LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072324\
Data File : BG062199.D
Acq On : 24 Jul 2024 00:03
Operator : MA/JU
Sample : P3244-04
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
YDZD5

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_G\Methods\SFAM-EPA-BG071924.MA.M
Title : SVOA CALIBRATION

Signal : TIC: BG062199.D\data.ms

| peak # | R.T. min | first scan | max scan | last scan | PK TY | peak height | corr. area | corr. % max. | % of total |
|---|---|---|---|---|---|---|---|---|---|
| 1 | 4.217 | 151 | 158 | 175 | rBV | 682437 | 1152469 | 19.87% | 1.427% |
| 2 | 5.556 | 379 | 386 | 392 | rVV | 94489 | 165436 | 2.85% | 0.205% |
| 3 | 5.685 | 401 | 408 | 423 | rBV | 276461 | 604023 | 10.41% | 0.748% |
| 4 | 5.944 | 444 | 452 | 459 | rBV2 | 131251 | 234556 | 4.04% | 0.290% |
| 5 | 6.061 | 465 | 472 | 480 | rBV | 292806 | 520786 | 8.98% | 0.645% |
| 6 | 7.142 | 650 | 656 | 661 | rBV | 242335 | 384229 | 6.62% | 0.476% |
| 7 | 7.201 | 661 | 666 | 671 | rVB2 | 138323 | 223763 | 3.86% | 0.277% |
| 8 | 7.277 | 671 | 679 | 685 | rBV2 | 202077 | 374791 | 6.46% | 0.464% |
| 9 | 7.447 | 703 | 708 | 714 | rVB | 173185 | 293689 | 5.06% | 0.364% |
| 10 | 7.612 | 729 | 736 | 740 | rBV5 | 106987 | 218084 | 3.76% | 0.270% |
| 11 | 7.706 | 746 | 752 | 759 | rBV | 603513 | 985378 | 16.99% | 1.220% |
| 12 | 8.023 | 800 | 806 | 810 | rVV | 363189 | 603732 | 10.41% | 0.747% |
| 13 | 8.070 | 810 | 814 | 826 | rVB2 | 176763 | 375813 | 6.48% | 0.465% |
| 14 | 8.176 | 826 | 832 | 844 | rVB3 | 405892 | 880440 | 15.18% | 1.090% |
| 15 | 8.364 | 853 | 864 | 871 | rBV2 | 370705 | 701088 | 12.09% | 0.868% |
| 16 | 8.434 | 871 | 876 | 883 | rBV4 | 100157 | 195204 | 3.36% | 0.242% |
| 17 | 8.534 | 889 | 893 | 900 | rVB4 | 148396 | 271205 | 4.68% | 0.336% |
| 18 | 8.693 | 914 | 920 | 933 | rBV3 | 256561 | 589494 | 10.16% | 0.730% |
| 19 | 8.804 | 933 | 939 | 943 | rBV2 | 95543 | 166500 | 2.87% | 0.206% |
| 20 | 8.863 | 943 | 949 | 957 | rVB4 | 85211 | 180470 | 3.11% | 0.223% |
| 21 | 9.063 | 979 | 983 | 989 | rVB | 99471 | 156972 | 2.71% | 0.194% |
| 22 | 9.169 | 989 | 1001 | 1008 | rVB4 | 149322 | 341592 | 5.89% | 0.423% |
| 23 | 9.245 | 1008 | 1014 | 1019 | rBV4 | 208990 | 416037 | 7.17% | 0.515% |
| 24 | 9.310 | 1019 | 1025 | 1035 | rVB | 761838 | 1320988 | 22.77% | 1.635% |
| 25 | 9.503 | 1048 | 1058 | 1063 | rBV3 | 218208 | 455848 | 7.86% | 0.564% |
| 26 | 9.750 | 1096 | 1100 | 1103 | rVV | 105278 | 164376 | 2.83% | 0.203% |
| 27 | 9.791 | 1103 | 1107 | 1116 | rVB4 | 157185 | 319240 | 5.50% | 0.395% |
| 28 | 9.891 | 1116 | 1124 | 1130 | rBV6 | 60027 | 178800 | 3.08% | 0.221% |
| 29 | 9.962 | 1130 | 1136 | 1145 | rVB7 | 115361 | 271851 | 4.69% | 0.336% |
| 30 | 10.067 | 1147 | 1154 | 1156 | rBV3 | 299186 | 509723 | 8.79% | 0.631% |
| 31 | 10.103 | 1156 | 1160 | 1166 | rVB2 | 547609 | 915340 | 15.78% | 1.133% |
| 32 | 10.250 | 1177 | 1185 | 1189 | rBV | 2190751 | 4040469 | 69.65% | 5.001% |
| 33 | 10.297 | 1189 | 1193 | 1201 | rVB3 | 1185464 | 2129347 | 36.71% | 2.636% |
| 34 | 10.379 | 1201 | 1207 | 1211 | rBV | 542087 | 878906 | 15.15% | 1.088% |
| 35 | 10.526 | 1221 | 1232 | 1238 | rBV5 | 477698 | 1205906 | 20.79% | 1.493% |
| 36 | 10.655 | 1248 | 1254 | 1262 | rBV2 | 143742 | 316963 | 5.46% | 0.392% |

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072324\
Data File : BG062199.D
Acq On : 24 Jul 2024 00:03
Operator : MA/JU
Sample : P3244-04
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
YDZD5

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_G\Methods\SFAM-EPA-BG071924.MA.M
Title : SVOA CALIBRATION

| | | | | | | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 37 | 10.766 | 1266 | 1273 | 1279 | rVV | 266804 | 512368 | 8.83% | 0.634% |
| 38 | 10.884 | 1283 | 1293 | 1297 | rVV | 302042 | 712064 | 12.27% | 0.881% |
| 39 | 10.937 | 1297 | 1302 | 1307 | rVV | 474194 | 829261 | 14.29% | 1.026% |
| 40 | 11.031 | 1313 | 1318 | 1324 | rVV5 | 186220 | 390165 | 6.73% | 0.483% |
| 41 | 11.242 | 1349 | 1354 | 1360 | rBV3 | 341219 | 571593 | 9.85% | 0.708% |
| 42 | 11.424 | 1376 | 1385 | 1389 | rBV3 | 85793 | 216889 | 3.74% | 0.268% |
| 43 | 11.724 | 1431 | 1436 | 1441 | rBV2 | 250318 | 397218 | 6.85% | 0.492% |
| 44 | 11.912 | 1463 | 1468 | 1473 | rBV2 | 185464 | 262660 | 4.53% | 0.325% |
| 45 | 11.971 | 1473 | 1478 | 1481 | rBV3 | 213800 | 373693 | 6.44% | 0.463% |
| 46 | 12.059 | 1489 | 1493 | 1499 | rVB5 | 89982 | 159850 | 2.76% | 0.198% |
| 47 | 12.188 | 1510 | 1515 | 1517 | rBV5 | 127688 | 206414 | 3.56% | 0.255% |
| 48 | 12.229 | 1517 | 1522 | 1530 | rVB | 610502 | 1017569 | 17.54% | 1.260% |
| 49 | 12.394 | 1543 | 1550 | 1556 | rBV9 | 66930 | 197781 | 3.41% | 0.245% |
| 50 | 12.482 | 1561 | 1565 | 1568 | rVV4 | 194409 | 295097 | 5.09% | 0.365% |
| 51 | 12.541 | 1568 | 1575 | 1580 | rVV | 3341915 | 5801099 | 100.00% | 7.180% |
| 52 | 12.588 | 1580 | 1583 | 1590 | rVB3 | 184403 | 329205 | 5.67% | 0.407% |
| 53 | 12.758 | 1604 | 1612 | 1620 | rVB | 2716487 | 4610962 | 79.48% | 5.707% |
| 54 | 12.852 | 1621 | 1628 | 1633 | rBV7 | 88332 | 218248 | 3.76% | 0.270% |
| 55 | 12.911 | 1633 | 1638 | 1640 | rBV2 | 95749 | 166017 | 2.86% | 0.205% |
| 56 | 13.052 | 1657 | 1662 | 1669 | rVV8 | 92150 | 215381 | 3.71% | 0.267% |
| 57 | 13.122 | 1670 | 1674 | 1681 | rVB2 | 179135 | 317562 | 5.47% | 0.393% |
| 58 | 13.292 | 1699 | 1703 | 1712 | rVB2 | 153933 | 248628 | 4.29% | 0.308% |
| 59 | 13.510 | 1736 | 1740 | 1744 | rBV5 | 113790 | 220543 | 3.80% | 0.273% |
| 60 | 13.727 | 1773 | 1777 | 1787 | rVB | 191498 | 454756 | 7.84% | 0.563% |
| 61 | 13.856 | 1791 | 1799 | 1800 | rBV5 | 390108 | 614866 | 10.60% | 0.761% |
| 62 | 14.021 | 1821 | 1827 | 1832 | rVV2 | 506525 | 1002125 | 17.27% | 1.240% |
| 63 | 14.074 | 1832 | 1836 | 1838 | rVV2 | 388254 | 592269 | 10.21% | 0.733% |
| 64 | 14.097 | 1838 | 1840 | 1846 | rVB | 428557 | 520686 | 8.98% | 0.644% |
| 65 | 14.256 | 1862 | 1867 | 1870 | rBV | 261676 | 403020 | 6.95% | 0.499% |
| 66 | 14.397 | 1886 | 1891 | 1902 | rVB3 | 418440 | 855548 | 14.75% | 1.059% |
| 67 | 14.555 | 1914 | 1918 | 1923 | rVB2 | 185186 | 288733 | 4.98% | 0.357% |
| 68 | 14.649 | 1930 | 1934 | 1937 | rBV2 | 120516 | 169006 | 2.91% | 0.209% |
| 69 | 14.696 | 1937 | 1942 | 1948 | rVB | 481750 | 730978 | 12.60% | 0.905% |
| 70 | 15.454 | 2065 | 2071 | 2078 | rVV | 820970 | 1304797 | 22.49% | 1.615% |
| 71 | 15.525 | 2079 | 2083 | 2089 | rVB | 325051 | 502716 | 8.67% | 0.622% |
| 72 | 15.689 | 2106 | 2111 | 2117 | rVB2 | 1221332 | 1793558 | 30.92% | 2.220% |
| 73 | 16.012 | 2162 | 2166 | 2172 | rVB4 | 86109 | 164714 | 2.84% | 0.204% |
| 74 | 16.124 | 2181 | 2185 | 2187 | rVV2 | 179251 | 256173 | 4.42% | 0.317% |
| 75 | 16.159 | 2187 | 2191 | 2195 | rVB2 | 176008 | 249357 | 4.30% | 0.309% |
| 76 | 16.424 | 2230 | 2236 | 2245 | rVV | 774990 | 1146580 | 19.76% | 1.419% |

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072324\
Data File : BG062199.D
Acq On : 24 Jul 2024 00:03
Operator : MA/JU
Sample : P3244-04
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
YDZD5

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_G\Methods\SFAM-EPA-BG071924.MA.M
Title : SVOA CALIBRATION

| | | | | | | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 77 | 16.518 | 2246 | 2252 | 2256 | rVV | 1634510 | 2343181 | 40.39% | 2.900% |
| 78 | 16.576 | 2256 | 2262 | 2268 | rVV3 | 2444629 | 4666093 | 80.43% | 5.776% |
| 79 | 16.635 | 2268 | 2272 | 2277 | rVV2 | 2124051 | 3216084 | 55.44% | 3.981% |
| 80 | 16.682 | 2277 | 2280 | 2284 | rVV | 1150155 | 1476903 | 25.46% | 1.828% |
| 81 | 16.729 | 2284 | 2288 | 2291 | rVV | 583257 | 1001997 | 17.27% | 1.240% |
| 82 | 16.758 | 2291 | 2293 | 2295 | rVV | 608857 | 695686 | 11.99% | 0.861% |
| 83 | 16.782 | 2295 | 2297 | 2301 | rVV | 492252 | 617241 | 10.64% | 0.764% |
| 84 | 16.841 | 2301 | 2307 | 2313 | rVV3 | 1422506 | 2674558 | 46.10% | 3.311% |
| 85 | 16.905 | 2313 | 2318 | 2321 | rVV | 1435179 | 2089020 | 36.01% | 2.586% |
| 86 | 16.940 | 2321 | 2324 | 2327 | rVV2 | 710608 | 1116490 | 19.25% | 1.382% |
| 87 | 16.976 | 2327 | 2330 | 2336 | rVV2 | 903513 | 1402988 | 24.18% | 1.737% |
| 88 | 17.034 | 2337 | 2340 | 2347 | rVV7 | 105861 | 198316 | 3.42% | 0.245% |
| 89 | 17.111 | 2348 | 2353 | 2361 | rVB4 | 71940 | 184773 | 3.19% | 0.229% |
| 90 | 17.246 | 2373 | 2376 | 2383 | rVB | 202169 | 299578 | 5.16% | 0.371% |
| 91 | 17.446 | 2405 | 2410 | 2415 | rVB | 556084 | 785901 | 13.55% | 0.973% |
| 92 | 17.540 | 2421 | 2426 | 2432 | rBV | 538192 | 838952 | 14.46% | 1.038% |
| 93 | 17.886 | 2480 | 2485 | 2492 | rBV3 | 201862 | 309710 | 5.34% | 0.383% |
| 94 | 19.490 | 2755 | 2758 | 2765 | rVB4 | 148804 | 220715 | 3.80% | 0.273% |
| 95 | 19.795 | 2805 | 2810 | 2812 | rBV | 450991 | 619987 | 10.69% | 0.767% |
| 96 | 19.825 | 2812 | 2815 | 2820 | rVB | 748335 | 1038700 | 17.91% | 1.286% |
| 97 | 20.007 | 2842 | 2846 | 2850 | rBV | 123954 | 167823 | 2.89% | 0.208% |
| 98 | 21.705 | 3129 | 3135 | 3141 | rBV | 595608 | 941784 | 16.23% | 1.166% |
| 99 | 24.718 | 3640 | 3648 | 3659 | rVB2 | 323339 | 870095 | 15.00% | 1.077% |
| 100 | 24.941 | 3679 | 3686 | 3700 | rVB2 | 324755 | 949647 | 16.37% | 1.175% |

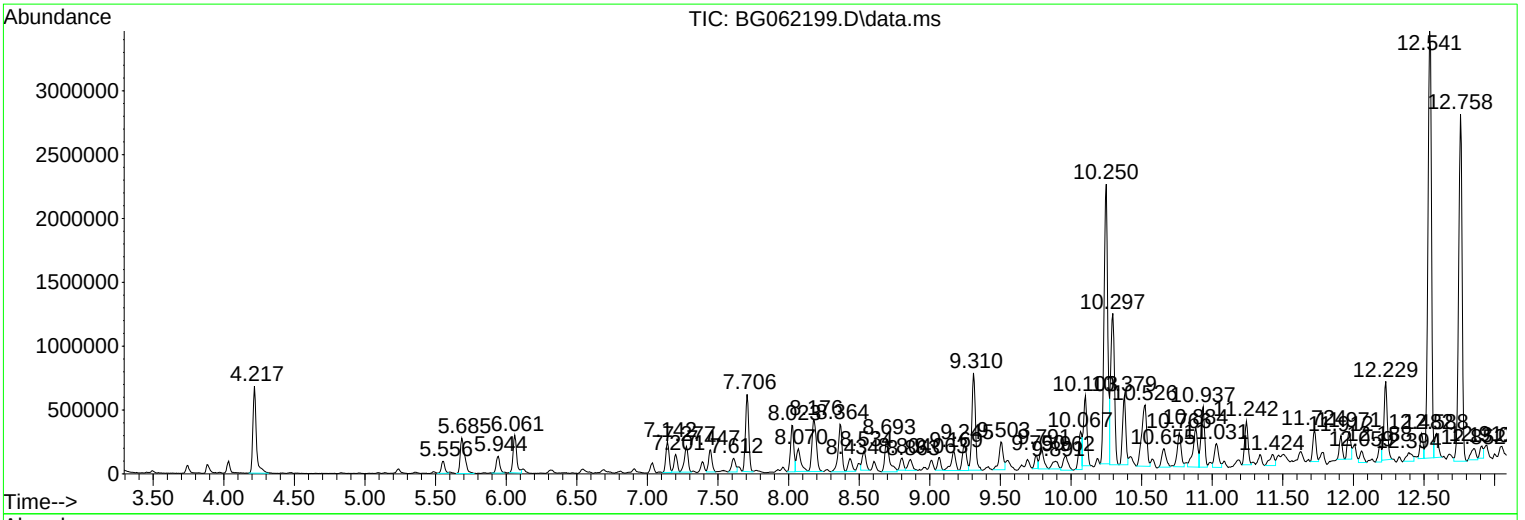
Sum of corrected areas: 80789879

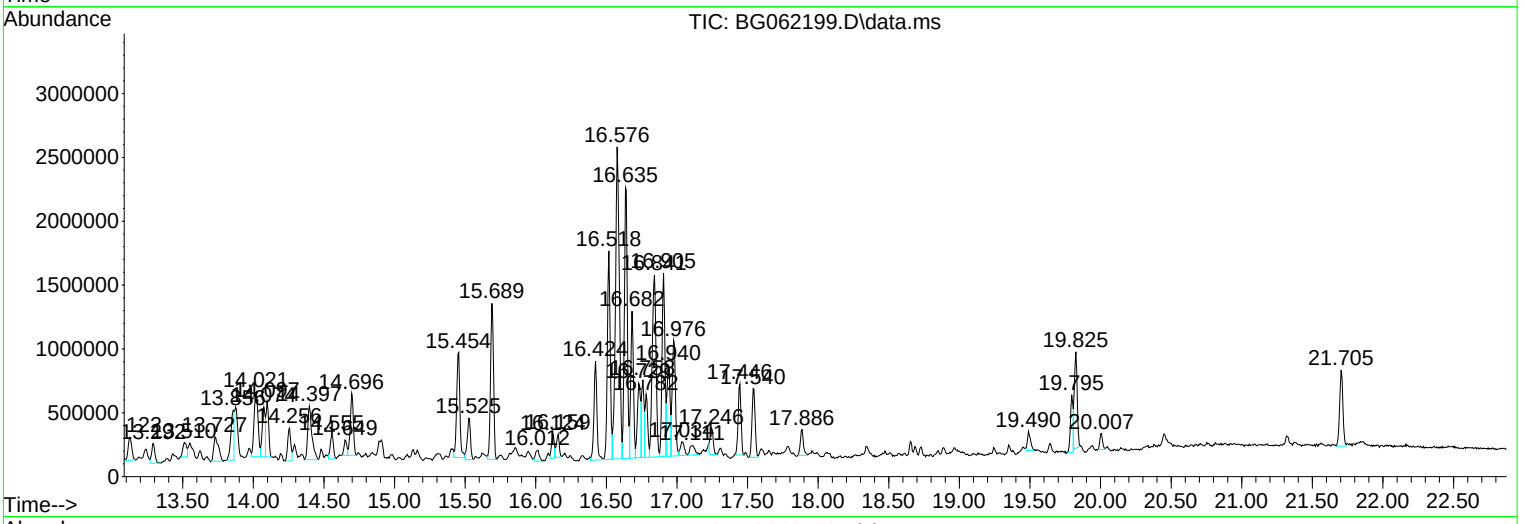
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072324\
Data File : BG062199.D
Acq On : 24 Jul 2024 00:03
Operator : MA/JU
Sample : P3244-04
Misc :
ALS Vial : 19 Sample Multiplier: 1

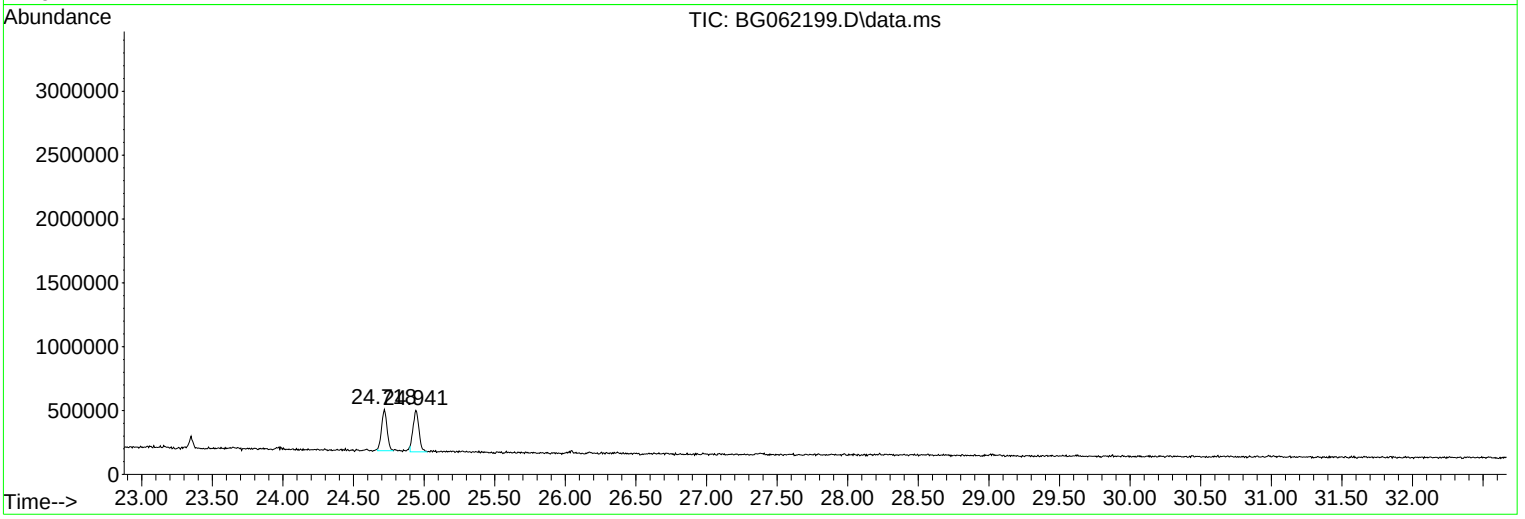
**Instrument :**
BNA_G
**ClientSampleId :**
YDZD5

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072324\
Data File : BG062199.D
Acq On : 24 Jul 2024 00:03
Operator : MA/JU
Sample : P3244-04
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
YDZD5

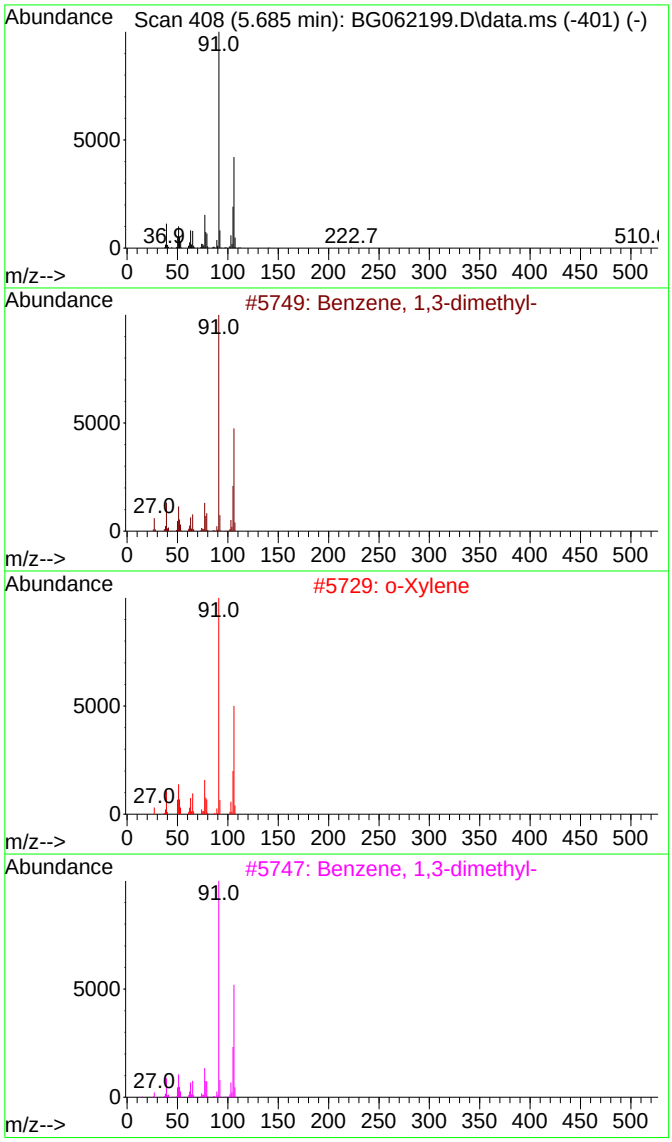
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

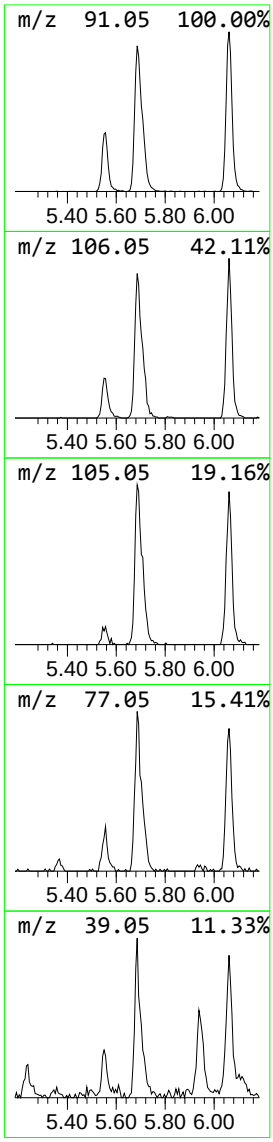
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 3 Benzene, 1,3-dimethyl- Concentration Rank 16

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 5.685 | 32.14 ng/ul | 604023 | 1,4-Dichlorobenzene-d4 | 8.070 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzene, 1,3-dimethyl- | 106 | C8H10 | 000108-38-3 | 97 |
| 2 | o-Xylene | 106 | C8H10 | 000095-47-6 | 95 |
| 3 | Benzene, 1,3-dimethyl- | 106 | C8H10 | 000108-38-3 | 95 |
| 4 | o-Xylene | 106 | C8H10 | 000095-47-6 | 94 |
| 5 | p-Xylene | 106 | C8H10 | 000106-42-3 | 94 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072324\
Data File : BG062199.D
Acq On : 24 Jul 2024 00:03
Operator : MA/JU
Sample : P3244-04
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
YDZD5

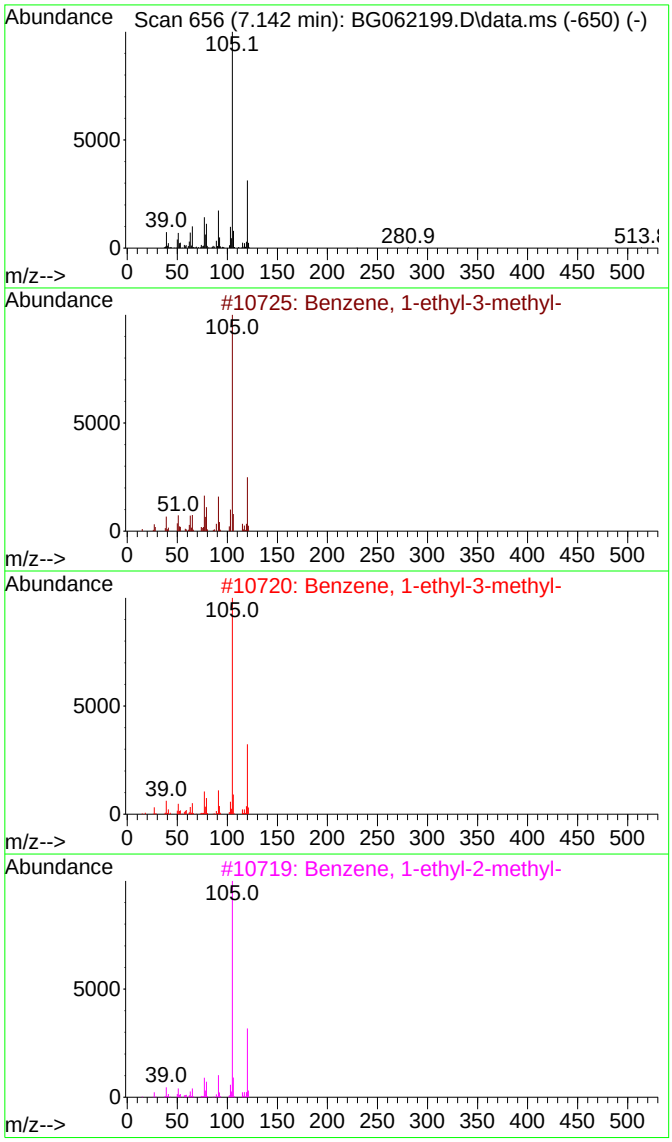
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

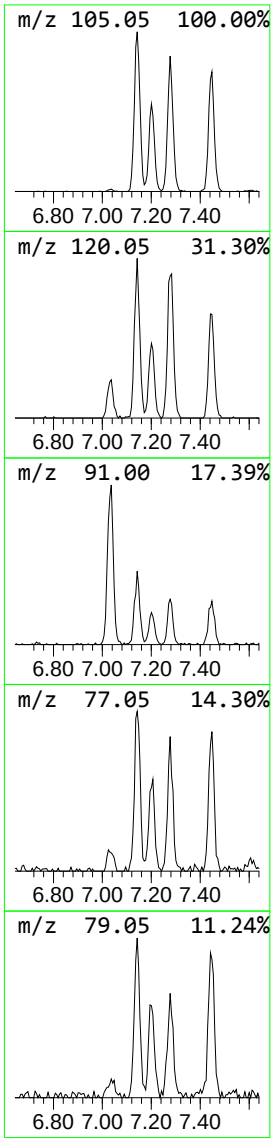
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 6 Benzene, 1-ethyl-3-methyl- Concentration Rank 24

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 7.142 | 20.45 ng/ul | 384229 | 1,4-Dichlorobenzene-d4 | 8.070 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzene, 1-ethyl-3-methyl- | 120 | C9H12 | 000620-14-4 | 93 |
| 2 | Benzene, 1-ethyl-3-methyl- | 120 | C9H12 | 000620-14-4 | 90 |
| 3 | Benzene, 1-ethyl-2-methyl- | 120 | C9H12 | 000611-14-3 | 90 |
| 4 | Mesitylene | 120 | C9H12 | 000108-67-8 | 90 |
| 5 | Benzene, 1-ethyl-2-methyl- | 120 | C9H12 | 000611-14-3 | 90 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072324\
Data File : BG062199.D
Acq On : 24 Jul 2024 00:03
Operator : MA/JU
Sample : P3244-04
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
YDZD5

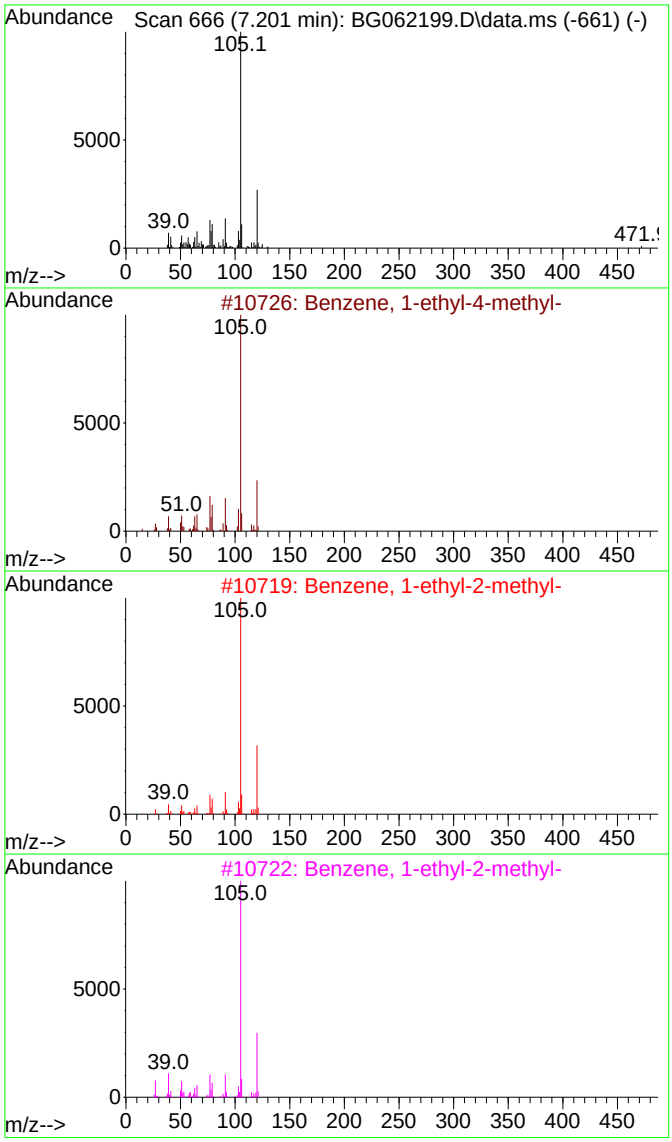
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

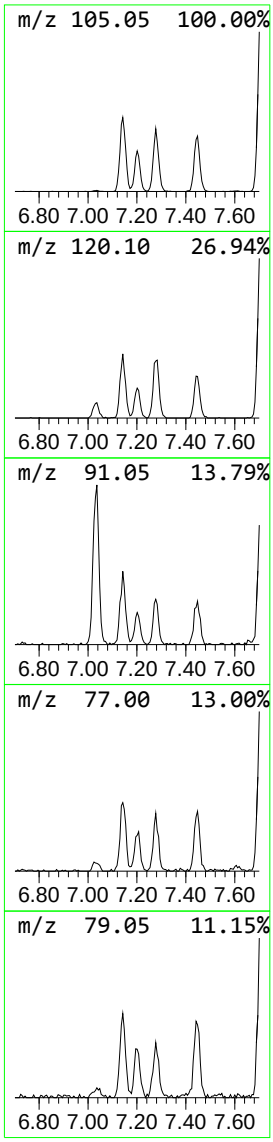
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 7 Benzene, 1-ethyl-4-methyl- Concentration Rank 36

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 7.201 | 11.91 ng/ul | 223763 | 1,4-Dichlorobenzene-d4 | 8.070 |

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072324\
Data File : BG062199.D
Acq On : 24 Jul 2024 00:03
Operator : MA/JU
Sample : P3244-04
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
YDZD5

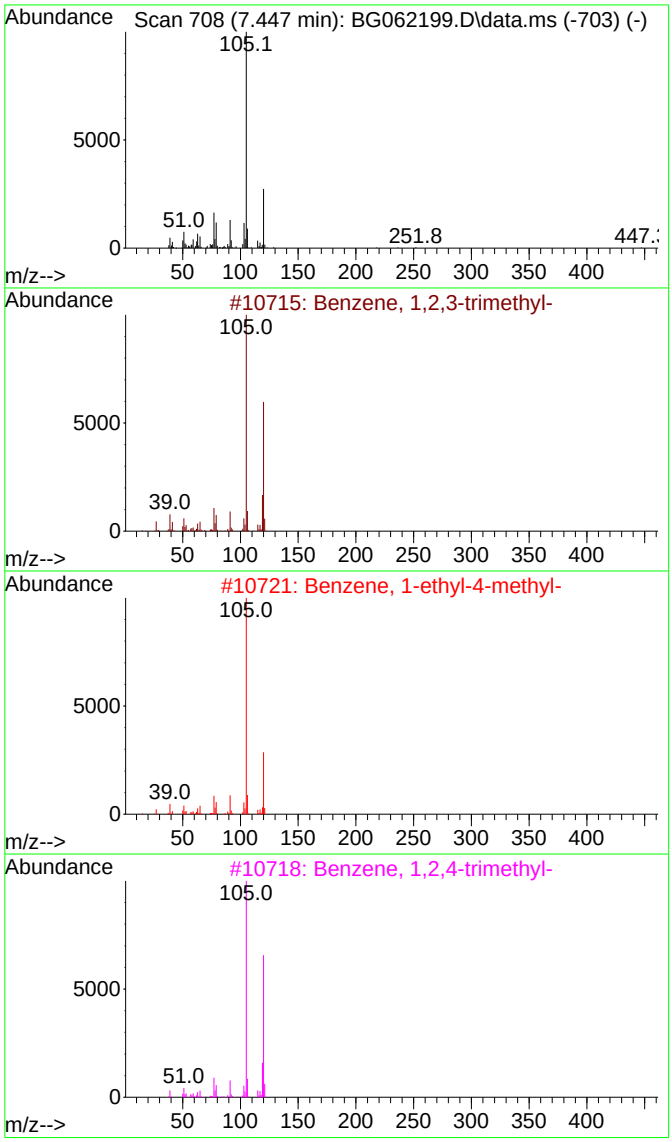
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

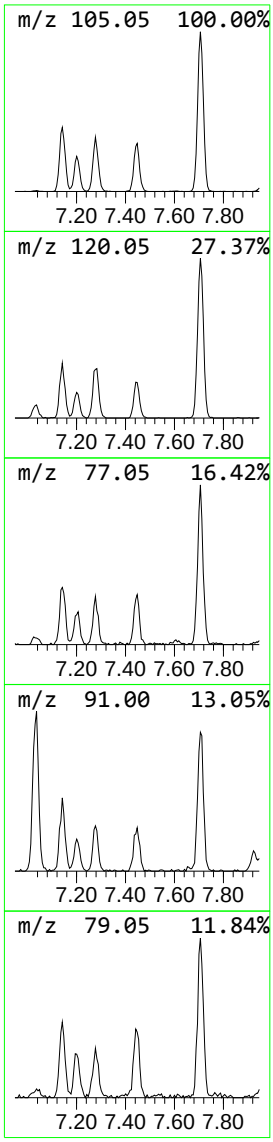
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 8 Benzene, 1,2,3-trimethyl- Concentration Rank 30

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 7.447 | 15.63 ng/ul | 293689 | 1,4-Dichlorobenzene-d4 | 8.070 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzene, 1,2,3-trimethyl- | 120 | C9H12 | 000526-73-8 | 91 |
| 2 | Benzene, 1-ethyl-4-methyl- | 120 | C9H12 | 000622-96-8 | 91 |
| 3 | Benzene, 1,2,4-trimethyl- | 120 | C9H12 | 000095-63-6 | 91 |
| 4 | Benzene, 1-ethyl-3-methyl- | 120 | C9H12 | 000620-14-4 | 90 |
| 5 | Benzene, 1-ethyl-3-methyl- | 120 | C9H12 | 000620-14-4 | 90 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072324\
Data File : BG062199.D
Acq On : 24 Jul 2024 00:03
Operator : MA/JU
Sample : P3244-04
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
YDZD5

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

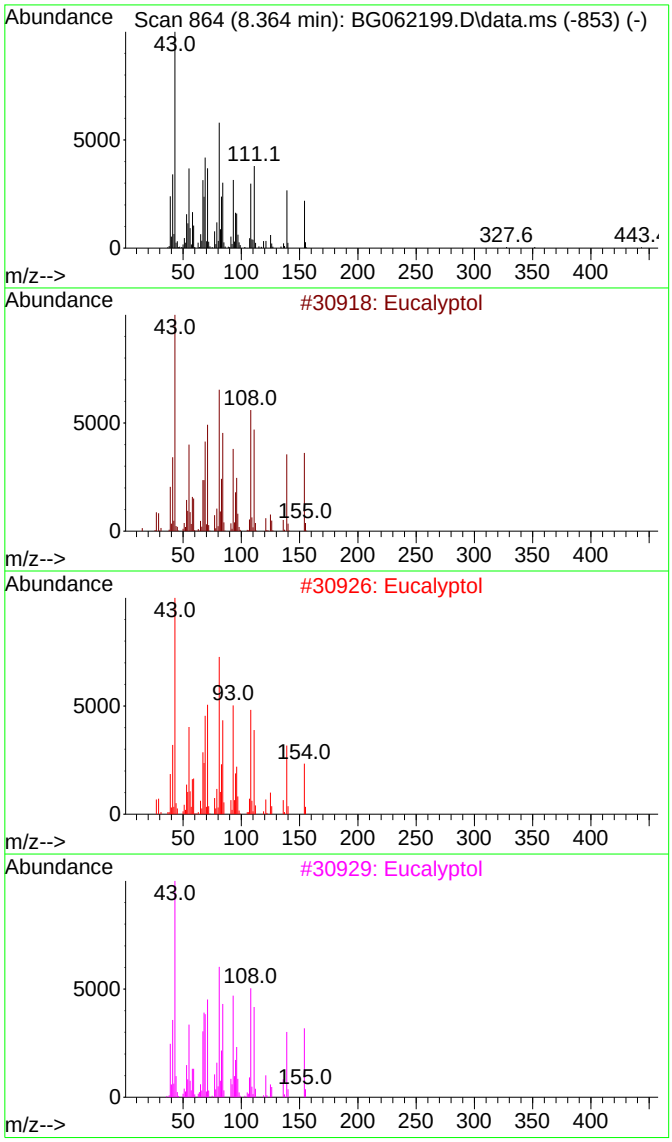
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

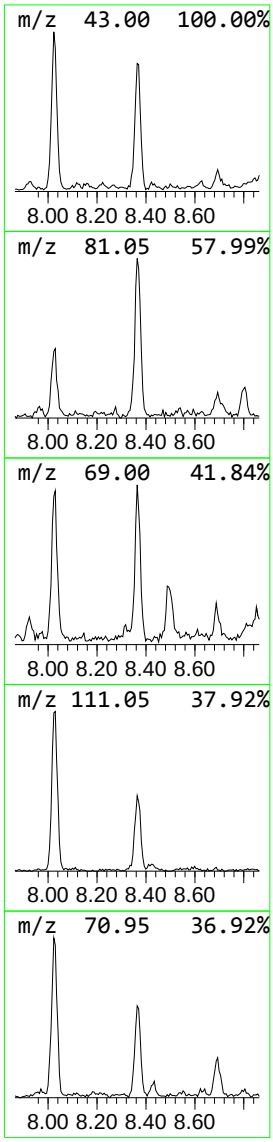
**********

Peak Number 11 Eucalyptol Concentration Rank 13

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 8.364 | 37.31 ng/ul | 701088 | 1,4-Dichlorobenzene-d4 | 8.070 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Eucalyptol | 154 | C10H18O | 000470-82-6 | 91 |
| 2 | Eucalyptol | 154 | C10H18O | 000470-82-6 | 91 |
| 3 | Eucalyptol | 154 | C10H18O | 000470-82-6 | 83 |
| 4 | Eucalyptol | 154 | C10H18O | 000470-82-6 | 81 |
| 5 | Eucalyptol | 154 | C10H18O | 000470-82-6 | 76 |



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Data File : BG062199.D
Acq On : 24 Jul 2024 00:03
Operator : MA/JU
Sample : P3244-04
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
YDZD5

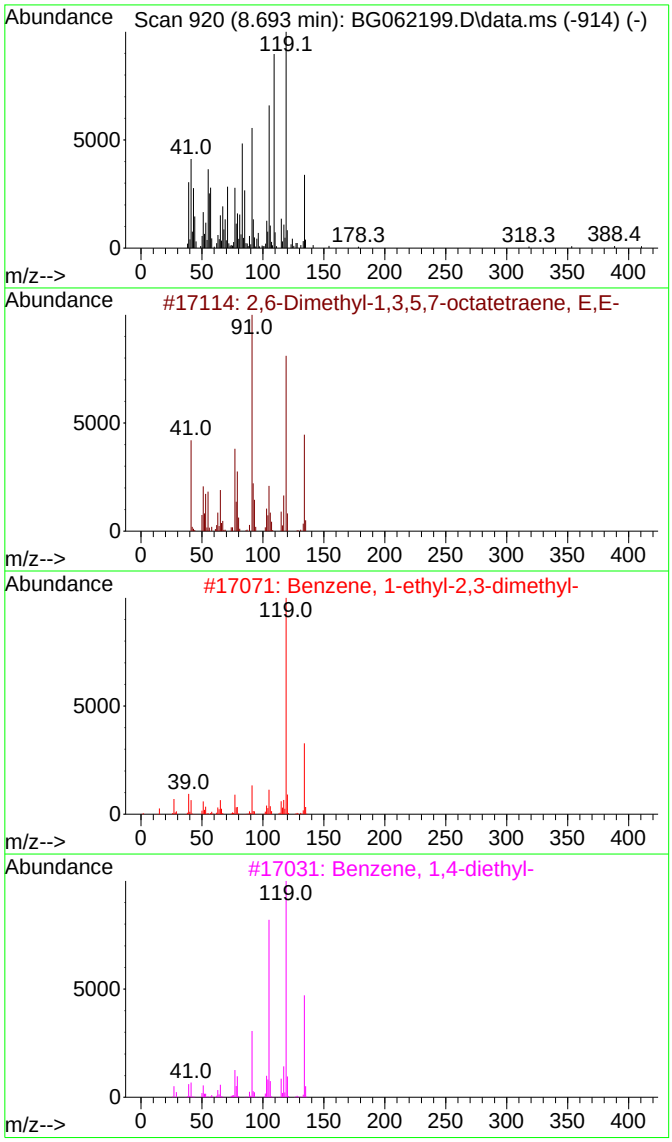
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

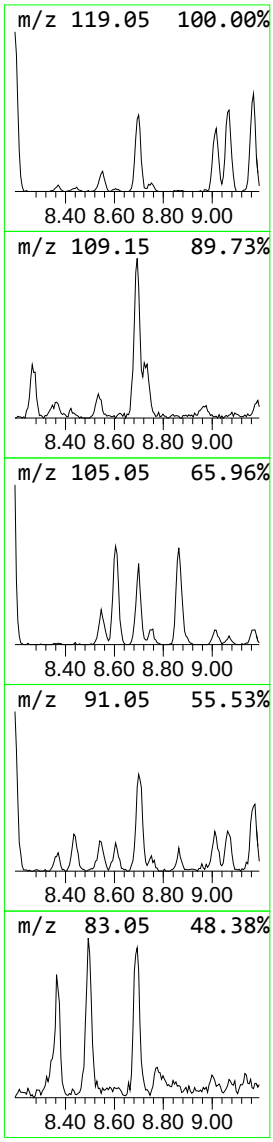
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 13 2,6-Dimethyl-1,3,5,7-octate... Concentration Rank 17

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 8.693 | 31.37 ng/ul | 589494 | 1,4-Dichlorobenzene-d4 | 8.070 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2,6-Dimethyl-1,3,5,7-octatetraen... | 134 | C10H14 | 000460-01-5 | 70 |
| 2 | Benzene, 1-ethyl-2,3-dimethyl- | 134 | C10H14 | 000933-98-2 | 49 |
| 3 | Benzene, 1,4-diethyl- | 134 | C10H14 | 000105-05-5 | 46 |
| 4 | 1,3,8-p-Menthatriene | 134 | C10H14 | 018368-95-1 | 46 |
| 5 | Benzene, 1,3-diethyl- | 134 | C10H14 | 000141-93-5 | 46 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072324\
Data File : BG062199.D
Acq On : 24 Jul 2024 00:03
Operator : MA/JU
Sample : P3244-04
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
YDZD5

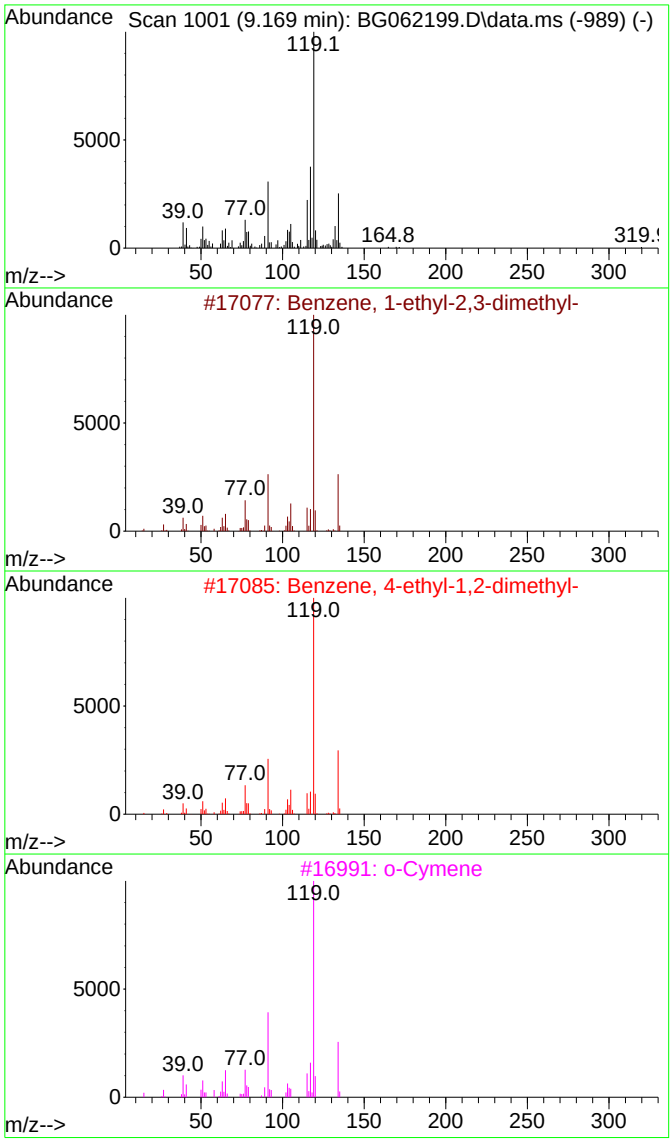
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

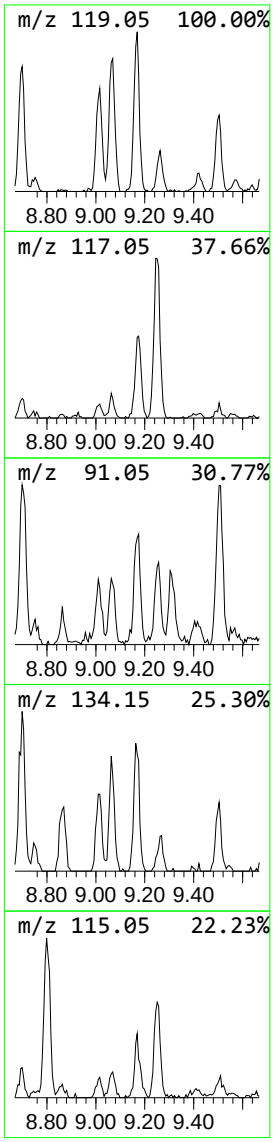
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 16 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 25

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 9.169 | 18.18 ng/ul | 341592 | 1,4-Dichlorobenzene-d4 | 8.070 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzene, 1-ethyl-2,3-dimethyl- | 134 | C10H14 | 000933-98-2 | 91 |
| 2 | Benzene, 4-ethyl-1,2-dimethyl- | 134 | C10H14 | 000934-80-5 | 91 |
| 3 | o-Cymene | 134 | C10H14 | 000527-84-4 | 70 |
| 4 | Benzene, 1-ethyl-3,5-dimethyl- | 134 | C10H14 | 000934-74-7 | 70 |
| 5 | Benzene, 2-ethyl-1,3-dimethyl- | 134 | C10H14 | 002870-04-4 | 64 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072324\
Data File : BG062199.D
Acq On : 24 Jul 2024 00:03
Operator : MA/JU
Sample : P3244-04
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
YDZD5

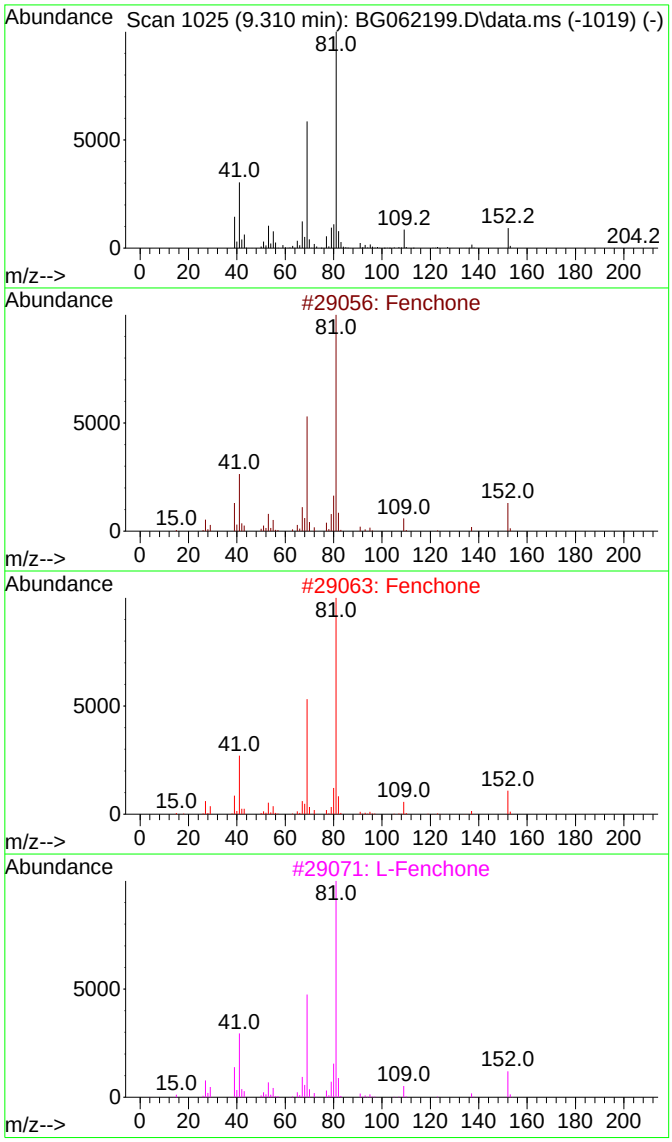
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

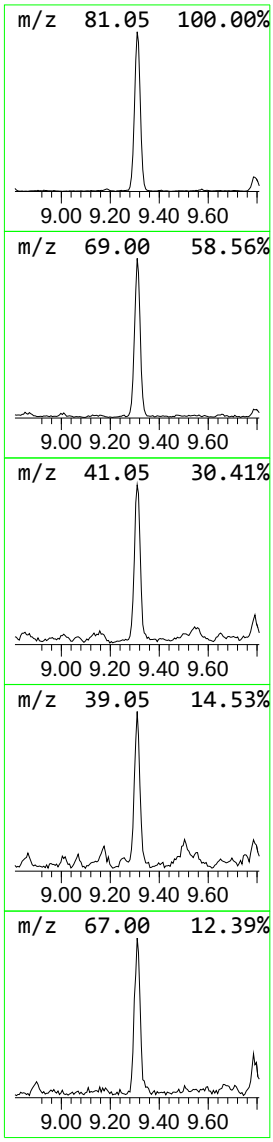
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 17 Fenchone Concentration Rank 4

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 9.310 | 70.30 ng/ul | 1320990 | 1,4-Dichlorobenzene-d4 | 8.070 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Fenchone | 152 | C10H16O | 001195-79-5 | 93 |
| 2 | Fenchone | 152 | C10H16O | 001195-79-5 | 90 |
| 3 | L-Fenchone | 152 | C10H16O | 007787-20-4 | 90 |
| 4 | Fenchone | 152 | C10H16O | 001195-79-5 | 87 |
| 5 | L-Fenchone | 152 | C10H16O | 007787-20-4 | 87 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072324\
Data File : BG062199.D
Acq On : 24 Jul 2024 00:03
Operator : MA/JU
Sample : P3244-04
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
YDZD5

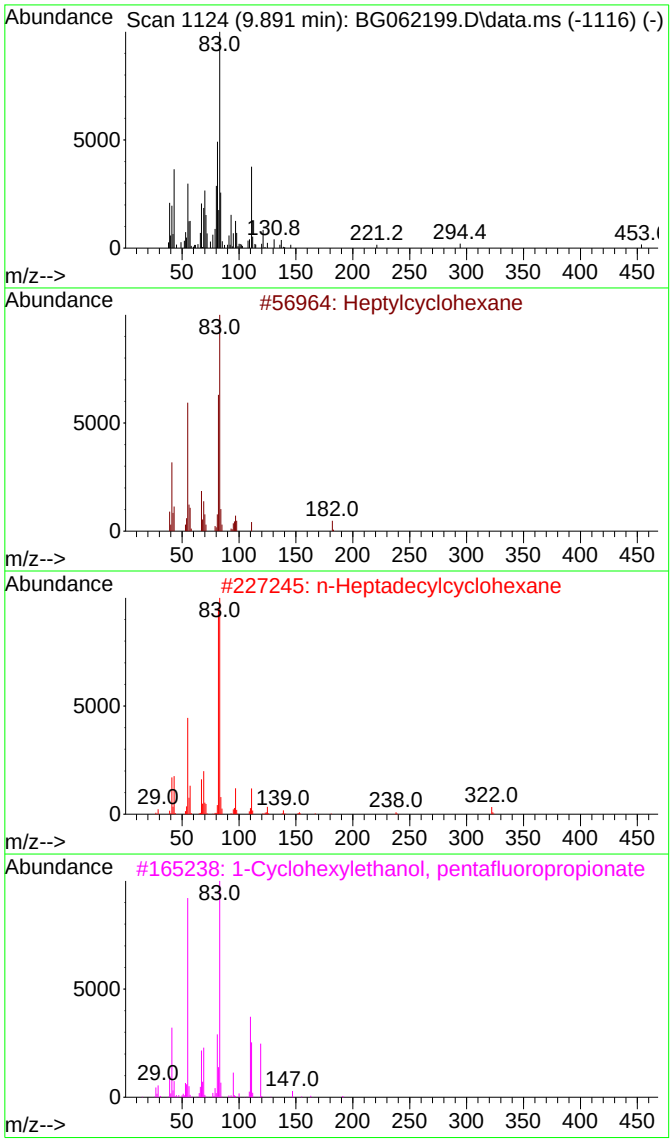
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

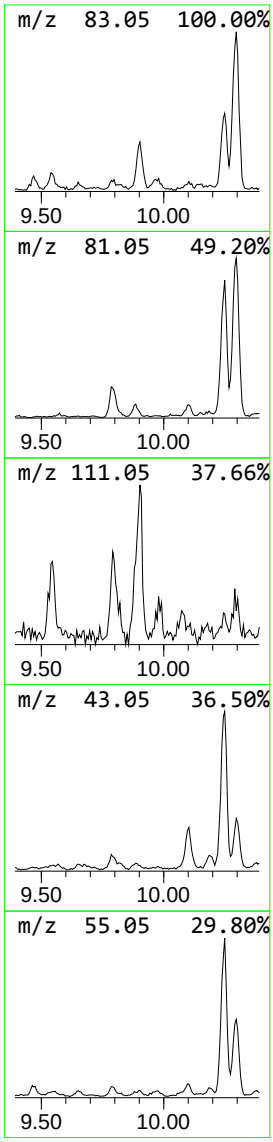
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 21 (DEL) Alkane: Cyclic9.891 Concentration Rank 61

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 9.891 | 5.02 ng/ul | 178800 | Naphthalene-d8 | 10.884 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Heptylcyclohexane | 182 | C13H26 | 005617-41-4 | 47 |
| 2 | n-Heptadecylcyclohexane | 322 | C23H46 | 019781-73-8 | 43 |
| 3 | 1-Cyclohexylethanol, pentafluoro... | 274 | C11H15F5O2 | 1000365-50-1 | 43 |
| 4 | Cyclohexane, tetradecyl- | 280 | C20H40 | 001795-18-2 | 40 |
| 5 | Cyclohexane, (4-methylpentyl)- | 168 | C12H24 | 061142-20-9 | 38 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072324\
Data File : BG062199.D
Acq On : 24 Jul 2024 00:03
Operator : MA/JU
Sample : P3244-04
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
YDZD5

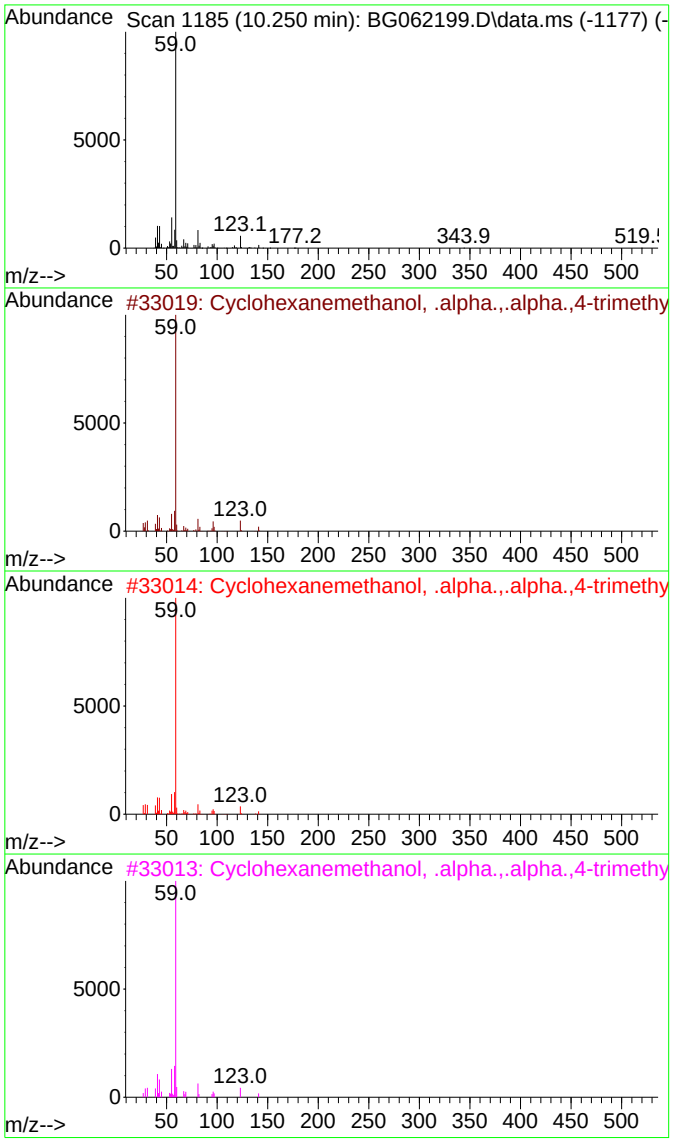
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

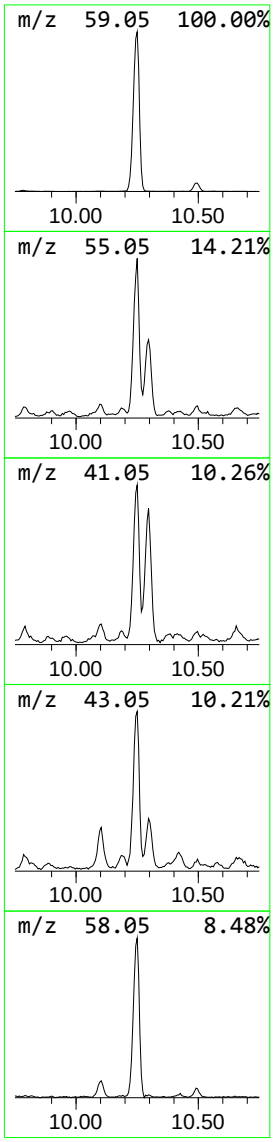
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 22 Cyclohexanemethanol, .alpha... Concentration Rank 2

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 10.249 | 113.49 ng/ul | 4040470 | Naphthalene-d8 | 10.884 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Cyclohexanemethanol, .alpha.,.al... | 156 | C10H20O | 007322-63-6 | 72 |
| 2 | Cyclohexanemethanol, .alpha.,.al... | 156 | C10H20O | 000498-81-7 | 64 |
| 3 | Cyclohexanemethanol, .alpha.,.al... | 156 | C10H20O | 000498-81-7 | 64 |
| 4 | Cyclohexanemethanol, .alpha.,.al... | 156 | C10H20O | 005114-00-1 | 56 |
| 5 | o-Menthan-8-ol | 156 | C10H20O | 343855-44-7 | 56 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072324\
Data File : BG062199.D
Acq On : 24 Jul 2024 00:03
Operator : MA/JU
Sample : P3244-04
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
YDZD5

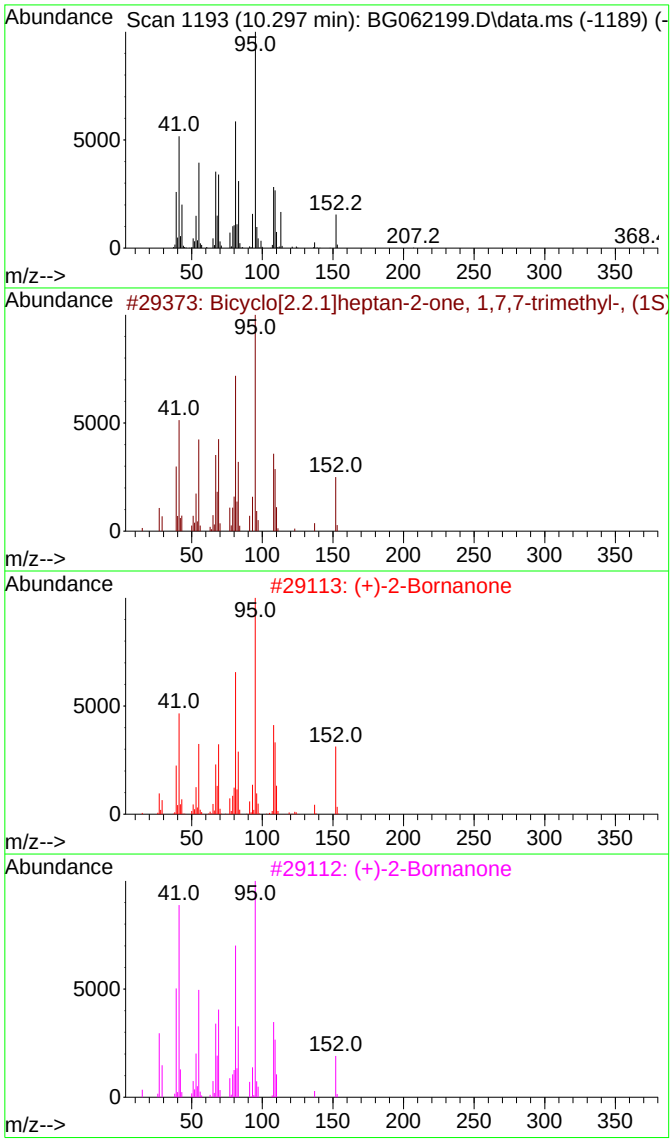
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

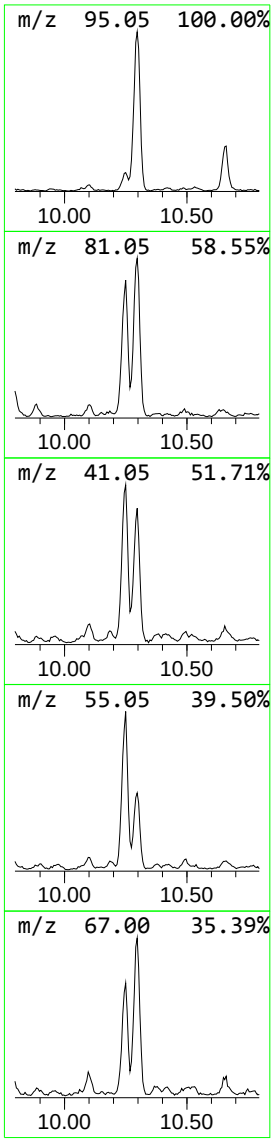
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 23 Bicyclo[2.2.1]heptan-2-one,... Concentration Rank 7

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 10.297 | 59.81 ng/ul | 2129350 | Naphthalene-d8 | 10.884 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Bicyclo[2.2.1]heptan-2-one, 1,7,... | 152 | C10H16O | 000464-48-2 | 93 |
| 2 | (+)-2-Bornanone | 152 | C10H16O | 000464-49-3 | 93 |
| 3 | (+)-2-Bornanone | 152 | C10H16O | 000464-49-3 | 90 |
| 4 | Camphor | 152 | C10H16O | 000076-22-2 | 86 |
| 5 | Bicyclo[2.2.1]heptan-2-one, 1,7,... | 152 | C10H16O | 000464-48-2 | 86 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072324\
Data File : BG062199.D
Acq On : 24 Jul 2024 00:03
Operator : MA/JU
Sample : P3244-04
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
YDZD5

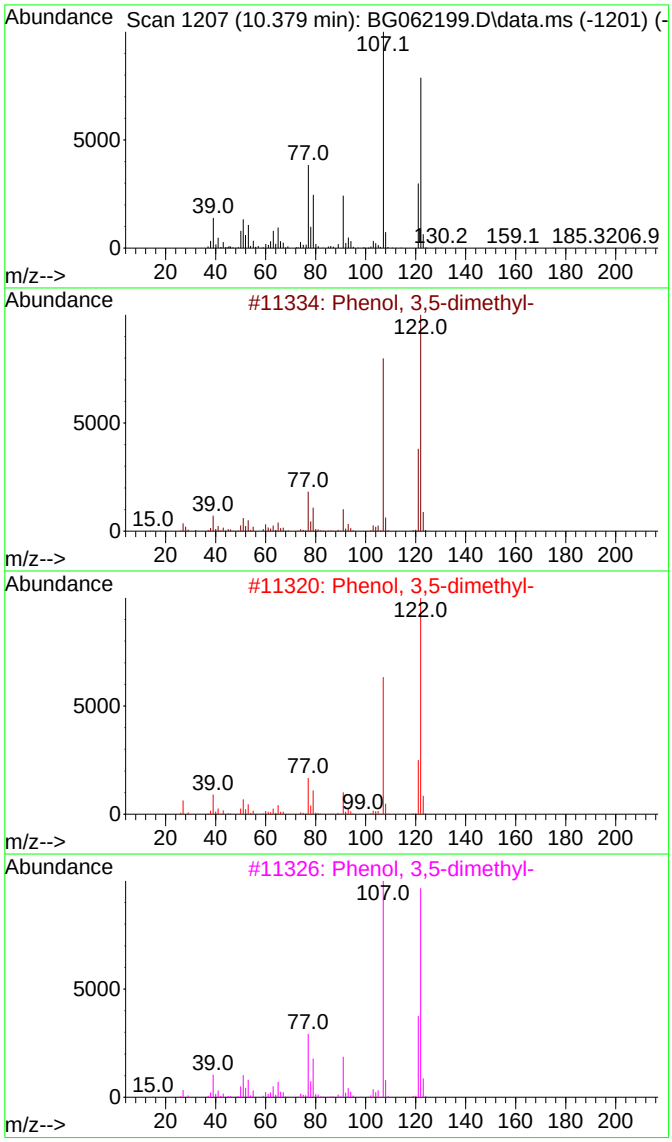
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

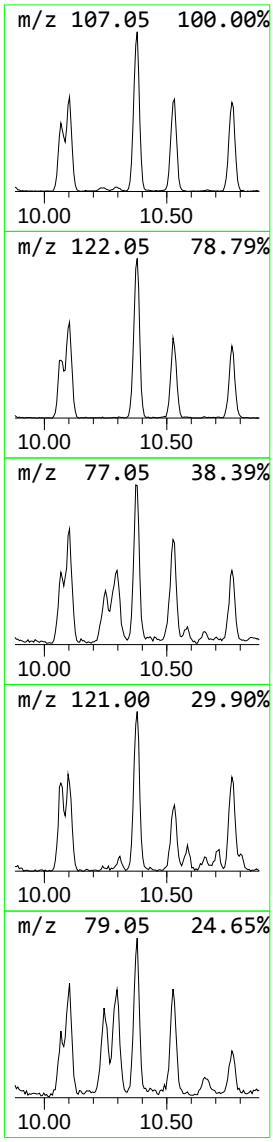
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 24 Phenol, 3,5-dimethyl- Concentration Rank 23

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 10.379 | 24.69 ng/ul | 878906 | Naphthalene-d8 | 10.884 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Phenol, 3,5-dimethyl- | 122 | C8H10O | 000108-68-9 | 96 |
| 2 | Phenol, 3,5-dimethyl- | 122 | C8H10O | 000108-68-9 | 95 |
| 3 | Phenol, 3,5-dimethyl- | 122 | C8H10O | 000108-68-9 | 95 |
| 4 | Phenol, 2,6-dimethyl- | 122 | C8H10O | 000576-26-1 | 91 |
| 5 | Phenol, 2-ethyl- | 122 | C8H10O | 000090-00-6 | 91 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072324\
Data File : BG062199.D
Acq On : 24 Jul 2024 00:03
Operator : MA/JU
Sample : P3244-04
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
YDZD5

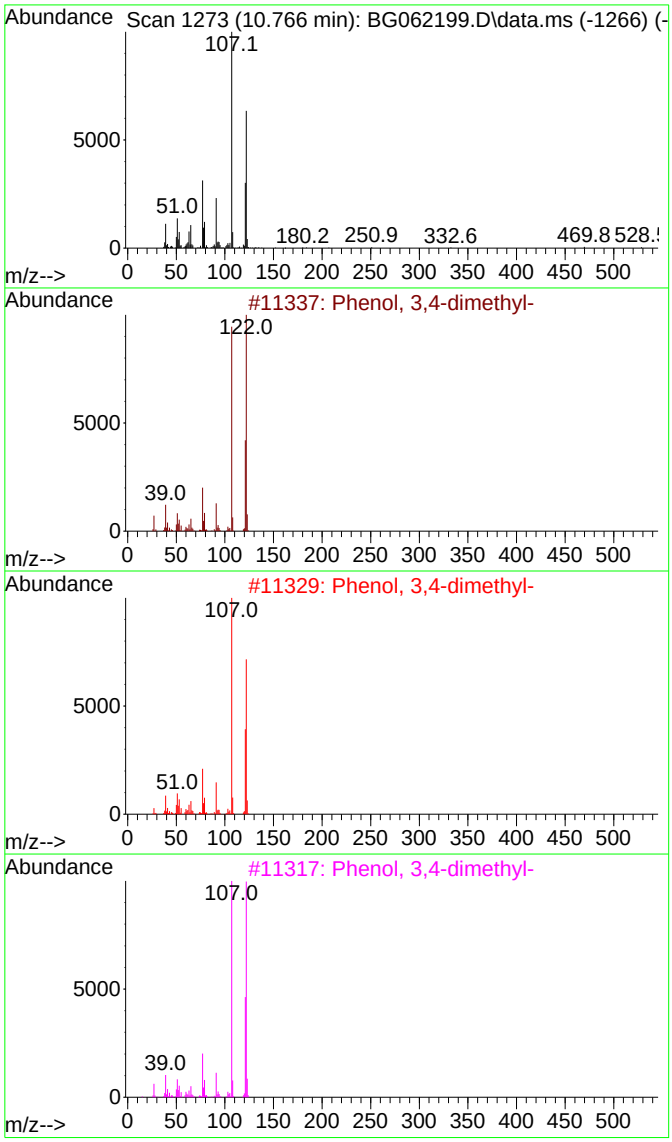
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

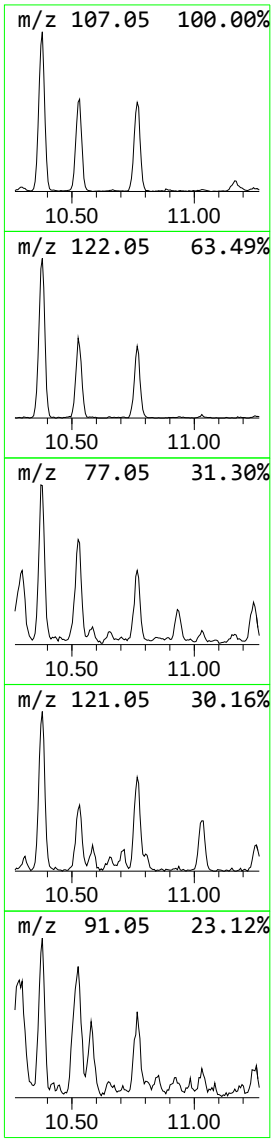
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 26 Phenol, 3,4-dimethyl- Concentration Rank 31

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 10.767 | 14.39 ng/ul | 512368 | Naphthalene-d8 | 10.884 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Phenol, 3,4-dimethyl- | 122 | C8H10O | 000095-65-8 | 95 |
| 2 | Phenol, 3,4-dimethyl- | 122 | C8H10O | 000095-65-8 | 95 |
| 3 | Phenol, 3,4-dimethyl- | 122 | C8H10O | 000095-65-8 | 94 |
| 4 | Phenol, 3,4-dimethyl- | 122 | C8H10O | 000095-65-8 | 93 |
| 5 | Phenol, 2,4-dimethyl- | 122 | C8H10O | 000105-67-9 | 91 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072324\
Data File : BG062199.D
Acq On : 24 Jul 2024 00:03
Operator : MA/JU
Sample : P3244-04
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
YDZD5

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

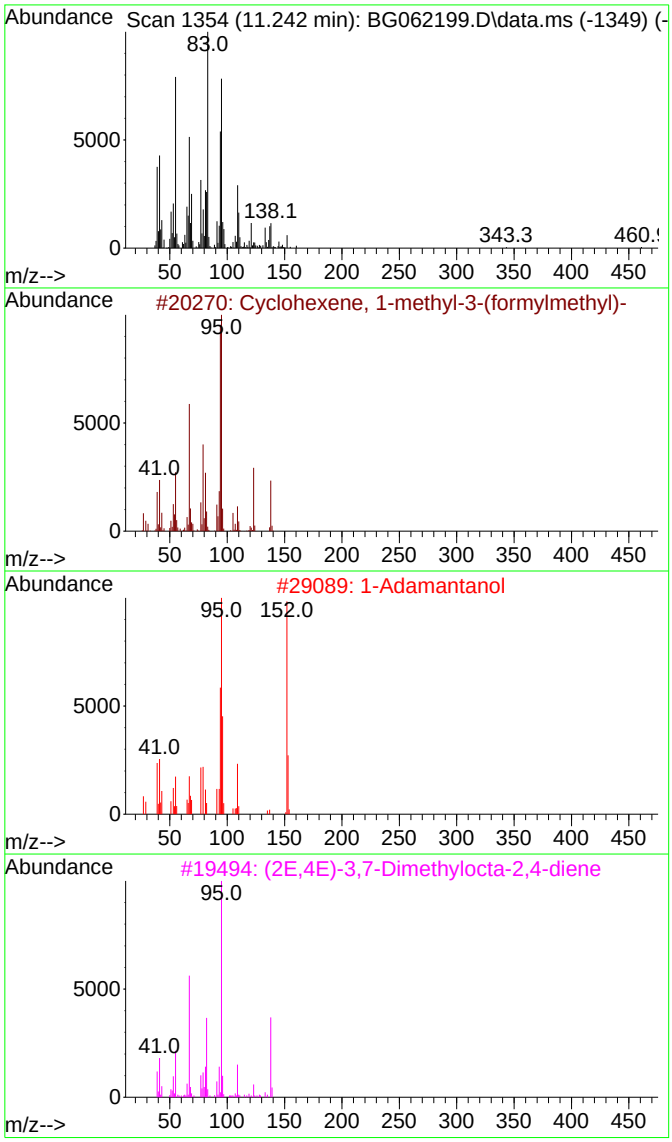
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

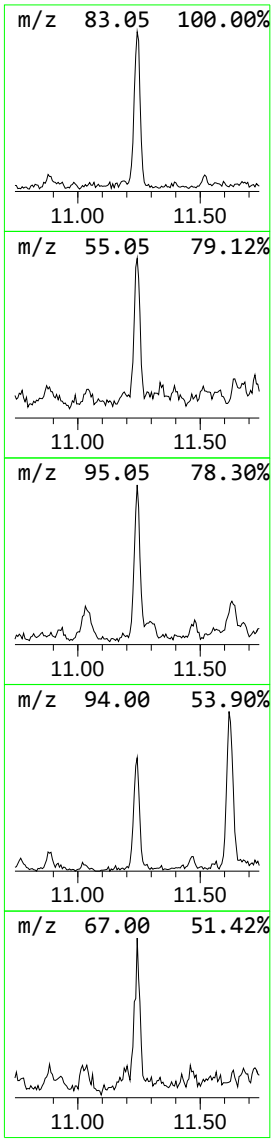
**********

Peak Number 27 unknown-01 Concentration Rank 28

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 11.242 | 16.05 ng/ul | 571593 | Naphthalene-d8 | 10.884 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Cyclohexene, 1-methyl-3-(formylm... | 138 | C9H14O | 129993-40-4 | 46 |
| 2 | 1-Adamantanol | 152 | C10H16O | 000768-95-6 | 46 |
| 3 | (2E,4E)-3,7-Dimethylocta-2,4-diene | 138 | C10H18 | 006874-39-1 | 42 |
| 4 | Cyclopentane, (3-methylbutylidene)- | 138 | C10H18 | 053366-51-1 | 38 |
| 5 | Bicyclo[2.2.1]heptane, 2,2,3-tri... | 138 | C10H18 | 000473-19-8 | 38 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072324\
Data File : BG062199.D
Acq On : 24 Jul 2024 00:03
Operator : MA/JU
Sample : P3244-04
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
YDZD5

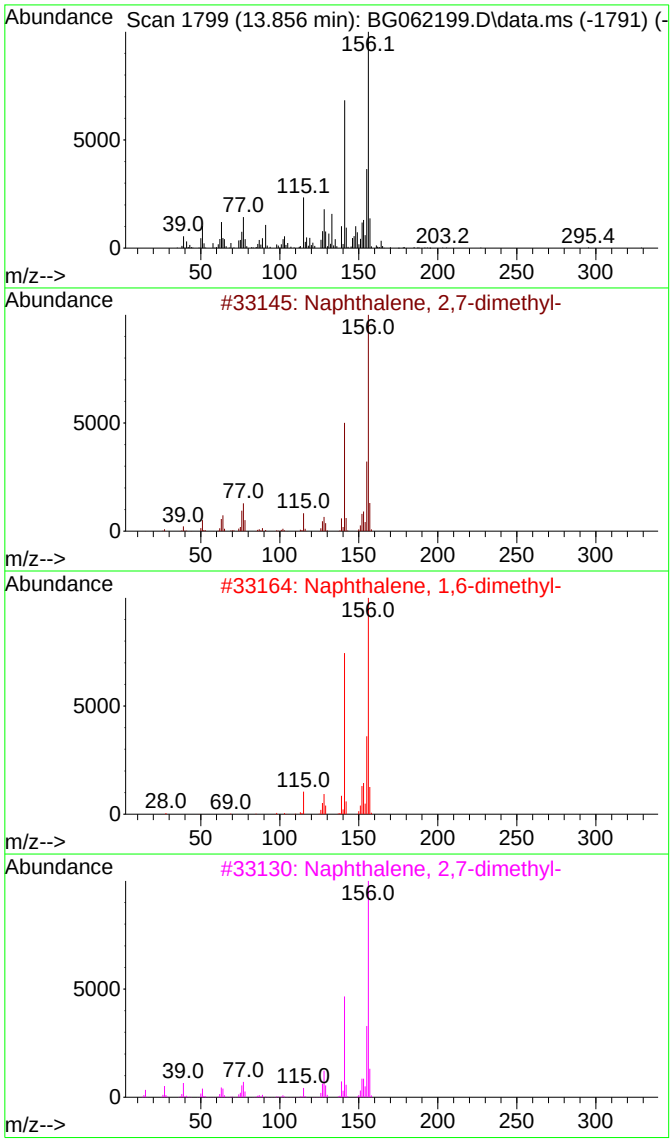
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

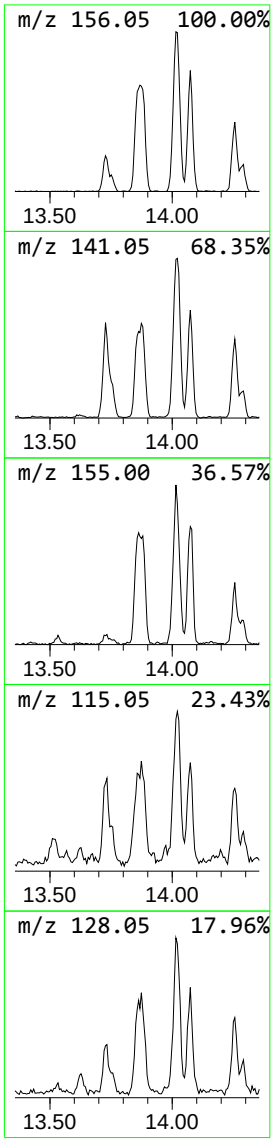
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 41 Naphthalene, 2,7-dimethyl- Concentration Rank 27

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 13.856 | 16.82 ng/ul | 614866 | Acenaphthene-d10 | 14.697 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Naphthalene, 2,7-dimethyl- | 156 | C12H12 | 000582-16-1 | 96 |
| 2 | Naphthalene, 1,6-dimethyl- | 156 | C12H12 | 000575-43-9 | 96 |
| 3 | Naphthalene, 2,7-dimethyl- | 156 | C12H12 | 000582-16-1 | 96 |
| 4 | Naphthalene, 1,7-dimethyl- | 156 | C12H12 | 000575-37-1 | 95 |
| 5 | Naphthalene, 1,6-dimethyl- | 156 | C12H12 | 000575-43-9 | 95 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072324\
Data File : BG062199.D
Acq On : 24 Jul 2024 00:03
Operator : MA/JU
Sample : P3244-04
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
YDZD5

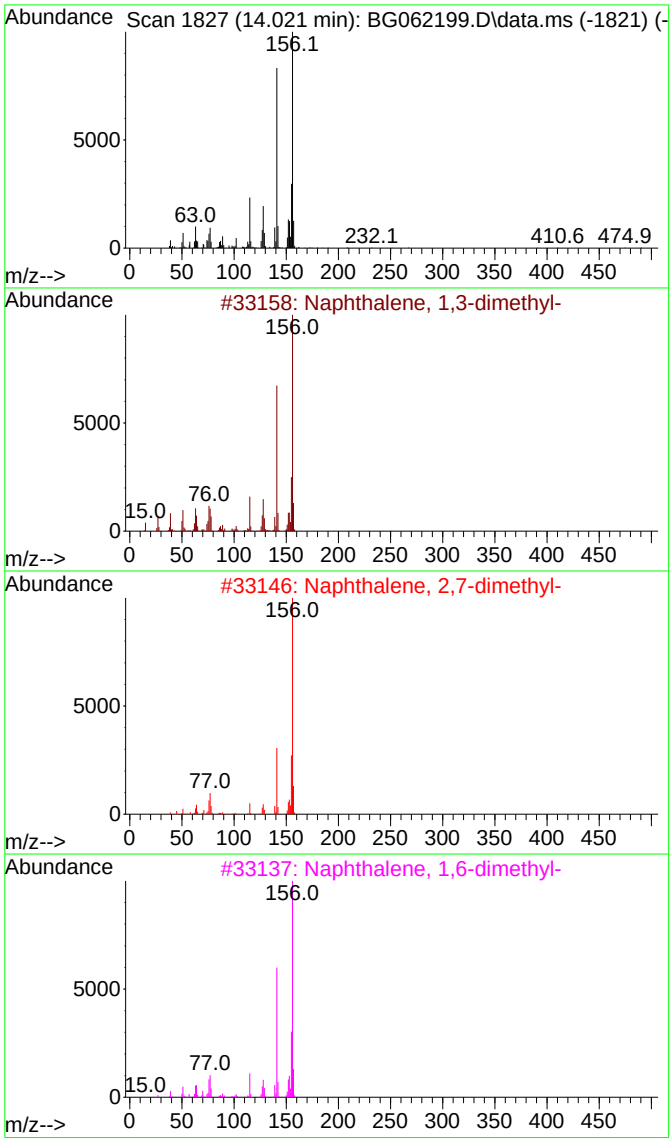
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

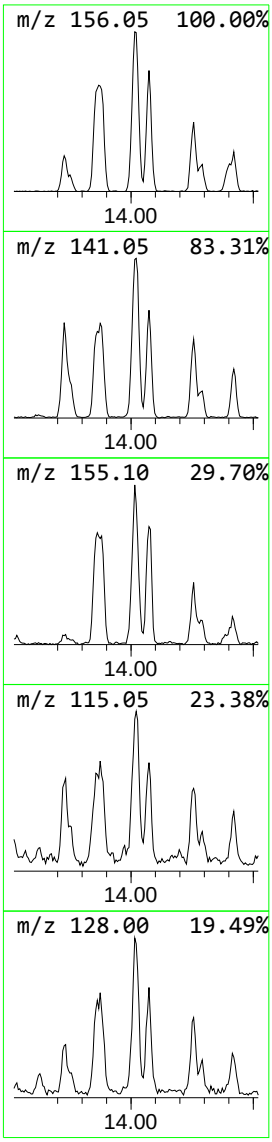
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 42 Naphthalene, 1,3-dimethyl- Concentration Rank 21

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 14.021 | 27.42 ng/ul | 1002130 | Acenaphthene-d10 | 14.697 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Naphthalene, 1,3-dimethyl- | 156 | C12H12 | 000575-41-7 | 97 |
| 2 | Naphthalene, 2,7-dimethyl- | 156 | C12H12 | 000582-16-1 | 96 |
| 3 | Naphthalene, 1,6-dimethyl- | 156 | C12H12 | 000575-43-9 | 95 |
| 4 | Naphthalene, 2,7-dimethyl- | 156 | C12H12 | 000582-16-1 | 95 |
| 5 | Naphthalene, 1,3-dimethyl- | 156 | C12H12 | 000575-41-7 | 95 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072324\
Data File : BG062199.D
Acq On : 24 Jul 2024 00:03
Operator : MA/JU
Sample : P3244-04
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
YDZD5

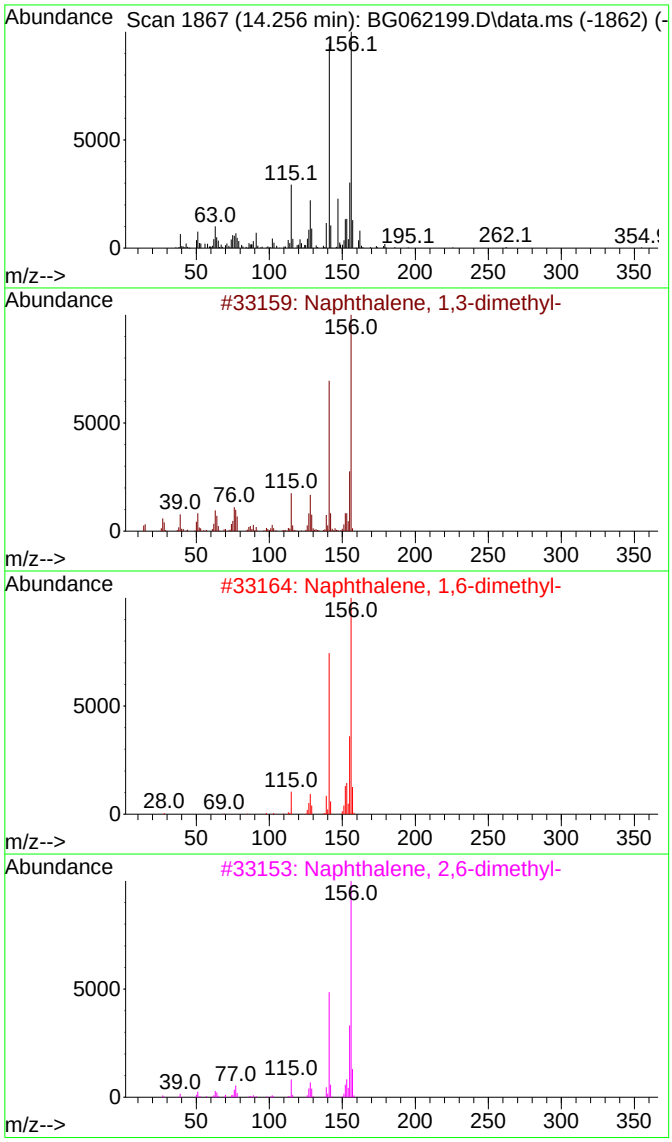
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

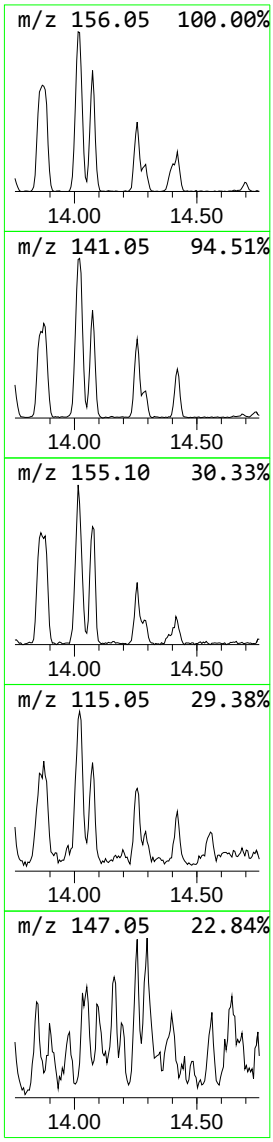
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 43 Naphthalene, 1,6-dimethyl- Concentration Rank 38

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 14.256 | 11.03 ng/ul | 403020 | Acenaphthene-d10 | 14.697 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Naphthalene, 1,3-dimethyl- | 156 | C12H12 | 000575-41-7 | 95 |
| 2 | Naphthalene, 1,6-dimethyl- | 156 | C12H12 | 000575-43-9 | 95 |
| 3 | Naphthalene, 2,6-dimethyl- | 156 | C12H12 | 000581-42-0 | 95 |
| 4 | Naphthalene, 1,3-dimethyl- | 156 | C12H12 | 000575-41-7 | 94 |
| 5 | Naphthalene, 1,6-dimethyl- | 156 | C12H12 | 000575-43-9 | 93 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072324\
Data File : BG062199.D
Acq On : 24 Jul 2024 00:03
Operator : MA/JU
Sample : P3244-04
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
YDZD5

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

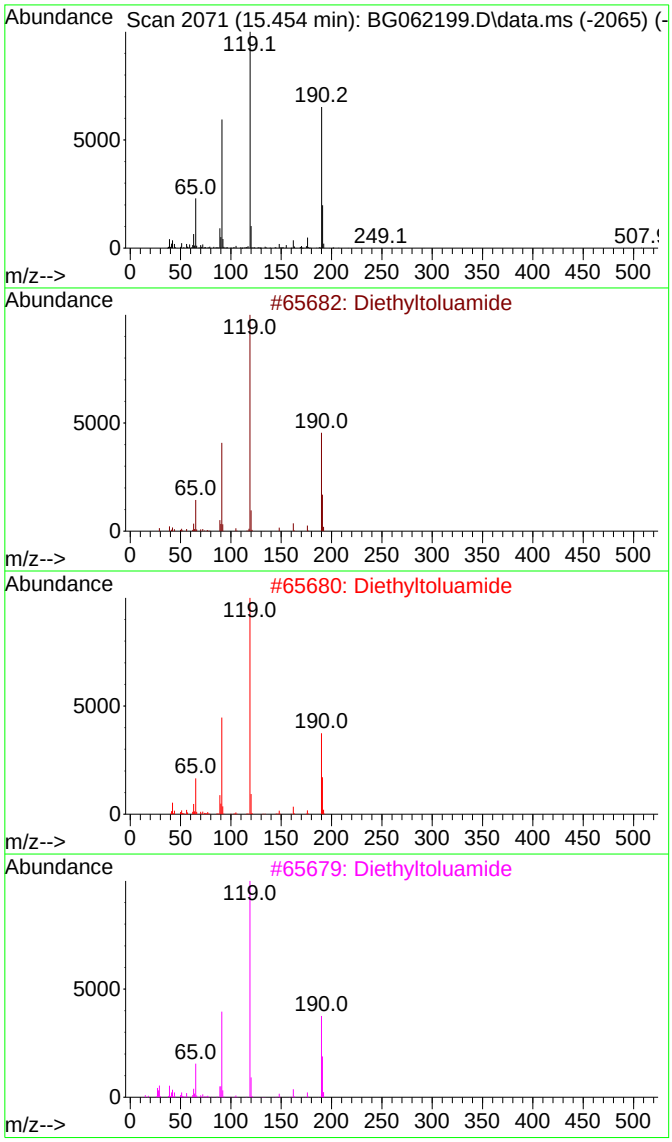
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

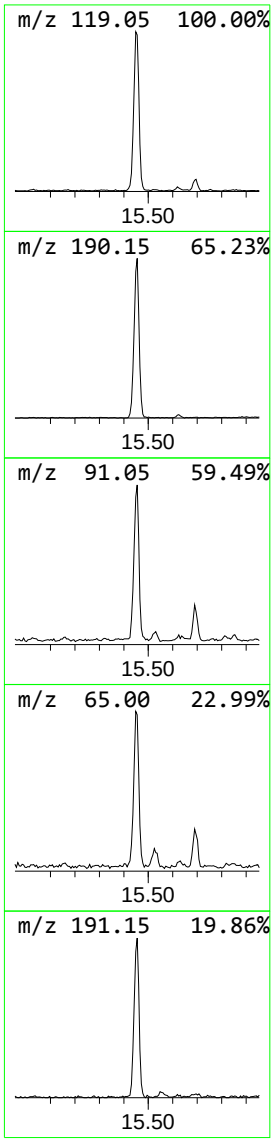
**********

Peak Number 45 Diethyltoluamide Concentration Rank 15

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 15.454 | 35.70 ng/ul | 1304800 | Acenaphthene-d10 | 14.697 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Diethyltoluamide | 191 | C12H17NO | 000134-62-3 | 95 |
| 2 | Diethyltoluamide | 191 | C12H17NO | 000134-62-3 | 95 |
| 3 | Diethyltoluamide | 191 | C12H17NO | 000134-62-3 | 81 |
| 4 | Benzamide, N,N-diethyl-4-methyl- | 191 | C12H17NO | 002728-05-4 | 72 |
| 5 | Diethyltoluamide | 191 | C12H17NO | 000134-62-3 | 68 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072324\
Data File : BG062199.D
Acq On : 24 Jul 2024 00:03
Operator : MA/JU
Sample : P3244-04
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
YDZD5

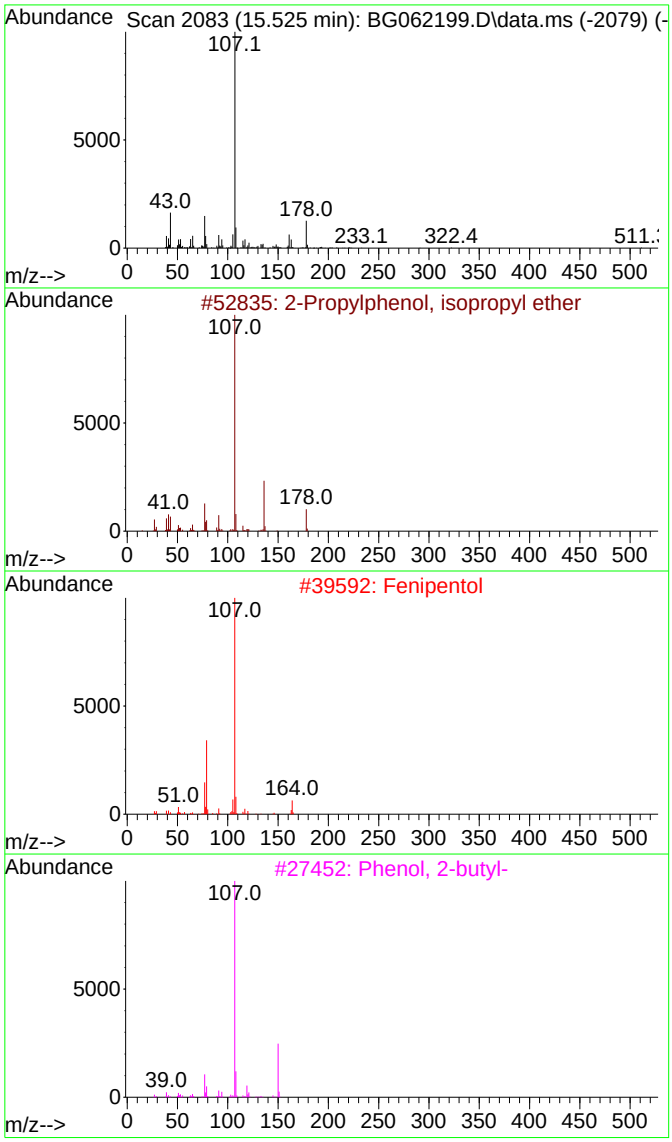
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

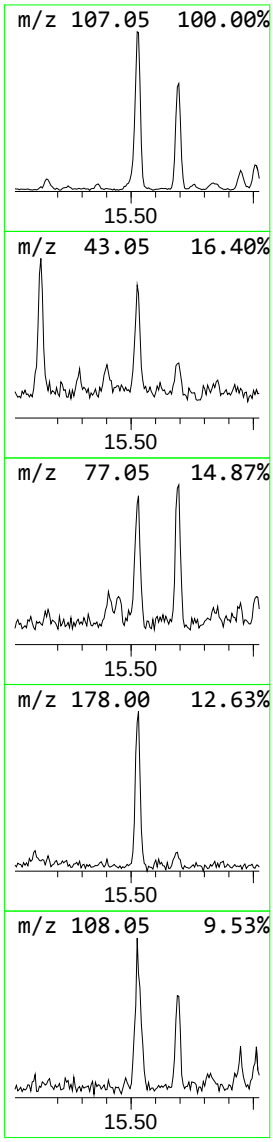
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 46 2-Propylphenol, isopropyl e... Concentration Rank 32

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 15.525 | 13.75 ng/ul | 502716 | Acenaphthene-d10 | 14.697 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2-Propylphenol, isopropyl ether | 178 | C12H18O | 1000395-18-7 | 72 |
| 2 | Fenipentol | 164 | C11H16O | 000583-03-9 | 64 |
| 3 | Phenol, 2-butyl- | 150 | C10H14O | 003180-09-4 | 64 |
| 4 | Phenol, 4-hexyl- | 178 | C12H18O | 002446-69-7 | 64 |
| 5 | L-Tyrosine, N-(methoxycarbonyl)-... | 253 | C12H15NO5 | 215300-34-8 | 64 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072324\
Data File : BG062199.D
Acq On : 24 Jul 2024 00:03
Operator : MA/JU
Sample : P3244-04
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
YDZD5

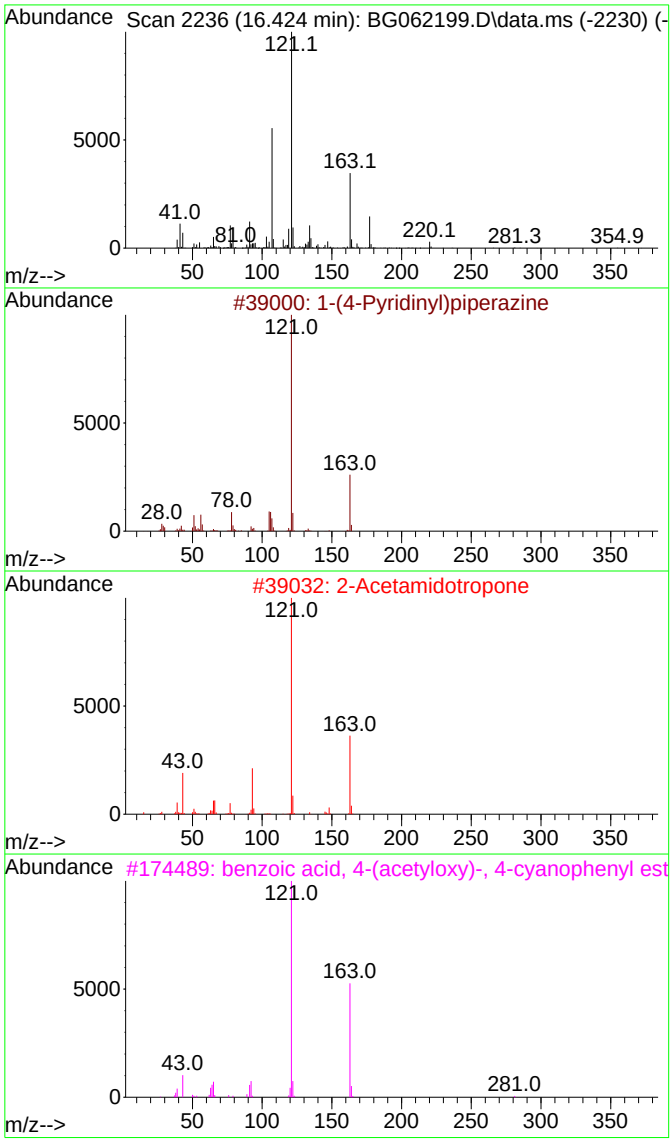
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

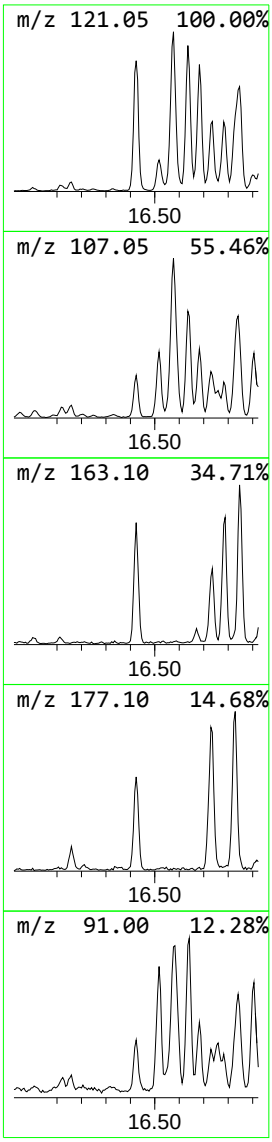
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 50 1-(4-Pyridinyl)piperazine Concentration Rank 18

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 16.424 | 29.18 ng/ul | 1146580 | Phenanthrene-d10 | 17.446 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1-(4-Pyridinyl)piperazine | 163 | C9H13N3 | 001008-91-9 | 50 |
| 2 | 2-Acetamidotropone | 163 | C9H9NO2 | 006422-12-4 | 49 |
| 3 | benzoic acid, 4-(acetyloxy)-, 4-... | 281 | C16H11NO4 | 1000401-91-0 | 47 |
| 4 | 1-Propyl-1,2,3,4-tetrahydropyrro... | 164 | C10H16N2 | 112758-86-8 | 47 |
| 5 | 4-Isopropoxybenzohydrazide, 3Ac ... | 320 | C16H20N2O5 | 1000486-70-2 | 47 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072324\
Data File : BG062199.D
Acq On : 24 Jul 2024 00:03
Operator : MA/JU
Sample : P3244-04
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
YDZD5

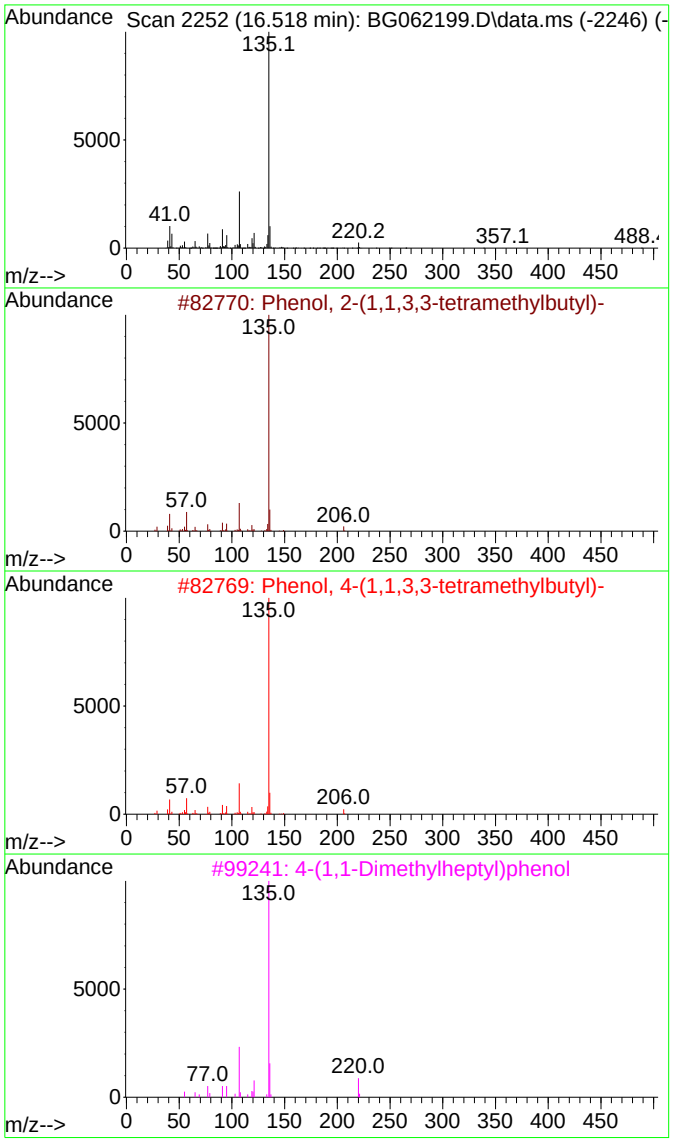
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

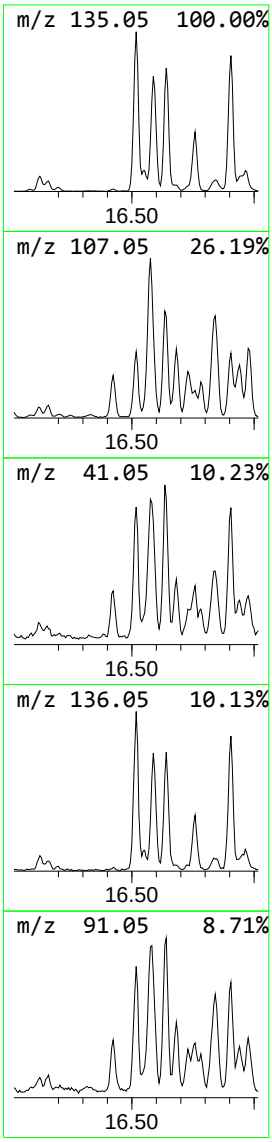
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 51 Phenol, 2-(1,1,3,3-tetramet... Concentration Rank 8

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 16.517 | 59.63 ng/ul | 2343180 | Phenanthrene-d10 | 17.446 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Phenol, 2-(1,1,3,3-tetramethylbu... | 206 | C14H22O | 003884-95-5 | 90 |
| 2 | Phenol, 4-(1,1,3,3-tetramethylbu... | 206 | C14H22O | 000140-66-9 | 90 |
| 3 | 4-(1,1-Dimethylheptyl)phenol | 220 | C15H24O | 1000384-80-3 | 87 |
| 4 | ((1,2-Diethylethylene)bis(p-phen... | 354 | C22H26O4 | 004547-76-6 | 78 |
| 5 | Phenol, 4-(1,1-dimethylpropyl)- | 164 | C11H16O | 000080-46-6 | 78 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072324\
Data File : BG062199.D
Acq On : 24 Jul 2024 00:03
Operator : MA/JU
Sample : P3244-04
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
YDZD5

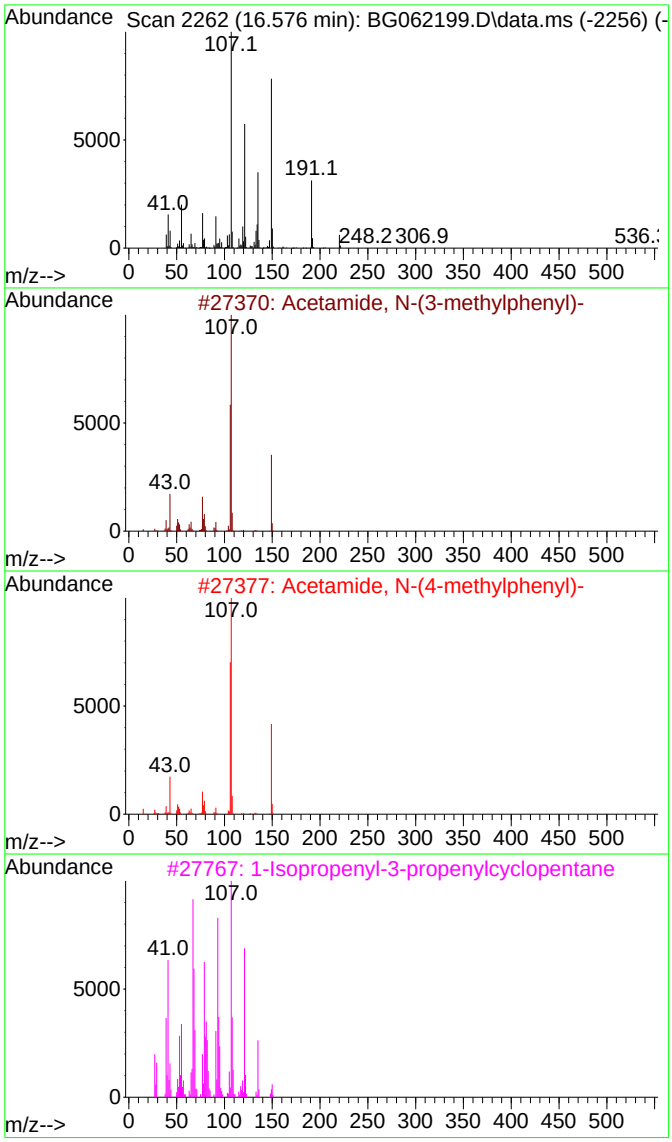
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

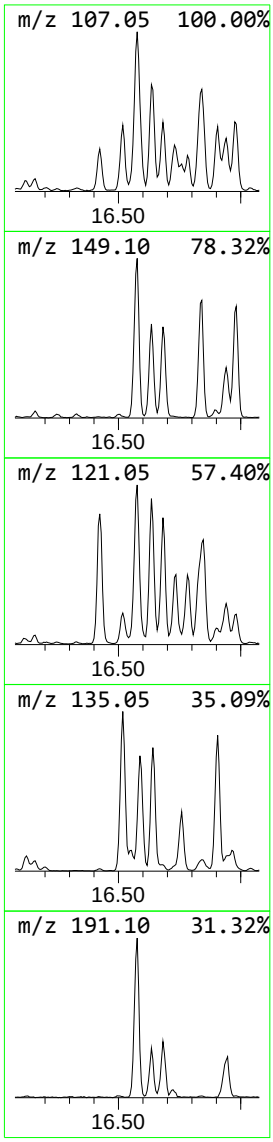
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 52 unknown-02 Concentration Rank 1

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 16.576 | 118.75 ng/ul | 4666090 | Phenanthrene-d10 | 17.446 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Acetamide, N-(3-methylphenyl)- | 149 | C9H11NO | 000537-92-8 | 43 |
| 2 | Acetamide, N-(4-methylphenyl)- | 149 | C9H11NO | 000103-89-9 | 38 |
| 3 | 1-Isopropenyl-3-propenylcyclopen... | 150 | C11H18 | 1000192-81-2 | 38 |
| 4 | Pyridine-3-carbonitrile, 1,2-dih... | 191 | C9H9N3O2 | 220098-92-0 | 35 |
| 5 | 4-Methyl-2-tert-octylphenol | 220 | C15H24O | 004979-46-8 | 30 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072324\
Data File : BG062199.D
Acq On : 24 Jul 2024 00:03
Operator : MA/JU
Sample : P3244-04
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
YDZD5

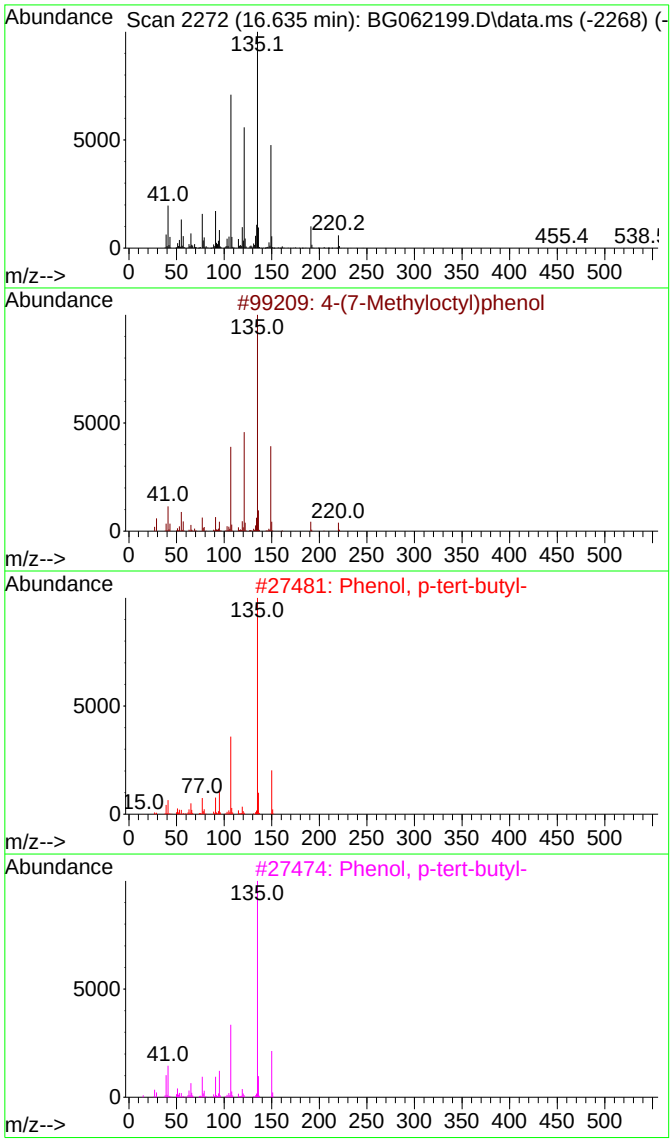
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

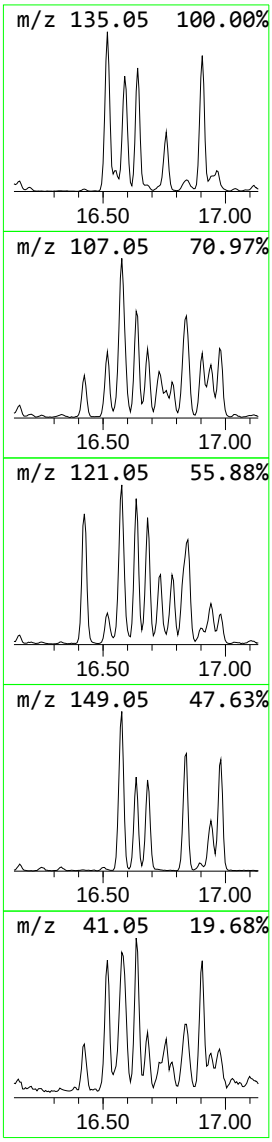
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 53 4-(7-Methyloctyl)phenol Concentration Rank 3

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 16.635 | 81.84 ng/ul | 3216080 | Phenanthrene-d10 | 17.446 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 4-(7-Methyloctyl)phenol | 220 | C15H24O | 024518-48-7 | 97 |
| 2 | Phenol, p-tert-butyl- | 150 | C10H14O | 000098-54-4 | 64 |
| 3 | Phenol, p-tert-butyl- | 150 | C10H14O | 000098-54-4 | 52 |
| 4 | 1,5,6,7-Tetrahydro-4-indolone | 135 | C8H9NO | 013754-86-4 | 49 |
| 5 | Phenol, 2-(1,1-dimethylethyl)- | 150 | C10H14O | 000088-18-6 | 49 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072324\
Data File : BG062199.D
Acq On : 24 Jul 2024 00:03
Operator : MA/JU
Sample : P3244-04
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
YDZD5

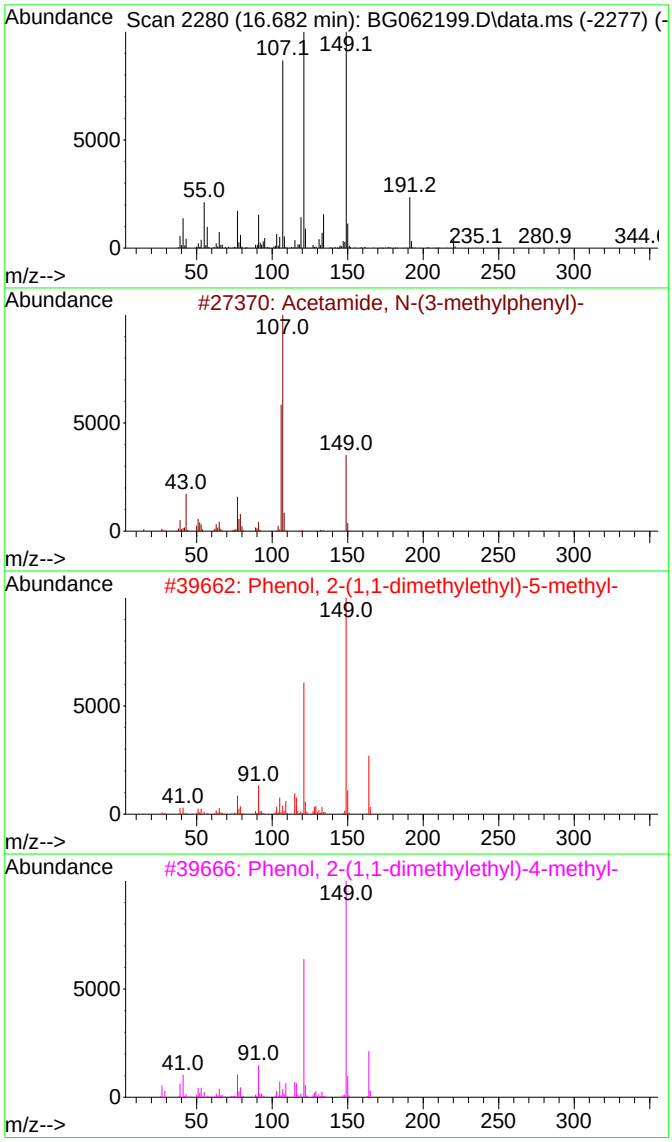
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

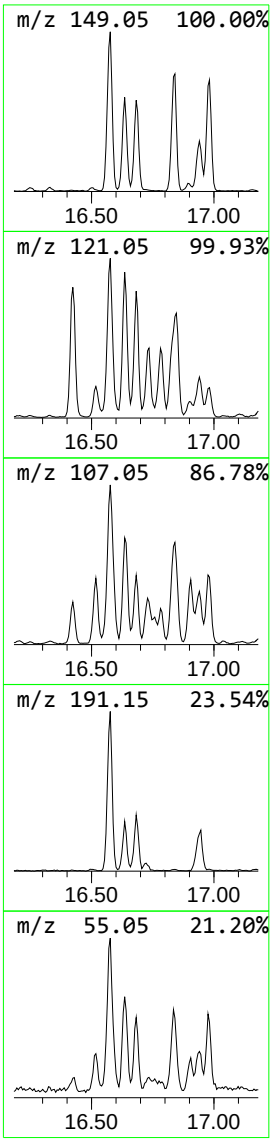
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 54 Acetamide, N-(3-methylphenyl)- Concentration Rank 12

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 16.682 | 37.59 ng/ul | 1476900 | Phenanthrene-d10 | 17.446 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Acetamide, N-(3-methylphenyl)- | 149 | C9H11NO | 000537-92-8 | 55 |
| 2 | Phenol, 2-(1,1-dimethylethyl)-5-... | 164 | C11H16O | 000088-60-8 | 47 |
| 3 | Phenol, 2-(1,1-dimethylethyl)-4-... | 164 | C11H16O | 002409-55-4 | 47 |
| 4 | Phenol, 2-(1,1-dimethylethyl)-5-... | 164 | C11H16O | 000088-60-8 | 43 |
| 5 | Phenol, 2-(1,1-dimethylethyl)-6-... | 164 | C11H16O | 002219-82-1 | 43 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072324\
Data File : BG062199.D
Acq On : 24 Jul 2024 00:03
Operator : MA/JU
Sample : P3244-04
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
YDZD5

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

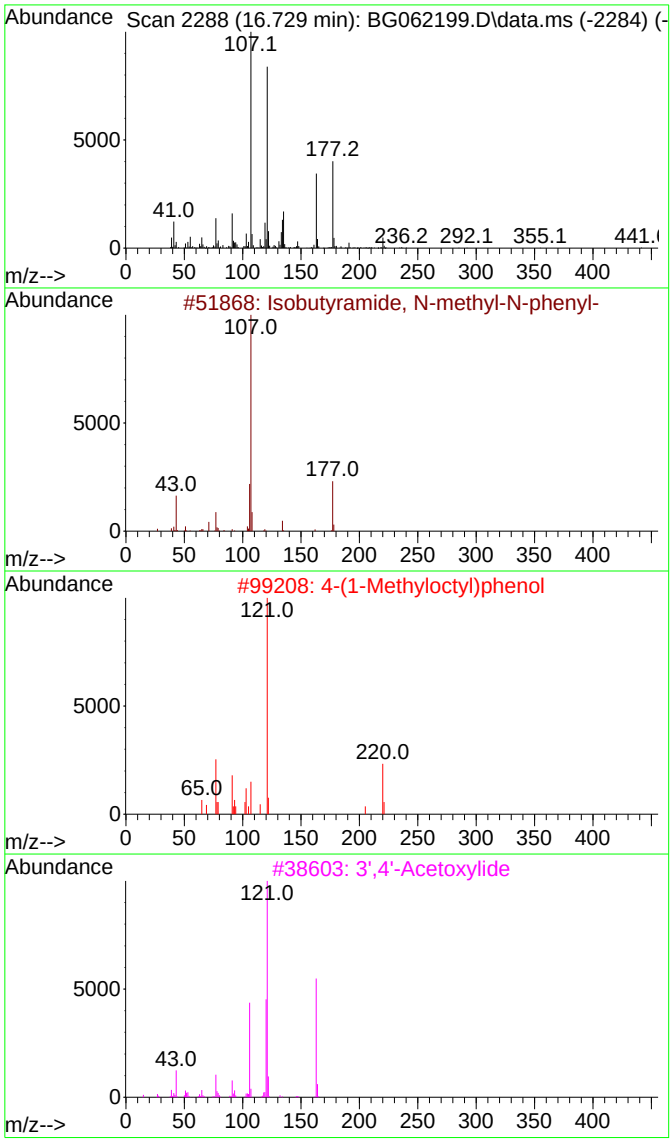
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

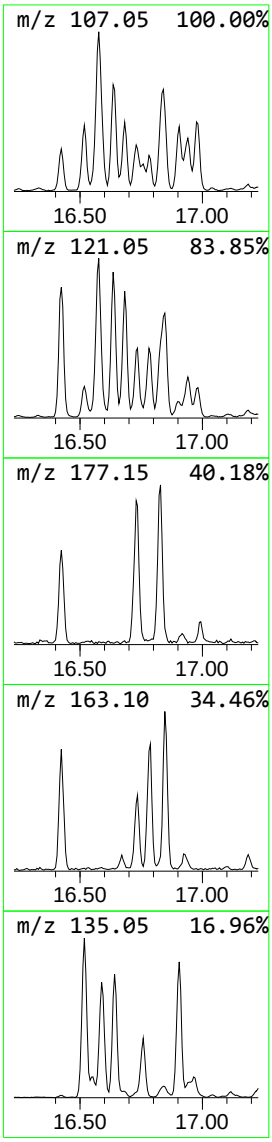
**********

Peak Number 55 unknown-03 Concentration Rank 22

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 16.729 | 25.50 ng/ul | 1002000 | Phenanthrene-d10 | 17.446 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Isobutyramide, N-methyl-N-phenyl- | 177 | C11H15NO | 1000339-96-1 | 38 |
| 2 | 4-(1-Methyloctyl)phenol | 220 | C15H24O | 1000384-80-2 | 35 |
| 3 | 3',4'-Acetoxylide | 163 | C10H13NO | 002198-54-1 | 30 |
| 4 | p-Propionotoluidide | 163 | C10H13NO | 002759-55-9 | 30 |
| 5 | Acetamide, N-methyl-N-(4-methylp... | 163 | C10H13NO | 000612-03-3 | 27 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072324\
Data File : BG062199.D
Acq On : 24 Jul 2024 00:03
Operator : MA/JU
Sample : P3244-04
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
YDZD5

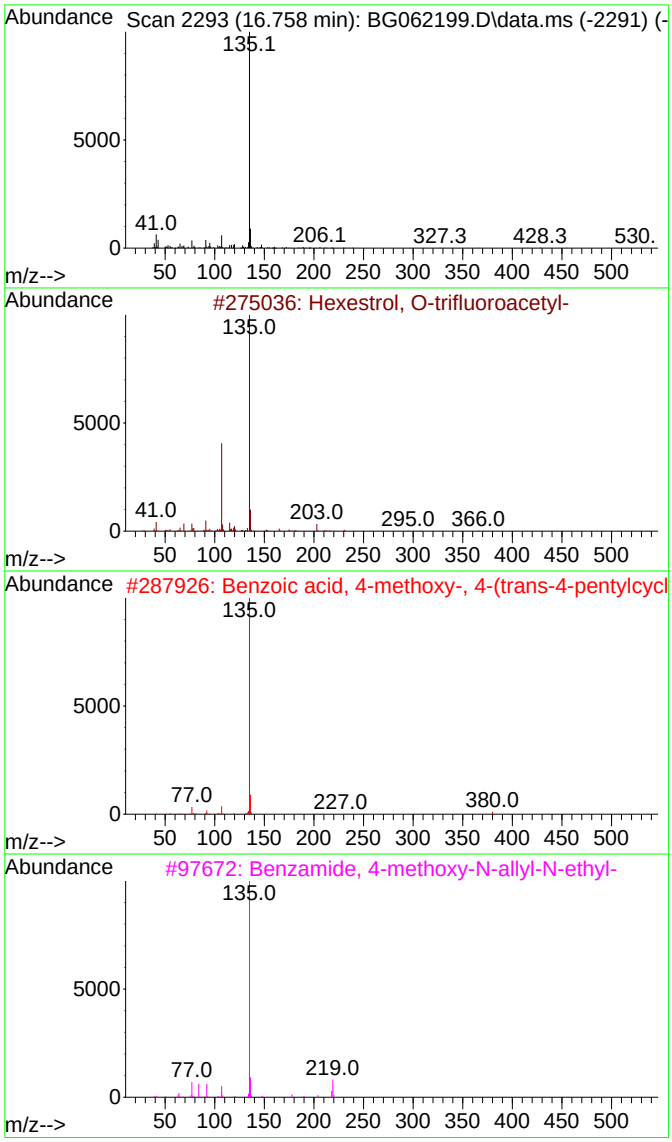
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

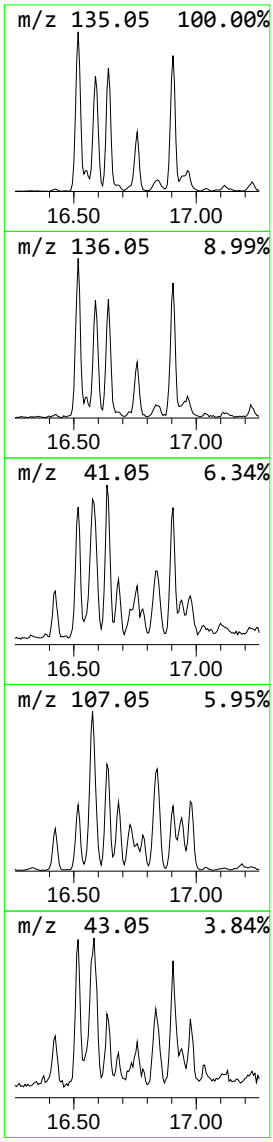
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 56 Hexestrol, O-trifluoroacetyl- Concentration Rank 26

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 16.758 | 17.70 ng/ul | 695686 | Phenanthrene-d10 | 17.446 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Hexestrol, O-trifluoroacetyl- | 366 | C20H21F3O3 | 1000365-31-7 | 74 |
| 2 | Benzoic acid, 4-methoxy-, 4-(tra... | 380 | C25H32O3 | 1229648-08-1 | 72 |
| 3 | Benzamide, 4-methoxy-N-allyl-N-e... | 219 | C13H17NO2 | 1000446-33-5 | 72 |
| 4 | 1-Adamantanecarboxylic acid, 3-m... | 270 | C18H22O2 | 1000307-65-9 | 64 |
| 5 | 1-Adamantanecarboxylic acid, 3,4... | 324 | C17H18Cl2O2 | 1000307-66-4 | 64 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072324\
Data File : BG062199.D
Acq On : 24 Jul 2024 00:03
Operator : MA/JU
Sample : P3244-04
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
YDZD5

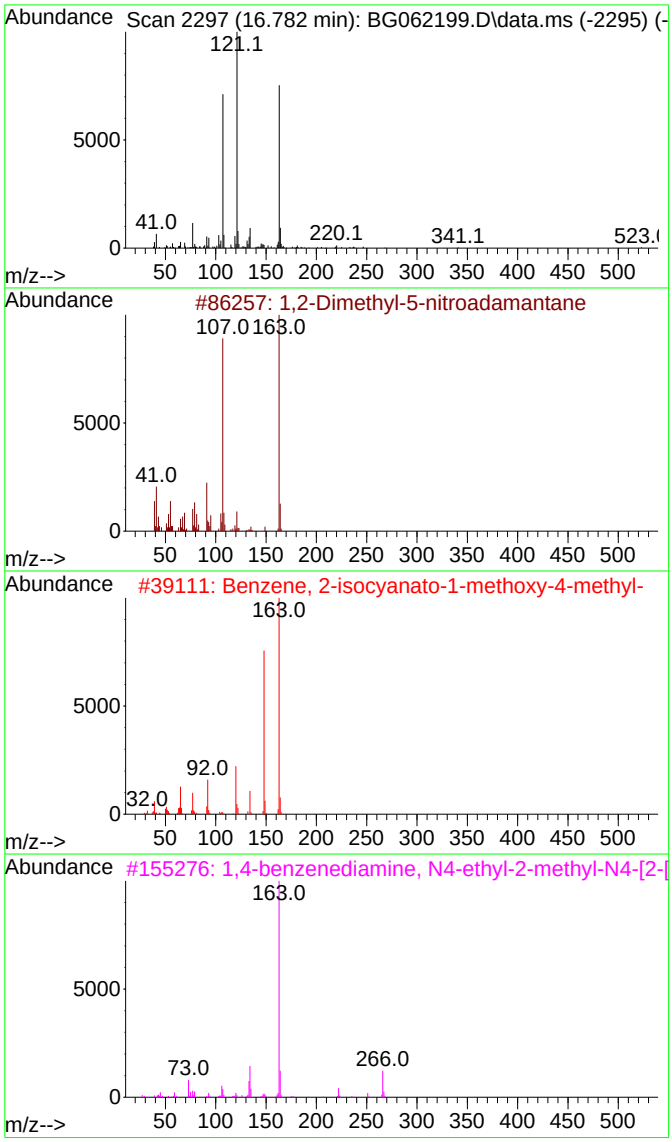
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

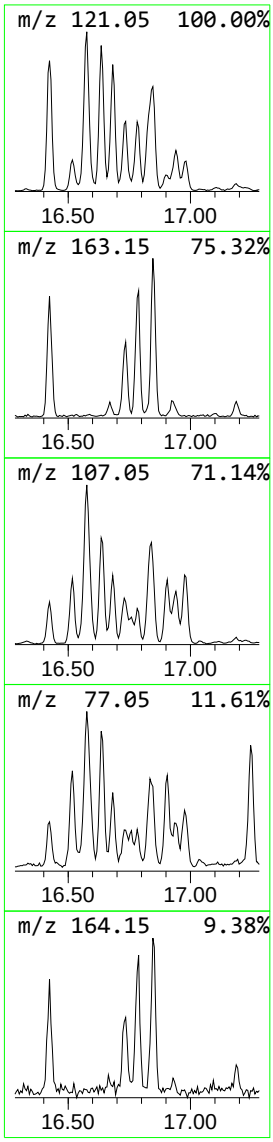
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 57 unknown-04 Concentration Rank 29

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 16.782 | 15.71 ng/ul | 617241 | Phenanthrene-d10 | 17.446 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1,2-Dimethyl-5-nitroadamantane | 209 | C12H19NO2 | 006588-68-7 | 43 |
| 2 | Benzene, 2-isocyanato-1-methoxy-... | 163 | C9H9NO2 | 1000319-47-4 | 35 |
| 3 | 1,4-benzenediamine, N4-ethyl-2-m... | 266 | C14H26N2OSi | 1000396-92-9 | 35 |
| 4 | 2-Propanone, 1-(4-methoxyphenyl)- | 164 | C10H12O2 | 000122-84-9 | 30 |
| 5 | 2-Propanone, 1-(4-methoxyphenyl)- | 164 | C10H12O2 | 000122-84-9 | 30 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072324\
Data File : BG062199.D
Acq On : 24 Jul 2024 00:03
Operator : MA/JU
Sample : P3244-04
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
YDZD5

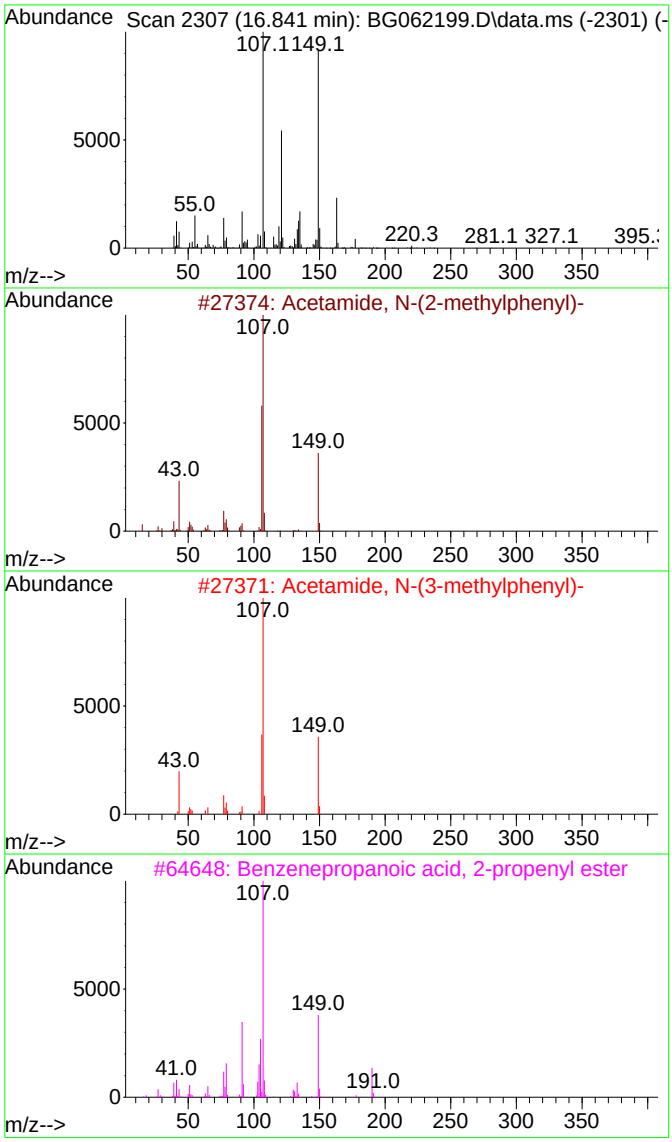
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

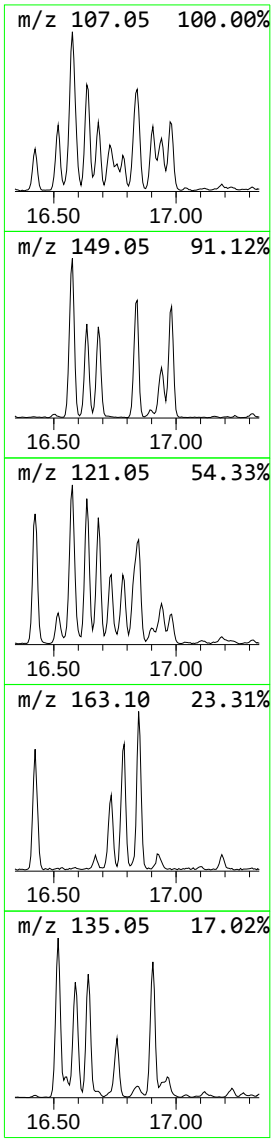
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 58 Acetamide, N-(2-methylphenyl)- Concentration Rank 5

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 16.841 | 68.06 ng/ul | 2674560 | Phenanthrene-d10 | 17.446 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Acetamide, N-(2-methylphenyl)- | 149 | C9H11NO | 000120-66-1 | 53 |
| 2 | Acetamide, N-(3-methylphenyl)- | 149 | C9H11NO | 000537-92-8 | 46 |
| 3 | Benzenepropanoic acid, 2-propeny... | 190 | C12H14O2 | 015814-45-6 | 43 |
| 4 | Acetamide, N-(3-methylphenyl)- | 149 | C9H11NO | 000537-92-8 | 43 |
| 5 | Acetamide, N-(4-methylphenyl)- | 149 | C9H11NO | 000103-89-9 | 43 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072324\
Data File : BG062199.D
Acq On : 24 Jul 2024 00:03
Operator : MA/JU
Sample : P3244-04
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
YDZD5

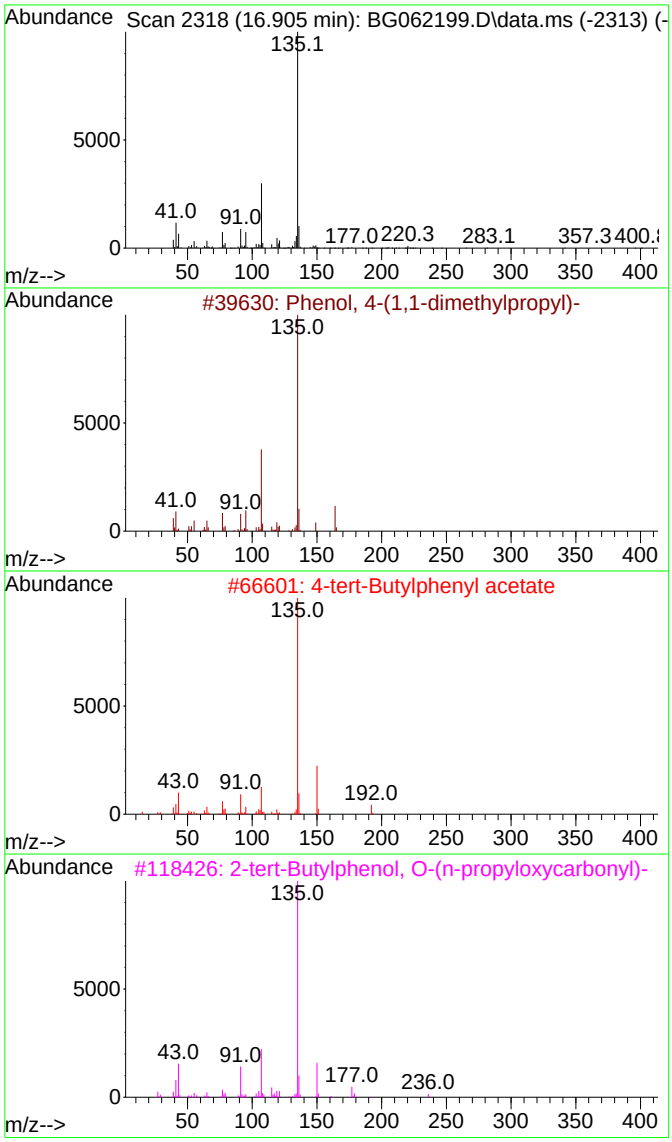
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

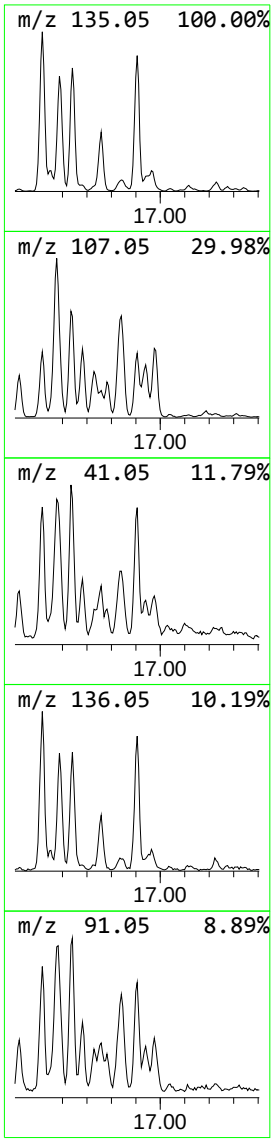
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 59 Phenol, 4-(1,1-dimethylprop... Concentration Rank 9

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 16.905 | 53.16 ng/ul | 2089020 | Phenanthrene-d10 | 17.446 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Phenol, 4-(1,1-dimethylpropyl)- | 164 | C11H16O | 000080-46-6 | 83 |
| 2 | 4-tert-Butylphenyl acetate | 192 | C12H16O2 | 003056-64-2 | 78 |
| 3 | 2-tert-Butylphenol, O-(n-propylo... | 236 | C14H20O3 | 1000487-37-9 | 78 |
| 4 | Phenol, 4-(1,1-dimethylpropyl)- | 164 | C11H16O | 000080-46-6 | 78 |
| 5 | Phenol, m-tert-butyl- | 150 | C10H14O | 000585-34-2 | 78 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072324\
Data File : BG062199.D
Acq On : 24 Jul 2024 00:03
Operator : MA/JU
Sample : P3244-04
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
YDZD5

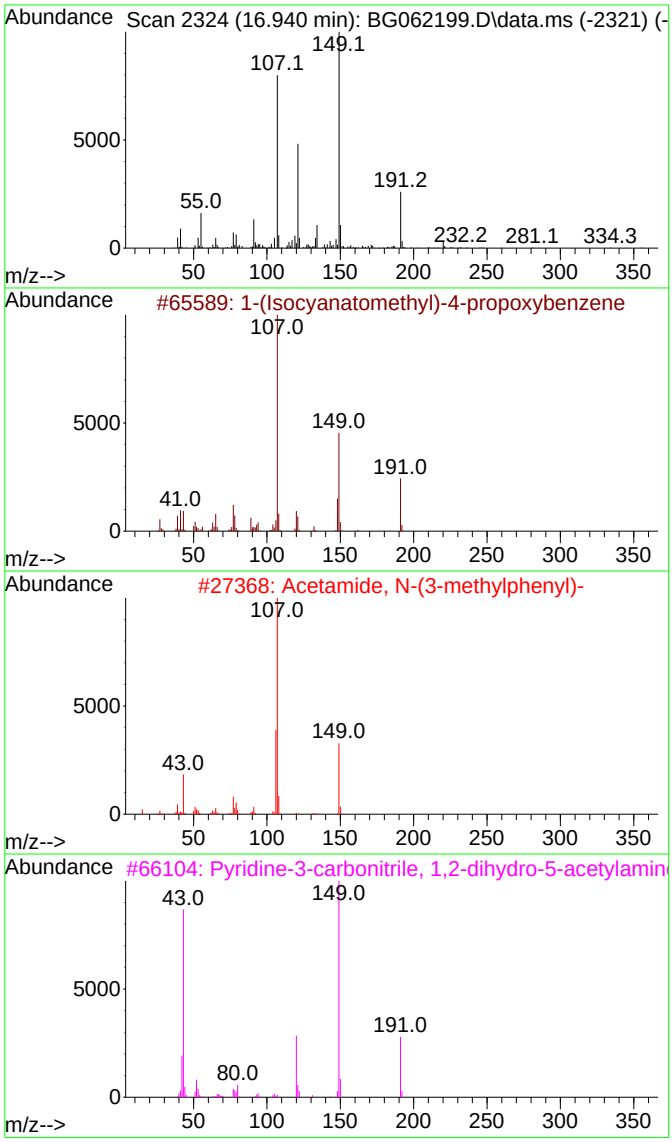
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

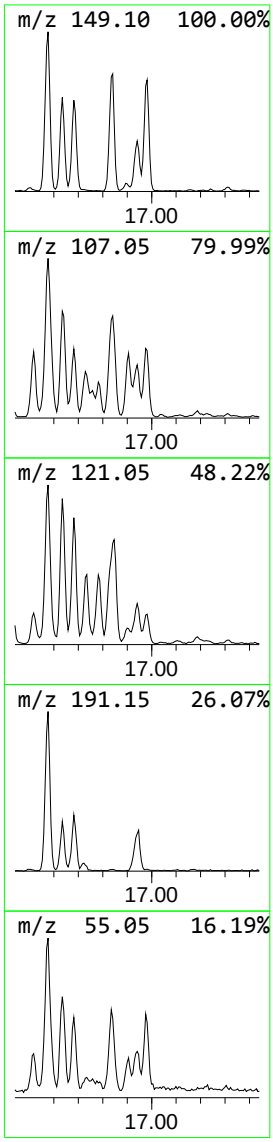
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 60 1-(Isocyanatomethyl)-4-prop... Concentration Rank 19

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 16.941 | 28.41 ng/ul | 1116490 | Phenanthrene-d10 | 17.446 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1-(Isocyanatomethyl)-4-propoxybe... | 191 | C11H13NO2 | 936095-07-7 | 72 |
| 2 | Acetamide, N-(3-methylphenyl)- | 149 | C9H11NO | 000537-92-8 | 64 |
| 3 | Pyridine-3-carbonitrile, 1,2-dih... | 191 | C9H9N3O2 | 220098-92-0 | 43 |
| 4 | Phthalic acid, 2,3-dimethylpheny... | 312 | C19H20O4 | 1000357-08-5 | 38 |
| 5 | Phthalic acid, 2-methylphenyl pr... | 298 | C18H18O4 | 1000314-95-1 | 38 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072324\
Data File : BG062199.D
Acq On : 24 Jul 2024 00:03
Operator : MA/JU
Sample : P3244-04
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
YDZD5

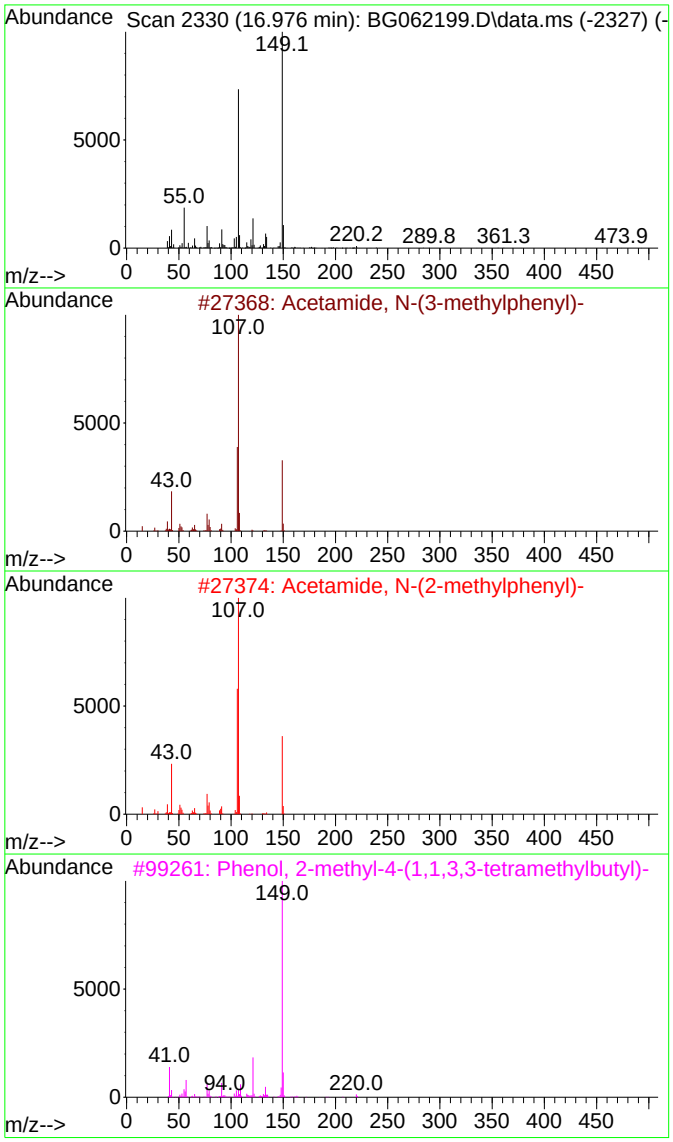
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

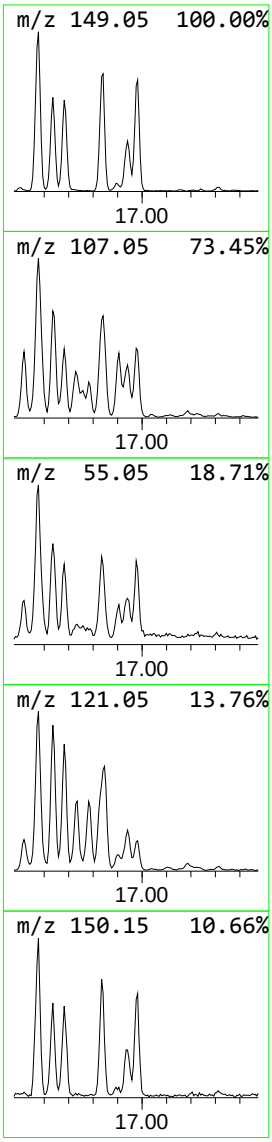
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 61 unknown-05 Concentration Rank 14

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 16.976 | 35.70 ng/ul | 1402990 | Phenanthrene-d10 | 17.446 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Acetamide, N-(3-methylphenyl)- | 149 | C9H11NO | 000537-92-8 | 50 |
| 2 | Acetamide, N-(2-methylphenyl)- | 149 | C9H11NO | 000120-66-1 | 49 |
| 3 | Phenol, 2-methyl-4-(1,1,3,3-tetr... | 220 | C15H24O | 002219-84-3 | 49 |
| 4 | Phenol, 2-methyl-4-(1,1,3,3-tetr... | 220 | C15H24O | 002219-84-3 | 43 |
| 5 | 3-(4-Hydroxyphenyl)propionic aci... | 263 | C13H13NO5 | 034071-95-9 | 43 |



Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072324\
Data File : BG062199.D
Acq On : 24 Jul 2024 00:03
Operator : MA/JU
Sample : P3244-04
Misc :
ALS Vial : 19 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
YDZD5

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071924.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

| | | | | | --Internal Standard--- | | | |
|---|---|---|---|---|---|---|---|---|
| TIC Top Hit name | RT | EstConc | Units | Response | # | RT | Resp | Conc |
| Benzene, 1,3-di... | 5.685 | 32.1 | ng/ul | 604023 | 1 | 8.070 | 375813 | 20.0 |
| Benzene, 1-ethy... | 7.142 | 20.4 | ng/ul | 384229 | 1 | 8.070 | 375813 | 20.0 |
| Benzene, 1-ethy... | 7.201 | 11.9 | ng/ul | 223763 | 1 | 8.070 | 375813 | 20.0 |
| Benzene, 1,2,3-... | 7.447 | 15.6 | ng/ul | 293689 | 1 | 8.070 | 375813 | 20.0 |
| Eucalyptol | 8.364 | 37.3 | ng/ul | 701088 | 1 | 8.070 | 375813 | 20.0 |
| 2,6-Dimethyl-1,... | 8.693 | 31.4 | ng/ul | 589494 | 1 | 8.070 | 375813 | 20.0 |
| Benzene, 1-ethy... | 9.169 | 18.2 | ng/ul | 341592 | 1 | 8.070 | 375813 | 20.0 |
| Fenchone | 9.310 | 70.3 | ng/ul | 1320990 | 1 | 8.070 | 375813 | 20.0 |
| (DEL) Alkane: C... | 9.891 | 5.0 | ng/ul | 178800 | 2 | 10.884 | 712064 | 20.0 |
| Cyclohexanemeth... | 10.249 | 113.5 | ng/ul | 4040470 | 2 | 10.884 | 712064 | 20.0 |
| Bicyclo[2.2.1]h... | 10.297 | 59.8 | ng/ul | 2129350 | 2 | 10.884 | 712064 | 20.0 |
| Phenol, 3,5-dim... | 10.379 | 24.7 | ng/ul | 878906 | 2 | 10.884 | 712064 | 20.0 |
| Phenol, 3,4-dim... | 10.767 | 14.4 | ng/ul | 512368 | 2 | 10.884 | 712064 | 20.0 |
| unknown-01 | 11.242 | 16.1 | ng/ul | 571593 | 2 | 10.884 | 712064 | 20.0 |
| Naphthalene, 2,... | 13.856 | 16.8 | ng/ul | 614866 | 3 | 14.697 | 730978 | 20.0 |
| Naphthalene, 1,... | 14.021 | 27.4 | ng/ul | 1002130 | 3 | 14.697 | 730978 | 20.0 |
| Naphthalene, 1,... | 14.256 | 11.0 | ng/ul | 403020 | 3 | 14.697 | 730978 | 20.0 |
| Diethyltoluamide | 15.454 | 35.7 | ng/ul | 1304800 | 3 | 14.697 | 730978 | 20.0 |
| 2-Propylphenol,... | 15.525 | 13.8 | ng/ul | 502716 | 3 | 14.697 | 730978 | 20.0 |
| 1-(4-Pyridinyl)... | 16.424 | 29.2 | ng/ul | 1146580 | 4 | 17.446 | 785901 | 20.0 |
| Phenol, 2-(1,1,... | 16.517 | 59.6 | ng/ul | 2343180 | 4 | 17.446 | 785901 | 20.0 |
| unknown-02 | 16.576 | 118.8 | ng/ul | 4666090 | 4 | 17.446 | 785901 | 20.0 |
| 4-(7-Methylocty... | 16.635 | 81.8 | ng/ul | 3216080 | 4 | 17.446 | 785901 | 20.0 |
| Acetamide, N-(3... | 16.682 | 37.6 | ng/ul | 1476900 | 4 | 17.446 | 785901 | 20.0 |
| unknown-03 | 16.729 | 25.5 | ng/ul | 1002000 | 4 | 17.446 | 785901 | 20.0 |
| Hexestrol, O-tr... | 16.758 | 17.7 | ng/ul | 695686 | 4 | 17.446 | 785901 | 20.0 |
| unknown-04 | 16.782 | 15.7 | ng/ul | 617241 | 4 | 17.446 | 785901 | 20.0 |
| Acetamide, N-(2... | 16.841 | 68.1 | ng/ul | 2674560 | 4 | 17.446 | 785901 | 20.0 |
| Phenol, 4-(1,1-... | 16.905 | 53.2 | ng/ul | 2089020 | 4 | 17.446 | 785901 | 20.0 |
| 1-(Isocyanatome... | 16.941 | 28.4 | ng/ul | 1116490 | 4 | 17.446 | 785901 | 20.0 |
| unknown-05 | 16.976 | 35.7 | ng/ul | 1402990 | 4 | 17.446 | 785901 | 20.0 |