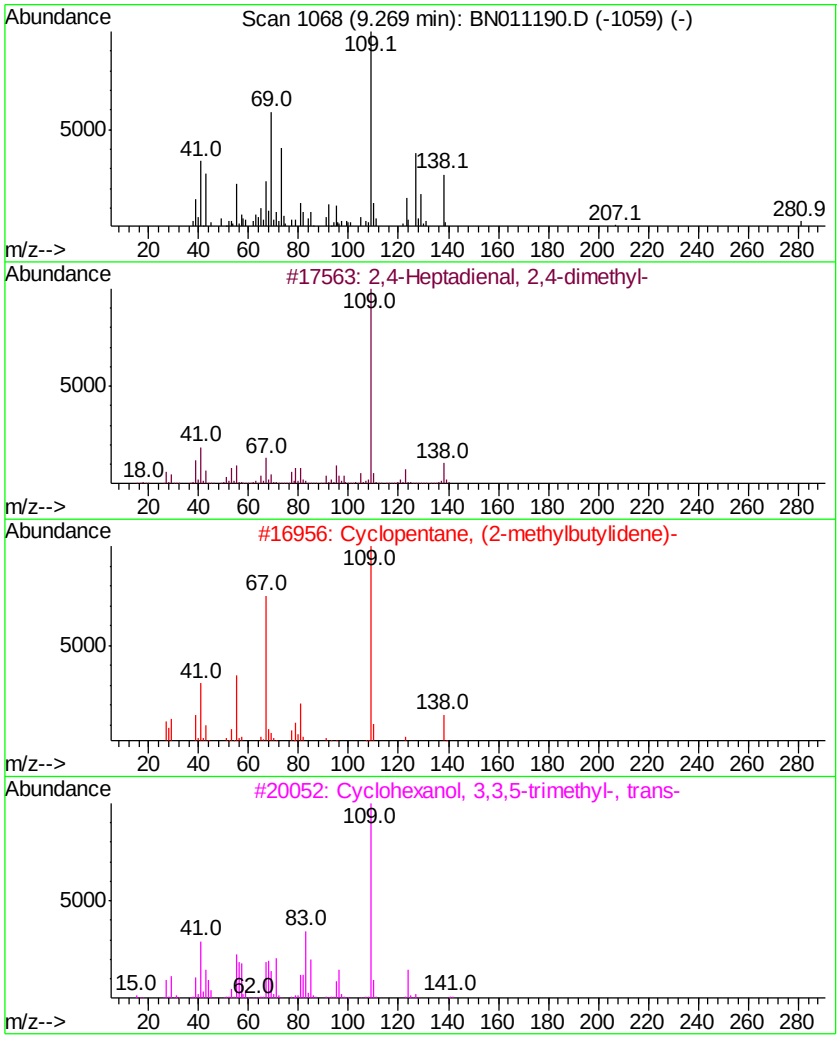
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN071520MA.M
Quant Title : SVOA CALIBRATION

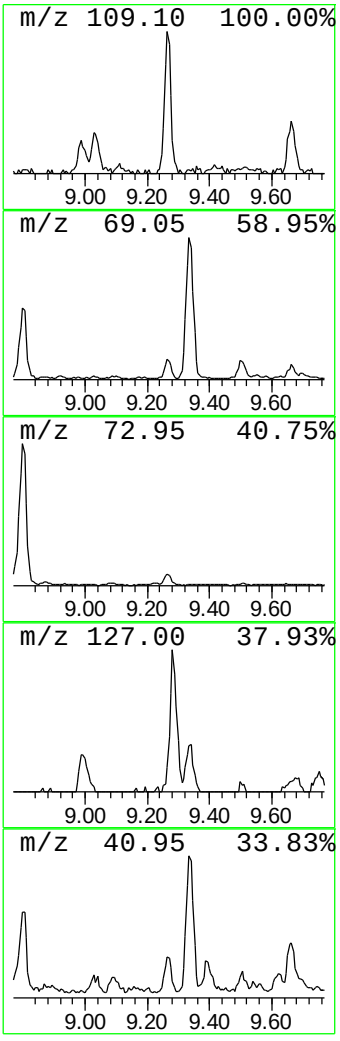
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 18 2,4-Heptadienal, 2,4-dimethyl- Concentration Rank 33

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 9.27 | 3.24 ng/ul | 230695 | Naphthalene-d8 | 10.23 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2,4-Heptadienal, 2,4-dimethyl- | 138 | C9H14O | 042452-48-2 | 58 |
| 2 | Cyclopentane, (2-methylbutylidene)- | 138 | C10H18 | 053366-54-4 | 50 |
| 3 | Cyclohexanol, 3,3,5-trimethyl-, ... | 142 | C9H18O | 000767-54-4 | 47 |
| 4 | Bicyclo[2.2.2]octane, 1-iodo- | 236 | C8H13I | 000931-98-6 | 47 |
| 5 | 1-Propanone, 1-(5-methyl-2-furan... | 138 | C8H10O2 | 010599-69-6 | 46 |



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Data File : BN011190.D
Acq On : 16 Jul 2020 15:23
Operator : CG/JU
Sample : L3261-07
Misc :
ALS Vial : 6 Sample Multiplier: 1

**Instrument :** BNA_N
**ClientSampleId :** BE506

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN071520MA.M
Quant Title : SVOA CALIBRATION

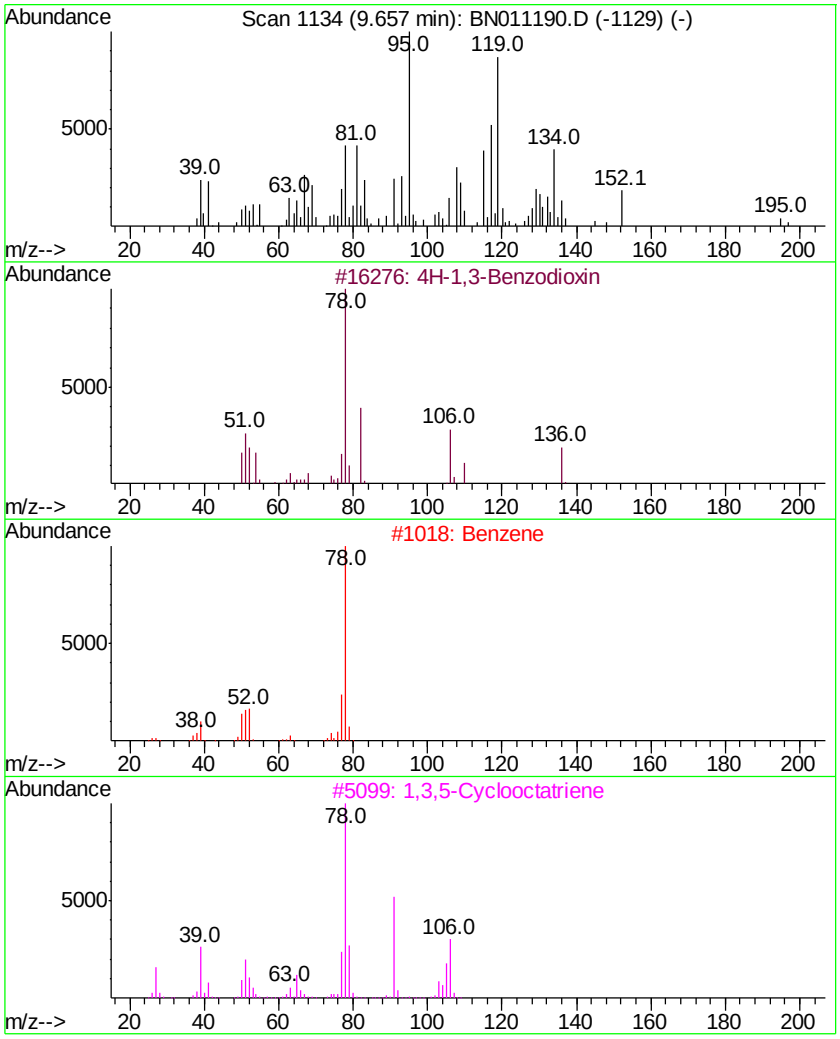
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

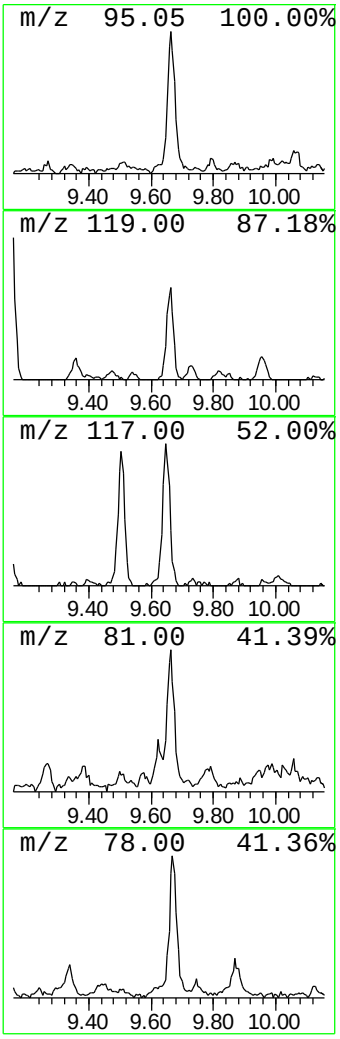
**********

Peak Number 19 unknown-03 Concentration Rank 21

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 9.66 | 6.83 ng/ul | 486626 | Naphthalene-d8 | 10.23 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 4H-1,3-Benzodioxin | 136 | C8H8O2 | 000254-27-3 | 49 |
| 2 | Benzene | 78 | C6H6 | 000071-43-2 | 45 |
| 3 | 1,3,5-Cyclooctatriene | 106 | C8H10 | 001871-52-9 | 43 |
| 4 | 2-Coumaranone | 134 | C8H6O2 | 000553-86-6 | 43 |
| 5 | 2,4,6-Cycloheptatrien-1-one | 106 | C7H6O | 000539-80-0 | 38 |



# Library Search Compound Report

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Data File : BN011190.D
Acq On : 16 Jul 2020 15:23
Operator : CG/JU
Sample : L3261-07
Misc :
ALS Vial : 6 Sample Multiplier: 1

**Instrument :** BNA_N
**ClientSampleId :** BE506

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN071520MA.M
Quant Title : SVOA CALIBRATION

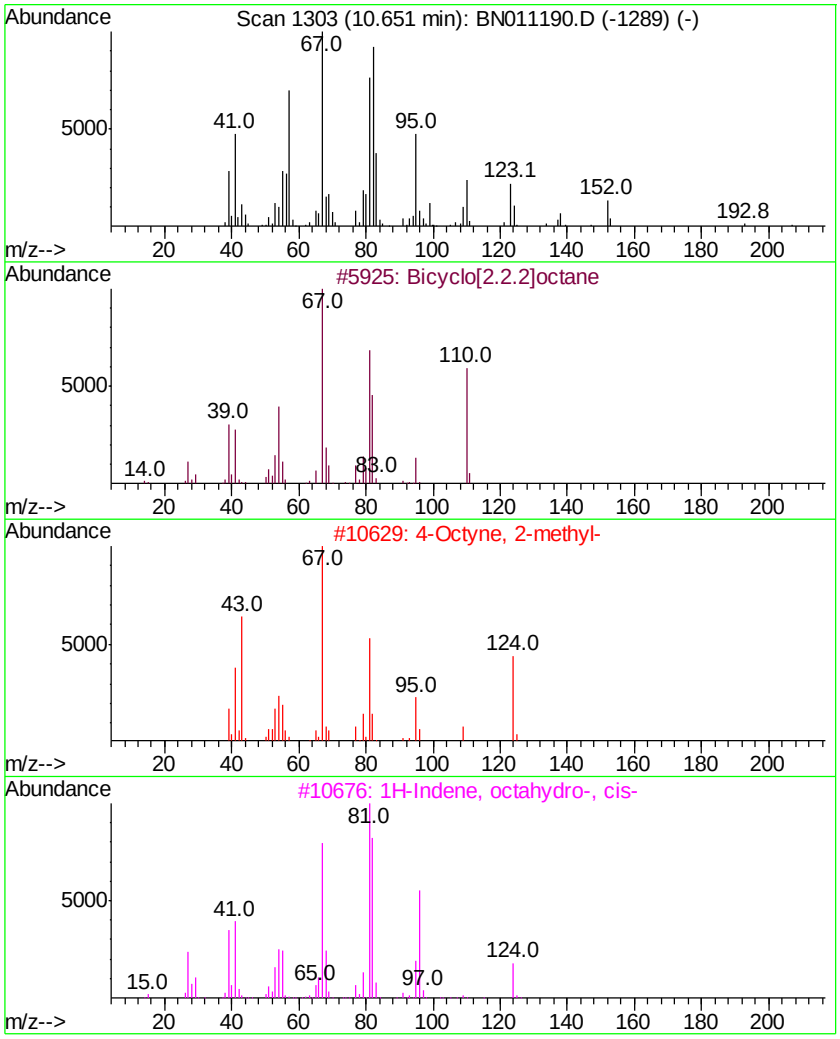
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

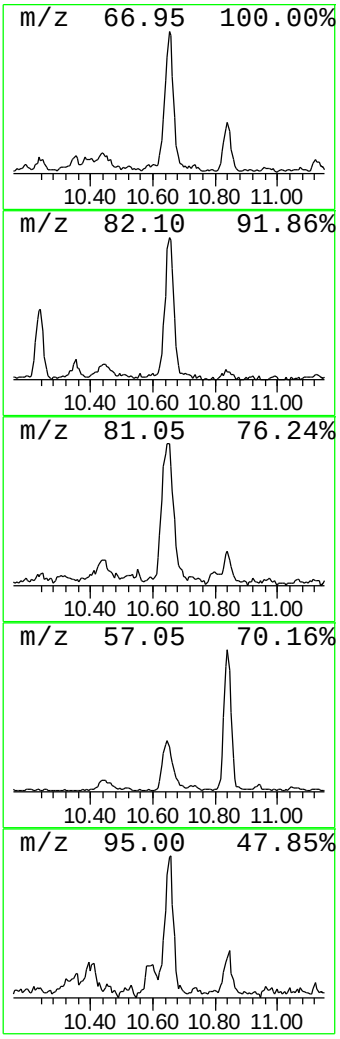
**********

Peak Number 20 Bicyclo[2.2.2]octane Concentration Rank 16

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 10.65 | 8.87 ng/ul | 631752 | Naphthalene-d8 | 10.23 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Bicyclo[2.2.2]octane | 110 | C8H14 | 000280-33-1 | 53 |
| 2 | 4-Octyne, 2-methyl- | 124 | C9H16 | 010306-94-2 | 46 |
| 3 | 1H-Indene, octahydro-, cis- | 124 | C9H16 | 004551-51-3 | 43 |
| 4 | Methyl ethyl cyclopentene | 110 | C8H14 | 019780-56-4 | 41 |
| 5 | 1-Ethyl-5-methylcyclopentene | 110 | C8H14 | 097797-57-4 | 41 |



# Library Search Compound Report

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Data File : BN011190.D
Acq On : 16 Jul 2020 15:23
Operator : CG/JU
Sample : L3261-07
Misc :
ALS Vial : 6 Sample Multiplier: 1

**Instrument :** BNA_N
**ClientSampleId :** BE506

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN071520MA.M
Quant Title : SVOA CALIBRATION

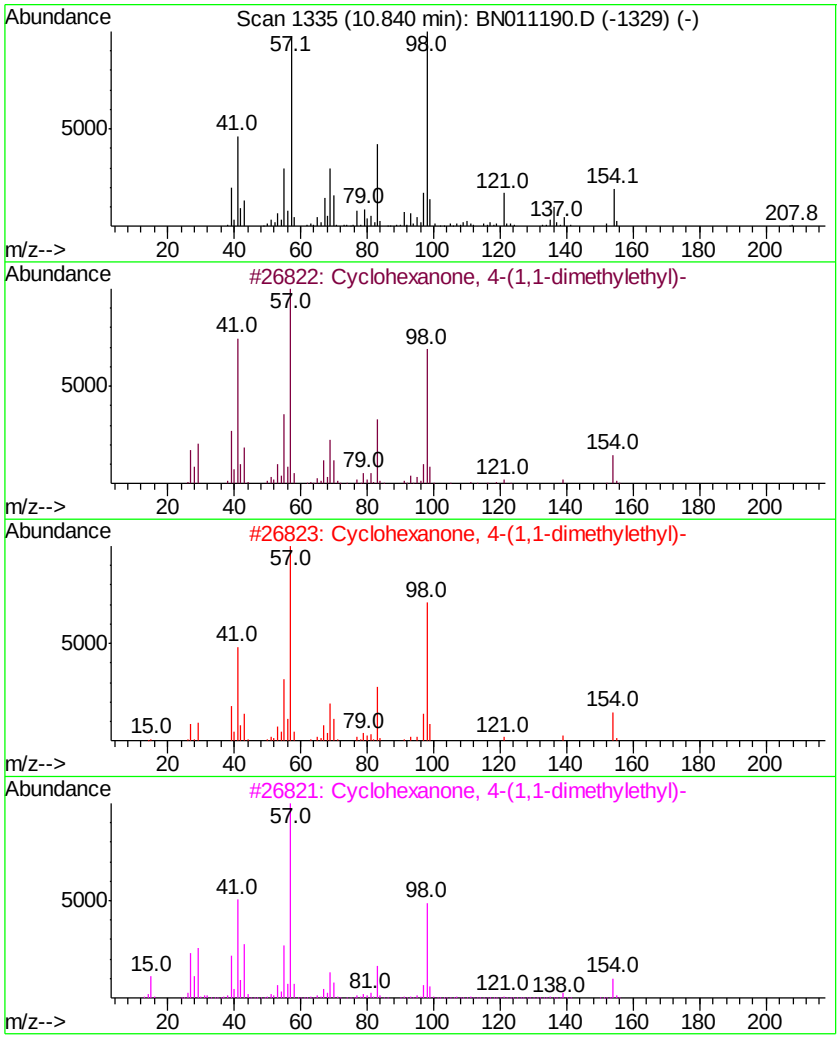
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

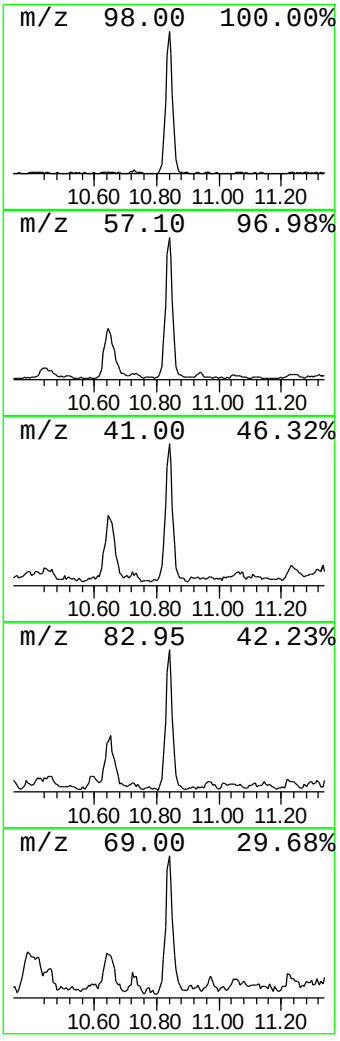
**********

Peak Number 21 Cyclohexanone, 4-(1,1-dimet... Concentration Rank 12

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 10.84 | 10.15 ng/ul | 722907 | Naphthalene-d8 | 10.23 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Cyclohexanone, 4-(1,1-dimethylet... | 154 | C10H18O | 000098-53-3 | 93 |
| 2 | Cyclohexanone, 4-(1,1-dimethylet... | 154 | C10H18O | 000098-53-3 | 81 |
| 3 | Cyclohexanone, 4-(1,1-dimethylet... | 154 | C10H18O | 000098-53-3 | 76 |
| 4 | Cyclohexanone, 4-(1,1-dimethylet... | 154 | C10H18O | 000098-53-3 | 64 |
| 5 | 1H-Inden-1-one, octahydro-7a-hyd... | 154 | C9H14O2 | 010421-80-4 | 58 |



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Data File : BN011190.D
Acq On : 16 Jul 2020 15:23
Operator : CG/JU
Sample : L3261-07
Misc :
ALS Vial : 6 Sample Multiplier: 1

**Instrument :** BNA_N
**ClientSampleId :** BE506

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN071520MA.M
Quant Title : SVOA CALIBRATION

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

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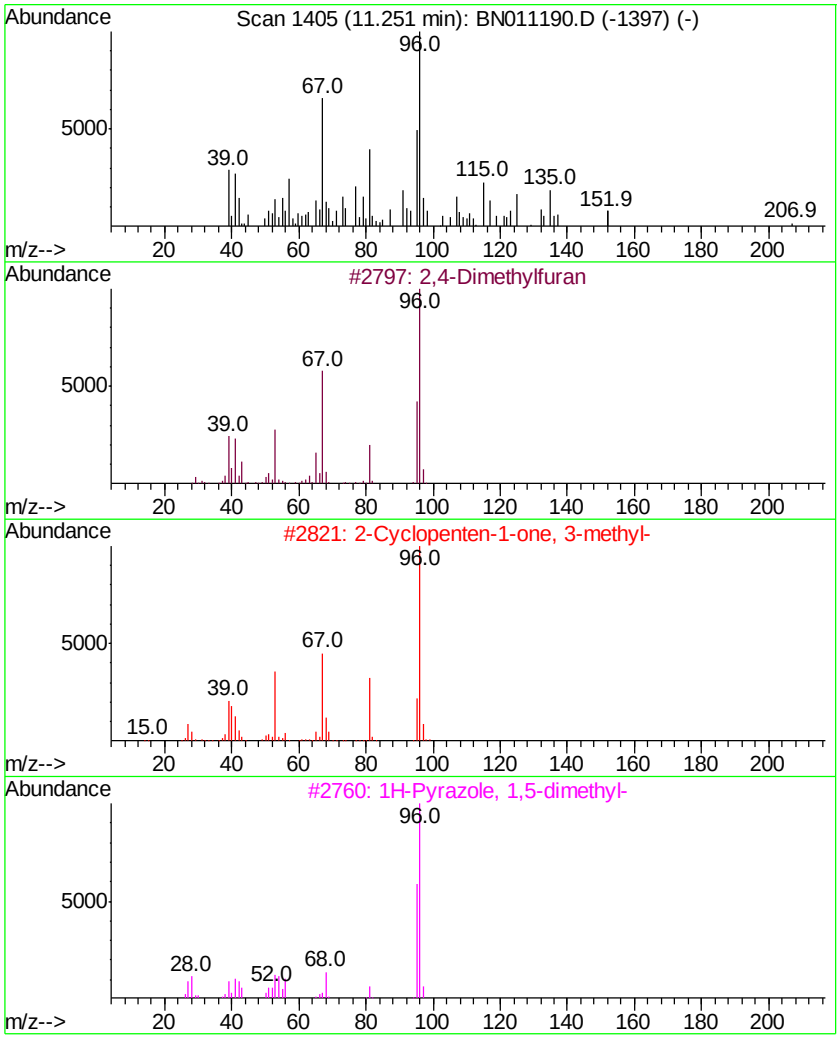
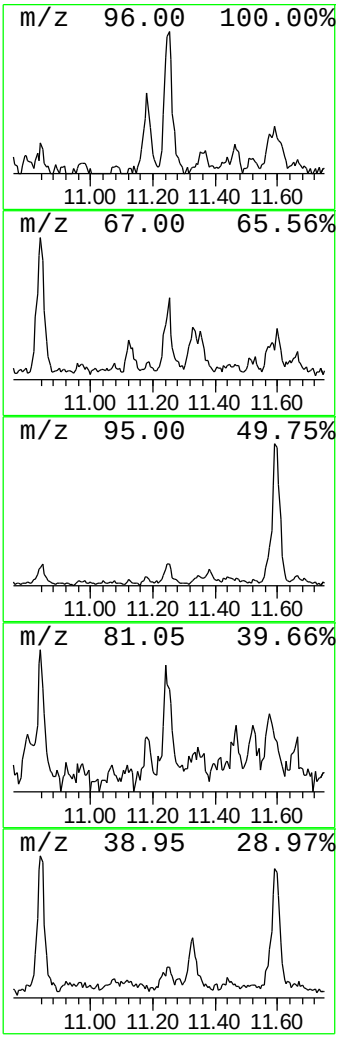
Peak Number 22 unknown-04 Concentration Rank 34

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 11.25 | 3.21 ng/ul | 228846 | Naphthalene-d8 | 10.23 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2,4-Dimethylfuran | 96 | C6H8O | 003710-43-8 | 47 |
| 2 | 2-Cyclopenten-1-one, 3-methyl- | 96 | C6H8O | 002758-18-1 | 38 |
| 3 | 1H-Pyrazole, 1,5-dimethyl- | 96 | C5H8N2 | 000694-31-5 | 38 |
| 4 | 1H-Pyrazole, 1,3-dimethyl- | 96 | C5H8N2 | 000694-48-4 | 35 |
| 5 | 1H-Imidazole, 1,2-dimethyl- | 96 | C5H8N2 | 001739-84-0 | 35 |



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Data File : BN011190.D
Acq On : 16 Jul 2020 15:23
Operator : CG/JU
Sample : L3261-07
Misc :
ALS Vial : 6 Sample Multiplier: 1

**Instrument :** BNA_N
**ClientSampleId :** BE506

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN071520MA.M
Quant Title : SVOA CALIBRATION

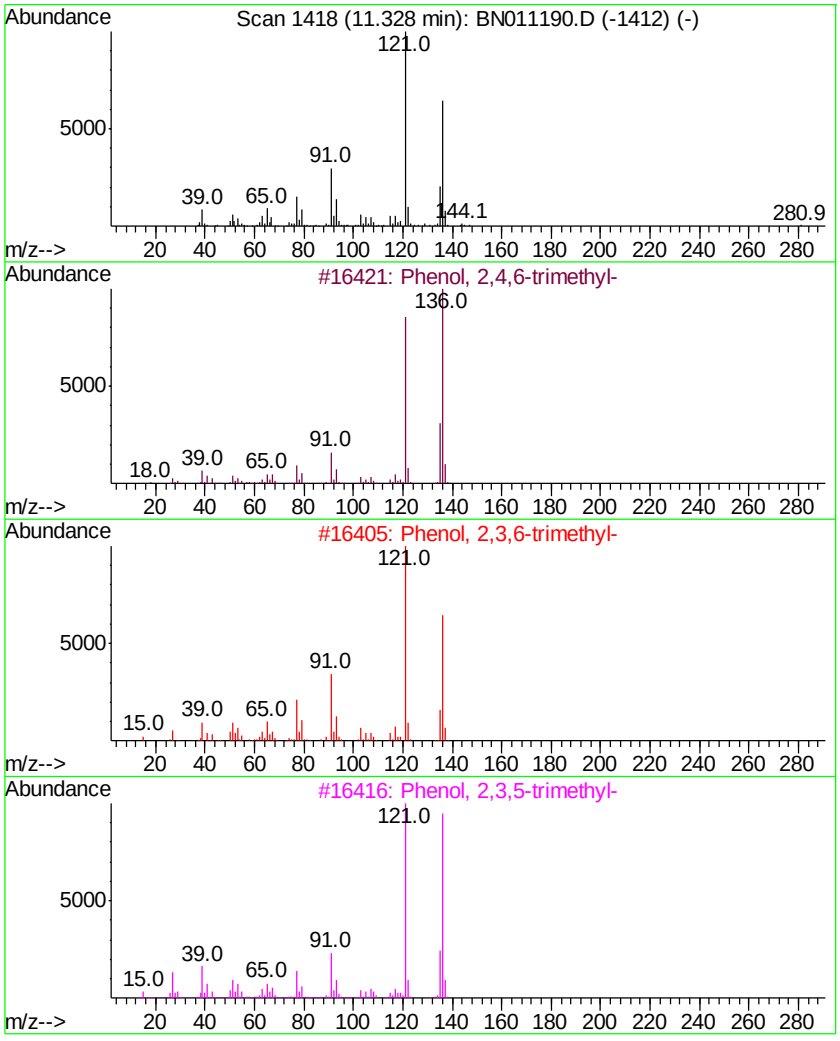
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

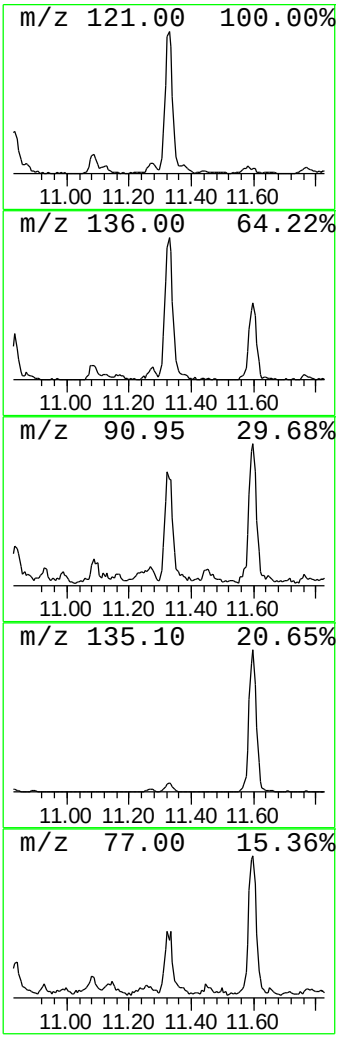
**********

Peak Number 23 Phenol, 2,4,6-trimethyl- Concentration Rank 19

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 11.33 | 7.18 ng/ul | 511414 | Naphthalene-d8 | 10.23 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Phenol, 2,4,6-trimethyl- | 136 | C9H12O | 000527-60-6 | 97 |
| 2 | Phenol, 2,3,6-trimethyl- | 136 | C9H12O | 002416-94-6 | 97 |
| 3 | Phenol, 2,3,5-trimethyl- | 136 | C9H12O | 000697-82-5 | 94 |
| 4 | Phenol, 2,4,6-trimethyl- | 136 | C9H12O | 000527-60-6 | 94 |
| 5 | Phenol, 2,3,6-trimethyl- | 136 | C9H12O | 002416-94-6 | 94 |



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Data File : BN011190.D
Acq On : 16 Jul 2020 15:23
Operator : CG/JU
Sample : L3261-07
Misc :
ALS Vial : 6 Sample Multiplier: 1

**Instrument :** BNA_N
**ClientSampleId :** BE506

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN071520MA.M
Quant Title : SVOA CALIBRATION

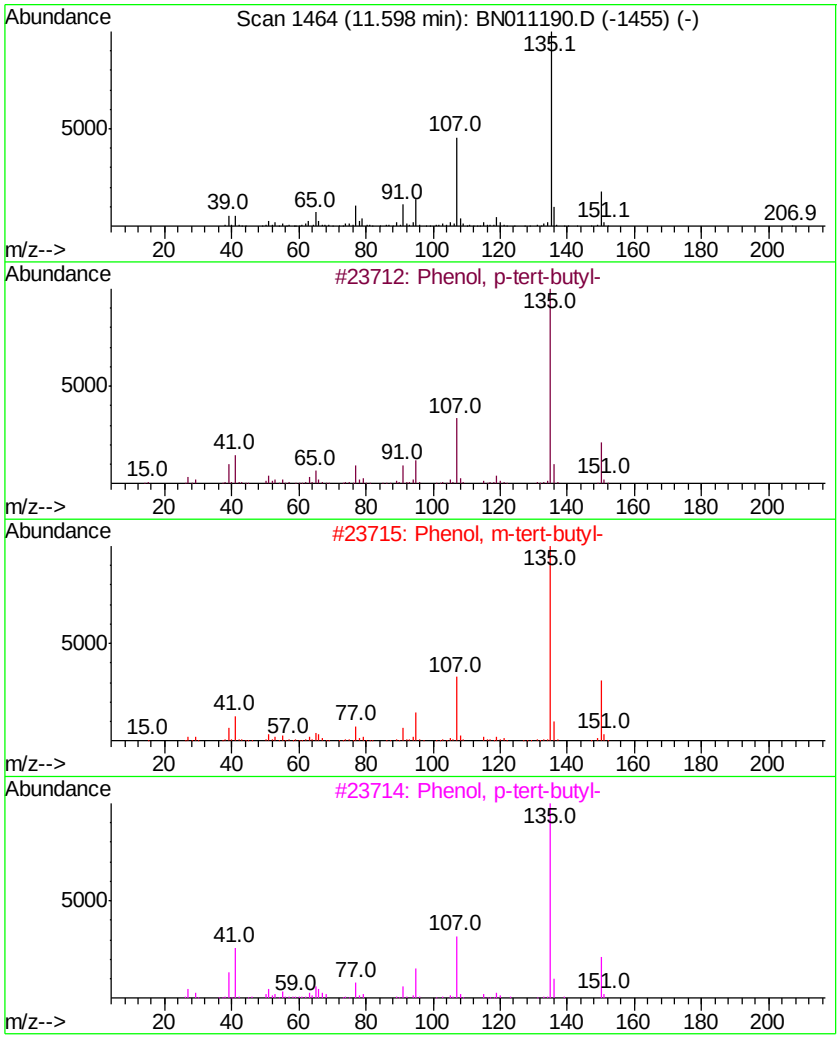
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

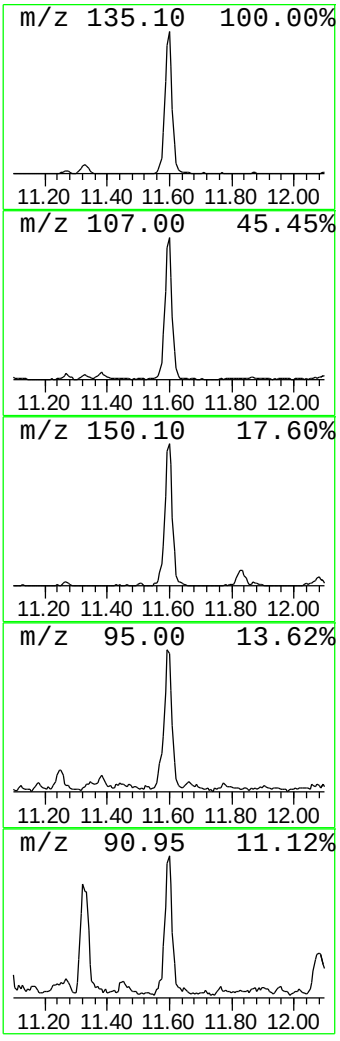
**********

Peak Number 24 Phenol, p-tert-butyl- Concentration Rank 9

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 11.60 | 17.53 ng/ul | 1248910 | Naphthalene-d8 | 10.23 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Phenol, p-tert-butyl- | 150 | C10H14O | 000098-54-4 | 94 |
| 2 | Phenol, m-tert-butyl- | 150 | C10H14O | 000585-34-2 | 91 |
| 3 | Phenol, p-tert-butyl- | 150 | C10H14O | 000098-54-4 | 91 |
| 4 | Phenol, m-tert-butyl- | 150 | C10H14O | 000585-34-2 | 81 |
| 5 | Phenol, 2-(1,1-dimethylethyl)- | 150 | C10H14O | 000088-18-6 | 58 |



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Data File : BN011190.D
Acq On : 16 Jul 2020 15:23
Operator : CG/JU
Sample : L3261-07
Misc :
ALS Vial : 6 Sample Multiplier: 1

**Instrument :** BNA_N
**ClientSampleId :** BE506

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN071520MA.M
Quant Title : SVOA CALIBRATION

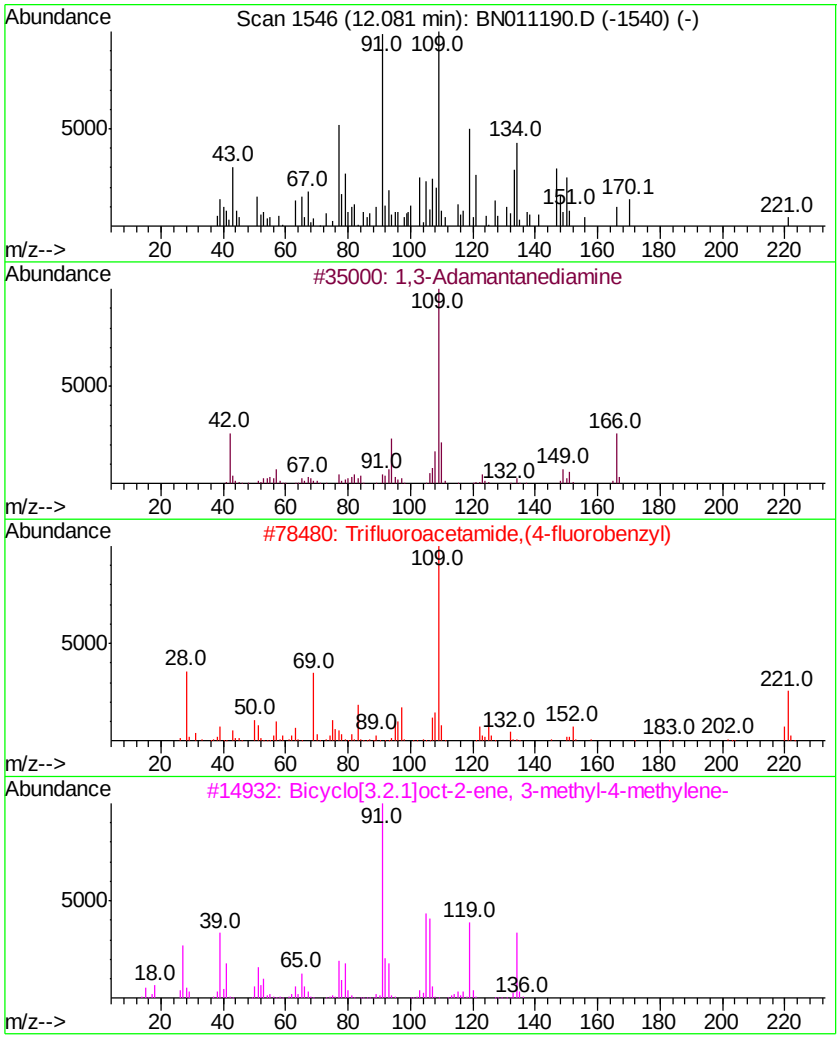
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

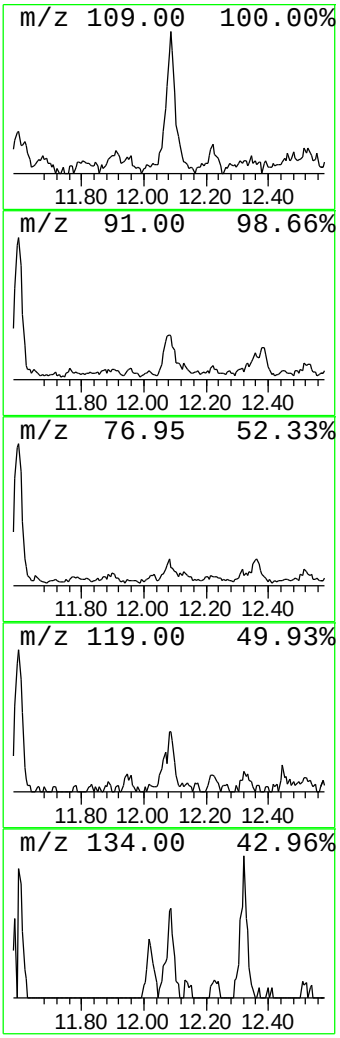
**********

Peak Number 25 unknown-05 Concentration Rank 39

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 12.08 | 2.59 ng/ul | 184759 | Naphthalene-d8 | 10.23 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1,3-Adamantanediamine | 166 | C10H18N2 | 026562-81-2 | 22 |
| 2 | Trifluoroacetamide,(4-fluorobenzyl) | 221 | C9H7F4NO | 170374-90-0 | 22 |
| 3 | Bicyclo[3.2.1]oct-2-ene, 3-methy... | 134 | C10H14 | 049826-53-1 | 18 |
| 4 | Tetracyclo[3.3.1.1(1,8).0(2,4)]d... | 134 | C10H14 | 1000185-58-7 | 18 |
| 5 | 2-Hydrazinopyridine | 109 | C5H7N3 | 004930-98-7 | 14 |



# Library Search Compound Report

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Data File : BN011190.D
Acq On : 16 Jul 2020 15:23
Operator : CG/JU
Sample : L3261-07
Misc :
ALS Vial : 6 Sample Multiplier: 1

**Instrument :** BNA_N
**ClientSampleId :** BE506

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN071520MA.M
Quant Title : SVOA CALIBRATION

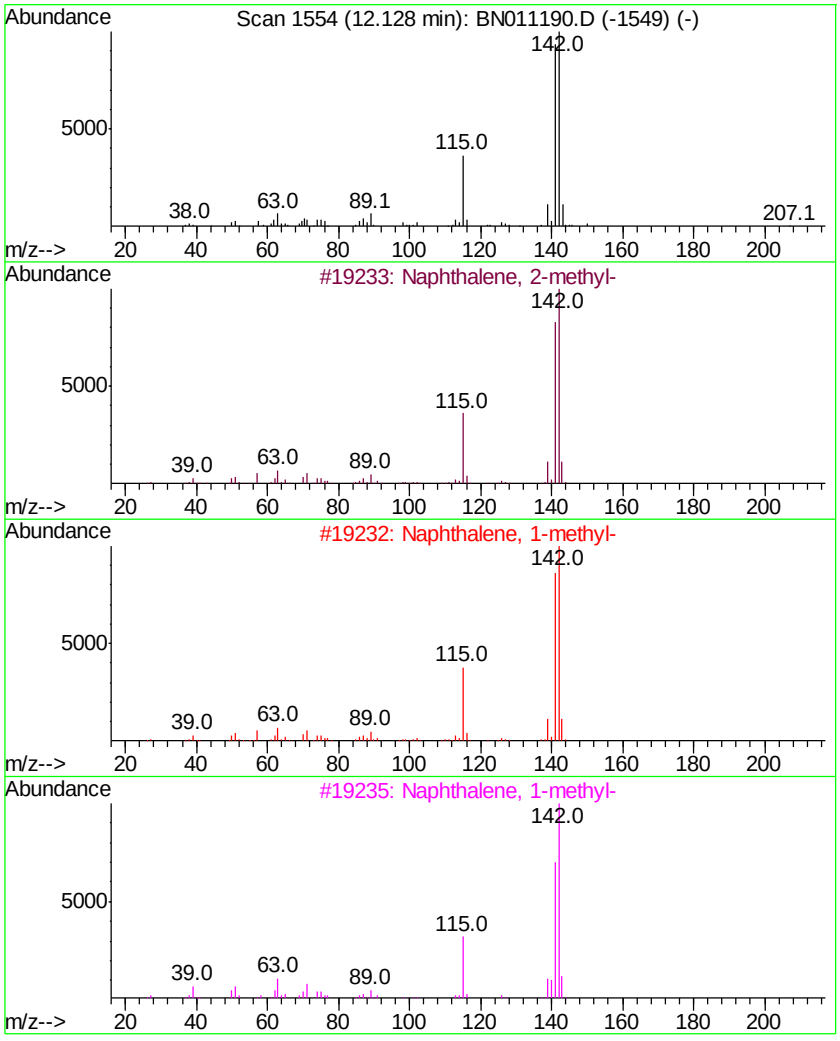
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

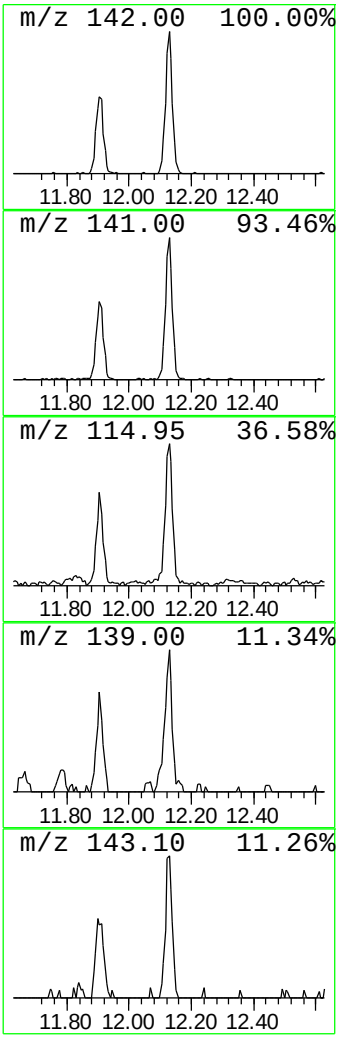
**********

Peak Number 26 Naphthalene, 1-methyl- Concentration Rank 23

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 12.13 | 5.29 ng/ul | 376934 | Naphthalene-d8 | 10.23 |

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



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN071620\
Data File : BN011190.D
Acq On : 16 Jul 2020 15:23
Operator : CG/JU
Sample : L3261-07
Misc :
ALS Vial : 6 Sample Multiplier: 1

**Instrument :** BNA_N
**ClientSampleId :** BE506

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN071520MA.M
Quant Title : SVOA CALIBRATION

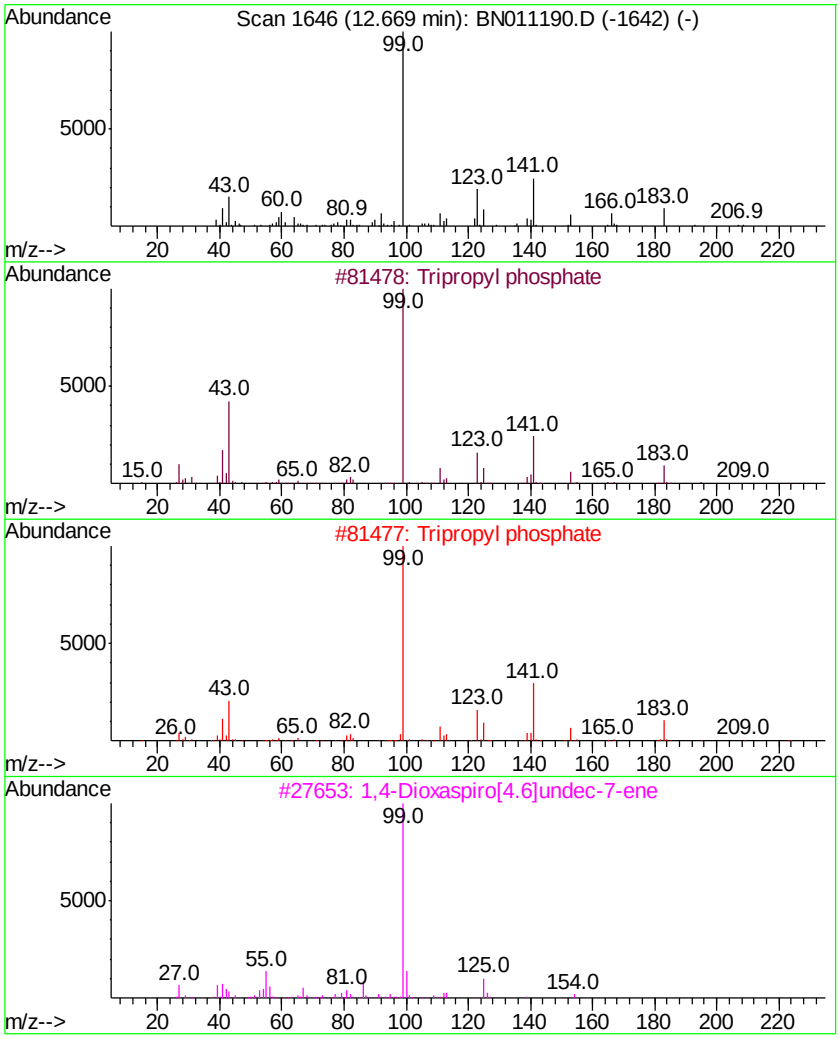
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

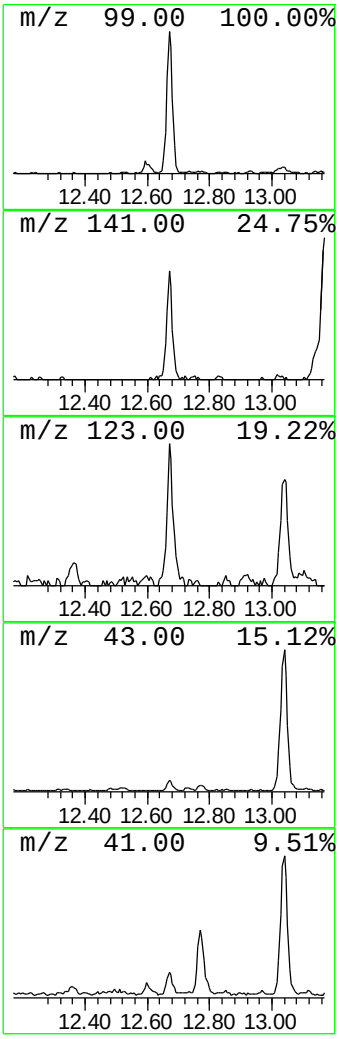
**********

Peak Number 27 Tripropyl phosphate Concentration Rank 36

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 12.67 | 2.88 ng/ul | 312851 | Acenaphthene-d10 | 14.12 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Tripropyl phosphate | 224 | C9H21O4P | 000513-08-6 | 64 |
| 2 | Tripropyl phosphate | 224 | C9H21O4P | 000513-08-6 | 59 |
| 3 | 1,4-Dioxaspiro[4.6]undec-7-ene | 154 | C9H14O2 | 007140-60-5 | 35 |
| 4 | Hexanoic acid, 2-tridecyl ester | 298 | C19H38O2 | 055193-20-9 | 25 |
| 5 | n-Butyl methylphosphonofluoridate | 154 | C5H12FO2P | 000352-63-6 | 17 |



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Data File : BN011190.D
Acq On : 16 Jul 2020 15:23
Operator : CG/JU
Sample : L3261-07
Misc :
ALS Vial : 6 Sample Multiplier: 1

**Instrument :** BNA_N
**ClientSampleId :** BE506

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN071520MA.M
Quant Title : SVOA CALIBRATION

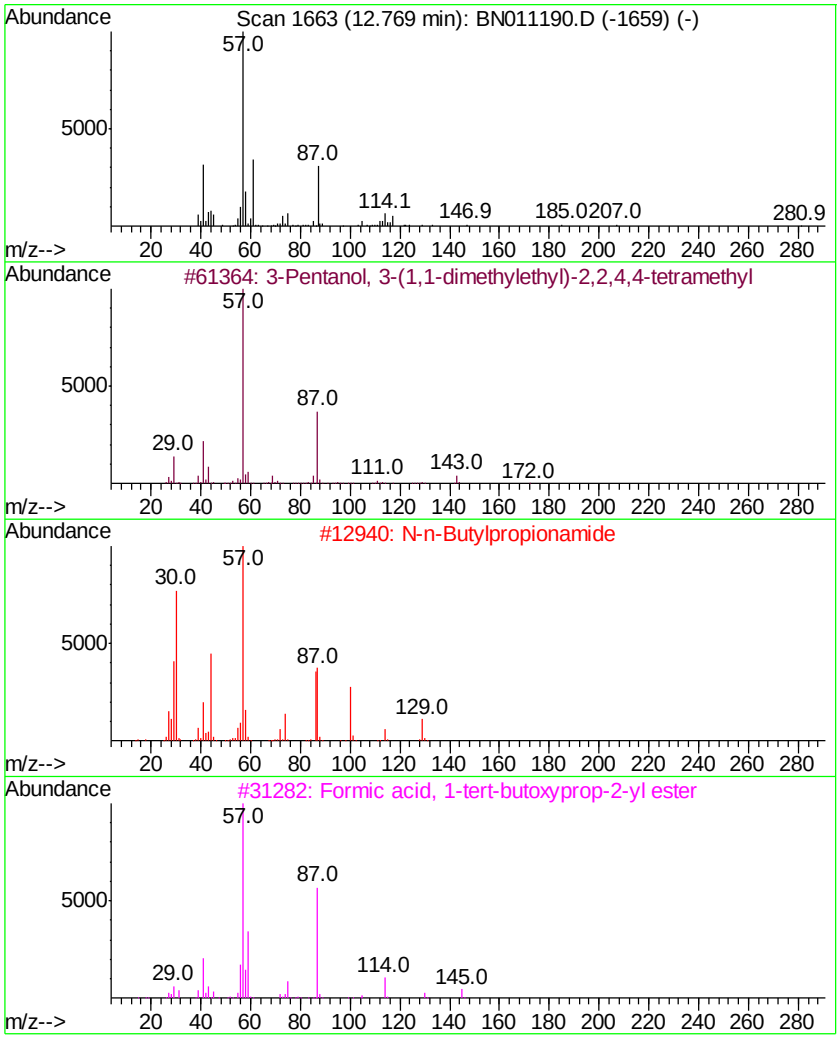
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

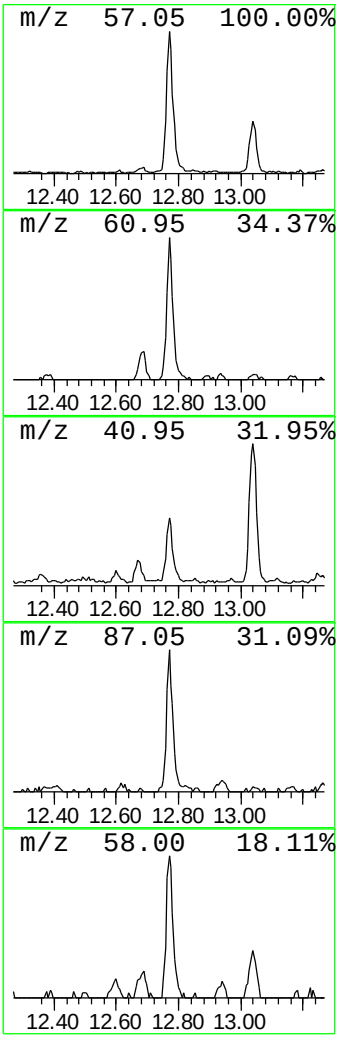
**********

Peak Number 28 unknown-06 Concentration Rank 43

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 12.77 | 2.08 ng/ul | 226333 | Acenaphthene-d10 | 14.12 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 3-Pentanol, 3-(1,1-dimethylethyl... | 200 | C13H28O | 041902-42-5 | 35 |
| 2 | N-n-Butylpropionamide | 129 | C7H15NO | 002955-67-1 | 28 |
| 3 | Formic acid, 1-tert-butoxyprop-2... | 160 | C8H16O3 | 1000368-95-6 | 25 |
| 4 | 3-Pentanone, 2,2-dimethyl- | 114 | C7H14O | 000564-04-5 | 16 |
| 5 | 1,2-Butanediol, 3,3-dimethyl- | 118 | C6H14O2 | 059562-82-2 | 12 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN071620\
Data File : BN011190.D
Acq On : 16 Jul 2020 15:23
Operator : CG/JU
Sample : L3261-07
Misc :
ALS Vial : 6 Sample Multiplier: 1

**Instrument :** BNA_N
**ClientSampleId :** BE506

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN071520MA.M
Quant Title : SVOA CALIBRATION

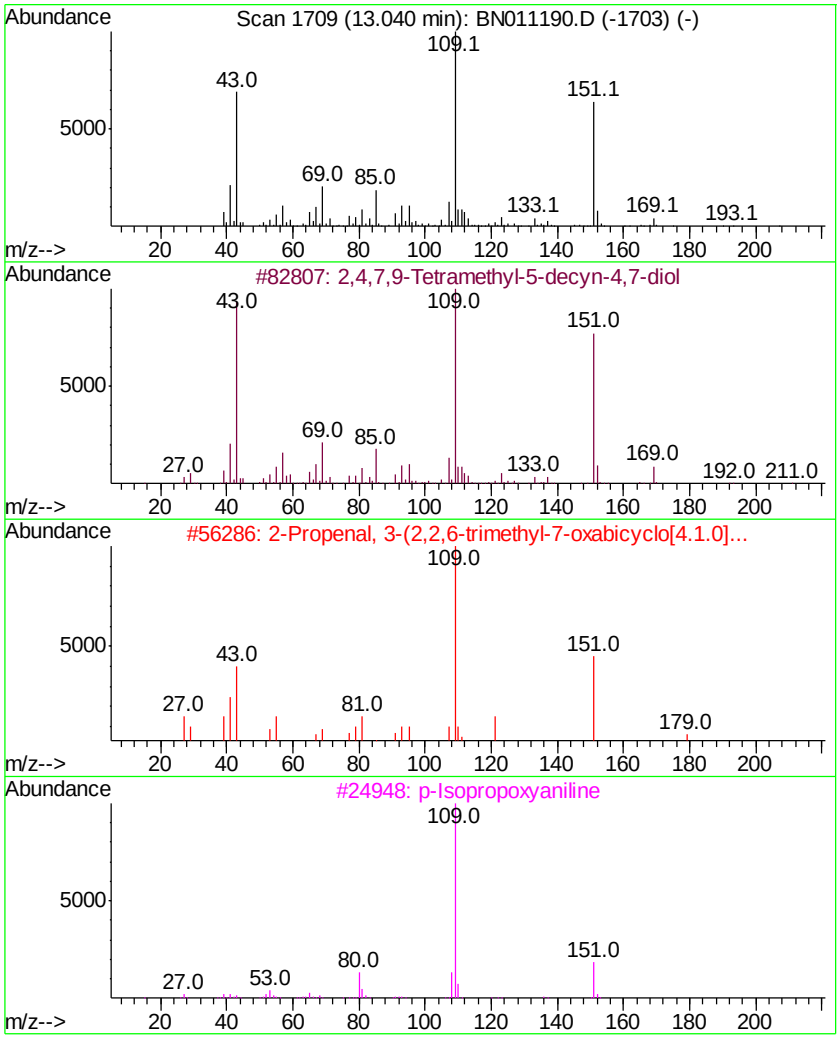
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

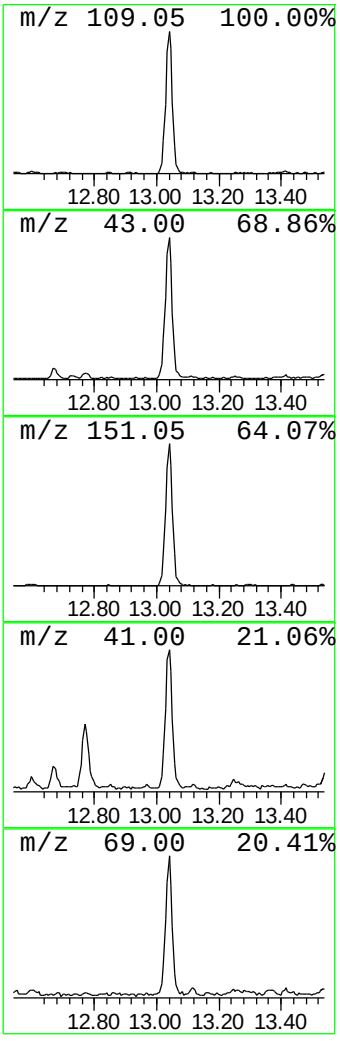
**********

Peak Number 29 2,4,7,9-Tetramethyl-5-decyn... Concentration Rank 11

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 13.04 | 13.00 ng/ul | 1411870 | Acenaphthene-d10 | 14.12 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2,4,7,9-Tetramethyl-5-decyn-4,7-... | 226 | C14H26O2 | 000126-86-3 | 91 |
| 2 | 2-Propenal, 3-(2,2,6-trimethyl-7... | 194 | C12H18O2 | 055759-91-6 | 64 |
| 3 | p-Isopropoxyaniline | 151 | C9H13NO | 007664-66-6 | 58 |
| 4 | m-Isopropoxyaniline | 151 | C9H13NO | 041406-00-2 | 53 |
| 5 | Acetaminophen | 151 | C8H9NO2 | 000103-90-2 | 53 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN071620\
Data File : BN011190.D
Acq On : 16 Jul 2020 15:23
Operator : CG/JU
Sample : L3261-07
Misc :
ALS Vial : 6 Sample Multiplier: 1

**Instrument :** BNA_N
**ClientSampleId :** BE506

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN071520MA.M
Quant Title : SVOA CALIBRATION

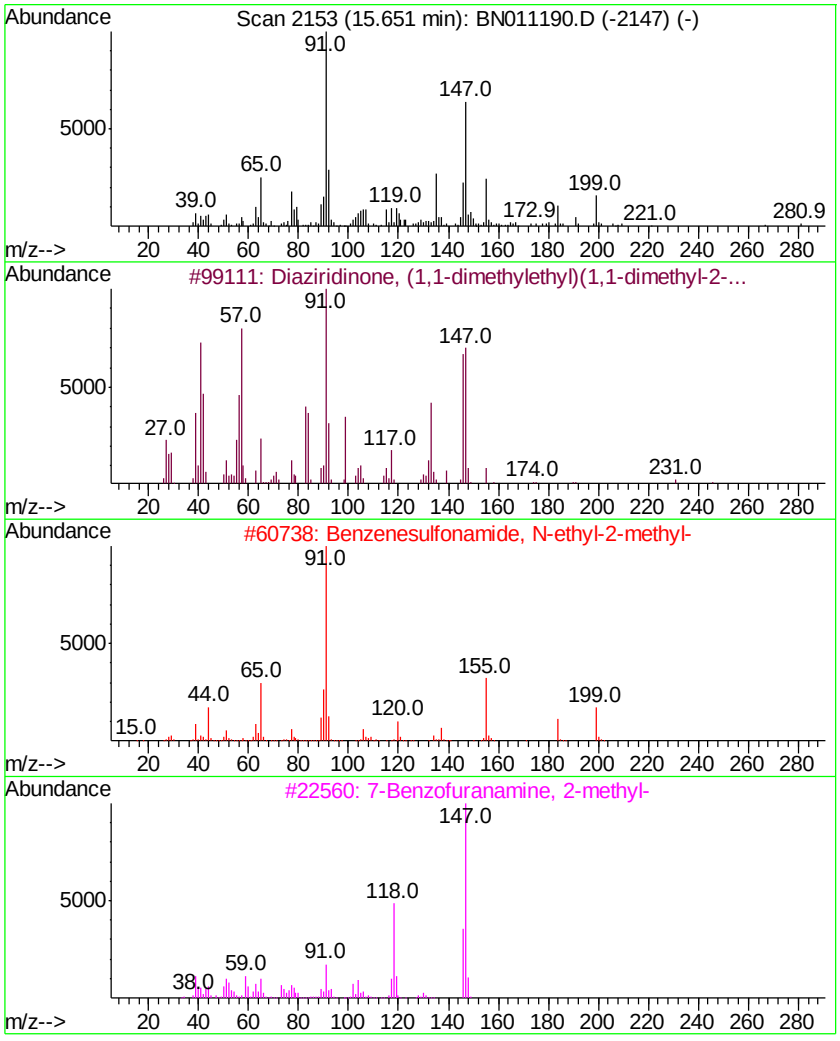
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

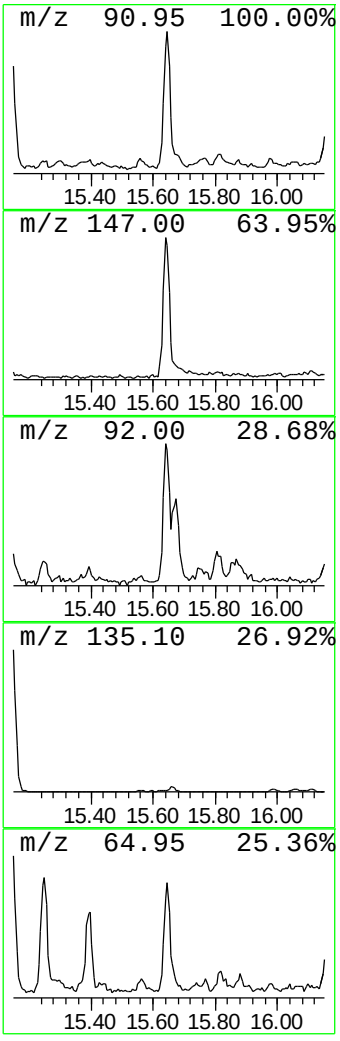
**********

Peak Number 30 Diaziridinone, (1,1-dimethy... Concentration Rank 27

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 15.65 | 3.91 ng/ul | 700547 | Phenanthrene-d10 | 16.87 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Diaziridinone, (1,1-dimethylethy... | 246 | C15H22N2O | 019656-67-8 | 52 |
| 2 | Benzenesulfonamide, N-ethyl-2-me... | 199 | C9H13NO2S | 001077-56-1 | 50 |
| 3 | 7-Benzofuranamine, 2-methyl- | 147 | C9H9NO | 1000327-46-4 | 49 |
| 4 | 1,3,2-Benzoxazaborole, 2-ethyl-2... | 147 | C8H10BNO | 129363-62-8 | 38 |
| 5 | Pyrido[3,2-d]pyrimidin-4-ol | 147 | C7H5N3O | 037538-67-3 | 38 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN071620\
Data File : BN011190.D
Acq On : 16 Jul 2020 15:23
Operator : CG/JU
Sample : L3261-07
Misc :
ALS Vial : 6 Sample Multiplier: 1

**Instrument :** BNA_N
**ClientSampleId :** BE506

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN071520MA.M
Quant Title : SVOA CALIBRATION

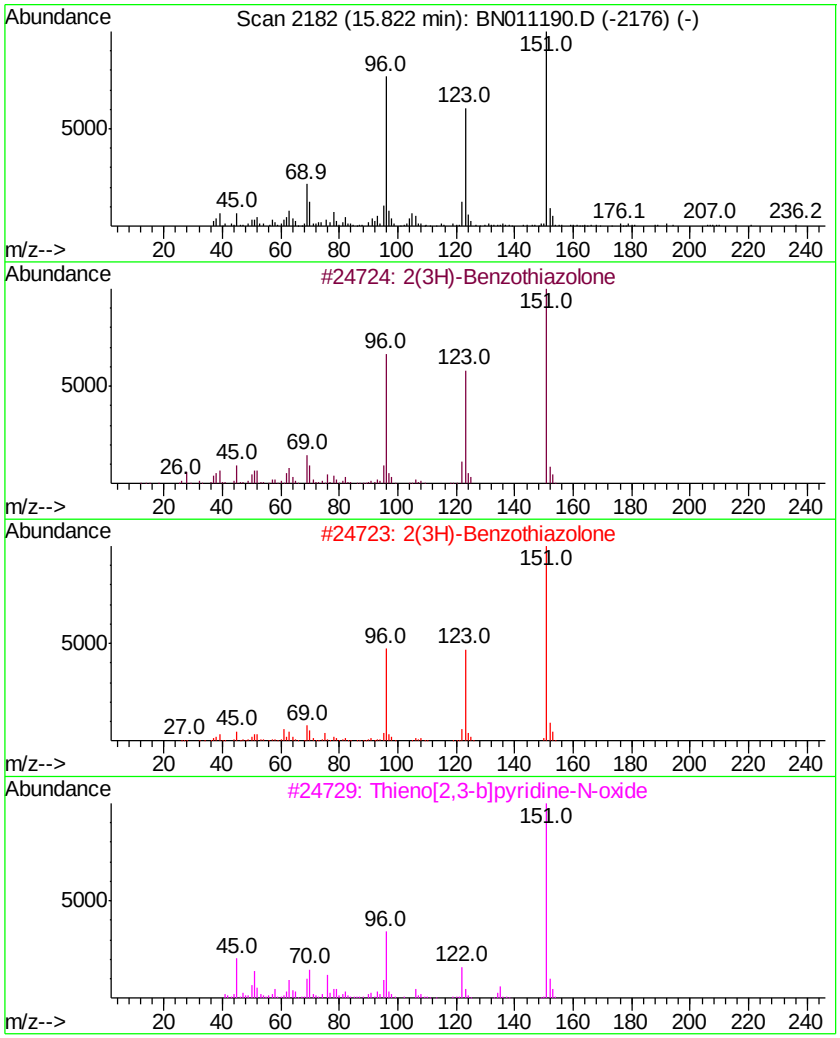
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

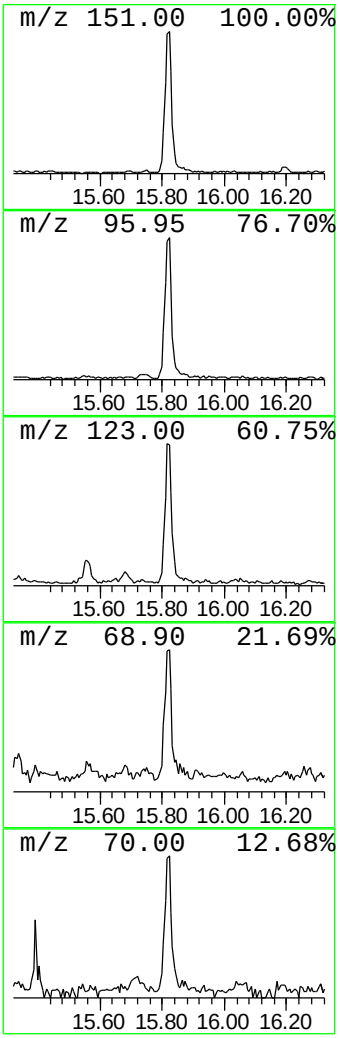
**********

Peak Number 31 2(3H)-Benzothiazolone Concentration Rank 32

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 15.82 | 3.26 ng/ul | 584216 | Phenanthrene-d10 | 16.87 |

| 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 | 91 |
| 3 | Thieno[2,3-b]pyridine-N-oxide | 151 | C7H5NOS | 025557-50-0 | 46 |
| 4 | 1,2,4-Triazolo[4,3-a]pyridine-3(... | 151 | C6H5N3S | 006952-68-7 | 42 |
| 5 | 2(3H)-Benzoxazolethione | 151 | C7H5NOS | 002382-96-9 | 38 |



# Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN071620\
Data File : BN011190.D
Acq On : 16 Jul 2020 15:23
Operator : CG/JU
Sample : L3261-07
Misc :
ALS Vial : 6 Sample Multiplier: 1

**Instrument :** BNA_N
**ClientSampleId :** BE506

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN071520MA.M
Quant Title : SVOA CALIBRATION

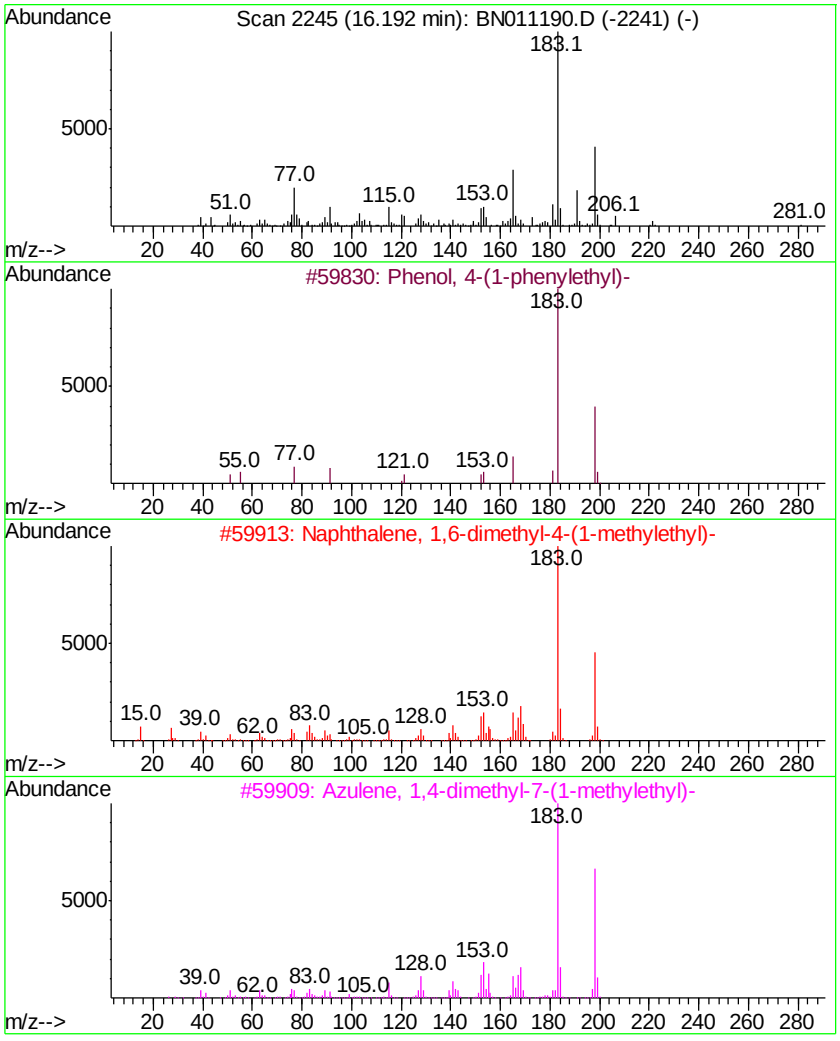
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

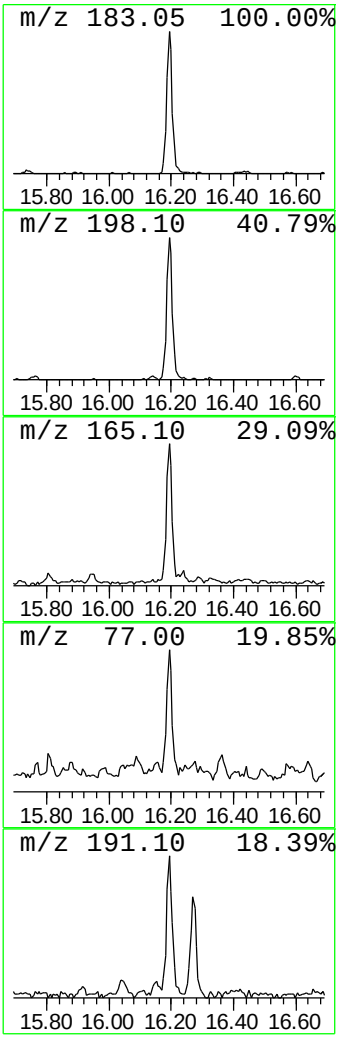
**********

Peak Number 32 Phenol, 4-(1-phenylethyl)- Concentration Rank 30

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 16.19 | 3.37 ng/ul | 603658 | Phenanthrene-d10 | 16.87 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Phenol, 4-(1-phenylethyl)- | 198 | C14H14O | 001988-89-2 | 76 |
| 2 | Naphthalene, 1,6-dimethyl-4-(1-m... | 198 | C15H18 | 000483-78-3 | 58 |
| 3 | Azulene, 1,4-dimethyl-7-(1-methy... | 198 | C15H18 | 000489-84-9 | 49 |
| 4 | Phenol, 2-(1-phenylethyl)- | 198 | C14H14O | 004237-44-9 | 46 |
| 5 | Naphthalene, 1,6-dimethyl-4-(1-m... | 198 | C15H18 | 000483-78-3 | 46 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN071620\
Data File : BN011190.D
Acq On : 16 Jul 2020 15:23
Operator : CG/JU
Sample : L3261-07
Misc :
ALS Vial : 6 Sample Multiplier: 1

**Instrument :** BNA_N
**ClientSampleId :** BE506

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN071520MA.M
Quant Title : SVOA CALIBRATION

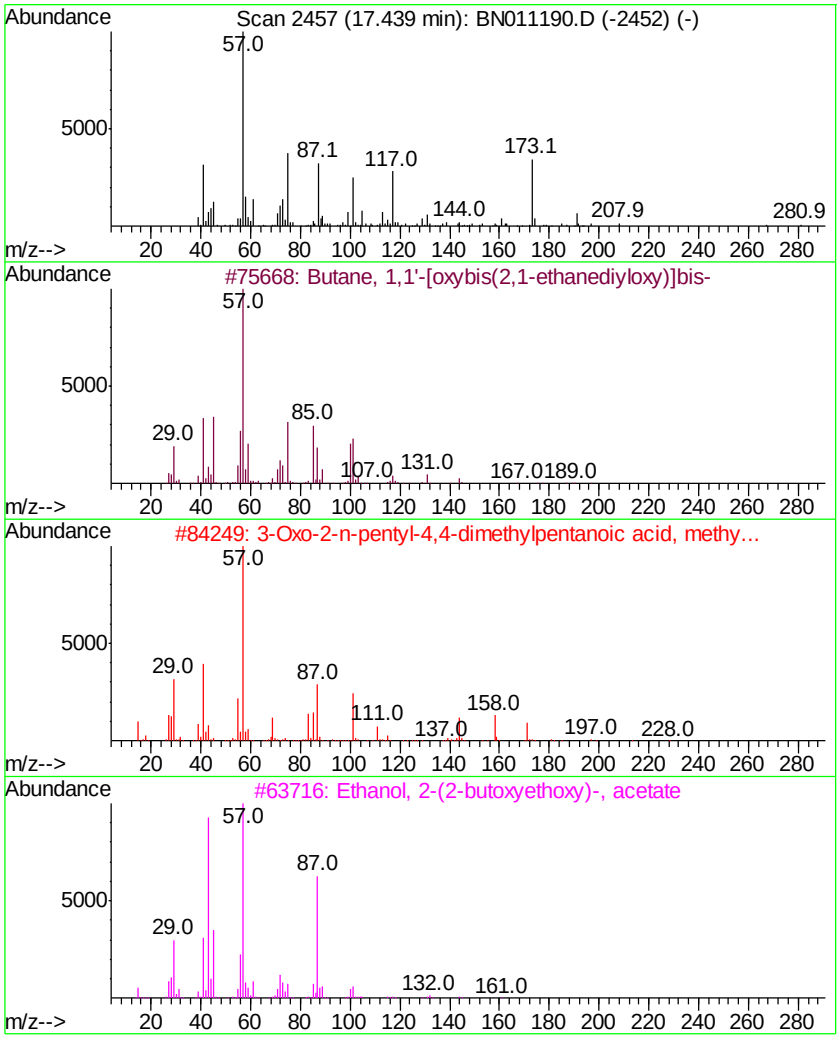
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

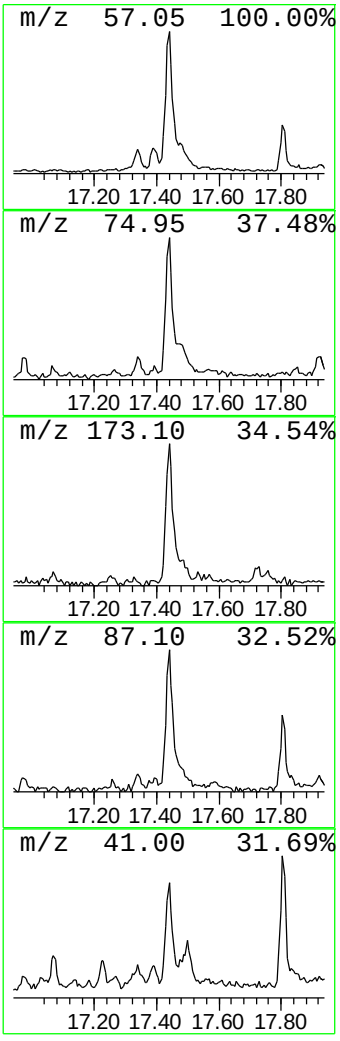
**********

Peak Number 33 unknown-07 Concentration Rank 42

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 17.44 | 2.31 ng/ul | 413145 | Phenanthrene-d10 | 16.87 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Butane, 1,1'-[oxybis(2,1-ethaned... | 218 | C12H26O3 | 000112-73-2 | 37 |
| 2 | 3-Oxo-2-n-pentyl-4,4-dimethylpen... | 228 | C13H24O3 | 1000202-26-0 | 27 |
| 3 | Ethanol, 2-(2-butoxyethoxy)-, ac... | 204 | C10H20O4 | 000124-17-4 | 25 |
| 4 | p-Dioxane-2,5-dimethanol | 148 | C6H12O4 | 014236-12-5 | 25 |
| 5 | Ethenol, 1-[bis(1,1-dimethylethy... | 230 | C12H23O2P | 050838-15-8 | 25 |



# Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN071620\
Data File : BN011190.D
Acq On : 16 Jul 2020 15:23
Operator : CG/JU
Sample : L3261-07
Misc :
ALS Vial : 6 Sample Multiplier: 1

**Instrument :** BNA_N
**ClientSampleId :** BE506

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN071520MA.M
Quant Title : SVOA CALIBRATION

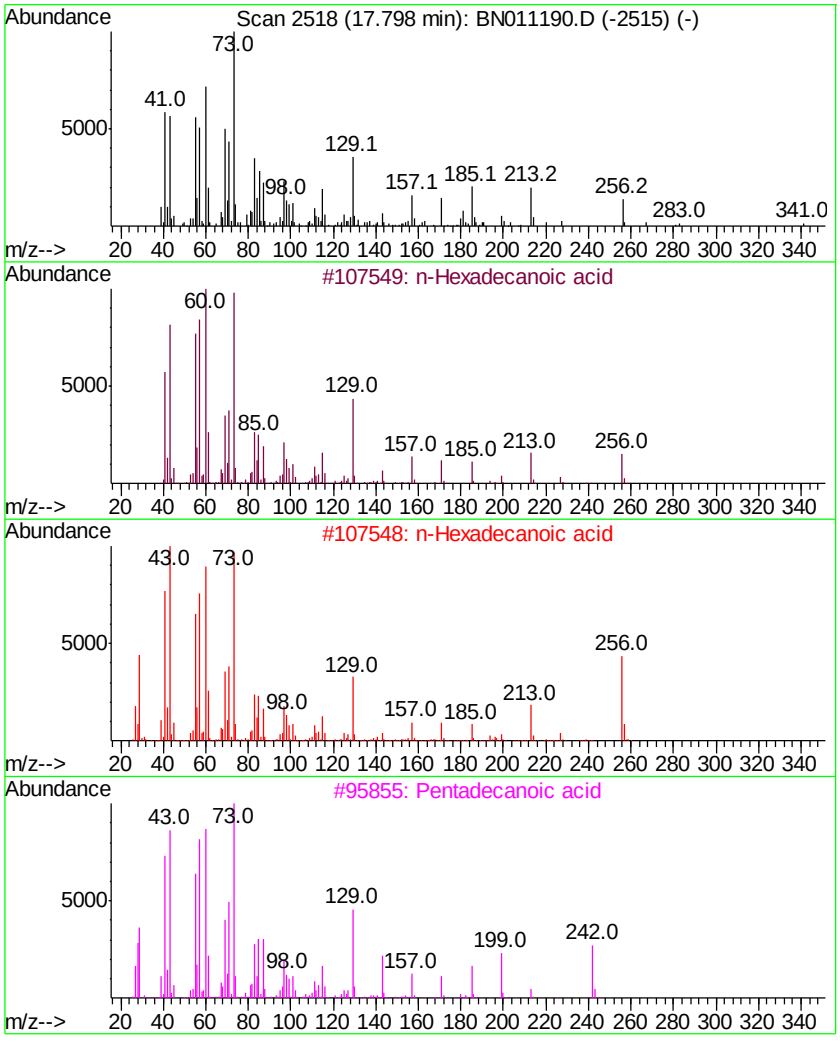
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

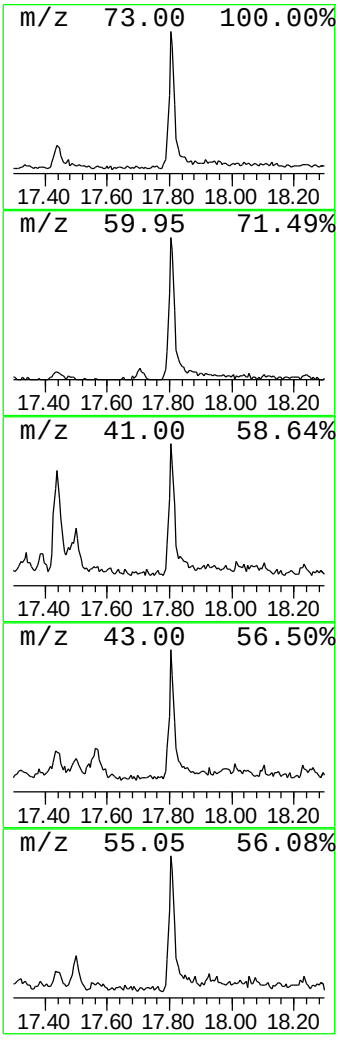
**********

Peak Number 34 n-Hexadecanoic acid Concentration Rank 37

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 17.80 | 2.83 ng/ul | 506159 | Phenanthrene-d10 | 16.87 |

| 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 | 95 |
| 3 | Pentadecanoic acid | 242 | C15H30O2 | 001002-84-2 | 83 |
| 4 | Tridecanoic acid | 214 | C13H26O2 | 000638-53-9 | 64 |
| 5 | Octadecanoic acid | 284 | C18H36O2 | 000057-11-4 | 52 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN071620\
Data File : BN011190.D
Acq On : 16 Jul 2020 15:23
Operator : CG/JU
Sample : L3261-07
Misc :
ALS Vial : 6 Sample Multiplier: 1

**Instrument :** BNA_N
**ClientSampleId :** BE506

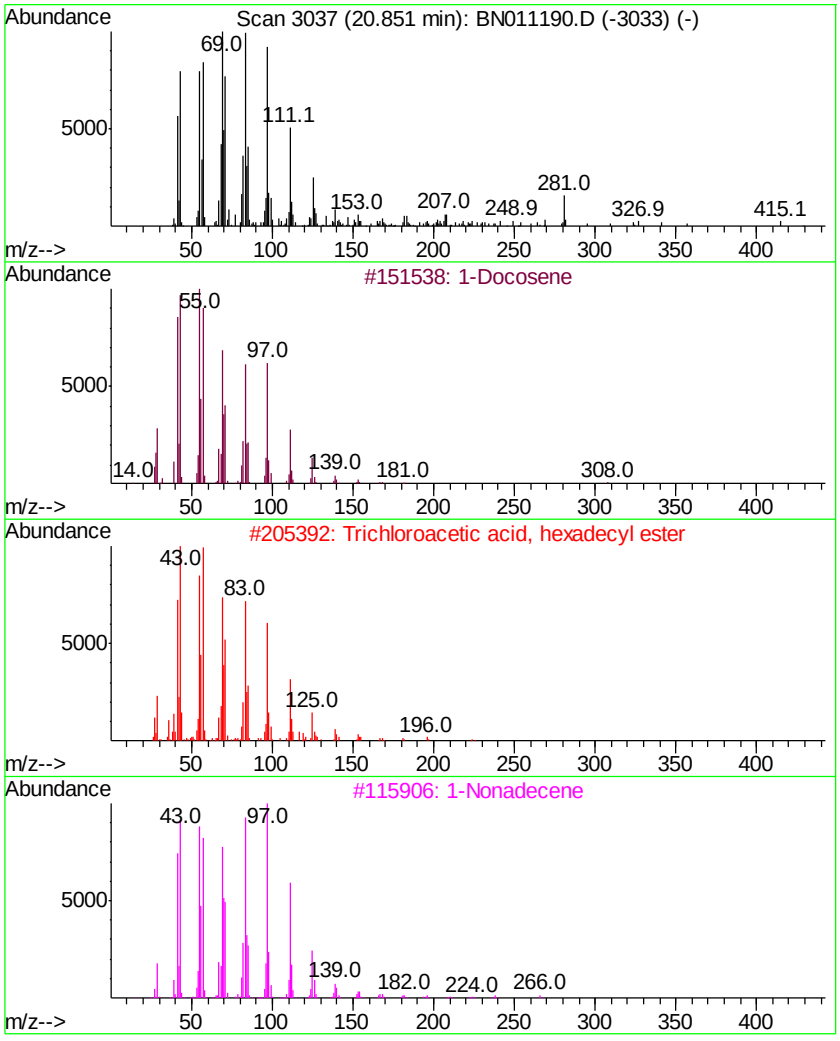
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN071520MA.M
Quant Title : SVOA CALIBRATION

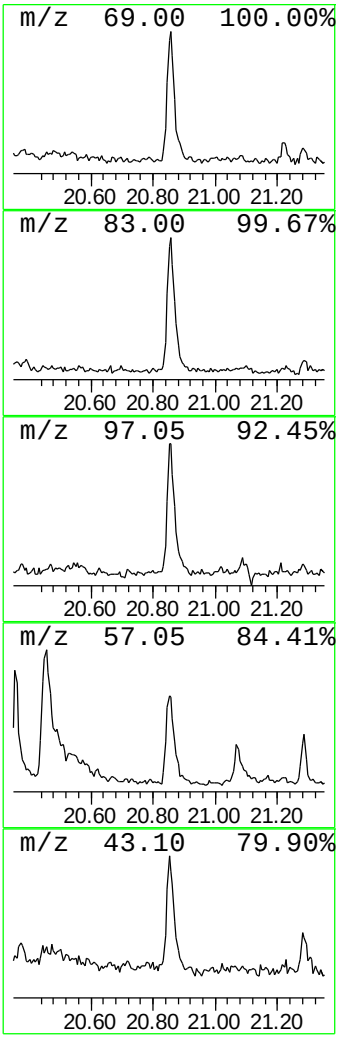
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 41 (DEL) Alkane: Straight-Chai... Concentration Rank 35

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 20.85 | 3.02 ng/ul | 539983 | Chrysene-d12 | 21.09 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1-Docosene | 308 | C22H44 | 001599-67-3 | 81 |
| 2 | Trichloroacetic acid, hexadecyl ... | 386 | C18H33Cl3O2 | 074339-54-1 | 76 |
| 3 | 1-Nonadecene | 266 | C19H38 | 018435-45-5 | 64 |
| 4 | Carbonic acid, hexadecyl isobutyl... | 342 | C21H42O3 | 1000314-61-3 | 62 |
| 5 | Tricosyl trifluoroacetate | 436 | C25H47F3O2 | 1000351-75-1 | 62 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN071620\
Data File : BN011190.D
Acq On : 16 Jul 2020 15:23
Operator : CG/JU
Sample : L3261-07
Misc :
ALS Vial : 6 Sample Multiplier: 1

**Instrument :** BNA_N
**ClientSampleId :** BE506

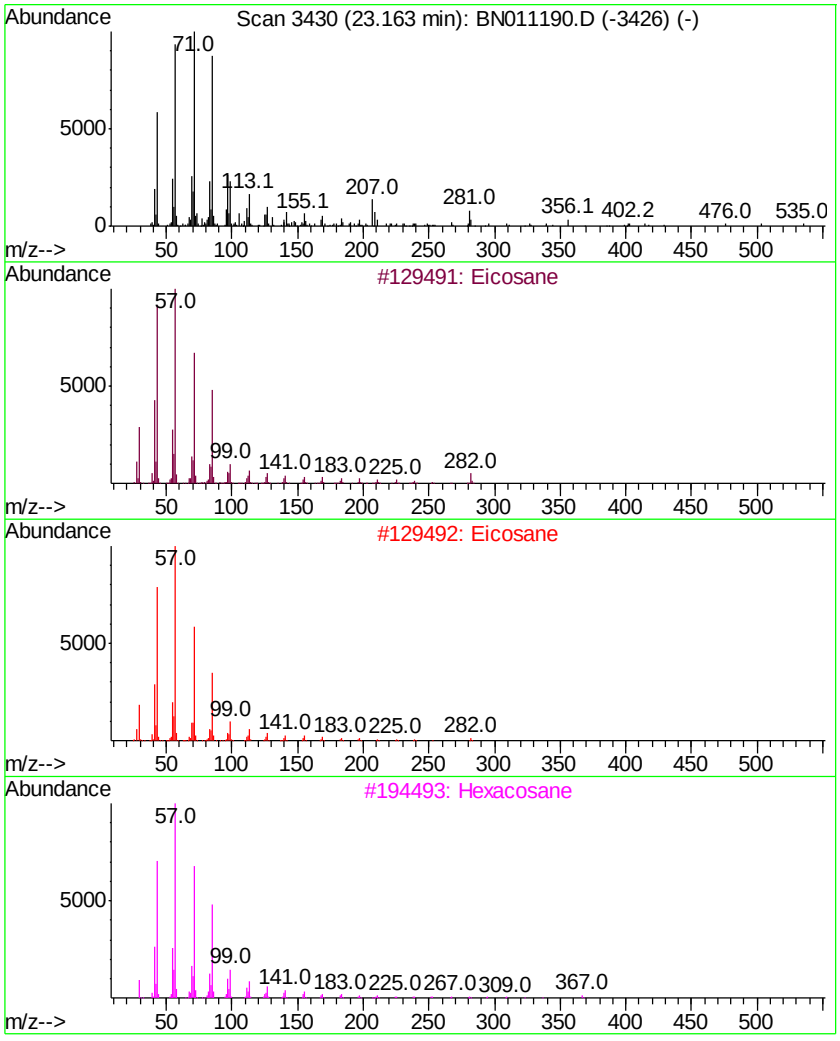
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN071520MA.M
Quant Title : SVOA CALIBRATION

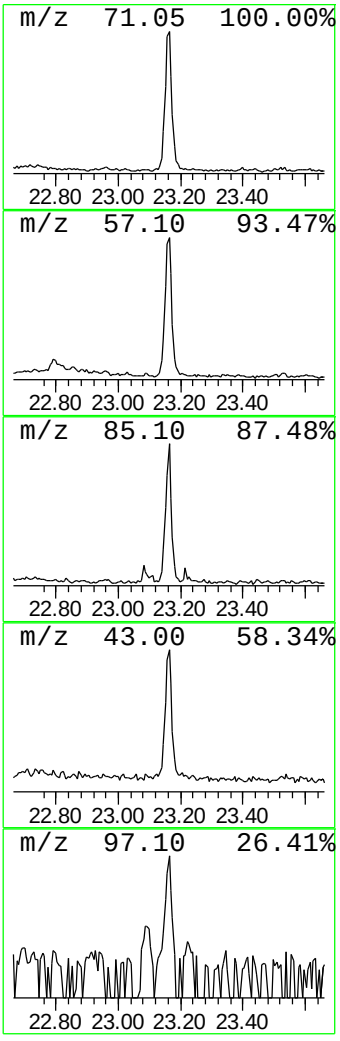
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 43 (DEL) Alkane: Straight-Chai... Concentration Rank 41

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 23.16 | 2.49 ng/ul | 432150 | Perylene-d12 | 23.22 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Eicosane | 282 | C20H42 | 000112-95-8 | 87 |
| 2 | Eicosane | 282 | C20H42 | 000112-95-8 | 87 |
| 3 | Hexacosane | 366 | C26H54 | 000630-01-3 | 83 |
| 4 | 2-methyloctacosane | 408 | C29H60 | 1000376-72-8 | 83 |
| 5 | Octacosane | 394 | C28H58 | 000630-02-4 | 83 |



Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN071620\
Data File : BN011190.D
Acq On : 16 Jul 2020 15:23
Operator : CG/JU
Sample : L3261-07
Misc :
ALS Vial : 6 Sample Multiplier: 1

**Instrument :** BNA_N
**ClientSampleId :** BE506

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN071520MA.M
Quant Title : SVOA CALIBRATION

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

| | | | | | Internal Standard | | | |
|---|---|---|---|---|---|---|---|---|
| TIC Top Hit name | RT | EstConc | Units | Response | # | RT | Resp | Conc |
| Benzene, 1,3-dime... | 5.20 | 175.4 | ng/ul | 8801190 | 1 | 7.49 | 1003790 | 20.0 |
| Benzene, (1-methy... | 6.02 | 13.6 | ng/ul | 683615 | 1 | 7.49 | 1003790 | 20.0 |
| 3-Heptanone, 5-me... | 6.20 | 3.8 | ng/ul | 190198 | 1 | 7.49 | 1003790 | 20.0 |
| Benzene, propyl- | 6.49 | 7.1 | ng/ul | 358328 | 1 | 7.49 | 1003790 | 20.0 |
| Benzene, 1-ethyl-... | 6.60 | 8.1 | ng/ul | 404828 | 1 | 7.49 | 1003790 | 20.0 |
| Benzene, 1,2,3-tr... | 7.15 | 27.9 | ng/ul | 1400580 | 1 | 7.49 | 1003790 | 20.0 |
| Naphthalene, 1,1'... | 7.22 | 3.9 | ng/ul | 196833 | 1 | 7.49 | 1003790 | 20.0 |
| Mesitylene | 7.60 | 10.0 | ng/ul | 502561 | 1 | 7.49 | 1003790 | 20.0 |
| unknown-01 | 7.70 | 4.5 | ng/ul | 224343 | 1 | 7.49 | 1003790 | 20.0 |
| Benzene, cyclopro... | 7.85 | 18.4 | ng/ul | 922699 | 1 | 7.49 | 1003790 | 20.0 |
| Cyclohexanol, 3,3... | 8.11 | 27.2 | ng/ul | 1365430 | 1 | 7.49 | 1003790 | 20.0 |
| unknown-02 | 8.70 | 4.3 | ng/ul | 216015 | 1 | 7.49 | 1003790 | 20.0 |
| 3-Octanol, 3,6-di... | 8.80 | 9.6 | ng/ul | 479550 | 1 | 7.49 | 1003790 | 20.0 |
| 2,4-Heptadienal, ... | 9.27 | 3.2 | ng/ul | 230695 | 2 | 10.23 | 1425010 | 20.0 |
| unknown-03 | 9.66 | 6.8 | ng/ul | 486626 | 2 | 10.23 | 1425010 | 20.0 |
| Bicyclo[2.2.2]octane | 10.65 | 8.9 | ng/ul | 631752 | 2 | 10.23 | 1425010 | 20.0 |
| Cyclohexanone, 4-... | 10.84 | 10.2 | ng/ul | 722907 | 2 | 10.23 | 1425010 | 20.0 |
| unknown-04 | 11.25 | 3.2 | ng/ul | 228846 | 2 | 10.23 | 1425010 | 20.0 |
| Phenol, 2,4,6-tri... | 11.33 | 7.2 | ng/ul | 511414 | 2 | 10.23 | 1425010 | 20.0 |
| Phenol, p-tert-bu... | 11.60 | 17.5 | ng/ul | 1248910 | 2 | 10.23 | 1425010 | 20.0 |
| unknown-05 | 12.08 | 2.6 | ng/ul | 184759 | 2 | 10.23 | 1425010 | 20.0 |
| Naphthalene, 1-me... | 12.13 | 5.3 | ng/ul | 376934 | 2 | 10.23 | 1425010 | 20.0 |
| Tripropyl phosphate | 12.67 | 2.9 | ng/ul | 312851 | 3 | 14.12 | 2171550 | 20.0 |
| unknown-06 | 12.77 | 2.1 | ng/ul | 226333 | 3 | 14.12 | 2171550 | 20.0 |
| 2,4,7,9-Tetrameth... | 13.04 | 13.0 | ng/ul | 1411870 | 3 | 14.12 | 2171550 | 20.0 |
| Diaziridinone, (1... | 15.65 | 3.9 | ng/ul | 700547 | 4 | 16.87 | 3583360 | 20.0 |
| 2(3H)-Benzothiazo... | 15.82 | 3.3 | ng/ul | 584216 | 4 | 16.87 | 3583360 | 20.0 |
| Phenol, 4-(1-phen... | 16.19 | 3.4 | ng/ul | 603658 | 4 | 16.87 | 3583360 | 20.0 |
| unknown-07 | 17.44 | 2.3 | ng/ul | 413145 | 4 | 16.87 | 3583360 | 20.0 |
| n-Hexadecanoic acid | 17.80 | 2.8 | ng/ul | 506159 | 4 | 16.87 | 3583360 | 20.0 |
| (DEL) Alkane: Str... | 20.85 | 3.0 | ng/ul | 539983 | 5 | 21.09 | 3570630 | 20.0 |
| (DEL) Alkane: Str... | 23.16 | 2.5 | ng/ul | 432150 | 6 | 23.22 | 3470290 | 20.0 |