LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062424\
Data File : BG061731.D
Acq On : 26 Jun 2024 11:43
Operator : MA/JU
Sample : P2873-09DL 10X
Misc :
ALS Vial : 30 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** C0SJ9DL

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-BG062024.MA.M
Title : SVOA CALIBRATION

Signal : TIC: BG061731.D\data.ms

| peak # | R.T. min | first scan | max scan | last scan | PK TY | peak height | corr. area | corr. % max. | % of total |
|---|---|---|---|---|---|---|---|---|---|
| 1 | 4.225 | 155 | 159 | 166 | rVB | 39326 | 60696 | 1.10% | 0.137% |
| 2 | 4.759 | 243 | 250 | 258 | rBV | 574929 | 920905 | 16.75% | 2.082% |
| 3 | 5.564 | 381 | 387 | 396 | rBV | 151720 | 237649 | 4.32% | 0.537% |
| 4 | 5.699 | 402 | 410 | 420 | rBV | 127647 | 292837 | 5.33% | 0.662% |
| 5 | 6.070 | 464 | 473 | 481 | rVB4 | 89619 | 196255 | 3.57% | 0.444% |
| 6 | 7.151 | 652 | 657 | 662 | rBV2 | 38706 | 56929 | 1.04% | 0.129% |
| 7 | 7.215 | 663 | 668 | 673 | rVB4 | 65231 | 112157 | 2.04% | 0.254% |
| 8 | 7.286 | 675 | 680 | 686 | rBV2 | 45940 | 73945 | 1.35% | 0.167% |
| 9 | 7.721 | 748 | 754 | 763 | rVV2 | 277665 | 529629 | 9.64% | 1.197% |
| 10 | 7.809 | 763 | 769 | 775 | rVB3 | 54974 | 94474 | 1.72% | 0.214% |
| 11 | 8.073 | 803 | 814 | 822 | rBV | 229166 | 399993 | 7.28% | 0.904% |
| 12 | 8.185 | 826 | 833 | 839 | rVV3 | 43610 | 84592 | 1.54% | 0.191% |
| 13 | 8.249 | 839 | 844 | 851 | rVB2 | 32260 | 53456 | 0.97% | 0.121% |
| 14 | 8.614 | 895 | 906 | 913 | rVV | 464478 | 833918 | 15.17% | 1.885% |
| 15 | 9.236 | 1007 | 1012 | 1016 | rBV3 | 26771 | 46890 | 0.85% | 0.106% |
| 16 | 9.342 | 1025 | 1030 | 1039 | rVV3 | 54582 | 104266 | 1.90% | 0.236% |
| 17 | 9.759 | 1097 | 1101 | 1108 | rVB2 | 31051 | 54439 | 0.99% | 0.123% |
| 18 | 9.842 | 1110 | 1115 | 1121 | rBV2 | 39189 | 63747 | 1.16% | 0.144% |
| 19 | 10.306 | 1183 | 1194 | 1200 | rBV3 | 140624 | 344187 | 6.26% | 0.778% |
| 20 | 10.376 | 1200 | 1206 | 1210 | rBV | 104814 | 205120 | 3.73% | 0.464% |
| 21 | 10.494 | 1221 | 1226 | 1232 | rVB3 | 35122 | 65818 | 1.20% | 0.149% |
| 22 | 10.588 | 1236 | 1242 | 1247 | rVB5 | 26209 | 50035 | 0.91% | 0.113% |
| 23 | 10.887 | 1281 | 1293 | 1296 | rBV2 | 321148 | 656801 | 11.95% | 1.485% |
| 24 | 10.940 | 1296 | 1302 | 1310 | rVV | 3028481 | 5496555 | 100.00% | 12.425% |
| 25 | 11.052 | 1317 | 1321 | 1327 | rVB | 59029 | 108048 | 1.97% | 0.244% |
| 26 | 12.192 | 1507 | 1515 | 1520 | rBV8 | 24197 | 53491 | 0.97% | 0.121% |
| 27 | 12.292 | 1525 | 1532 | 1537 | rVV6 | 25109 | 57110 | 1.04% | 0.129% |
| 28 | 12.392 | 1543 | 1549 | 1552 | rVV6 | 24918 | 54031 | 0.98% | 0.122% |
| 29 | 12.433 | 1552 | 1556 | 1561 | rVV4 | 37759 | 68006 | 1.24% | 0.154% |
| 30 | 12.538 | 1565 | 1574 | 1582 | rVB | 1469436 | 2440007 | 44.39% | 5.516% |
| 31 | 12.650 | 1587 | 1593 | 1598 | rVV3 | 25791 | 51091 | 0.93% | 0.115% |
| 32 | 12.756 | 1598 | 1611 | 1620 | rVB | 888247 | 1530029 | 27.84% | 3.459% |
| 33 | 13.537 | 1736 | 1744 | 1750 | rBV | 195944 | 333355 | 6.06% | 0.754% |
| 34 | 13.678 | 1763 | 1768 | 1772 | rBV6 | 31569 | 62123 | 1.13% | 0.140% |
| 35 | 13.731 | 1772 | 1777 | 1780 | rBV2 | 55017 | 81394 | 1.48% | 0.184% |
| 36 | 13.866 | 1793 | 1800 | 1801 | rBV2 | 205099 | 312623 | 5.69% | 0.707% |

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062424\
Data File : BG061731.D
Acq On : 26 Jun 2024 11:43
Operator : MA/JU
Sample : P2873-09DL 10X
Misc :
ALS Vial : 30 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** C0SJ9DL

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-BG062024.MA.M
Title : SVOA CALIBRATION

| | | | | | | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 37 | 13.943 | 1809 | 1813 | 1819 | rVB6 | 37665 | 56765 | 1.03% | 0.128% |
| 38 | 14.019 | 1819 | 1826 | 1831 | rVV2 | 362602 | 657174 | 11.96% | 1.486% |
| 39 | 14.072 | 1831 | 1835 | 1844 | rVV2 | 212835 | 477758 | 8.69% | 1.080% |
| 40 | 14.154 | 1844 | 1849 | 1856 | rVB | 269056 | 425906 | 7.75% | 0.963% |
| 41 | 14.260 | 1861 | 1867 | 1870 | rBV | 164372 | 261244 | 4.75% | 0.591% |
| 42 | 14.424 | 1884 | 1895 | 1902 | rBV | 805170 | 1414514 | 25.73% | 3.198% |
| 43 | 14.701 | 1930 | 1942 | 1947 | rBV2 | 592395 | 1087658 | 19.79% | 2.459% |
| 44 | 14.765 | 1947 | 1953 | 1960 | rVV4 | 85004 | 200371 | 3.65% | 0.453% |
| 45 | 14.906 | 1972 | 1977 | 1986 | rVB5 | 44170 | 106436 | 1.94% | 0.241% |
| 46 | 15.094 | 2005 | 2009 | 2017 | rVB2 | 117918 | 185811 | 3.38% | 0.420% |
| 47 | 15.171 | 2017 | 2022 | 2026 | rVB | 77096 | 109808 | 2.00% | 0.248% |
| 48 | 15.300 | 2037 | 2044 | 2049 | rBV2 | 129838 | 286713 | 5.22% | 0.648% |
| 49 | 15.365 | 2052 | 2055 | 2058 | rVB2 | 56012 | 67401 | 1.23% | 0.152% |
| 50 | 15.459 | 2066 | 2071 | 2073 | rBV2 | 56612 | 81482 | 1.48% | 0.184% |
| 51 | 15.564 | 2085 | 2089 | 2095 | rVB2 | 144907 | 234112 | 4.26% | 0.529% |
| 52 | 15.688 | 2106 | 2110 | 2115 | rVV2 | 225817 | 344561 | 6.27% | 0.779% |
| 53 | 15.746 | 2115 | 2120 | 2132 | rVB2 | 529062 | 990924 | 18.03% | 2.240% |
| 54 | 15.858 | 2132 | 2139 | 2144 | rBV6 | 48710 | 134025 | 2.44% | 0.303% |
| 55 | 15.946 | 2150 | 2154 | 2160 | rBV4 | 64107 | 117604 | 2.14% | 0.266% |
| 56 | 16.034 | 2164 | 2169 | 2174 | rBV2 | 50332 | 89991 | 1.64% | 0.203% |
| 57 | 16.158 | 2184 | 2190 | 2203 | rVB2 | 192163 | 388997 | 7.08% | 0.879% |
| 58 | 16.405 | 2228 | 2232 | 2243 | rVB4 | 62693 | 129942 | 2.36% | 0.294% |
| 59 | 16.739 | 2282 | 2289 | 2295 | rBV3 | 137966 | 307667 | 5.60% | 0.695% |
| 60 | 16.792 | 2295 | 2298 | 2304 | rVB2 | 92819 | 141228 | 2.57% | 0.319% |
| 61 | 16.892 | 2311 | 2315 | 2320 | rBV2 | 82300 | 125433 | 2.28% | 0.284% |
| 62 | 17.104 | 2347 | 2351 | 2357 | rVB2 | 68421 | 103555 | 1.88% | 0.234% |
| 63 | 17.262 | 2371 | 2378 | 2383 | rBV | 220247 | 366524 | 6.67% | 0.829% |
| 64 | 17.444 | 2403 | 2409 | 2412 | rBV | 720531 | 1118053 | 20.34% | 2.527% |
| 65 | 17.486 | 2412 | 2416 | 2422 | rVV | 2108964 | 3245097 | 59.04% | 7.336% |
| 66 | 17.538 | 2422 | 2425 | 2427 | rVV | 134260 | 179157 | 3.26% | 0.405% |
| 67 | 17.574 | 2427 | 2431 | 2437 | rVV | 554778 | 842139 | 15.32% | 1.904% |
| 68 | 17.627 | 2437 | 2440 | 2445 | rVV3 | 41404 | 77544 | 1.41% | 0.175% |
| 69 | 18.020 | 2503 | 2507 | 2514 | rVB2 | 122681 | 185547 | 3.38% | 0.419% |
| 70 | 18.167 | 2527 | 2532 | 2539 | rVB2 | 79183 | 169000 | 3.07% | 0.382% |
| 71 | 18.320 | 2552 | 2558 | 2562 | rBV | 429767 | 650436 | 11.83% | 1.470% |
| 72 | 18.367 | 2562 | 2566 | 2573 | rVB | 468038 | 687892 | 12.51% | 1.555% |
| 73 | 18.449 | 2575 | 2580 | 2584 | rBV | 191309 | 266971 | 4.86% | 0.604% |
| 74 | 18.508 | 2584 | 2590 | 2595 | rBV3 | 499975 | 1050215 | 19.11% | 2.374% |
| 75 | 18.625 | 2606 | 2610 | 2613 | rBV5 | 37186 | 51305 | 0.93% | 0.116% |
| 76 | 18.813 | 2637 | 2642 | 2646 | rBV | 202260 | 287030 | 5.22% | 0.649% |

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062424\
Data File : BG061731.D
Acq On : 26 Jun 2024 11:43
Operator : MA/JU
Sample : P2873-09DL 10X
Misc :
ALS Vial : 30 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** C0SJ9DL

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-BG062024.MA.M
Title : SVOA CALIBRATION

| | | | | | | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 77 | 19.054 | 2681 | 2683 | 2688 | rVB6 | 41726 | 52067 | 0.95% | 0.118% |
| 78 | 19.107 | 2689 | 2692 | 2695 | rVB | 64470 | 63099 | 1.15% | 0.143% |
| 79 | 19.142 | 2696 | 2698 | 2703 | rVB2 | 65958 | 73881 | 1.34% | 0.167% |
| 80 | 19.225 | 2707 | 2712 | 2717 | rBV2 | 224325 | 378632 | 6.89% | 0.856% |
| 81 | 19.366 | 2733 | 2736 | 2738 | rBV4 | 37681 | 54657 | 0.99% | 0.124% |
| 82 | 19.489 | 2751 | 2757 | 2764 | rBV | 936458 | 1295052 | 23.56% | 2.928% |
| 83 | 19.548 | 2764 | 2767 | 2770 | rVV3 | 50393 | 70032 | 1.27% | 0.158% |
| 84 | 19.583 | 2770 | 2773 | 2777 | rVV4 | 51634 | 75624 | 1.38% | 0.171% |
| 85 | 19.636 | 2778 | 2782 | 2793 | rVB | 339561 | 526057 | 9.57% | 1.189% |
| 86 | 19.853 | 2809 | 2819 | 2826 | rVB | 1057658 | 1833925 | 33.36% | 4.146% |
| 87 | 19.924 | 2828 | 2831 | 2836 | rVB4 | 44881 | 63616 | 1.16% | 0.144% |
| 88 | 20.171 | 2868 | 2873 | 2880 | rBV3 | 109876 | 255535 | 4.65% | 0.578% |
| 89 | 20.312 | 2893 | 2897 | 2898 | rBV3 | 82065 | 120236 | 2.19% | 0.272% |
| 90 | 20.335 | 2899 | 2901 | 2909 | rVB | 269568 | 390023 | 7.10% | 0.882% |
| 91 | 20.441 | 2915 | 2919 | 2923 | rVB | 225843 | 291125 | 5.30% | 0.658% |
| 92 | 20.500 | 2925 | 2929 | 2933 | rBV | 159952 | 196822 | 3.58% | 0.445% |
| 93 | 20.635 | 2949 | 2952 | 2956 | rBV | 104006 | 158120 | 2.88% | 0.357% |
| 94 | 20.682 | 2956 | 2960 | 2967 | rVB | 196686 | 302801 | 5.51% | 0.684% |
| 95 | 21.328 | 3067 | 3070 | 3074 | rVB2 | 83335 | 100807 | 1.83% | 0.228% |
| 96 | 21.704 | 3125 | 3134 | 3139 | rBV3 | 863884 | 2047683 | 37.25% | 4.629% |
| 97 | 21.751 | 3139 | 3142 | 3150 | rVB | 351630 | 575239 | 10.47% | 1.300% |
| 98 | 22.756 | 3311 | 3313 | 3322 | rVB3 | 99427 | 155285 | 2.83% | 0.351% |
| 99 | 23.907 | 3505 | 3509 | 3516 | rBV4 | 76618 | 210207 | 3.82% | 0.475% |
| 100 | 24.930 | 3676 | 3683 | 3693 | rVB2 | 397921 | 1055818 | 19.21% | 2.387% |

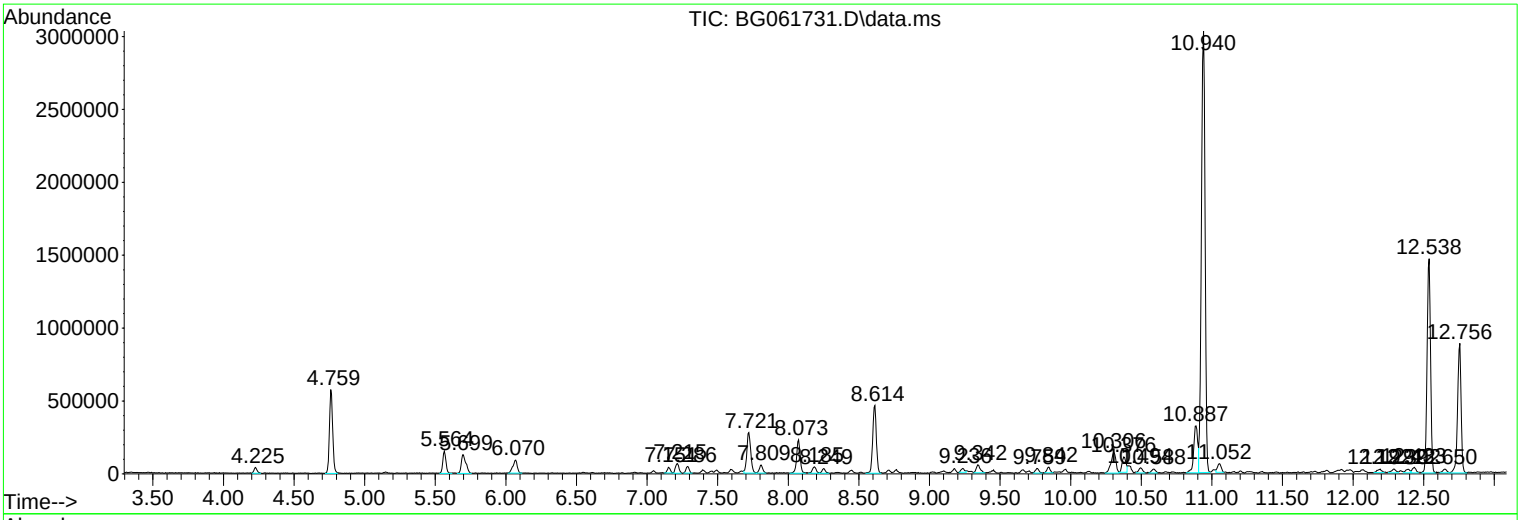
Sum of corrected areas: 44236934

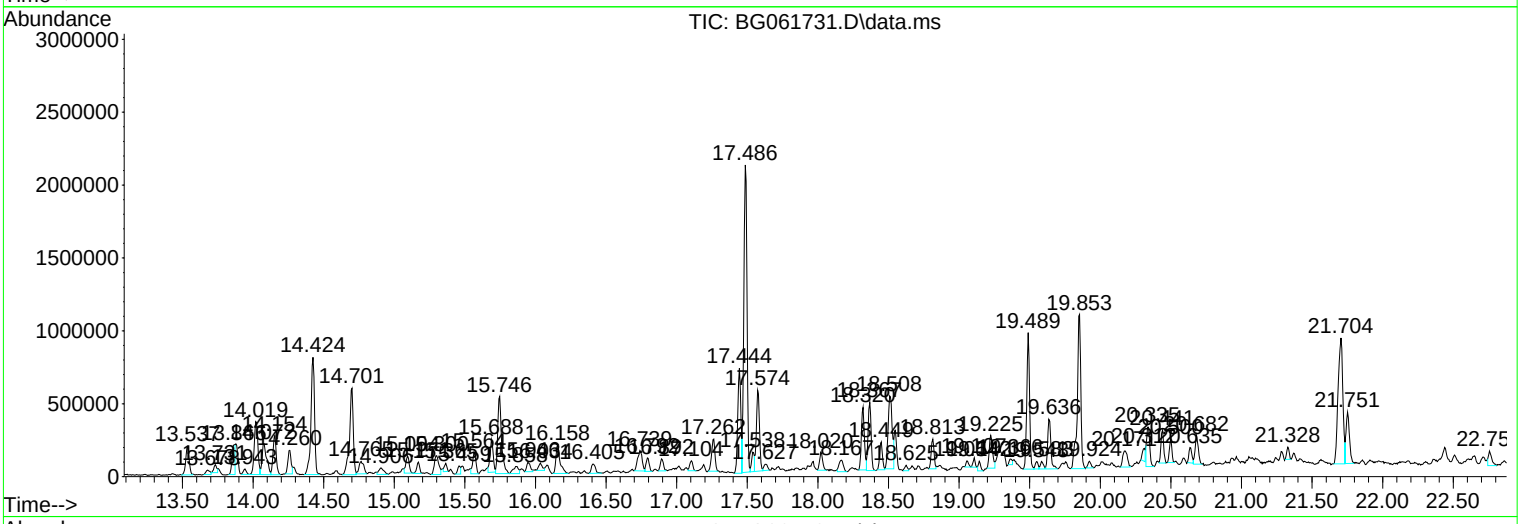
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062424\
Data File : BG061731.D
Acq On : 26 Jun 2024 11:43
Operator : MA/JU
Sample : P2873-09DL 10X
Misc :
ALS Vial : 30 Sample Multiplier: 1

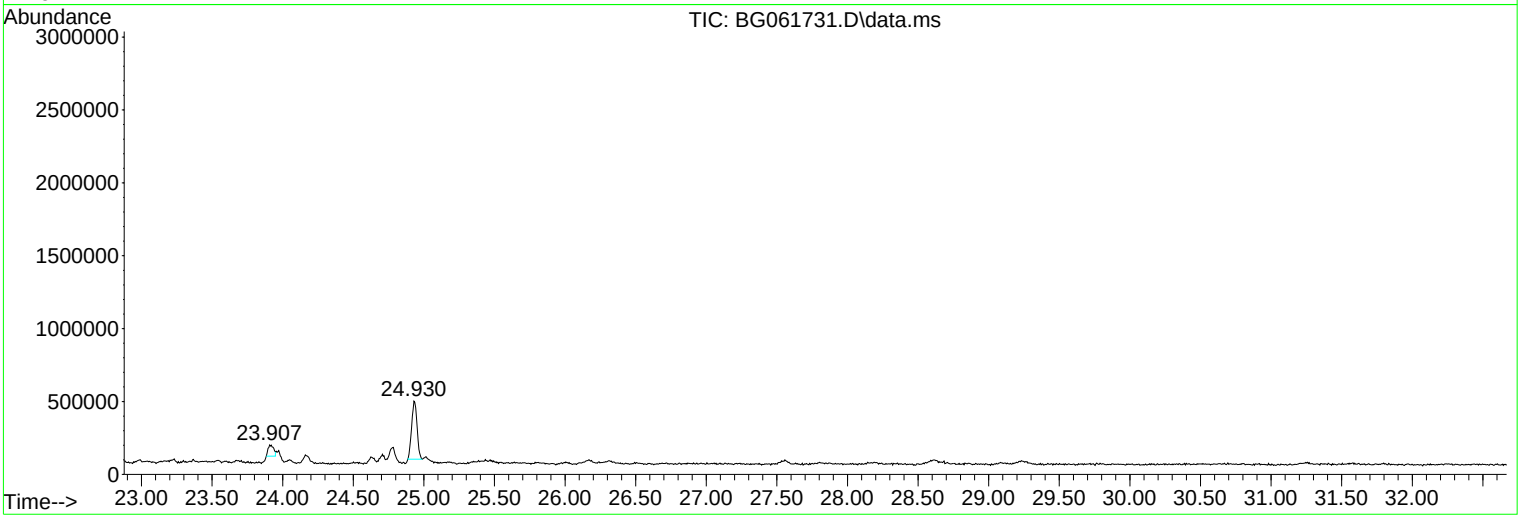
**Instrument :**
BNA_G
**ClientSampleId :**
C0SJ9DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.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\BG062424\
Data File : BG061731.D
Acq On : 26 Jun 2024 11:43
Operator : MA/JU
Sample : P2873-09DL 10X
Misc :
ALS Vial : 30 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
C0SJ9DL

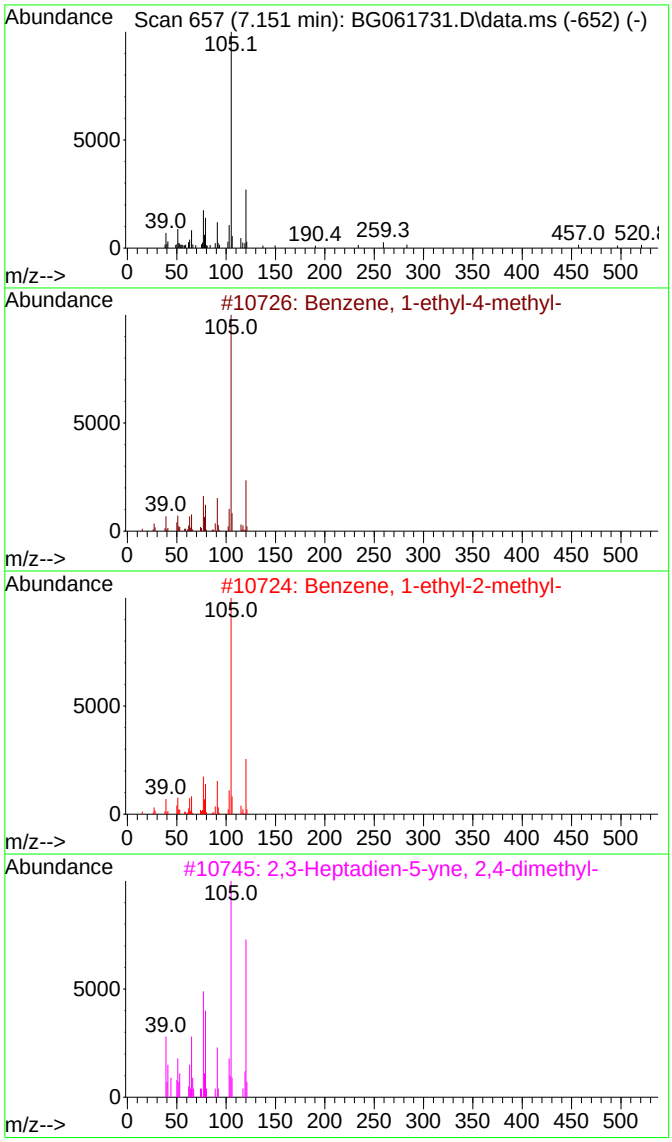
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

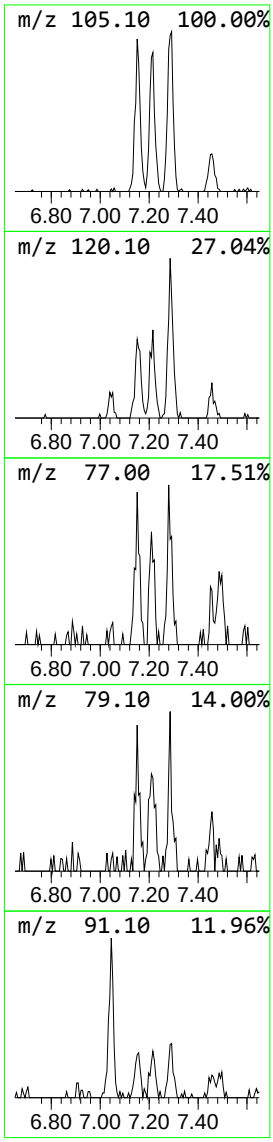
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 6 Benzene, 1-ethyl-4-methyl- Concentration Rank 34

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 7.151 | 2.85 ng/ul | 56929 | 1,4-Dichlorobenzene-d4 | 8.073 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzene, 1-ethyl-4-methyl- | 120 | C9H12 | 000622-96-8 | 87 |
| 2 | Benzene, 1-ethyl-2-methyl- | 120 | C9H12 | 000611-14-3 | 87 |
| 3 | 2,3-Heptadien-5-yne, 2,4-dimethyl- | 120 | C9H12 | 041898-89-9 | 76 |
| 4 | Benzene, 1-ethyl-2-methyl- | 120 | C9H12 | 000611-14-3 | 74 |
| 5 | Benzene, (1-methylethyl)- | 120 | C9H12 | 000098-82-8 | 72 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062424\
Data File : BG061731.D
Acq On : 26 Jun 2024 11:43
Operator : MA/JU
Sample : P2873-09DL 10X
Misc :
ALS Vial : 30 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
C0SJ9DL

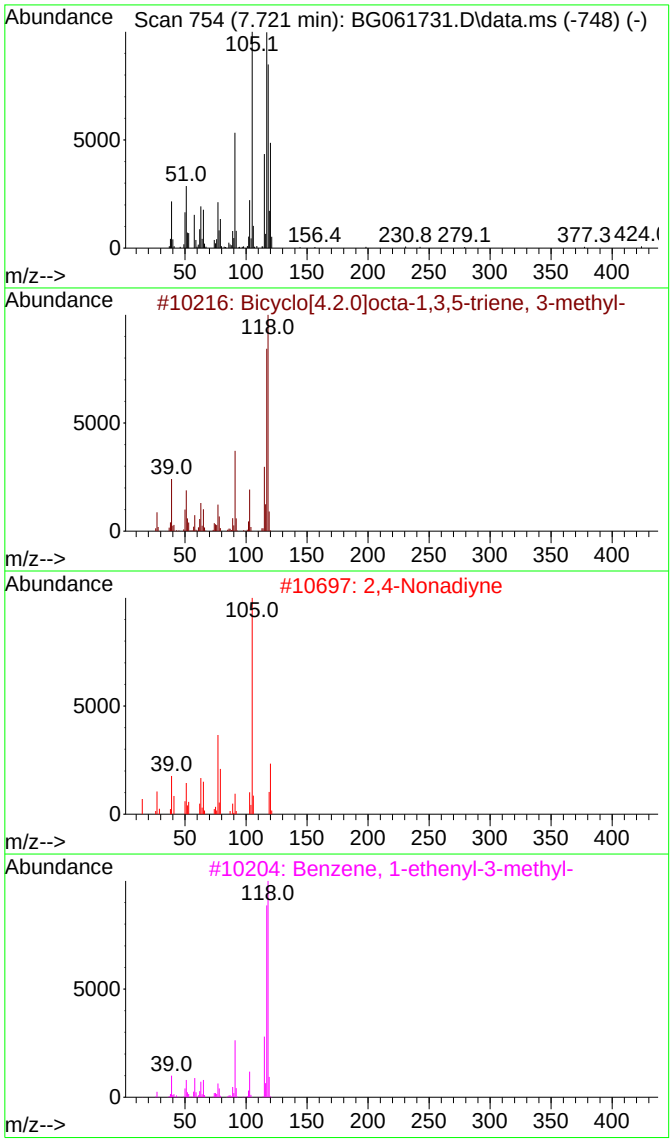
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

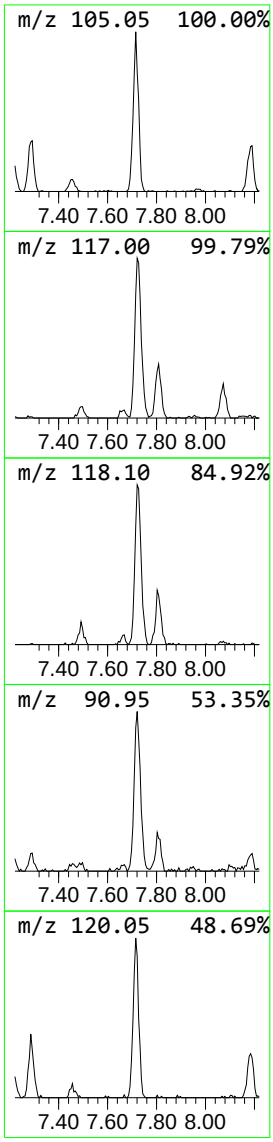
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 8 Bicyclo[4.2.0]octa-1,3,5-tr... Concentration Rank 3

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 7.721 | 26.48 ng/ul | 529629 | 1,4-Dichlorobenzene-d4 | 8.073 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Bicyclo[4.2.0]octa-1,3,5-triene,... | 118 | C9H10 | 022250-74-4 | 83 |
| 2 | 2,4-Nonadiyne | 120 | C9H12 | 063621-15-8 | 70 |
| 3 | Benzene, 1-ethenyl-3-methyl- | 118 | C9H10 | 000100-80-1 | 70 |
| 4 | Benzene, 1-ethenyl-4-methyl- | 118 | C9H10 | 000622-97-9 | 70 |
| 5 | Benzene, 1-ethenyl-4-methyl- | 118 | C9H10 | 000622-97-9 | 70 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062424\
Data File : BG061731.D
Acq On : 26 Jun 2024 11:43
Operator : MA/JU
Sample : P2873-09DL 10X
Misc :
ALS Vial : 30 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
C0SJ9DL

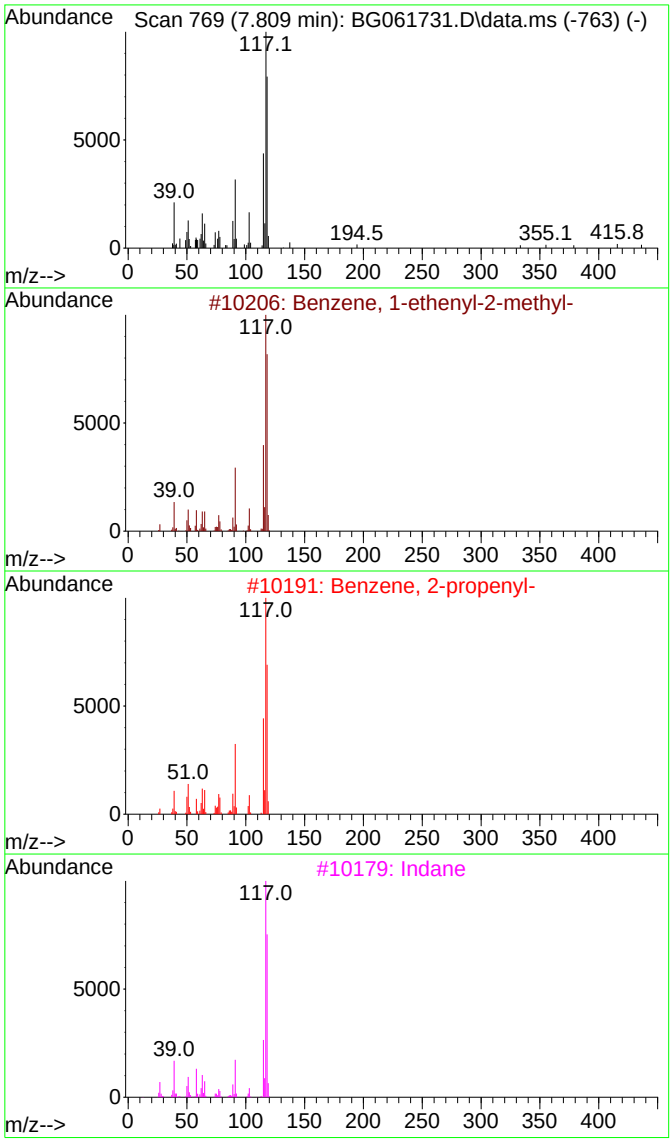
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

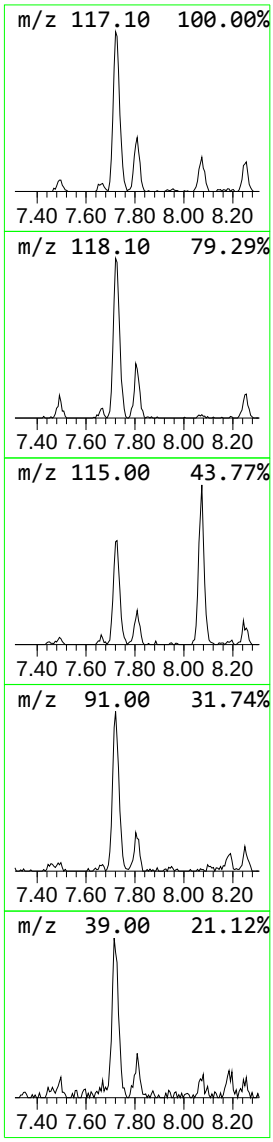
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 9 Benzene, 1-ethenyl-2-methyl- Concentration Rank 25

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 7.809 | 4.72 ng/ul | 94474 | 1,4-Dichlorobenzene-d4 | 8.073 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzene, 1-ethenyl-2-methyl- | 118 | C9H10 | 000611-15-4 | 93 |
| 2 | Benzene, 2-propenyl- | 118 | C9H10 | 000300-57-2 | 89 |
| 3 | Indane | 118 | C9H10 | 000496-11-7 | 89 |
| 4 | Bicyclo[4.2.0]octa-1,3,5-triene,... | 118 | C9H10 | 055337-80-9 | 89 |
| 5 | (Z)-1-Phenylpropene | 118 | C9H10 | 000766-90-5 | 89 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062424\
Data File : BG061731.D
Acq On : 26 Jun 2024 11:43
Operator : MA/JU
Sample : P2873-09DL 10X
Misc :
ALS Vial : 30 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
C0SJ9DL

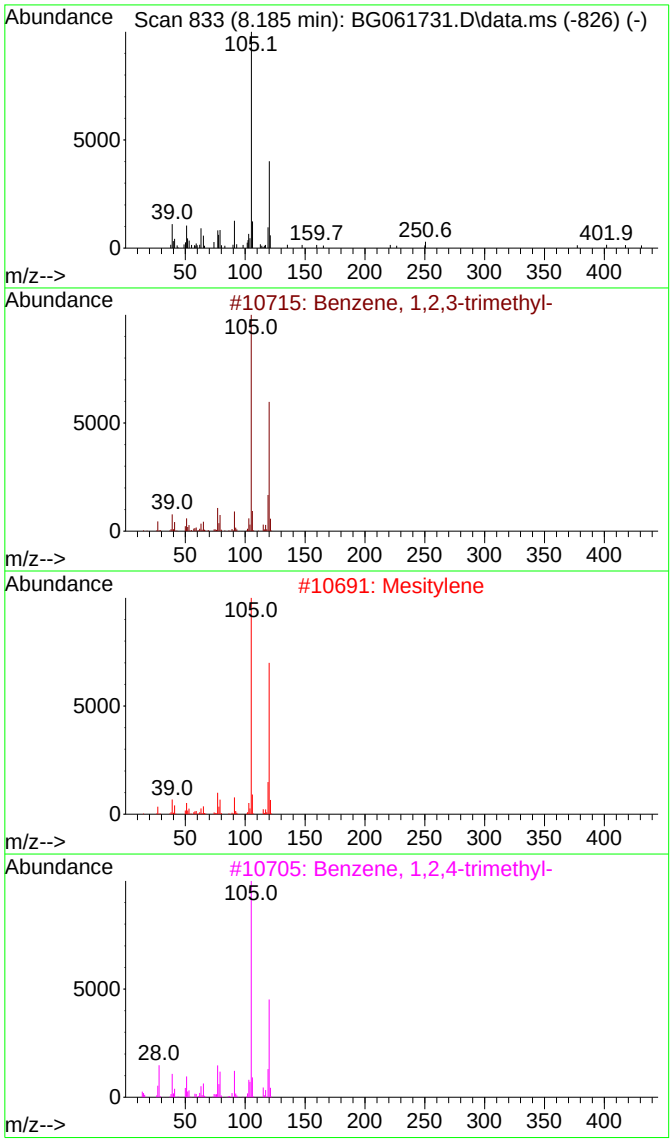
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

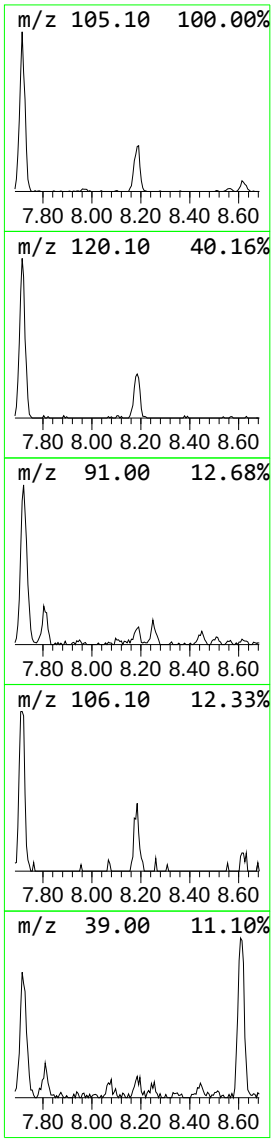
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 10 Benzene, 1,2,3-trimethyl- Concentration Rank 27

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 8.185 | 4.23 ng/ul | 84592 | 1,4-Dichlorobenzene-d4 | 8.073 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzene, 1,2,3-trimethyl- | 120 | C9H12 | 000526-73-8 | 87 |
| 2 | Mesitylene | 120 | C9H12 | 000108-67-8 | 87 |
| 3 | Benzene, 1,2,4-trimethyl- | 120 | C9H12 | 000095-63-6 | 83 |
| 4 | Benzene, 1,2,3-trimethyl- | 120 | C9H12 | 000526-73-8 | 83 |
| 5 | Mesitylene | 120 | C9H12 | 000108-67-8 | 83 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062424\
Data File : BG061731.D
Acq On : 26 Jun 2024 11:43
Operator : MA/JU
Sample : P2873-09DL 10X
Misc :
ALS Vial : 30 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
C0SJ9DL

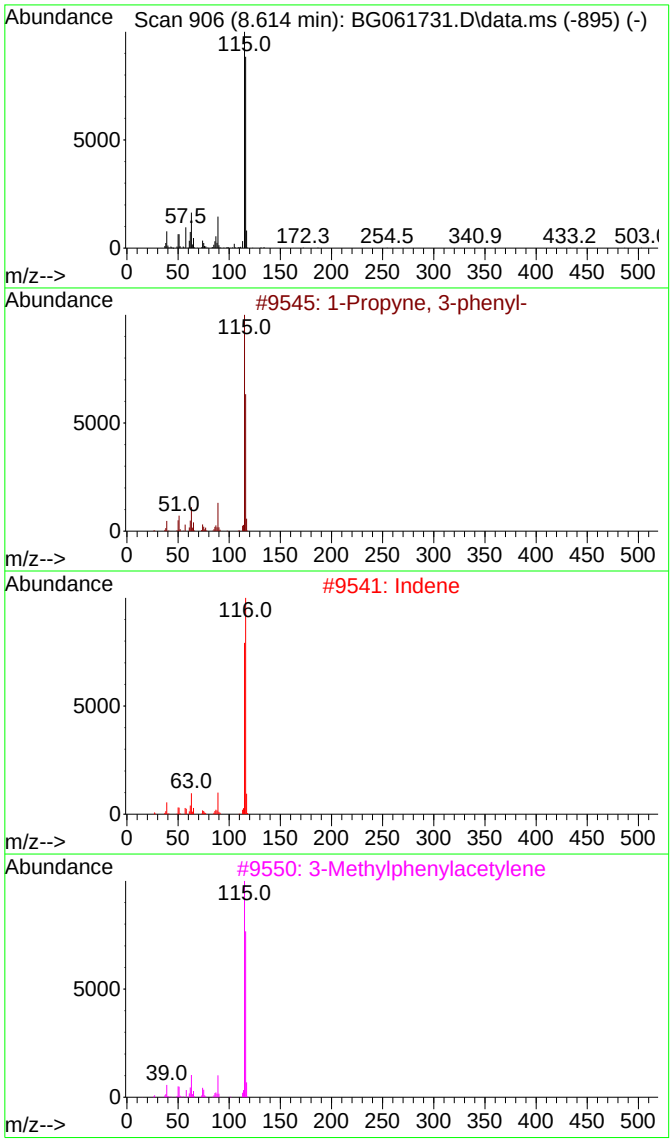
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

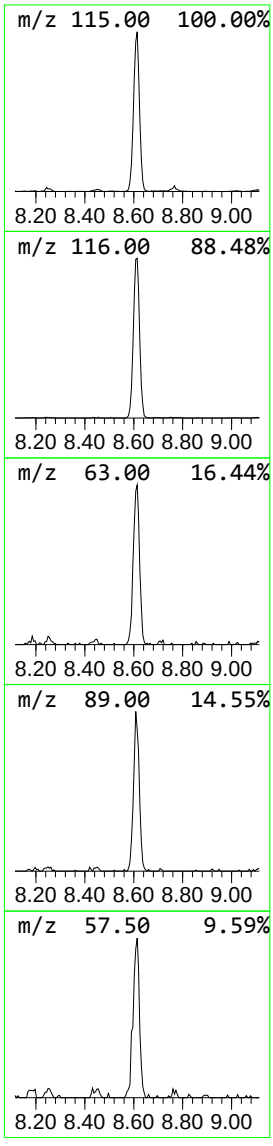
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 12 1-Propyne, 3-phenyl- Concentration Rank 2

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 8.614 | 41.70 ng/ul | 833918 | 1,4-Dichlorobenzene-d4 | 8.073 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1-Propyne, 3-phenyl- | 116 | C9H8 | 010147-11-2 | 94 |
| 2 | Indene | 116 | C9H8 | 000095-13-6 | 91 |
| 3 | 3-Methylphenylacetylene | 116 | C9H8 | 000766-82-5 | 91 |
| 4 | Indene | 116 | C9H8 | 000095-13-6 | 91 |
| 5 | Benzene, 1-propynyl- | 116 | C9H8 | 000673-32-5 | 91 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062424\
Data File : BG061731.D
Acq On : 26 Jun 2024 11:43
Operator : MA/JU
Sample : P2873-09DL 10X
Misc :
ALS Vial : 30 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
C0SJ9DL

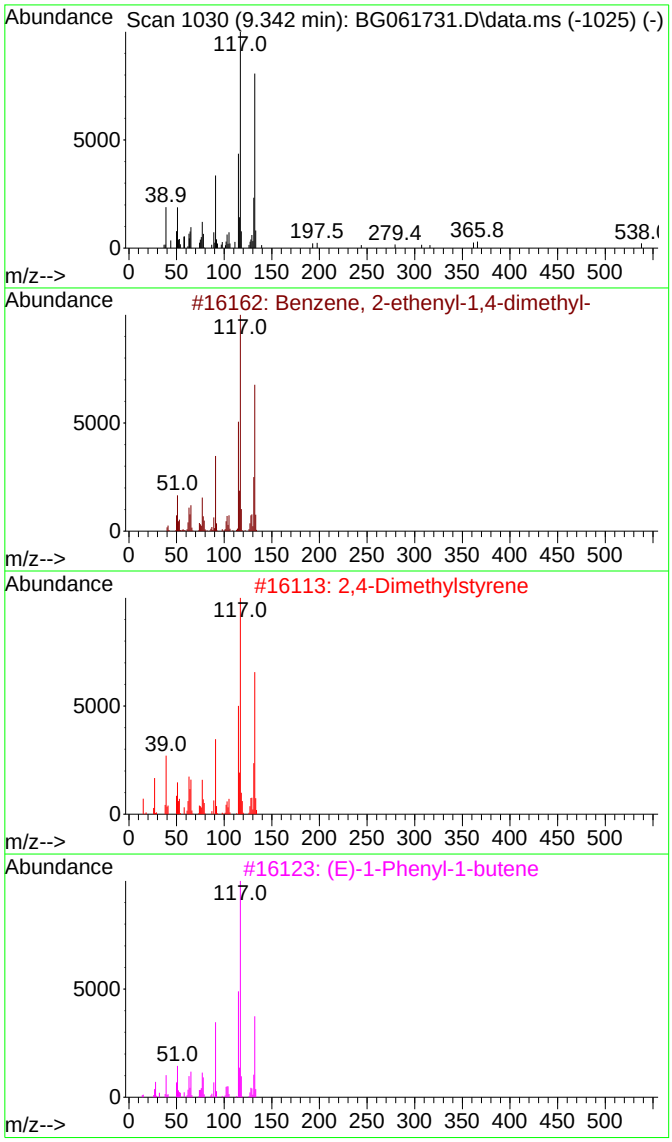
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

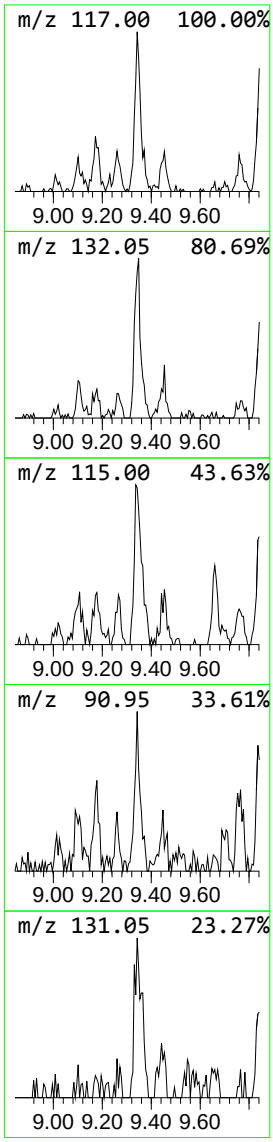
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 13 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 20

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 9.342 | 5.21 ng/ul | 104266 | 1,4-Dichlorobenzene-d4 | 8.073 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzene, 2-ethenyl-1,4-dimethyl- | 132 | C10H12 | 002039-89-6 | 95 |
| 2 | 2,4-Dimethylstyrene | 132 | C10H12 | 002234-20-0 | 95 |
| 3 | (E)-1-Phenyl-1-butene | 132 | C10H12 | 001005-64-7 | 95 |
| 4 | Benzene, (1-methyl-1-propenyl)-,... | 132 | C10H12 | 000767-99-7 | 95 |
| 5 | Benzene, (2-methyl-1-propenyl)- | 132 | C10H12 | 000768-49-0 | 95 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062424\
Data File : BG061731.D
Acq On : 26 Jun 2024 11:43
Operator : MA/JU
Sample : P2873-09DL 10X
Misc :
ALS Vial : 30 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
C0SJ9DL

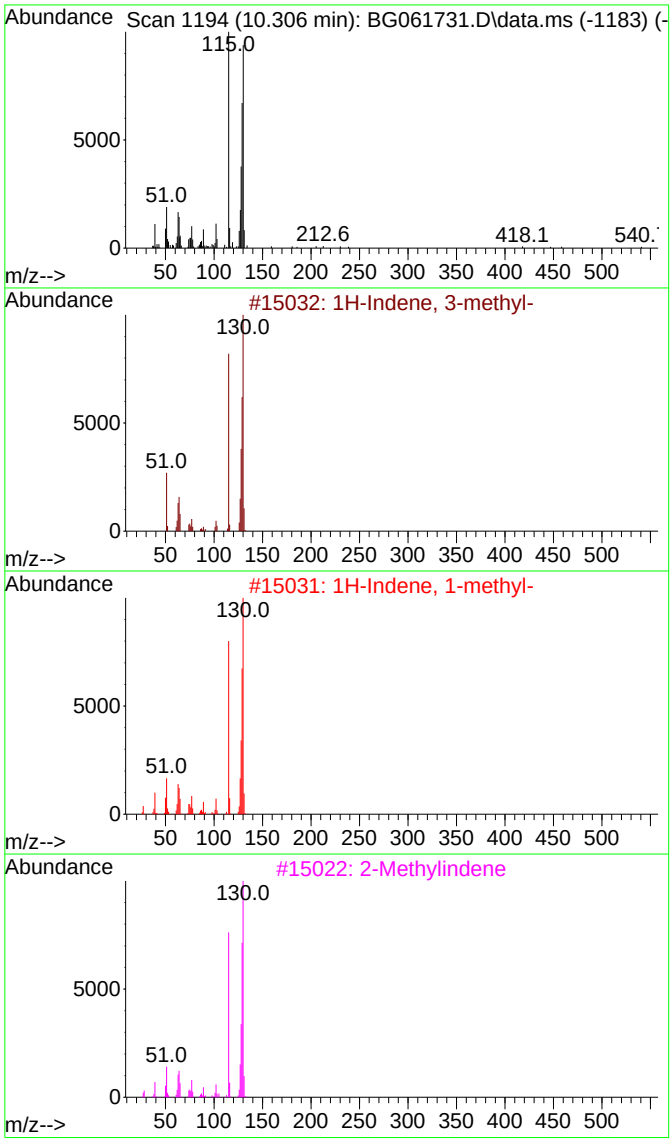
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

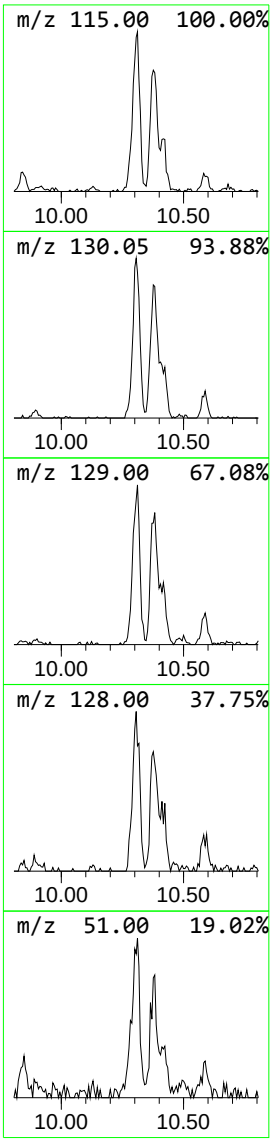
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 14 1H-Indene, 3-methyl- Concentration Rank 10

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 10.306 | 10.48 ng/ul | 344187 | Naphthalene-d8 | 10.887 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1H-Indene, 3-methyl- | 130 | C10H10 | 000767-60-2 | 94 |
| 2 | 1H-Indene, 1-methyl- | 130 | C10H10 | 000767-59-9 | 94 |
| 3 | 2-Methylindene | 130 | C10H10 | 002177-47-1 | 93 |
| 4 | 1H-Indene, 1-methyl- | 130 | C10H10 | 000767-59-9 | 93 |
| 5 | Benzene, (1-methyl-2-cyclopropen... | 130 | C10H10 | 065051-83-4 | 93 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062424\
Data File : BG061731.D
Acq On : 26 Jun 2024 11:43
Operator : MA/JU
Sample : P2873-09DL 10X
Misc :
ALS Vial : 30 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
C0SJ9DL

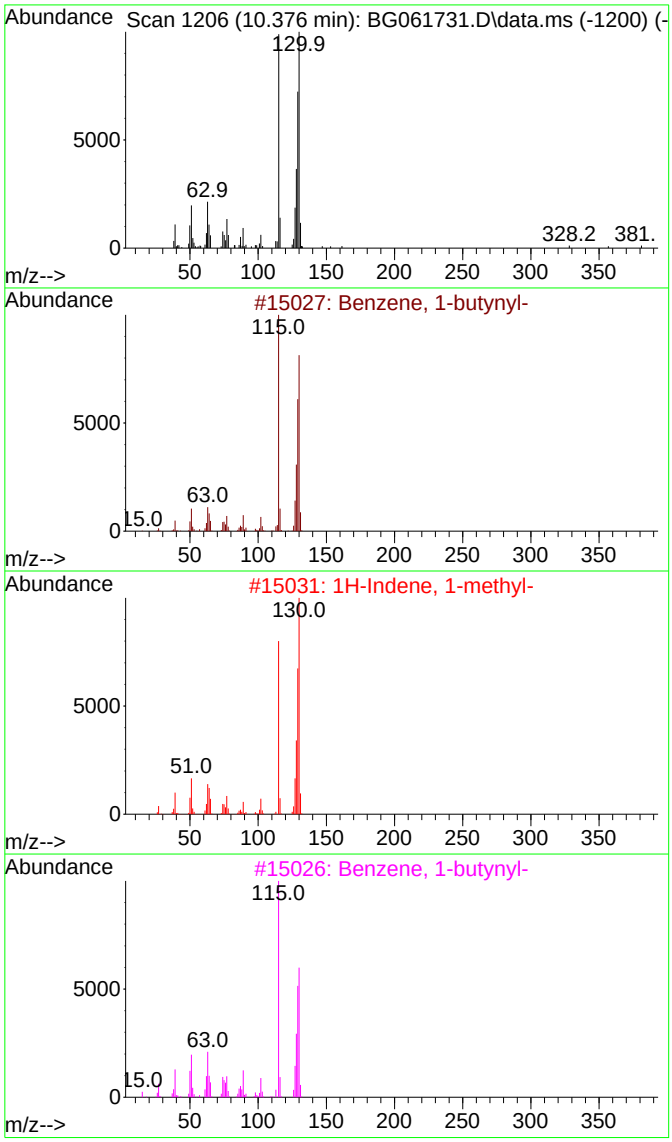
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

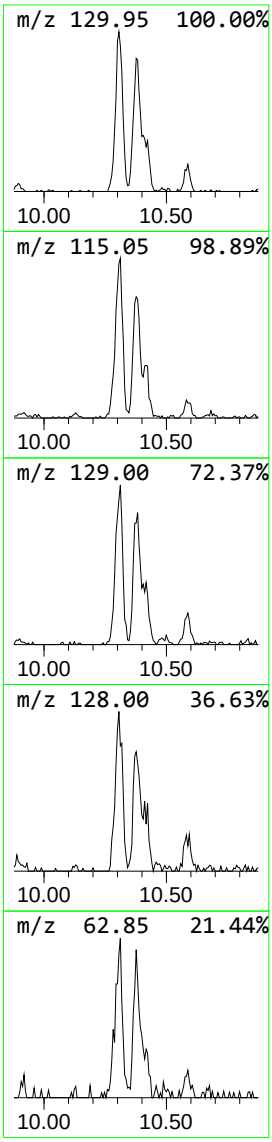
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 15 Benzene, 1-butynyl- Concentration Rank 16

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 10.376 | 6.25 ng/ul | 205120 | Naphthalene-d8 | 10.887 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzene, 1-butynyl- | 130 | C10H10 | 000622-76-4 | 95 |
| 2 | 1H-Indene, 1-methyl- | 130 | C10H10 | 000767-59-9 | 94 |
| 3 | Benzene, 1-butynyl- | 130 | C10H10 | 000622-76-4 | 93 |
| 4 | Benzene,1-methyl-1,2-propadienyl- | 130 | C10H10 | 022433-39-2 | 91 |
| 5 | Benzene, (1-methyl-2-cyclopropen... | 130 | C10H10 | 065051-83-4 | 91 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062424\
Data File : BG061731.D
Acq On : 26 Jun 2024 11:43
Operator : MA/JU
Sample : P2873-09DL 10X
Misc :
ALS Vial : 30 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
C0SJ9DL

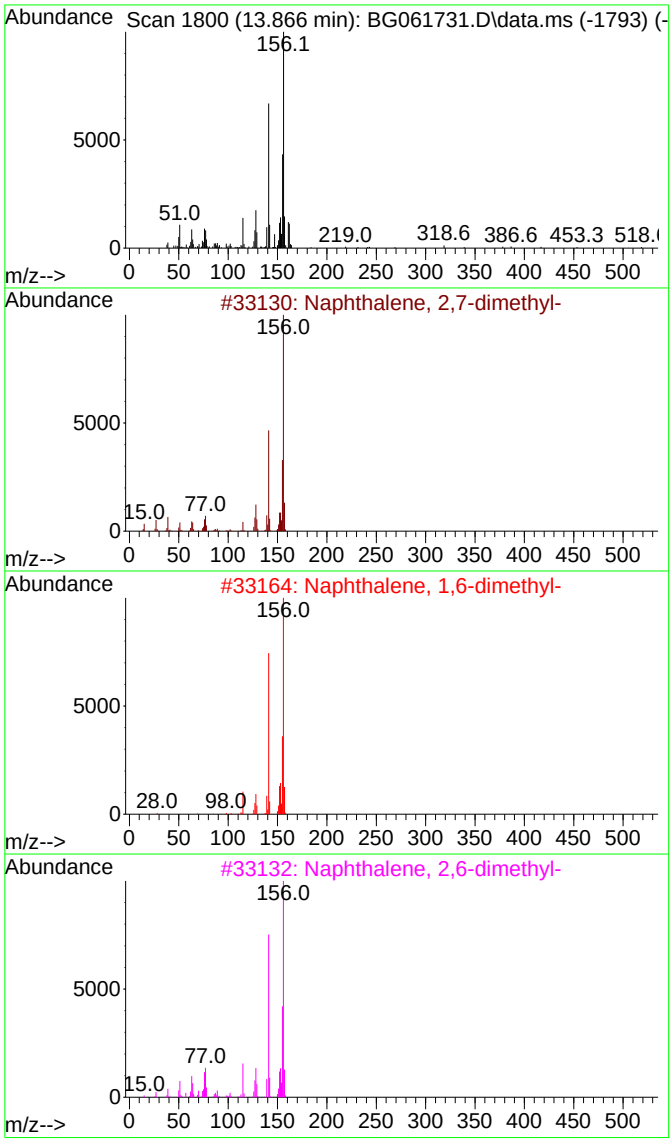
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

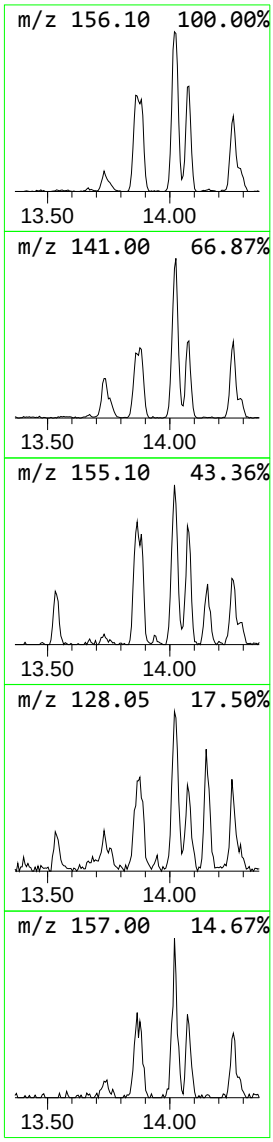
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 17 Naphthalene, 2,7-dimethyl- Concentration Rank 17

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 13.866 | 5.75 ng/ul | 312623 | Acenaphthene-d10 | 14.701 |

| 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, 2,6-dimethyl- | 156 | C12H12 | 000581-42-0 | 95 |
| 4 | Naphthalene, 1,5-dimethyl- | 156 | C12H12 | 000571-61-9 | 95 |
| 5 | Naphthalene, 2,3-dimethyl- | 156 | C12H12 | 000581-40-8 | 95 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062424\
Data File : BG061731.D
Acq On : 26 Jun 2024 11:43
Operator : MA/JU
Sample : P2873-09DL 10X
Misc :
ALS Vial : 30 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
C0SJ9DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

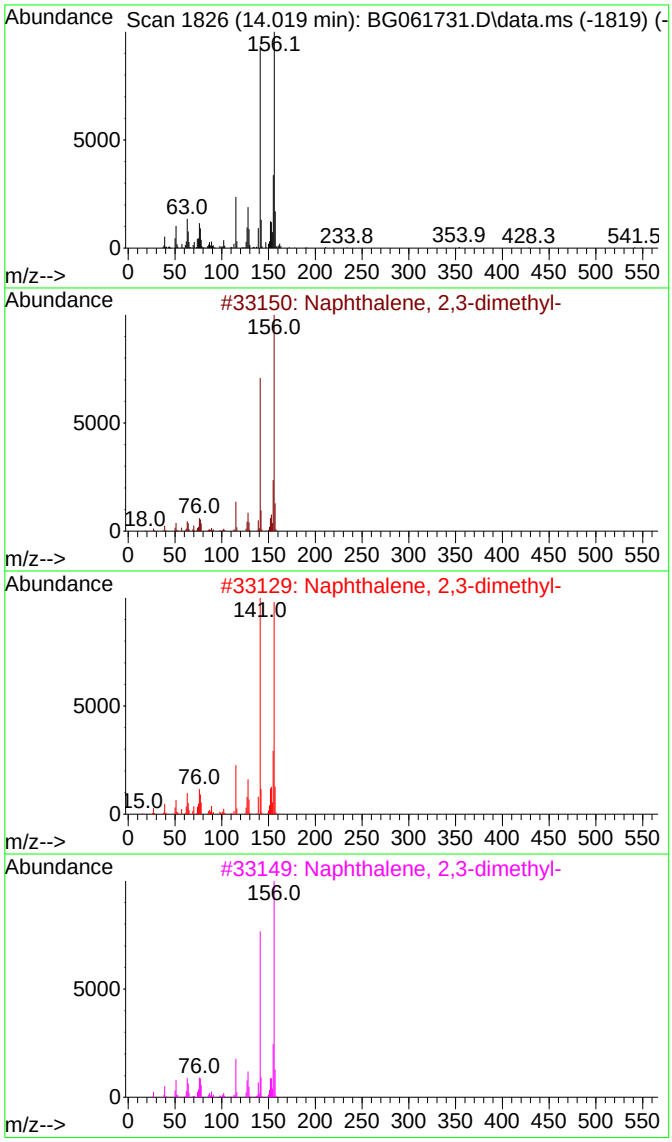
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

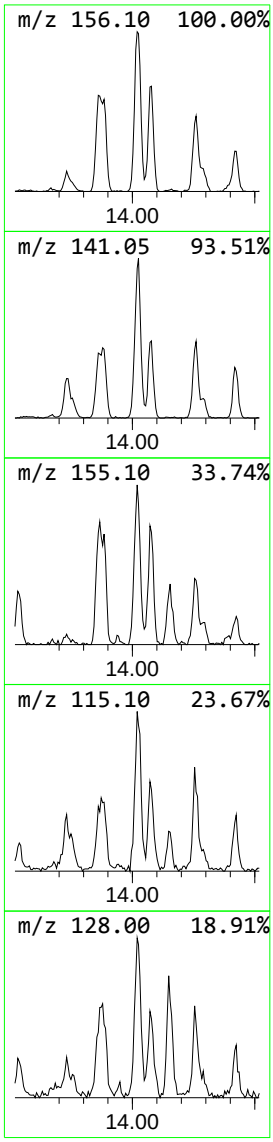
**********

Peak Number 18 Naphthalene, 2,3-dimethyl- Concentration Rank 7

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 14.019 | 12.08 ng/ul | 657174 | Acenaphthene-d10 | 14.701 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Naphthalene, 2,3-dimethyl- | 156 | C12H12 | 000581-40-8 | 97 |
| 2 | Naphthalene, 2,3-dimethyl- | 156 | C12H12 | 000581-40-8 | 97 |
| 3 | Naphthalene, 2,3-dimethyl- | 156 | C12H12 | 000581-40-8 | 96 |
| 4 | Naphthalene, 1,3-dimethyl- | 156 | C12H12 | 000575-41-7 | 96 |
| 5 | Naphthalene, 2,6-dimethyl- | 156 | C12H12 | 000581-42-0 | 96 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062424\
Data File : BG061731.D
Acq On : 26 Jun 2024 11:43
Operator : MA/JU
Sample : P2873-09DL 10X
Misc :
ALS Vial : 30 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
C0SJ9DL

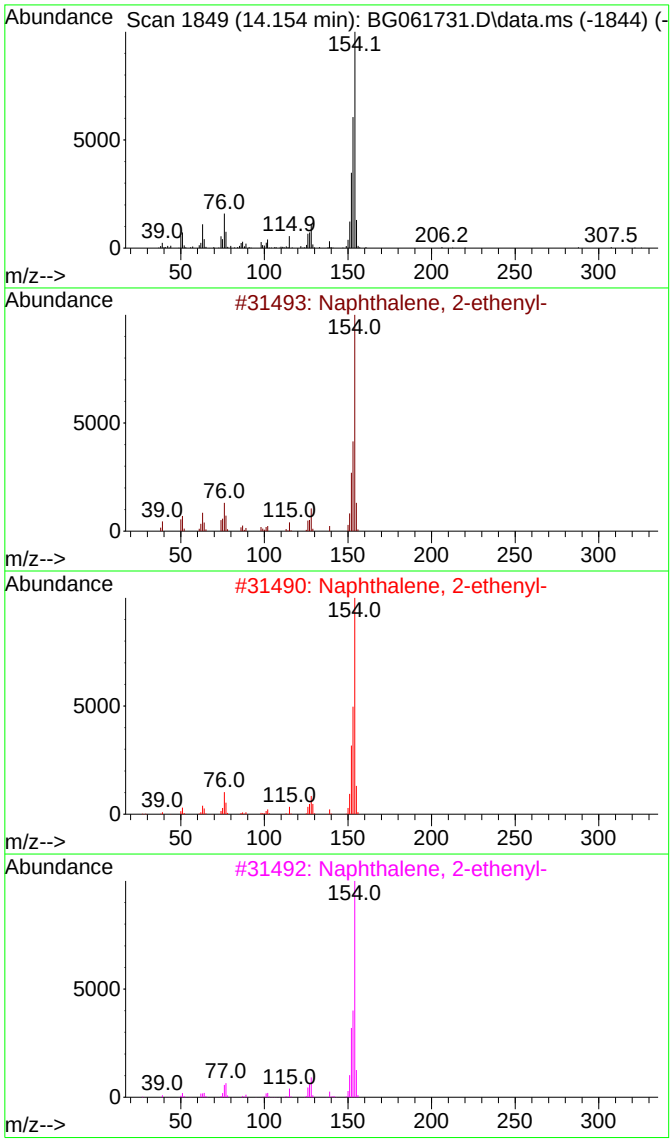
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

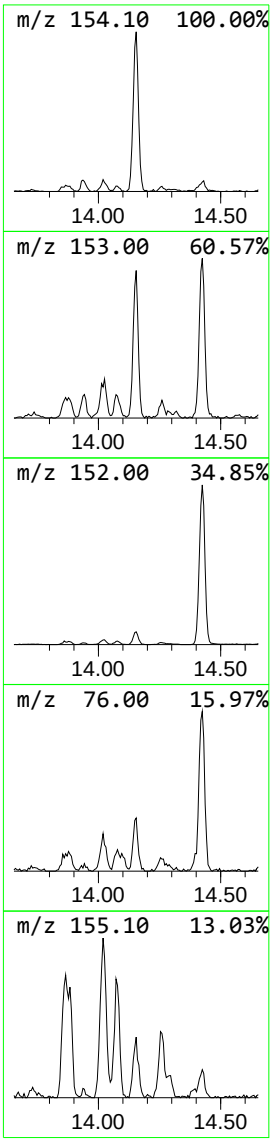
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 19 Naphthalene, 2-ethenyl- Concentration Rank 12

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 14.154 | 7.83 ng/ul | 425906 | Acenaphthene-d10 | 14.701 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Naphthalene, 2-ethenyl- | 154 | C12H10 | 000827-54-3 | 96 |
| 2 | Naphthalene, 2-ethenyl- | 154 | C12H10 | 000827-54-3 | 94 |
| 3 | Naphthalene, 2-ethenyl- | 154 | C12H10 | 000827-54-3 | 93 |
| 4 | Biphenyl | 154 | C12H10 | 000092-52-4 | 91 |
| 5 | Biphenyl | 154 | C12H10 | 000092-52-4 | 83 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062424\
Data File : BG061731.D
Acq On : 26 Jun 2024 11:43
Operator : MA/JU
Sample : P2873-09DL 10X
Misc :
ALS Vial : 30 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
C0SJ9DL

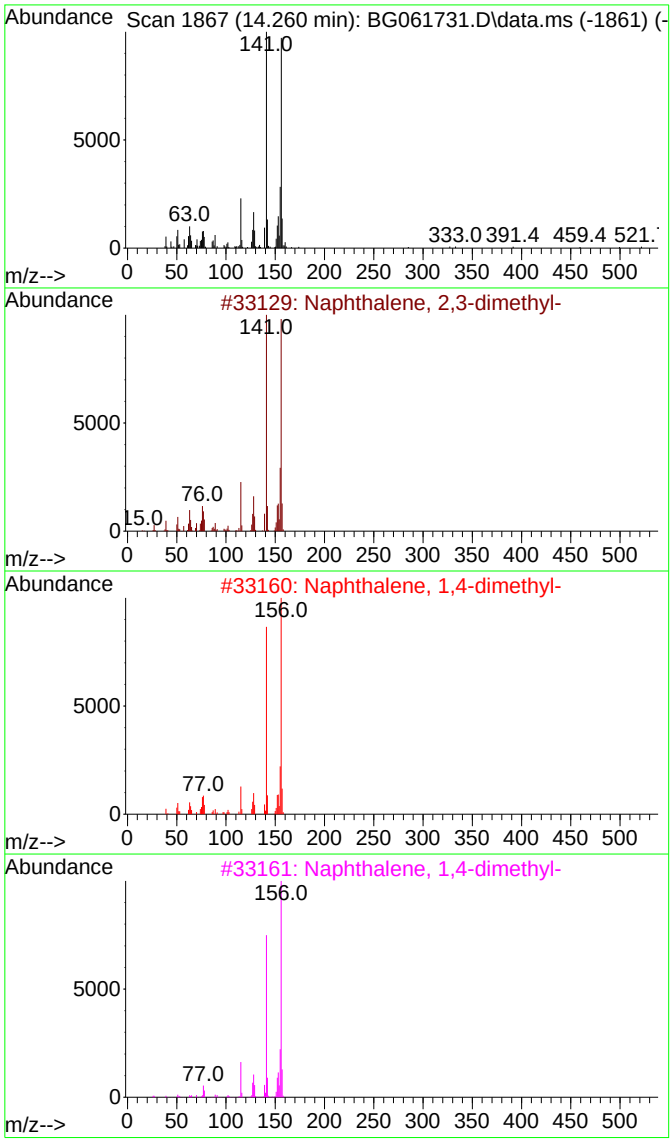
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

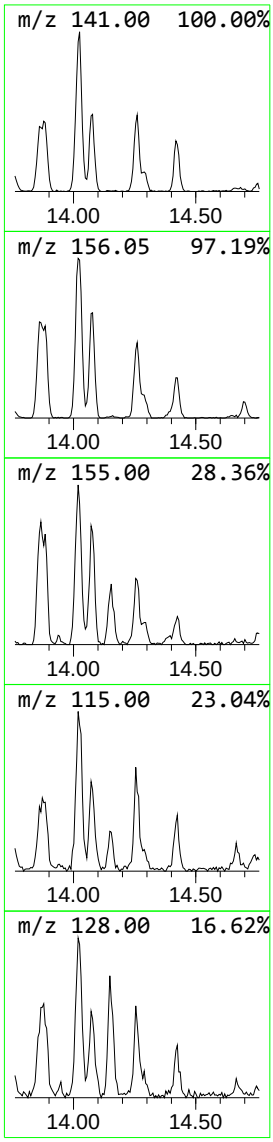
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 20 Naphthalene, 1,4-dimethyl- Concentration Rank 23

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 14.260 | 4.80 ng/ul | 261244 | Acenaphthene-d10 | 14.701 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Naphthalene, 2,3-dimethyl- | 156 | C12H12 | 000581-40-8 | 97 |
| 2 | Naphthalene, 1,4-dimethyl- | 156 | C12H12 | 000571-58-4 | 97 |
| 3 | Naphthalene, 1,4-dimethyl- | 156 | C12H12 | 000571-58-4 | 97 |
| 4 | Naphthalene, 1,4-dimethyl- | 156 | C12H12 | 000571-58-4 | 97 |
| 5 | Naphthalene, 2,3-dimethyl- | 156 | C12H12 | 000581-40-8 | 96 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062424\
Data File : BG061731.D
Acq On : 26 Jun 2024 11:43
Operator : MA/JU
Sample : P2873-09DL 10X
Misc :
ALS Vial : 30 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
C0SJ9DL

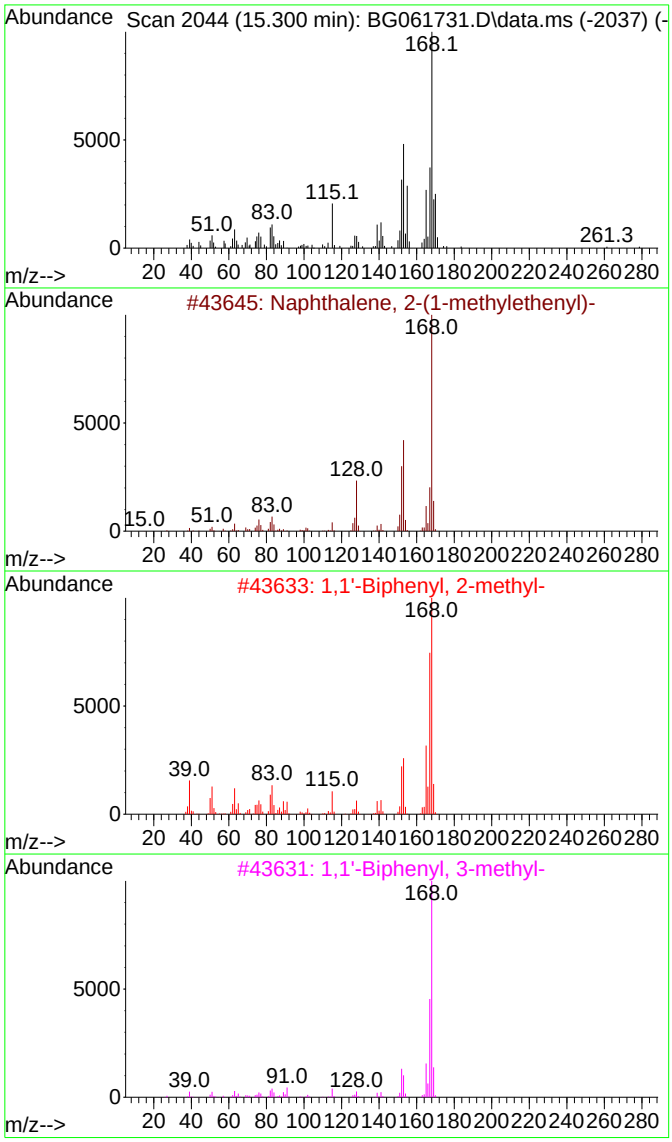
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

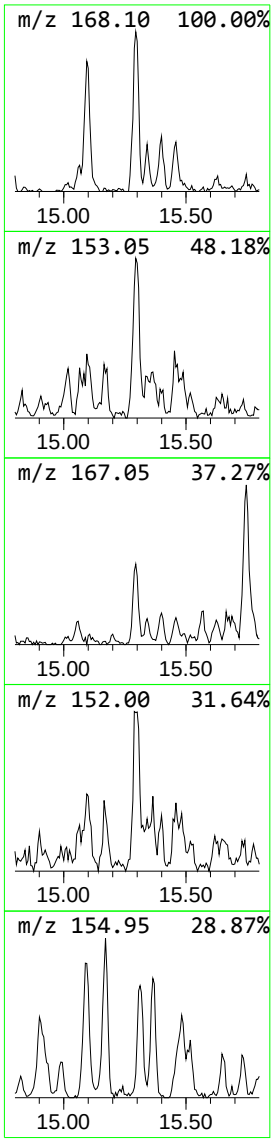
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 22 Naphthalene, 2-(1-methyleth... Concentration Rank 19

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 15.300 | 5.27 ng/ul | 286713 | Acenaphthene-d10 | 14.701 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Naphthalene, 2-(1-methylethenyl)- | 168 | C13H12 | 003710-23-4 | 58 |
| 2 | 1,1'-Biphenyl, 2-methyl- | 168 | C13H12 | 000643-58-3 | 52 |
| 3 | 1,1'-Biphenyl, 3-methyl- | 168 | C13H12 | 000643-93-6 | 49 |
| 4 | Naphthalene, 1-(2-propenyl)- | 168 | C13H12 | 002489-86-3 | 47 |
| 5 | 1,1'-Biphenyl, 3-methyl- | 168 | C13H12 | 000643-93-6 | 43 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062424\
Data File : BG061731.D
Acq On : 26 Jun 2024 11:43
Operator : MA/JU
Sample : P2873-09DL 10X
Misc :
ALS Vial : 30 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
C0SJ9DL

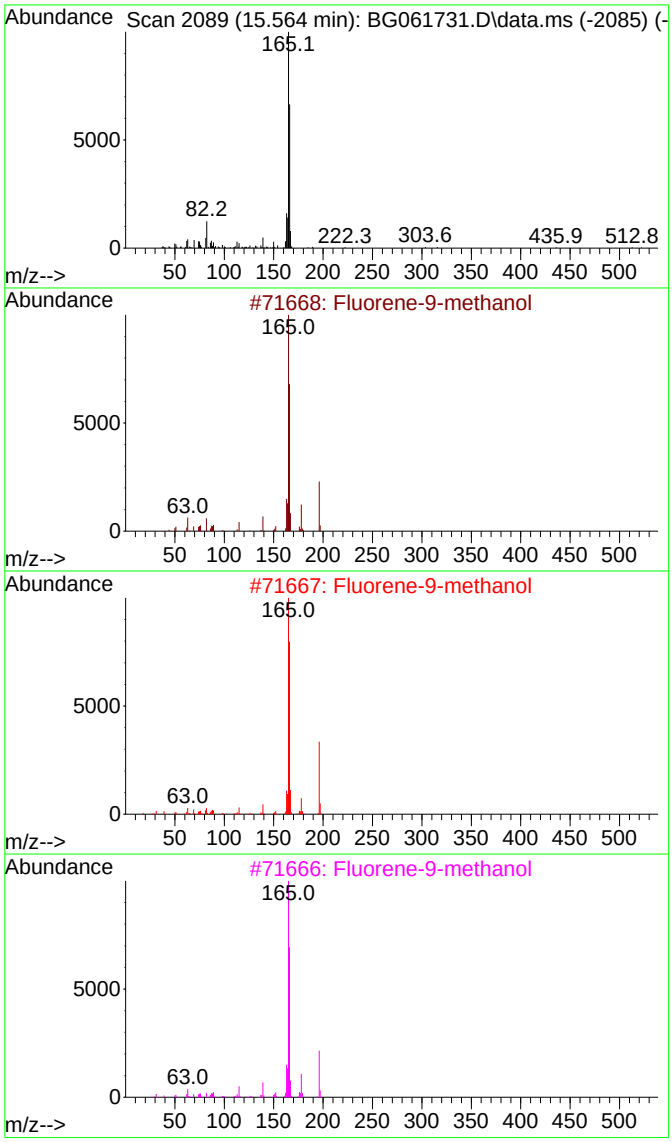
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

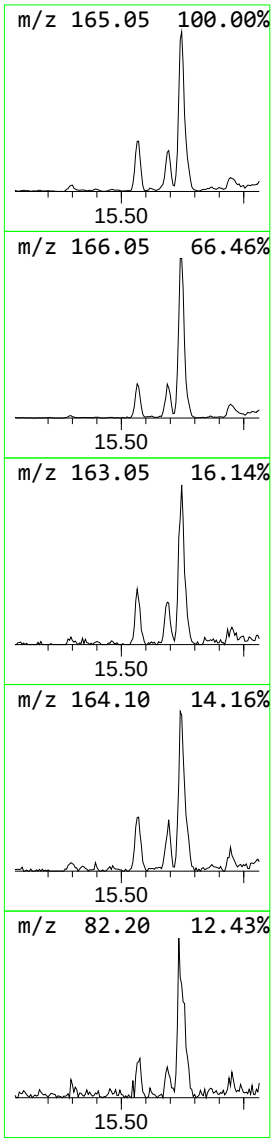
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 23 Fluorene-9-methanol Concentration Rank 26

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 15.564 | 4.30 ng/ul | 234112 | Acenaphthene-d10 | 14.701 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Fluorene-9-methanol | 196 | C14H12O | 024324-17-2 | 78 |
| 2 | Fluorene-9-methanol | 196 | C14H12O | 024324-17-2 | 78 |
| 3 | Fluorene-9-methanol | 196 | C14H12O | 024324-17-2 | 78 |
| 4 | Fluorene-9-methanol | 196 | C14H12O | 024324-17-2 | 64 |
| 5 | Benzene, 5-heptene-1,3-diyn-1-yl- | 166 | C13H10 | 013678-98-3 | 64 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062424\
Data File : BG061731.D
Acq On : 26 Jun 2024 11:43
Operator : MA/JU
Sample : P2873-09DL 10X
Misc :
ALS Vial : 30 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
C0SJ9DL

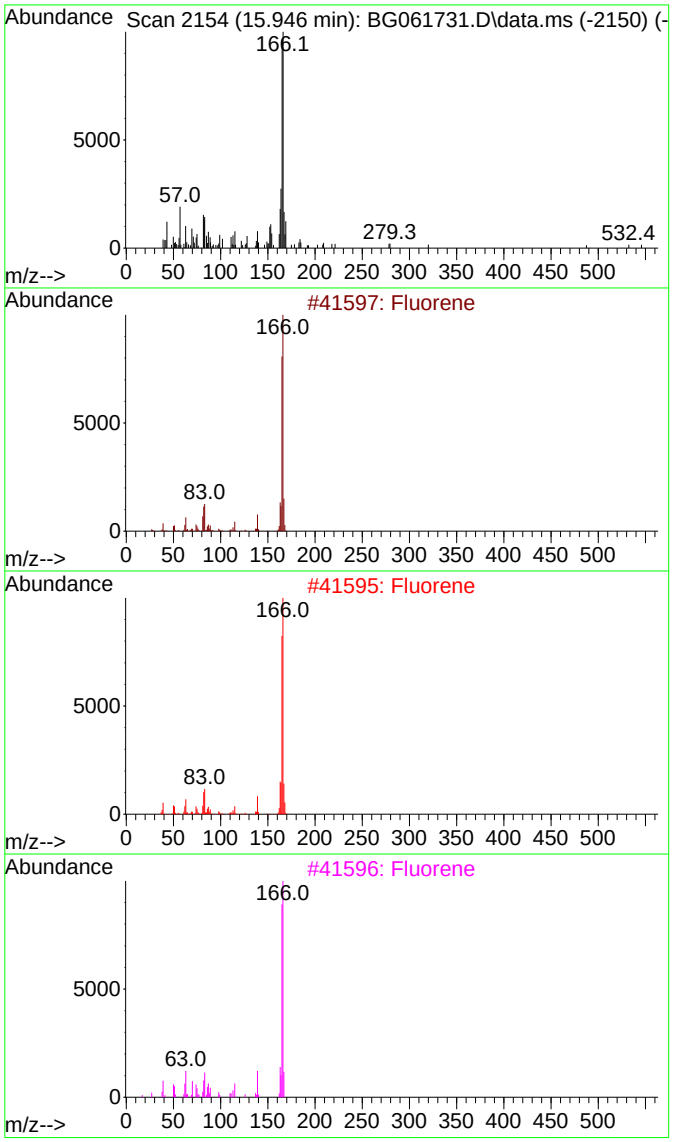
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

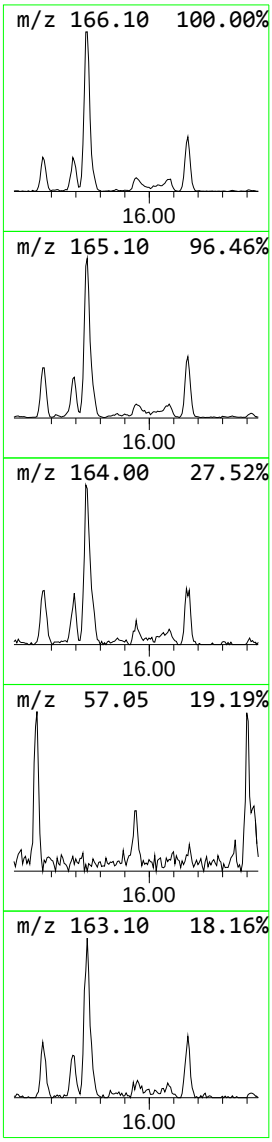
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 25 Benzene, 5-heptene-1,3-diyn... Concentration Rank 42

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 15.946 | 2.16 ng/ul | 117604 | Acenaphthene-d10 | 14.701 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Fluorene | 166 | C13H10 | 000086-73-7 | 87 |
| 2 | Fluorene | 166 | C13H10 | 000086-73-7 | 83 |
| 3 | Fluorene | 166 | C13H10 | 000086-73-7 | 80 |
| 4 | Fluorene | 166 | C13H10 | 000086-73-7 | 80 |
| 5 | Benzene, 5-heptene-1,3-diyn-1-yl- | 166 | C13H10 | 013678-98-3 | 74 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062424\
Data File : BG061731.D
Acq On : 26 Jun 2024 11:43
Operator : MA/JU
Sample : P2873-09DL 10X
Misc :
ALS Vial : 30 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
C0SJ9DL

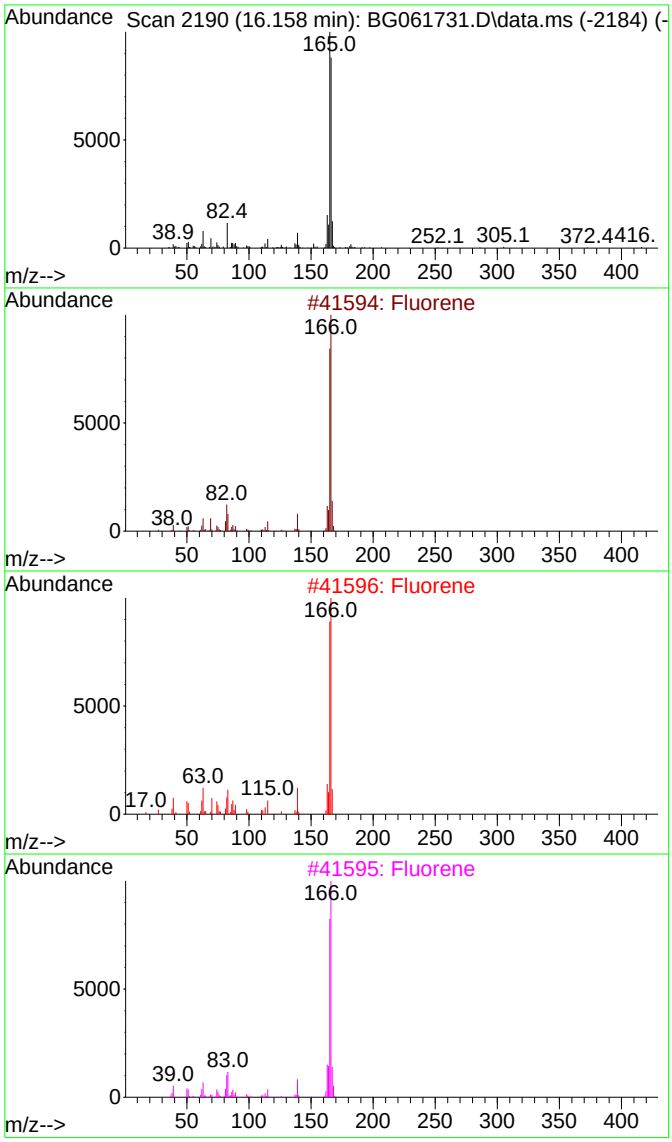
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

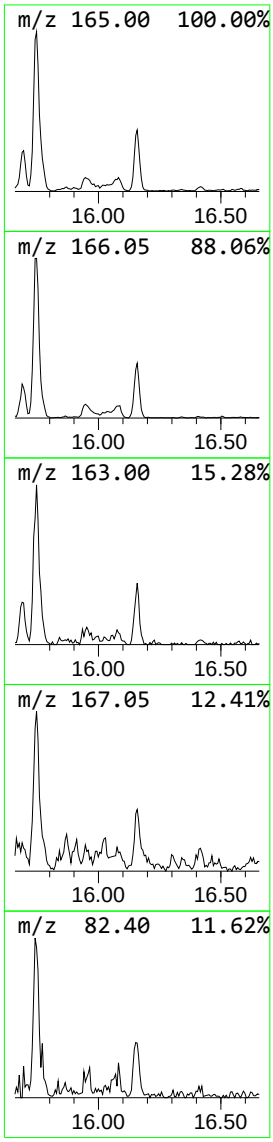
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 26 Benzaldehyde, 4,6-dihydroxy... Concentration Rank 13

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 16.158 | 6.96 ng/ul | 388997 | Phenanthrene-d10 | 17.445 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Fluorene | 166 | C13H10 | 000086-73-7 | 91 |
| 2 | Fluorene | 166 | C13H10 | 000086-73-7 | 90 |
| 3 | Fluorene | 166 | C13H10 | 000086-73-7 | 86 |
| 4 | Fluorene | 166 | C13H10 | 000086-73-7 | 76 |
| 5 | Benzaldehyde, 4,6-dihydroxy-2,3-... | 166 | C9H10O3 | 002990-31-0 | 72 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062424\
Data File : BG061731.D
Acq On : 26 Jun 2024 11:43
Operator : MA/JU
Sample : P2873-09DL 10X
Misc :
ALS Vial : 30 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
C0SJ9DL

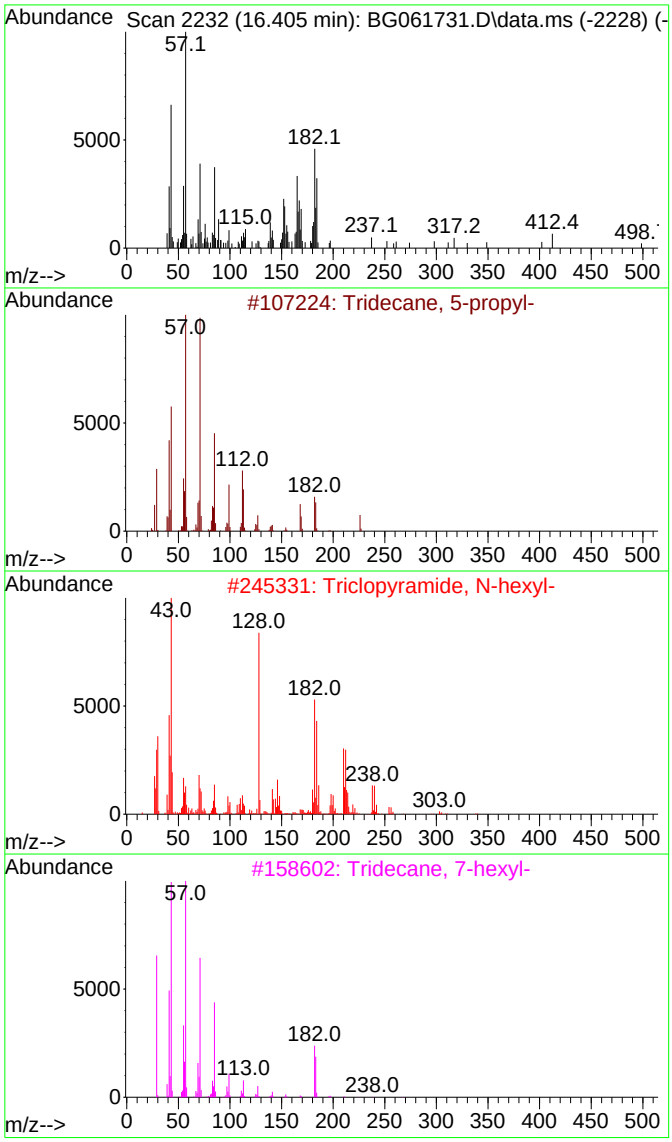
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

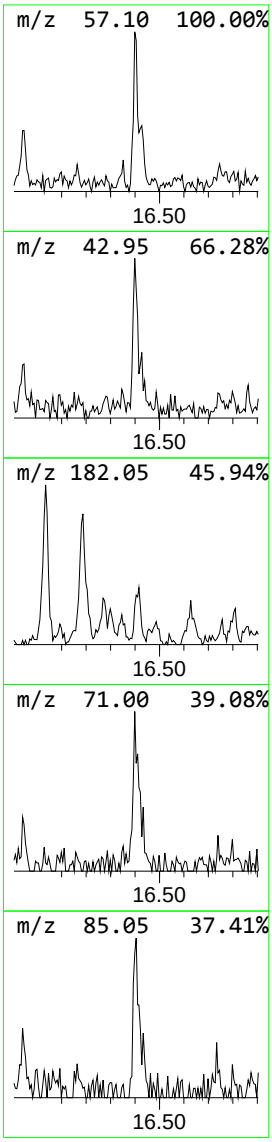
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 27 (DEL) Alkane: Straight-Chai... Concentration Rank 40

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 16.404 | 2.32 ng/ul | 129942 | Phenanthrene-d10 | 17.445 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Tridecane, 5-propyl- | 226 | C16H34 | 055045-11-9 | 14 |
| 2 | Triclopyramide, N-hexyl- | 338 | C13H17Cl3N2O2 | 1000420-57-8 | 14 |
| 3 | Tridecane, 7-hexyl- | 268 | C19H40 | 007225-66-3 | 12 |
| 4 | 3-(Boc-amino)propyl bromide | 237 | C8H16BrNO2 | 083948-53-2 | 10 |
| 5 | 2,2-Diphenylethylamine | 197 | C14H15N | 003963-62-0 | 10 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062424\
Data File : BG061731.D
Acq On : 26 Jun 2024 11:43
Operator : MA/JU
Sample : P2873-09DL 10X
Misc :
ALS Vial : 30 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
C0SJ9DL

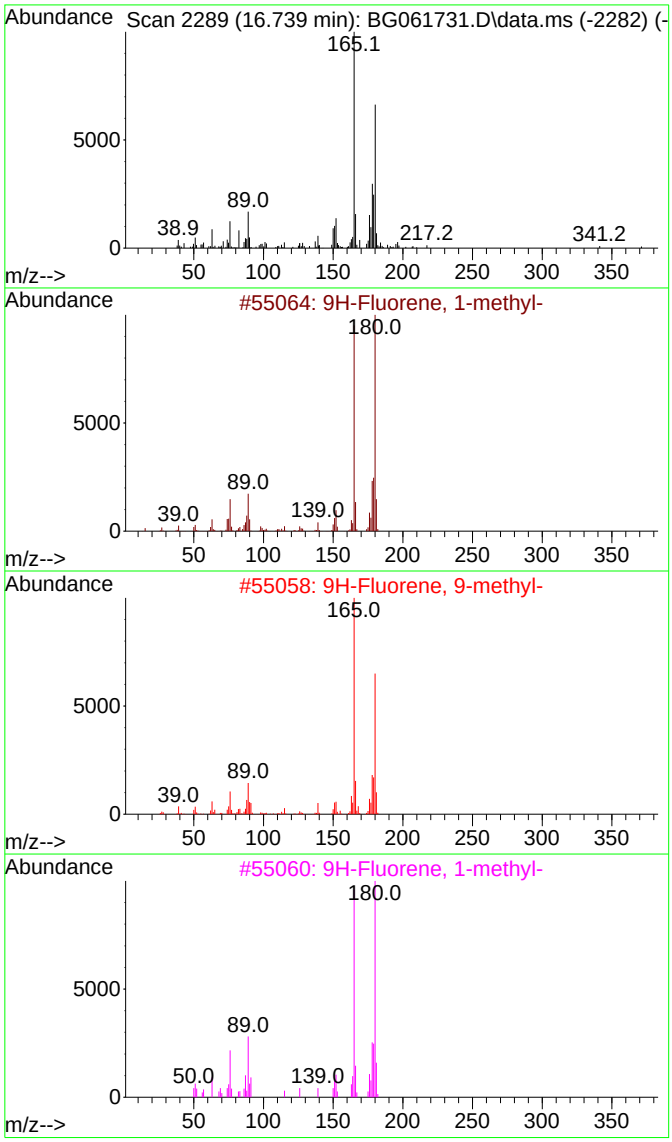
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

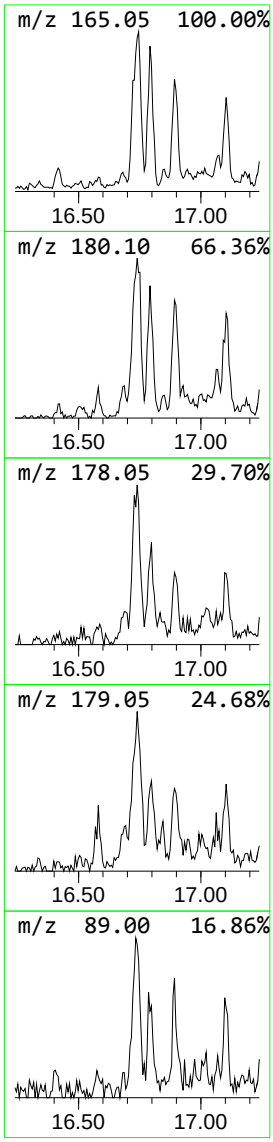
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 28 9H-Fluorene, 1-methyl- Concentration Rank 18

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 16.739 | 5.50 ng/ul | 307667 | Phenanthrene-d10 | 17.445 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 9H-Fluorene, 1-methyl- | 180 | C14H12 | 001730-37-6 | 90 |
| 2 | 9H-Fluorene, 9-methyl- | 180 | C14H12 | 002523-37-7 | 90 |
| 3 | 9H-Fluorene, 1-methyl- | 180 | C14H12 | 001730-37-6 | 87 |
| 4 | 9H-Fluorene, 1-methyl- | 180 | C14H12 | 001730-37-6 | 83 |
| 5 | 9H-Fluorene, 2-methyl- | 180 | C14H12 | 001430-97-3 | 76 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062424\
Data File : BG061731.D
Acq On : 26 Jun 2024 11:43
Operator : MA/JU
Sample : P2873-09DL 10X
Misc :
ALS Vial : 30 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
C0SJ9DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

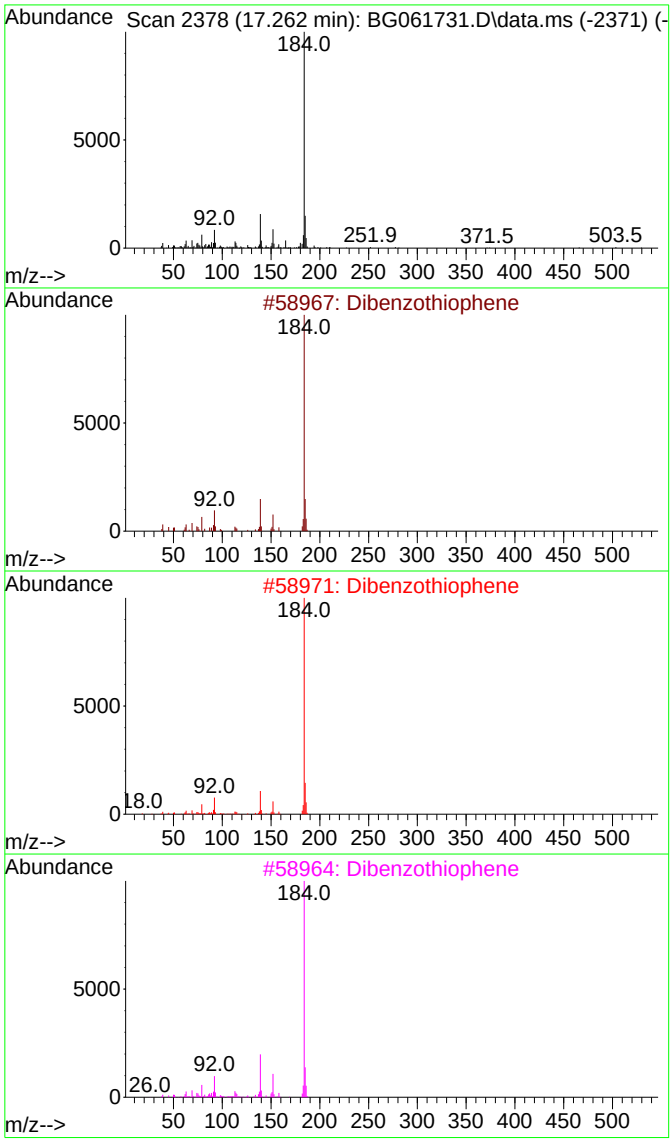
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

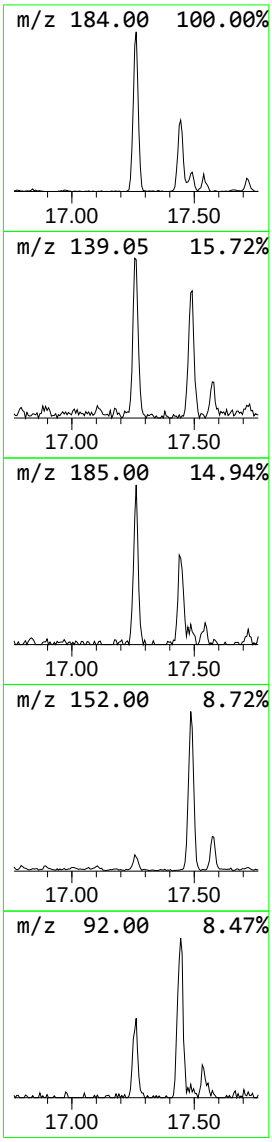
**********

Peak Number 31 Dibenzothiophene Concentration Rank 15

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 17.262 | 6.56 ng/ul | 366524 | Phenanthrene-d10 | 17.445 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Dibenzothiophene | 184 | C12H8S | 000132-65-0 | 96 |
| 2 | Dibenzothiophene | 184 | C12H8S | 000132-65-0 | 93 |
| 3 | Dibenzothiophene | 184 | C12H8S | 000132-65-0 | 93 |
| 4 | Dibenzothiophene | 184 | C12H8S | 000132-65-0 | 93 |
| 5 | Dibenzothiophene | 184 | C12H8S | 000132-65-0 | 93 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062424\
Data File : BG061731.D
Acq On : 26 Jun 2024 11:43
Operator : MA/JU
Sample : P2873-09DL 10X
Misc :
ALS Vial : 30 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
C0SJ9DL

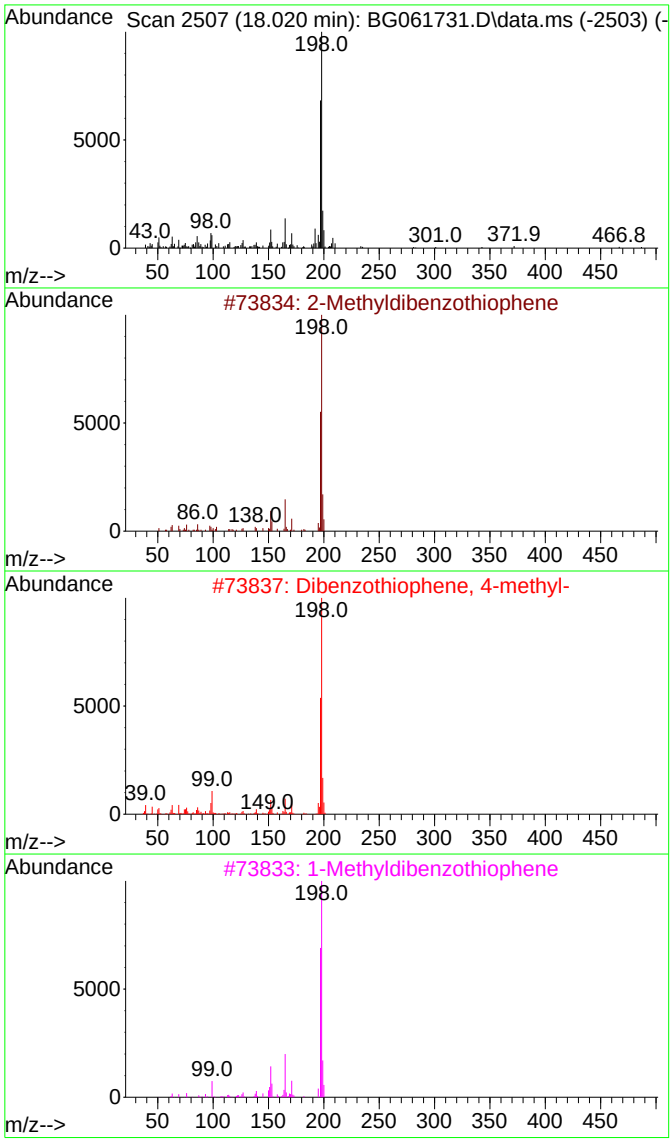
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

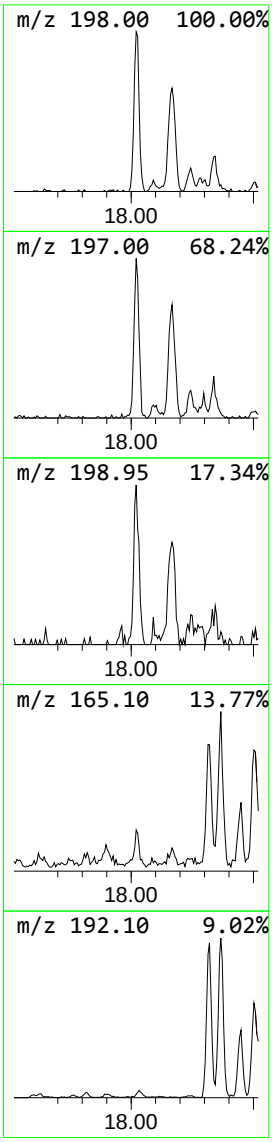
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 32 2-Methyldibenzothiophene Concentration Rank 30

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 18.020 | 3.32 ng/ul | 185547 | Phenanthrene-d10 | 17.445 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2-Methyldibenzothiophene | 198 | C13H10S | 1000383-55-1 | 93 |
| 2 | Dibenzothiophene, 4-methyl- | 198 | C13H10S | 007372-88-5 | 93 |
| 3 | 1-Methyldibenzothiophene | 198 | C13H10S | 031317-07-4 | 91 |
| 4 | Dibenzothiophene, 3-methyl- | 198 | C13H10S | 016587-52-3 | 90 |
| 5 | 4-Methylnaphtho[1,2-b]thiophene | 198 | C13H10S | 067388-11-8 | 81 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062424\
Data File : BG061731.D
Acq On : 26 Jun 2024 11:43
Operator : MA/JU
Sample : P2873-09DL 10X
Misc :
ALS Vial : 30 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
C0SJ9DL

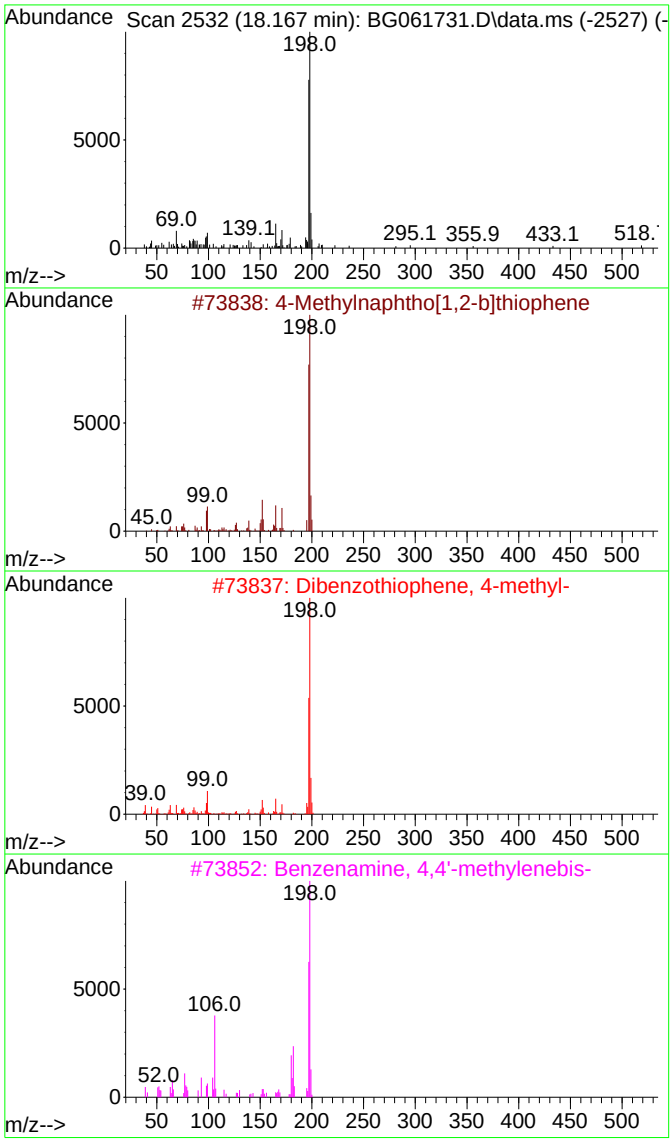
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

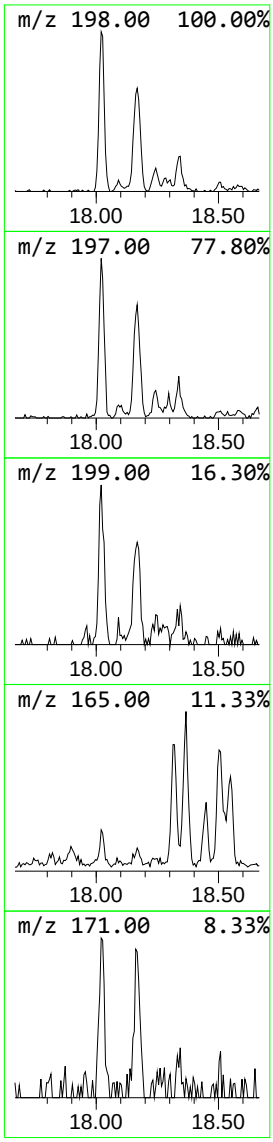
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 33 4-Methylnaphtho[1,2-b]thiop... Concentration Rank 32

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 18.167 | 3.02 ng/ul | 169000 | Phenanthrene-d10 | 17.445 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 4-Methylnaphtho[1,2-b]thiophene | 198 | C13H10S | 067388-11-8 | 87 |
| 2 | Dibenzothiophene, 4-methyl- | 198 | C13H10S | 007372-88-5 | 86 |
| 3 | Benzenamine, 4,4'-methylenebis- | 198 | C13H14N2 | 000101-77-9 | 80 |
| 4 | Benzenamine, 4,4'-methylenebis- | 198 | C13H14N2 | 000101-77-9 | 72 |
| 5 | Tacrine | 198 | C13H14N2 | 000321-64-2 | 72 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062424\
Data File : BG061731.D
Acq On : 26 Jun 2024 11:43
Operator : MA/JU
Sample : P2873-09DL 10X
Misc :
ALS Vial : 30 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
C0SJ9DL

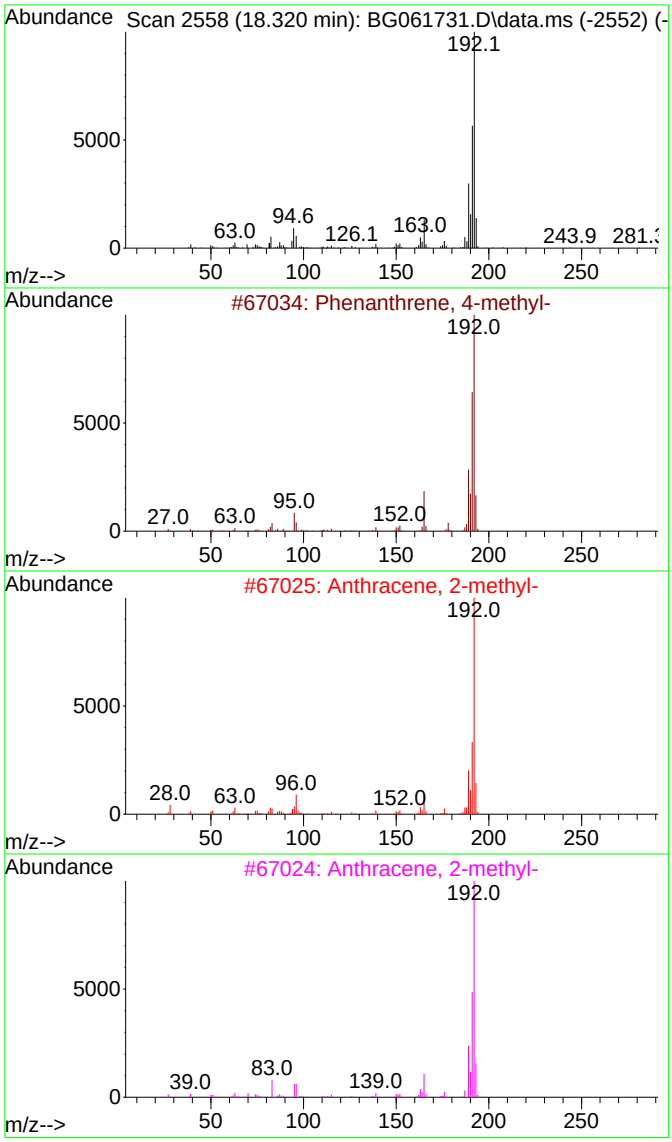
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

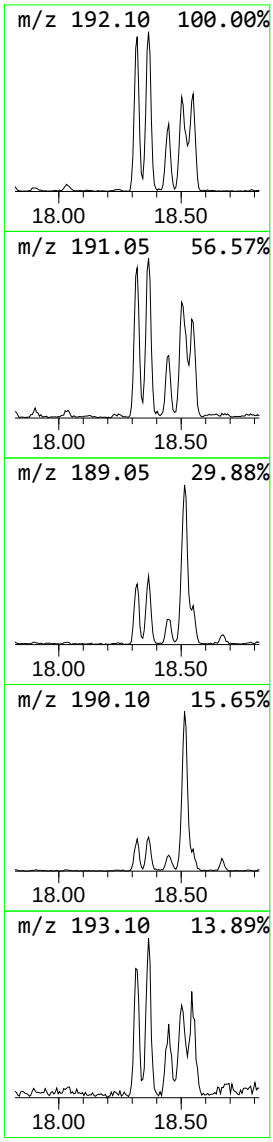
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 34 Phenanthrene, 4-methyl- Concentration Rank 9

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 18.320 | 11.64 ng/ul | 650436 | Phenanthrene-d10 | 17.445 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Phenanthrene, 4-methyl- | 192 | C15H12 | 000832-64-4 | 95 |
| 2 | Anthracene, 2-methyl- | 192 | C15H12 | 000613-12-7 | 95 |
| 3 | Anthracene, 2-methyl- | 192 | C15H12 | 000613-12-7 | 93 |
| 4 | Phenanthrene, 2-methyl- | 192 | C15H12 | 002531-84-2 | 92 |
| 5 | Phenanthrene, 2-methyl- | 192 | C15H12 | 002531-84-2 | 91 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062424\
Data File : BG061731.D
Acq On : 26 Jun 2024 11:43
Operator : MA/JU
Sample : P2873-09DL 10X
Misc :
ALS Vial : 30 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
C0SJ9DL

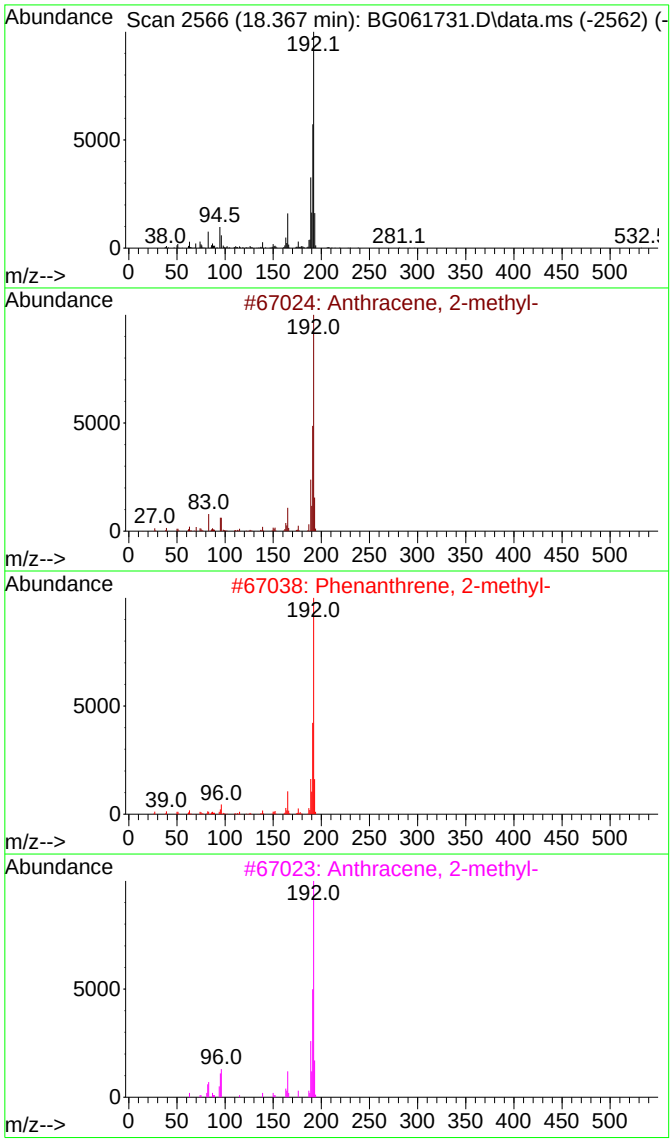
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

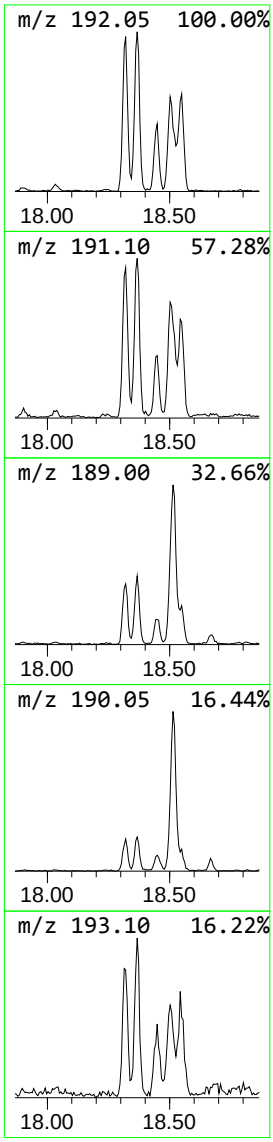
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 35 Anthracene, 2-methyl- Concentration Rank 6

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 18.367 | 12.31 ng/ul | 687892 | Phenanthrene-d10 | 17.445 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Anthracene, 2-methyl- | 192 | C15H12 | 000613-12-7 | 95 |
| 2 | Phenanthrene, 2-methyl- | 192 | C15H12 | 002531-84-2 | 93 |
| 3 | Anthracene, 2-methyl- | 192 | C15H12 | 000613-12-7 | 93 |
| 4 | Phenanthrene, 2-methyl- | 192 | C15H12 | 002531-84-2 | 93 |
| 5 | Phenanthrene, 1-methyl- | 192 | C15H12 | 000832-69-9 | 91 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062424\
Data File : BG061731.D
Acq On : 26 Jun 2024 11:43
Operator : MA/JU
Sample : P2873-09DL 10X
Misc :
ALS Vial : 30 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
C0SJ9DL

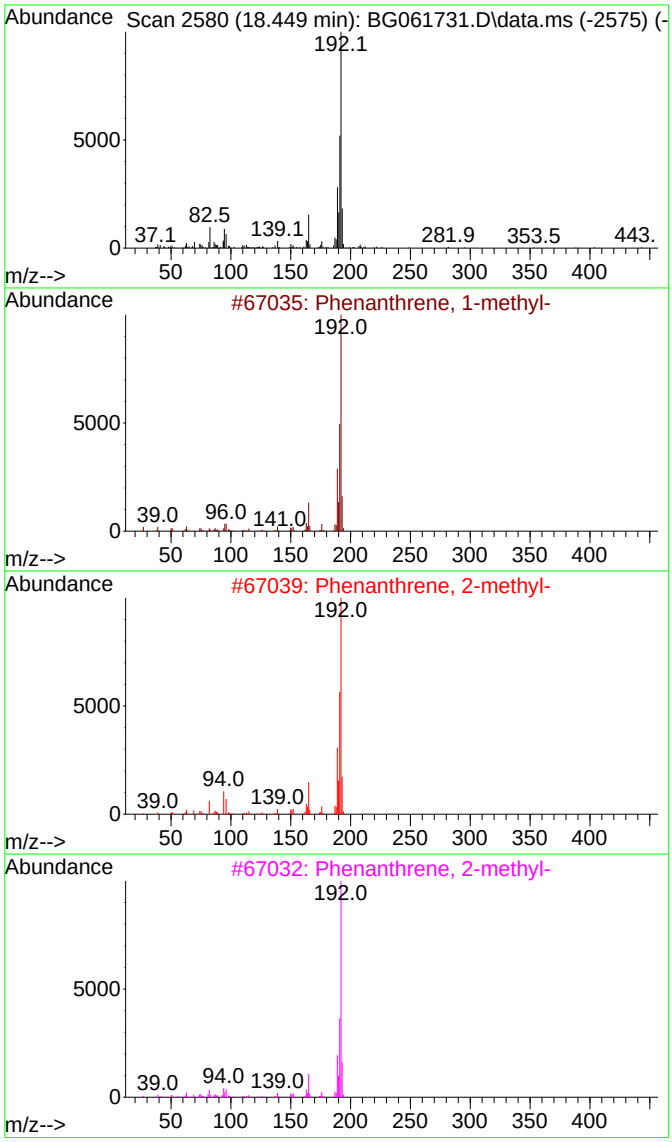
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

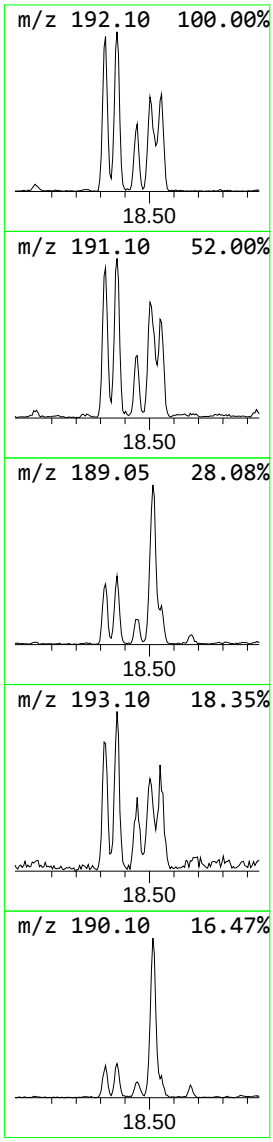
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 36 Phenanthrene, 1-methyl- Concentration Rank 24

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 18.449 | 4.78 ng/ul | 266971 | Phenanthrene-d10 | 17.445 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Phenanthrene, 1-methyl- | 192 | C15H12 | 000832-69-9 | 95 |
| 2 | Phenanthrene, 2-methyl- | 192 | C15H12 | 002531-84-2 | 93 |
| 3 | Phenanthrene, 2-methyl- | 192 | C15H12 | 002531-84-2 | 91 |
| 4 | Naphtho[2,3-b]norbornadiene | 192 | C15H12 | 107426-38-0 | 90 |
| 5 | 1H-Cyclopropa[l]phenanthrene,1a,... | 192 | C15H12 | 000949-41-7 | 90 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062424\
Data File : BG061731.D
Acq On : 26 Jun 2024 11:43
Operator : MA/JU
Sample : P2873-09DL 10X
Misc :
ALS Vial : 30 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
C0SJ9DL

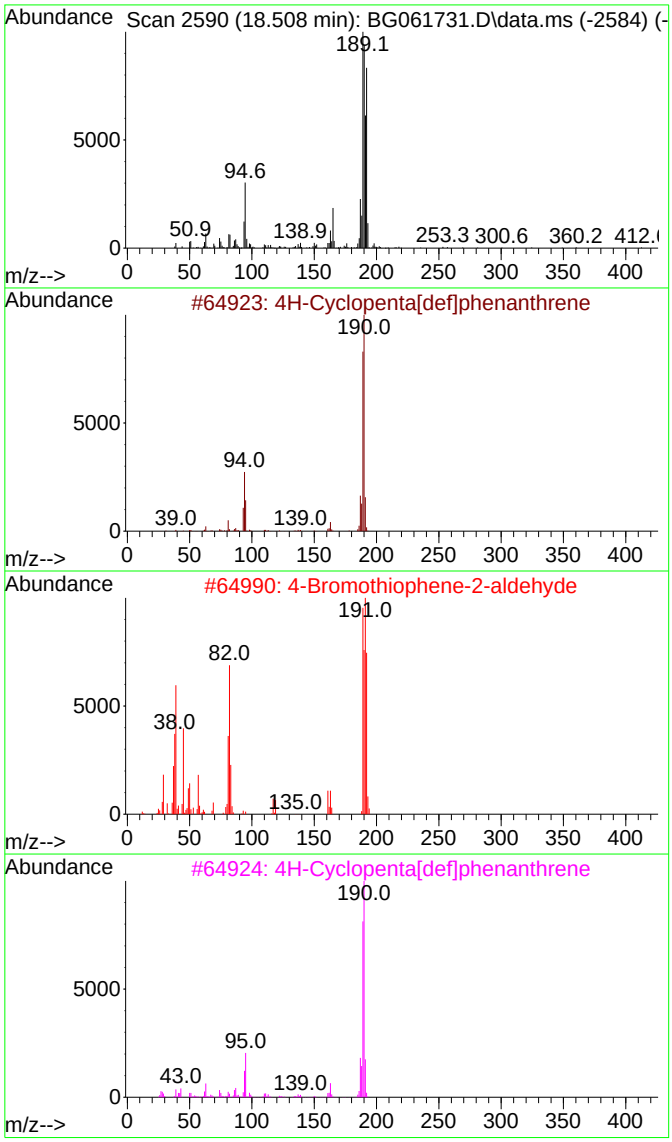
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

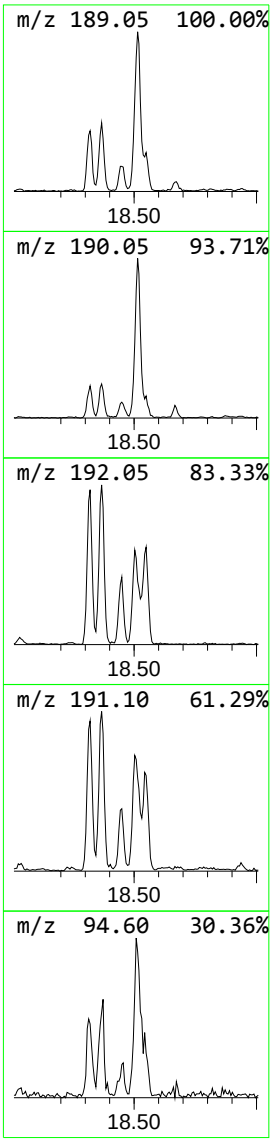
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 37 4H-Cyclopenta[def]phenanthrene Concentration Rank 4

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 18.508 | 18.79 ng/ul | 1050220 | Phenanthrene-d10 | 17.445 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 4H-Cyclopenta[def]phenanthrene | 190 | C15H10 | 000203-64-5 | 64 |
| 2 | 4-Bromothiophene-2-aldehyde | 190 | C5H3BrOS | 018791-75-8 | 58 |
| 3 | 4H-Cyclopenta[def]phenanthrene | 190 | C15H10 | 000203-64-5 | 55 |
| 4 | 6H-Cyclobuta[jk]phenanthrene | 190 | C15H10 | 083469-43-6 | 49 |
| 5 | Phenanthrene, 1-methyl- | 192 | C15H12 | 000832-69-9 | 38 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062424\
Data File : BG061731.D
Acq On : 26 Jun 2024 11:43
Operator : MA/JU
Sample : P2873-09DL 10X
Misc :
ALS Vial : 30 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
C0SJ9DL

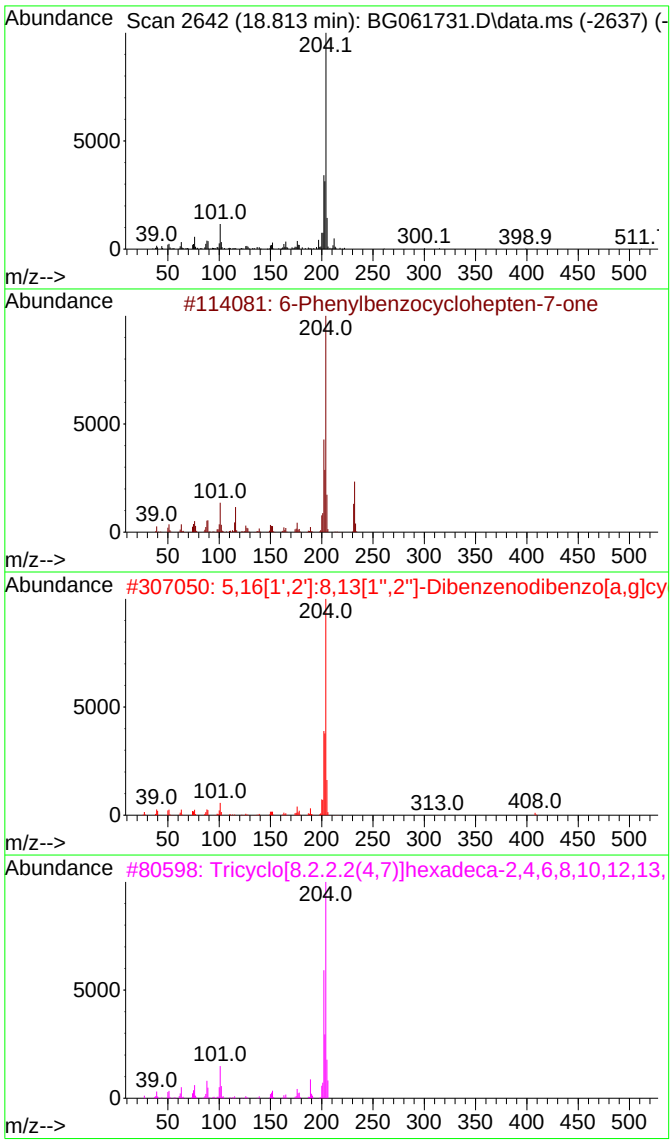
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

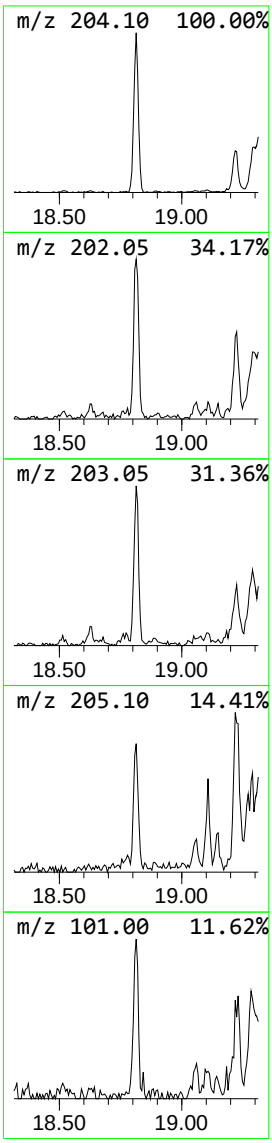
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 38 6-Phenylbenzocyclohepten-7-one Concentration Rank 22

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 18.814 | 5.13 ng/ul | 287030 | Phenanthrene-d10 | 17.445 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 6-Phenylbenzocyclohepten-7-one | 232 | C17H12O | 093327-56-1 | 90 |
| 2 | 5,16[1',2']:8,13[1'',2'']-Dibenz... | 408 | C32H24 | 005672-97-9 | 87 |
| 3 | Tricyclo[8.2.2.2(4,7)]hexadeca-2... | 204 | C16H12 | 006572-60-7 | 74 |
| 4 | Naphthalene, 2-phenyl- | 204 | C16H12 | 000612-94-2 | 72 |
| 5 | 9,10-Bis(bromomethyl)anthracene | 362 | C16H12Br2 | 034373-96-1 | 70 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062424\
Data File : BG061731.D
Acq On : 26 Jun 2024 11:43
Operator : MA/JU
Sample : P2873-09DL 10X
Misc :
ALS Vial : 30 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
C0SJ9DL

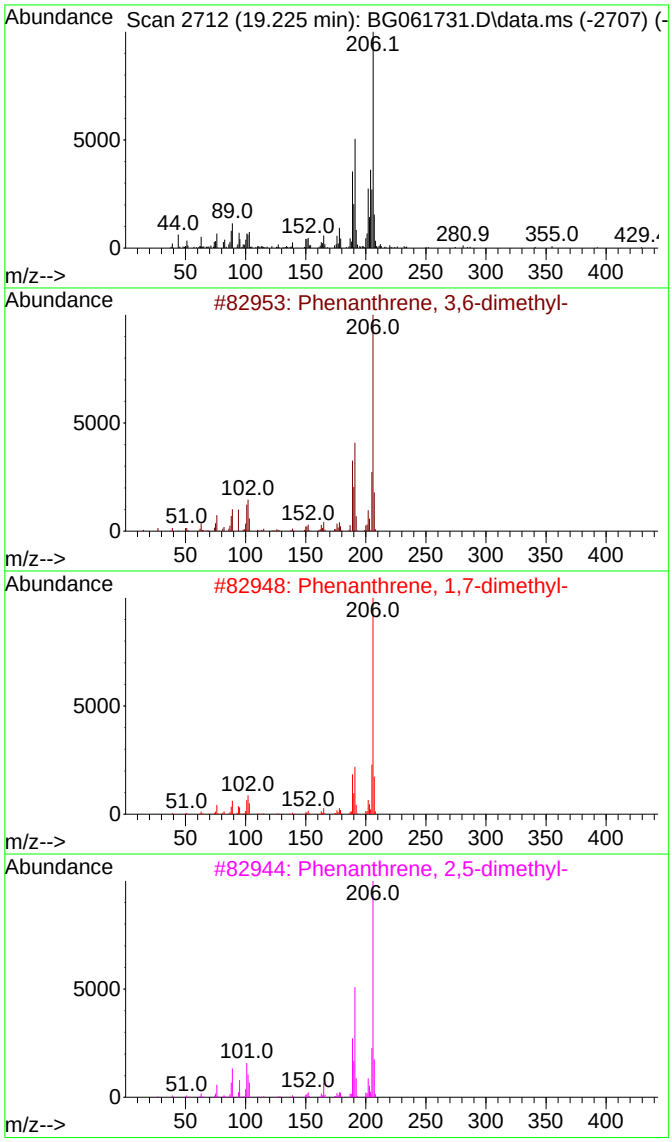
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

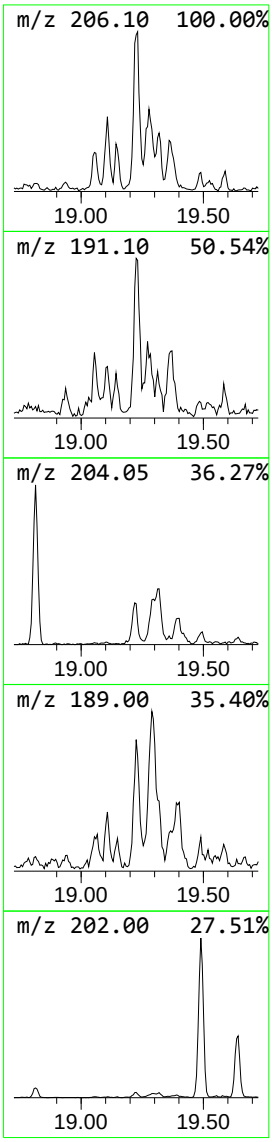
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 39 Phenanthrene, 3,6-dimethyl- Concentration Rank 14

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 19.225 | 6.77 ng/ul | 378632 | Phenanthrene-d10 | 17.445 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Phenanthrene, 3,6-dimethyl- | 206 | C16H14 | 001576-67-6 | 93 |
| 2 | Phenanthrene, 1,7-dimethyl- | 206 | C16H14 | 000483-87-4 | 90 |
| 3 | Phenanthrene, 2,5-dimethyl- | 206 | C16H14 | 003674-66-6 | 90 |
| 4 | Phenanthrene, 3,6-dimethyl- | 206 | C16H14 | 001576-67-6 | 90 |
| 5 | Anthracene, 1,4-dimethyl- | 206 | C16H14 | 000781-92-0 | 90 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062424\
Data File : BG061731.D
Acq On : 26 Jun 2024 11:43
Operator : MA/JU
Sample : P2873-09DL 10X
Misc :
ALS Vial : 30 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
C0SJ9DL

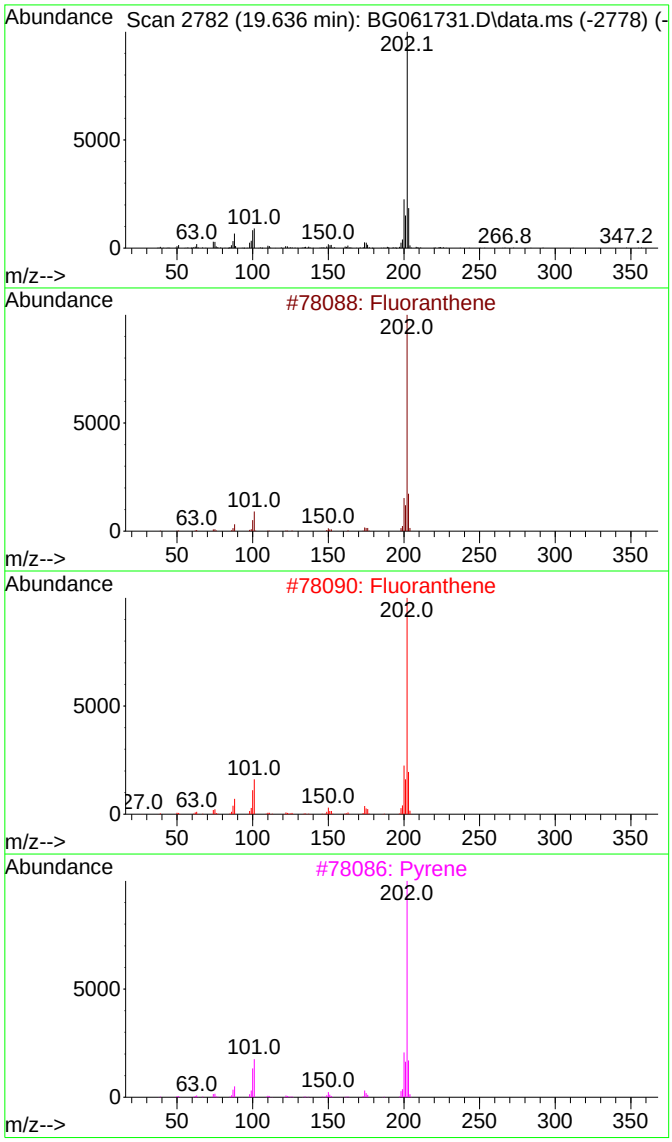
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

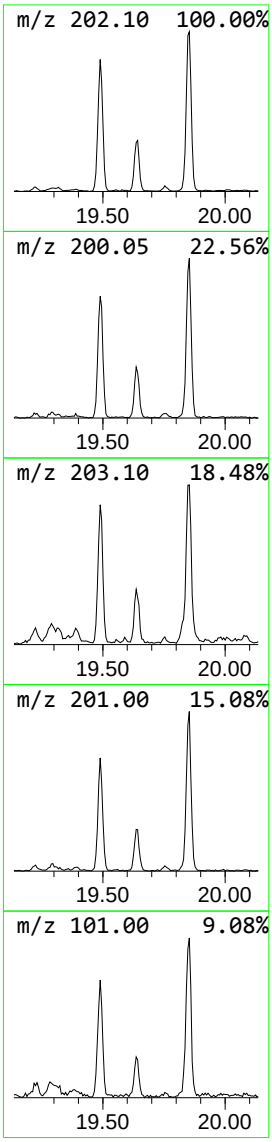
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 40 Benzene, 1,1'-(1,3-butadiyn... Concentration Rank 21

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 19.636 | 5.14 ng/ul | 526057 | Chrysene-d12 | 21.710 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Fluoranthene | 202 | C16H10 | 000206-44-0 | 94 |
| 2 | Fluoranthene | 202 | C16H10 | 000206-44-0 | 94 |
| 3 | Pyrene | 202 | C16H10 | 000129-00-0 | 89 |
| 4 | Fluoranthene | 202 | C16H10 | 000206-44-0 | 87 |
| 5 | Benzene, 1,1'-(1,3-butadiyne-1,4... | 202 | C16H10 | 000886-66-8 | 81 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062424\
Data File : BG061731.D
Acq On : 26 Jun 2024 11:43
Operator : MA/JU
Sample : P2873-09DL 10X
Misc :
ALS Vial : 30 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
C0SJ9DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

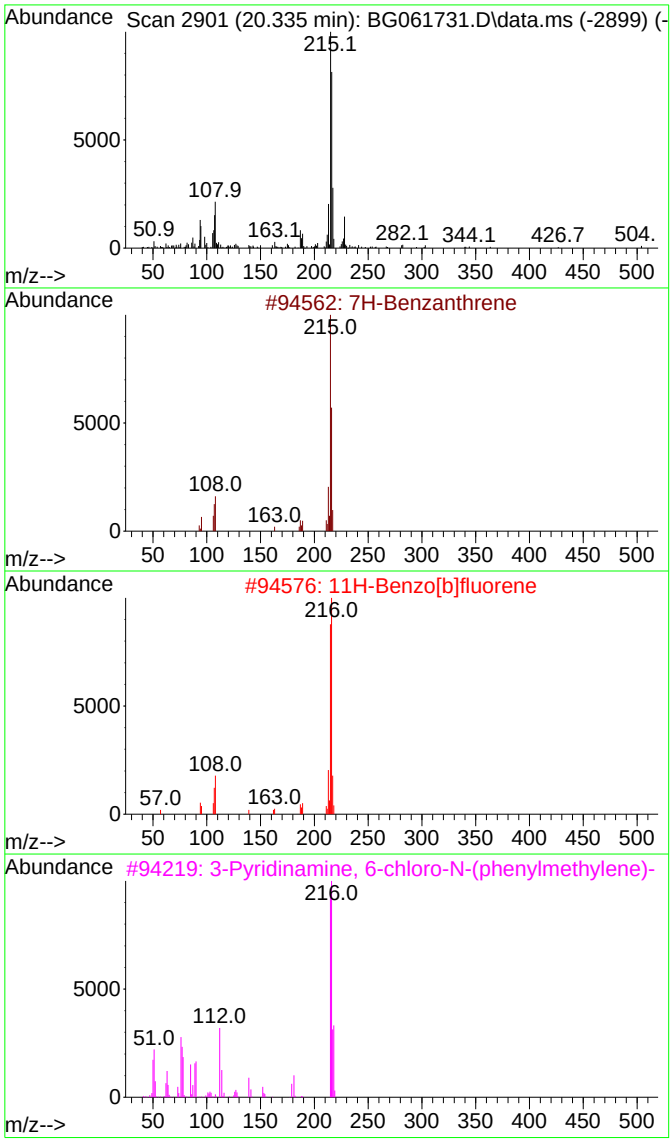
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

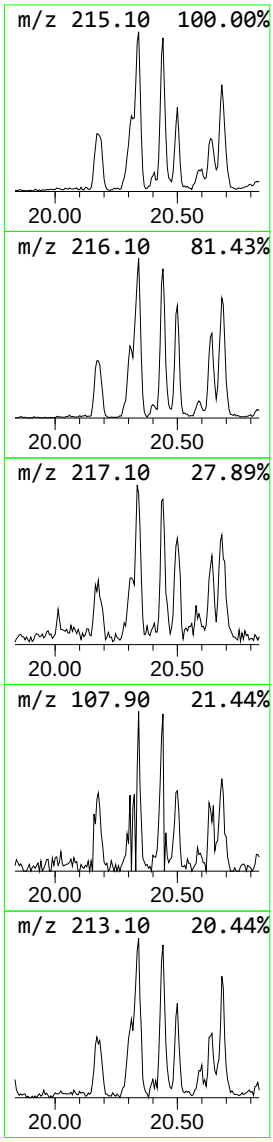
**********

Peak Number 42 7H-Benzanthrene Concentration Rank 28

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 20.335 | 3.81 ng/ul | 390023 | Chrysene-d12 | 21.710 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 7H-Benzanthrene | 216 | C17H12 | 000199-94-0 | 76 |
| 2 | 11H-Benzo[b]fluorene | 216 | C17H12 | 000243-17-4 | 53 |
| 3 | 3-Pyridinamine, 6-chloro-N-(phen... | 216 | C12H9ClN2 | 005325-71-3 | 50 |
| 4 | Fluoranthene, 2-methyl- | 216 | C17H12 | 033543-31-6 | 47 |
| 5 | Pyrene, 1-methyl- | 216 | C17H12 | 002381-21-7 | 47 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062424\
Data File : BG061731.D
Acq On : 26 Jun 2024 11:43
Operator : MA/JU
Sample : P2873-09DL 10X
Misc :
ALS Vial : 30 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
C0SJ9DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

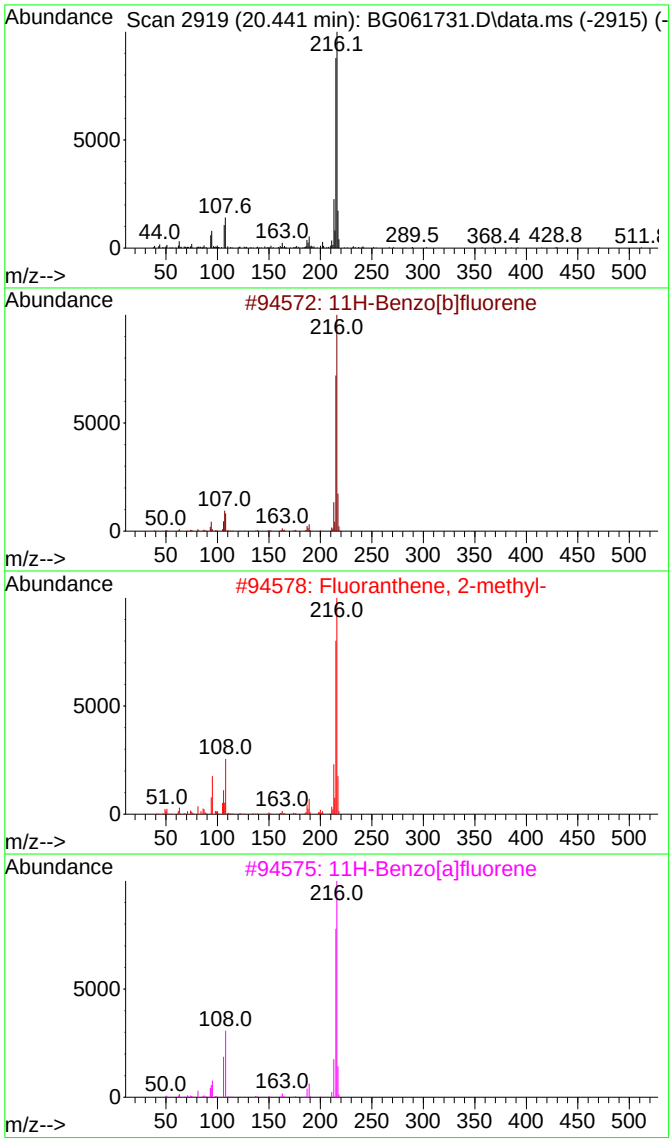
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

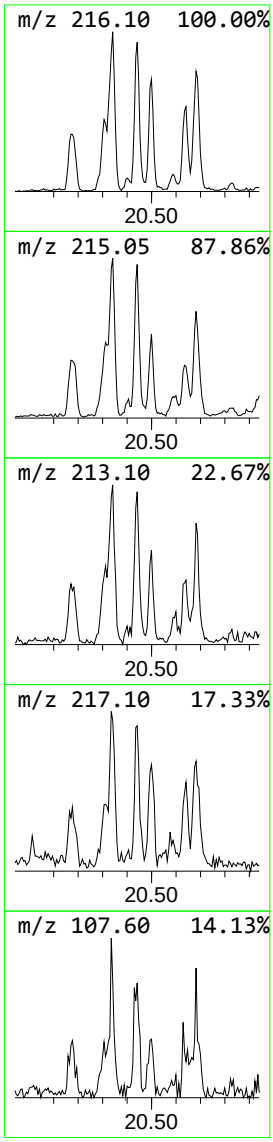
**********

Peak Number 43 11H-Benzo[b]fluorene Concentration Rank 35

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 20.441 | 2.84 ng/ul | 291125 | Chrysene-d12 | 21.710 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 11H-Benzo[b]fluorene | 216 | C17H12 | 000243-17-4 | 95 |
| 2 | Fluoranthene, 2-methyl- | 216 | C17H12 | 033543-31-6 | 87 |
| 3 | 11H-Benzo[a]fluorene | 216 | C17H12 | 000238-84-6 | 87 |
| 4 | Pyrene, 1-methyl- | 216 | C17H12 | 002381-21-7 | 83 |
| 5 | 11H-Benzo[a]fluorene | 216 | C17H12 | 000238-84-6 | 81 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062424\
Data File : BG061731.D
Acq On : 26 Jun 2024 11:43
Operator : MA/JU
Sample : P2873-09DL 10X
Misc :
ALS Vial : 30 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
C0SJ9DL

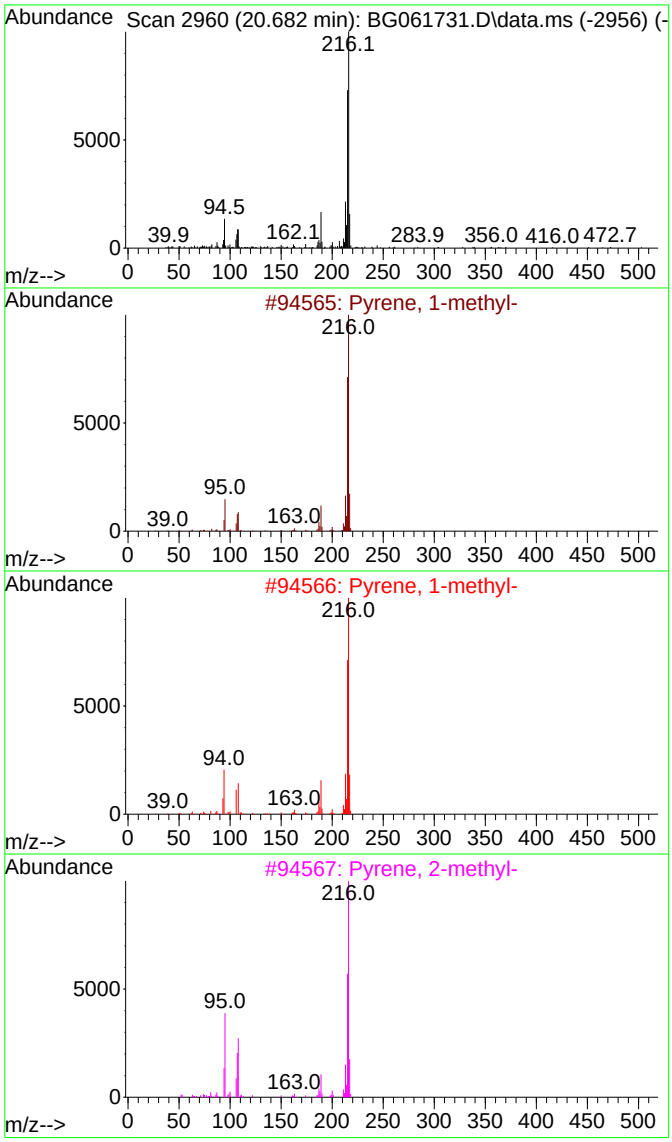
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
Quant Title : SVOA CALIBRATION

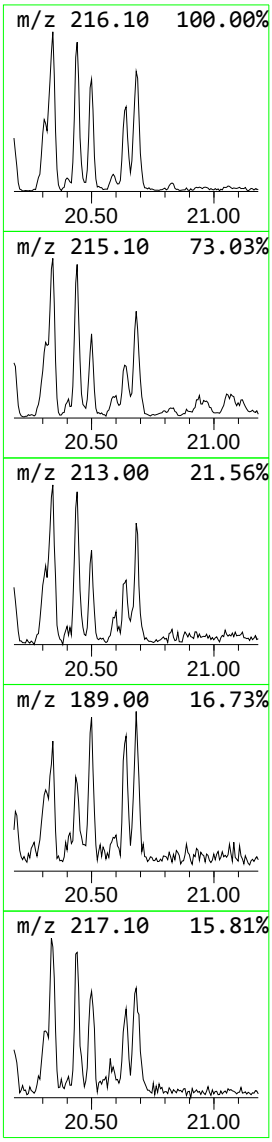
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 44 Pyrene, 1-methyl- Concentration Rank 33

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 20.682 | 2.96 ng/ul | 302801 | Chrysene-d12 | 21.710 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Pyrene, 1-methyl- | 216 | C17H12 | 002381-21-7 | 95 |
| 2 | Pyrene, 1-methyl- | 216 | C17H12 | 002381-21-7 | 91 |
| 3 | Pyrene, 2-methyl- | 216 | C17H12 | 003442-78-2 | 90 |
| 4 | Pyrene, 1-methyl- | 216 | C17H12 | 002381-21-7 | 90 |
| 5 | Pyrene, 1-methyl- | 216 | C17H12 | 002381-21-7 | 87 |



Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062424\
Data File : BG061731.D
Acq On : 26 Jun 2024 11:43
Operator : MA/JU
Sample : P2873-09DL 10X
Misc :
ALS Vial : 30 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
C0SJ9DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.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-ethy... | 7.151 | 2.9 | ng/ul | 56929 | 1 | 8.073 | 399993 | 20.0 |
| Bicyclo[4.2.0]o... | 7.721 | 26.5 | ng/ul | 529629 | 1 | 8.073 | 399993 | 20.0 |
| Benzene, 1-ethe... | 7.809 | 4.7 | ng/ul | 94474 | 1 | 8.073 | 399993 | 20.0 |
| Benzene, 1,2,3-... | 8.185 | 4.2 | ng/ul | 84592 | 1 | 8.073 | 399993 | 20.0 |
| 1-Propyne, 3-ph... | 8.614 | 41.7 | ng/ul | 833918 | 1 | 8.073 | 399993 | 20.0 |
| Benzene, 2-ethe... | 9.342 | 5.2 | ng/ul | 104266 | 1 | 8.073 | 399993 | 20.0 |
| 1H-Indene, 3-me... | 10.306 | 10.5 | ng/ul | 344187 | 2 | 10.887 | 656801 | 20.0 |
| Benzene, 1-buty... | 10.376 | 6.3 | ng/ul | 205120 | 2 | 10.887 | 656801 | 20.0 |
| Naphthalene, 2,... | 13.866 | 5.8 | ng/ul | 312623 | 3 | 14.701 | 1087660 | 20.0 |
| Naphthalene, 2,... | 14.019 | 12.1 | ng/ul | 657174 | 3 | 14.701 | 1087660 | 20.0 |
| Naphthalene, 2-... | 14.154 | 7.8 | ng/ul | 425906 | 3 | 14.701 | 1087660 | 20.0 |
| Naphthalene, 1,... | 14.260 | 4.8 | ng/ul | 261244 | 3 | 14.701 | 1087660 | 20.0 |
| Naphthalene, 2-... | 15.300 | 5.3 | ng/ul | 286713 | 3 | 14.701 | 1087660 | 20.0 |
| Fluorene-9-meth... | 15.564 | 4.3 | ng/ul | 234112 | 3 | 14.701 | 1087660 | 20.0 |
| Benzene, 5-hept... | 15.946 | 2.2 | ng/ul | 117604 | 3 | 14.701 | 1087660 | 20.0 |
| Benzaldehyde, 4... | 16.158 | 7.0 | ng/ul | 388997 | 4 | 17.445 | 1118050 | 20.0 |
| (DEL) Alkane: S... | 16.404 | 2.3 | ng/ul | 129942 | 4 | 17.445 | 1118050 | 20.0 |
| 9H-Fluorene, 1-... | 16.739 | 5.5 | ng/ul | 307667 | 4 | 17.445 | 1118050 | 20.0 |
| Dibenzothiophene | 17.262 | 6.6 | ng/ul | 366524 | 4 | 17.445 | 1118050 | 20.0 |
| 2-Methyldibenzo... | 18.020 | 3.3 | ng/ul | 185547 | 4 | 17.445 | 1118050 | 20.0 |
| 4-Methylnaphtho... | 18.167 | 3.0 | ng/ul | 169000 | 4 | 17.445 | 1118050 | 20.0 |
| Phenanthrene, 4... | 18.320 | 11.6 | ng/ul | 650436 | 4 | 17.445 | 1118050 | 20.0 |
| Anthracene, 2-m... | 18.367 | 12.3 | ng/ul | 687892 | 4 | 17.445 | 1118050 | 20.0 |
| Phenanthrene, 1... | 18.449 | 4.8 | ng/ul | 266971 | 4 | 17.445 | 1118050 | 20.0 |
| 4H-Cyclopenta[d... | 18.508 | 18.8 | ng/ul | 1050220 | 4 | 17.445 | 1118050 | 20.0 |
| 6-Phenylbenzocy... | 18.814 | 5.1 | ng/ul | 287030 | 4 | 17.445 | 1118050 | 20.0 |
| Phenanthrene, 3... | 19.225 | 6.8 | ng/ul | 378632 | 4 | 17.445 | 1118050 | 20.0 |
| Benzene, 1,1'-(... | 19.636 | 5.1 | ng/ul | 526057 | 5 | 21.710 | 2047680 | 20.0 |
| 7H-Benzanthrene | 20.335 | 3.8 | ng/ul | 390023 | 5 | 21.710 | 2047680 | 20.0 |
| 11H-Benzo[b]flu... | 20.441 | 2.8 | ng/ul | 291125 | 5 | 21.710 | 2047680 | 20.0 |
| Pyrene, 1-methyl- | 20.682 | 3.0 | ng/ul | 302801 | 5 | 21.710 | 2047680 | 20.0 |