

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG121721\  
 Data File : BG051641.D  
 Acq On : 18 Dec 2021 22:49  
 Operator : CG/JU  
 Sample : M5153-06  
 Misc :  
 ALS Vial : 57 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 BGL71

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
 Title : SVOA CALIBRATION

Signal : TIC: BG051641.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         | 7.408       | 660           | 667         | 676          | rVV      | 94562          | 181437        | 3.21%           | 0.632%        |
| 2         | 7.578       | 690           | 696         | 703          | rBV      | 62699          | 112819        | 2.00%           | 0.393%        |
| 3         | 7.801       | 729           | 734         | 741          | rVV      | 83797          | 147209        | 2.61%           | 0.513%        |
| 4         | 8.277       | 805           | 815         | 821          | rBV      | 134043         | 250882        | 4.44%           | 0.874%        |
| 5         | 8.471       | 843           | 848         | 853          | rBV2     | 29464          | 49968         | 0.88%           | 0.174%        |
| 6         | 8.965       | 926           | 932         | 941          | rVB2     | 114837         | 202983        | 3.59%           | 0.707%        |
| 7         | 9.446       | 1008          | 1014        | 1020         | rVV      | 84468          | 155960        | 2.76%           | 0.544%        |
| 8         | 10.175      | 1131          | 1138        | 1144         | rBV2     | 73825          | 143364        | 2.54%           | 0.500%        |
| 9         | 10.721      | 1223          | 1231        | 1240         | rVB2     | 107381         | 210365        | 3.72%           | 0.733%        |
| 10        | 11.115      | 1285          | 1298        | 1302         | rBV      | 183648         | 395956        | 7.01%           | 1.380%        |
| 11        | 11.162      | 1302          | 1306        | 1312         | rVV      | 135454         | 251772        | 4.46%           | 0.877%        |
| 12        | 11.232      | 1312          | 1318        | 1321         | rVV      | 37354          | 68039         | 1.20%           | 0.237%        |
| 13        | 12.119      | 1463          | 1469        | 1474         | rBV4     | 29350          | 53382         | 0.94%           | 0.186%        |
| 14        | 12.507      | 1527          | 1535        | 1540         | rBV3     | 53110          | 89892         | 1.59%           | 0.313%        |
| 15        | 12.730      | 1567          | 1573        | 1575         | rVV      | 50399          | 81160         | 1.44%           | 0.283%        |
| 16        | 12.754      | 1575          | 1577        | 1583         | rVB      | 50189          | 62718         | 1.11%           | 0.219%        |
| 17        | 12.971      | 1604          | 1614        | 1622         | rBV2     | 139685         | 253269        | 4.48%           | 0.883%        |
| 18        | 13.353      | 1674          | 1679        | 1683         | rVV2     | 46185          | 76831         | 1.36%           | 0.268%        |
| 19        | 13.447      | 1691          | 1695        | 1702         | rVV2     | 31993          | 49507         | 0.88%           | 0.173%        |
| 20        | 13.723      | 1736          | 1742        | 1750         | rBV3     | 81186          | 181299        | 3.21%           | 0.632%        |
| 21        | 13.946      | 1776          | 1780        | 1787         | rVB3     | 27659          | 68269         | 1.21%           | 0.238%        |
| 22        | 14.069      | 1796          | 1801        | 1803         | rBV3     | 57115          | 85022         | 1.50%           | 0.296%        |
| 23        | 14.234      | 1823          | 1829        | 1834         | rBV3     | 107782         | 196802        | 3.48%           | 0.686%        |
| 24        | 14.298      | 1834          | 1840        | 1845         | rVB2     | 244938         | 389494        | 6.89%           | 1.357%        |
| 25        | 14.381      | 1851          | 1854        | 1858         | rVB2     | 39990          | 49645         | 0.88%           | 0.173%        |
| 26        | 14.469      | 1864          | 1869        | 1872         | rBV      | 39846          | 64699         | 1.15%           | 0.225%        |
| 27        | 14.604      | 1887          | 1892        | 1900         | rBV      | 211108         | 340876        | 6.03%           | 1.188%        |
| 28        | 14.792      | 1920          | 1924        | 1932         | rVV2     | 34846          | 66182         | 1.17%           | 0.231%        |
| 29        | 14.909      | 1939          | 1944        | 1950         | rVB      | 311437         | 485741        | 8.60%           | 1.693%        |
| 30        | 15.068      | 1966          | 1971        | 1975         | rBV2     | 79683          | 127488        | 2.26%           | 0.444%        |
| 31        | 15.115      | 1975          | 1979        | 1986         | rVB2     | 30698          | 56924         | 1.01%           | 0.198%        |
| 32        | 15.297      | 2005          | 2010        | 2016         | rVB3     | 59146          | 110818        | 1.96%           | 0.386%        |
| 33        | 15.374      | 2016          | 2023        | 2027         | rBV2     | 36335          | 72834         | 1.29%           | 0.254%        |
| 34        | 15.573      | 2053          | 2057        | 2063         | rVB4     | 31698          | 44273         | 0.78%           | 0.154%        |
| 35        | 15.691      | 2071          | 2077        | 2083         | rBV2     | 148169         | 266811        | 4.72%           | 0.930%        |
| 36        | 15.843      | 2099          | 2103        | 2107         | rVV      | 30870          | 52714         | 0.93%           | 0.184%        |

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 Sample : M5153-06  
 Misc :  
 ALS Vial : 57 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 BGL71

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
 Title : SVOA CALIBRATION

|    |        |      |      |      |      |         |         |        |        |
|----|--------|------|------|------|------|---------|---------|--------|--------|
| 37 | 15.896 | 2107 | 2112 | 2117 | rVB  | 403541  | 579503  | 10.26% | 2.020% |
| 38 | 15.996 | 2125 | 2129 | 2133 | rBV2 | 59125   | 77918   | 1.38%  | 0.272% |
| 39 | 16.084 | 2141 | 2144 | 2148 | rVV3 | 50019   | 66876   | 1.18%  | 0.233% |
| 40 | 16.149 | 2150 | 2155 | 2159 | rVV  | 114225  | 179148  | 3.17%  | 0.624% |
| 41 | 16.184 | 2159 | 2161 | 2166 | rVB3 | 39176   | 52204   | 0.92%  | 0.182% |
| 42 | 16.243 | 2167 | 2171 | 2178 | rVB8 | 25276   | 53904   | 0.95%  | 0.188% |
| 43 | 16.313 | 2179 | 2183 | 2188 | rBV4 | 30269   | 55556   | 0.98%  | 0.194% |
| 44 | 16.396 | 2194 | 2197 | 2201 | rVB2 | 52786   | 61237   | 1.08%  | 0.213% |
| 45 | 16.625 | 2228 | 2236 | 2241 | rBV3 | 132796  | 257937  | 4.57%  | 0.899% |
| 46 | 16.725 | 2248 | 2253 | 2257 | rBV  | 182483  | 277975  | 4.92%  | 0.969% |
| 47 | 16.777 | 2258 | 2262 | 2269 | rVB4 | 141639  | 278539  | 4.93%  | 0.971% |
| 48 | 16.842 | 2269 | 2273 | 2276 | rBV  | 128990  | 168660  | 2.99%  | 0.588% |
| 49 | 16.883 | 2276 | 2280 | 2284 | rVB2 | 81287   | 107080  | 1.90%  | 0.373% |
| 50 | 16.924 | 2284 | 2287 | 2291 | rBV4 | 27223   | 54733   | 0.97%  | 0.191% |
| 51 | 17.048 | 2302 | 2308 | 2314 | rVB5 | 101386  | 225194  | 3.99%  | 0.785% |
| 52 | 17.106 | 2314 | 2318 | 2322 | rVV  | 117123  | 160178  | 2.84%  | 0.558% |
| 53 | 17.148 | 2322 | 2325 | 2327 | rVV4 | 35540   | 52979   | 0.94%  | 0.185% |
| 54 | 17.183 | 2327 | 2331 | 2336 | rVV2 | 71072   | 128876  | 2.28%  | 0.449% |
| 55 | 17.230 | 2336 | 2339 | 2345 | rVB  | 39279   | 61597   | 1.09%  | 0.215% |
| 56 | 17.453 | 2373 | 2377 | 2380 | rVB4 | 31256   | 48792   | 0.86%  | 0.170% |
| 57 | 17.541 | 2387 | 2392 | 2398 | rVB  | 58491   | 101515  | 1.80%  | 0.354% |
| 58 | 17.659 | 2406 | 2412 | 2416 | rBV  | 364139  | 580674  | 10.28% | 2.024% |
| 59 | 17.700 | 2416 | 2419 | 2422 | rVV  | 43144   | 58904   | 1.04%  | 0.205% |
| 60 | 17.753 | 2423 | 2428 | 2434 | rVB  | 310645  | 492197  | 8.71%  | 1.715% |
| 61 | 17.935 | 2455 | 2459 | 2466 | rBV2 | 45056   | 84598   | 1.50%  | 0.295% |
| 62 | 18.005 | 2466 | 2471 | 2484 | rBV  | 1006124 | 1542648 | 27.30% | 5.376% |
| 63 | 18.134 | 2490 | 2493 | 2499 | rVB4 | 43251   | 62403   | 1.10%  | 0.217% |
| 64 | 18.305 | 2518 | 2522 | 2526 | rVB3 | 35533   | 49786   | 0.88%  | 0.174% |
| 65 | 18.522 | 2556 | 2559 | 2564 | rVB4 | 41418   | 58856   | 1.04%  | 0.205% |
| 66 | 18.663 | 2579 | 2583 | 2587 | rBV  | 40887   | 74053   | 1.31%  | 0.258% |
| 67 | 18.716 | 2589 | 2592 | 2599 | rVB6 | 31190   | 62396   | 1.10%  | 0.217% |
| 68 | 19.010 | 2636 | 2642 | 2647 | rBV3 | 74120   | 133041  | 2.35%  | 0.464% |
| 69 | 19.198 | 2669 | 2674 | 2681 | rBV  | 241580  | 312471  | 5.53%  | 1.089% |
| 70 | 19.627 | 2743 | 2747 | 2751 | rVB2 | 72987   | 93429   | 1.65%  | 0.326% |
| 71 | 19.750 | 2764 | 2768 | 2771 | rBV4 | 38604   | 53030   | 0.94%  | 0.185% |
| 72 | 19.897 | 2789 | 2793 | 2802 | rVB3 | 107714  | 207671  | 3.68%  | 0.724% |
| 73 | 20.032 | 2810 | 2816 | 2826 | rVB  | 900031  | 1280933 | 22.67% | 4.464% |
| 74 | 20.807 | 2944 | 2948 | 2952 | rVB  | 73478   | 85504   | 1.51%  | 0.298% |
| 75 | 20.931 | 2965 | 2969 | 2977 | rVB  | 141400  | 204887  | 3.63%  | 0.714% |
| 76 | 21.260 | 3021 | 3025 | 3028 | rBV  | 82499   | 106097  | 1.88%  | 0.370% |

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Instrument :  
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 ClientSampleId :  
 BGL71

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
 Title : SVOA CALIBRATION

|     |        |      |      |      |      |         |         |         |         |
|-----|--------|------|------|------|------|---------|---------|---------|---------|
| 77  | 21.301 | 3028 | 3032 | 3040 | rVB  | 227409  | 309544  | 5.48%   | 1.079%  |
| 78  | 21.424 | 3044 | 3053 | 3057 | rBV2 | 248242  | 466635  | 8.26%   | 1.626%  |
| 79  | 21.471 | 3057 | 3061 | 3067 | rVB  | 141407  | 218064  | 3.86%   | 0.760%  |
| 80  | 21.589 | 3077 | 3081 | 3088 | rVB3 | 71184   | 130261  | 2.31%   | 0.454%  |
| 81  | 21.706 | 3098 | 3101 | 3106 | rVB6 | 45350   | 69018   | 1.22%   | 0.241%  |
| 82  | 21.794 | 3111 | 3116 | 3119 | rVV  | 74569   | 129322  | 2.29%   | 0.451%  |
| 83  | 21.853 | 3119 | 3126 | 3140 | rVV  | 3290875 | 5649933 | 100.00% | 19.690% |
| 84  | 21.970 | 3140 | 3146 | 3154 | rVB3 | 433890  | 863464  | 15.28%  | 3.009%  |
| 85  | 22.147 | 3172 | 3176 | 3178 | rBV2 | 54363   | 78908   | 1.40%   | 0.275%  |
| 86  | 22.188 | 3179 | 3183 | 3195 | rVB5 | 118519  | 307558  | 5.44%   | 1.072%  |
| 87  | 22.393 | 3216 | 3218 | 3225 | rVB3 | 45203   | 72765   | 1.29%   | 0.254%  |
| 88  | 22.540 | 3237 | 3243 | 3246 | rBV2 | 37266   | 82245   | 1.46%   | 0.287%  |
| 89  | 22.593 | 3247 | 3252 | 3258 | rVV  | 154435  | 315844  | 5.59%   | 1.101%  |
| 90  | 22.669 | 3259 | 3265 | 3273 | rVB  | 339461  | 623871  | 11.04%  | 2.174%  |
| 91  | 22.834 | 3291 | 3293 | 3300 | rVB3 | 41235   | 65733   | 1.16%   | 0.229%  |
| 92  | 23.169 | 3342 | 3350 | 3364 | rVB  | 450628  | 1215190 | 21.51%  | 4.235%  |
| 93  | 23.545 | 3410 | 3414 | 3419 | rBV3 | 44417   | 76828   | 1.36%   | 0.268%  |
| 94  | 23.598 | 3420 | 3423 | 3430 | rVB3 | 46231   | 77160   | 1.37%   | 0.269%  |
| 95  | 24.103 | 3502 | 3509 | 3521 | rVB  | 105417  | 273766  | 4.85%   | 0.954%  |
| 96  | 24.273 | 3530 | 3538 | 3546 | rVB2 | 187067  | 482326  | 8.54%   | 1.681%  |
| 97  | 24.890 | 3635 | 3643 | 3660 | rVB  | 333007  | 1264802 | 22.39%  | 4.408%  |
| 98  | 25.231 | 3693 | 3701 | 3713 | rVB2 | 188862  | 593908  | 10.51%  | 2.070%  |
| 99  | 25.472 | 3733 | 3742 | 3752 | rBV2 | 192633  | 583912  | 10.33%  | 2.035%  |
| 100 | 27.287 | 4042 | 4051 | 4067 | rVB  | 166766  | 651455  | 11.53%  | 2.270%  |

Sum of corrected areas: 28693895



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Sample : M5153-06  
Misc :  
ALS Vial : 57 Sample Multiplier: 1

Instrument :  
BNA\_G  
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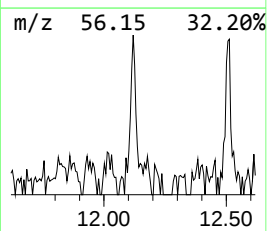
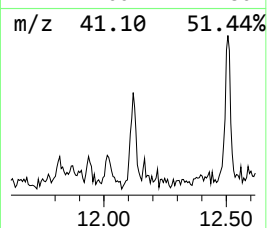
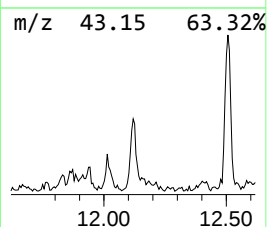
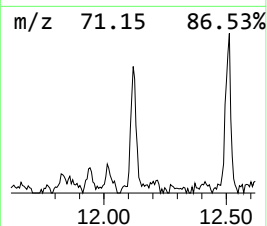
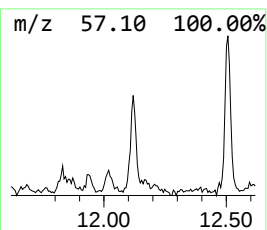
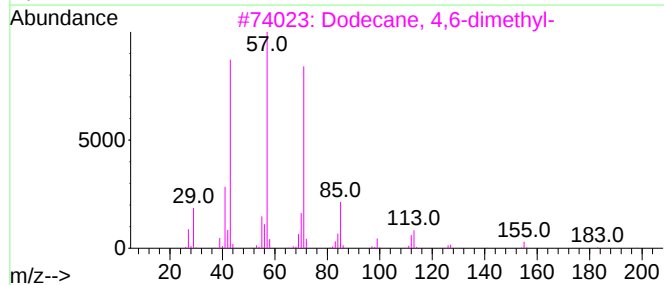
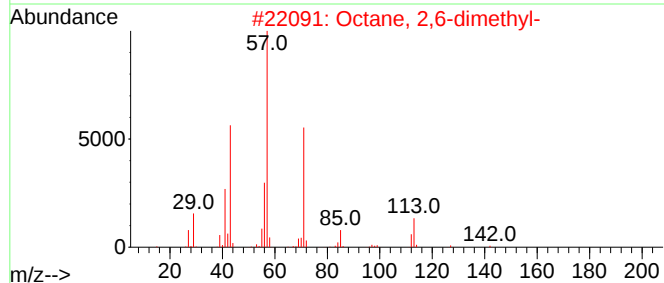
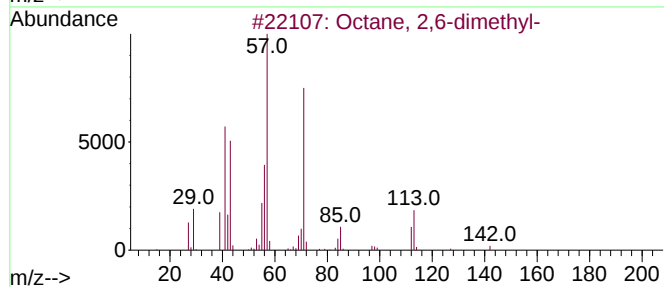
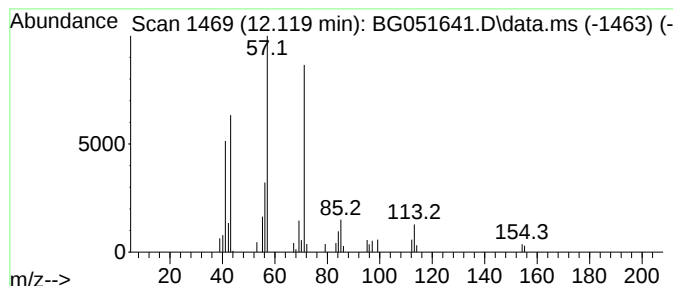
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 39

| R.T.   | EstConc    | Area  | Relative to ISTD | R.T.   |
|--------|------------|-------|------------------|--------|
| 12.119 | 2.70 ng/ul | 53382 | Naphthalene-d8   | 11.115 |

| Hit# | of | 5 | Tentative ID                   | MW  | MolForm  | CAS#         | Qual |
|------|----|---|--------------------------------|-----|----------|--------------|------|
| 1    |    |   | Octane, 2,6-dimethyl-          | 142 | C10H22   | 002051-30-1  | 78   |
| 2    |    |   | Octane, 2,6-dimethyl-          | 142 | C10H22   | 002051-30-1  | 72   |
| 3    |    |   | Dodecane, 4,6-dimethyl-        | 198 | C14H30   | 061141-72-8  | 64   |
| 4    |    |   | Oxalic acid, allyl octyl ester | 242 | C13H22O4 | 1000309-23-6 | 59   |
| 5    |    |   | Oxalic acid, allyl nonyl ester | 256 | C14H24O4 | 1000309-23-7 | 59   |



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Quant Title : SVOA CALIBRATION

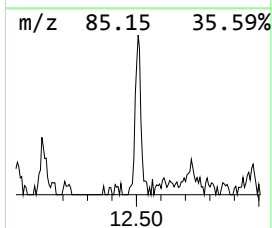
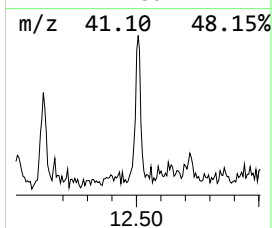
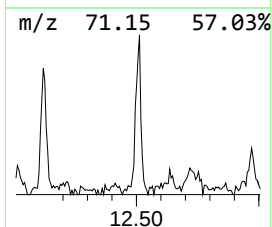
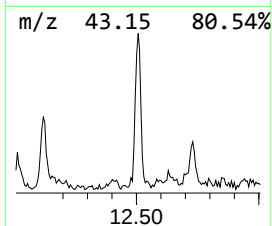
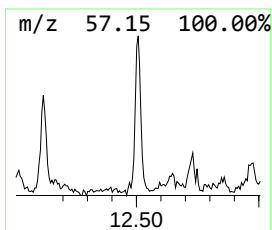
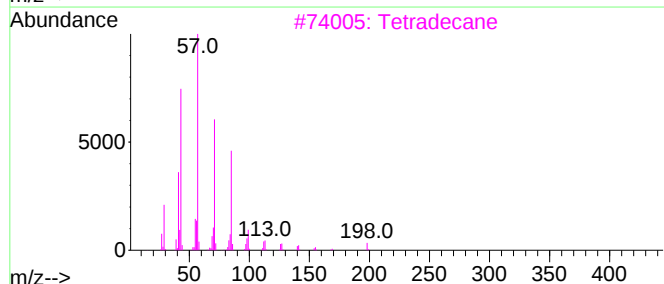
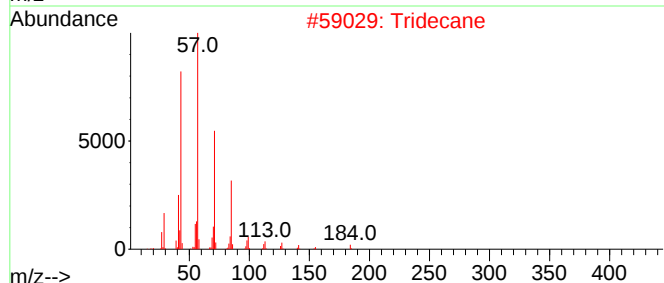
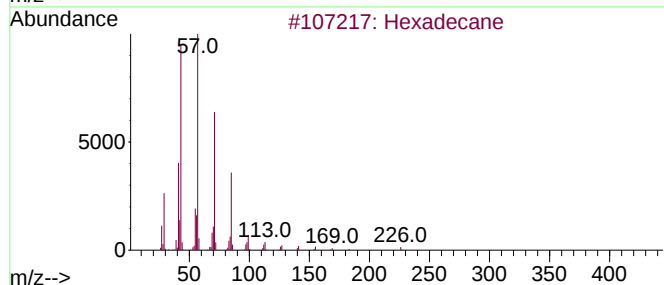
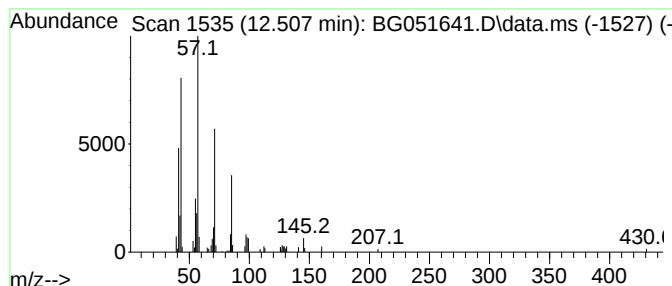
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 24

| R.T.   | EstConc    | Area  | Relative to ISTD | R.T.   |
|--------|------------|-------|------------------|--------|
| 12.507 | 4.54 ng/ul | 89892 | Naphthalene-d8   | 11.115 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Hexadecane                          | 226 | C16H34   | 000544-76-3  | 80   |
| 2    |    |   | Tridecane                           | 184 | C13H28   | 000629-50-5  | 72   |
| 3    |    |   | Tetradecane                         | 198 | C14H30   | 000629-59-4  | 72   |
| 4    |    |   | Carbonic acid, dodecyl vinyl ester  | 256 | C15H28O3 | 1000382-54-8 | 72   |
| 5    |    |   | Carbonic acid, octadecyl prop-1-... | 354 | C22H42O3 | 1000383-11-5 | 72   |



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Misc :  
ALS Vial : 57 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL71

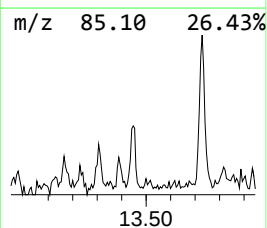
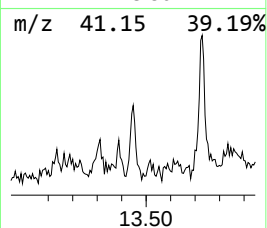
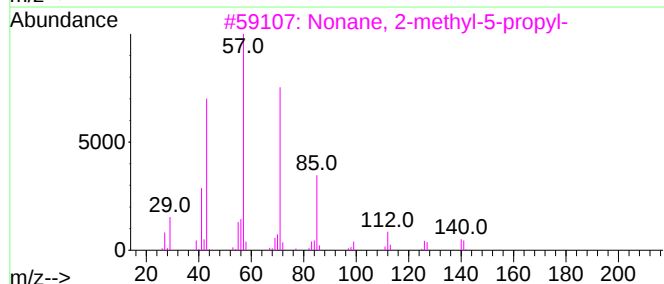
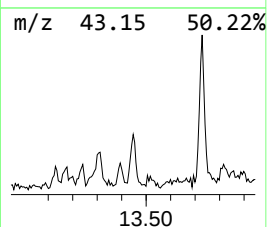
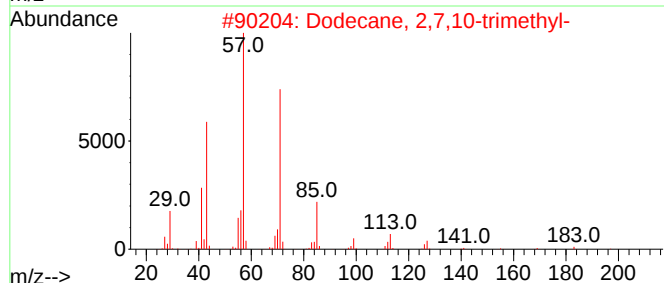
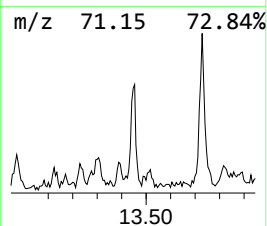
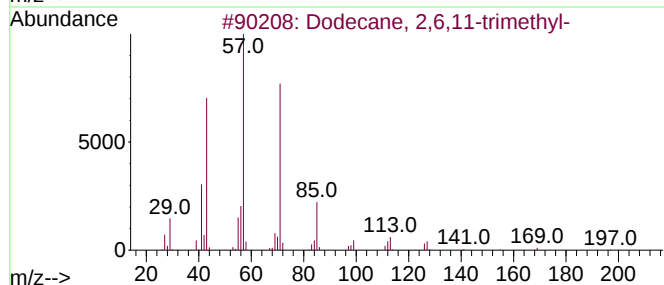
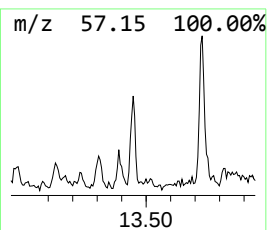
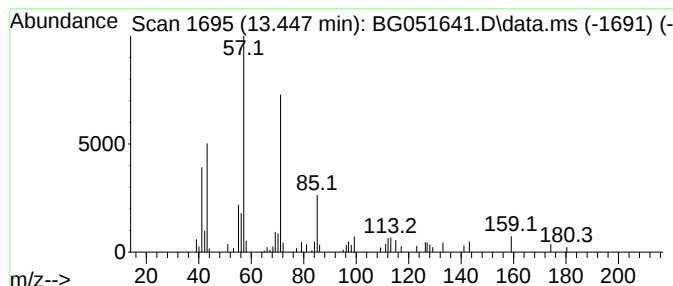
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Quant Title : SVOA CALIBRATION

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

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Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 50

| R.T.   | EstConc    | Area  | Relative to ISTD | R.T.   |
|--------|------------|-------|------------------|--------|
| 13.447 | 2.04 ng/ul | 49507 | Acenaphthene-d10 | 14.909 |

| Hit# | of | 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------------|-----|---------|-------------|------|
| 1    |    |   | Dodecane, 2,6,11-trimethyl- | 212 | C15H32  | 031295-56-4 | 64   |
| 2    |    |   | Dodecane, 2,7,10-trimethyl- | 212 | C15H32  | 074645-98-0 | 64   |
| 3    |    |   | Nonane, 2-methyl-5-propyl-  | 184 | C13H28  | 031081-17-1 | 59   |
| 4    |    |   | Heptacosane                 | 380 | C27H56  | 000593-49-7 | 58   |
| 5    |    |   | Pentadecane                 | 212 | C15H32  | 000629-62-9 | 58   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG121721\  
Data File : BG051641.D  
Acq On : 18 Dec 2021 22:49  
Operator : CG/JU  
Sample : M5153-06  
Misc :  
ALS Vial : 57 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL71

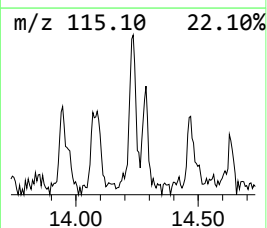
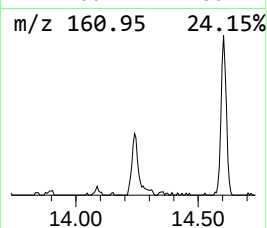
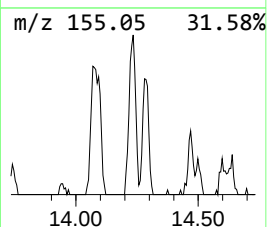
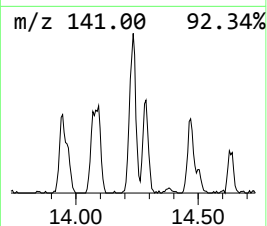
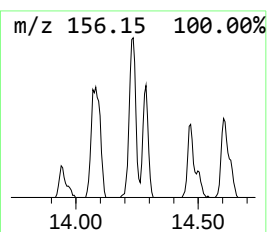
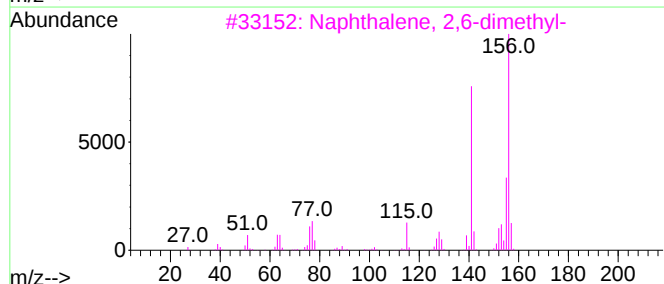
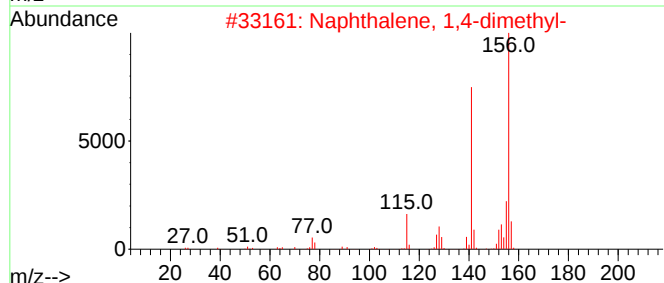
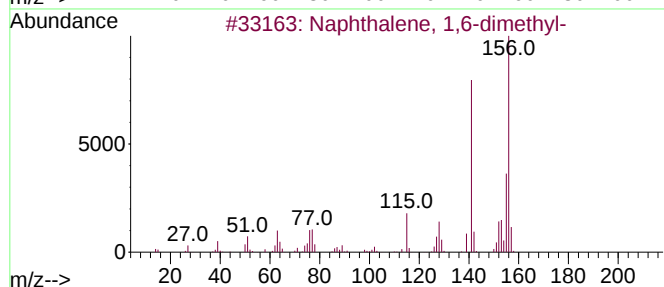
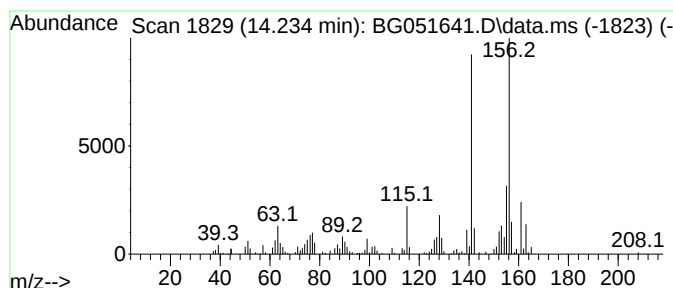
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Quant Title : SVOA CALIBRATION

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

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Peak Number 8 Naphthalene, 1,6-dimethyl- Concentration Rank 13

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 14.234 | 8.10 ng/ul | 196802 | Acenaphthene-d10 | 14.909 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, 1,6-dimethyl- | 156 | C12H12  | 000575-43-9 | 98   |
| 2    |    |   | Naphthalene, 1,4-dimethyl- | 156 | C12H12  | 000571-58-4 | 96   |
| 3    |    |   | Naphthalene, 2,6-dimethyl- | 156 | C12H12  | 000581-42-0 | 96   |
| 4    |    |   | Naphthalene, 1,6-dimethyl- | 156 | C12H12  | 000575-43-9 | 96   |
| 5    |    |   | Naphthalene, 1,7-dimethyl- | 156 | C12H12  | 000575-37-1 | 96   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG121721\  
Data File : BG051641.D  
Acq On : 18 Dec 2021 22:49  
Operator : CG/JU  
Sample : M5153-06  
Misc :  
ALS Vial : 57 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL71

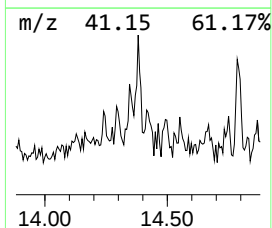
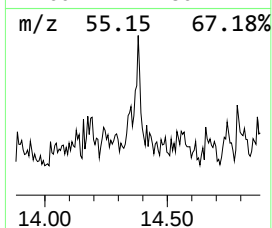
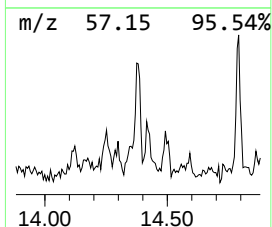
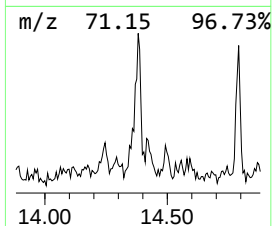
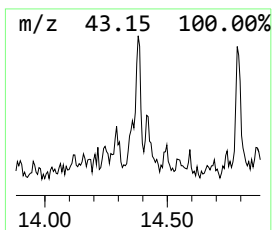
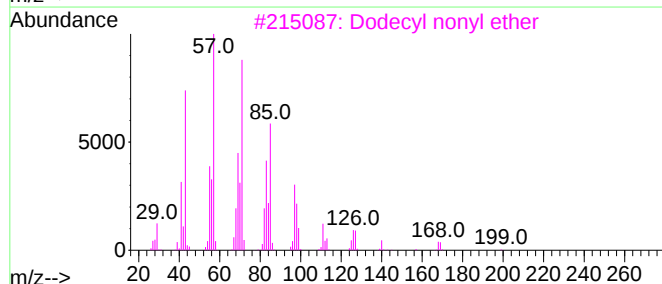
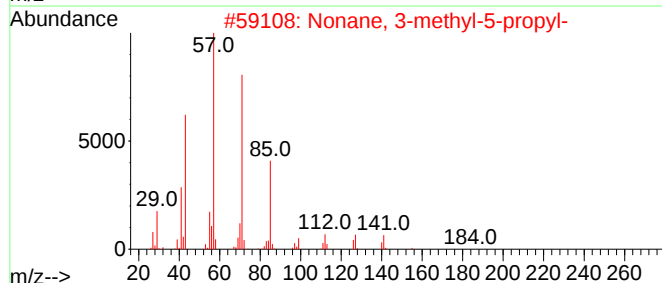
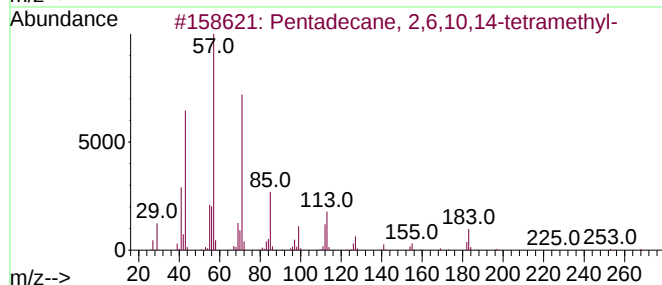
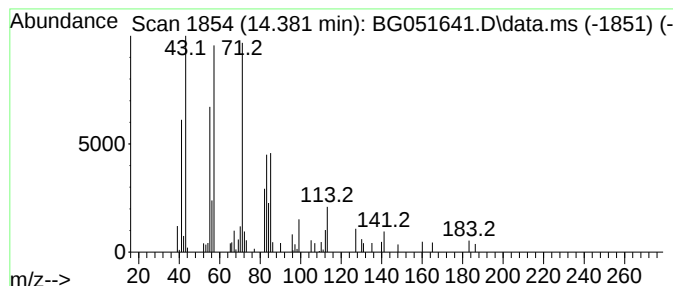
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Quant Title : SVOA CALIBRATION

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

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Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 49

| R.T.      | EstConc                             | Area  | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|-------|------------------|--------------|------|
| 14.381    | 2.04 ng/ul                          | 49645 | Acenaphthene-d10 | 14.909       |      |
| Hit# of 5 | Tentative ID                        | MW    | MolForm          | CAS#         | Qual |
| 1         | Pentadecane, 2,6,10,14-tetramethyl- | 268   | C19H40           | 001921-70-6  | 50   |
| 2         | Nonane, 3-methyl-5-propyl-          | 184   | C13H28           | 031081-18-2  | 50   |
| 3         | Dodecyl nonyl ether                 | 312   | C21H44O          | 1000406-37-5 | 50   |
| 4         | Nonane, 5-methyl-5-propyl-          | 184   | C13H28           | 017312-75-3  | 47   |
| 5         | Succinic acid, nonyl tetrahydrof... | 328   | C18H32O5         | 1000329-86-7 | 43   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG121721\  
Data File : BG051641.D  
Acq On : 18 Dec 2021 22:49  
Operator : CG/JU  
Sample : M5153-06  
Misc :  
ALS Vial : 57 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL71

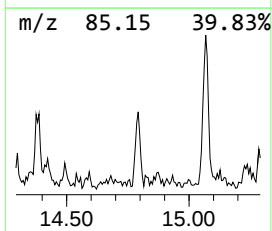
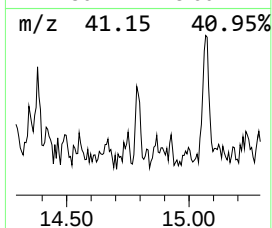
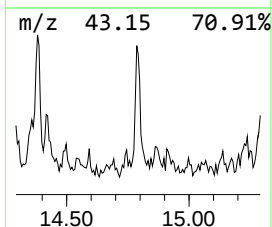
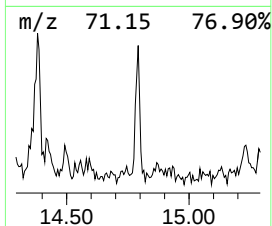
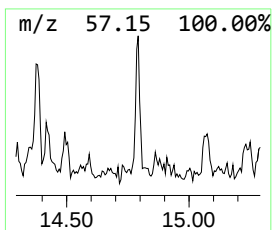
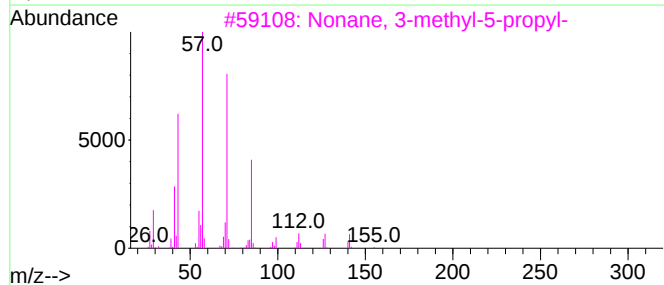
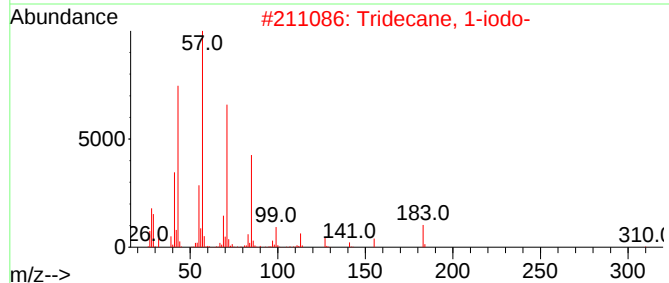
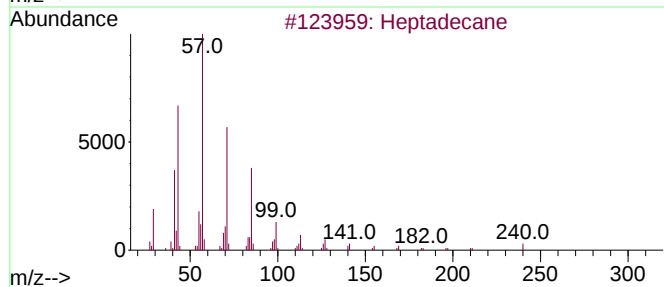
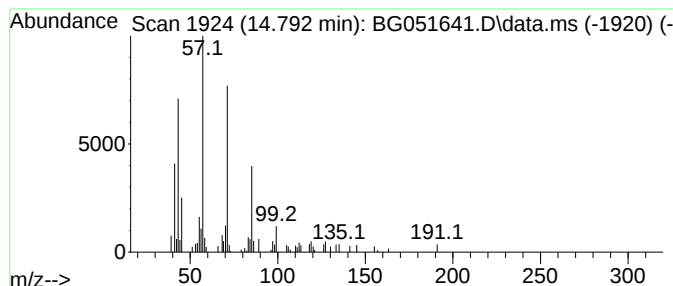
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Quant Title : SVOA CALIBRATION

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

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Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 38

| R.T.      | EstConc                    | Area  | Relative to ISTD | R.T.        |      |
|-----------|----------------------------|-------|------------------|-------------|------|
| 14.792    | 2.72 ng/ul                 | 66182 | Acenaphthene-d10 | 14.909      |      |
| Hit# of 5 | Tentative ID               | MW    | MolForm          | CAS#        | Qual |
| 1         | Heptadecane                | 240   | C17H36           | 000629-78-7 | 72   |
| 2         | Tridecane, 1-iodo-         | 310   | C13H27I          | 035599-77-0 | 72   |
| 3         | Nonane, 3-methyl-5-propyl- | 184   | C13H28           | 031081-18-2 | 64   |
| 4         | Heptacosane                | 380   | C27H56           | 000593-49-7 | 64   |
| 5         | Heneicosane                | 296   | C21H44           | 000629-94-7 | 64   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG121721\  
Data File : BG051641.D  
Acq On : 18 Dec 2021 22:49  
Operator : CG/JU  
Sample : M5153-06  
Misc :  
ALS Vial : 57 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL71

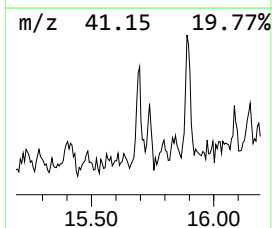
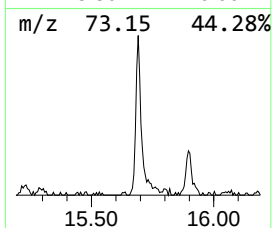
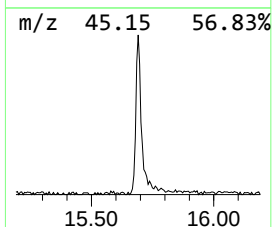
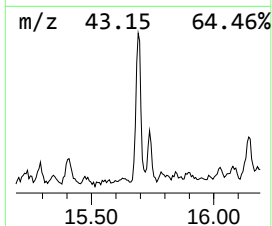
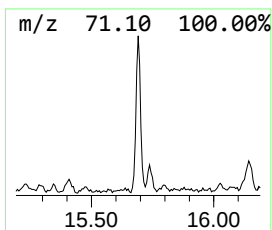
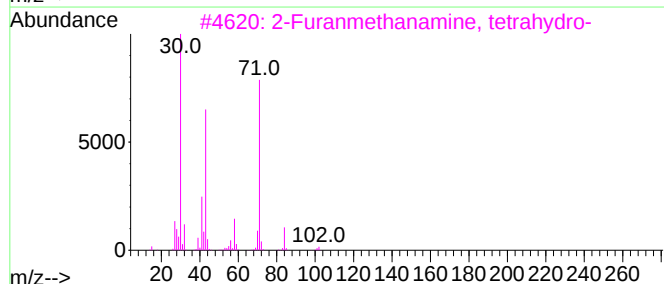
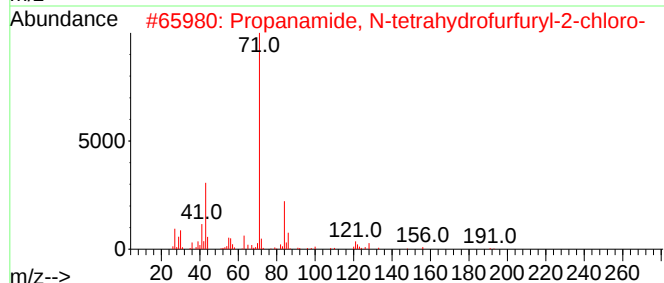
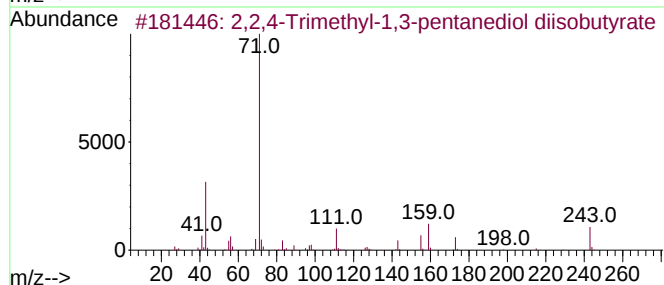
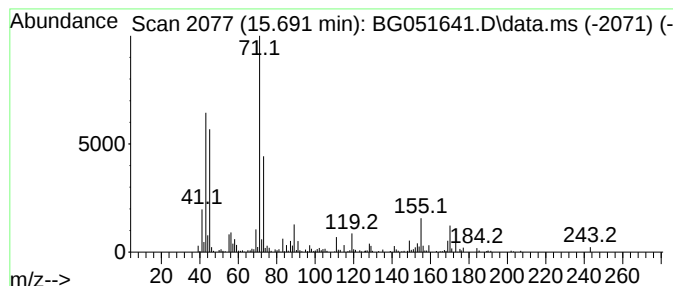
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Quant Title : SVOA CALIBRATION

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

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Peak Number 14 unknown-01 Concentration Rank 6

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 15.691 | 10.99 ng/ul | 266811 | Acenaphthene-d10 | 14.909 |

| Hit# | of                                    | 5   | Tentative ID | MW           | MolForm | CAS# | Qual |
|------|---------------------------------------|-----|--------------|--------------|---------|------|------|
| 1    | 2,2,4-Trimethyl-1,3-pentanediol ...   | 286 | C16H30O4     | 006846-50-0  | 38      |      |      |
| 2    | Propanamide, N-tetrahydrofurfuryl...  | 191 | C8H14ClNO2   | 1000307-48-2 | 35      |      |      |
| 3    | 2-Furanmethanamine, tetrahydro-       | 101 | C5H11NO      | 004795-29-3  | 35      |      |      |
| 4    | Methylamine, N-(1-methylhexylidene... | 127 | C8H17N       | 022058-71-5  | 35      |      |      |
| 5    | Acetamide, N-tetrahydrofurfuryl-      | 211 | C7H11Cl2NO2  | 1000307-29-6 | 35      |      |      |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG121721\  
Data File : BG051641.D  
Acq On : 18 Dec 2021 22:49  
Operator : CG/JU  
Sample : M5153-06  
Misc :  
ALS Vial : 57 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
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Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

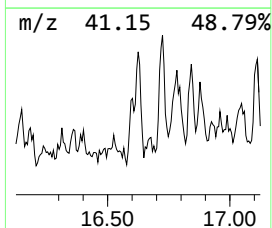
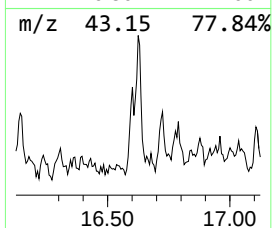
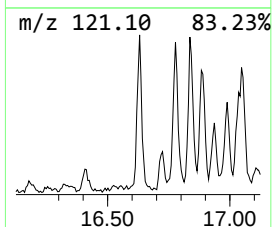
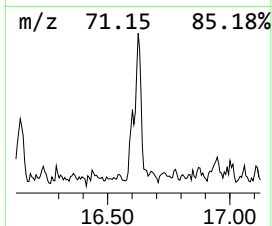
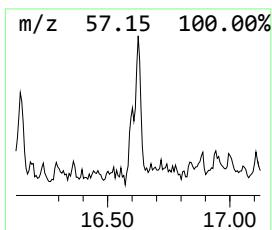
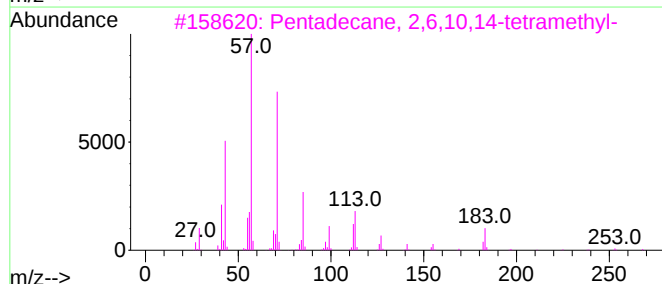
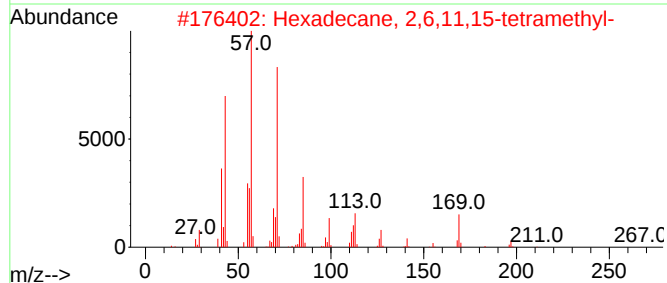
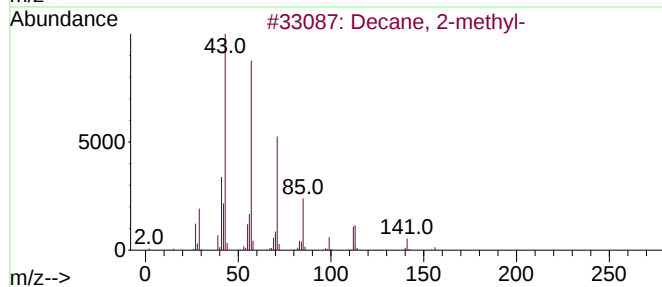
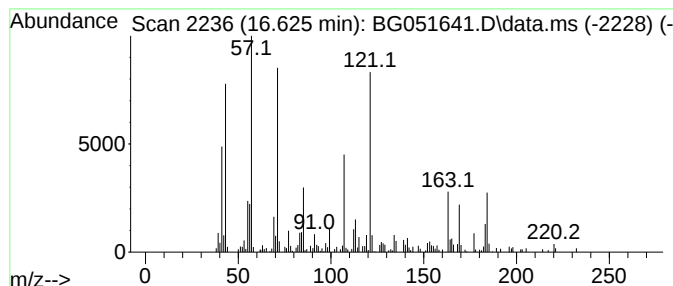
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TIC Integration Parameters: LSCINT.P

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Peak Number 19 (DEL) Alkane: Straight-Chai... Concentration Rank 12

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 16.625 | 8.88 ng/ul | 257937 | Phenanthrene-d10 | 17.659 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Decane, 2-methyl-                   | 156 | C11H24  | 006975-98-0 | 38   |
| 2    |    |   | Hexadecane, 2,6,11,15-tetramethyl-  | 282 | C20H42  | 000504-44-9 | 30   |
| 3    |    |   | Pentadecane, 2,6,10,14-tetramethyl- | 268 | C19H40  | 001921-70-6 | 30   |
| 4    |    |   | Tridecane                           | 184 | C13H28  | 000629-50-5 | 25   |
| 5    |    |   | Decane, 2-methyl-                   | 156 | C11H24  | 006975-98-0 | 25   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG121721\  
Data File : BG051641.D  
Acq On : 18 Dec 2021 22:49  
Operator : CG/JU  
Sample : M5153-06  
Misc :  
ALS Vial : 57 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL71

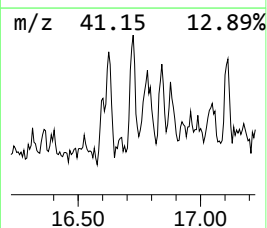
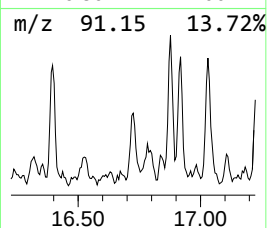
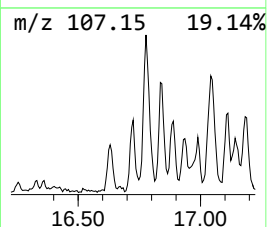
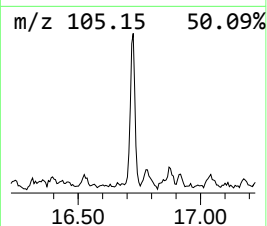
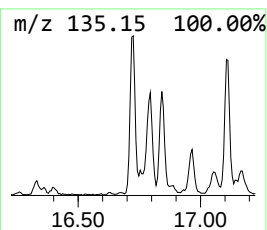
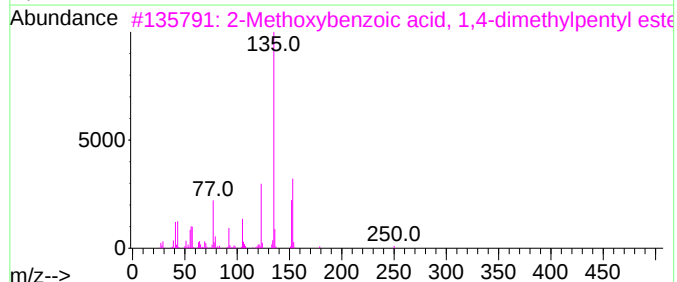
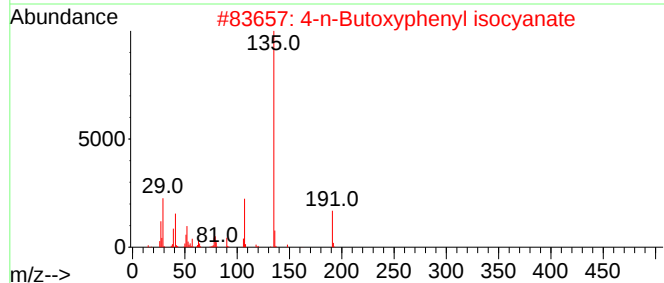
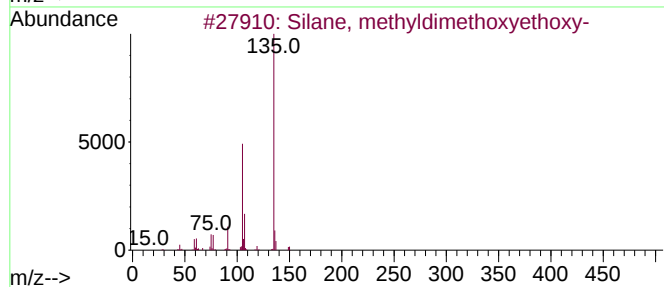
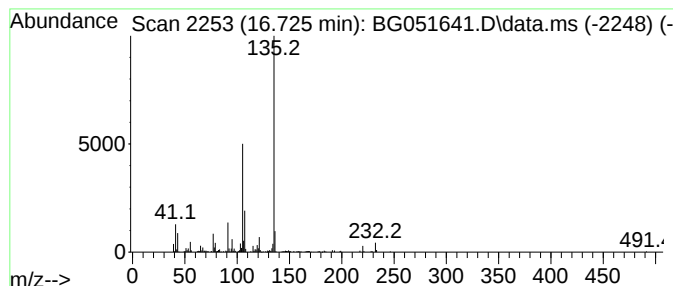
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Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 20 Silane, methylmethoxyethoxy- Concentration Rank 10

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 16.725 | 9.57 ng/ul | 277975 | Phenanthrene-d10 | 17.659 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Silane, methylmethoxyethoxy-        | 150 | C5H14O3Si | 1010416-18-1 | 72   |
| 2    |    |   | 4-n-Butoxyphenyl isocyanate         | 207 | C11H13NOS | 1000342-44-4 | 64   |
| 3    |    |   | 2-Methoxybenzoic acid, 1,4-dimet... | 250 | C15H22O3  | 1000374-94-2 | 64   |
| 4    |    |   | Phenol, 4-(1,1,3,3-tetramethylbu... | 206 | C14H22O   | 000140-66-9  | 64   |
| 5    |    |   | Phenol, 4-(1,1-dimethylpropyl)-     | 164 | C11H16O   | 000080-46-6  | 64   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG121721\  
Data File : BG051641.D  
Acq On : 18 Dec 2021 22:49  
Operator : CG/JU  
Sample : M5153-06  
Misc :  
ALS Vial : 57 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL71

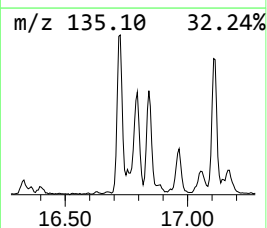
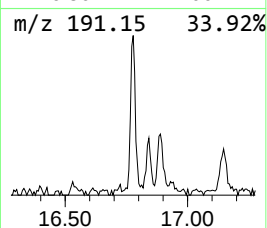
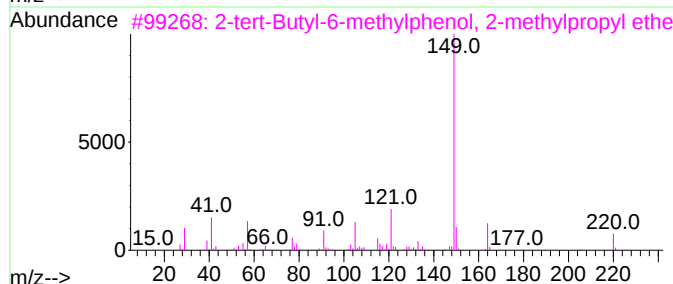
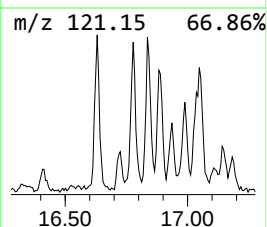
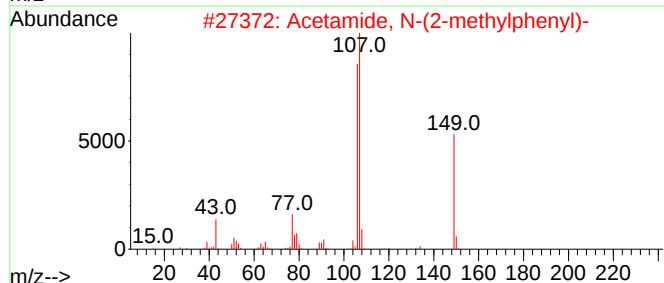
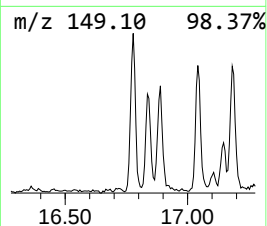
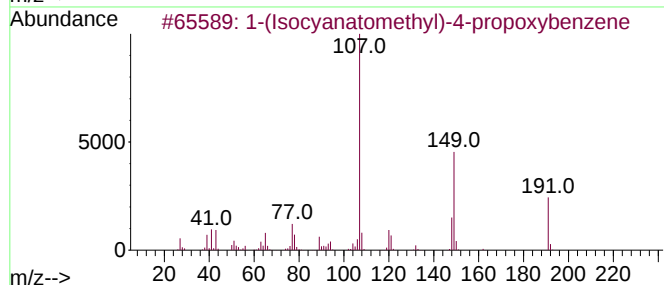
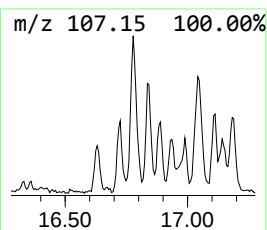
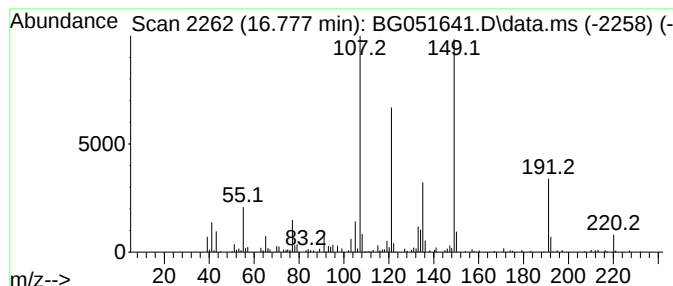
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Quant Title : SVOA CALIBRATION

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

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Peak Number 21 1-(Isocyanatomethyl)-4-prop... Concentration Rank 9

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 16.778 | 9.59 ng/ul | 278539 | Phenanthrene-d10 | 17.659 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | 1-(Isocyanatomethyl)-4-propoxybe... | 191 | C11H13NO2 | 936095-07-7  | 64   |
| 2    |    |   | Acetamide, N-(2-methylphenyl)-      | 149 | C9H11NO   | 000120-66-1  | 38   |
| 3    |    |   | 2-tert-Butyl-6-methylphenol, 2-m... | 220 | C15H24O   | 1000395-19-4 | 38   |
| 4    |    |   | Phthalic acid, 2-(methylthio)phe... | 330 | C18H18O4S | 1000415-55-9 | 35   |
| 5    |    |   | Butanamide, N-(2-methylphenyl)-3... | 191 | C11H13NO2 | 000093-68-5  | 35   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG121721\  
Data File : BG051641.D  
Acq On : 18 Dec 2021 22:49  
Operator : CG/JU  
Sample : M5153-06  
Misc :  
ALS Vial : 57 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL71

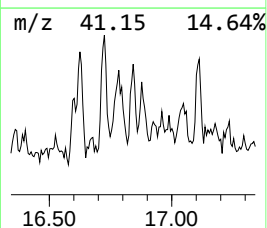
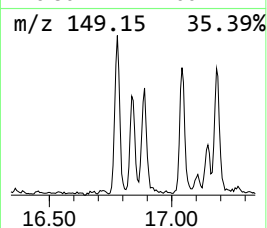
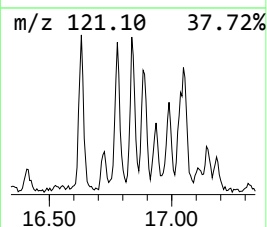
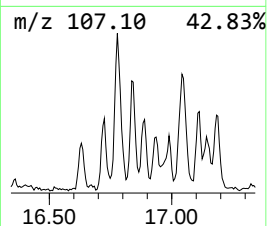
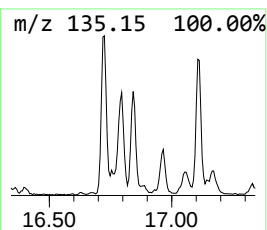
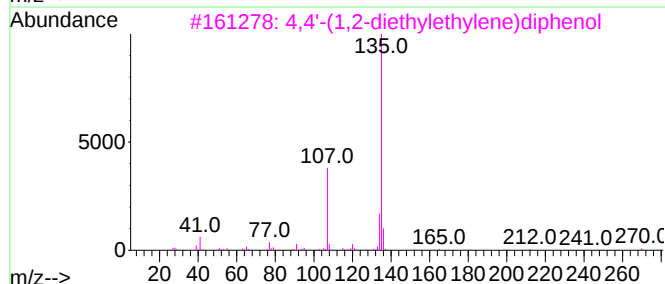
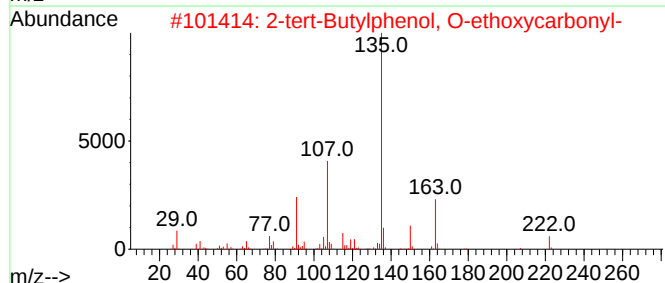
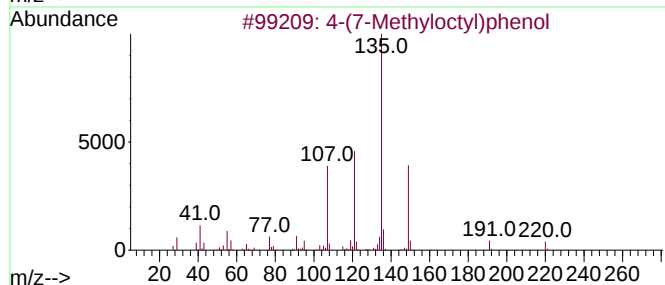
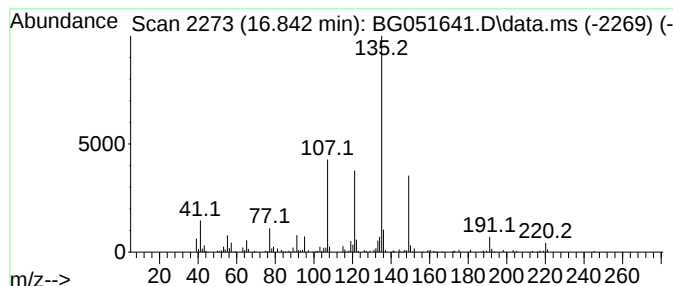
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Quant Title : SVOA CALIBRATION

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

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Peak Number 22 4-(7-Methyloctyl)phenol Concentration Rank 18

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 16.842 | 5.81 ng/ul | 168660 | Phenanthrene-d10 | 17.659 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|------------|--------------|------|
| 1    |    |   | 4-(7-Methyloctyl)phenol             | 220 | C15H24O    | 024518-48-7  | 95   |
| 2    |    |   | 2-tert-Butylphenol, O-ethoxycarb... | 222 | C13H18O3   | 1000487-37-8 | 59   |
| 3    |    |   | 4,4'-(1,2-diethylethylene)diphenol  | 270 | C18H22O2   | 005635-50-7  | 59   |
| 4    |    |   | Phenol, 4-(1,1-dimethylpropyl)-     | 164 | C11H16O    | 000080-46-6  | 59   |
| 5    |    |   | Hexestrol, O-heptafluorobutryl-     | 466 | C22H21F7O3 | 1000365-31-9 | 59   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG121721\  
Data File : BG051641.D  
Acq On : 18 Dec 2021 22:49  
Operator : CG/JU  
Sample : M5153-06  
Misc :  
ALS Vial : 57 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL71

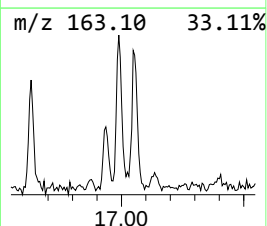
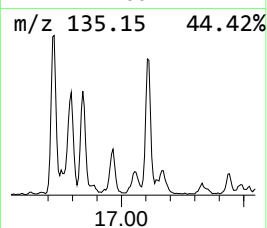
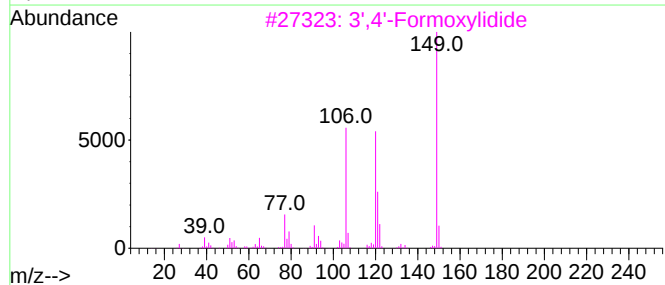
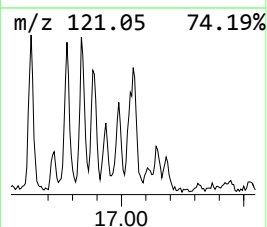
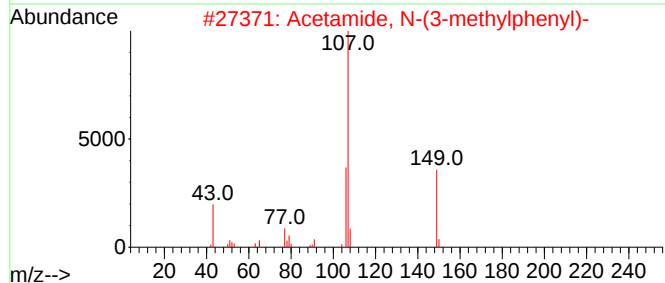
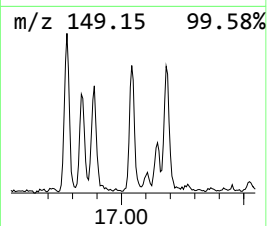
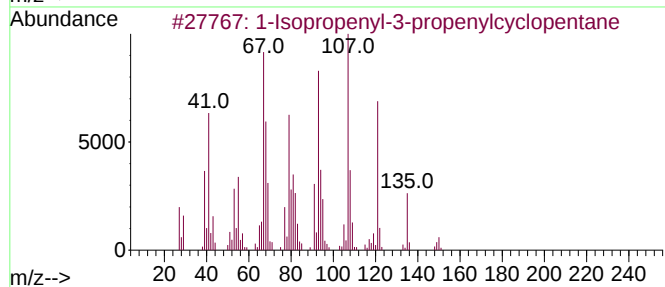
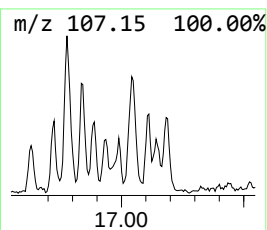
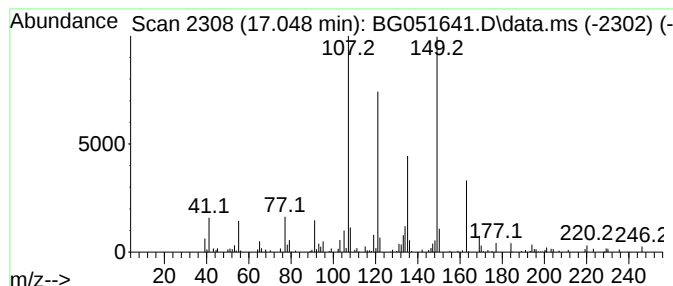
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Quant Title : SVOA CALIBRATION

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

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Peak Number 24 unknown-02 Concentration Rank 14

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 17.048 | 7.76 ng/ul | 225194 | Phenanthrene-d10 | 17.659 |

| Hit# | of                                   | 5   | Tentative ID | MW           | MolForm | CAS# | Qual |
|------|--------------------------------------|-----|--------------|--------------|---------|------|------|
| 1    | 1-Isopropenyl-3-propenylcyclopentane | 150 | C11H18       | 1000192-81-2 | 46      |      |      |
| 2    | Acetamide, N-(3-methylphenyl)-       | 149 | C9H11NO      | 000537-92-8  | 43      |      |      |
| 3    | 3',4'-Formoxylidide                  | 149 | C9H11NO      | 006639-60-7  | 18      |      |      |
| 4    | 4,5,7-Trimethyl-4,5,6,7-tetrahyd...  | 164 | C10H16N2     | 113849-86-8  | 18      |      |      |
| 5    | Acetamide, N-(2-methylphenyl)-       | 149 | C9H11NO      | 000120-66-1  | 18      |      |      |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG121721\  
Data File : BG051641.D  
Acq On : 18 Dec 2021 22:49  
Operator : CG/JU  
Sample : M5153-06  
Misc :  
ALS Vial : 57 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL71

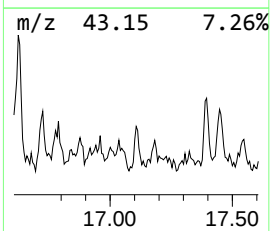
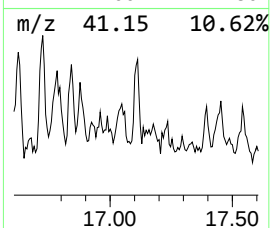
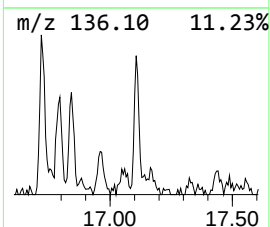
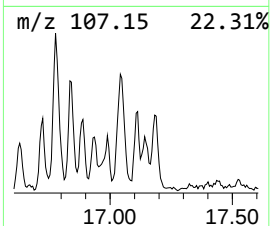
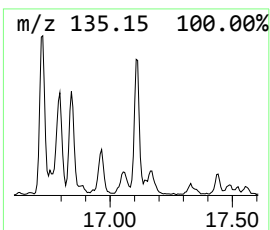
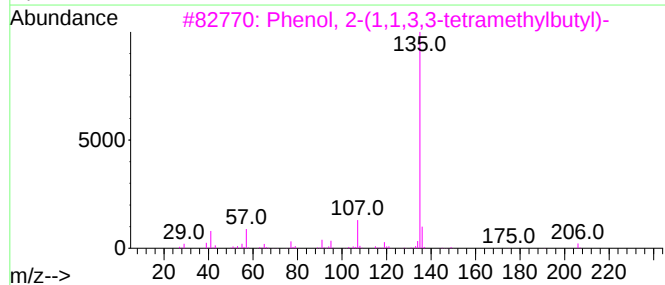
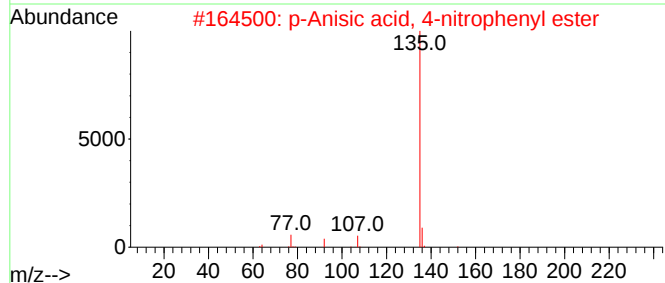
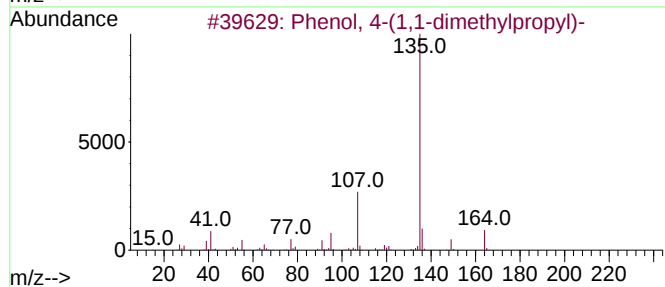
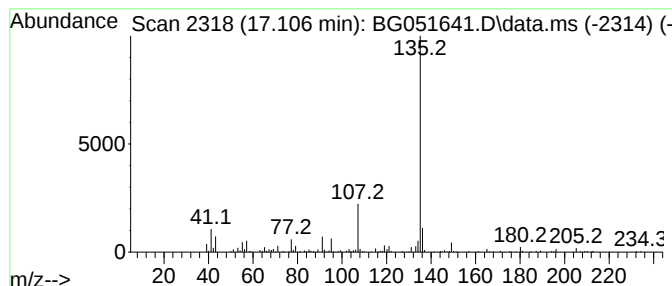
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Quant Title : SVOA CALIBRATION

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

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Peak Number 25 Phenol, 4-(1,1-dimethylprop... Concentration Rank 19

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 17.107 | 5.52 ng/ul | 160178 | Phenanthrene-d10 | 17.659 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Phenol, 4-(1,1-dimethylpropyl)-     | 164 | C11H16O   | 000080-46-6  | 78   |
| 2    |    |   | p-Anisic acid, 4-nitrophenyl ester  | 273 | C14H11NO5 | 1000307-63-9 | 64   |
| 3    |    |   | Phenol, 2-(1,1,3,3-tetramethylbu... | 206 | C14H22O   | 003884-95-5  | 64   |
| 4    |    |   | Phenol, 4-(1,1,3,3-tetramethylbu... | 206 | C14H22O   | 000140-66-9  | 64   |
| 5    |    |   | p-Anisic acid, 4-cyanophenyl ester  | 253 | C15H11NO3 | 1000307-63-7 | 64   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG121721\  
Data File : BG051641.D  
Acq On : 18 Dec 2021 22:49  
Operator : CG/JU  
Sample : M5153-06  
Misc :  
ALS Vial : 57 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL71

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

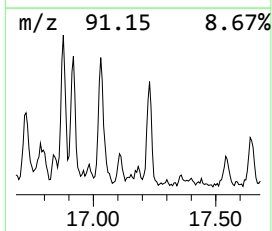
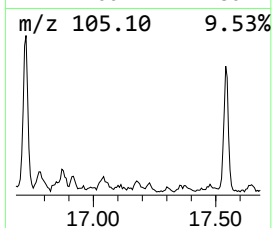
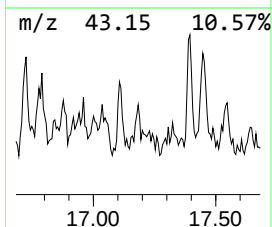
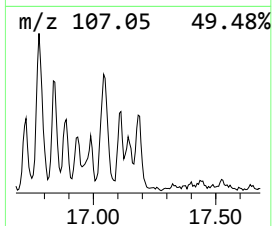
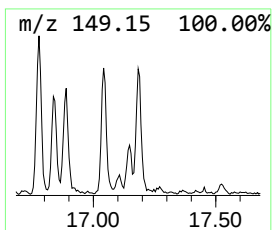
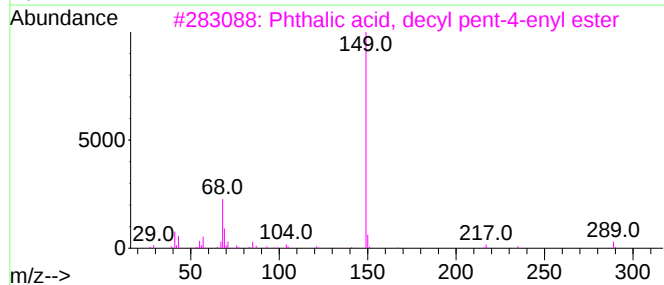
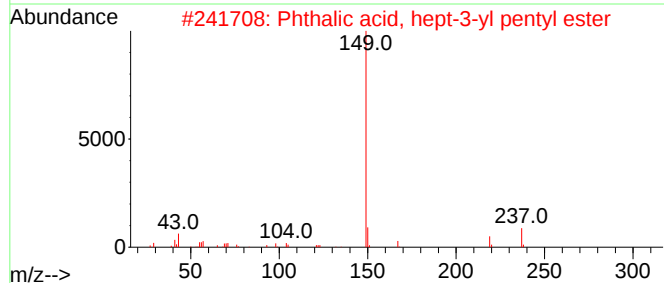
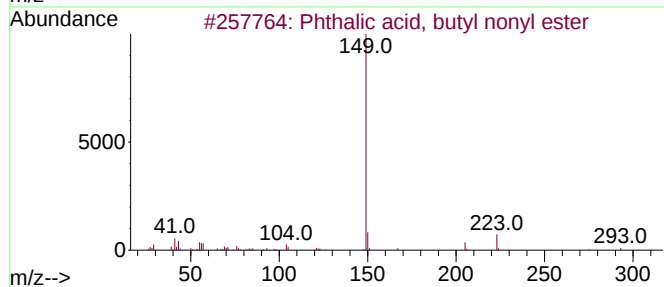
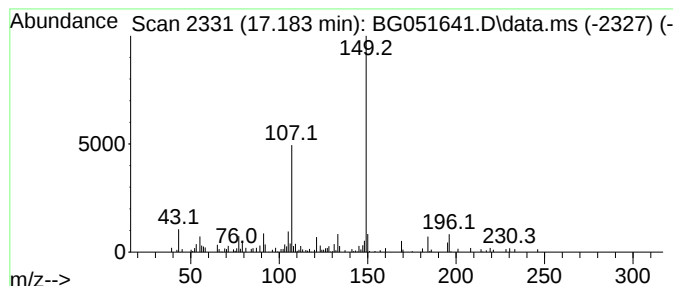
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TIC Integration Parameters: LSCINT.P

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Peak Number 26 unknown-03 Concentration Rank 25

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 17.183 | 4.44 ng/ul | 128876 | Phenanthrene-d10 | 17.659 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Phthalic acid, butyl nonyl ester    | 348 | C21H32O4 | 1000308-96-8 | 43   |
| 2    |    |   | Phthalic acid, hept-3-yl pentyl ... | 334 | C20H30O4 | 1010356-98-8 | 43   |
| 3    |    |   | Phthalic acid, decyl pent-4-enyl... | 374 | C23H34O4 | 1000360-46-9 | 43   |
| 4    |    |   | Phthalic acid, hexyl heptyl ester   | 348 | C21H32O4 | 1000308-93-8 | 43   |
| 5    |    |   | Phthalic acid, nonyl pentyl ester   | 362 | C22H34O4 | 1000308-93-1 | 43   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG121721\  
Data File : BG051641.D  
Acq On : 18 Dec 2021 22:49  
Operator : CG/JU  
Sample : M5153-06  
Misc :  
ALS Vial : 57 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL71

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

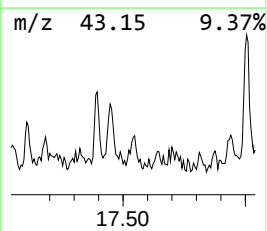
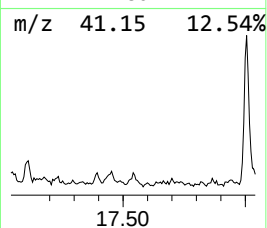
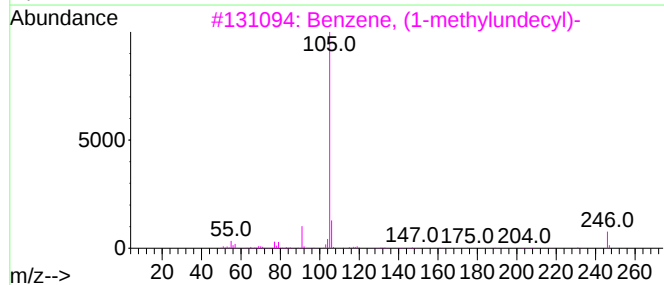
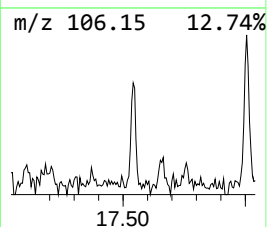
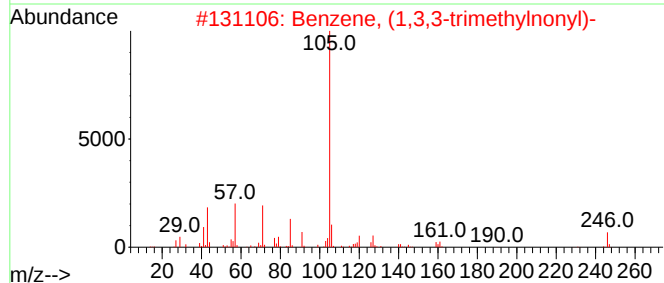
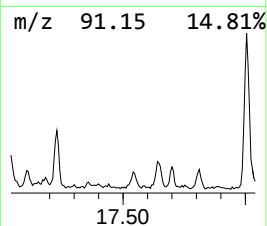
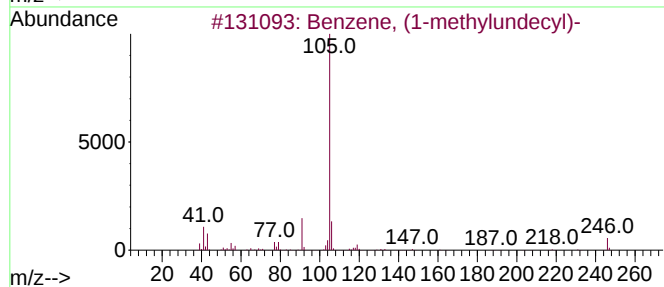
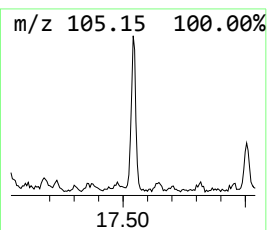
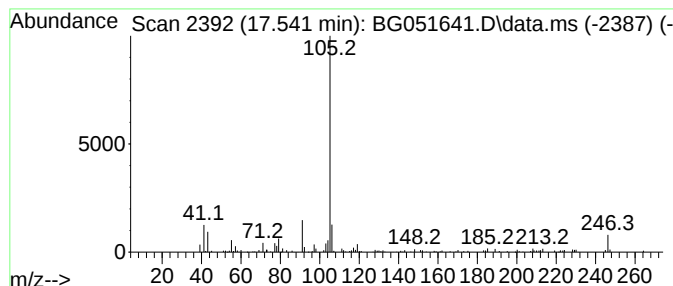
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 28 Benzene, (1-methylundecyl)- Concentration Rank 29

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 17.541 | 3.50 ng/ul | 101515 | Phenanthrene-d10 | 17.659 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|-----------|-------------|------|
| 1    |    |   | Benzene, (1-methylundecyl)-         | 246 | C18H30    | 002719-61-1 | 95   |
| 2    |    |   | Benzene, (1,3,3-trimethylnonyl)-    | 246 | C18H30    | 054986-44-6 | 64   |
| 3    |    |   | Benzene, (1-methylundecyl)-         | 246 | C18H30    | 002719-61-1 | 62   |
| 4    |    |   | Benzene, (1-methylundecyl)-         | 246 | C18H30    | 002719-61-1 | 62   |
| 5    |    |   | 2-Cyano-3-methyl-2-(O-methylbenz... | 231 | C14H17NO2 | 029277-53-0 | 47   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG121721\  
Data File : BG051641.D  
Acq On : 18 Dec 2021 22:49  
Operator : CG/JU  
Sample : M5153-06  
Misc :  
ALS Vial : 57 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL71

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

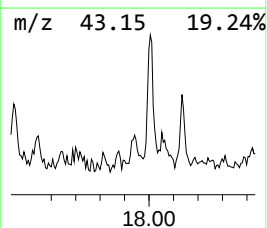
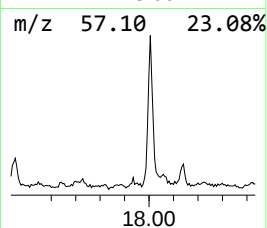
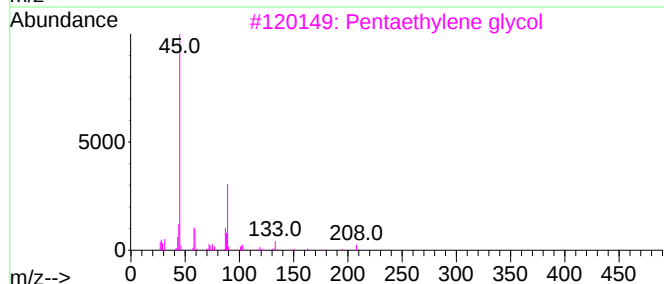
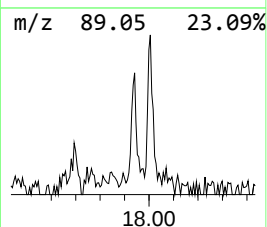
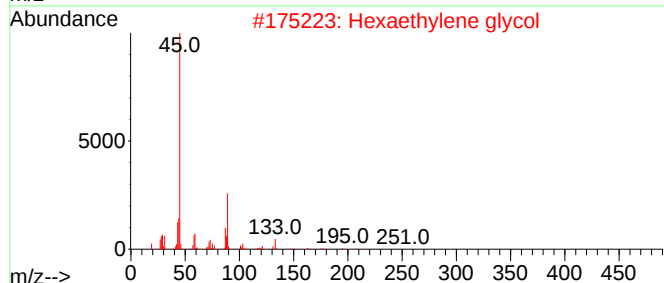
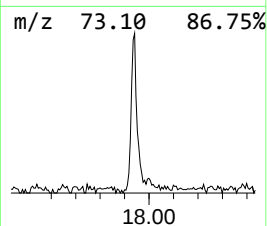
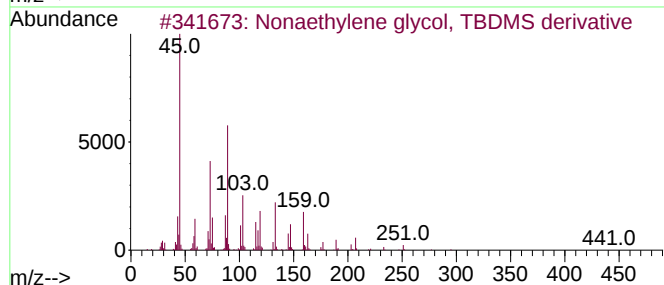
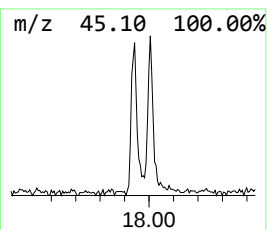
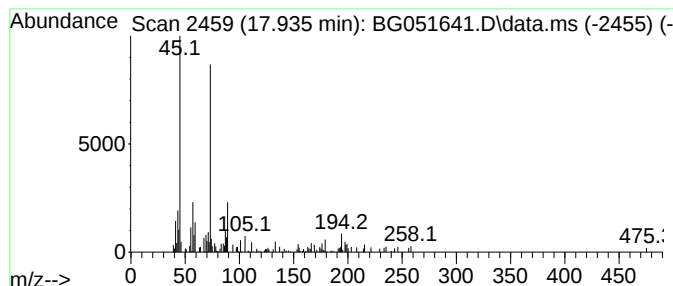
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 29 Nonaethylene glycol, TBDMS ... Concentration Rank 35

| R.T.   | EstConc    | Area  | Relative to ISTD | R.T.   |
|--------|------------|-------|------------------|--------|
| 17.935 | 2.91 ng/ul | 84598 | Phenanthrene-d10 | 17.659 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm     | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-------------|--------------|------|
| 1    |    |   | Nonaethylene glycol, TBDMS deriv... | 528 | C24H52O10Si | 1010352-11-3 | 78   |
| 2    |    |   | Hexaethylene glycol                 | 282 | C12H26O7    | 002615-15-8  | 64   |
| 3    |    |   | Pentaethylene glycol                | 238 | C10H22O6    | 004792-15-8  | 64   |
| 4    |    |   | Pentaethylene glycol                | 238 | C10H22O6    | 004792-15-8  | 59   |
| 5    |    |   | Pentaethylene glycol                | 238 | C10H22O6    | 004792-15-8  | 59   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG121721\  
Data File : BG051641.D  
Acq On : 18 Dec 2021 22:49  
Operator : CG/JU  
Sample : M5153-06  
Misc :  
ALS Vial : 57 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL71

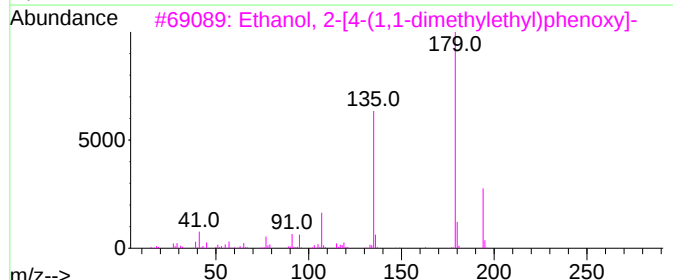
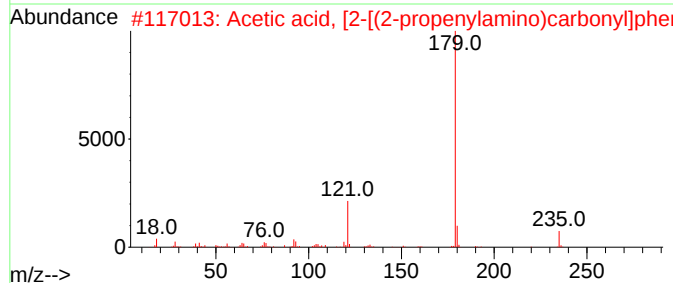
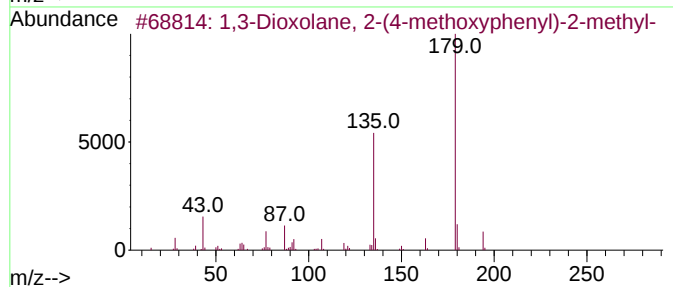
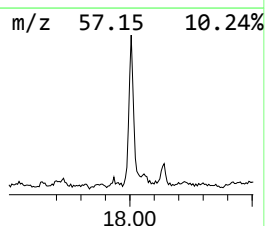
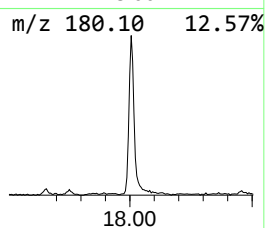
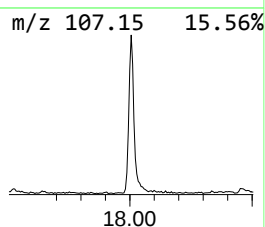
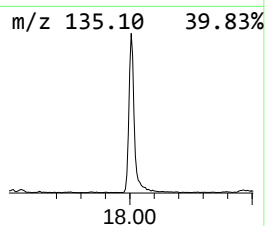
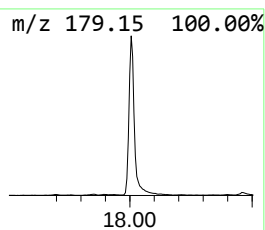
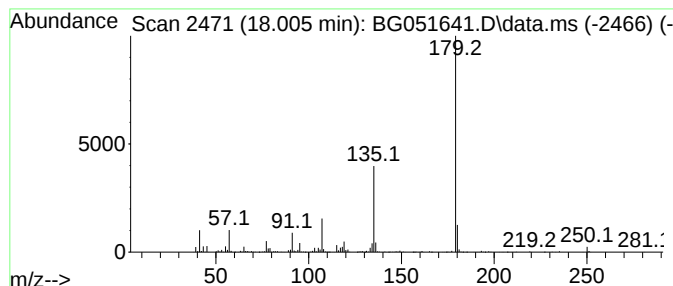
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Quant Title : SVOA CALIBRATION

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

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Peak Number 30 1,3-Dioxolane, 2-(4-methoxy... Concentration Rank 1

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 18.005 | 53.13 ng/ul | 1542650 | Phenanthrene-d10 | 17.659 |

| Hit# | of                                  | 5   | Tentative ID | MW          | MolForm | CAS# | Qual |
|------|-------------------------------------|-----|--------------|-------------|---------|------|------|
| 1    | 1,3-Dioxolane, 2-(4-methoxypheny... | 194 | C11H14O3     | 036881-00-2 | 50      |      |      |
| 2    | Acetic acid, [2-[(2-propenylamin... | 235 | C12H13NO4    | 000119-45-9 | 43      |      |      |
| 3    | Ethanol, 2-[4-(1,1-dimethylethyl... | 194 | C12H18O2     | 000713-46-2 | 40      |      |      |
| 4    | 2-(4-Formyl-phenoxy)-acetamide      | 179 | C9H9NO3      | 135857-20-4 | 38      |      |      |
| 5    | 2H-1-Benzopyran, 3,4,4a,5,6,8a-h... | 194 | C13H22O      | 041678-32-4 | 28      |      |      |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG121721\  
Data File : BG051641.D  
Acq On : 18 Dec 2021 22:49  
Operator : CG/JU  
Sample : M5153-06  
Misc :  
ALS Vial : 57 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL71

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

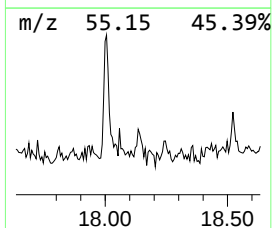
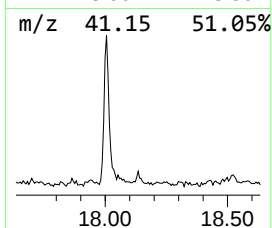
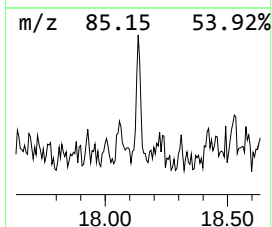
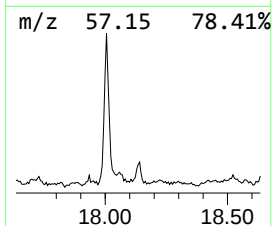
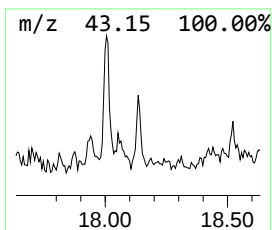
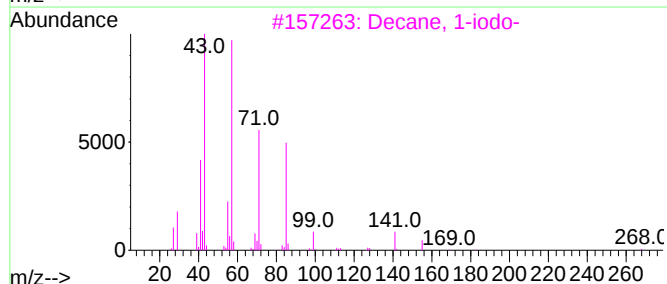
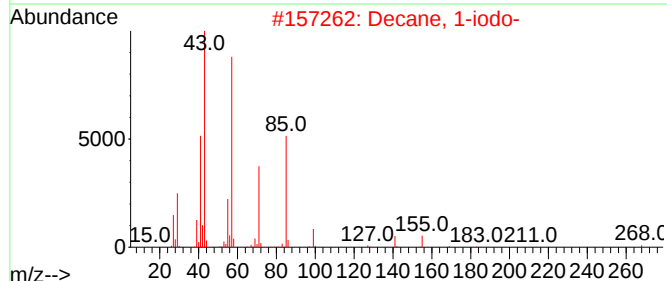
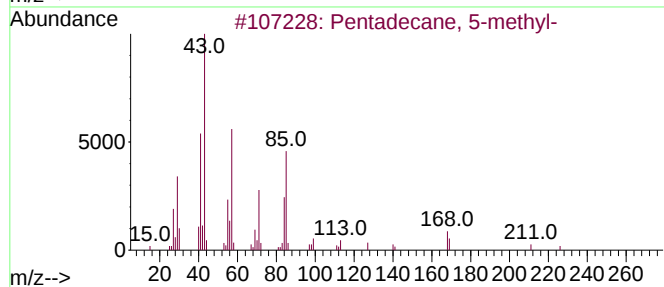
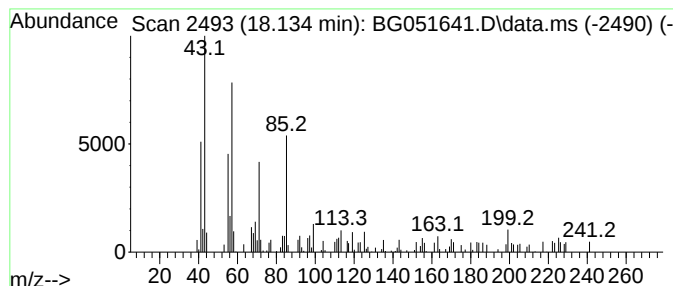
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 31 (DEL) Alkane: Straight-Chai... Concentration Rank 45

| R.T.   | EstConc    | Area  | Relative to ISTD | R.T.   |
|--------|------------|-------|------------------|--------|
| 18.134 | 2.15 ng/ul | 62403 | Phenanthrene-d10 | 17.659 |

| Hit# | of 5 | Tentative ID                      | MW  | MolForm   | CAS#         | Qual |
|------|------|-----------------------------------|-----|-----------|--------------|------|
| 1    |      | Pentadecane, 5-methyl-            | 226 | C16H34    | 025117-33-3  | 49   |
| 2    |      | Decane, 1-iodo-                   | 268 | C10H21I   | 002050-77-3  | 47   |
| 3    |      | Decane, 1-iodo-                   | 268 | C10H21I   | 002050-77-3  | 47   |
| 4    |      | Sulfurous acid, hexyl octyl ester | 278 | C14H30O3S | 1000309-13-0 | 47   |
| 5    |      | Tridecane                         | 184 | C13H28    | 000629-50-5  | 45   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG121721\  
Data File : BG051641.D  
Acq On : 18 Dec 2021 22:49  
Operator : CG/JU  
Sample : M5153-06  
Misc :  
ALS Vial : 57 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL71

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

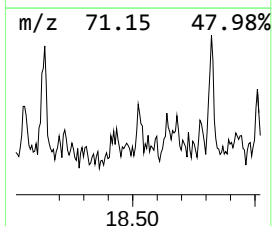
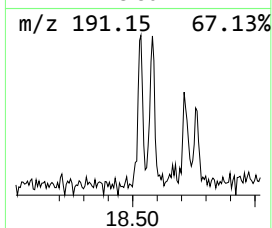
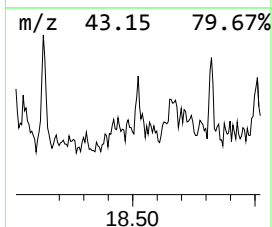
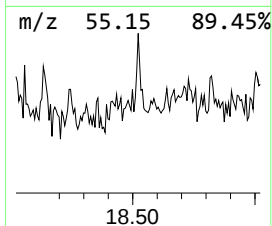
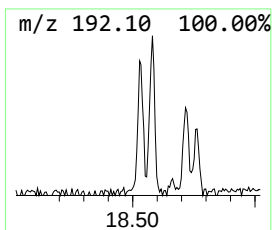
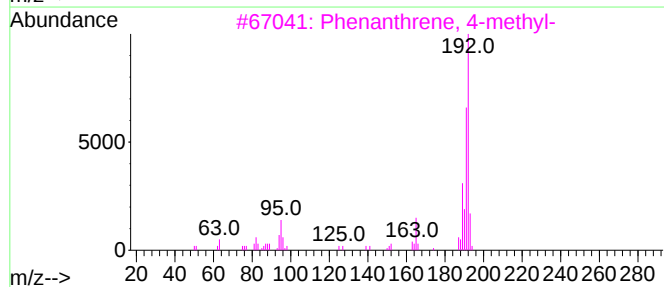
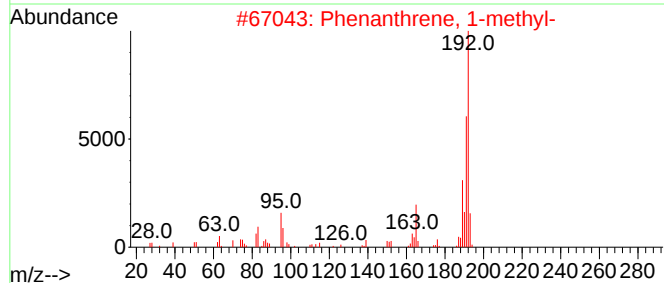
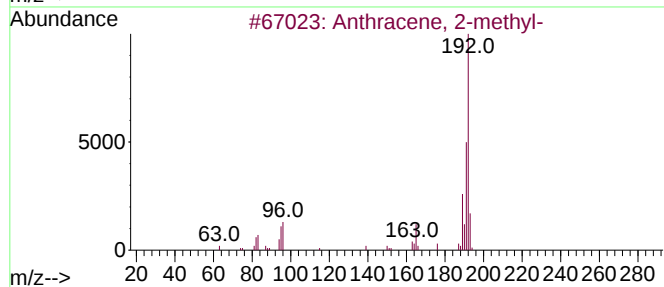
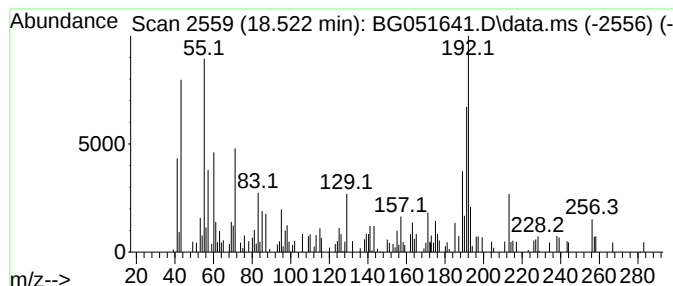
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 32 unknown-04 Concentration Rank 51

| R.T.   | EstConc    | Area  | Relative to ISTD | R.T.   |
|--------|------------|-------|------------------|--------|
| 18.522 | 2.03 ng/ul | 58856 | Phenanthrene-d10 | 17.659 |

| Hit# | of | 5 | Tentative ID                | MW  | MolForm | CAS#         | Qual |
|------|----|---|-----------------------------|-----|---------|--------------|------|
| 1    |    |   | Anthracene, 2-methyl-       | 192 | C15H12  | 000613-12-7  | 43   |
| 2    |    |   | Phenanthrene, 1-methyl-     | 192 | C15H12  | 000832-69-9  | 43   |
| 3    |    |   | Phenanthrene, 4-methyl-     | 192 | C15H12  | 000832-64-4  | 43   |
| 4    |    |   | Phenanthrene, 4-methyl-     | 192 | C15H12  | 000832-64-4  | 38   |
| 5    |    |   | Naphtho[1,2-b]norbornadiene | 192 | C15H12  | 1000210-14-8 | 38   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG121721\  
Data File : BG051641.D  
Acq On : 18 Dec 2021 22:49  
Operator : CG/JU  
Sample : M5153-06  
Misc :  
ALS Vial : 57 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL71

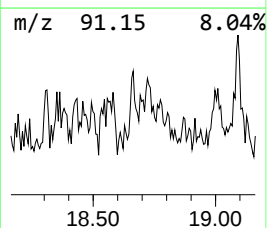
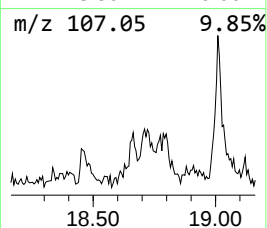
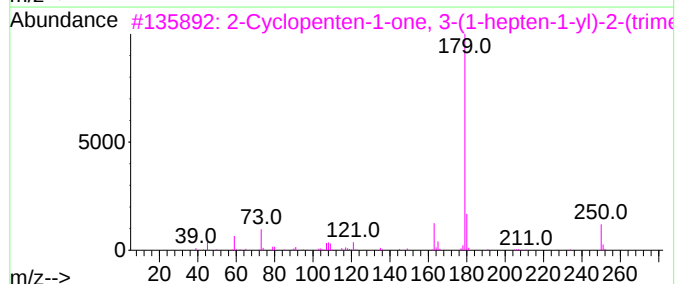
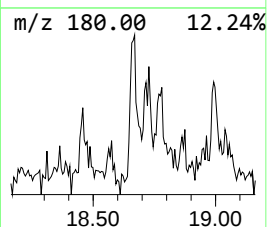
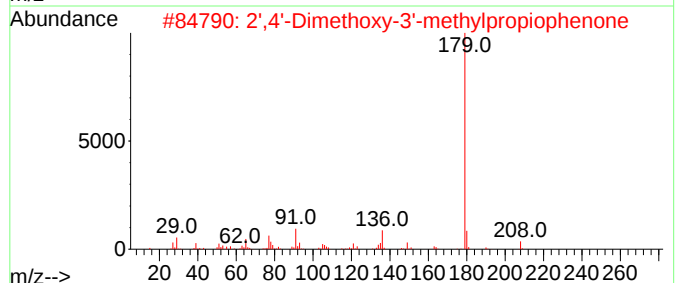
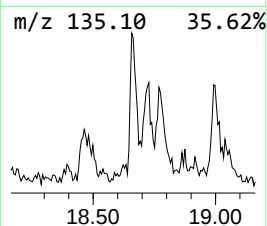
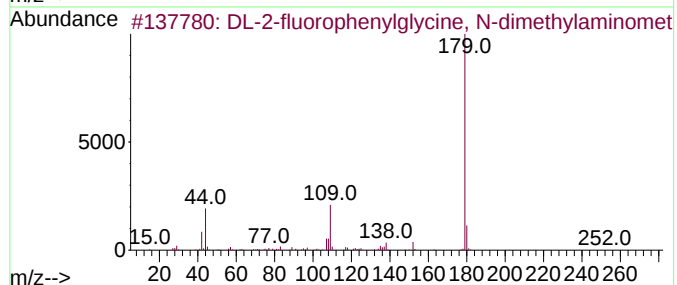
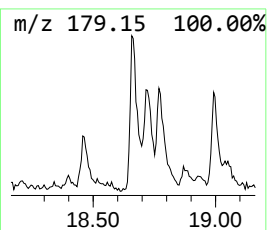
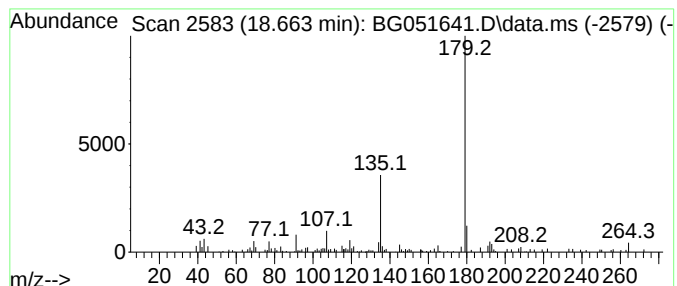
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Quant Title : SVOA CALIBRATION

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

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Peak Number 33 unknown-05 Concentration Rank 41

| R.T.   | EstConc    | Area  | Relative to ISTD | R.T.   |
|--------|------------|-------|------------------|--------|
| 18.663 | 2.55 ng/ul | 74053 | Phenanthrene-d10 | 17.659 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm     | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-------------|--------------|------|
| 1    |    |   | DL-2-fluorophenylglycine, N-dime... | 252 | C13H17FN2O2 | 1000375-81-0 | 42   |
| 2    |    |   | 2',4'-Dimethoxy-3'-methylpropiop... | 208 | C12H16O3    | 077942-13-3  | 40   |
| 3    |    |   | 2-Cyclopenten-1-one, 3-(1-hepten... | 250 | C15H26OSi   | 1000159-28-7 | 40   |
| 4    |    |   | Benzoic acid, 4-(dimethoxymethyl... | 210 | C11H14O4    | 042228-16-0  | 40   |
| 5    |    |   | 2-(4-Formyl-phenoxy)-acetamide      | 179 | C9H9NO3     | 135857-20-4  | 40   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG121721\  
Data File : BG051641.D  
Acq On : 18 Dec 2021 22:49  
Operator : CG/JU  
Sample : M5153-06  
Misc :  
ALS Vial : 57 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL71

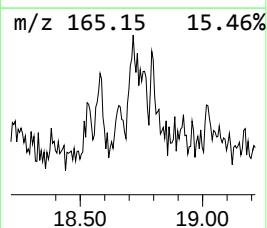
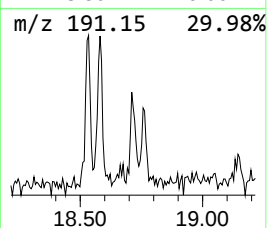
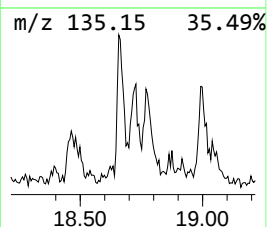
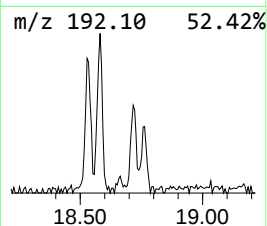
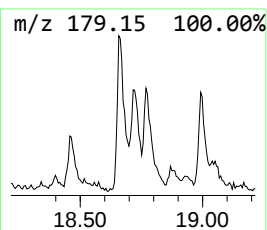
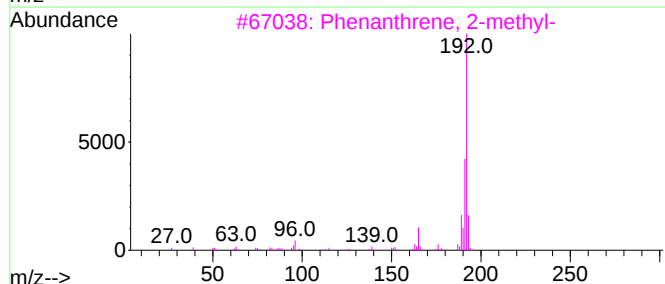
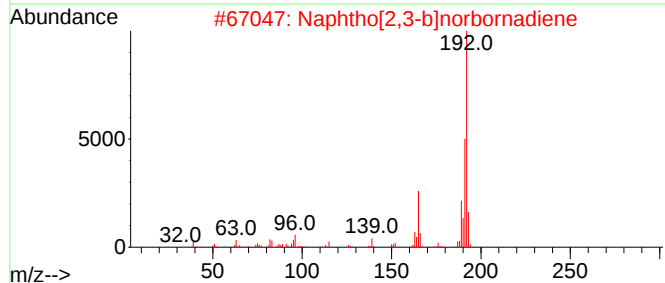
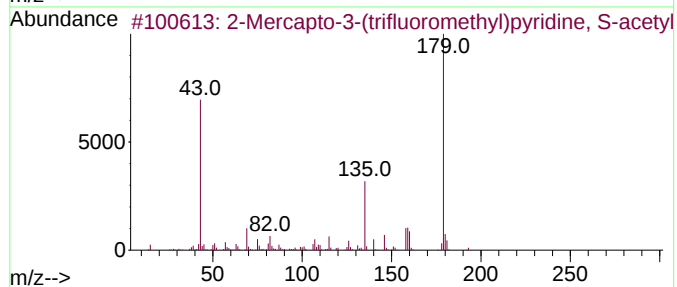
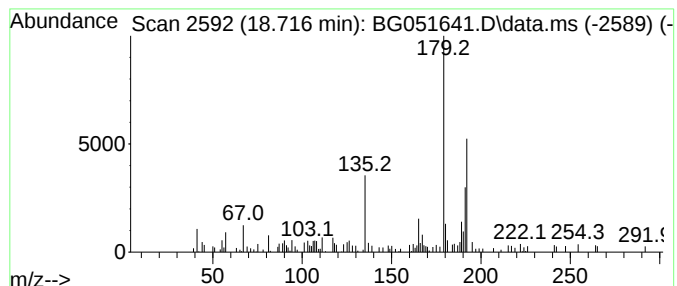
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Quant Title : SVOA CALIBRATION

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

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Peak Number 34 unknown-06 Concentration Rank 46

| R.T.   | EstConc    | Area  | Relative to ISTD | R.T.   |
|--------|------------|-------|------------------|--------|
| 18.716 | 2.15 ng/ul | 62396 | Phenanthrene-d10 | 17.659 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|------|-------------------------------------|-----|-----------|--------------|------|
| 1    |      | 2-Mercapto-3-(trifluoromethyl)py... | 221 | C8H6F3NOS | 1000463-18-0 | 38   |
| 2    |      | Naphtho[2,3-b]norbornadiene         | 192 | C15H12    | 107426-38-0  | 30   |
| 3    |      | Phenanthrene, 2-methyl-             | 192 | C15H12    | 002531-84-2  | 30   |
| 4    |      | Phenanthrene, 1-methyl-             | 192 | C15H12    | 000832-69-9  | 30   |
| 5    |      | Benzothiazole, 6-methoxy-2-methyl-  | 179 | C9H9NOS   | 002941-72-2  | 30   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG121721\  
Data File : BG051641.D  
Acq On : 18 Dec 2021 22:49  
Operator : CG/JU  
Sample : M5153-06  
Misc :  
ALS Vial : 57 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL71

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

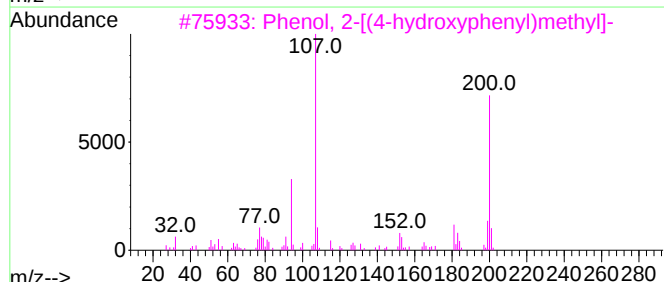
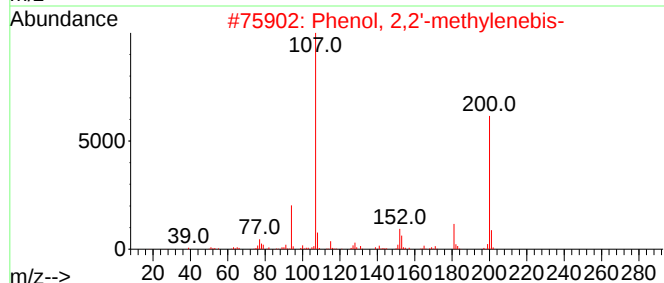
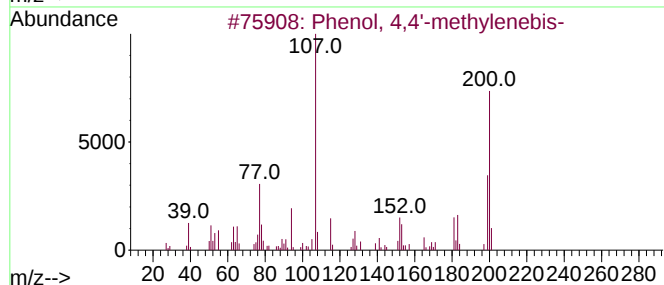
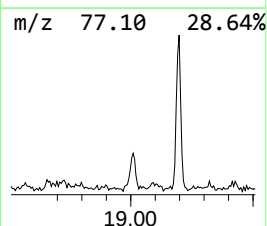
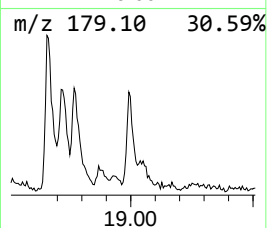
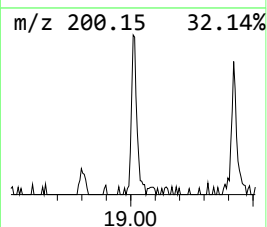
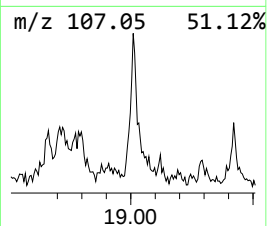
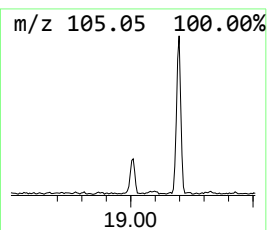
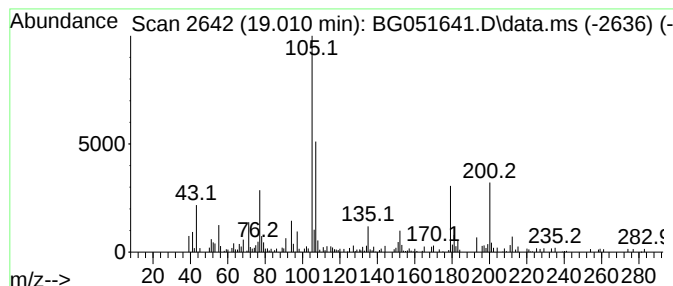
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TIC Integration Parameters: LSCINT.P

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Peak Number 35 unknown-07 Concentration Rank 22

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 19.010 | 4.58 ng/ul | 133041 | Phenanthrene-d10 | 17.659 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | Phenol, 4,4'-methylenebis-          | 200 | C13H12O2 | 000620-92-8 | 46   |
| 2    |    |   | Phenol, 2,2'-methylenebis-          | 200 | C13H12O2 | 002467-02-9 | 45   |
| 3    |    |   | Phenol, 2-[(4-hydroxyphenyl)meth... | 200 | C13H12O2 | 002467-03-0 | 41   |
| 4    |    |   | Phenol, 4,4'-methylenebis-          | 200 | C13H12O2 | 000620-92-8 | 38   |
| 5    |    |   | Phenol, 2-[(4-hydroxyphenyl)meth... | 200 | C13H12O2 | 002467-03-0 | 38   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG121721\  
Data File : BG051641.D  
Acq On : 18 Dec 2021 22:49  
Operator : CG/JU  
Sample : M5153-06  
Misc :  
ALS Vial : 57 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL71

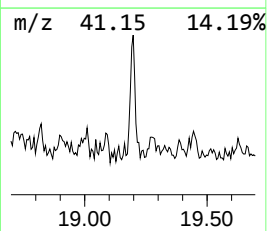
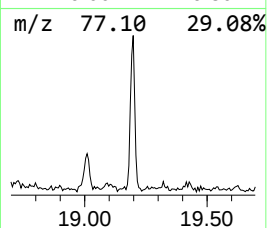
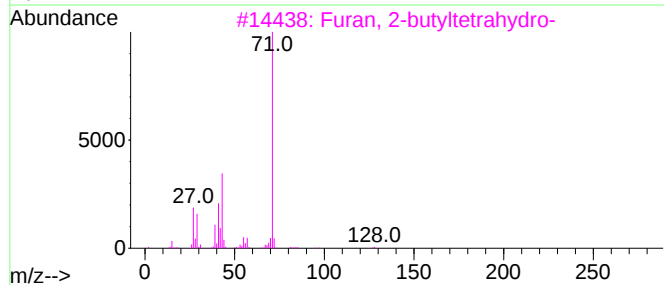
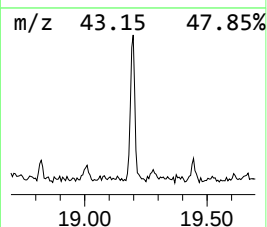
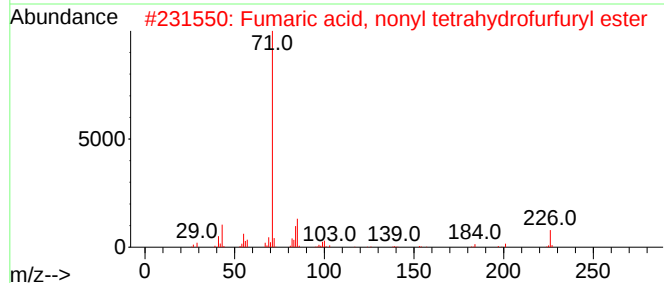
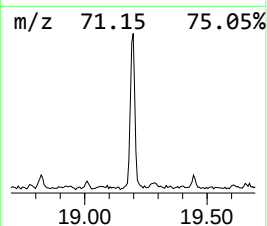
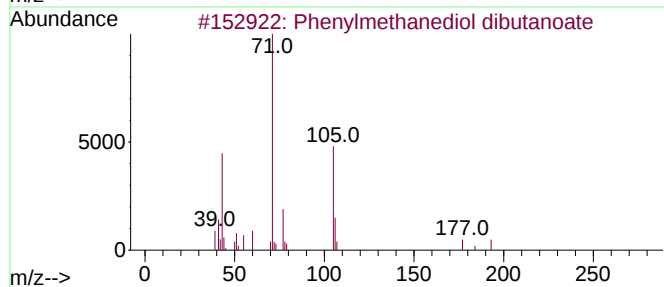
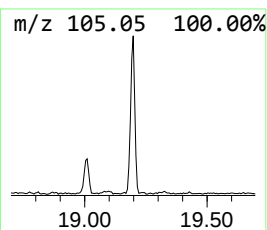
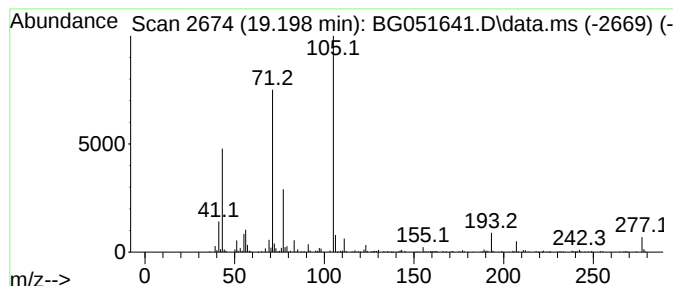
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Quant Title : SVOA CALIBRATION

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

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Peak Number 36 Phenylmethanediol dibutanoate Concentration Rank 8

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|--------|------------------|--------------|------|
| 19.198    | 10.76 ng/ul                         | 312471 | Phenanthrene-d10 | 17.659       |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#         | Qual |
| 1         | Phenylmethanediol dibutanoate       | 264    | C15H20O4         | 002929-77-3  | 64   |
| 2         | Fumaric acid, nonyl tetrahydrofu... | 326    | C18H30O5         | 1000330-58-3 | 43   |
| 3         | Furan, 2-butyltetrahydro-           | 128    | C8H16O           | 001004-29-1  | 43   |
| 4         | Propanoic acid, 2-methyl-, 1-met... | 158    | C9H18O2          | 054340-93-1  | 43   |
| 5         | N-(3-Methyl-1-phenyl-1H-pyrazol-... | 277    | C17H15N3O        | 064664-15-9  | 22   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG121721\  
Data File : BG051641.D  
Acq On : 18 Dec 2021 22:49  
Operator : CG/JU  
Sample : M5153-06  
Misc :  
ALS Vial : 57 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL71

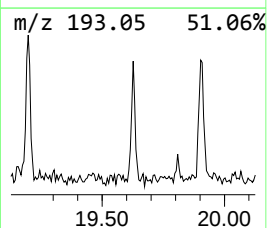
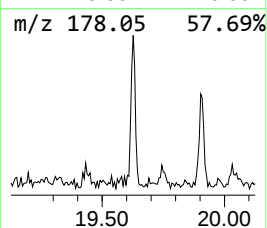
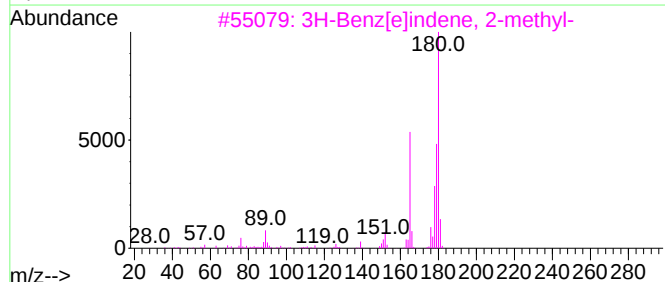
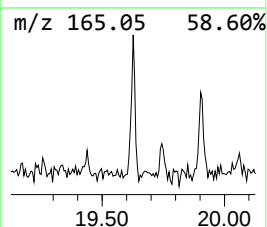
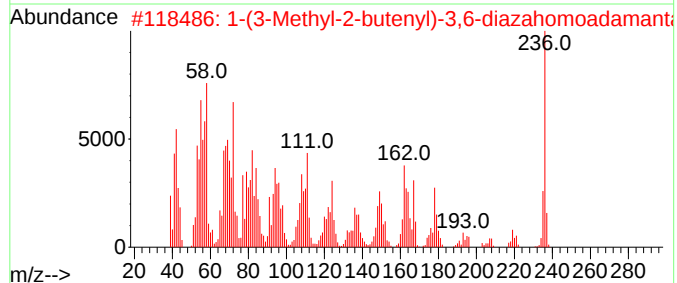
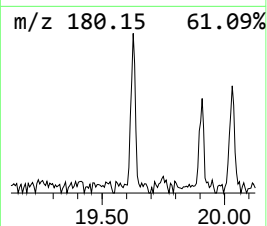
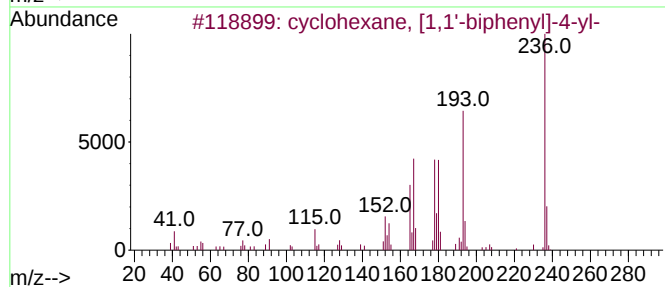
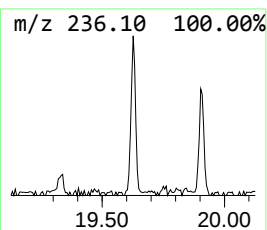
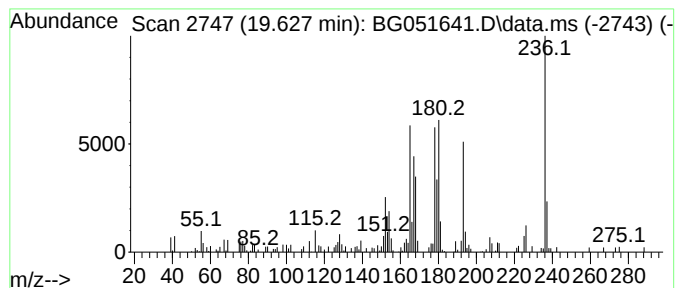
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Quant Title : SVOA CALIBRATION

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

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Peak Number 37 cyclohexane, [1,1'-biphenyl]... Concentration Rank 30

| R.T.   | EstConc    | Area  | Relative to ISTD | R.T.   |
|--------|------------|-------|------------------|--------|
| 19.627 | 3.22 ng/ul | 93429 | Phenanthrene-d10 | 17.659 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | cyclohexane, [1,1'-biphenyl]-4-yl-  | 236 | C18H20    | 1000401-12-4 | 76   |
| 2    |    |   | 1-(3-Methyl-2-butenyl)-3,6-diaza... | 236 | C14H24N2O | 1000216-26-4 | 46   |
| 3    |    |   | 3H-Benz[e]indene, 2-methyl-         | 180 | C14H12    | 150096-60-9  | 44   |
| 4    |    |   | 1-Diisopropylsilyloxy-3,5-dimeth... | 236 | C14H24OSi | 1000307-97-3 | 43   |
| 5    |    |   | 9,10-Anthracenedione, 2-ethyl-      | 236 | C16H12O2  | 000084-51-5  | 43   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG121721\  
Data File : BG051641.D  
Acq On : 18 Dec 2021 22:49  
Operator : CG/JU  
Sample : M5153-06  
Misc :  
ALS Vial : 57 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL71

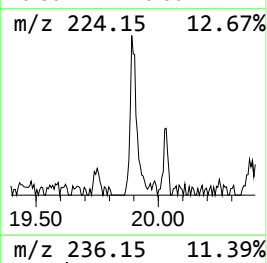
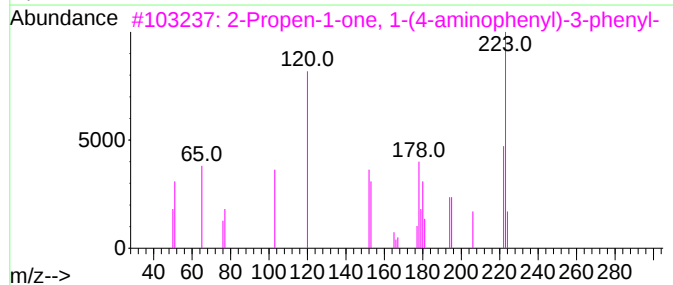
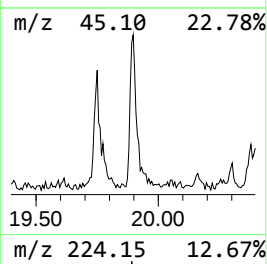
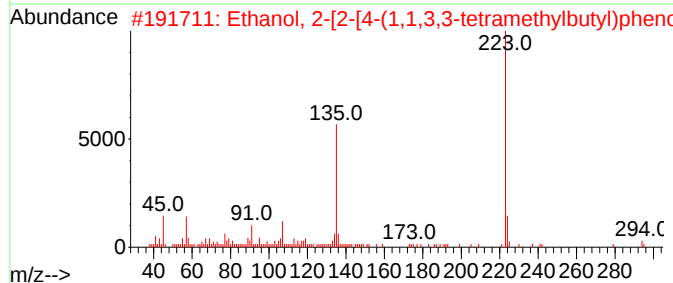
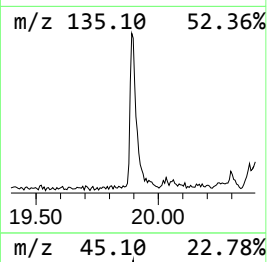
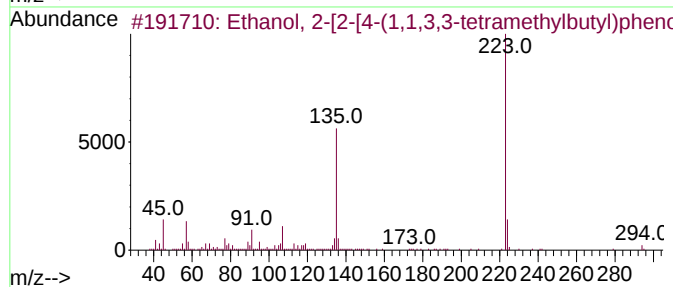
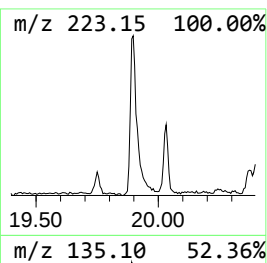
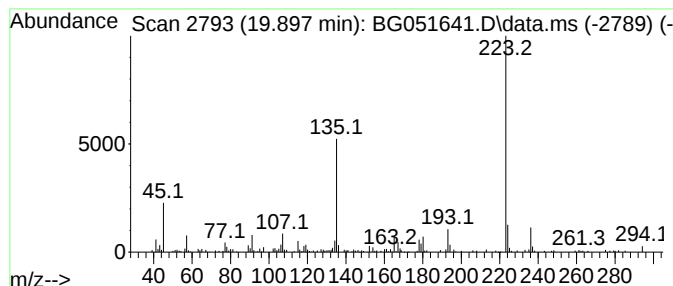
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Quant Title : SVOA CALIBRATION

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

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Peak Number 38 Ethanol, 2-[2-[4-(1,1,3,3-t... Concentration Rank 21

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------------|--------|------------------|-------------|------|
| 19.897    | 4.81 ng/ul                          | 207671 | Chrysene-d12     | 21.976      |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#        | Qual |
| 1         | Ethanol, 2-[2-[4-(1,1,3,3-tetram... | 294    | C18H30O3         | 002315-61-9 | 90   |
| 2         | Ethanol, 2-[2-[4-(1,1,3,3-tetram... | 294    | C18H30O3         | 002315-61-9 | 64   |
| 3         | 2-Propen-1-one, 1-(4-aminophenyl... | 223    | C15H13NO         | 002403-30-7 | 55   |
| 4         | 1,4,7-Trimethyl-2-azafluorenone     | 223    | C15H13NO         | 064322-98-1 | 47   |
| 5         | 1H-Isindole-1,3(2H)-dione, 2-ph...  | 223    | C14H9NO2         | 000520-03-6 | 43   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG121721\  
Data File : BG051641.D  
Acq On : 18 Dec 2021 22:49  
Operator : CG/JU  
Sample : M5153-06  
Misc :  
ALS Vial : 57 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL71

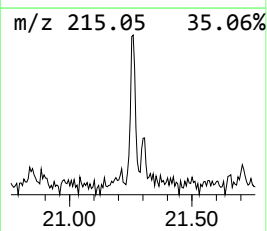
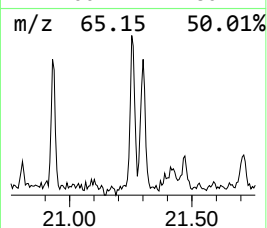
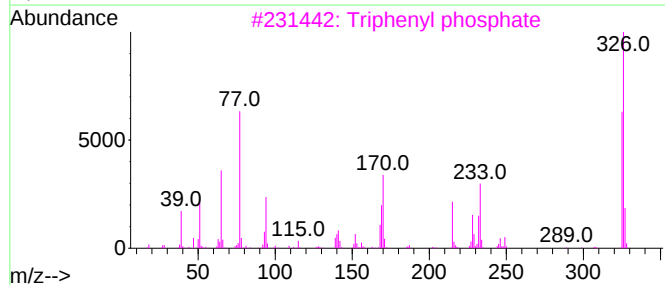
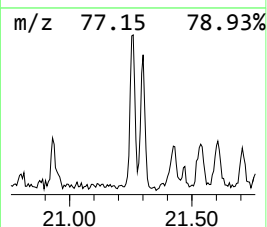
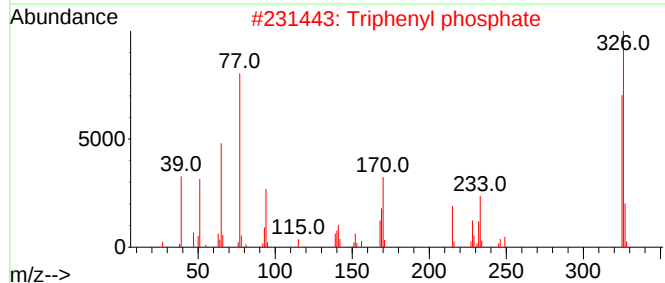
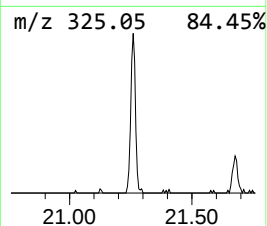
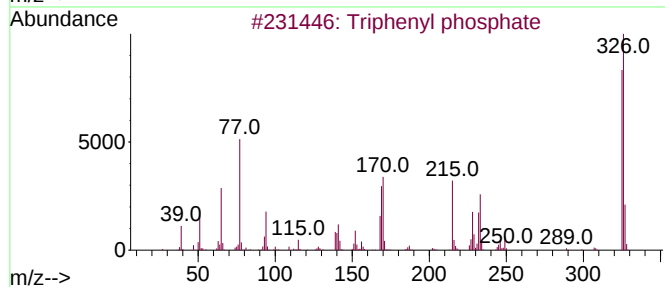
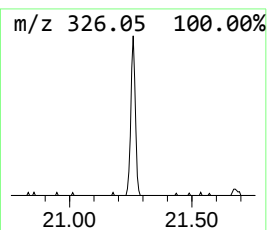
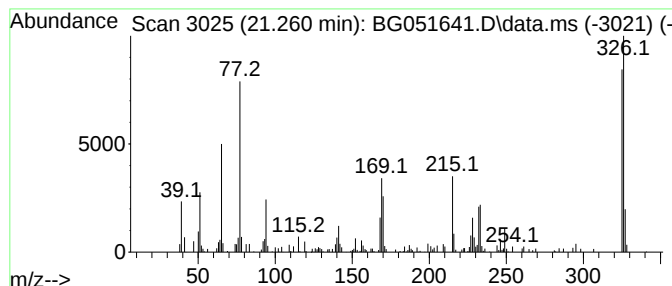
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Quant Title : SVOA CALIBRATION

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

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Peak Number 39 Triphenyl phosphate Concentration Rank 42

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 21.260 | 2.46 ng/ul | 106097 | Chrysene-d12     | 21.976 |

| Hit# | of | 5 | Tentative ID        | MW  | MolForm   | CAS#        | Qual |
|------|----|---|---------------------|-----|-----------|-------------|------|
| 1    |    |   | Triphenyl phosphate | 326 | C18H15O4P | 000115-86-6 | 96   |
| 2    |    |   | Triphenyl phosphate | 326 | C18H15O4P | 000115-86-6 | 93   |
| 3    |    |   | Triphenyl phosphate | 326 | C18H15O4P | 000115-86-6 | 86   |
| 4    |    |   | Triphenyl phosphate | 326 | C18H15O4P | 000115-86-6 | 70   |
| 5    |    |   | Triphenyl phosphate | 326 | C18H15O4P | 000115-86-6 | 70   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG121721\  
Data File : BG051641.D  
Acq On : 18 Dec 2021 22:49  
Operator : CG/JU  
Sample : M5153-06  
Misc :  
ALS Vial : 57 Sample Multiplier: 1

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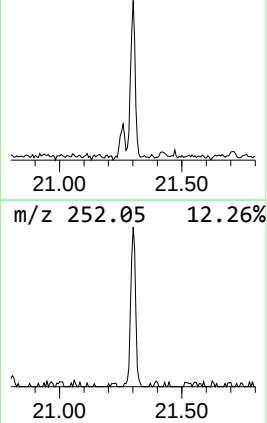
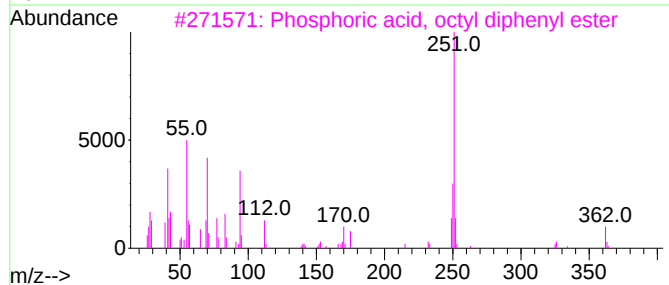
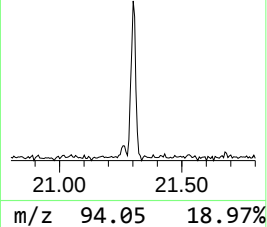
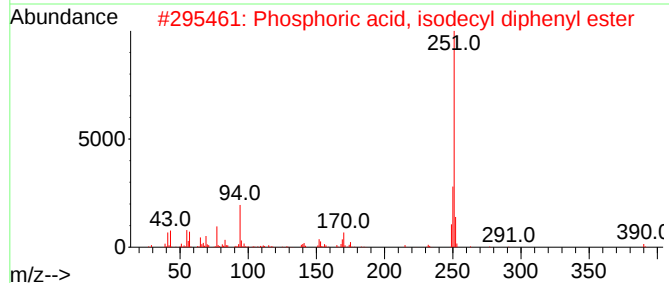
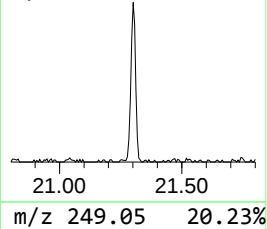
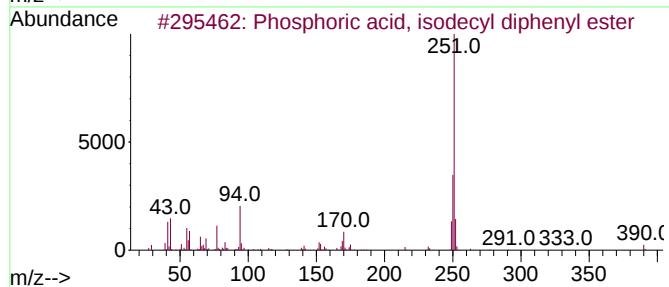
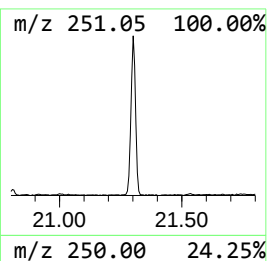
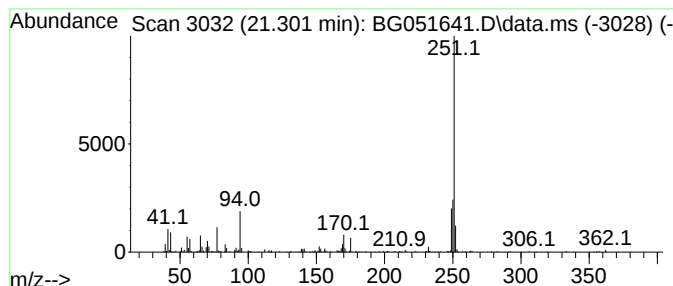
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Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 40 unknown-08 Concentration Rank 16

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 21.301 | 7.17 ng/ul | 309544 | Chrysene-d12     | 21.976 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|------------|--------------|------|
| 1    |    |   | Phosphoric acid, isodecyl diphen... | 390 | C22H31O4P  | 1346599-15-2 | 47   |
| 2    |    |   | Phosphoric acid, isodecyl diphen... | 390 | C22H31O4P  | 1346599-15-2 | 43   |
| 3    |    |   | Phosphoric acid, octyl diphenyl ... | 362 | C20H27O4P  | 000115-88-8  | 37   |
| 4    |    |   | 7-Amino-2,3-dihydro-5-phenyl-1H-... | 251 | C15H13N3O  | 004928-02-3  | 36   |
| 5    |    |   | Ethyl 2-(0-nitrophenylhydrazono)... | 251 | C11H13N3O4 | 006954-88-7  | 25   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG121721\  
Data File : BG051641.D  
Acq On : 18 Dec 2021 22:49  
Operator : CG/JU  
Sample : M5153-06  
Misc :  
ALS Vial : 57 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL71

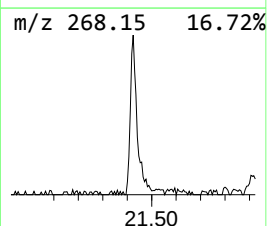
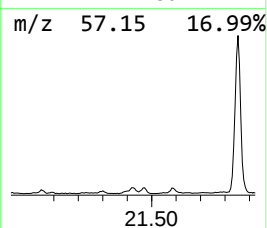
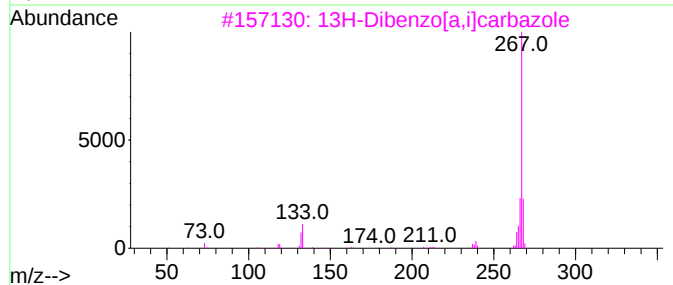
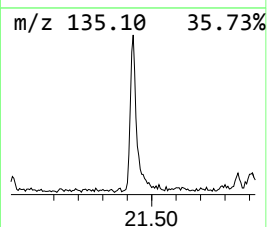
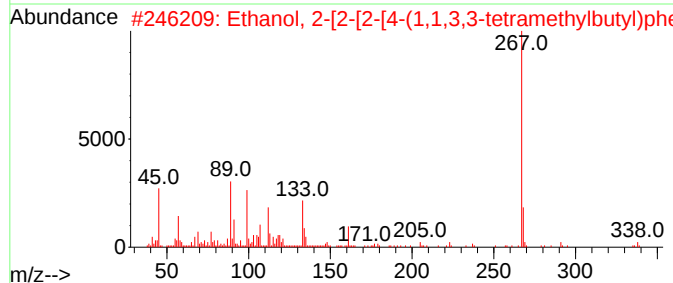
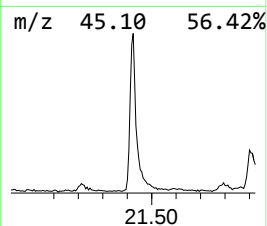
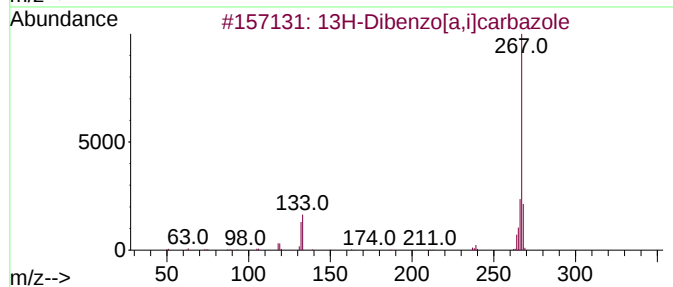
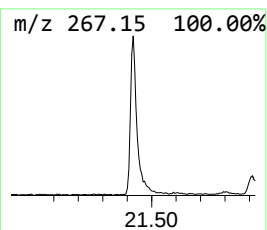
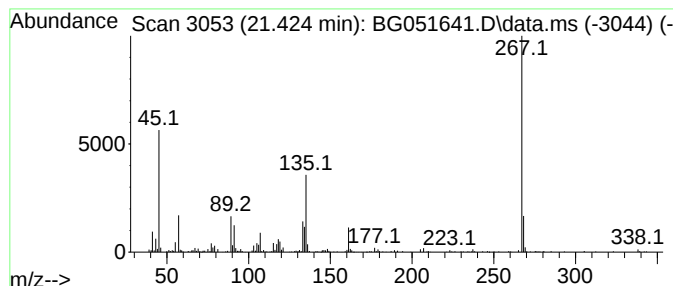
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Quant Title : SVOA CALIBRATION

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

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Peak Number 41 13H-Dibenzo[a,i]carbazole Concentration Rank 7

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 21.424 | 10.81 ng/ul | 466635 | Chrysene-d12     | 21.976 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | 13H-Dibenzo[a,i]carbazole           | 267 | C20H13N  | 000239-64-5 | 50   |
| 2    |    |   | Ethanol, 2-[2-[2-[4-(1,1,3,3-tet... | 338 | C20H34O4 | 002315-62-0 | 50   |
| 3    |    |   | 13H-Dibenzo[a,i]carbazole           | 267 | C20H13N  | 000239-64-5 | 50   |
| 4    |    |   | 13H-Dibenzo[a,i]carbazole           | 267 | C20H13N  | 000239-64-5 | 43   |
| 5    |    |   | 7H-Dibenzo(a,g)carbazole            | 267 | C20H13N  | 000207-84-1 | 38   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG121721\  
Data File : BG051641.D  
Acq On : 18 Dec 2021 22:49  
Operator : CG/JU  
Sample : M5153-06  
Misc :  
ALS Vial : 57 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL71

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

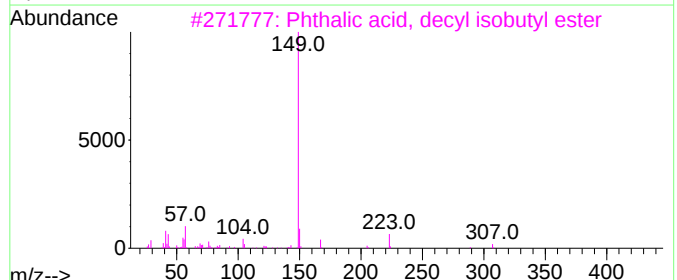
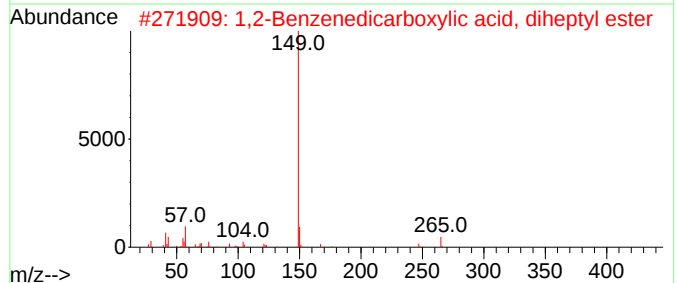
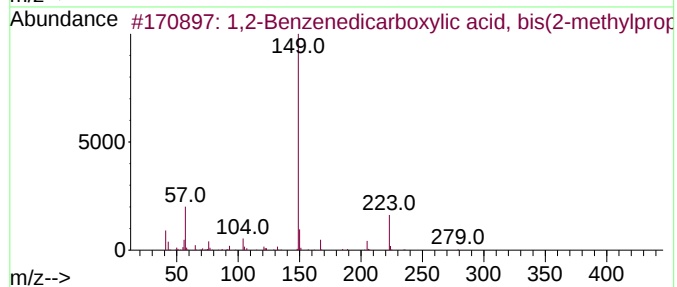
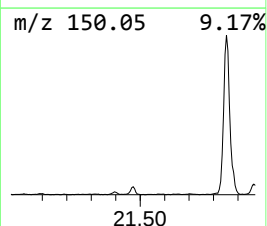
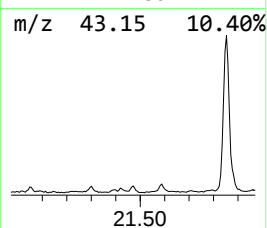
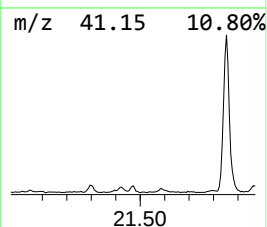
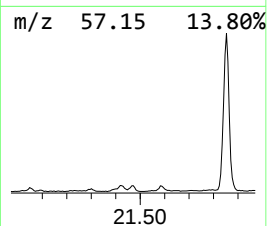
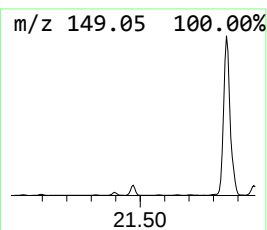
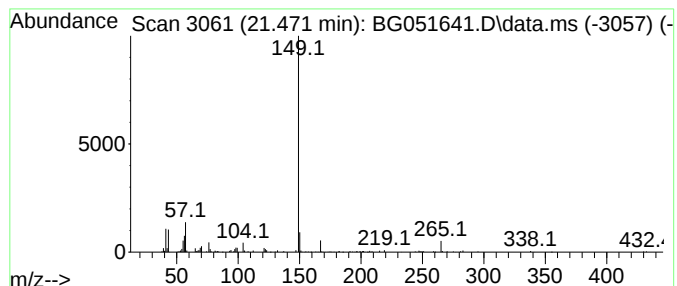
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 42 1,2-Benzenedicarboxylic aci... Concentration Rank 20

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 21.471 | 5.05 ng/ul | 218064 | Chrysene-d12     | 21.976 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm     | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-------------|--------------|------|
| 1    |    |   | 1,2-Benzenedicarboxylic acid, bi... | 278 | C16H22O4    | 000084-69-5  | 78   |
| 2    |    |   | 1,2-Benzenedicarboxylic acid, di... | 362 | C22H34O4    | 003648-21-3  | 78   |
| 3    |    |   | Phthalic acid, decyl isobutyl ester | 362 | C22H34O4    | 1000308-94-2 | 78   |
| 4    |    |   | Phthalic acid, 4-chlorophenyl he... | 374 | C21H23ClO4  | 1000315-25-1 | 72   |
| 5    |    |   | Phthalic acid, 3,4-dichloropheny... | 464 | C25H30Cl2O4 | 1000315-28-0 | 72   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG121721\  
Data File : BG051641.D  
Acq On : 18 Dec 2021 22:49  
Operator : CG/JU  
Sample : M5153-06  
Misc :  
ALS Vial : 57 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL71

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

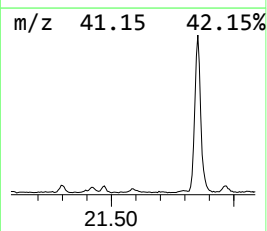
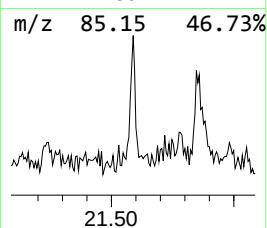
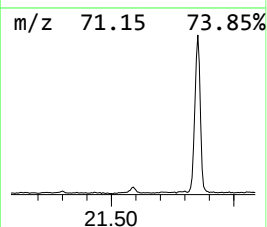
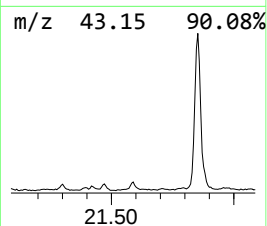
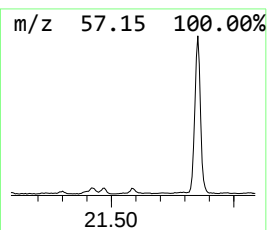
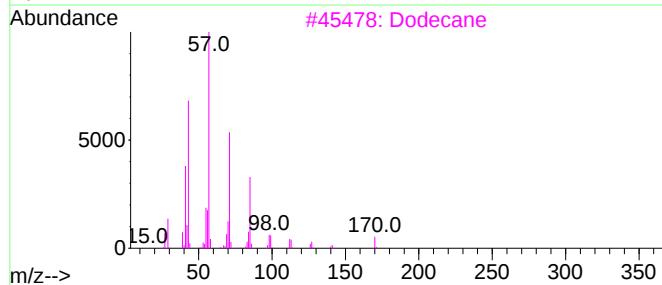
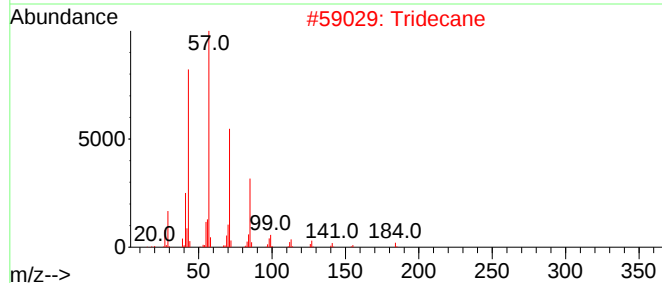
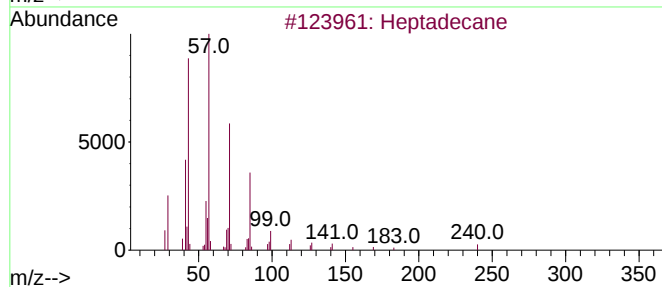
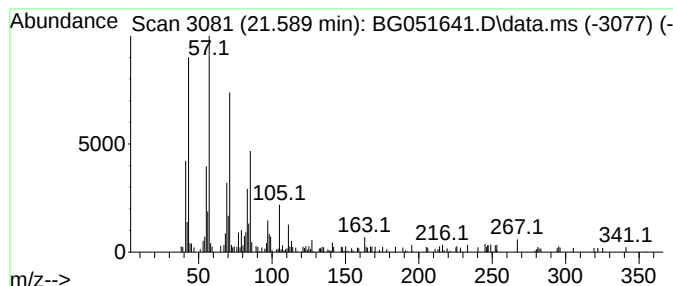
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 43 (DEL) Alkane: Straight-Chai... Concentration Rank 32

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 21.589 | 3.02 ng/ul | 130261 | Chrysene-d12     | 21.976 |

| Hit# | of 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|------|--------------|-----|---------|-------------|------|
| 1    |      | Heptadecane  | 240 | C17H36  | 000629-78-7 | 64   |
| 2    |      | Tridecane    | 184 | C13H28  | 000629-50-5 | 64   |
| 3    |      | Dodecane     | 170 | C12H26  | 000112-40-3 | 64   |
| 4    |      | Hexadecane   | 226 | C16H34  | 000544-76-3 | 64   |
| 5    |      | Hexadecane   | 226 | C16H34  | 000544-76-3 | 64   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG121721\  
Data File : BG051641.D  
Acq On : 18 Dec 2021 22:49  
Operator : CG/JU  
Sample : M5153-06  
Misc :  
ALS Vial : 57 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL71

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

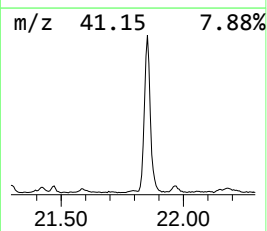
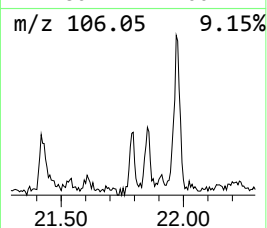
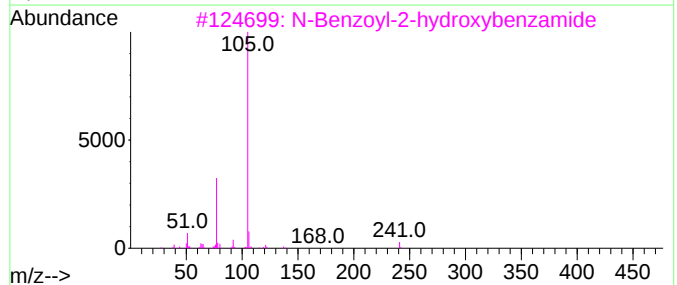
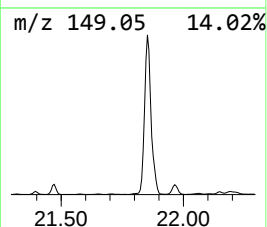
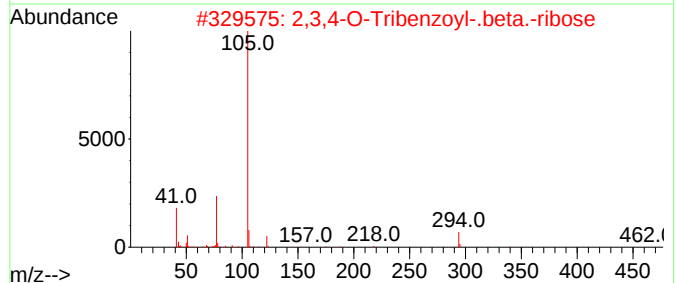
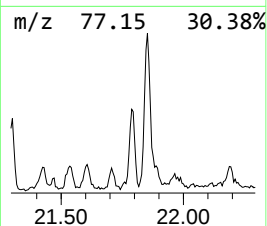
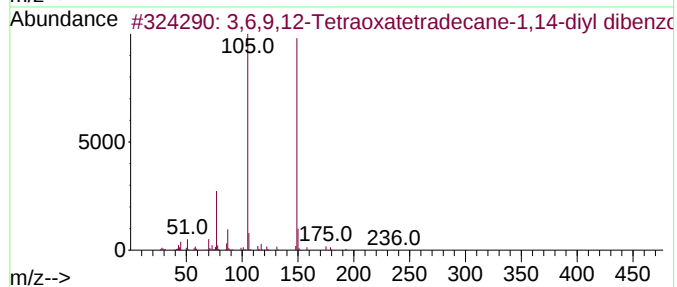
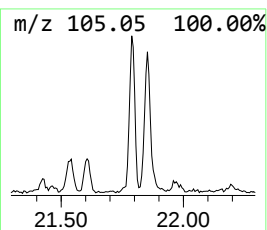
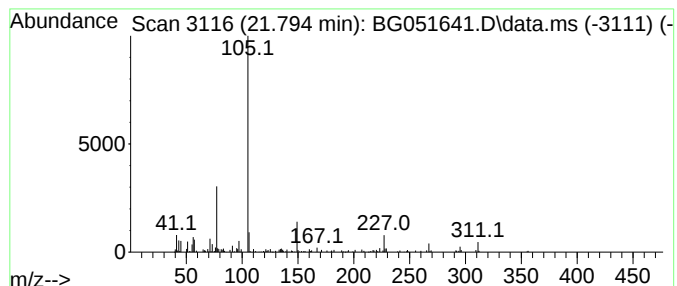
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 44 3,6,9,12-Tetraoxatetradecan... Concentration Rank 34

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 21.794 | 3.00 ng/ul | 129322 | Chrysene-d12     | 21.976 |

| Hit# | of                                  | 5   | Tentative ID | MW           | MolForm | CAS# | Qual |
|------|-------------------------------------|-----|--------------|--------------|---------|------|------|
| 1    | 3,6,9,12-Tetraoxatetradecane-1,1... | 446 | C24H30O8     | 1000366-97-0 | 53      |      |      |
| 2    | 2,3,4-O-Tribenzoyl-.beta.-ribose    | 462 | C26H22O8     | 1000129-80-8 | 47      |      |      |
| 3    | N-Benzoyl-2-hydroxybenzamide        | 241 | C14H11NO3    | 005663-74-1  | 46      |      |      |
| 4    | N-(2,4-Dihydroxy-6-methyl-5-pyri... | 245 | C12H11N3O3   | 1000472-87-2 | 43      |      |      |
| 5    | 3-Cyclopenten-1-ol, benzoate        | 188 | C12H12O2     | 043019-84-7  | 43      |      |      |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG121721\  
Data File : BG051641.D  
Acq On : 18 Dec 2021 22:49  
Operator : CG/JU  
Sample : M5153-06  
Misc :  
ALS Vial : 57 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL71

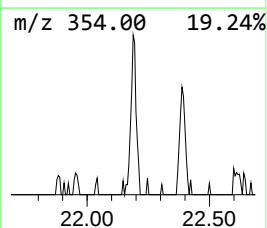
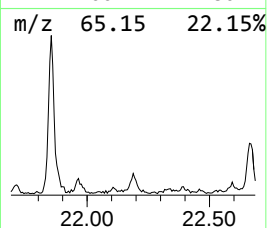
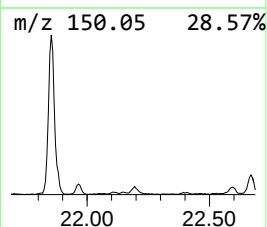
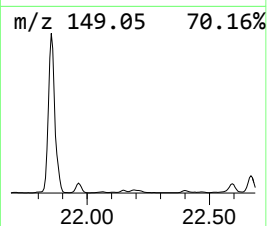
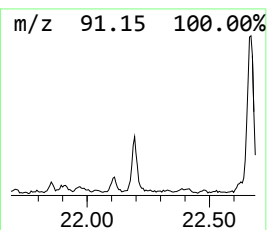
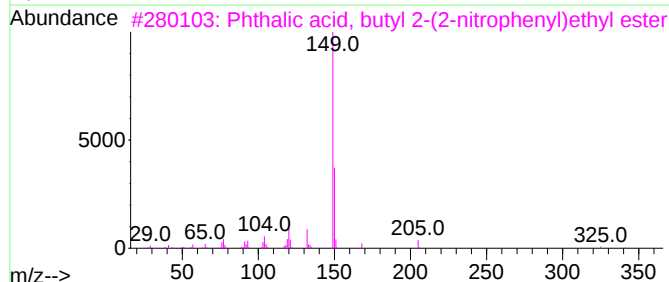
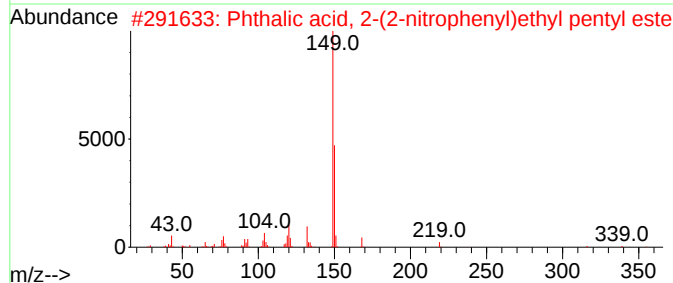
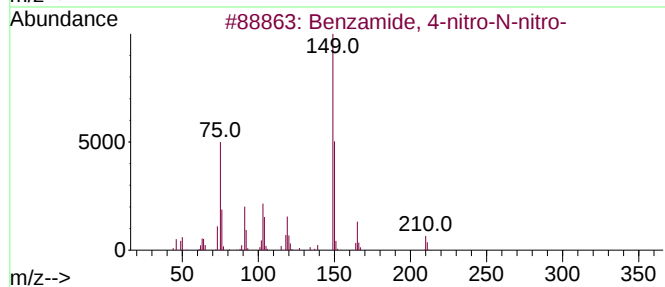
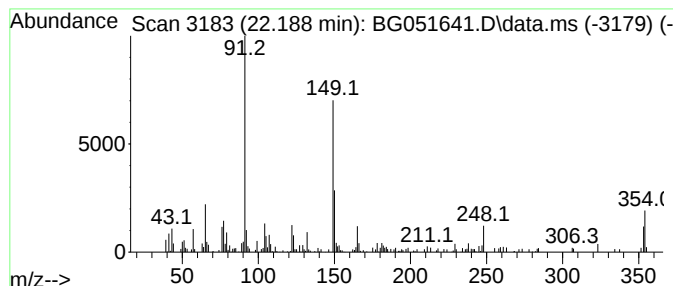
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Quant Title : SVOA CALIBRATION

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

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Peak Number 45 Benzamide, 4-nitro-N-nitro- Concentration Rank 17

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 22.188 | 7.12 ng/ul | 307558 | Chrysene-d12     | 21.976 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Benzamide, 4-nitro-N-nitro-         | 211 | C7H5N3O5  | 069719-32-0  | 50   |
| 2    |    |   | Phthalic acid, 2-(2-nitrophenyl)... | 385 | C21H23N06 | 1000377-99-5 | 38   |
| 3    |    |   | Phthalic acid, butyl 2-(2-nitrop... | 371 | C20H21N06 | 1000377-99-4 | 38   |
| 4    |    |   | (E)-2-((But-2-enyloxy)carbonyl)b... | 220 | C12H12O4  | 855233-54-4  | 38   |
| 5    |    |   | Phthalic acid, 3-methylbenzyl te... | 466 | C30H42O4  | 1000371-09-7 | 35   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG121721\  
Data File : BG051641.D  
Acq On : 18 Dec 2021 22:49  
Operator : CG/JU  
Sample : M5153-06  
Misc :  
ALS Vial : 57 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL71

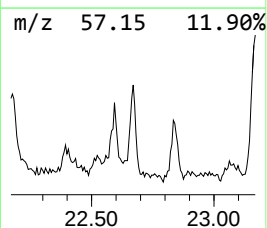
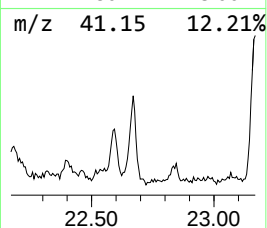
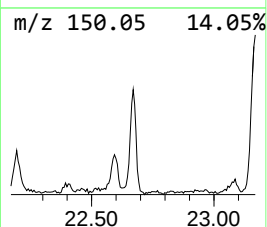
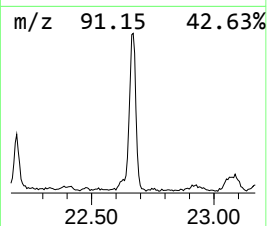
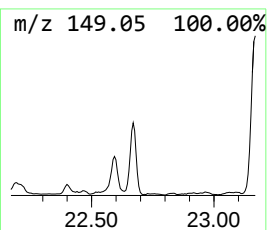
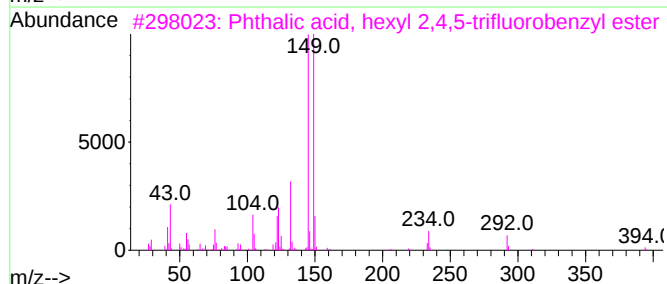
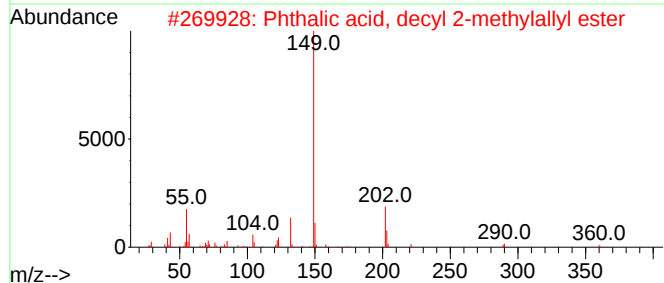
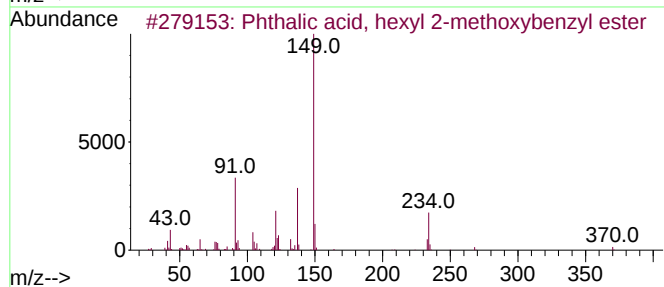
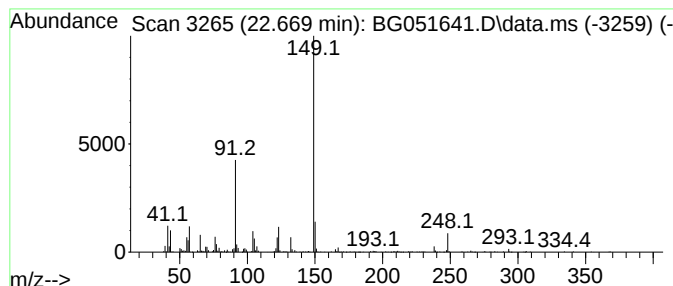
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Quant Title : SVOA CALIBRATION

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

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Peak Number 47 Phthalic acid, hexyl 2-meth... Concentration Rank 5

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 22.669 | 14.45 ng/ul | 623871 | Chrysene-d12     | 21.976 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|------------|--------------|------|
| 1    |    |   | Phthalic acid, hexyl 2-methoxybe... | 370 | C22H26O5   | 1000382-49-5 | 78   |
| 2    |    |   | Phthalic acid, decyl 2-methylall... | 360 | C22H32O4   | 1000315-83-0 | 72   |
| 3    |    |   | Phthalic acid, hexyl 2,4,5-trifl... | 394 | C21H21F3O4 | 1000415-49-6 | 64   |
| 4    |    |   | 1,2-Benzenedicarboxylic acid, di... | 250 | C14H18O4   | 000131-16-8  | 59   |
| 5    |    |   | Phthalic acid, 2-isopropoxypheny... | 370 | C22H26O5   | 1000309-83-1 | 59   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG121721\  
Data File : BG051641.D  
Acq On : 18 Dec 2021 22:49  
Operator : CG/JU  
Sample : M5153-06  
Misc :  
ALS Vial : 57 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL71

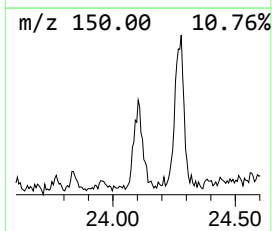
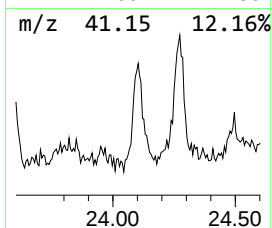
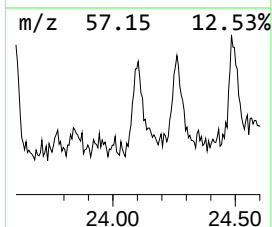
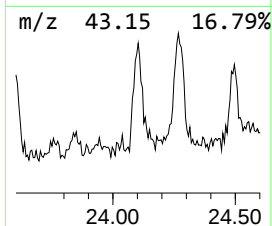
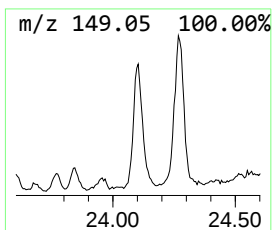
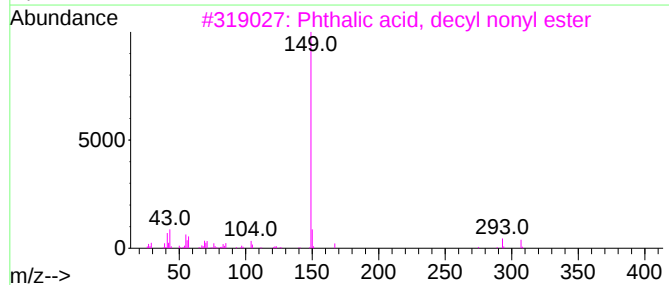
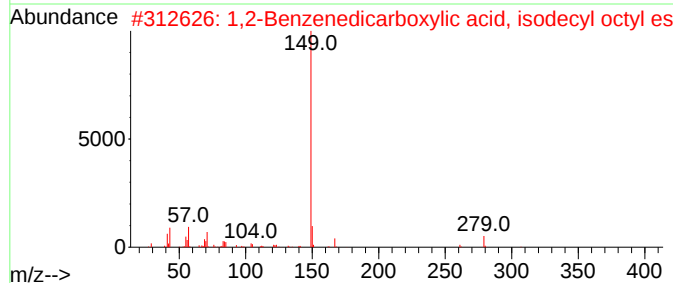
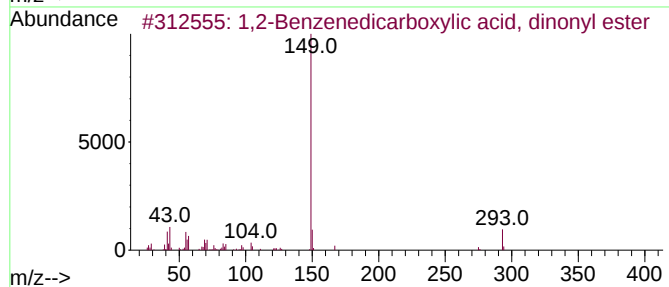
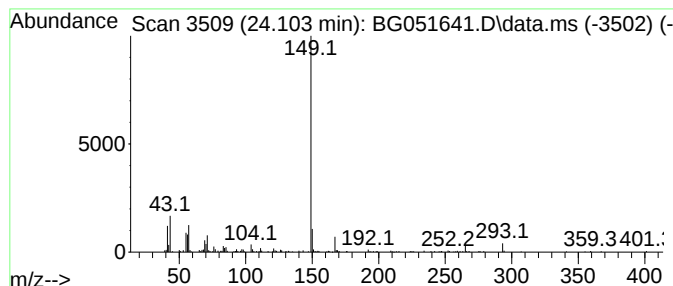
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Quant Title : SVOA CALIBRATION

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

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Peak Number 48 1,2-Benzenedicarboxylic aci... Concentration Rank 11

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|--------|------------------|--------------|------|
| 24.103    | 9.38 ng/ul                          | 273766 | Perylene-d12     | 25.477       |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#         | Qual |
| 1         | 1,2-Benzenedicarboxylic acid, di... | 418    | C26H42O4         | 000084-76-4  | 80   |
| 2         | 1,2-Benzenedicarboxylic acid, is... | 418    | C26H42O4         | 001330-96-7  | 80   |
| 3         | Phthalic acid, decyl nonyl ester    | 432    | C27H44O4         | 1000308-89-8 | 80   |
| 4         | Phthalic acid, dodecyl nonyl ester  | 460    | C29H48O4         | 1000308-92-6 | 80   |
| 5         | Phthalic acid, hept-2-yl tetrade... | 460    | C29H48O4         | 1000356-98-1 | 72   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG121721\  
Data File : BG051641.D  
Acq On : 18 Dec 2021 22:49  
Operator : CG/JU  
Sample : M5153-06  
Misc :  
ALS Vial : 57 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL71

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

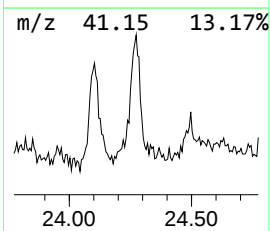
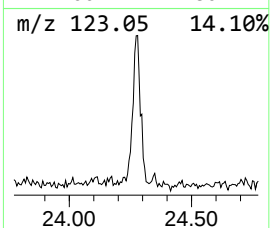
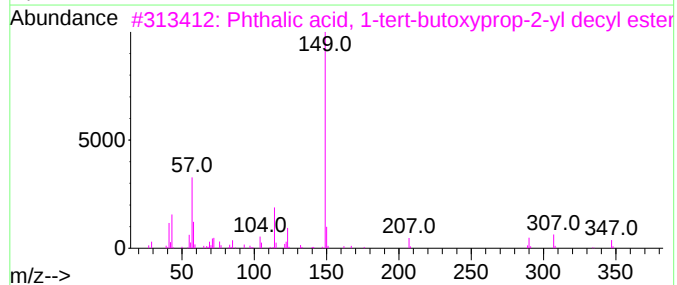
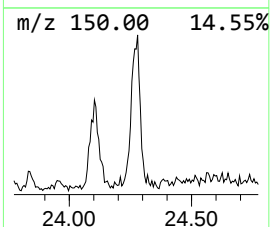
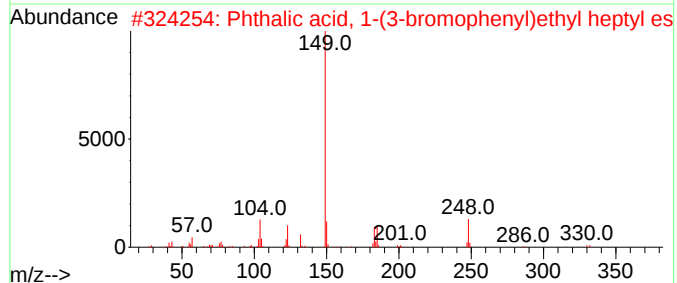
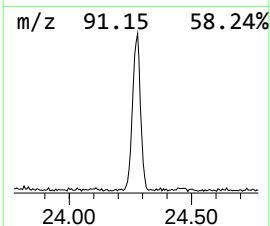
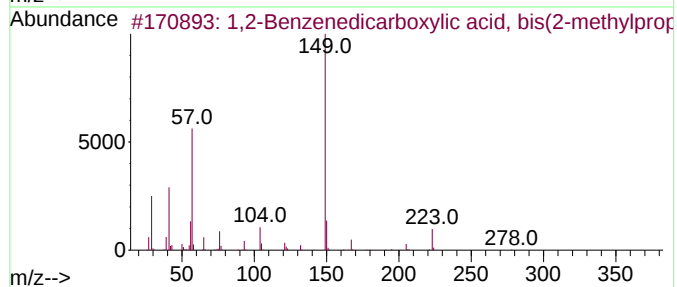
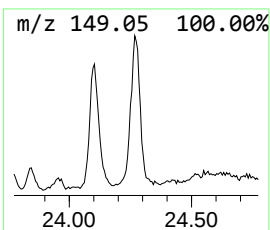
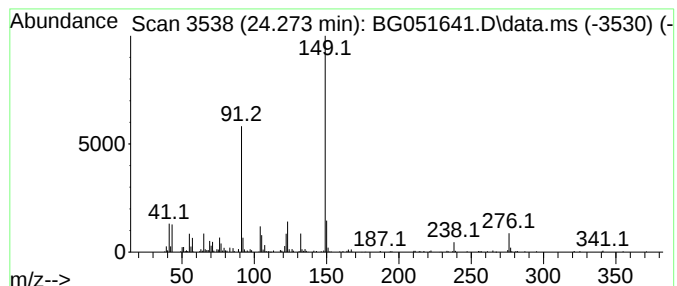
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 49 1,2-Benzenedicarboxylic aci... Concentration Rank 4

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 24.273 | 16.52 ng/ul | 482326 | Perylene-d12     | 25.477 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|------------|--------------|------|
| 1    |    |   | 1,2-Benzenedicarboxylic acid, bi... | 278 | C16H22O4   | 000084-69-5  | 64   |
| 2    |    |   | Phthalic acid, 1-(3-bromophenyl)... | 446 | C23H27BrO4 | 1000377-88-0 | 64   |
| 3    |    |   | Phthalic acid, 1-tert-butoxyprop... | 420 | C25H40O5   | 1000371-01-5 | 64   |
| 4    |    |   | 1,2-Benzenedicarboxylic acid, mo... | 278 | C16H22O4   | 068296-97-9  | 59   |
| 5    |    |   | Phthalic acid, 2-hexyl ester        | 250 | C14H18O4   | 079107-80-5  | 59   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG121721\  
Data File : BG051641.D  
Acq On : 18 Dec 2021 22:49  
Operator : CG/JU  
Sample : M5153-06  
Misc :  
ALS Vial : 57 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL71

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

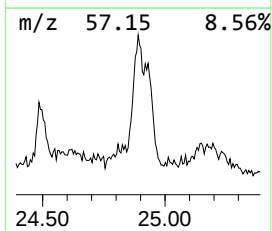
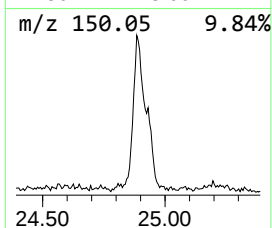
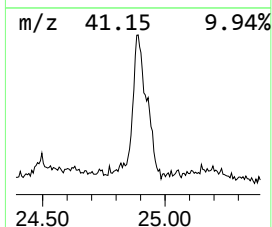
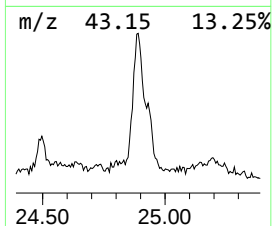
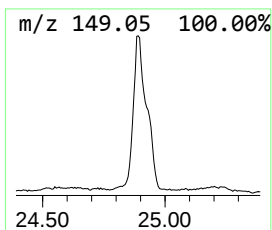
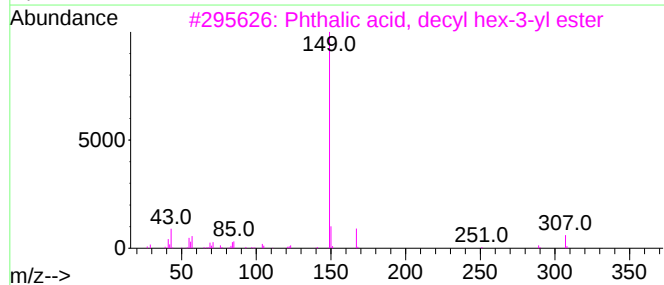
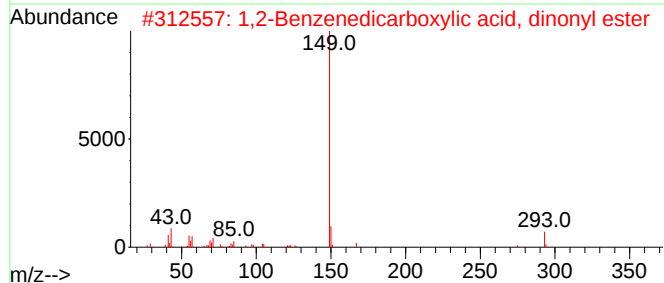
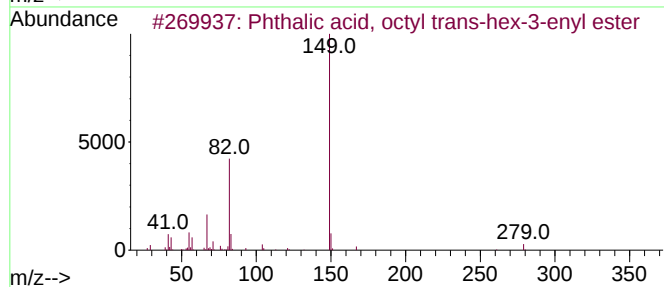
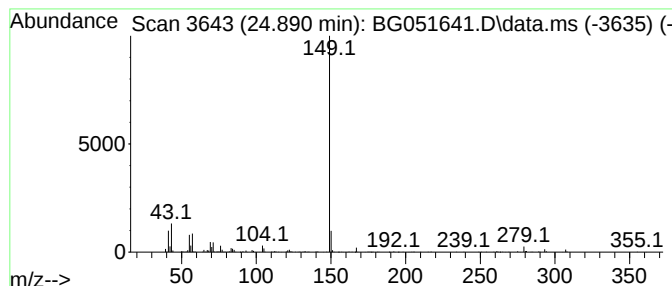
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 50 Phthalic acid, octyl trans-... Concentration Rank 2

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 24.890 | 43.32 ng/ul | 1264800 | Perylene-d12     | 25.477 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Phthalic acid, octyl trans-hex-3... | 360 | C22H32O4 | 1000360-48-6 | 80   |
| 2    |    |   | 1,2-Benzenedicarboxylic acid, di... | 418 | C26H42O4 | 000084-76-4  | 80   |
| 3    |    |   | Phthalic acid, decyl hex-3-yl ester | 390 | C24H38O4 | 1000356-96-1 | 80   |
| 4    |    |   | Di-n-octyl phthalate                | 390 | C24H38O4 | 000117-84-0  | 72   |
| 5    |    |   | Phthalic acid, 3,3-dimethylbut-2... | 376 | C23H36O4 | 1000357-00-8 | 72   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG121721\  
Data File : BG051641.D  
Acq On : 18 Dec 2021 22:49  
Operator : CG/JU  
Sample : M5153-06  
Misc :  
ALS Vial : 57 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL71

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

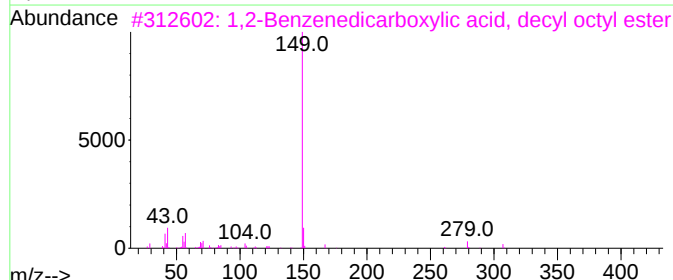
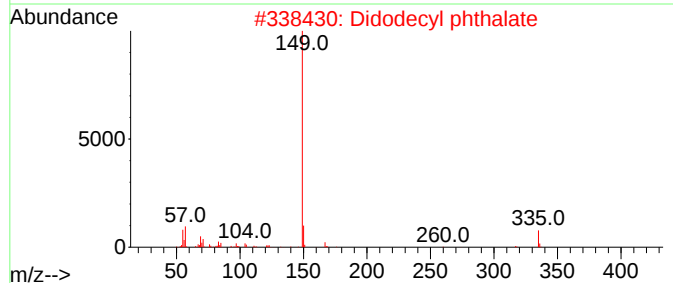
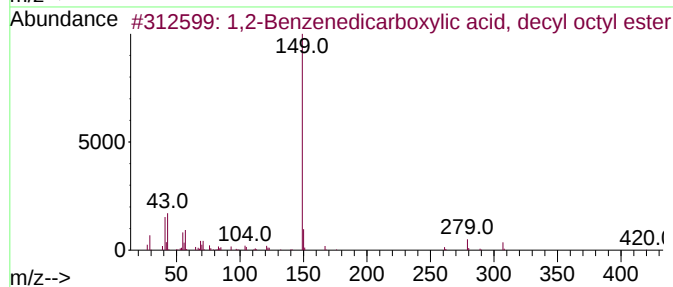
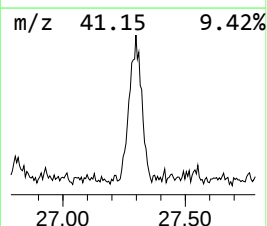
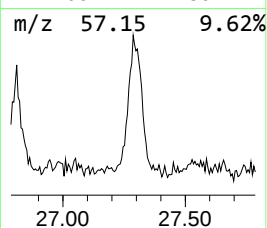
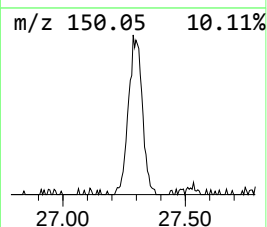
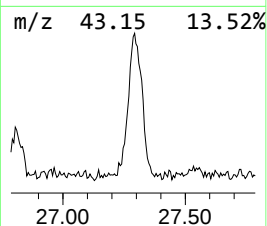
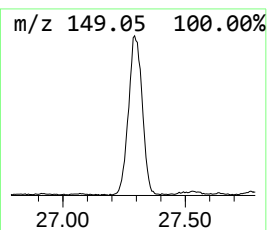
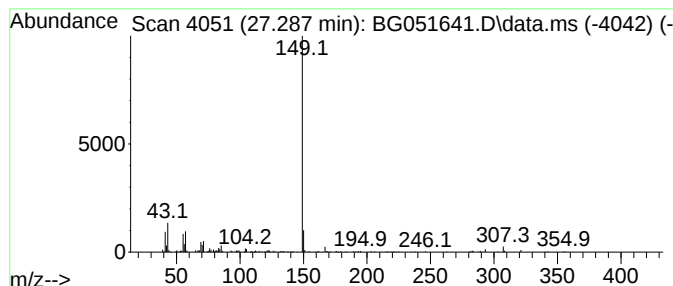
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 51 1,2-Benzenedicarboxylic aci... Concentration Rank 3

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 27.287 | 22.31 ng/ul | 651455 | Perylene-d12     | 25.477 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | 1,2-Benzenedicarboxylic acid, de... | 418 | C26H42O4 | 000119-07-3  | 87   |
| 2    |    |   | Didodecyl phthalate                 | 502 | C32H54O4 | 002432-90-8  | 83   |
| 3    |    |   | 1,2-Benzenedicarboxylic acid, de... | 418 | C26H42O4 | 000119-07-3  | 80   |
| 4    |    |   | Phthalic acid, decyl dec-2-yl ester | 446 | C28H46O4 | 1000356-94-5 | 80   |
| 5    |    |   | 1,2-Benzenedicarboxylic acid, di... | 474 | C30H50O4 | 003648-20-2  | 80   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG121721\  
 Data File : BG051641.D  
 Acq On : 18 Dec 2021 22:49  
 Operator : CG/JU  
 Sample : M5153-06  
 Misc :  
 ALS Vial : 57 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 BGL71

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.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 |
| (DEL) Alkane: S...  | 12.119 | 2.7     | ng/ul | 53382    | 2                     | 11.115 | 395956 | 20.0 |
| (DEL) Alkane: S...  | 12.507 | 4.5     | ng/ul | 89892    | 2                     | 11.115 | 395956 | 20.0 |
| (DEL) Alkane: S...  | 13.447 | 2.0     | ng/ul | 49507    | 3                     | 14.909 | 485741 | 20.0 |
| Naphthalene, 1,...  | 14.234 | 8.1     | ng/ul | 196802   | 3                     | 14.909 | 485741 | 20.0 |
| (DEL) Alkane: S...  | 14.381 | 2.0     | ng/ul | 49645    | 3                     | 14.909 | 485741 | 20.0 |
| (DEL) Alkane: S...  | 14.792 | 2.7     | ng/ul | 66182    | 3                     | 14.909 | 485741 | 20.0 |
| unknown-01          | 15.691 | 11.0    | ng/ul | 266811   | 3                     | 14.909 | 485741 | 20.0 |
| (DEL) Alkane: S...  | 16.625 | 8.9     | ng/ul | 257937   | 4                     | 17.659 | 580674 | 20.0 |
| Silane, methyl...   | 16.725 | 9.6     | ng/ul | 277975   | 4                     | 17.659 | 580674 | 20.0 |
| 1-(Isocyanatome...  | 16.778 | 9.6     | ng/ul | 278539   | 4                     | 17.659 | 580674 | 20.0 |
| 4-(7-Methylocty...  | 16.842 | 5.8     | ng/ul | 168660   | 4                     | 17.659 | 580674 | 20.0 |
| unknown-02          | 17.048 | 7.8     | ng/ul | 225194   | 4                     | 17.659 | 580674 | 20.0 |
| Phenol, 4-(1,1-...  | 17.107 | 5.5     | ng/ul | 160178   | 4                     | 17.659 | 580674 | 20.0 |
| unknown-03          | 17.183 | 4.4     | ng/ul | 128876   | 4                     | 17.659 | 580674 | 20.0 |
| Benzene, (1-met...  | 17.541 | 3.5     | ng/ul | 101515   | 4                     | 17.659 | 580674 | 20.0 |
| Nonaethylene gl...  | 17.935 | 2.9     | ng/ul | 84598    | 4                     | 17.659 | 580674 | 20.0 |
| 1,3-Dioxolane, ...  | 18.005 | 53.1    | ng/ul | 1542650  | 4                     | 17.659 | 580674 | 20.0 |
| (DEL) Alkane: S...  | 18.134 | 2.1     | ng/ul | 62403    | 4                     | 17.659 | 580674 | 20.0 |
| unknown-04          | 18.522 | 2.0     | ng/ul | 58856    | 4                     | 17.659 | 580674 | 20.0 |
| unknown-05          | 18.663 | 2.5     | ng/ul | 74053    | 4                     | 17.659 | 580674 | 20.0 |
| unknown-06          | 18.716 | 2.1     | ng/ul | 62396    | 4                     | 17.659 | 580674 | 20.0 |
| unknown-07          | 19.010 | 4.6     | ng/ul | 133041   | 4                     | 17.659 | 580674 | 20.0 |
| Phenylmethanedi...  | 19.198 | 10.8    | ng/ul | 312471   | 4                     | 17.659 | 580674 | 20.0 |
| cyclohexane, [1...  | 19.627 | 3.2     | ng/ul | 93429    | 4                     | 17.659 | 580674 | 20.0 |
| Ethanol, 2-[2-[...  | 19.897 | 4.8     | ng/ul | 207671   | 5                     | 21.976 | 863464 | 20.0 |
| Triphenyl phosph... | 21.260 | 2.5     | ng/ul | 106097   | 5                     | 21.976 | 863464 | 20.0 |
| unknown-08          | 21.301 | 7.2     | ng/ul | 309544   | 5                     | 21.976 | 863464 | 20.0 |
| 13H-Dibenzo[a,i...  | 21.424 | 10.8    | ng/ul | 466635   | 5                     | 21.976 | 863464 | 20.0 |
| 1,2-Benzenedica...  | 21.471 | 5.0     | ng/ul | 218064   | 5                     | 21.976 | 863464 | 20.0 |
| (DEL) Alkane: S...  | 21.589 | 3.0     | ng/ul | 130261   | 5                     | 21.976 | 863464 | 20.0 |
| 3,6,9,12-Tetrao...  | 21.794 | 3.0     | ng/ul | 129322   | 5                     | 21.976 | 863464 | 20.0 |
| Benzamide, 4-ni...  | 22.188 | 7.1     | ng/ul | 307558   | 5                     | 21.976 | 863464 | 20.0 |
| Phthalic acid, ...  | 22.669 | 14.4    | ng/ul | 623871   | 5                     | 21.976 | 863464 | 20.0 |
| 1,2-Benzenedica...  | 24.103 | 9.4     | ng/ul | 273766   | 6                     | 25.477 | 583912 | 20.0 |
| 1,2-Benzenedica...  | 24.273 | 16.5    | ng/ul | 482326   | 6                     | 25.477 | 583912 | 20.0 |
| Phthalic acid, ...  | 24.890 | 43.3    | ng/ul | 1264800  | 6                     | 25.477 | 583912 | 20.0 |
| 1,2-Benzenedica...  | 27.287 | 22.3    | ng/ul | 651455   | 6                     | 25.477 | 583912 | 20.0 |