

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM010224\  
 Data File : BM043693.D  
 Acq On : 02 Jan 2024 13:53  
 Operator : MA/JU  
 Sample : 05918-19  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 GCF32

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\_M\Methods\SFAM-EPA-BM121523.MA.M  
 Title : SVOA CALIBRATION

Signal : TIC: BM043693.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         | 3.369       | 28            | 31          | 39           | rVB      | 50724          | 70091         | 0.83%           | 0.075%        |
| 2         | 3.669       | 80            | 82          | 91           | rVB3     | 17606          | 34243         | 0.40%           | 0.037%        |
| 3         | 3.799       | 101           | 104         | 120          | rBV      | 412728         | 713179        | 8.41%           | 0.764%        |
| 4         | 4.346       | 193           | 197         | 203          | rVB      | 35203          | 46197         | 0.54%           | 0.049%        |
| 5         | 5.304       | 353           | 360         | 371          | rBV      | 93424          | 168298        | 1.98%           | 0.180%        |
| 6         | 6.463       | 550           | 557         | 567          | rVB      | 98692          | 172391        | 2.03%           | 0.185%        |
| 7         | 6.622       | 579           | 584         | 589          | rVB2     | 31545          | 50195         | 0.59%           | 0.054%        |
| 8         | 7.128       | 662           | 670         | 681          | rBV      | 776257         | 1254399       | 14.79%          | 1.344%        |
| 9         | 7.251       | 685           | 691         | 694          | rBV      | 154478         | 241838        | 2.85%           | 0.259%        |
| 10        | 7.298       | 694           | 699         | 710          | rVB      | 584175         | 925338        | 10.91%          | 0.991%        |
| 11        | 7.487       | 725           | 731         | 741          | rVB      | 741791         | 1197058       | 14.12%          | 1.283%        |
| 12        | 7.957       | 804           | 811         | 818          | rBV      | 754454         | 1227682       | 14.48%          | 1.315%        |
| 13        | 8.163       | 839           | 846         | 854          | rBV8     | 32058          | 81215         | 0.96%           | 0.087%        |
| 14        | 8.622       | 918           | 924         | 927          | rBV2     | 36639          | 56532         | 0.67%           | 0.061%        |
| 15        | 8.675       | 927           | 933         | 942          | rVB      | 822416         | 1341951       | 15.83%          | 1.438%        |
| 16        | 9.133       | 1004          | 1011        | 1021         | rBV      | 677181         | 1141706       | 13.47%          | 1.223%        |
| 17        | 9.716       | 1101          | 1110        | 1120         | rVB4     | 63002          | 130911        | 1.54%           | 0.140%        |
| 18        | 9.851       | 1126          | 1133        | 1142         | rBV      | 569665         | 955147        | 11.27%          | 1.023%        |
| 19        | 10.386      | 1216          | 1224        | 1237         | rBV      | 773465         | 1391497       | 16.41%          | 1.491%        |
| 20        | 10.769      | 1276          | 1289        | 1297         | rBV      | 933179         | 1638978       | 19.33%          | 1.756%        |
| 21        | 10.916      | 1306          | 1314        | 1328         | rBV      | 463004         | 889442        | 10.49%          | 0.953%        |
| 22        | 11.057      | 1333          | 1338        | 1346         | rVB8     | 19688          | 39978         | 0.47%           | 0.043%        |
| 23        | 11.139      | 1347          | 1352        | 1358         | rBV2     | 23424          | 42448         | 0.50%           | 0.045%        |
| 24        | 12.345      | 1550          | 1557        | 1564         | rBV7     | 22645          | 51651         | 0.61%           | 0.055%        |
| 25        | 13.098      | 1680          | 1685        | 1694         | rVB4     | 44512          | 88528         | 1.04%           | 0.095%        |
| 26        | 13.374      | 1727          | 1732        | 1737         | rVB      | 44352          | 65715         | 0.78%           | 0.070%        |
| 27        | 13.468      | 1742          | 1748        | 1750         | rVV6     | 39897          | 73557         | 0.87%           | 0.079%        |
| 28        | 13.492      | 1750          | 1752        | 1756         | rVV      | 39686          | 50784         | 0.60%           | 0.054%        |
| 29        | 13.539      | 1756          | 1760        | 1763         | rVV4     | 19017          | 33502         | 0.40%           | 0.036%        |
| 30        | 13.598      | 1765          | 1770        | 1775         | rVB2     | 144869         | 210275        | 2.48%           | 0.225%        |
| 31        | 13.727      | 1787          | 1792        | 1796         | rVV3     | 54671          | 92112         | 1.09%           | 0.099%        |
| 32        | 13.774      | 1796          | 1800        | 1807         | rVB3     | 91005          | 156470        | 1.85%           | 0.168%        |
| 33        | 13.857      | 1807          | 1814        | 1819         | rBV      | 1632461        | 2415776       | 28.49%          | 2.588%        |
| 34        | 13.904      | 1819          | 1822        | 1832         | rVB2     | 258695         | 405015        | 4.78%           | 0.434%        |
| 35        | 14.016      | 1832          | 1841        | 1847         | rBV      | 1237363        | 1776973       | 20.96%          | 1.904%        |
| 36        | 14.110      | 1852          | 1857        | 1865         | rVB      | 192004         | 290133        | 3.42%           | 0.311%        |

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 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

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 ClientSampleId :  
 GCF32

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\_M\Methods\SFAM-EPA-BM121523.MA.M  
 Title : SVOA CALIBRATION

|    |        |      |      |      |      |         |         |        |        |
|----|--------|------|------|------|------|---------|---------|--------|--------|
| 37 | 14.245 | 1875 | 1880 | 1882 | rBV5 | 20992   | 34054   | 0.40%  | 0.036% |
| 38 | 14.292 | 1882 | 1888 | 1895 | rVV  | 1431840 | 2227903 | 26.28% | 2.387% |
| 39 | 14.374 | 1898 | 1902 | 1903 | rVV4 | 28066   | 37831   | 0.45%  | 0.041% |
| 40 | 14.410 | 1903 | 1908 | 1914 | rVV  | 322891  | 485743  | 5.73%  | 0.520% |
| 41 | 14.480 | 1914 | 1920 | 1926 | rVB  | 567103  | 813146  | 9.59%  | 0.871% |
| 42 | 14.592 | 1933 | 1939 | 1946 | rVV  | 1276514 | 2079593 | 24.53% | 2.228% |
| 43 | 14.657 | 1946 | 1950 | 1956 | rVV2 | 132181  | 201726  | 2.38%  | 0.216% |
| 44 | 14.739 | 1960 | 1964 | 1969 | rVB2 | 113665  | 161328  | 1.90%  | 0.173% |
| 45 | 14.810 | 1969 | 1976 | 1991 | rBV2 | 522780  | 1238793 | 14.61% | 1.327% |
| 46 | 14.957 | 1996 | 2001 | 2006 | rBV4 | 74584   | 140302  | 1.65%  | 0.150% |
| 47 | 15.010 | 2007 | 2010 | 2013 | rVV2 | 44775   | 73480   | 0.87%  | 0.079% |
| 48 | 15.045 | 2013 | 2016 | 2019 | rVV  | 111193  | 173806  | 2.05%  | 0.186% |
| 49 | 15.074 | 2019 | 2021 | 2026 | rVV2 | 84451   | 104085  | 1.23%  | 0.112% |
| 50 | 15.145 | 2028 | 2033 | 2037 | rVB3 | 115787  | 169765  | 2.00%  | 0.182% |
| 51 | 15.586 | 2102 | 2108 | 2116 | rBV  | 1865569 | 2702620 | 31.88% | 2.896% |
| 52 | 15.710 | 2124 | 2129 | 2135 | rBV  | 435434  | 634525  | 7.48%  | 0.680% |
| 53 | 15.757 | 2135 | 2137 | 2141 | rVV4 | 34498   | 49202   | 0.58%  | 0.053% |
| 54 | 15.821 | 2144 | 2148 | 2152 | rVV5 | 53564   | 87907   | 1.04%  | 0.094% |
| 55 | 15.874 | 2152 | 2157 | 2163 | rVB  | 914187  | 1253981 | 14.79% | 1.344% |
| 56 | 16.045 | 2183 | 2186 | 2190 | rVB4 | 30136   | 37070   | 0.44%  | 0.040% |
| 57 | 16.215 | 2210 | 2215 | 2225 | rBV4 | 77536   | 179351  | 2.12%  | 0.192% |
| 58 | 16.315 | 2227 | 2232 | 2239 | rVB8 | 30154   | 75771   | 0.89%  | 0.081% |
| 59 | 16.527 | 2264 | 2268 | 2272 | rBV4 | 38041   | 62814   | 0.74%  | 0.067% |
| 60 | 16.739 | 2300 | 2304 | 2313 | rBV2 | 152951  | 263461  | 3.11%  | 0.282% |
| 61 | 16.821 | 2313 | 2318 | 2320 | rBV  | 146083  | 201847  | 2.38%  | 0.216% |
| 62 | 16.986 | 2341 | 2346 | 2356 | rBV  | 561821  | 829248  | 9.78%  | 0.889% |
| 63 | 17.086 | 2358 | 2363 | 2373 | rVV  | 492315  | 737270  | 8.70%  | 0.790% |
| 64 | 17.239 | 2385 | 2389 | 2396 | rBV2 | 119536  | 197461  | 2.33%  | 0.212% |
| 65 | 17.304 | 2397 | 2400 | 2401 | rVV  | 71348   | 78988   | 0.93%  | 0.085% |
| 66 | 17.339 | 2401 | 2406 | 2411 | rVV  | 1490973 | 2219226 | 26.17% | 2.378% |
| 67 | 17.380 | 2411 | 2413 | 2417 | rVV  | 111724  | 150101  | 1.77%  | 0.161% |
| 68 | 17.439 | 2417 | 2423 | 2432 | rVB2 | 1930169 | 2760016 | 32.55% | 2.957% |
| 69 | 17.568 | 2441 | 2445 | 2452 | rVB4 | 113990  | 236269  | 2.79%  | 0.253% |
| 70 | 17.633 | 2453 | 2456 | 2464 | rVB6 | 39926   | 66047   | 0.78%  | 0.071% |
| 71 | 18.121 | 2534 | 2539 | 2545 | rBV2 | 144872  | 264967  | 3.13%  | 0.284% |
| 72 | 18.186 | 2545 | 2550 | 2554 | rBV  | 702637  | 1034362 | 12.20% | 1.108% |
| 73 | 18.245 | 2555 | 2560 | 2576 | rVB  | 1891583 | 2910954 | 34.33% | 3.119% |
| 74 | 18.539 | 2606 | 2610 | 2614 | rBV2 | 106117  | 191763  | 2.26%  | 0.205% |
| 75 | 19.203 | 2719 | 2723 | 2727 | rVB  | 272651  | 327639  | 3.86%  | 0.351% |
| 76 | 19.256 | 2729 | 2732 | 2736 | rVB  | 93614   | 114200  | 1.35%  | 0.122% |

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Instrument :  
 BNA\_M  
 ClientSampleId :  
 GCF32

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\_M\Methods\SFAM-EPA-BM121523.MA.M  
 Title : SVOA CALIBRATION

|     |        |      |      |      |      |         |         |         |        |
|-----|--------|------|------|------|------|---------|---------|---------|--------|
| 77  | 19.374 | 2749 | 2752 | 2754 | rBV2 | 91832   | 108980  | 1.29%   | 0.117% |
| 78  | 19.403 | 2754 | 2757 | 2759 | rBV  | 200194  | 257617  | 3.04%   | 0.276% |
| 79  | 19.521 | 2773 | 2777 | 2784 | rBV  | 521036  | 795008  | 9.38%   | 0.852% |
| 80  | 19.598 | 2786 | 2790 | 2796 | rVB  | 1352384 | 1695593 | 20.00%  | 1.817% |
| 81  | 19.727 | 2807 | 2812 | 2819 | rBV  | 2210595 | 3118360 | 36.78%  | 3.341% |
| 82  | 19.903 | 2839 | 2842 | 2847 | rBV2 | 143442  | 190121  | 2.24%   | 0.204% |
| 83  | 20.250 | 2897 | 2901 | 2910 | rBV4 | 227482  | 564896  | 6.66%   | 0.605% |
| 84  | 20.439 | 2929 | 2933 | 2941 | rVB  | 874347  | 1183271 | 13.96%  | 1.268% |
| 85  | 20.521 | 2945 | 2947 | 2952 | rVB4 | 209999  | 271643  | 3.20%   | 0.291% |
| 86  | 20.580 | 2954 | 2957 | 2958 | rBV  | 174455  | 197336  | 2.33%   | 0.211% |
| 87  | 20.609 | 2958 | 2962 | 2966 | rVV  | 6527622 | 8478590 | 100.00% | 9.085% |
| 88  | 20.650 | 2966 | 2969 | 2985 | rVB2 | 1643987 | 2831530 | 33.40%  | 3.034% |
| 89  | 21.115 | 3045 | 3048 | 3057 | rVB3 | 477704  | 689860  | 8.14%   | 0.739% |
| 90  | 21.239 | 3065 | 3069 | 3073 | rBV  | 1140388 | 1319742 | 15.57%  | 1.414% |
| 91  | 21.339 | 3076 | 3086 | 3097 | rVV5 | 1472826 | 5457438 | 64.37%  | 5.848% |
| 92  | 21.521 | 3113 | 3117 | 3133 | rVB2 | 1754825 | 2655529 | 31.32%  | 2.845% |
| 93  | 22.103 | 3214 | 3216 | 3220 | rVB  | 365896  | 409187  | 4.83%   | 0.438% |
| 94  | 23.227 | 3404 | 3407 | 3411 | rVV  | 402618  | 550853  | 6.50%   | 0.590% |
| 95  | 23.274 | 3411 | 3415 | 3424 | rVB  | 794404  | 1346815 | 15.88%  | 1.443% |
| 96  | 23.780 | 3495 | 3501 | 3508 | rVB2 | 1554759 | 3021824 | 35.64%  | 3.238% |
| 97  | 23.938 | 3522 | 3528 | 3536 | rVB  | 1249627 | 2407165 | 28.39%  | 2.579% |
| 98  | 24.609 | 3636 | 3642 | 3649 | rVB  | 1207053 | 2586969 | 30.51%  | 2.772% |
| 99  | 24.697 | 3651 | 3657 | 3669 | rVB  | 1969453 | 4198835 | 49.52%  | 4.499% |
| 100 | 27.403 | 4110 | 4117 | 4131 | rVB2 | 1152912 | 3860621 | 45.53%  | 4.137% |

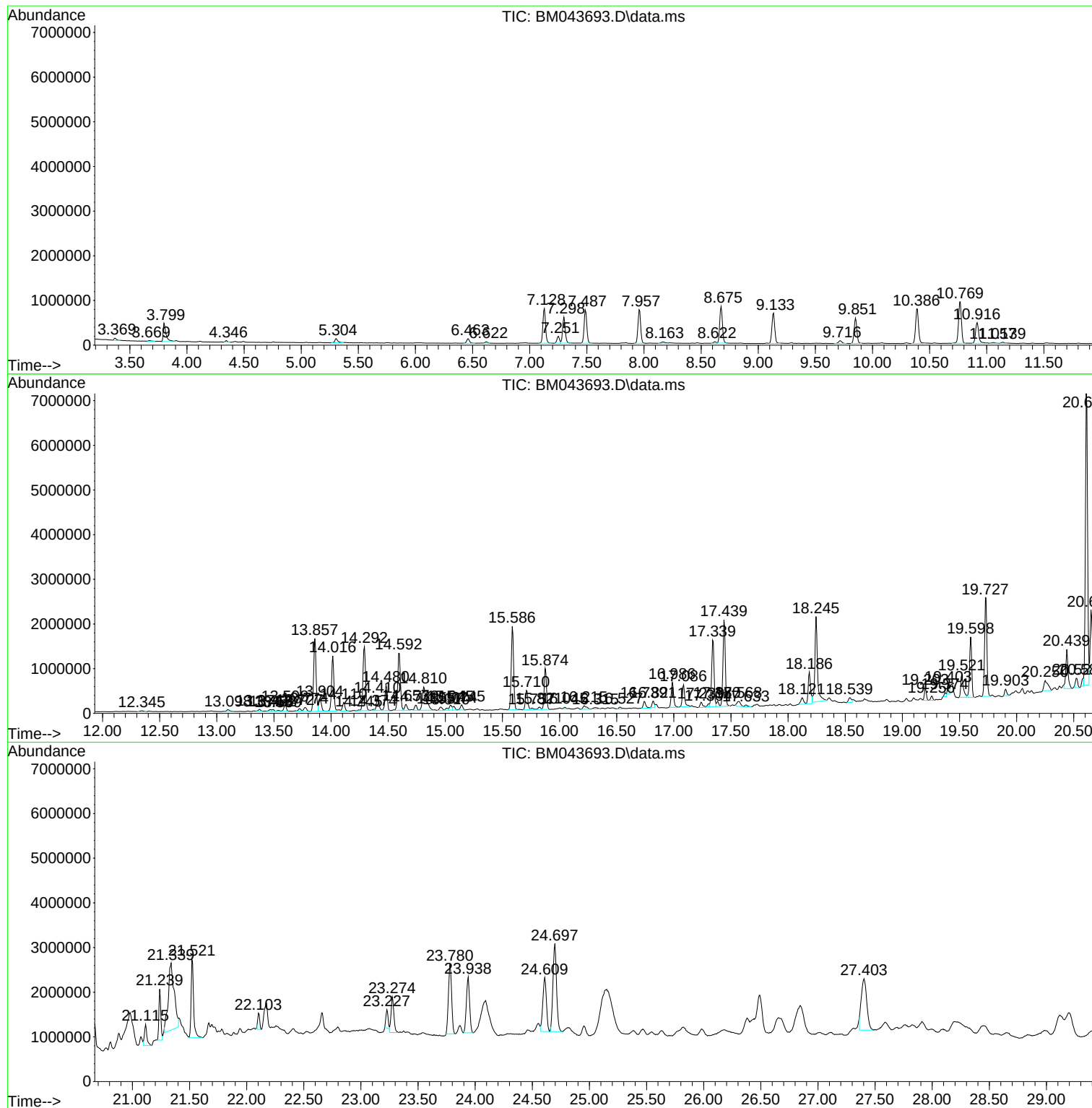
Sum of corrected areas: 93327603

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Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM121523.MA.M  
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM010224\  
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Instrument :  
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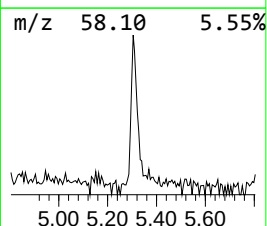
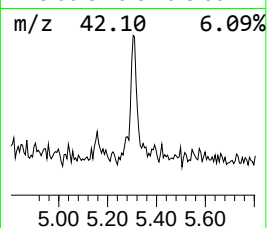
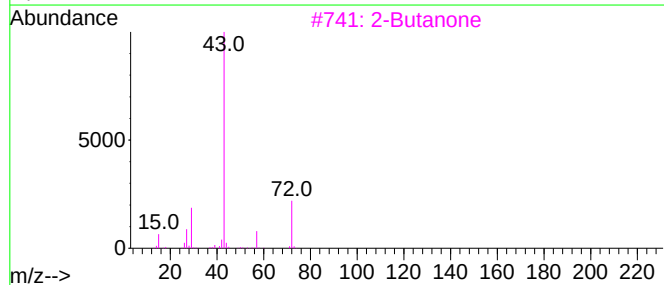
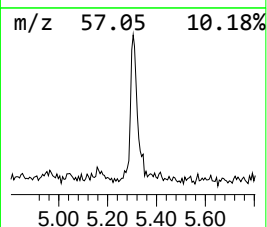
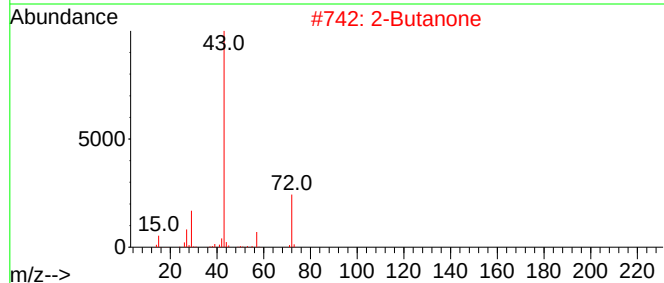
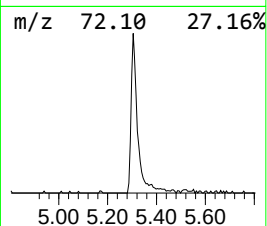
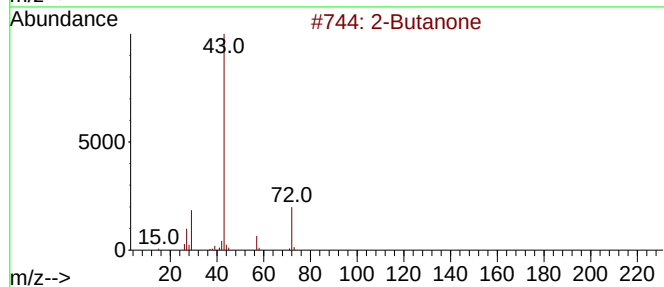
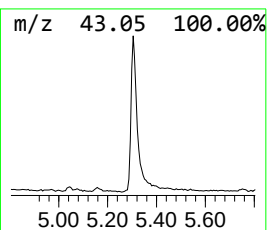
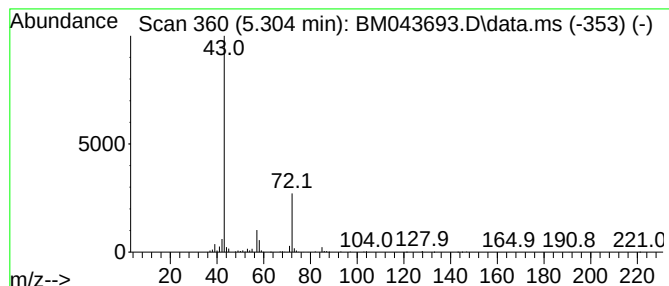
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM121523.MA.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 1 2-Butanone Concentration Rank 27

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 5.304 | 2.74 ng/ul | 168298 | 1,4-Dichlorobenzene-d4 | 7.957 |

| Hit# | of                     | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|------------------------|---|--------------|-----|---------|-------------|------|
| 1    | 2-Butanone             |   |              | 72  | C4H8O   | 000078-93-3 | 72   |
| 2    | 2-Butanone             |   |              | 72  | C4H8O   | 000078-93-3 | 64   |
| 3    | 2-Butanone             |   |              | 72  | C4H8O   | 000078-93-3 | 64   |
| 4    | 2-Butanone             |   |              | 72  | C4H8O   | 000078-93-3 | 59   |
| 5    | 2-Heptanone, 3-methyl- |   |              | 128 | C8H16O  | 002371-19-9 | 50   |



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Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM121523.MA.M  
Quant Title : SVOA CALIBRATION

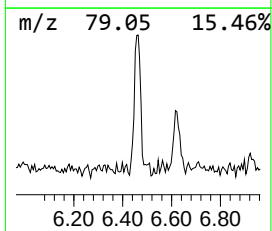
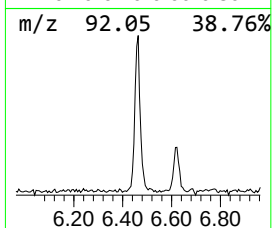
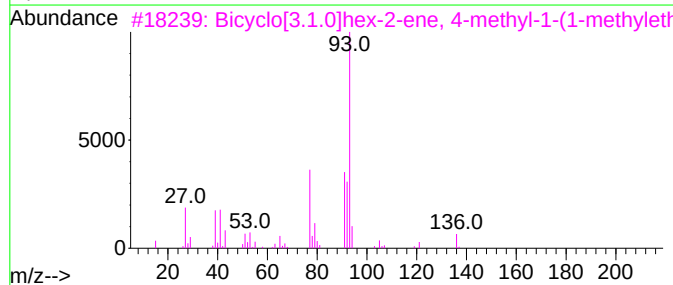
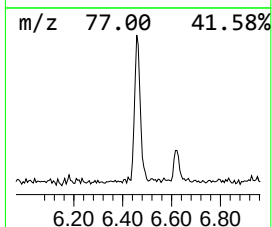
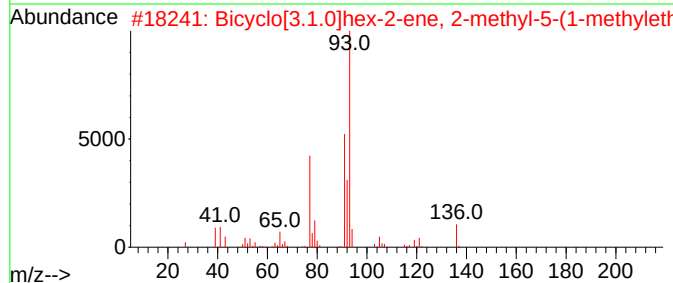
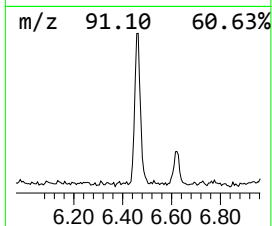
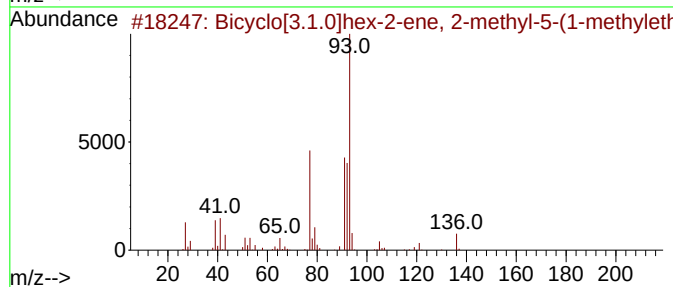
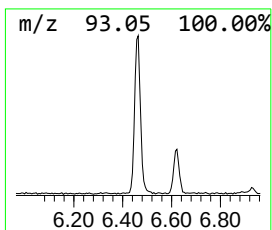
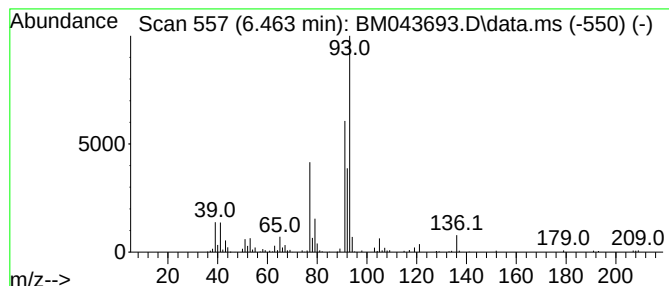
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 2 Bicyclo[3.1.0]hex-2-ene, 2-... Concentration Rank 25

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 6.463 | 2.81 ng/ul | 172391 | 1,4-Dichlorobenzene-d4 | 7.957 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Bicyclo[3.1.0]hex-2-ene, 2-methy... | 136 | C10H16  | 002867-05-2 | 91   |
| 2    |    |   | Bicyclo[3.1.0]hex-2-ene, 2-methy... | 136 | C10H16  | 002867-05-2 | 91   |
| 3    |    |   | Bicyclo[3.1.0]hex-2-ene, 4-methy... | 136 | C10H16  | 028634-89-1 | 90   |
| 4    |    |   | .alpha.-Phellandrene                | 136 | C10H16  | 000099-83-2 | 90   |
| 5    |    |   | Bicyclo[3.1.0]hex-2-ene, 2-methy... | 136 | C10H16  | 002867-05-2 | 86   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM010224\  
Data File : BM043693.D  
Acq On : 02 Jan 2024 13:53  
Operator : MA/JU  
Sample : 05918-19  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
GCF32

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM121523.MA.M  
Quant Title : SVOA CALIBRATION

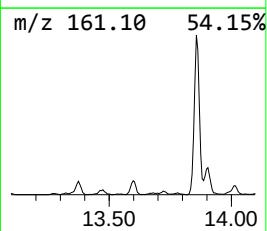
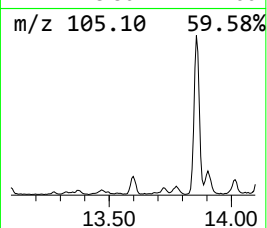
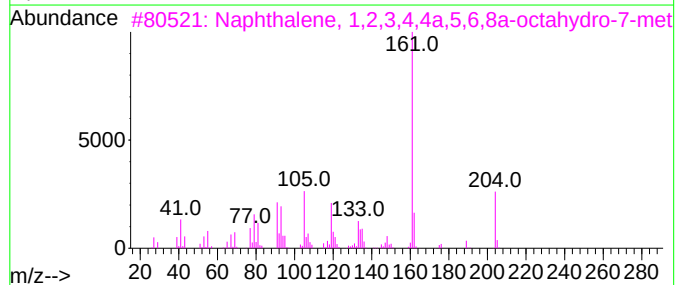
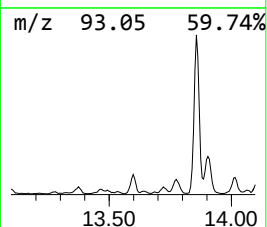
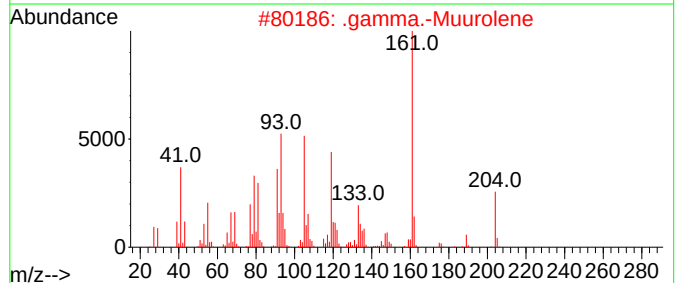
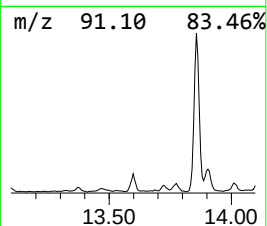
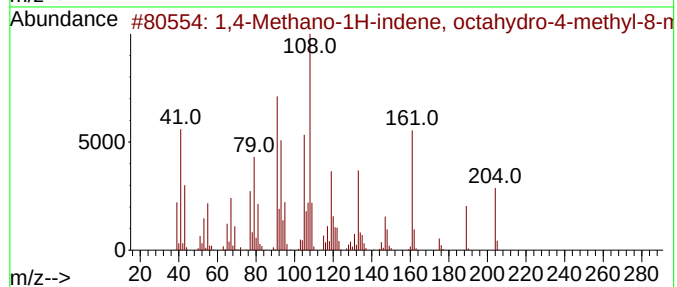
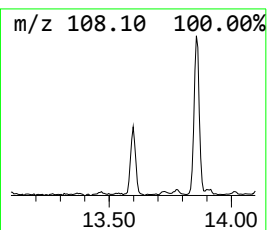
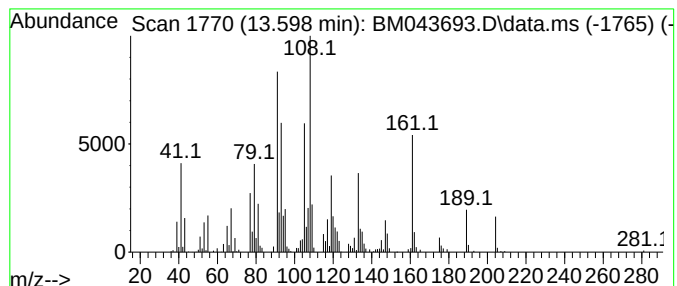
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TIC Integration Parameters: LSCINT.P

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Peak Number 3 1,4-Methano-1H-indene, octa... Concentration Rank 33

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 13.598 | 2.02 ng/ul | 210275 | Acenaphthene-d10 | 14.592 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | 1,4-Methano-1H-indene, octahydro... | 204 | C15H24  | 003650-28-0 | 99   |
| 2    |    |   | .gamma.-Muurolene                   | 204 | C15H24  | 030021-74-0 | 95   |
| 3    |    |   | Naphthalene, 1,2,3,4,4a,5,6,8a-o... | 204 | C15H24  | 039029-41-9 | 95   |
| 4    |    |   | Naphthalene, 1,2,4a,5,6,8a-hexah... | 204 | C15H24  | 000483-75-0 | 94   |
| 5    |    |   | 1,4-Methano-1H-indene, octahydro... | 204 | C15H24  | 003650-28-0 | 93   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM010224\  
Data File : BM043693.D  
Acq On : 02 Jan 2024 13:53  
Operator : MA/JU  
Sample : 05918-19  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
GCF32

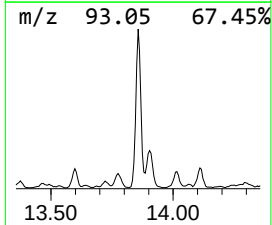
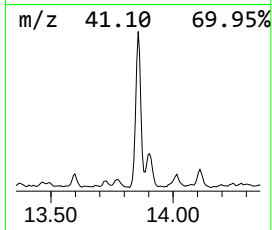
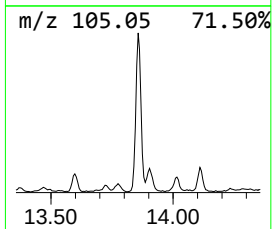
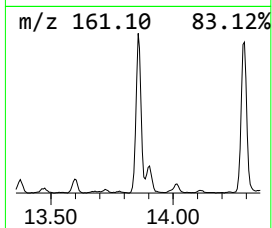
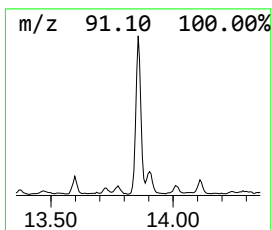
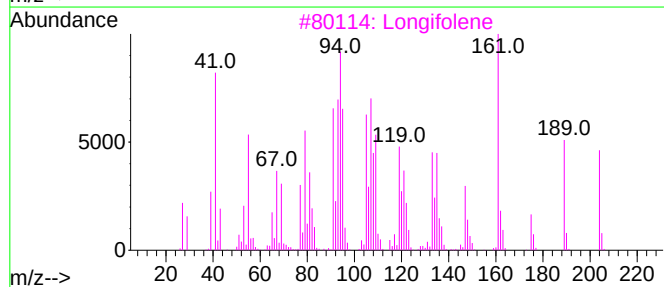
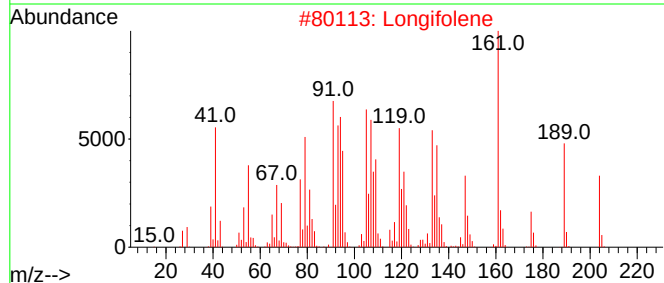
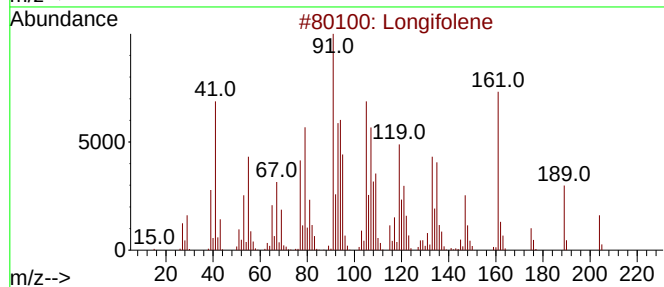
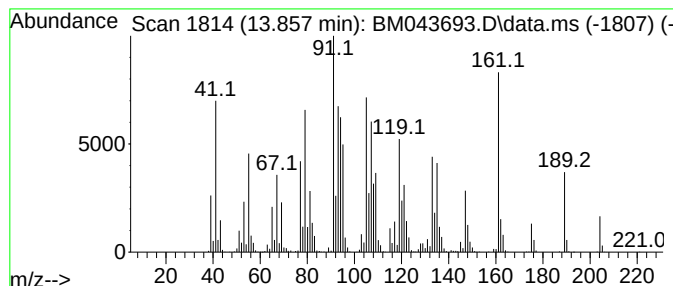
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Quant Title : SVOA CALIBRATION

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

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Peak Number 4 Longifolene Concentration Rank 6

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 13.857 | 23.23 ng/ul | 2415780 | Acenaphthene-d10 | 14.592 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Longifolene                         | 204 | C15H24  | 000475-20-7 | 99   |
| 2    |    |   | Longifolene                         | 204 | C15H24  | 000475-20-7 | 99   |
| 3    |    |   | Longifolene                         | 204 | C15H24  | 000475-20-7 | 98   |
| 4    |    |   | (3R,4aS,5R)-4a,5-Dimethyl-3-(pro... | 204 | C15H24  | 024741-64-8 | 96   |
| 5    |    |   | Longifolene                         | 204 | C15H24  | 000475-20-7 | 94   |





Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM010224\  
Data File : BM043693.D  
Acq On : 02 Jan 2024 13:53  
Operator : MA/JU  
Sample : 05918-19  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
GCF32

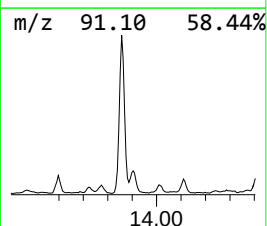
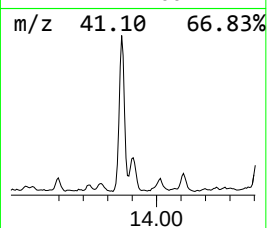
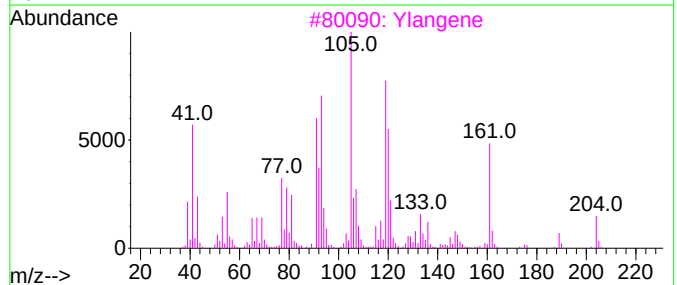
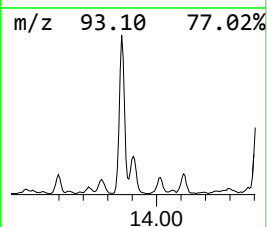
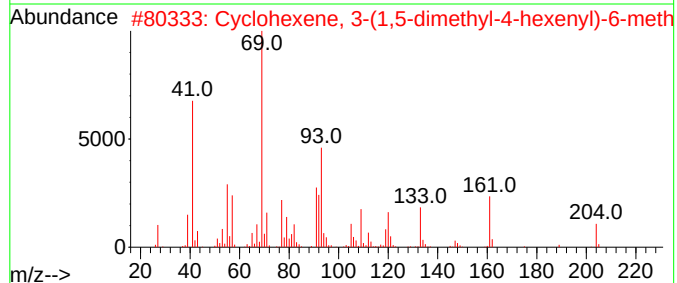
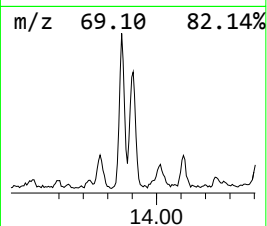
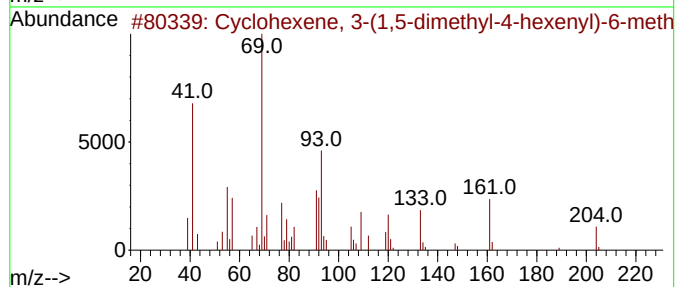
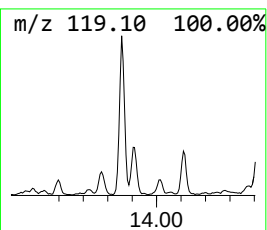
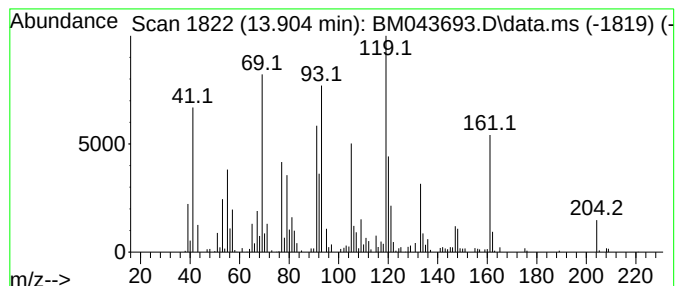
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Quant Title : SVOA CALIBRATION

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

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Peak Number 5 Cyclohexene, 3-(1,5-dimethy... Concentration Rank 22

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------------|--------|------------------|-------------|------|
| 13.904    | 3.90 ng/ul                          | 405015 | Acenaphthene-d10 | 14.592      |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#        | Qual |
| 1         | Cyclohexene, 3-(1,5-dimethyl-4-h... | 204    | C15H24           | 020307-83-9 | 62   |
| 2         | Cyclohexene, 3-(1,5-dimethyl-4-h... | 204    | C15H24           | 020307-83-9 | 62   |
| 3         | Ylangene                            | 204    | C15H24           | 014912-44-8 | 52   |
| 4         | (1S,5S)-4-Methylene-1-((R)-6-met... | 204    | C15H24           | 058319-04-3 | 49   |
| 5         | (1R,4aR,8aR)-2,5,5,8a-Tetramethy... | 204    | C15H24           | 079562-96-2 | 42   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM010224\  
Data File : BM043693.D  
Acq On : 02 Jan 2024 13:53  
Operator : MA/JU  
Sample : 05918-19  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
GCF32

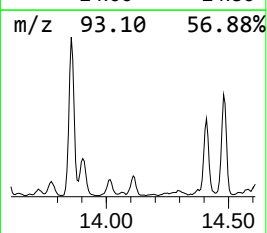
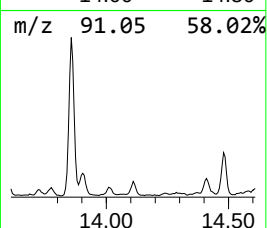
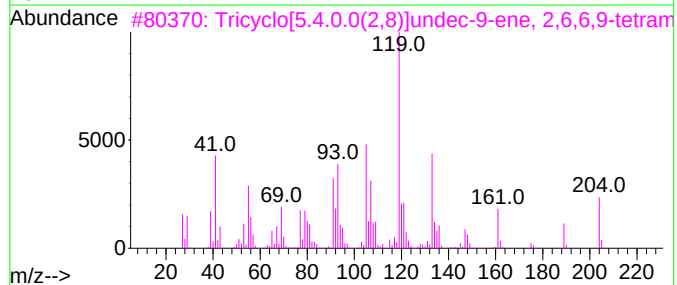
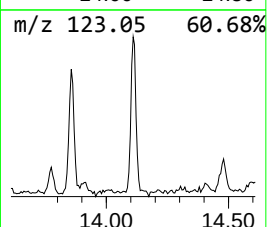
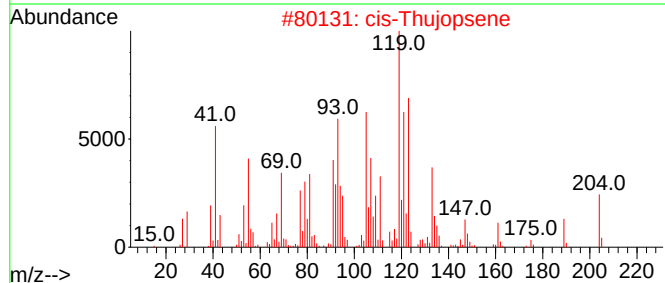
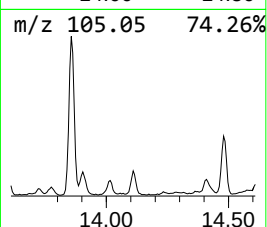
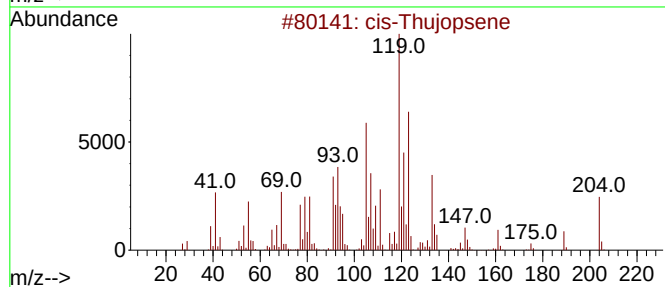
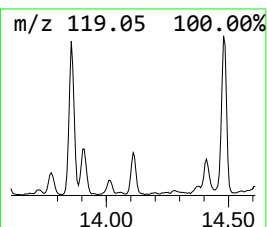
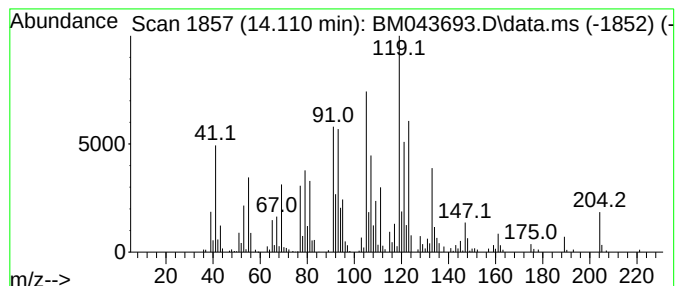
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Quant Title : SVOA CALIBRATION

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

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Peak Number 6 cis-Thujopsene Concentration Rank 26

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 14.110 | 2.79 ng/ul | 290133 | Acenaphthene-d10 | 14.592 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | cis-Thujopsene                      | 204 | C15H24  | 000470-40-6 | 99   |
| 2    |    |   | cis-Thujopsene                      | 204 | C15H24  | 000470-40-6 | 94   |
| 3    |    |   | Tricyclo[5.4.0.0(2,8)]undec-9-en... | 204 | C15H24  | 005989-08-2 | 89   |
| 4    |    |   | cis-Thujopsene                      | 204 | C15H24  | 000470-40-6 | 87   |
| 5    |    |   | 1H-Benzocycloheptene, 2,4a,5,6,7... | 204 | C15H24  | 001461-03-6 | 64   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM010224\  
Data File : BM043693.D  
Acq On : 02 Jan 2024 13:53  
Operator : MA/JU  
Sample : 05918-19  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
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Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM121523.MA.M  
Quant Title : SVOA CALIBRATION

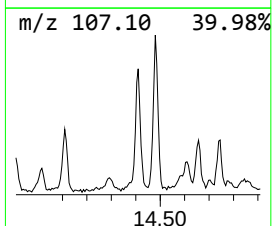
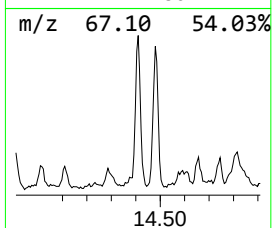
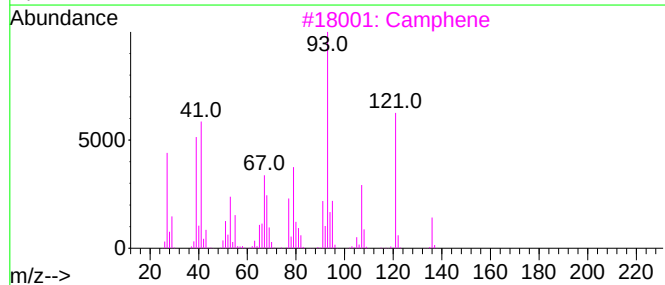
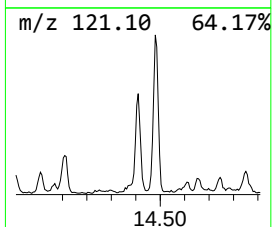
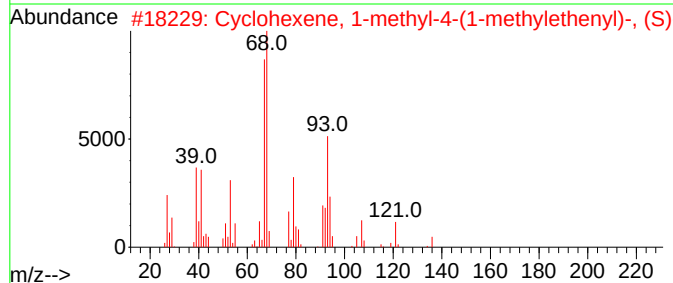
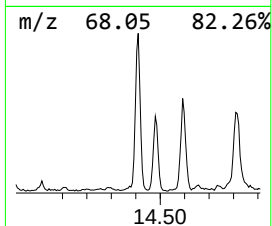
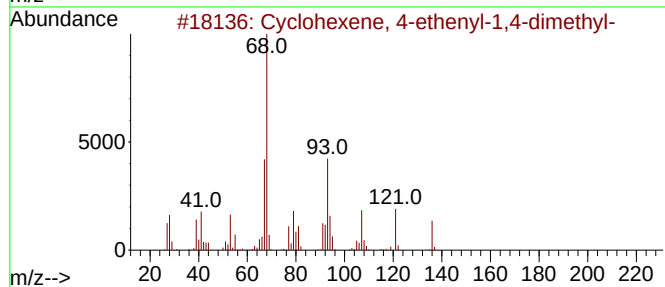
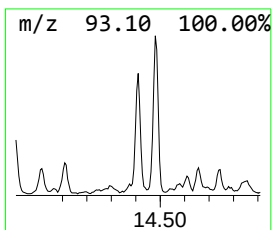
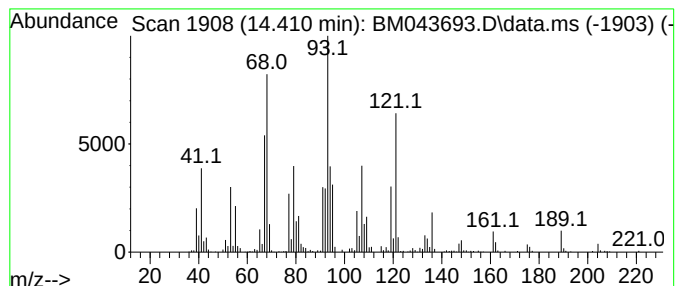
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TIC Integration Parameters: LSCINT.P

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Peak Number 7 Cyclohexene, 4-ethenyl-1,4-... Concentration Rank 19

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 14.410 | 4.67 ng/ul | 485743 | Acenaphthene-d10 | 14.592 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Cyclohexene, 4-ethenyl-1,4-dimet... | 136 | C10H16  | 001743-61-9 | 81   |
| 2    |    |   | Cyclohexene, 1-methyl-4-(1-methy... | 136 | C10H16  | 005989-54-8 | 74   |
| 3    |    |   | Camphene                            | 136 | C10H16  | 000079-92-5 | 60   |
| 4    |    |   | Camphene                            | 136 | C10H16  | 000079-92-5 | 55   |
| 5    |    |   | Limonene                            | 136 | C10H16  | 000138-86-3 | 53   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM010224\  
Data File : BM043693.D  
Acq On : 02 Jan 2024 13:53  
Operator : MA/JU  
Sample : 05918-19  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
GCF32

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM121523.MA.M  
Quant Title : SVOA CALIBRATION

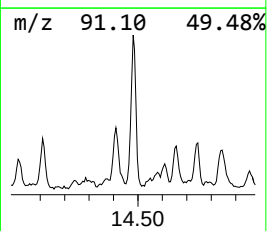
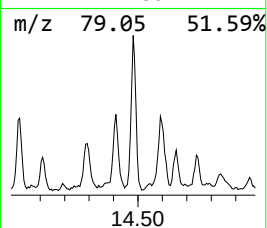
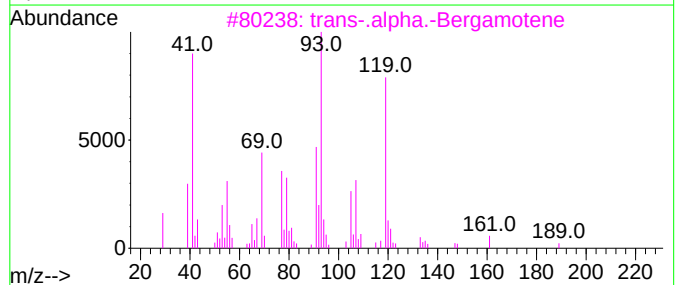
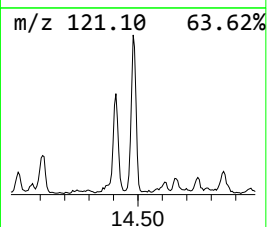
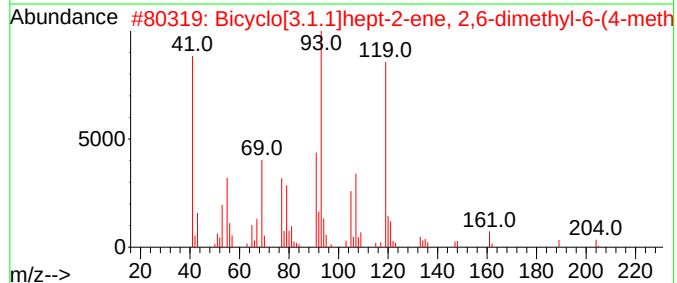
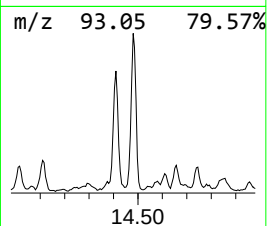
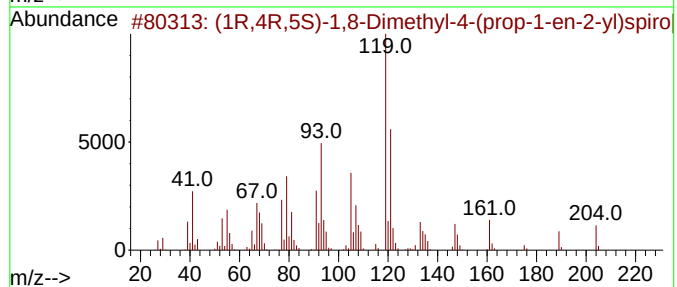
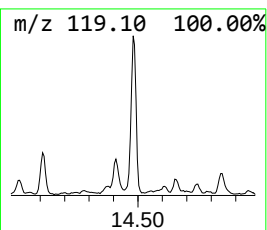
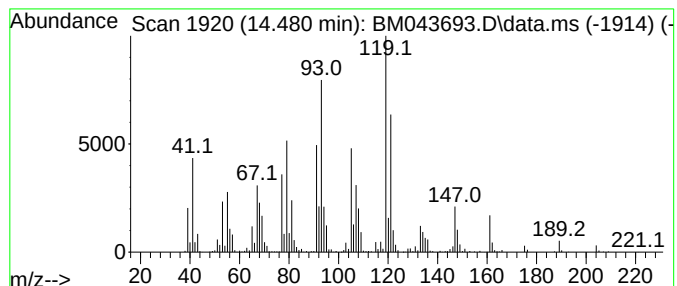
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TIC Integration Parameters: LSCINT.P

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Peak Number 8 (1R,4R,5S)-1,8-Dimethyl-4-(... Concentration Rank 15

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 14.480 | 7.82 ng/ul | 813146 | Acenaphthene-d10 | 14.592 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | (1R,4R,5S)-1,8-Dimethyl-4-(prop-... | 204 | C15H24  | 729602-94-2 | 95   |
| 2    |    |   | Bicyclo[3.1.1]hept-2-ene, 2,6-di... | 204 | C15H24  | 017699-05-7 | 70   |
| 3    |    |   | trans-.alpha.-Bergamotene           | 204 | C15H24  | 013474-59-4 | 64   |
| 4    |    |   | cis-.alpha.-Bergamotene             | 204 | C15H24  | 018252-46-5 | 64   |
| 5    |    |   | Bicyclo[2.2.1]heptane, 2,2-dimet... | 136 | C10H16  | 005794-03-6 | 49   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM010224\  
Data File : BM043693.D  
Acq On : 02 Jan 2024 13:53  
Operator : MA/JU  
Sample : 05918-19  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
GCF32

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Quant Title : SVOA CALIBRATION

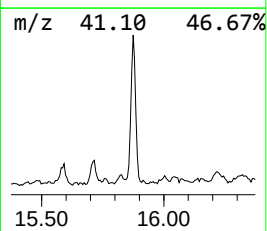
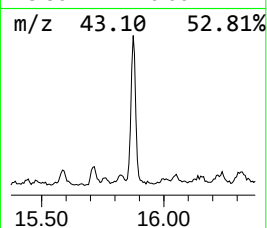
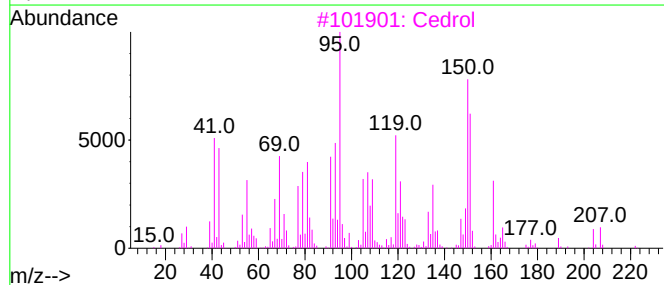
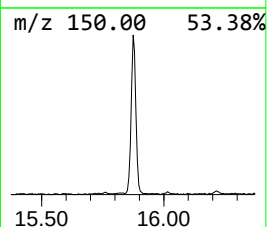
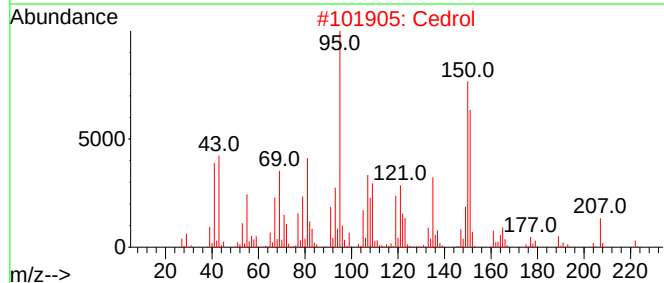
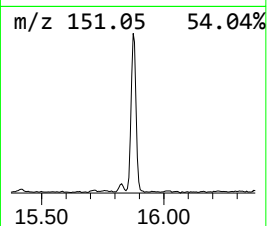
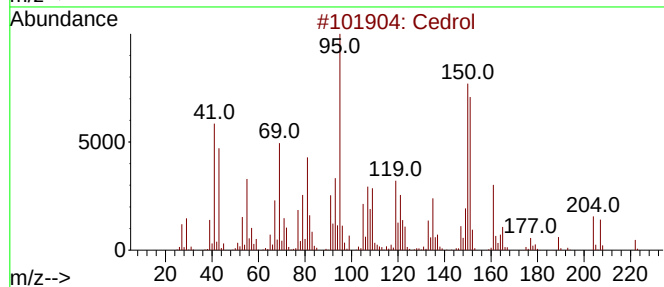
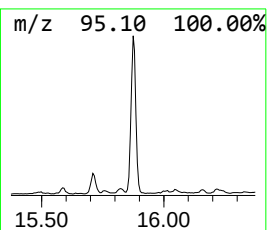
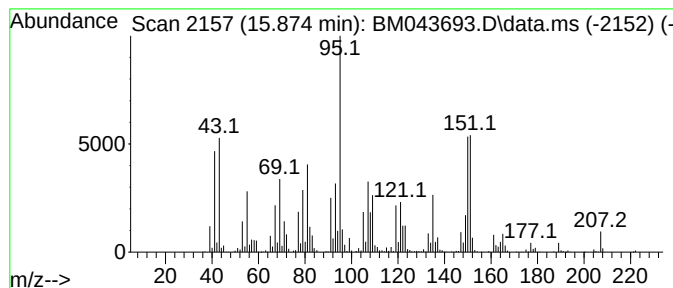
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TIC Integration Parameters: LSCINT.P

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Peak Number 9 Cedrol Concentration Rank 10

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 15.874 | 12.06 ng/ul | 1253980 | Acenaphthene-d10 | 14.592 |

| Hit# | of                                  | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|-------------------------------------|---|--------------|-----|---------|-------------|------|
| 1    | Cedrol                              |   |              | 222 | C15H26O | 000077-53-2 | 93   |
| 2    | Cedrol                              |   |              | 222 | C15H26O | 000077-53-2 | 91   |
| 3    | Cedrol                              |   |              | 222 | C15H26O | 000077-53-2 | 89   |
| 4    | Cedrol                              |   |              | 222 | C15H26O | 000077-53-2 | 87   |
| 5    | (3R,3aS,6S,7R)-3,6,8,8-Tetrameth... |   |              | 222 | C15H26O | 019903-73-2 | 83   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM010224\  
Data File : BM043693.D  
Acq On : 02 Jan 2024 13:53  
Operator : MA/JU  
Sample : 05918-19  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
GCF32

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM121523.MA.M  
Quant Title : SVOA CALIBRATION

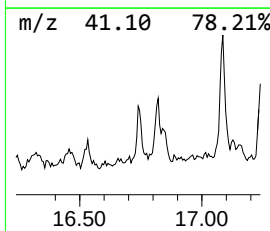
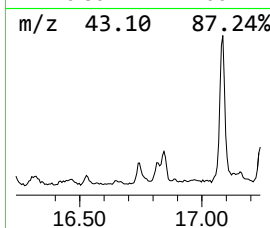
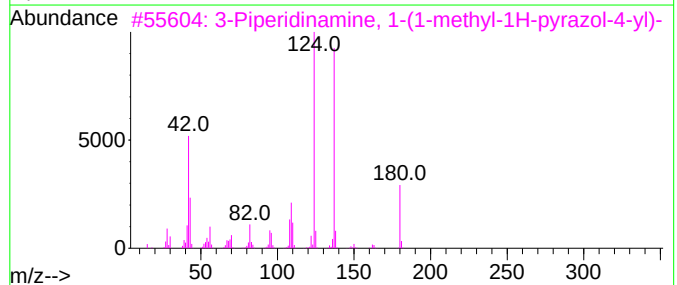
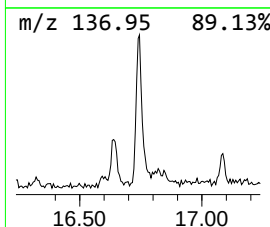
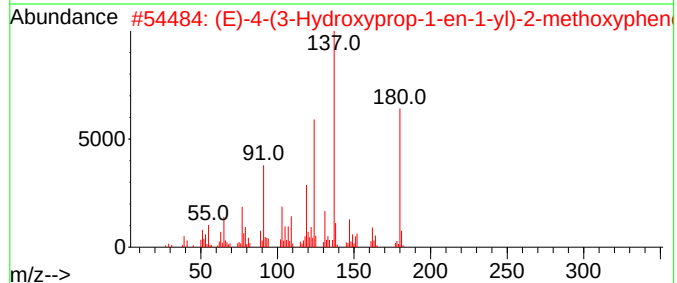
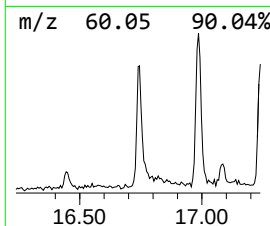
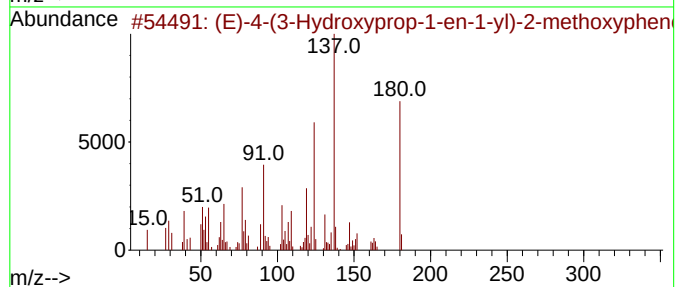
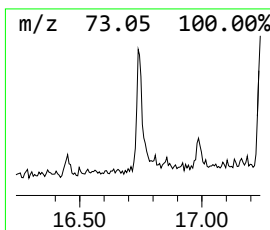
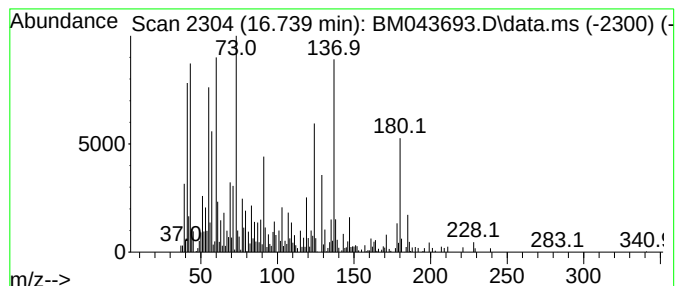
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TIC Integration Parameters: LSCINT.P

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Peak Number 10 (E)-4-(3-Hydroxyprop-1-en-1-yl) Concentration Rank 29

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 16.739 | 2.37 ng/ul | 263461 | Phenanthrene-d10 | 17.339 |

| Hit# | of                                  | 5   | Tentative ID | MW           | MolForm | CAS# | Qual |
|------|-------------------------------------|-----|--------------|--------------|---------|------|------|
| 1    | (E)-4-(3-Hydroxyprop-1-en-1-yl)-... | 180 | C10H12O3     | 032811-40-8  | 95      |      |      |
| 2    | (E)-4-(3-Hydroxyprop-1-en-1-yl)-... | 180 | C10H12O3     | 032811-40-8  | 49      |      |      |
| 3    | 3-Piperidinamine, 1-(1-methyl-1H... | 180 | C9H16N4      | 1251330-45-6 | 46      |      |      |
| 4    | 1-Hydroxy-1-(4-methoxyphenyl)pro... | 180 | C10H12O3     | 015482-29-8  | 38      |      |      |
| 5    | 3,7-Benzofurandiyl, 2,3-dihydro-... | 180 | C10H12O3     | 017781-15-6  | 30      |      |      |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM010224\  
Data File : BM043693.D  
Acq On : 02 Jan 2024 13:53  
Operator : MA/JU  
Sample : 05918-19  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
GCF32

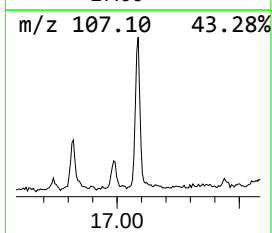
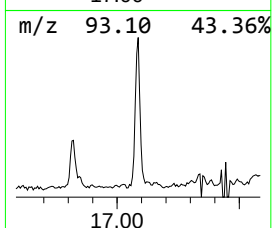
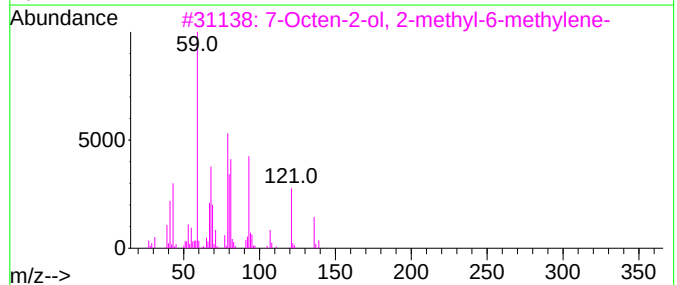
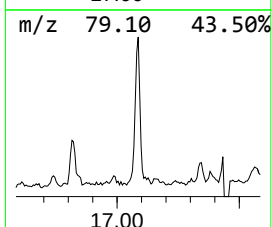
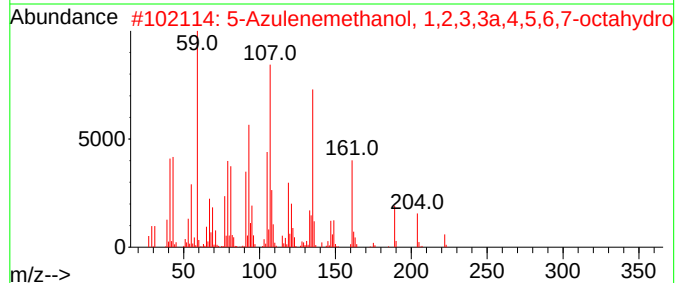
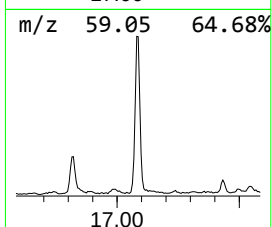
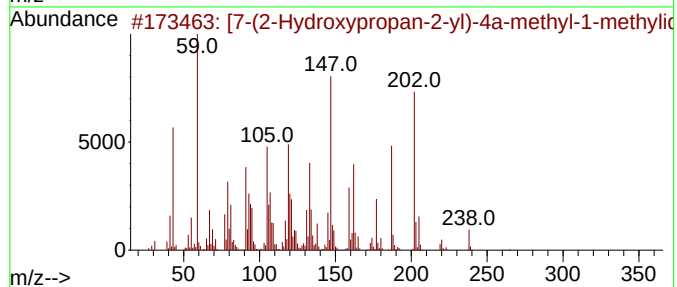
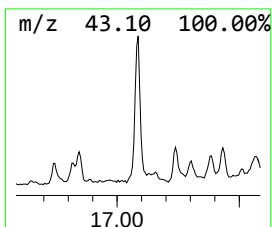
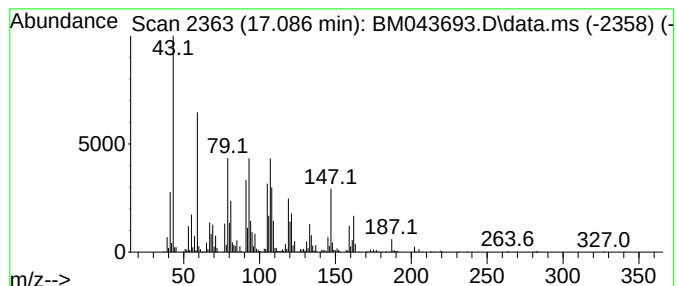
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Quant Title : SVOA CALIBRATION

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

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Peak Number 11 unknown-01 Concentration Rank 16

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|--------|------------------|--------------|------|
| 17.086    | 6.64 ng/ul                          | 737270 | Phenanthrene-d10 | 17.339       |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#         | Qual |
| 1         | [7-(2-Hydroxypropan-2-yl)-4a-met... | 280    | C17H28O3         | 1000502-54-5 | 41   |
| 2         | 5-Azulenemethanol, 1,2,3,3a,4,5,... | 222    | C15H26O          | 022451-73-6  | 37   |
| 3         | 7-Octen-2-ol, 2-methyl-6-methylene- | 154    | C10H18O          | 000543-39-5  | 32   |
| 4         | 1,1,4,7-Tetramethyldecahydro-1H-... | 238    | C15H26O2         | 1212211-43-2 | 32   |
| 5         | 2-Oxa-3-methylbicyclo(3.3.0)octane  | 140    | C9H16O           | 029183-69-5  | 25   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM010224\  
Data File : BM043693.D  
Acq On : 02 Jan 2024 13:53  
Operator : MA/JU  
Sample : 05918-19  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
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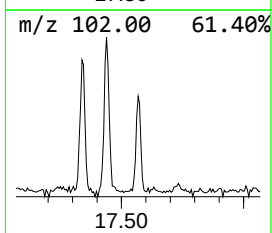
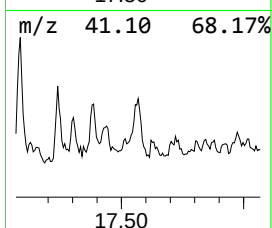
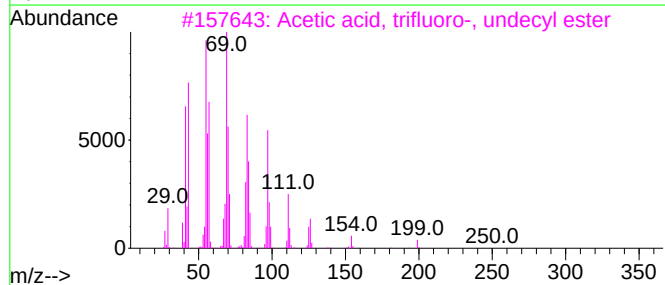
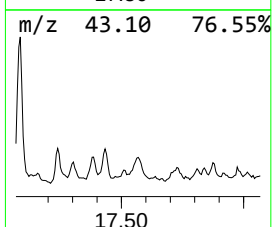
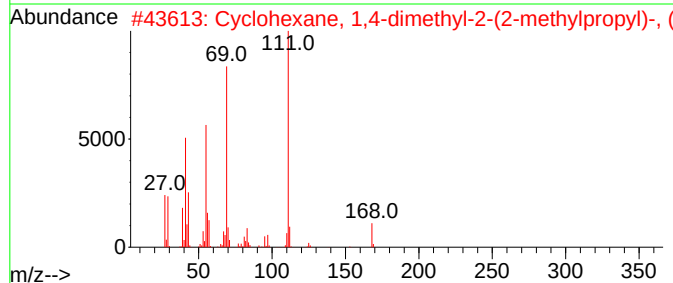
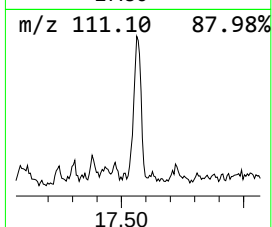
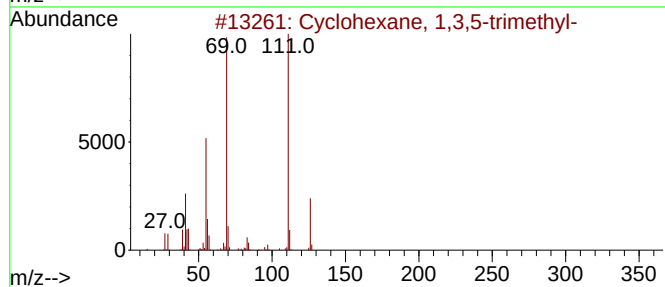
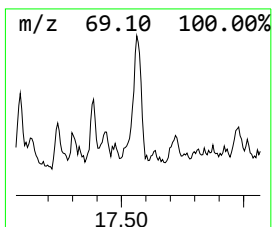
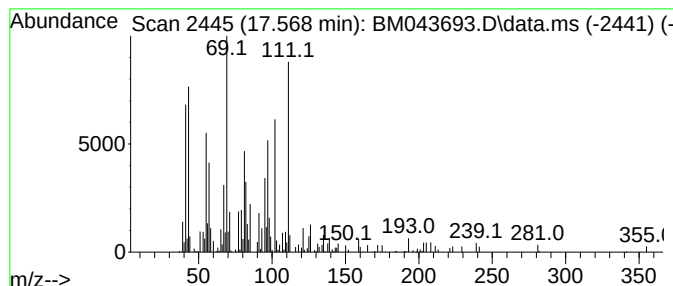
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Quant Title : SVOA CALIBRATION

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

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Peak Number 12 (DEL) Alkane: Cyclic17.568 Concentration Rank 31

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 17.568 | 2.13 ng/ul | 236269 | Phenanthrene-d10 | 17.339 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|------------|--------------|------|
| 1    |    |   | Cyclohexane, 1,3,5-trimethyl-       | 126 | C9H18      | 001839-63-0  | 46   |
| 2    |    |   | Cyclohexane, 1,4-dimethyl-2-(2-m... | 168 | C12H24     | 054411-17-5  | 38   |
| 3    |    |   | Acetic acid, trifluoro-, undecyl... | 268 | C13H23F3O2 | 053800-01-4  | 35   |
| 4    |    |   | 3-methyl-5-(2,6-dimethylheptyl)-... | 238 | C15H26O2   | 1000222-09-3 | 22   |
| 5    |    |   | 4-Pyridinol-1-oxide                 | 111 | C5H5NO2    | 006890-62-6  | 14   |





Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM010224\  
Data File : BM043693.D  
Acq On : 02 Jan 2024 13:53  
Operator : MA/JU  
Sample : 05918-19  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
GCF32

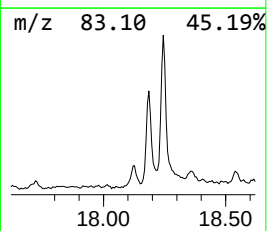
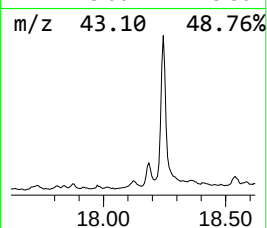
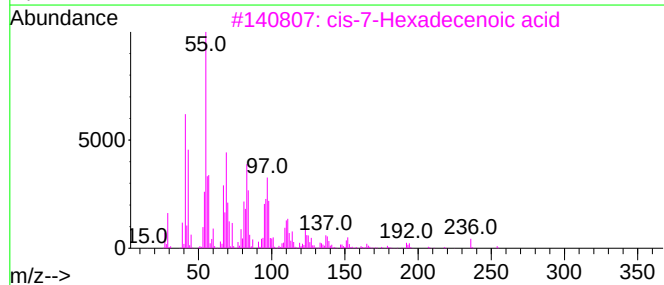
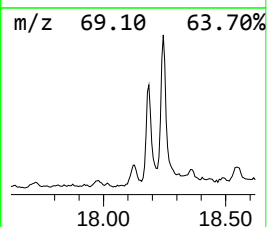
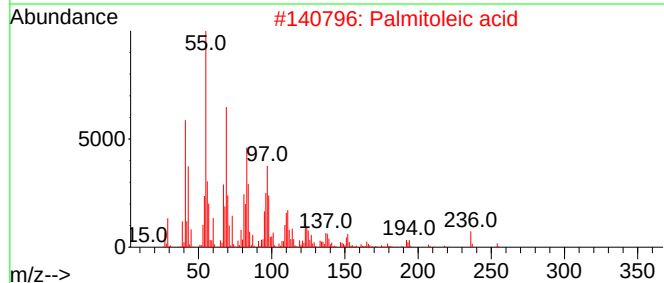
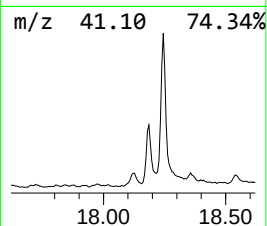
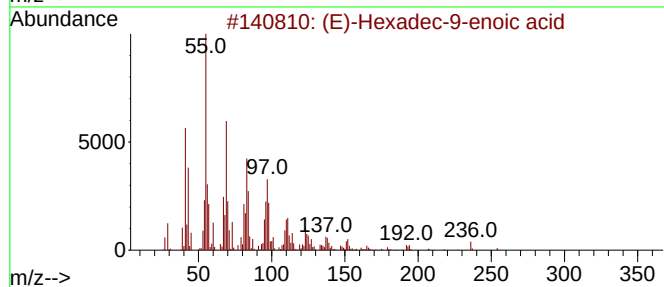
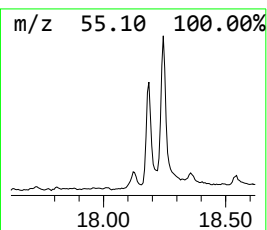
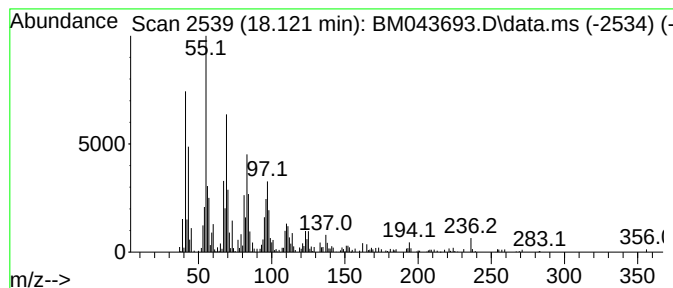
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Quant Title : SVOA CALIBRATION

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

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Peak Number 13 (E)-Hexadec-9-enoic acid Concentration Rank 28

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 18.121 | 2.39 ng/ul | 264967 | Phenanthrene-d10 | 17.339 |

| Hit# | of                       | 5 | Tentative ID | MW  | MolForm  | CAS#        | Qual |
|------|--------------------------|---|--------------|-----|----------|-------------|------|
| 1    | (E)-Hexadec-9-enoic acid |   |              | 254 | C16H30O2 | 010030-73-6 | 99   |
| 2    | Palmitoleic acid         |   |              | 254 | C16H30O2 | 000373-49-9 | 99   |
| 3    | cis-7-Hexadecenoic acid  |   |              | 254 | C16H30O2 | 002416-19-5 | 99   |
| 4    | Palmitoleic acid         |   |              | 254 | C16H30O2 | 000373-49-9 | 97   |
| 5    | Hexadecenoic acid, Z-11- |   |              | 254 | C16H30O2 | 002416-20-8 | 93   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM010224\  
Data File : BM043693.D  
Acq On : 02 Jan 2024 13:53  
Operator : MA/JU  
Sample : 05918-19  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
GCF32

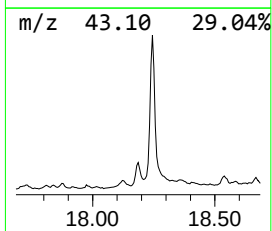
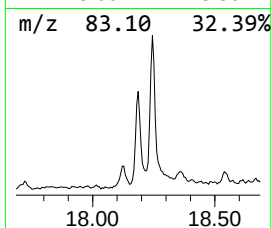
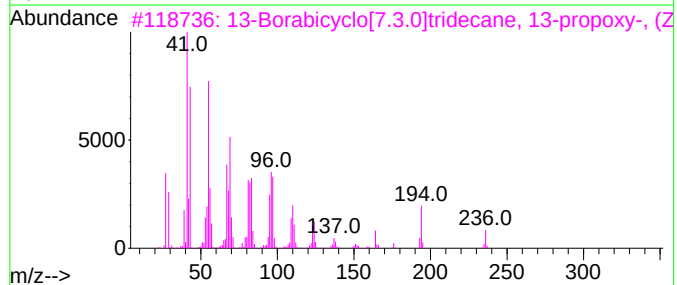
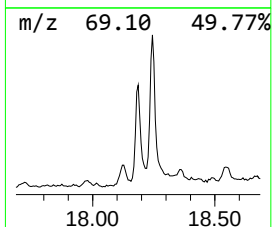
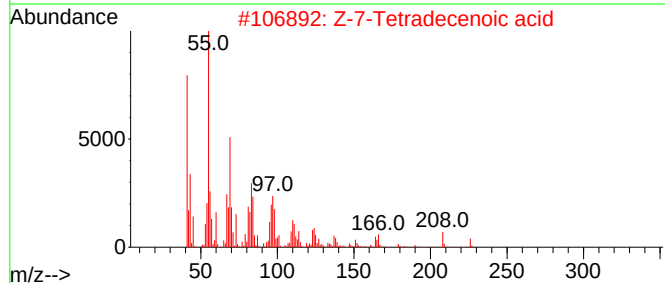
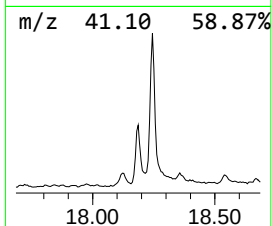
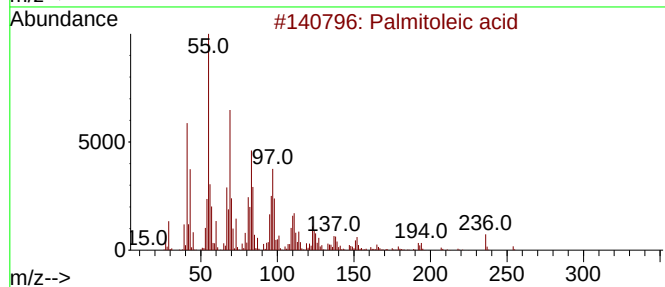
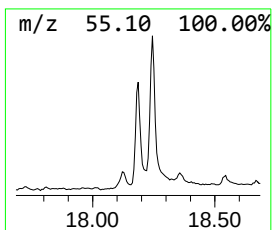
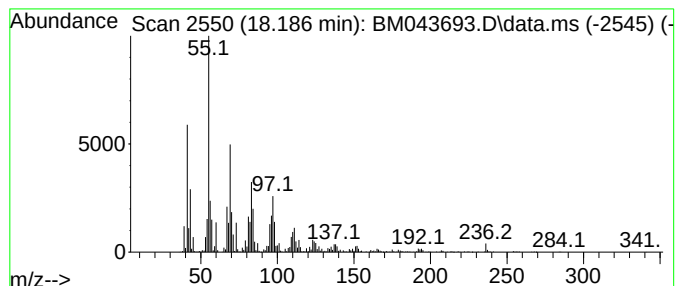
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Quant Title : SVOA CALIBRATION

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

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Peak Number 14 Palmitoleic acid Concentration Rank 13

| R.T.   | EstConc    | Area    | Relative to ISTD | R.T.   |
|--------|------------|---------|------------------|--------|
| 18.186 | 9.32 ng/ul | 1034360 | Phenanthrene-d10 | 17.339 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Palmitoleic acid                    | 254 | C16H30O2 | 000373-49-9  | 91   |
| 2    |    |   | Z-7-Tetradecenoic acid              | 226 | C14H26O2 | 1000130-98-4 | 91   |
| 3    |    |   | 13-Borabicyclo[7.3.0]tridecane, ... | 236 | C15H29BO | 1000156-41-7 | 89   |
| 4    |    |   | (E)-Hexadec-9-enoic acid            | 254 | C16H30O2 | 010030-73-6  | 87   |
| 5    |    |   | 9-Hexadecenoic acid                 | 254 | C16H30O2 | 002091-29-4  | 86   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM010224\  
Data File : BM043693.D  
Acq On : 02 Jan 2024 13:53  
Operator : MA/JU  
Sample : 05918-19  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
GCF32

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM121523.MA.M  
Quant Title : SVOA CALIBRATION

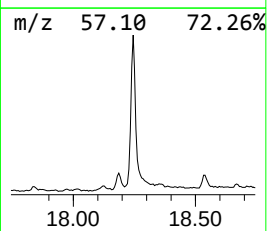
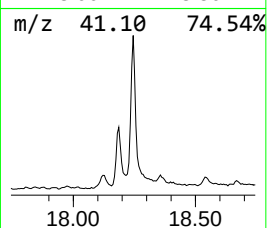
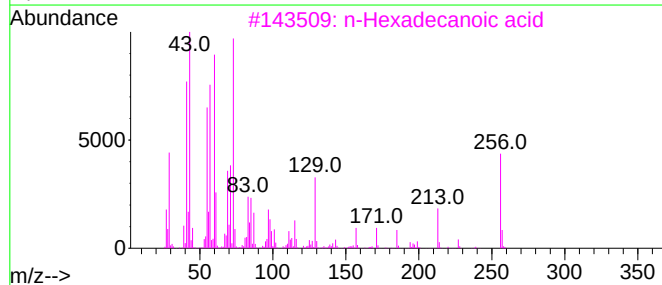
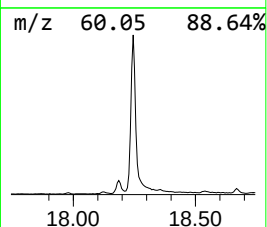
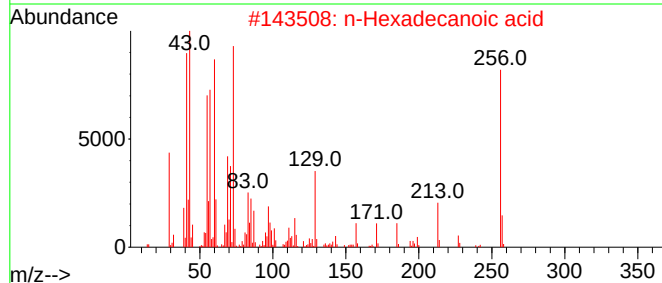
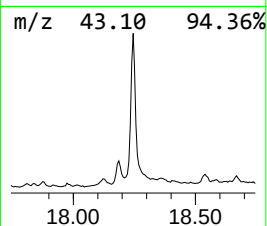
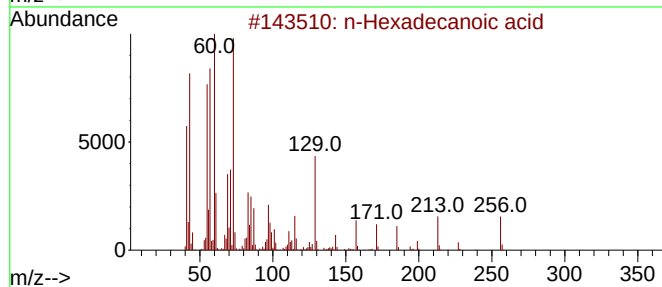
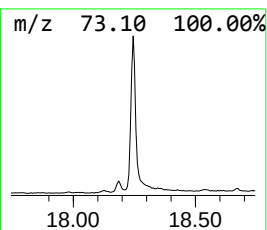
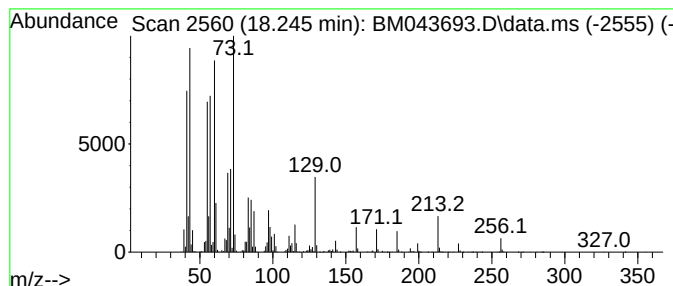
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 15 n-Hexadecanoic acid Concentration Rank 5

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 18.245 | 26.23 ng/ul | 2910950 | Phenanthrene-d10 | 17.339 |

| 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\_M\Data\BM010224\  
Data File : BM043693.D  
Acq On : 02 Jan 2024 13:53  
Operator : MA/JU  
Sample : 05918-19  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
GCF32

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM121523.MA.M  
Quant Title : SVOA CALIBRATION

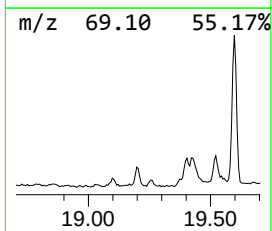
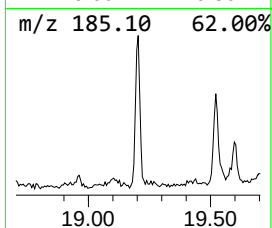
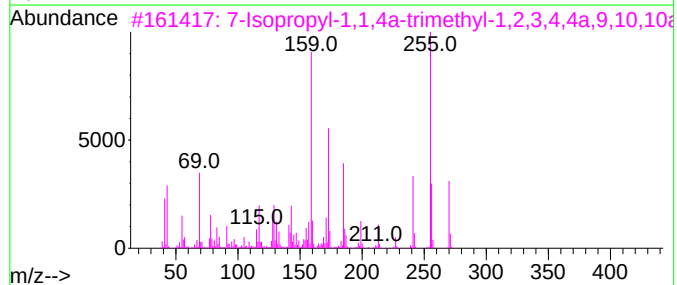
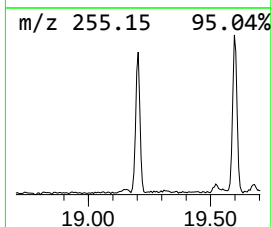
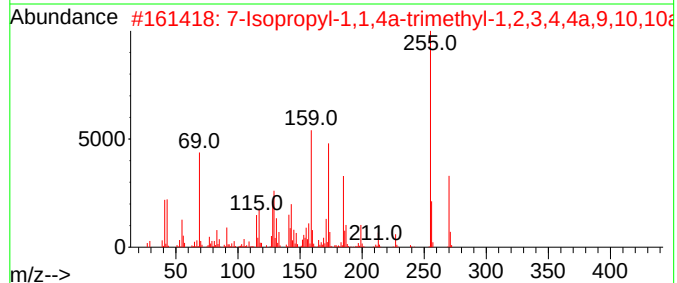
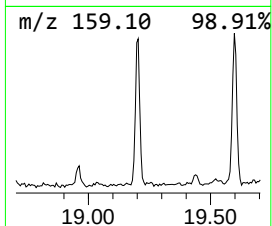
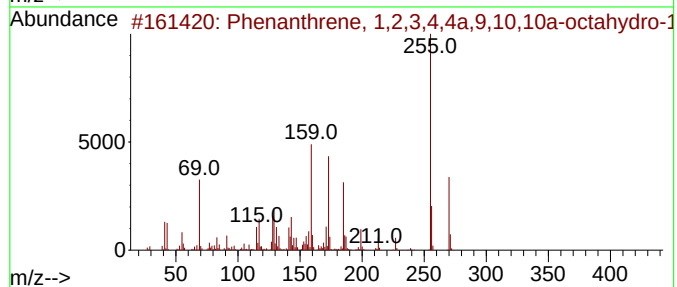
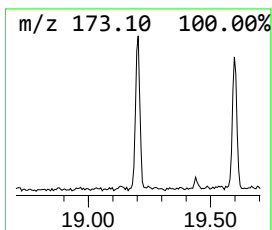
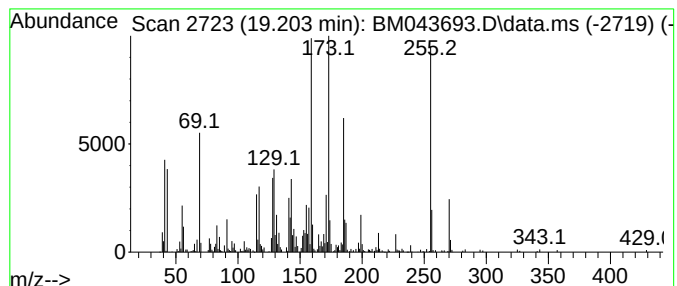
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 16 Phenanthrene, 1,2,3,4,4a,9,... Concentration Rank 24

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 19.203 | 2.95 ng/ul | 327639 | Phenanthrene-d10 | 17.339 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Phenanthrene, 1,2,3,4,4a,9,10,10... | 270 | C20H30    | 019407-28-4  | 99   |
| 2    |    |   | 7-Isopropyl-1,1,4a-trimethyl-1,2... | 270 | C20H30    | 109680-01-5  | 94   |
| 3    |    |   | 7-Isopropyl-1,1,4a-trimethyl-1,2... | 270 | C20H30    | 109680-01-5  | 89   |
| 4    |    |   | 2-Thioxo-imidazolidin-4-one-5-ac... | 173 | C5H7N3O2S | 1010146-03-1 | 35   |
| 5    |    |   | 2(5H)-Furanone, 5-(phenylimino)-    | 173 | C10H7NO2  | 019990-26-2  | 20   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM010224\  
Data File : BM043693.D  
Acq On : 02 Jan 2024 13:53  
Operator : MA/JU  
Sample : 05918-19  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

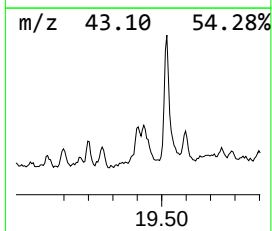
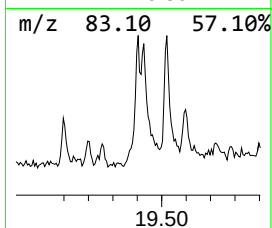
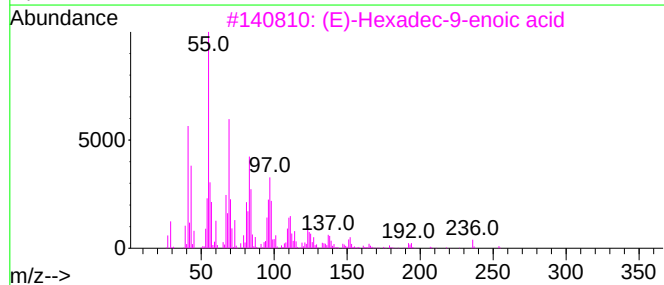
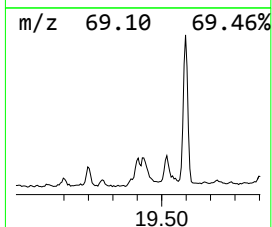
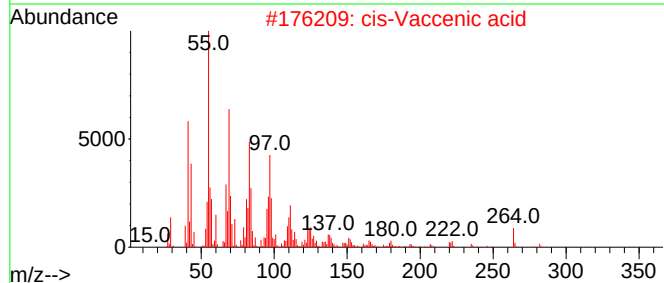
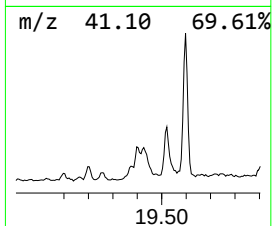
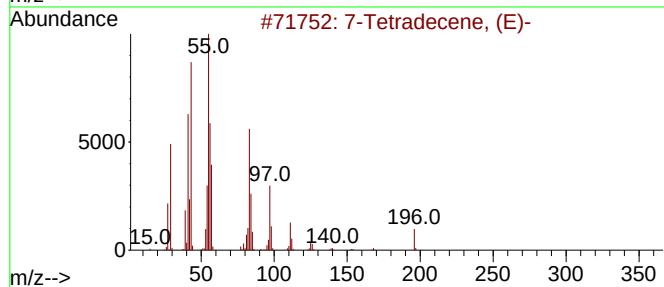
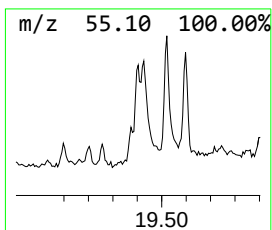
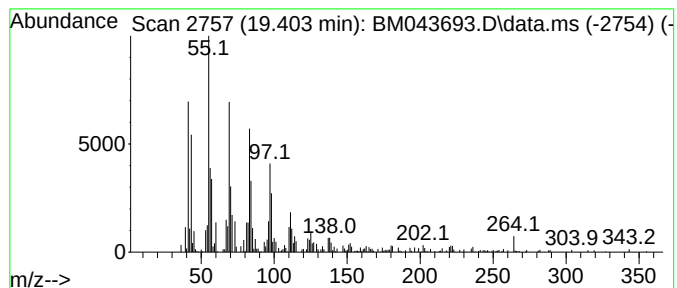
Instrument :  
BNA\_M  
ClientSampleId :  
GCF32

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM121523.MA.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 17 (DEL) Alkane: Straight-Chai... Concentration Rank 30

| R.T.      | EstConc                             | Area   | Relative to ISTD |           | R.T.         |      |
|-----------|-------------------------------------|--------|------------------|-----------|--------------|------|
| 19.403    | 2.32 ng/ul                          | 257617 | Phenanthrene-d10 |           | 17.339       |      |
| Hit# of 5 | Tentative ID                        |        | MW               | MolForm   | CAS#         | Qual |
| 1         | 7-Tetradecene, (E)-                 |        | 196              | C14H28    | 041446-63-3  | 91   |
| 2         | cis-Vaccenic acid                   |        | 282              | C18H34O2  | 000506-17-2  | 91   |
| 3         | (E)-Hexadec-9-enoic acid            |        | 254              | C16H30O2  | 010030-73-6  | 87   |
| 4         | trans-2-methyl-4-n-pentylthiane,... |        | 218              | C11H22O2S | 1000215-75-3 | 86   |
| 5         | Heptadecanolide                     |        | 268              | C17H32O2  | 005637-97-8  | 86   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM010224\  
Data File : BM043693.D  
Acq On : 02 Jan 2024 13:53  
Operator : MA/JU  
Sample : 05918-19  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
GCF32

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM121523.MA.M  
Quant Title : SVOA CALIBRATION

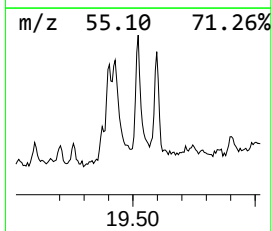
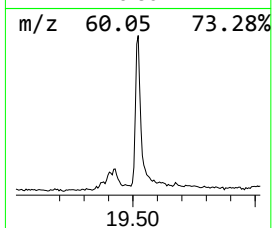
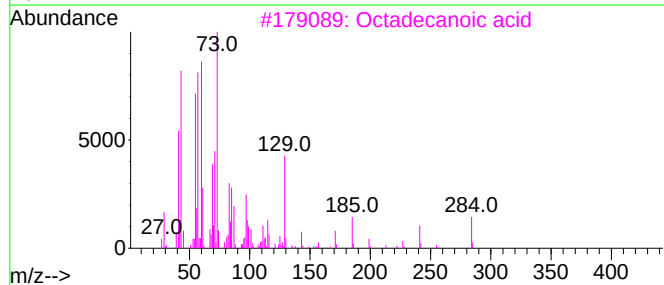
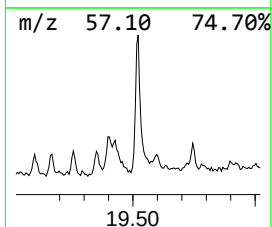
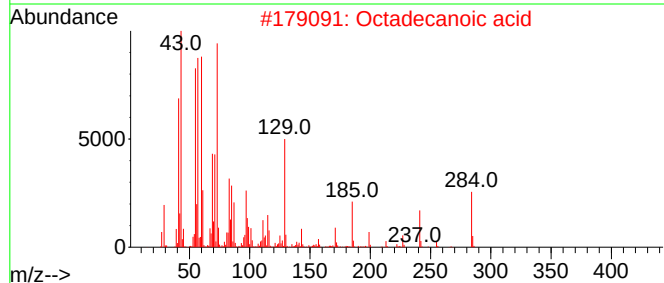
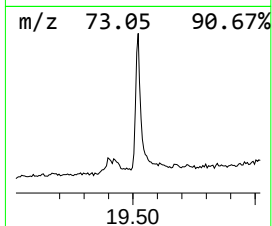
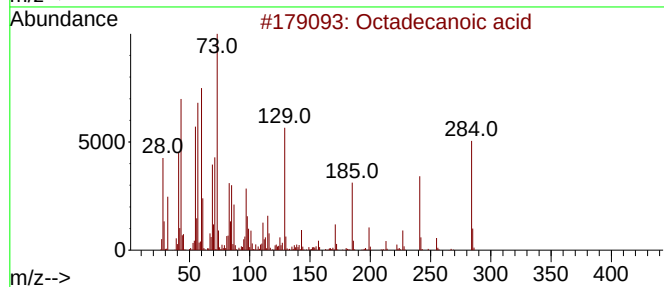
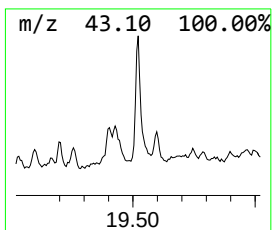
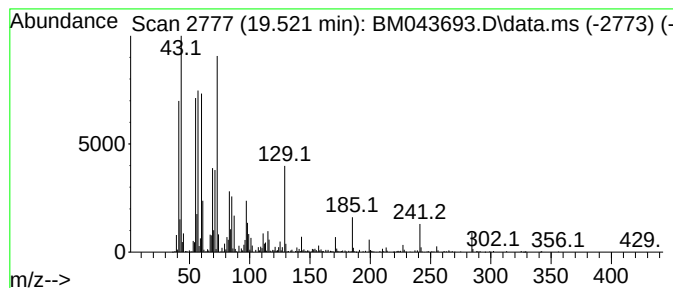
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 18 Octadecanoic acid Concentration Rank 17

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 19.521 | 5.99 ng/ul | 795008 | Chrysene-d12     | 21.521 |

| 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 | 96   |
| 4    |      | Octadecanoic acid  | 284 | C18H36O2 | 000057-11-4 | 95   |
| 5    |      | Pentadecanoic acid | 242 | C15H30O2 | 001002-84-2 | 90   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM010224\  
Data File : BM043693.D  
Acq On : 02 Jan 2024 13:53  
Operator : MA/JU  
Sample : 05918-19  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
GCF32

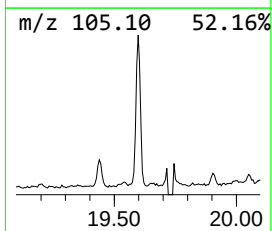
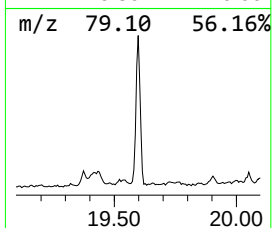
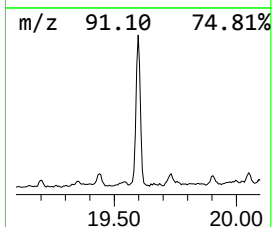
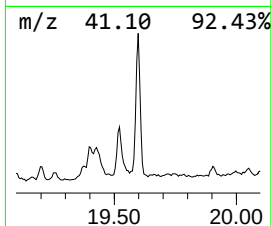
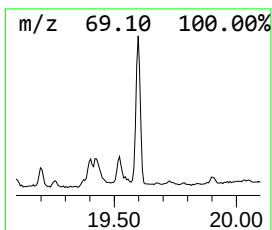
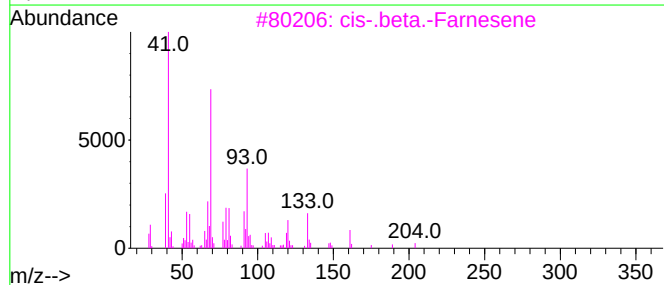
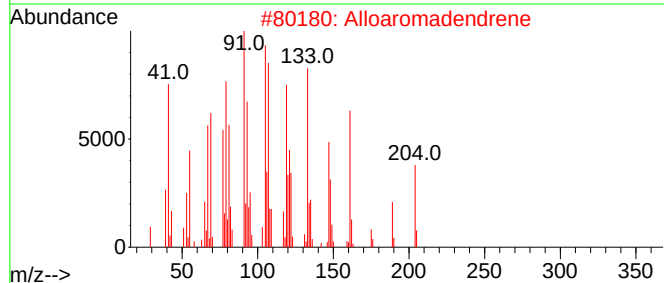
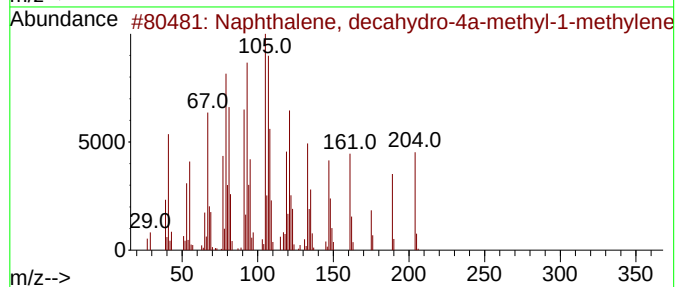
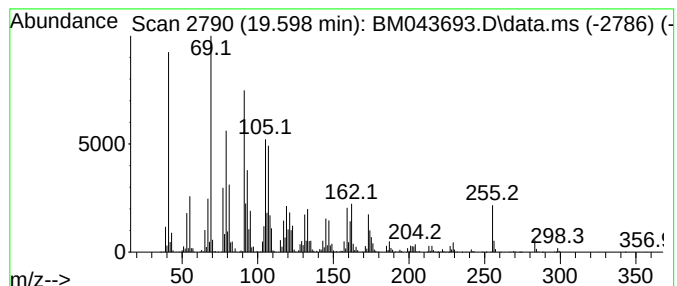
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Quant Title : SVOA CALIBRATION

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

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Peak Number 19 unknown-02 Concentration Rank 9

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 19.598 | 12.77 ng/ul | 1695590 | Chrysene-d12     | 21.521 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, decahydro-4a-methyl... | 204 | C15H24  | 017066-67-0 | 46   |
| 2    |    |   | Alloaromadendrene                   | 204 | C15H24  | 025246-27-9 | 42   |
| 3    |    |   | cis-.beta.-Farnesene                | 204 | C15H24  | 028973-97-9 | 38   |
| 4    |    |   | 1,6,10-Dodecatriene, 7,11-dimeth... | 204 | C15H24  | 077129-48-7 | 30   |
| 5    |    |   | 2,6,10-Dodecatrien-1-ol, 3,7,11-... | 222 | C15H26O | 004602-84-0 | 25   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM010224\  
Data File : BM043693.D  
Acq On : 02 Jan 2024 13:53  
Operator : MA/JU  
Sample : 05918-19  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
GCF32

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM121523.MA.M  
Quant Title : SVOA CALIBRATION

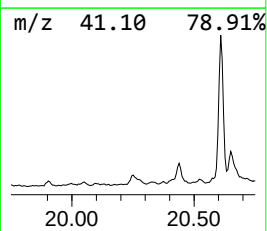
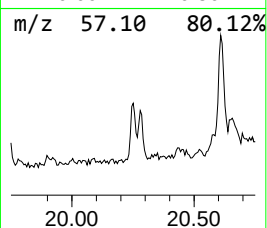
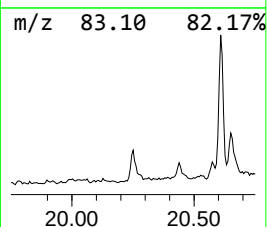
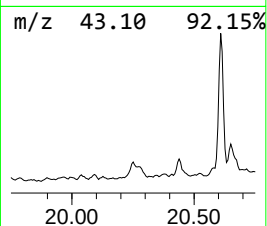
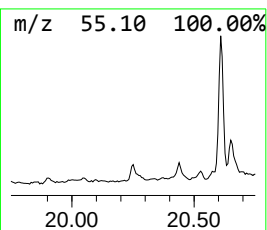
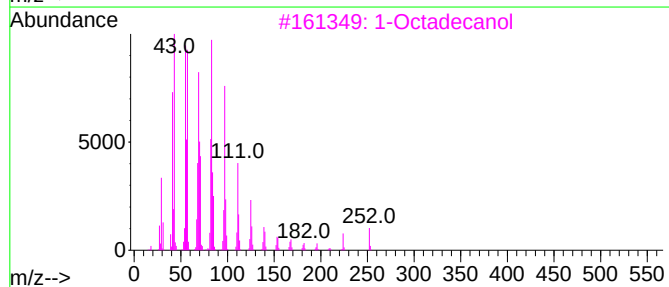
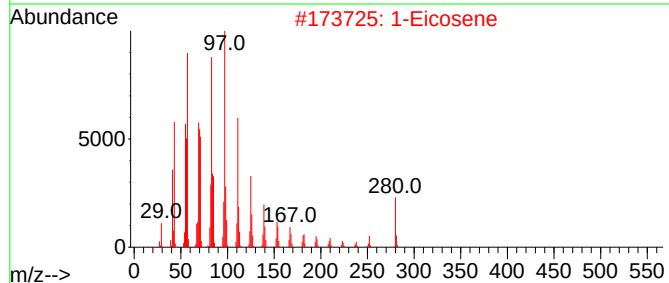
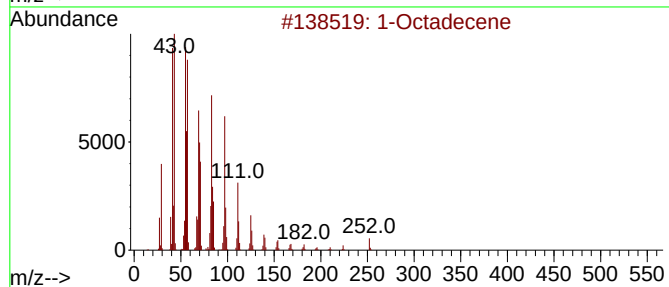
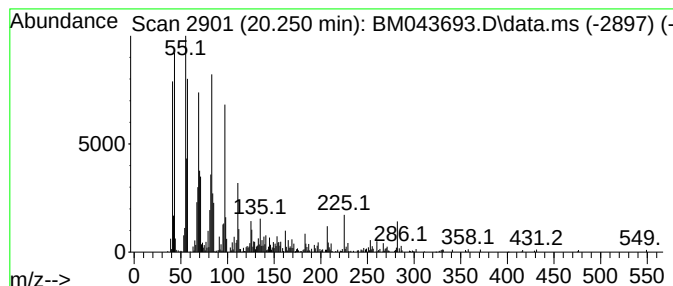
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 20 (DEL) Alkane: Straight-Chai... Concentration Rank 21

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 20.250 | 4.25 ng/ul | 564896 | Chrysene-d12     | 21.521 |

| Hit# | of 5 | Tentative ID     | MW  | MolForm | CAS#        | Qual |
|------|------|------------------|-----|---------|-------------|------|
| 1    |      | 1-Octadecene     | 252 | C18H36  | 000112-88-9 | 98   |
| 2    |      | 1-Eicosene       | 280 | C20H40  | 003452-07-1 | 96   |
| 3    |      | 1-Octadecanol    | 270 | C18H38O | 000112-92-5 | 95   |
| 4    |      | n-Nonadecanol-1  | 284 | C19H40O | 001454-84-8 | 95   |
| 5    |      | n-Heptadecanol-1 | 256 | C17H36O | 001454-85-9 | 94   |





Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM010224\  
Data File : BM043693.D  
Acq On : 02 Jan 2024 13:53  
Operator : MA/JU  
Sample : 05918-19  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
GCF32

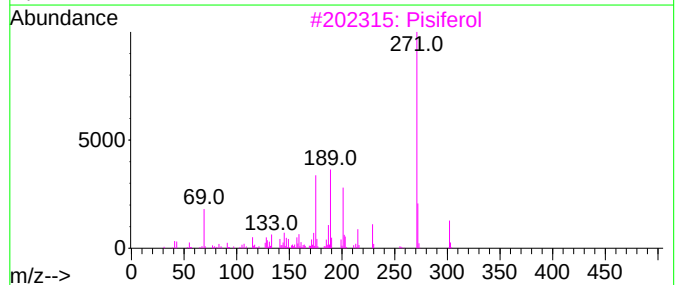
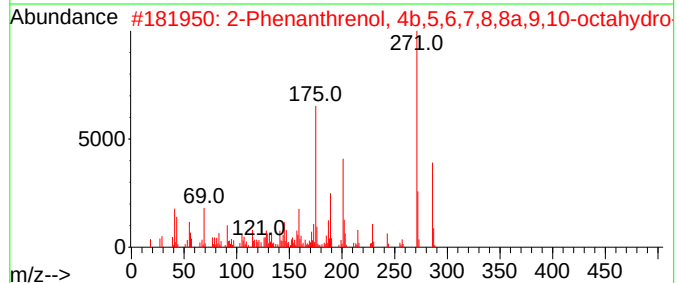
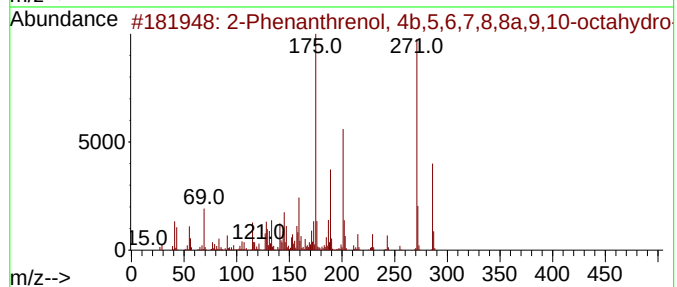
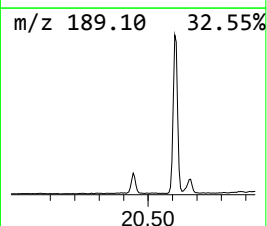
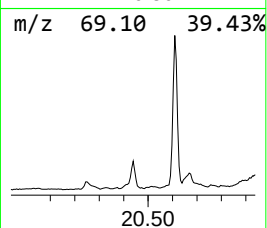
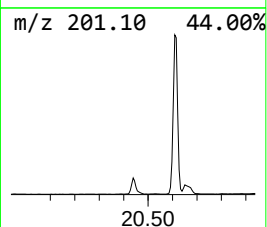
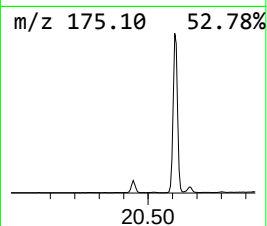
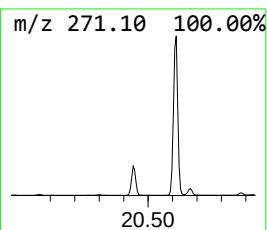
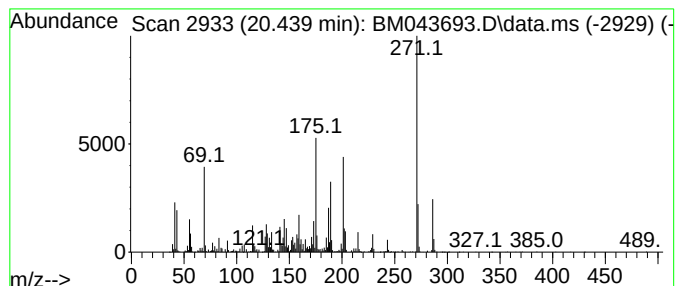
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Quant Title : SVOA CALIBRATION

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

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Peak Number 21 2-Phenanthrenol, 4b,5,6,7,8... Concentration Rank 14

| R.T.   | EstConc    | Area    | Relative to ISTD | R.T.   |
|--------|------------|---------|------------------|--------|
| 20.439 | 8.91 ng/ul | 1183270 | Chrysene-d12     | 21.521 |

| Hit# | of                                  | 5   | Tentative ID | MW           | MolForm | CAS# | Qual |
|------|-------------------------------------|-----|--------------|--------------|---------|------|------|
| 1    | 2-Phenanthrenol, 4b,5,6,7,8,8a,9... | 286 | C20H30O      | 000511-15-9  | 98      |      |      |
| 2    | 2-Phenanthrenol, 4b,5,6,7,8,8a,9... | 286 | C20H30O      | 000511-15-9  | 55      |      |      |
| 3    | Pisiferol                           | 302 | C20H30O2     | 024035-36-7  | 52      |      |      |
| 4    | 13-Isopropylpodocarpene-12-ol-20-al | 300 | C20H28O2     | 024035-37-8  | 50      |      |      |
| 5    | Silane, dimethyl(2-chlorophenoxy... | 286 | C14H23ClO2Si | 1000347-07-3 | 49      |      |      |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM010224\  
Data File : BM043693.D  
Acq On : 02 Jan 2024 13:53  
Operator : MA/JU  
Sample : 05918-19  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
GCF32

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM121523.MA.M  
Quant Title : SVOA CALIBRATION

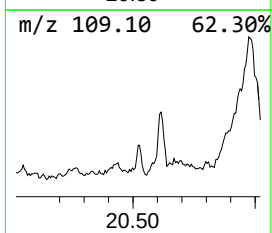
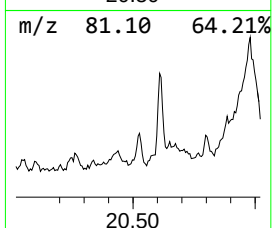
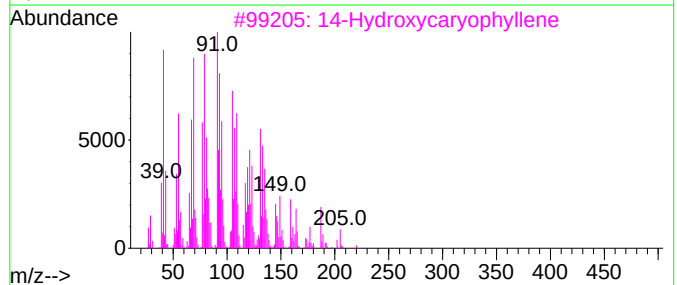
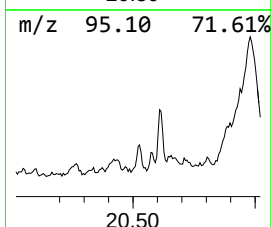
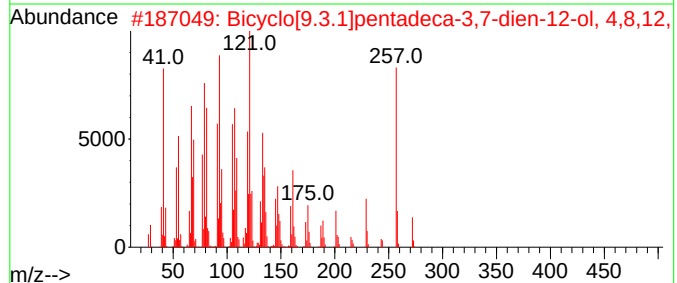
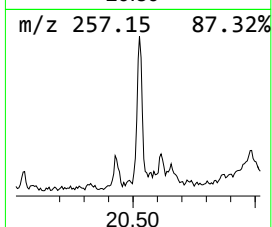
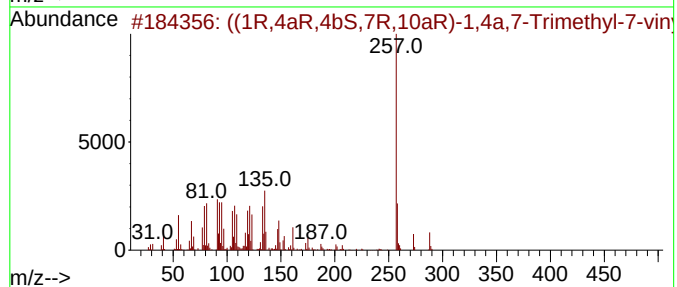
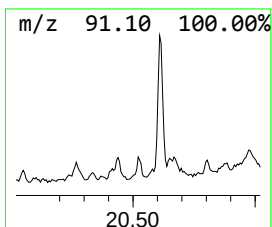
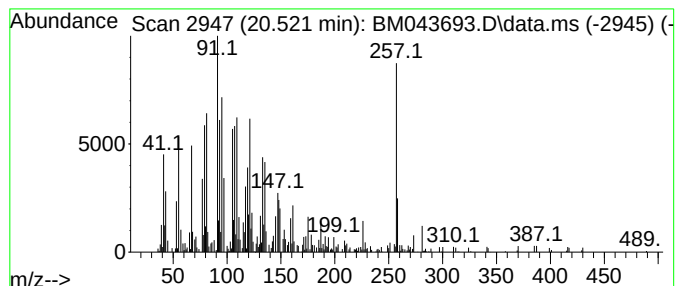
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TIC Integration Parameters: LSCINT.P

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Peak Number 22 ((1R,4aR,4bS,7R,10aR)-1,4a,... Concentration Rank 32

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 20.521 | 2.05 ng/ul | 271643 | Chrysene-d12     | 21.521 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | ((1R,4aR,4bS,7R,10aR)-1,4a,7-Tri... | 288 | C20H32O   | 024563-84-6  | 80   |
| 2    |    |   | Bicyclo[9.3.1]pentadeca-3,7-dien... | 290 | C20H34O   | 070000-19-0  | 49   |
| 3    |    |   | 14-Hydroxycaryophyllene             | 220 | C15H24O   | 050277-33-3  | 41   |
| 4    |    |   | 1R,3Z,9S-4,11,11-Trimethyl-8-met... | 204 | C15H24    | 1000159-39-4 | 27   |
| 5    |    |   | Acetamide, N-(3-methylphenyl)-2-... | 257 | C15H15NOS | 1000307-20-6 | 25   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM010224\  
Data File : BM043693.D  
Acq On : 02 Jan 2024 13:53  
Operator : MA/JU  
Sample : 05918-19  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
GCF32

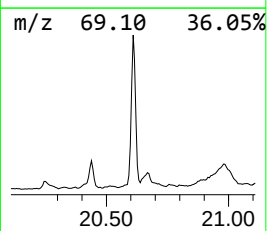
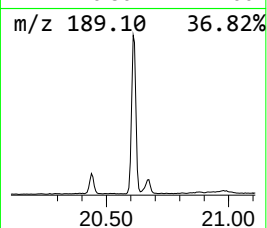
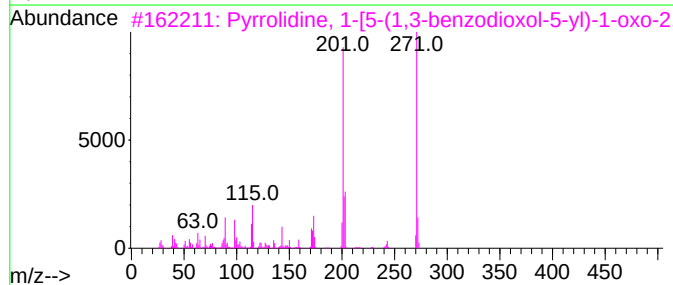
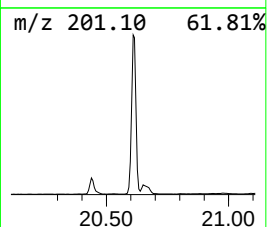
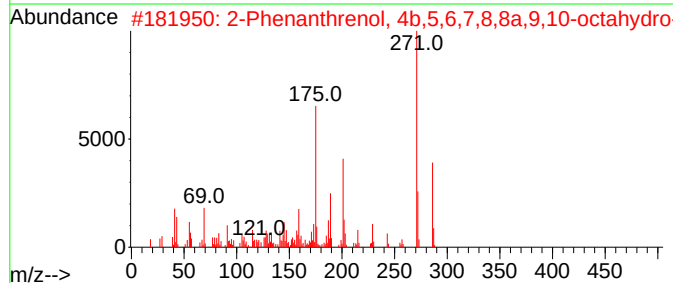
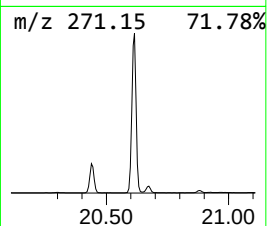
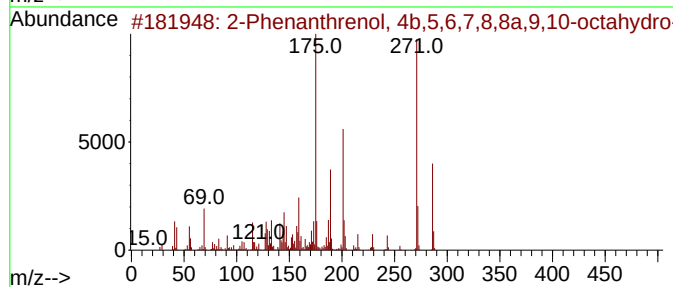
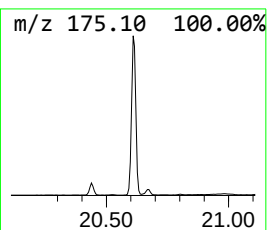
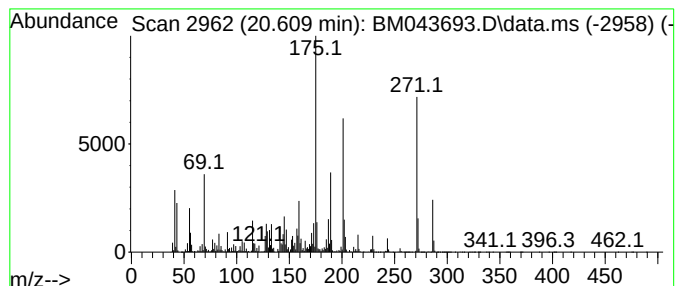
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Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 23 unknown-03 Concentration Rank 1

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 20.609 | 63.86 ng/ul | 8478590 | Chrysene-d12     | 21.521 |

| Hit# | of                                  | 5   | Tentative ID | MW          | MolForm | CAS# | Qual |
|------|-------------------------------------|-----|--------------|-------------|---------|------|------|
| 1    | 2-Phenanthrenol, 4b,5,6,7,8,8a,9... | 286 | C20H30O      | 000511-15-9 | 99      |      |      |
| 2    | 2-Phenanthrenol, 4b,5,6,7,8,8a,9... | 286 | C20H30O      | 000511-15-9 | 95      |      |      |
| 3    | Pyrrolidine, 1-[5-(1,3-benzodiox... | 271 | C16H17NO3    | 025924-78-1 | 38      |      |      |
| 4    | m-Cymene, 5-tert-butyl-             | 190 | C14H22       | 029577-19-3 | 18      |      |      |
| 5    | Benzene, 1,4-dimethyl-2,5-bis(1-... | 190 | C14H22       | 010375-96-9 | 18      |      |      |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM010224\  
Data File : BM043693.D  
Acq On : 02 Jan 2024 13:53  
Operator : MA/JU  
Sample : 05918-19  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
GCF32

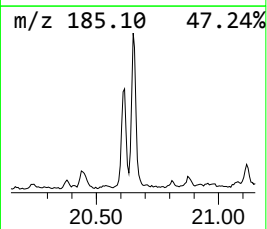
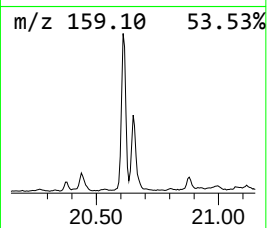
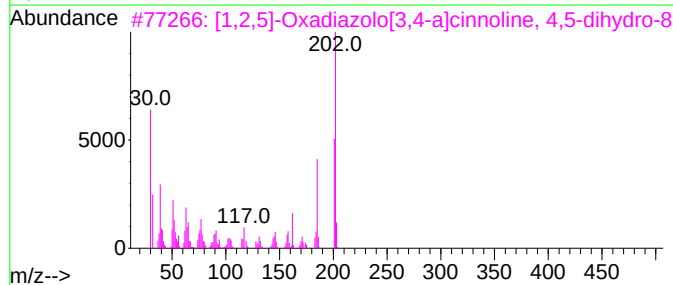
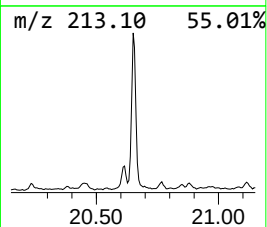
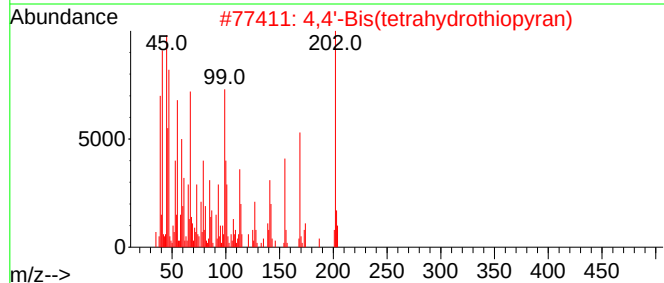
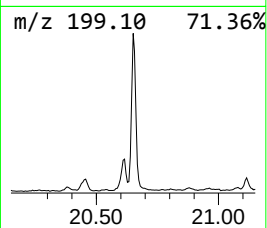
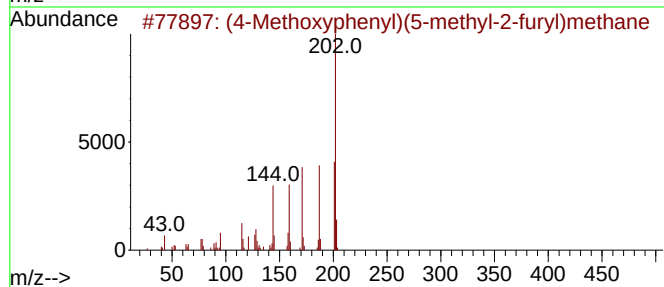
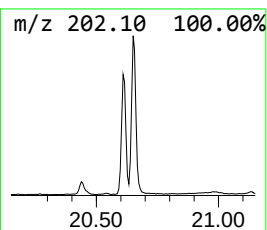
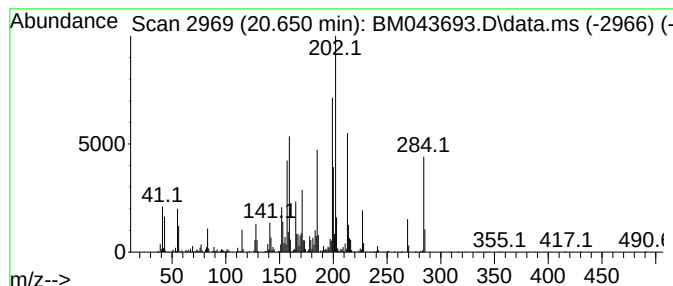
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Quant Title : SVOA CALIBRATION

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

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Peak Number 24 unknown-04 Concentration Rank 8

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 20.650 | 21.33 ng/ul | 2831530 | Chrysene-d12     | 21.521 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | (4-Methoxyphenyl)(5-methyl-2-fur... | 202 | C13H14O2  | 1000440-64-7 | 25   |
| 2    |    |   | 4,4'-Bis(tetrahydrothiopyran)       | 202 | C10H18S2  | 116196-83-9  | 25   |
| 3    |    |   | [1,2,5]-Oxadiazolo[3,4-a]cinnoli... | 202 | C10H10N4O | 1000260-59-4 | 20   |
| 4    |    |   | 9-Borabicyclo[3.3.1]nonane, 9-br... | 200 | C8H14BBr  | 022086-45-9  | 18   |
| 5    |    |   | 1-Naphthalenecarboxylic acid, 2-... | 202 | C12H10O3  | 000947-62-6  | 18   |





Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM010224\  
Data File : BM043693.D  
Acq On : 02 Jan 2024 13:53  
Operator : MA/JU  
Sample : 05918-19  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
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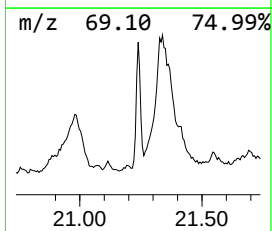
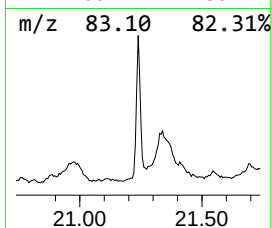
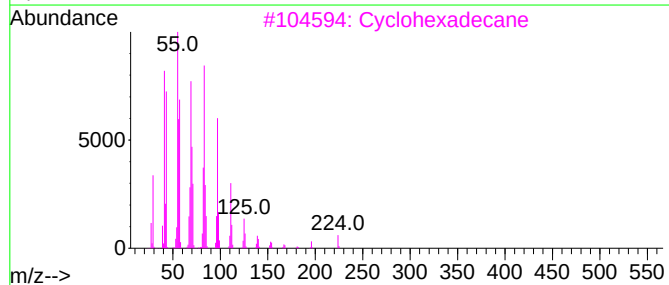
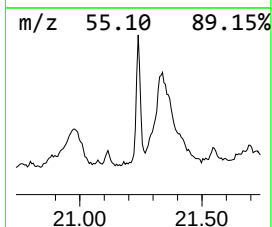
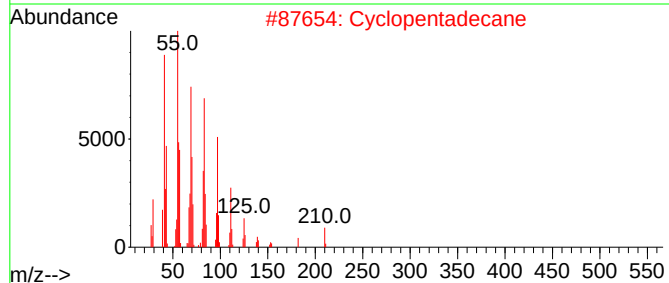
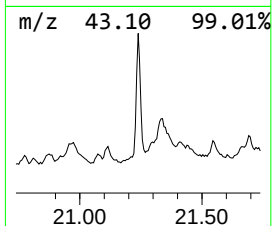
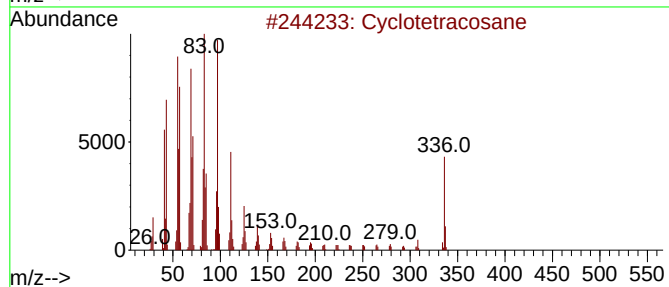
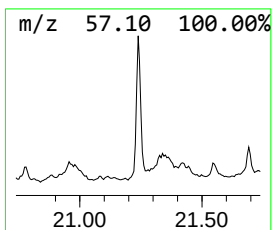
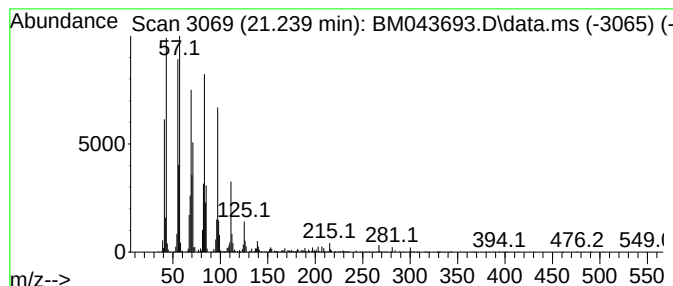
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Quant Title : SVOA CALIBRATION

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

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Peak Number 26 (DEL) Alkane: Cyclic21.239 Concentration Rank 12

| R.T.   | EstConc    | Area    | Relative to ISTD | R.T.   |
|--------|------------|---------|------------------|--------|
| 21.239 | 9.94 ng/ul | 1319740 | Chrysene-d12     | 21.521 |

| Hit# | of | 5 | Tentative ID         | MW  | MolForm  | CAS#        | Qual |
|------|----|---|----------------------|-----|----------|-------------|------|
| 1    |    |   | Cyclotetracosane     | 336 | C24H48   | 000297-03-0 | 98   |
| 2    |    |   | Cyclopentadecane     | 210 | C15H30   | 000295-48-7 | 97   |
| 3    |    |   | Cyclohexadecane      | 224 | C16H32   | 000295-65-8 | 94   |
| 4    |    |   | 1-Hexacosene         | 364 | C26H52   | 018835-33-1 | 94   |
| 5    |    |   | 1-Heneicosyl formate | 340 | C22H44O2 | 077899-03-7 | 94   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM010224\  
Data File : BM043693.D  
Acq On : 02 Jan 2024 13:53  
Operator : MA/JU  
Sample : 05918-19  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

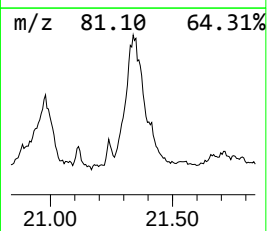
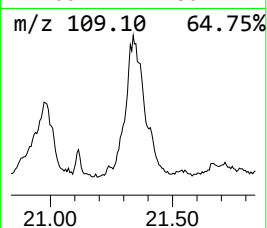
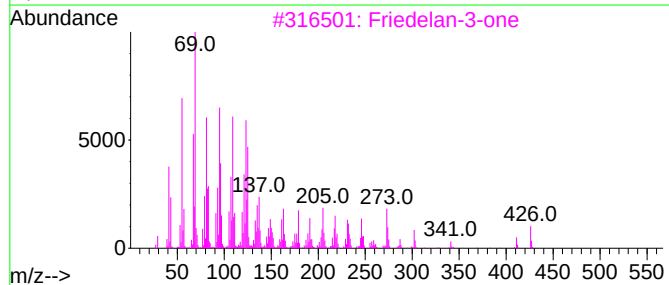
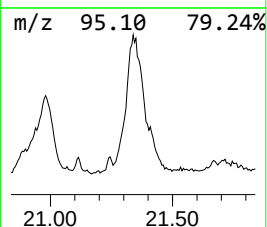
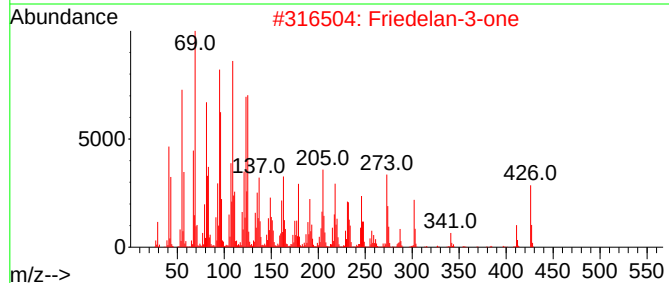
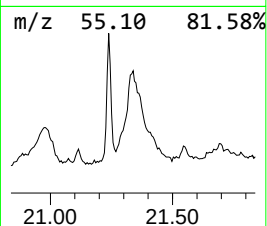
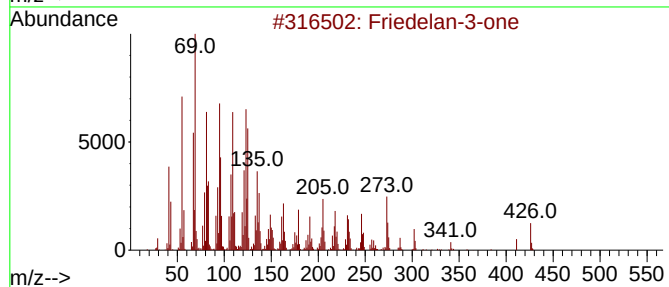
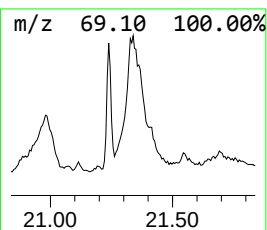
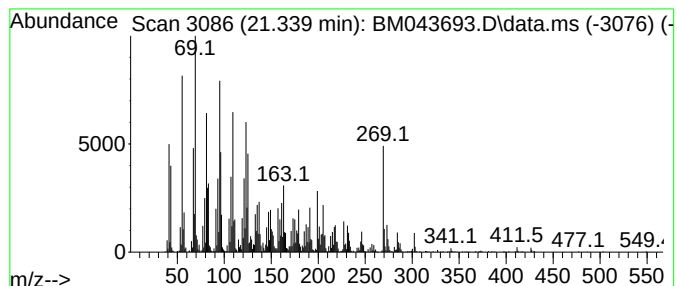
Instrument :  
BNA\_M  
ClientSampleId :  
GCF32

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM121523.MA.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 27 unknown-05 Concentration Rank 2

| R.T.      | EstConc                             | Area    | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|---------|------------------|--------------|------|
| 21.339    | 41.10 ng/ul                         | 5457440 | Chrysene-d12     | 21.521       |      |
| Hit# of 5 | Tentative ID                        | MW      | MolForm          | CAS#         | Qual |
| 1         | Friedelan-3-one                     | 426     | C30H50O          | 000559-74-0  | 96   |
| 2         | Friedelan-3-one                     | 426     | C30H50O          | 000559-74-0  | 90   |
| 3         | Friedelan-3-one                     | 426     | C30H50O          | 000559-74-0  | 64   |
| 4         | (1,2,4-Trimethylcyclohexyl)metha... | 198     | C12H22O2         | 1000499-04-5 | 25   |
| 5         | Cyclohexanone, 2-(1-mercapto-1-m... | 186     | C10H18OS         | 033281-91-3  | 22   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM010224\  
Data File : BM043693.D  
Acq On : 02 Jan 2024 13:53  
Operator : MA/JU  
Sample : 05918-19  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
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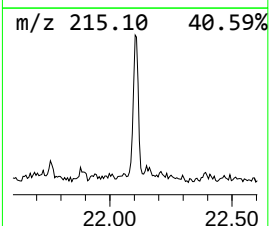
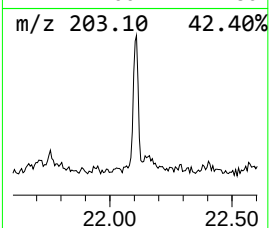
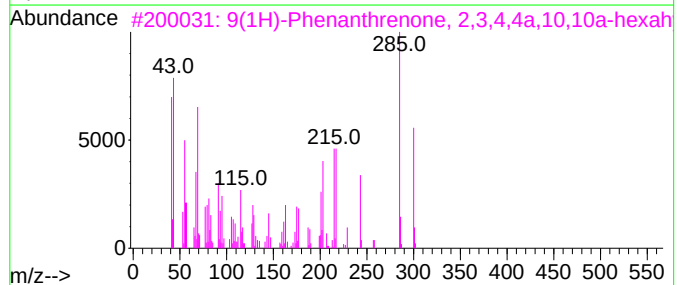
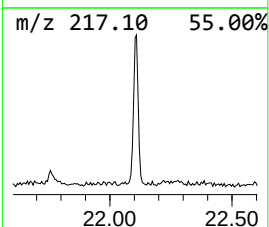
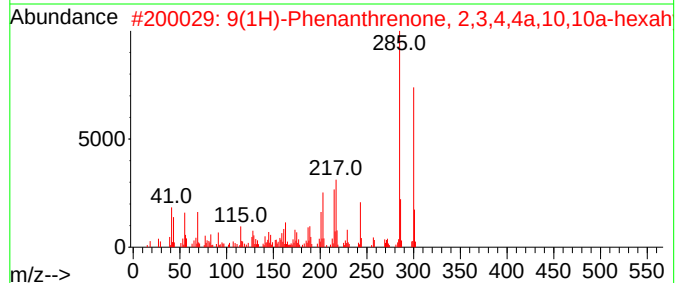
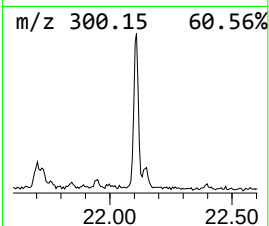
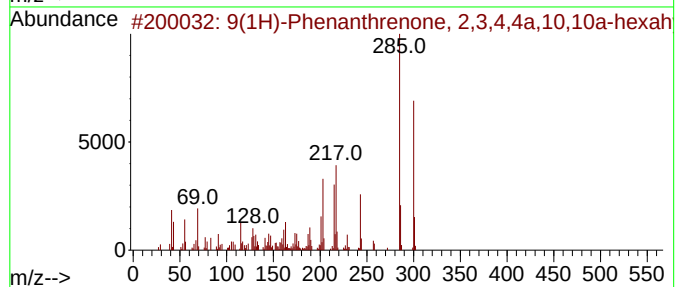
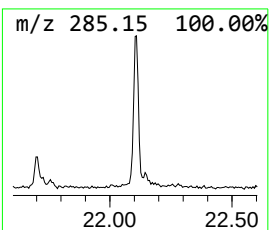
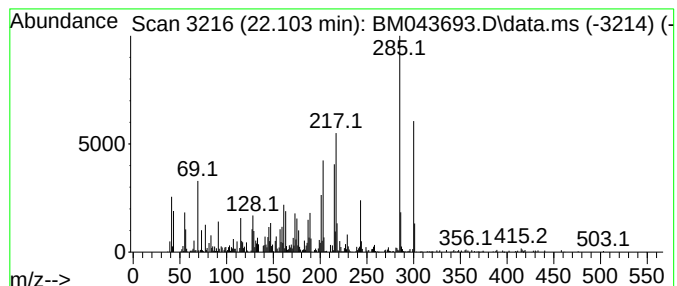
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Quant Title : SVOA CALIBRATION

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

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Peak Number 28 9(1H)-Phenanthrene, 2,3,4... Concentration Rank 23

| R.T.   | EstConc                             | Area         | Relative to ISTD | R.T.      |
|--------|-------------------------------------|--------------|------------------|-----------|
| 22.103 | 3.08 ng/ul                          | 409187       | Chrysene-d12     | 21.521    |
| Hit#   | of 5                                | Tentative ID | MW MolForm       | CAS# Qual |
| 1      | 9(1H)-Phenanthrene, 2,3,4,4a,1...   | 300 C20H28O2 | 000511-05-7      | 96        |
| 2      | 9(1H)-Phenanthrene, 2,3,4,4a,1...   | 300 C20H28O2 | 000511-05-7      | 94        |
| 3      | 9(1H)-Phenanthrene, 2,3,4,4a,1...   | 300 C20H28O2 | 000511-05-7      | 90        |
| 4      | Phenanthrene, 1,2,3,4,4a,9,10,10... | 300 C21H32O  | 015340-83-7      | 46        |
| 5      | 2(1H)-Phenanthrene, 3,4,4a,9,1...   | 300 C20H28O2 | 006755-93-7      | 43        |





Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM010224\  
Data File : BM043693.D  
Acq On : 02 Jan 2024 13:53  
Operator : MA/JU  
Sample : 05918-19  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

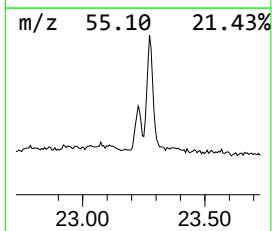
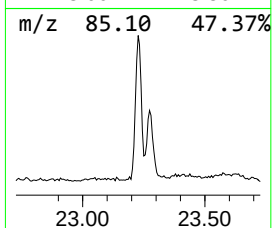
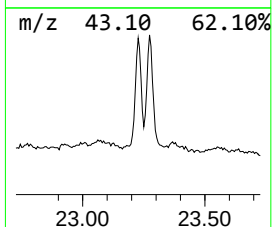
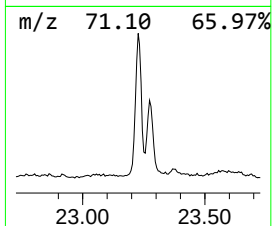
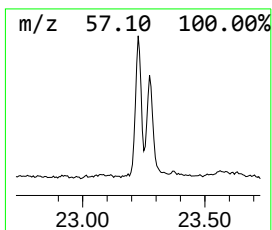
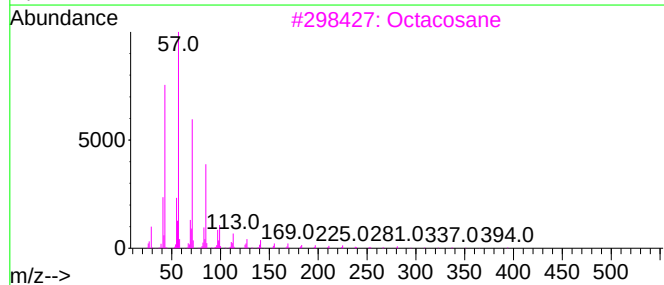
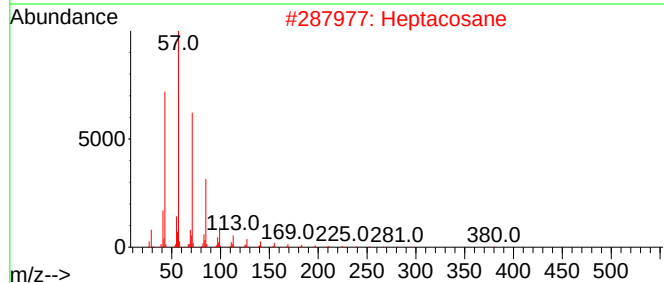
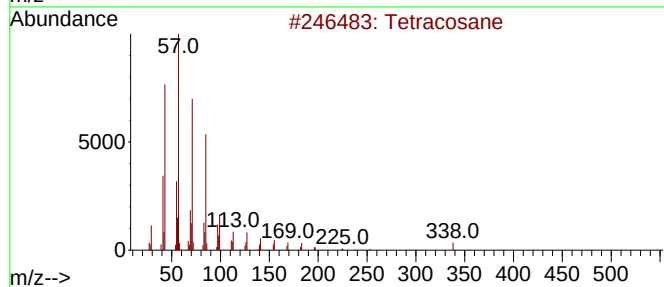
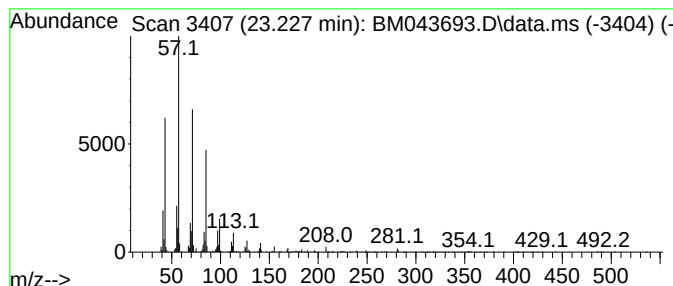
Instrument :  
BNA\_M  
ClientSampleId :  
GCF32

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM121523.MA.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 29 (DEL) Alkane: Straight-Chai... Concentration Rank 20

| R.T.      | EstConc              | Area   | Relative to ISTD |         | R.T.        |      |
|-----------|----------------------|--------|------------------|---------|-------------|------|
| 23.227    | 4.58 ng/ul           | 550853 | Perylene-d12     |         | 23.938      |      |
| Hit# of 5 | Tentative ID         |        | MW               | MolForm | CAS#        | Qual |
| 1         | Tetracosane          |        | 338              | C24H50  | 000646-31-1 | 91   |
| 2         | Heptacosane          |        | 380              | C27H56  | 000593-49-7 | 90   |
| 3         | Octacosane           |        | 394              | C28H58  | 000630-02-4 | 90   |
| 4         | Tetracosane          |        | 338              | C24H50  | 000646-31-1 | 87   |
| 5         | Tetradecane, 1-iodo- |        | 324              | C14H29I | 019218-94-1 | 83   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM010224\  
Data File : BM043693.D  
Acq On : 02 Jan 2024 13:53  
Operator : MA/JU  
Sample : 05918-19  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
GCF32

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM121523.MA.M  
Quant Title : SVOA CALIBRATION

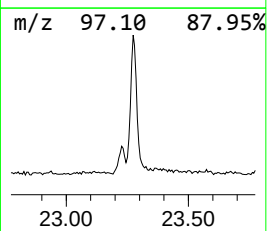
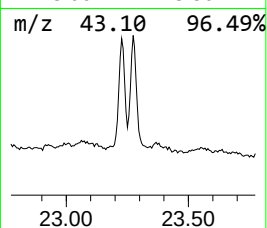
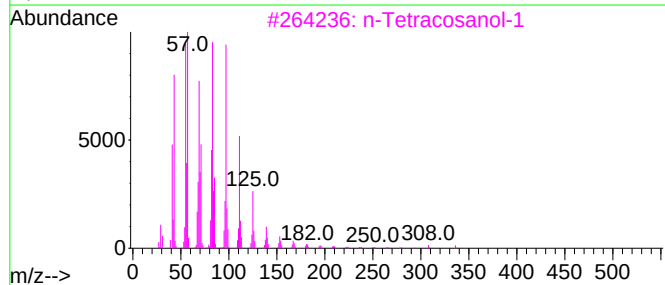
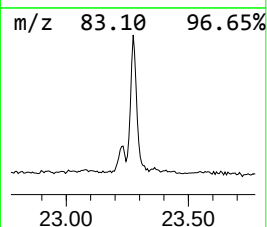
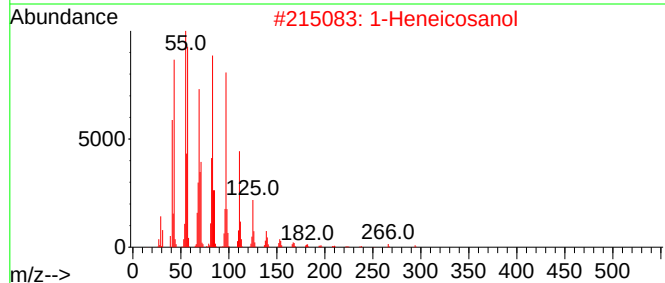
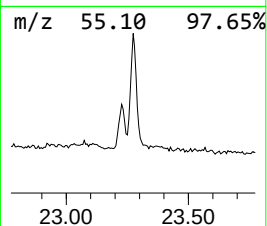
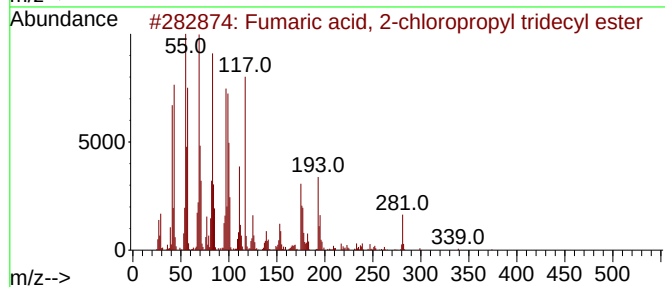
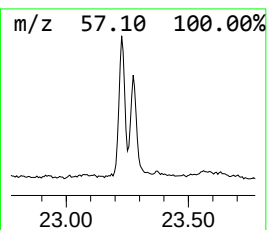
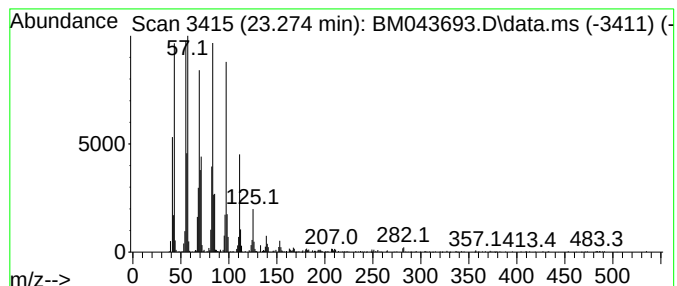
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 30 Fumaric acid, 2-chloropropyl... Concentration Rank 11

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 23.274 | 11.19 ng/ul | 1346820 | Perylene-d12     | 23.938 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|------------|--------------|------|
| 1    |    |   | Fumaric acid, 2-chloropropyl tri... | 374 | C20H35ClO4 | 1010348-57-1 | 98   |
| 2    |    |   | 1-Heneicosanol                      | 312 | C21H44O    | 015594-90-8  | 95   |
| 3    |    |   | n-Tetracosanol-1                    | 354 | C24H50O    | 000506-51-4  | 94   |
| 4    |    |   | Behenic alcohol                     | 326 | C22H46O    | 000661-19-8  | 94   |
| 5    |    |   | 1-Tetracosene                       | 336 | C24H48     | 010192-32-2  | 94   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM010224\  
Data File : BM043693.D  
Acq On : 02 Jan 2024 13:53  
Operator : MA/JU  
Sample : 05918-19  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
GCF32

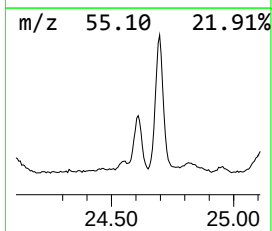
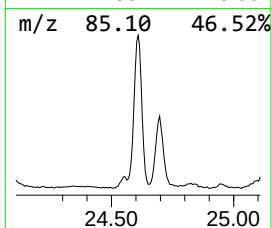
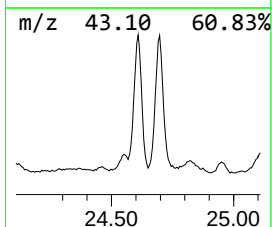
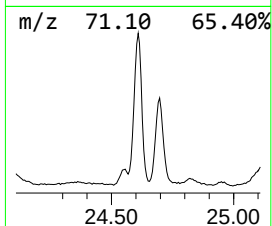
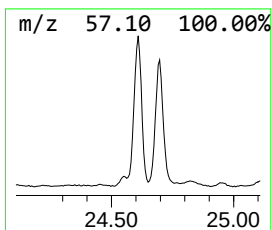
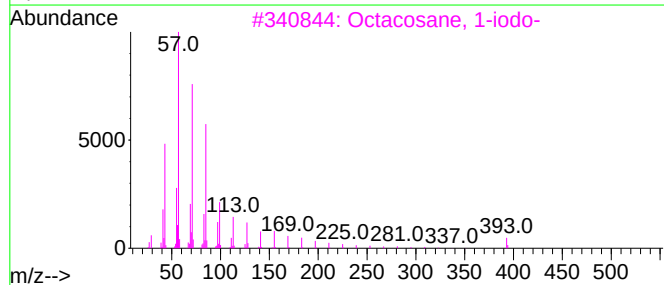
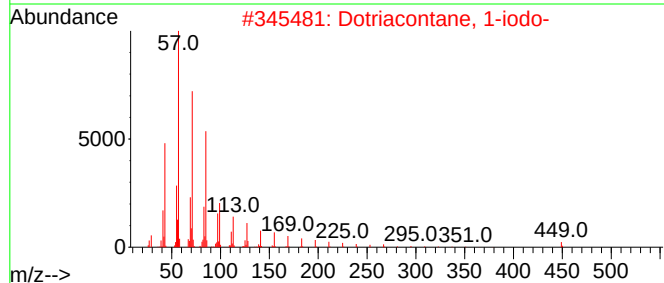
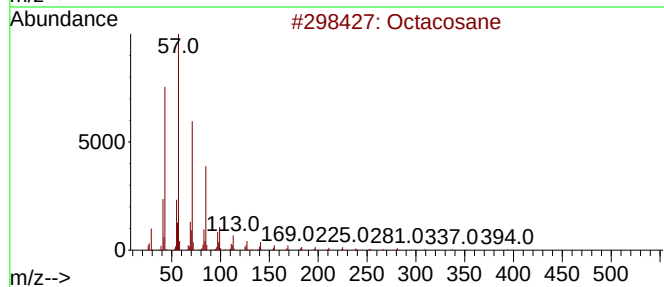
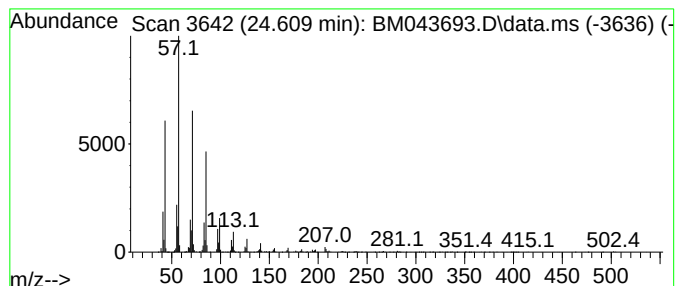
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Quant Title : SVOA CALIBRATION

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

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Peak Number 31 (DEL) Alkane: Straight-Chai... Concentration Rank 7

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 24.609 | 21.49 ng/ul | 2586970 | Perylene-d12     | 23.938 |

| Hit# | of 5 | Tentative ID           | MW  | MolForm | CAS#         | Qual |
|------|------|------------------------|-----|---------|--------------|------|
| 1    |      | Octacosane             | 394 | C28H58  | 000630-02-4  | 91   |
| 2    |      | Dotriacontane, 1-iodo- | 576 | C32H65I | 1000406-32-4 | 91   |
| 3    |      | Octacosane, 1-iodo-    | 520 | C28H57I | 1000406-32-2 | 91   |
| 4    |      | Tetratetracontane      | 619 | C44H90  | 007098-22-8  | 91   |
| 5    |      | Triacontane, 1-iodo-   | 548 | C30H61I | 1000406-32-3 | 91   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM010224\  
Data File : BM043693.D  
Acq On : 02 Jan 2024 13:53  
Operator : MA/JU  
Sample : 05918-19  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

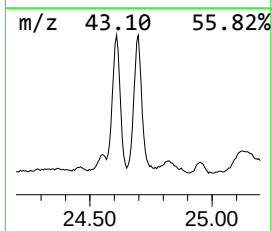
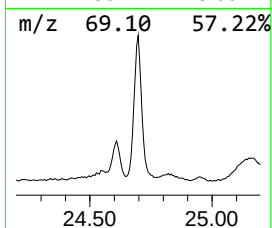
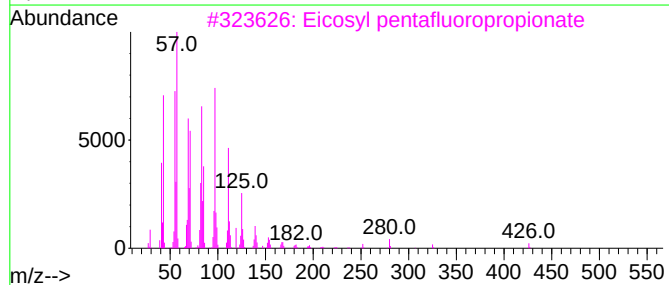
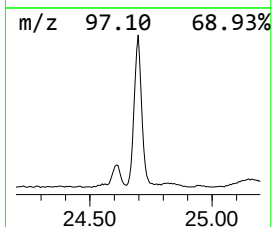
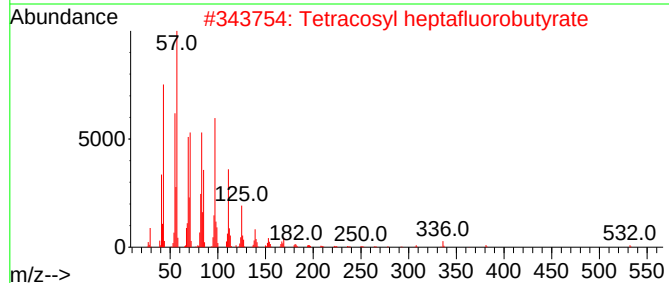
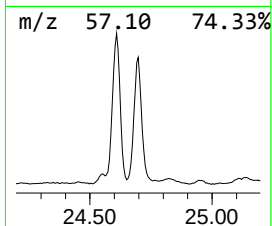
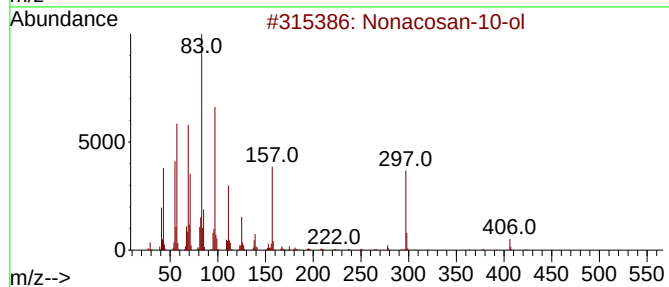
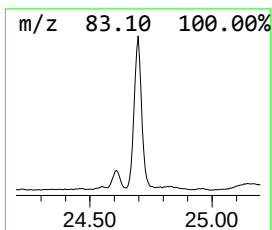
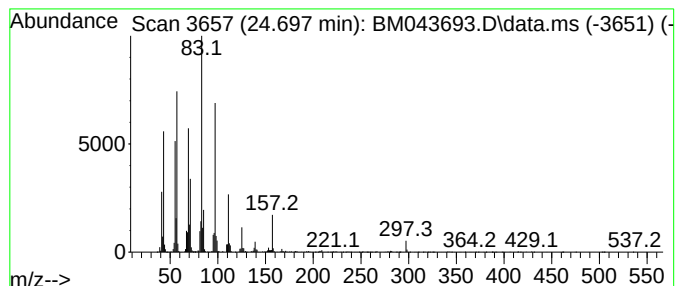
Instrument :  
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ClientSampleId :  
GCF32

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM121523.MA.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 32 Nonacosan-10-ol Concentration Rank 3

| R.T.      | EstConc                        | Area    | Relative to ISTD |            | R.T.         |      |
|-----------|--------------------------------|---------|------------------|------------|--------------|------|
| 24.697    | 34.89 ng/ul                    | 4198840 | Perylene-d12     |            | 23.938       |      |
| Hit# of 5 | Tentative ID                   |         | MW               | MolForm    | CAS#         | Qual |
| 1         | Nonacosan-10-ol                |         | 424              | C29H60O    | 000504-55-2  | 86   |
| 2         | Tetracosyl heptafluorobutyrate |         | 550              | C28H49F7O2 | 1000351-83-7 | 49   |
| 3         | Eicosyl pentafluoropropionate  |         | 444              | C23H41F5O2 | 1000351-80-8 | 46   |
| 4         | Eicosyl heptafluorobutyrate    |         | 494              | C24H41F7O2 | 1000351-83-5 | 46   |
| 5         | Tetracosyl trifluoroacetate    |         | 450              | C26H49F3O2 | 1000351-76-4 | 43   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM010224\  
Data File : BM043693.D  
Acq On : 02 Jan 2024 13:53  
Operator : MA/JU  
Sample : 05918-19  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

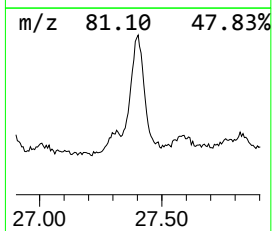
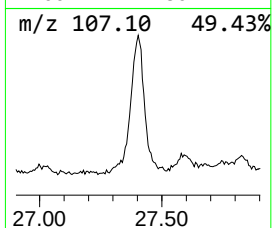
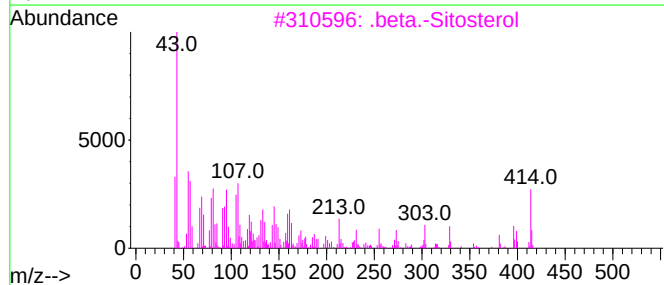
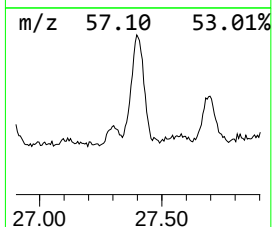
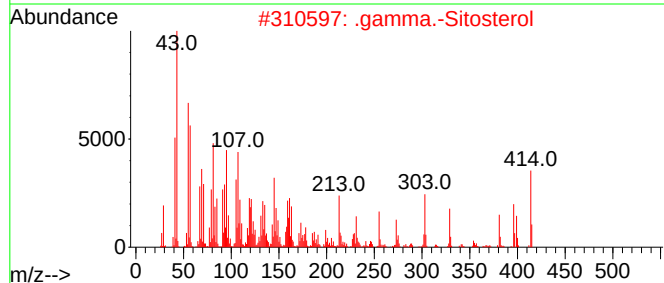
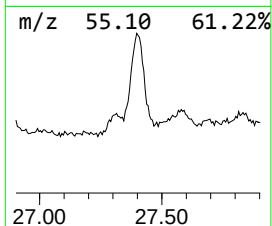
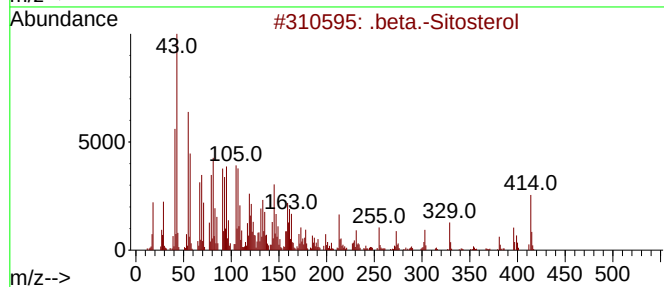
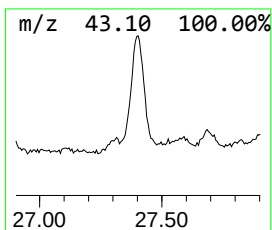
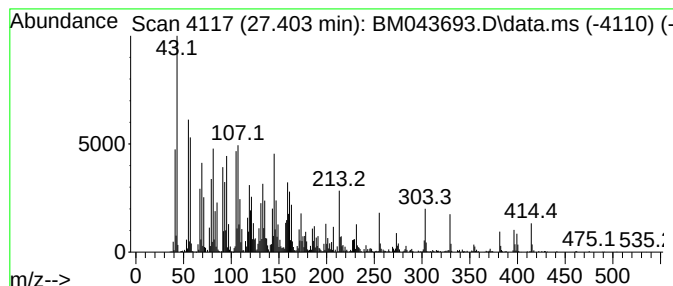
Instrument :  
BNA\_M  
ClientSampleId :  
GCF32

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM121523.MA.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 33 .beta.-Sitosterol Concentration Rank 4

| R.T.      | EstConc            | Area    | Relative to ISTD | R.T.        |      |
|-----------|--------------------|---------|------------------|-------------|------|
| 27.403    | 32.08 ng/ul        | 3860620 | Perylene-d12     | 23.938      |      |
| Hit# of 5 | Tentative ID       | MW      | MolForm          | CAS#        | Qual |
| 1         | .beta.-Sitosterol  | 414     | C29H50O          | 000083-46-5 | 96   |
| 2         | .gamma.-Sitosterol | 414     | C29H50O          | 000083-47-6 | 95   |
| 3         | .beta.-Sitosterol  | 414     | C29H50O          | 000083-46-5 | 91   |
| 4         | .gamma.-Sitosterol | 414     | C29H50O          | 000083-47-6 | 83   |
| 5         | Campesterol        | 400     | C28H48O          | 000474-62-4 | 64   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM010224\  
 Data File : BM043693.D  
 Acq On : 02 Jan 2024 13:53  
 Operator : MA/JU  
 Sample : 05918-19  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 GCF32

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\SFAM-EPA-BM121523.MA.M  
 Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST0.L  
 TIC Integration Parameters: LSCINT.P

| TIC Top Hit name   | RT     | EstConc | Units | Response | --Internal Standard-- |        |         |      |
|--------------------|--------|---------|-------|----------|-----------------------|--------|---------|------|
|                    |        |         |       |          | #                     | RT     | Resp    | Conc |
| 2-Butanone         | 5.304  | 2.7     | ng/ul | 168298   | 1                     | 7.957  | 1227680 | 20.0 |
| Bicyclo[3.1.0]h... | 6.463  | 2.8     | ng/ul | 172391   | 1                     | 7.957  | 1227680 | 20.0 |
| 1,4-Methano-1H-... | 13.598 | 2.0     | ng/ul | 210275   | 3                     | 14.592 | 2079590 | 20.0 |
| Longifolene        | 13.857 | 23.2    | ng/ul | 2415780  | 3                     | 14.592 | 2079590 | 20.0 |
| Cyclohexene, 3-... | 13.904 | 3.9     | ng/ul | 405015   | 3                     | 14.592 | 2079590 | 20.0 |
| cis-Thujopsene     | 14.110 | 2.8     | ng/ul | 290133   | 3                     | 14.592 | 2079590 | 20.0 |
| Cyclohexene, 4-... | 14.410 | 4.7     | ng/ul | 485743   | 3                     | 14.592 | 2079590 | 20.0 |
| (1R,4R,5S)-1,8-... | 14.480 | 7.8     | ng/ul | 813146   | 3                     | 14.592 | 2079590 | 20.0 |
| Cedrol             | 15.874 | 12.1    | ng/ul | 1253980  | 3                     | 14.592 | 2079590 | 20.0 |
| (E)-4-(3-Hydrox... | 16.739 | 2.4     | ng/ul | 263461   | 4                     | 17.339 | 2219230 | 20.0 |
| unknown-01         | 17.086 | 6.6     | ng/ul | 737270   | 4                     | 17.339 | 2219230 | 20.0 |
| (DEL) Alkane: C... | 17.568 | 2.1     | ng/ul | 236269   | 4                     | 17.339 | 2219230 | 20.0 |
| (E)-Hexadec-9-e... | 18.121 | 2.4     | ng/ul | 264967   | 4                     | 17.339 | 2219230 | 20.0 |
| Palmitoleic acid   | 18.186 | 9.3     | ng/ul | 1034360  | 4                     | 17.339 | 2219230 | 20.0 |
| n-Hexadecanoic ... | 18.245 | 26.2    | ng/ul | 2910950  | 4                     | 17.339 | 2219230 | 20.0 |
| Phenanthrene, 1... | 19.203 | 3.0     | ng/ul | 327639   | 4                     | 17.339 | 2219230 | 20.0 |
| (DEL) Alkane: S... | 19.403 | 2.3     | ng/ul | 257617   | 4                     | 17.339 | 2219230 | 20.0 |
| Octadecanoic acid  | 19.521 | 6.0     | ng/ul | 795008   | 5                     | 21.521 | 2655530 | 20.0 |
| unknown-02         | 19.598 | 12.8    | ng/ul | 1695590  | 5                     | 21.521 | 2655530 | 20.0 |
| (DEL) Alkane: S... | 20.250 | 4.3     | ng/ul | 564896   | 5                     | 21.521 | 2655530 | 20.0 |
| 2-Phenanthrenol... | 20.439 | 8.9     | ng/ul | 1183270  | 5                     | 21.521 | 2655530 | 20.0 |
| ((1R,4aR,4bS,7R... | 20.521 | 2.0     | ng/ul | 271643   | 5                     | 21.521 | 2655530 | 20.0 |
| unknown-03         | 20.609 | 63.9    | ng/ul | 8478590  | 5                     | 21.521 | 2655530 | 20.0 |
| unknown-04         | 20.650 | 21.3    | ng/ul | 2831530  | 5                     | 21.521 | 2655530 | 20.0 |
| ((1R,4aR,4bR,10... | 21.115 | 5.2     | ng/ul | 689860   | 5                     | 21.521 | 2655530 | 20.0 |
| (DEL) Alkane: C... | 21.239 | 9.9     | ng/ul | 1319740  | 5                     | 21.521 | 2655530 | 20.0 |
| unknown-05         | 21.339 | 41.1    | ng/ul | 5457440  | 5                     | 21.521 | 2655530 | 20.0 |
| 9(1H)-Phenanthr... | 22.103 | 3.1     | ng/ul | 409187   | 5                     | 21.521 | 2655530 | 20.0 |
| (DEL) Alkane: S... | 23.227 | 4.6     | ng/ul | 550853   | 6                     | 23.938 | 2407170 | 20.0 |
| Fumaric acid, 2... | 23.274 | 11.2    | ng/ul | 1346820  | 6                     | 23.938 | 2407170 | 20.0 |
| (DEL) Alkane: S... | 24.609 | 21.5    | ng/ul | 2586970  | 6                     | 23.938 | 2407170 | 20.0 |
| Nonacosan-10-ol    | 24.697 | 34.9    | ng/ul | 4198840  | 6                     | 23.938 | 2407170 | 20.0 |
| .beta.-Sitosterol  | 27.403 | 32.1    | ng/ul | 3860620  | 6                     | 23.938 | 2407170 | 20.0 |