

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP061423\  
 Data File : BP015546.D  
 Acq On : 15 Jun 2023 07:45  
 Operator : MA/JU  
 Sample : 03088-05  
 Misc :  
 ALS Vial : 41 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 DCFR1

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\_P\Methods\SFAM-EPA-BP060723.MA.M  
 Title : SVOA CALIBRATION

Signal : TIC: BP015546.D\data.ms

| peak<br># | R.T.<br>min | first<br>scan | max<br>scan | last<br>scan | PK<br>TY | peak<br>height | corr.<br>area | corr.<br>% max. | % of<br>total |
|-----------|-------------|---------------|-------------|--------------|----------|----------------|---------------|-----------------|---------------|
| 1         | 3.958       | 128           | 131         | 139          | rVB      | 25567          | 35616         | 0.13%           | 0.046%        |
| 2         | 4.081       | 148           | 152         | 160          | rVB      | 19403          | 35715         | 0.14%           | 0.046%        |
| 3         | 6.740       | 598           | 604         | 611          | rVB      | 31758          | 51239         | 0.19%           | 0.066%        |
| 4         | 7.240       | 681           | 689         | 706          | rBV      | 263979         | 538373        | 2.04%           | 0.693%        |
| 5         | 7.411       | 712           | 718         | 730          | rVB      | 210227         | 343882        | 1.30%           | 0.442%        |
| 6         | 7.511       | 730           | 735         | 745          | rVB3     | 36019          | 68185         | 0.26%           | 0.088%        |
| 7         | 7.611       | 745           | 752         | 762          | rBV      | 271291         | 464865        | 1.76%           | 0.598%        |
| 8         | 8.087       | 823           | 833         | 844          | rBV      | 314048         | 515974        | 1.95%           | 0.664%        |
| 9         | 8.805       | 948           | 955         | 966          | rBV      | 263406         | 488546        | 1.85%           | 0.628%        |
| 10        | 8.922       | 969           | 975         | 986          | rVV      | 183227         | 317421        | 1.20%           | 0.408%        |
| 11        | 9.263       | 1025          | 1033        | 1044         | rBV      | 271297         | 483724        | 1.83%           | 0.622%        |
| 12        | 9.993       | 1148          | 1157        | 1172         | rBV      | 239253         | 432779        | 1.64%           | 0.557%        |
| 13        | 10.175      | 1180          | 1188        | 1191         | rBV4     | 21830          | 43253         | 0.16%           | 0.056%        |
| 14        | 10.357      | 1214          | 1219        | 1226         | rVB3     | 19439          | 35641         | 0.13%           | 0.046%        |
| 15        | 10.540      | 1243          | 1250        | 1264         | rBV      | 290692         | 602803        | 2.28%           | 0.775%        |
| 16        | 10.910      | 1306          | 1313        | 1321         | rBV      | 477350         | 795022        | 3.01%           | 1.023%        |
| 17        | 11.063      | 1333          | 1339        | 1351         | rBV      | 70302          | 165105        | 0.62%           | 0.212%        |
| 18        | 13.457      | 1741          | 1746        | 1749         | rBV6     | 18044          | 29308         | 0.11%           | 0.038%        |
| 19        | 13.610      | 1763          | 1772        | 1777         | rBV4     | 36460          | 68945         | 0.26%           | 0.089%        |
| 20        | 13.710      | 1786          | 1789        | 1796         | rVB      | 26669          | 38638         | 0.15%           | 0.050%        |
| 21        | 14.046      | 1841          | 1846        | 1851         | rBV4     | 35535          | 49137         | 0.19%           | 0.063%        |
| 22        | 14.134      | 1853          | 1861        | 1866         | rBV      | 706903         | 1000763       | 3.79%           | 1.287%        |
| 23        | 14.181      | 1866          | 1869        | 1874         | rVV3     | 47640          | 78223         | 0.30%           | 0.101%        |
| 24        | 14.246      | 1877          | 1880        | 1884         | rVB2     | 29424          | 37902         | 0.14%           | 0.049%        |
| 25        | 14.422      | 1904          | 1910        | 1918         | rBV      | 754644         | 1117387       | 4.23%           | 1.437%        |
| 26        | 14.604      | 1937          | 1941        | 1948         | rVV6     | 14803          | 32257         | 0.12%           | 0.041%        |
| 27        | 14.675      | 1950          | 1953        | 1956         | rVV3     | 21512          | 30399         | 0.12%           | 0.039%        |
| 28        | 14.728      | 1956          | 1962        | 1968         | rVV      | 763973         | 1143835       | 4.33%           | 1.471%        |
| 29        | 14.787      | 1968          | 1972        | 1978         | rVV4     | 55765          | 100665        | 0.38%           | 0.129%        |
| 30        | 14.957      | 1994          | 2001        | 2012         | rBV3     | 84421          | 266526        | 1.01%           | 0.343%        |
| 31        | 15.134      | 2024          | 2031        | 2035         | rVB8     | 21574          | 41511         | 0.16%           | 0.053%        |
| 32        | 15.257      | 2048          | 2052        | 2060         | rVB4     | 26679          | 45308         | 0.17%           | 0.058%        |
| 33        | 15.487      | 2087          | 2091        | 2097         | rBV6     | 18356          | 45134         | 0.17%           | 0.058%        |
| 34        | 15.722      | 2124          | 2131        | 2137         | rBV      | 982323         | 1488510       | 5.63%           | 1.915%        |
| 35        | 15.851      | 2148          | 2153        | 2165         | rBV      | 238291         | 394995        | 1.49%           | 0.508%        |
| 36        | 16.087      | 2188          | 2193        | 2199         | rBV      | 122451         | 182923        | 0.69%           | 0.235%        |

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 Sample : 03088-05  
 Misc :  
 ALS Vial : 41 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 DCFR1

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\_P\Methods\SFAM-EPA-BP060723.MA.M  
 Title : SVOA CALIBRATION

|    |        |      |      |      |       |         |         |       |        |
|----|--------|------|------|------|-------|---------|---------|-------|--------|
| 37 | 16.445 | 2247 | 2254 | 2260 | rBV8  | 36997   | 90887   | 0.34% | 0.117% |
| 38 | 17.357 | 2406 | 2409 | 2414 | rBV7  | 22677   | 33824   | 0.13% | 0.044% |
| 39 | 17.486 | 2425 | 2431 | 2436 | rBV   | 969227  | 1420804 | 5.38% | 1.828% |
| 40 | 17.528 | 2436 | 2438 | 2441 | rVV   | 74759   | 87279   | 0.33% | 0.112% |
| 41 | 17.587 | 2442 | 2448 | 2455 | rVB   | 920068  | 1341983 | 5.08% | 1.726% |
| 42 | 17.951 | 2507 | 2510 | 2515 | rBV3  | 20257   | 30240   | 0.11% | 0.039% |
| 43 | 18.234 | 2554 | 2558 | 2561 | rBV2  | 48783   | 66392   | 0.25% | 0.085% |
| 44 | 18.281 | 2561 | 2566 | 2572 | rVV2  | 405726  | 689624  | 2.61% | 0.887% |
| 45 | 18.345 | 2572 | 2577 | 2587 | rVV   | 1369483 | 1744392 | 6.60% | 2.244% |
| 46 | 18.416 | 2587 | 2589 | 2592 | rVB   | 41379   | 41799   | 0.16% | 0.054% |
| 47 | 18.622 | 2620 | 2624 | 2632 | rBV2  | 44717   | 82811   | 0.31% | 0.107% |
| 48 | 18.910 | 2670 | 2673 | 2677 | rVB   | 55780   | 64306   | 0.24% | 0.083% |
| 49 | 19.004 | 2682 | 2689 | 2693 | rBV10 | 53145   | 159669  | 0.60% | 0.205% |
| 50 | 19.228 | 2723 | 2727 | 2732 | rVB   | 61232   | 78214   | 0.30% | 0.101% |
| 51 | 19.428 | 2758 | 2761 | 2763 | rBV   | 150024  | 169995  | 0.64% | 0.219% |
| 52 | 19.457 | 2763 | 2766 | 2767 | rVV   | 242892  | 282060  | 1.07% | 0.363% |
| 53 | 19.481 | 2767 | 2770 | 2777 | rVB   | 387243  | 558040  | 2.11% | 0.718% |
| 54 | 19.569 | 2781 | 2785 | 2797 | rBV2  | 526644  | 716524  | 2.71% | 0.922% |
| 55 | 19.810 | 2819 | 2826 | 2835 | rBV2  | 989961  | 1631647 | 6.18% | 2.099% |
| 56 | 20.298 | 2906 | 2909 | 2915 | rBV2  | 95788   | 110619  | 0.42% | 0.142% |
| 57 | 20.616 | 2959 | 2963 | 2973 | rVB3  | 194848  | 302160  | 1.14% | 0.389% |
| 58 | 20.780 | 2989 | 2991 | 2995 | rBV   | 76998   | 108169  | 0.41% | 0.139% |
| 59 | 21.245 | 3065 | 3070 | 3085 | rVB3  | 571889  | 1140497 | 4.32% | 1.467% |
| 60 | 21.445 | 3101 | 3104 | 3109 | rVB   | 179861  | 197878  | 0.75% | 0.255% |
| 61 | 21.563 | 3118 | 3124 | 3129 | rBV   | 1165029 | 2161410 | 8.18% | 2.780% |
| 62 | 21.892 | 3177 | 3180 | 3185 | rVB5  | 140882  | 163456  | 0.62% | 0.210% |
| 63 | 22.151 | 3221 | 3224 | 3230 | rBV4  | 436523  | 954622  | 3.61% | 1.228% |
| 64 | 22.204 | 3230 | 3233 | 3242 | rVB4  | 625691  | 1166630 | 4.42% | 1.501% |
| 65 | 22.286 | 3242 | 3247 | 3258 | rVB2  | 916854  | 1489218 | 5.64% | 1.916% |
| 66 | 22.404 | 3260 | 3267 | 3272 | rBV4  | 448475  | 938873  | 3.55% | 1.208% |
| 67 | 22.551 | 3287 | 3292 | 3300 | rVB3  | 486245  | 842478  | 3.19% | 1.084% |
| 68 | 22.663 | 3307 | 3311 | 3314 | rBV2  | 370670  | 569986  | 2.16% | 0.733% |
| 69 | 22.969 | 3359 | 3363 | 3373 | rBV4  | 291935  | 687453  | 2.60% | 0.884% |
| 70 | 23.133 | 3386 | 3391 | 3396 | rVB   | 1180757 | 1926248 | 7.29% | 2.478% |
| 71 | 23.286 | 3411 | 3417 | 3422 | rBV   | 1156275 | 1941043 | 7.35% | 2.497% |
| 72 | 23.339 | 3422 | 3426 | 3440 | rVB2  | 558294  | 1834806 | 6.94% | 2.360% |
| 73 | 23.951 | 3524 | 3530 | 3537 | rBV3  | 506330  | 1071542 | 4.06% | 1.378% |
| 74 | 24.116 | 3552 | 3558 | 3568 | rVB   | 627290  | 1298795 | 4.92% | 1.671% |
| 75 | 24.739 | 3657 | 3664 | 3673 | rVB   | 969612  | 2108609 | 7.98% | 2.712% |
| 76 | 25.651 | 3814 | 3819 | 3829 | rVB2  | 288922  | 741119  | 2.80% | 0.953% |

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Sample : 03088-05  
Misc :  
ALS Vial : 41 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
DCFR1

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\_P\Methods\SFAM-EPA-BP060723.MA.M  
Title : SVOA CALIBRATION

|    |        |      |      |      |      |         |          |         |         |
|----|--------|------|------|------|------|---------|----------|---------|---------|
| 77 | 26.204 | 3907 | 3913 | 3930 | rVB2 | 449993  | 1299838  | 4.92%   | 1.672%  |
| 78 | 26.727 | 3994 | 4002 | 4027 | rVB  | 770006  | 3278061  | 12.41%  | 4.217%  |
| 79 | 27.804 | 4175 | 4185 | 4206 | rVB3 | 872421  | 4084042  | 15.46%  | 5.253%  |
| 80 | 28.886 | 4351 | 4369 | 4387 | rVB2 | 6179783 | 26421599 | 100.00% | 33.987% |

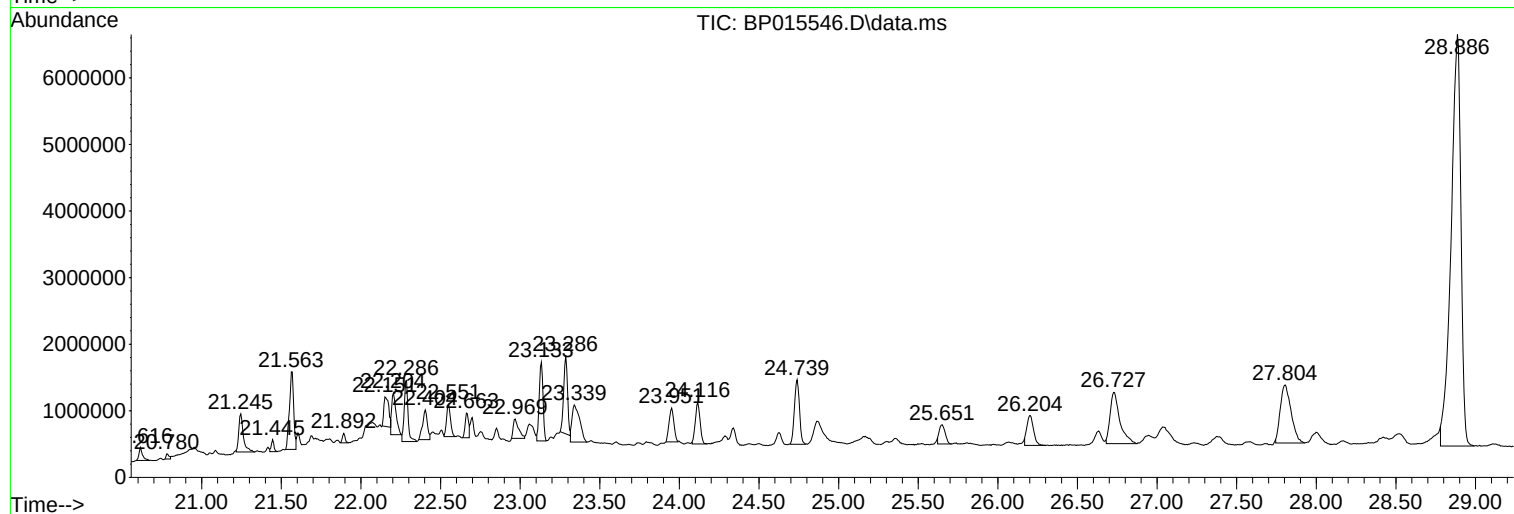
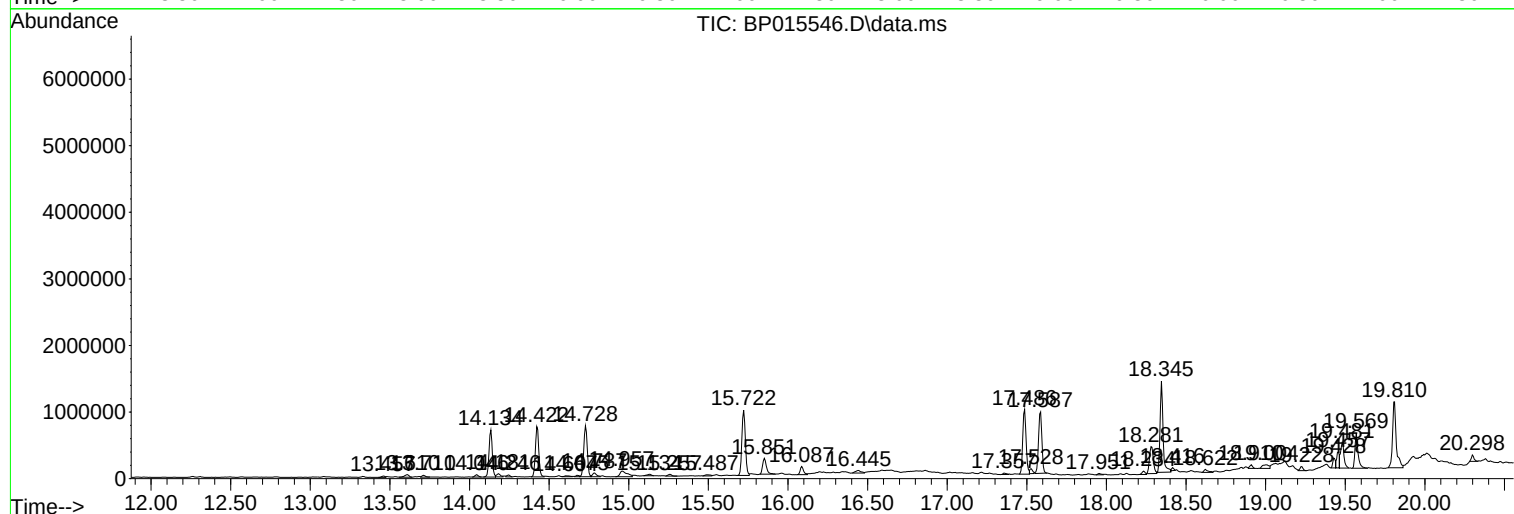
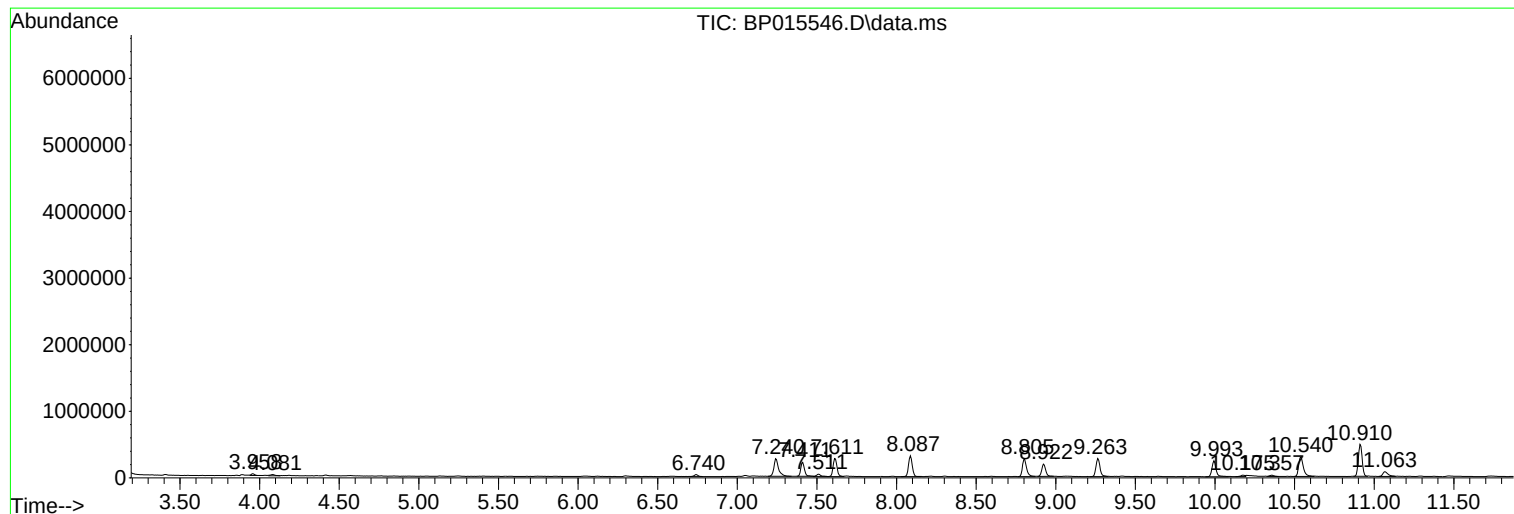
Sum of corrected areas: 77740150

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Instrument :  
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Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP060723.MA.M  
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP061423\  
Data File : BP015546.D  
Acq On : 15 Jun 2023 07:45  
Operator : MA/JU  
Sample : 03088-05  
Misc :  
ALS Vial : 41 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
DCFR1

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP060723.MA.M  
Quant Title : SVOA CALIBRATION

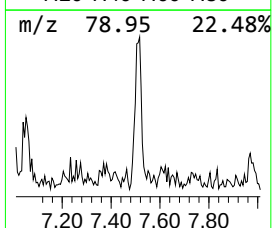
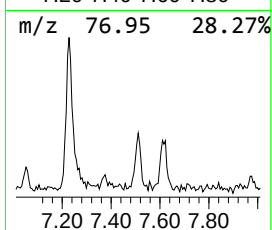
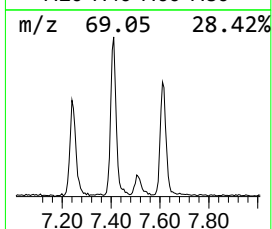
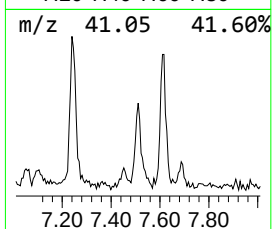
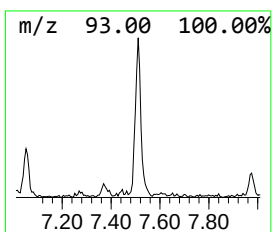
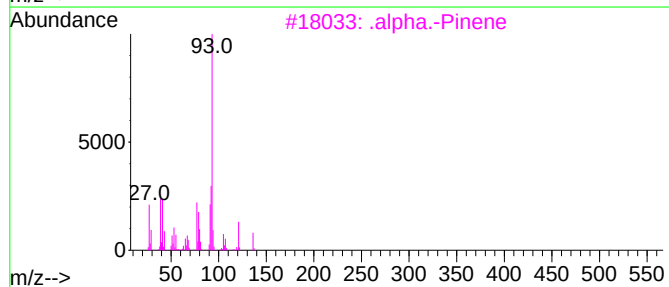
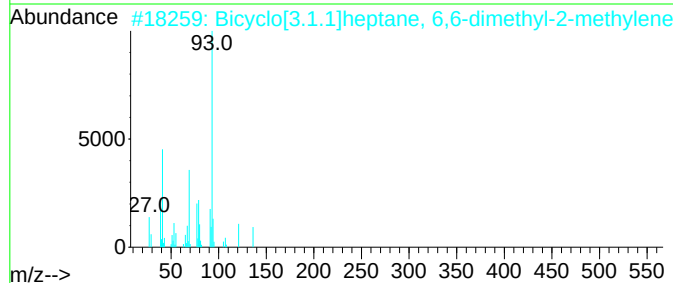
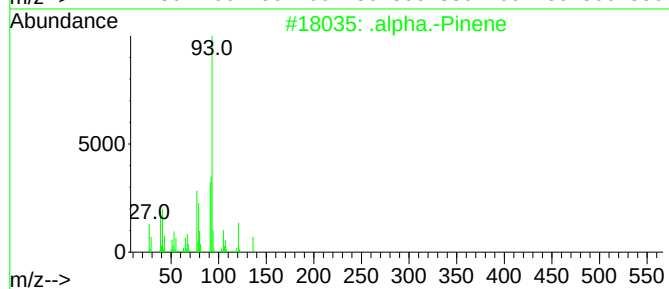
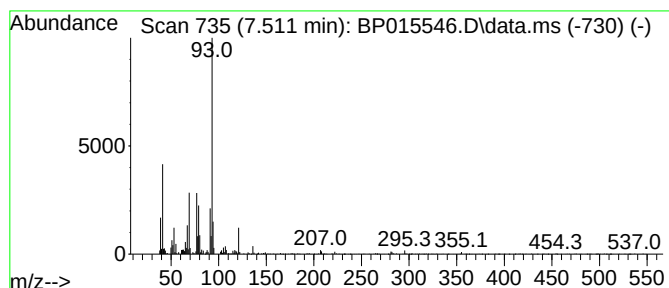
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 1 .alpha.-Pinene Concentration Rank 22

| R.T.  | EstConc    | Area  | Relative to ISTD       | R.T.  |
|-------|------------|-------|------------------------|-------|
| 7.511 | 2.64 ng/ul | 68185 | 1,4-Dichlorobenzene-d4 | 8.087 |

| Hit# | of                                  | 5 | Tentative ID   | MW  | MolForm | CAS#        | Qual |
|------|-------------------------------------|---|----------------|-----|---------|-------------|------|
| 1    | .                                   |   | .alpha.-Pinene | 136 | C10H16  | 000080-56-8 | 87   |
| 2    | Bicyclo[3.1.1]heptane, 6,6-dimet... |   |                | 136 | C10H16  | 018172-67-3 | 87   |
| 3    | .                                   |   | .alpha.-Pinene | 136 | C10H16  | 000080-56-8 | 87   |
| 4    | .                                   |   | .beta.-Pinene  | 136 | C10H16  | 000127-91-3 | 87   |
| 5    | Tricyclo[2.2.1.0(2,6)]heptane, 1... |   |                | 136 | C10H16  | 000488-97-1 | 86   |



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ALS Vial : 41 Sample Multiplier: 1

Instrument :  
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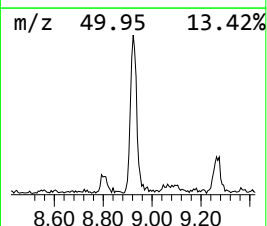
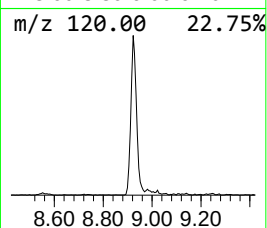
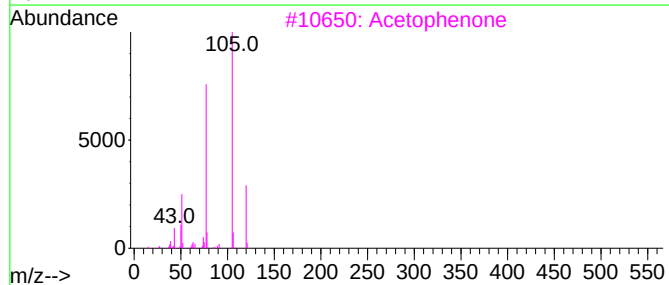
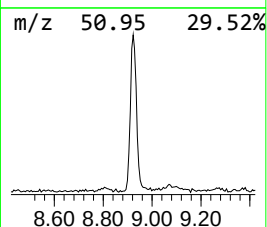
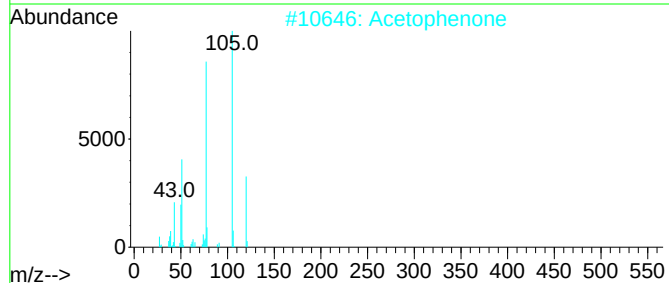
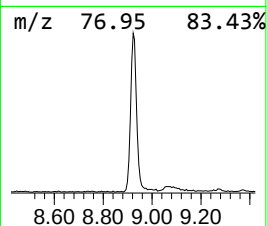
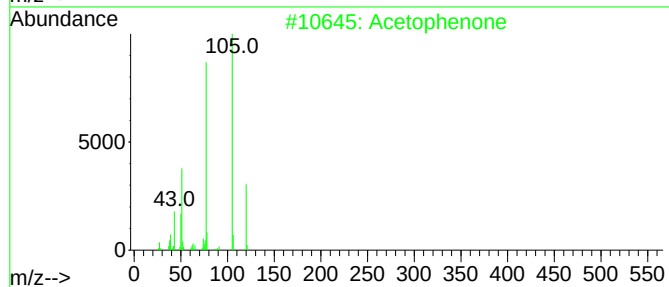
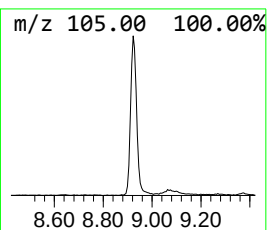
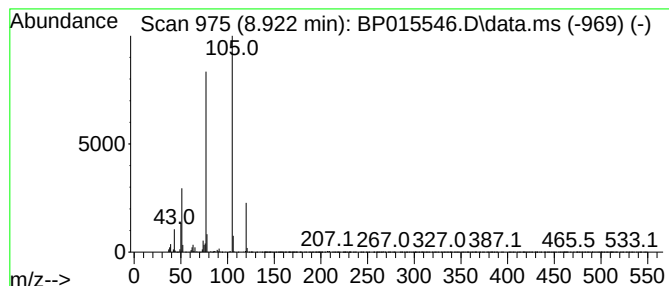
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP060723.MA.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 2 Aceto phenone Concentration Rank 9

| R.T.  | EstConc     | Area   | Relative to ISTD       | R.T.  |
|-------|-------------|--------|------------------------|-------|
| 8.922 | 12.30 ng/ul | 317421 | 1,4-Dichlorobenzene-d4 | 8.087 |

| Hit# | of           | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|--------------|---|--------------|-----|---------|-------------|------|
| 1    | Acetophenone |   |              | 120 | C8H8O   | 000098-86-2 | 97   |
| 2    | Acetophenone |   |              | 120 | C8H8O   | 000098-86-2 | 96   |
| 3    | Acetophenone |   |              | 120 | C8H8O   | 000098-86-2 | 94   |
| 4    | Acetophenone |   |              | 120 | C8H8O   | 000098-86-2 | 91   |
| 5    | Acetophenone |   |              | 120 | C8H8O   | 000098-86-2 | 91   |



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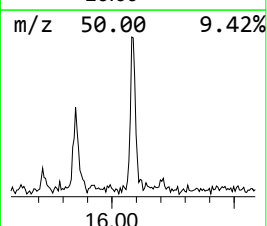
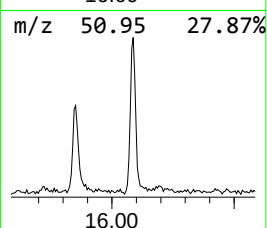
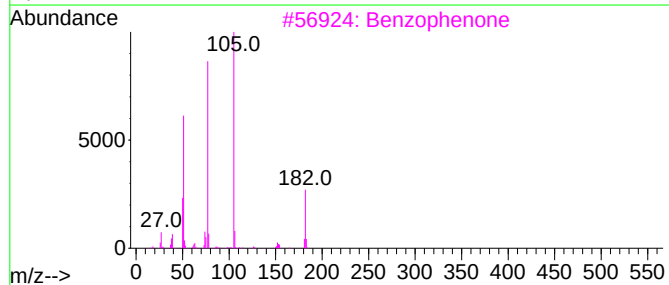
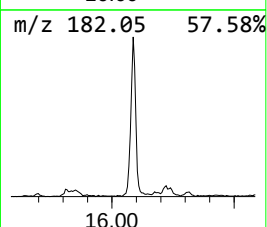
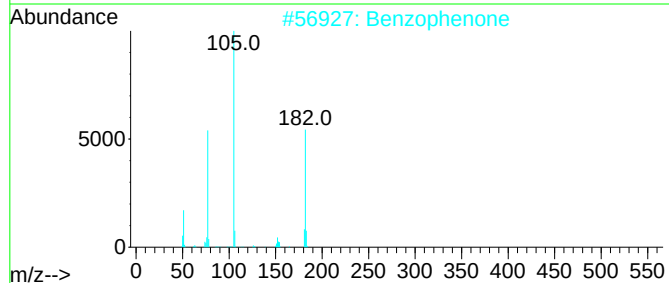
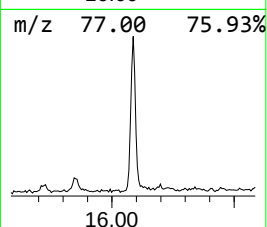
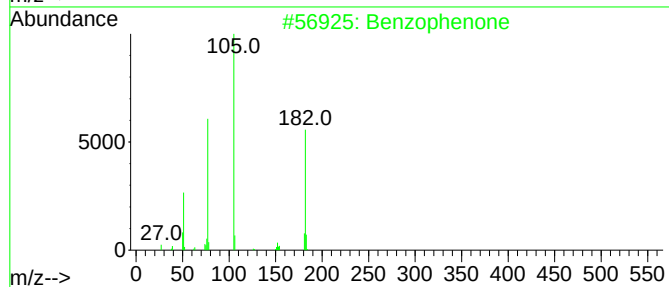
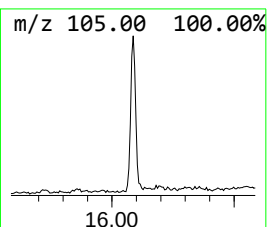
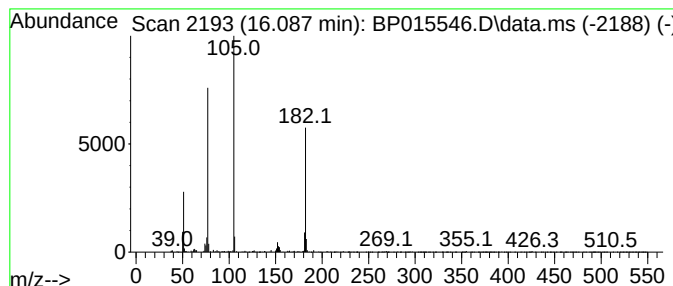
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Quant Title : SVOA CALIBRATION

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

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Peak Number 3 Benzophenone Concentration Rank 20

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 16.087 | 3.20 ng/ul | 182923 | Acenaphthene-d10 | 14.728 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | Benzophenone                        | 182 | C13H10O  | 000119-61-9 | 96   |
| 2    |    |   | Benzophenone                        | 182 | C13H10O  | 000119-61-9 | 95   |
| 3    |    |   | Benzophenone                        | 182 | C13H10O  | 000119-61-9 | 83   |
| 4    |    |   | 1,2,4,5-Tetroxane, 3,3,6,6-tetra... | 396 | C26H20O4 | 016204-36-7 | 78   |
| 5    |    |   | Benzophenone                        | 182 | C13H10O  | 000119-61-9 | 76   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP061423\  
Data File : BP015546.D  
Acq On : 15 Jun 2023 07:45  
Operator : MA/JU  
Sample : 03088-05  
Misc :  
ALS Vial : 41 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
DCFR1

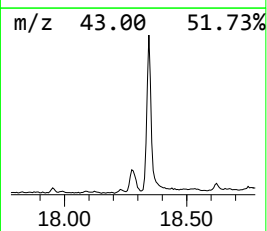
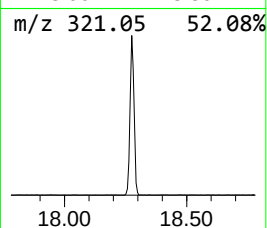
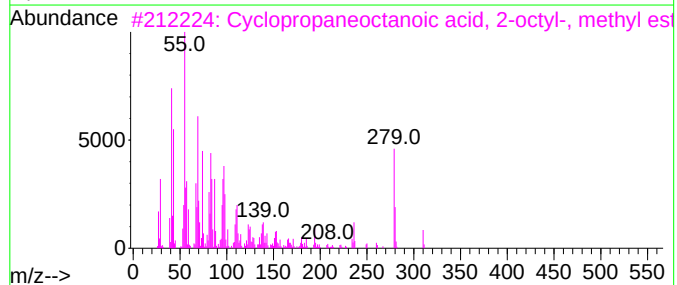
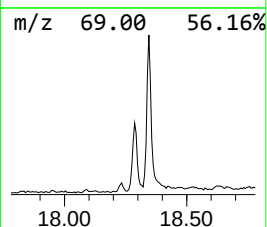
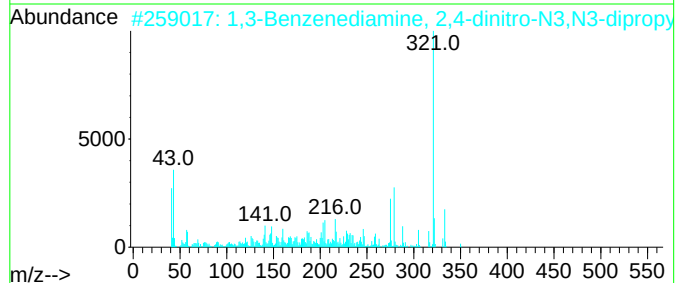
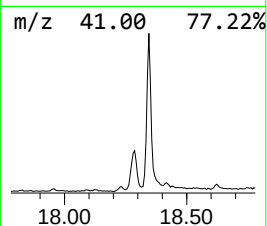
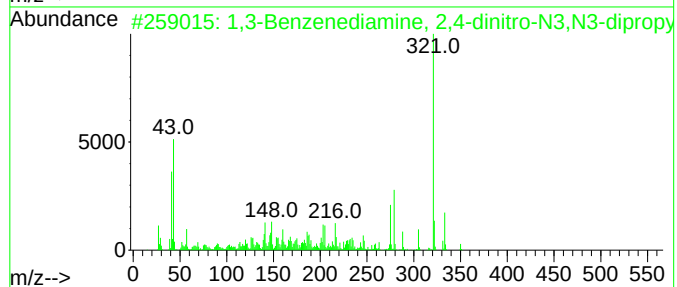
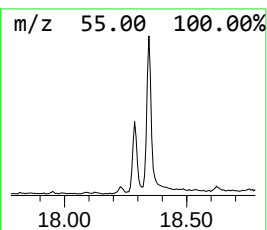
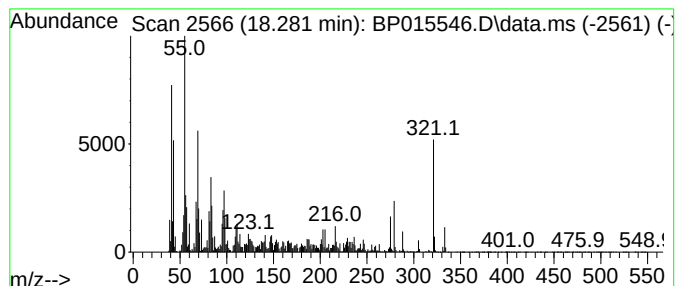
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Quant Title : SVOA CALIBRATION

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

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Peak Number 4 1,3-Benzenediamine, 2,4-din... Concentration Rank 14

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 18.281 | 9.71 ng/ul | 689624 | Phenanthrene-d10 | 17.486 |

| Hit# | of                                  | 5   | Tentative ID | MW           | MolForm | CAS# | Qual |
|------|-------------------------------------|-----|--------------|--------------|---------|------|------|
| 1    | 1,3-Benzenediamine, 2,4-dinitro-... | 350 | C13H17F3N4O4 | 029091-21-2  | 87      |      |      |
| 2    | 1,3-Benzenediamine, 2,4-dinitro-... | 350 | C13H17F3N4O4 | 029091-21-2  | 55      |      |      |
| 3    | Cyclopropaneoctanoic acid, 2-oct... | 310 | C20H38O2     | 005135-07-9  | 38      |      |      |
| 4    | 9-Tricosene, (Z)-                   | 322 | C23H46       | 027519-02-4  | 25      |      |      |
| 5    | 4-Amino-5-nitro-6-[3,4,5-trimeth... | 321 | C13H15N5O5   | 1000254-14-1 | 25      |      |      |





Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP061423\  
Data File : BP015546.D  
Acq On : 15 Jun 2023 07:45  
Operator : MA/JU  
Sample : 03088-05  
Misc :  
ALS Vial : 41 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
DCFR1

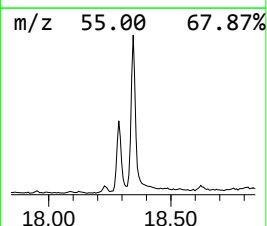
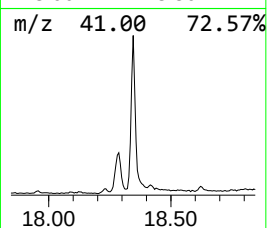
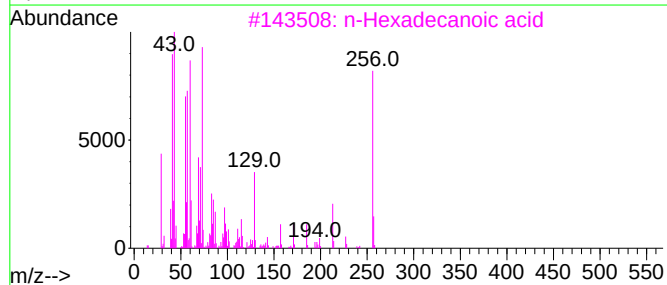
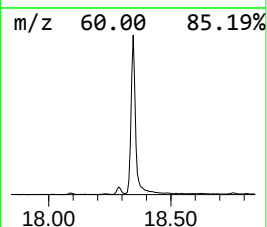
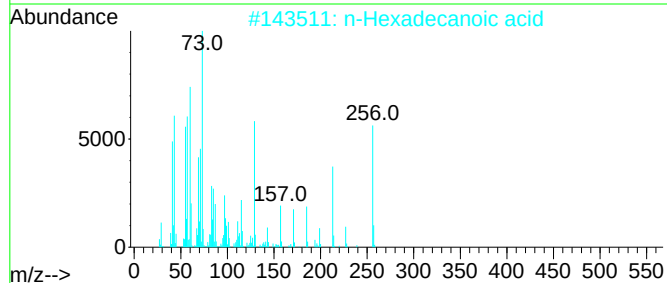
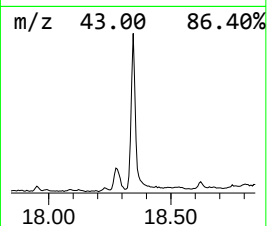
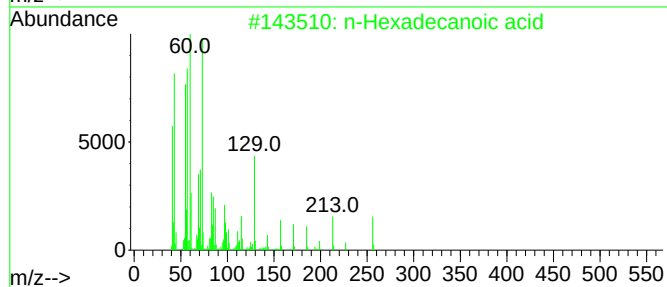
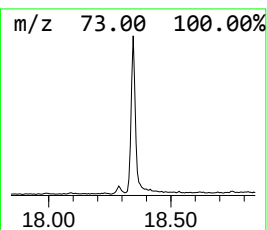
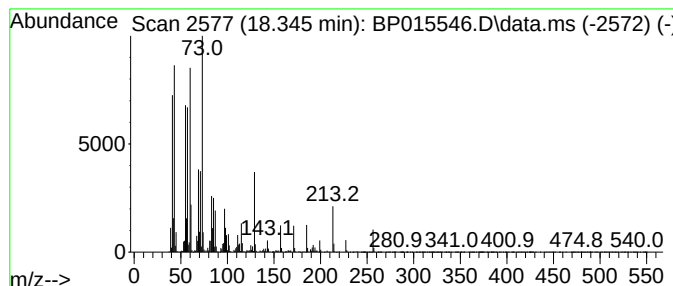
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Quant Title : SVOA CALIBRATION

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

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Peak Number 5 n-Hexadecanoic acid Concentration Rank 6

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 18.345 | 24.56 ng/ul | 1744390 | Phenanthrene-d10 | 17.486 |

| 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 | 99   |
| 3    |    |   | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 98   |
| 4    |    |   | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 98   |
| 5    |    |   | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 91   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP061423\  
Data File : BP015546.D  
Acq On : 15 Jun 2023 07:45  
Operator : MA/JU  
Sample : 03088-05  
Misc :  
ALS Vial : 41 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
DCFR1

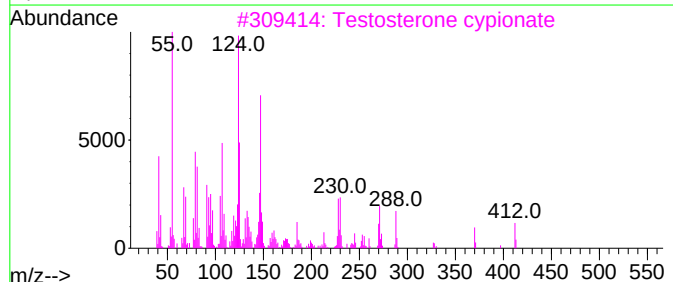
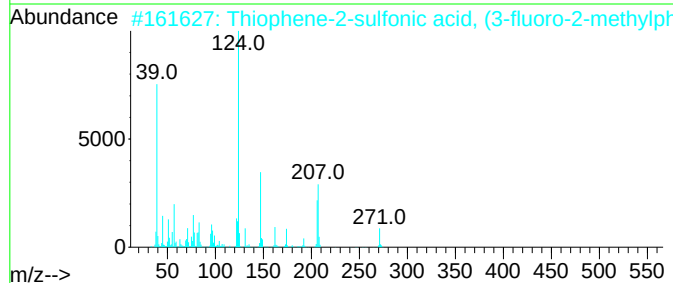
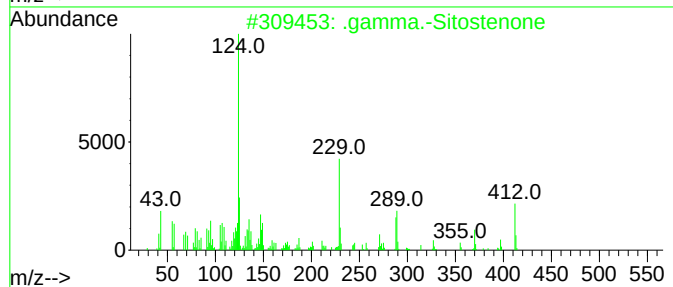
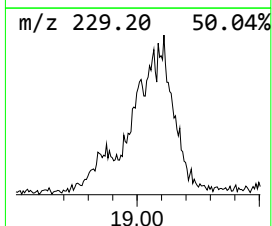
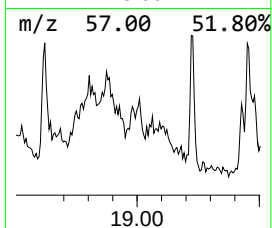
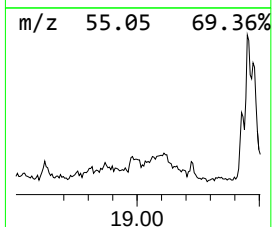
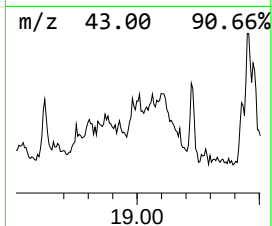
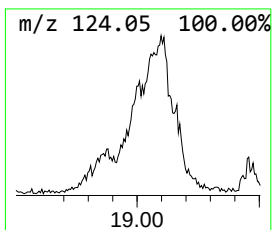
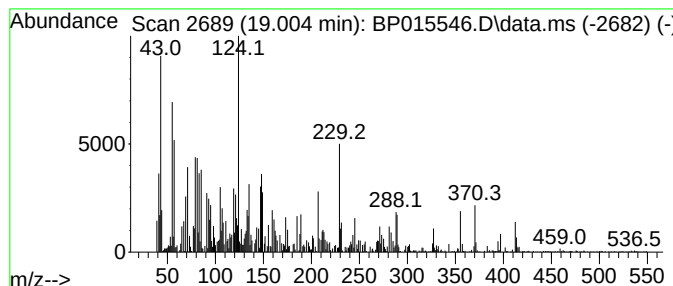
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Quant Title : SVOA CALIBRATION

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

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Peak Number 6 .gamma.-Sitostenone Concentration Rank 24

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 19.004 | 2.25 ng/ul | 159669 | Phenanthrene-d10 | 17.486 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm      | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|--------------|--------------|------|
| 1    |    |   | .gamma.-Sitostenone                 | 412 | C29H48O      | 084924-96-9  | 70   |
| 2    |    |   | Thiophene-2-sulfonic acid, (3-fl... | 271 | C11H10FNO2S2 | 1000311-21-1 | 51   |
| 3    |    |   | Testosterone cypionate              | 412 | C27H40O3     | 000058-20-8  | 46   |
| 4    |    |   | cis-Cyclohex-4-en-1,2-dicarboxyl... | 380 | C23H40O4     | 1000382-76-6 | 43   |
| 5    |    |   | Testosterone cypionate              | 412 | C27H40O3     | 000058-20-8  | 38   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP061423\  
Data File : BP015546.D  
Acq On : 15 Jun 2023 07:45  
Operator : MA/JU  
Sample : 03088-05  
Misc :  
ALS Vial : 41 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
DCFR1

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP060723.MA.M  
Quant Title : SVOA CALIBRATION

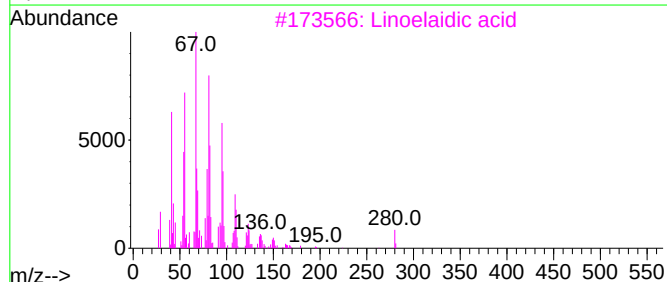
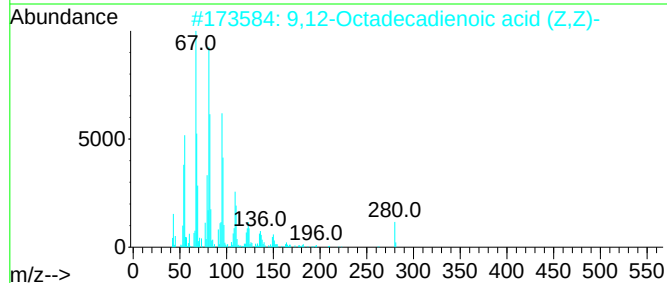
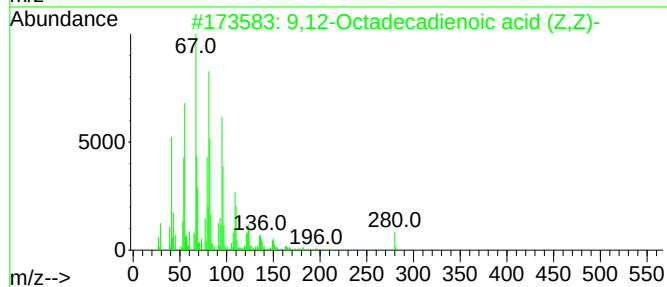
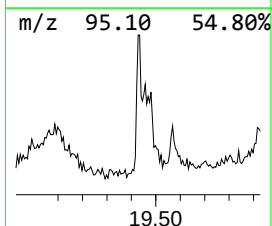
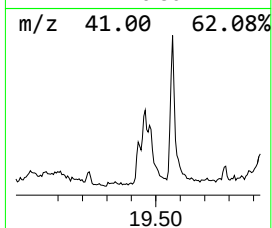
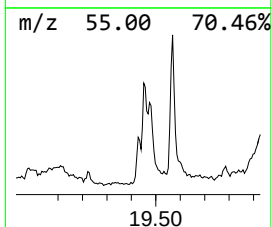
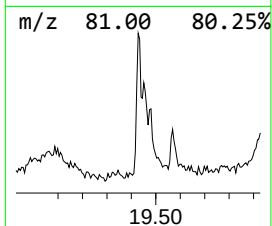
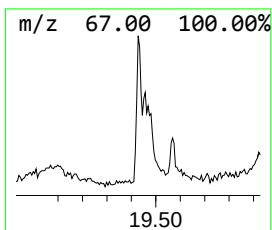
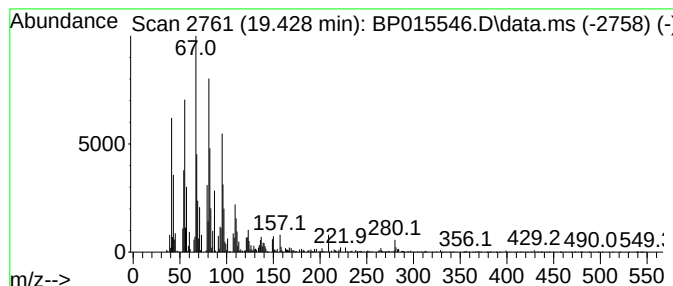
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TIC Integration Parameters: LSCINT.P

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Peak Number 7 9,12-Octadecadienoic acid (... Concentration Rank 23

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 19.428 | 2.39 ng/ul | 169995 | Phenanthrene-d10 | 17.486 |

| Hit# | of | 5 | Tentative ID                     | MW  | MolForm  | CAS#        | Qual |
|------|----|---|----------------------------------|-----|----------|-------------|------|
| 1    |    |   | 9,12-Octadecadienoic acid (Z,Z)- | 280 | C18H32O2 | 000060-33-3 | 99   |
| 2    |    |   | 9,12-Octadecadienoic acid (Z,Z)- | 280 | C18H32O2 | 000060-33-3 | 99   |
| 3    |    |   | Linoleic acid                    | 280 | C18H32O2 | 000506-21-8 | 98   |
| 4    |    |   | 11,14-Eicosadienoic acid         | 308 | C20H36O2 | 002091-39-6 | 97   |
| 5    |    |   | 10E,12Z-Octadecadienoic acid     | 280 | C18H32O2 | 002420-56-6 | 97   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP061423\  
Data File : BP015546.D  
Acq On : 15 Jun 2023 07:45  
Operator : MA/JU  
Sample : 03088-05  
Misc :  
ALS Vial : 41 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
DCFR1

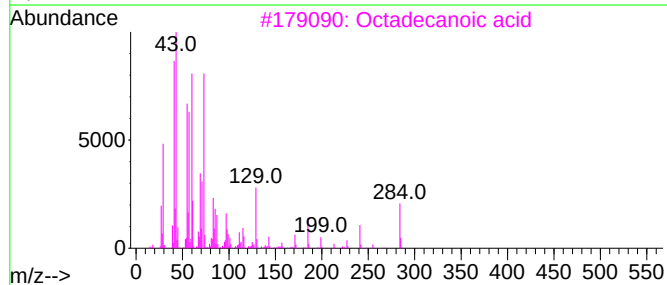
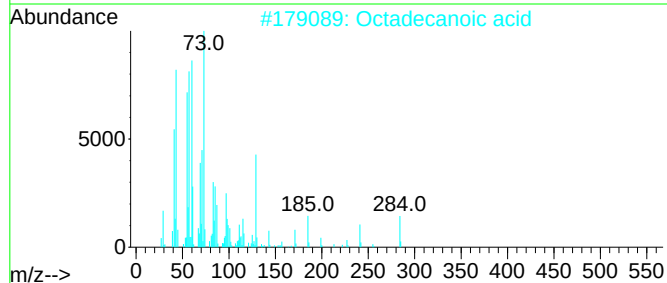
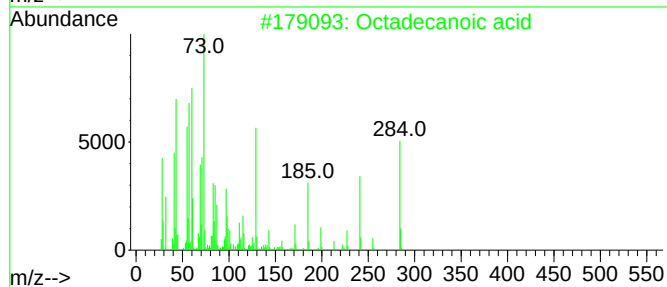
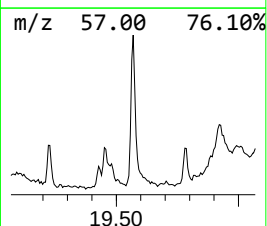
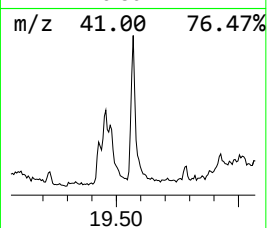
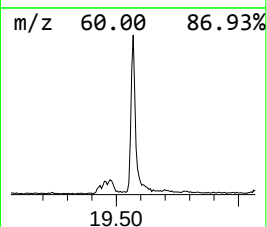
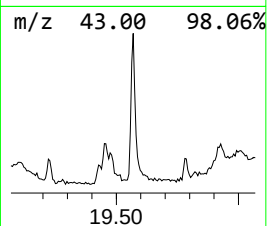
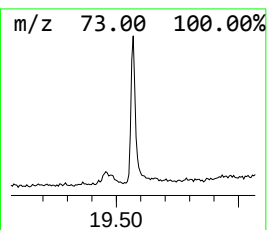
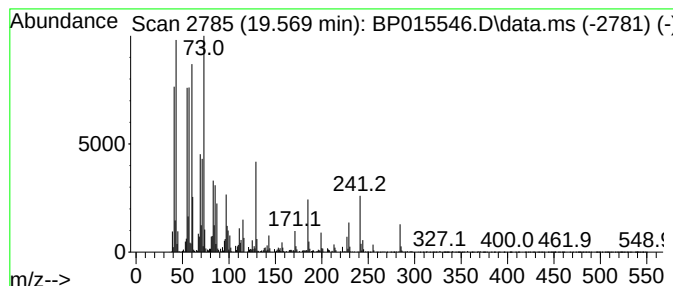
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Quant Title : SVOA CALIBRATION

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

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Peak Number 8 Octadecanoic acid Concentration Rank 18

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 19.569 | 6.63 ng/ul | 716524 | Chrysene-d12     | 21.569 |

| Hit# | of | 5 | Tentative ID      | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------|-----|----------|-------------|------|
| 1    |    |   | Octadecanoic acid | 284 | C18H36O2 | 000057-11-4 | 99   |
| 2    |    |   | Octadecanoic acid | 284 | C18H36O2 | 000057-11-4 | 99   |
| 3    |    |   | Octadecanoic acid | 284 | C18H36O2 | 000057-11-4 | 99   |
| 4    |    |   | Octadecanoic acid | 284 | C18H36O2 | 000057-11-4 | 99   |
| 5    |    |   | Octadecanoic acid | 284 | C18H36O2 | 000057-11-4 | 93   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP061423\  
Data File : BP015546.D  
Acq On : 15 Jun 2023 07:45  
Operator : MA/JU  
Sample : 03088-05  
Misc :  
ALS Vial : 41 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
DCFR1

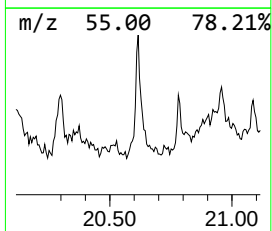
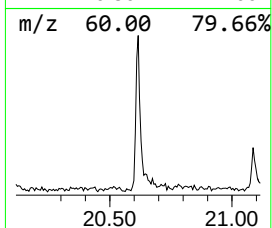
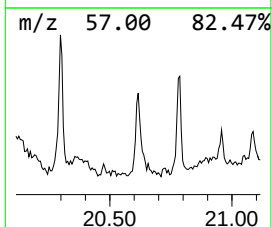
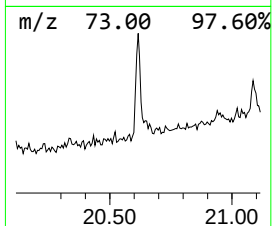
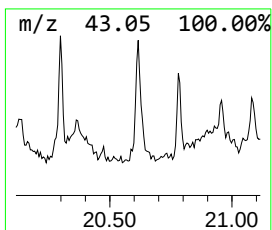
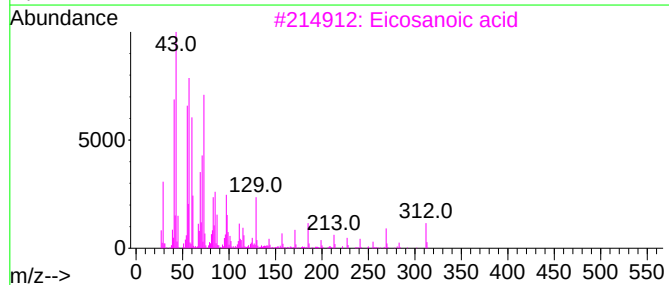
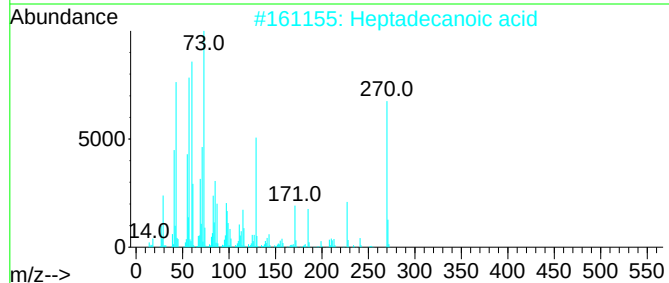
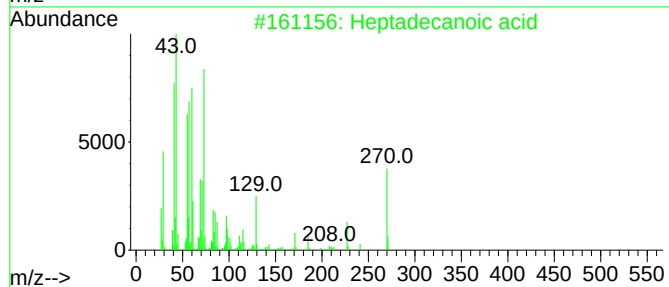
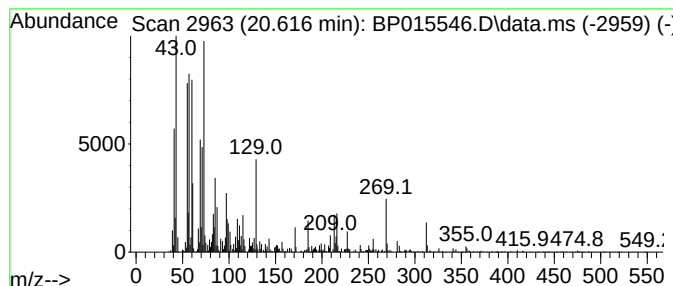
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Quant Title : SVOA CALIBRATION

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

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Peak Number 9 Heptadecanoic acid Concentration Rank 21

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 20.616 | 2.80 ng/ul | 302160 | Chrysene-d12     | 21.569 |

| Hit# | of | 5 | Tentative ID       | MW  | MolForm  | CAS#        | Qual |
|------|----|---|--------------------|-----|----------|-------------|------|
| 1    |    |   | Heptadecanoic acid | 270 | C17H34O2 | 000506-12-7 | 97   |
| 2    |    |   | Heptadecanoic acid | 270 | C17H34O2 | 000506-12-7 | 97   |
| 3    |    |   | Eicosanoic acid    | 312 | C20H40O2 | 000506-30-9 | 93   |
| 4    |    |   | Octadecanoic acid  | 284 | C18H36O2 | 000057-11-4 | 87   |
| 5    |    |   | Tetradecanoic acid | 228 | C14H28O2 | 000544-63-8 | 86   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP061423\  
Data File : BP015546.D  
Acq On : 15 Jun 2023 07:45  
Operator : MA/JU  
Sample : 03088-05  
Misc :  
ALS Vial : 41 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
DCFR1

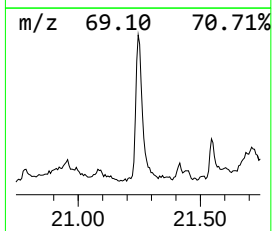
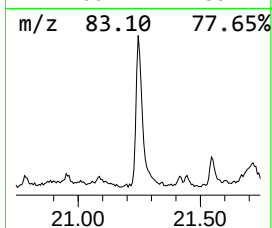
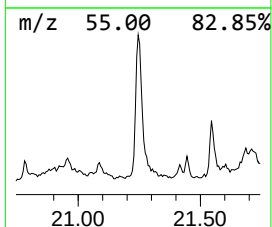
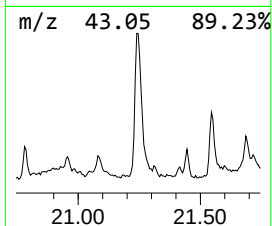
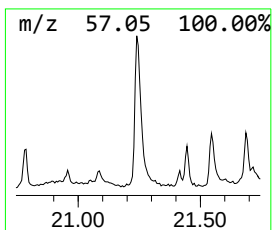
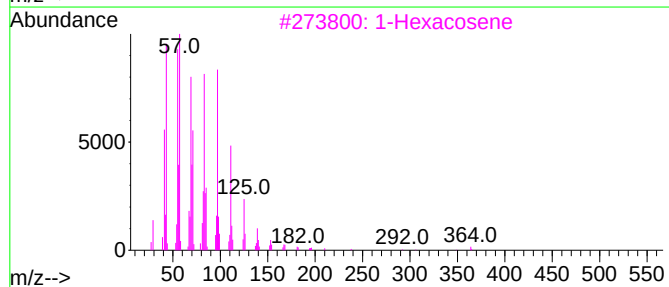
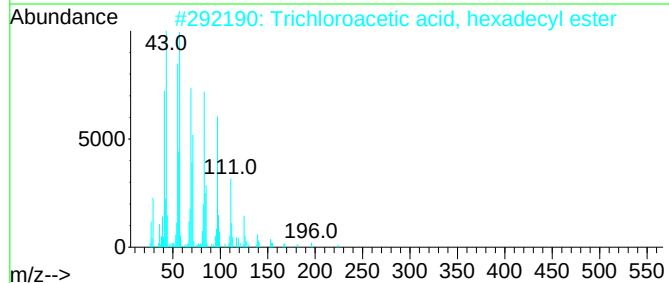
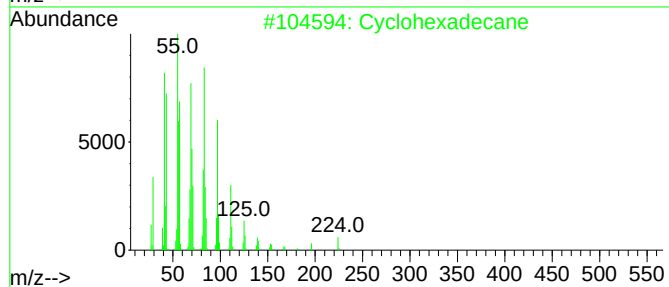
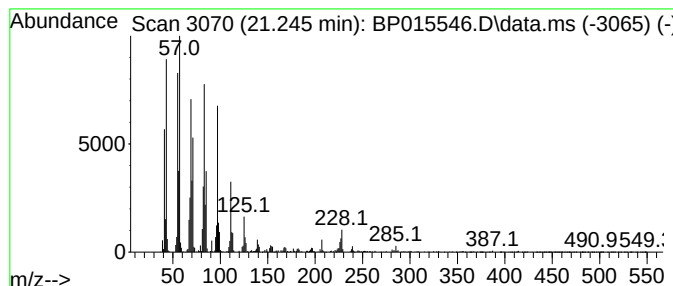
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Quant Title : SVOA CALIBRATION

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

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Peak Number 10 (DEL) Alkane: Cyclic21.245 Concentration Rank 13

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 21.245 | 10.55 ng/ul | 1140500 | Chrysene-d12     | 21.569 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm     | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|-------------|-------------|------|
| 1    |    |   | Cyclohexadecane                     | 224 | C16H32      | 000295-65-8 | 97   |
| 2    |    |   | Trichloroacetic acid, hexadecyl ... | 386 | C18H33Cl3O2 | 074339-54-1 | 94   |
| 3    |    |   | 1-Hexacosene                        | 364 | C26H52      | 018835-33-1 | 94   |
| 4    |    |   | Heptadecyl heptafluorobutyrate      | 452 | C21H35F7O2  | 959085-66-6 | 94   |
| 5    |    |   | 1-Heneicosyl formate                | 340 | C22H44O2    | 077899-03-7 | 93   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP061423\  
Data File : BP015546.D  
Acq On : 15 Jun 2023 07:45  
Operator : MA/JU  
Sample : 03088-05  
Misc :  
ALS Vial : 41 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
DCFR1

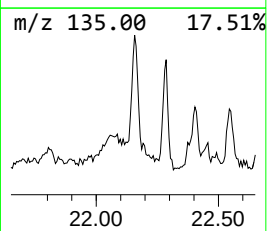
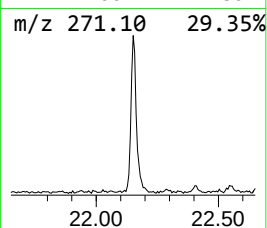
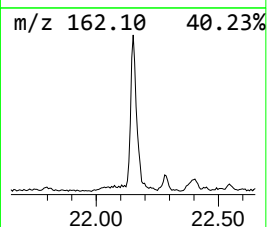
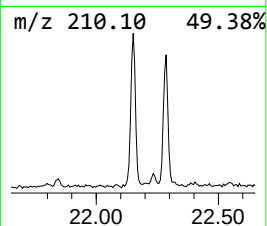
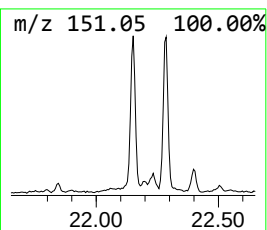
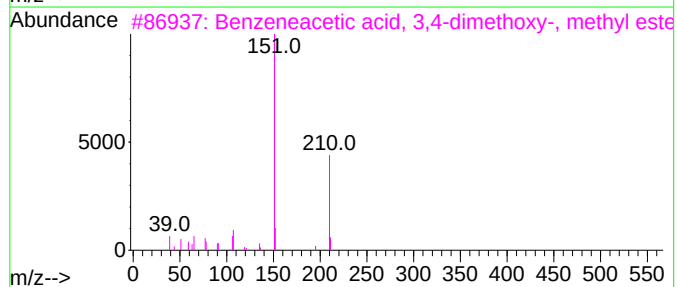
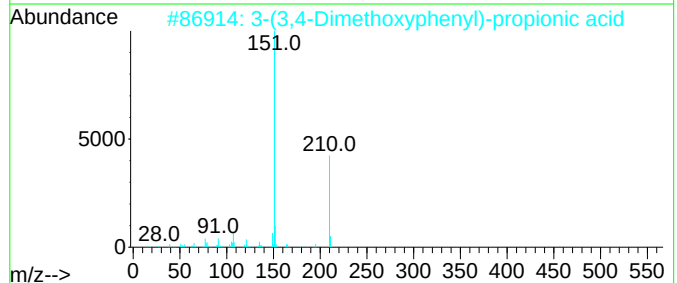
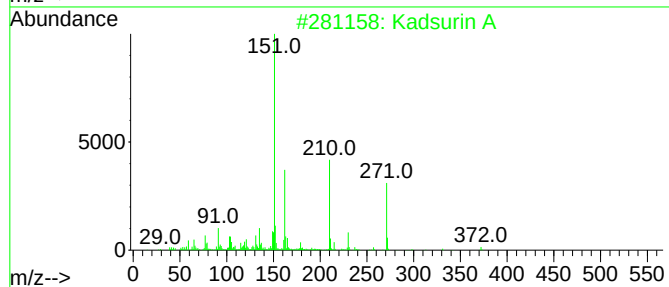
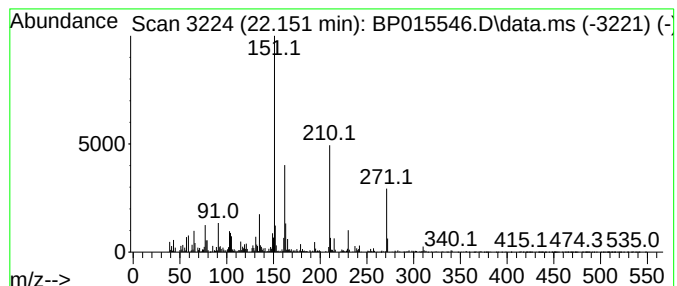
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Quant Title : SVOA CALIBRATION

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

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Peak Number 11 Kadsurin A Concentration Rank 15

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 22.151 | 8.83 ng/ul | 954622 | Chrysene-d12     | 21.569 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | Kadsurin A                          | 372 | C21H24O6 | 099340-07-5 | 83   |
| 2    |    |   | 3-(3,4-Dimethoxyphenyl)-propioni... | 210 | C11H14O4 | 002107-70-2 | 46   |
| 3    |    |   | Benzeneacetic acid, 3,4-dimethox... | 210 | C11H14O4 | 015964-79-1 | 46   |
| 4    |    |   | Benzeneacetic acid, 3,4-dimethox... | 210 | C11H14O4 | 015964-79-1 | 46   |
| 5    |    |   | 3-(3,4-Dimethoxyphenyl)-propioni... | 210 | C11H14O4 | 002107-70-2 | 43   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP061423\  
Data File : BP015546.D  
Acq On : 15 Jun 2023 07:45  
Operator : MA/JU  
Sample : 03088-05  
Misc :  
ALS Vial : 41 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
DCFR1

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP060723.MA.M  
Quant Title : SVOA CALIBRATION

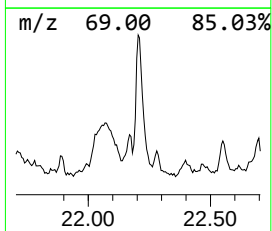
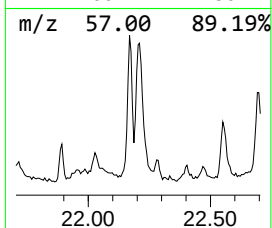
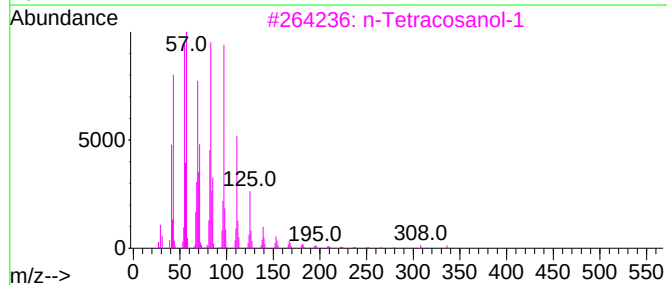
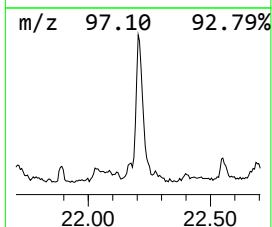
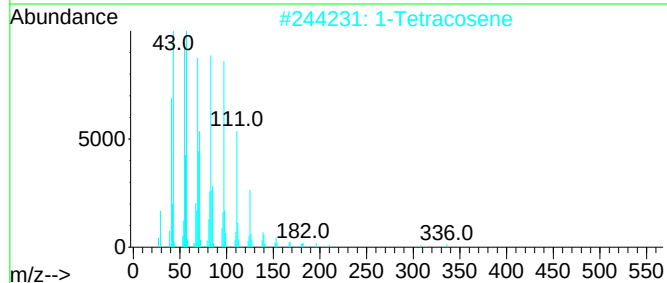
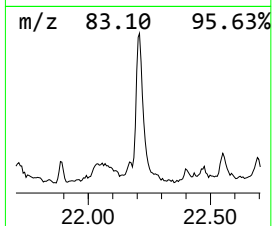
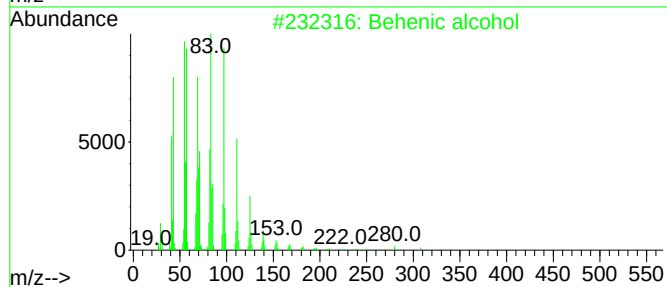
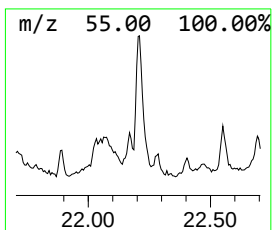
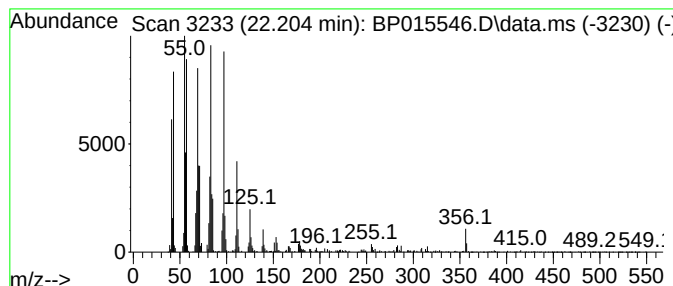
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TIC Integration Parameters: LSCINT.P

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Peak Number 12 Behenic alcohol Concentration Rank 11

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 22.204 | 10.80 ng/ul | 1166630 | Chrysene-d12     | 21.569 |

| Hit# | of | 5 | Tentative ID     | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------|-----|---------|-------------|------|
| 1    |    |   | Behenic alcohol  | 326 | C22H46O | 000661-19-8 | 95   |
| 2    |    |   | 1-Tetracosene    | 336 | C24H48  | 010192-32-2 | 94   |
| 3    |    |   | n-Tetracosanol-1 | 354 | C24H50O | 000506-51-4 | 93   |
| 4    |    |   | 1-Tricosene      | 322 | C23H46  | 018835-32-0 | 91   |
| 5    |    |   | 1-Nonadecene     | 266 | C19H38  | 018435-45-5 | 91   |





Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP061423\  
Data File : BP015546.D  
Acq On : 15 Jun 2023 07:45  
Operator : MA/JU  
Sample : 03088-05  
Misc :  
ALS Vial : 41 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
DCFR1

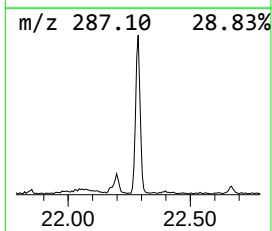
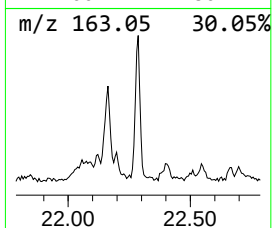
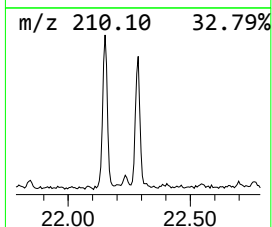
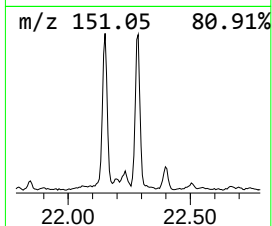
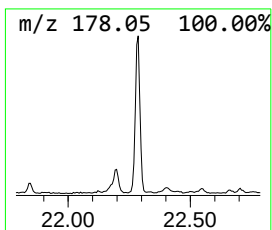
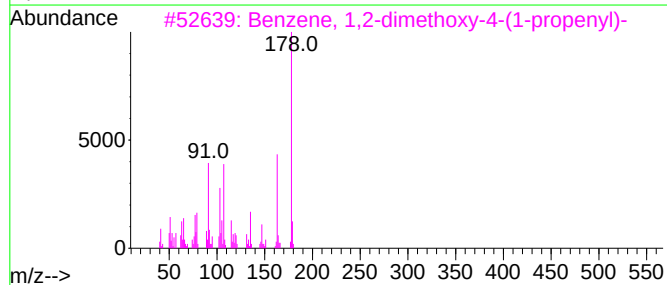
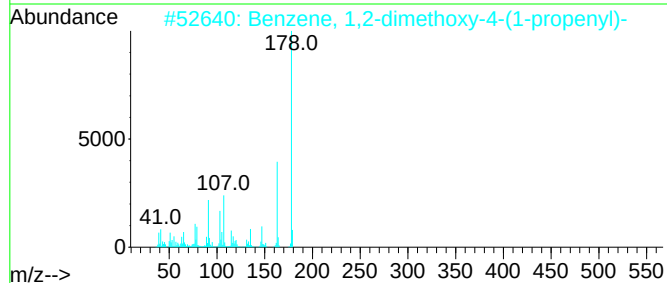
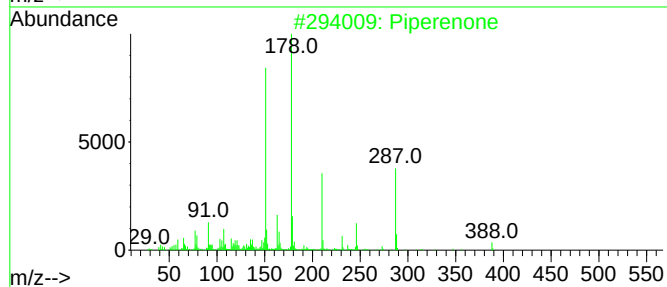
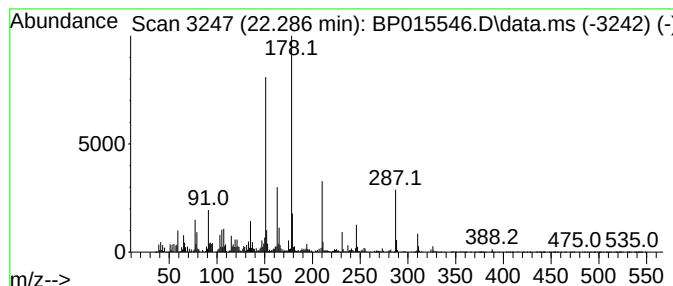
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Quant Title : SVOA CALIBRATION

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

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Peak Number 13 Piperone Concentration Rank 8

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 22.286 | 13.78 ng/ul | 1489220 | Chrysene-d12     | 21.569 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | Piperone                            | 388 | C22H28O6 | 057625-31-7 | 76   |
| 2    |    |   | Benzene, 1,2-dimethoxy-4-(1-prop... | 178 | C11H14O2 | 000093-16-3 | 42   |
| 3    |    |   | Benzene, 1,2-dimethoxy-4-(1-prop... | 178 | C11H14O2 | 000093-16-3 | 42   |
| 4    |    |   | Benzene, 1,2-dimethoxy-4-(1-prop... | 178 | C11H14O2 | 000093-16-3 | 42   |
| 5    |    |   | Benzeneacetic acid, 3,4-dimethox... | 210 | C11H14O4 | 015964-79-1 | 38   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP061423\  
Data File : BP015546.D  
Acq On : 15 Jun 2023 07:45  
Operator : MA/JU  
Sample : 03088-05  
Misc :  
ALS Vial : 41 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
DCFR1

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP060723.MA.M  
Quant Title : SVOA CALIBRATION

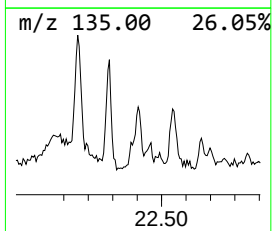
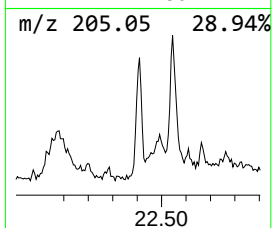
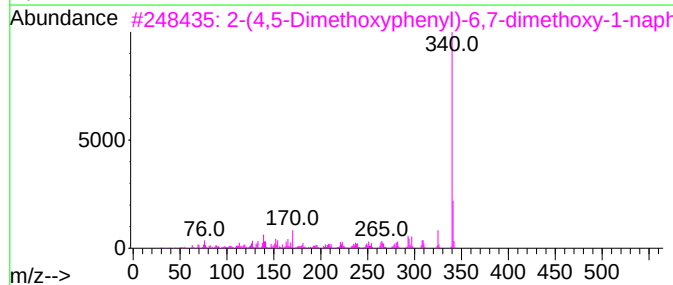
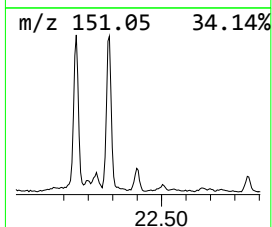
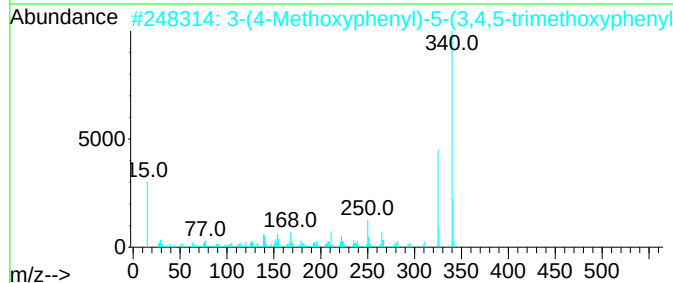
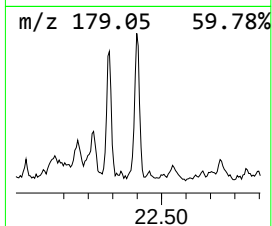
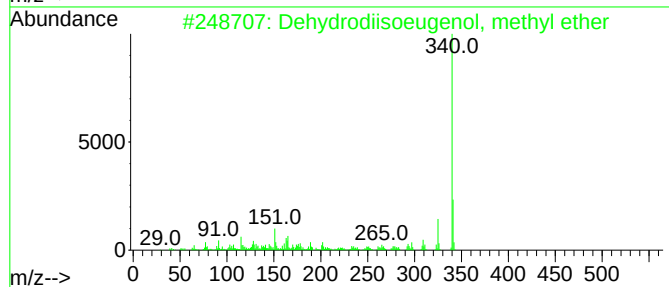
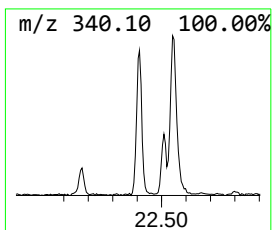
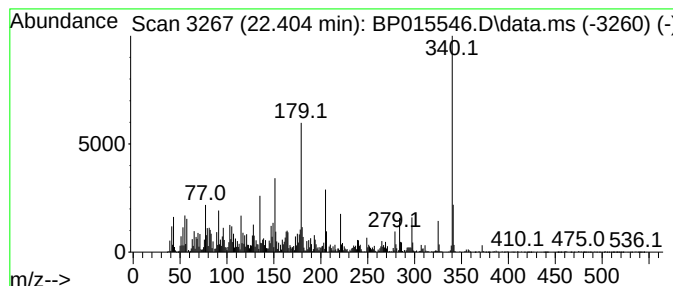
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TIC Integration Parameters: LSCINT.P

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Peak Number 14 Dehydrodiisoeugenol, methyl... Concentration Rank 16

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 22.404 | 8.69 ng/ul | 938873 | Chrysene-d12     | 21.569 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|------------|--------------|------|
| 1    |    |   | Dehydrodiisoeugenol, methyl ether   | 340 | C21H24O4   | 006544-71-4  | 64   |
| 2    |    |   | 3-(4-Methoxyphenyl)-5-(3,4,5-tri... | 340 | C19H20N2O4 | 1000460-31-8 | 46   |
| 3    |    |   | 2-(4,5-Dimethoxyphenyl)-6,7-dime... | 340 | C20H20O5   | 065047-63-4  | 45   |
| 4    |    |   | Dibenzofurane-1,3-dicarbonitrile... | 340 | C20H12N4O2 | 328286-70-0  | 41   |
| 5    |    |   | 6-Methyl-2-(4-nitrophenyl)-4-phe... | 340 | C22H16N2O2 | 073402-99-0  | 41   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP061423\  
Data File : BP015546.D  
Acq On : 15 Jun 2023 07:45  
Operator : MA/JU  
Sample : 03088-05  
Misc :  
ALS Vial : 41 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
DCFR1

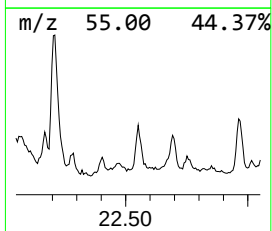
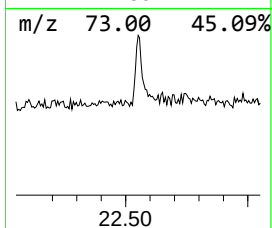
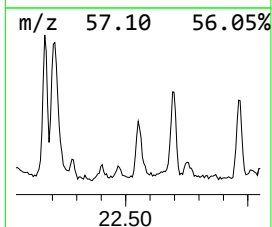
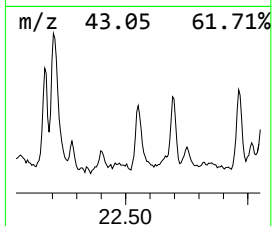
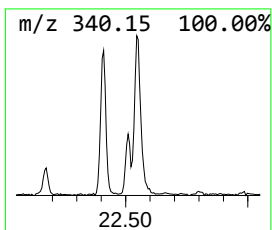
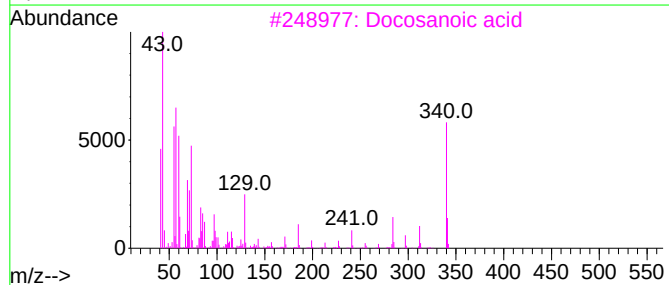
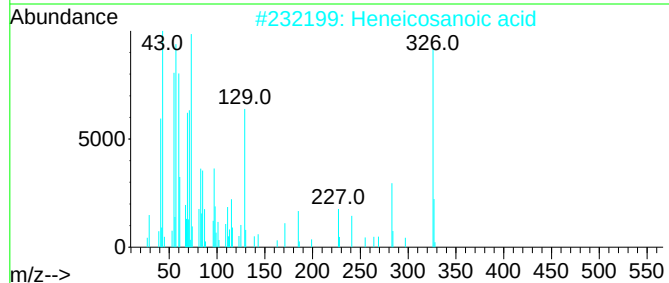
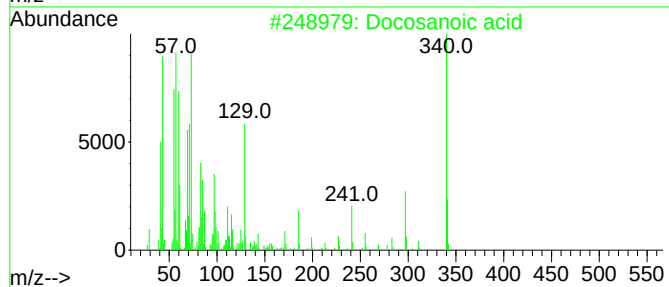
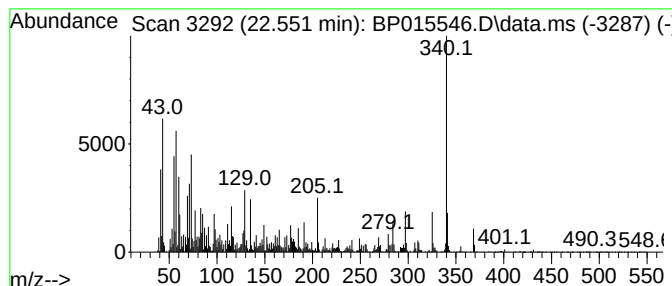
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Quant Title : SVOA CALIBRATION

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

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Peak Number 15 Docosanoic acid Concentration Rank 17

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------------|--------|------------------|-------------|------|
| 22.551    | 7.80 ng/ul                          | 842478 | Chrysene-d12     | 21.569      |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#        | Qual |
| 1         | Docosanoic acid                     | 340    | C22H44O2         | 000112-85-6 | 93   |
| 2         | Heneicosanoic acid                  | 326    | C21H42O2         | 002363-71-5 | 90   |
| 3         | Docosanoic acid                     | 340    | C22H44O2         | 000112-85-6 | 58   |
| 4         | Tetracosanoic acid                  | 368    | C24H48O2         | 000557-59-5 | 55   |
| 5         | Dibenzofurane-1,3-dicarbonitrile... | 340    | C20H12N4O2       | 330158-97-9 | 50   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP061423\  
Data File : BP015546.D  
Acq On : 15 Jun 2023 07:45  
Operator : MA/JU  
Sample : 03088-05  
Misc :  
ALS Vial : 41 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
DCFR1

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP060723.MA.M  
Quant Title : SVOA CALIBRATION

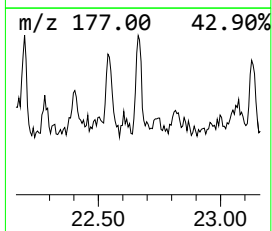
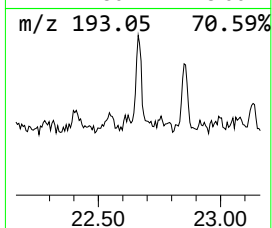
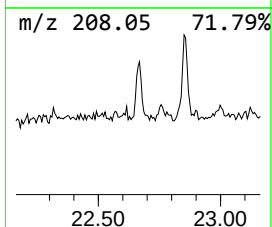
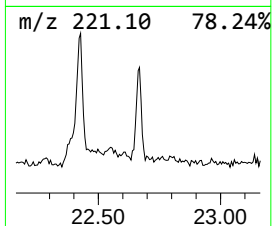
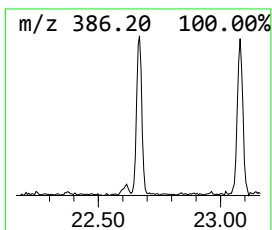
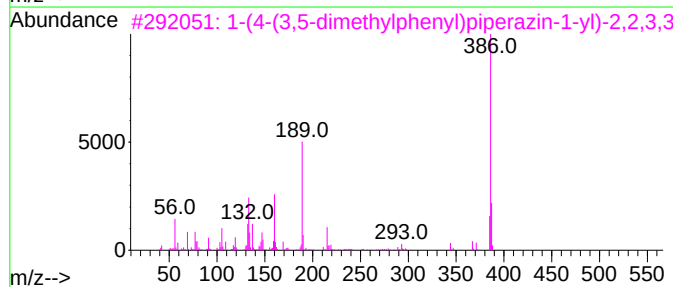
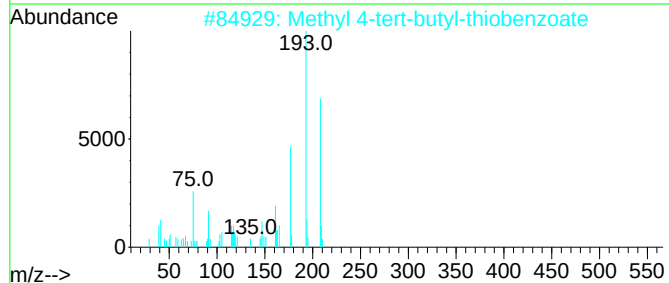
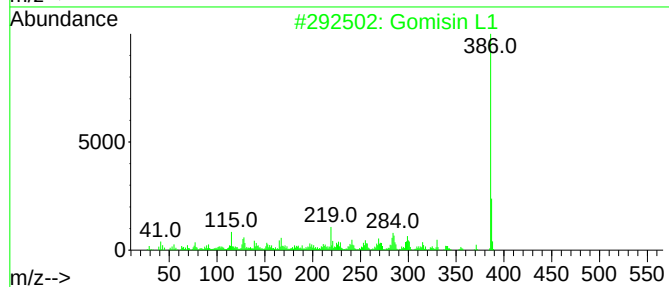
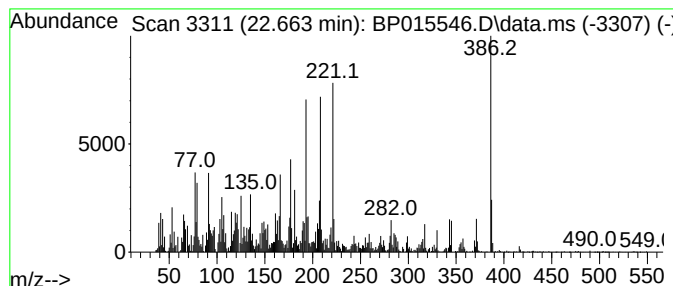
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TIC Integration Parameters: LSCINT.P

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Peak Number 16 unknown-01 Concentration Rank 19

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 22.663 | 5.27 ng/ul | 569986 | Chrysene-d12     | 21.569 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm     | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-------------|--------------|------|
| 1    |    |   | Gomisin L1                          | 386 | C22H26O6    | 082425-43-2  | 44   |
| 2    |    |   | Methyl 4-tert-butyl-thiobenzoate    | 208 | C12H16OS    | 078906-90-8  | 38   |
| 3    |    |   | 1-(4-(3,5-dimethylphenyl)piperaz... | 386 | C16H17F7N2O | 1000455-14-9 | 35   |
| 4    |    |   | 1-(4-(2,5-dimethylphenyl)piperaz... | 386 | C16H17F7N2O | 1000455-15-2 | 35   |
| 5    |    |   | (E)-methyl 3-(3-hydroxy-4-methox... | 208 | C11H12O4    | 1000441-28-8 | 25   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP061423\  
Data File : BP015546.D  
Acq On : 15 Jun 2023 07:45  
Operator : MA/JU  
Sample : 03088-05  
Misc :  
ALS Vial : 41 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
DCFR1

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP060723.MA.M  
Quant Title : SVOA CALIBRATION

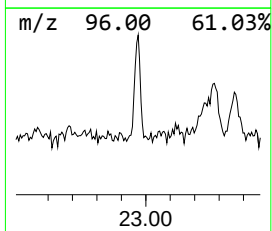
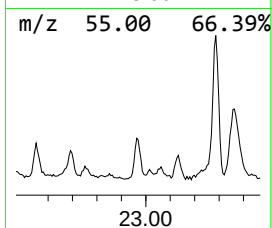
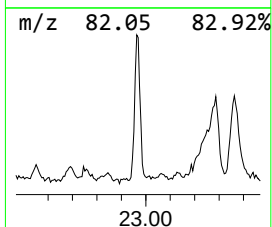
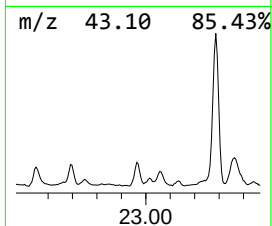
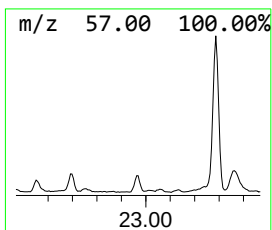
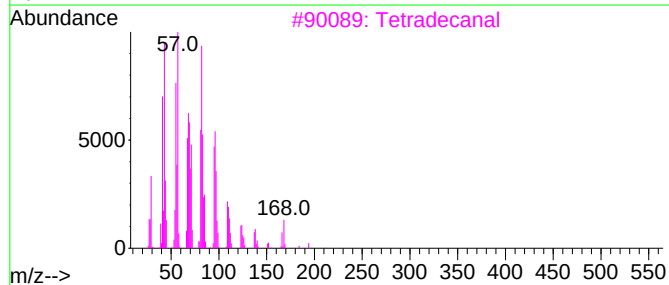
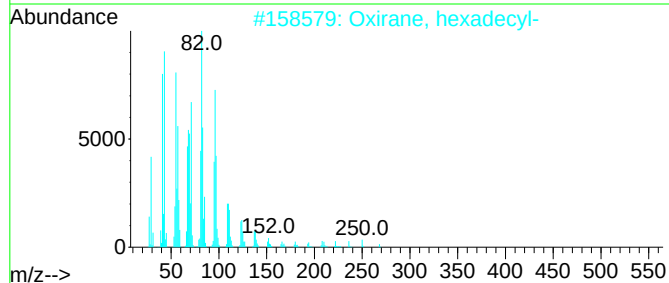
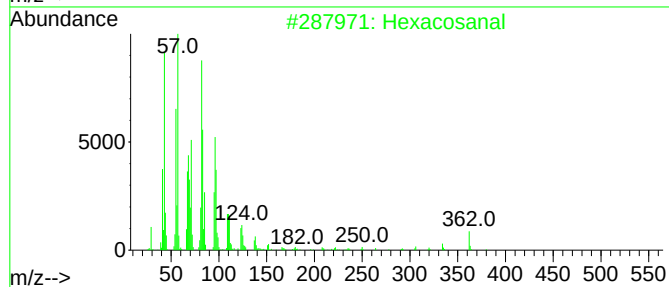
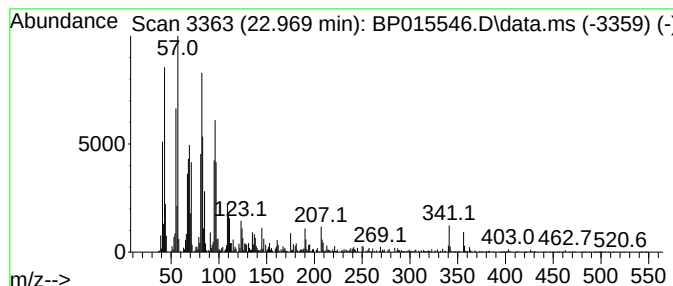
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 17 Hexacosanal Concentration Rank 12

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 22.969 | 10.59 ng/ul | 687453 | Perylene-d12     | 24.116 |

| Hit# | of | 5 | Tentative ID        | MW  | MolForm | CAS#        | Qual |
|------|----|---|---------------------|-----|---------|-------------|------|
| 1    |    |   | Hexacosanal         | 380 | C26H52O | 026627-85-0 | 95   |
| 2    |    |   | Oxirane, hexadecyl- | 268 | C18H36O | 007390-81-0 | 95   |
| 3    |    |   | Tetradecanal        | 212 | C14H28O | 000124-25-4 | 95   |
| 4    |    |   | Tetracosanal        | 352 | C24H48O | 057866-08-7 | 93   |
| 5    |    |   | Nonacosanal         | 422 | C29H58O | 072934-04-4 | 91   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP061423\  
Data File : BP015546.D  
Acq On : 15 Jun 2023 07:45  
Operator : MA/JU  
Sample : 03088-05  
Misc :  
ALS Vial : 41 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
DCFR1

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP060723.MA.M  
Quant Title : SVOA CALIBRATION

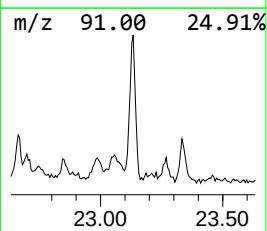
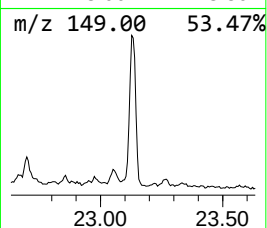
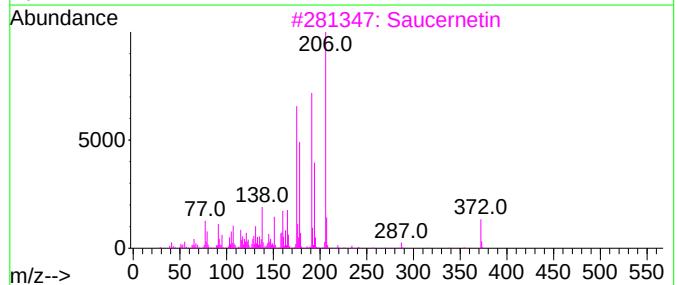
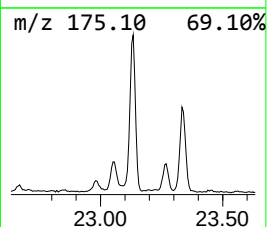
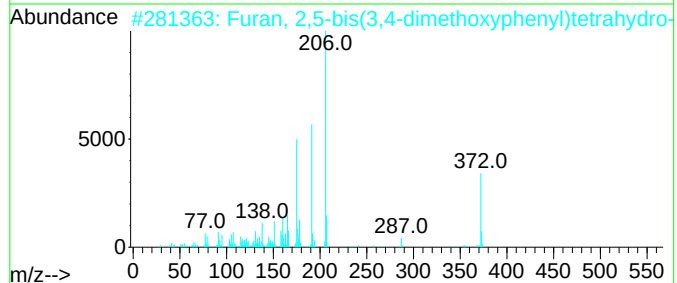
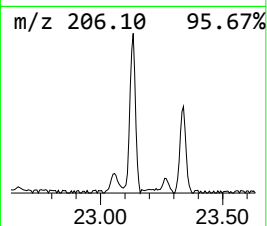
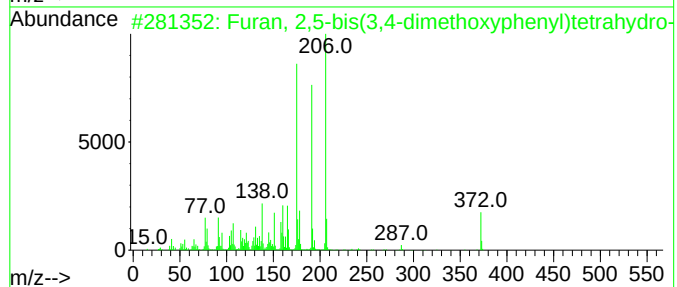
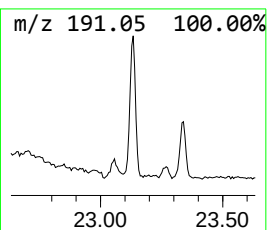
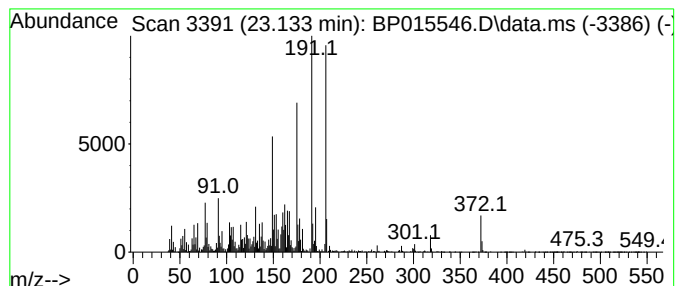
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 18 Furan, 2,5-bis(3,4-dimethox... Concentration Rank 5

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 23.133 | 29.66 ng/ul | 1926250 | Perylene-d12     | 24.116 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | Furan, 2,5-bis(3,4-dimethoxyphen... | 372 | C22H28O5 | 102518-40-1 | 96   |
| 2    |    |   | Furan, 2,5-bis(3,4-dimethoxyphen... | 372 | C22H28O5 | 000528-63-2 | 95   |
| 3    |    |   | Saucernetin                         | 372 | C22H28O5 | 083377-13-3 | 94   |
| 4    |    |   | Furan, 2,5-bis(3,4-dimethoxyphen... | 372 | C22H28O5 | 000528-63-2 | 90   |
| 5    |    |   | Furan, 2,5-bis(3,4-dimethoxyphen... | 372 | C22H28O5 | 102518-40-1 | 70   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP061423\  
Data File : BP015546.D  
Acq On : 15 Jun 2023 07:45  
Operator : MA/JU  
Sample : 03088-05  
Misc :  
ALS Vial : 41 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
DCFR1

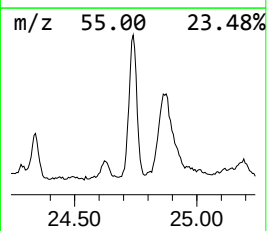
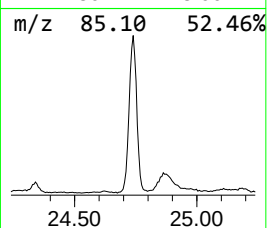
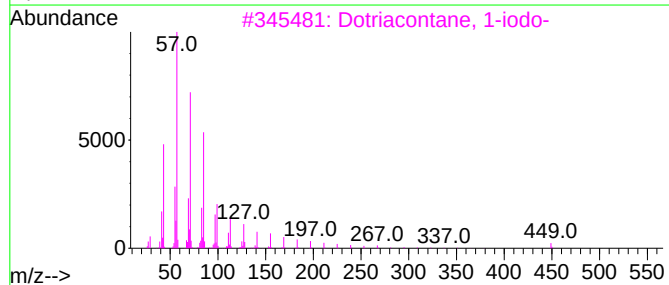
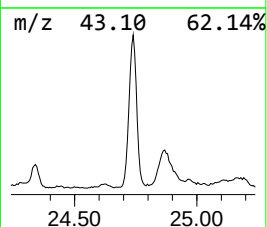
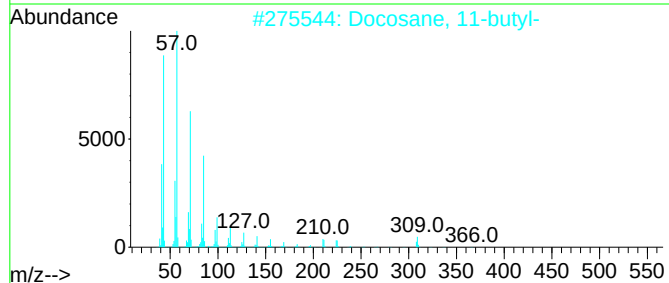
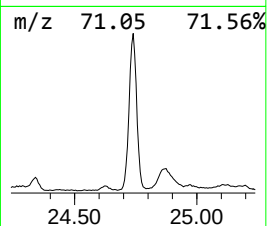
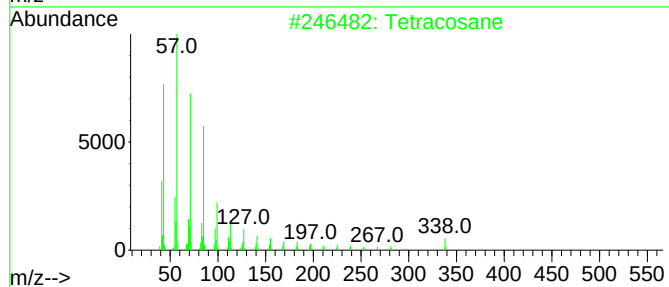
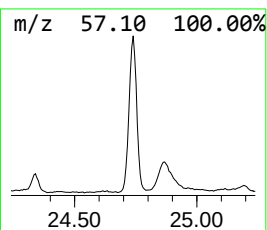
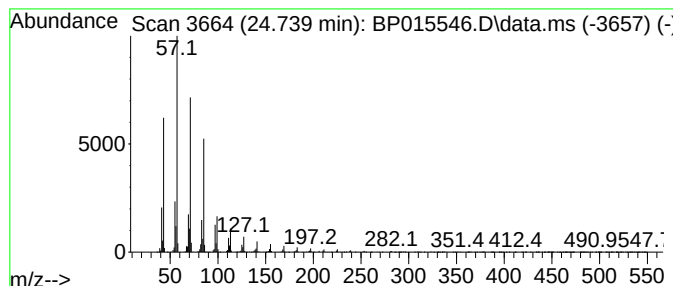
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Quant Title : SVOA CALIBRATION

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

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Peak Number 19 (DEL) Alkane: Straight-Chai... Concentration Rank 4

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 24.739 | 32.47 ng/ul | 2108610 | Perylene-d12     | 24.116 |

| Hit# | of 5 | Tentative ID           | MW  | MolForm | CAS#         | Qual |
|------|------|------------------------|-----|---------|--------------|------|
| 1    |      | Tetracosane            | 338 | C24H50  | 000646-31-1  | 98   |
| 2    |      | Docosane, 11-butyl-    | 366 | C26H54  | 013475-76-8  | 94   |
| 3    |      | Dotriacontane, 1-iodo- | 576 | C32H65I | 1000406-32-4 | 91   |
| 4    |      | Octacosane, 2-methyl-  | 408 | C29H60  | 001560-98-1  | 91   |
| 5    |      | Heneicosane            | 296 | C21H44  | 000629-94-7  | 91   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP061423\  
Data File : BP015546.D  
Acq On : 15 Jun 2023 07:45  
Operator : MA/JU  
Sample : 03088-05  
Misc :  
ALS Vial : 41 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
DCFR1

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP060723.MA.M  
Quant Title : SVOA CALIBRATION

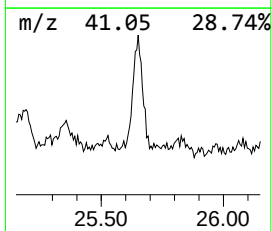
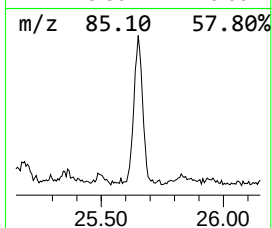
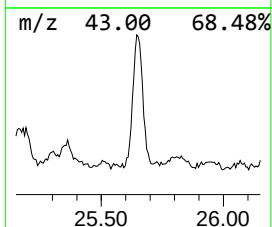
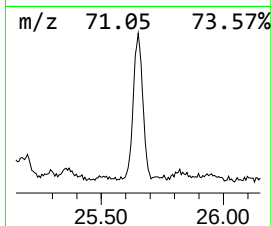
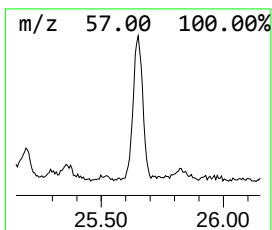
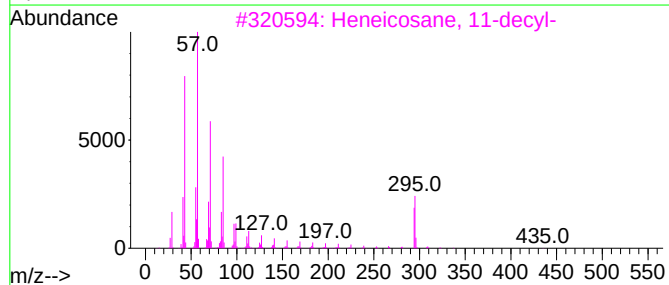
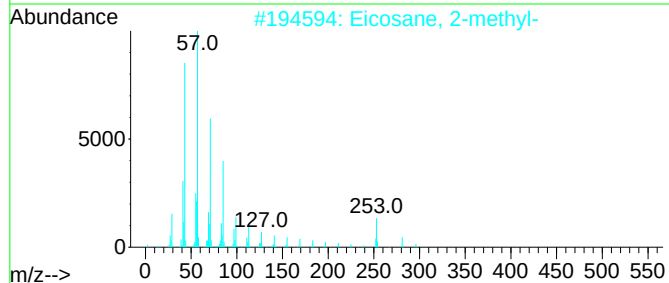
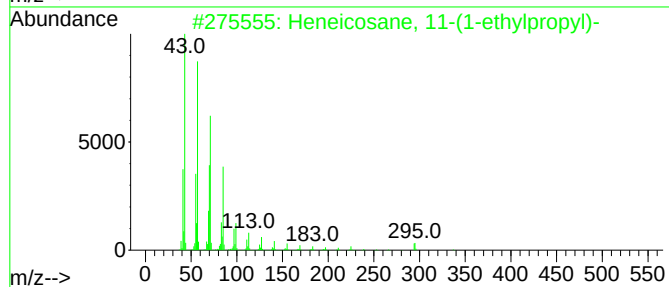
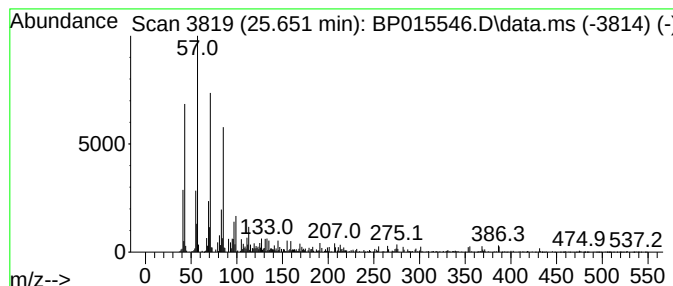
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 20 (DEL) Alkane: Straight-Chai... Concentration Rank 10

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 25.651 | 11.41 ng/ul | 741119 | Perylene-d12     | 24.116 |

| Hit# | of 5 | Tentative ID                     | MW  | MolForm | CAS#        | Qual |
|------|------|----------------------------------|-----|---------|-------------|------|
| 1    |      | Heneicosane, 11-(1-ethylpropyl)- | 366 | C26H54  | 055282-11-6 | 91   |
| 2    |      | Eicosane, 2-methyl-              | 296 | C21H44  | 001560-84-5 | 90   |
| 3    |      | Heneicosane, 11-decyl-           | 437 | C31H64  | 055320-06-4 | 90   |
| 4    |      | Tetracosane, 9-octyl-            | 451 | C32H66  | 055401-54-2 | 90   |
| 5    |      | Docosane, 7-hexyl-               | 394 | C28H58  | 055373-86-9 | 87   |





Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP061423\  
Data File : BP015546.D  
Acq On : 15 Jun 2023 07:45  
Operator : MA/JU  
Sample : 03088-05  
Misc :  
ALS Vial : 41 Sample Multiplier: 1

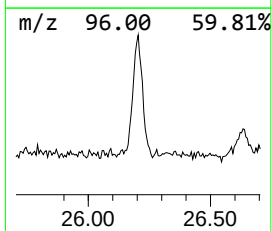
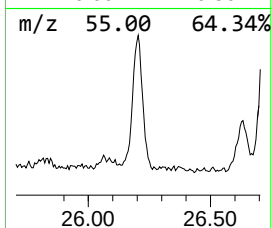
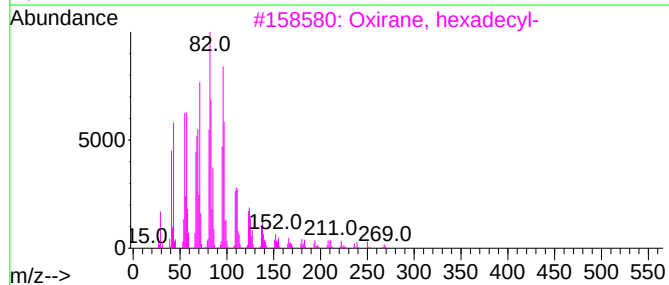
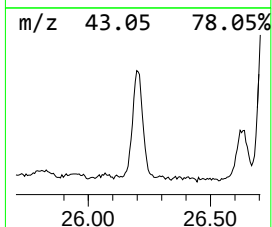
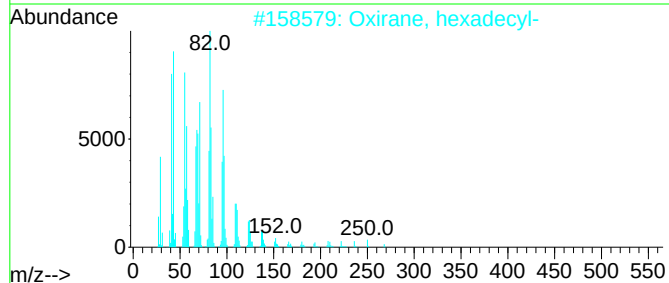
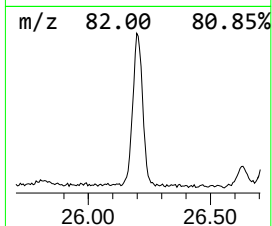
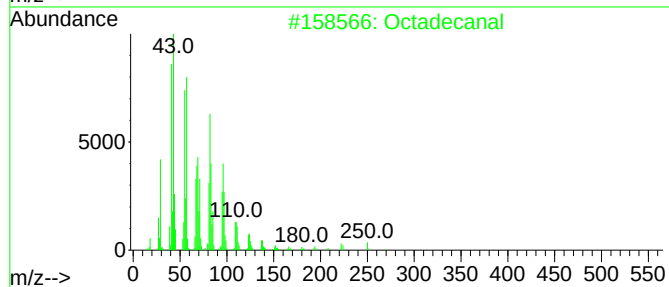
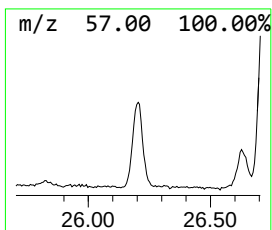
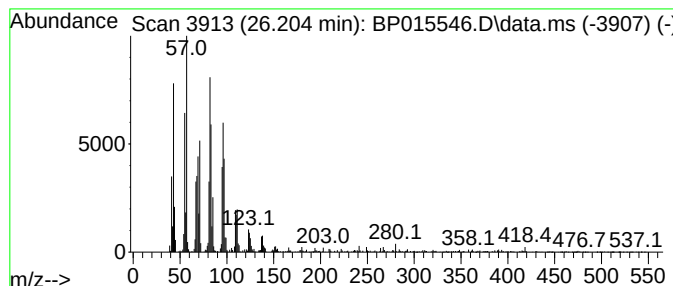
Instrument :  
BNA\_P  
ClientSampleId :  
DCFR1

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP060723.MA.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 21 Octadecanal Concentration Rank 7

| R.T.      | EstConc                   | Area    | Relative to ISTD |          | R.T.        |      |
|-----------|---------------------------|---------|------------------|----------|-------------|------|
| 26.204    | 20.02 ng/ul               | 1299840 | Perylene-d12     |          | 24.116      |      |
| Hit# of 5 | Tentative ID              |         | MW               | MolForm  | CAS#        | Qual |
| 1         | Octadecanal               |         | 268              | C18H36O  | 000638-66-4 | 94   |
| 2         | Oxirane, hexadecyl-       |         | 268              | C18H36O  | 007390-81-0 | 93   |
| 3         | Oxirane, hexadecyl-       |         | 268              | C18H36O  | 007390-81-0 | 91   |
| 4         | Dotriacontanal            |         | 464              | C32H64O  | 057878-00-9 | 91   |
| 5         | (Z)-14-Tricosenyl formate |         | 366              | C24H46O2 | 077899-10-6 | 89   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP061423\  
Data File : BP015546.D  
Acq On : 15 Jun 2023 07:45  
Operator : MA/JU  
Sample : 03088-05  
Misc :  
ALS Vial : 41 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
DCFR1

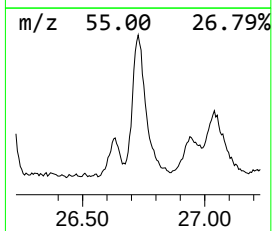
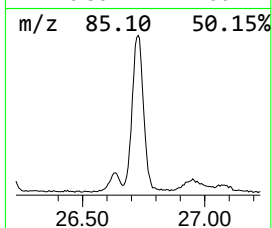
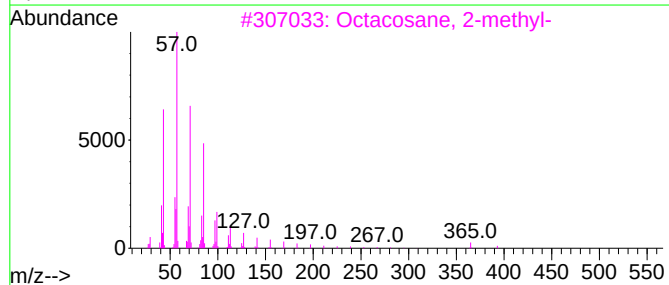
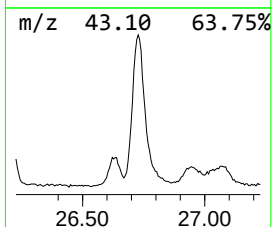
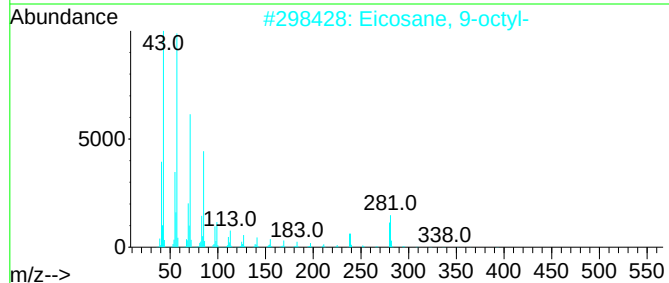
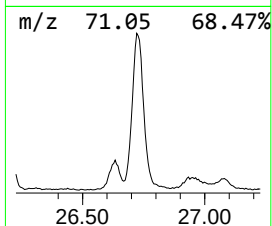
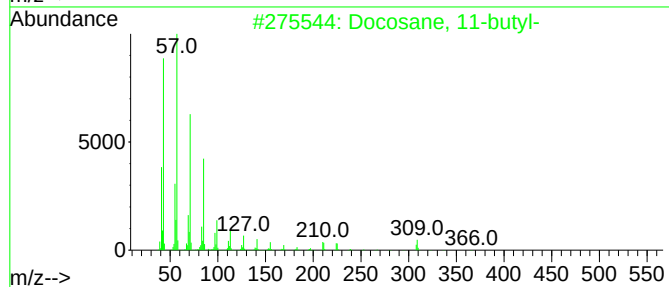
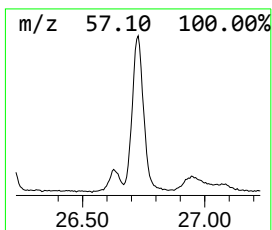
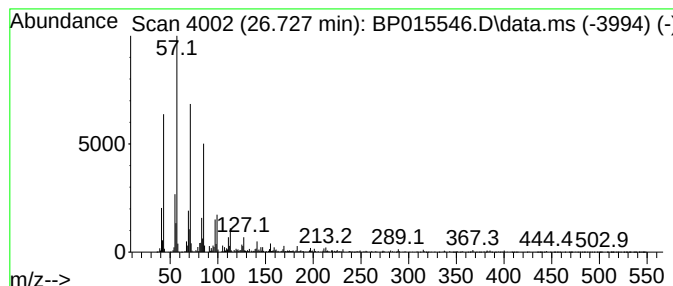
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Quant Title : SVOA CALIBRATION

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

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Peak Number 22 (DEL) Alkane: Straight-Chai... Concentration Rank 3

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 26.727 | 50.48 ng/ul | 3278060 | Perylene-d12     | 24.116 |

| Hit# | of | 5 | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------------|-----|---------|-------------|------|
| 1    |    |   | Docosane, 11-butyl-    | 366 | C26H54  | 013475-76-8 | 95   |
| 2    |    |   | Eicosane, 9-octyl-     | 394 | C28H58  | 013475-77-9 | 94   |
| 3    |    |   | Octacosane, 2-methyl-  | 408 | C29H60  | 001560-98-1 | 91   |
| 4    |    |   | Tetracosane, 11-decyl- | 479 | C34H70  | 055429-84-0 | 91   |
| 5    |    |   | Octacosane             | 394 | C28H58  | 000630-02-4 | 91   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP061423\  
Data File : BP015546.D  
Acq On : 15 Jun 2023 07:45  
Operator : MA/JU  
Sample : 03088-05  
Misc :  
ALS Vial : 41 Sample Multiplier: 1

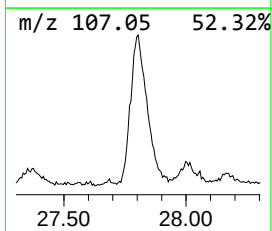
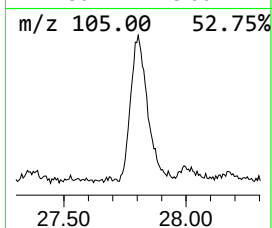
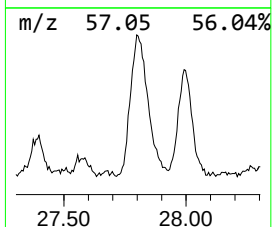
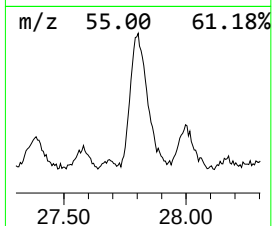
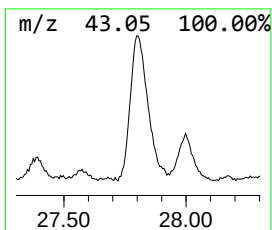
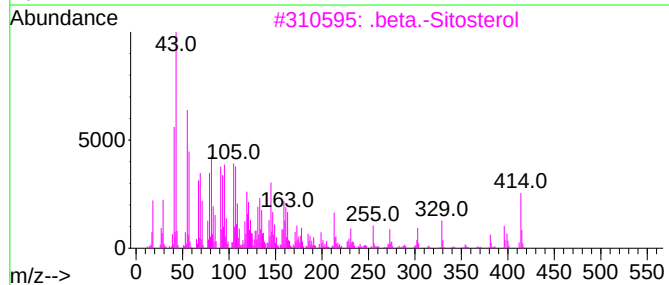
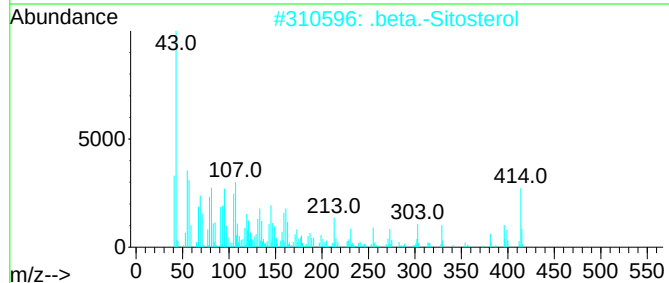
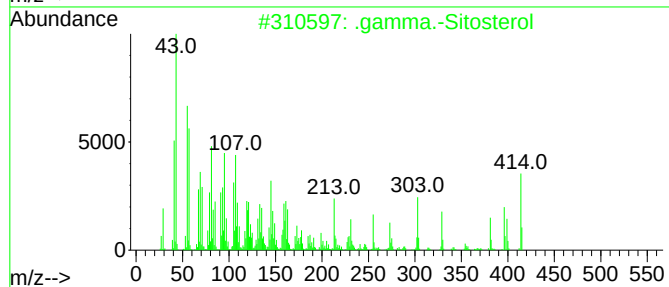
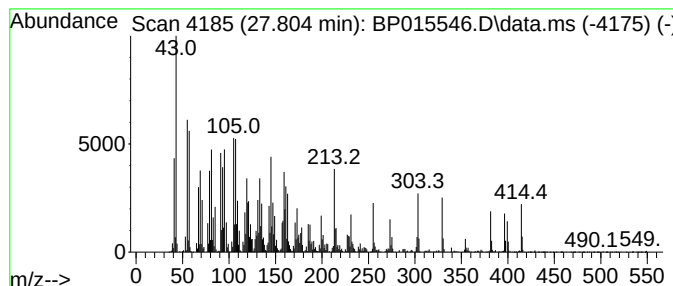
Instrument :  
BNA\_P  
ClientSampleId :  
DCFR1

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP060723.MA.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 23 .gamma.-Sitosterol Concentration Rank 2

| R.T.      | EstConc                             | Area    | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------------|---------|------------------|-------------|------|
| 27.804    | 62.89 ng/ul                         | 4084040 | Perylene-d12     | 24.116      |      |
| Hit# of 5 | Tentative ID                        | MW      | MolForm          | CAS#        | Qual |
| 1         | .gamma.-Sitosterol                  | 414     | C29H50O          | 000083-47-6 | 99   |
| 2         | .beta.-Sitosterol                   | 414     | C29H50O          | 000083-46-5 | 95   |
| 3         | .beta.-Sitosterol                   | 414     | C29H50O          | 000083-46-5 | 93   |
| 4         | Pregn-5-en-3-ol, 21-bromo-20-met... | 394     | C22H35BrO        | 055103-80-5 | 91   |
| 5         | (3S,8S,9S,10R,13R,14S,17R)-17-((... | 414     | C29H50O          | 029365-29-5 | 64   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP061423\  
Data File : BP015546.D  
Acq On : 15 Jun 2023 07:45  
Operator : MA/JU  
Sample : 03088-05  
Misc :  
ALS Vial : 41 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
DCFR1

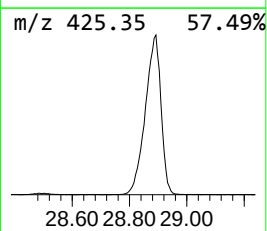
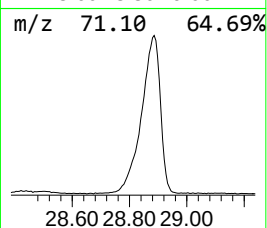
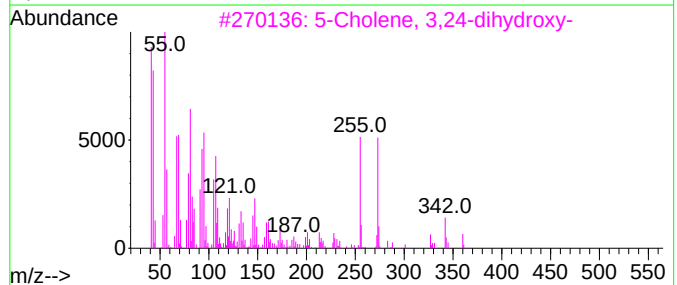
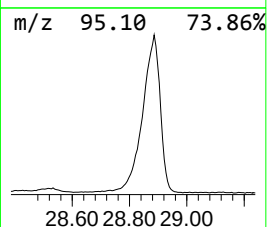
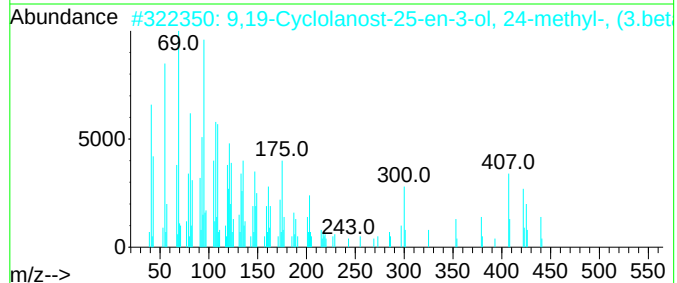
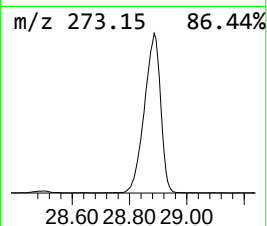
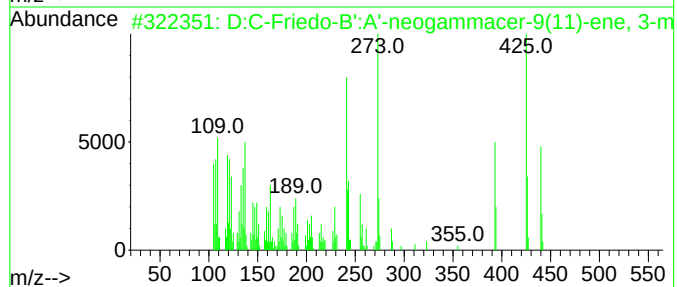
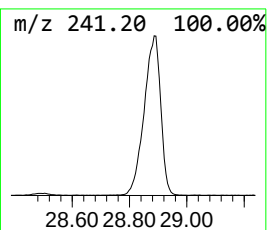
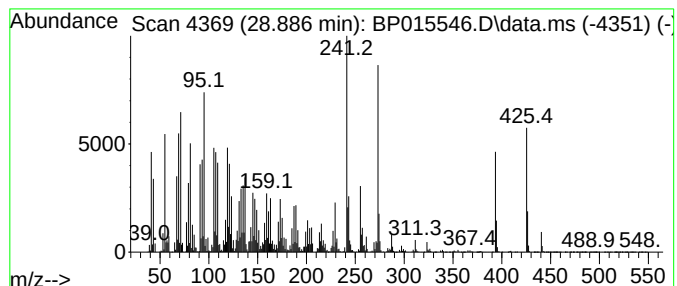
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Quant Title : SVOA CALIBRATION

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

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Peak Number 24 unknown-02 Concentration Rank 1

| R.T.   | EstConc      | Area     | Relative to ISTD | R.T.   |
|--------|--------------|----------|------------------|--------|
| 28.886 | 406.86 ng/ul | 26421600 | Perylene-d12     | 24.116 |

| Hit# | of | 5 | Tentative ID                         | MW  | MolForm  | CAS#         | Qual |
|------|----|---|--------------------------------------|-----|----------|--------------|------|
| 1    |    |   | D:C-Friedo-B':A'-neogammacer-9(1...  | 440 | C31H52O  | 004555-56-0  | 46   |
| 2    |    |   | 9,19-Cyclolanost-25-en-3-ol, 24-...  | 440 | C31H52O  | 000511-61-5  | 30   |
| 3    |    |   | 5-Cholene, 3,24-dihydroxy-           | 360 | C24H40O2 | 1000251-69-2 | 27   |
| 4    |    |   | 16,22-Epoxycooprostan-3.beta.,26,... | 434 | C27H46O4 | 122337-91-1  | 16   |
| 5    |    |   | Undec-10-ynoic acid, 2,7-dimethy...  | 316 | C21H32O2 | 1000406-96-7 | 14   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP061423\  
 Data File : BP015546.D  
 Acq On : 15 Jun 2023 07:45  
 Operator : MA/JU  
 Sample : 03088-05  
 Misc :  
 ALS Vial : 41 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 DCFR1

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP060723.MA.M  
 Quant Title : SVOA CALIBRATION

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

| TIC Top Hit name   | RT     | EstConc | Units | Response | --Internal Standard-- |        |         |      |
|--------------------|--------|---------|-------|----------|-----------------------|--------|---------|------|
|                    |        |         |       |          | #                     | RT     | Resp    | Conc |
| .alpha.-Pinene     | 7.511  | 2.6     | ng/ul | 68185    | 1                     | 8.087  | 515974  | 20.0 |
| Aceto phenone      | 8.922  | 12.3    | ng/ul | 317421   | 1                     | 8.087  | 515974  | 20.0 |
| Benzophenone       | 16.087 | 3.2     | ng/ul | 182923   | 3                     | 14.728 | 1143840 | 20.0 |
| 1,3-Benzenediam... | 18.281 | 9.7     | ng/ul | 689624   | 4                     | 17.486 | 1420800 | 20.0 |
| n-Hexadecanoic ... | 18.345 | 24.6    | ng/ul | 1744390  | 4                     | 17.486 | 1420800 | 20.0 |
| .gamma.-Sitoste... | 19.004 | 2.3     | ng/ul | 159669   | 4                     | 17.486 | 1420800 | 20.0 |
| 9,12-Octadecadi... | 19.428 | 2.4     | ng/ul | 169995   | 4                     | 17.486 | 1420800 | 20.0 |
| Octadecanoic acid  | 19.569 | 6.6     | ng/ul | 716524   | 5                     | 21.569 | 2161410 | 20.0 |
| Heptadecanoic acid | 20.616 | 2.8     | ng/ul | 302160   | 5                     | 21.569 | 2161410 | 20.0 |
| (DEL) Alkane: C... | 21.245 | 10.6    | ng/ul | 1140500  | 5                     | 21.569 | 2161410 | 20.0 |
| Kadsurin A         | 22.151 | 8.8     | ng/ul | 954622   | 5                     | 21.569 | 2161410 | 20.0 |
| Behenic alcohol    | 22.204 | 10.8    | ng/ul | 1166630  | 5                     | 21.569 | 2161410 | 20.0 |
| Piperenone         | 22.286 | 13.8    | ng/ul | 1489220  | 5                     | 21.569 | 2161410 | 20.0 |
| Dehydrodiisoeug... | 22.404 | 8.7     | ng/ul | 938873   | 5                     | 21.569 | 2161410 | 20.0 |
| Docosanoic acid    | 22.551 | 7.8     | ng/ul | 842478   | 5                     | 21.569 | 2161410 | 20.0 |
| unknown-01         | 22.663 | 5.3     | ng/ul | 569986   | 5                     | 21.569 | 2161410 | 20.0 |
| Hexacosanal        | 22.969 | 10.6    | ng/ul | 687453   | 6                     | 24.116 | 1298800 | 20.0 |
| Furan, 2,5-bis(... | 23.133 | 29.7    | ng/ul | 1926250  | 6                     | 24.116 | 1298800 | 20.0 |
| (DEL) Alkane: S... | 24.739 | 32.5    | ng/ul | 2108610  | 6                     | 24.116 | 1298800 | 20.0 |
| (DEL) Alkane: S... | 25.651 | 11.4    | ng/ul | 741119   | 6                     | 24.116 | 1298800 | 20.0 |
| Octadecanal        | 26.204 | 20.0    | ng/ul | 1299840  | 6                     | 24.116 | 1298800 | 20.0 |
| (DEL) Alkane: S... | 26.727 | 50.5    | ng/ul | 3278060  | 6                     | 24.116 | 1298800 | 20.0 |
| .gamma.-Sitosterol | 27.804 | 62.9    | ng/ul | 4084040  | 6                     | 24.116 | 1298800 | 20.0 |
| unknown-02         | 28.886 | 406.9   | ng/ul | 26421600 | 6                     | 24.116 | 1298800 | 20.0 |