LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049714.D
Acq On : 7 Aug 2021 21:35
Operator : CG/JU
Sample : M3244-05
Misc :
ALS Vial : 47 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** JCE05

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-BG080421.M
Title : SVOA CALIBRATION

Signal : TIC: BG049714.D\data.ms

| peak # | R.T. min | first scan | max scan | last scan | PK TY | peak height | corr. area | corr. % max. | % of total |
|---|---|---|---|---|---|---|---|---|---|
| 1 | 3.078 | 3 | 7 | 15 | rVV3 | 76826 | 184169 | 4.79% | 0.619% |
| 2 | 3.161 | 15 | 21 | 34 | rVB | 323477 | 664338 | 17.26% | 2.232% |
| 3 | 3.425 | 61 | 66 | 72 | rBV6 | 9042 | 17936 | 0.47% | 0.060% |
| 4 | 3.860 | 136 | 140 | 150 | rBV | 43318 | 89025 | 2.31% | 0.299% |
| 5 | 3.965 | 153 | 158 | 163 | rVB3 | 11259 | 18257 | 0.47% | 0.061% |
| 6 | 4.412 | 228 | 234 | 239 | rBV3 | 15842 | 26923 | 0.70% | 0.090% |
| 7 | 4.547 | 255 | 257 | 268 | rVB5 | 8737 | 18127 | 0.47% | 0.061% |
| 8 | 4.829 | 300 | 305 | 312 | rVB3 | 12413 | 19544 | 0.51% | 0.066% |
| 9 | 5.358 | 388 | 395 | 398 | rBV5 | 5039 | 10468 | 0.27% | 0.035% |
| 10 | 5.522 | 418 | 423 | 430 | rVB3 | 7007 | 14922 | 0.39% | 0.050% |
| 11 | 6.027 | 504 | 509 | 513 | rBV6 | 5366 | 10407 | 0.27% | 0.035% |
| 12 | 6.715 | 621 | 626 | 632 | rBV4 | 22845 | 40312 | 1.05% | 0.135% |
| 13 | 7.196 | 700 | 708 | 719 | rVB | 202312 | 384997 | 10.00% | 1.294% |
| 14 | 7.361 | 730 | 736 | 745 | rBV | 132254 | 233829 | 6.08% | 0.786% |
| 15 | 7.478 | 750 | 756 | 765 | rVV6 | 7956 | 16642 | 0.43% | 0.056% |
| 16 | 7.572 | 765 | 772 | 780 | rVB | 203978 | 350363 | 9.10% | 1.177% |
| 17 | 8.042 | 845 | 852 | 859 | rBV | 121567 | 214022 | 5.56% | 0.719% |
| 18 | 8.171 | 870 | 874 | 880 | rVB3 | 7354 | 10961 | 0.28% | 0.037% |
| 19 | 8.318 | 893 | 899 | 905 | rBV3 | 16289 | 29037 | 0.75% | 0.098% |
| 20 | 8.747 | 966 | 972 | 985 | rVV | 187601 | 341791 | 8.88% | 1.148% |
| 21 | 8.870 | 987 | 993 | 998 | rVV3 | 6721 | 14137 | 0.37% | 0.047% |
| 22 | 9.211 | 1043 | 1051 | 1059 | rBV | 205684 | 371134 | 9.64% | 1.247% |
| 23 | 9.364 | 1072 | 1077 | 1082 | rVB6 | 12765 | 23128 | 0.60% | 0.078% |
| 24 | 9.940 | 1166 | 1175 | 1183 | rBV | 188015 | 349717 | 9.09% | 1.175% |
| 25 | 10.480 | 1260 | 1267 | 1279 | rBV | 253598 | 463991 | 12.06% | 1.559% |
| 26 | 10.627 | 1286 | 1292 | 1297 | rBV3 | 13759 | 23944 | 0.62% | 0.080% |
| 27 | 10.862 | 1321 | 1332 | 1339 | rBV | 165625 | 311680 | 8.10% | 1.047% |
| 28 | 11.003 | 1349 | 1356 | 1361 | rBV4 | 11425 | 20381 | 0.53% | 0.068% |
| 29 | 11.655 | 1463 | 1467 | 1474 | rBV5 | 8263 | 15516 | 0.40% | 0.052% |
| 30 | 11.796 | 1487 | 1491 | 1496 | rVV7 | 8211 | 14416 | 0.37% | 0.048% |
| 31 | 12.272 | 1565 | 1572 | 1575 | rBV4 | 6129 | 8709 | 0.23% | 0.029% |
| 32 | 13.041 | 1699 | 1703 | 1709 | rBV7 | 4323 | 8807 | 0.23% | 0.030% |
| 33 | 13.141 | 1713 | 1720 | 1724 | rVB3 | 10081 | 16303 | 0.42% | 0.055% |
| 34 | 13.511 | 1778 | 1783 | 1786 | rVV4 | 7015 | 12016 | 0.31% | 0.040% |
| 35 | 13.576 | 1789 | 1794 | 1796 | rVV6 | 9241 | 15591 | 0.41% | 0.052% |
| 36 | 13.605 | 1796 | 1799 | 1809 | rVV5 | 16848 | 32607 | 0.85% | 0.110% |

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049714.D
Acq On : 7 Aug 2021 21:35
Operator : CG/JU
Sample : M3244-05
Misc :
ALS Vial : 47 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** JCE05

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-BG080421.M
Title : SVOA CALIBRATION

| | | | | | | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 37 | 14.093 | 1875 | 1882 | 1888 | rVV | 470311 | 698921 | 18.16% | 2.348% |
| 38 | 14.328 | 1918 | 1922 | 1926 | rBV5 | 8341 | 17470 | 0.45% | 0.059% |
| 39 | 14.387 | 1926 | 1932 | 1940 | rVB | 476082 | 775064 | 20.14% | 2.604% |
| 40 | 14.574 | 1961 | 1964 | 1969 | rVB3 | 7033 | 9932 | 0.26% | 0.033% |
| 41 | 14.692 | 1975 | 1984 | 1993 | rBV | 257738 | 406395 | 10.56% | 1.365% |
| 42 | 14.892 | 2011 | 2018 | 2029 | rBV | 147270 | 289255 | 7.52% | 0.972% |
| 43 | 15.039 | 2037 | 2043 | 2049 | rBV2 | 31590 | 50829 | 1.32% | 0.171% |
| 44 | 15.091 | 2049 | 2052 | 2057 | rVB5 | 9172 | 11482 | 0.30% | 0.039% |
| 45 | 15.426 | 2105 | 2109 | 2113 | rBV | 26846 | 36150 | 0.94% | 0.121% |
| 46 | 15.491 | 2113 | 2120 | 2124 | rBV4 | 14689 | 35679 | 0.93% | 0.120% |
| 47 | 15.685 | 2146 | 2153 | 2161 | rBV | 609744 | 967478 | 25.14% | 3.251% |
| 48 | 15.802 | 2167 | 2173 | 2181 | rBV | 184054 | 291133 | 7.56% | 0.978% |
| 49 | 15.926 | 2190 | 2194 | 2199 | rVB4 | 7657 | 14121 | 0.37% | 0.047% |
| 50 | 16.237 | 2242 | 2247 | 2252 | rBV5 | 16882 | 29247 | 0.76% | 0.098% |
| 51 | 16.390 | 2267 | 2273 | 2280 | rBV3 | 12880 | 24138 | 0.63% | 0.081% |
| 52 | 16.572 | 2296 | 2304 | 2309 | rBV4 | 28803 | 44112 | 1.15% | 0.148% |
| 53 | 16.631 | 2309 | 2314 | 2318 | rBV8 | 5485 | 11673 | 0.30% | 0.039% |
| 54 | 16.830 | 2343 | 2348 | 2353 | rVB6 | 15198 | 22393 | 0.58% | 0.075% |
| 55 | 17.194 | 2404 | 2410 | 2419 | rBV7 | 27390 | 67665 | 1.76% | 0.227% |
| 56 | 17.324 | 2427 | 2432 | 2440 | rBV7 | 20890 | 41807 | 1.09% | 0.140% |
| 57 | 17.388 | 2440 | 2443 | 2446 | rVV4 | 8121 | 8726 | 0.23% | 0.029% |
| 58 | 17.441 | 2446 | 2452 | 2458 | rVV | 291299 | 452913 | 11.77% | 1.522% |
| 59 | 17.547 | 2462 | 2470 | 2477 | rVV | 564965 | 898854 | 23.35% | 3.020% |
| 60 | 17.653 | 2484 | 2488 | 2492 | rVB6 | 11955 | 16247 | 0.42% | 0.055% |
| 61 | 17.823 | 2512 | 2517 | 2519 | rBV5 | 8373 | 11794 | 0.31% | 0.040% |
| 62 | 17.923 | 2530 | 2534 | 2540 | rBV | 23892 | 30040 | 0.78% | 0.101% |
| 63 | 18.099 | 2560 | 2564 | 2568 | rVB3 | 24053 | 29122 | 0.76% | 0.098% |
| 64 | 18.211 | 2577 | 2583 | 2587 | rBV5 | 15374 | 27380 | 0.71% | 0.092% |
| 65 | 18.269 | 2590 | 2593 | 2598 | rVB4 | 27061 | 36621 | 0.95% | 0.123% |
| 66 | 18.328 | 2598 | 2603 | 2612 | rBV2 | 183248 | 280265 | 7.28% | 0.942% |
| 67 | 18.616 | 2648 | 2652 | 2655 | rBV3 | 21865 | 31032 | 0.81% | 0.104% |
| 68 | 18.974 | 2709 | 2713 | 2717 | rBV4 | 33830 | 50088 | 1.30% | 0.168% |
| 69 | 19.016 | 2717 | 2720 | 2724 | rVB4 | 54037 | 70721 | 1.84% | 0.238% |
| 70 | 19.115 | 2733 | 2737 | 2743 | rVB8 | 14396 | 23097 | 0.60% | 0.078% |
| 71 | 19.186 | 2745 | 2749 | 2752 | rBV4 | 24445 | 45138 | 1.17% | 0.152% |
| 72 | 19.256 | 2755 | 2761 | 2765 | rVV3 | 90060 | 192358 | 5.00% | 0.646% |
| 73 | 19.297 | 2765 | 2768 | 2770 | rVB2 | 23081 | 24351 | 0.63% | 0.082% |
| 74 | 19.480 | 2791 | 2799 | 2805 | rBV4 | 88630 | 184445 | 4.79% | 0.620% |
| 75 | 19.532 | 2806 | 2808 | 2811 | rVV3 | 32410 | 31464 | 0.82% | 0.106% |
| 76 | 19.597 | 2815 | 2819 | 2827 | rBV6 | 42741 | 68367 | 1.78% | 0.230% |

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049714.D
Acq On : 7 Aug 2021 21:35
Operator : CG/JU
Sample : M3244-05
Misc :
ALS Vial : 47 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** JCE05

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-BG080421.M
Title : SVOA CALIBRATION

| | | | | | | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 77 | 19.832 | 2853 | 2859 | 2865 | rVB | 752247 | 1087362 | 28.25% | 3.653% |
| 78 | 20.326 | 2938 | 2943 | 2944 | rBV | 123505 | 160664 | 4.17% | 0.540% |
| 79 | 20.349 | 2944 | 2947 | 2956 | rVB | 196671 | 258938 | 6.73% | 0.870% |
| 80 | 20.525 | 2974 | 2977 | 2981 | rVB6 | 38657 | 48666 | 1.26% | 0.164% |
| 81 | 20.672 | 2999 | 3002 | 3010 | rBV4 | 73836 | 124535 | 3.24% | 0.418% |
| 82 | 20.790 | 3019 | 3022 | 3027 | rBV3 | 26437 | 39373 | 1.02% | 0.132% |
| 83 | 20.842 | 3027 | 3031 | 3038 | rVB2 | 62753 | 92769 | 2.41% | 0.312% |
| 84 | 21.030 | 3058 | 3063 | 3068 | rBV3 | 81169 | 131827 | 3.43% | 0.443% |
| 85 | 21.354 | 3113 | 3118 | 3131 | rBV | 1014618 | 1560091 | 40.54% | 5.242% |
| 86 | 21.730 | 3176 | 3182 | 3191 | rVB2 | 312480 | 535171 | 13.91% | 1.798% |
| 87 | 22.158 | 3250 | 3255 | 3261 | rVB | 202801 | 309569 | 8.04% | 1.040% |
| 88 | 22.558 | 3312 | 3323 | 3337 | rBV2 | 680967 | 2188957 | 56.88% | 7.354% |
| 89 | 22.822 | 3364 | 3368 | 3374 | rVB2 | 82356 | 129922 | 3.38% | 0.437% |
| 90 | 23.304 | 3443 | 3450 | 3458 | rBV4 | 118341 | 276071 | 7.17% | 0.928% |
| 91 | 23.592 | 3492 | 3499 | 3508 | rVB | 598440 | 1220256 | 31.71% | 4.100% |
| 92 | 24.062 | 3571 | 3579 | 3584 | rBV | 517232 | 1170071 | 30.40% | 3.931% |
| 93 | 24.126 | 3584 | 3590 | 3608 | rVB2 | 639160 | 1673868 | 43.49% | 5.624% |
| 94 | 24.490 | 3647 | 3652 | 3658 | rVB3 | 76632 | 154979 | 4.03% | 0.521% |
| 95 | 24.790 | 3694 | 3703 | 3714 | rVB2 | 321319 | 905805 | 23.54% | 3.043% |
| 96 | 25.013 | 3732 | 3741 | 3751 | rVB2 | 207186 | 618491 | 16.07% | 2.078% |
| 97 | 25.131 | 3755 | 3761 | 3771 | rVB3 | 105424 | 280564 | 7.29% | 0.943% |
| 98 | 25.560 | 3823 | 3834 | 3845 | rBV2 | 485614 | 1482045 | 38.51% | 4.979% |
| 99 | 26.188 | 3932 | 3941 | 3951 | rVB | 218490 | 709020 | 18.42% | 2.382% |
| 100 | 26.341 | 3953 | 3967 | 3987 | rBV2 | 1006374 | 3848662 | 100.00% | 12.931% |

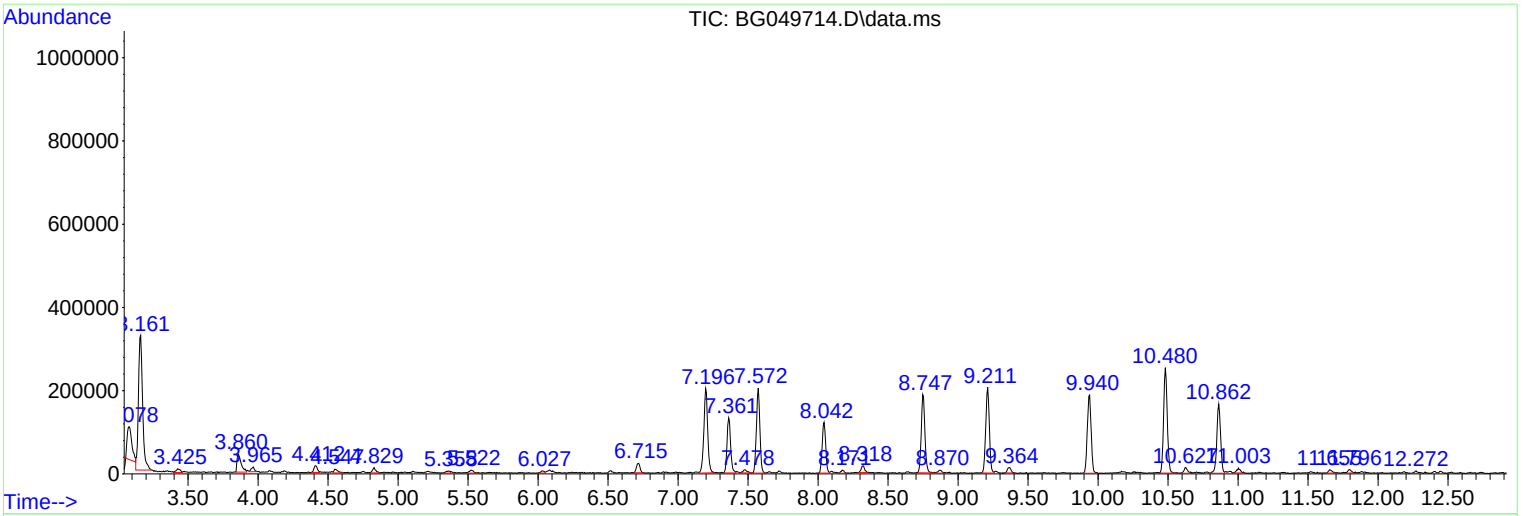
Sum of corrected areas: 29763790

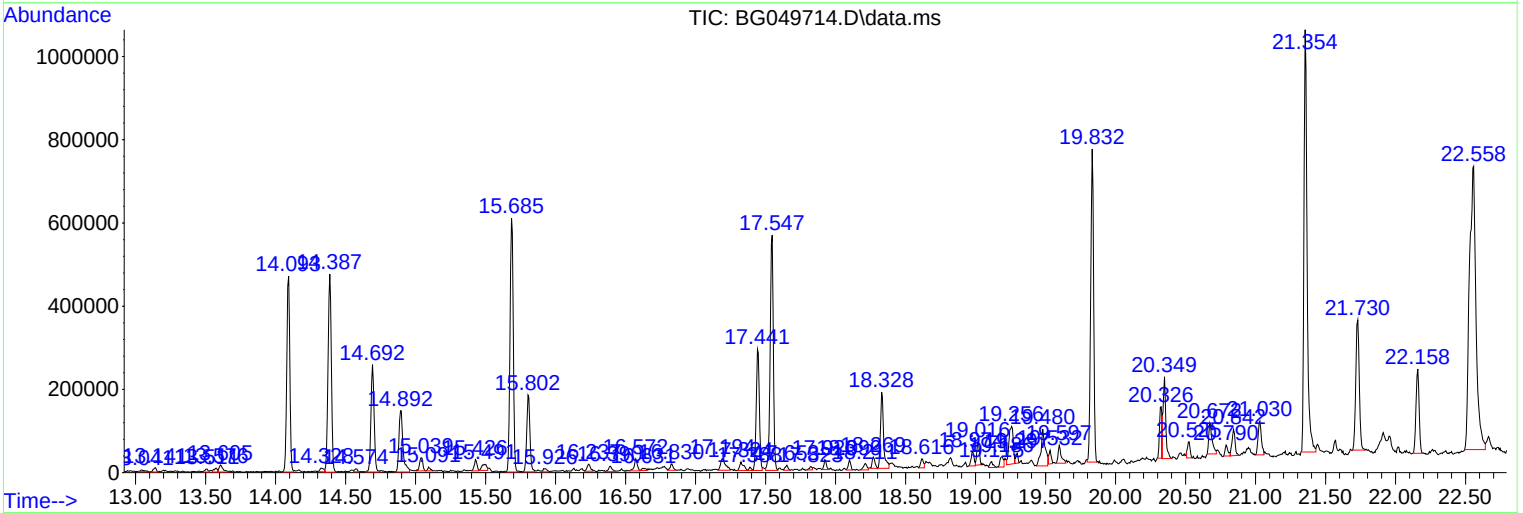
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049714.D
Acq On : 7 Aug 2021 21:35
Operator : CG/JU
Sample : M3244-05
Misc :
ALS Vial : 47 Sample Multiplier: 1

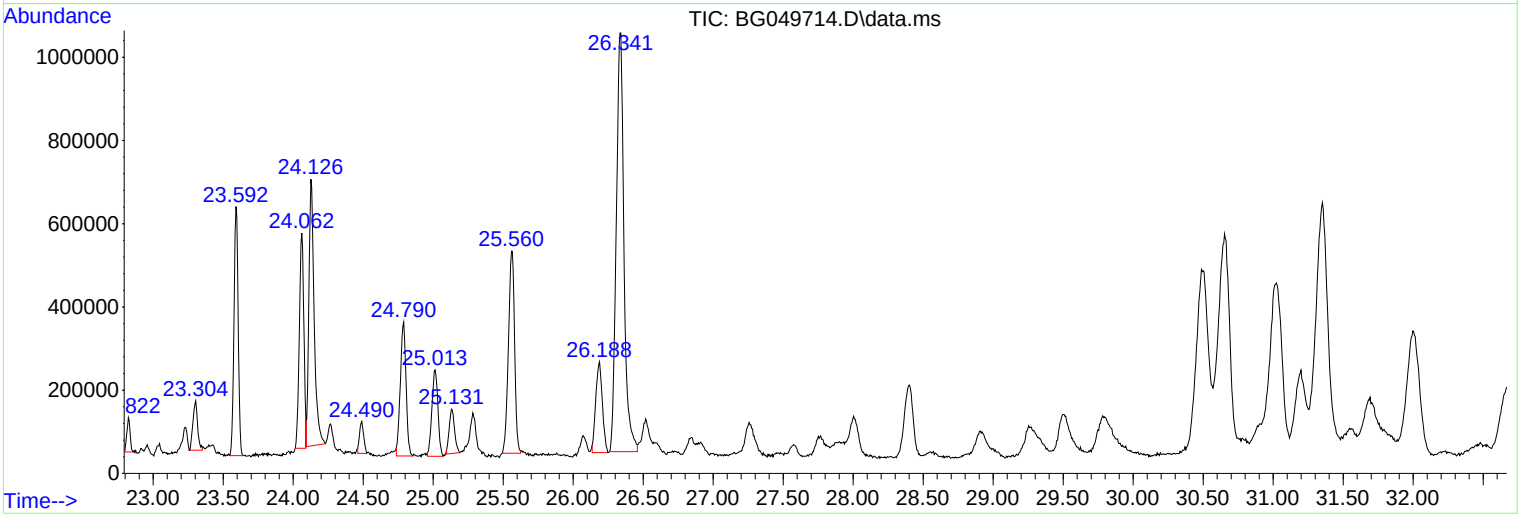
**Instrument :**
BNA_G
**ClientSampleId :**
JCE05

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049714.D
Acq On : 7 Aug 2021 21:35
Operator : CG/JU
Sample : M3244-05
Misc :
ALS Vial : 47 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** JCE05

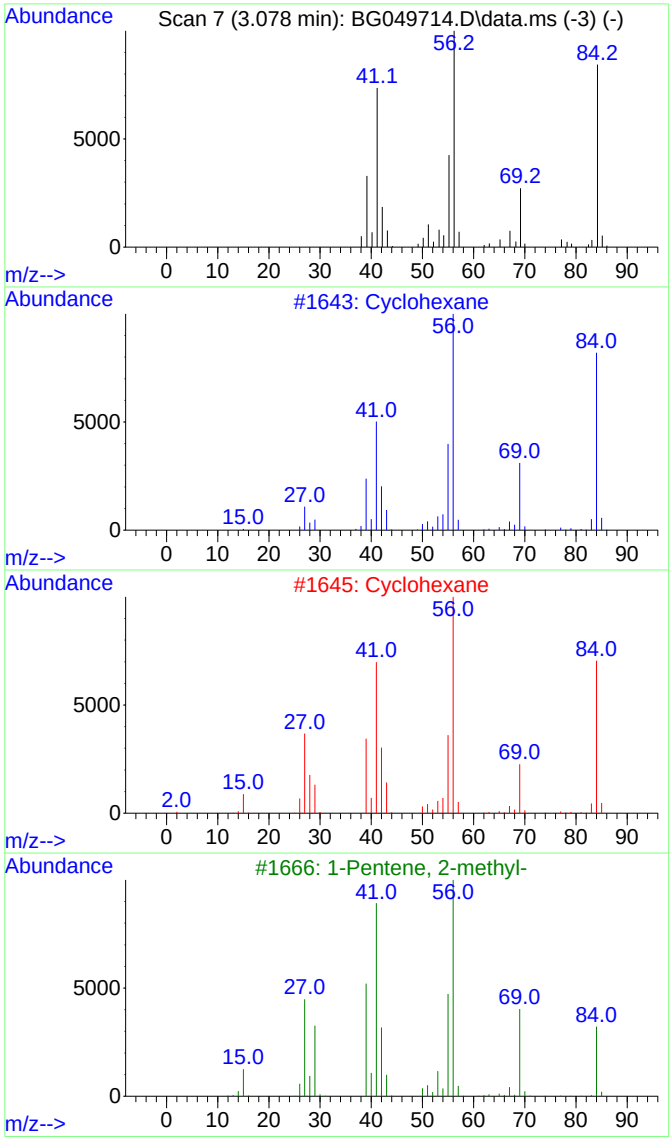
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

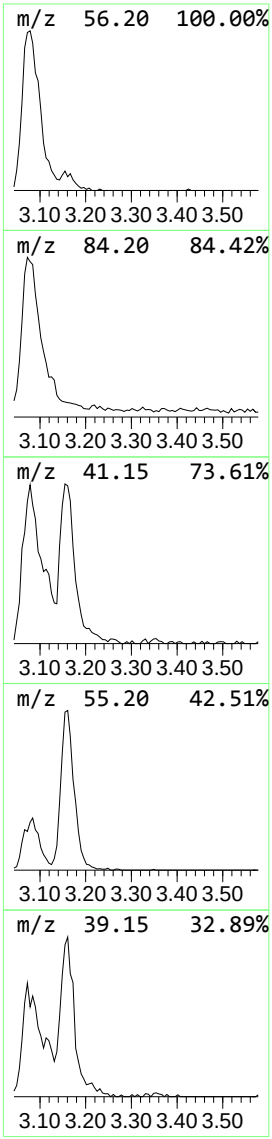
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 1 (DEL) Alkane: Cyclic3.078 Concentration Rank 10

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 3.078 | 17.21 ng/ul | 184169 | 1,4-Dichlorobenzene-d4 | 8.042 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Cyclohexane | 84 | C6H12 | 000110-82-7 | 87 |
| 2 | Cyclohexane | 84 | C6H12 | 000110-82-7 | 86 |
| 3 | 1-Pentene, 2-methyl- | 84 | C6H12 | 000763-29-1 | 86 |
| 4 | 1-Pentene, 2-methyl- | 84 | C6H12 | 000763-29-1 | 86 |
| 5 | Cyclohexane | 84 | C6H12 | 000110-82-7 | 83 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049714.D
Acq On : 7 Aug 2021 21:35
Operator : CG/JU
Sample : M3244-05
Misc :
ALS Vial : 47 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** JCE05

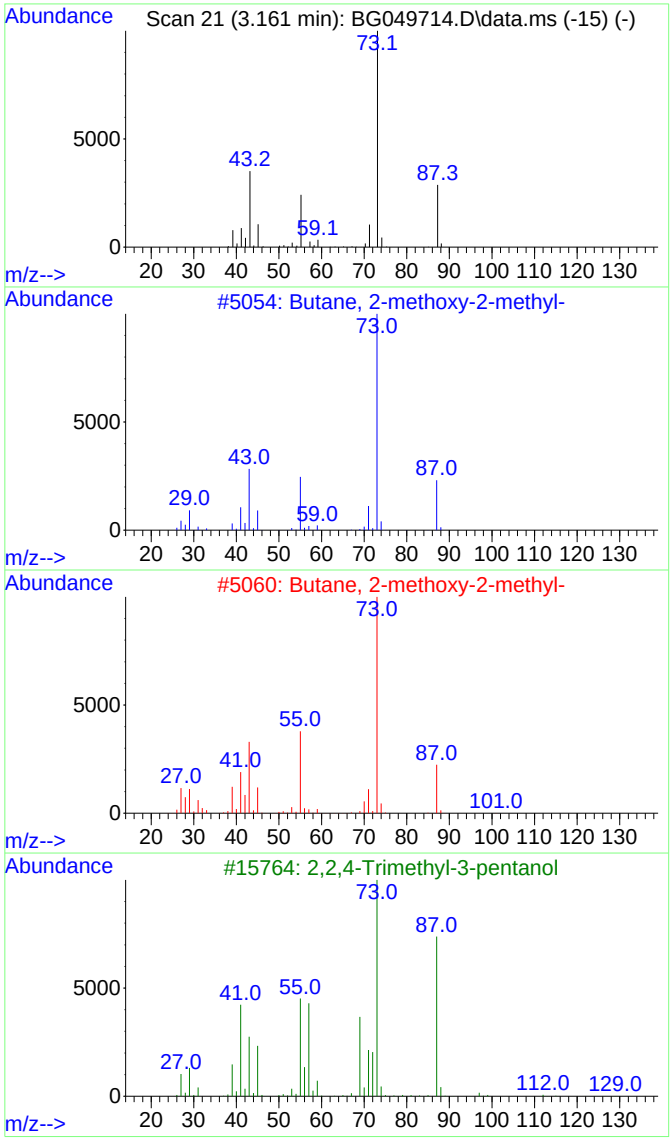
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

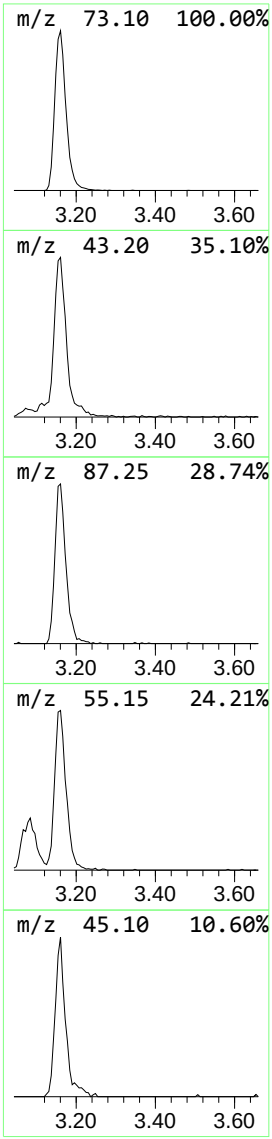
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 3

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 3.161 | 62.08 ng/ul | 664338 | 1,4-Dichlorobenzene-d4 | 8.042 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Butane, 2-methoxy-2-methyl- | 102 | C6H14O | 000994-05-8 | 78 |
| 2 | Butane, 2-methoxy-2-methyl- | 102 | C6H14O | 000994-05-8 | 74 |
| 3 | 2,2,4-Trimethyl-3-pentanol | 130 | C8H18O | 005162-48-1 | 45 |
| 4 | Acetamide, N-ethyl- | 87 | C4H9NO | 000625-50-3 | 27 |
| 5 | Pentane, 3-methoxy- | 102 | C6H14O | 036839-67-5 | 23 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049714.D
Acq On : 7 Aug 2021 21:35
Operator : CG/JU
Sample : M3244-05
Misc :
ALS Vial : 47 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** JCE05

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

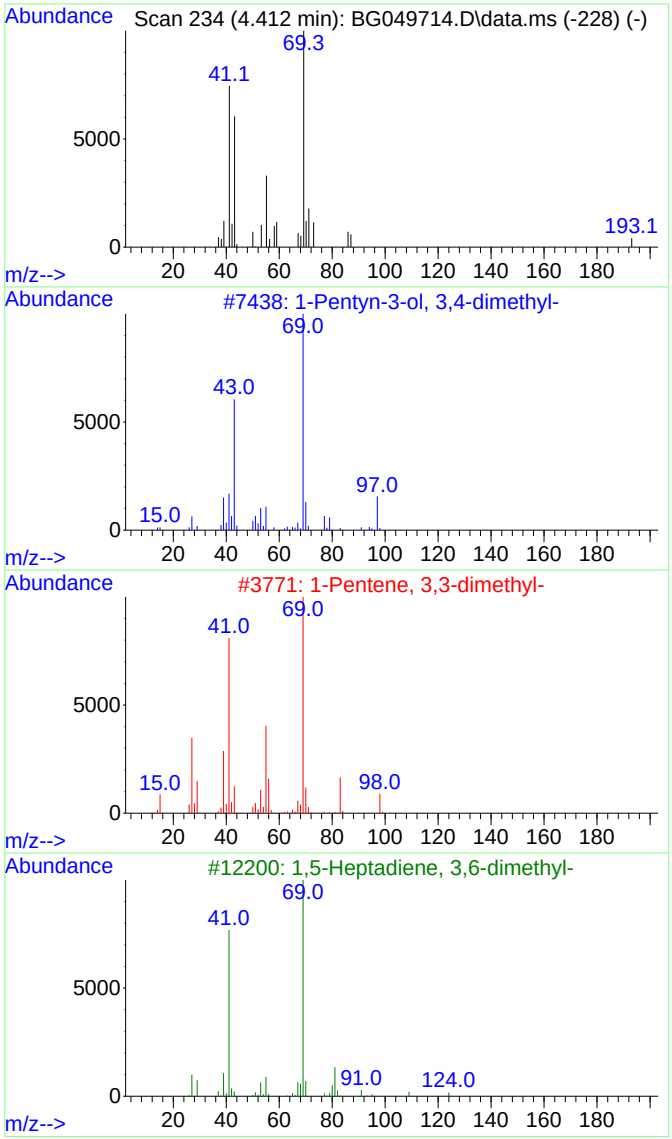
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

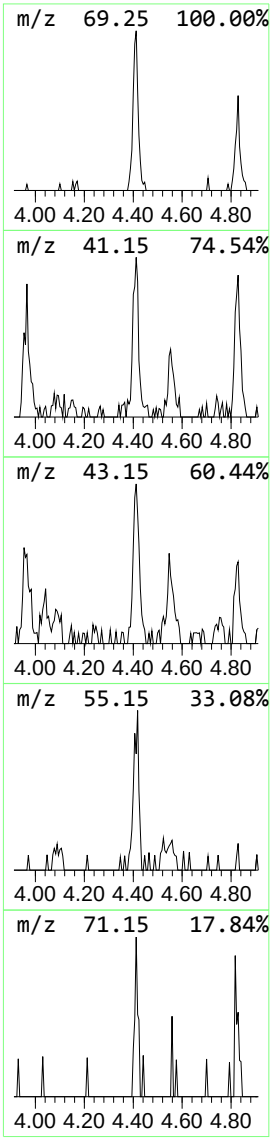
**********

Peak Number 3 unknown-01 Concentration Rank 29

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 4.412 | 2.52 ng/ul | 26923 | 1,4-Dichlorobenzene-d4 | 8.042 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1-Pentyn-3-ol, 3,4-dimethyl- | 112 | C7H12O | 001482-15-1 | 40 |
| 2 | 1-Pentene, 3,3-dimethyl- | 98 | C7H14 | 003404-73-7 | 40 |
| 3 | 1,5-Heptadiene, 3,6-dimethyl- | 124 | C9H16 | 034891-10-6 | 40 |
| 4 | 3-Butyn-2-ol, 2-methyl- | 84 | C5H8O | 000115-19-5 | 38 |
| 5 | 2-Propenoic acid, 2-methyl-, 1-m... | 128 | C7H12O2 | 004655-34-9 | 38 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049714.D
Acq On : 7 Aug 2021 21:35
Operator : CG/JU
Sample : M3244-05
Misc :
ALS Vial : 47 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** JCE05

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

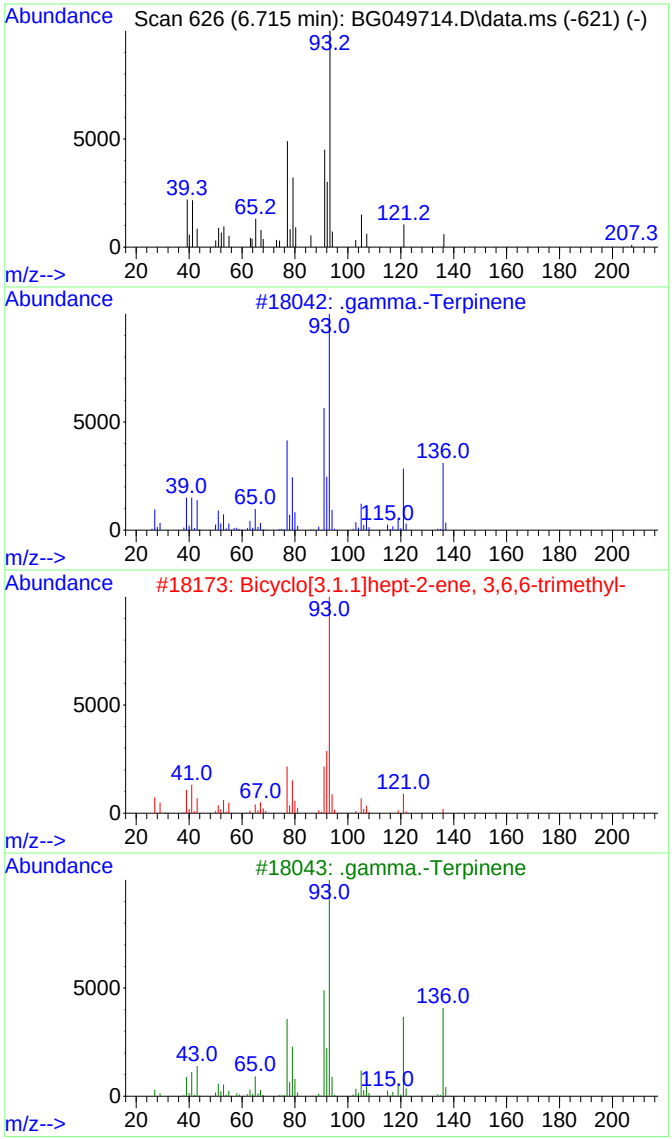
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

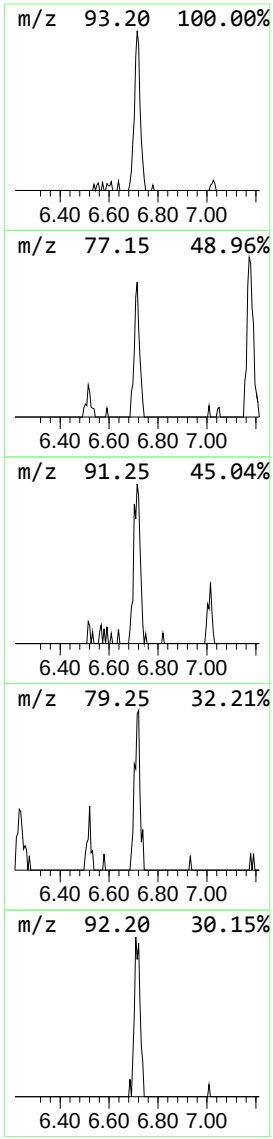
**********

Peak Number 4 .gamma.-Terpinene Concentration Rank 23

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 6.715 | 3.77 ng/ul | 40312 | 1,4-Dichlorobenzene-d4 | 8.042 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | .gamma.-Terpinene | 136 | C10H16 | 000099-85-4 | 94 |
| 2 | Bicyclo[3.1.1]hept-2-ene, 3,6,6-... | 136 | C10H16 | 004889-83-2 | 90 |
| 3 | .gamma.-Terpinene | 136 | C10H16 | 000099-85-4 | 87 |
| 4 | .alpha.-Pinene | 136 | C10H16 | 000080-56-8 | 83 |
| 5 | .alpha.-Pinene | 136 | C10H16 | 000080-56-8 | 72 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049714.D
Acq On : 7 Aug 2021 21:35
Operator : CG/JU
Sample : M3244-05
Misc :
ALS Vial : 47 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** JCE05

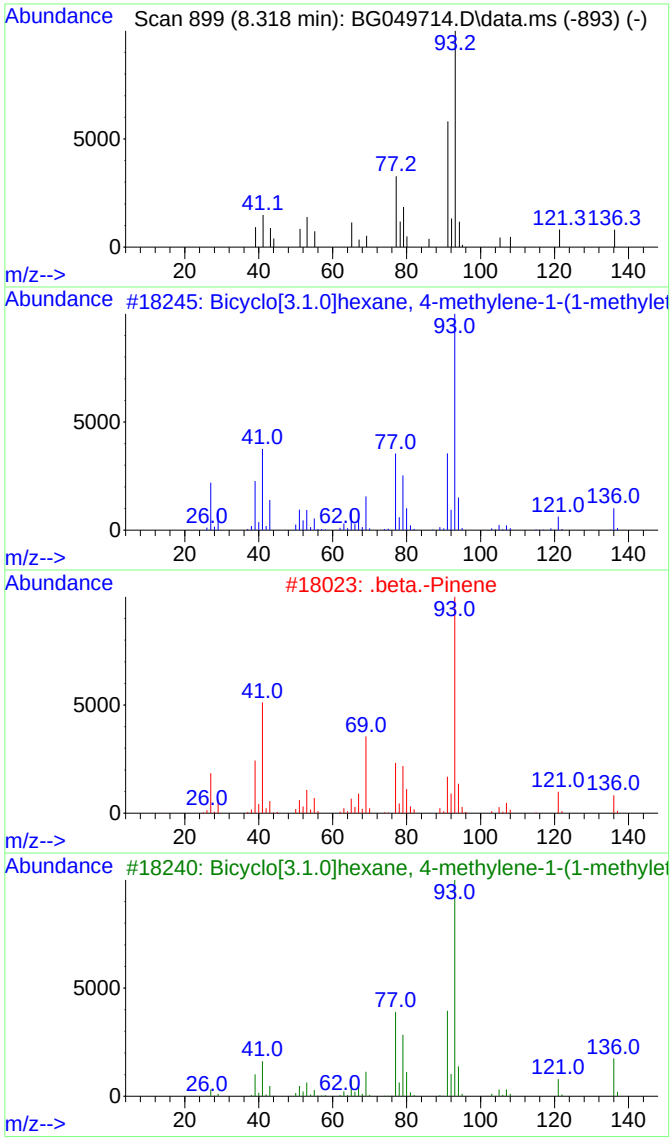
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

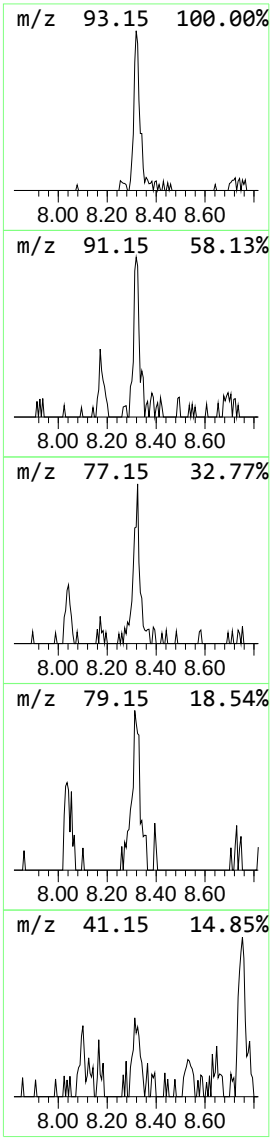
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 5 Bicyclo[3.1.0]hexane, 4-met... Concentration Rank 27

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 8.318 | 2.71 ng/ul | 29037 | 1,4-Dichlorobenzene-d4 | 8.042 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Bicyclo[3.1.0]hexane, 4-methylen... | 136 | C10H16 | 003387-41-5 | 87 |
| 2 | .beta.-Pinene | 136 | C10H16 | 000127-91-3 | 86 |
| 3 | Bicyclo[3.1.0]hexane, 4-methylen... | 136 | C10H16 | 003387-41-5 | 80 |
| 4 | .beta.-Pinene | 136 | C10H16 | 000127-91-3 | 80 |
| 5 | Bicyclo[3.1.0]hexane, 4-methylen... | 136 | C10H16 | 003387-41-5 | 80 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049714.D
Acq On : 7 Aug 2021 21:35
Operator : CG/JU
Sample : M3244-05
Misc :
ALS Vial : 47 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** JCE05

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

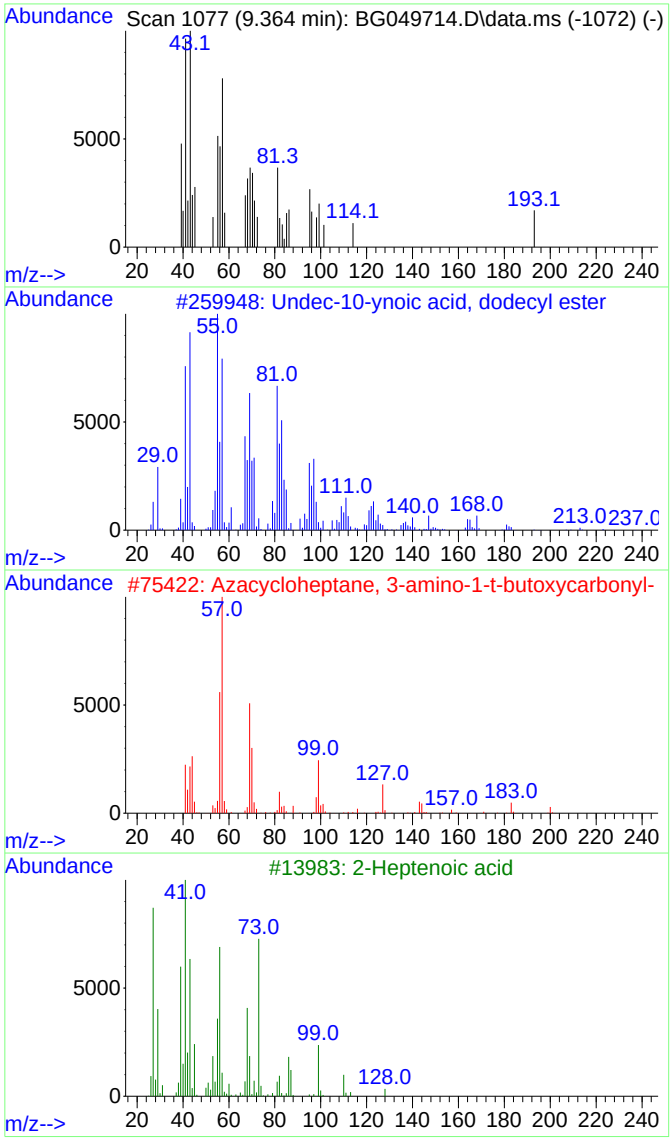
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

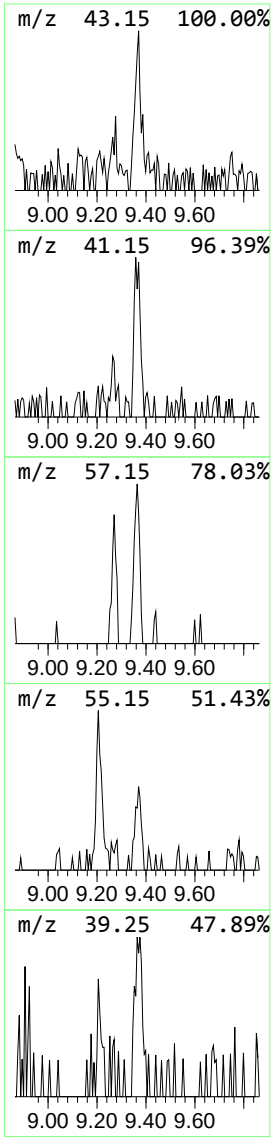
**********

Peak Number 6 unknown-02 Concentration Rank 32

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 9.364 | 2.16 ng/ul | 23128 | 1,4-Dichlorobenzene-d4 | 8.042 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Undec-10-ynoic acid, dodecyl ester | 350 | C23H42O2 | 1010406-16-5 | 38 |
| 2 | Azacycloheptane, 3-amino-1-t-but... | 200 | C10H20N2O2 | 1000126-78-8 | 35 |
| 3 | 2-Heptenoic acid | 128 | C7H12O2 | 018999-28-5 | 35 |
| 4 | trans-2,3-Epoxynonane | 142 | C9H18O | 056820-01-0 | 27 |
| 5 | 4-Tridecene, (Z)- | 182 | C13H26 | 041446-54-2 | 27 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049714.D
Acq On : 7 Aug 2021 21:35
Operator : CG/JU
Sample : M3244-05
Misc :
ALS Vial : 47 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** JCE05

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

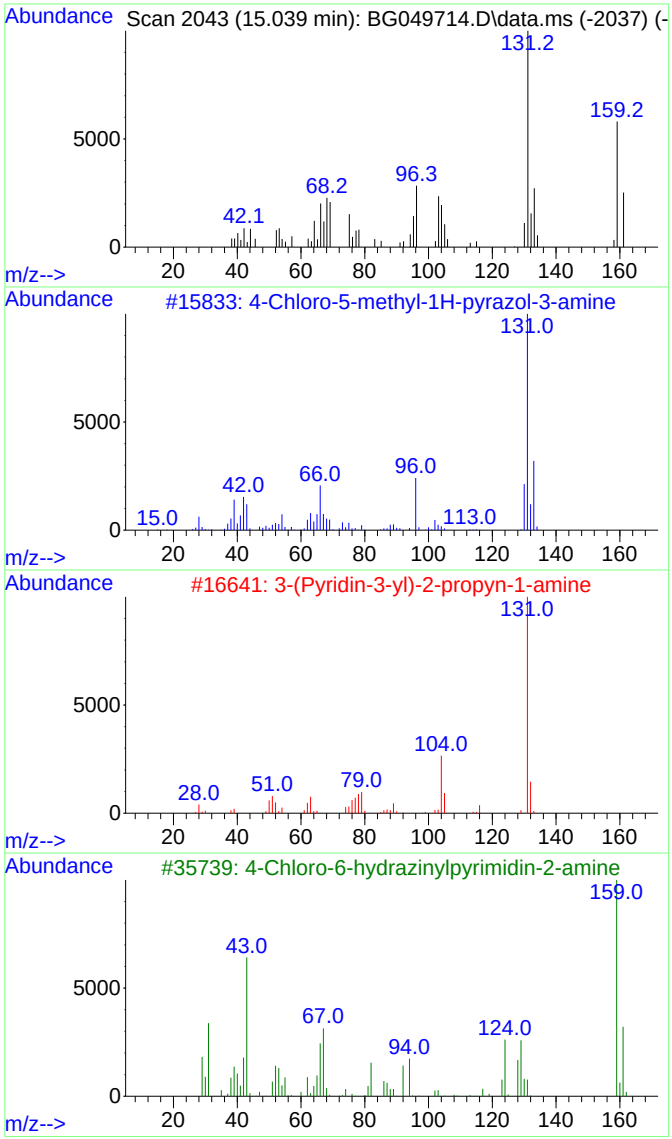
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

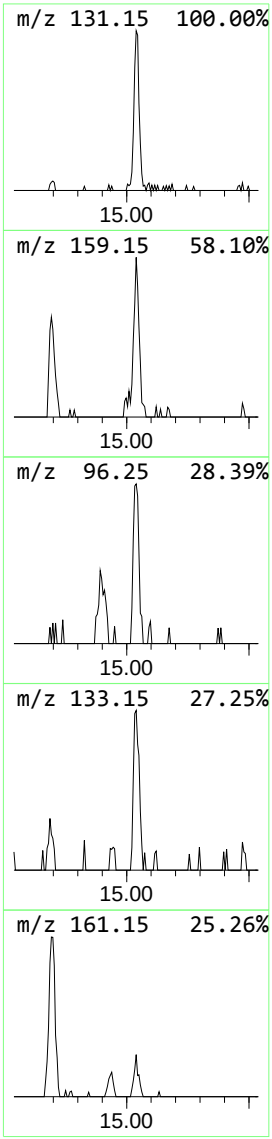
**********

Peak Number 7 unknown-03 Concentration Rank 30

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 15.039 | 2.50 ng/ul | 50829 | Acenaphthene-d10 | 14.692 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 4-Chloro-5-methyl-1H-pyrazol-3-a... | 131 | C4H6ClN3 | 110580-44-4 | 38 |
| 2 | 3-(Pyridin-3-yl)-2-propyn-1-amine | 132 | C8H8N2 | 777856-62-9 | 25 |
| 3 | 4-Chloro-6-hydrazinylpyrimidin-2... | 159 | C4H6ClN5 | 089124-04-9 | 17 |
| 4 | 2-(1H-1,2,4-Triazol-1-yl)benzyla... | 174 | C9H10N4 | 449756-97-2 | 16 |
| 5 | Isoxazole, 5-amino-4-bromo-3-phe... | 238 | C9H7BrN2O | 028884-13-1 | 12 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049714.D
Acq On : 7 Aug 2021 21:35
Operator : CG/JU
Sample : M3244-05
Misc :
ALS Vial : 47 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** JCE05

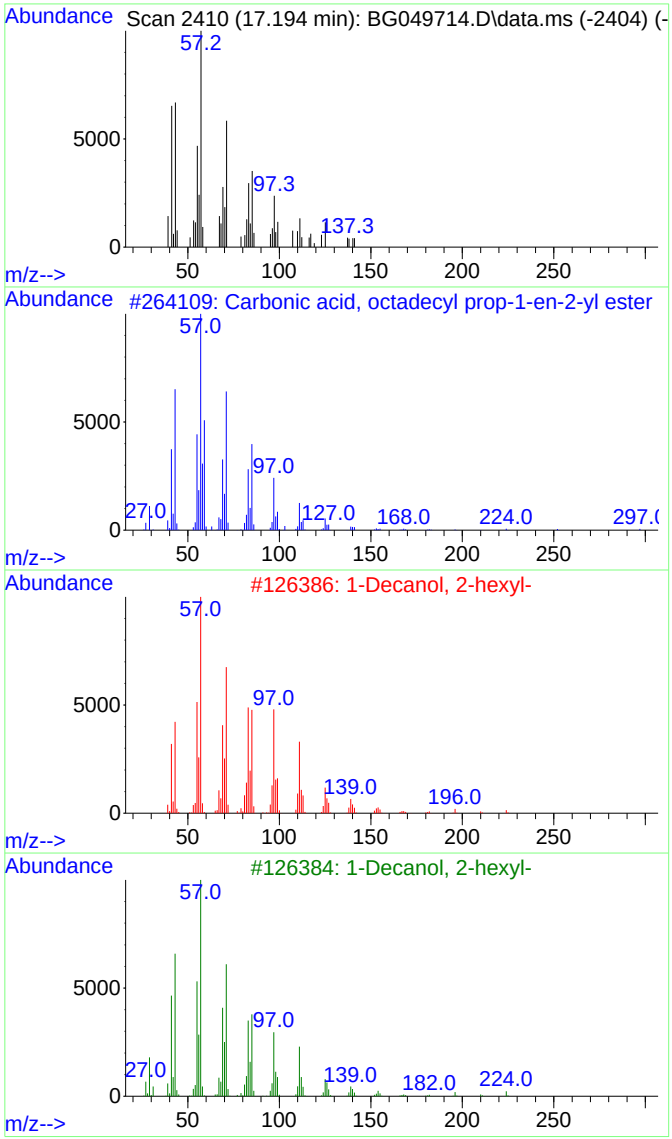
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

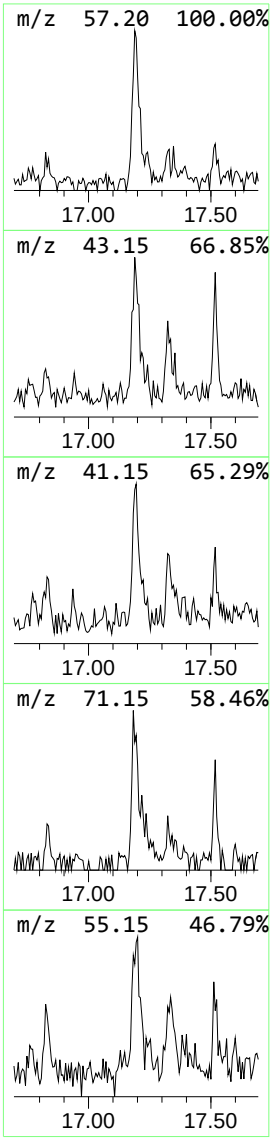
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 8 Carbonic acid, octadecyl pr... Concentration Rank 26

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 17.195 | 2.99 ng/ul | 67665 | Phenanthrene-d10 | 17.441 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Carbonic acid, octadecyl prop-1-... | 354 | C22H42O3 | 1000383-11-5 | 91 |
| 2 | 1-Decanol, 2-hexyl- | 242 | C16H34O | 002425-77-6 | 90 |
| 3 | 1-Decanol, 2-hexyl- | 242 | C16H34O | 002425-77-6 | 90 |
| 4 | Oxalic acid, allyl hexadecyl ester | 354 | C21H38O4 | 1000309-24-4 | 86 |
| 5 | Carbonic acid, eicosyl vinyl ester | 368 | C23H44O3 | 1000382-54-3 | 86 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049714.D
Acq On : 7 Aug 2021 21:35
Operator : CG/JU
Sample : M3244-05
Misc :
ALS Vial : 47 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** JCE05

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

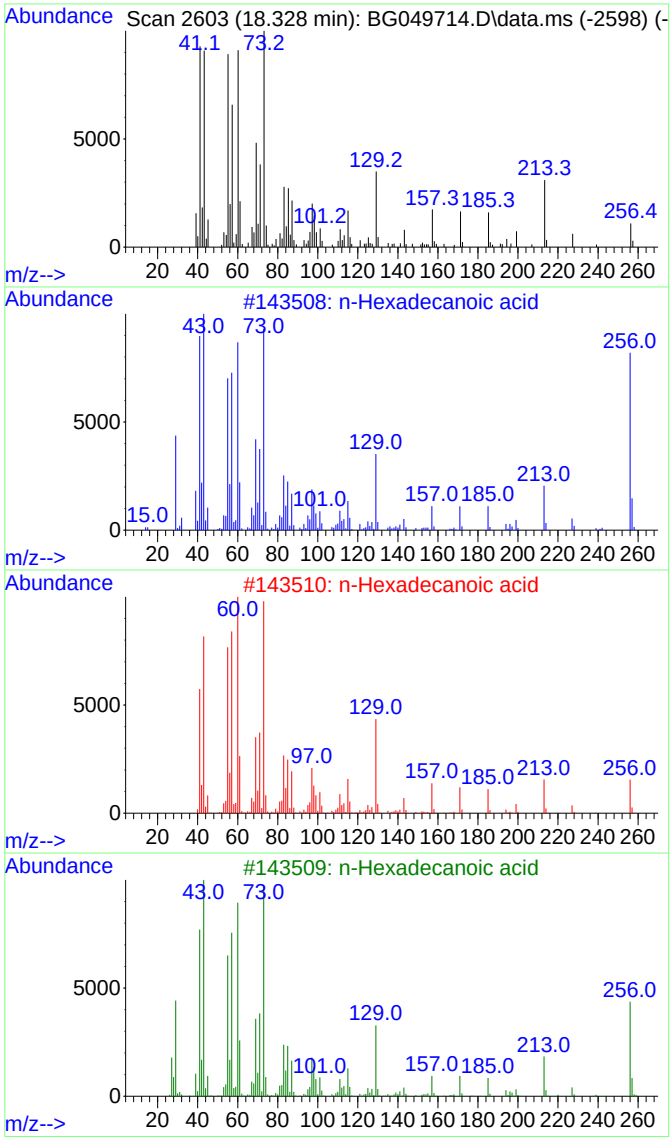
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

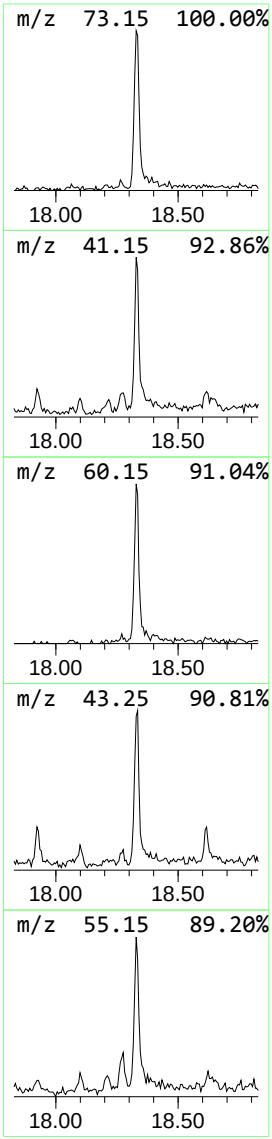
**********

Peak Number 9 n-Hexadecanoic acid Concentration Rank 11

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 18.328 | 12.38 ng/ul | 280265 | Phenanthrene-d10 | 17.441 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 99 |
| 2 | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 98 |
| 3 | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 96 |
| 4 | Tetradecanoic acid | 228 | C14H28O2 | 000544-63-8 | 83 |
| 5 | Pentadecanoic acid | 242 | C15H30O2 | 001002-84-2 | 83 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049714.D
Acq On : 7 Aug 2021 21:35
Operator : CG/JU
Sample : M3244-05
Misc :
ALS Vial : 47 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** JCE05

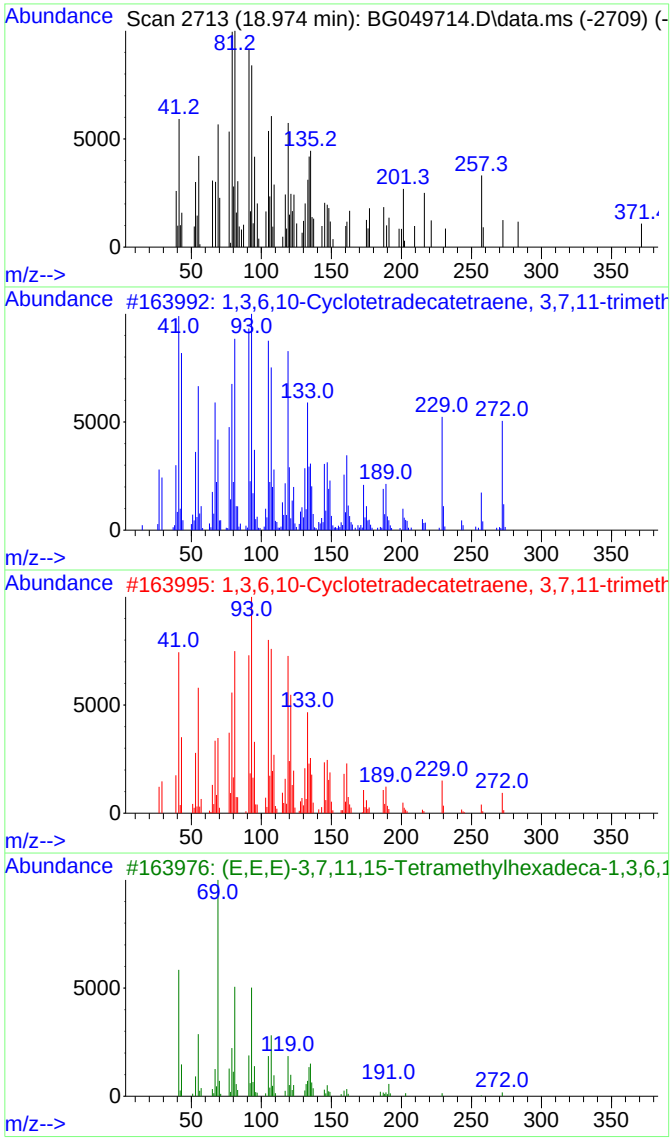
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

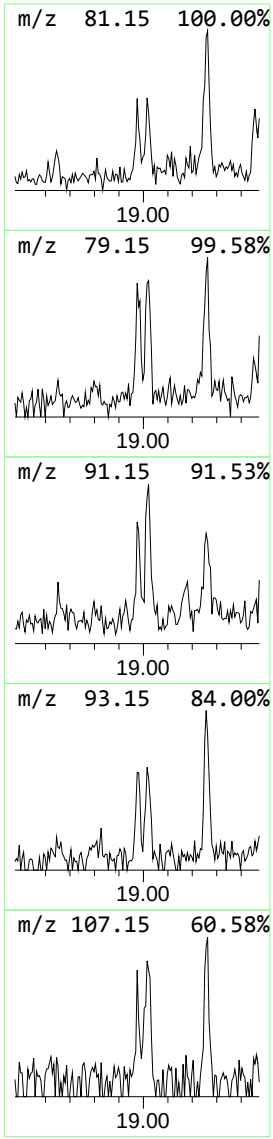
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 10 1,3,6,10-Cyclotetradecatetr... Concentration Rank 31

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 18.974 | 2.21 ng/ul | 50088 | Phenanthrene-d10 | 17.441 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1,3,6,10-Cyclotetradecatetraene,... | 272 | C20H32 | 001898-13-1 | 55 |
| 2 | 1,3,6,10-Cyclotetradecatetraene,... | 272 | C20H32 | 001898-13-1 | 49 |
| 3 | (E,E,E)-3,7,11,15-Tetramethylhex... | 272 | C20H32 | 077898-97-6 | 43 |
| 4 | Bicyclo[5.2.0]nonane, 4-methylen... | 204 | C15H24 | 1000159-38-2 | 38 |
| 5 | 9,10-Diazatricyclo[4.4.0.0(2,8)]... | 283 | C16H17N3O2 | 1000142-26-0 | 35 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049714.D
Acq On : 7 Aug 2021 21:35
Operator : CG/JU
Sample : M3244-05
Misc :
ALS Vial : 47 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** JCE05

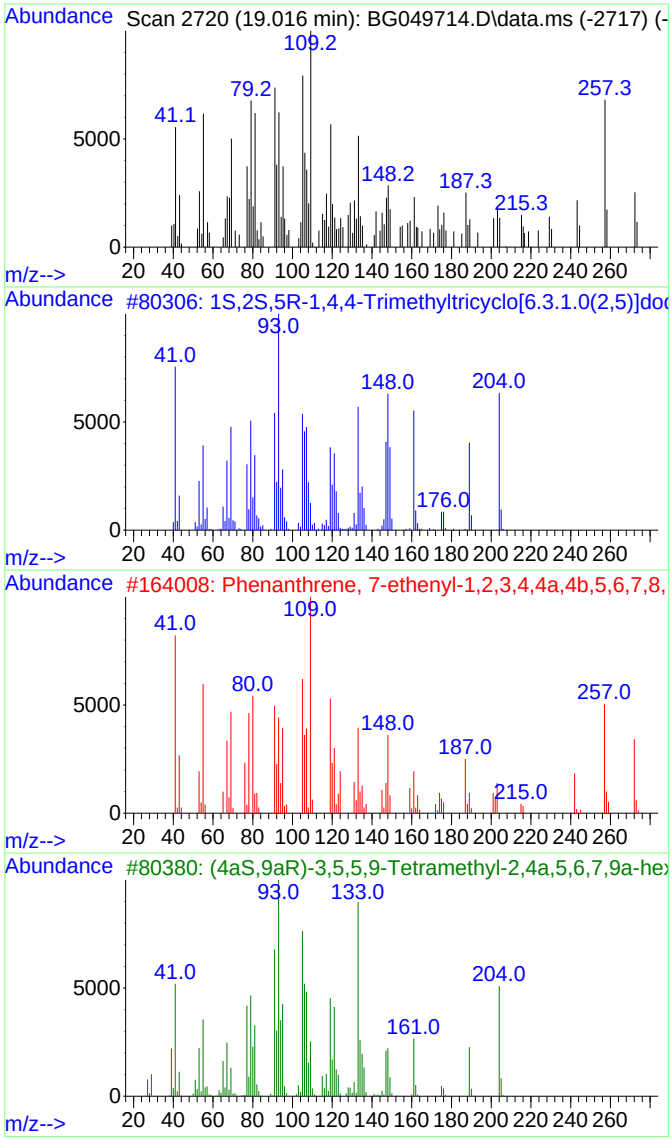
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

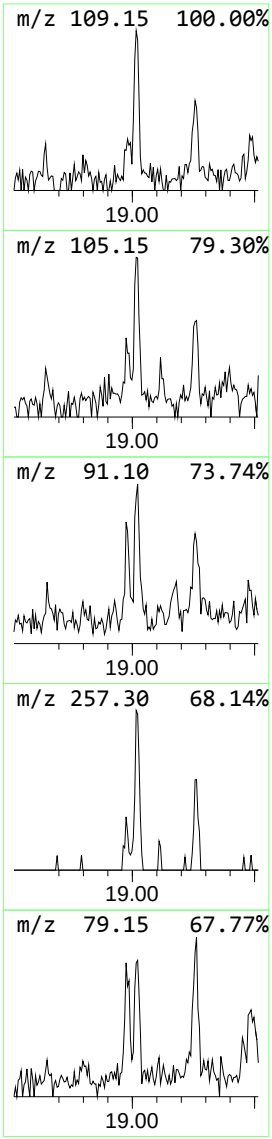
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 11 1S,2S,5R-1,4,4-Trimethyltri... Concentration Rank 25

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 19.015 | 3.12 ng/ul | 70721 | Phenanthrene-d10 | 17.441 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1S,2S,5R-1,4,4-Trimethyltricyclo... | 204 | C15H24 | 1000140-07-5 | 53 |
| 2 | Phenanthrene, 7-ethenyl-1,2,3,4,... | 272 | C20H32 | 001686-66-4 | 53 |
| 3 | (4aS,9aR)-3,5,5,9-Tetramethyl-2,... | 204 | C15H24 | 053111-25-4 | 52 |
| 4 | 1,3,6,10-Cyclotetradecatetraene,... | 272 | C20H32 | 001898-13-1 | 44 |
| 5 | Alloaromadendrene | 204 | C15H24 | 025246-27-9 | 41 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049714.D
Acq On : 7 Aug 2021 21:35
Operator : CG/JU
Sample : M3244-05
Misc :
ALS Vial : 47 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** JCE05

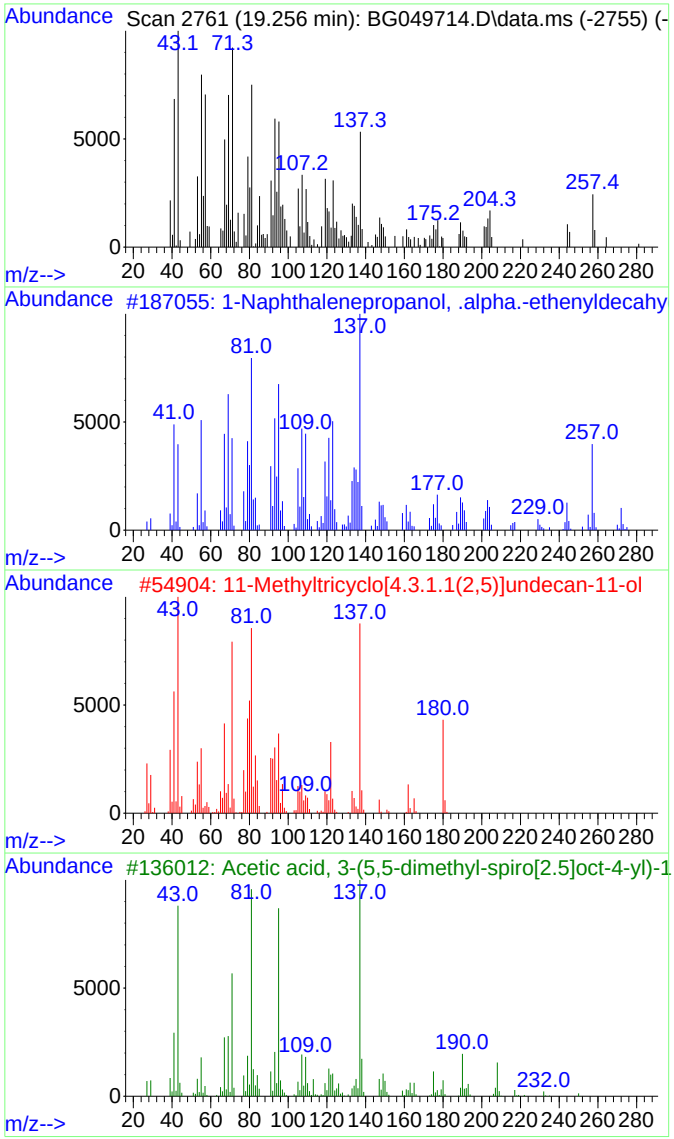
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

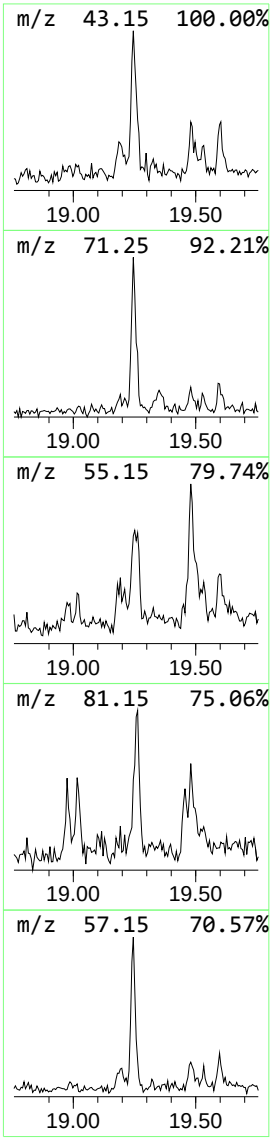
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 12 1-Naphthalenepropanol, .alp... Concentration Rank 16

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 19.256 | 8.49 ng/ul | 192358 | Phenanthrene-d10 | 17.441 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1-Naphthalenepropanol, .alpha.-e... | 290 | C20H34O | 001438-62-6 | 76 |
| 2 | 11-Methyltricyclo[4.3.1.1(2,5)]u... | 180 | C12H20O | 174226-52-9 | 58 |
| 3 | Acetic acid, 3-(5,5-dimethyl-spi... | 250 | C16H26O2 | 077143-33-0 | 43 |
| 4 | Acetic acid, 1-methyl-3-(1,3,3-t... | 250 | C16H26O2 | 1000192-22-3 | 43 |
| 5 | 2,6,10,10-Tetramethylbicyclo[7.2... | 204 | C15H24 | 136296-37-2 | 38 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049714.D
Acq On : 7 Aug 2021 21:35
Operator : CG/JU
Sample : M3244-05
Misc :
ALS Vial : 47 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** JCE05

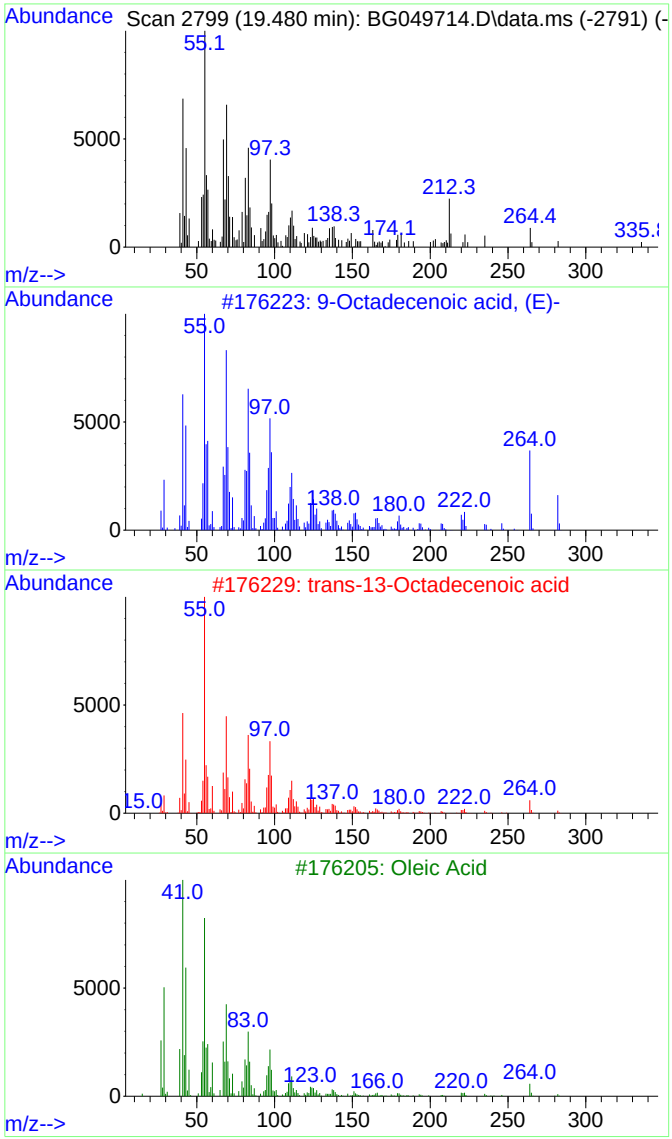
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

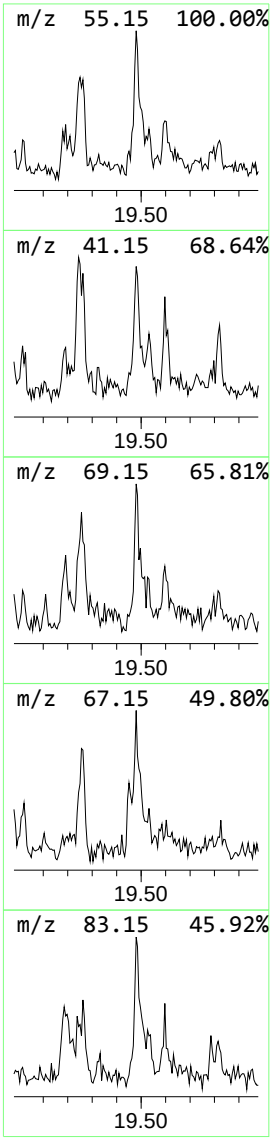
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 13 9-Octadecenoic acid, (E)- Concentration Rank 17

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 19.480 | 8.14 ng/ul | 184445 | Phenanthrene-d10 | 17.441 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 9-Octadecenoic acid, (E)- | 282 | C18H34O2 | 000112-79-8 | 86 |
| 2 | trans-13-Octadecenoic acid | 282 | C18H34O2 | 000693-71-0 | 78 |
| 3 | Oleic Acid | 282 | C18H34O2 | 000112-80-1 | 76 |
| 4 | Cetene | 224 | C16H32 | 000629-73-2 | 68 |
| 5 | 6-Octadecenoic acid | 282 | C18H34O2 | 1000336-66-8 | 64 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049714.D
Acq On : 7 Aug 2021 21:35
Operator : CG/JU
Sample : M3244-05
Misc :
ALS Vial : 47 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** JCE05

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

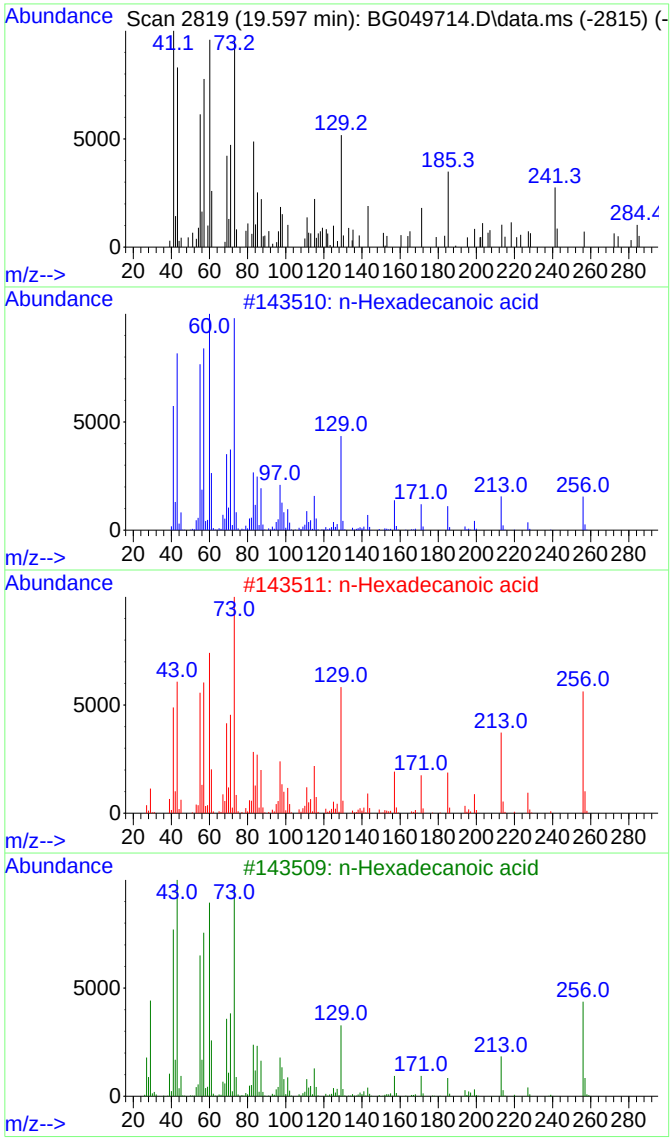
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

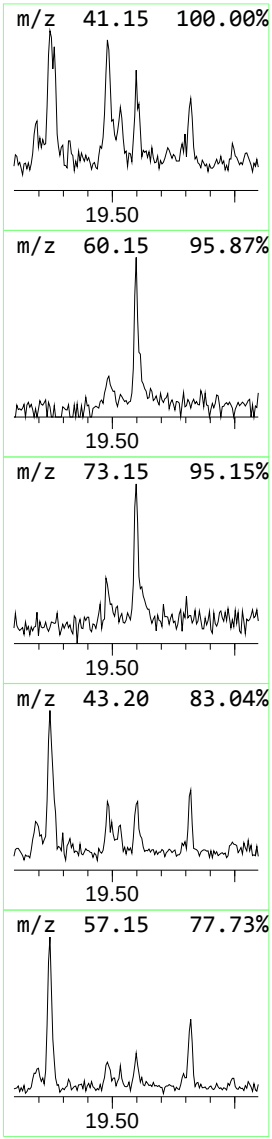
**********

Peak Number 14 Pentadecanoic acid Concentration Rank 28

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 19.597 | 2.55 ng/ul | 68367 | Chrysene-d12 | 21.724 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 96 |
| 2 | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 95 |
| 3 | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 91 |
| 4 | Pentadecanoic acid | 242 | C15H30O2 | 001002-84-2 | 89 |
| 5 | Octadecanoic acid | 284 | C18H36O2 | 000057-11-4 | 83 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049714.D
Acq On : 7 Aug 2021 21:35
Operator : CG/JU
Sample : M3244-05
Misc :
ALS Vial : 47 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** JCE05

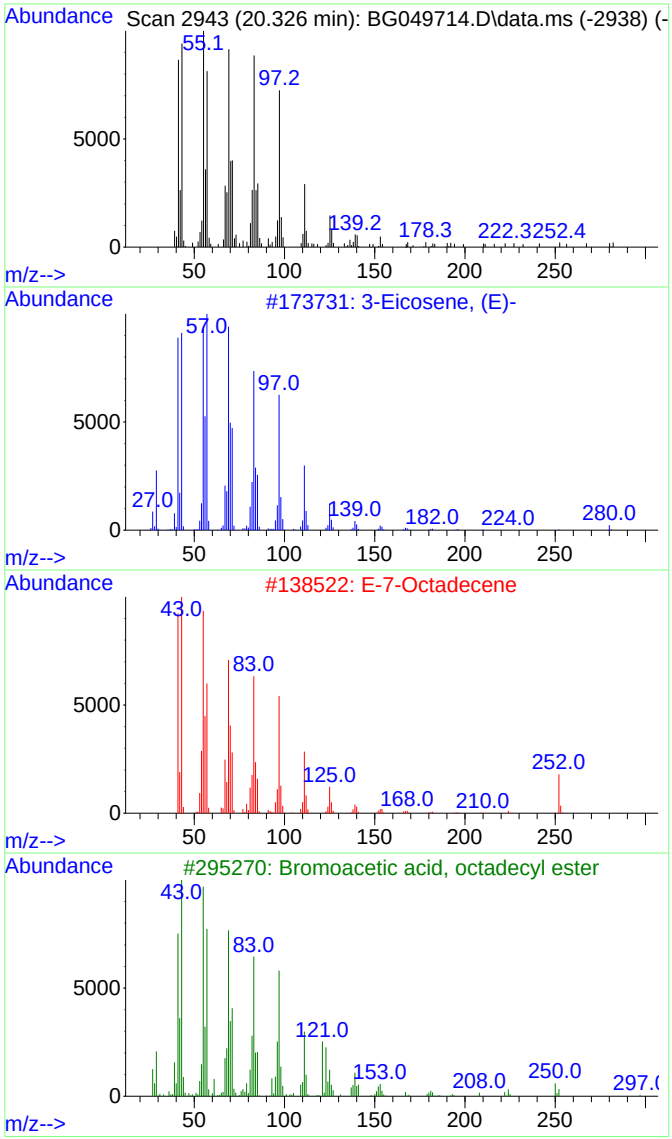
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

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

**********
Peak Number 15 (DEL) Alkane: Straight-Chai... Concentration Rank 18

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 20.326 | 6.00 ng/ul | 160664 | Chrysene-d12 | 21.724 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 3-Eicosene, (E)- | 280 | C20H40 | 074685-33-9 | 98 |
| 2 | E-7-Octadecene | 252 | C18H36 | 1000130-92-0 | 93 |
| 3 | Bromoacetic acid, octadecyl ester | 390 | C20H39BrO2 | 018992-03-5 | 91 |
| 4 | 5-Eicosene, (E)- | 280 | C20H40 | 074685-30-6 | 91 |
| 5 | 1-Octadecene | 252 | C18H36 | 000112-88-9 | 90 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049714.D
Acq On : 7 Aug 2021 21:35
Operator : CG/JU
Sample : M3244-05
Misc :
ALS Vial : 47 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** JCE05

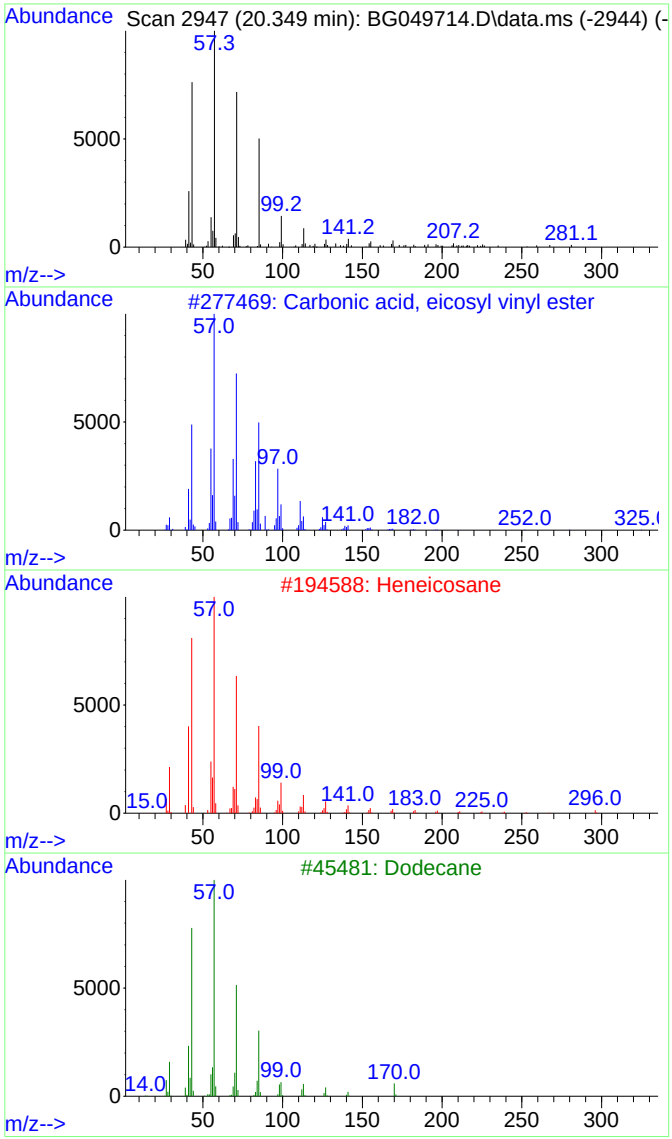
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

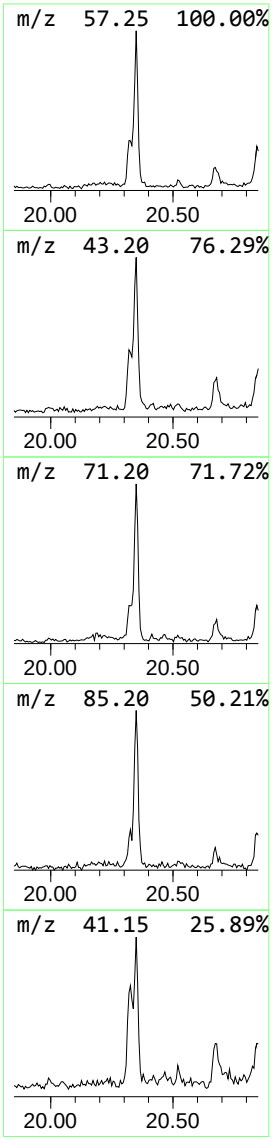
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 16 Carbonic acid, eicosyl viny... Concentration Rank 14

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 20.349 | 9.68 ng/ul | 258938 | Chrysene-d12 | 21.724 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Carbonic acid, eicosyl vinyl ester | 368 | C23H44O3 | 1000382-54-3 | 83 |
| 2 | Heneicosane | 296 | C21H44 | 000629-94-7 | 78 |
| 3 | Dodecane | 170 | C12H26 | 000112-40-3 | 78 |
| 4 | Tridecanol, 2-ethyl-2-methyl- | 242 | C16H34O | 1010115-66-1 | 78 |
| 5 | Undecane, 2,10-dimethyl- | 184 | C13H28 | 017301-27-8 | 72 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049714.D
Acq On : 7 Aug 2021 21:35
Operator : CG/JU
Sample : M3244-05
Misc :
ALS Vial : 47 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** JCE05

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

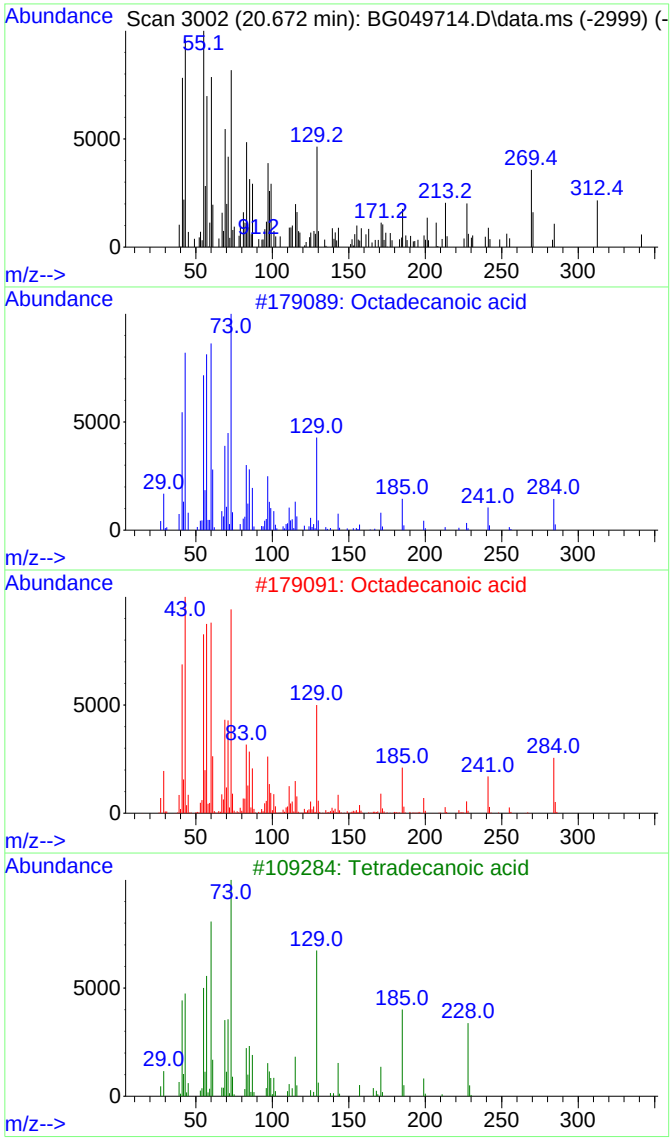
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

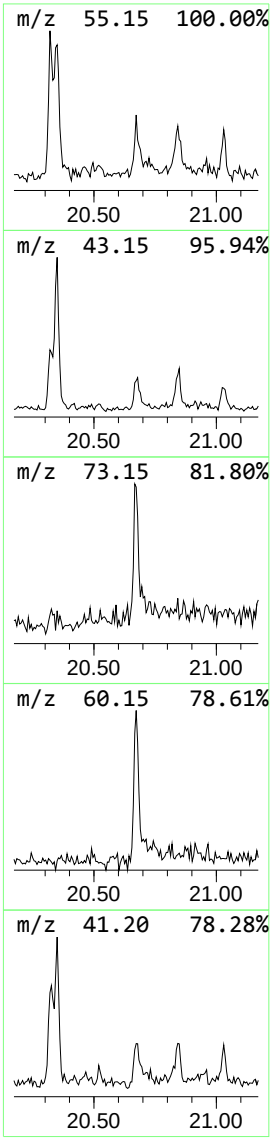
**********

Peak Number 17 Octadecanoic acid Concentration Rank 22

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 20.672 | 4.65 ng/ul | 124535 | Chrysene-d12 | 21.724 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Octadecanoic acid | 284 | C18H36O2 | 000057-11-4 | 90 |
| 2 | Octadecanoic acid | 284 | C18H36O2 | 000057-11-4 | 86 |
| 3 | Tetradecanoic acid | 228 | C14H28O2 | 000544-63-8 | 83 |
| 4 | Eicosanoic acid | 312 | C20H40O2 | 000506-30-9 | 55 |
| 5 | Tetradecanoic acid | 228 | C14H28O2 | 000544-63-8 | 52 |



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Data File : BG049714.D
Acq On : 7 Aug 2021 21:35
Operator : CG/JU
Sample : M3244-05
Misc :
ALS Vial : 47 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** JCE05

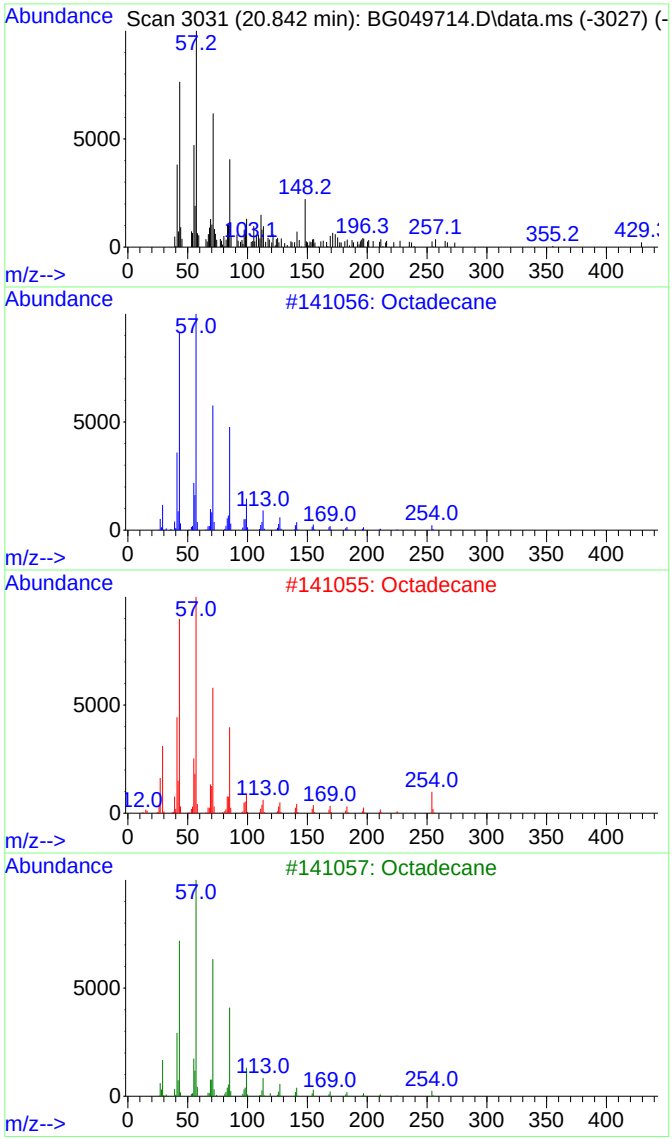
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

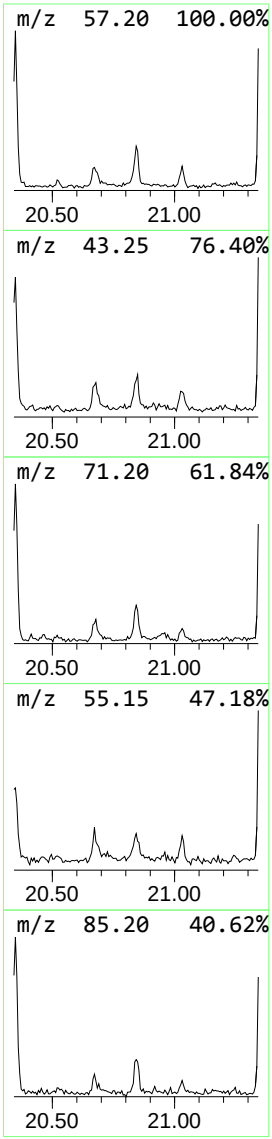
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 18 (DEL) Alkane: Straight-Chai... Concentration Rank 24

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 20.843 | 3.47 ng/ul | 92769 | Chrysene-d12 | 21.724 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Octadecane | 254 | C18H38 | 000593-45-3 | 94 |
| 2 | Octadecane | 254 | C18H38 | 000593-45-3 | 89 |
| 3 | Octadecane | 254 | C18H38 | 000593-45-3 | 76 |
| 4 | Octacosane | 394 | C28H58 | 000630-02-4 | 72 |
| 5 | Octacosane | 394 | C28H58 | 000630-02-4 | 64 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049714.D
Acq On : 7 Aug 2021 21:35
Operator : CG/JU
Sample : M3244-05
Misc :
ALS Vial : 47 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** JCE05

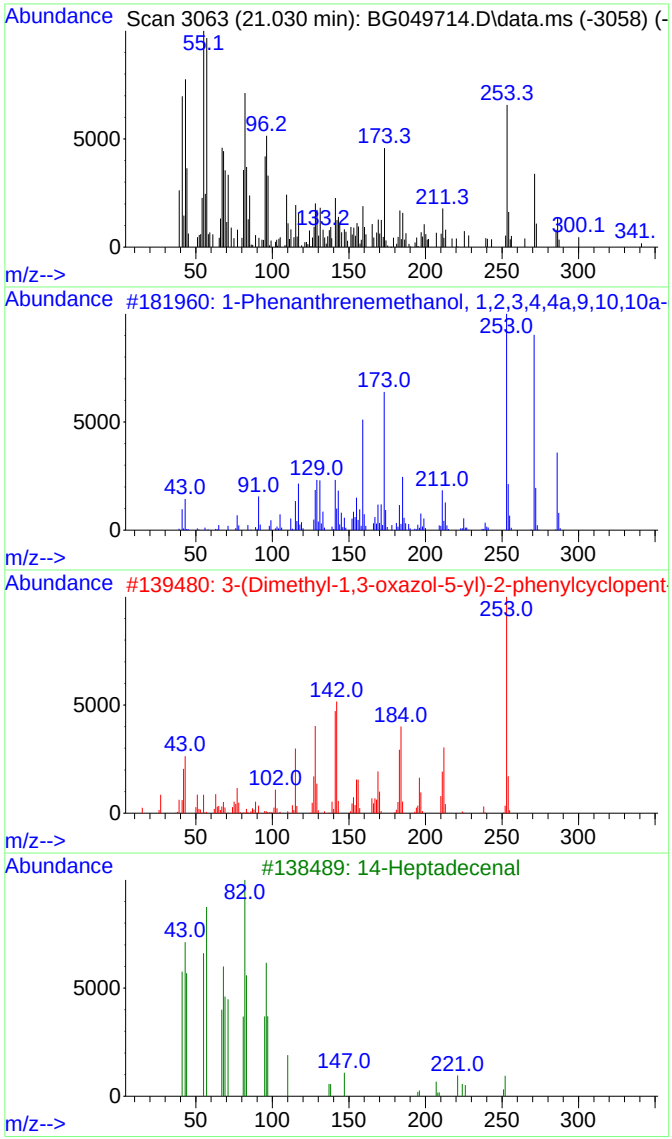
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

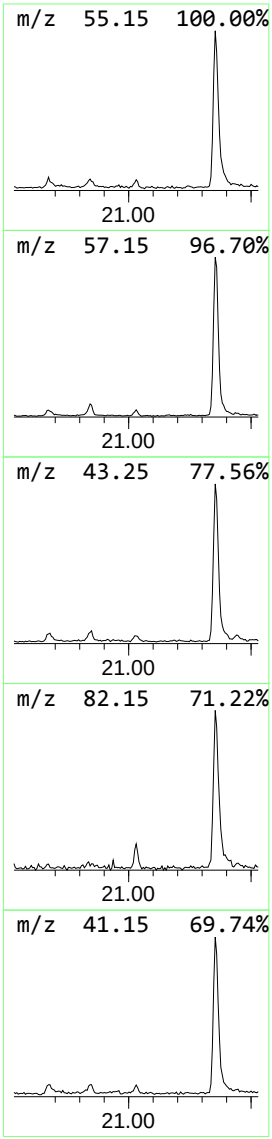
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 19 1-Phenanthrenemethanol, 1,2... Concentration Rank 20

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 21.031 | 4.93 ng/ul | 131827 | Chrysene-d12 | 21.724 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1-Phenanthrenemethanol, 1,2,3,4,... | 286 | C20H30O | 003772-55-2 | 70 |
| 2 | 3-(Dimethyl-1,3-oxazol-5-yl)-2-p... | 253 | C16H15NO2 | 1000445-13-1 | 60 |
| 3 | 14-Heptadecenal | 252 | C17H32O | 1010144-58-0 | 53 |
| 4 | E-2-Tetradecen-1-ol | 212 | C14H28O | 1000130-83-7 | 48 |
| 5 | 1-Phenanthrenemethanol, 1,2,3,4,... | 286 | C20H30O | 003772-55-2 | 42 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049714.D
Acq On : 7 Aug 2021 21:35
Operator : CG/JU
Sample : M3244-05
Misc :
ALS Vial : 47 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** JCE05

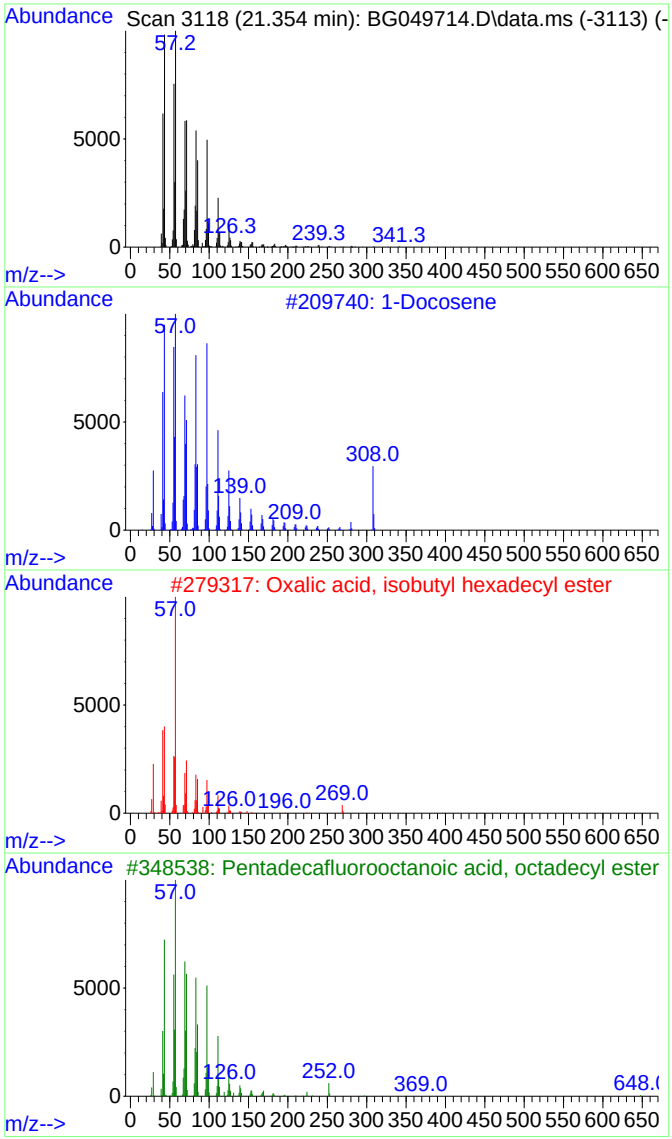
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

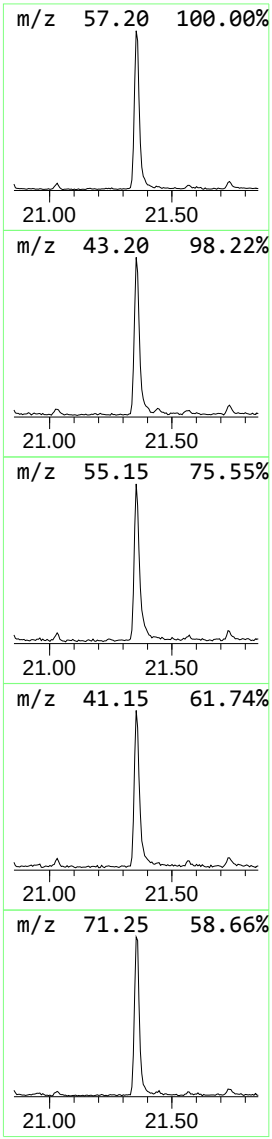
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 20 (DEL) Alkane: Straight-Chai... Concentration Rank 4

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 21.354 | 58.30 ng/ul | 1560090 | Chrysene-d12 | 21.724 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1-Docosene | 308 | C22H44 | 001599-67-3 | 96 |
| 2 | Oxalic acid, isobutyl hexadecyl ... | 370 | C22H42O4 | 1000309-38-1 | 91 |
| 3 | Pentadecafluorooctanoic acid, oc... | 666 | C26H37F15O2 | 1000406-04-8 | 90 |
| 4 | 1-Eicosene | 280 | C20H40 | 003452-07-1 | 89 |
| 5 | Oxalic acid, isobutyl octadecyl ... | 398 | C24H46O4 | 1000309-38-3 | 87 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049714.D
Acq On : 7 Aug 2021 21:35
Operator : CG/JU
Sample : M3244-05
Misc :
ALS Vial : 47 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** JCE05

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

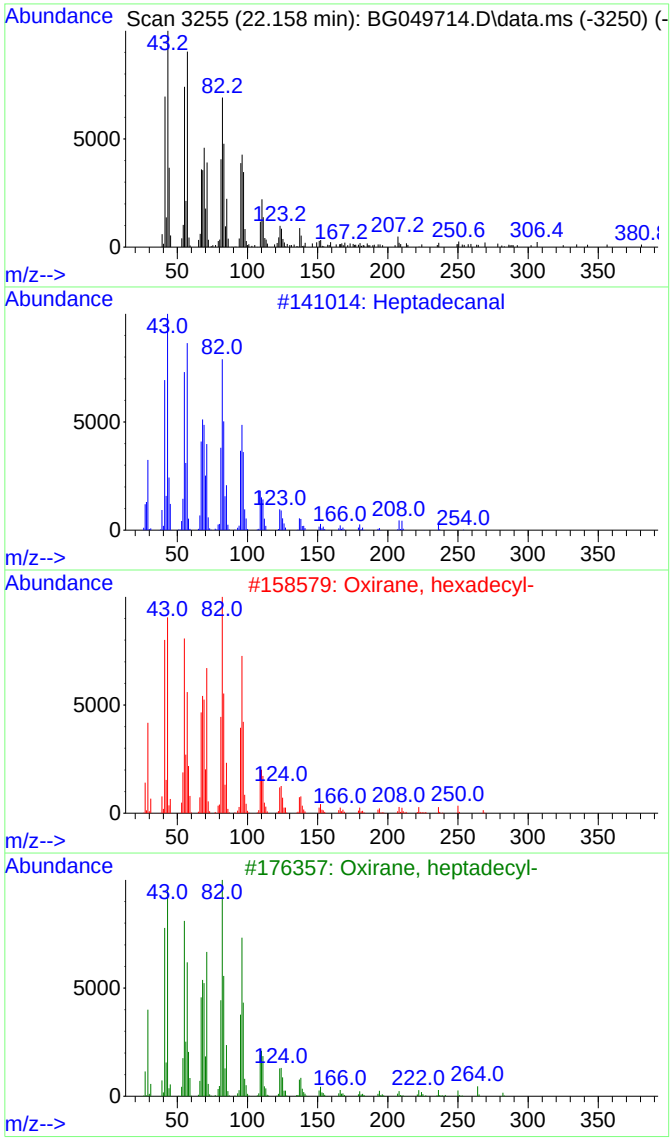
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

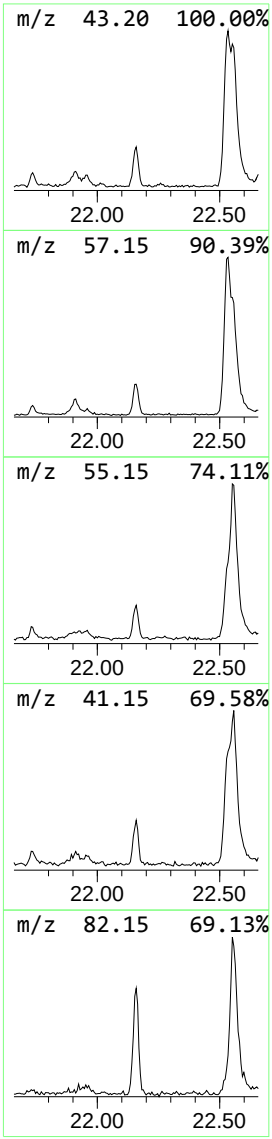
**********

Peak Number 21 Heptadecanal Concentration Rank 12

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 22.158 | 11.57 ng/ul | 309569 | Chrysene-d12 | 21.724 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Heptadecanal | 254 | C17H34O | 1000376-70-0 | 91 |
| 2 | Oxirane, hexadecyl- | 268 | C18H36O | 007390-81-0 | 87 |
| 3 | Oxirane, heptadecyl- | 282 | C19H38O | 067860-04-2 | 83 |
| 4 | 1,21-Docosadiene | 306 | C22H42 | 053057-53-7 | 83 |
| 5 | Eicosanal- | 296 | C20H40O | 002400-66-0 | 83 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049714.D
Acq On : 7 Aug 2021 21:35
Operator : CG/JU
Sample : M3244-05
Misc :
ALS Vial : 47 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** JCE05

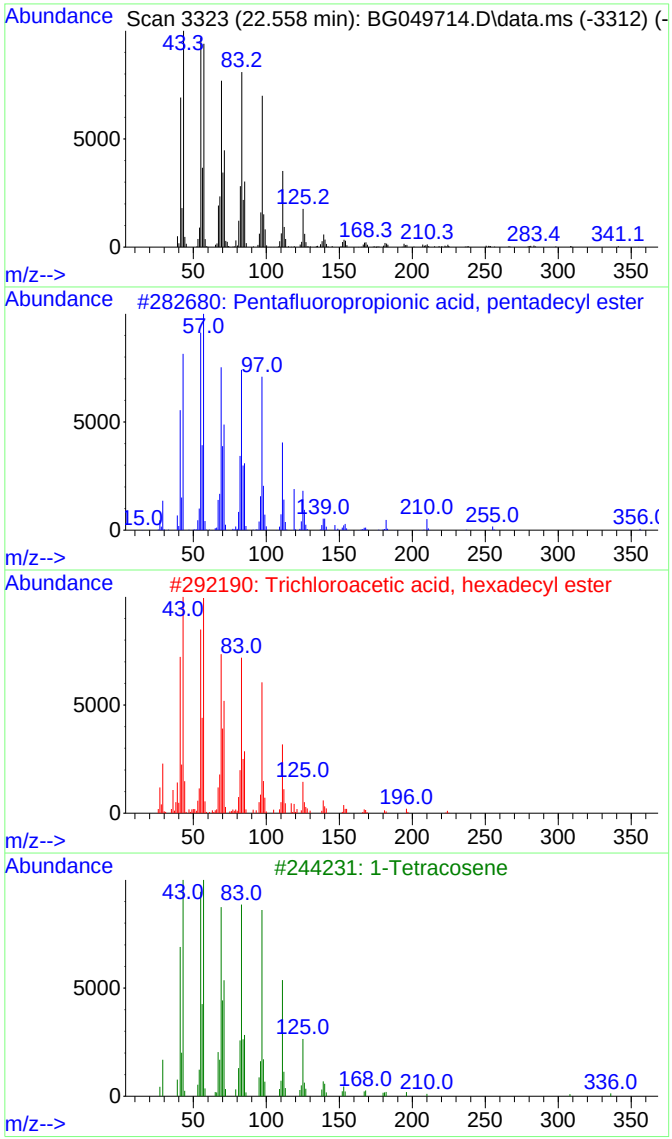
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

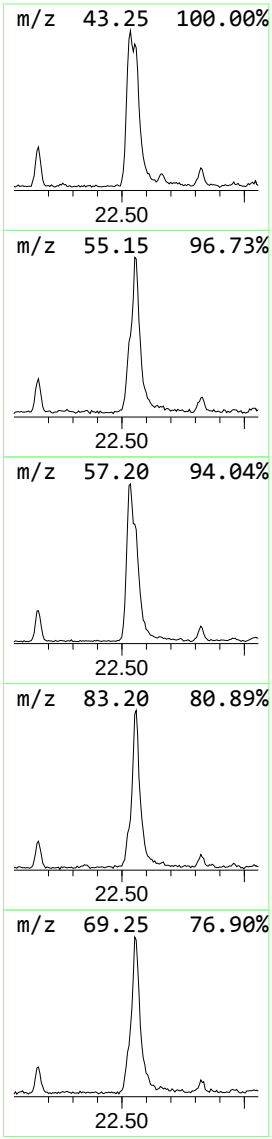
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 22 Pentafluoropropionic acid, ... Concentration Rank 2

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 22.558 | 81.80 ng/ul | 2188960 | Chrysene-d12 | 21.724 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Pentafluoropropionic acid, penta... | 374 | C18H31F5O2 | 959092-08-1 | 94 |
| 2 | Trichloroacetic acid, hexadecyl ... | 386 | C18H33Cl3O2 | 074339-54-1 | 94 |
| 3 | 1-Tetracosene | 336 | C24H48 | 010192-32-2 | 94 |
| 4 | Bromoacetic acid, hexadecyl ester | 362 | C18H35BrO2 | 005454-48-8 | 93 |
| 5 | E-15-Heptadecenal | 252 | C17H32O | 1000130-97-9 | 92 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049714.D
Acq On : 7 Aug 2021 21:35
Operator : CG/JU
Sample : M3244-05
Misc :
ALS Vial : 47 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** JCE05

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

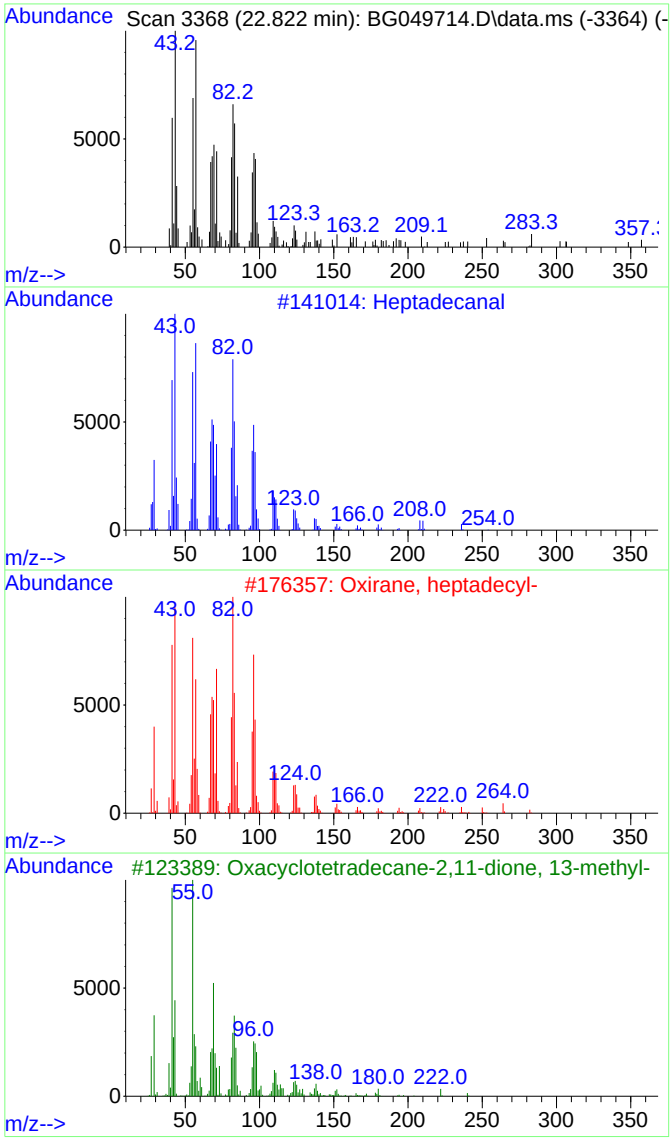
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

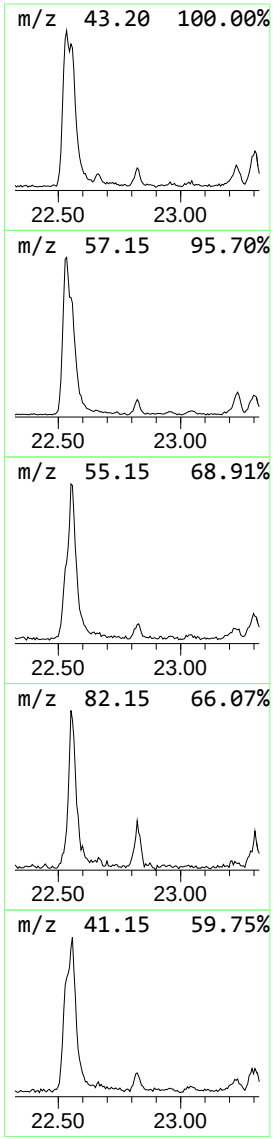
**********

Peak Number 23 Oxirane, heptadecyl- Concentration Rank 21

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 22.822 | 4.86 ng/ul | 129922 | Chrysene-d12 | 21.724 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Heptadecanal | 254 | C17H34O | 1000376-70-0 | 91 |
| 2 | Oxirane, heptadecyl- | 282 | C19H38O | 067860-04-2 | 74 |
| 3 | Oxacyclotetradecane-2,11-dione, ... | 240 | C14H24O3 | 074685-36-2 | 72 |
| 4 | 18-Nonadecen-1-ol | 282 | C19H38O | 1000142-89-2 | 58 |
| 5 | 17-Octadecenal | 266 | C18H34O | 056554-86-0 | 52 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049714.D
Acq On : 7 Aug 2021 21:35
Operator : CG/JU
Sample : M3244-05
Misc :
ALS Vial : 47 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** JCE05

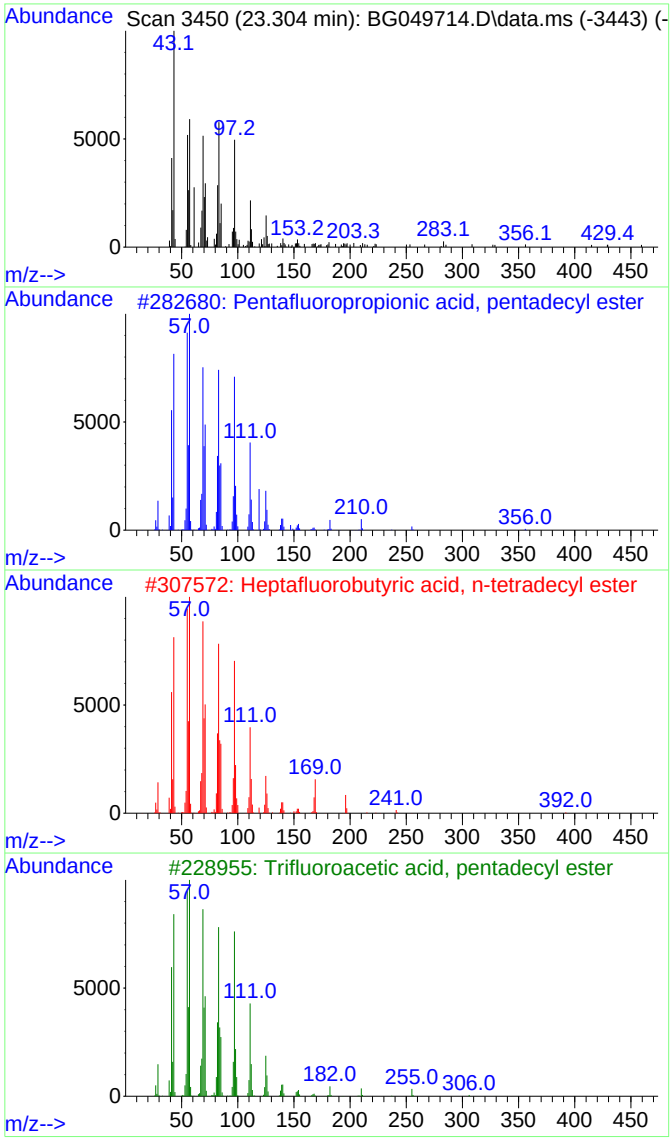
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

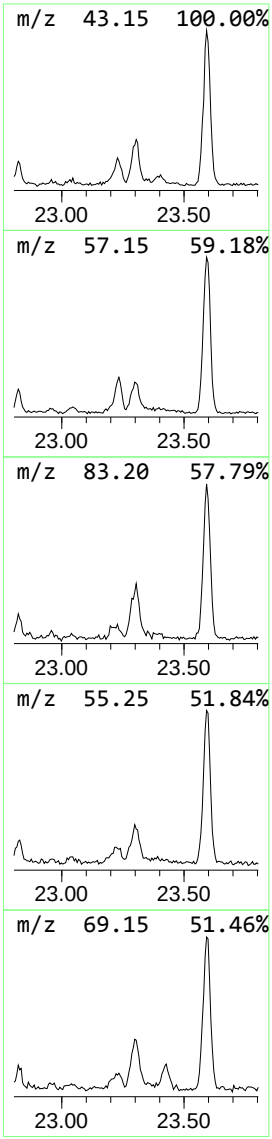
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 24 Heptafluorobutyric acid, n-... Concentration Rank 13

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 23.304 | 10.32 ng/ul | 276071 | Chrysene-d12 | 21.724 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Pentafluoropropionic acid, penta... | 374 | C18H31F5O2 | 959092-08-1 | 68 |
| 2 | Heptafluorobutyric acid, n-tetra... | 410 | C18H29F7O2 | 007365-36-8 | 68 |
| 3 | Trifluoroacetic acid, pentadecyl... | 324 | C17H31F3O2 | 959010-23-2 | 68 |
| 4 | Heptafluorobutyric acid, pentade... | 424 | C19H31F7O2 | 959261-23-5 | 68 |
| 5 | Heneicosyl heptafluorobutyrate | 508 | C25H43F7O2 | 1000351-83-8 | 62 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049714.D
Acq On : 7 Aug 2021 21:35
Operator : CG/JU
Sample : M3244-05
Misc :
ALS Vial : 47 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** JCE05

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

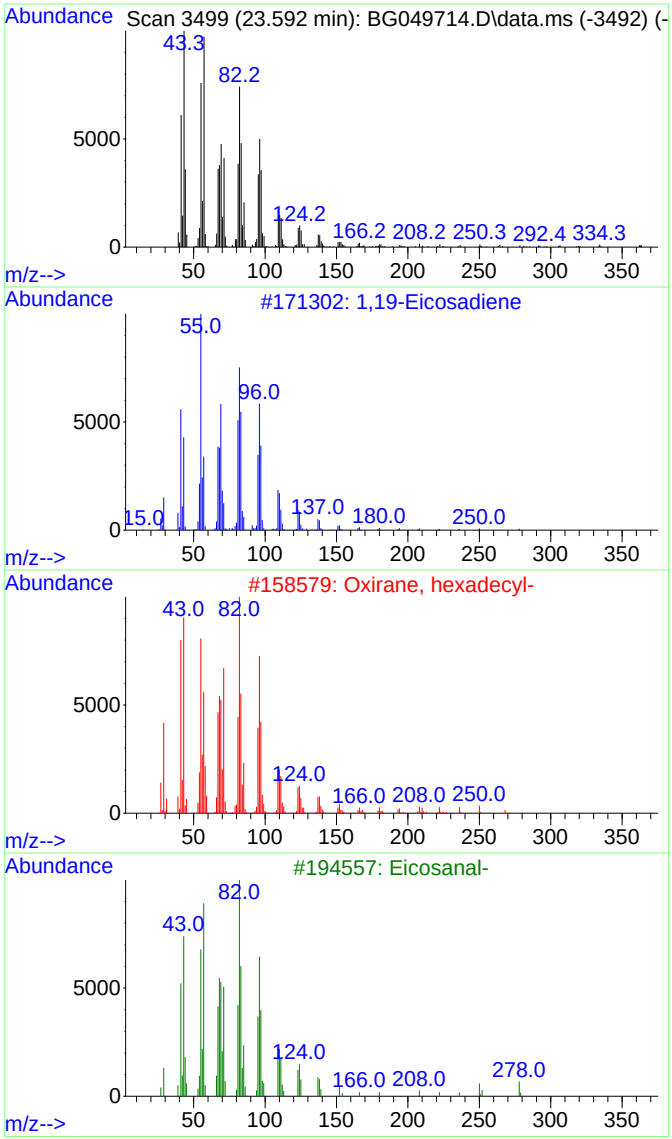
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

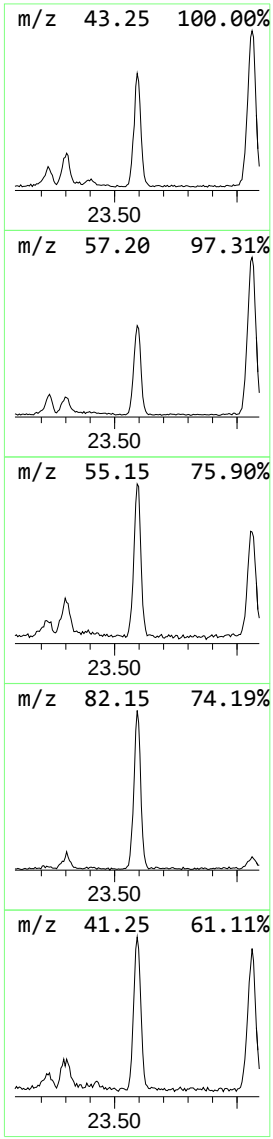
**********

Peak Number 25 1,19-Eicosadiene Concentration Rank 7

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 23.592 | 39.46 ng/ul | 1220260 | Perylene-d12 | 25.007 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1,19-Eicosadiene | 278 | C20H38 | 014811-95-1 | 95 |
| 2 | Oxirane, hexadecyl- | 268 | C18H36O | 007390-81-0 | 94 |
| 3 | Eicosanal- | 296 | C20H40O | 002400-66-0 | 94 |
| 4 | Hexacosanal | 380 | C26H52O | 026627-85-0 | 91 |
| 5 | Octadecanal | 268 | C18H36O | 000638-66-4 | 91 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049714.D
Acq On : 7 Aug 2021 21:35
Operator : CG/JU
Sample : M3244-05
Misc :
ALS Vial : 47 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** JCE05

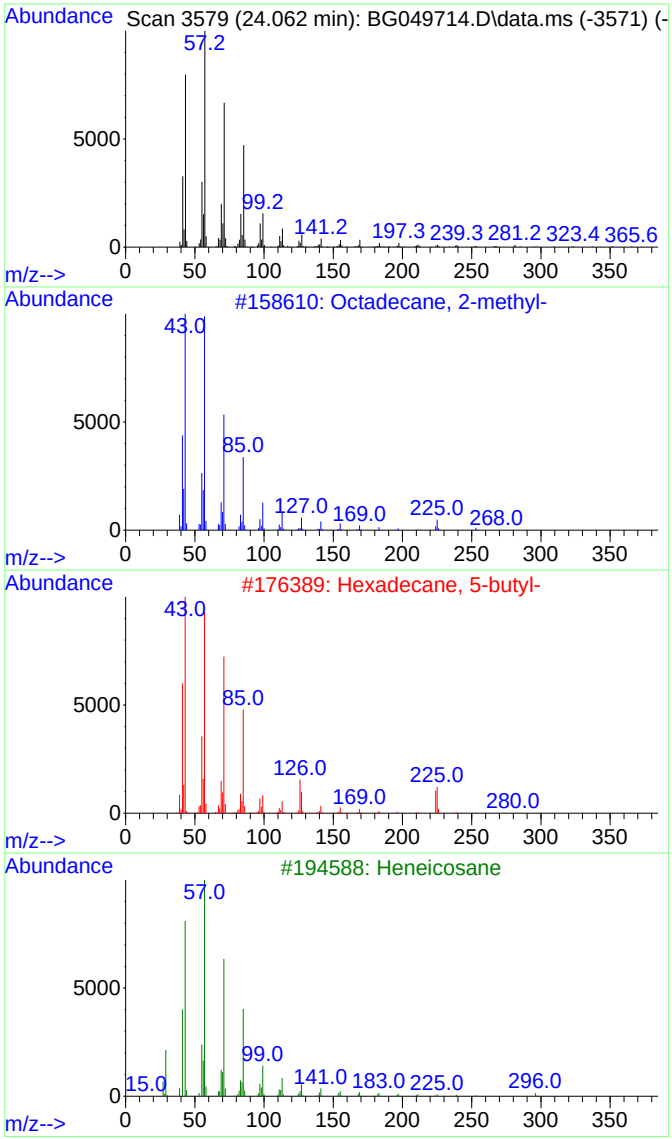
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

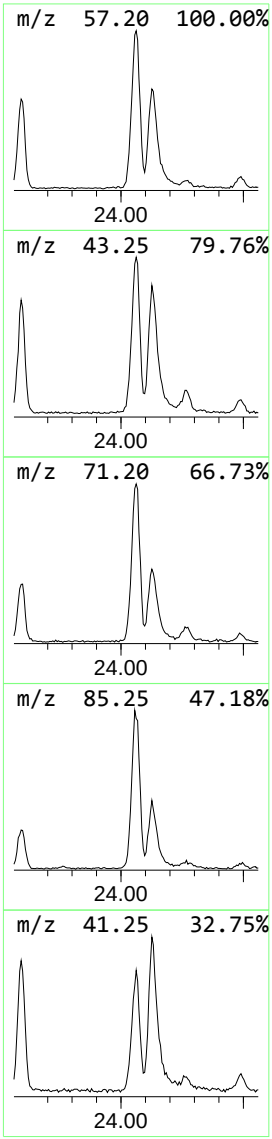
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 26 (DEL) Alkane: Straight-Chai... Concentration Rank 8

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 24.062 | 37.84 ng/ul | 1170070 | Perylene-d12 | 25.007 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Octadecane, 2-methyl- | 268 | C19H40 | 001560-88-9 | 94 |
| 2 | Hexadecane, 5-butyl- | 282 | C20H42 | 006912-07-8 | 93 |
| 3 | Heneicosane | 296 | C21H44 | 000629-94-7 | 91 |
| 4 | Octacosane | 394 | C28H58 | 000630-02-4 | 91 |
| 5 | Octacosane, 2-methyl- | 408 | C29H60 | 001560-98-1 | 91 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049714.D
Acq On : 7 Aug 2021 21:35
Operator : CG/JU
Sample : M3244-05
Misc :
ALS Vial : 47 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** JCE05

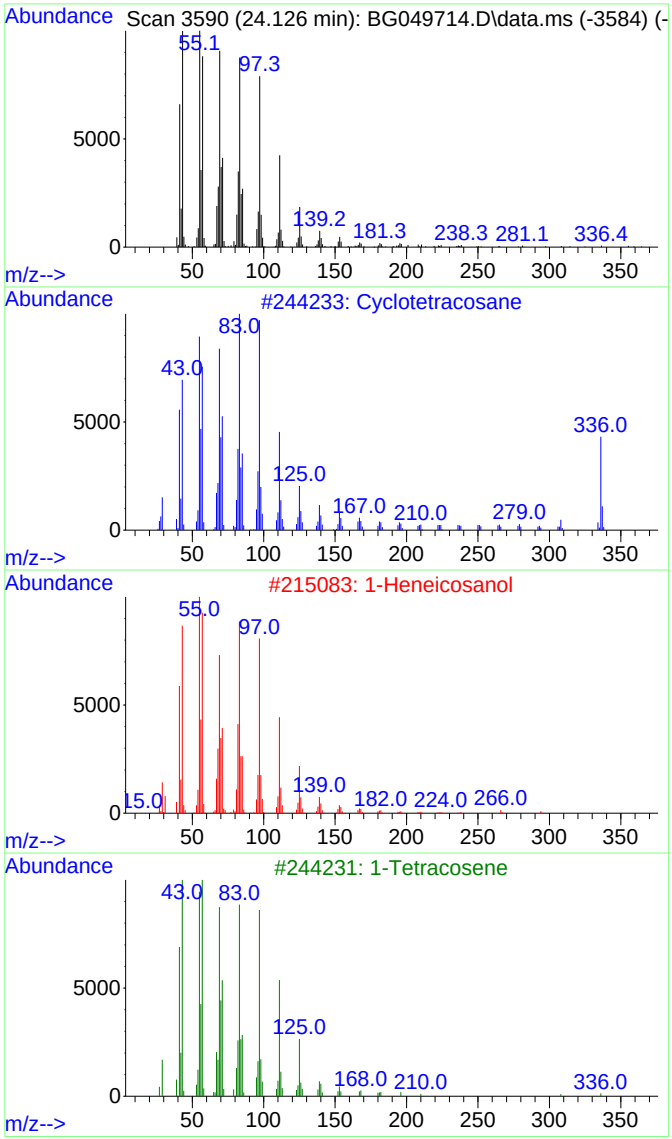
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

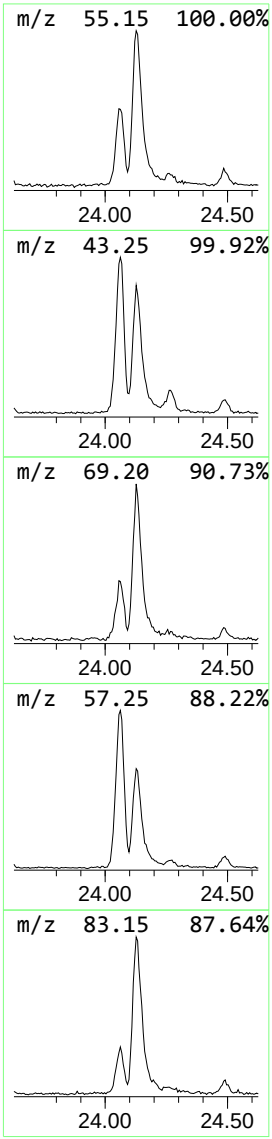
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 27 (DEL) Alkane: Cyclic24.126 Concentration Rank 5

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 24.126 | 54.13 ng/ul | 1673870 | Perylene-d12 | 25.007 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Cyclotetracosane | 336 | C24H48 | 000297-03-0 | 97 |
| 2 | 1-Heneicosanol | 312 | C21H44O | 015594-90-8 | 94 |
| 3 | 1-Tetracosene | 336 | C24H48 | 010192-32-2 | 94 |
| 4 | 1-Tricosene | 322 | C23H46 | 018835-32-0 | 93 |
| 5 | Cyclopentadecane | 210 | C15H30 | 000295-48-7 | 93 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049714.D
Acq On : 7 Aug 2021 21:35
Operator : CG/JU
Sample : M3244-05
Misc :
ALS Vial : 47 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** JCE05

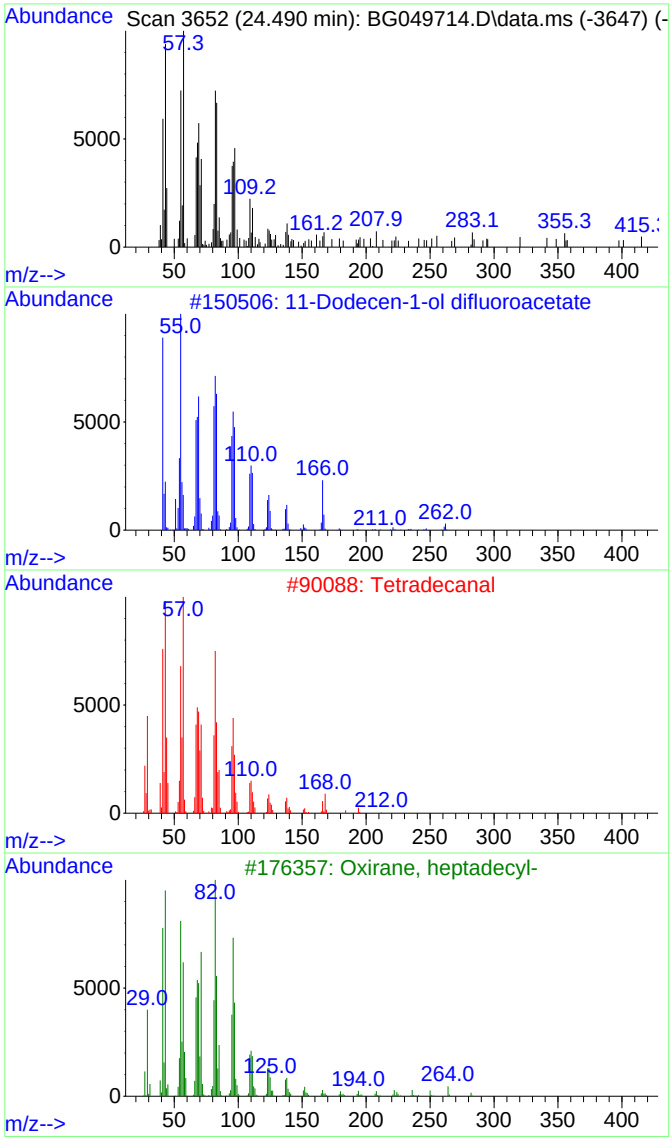
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

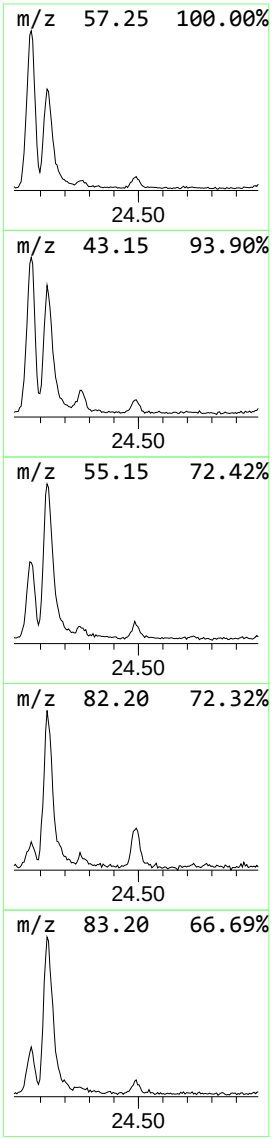
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 28 11-Dodecen-1-ol difluoroace... Concentration Rank 19

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 24.491 | 5.01 ng/ul | 154979 | Perylene-d12 | 25.007 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 11-Dodecen-1-ol difluoroacetate | 262 | C14H24F2O2 | 128792-45-0 | 80 |
| 2 | Tetradecanal | 212 | C14H28O | 000124-25-4 | 72 |
| 3 | Oxirane, heptadecyl- | 282 | C19H38O | 067860-04-2 | 70 |
| 4 | Z-11,13-Dimethyl-11-tetradecen-1... | 282 | C18H34O2 | 1000131-36-6 | 64 |
| 5 | Z-9-Hexadecen-1-ol acetate | 282 | C18H34O2 | 1000130-89-2 | 60 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049714.D
Acq On : 7 Aug 2021 21:35
Operator : CG/JU
Sample : M3244-05
Misc :
ALS Vial : 47 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** JCE05

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

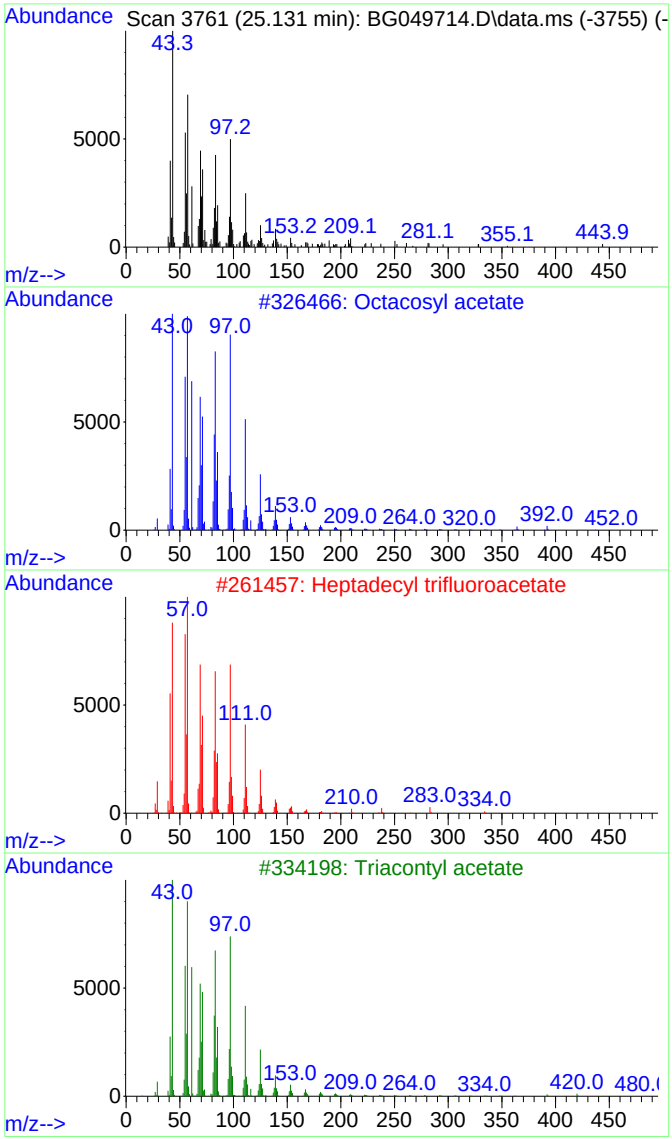
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

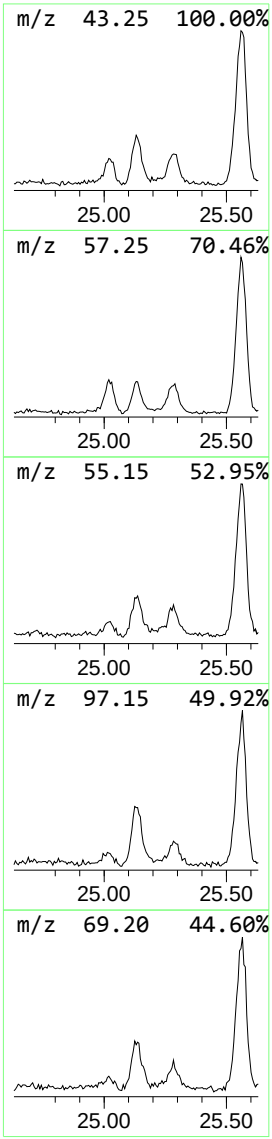
**********

Peak Number 29 Octacosyl acetate Concentration Rank 15

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 25.131 | 9.07 ng/ul | 280564 | Perylene-d12 | 25.007 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Octacosyl acetate | 452 | C30H60O2 | 018206-97-8 | 91 |
| 2 | Heptadecyl trifluoroacetate | 352 | C19H35F3O2 | 1010351-87-0 | 91 |
| 3 | Triacontyl acetate | 480 | C32H64O2 | 041755-58-2 | 91 |
| 4 | Heptacosyl acetate | 438 | C29H58O2 | 1000351-78-2 | 91 |
| 5 | Eicosyl acetate | 340 | C22H44O2 | 000822-24-2 | 90 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049714.D
Acq On : 7 Aug 2021 21:35
Operator : CG/JU
Sample : M3244-05
Misc :
ALS Vial : 47 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** JCE05

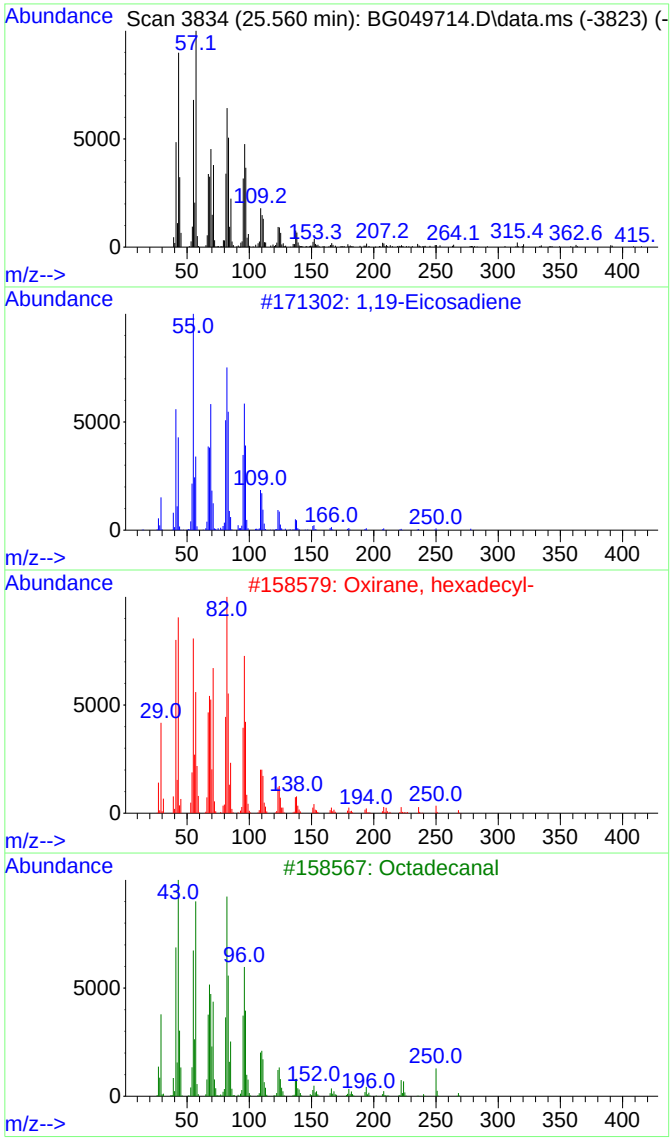
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

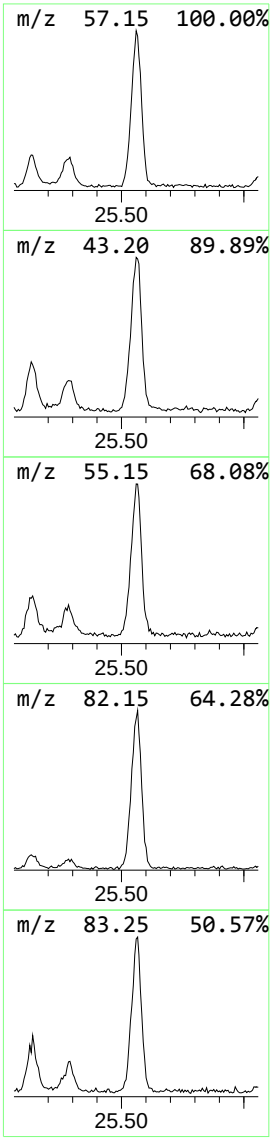
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 30 Oxirane, hexadecyl- Concentration Rank 6

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 25.560 | 47.92 ng/ul | 1482050 | Perylene-d12 | 25.007 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1,19-Eicosadiene | 278 | C20H38 | 014811-95-1 | 95 |
| 2 | Oxirane, hexadecyl- | 268 | C18H36O | 007390-81-0 | 91 |
| 3 | Octadecanal | 268 | C18H36O | 000638-66-4 | 91 |
| 4 | Eicosanal- | 296 | C20H40O | 002400-66-0 | 91 |
| 5 | Hexacosanal | 380 | C26H52O | 026627-85-0 | 91 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049714.D
Acq On : 7 Aug 2021 21:35
Operator : CG/JU
Sample : M3244-05
Misc :
ALS Vial : 47 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** JCE05

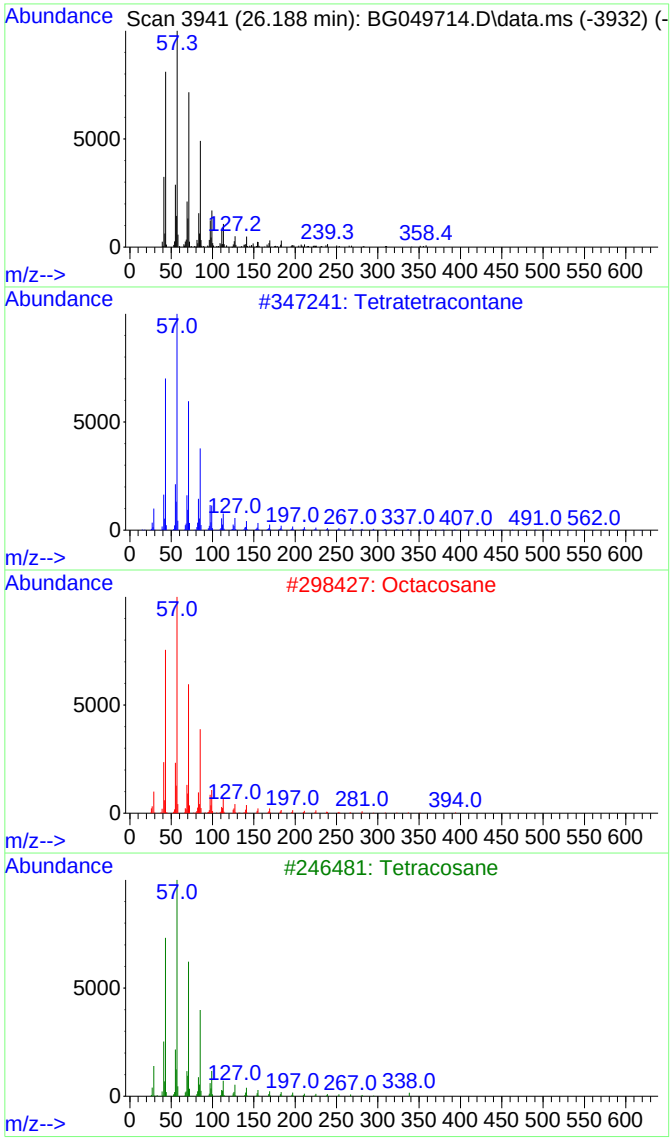
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

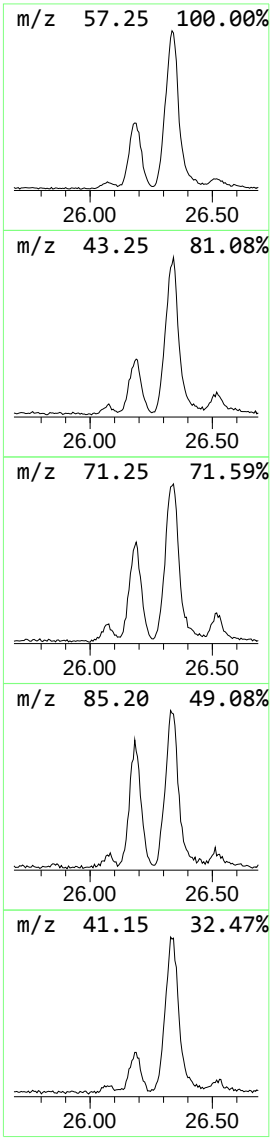
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 31 (DEL) Alkane: Straight-Chai... Concentration Rank 9

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 26.188 | 22.93 ng/ul | 709020 | Perylene-d12 | 25.007 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Tetratetracontane | 619 | C44H90 | 007098-22-8 | 91 |
| 2 | Octacosane | 394 | C28H58 | 000630-02-4 | 91 |
| 3 | Tetracosane | 338 | C24H50 | 000646-31-1 | 91 |
| 4 | Tridecane, 1-iodo- | 310 | C13H27I | 035599-77-0 | 90 |
| 5 | Nonadecane | 268 | C19H40 | 000629-92-5 | 90 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049714.D
Acq On : 7 Aug 2021 21:35
Operator : CG/JU
Sample : M3244-05
Misc :
ALS Vial : 47 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** JCE05

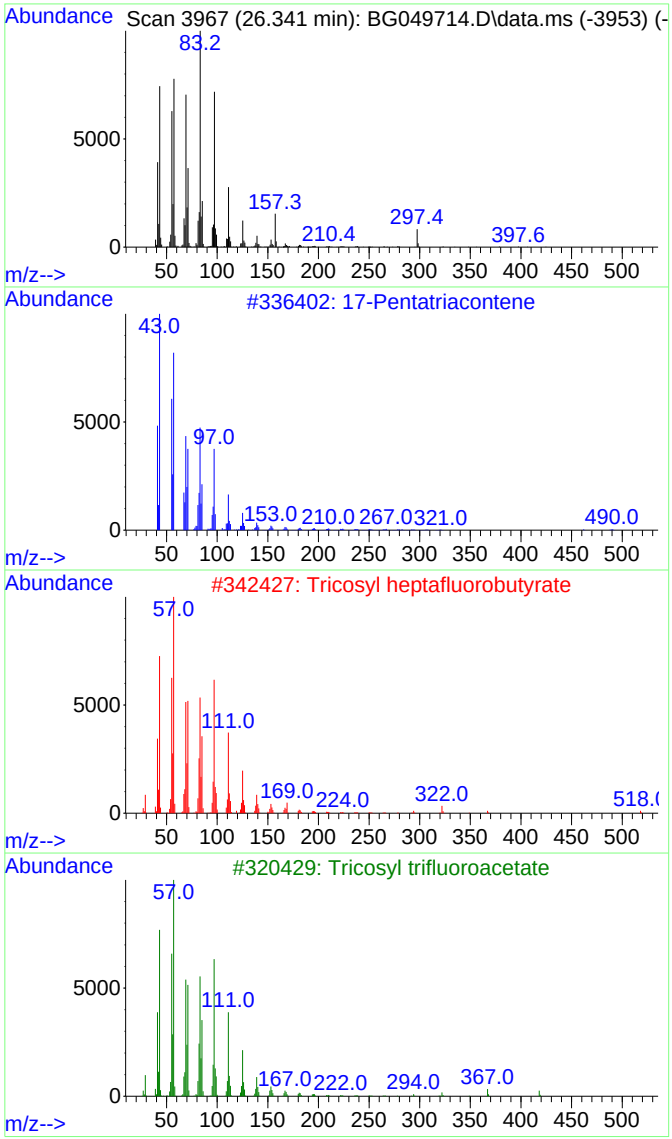
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.M
Quant Title : SVOA CALIBRATION

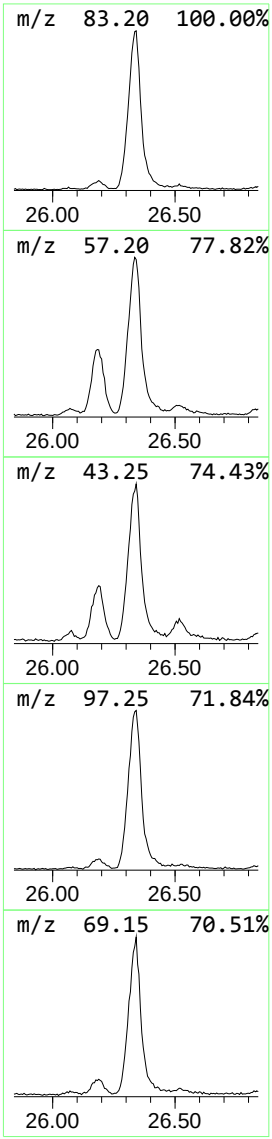
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 32 (DEL) Alkane: Straight-Chai... Concentration Rank 1

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 26.341 | 124.45 ng/ul | 3848660 | Perylene-d12 | 25.007 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 17-Pentatriacontene | 491 | C35H70 | 006971-40-0 | 50 |
| 2 | Tricosyl heptafluorobutyrate | 536 | C27H47F7O2 | 1000351-83-4 | 45 |
| 3 | Tricosyl trifluoroacetate | 436 | C25H47F3O2 | 1000351-75-1 | 43 |
| 4 | Octacosyl acetate | 452 | C30H60O2 | 018206-97-8 | 38 |
| 5 | Pentadecafluorooctanoic acid, oc... | 666 | C26H37F15O2 | 1000406-04-8 | 38 |



Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG080521\
Data File : BG049714.D
Acq On : 7 Aug 2021 21:35
Operator : CG/JU
Sample : M3244-05
Misc :
ALS Vial : 47 Sample Multiplier: 1

**Instrument :** BNA_G
**ClientSampleId :** JCE05

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG080421.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 |
| (DEL) Alkane: C... | 3.078 | 17.2 | ng/ul | 184169 | 1 | 8.042 | 214022 | 20.0 |
| Butane, 2-metho... | 3.161 | 62.1 | ng/ul | 664338 | 1 | 8.042 | 214022 | 20.0 |
| unknown-01 | 4.412 | 2.5 | ng/ul | 26923 | 1 | 8.042 | 214022 | 20.0 |
| .gamma.-Terpinene | 6.715 | 3.8 | ng/ul | 40312 | 1 | 8.042 | 214022 | 20.0 |
| Bicyclo[3.1.0]h... | 8.318 | 2.7 | ng/ul | 29037 | 1 | 8.042 | 214022 | 20.0 |
| unknown-02 | 9.364 | 2.2 | ng/ul | 23128 | 1 | 8.042 | 214022 | 20.0 |
| unknown-03 | 15.039 | 2.5 | ng/ul | 50829 | 3 | 14.692 | 406395 | 20.0 |
| Carbonic acid, ... | 17.195 | 3.0 | ng/ul | 67665 | 4 | 17.441 | 452913 | 20.0 |
| n-Hexadecanoic ... | 18.328 | 12.4 | ng/ul | 280265 | 4 | 17.441 | 452913 | 20.0 |
| 1,3,6,10-Cyclot... | 18.974 | 2.2 | ng/ul | 50088 | 4 | 17.441 | 452913 | 20.0 |
| 1S,2S,5R-1,4,4-... | 19.015 | 3.1 | ng/ul | 70721 | 4 | 17.441 | 452913 | 20.0 |
| 1-Naphthalenepr... | 19.256 | 8.5 | ng/ul | 192358 | 4 | 17.441 | 452913 | 20.0 |
| 9-Octadecenoic ... | 19.480 | 8.1 | ng/ul | 184445 | 4 | 17.441 | 452913 | 20.0 |
| Pentadecanoic acid | 19.597 | 2.5 | ng/ul | 68367 | 5 | 21.724 | 535171 | 20.0 |
| (DEL) Alkane: S... | 20.326 | 6.0 | ng/ul | 160664 | 5 | 21.724 | 535171 | 20.0 |
| Carbonic acid, ... | 20.349 | 9.7 | ng/ul | 258938 | 5 | 21.724 | 535171 | 20.0 |
| Octadecanoic acid | 20.672 | 4.7 | ng/ul | 124535 | 5 | 21.724 | 535171 | 20.0 |
| (DEL) Alkane: S... | 20.843 | 3.5 | ng/ul | 92769 | 5 | 21.724 | 535171 | 20.0 |
| 1-Phenanthrenem... | 21.031 | 4.9 | ng/ul | 131827 | 5 | 21.724 | 535171 | 20.0 |
| (DEL) Alkane: S... | 21.354 | 58.3 | ng/ul | 1560090 | 5 | 21.724 | 535171 | 20.0 |
| Heptadecanal | 22.158 | 11.6 | ng/ul | 309569 | 5 | 21.724 | 535171 | 20.0 |
| Pentafluoroprop... | 22.558 | 81.8 | ng/ul | 2188960 | 5 | 21.724 | 535171 | 20.0 |
| Oxirane, heptad... | 22.822 | 4.9 | ng/ul | 129922 | 5 | 21.724 | 535171 | 20.0 |
| Heptafluorobuty... | 23.304 | 10.3 | ng/ul | 276071 | 5 | 21.724 | 535171 | 20.0 |
| 1,19-Eicosadiene | 23.592 | 39.5 | ng/ul | 1220260 | 6 | 25.007 | 618491 | 20.0 |
| (DEL) Alkane: S... | 24.062 | 37.8 | ng/ul | 1170070 | 6 | 25.007 | 618491 | 20.0 |
| (DEL) Alkane: C... | 24.126 | 54.1 | ng/ul | 1673870 | 6 | 25.007 | 618491 | 20.0 |
| 11-Dodecen-1-ol... | 24.491 | 5.0 | ng/ul | 154979 | 6 | 25.007 | 618491 | 20.0 |
| Octacosyl acetate | 25.131 | 9.1 | ng/ul | 280564 | 6 | 25.007 | 618491 | 20.0 |
| Oxirane, hexade... | 25.560 | 47.9 | ng/ul | 1482050 | 6 | 25.007 | 618491 | 20.0 |
| (DEL) Alkane: S... | 26.188 | 22.9 | ng/ul | 709020 | 6 | 25.007 | 618491 | 20.0 |
| (DEL) Alkane: S... | 26.341 | 124.5 | ng/ul | 3848660 | 6 | 25.007 | 618491 | 20.0 |