

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122021\  
 Data File : BG051727.D  
 Acq On : 21 Dec 2021 22:20  
 Operator : CG/JU  
 Sample : M5121-04DL 10X  
 Misc :  
 ALS Vial : 54 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 BGL53DL

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: BG051727.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         | 5.874       | 402           | 406         | 414          | rBV2     | 7671           | 14162         | 0.19%           | 0.045%        |
| 2         | 7.407       | 661           | 667         | 675          | rVB4     | 19156          | 38052         | 0.50%           | 0.121%        |
| 3         | 7.577       | 688           | 696         | 703          | rBV      | 11683          | 20727         | 0.27%           | 0.066%        |
| 4         | 7.795       | 727           | 733         | 740          | rVB2     | 16141          | 28375         | 0.37%           | 0.090%        |
| 5         | 8.270       | 805           | 814         | 821          | rBV      | 218832         | 377425        | 4.93%           | 1.201%        |
| 6         | 8.964       | 924           | 932         | 938          | rVB5     | 19043          | 37262         | 0.49%           | 0.119%        |
| 7         | 9.439       | 1008          | 1013        | 1019         | rVB3     | 15970          | 25200         | 0.33%           | 0.080%        |
| 8         | 9.510       | 1019          | 1025        | 1029         | rBV2     | 10881          | 18168         | 0.24%           | 0.058%        |
| 9         | 10.174      | 1131          | 1138        | 1144         | rBV4     | 11433          | 24306         | 0.32%           | 0.077%        |
| 10        | 10.720      | 1225          | 1231        | 1237         | rVB2     | 18551          | 32861         | 0.43%           | 0.105%        |
| 11        | 10.967      | 1268          | 1273        | 1278         | rBV2     | 13753          | 23303         | 0.30%           | 0.074%        |
| 12        | 11.108      | 1284          | 1297        | 1303         | rBV      | 296746         | 576801        | 7.54%           | 1.835%        |
| 13        | 11.161      | 1303          | 1306        | 1312         | rVV      | 95327          | 151631        | 1.98%           | 0.482%        |
| 14        | 11.237      | 1312          | 1319        | 1321         | rVV2     | 11456          | 23517         | 0.31%           | 0.075%        |
| 15        | 11.255      | 1321          | 1322        | 1328         | rVB3     | 13284          | 15149         | 0.20%           | 0.048%        |
| 16        | 12.112      | 1463          | 1468        | 1472         | rBV3     | 8889           | 15784         | 0.21%           | 0.050%        |
| 17        | 12.500      | 1526          | 1534        | 1541         | rBV      | 18469          | 28975         | 0.38%           | 0.092%        |
| 18        | 12.747      | 1573          | 1576        | 1585         | rVB      | 11164          | 17479         | 0.23%           | 0.056%        |
| 19        | 12.970      | 1610          | 1614        | 1620         | rVB3     | 7976           | 12538         | 0.16%           | 0.040%        |
| 20        | 13.440      | 1690          | 1694        | 1700         | rVB4     | 18222          | 29191         | 0.38%           | 0.093%        |
| 21        | 13.722      | 1733          | 1742        | 1753         | rBV2     | 61063          | 114004        | 1.49%           | 0.363%        |
| 22        | 14.068      | 1797          | 1801        | 1805         | rBV7     | 5819           | 11634         | 0.15%           | 0.037%        |
| 23        | 14.233      | 1825          | 1829        | 1835         | rVB7     | 12404          | 22266         | 0.29%           | 0.071%        |
| 24        | 14.292      | 1835          | 1839        | 1844         | rVB      | 40989          | 57289         | 0.75%           | 0.182%        |
| 25        | 14.345      | 1844          | 1848        | 1850         | rBV4     | 16580          | 24744         | 0.32%           | 0.079%        |
| 26        | 14.380      | 1850          | 1854        | 1857         | rVV      | 39722          | 52323         | 0.68%           | 0.166%        |
| 27        | 14.603      | 1886          | 1892        | 1897         | rVV      | 39710          | 62888         | 0.82%           | 0.200%        |
| 28        | 14.668      | 1899          | 1903        | 1906         | rVV5     | 10362          | 16862         | 0.22%           | 0.054%        |
| 29        | 14.709      | 1907          | 1910        | 1913         | rVV4     | 10946          | 17295         | 0.23%           | 0.055%        |
| 30        | 14.744      | 1913          | 1916        | 1919         | rVV3     | 20608          | 25060         | 0.33%           | 0.080%        |
| 31        | 14.785      | 1919          | 1923        | 1928         | rVB      | 74834          | 92786         | 1.21%           | 0.295%        |
| 32        | 14.909      | 1934          | 1944        | 1953         | rBV      | 493347         | 791628        | 10.35%          | 2.519%        |
| 33        | 14.991      | 1955          | 1958        | 1963         | rBV7     | 7967           | 13979         | 0.18%           | 0.044%        |
| 34        | 15.073      | 1969          | 1972        | 1975         | rVB4     | 11749          | 12369         | 0.16%           | 0.039%        |
| 35        | 15.185      | 1987          | 1991        | 1994         | rBV2     | 14949          | 18638         | 0.24%           | 0.059%        |
| 36        | 15.226      | 1995          | 1998        | 2003         | rVV7     | 9103           | 17756         | 0.23%           | 0.056%        |

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 BNA\_G  
 ClientSampleId :  
 BGL53DL

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

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Start Thrs: 0.2

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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 | 15.290 | 2005 | 2009 | 2014 | rVB6 | 22956 | 38541 | 0.50% | 0.123% |
| 38 | 15.402 | 2024 | 2028 | 2035 | rVB8 | 23036 | 47737 | 0.62% | 0.152% |
| 39 | 15.572 | 2053 | 2057 | 2062 | rBV7 | 17712 | 34775 | 0.45% | 0.111% |
| 40 | 15.672 | 2071 | 2074 | 2075 | rBV3 | 18414 | 20225 | 0.26% | 0.064% |

|    |        |      |      |      |      |       |        |       |        |
|----|--------|------|------|------|------|-------|--------|-------|--------|
| 41 | 15.696 | 2076 | 2078 | 2081 | rVV3 | 36079 | 49579  | 0.65% | 0.158% |
| 42 | 15.737 | 2081 | 2085 | 2091 | rVB  | 85973 | 121307 | 1.59% | 0.386% |
| 43 | 15.843 | 2099 | 2103 | 2108 | rBV3 | 26005 | 45391  | 0.59% | 0.144% |
| 44 | 15.895 | 2108 | 2112 | 2116 | rBV2 | 71864 | 107430 | 1.40% | 0.342% |
| 45 | 16.048 | 2136 | 2138 | 2140 | rBV2 | 20459 | 21613  | 0.28% | 0.069% |

|    |        |      |      |      |      |       |       |       |        |
|----|--------|------|------|------|------|-------|-------|-------|--------|
| 46 | 16.078 | 2140 | 2143 | 2147 | rVB  | 40721 | 53513 | 0.70% | 0.170% |
| 47 | 16.142 | 2150 | 2154 | 2158 | rBV3 | 55779 | 92032 | 1.20% | 0.293% |
| 48 | 16.236 | 2167 | 2170 | 2173 | rBV3 | 28004 | 35044 | 0.46% | 0.112% |
| 49 | 16.312 | 2179 | 2183 | 2186 | rBV  | 39838 | 58295 | 0.76% | 0.185% |
| 50 | 16.395 | 2194 | 2197 | 2200 | rBV  | 24050 | 22301 | 0.29% | 0.071% |

|    |        |      |      |      |      |        |        |       |        |
|----|--------|------|------|------|------|--------|--------|-------|--------|
| 51 | 16.524 | 2216 | 2219 | 2227 | rVB4 | 35408  | 70666  | 0.92% | 0.225% |
| 52 | 16.594 | 2227 | 2231 | 2233 | rBV  | 122732 | 156356 | 2.04% | 0.497% |
| 53 | 16.624 | 2233 | 2236 | 2241 | rVB2 | 143993 | 199150 | 2.60% | 0.634% |
| 54 | 16.724 | 2249 | 2253 | 2259 | rBV3 | 48462  | 96267  | 1.26% | 0.306% |
| 55 | 16.871 | 2275 | 2278 | 2282 | rVV  | 54802  | 70288  | 0.92% | 0.224% |

|    |        |      |      |      |      |        |        |        |        |
|----|--------|------|------|------|------|--------|--------|--------|--------|
| 56 | 16.912 | 2282 | 2285 | 2294 | rVB2 | 39547  | 74158  | 0.97%  | 0.236% |
| 57 | 17.393 | 2363 | 2367 | 2371 | rBV  | 94891  | 114745 | 1.50%  | 0.365% |
| 58 | 17.446 | 2373 | 2376 | 2380 | rVB2 | 85256  | 106375 | 1.39%  | 0.338% |
| 59 | 17.658 | 2405 | 2412 | 2417 | rBV  | 538254 | 923949 | 12.08% | 2.940% |
| 60 | 18.134 | 2489 | 2493 | 2499 | rVB  | 105475 | 156085 | 2.04%  | 0.497% |

|    |        |      |      |      |      |        |        |       |        |
|----|--------|------|------|------|------|--------|--------|-------|--------|
| 61 | 18.298 | 2519 | 2521 | 2528 | rVB4 | 62492  | 84672  | 1.11% | 0.269% |
| 62 | 18.821 | 2606 | 2610 | 2616 | rBV  | 86736  | 108790 | 1.42% | 0.346% |
| 63 | 19.085 | 2652 | 2655 | 2665 | rVB2 | 111089 | 161729 | 2.11% | 0.515% |
| 64 | 19.191 | 2670 | 2673 | 2682 | rVB2 | 114934 | 175884 | 2.30% | 0.560% |
| 65 | 19.438 | 2710 | 2715 | 2719 | rVB4 | 94900  | 159109 | 2.08% | 0.506% |

|    |        |      |      |      |      |        |        |       |        |
|----|--------|------|------|------|------|--------|--------|-------|--------|
| 66 | 19.626 | 2743 | 2747 | 2751 | rVB  | 85657  | 112262 | 1.47% | 0.357% |
| 67 | 19.902 | 2792 | 2794 | 2798 | rVB2 | 62050  | 67767  | 0.89% | 0.216% |
| 68 | 20.013 | 2810 | 2813 | 2819 | rVB3 | 131444 | 230221 | 3.01% | 0.733% |
| 69 | 20.542 | 2899 | 2903 | 2907 | rBV  | 207479 | 230084 | 3.01% | 0.732% |
| 70 | 20.806 | 2945 | 2948 | 2953 | rVB  | 79894  | 94531  | 1.24% | 0.301% |

|    |        |      |      |      |      |        |        |        |        |
|----|--------|------|------|------|------|--------|--------|--------|--------|
| 71 | 20.936 | 2965 | 2970 | 2974 | rVB  | 661801 | 864855 | 11.31% | 2.752% |
| 72 | 21.047 | 2986 | 2989 | 2993 | rVB  | 258001 | 291689 | 3.81%  | 0.928% |
| 73 | 21.312 | 3028 | 3034 | 3041 | rVB3 | 386872 | 776592 | 10.15% | 2.471% |
| 74 | 21.388 | 3045 | 3047 | 3053 | rVV  | 71984  | 88229  | 1.15%  | 0.281% |
| 75 | 21.464 | 3057 | 3060 | 3065 | rVB  | 203291 | 258369 | 3.38%  | 0.822% |

|    |        |      |      |      |     |        |        |       |        |
|----|--------|------|------|------|-----|--------|--------|-------|--------|
| 76 | 21.582 | 3076 | 3080 | 3086 | rVB | 292623 | 426324 | 5.57% | 1.356% |
|----|--------|------|------|------|-----|--------|--------|-------|--------|

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Integrator: RTE

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 Title : SVOA CALIBRATION

|     |        |      |      |      |      |         |         |         |         |
|-----|--------|------|------|------|------|---------|---------|---------|---------|
| 77  | 21.852 | 3119 | 3126 | 3136 | rVB  | 4796659 | 7649813 | 100.00% | 24.340% |
| 78  | 21.975 | 3140 | 3147 | 3154 | rVB3 | 476432  | 956749  | 12.51%  | 3.044%  |
| 79  | 22.175 | 3172 | 3181 | 3186 | rBV2 | 366483  | 800450  | 10.46%  | 2.547%  |
| 80  | 22.398 | 3215 | 3219 | 3223 | rVV2 | 84841   | 142123  | 1.86%   | 0.452%  |
| 81  | 22.463 | 3225 | 3230 | 3236 | rVB2 | 165785  | 280619  | 3.67%   | 0.893%  |
| 82  | 22.592 | 3246 | 3252 | 3257 | rVB  | 280655  | 543758  | 7.11%   | 1.730%  |
| 83  | 22.669 | 3259 | 3265 | 3271 | rVB3 | 313121  | 640205  | 8.37%   | 2.037%  |
| 84  | 22.839 | 3287 | 3294 | 3301 | rBV  | 303930  | 561872  | 7.34%   | 1.788%  |
| 85  | 23.168 | 3342 | 3350 | 3364 | rVB  | 726297  | 1787869 | 23.37%  | 5.689%  |
| 86  | 23.597 | 3417 | 3423 | 3436 | rVB2 | 323620  | 692102  | 9.05%   | 2.202%  |
| 87  | 23.949 | 3478 | 3483 | 3490 | rVB2 | 130752  | 265910  | 3.48%   | 0.846%  |
| 88  | 24.096 | 3502 | 3508 | 3520 | rVB  | 340113  | 920892  | 12.04%  | 2.930%  |
| 89  | 24.490 | 3569 | 3575 | 3582 | rVB2 | 214510  | 473825  | 6.19%   | 1.508%  |
| 90  | 24.889 | 3635 | 3643 | 3646 | rBV  | 446956  | 1110424 | 14.52%  | 3.533%  |
| 91  | 25.482 | 3736 | 3744 | 3749 | rBV2 | 265505  | 764983  | 10.00%  | 2.434%  |
| 92  | 26.129 | 3847 | 3854 | 3861 | rBV2 | 103952  | 326696  | 4.27%   | 1.039%  |
| 93  | 26.205 | 3862 | 3867 | 3879 | rVB2 | 123316  | 357580  | 4.67%   | 1.138%  |
| 94  | 26.804 | 3960 | 3969 | 3979 | rVB  | 172818  | 542348  | 7.09%   | 1.726%  |
| 95  | 27.298 | 4041 | 4053 | 4068 | rVB  | 426859  | 1492148 | 19.51%  | 4.748%  |
| 96  | 28.320 | 4217 | 4227 | 4237 | rBV2 | 109157  | 410883  | 5.37%   | 1.307%  |
| 97  | 29.066 | 4348 | 4354 | 4367 | rVB3 | 79528   | 273104  | 3.57%   | 0.869%  |
| 98  | 30.147 | 4524 | 4538 | 4551 | rBV4 | 91291   | 420232  | 5.49%   | 1.337%  |
| 99  | 30.728 | 4627 | 4637 | 4638 | rBV4 | 89116   | 209300  | 2.74%   | 0.666%  |
| 100 | 32.355 | 4905 | 4914 | 4916 | rBV7 | 34944   | 95324   | 1.25%   | 0.303%  |

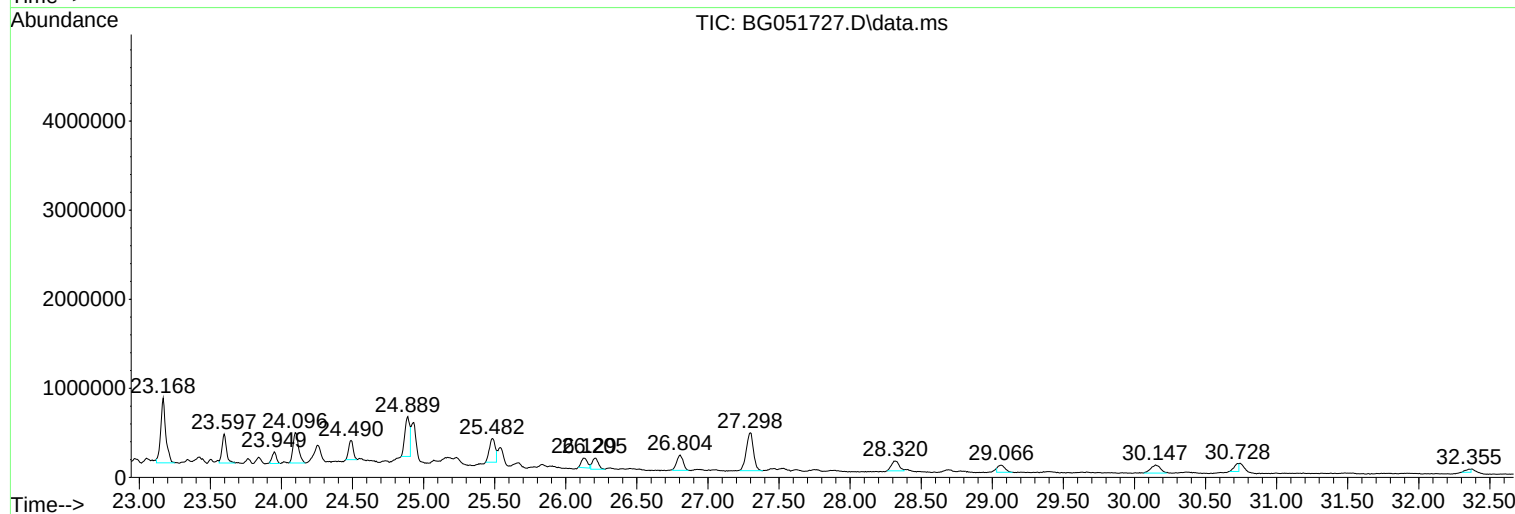
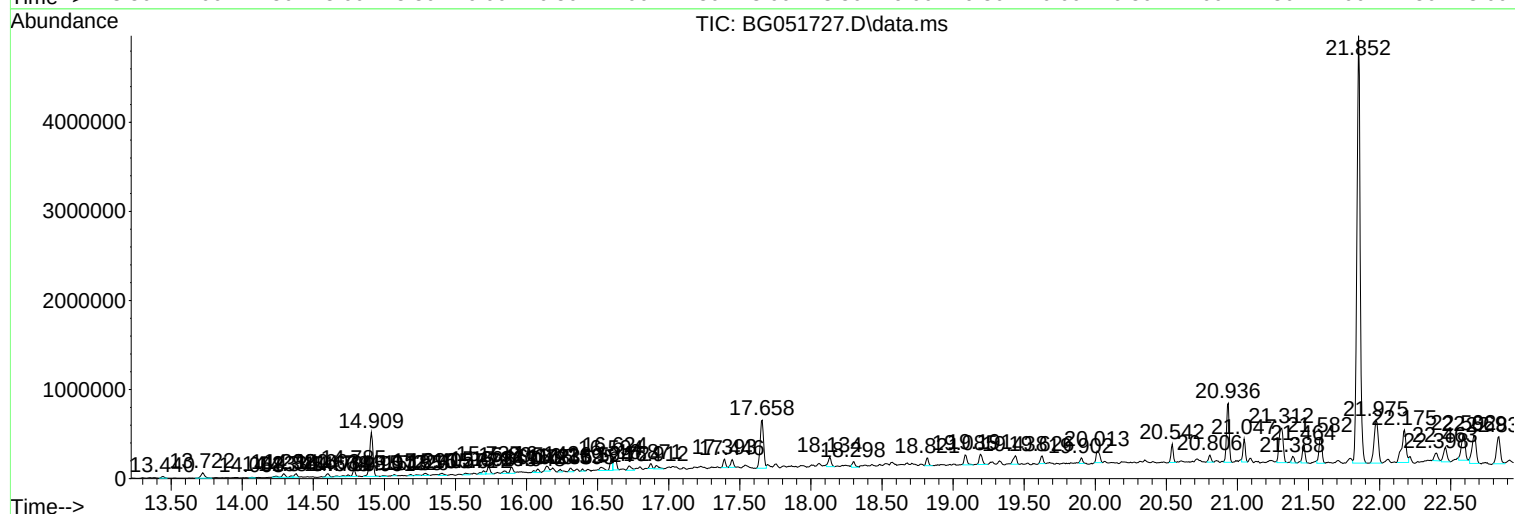
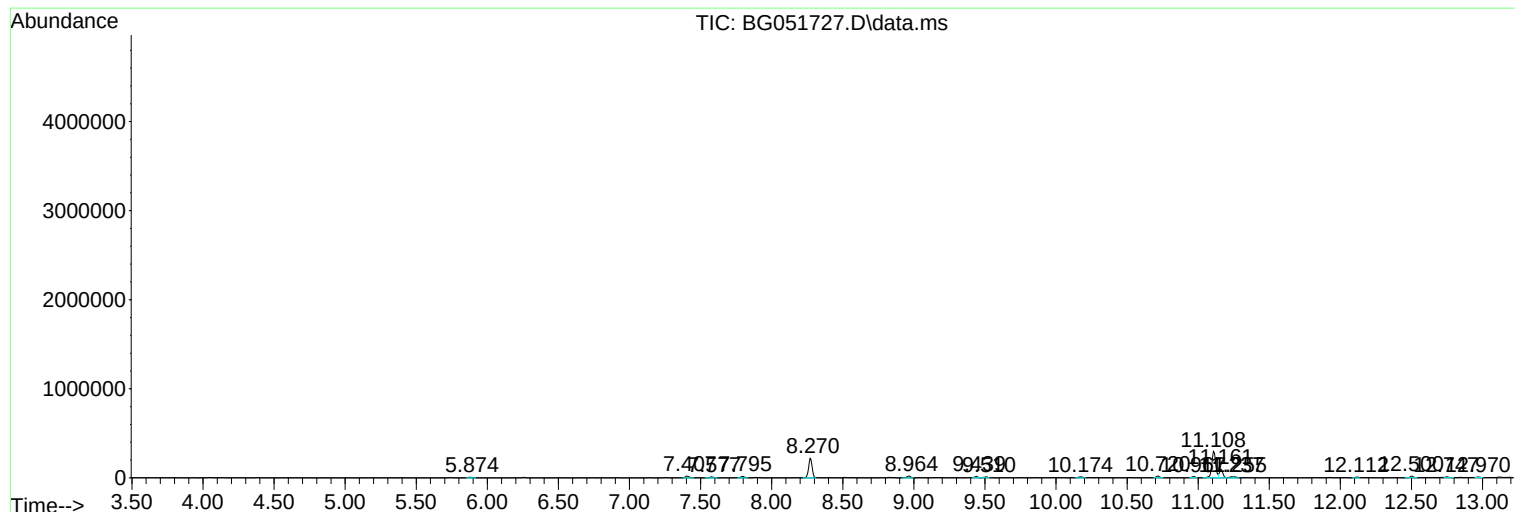
Sum of corrected areas: 31428565

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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



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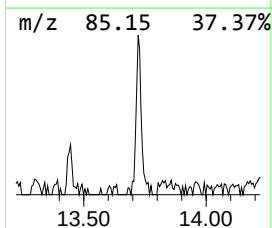
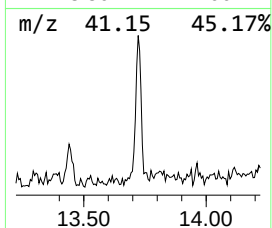
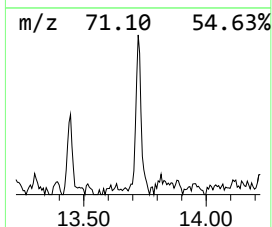
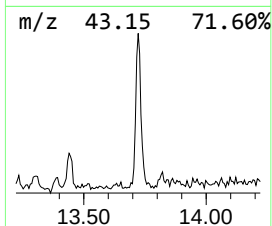
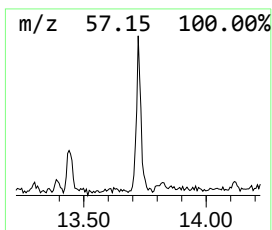
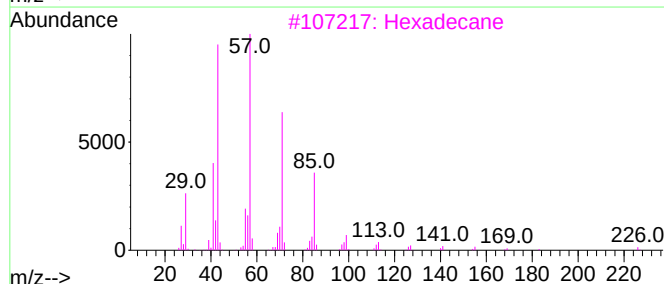
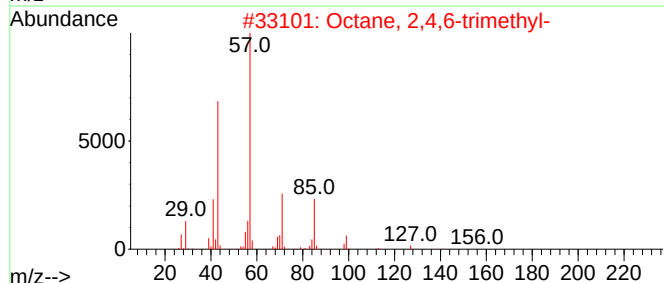
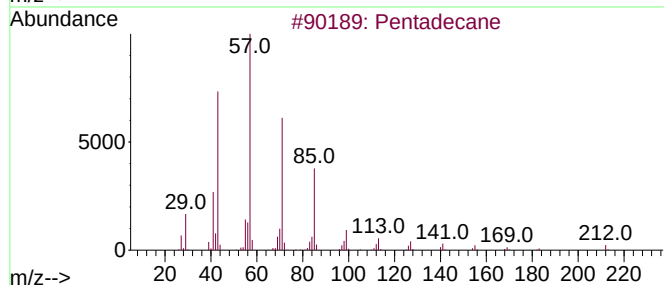
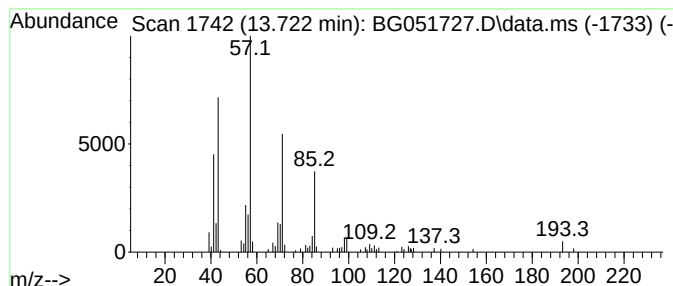
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

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

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

| R.T.      | EstConc                  | Area   | Relative to ISTD |         | R.T.        |      |
|-----------|--------------------------|--------|------------------|---------|-------------|------|
| 13.722    | 2.88 ng/ul               | 114004 | Acenaphthene-d10 |         | 14.909      |      |
| Hit# of 5 | Tentative ID             |        | MW               | MolForm | CAS#        | Qual |
| 1         | Pentadecane              |        | 212              | C15H32  | 000629-62-9 | 86   |
| 2         | Octane, 2,4,6-trimethyl- |        | 156              | C11H24  | 062016-37-9 | 80   |
| 3         | Hexadecane               |        | 226              | C16H34  | 000544-76-3 | 78   |
| 4         | Tridecane                |        | 184              | C13H28  | 000629-50-5 | 78   |
| 5         | Decane, 2,3,5-trimethyl- |        | 184              | C13H28  | 062238-11-3 | 78   |



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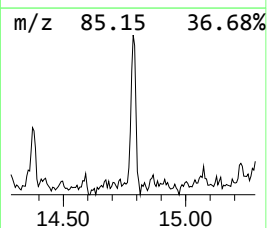
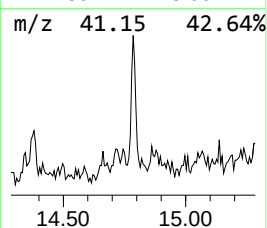
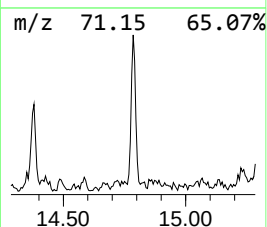
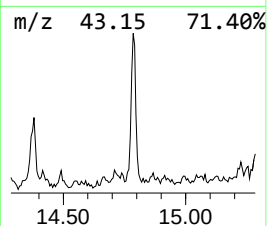
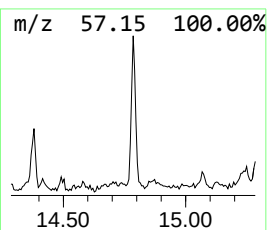
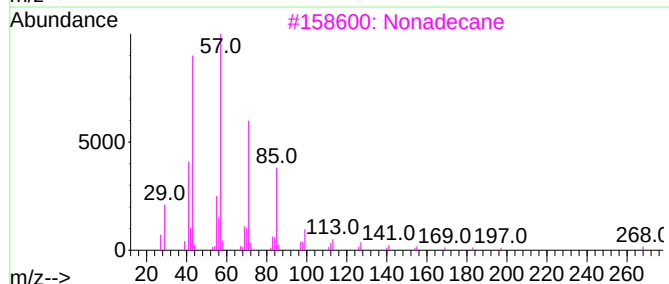
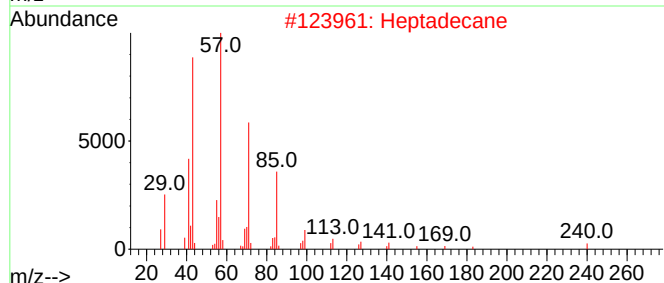
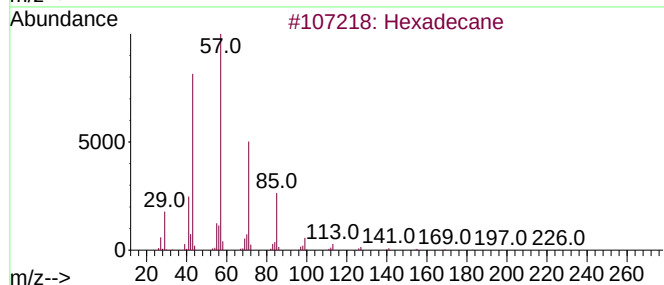
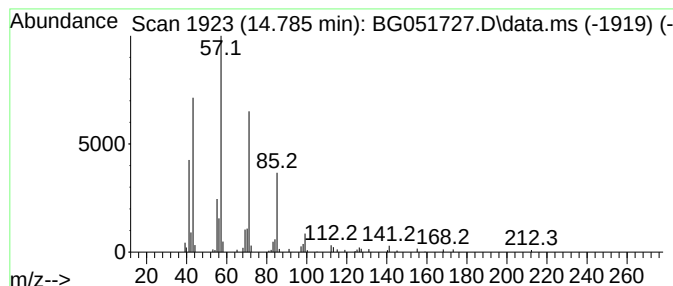
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 37

| R.T.   | EstConc    | Area  | Relative to ISTD | R.T.   |
|--------|------------|-------|------------------|--------|
| 14.785 | 2.34 ng/ul | 92786 | Acenaphthene-d10 | 14.909 |

| Hit# | of | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------|-----|---------|-------------|------|
| 1    |    |   | Hexadecane   | 226 | C16H34  | 000544-76-3 | 90   |
| 2    |    |   | Heptadecane  | 240 | C17H36  | 000629-78-7 | 90   |
| 3    |    |   | Nonadecane   | 268 | C19H40  | 000629-92-5 | 83   |
| 4    |    |   | Octacosane   | 394 | C28H58  | 000630-02-4 | 78   |
| 5    |    |   | Tridecane    | 184 | C13H28  | 000629-50-5 | 72   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122021\  
Data File : BG051727.D  
Acq On : 21 Dec 2021 22:20  
Operator : CG/JU  
Sample : M5121-04DL 10X  
Misc :  
ALS Vial : 54 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL53DL

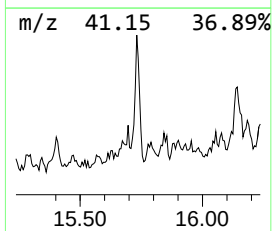
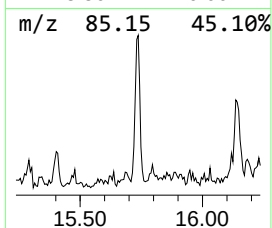
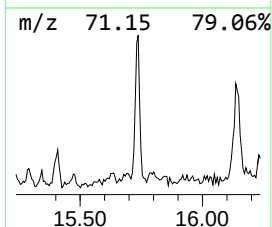
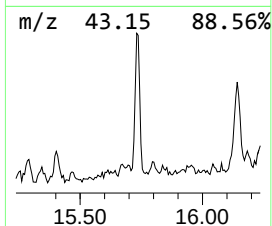
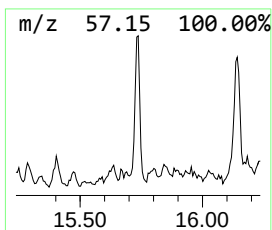
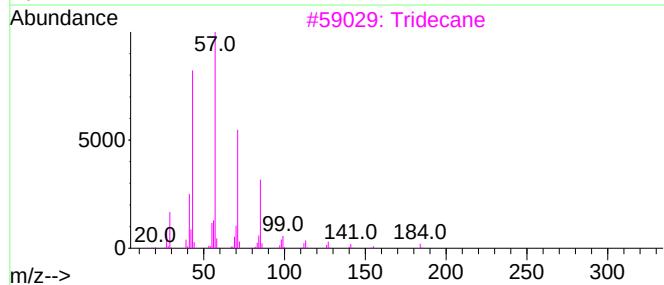
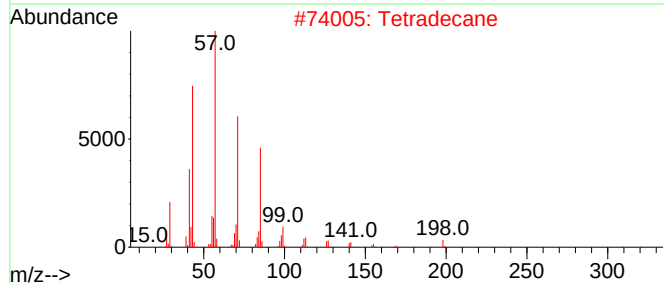
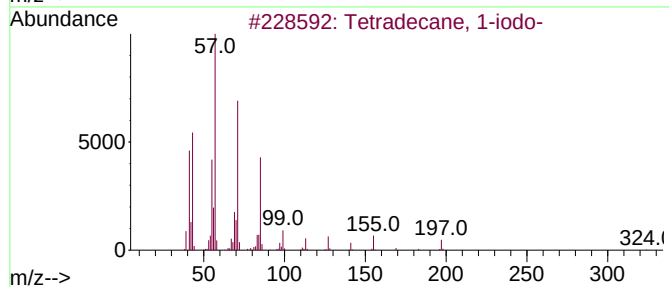
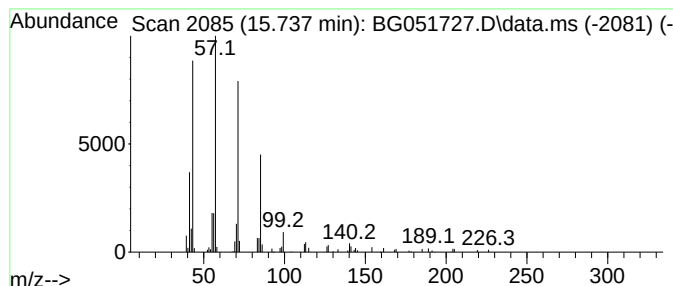
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Quant Title : SVOA CALIBRATION

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

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Peak Number 3 Tetradecane, 1-iodo- Concentration Rank 30

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 15.737 | 3.06 ng/ul | 121307 | Acenaphthene-d10 | 14.909 |

| Hit# | of | 5 | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------|-----|---------|-------------|------|
| 1    |    |   | Tetradecane, 1-iodo-    | 324 | C14H29I | 019218-94-1 | 90   |
| 2    |    |   | Tetradecane             | 198 | C14H30  | 000629-59-4 | 86   |
| 3    |    |   | Tridecane               | 184 | C13H28  | 000629-50-5 | 86   |
| 4    |    |   | Undecane, 2,8-dimethyl- | 184 | C13H28  | 017301-25-6 | 83   |
| 5    |    |   | Heptadecane             | 240 | C17H36  | 000629-78-7 | 80   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122021\  
Data File : BG051727.D  
Acq On : 21 Dec 2021 22:20  
Operator : CG/JU  
Sample : M5121-04DL 10X  
Misc :  
ALS Vial : 54 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL53DL

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

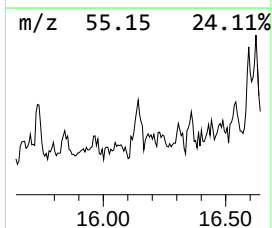
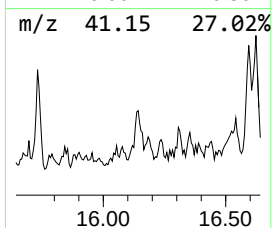
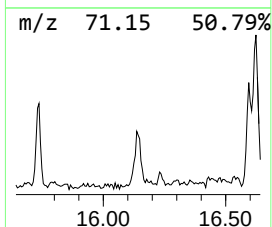
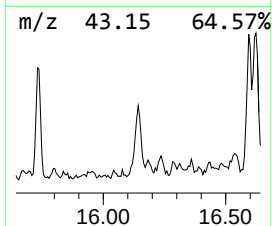
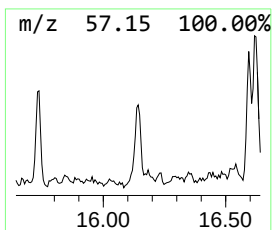
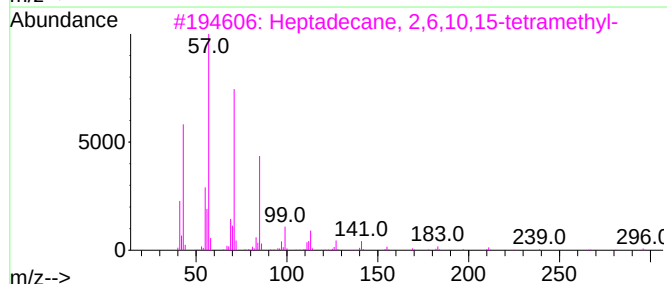
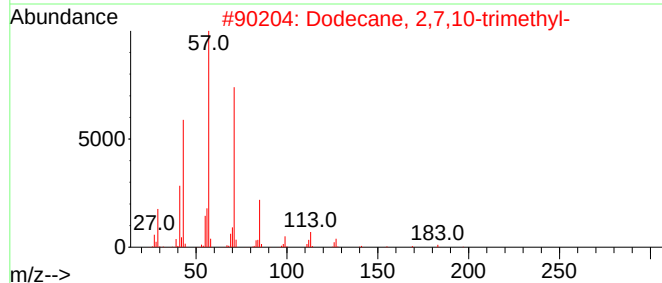
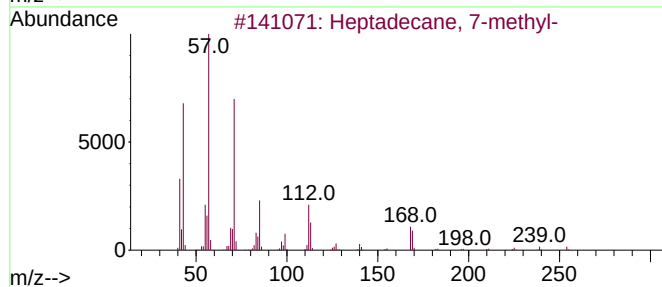
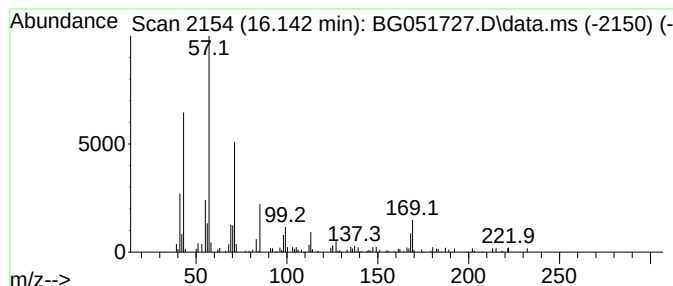
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TIC Integration Parameters: LSCINT.P

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Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 38

| R.T.   | EstConc    | Area  | Relative to ISTD | R.T.   |
|--------|------------|-------|------------------|--------|
| 16.142 | 2.33 ng/ul | 92032 | Acenaphthene-d10 | 14.909 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Heptadecane, 7-methyl-              | 254 | C18H38  | 020959-33-5 | 78   |
| 2    |    |   | Dodecane, 2,7,10-trimethyl-         | 212 | C15H32  | 074645-98-0 | 64   |
| 3    |    |   | Heptadecane, 2,6,10,15-tetramethyl- | 296 | C21H44  | 054833-48-6 | 59   |
| 4    |    |   | Undecane                            | 156 | C11H24  | 001120-21-4 | 53   |
| 5    |    |   | Dodecane, 2,6,10-trimethyl-         | 212 | C15H32  | 003891-98-3 | 50   |





Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122021\  
Data File : BG051727.D  
Acq On : 21 Dec 2021 22:20  
Operator : CG/JU  
Sample : M5121-04DL 10X  
Misc :  
ALS Vial : 54 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL53DL

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

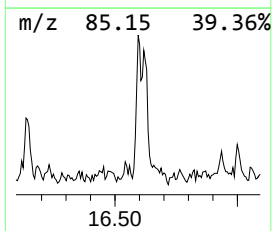
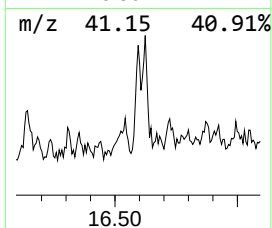
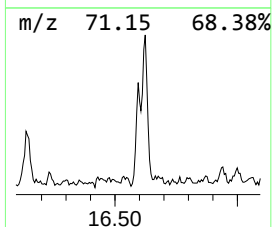
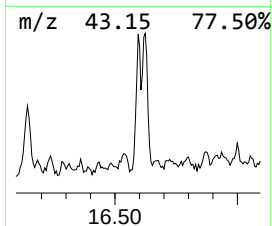
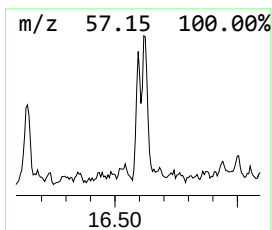
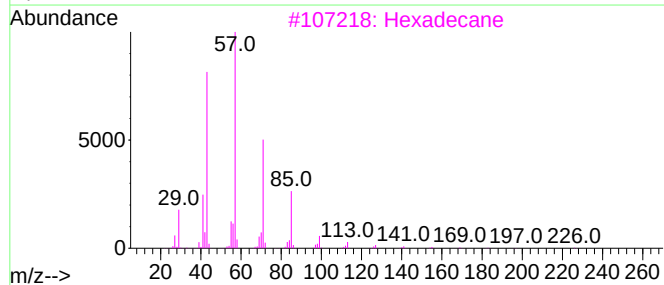
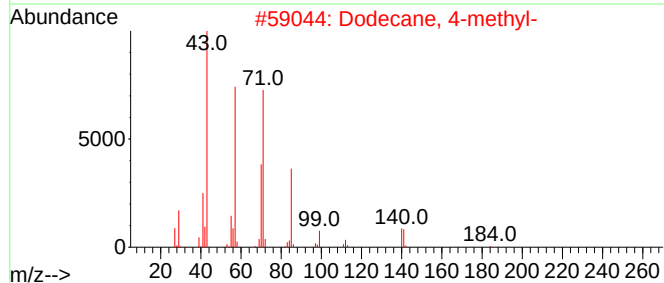
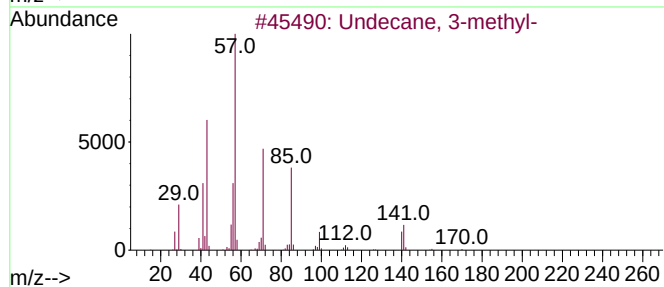
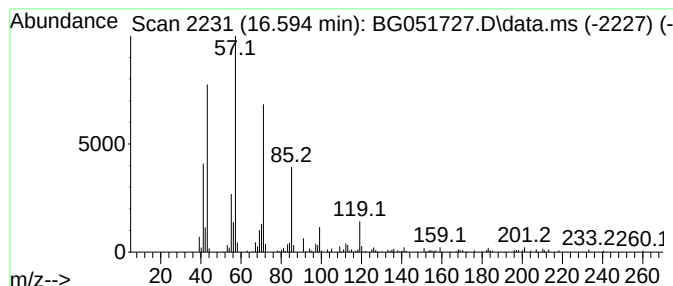
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 28

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 16.595 | 3.38 ng/ul | 156356 | Phenanthrene-d10 | 17.658 |

| Hit# | of | 5 | Tentative ID         | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------|-----|---------|-------------|------|
| 1    |    |   | Undecane, 3-methyl-  | 170 | C12H26  | 001002-43-3 | 83   |
| 2    |    |   | Dodecane, 4-methyl-  | 184 | C13H28  | 006117-97-1 | 80   |
| 3    |    |   | Hexadecane           | 226 | C16H34  | 000544-76-3 | 72   |
| 4    |    |   | Tridecane, 7-propyl- | 226 | C16H34  | 055045-09-5 | 72   |
| 5    |    |   | Eicosane             | 282 | C20H42  | 000112-95-8 | 72   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122021\  
Data File : BG051727.D  
Acq On : 21 Dec 2021 22:20  
Operator : CG/JU  
Sample : M5121-04DL 10X  
Misc :  
ALS Vial : 54 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL53DL

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

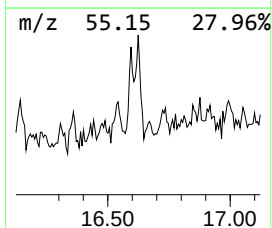
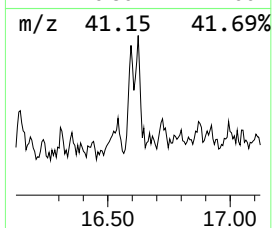
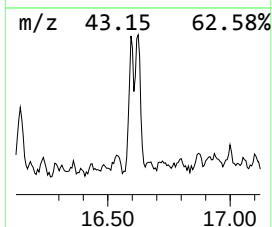
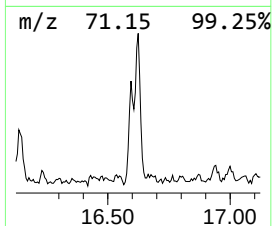
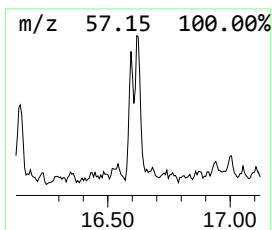
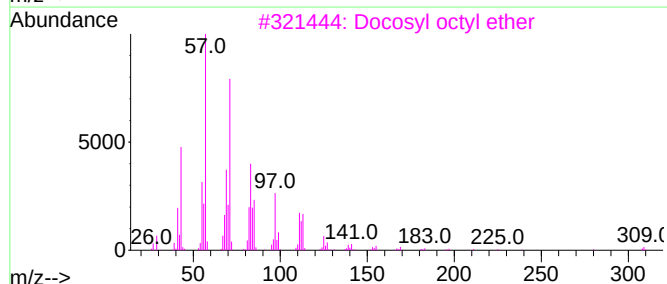
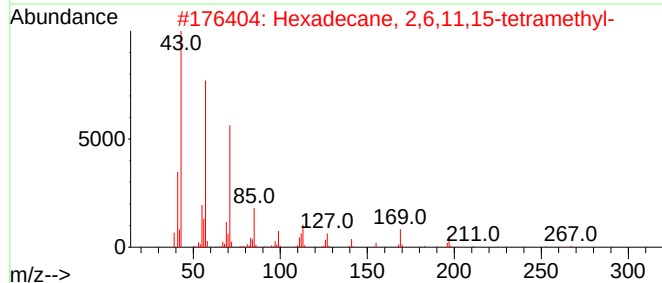
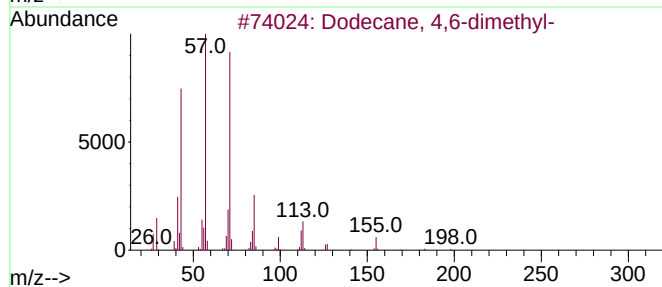
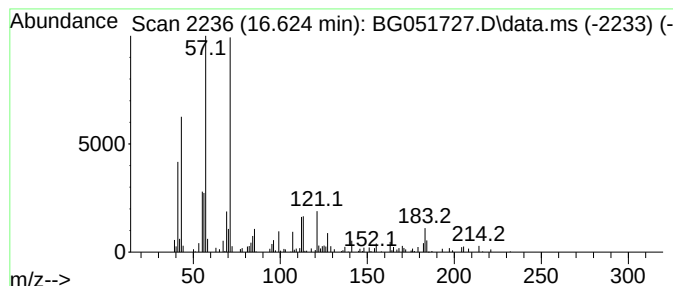
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 24

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 16.624 | 4.31 ng/ul | 199150 | Phenanthrene-d10 | 17.658 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Dodecane, 4,6-dimethyl-             | 198 | C14H30    | 061141-72-8  | 59   |
| 2    |    |   | Hexadecane, 2,6,11,15-tetramethyl-  | 282 | C20H42    | 000504-44-9  | 59   |
| 3    |    |   | Docosyl octyl ether                 | 438 | C30H62O   | 1000406-38-9 | 59   |
| 4    |    |   | Sulfurous acid, decyl 2-ethylhex... | 334 | C18H38O3S | 1000309-19-3 | 59   |
| 5    |    |   | Tridecane, 7-methyl-                | 198 | C14H30    | 026730-14-3  | 53   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122021\  
Data File : BG051727.D  
Acq On : 21 Dec 2021 22:20  
Operator : CG/JU  
Sample : M5121-04DL 10X  
Misc :  
ALS Vial : 54 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL53DL

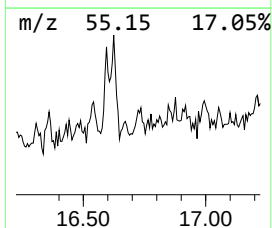
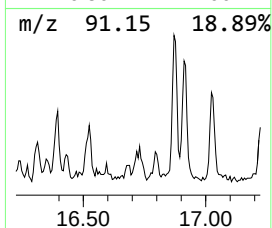
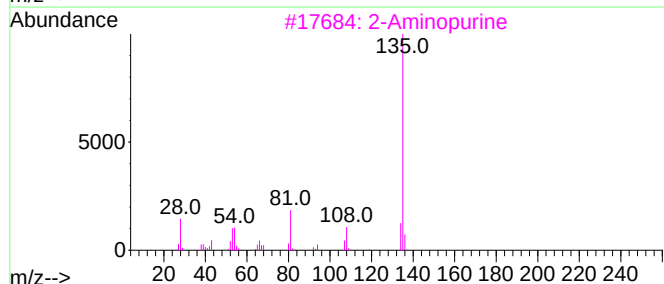
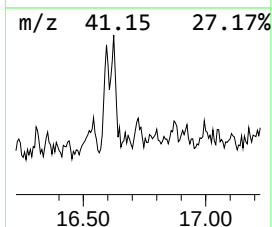
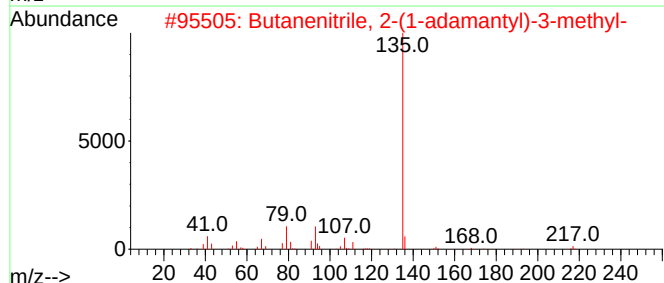
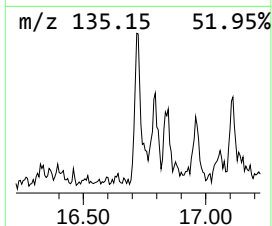
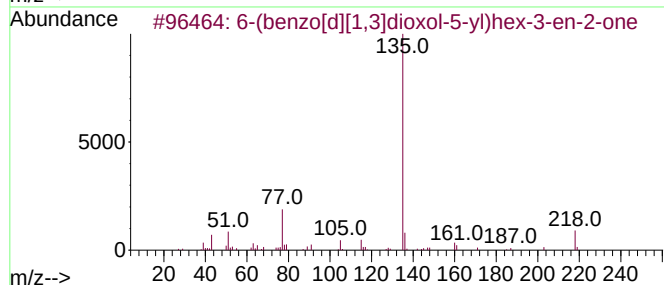
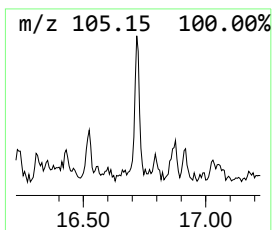
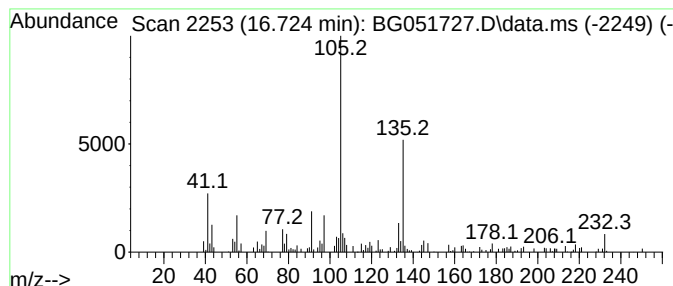
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Quant Title : SVOA CALIBRATION

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

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Peak Number 7 unknown-01 Concentration Rank 40

| R.T.   | EstConc    | Area  | Relative to ISTD | R.T.   |
|--------|------------|-------|------------------|--------|
| 16.724 | 2.08 ng/ul | 96267 | Phenanthrene-d10 | 17.658 |

| Hit# | of                                  | 5   | Tentative ID | MW           | MolForm | CAS# | Qual |
|------|-------------------------------------|-----|--------------|--------------|---------|------|------|
| 1    | 6-(benzo[d][1,3]dioxol-5-yl)hex-... | 218 | C13H14O3     | 075787-97-2  | 35      |      |      |
| 2    | Butanenitrile, 2-(1-adamantyl)-3... | 217 | C15H23N      | 1000266-44-6 | 30      |      |      |
| 3    | 2-Aminopurine                       | 135 | C5H5N5       | 000452-06-2  | 30      |      |      |
| 4    | Cyclopropane, 1-bromo-2,2,3,3-te... | 214 | C10H15Br     | 138777-60-3  | 11      |      |      |
| 5    | (.+/-)-2-Phenylpropanoic Acid, ...  | 192 | C12H16O2     | 1000453-23-1 | 11      |      |      |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122021\  
Data File : BG051727.D  
Acq On : 21 Dec 2021 22:20  
Operator : CG/JU  
Sample : M5121-04DL 10X  
Misc :  
ALS Vial : 54 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL53DL

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

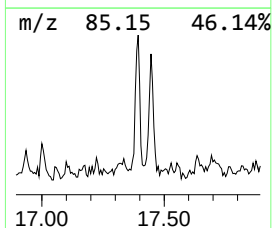
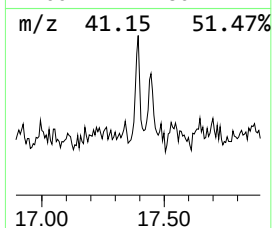
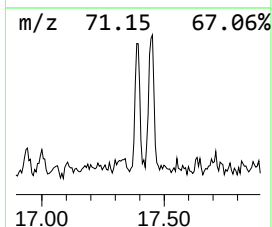
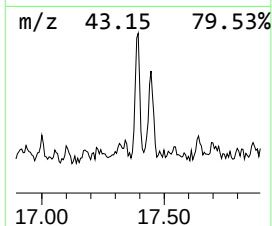
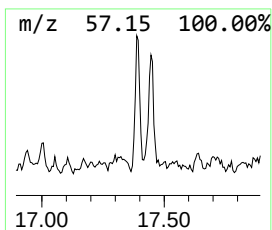
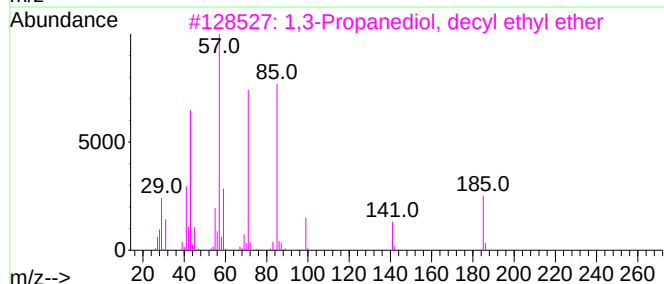
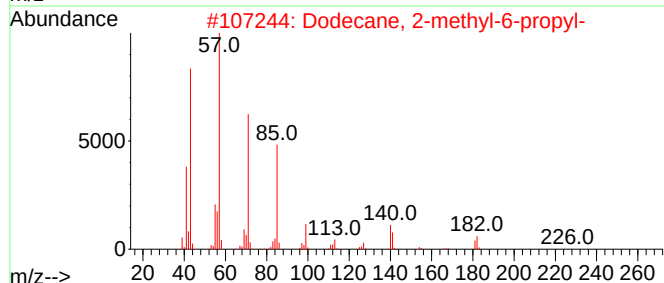
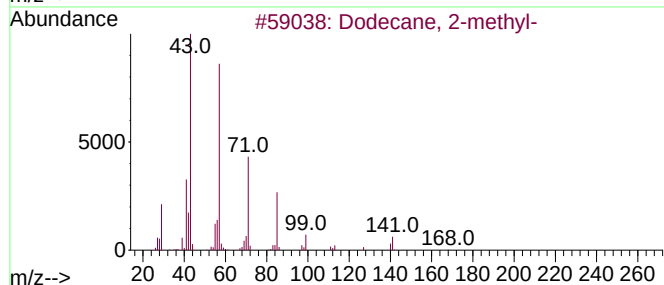
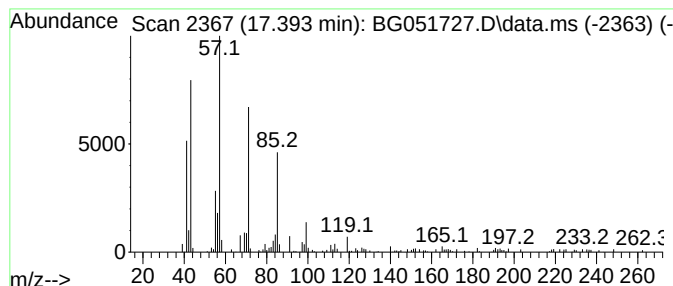
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TIC Integration Parameters: LSCINT.P

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Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 34

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 17.393 | 2.48 ng/ul | 114745 | Phenanthrene-d10 | 17.658 |

| Hit# | of | 5 | Tentative ID                       | MW  | MolForm  | CAS#         | Qual |
|------|----|---|------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Dodecane, 2-methyl-                | 184 | C13H28   | 001560-97-0  | 72   |
| 2    |    |   | Dodecane, 2-methyl-6-propyl-       | 226 | C16H34   | 055045-08-4  | 72   |
| 3    |    |   | 1,3-Propanediol, decyl ethyl ether | 244 | C15H32O2 | 1000406-35-0 | 72   |
| 4    |    |   | Undecane, 4,8-dimethyl-            | 184 | C13H28   | 017301-33-6  | 72   |
| 5    |    |   | Tridecane, 7-propyl-               | 226 | C16H34   | 055045-09-5  | 72   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122021\  
Data File : BG051727.D  
Acq On : 21 Dec 2021 22:20  
Operator : CG/JU  
Sample : M5121-04DL 10X  
Misc :  
ALS Vial : 54 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL53DL

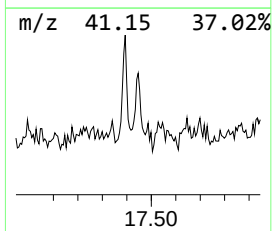
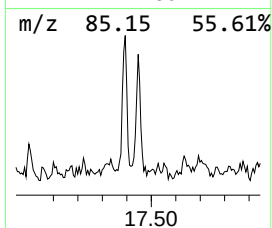
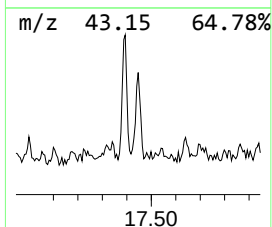
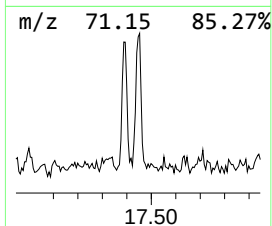
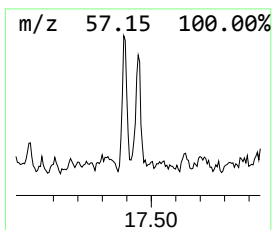
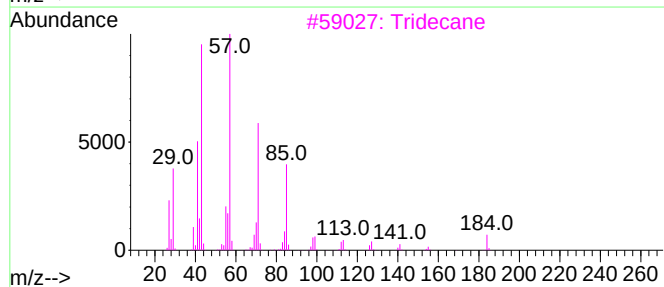
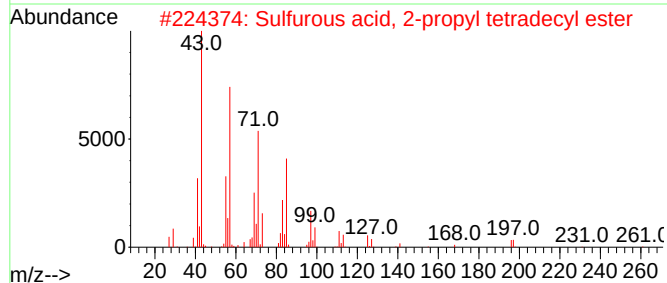
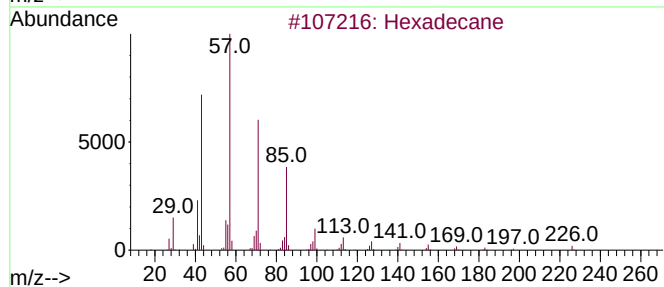
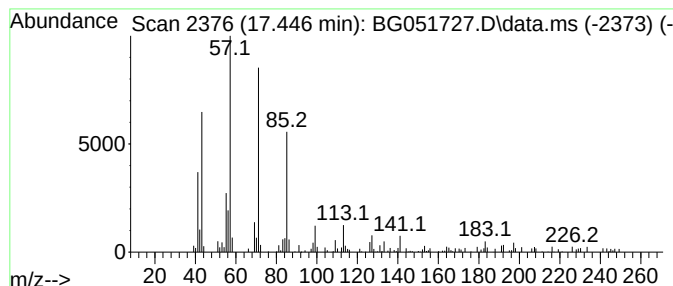
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Quant Title : SVOA CALIBRATION

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

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Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 39

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 17.446 | 2.30 ng/ul | 106375 | Phenanthrene-d10 | 17.658 |

| Hit# | of | 5 | Tentative ID                         | MW  | MolForm   | CAS#         | Qual |
|------|----|---|--------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Hexadecane                           | 226 | C16H34    | 000544-76-3  | 90   |
| 2    |    |   | Sulfurous acid, 2-propyl tetradec... | 320 | C17H36O3S | 1010309-12-5 | 90   |
| 3    |    |   | Tridecane                            | 184 | C13H28    | 000629-50-5  | 90   |
| 4    |    |   | Hexadecane                           | 226 | C16H34    | 000544-76-3  | 87   |
| 5    |    |   | Dodecane, 2-methyl-                  | 184 | C13H28    | 001560-97-0  | 87   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122021\  
Data File : BG051727.D  
Acq On : 21 Dec 2021 22:20  
Operator : CG/JU  
Sample : M5121-04DL 10X  
Misc :  
ALS Vial : 54 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL53DL

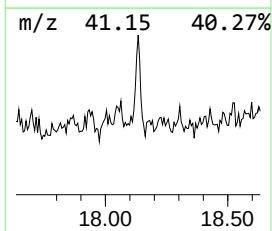
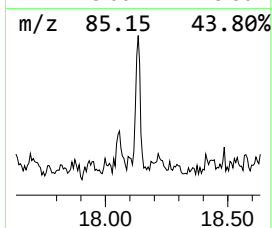
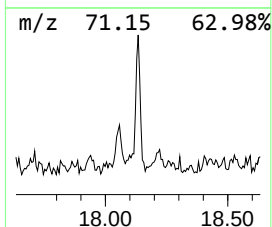
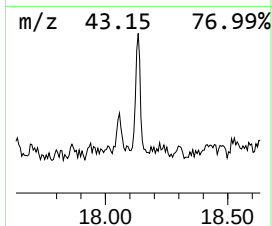
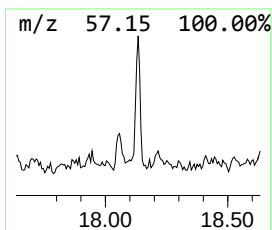
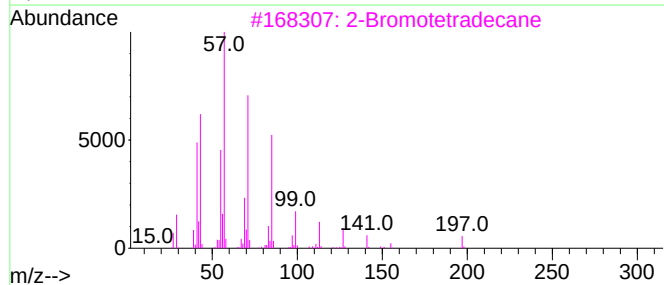
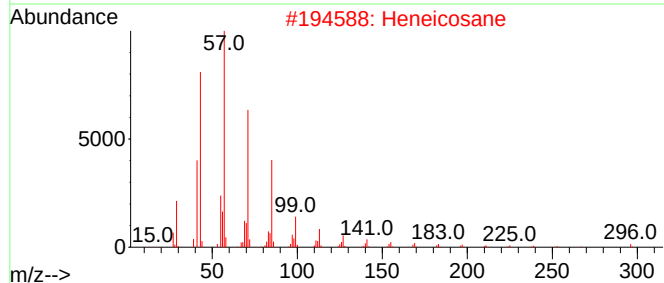
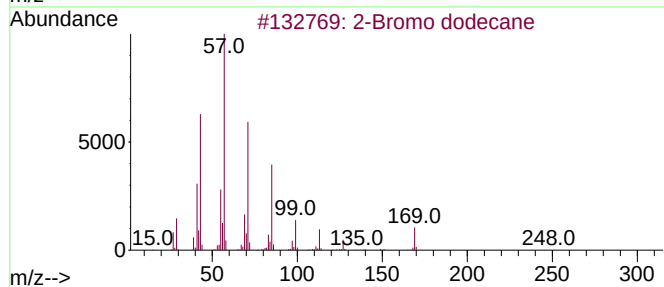
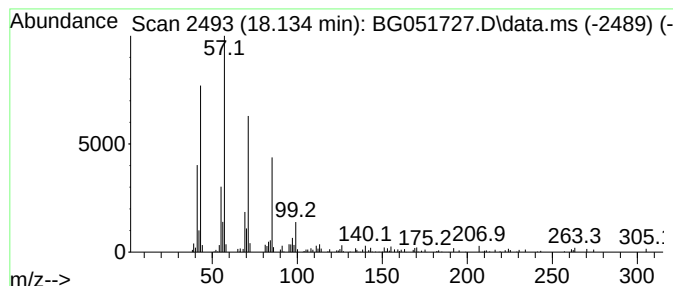
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Quant Title : SVOA CALIBRATION

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

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Peak Number 10 2-Bromo dodecane Concentration Rank 29

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 18.134 | 3.38 ng/ul | 156085 | Phenanthrene-d10 | 17.658 |

| Hit# | of 5 | Tentative ID        | MW  | MolForm  | CAS#        | Qual |
|------|------|---------------------|-----|----------|-------------|------|
| 1    |      | 2-Bromo dodecane    | 248 | C12H25Br | 013187-99-0 | 83   |
| 2    |      | Heneicosane         | 296 | C21H44   | 000629-94-7 | 80   |
| 3    |      | 2-Bromotetradecane  | 276 | C14H29Br | 074036-95-6 | 80   |
| 4    |      | Heptadecane         | 240 | C17H36   | 000629-78-7 | 80   |
| 5    |      | Octadecane, 1-iodo- | 380 | C18H37I  | 000629-93-6 | 80   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122021\  
Data File : BG051727.D  
Acq On : 21 Dec 2021 22:20  
Operator : CG/JU  
Sample : M5121-04DL 10X  
Misc :  
ALS Vial : 54 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL53DL

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

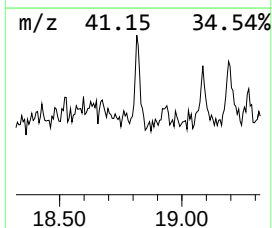
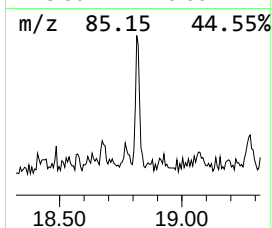
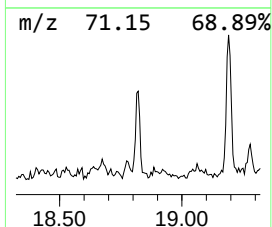
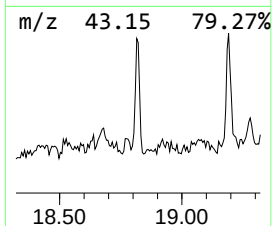
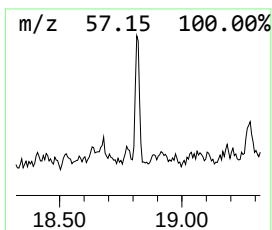
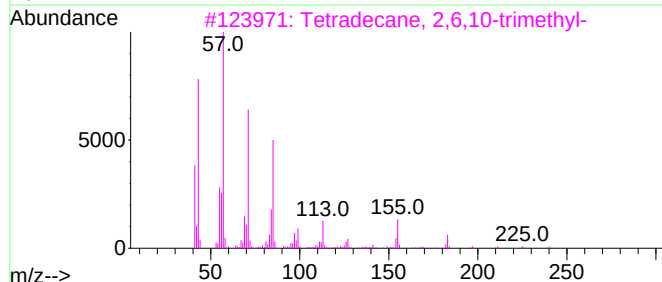
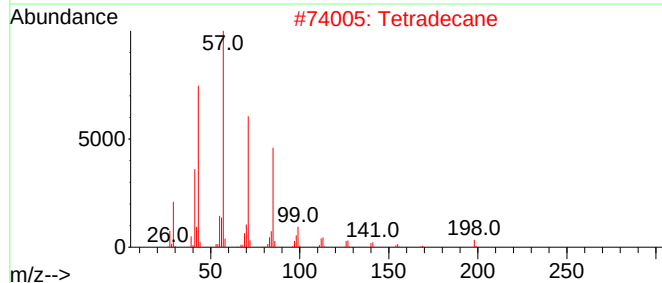
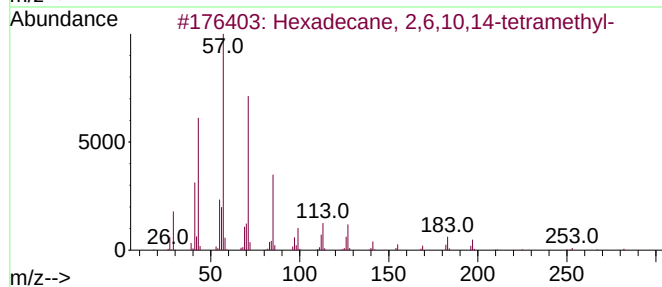
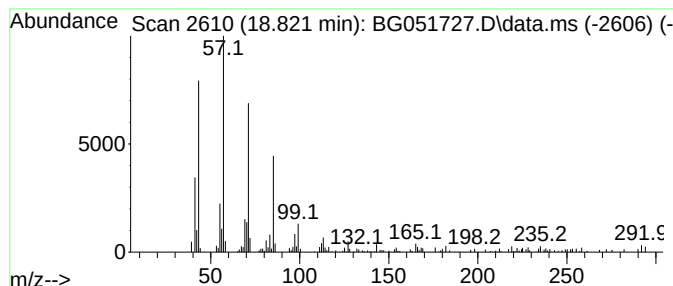
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TIC Integration Parameters: LSCINT.P

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Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 36

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 18.821 | 2.35 ng/ul | 108790 | Phenanthrene-d10 | 17.658 |

| Hit# | of | 5 | Tentative ID                       | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Hexadecane, 2,6,10,14-tetramethyl- | 282 | C20H42  | 000638-36-8 | 90   |
| 2    |    |   | Tetradecane                        | 198 | C14H30  | 000629-59-4 | 87   |
| 3    |    |   | Tetradecane, 2,6,10-trimethyl-     | 240 | C17H36  | 014905-56-7 | 87   |
| 4    |    |   | Tetradecane, 1-iodo-               | 324 | C14H29I | 019218-94-1 | 83   |
| 5    |    |   | Tridecane                          | 184 | C13H28  | 000629-50-5 | 83   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122021\  
Data File : BG051727.D  
Acq On : 21 Dec 2021 22:20  
Operator : CG/JU  
Sample : M5121-04DL 10X  
Misc :  
ALS Vial : 54 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL53DL

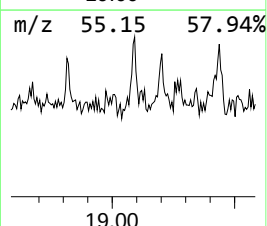
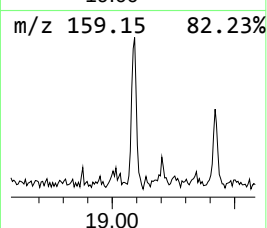
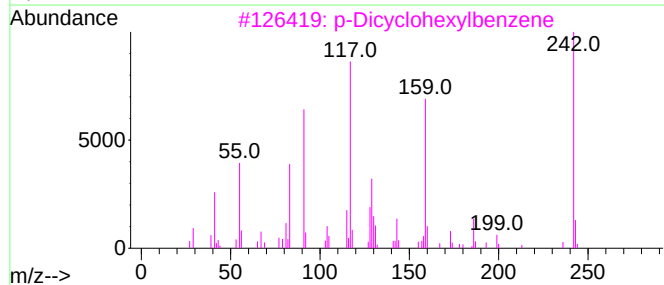
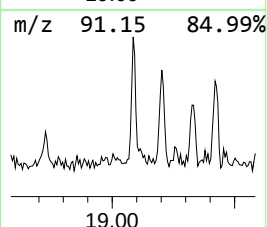
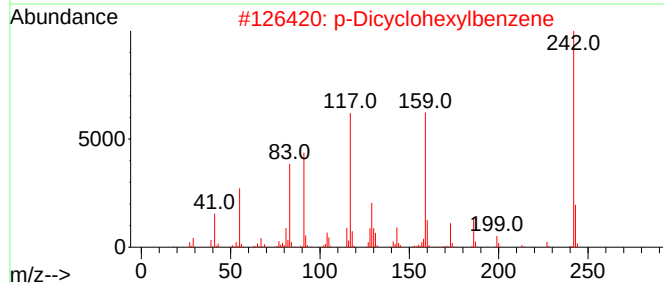
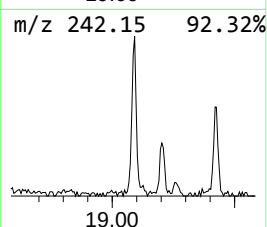
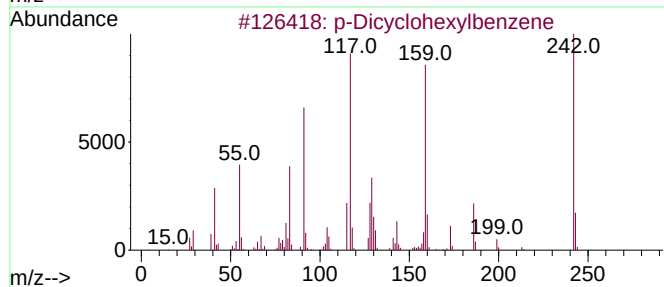
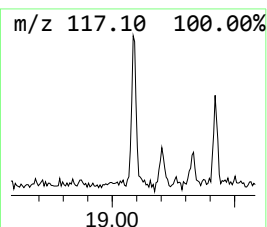
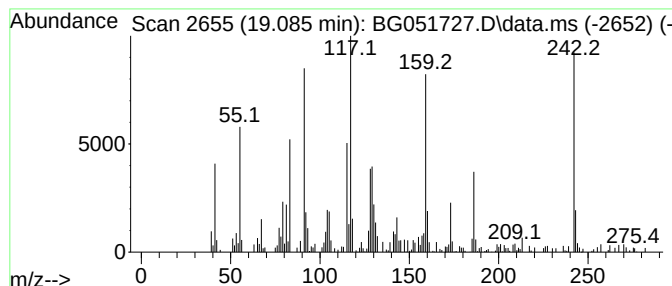
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Quant Title : SVOA CALIBRATION

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

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Peak Number 12 p-Dicyclohexylbenzene Concentration Rank 26

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 19.085 | 3.50 ng/ul | 161729 | Phenanthrene-d10 | 17.658 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | p-Dicyclohexylbenzene               | 242 | C18H26   | 001087-02-1  | 93   |
| 2    |    |   | p-Dicyclohexylbenzene               | 242 | C18H26   | 001087-02-1  | 91   |
| 3    |    |   | p-Dicyclohexylbenzene               | 242 | C18H26   | 001087-02-1  | 91   |
| 4    |    |   | Bicyclohexyl, 4-phenyl-             | 242 | C18H26   | 020273-27-2  | 76   |
| 5    |    |   | 2-Phenylpropionic acid, 4'-(2-me... | 218 | C14H18O2 | 1010454-94-3 | 70   |





Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122021\  
Data File : BG051727.D  
Acq On : 21 Dec 2021 22:20  
Operator : CG/JU  
Sample : M5121-04DL 10X  
Misc :  
ALS Vial : 54 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL53DL

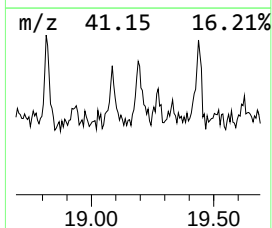
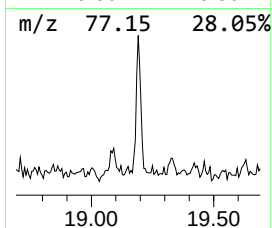
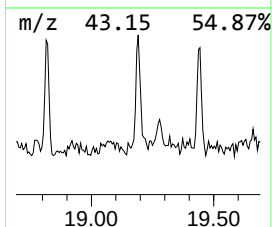
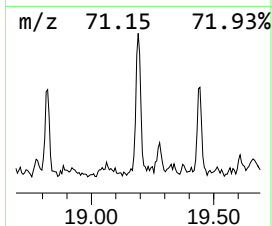
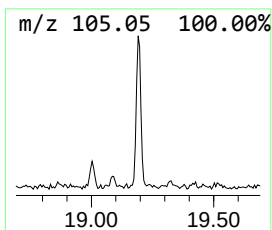
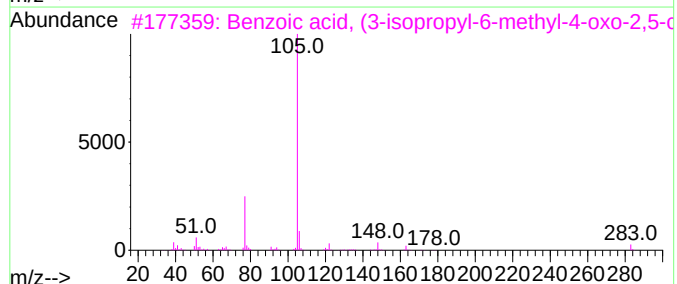
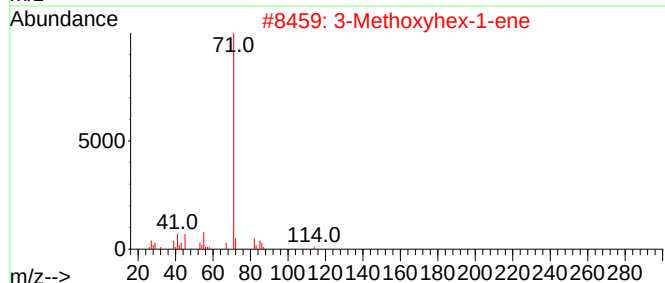
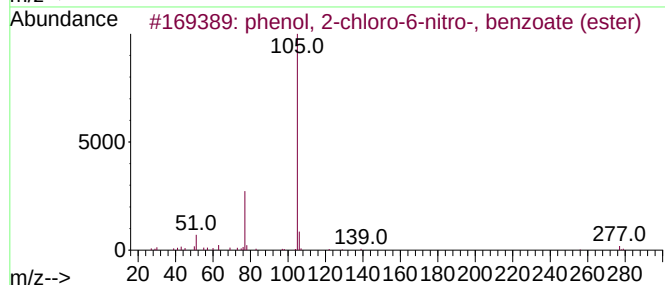
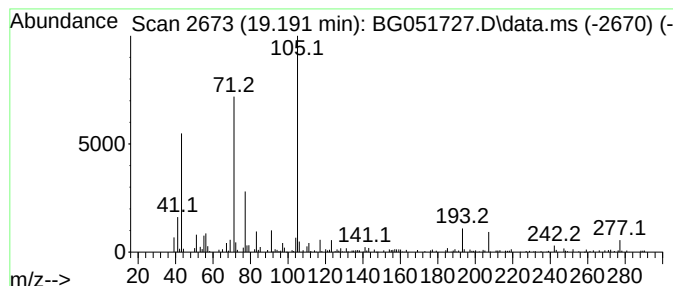
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Quant Title : SVOA CALIBRATION

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

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Peak Number 13 unknown-02 Concentration Rank 25

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 19.191 | 3.81 ng/ul | 175884 | Phenanthrene-d10 | 17.658 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|------------|--------------|------|
| 1    |    |   | phenol, 2-chloro-6-nitro-, benzo... | 277 | C13H8ClNO4 | 1010397-11-4 | 27   |
| 2    |    |   | 3-Methoxyhex-1-ene                  | 114 | C7H14O     | 108811-41-2  | 22   |
| 3    |    |   | Benzoic acid, (3-isopropyl-6-met... | 283 | C17H17NO3  | 039870-46-7  | 22   |
| 4    |    |   | Glutaric acid, propyl tetrahydro... | 258 | C13H22O5   | 1000359-65-7 | 22   |
| 5    |    |   | Cyclohexanol, 2,2-dichloro-1-met... | 182 | C7H12Cl2O  | 1000115-65-2 | 22   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122021\  
Data File : BG051727.D  
Acq On : 21 Dec 2021 22:20  
Operator : CG/JU  
Sample : M5121-04DL 10X  
Misc :  
ALS Vial : 54 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL53DL

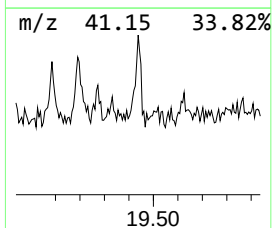
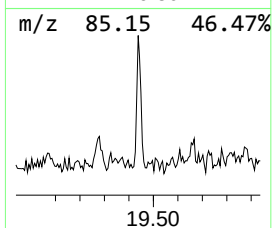
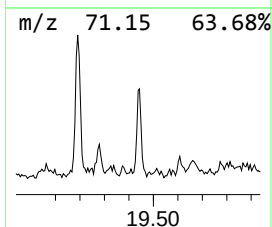
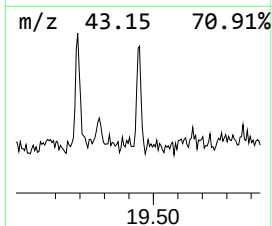
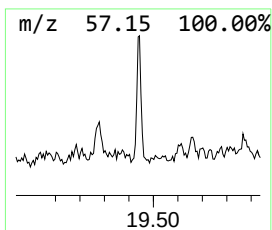
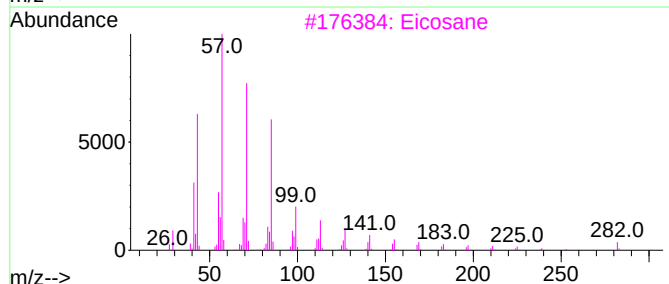
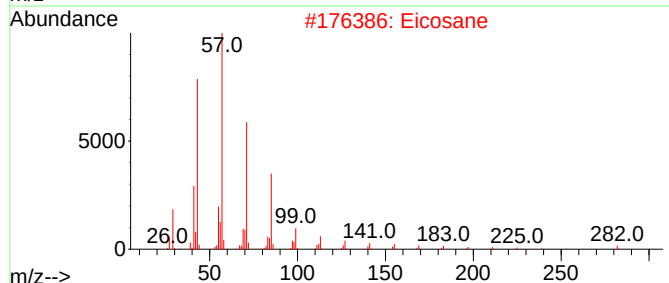
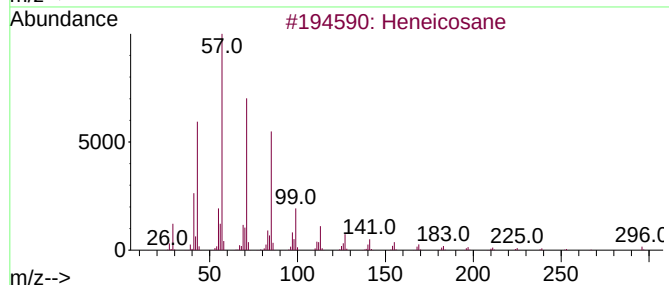
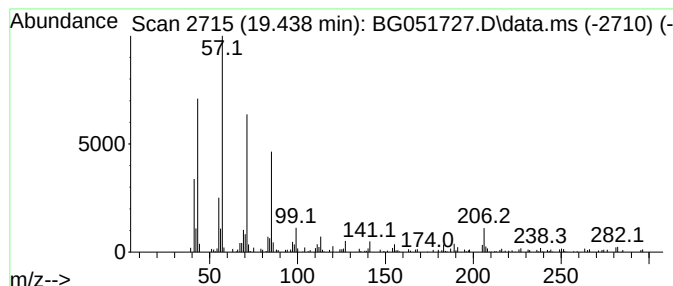
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Quant Title : SVOA CALIBRATION

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

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Peak Number 14 (DEL) Alkane: Straight-Chai... Concentration Rank 27

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 19.438 | 3.44 ng/ul | 159109 | Phenanthrene-d10 | 17.658 |

| Hit# | of 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|------|--------------|-----|---------|-------------|------|
| 1    |      | Heneicosane  | 296 | C21H44  | 000629-94-7 | 97   |
| 2    |      | Eicosane     | 282 | C20H42  | 000112-95-8 | 96   |
| 3    |      | Eicosane     | 282 | C20H42  | 000112-95-8 | 94   |
| 4    |      | Heneicosane  | 296 | C21H44  | 000629-94-7 | 92   |
| 5    |      | Heptacosane  | 380 | C27H56  | 000593-49-7 | 81   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122021\  
Data File : BG051727.D  
Acq On : 21 Dec 2021 22:20  
Operator : CG/JU  
Sample : M5121-04DL 10X  
Misc :  
ALS Vial : 54 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL53DL

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

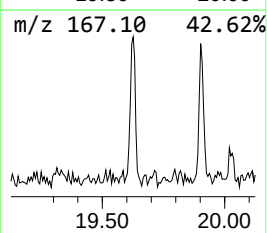
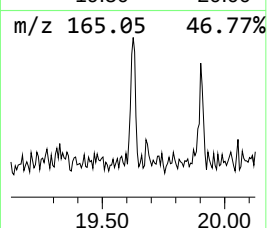
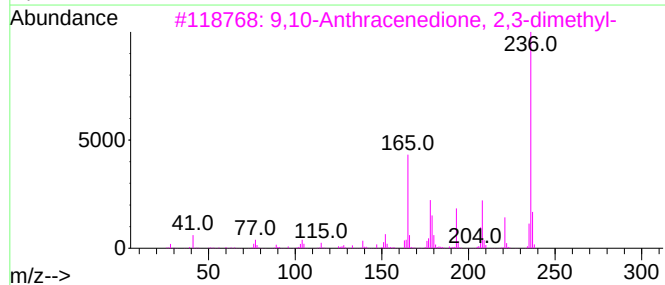
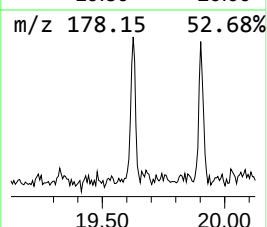
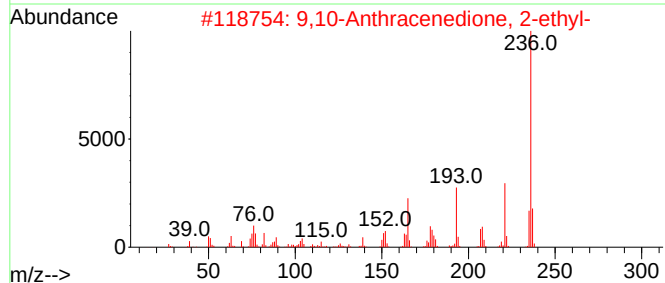
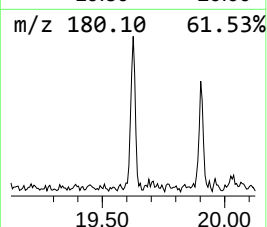
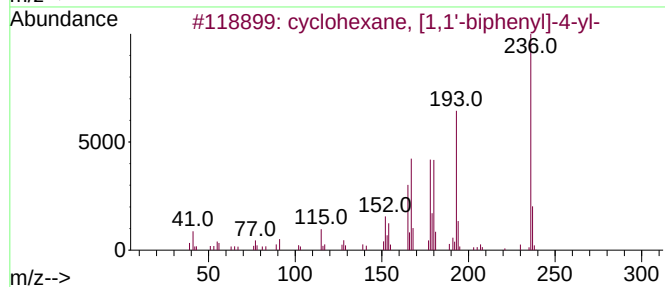
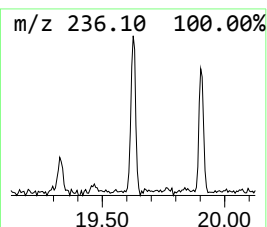
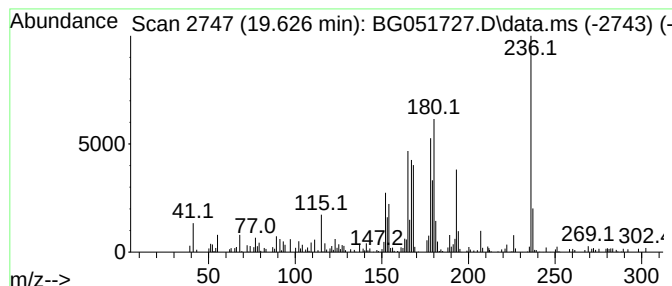
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 15 cyclohexane, [1,1'-biphenyl]... Concentration Rank 35

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 19.626 | 2.43 ng/ul | 112262 | Phenanthrene-d10 | 17.658 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | cyclohexane, [1,1'-biphenyl]-4-yl-  | 236 | C18H20   | 1000401-12-4 | 83   |
| 2    |    |   | 9,10-Anthracenedione, 2-ethyl-      | 236 | C16H12O2 | 000084-51-5  | 43   |
| 3    |    |   | 9,10-Anthracenedione, 2,3-dimethyl- | 236 | C16H12O2 | 006531-35-7  | 43   |
| 4    |    |   | 9,10-Anthracenedione, 2,7-dimethyl- | 236 | C16H12O2 | 003286-01-9  | 38   |
| 5    |    |   | 2,4-Diethoxy-5-methoxy-1-(2-prop... | 236 | C14H20O3 | 1000122-72-0 | 38   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122021\  
Data File : BG051727.D  
Acq On : 21 Dec 2021 22:20  
Operator : CG/JU  
Sample : M5121-04DL 10X  
Misc :  
ALS Vial : 54 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL53DL

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

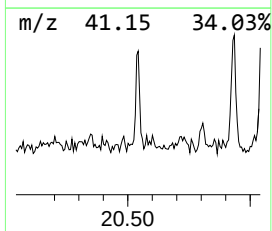
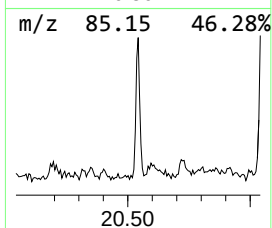
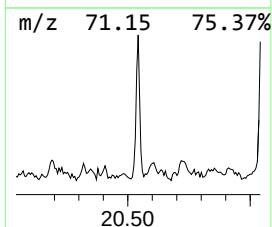
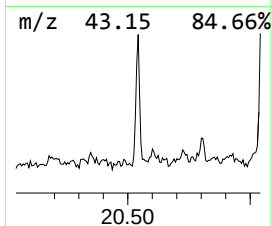
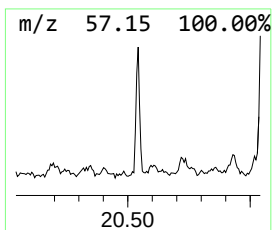
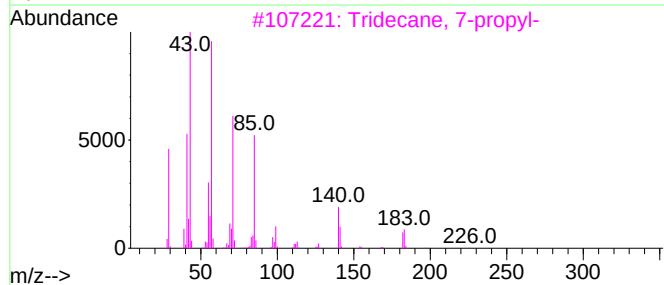
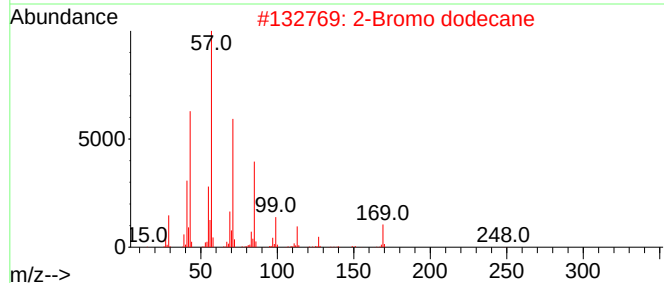
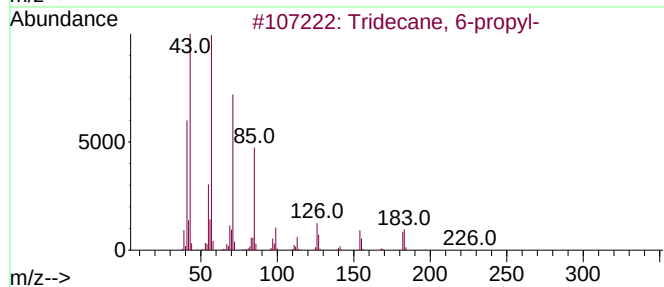
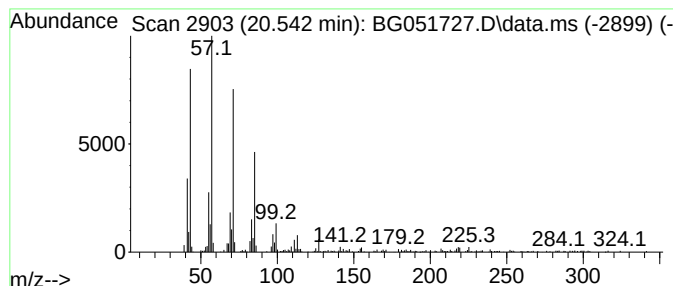
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 16 (DEL) Alkane: Straight-Chai... Concentration Rank 23

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 20.542 | 4.81 ng/ul | 230084 | Chrysene-d12     | 21.981 |

| Hit# | of 5 | Tentative ID         | MW  | MolForm  | CAS#        | Qual |
|------|------|----------------------|-----|----------|-------------|------|
| 1    |      | Tridecane, 6-propyl- | 226 | C16H34   | 055045-10-8 | 94   |
| 2    |      | 2-Bromo dodecane     | 248 | C12H25Br | 013187-99-0 | 93   |
| 3    |      | Tridecane, 7-propyl- | 226 | C16H34   | 055045-09-5 | 91   |
| 4    |      | Hexatriacontane      | 507 | C36H74   | 000630-06-8 | 91   |
| 5    |      | Heptacosane          | 380 | C27H56   | 000593-49-7 | 91   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122021\  
Data File : BG051727.D  
Acq On : 21 Dec 2021 22:20  
Operator : CG/JU  
Sample : M5121-04DL 10X  
Misc :  
ALS Vial : 54 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL53DL

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

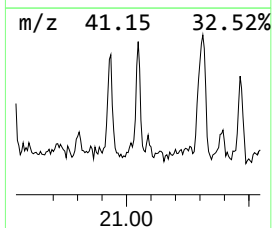
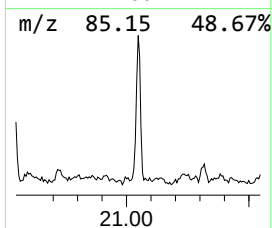
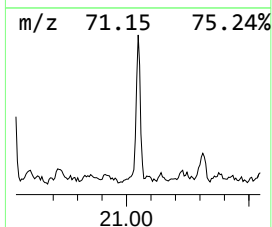
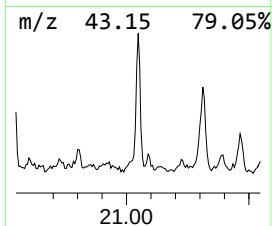
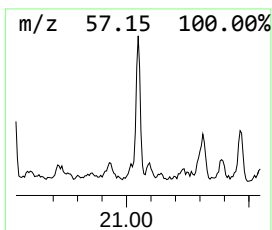
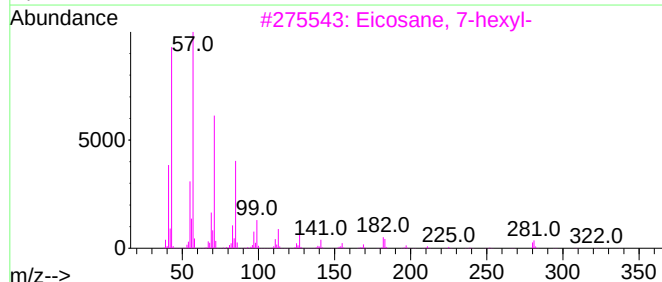
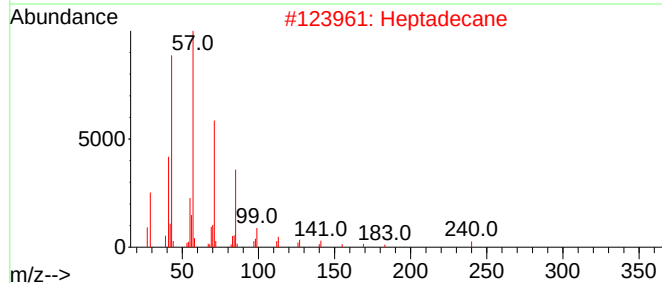
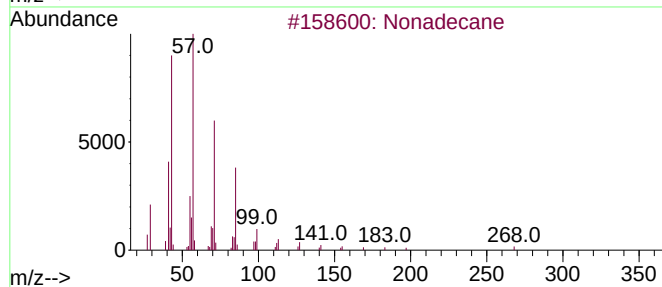
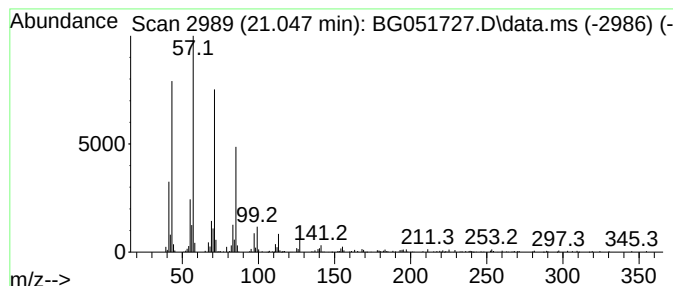
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 17 (DEL) Alkane: Straight-Chai... Concentration Rank 19

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 21.047 | 6.10 ng/ul | 291689 | Chrysene-d12     | 21.981 |

| Hit# | of | 5 | Tentative ID         | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------|-----|---------|-------------|------|
| 1    |    |   | Nonadecane           | 268 | C19H40  | 000629-92-5 | 93   |
| 2    |    |   | Heptadecane          | 240 | C17H36  | 000629-78-7 | 93   |
| 3    |    |   | Eicosane, 7-hexyl-   | 366 | C26H54  | 055333-99-8 | 90   |
| 4    |    |   | Decane, 5-propyl-    | 184 | C13H28  | 017312-62-8 | 87   |
| 5    |    |   | Tetradecane, 1-iodo- | 324 | C14H29I | 019218-94-1 | 86   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122021\  
Data File : BG051727.D  
Acq On : 21 Dec 2021 22:20  
Operator : CG/JU  
Sample : M5121-04DL 10X  
Misc :  
ALS Vial : 54 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL53DL

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

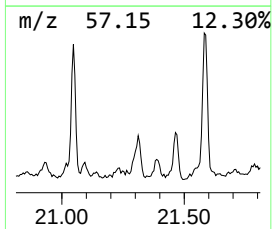
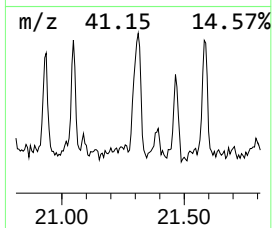
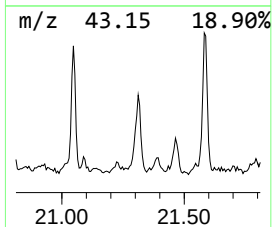
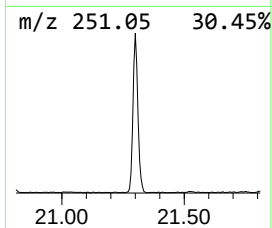
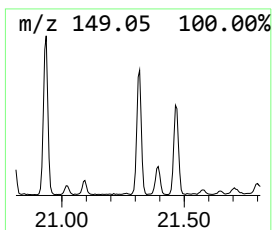
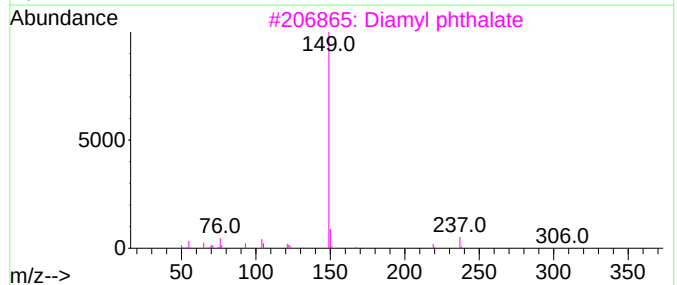
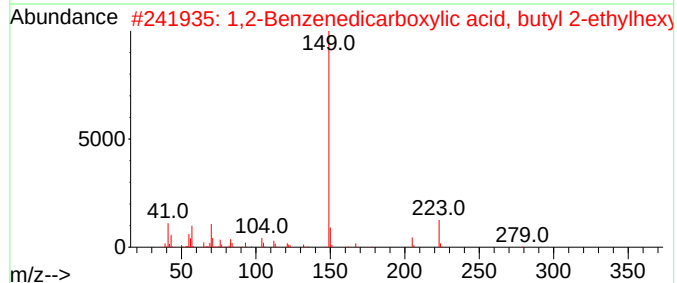
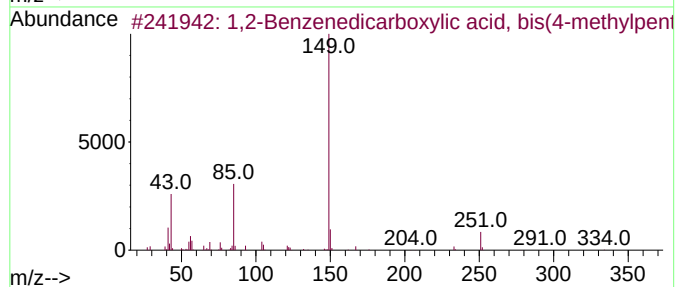
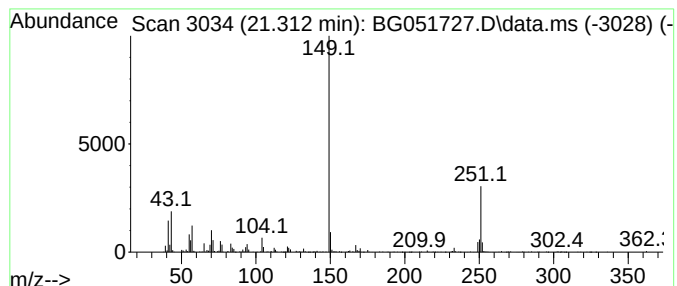
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TIC Integration Parameters: LSCINT.P

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Peak Number 18 1,2-Benzenedicarboxylic aci... Concentration Rank 5

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 21.312 | 16.23 ng/ul | 776592 | Chrysene-d12     | 21.981 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|------|-------------------------------------|-----|----------|--------------|------|
| 1    |      | 1,2-Benzenedicarboxylic acid, bi... | 334 | C20H3004 | 000146-50-9  | 64   |
| 2    |      | 1,2-Benzenedicarboxylic acid, bu... | 334 | C20H3004 | 000085-69-8  | 58   |
| 3    |      | Diamyl phthalate                    | 306 | C18H2604 | 000131-18-0  | 58   |
| 4    |      | Phthalic acid, 2-ethylhexyl hexy... | 362 | C22H3404 | 1000309-02-5 | 58   |
| 5    |      | Phthalic acid, 2-ethylhexyl prop... | 320 | C19H2804 | 1000309-00-2 | 58   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122021\  
Data File : BG051727.D  
Acq On : 21 Dec 2021 22:20  
Operator : CG/JU  
Sample : M5121-04DL 10X  
Misc :  
ALS Vial : 54 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL53DL

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

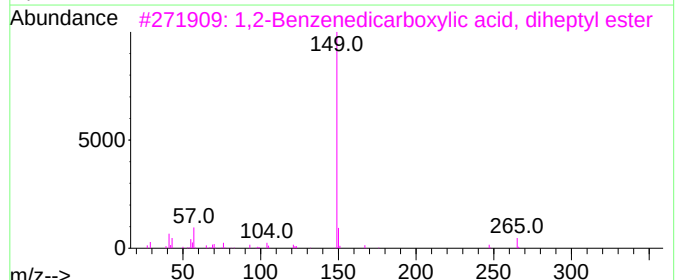
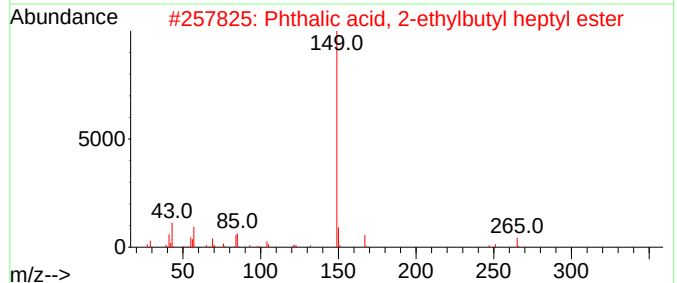
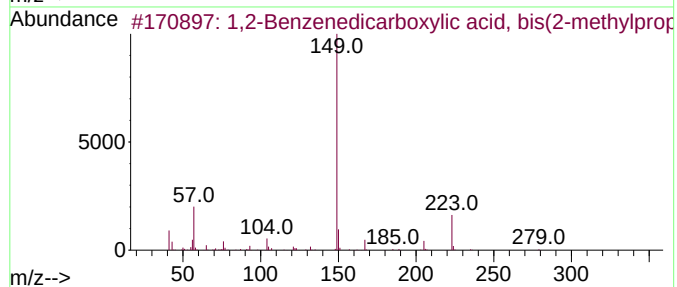
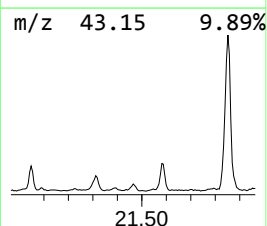
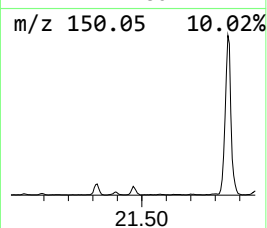
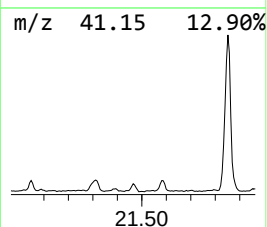
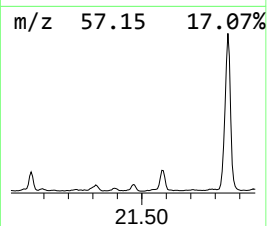
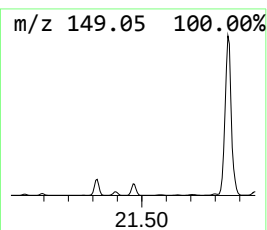
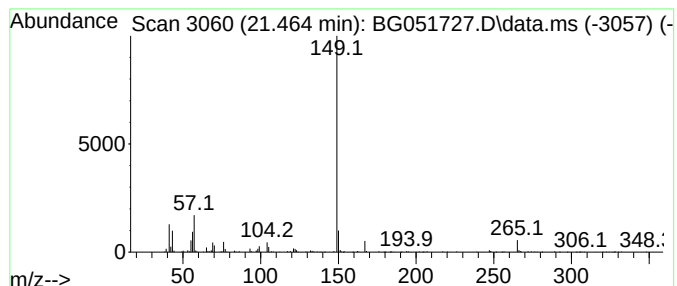
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 19 Phthalic acid, 2-ethylbutyl... Concentration Rank 22

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 21.464 | 5.40 ng/ul | 258369 | Chrysene-d12     | 21.981 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | 1,2-Benzenedicarboxylic acid, bi... | 278 | C16H22O4 | 000084-69-5  | 87   |
| 2    |    |   | Phthalic acid, 2-ethylbutyl hept... | 348 | C21H32O4 | 1000356-89-2 | 86   |
| 3    |    |   | 1,2-Benzenedicarboxylic acid, di... | 362 | C22H34O4 | 003648-21-3  | 80   |
| 4    |    |   | Diamyl phthalate                    | 306 | C18H26O4 | 000131-18-0  | 76   |
| 5    |    |   | Phthalic acid, butyl hexyl ester    | 306 | C18H26O4 | 1010308-99-5 | 72   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122021\  
Data File : BG051727.D  
Acq On : 21 Dec 2021 22:20  
Operator : CG/JU  
Sample : M5121-04DL 10X  
Misc :  
ALS Vial : 54 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL53DL

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

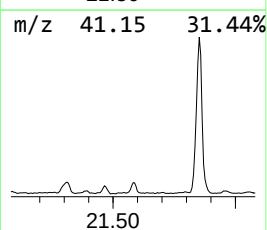
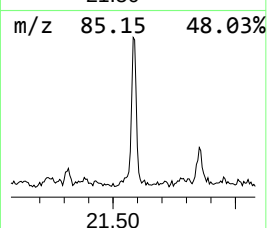
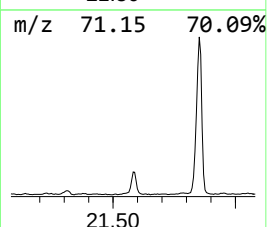
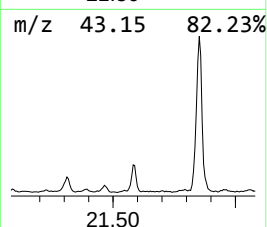
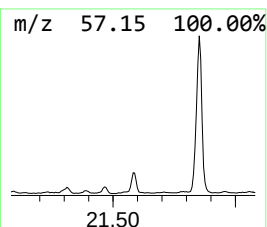
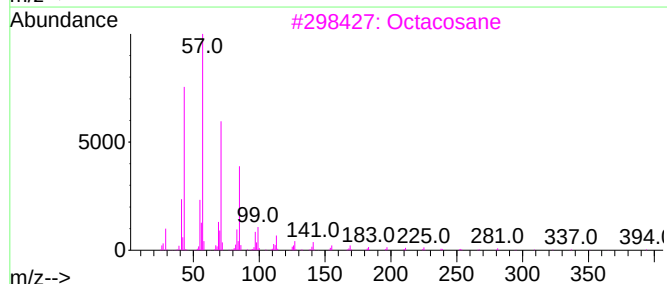
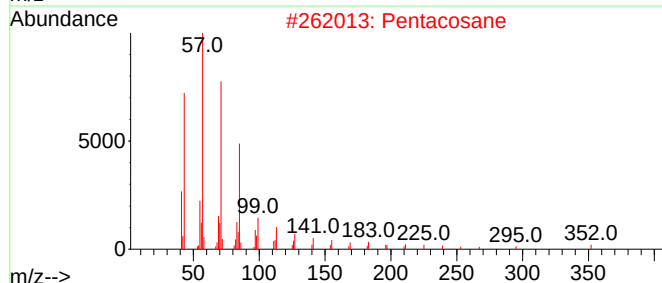
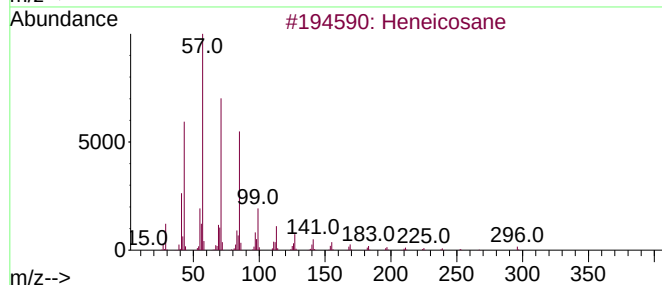
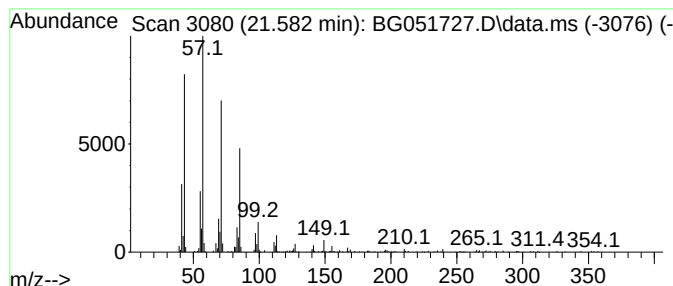
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TIC Integration Parameters: LSCINT.P

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Peak Number 20 (DEL) Alkane: Straight-Chai... Concentration Rank 15

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 21.582 | 8.91 ng/ul | 426324 | Chrysene-d12     | 21.981 |

| Hit# | of 5 | Tentative ID         | MW  | MolForm | CAS#        | Qual |
|------|------|----------------------|-----|---------|-------------|------|
| 1    |      | Heneicosane          | 296 | C21H44  | 000629-94-7 | 96   |
| 2    |      | Pentacosane          | 352 | C25H52  | 000629-99-2 | 93   |
| 3    |      | Octacosane           | 394 | C28H58  | 000630-02-4 | 87   |
| 4    |      | Pentadecane          | 212 | C15H32  | 000629-62-9 | 87   |
| 5    |      | Tetradecane, 1-iodo- | 324 | C14H29I | 019218-94-1 | 86   |





Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122021\  
Data File : BG051727.D  
Acq On : 21 Dec 2021 22:20  
Operator : CG/JU  
Sample : M5121-04DL 10X  
Misc :  
ALS Vial : 54 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL53DL

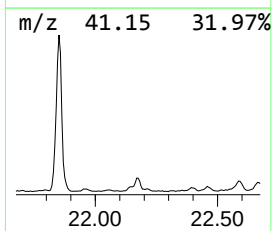
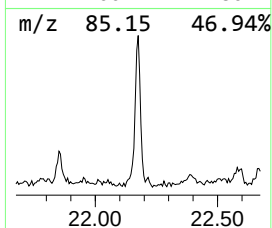
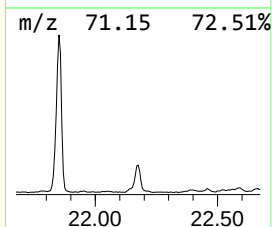
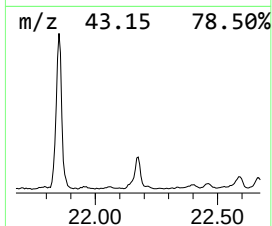
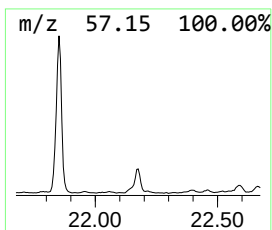
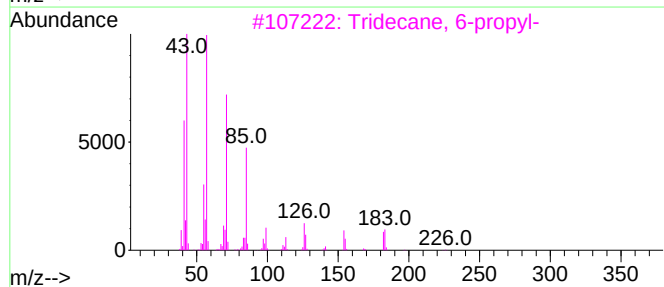
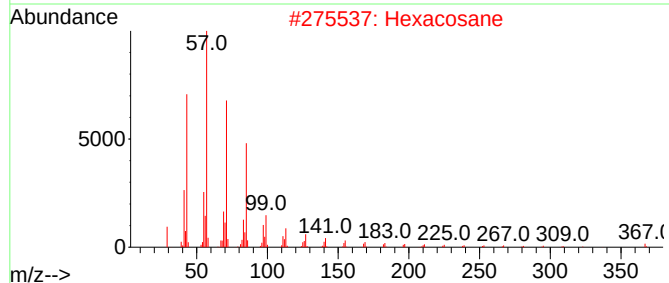
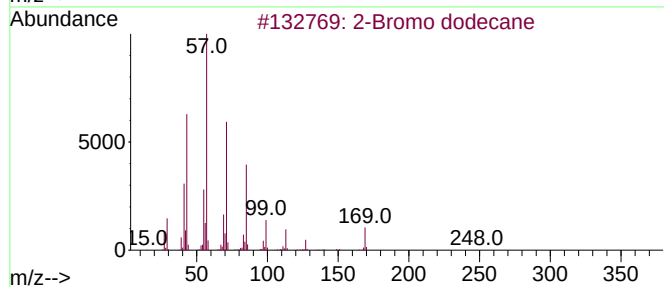
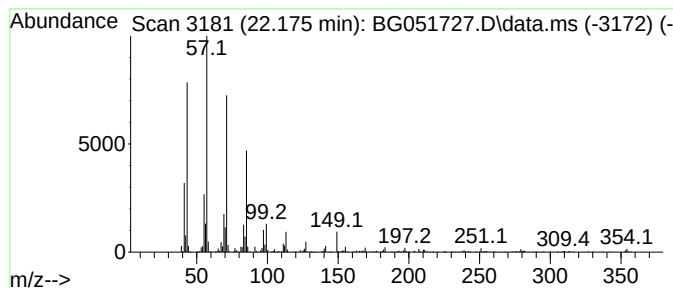
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Quant Title : SVOA CALIBRATION

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

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Peak Number 21 Hexacosane Concentration Rank 4

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 22.175 | 16.73 ng/ul | 800450 | Chrysene-d12     | 21.981 |

| Hit# | of 5 | Tentative ID         | MW  | MolForm  | CAS#        | Qual |
|------|------|----------------------|-----|----------|-------------|------|
| 1    |      | 2-Bromo dodecane     | 248 | C12H25Br | 013187-99-0 | 91   |
| 2    |      | Hexacosane           | 366 | C26H54   | 000630-01-3 | 87   |
| 3    |      | Tridecane, 6-propyl- | 226 | C16H34   | 055045-10-8 | 87   |
| 4    |      | Heptacosane          | 380 | C27H56   | 000593-49-7 | 87   |
| 5    |      | Octacosane           | 394 | C28H58   | 000630-02-4 | 87   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122021\  
Data File : BG051727.D  
Acq On : 21 Dec 2021 22:20  
Operator : CG/JU  
Sample : M5121-04DL 10X  
Misc :  
ALS Vial : 54 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL53DL

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

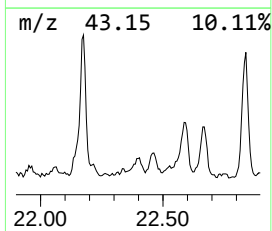
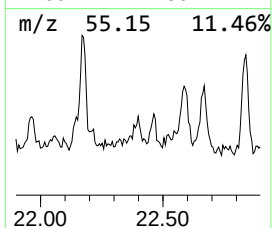
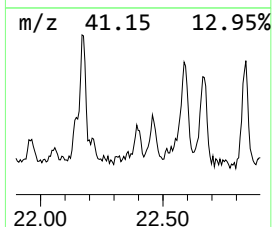
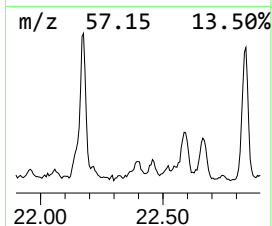
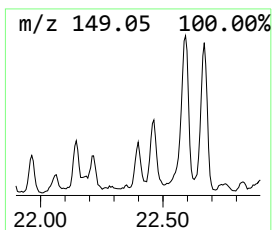
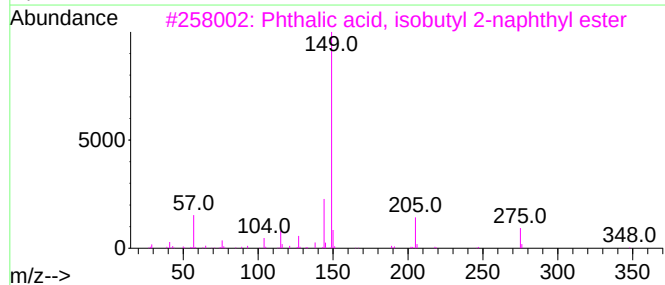
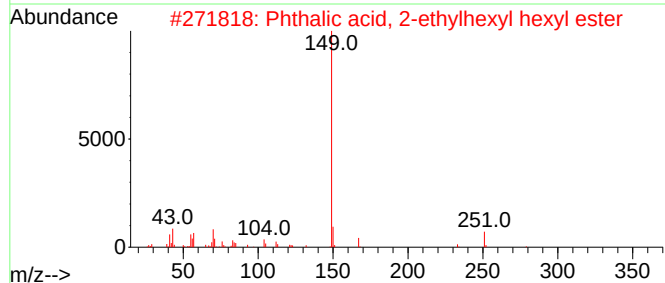
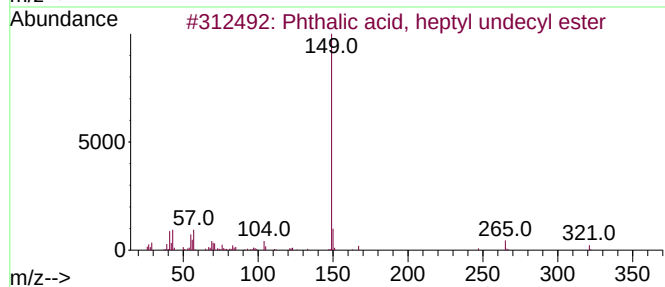
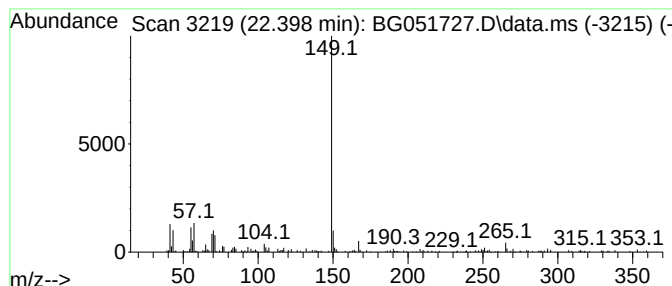
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 22 Phthalic acid, heptyl undec... Concentration Rank 31

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 22.398 | 2.97 ng/ul | 142123 | Chrysene-d12     | 21.981 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Phthalic acid, heptyl undecyl ester | 418 | C26H42O4 | 1010308-94-0 | 80   |
| 2    |    |   | Phthalic acid, 2-ethylhexyl hexy... | 362 | C22H34O4 | 1000309-02-5 | 72   |
| 3    |    |   | Phthalic acid, isobutyl 2-naphth... | 348 | C22H20O4 | 1000315-24-4 | 72   |
| 4    |    |   | 1,2-Benzenedicarboxylic acid, bu... | 334 | C20H30O4 | 000084-78-6  | 72   |
| 5    |    |   | Phthalic acid, di(oct-3-yl) ester   | 390 | C24H38O4 | 1000377-72-3 | 64   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122021\  
Data File : BG051727.D  
Acq On : 21 Dec 2021 22:20  
Operator : CG/JU  
Sample : M5121-04DL 10X  
Misc :  
ALS Vial : 54 Sample Multiplier: 1

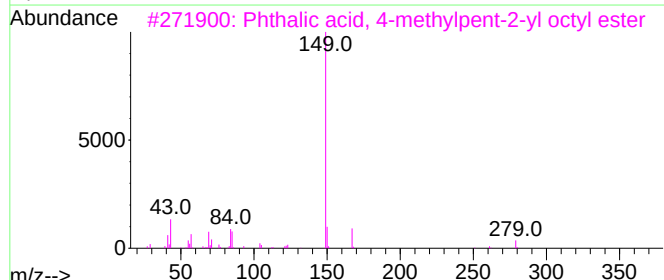
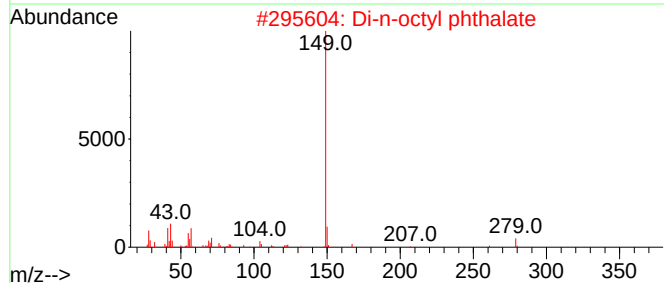
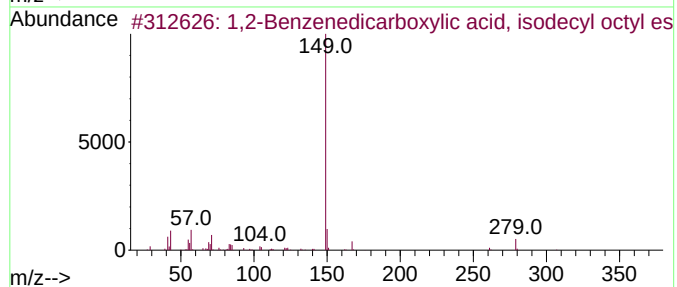
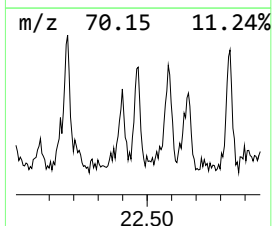
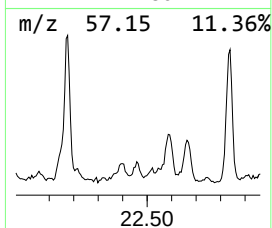
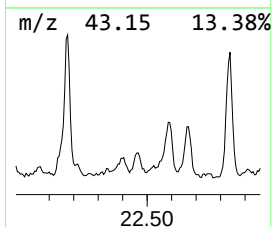
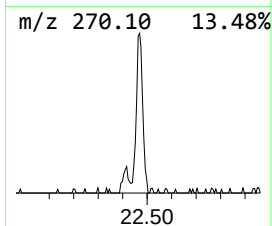
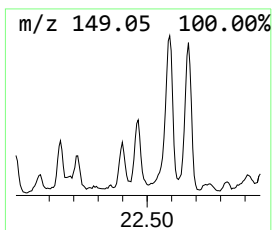
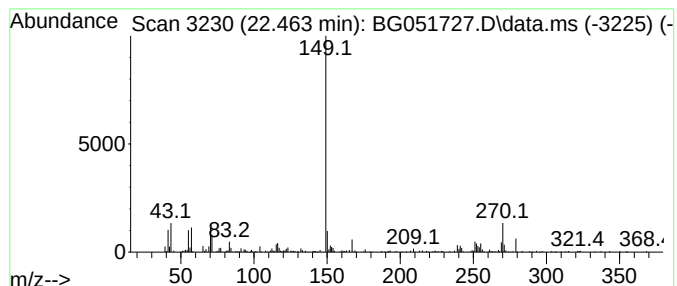
Instrument :  
BNA\_G  
ClientSampleId :  
BGL53DL

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 23 1,2-Benzenedicarboxylic aci... Concentration Rank 20

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|--------|------------------|--------------|------|
| 22.463    | 5.87 ng/ul                          | 280619 | Chrysene-d12     | 21.981       |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#         | Qual |
| 1         | 1,2-Benzenedicarboxylic acid, is... | 418    | C26H42O4         | 001330-96-7  | 64   |
| 2         | Di-n-octyl phthalate                | 390    | C24H38O4         | 000117-84-0  | 64   |
| 3         | Phthalic acid, 4-methylpent-2-yl... | 362    | C22H34O4         | 1000356-87-8 | 59   |
| 4         | Benzaldehyde, 3-(4-ethylphenoxy...  | 270    | C17H18O3         | 1000273-80-8 | 59   |
| 5         | m-Tolyl isothiocyanate              | 149    | C8H7NS           | 000621-30-7  | 58   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122021\  
Data File : BG051727.D  
Acq On : 21 Dec 2021 22:20  
Operator : CG/JU  
Sample : M5121-04DL 10X  
Misc :  
ALS Vial : 54 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL53DL

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

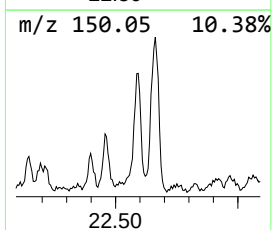
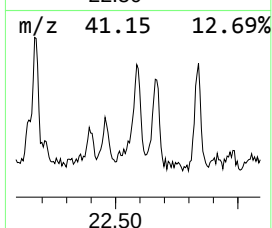
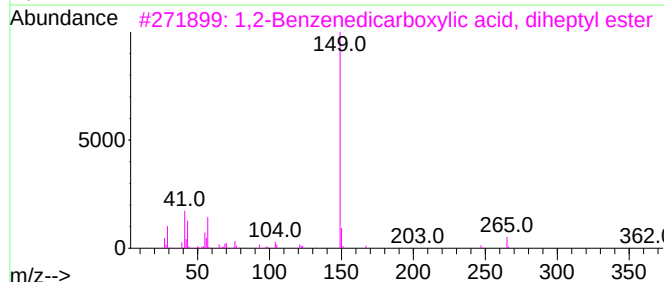
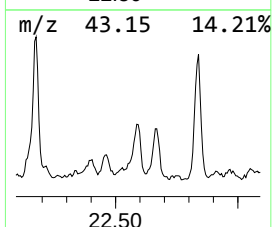
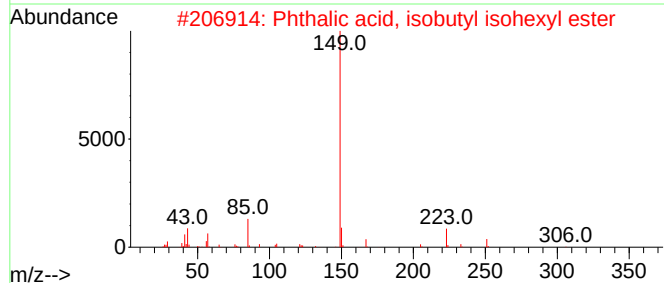
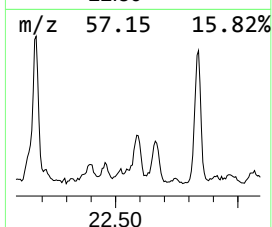
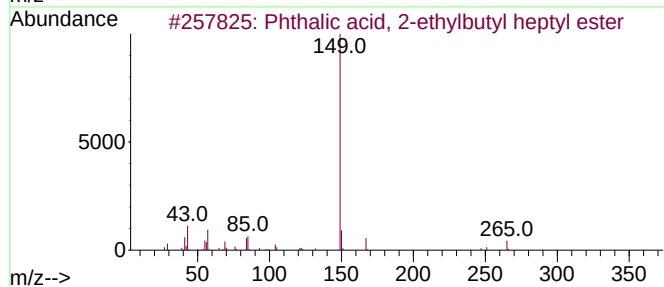
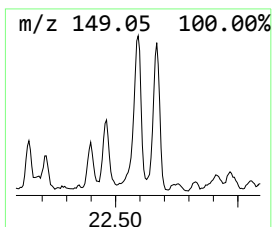
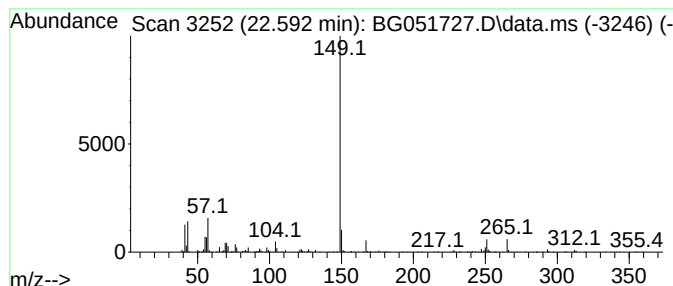
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 24 Phthalic acid, isobutyl iso... Concentration Rank 11

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 22.592 | 11.37 ng/ul | 543758 | Chrysene-d12     | 21.981 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Phthalic acid, 2-ethylbutyl hept... | 348 | C21H32O4 | 1000356-89-2 | 86   |
| 2    |    |   | Phthalic acid, isobutyl isohexyl... | 306 | C18H26O4 | 1000308-98-1 | 80   |
| 3    |    |   | 1,2-Benzenedicarboxylic acid, di... | 362 | C22H34O4 | 003648-21-3  | 80   |
| 4    |    |   | Phthalic acid, hexyl nonyl ester    | 376 | C23H36O4 | 1000309-01-1 | 80   |
| 5    |    |   | Phthalic acid, heptyl undecyl ester | 418 | C26H42O4 | 1010308-94-0 | 80   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122021\  
Data File : BG051727.D  
Acq On : 21 Dec 2021 22:20  
Operator : CG/JU  
Sample : M5121-04DL 10X  
Misc :  
ALS Vial : 54 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL53DL

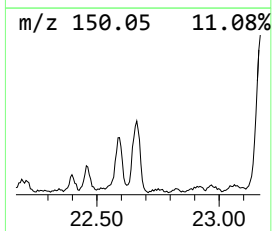
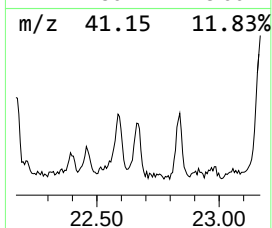
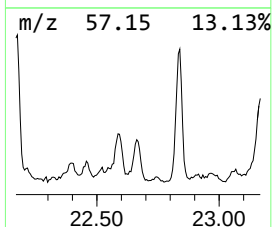
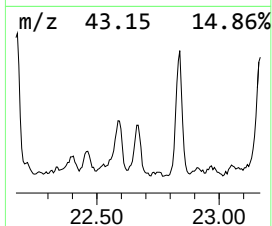
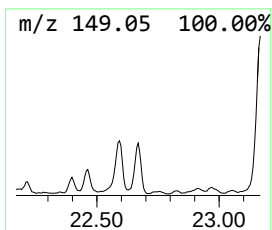
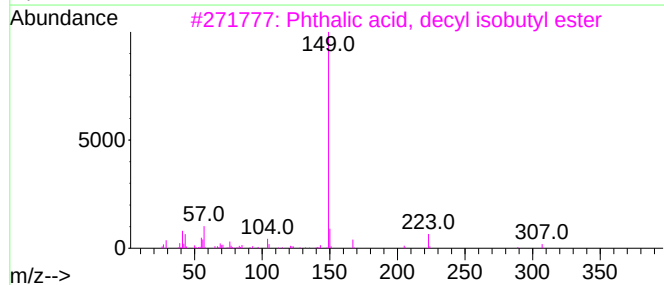
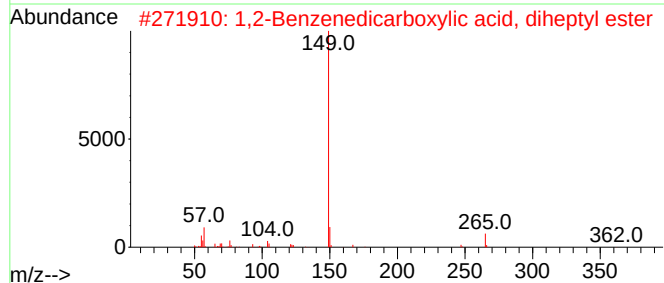
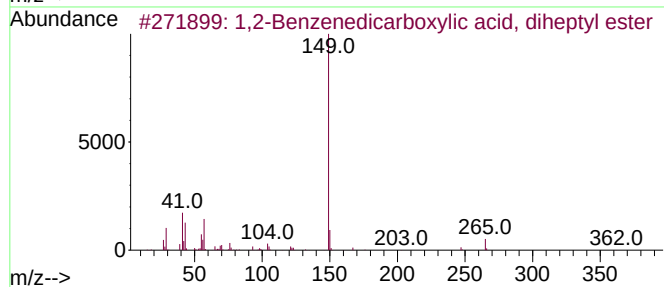
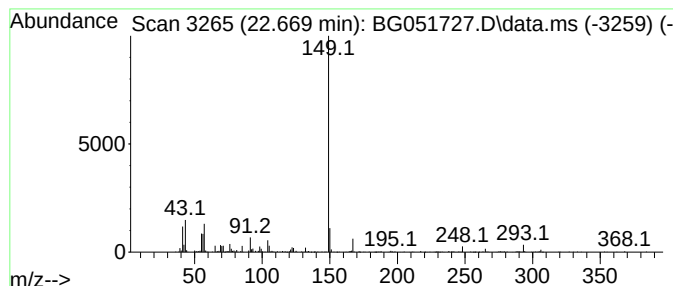
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Quant Title : SVOA CALIBRATION

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

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Peak Number 25 1,2-Benzenedicarboxylic aci... Concentration Rank 8

| R.T.   | EstConc     | Area                                | Relative to ISTD | R.T.     |              |      |
|--------|-------------|-------------------------------------|------------------|----------|--------------|------|
| 22.669 | 13.38 ng/ul | 640205                              | Chrysene-d12     | 21.981   |              |      |
| Hit#   | of 5        | Tentative ID                        | MW               | MolForm  | CAS#         | Qual |
| 1      |             | 1,2-Benzenedicarboxylic acid, di... | 362              | C22H34O4 | 003648-21-3  | 87   |
| 2      |             | 1,2-Benzenedicarboxylic acid, di... | 362              | C22H34O4 | 003648-21-3  | 87   |
| 3      |             | Phthalic acid, decyl isobutyl ester | 362              | C22H34O4 | 1000308-94-2 | 86   |
| 4      |             | Phthalic acid, isobutyl 2-methyl... | 306              | C18H26O4 | 1000356-90-7 | 78   |
| 5      |             | Diisooheptyl phthalate              | 362              | C22H34O4 | 090937-19-2  | 72   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122021\  
Data File : BG051727.D  
Acq On : 21 Dec 2021 22:20  
Operator : CG/JU  
Sample : M5121-04DL 10X  
Misc :  
ALS Vial : 54 Sample Multiplier: 1

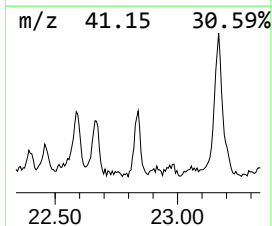
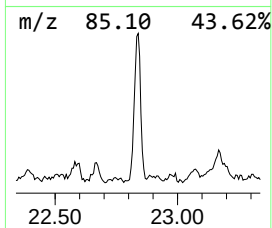
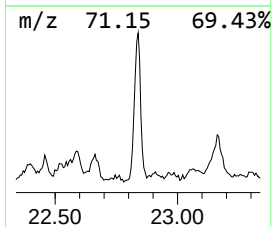
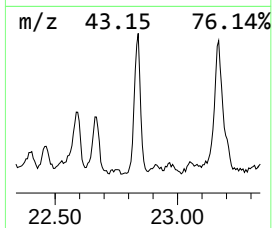
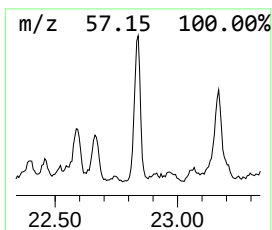
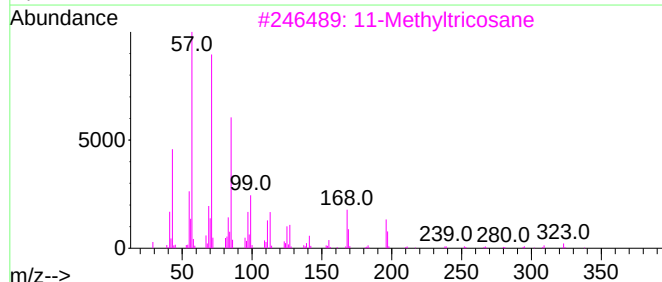
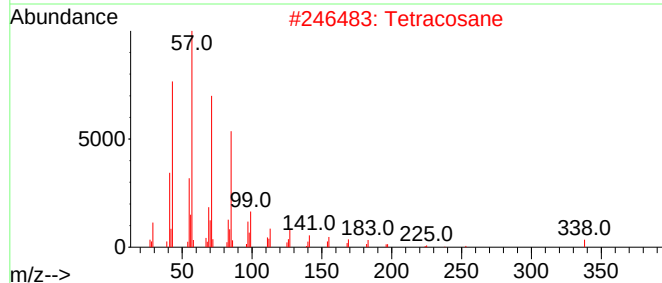
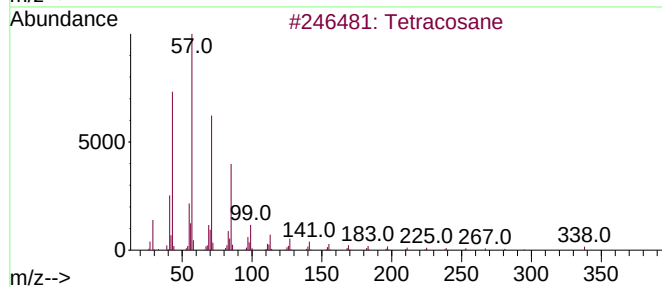
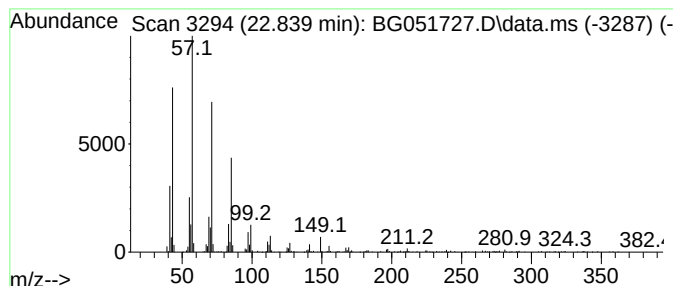
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BNA\_G  
ClientSampleId :  
BGL53DL

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 26 (DEL) Alkane: Straight-Chai... Concentration Rank 10

| R.T.      | EstConc            | Area   | Relative to ISTD | R.T.        |      |
|-----------|--------------------|--------|------------------|-------------|------|
| 22.839    | 11.75 ng/ul        | 561872 | Chrysene-d12     | 21.981      |      |
| Hit# of 5 | Tentative ID       | MW     | MolForm          | CAS#        | Qual |
| 1         | Tetracosane        | 338    | C24H50           | 000646-31-1 | 97   |
| 2         | Tetracosane        | 338    | C24H50           | 000646-31-1 | 95   |
| 3         | 11-Methyltricosane | 338    | C24H50           | 027538-41-6 | 90   |
| 4         | Octacosane         | 394    | C28H58           | 000630-02-4 | 90   |
| 5         | Heptacosane        | 380    | C27H56           | 000593-49-7 | 90   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122021\  
Data File : BG051727.D  
Acq On : 21 Dec 2021 22:20  
Operator : CG/JU  
Sample : M5121-04DL 10X  
Misc :  
ALS Vial : 54 Sample Multiplier: 1

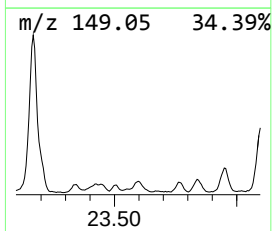
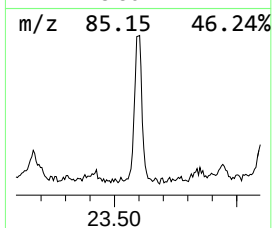
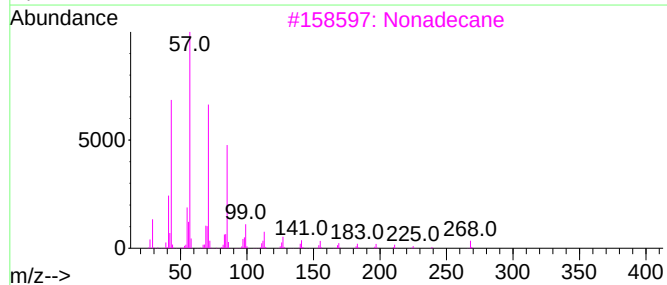
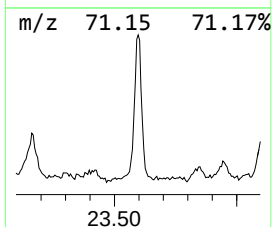
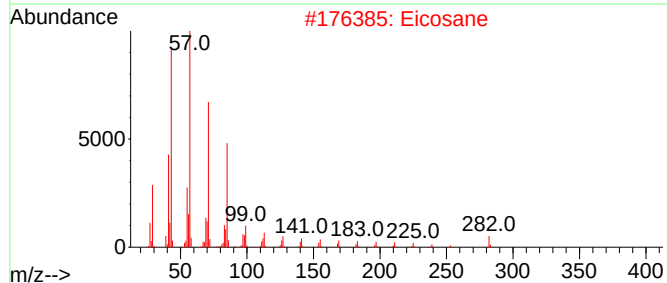
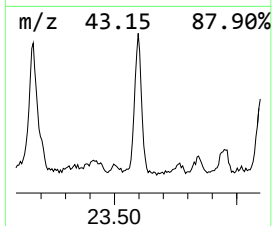
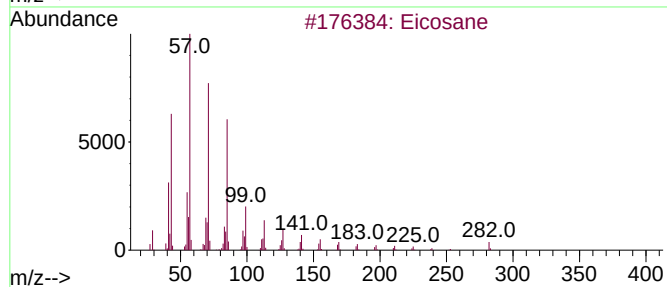
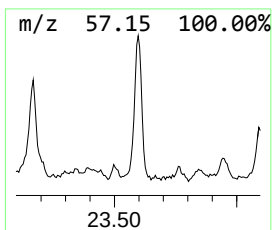
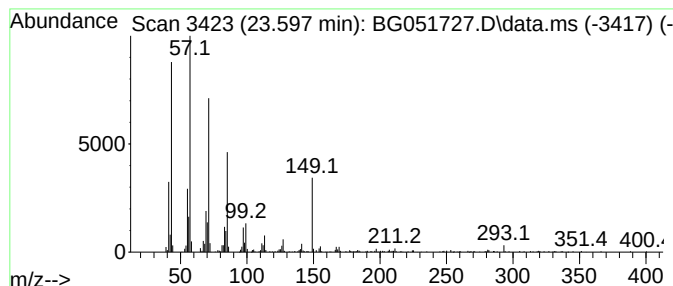
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BNA\_G  
ClientSampleId :  
BGL53DL

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 27 (DEL) Alkane: Straight-Chai... Concentration Rank 6

| R.T.      | EstConc      | Area   | Relative to ISTD |         | R.T.        |      |
|-----------|--------------|--------|------------------|---------|-------------|------|
| 23.597    | 14.47 ng/ul  | 692102 | Chrysene-d12     |         | 21.981      |      |
| Hit# of 5 | Tentative ID |        | MW               | MolForm | CAS#        | Qual |
| 1         | Eicosane     |        | 282              | C20H42  | 000112-95-8 | 95   |
| 2         | Eicosane     |        | 282              | C20H42  | 000112-95-8 | 95   |
| 3         | Nonadecane   |        | 268              | C19H40  | 000629-92-5 | 93   |
| 4         | Hexacosane   |        | 366              | C26H54  | 000630-01-3 | 93   |
| 5         | Heneicosane  |        | 296              | C21H44  | 000629-94-7 | 87   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122021\  
Data File : BG051727.D  
Acq On : 21 Dec 2021 22:20  
Operator : CG/JU  
Sample : M5121-04DL 10X  
Misc :  
ALS Vial : 54 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL53DL

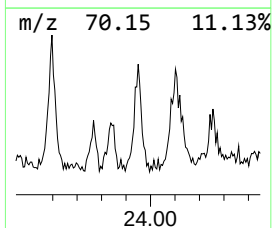
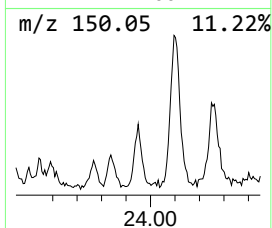
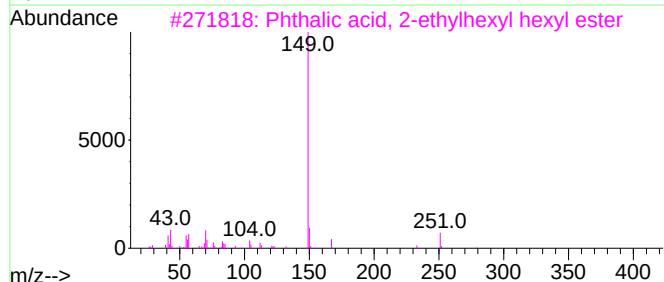
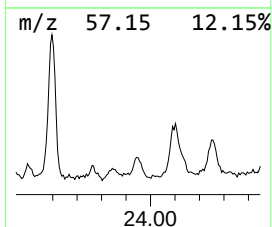
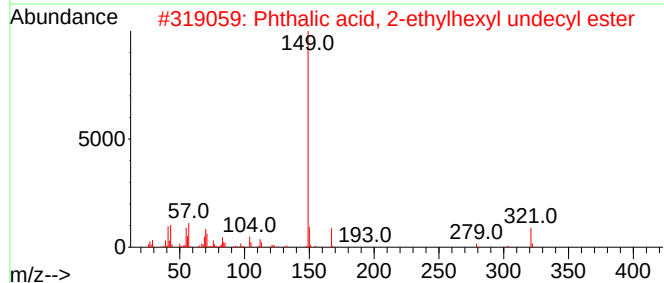
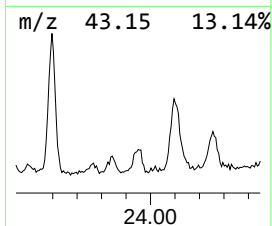
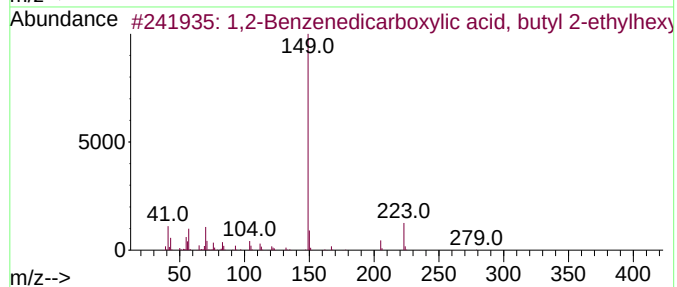
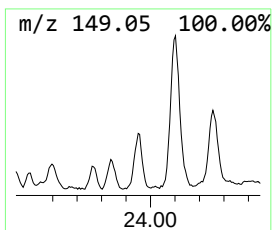
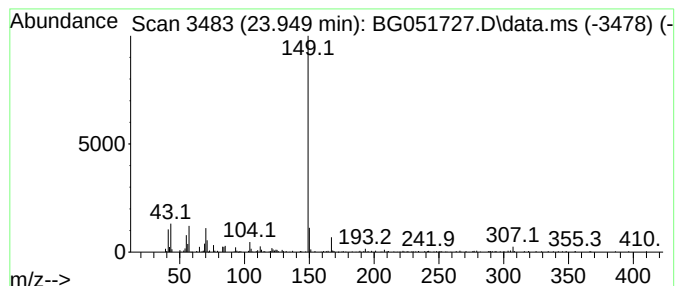
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Quant Title : SVOA CALIBRATION

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

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Peak Number 28 1,2-Benzenedicarboxylic aci... Concentration Rank 18

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 23.949 | 6.95 ng/ul | 265910 | Perylene-d12     | 25.482 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | 1,2-Benzenedicarboxylic acid, bu... | 334 | C20H30O4 | 000085-69-8  | 86   |
| 2    |    |   | Phthalic acid, 2-ethylhexyl unde... | 432 | C27H44O4 | 1000308-97-4 | 86   |
| 3    |    |   | Phthalic acid, 2-ethylhexyl hexy... | 362 | C22H34O4 | 1000309-02-5 | 83   |
| 4    |    |   | 1,2-Benzenedicarboxylic acid, di... | 362 | C22H34O4 | 003648-21-3  | 80   |
| 5    |    |   | Phthalic acid, 5-methylhex-2-yl ... | 320 | C19H28O4 | 1000371-09-1 | 80   |





Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122021\  
Data File : BG051727.D  
Acq On : 21 Dec 2021 22:20  
Operator : CG/JU  
Sample : M5121-04DL 10X  
Misc :  
ALS Vial : 54 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL53DL

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

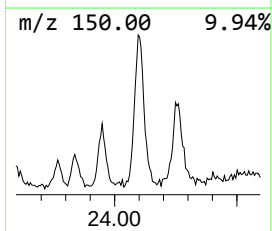
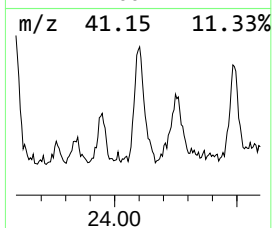
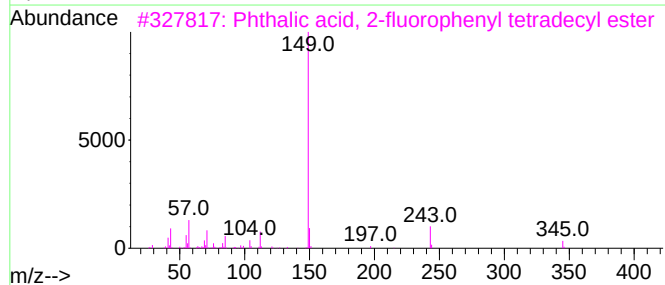
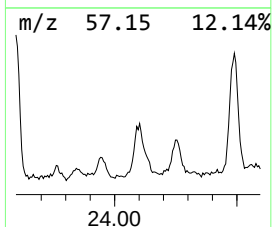
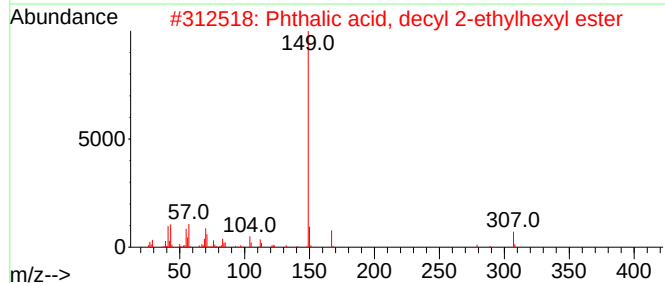
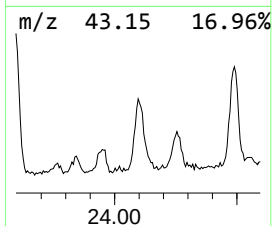
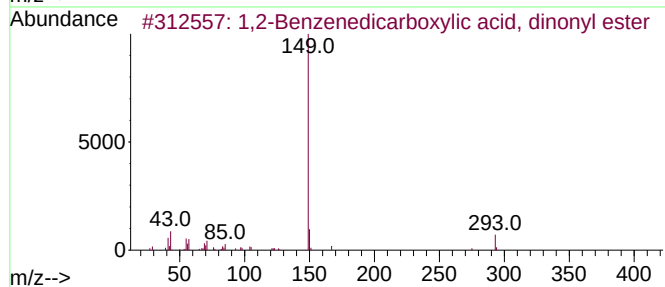
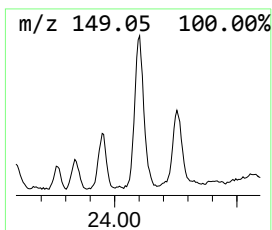
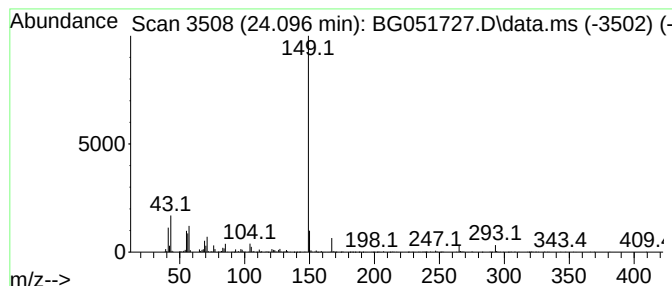
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 29 1,2-Benzenedicarboxylic aci... Concentration Rank 3

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 24.096 | 24.08 ng/ul | 920892 | Perylene-d12     | 25.482 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | 1,2-Benzenedicarboxylic acid, di... | 418 | C26H42O4  | 000084-76-4  | 86   |
| 2    |    |   | Phthalic acid, decyl 2-ethylhexy... | 418 | C26H42O4  | 1000309-00-0 | 86   |
| 3    |    |   | Phthalic acid, 2-fluorophenyl te... | 456 | C28H37FO4 | 1000356-50-3 | 72   |
| 4    |    |   | Phthalic acid, isobutyl nonyl ester | 348 | C21H32O4  | 1000309-04-4 | 72   |
| 5    |    |   | Phthalic acid, 3-fluorophenyl un... | 414 | C25H31FO4 | 1000356-66-9 | 72   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122021\  
Data File : BG051727.D  
Acq On : 21 Dec 2021 22:20  
Operator : CG/JU  
Sample : M5121-04DL 10X  
Misc :  
ALS Vial : 54 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL53DL

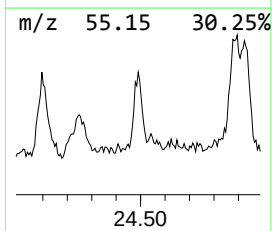
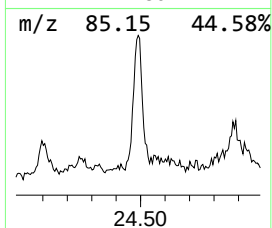
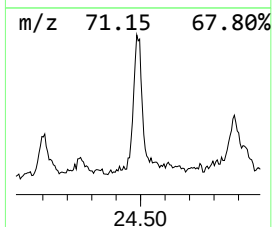
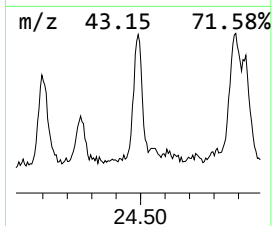
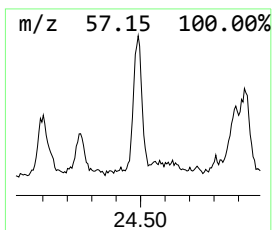
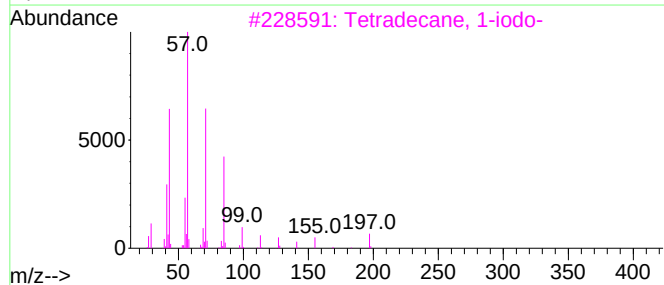
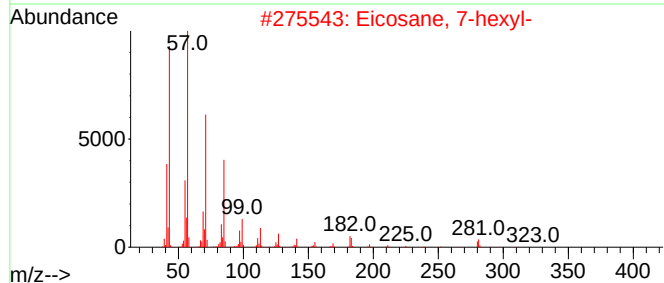
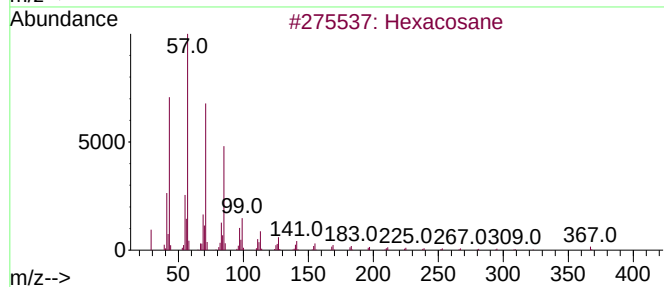
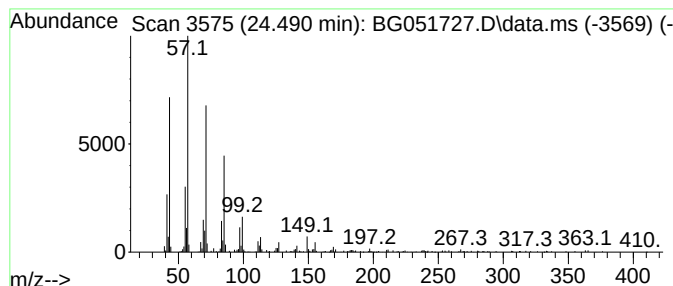
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Quant Title : SVOA CALIBRATION

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

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Peak Number 30 (DEL) Alkane: Straight-Chai... Concentration Rank 9

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 24.490 | 12.39 ng/ul | 473825 | Perylene-d12     | 25.482 |

| Hit# | of 5 | Tentative ID         | MW  | MolForm | CAS#        | Qual |
|------|------|----------------------|-----|---------|-------------|------|
| 1    |      | Hexacosane           | 366 | C26H54  | 000630-01-3 | 87   |
| 2    |      | Eicosane, 7-hexyl-   | 366 | C26H54  | 055333-99-8 | 86   |
| 3    |      | Tetradecane, 1-iodo- | 324 | C14H29I | 019218-94-1 | 83   |
| 4    |      | Heptadecane          | 240 | C17H36  | 000629-78-7 | 83   |
| 5    |      | Heptadecane          | 240 | C17H36  | 000629-78-7 | 83   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122021\  
Data File : BG051727.D  
Acq On : 21 Dec 2021 22:20  
Operator : CG/JU  
Sample : M5121-04DL 10X  
Misc :  
ALS Vial : 54 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL53DL

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

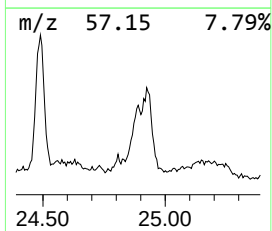
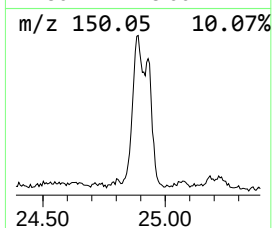
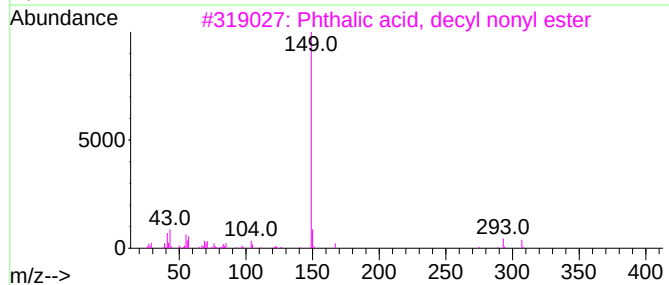
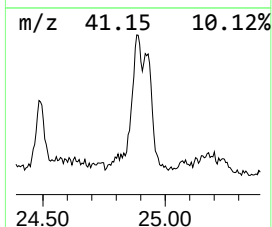
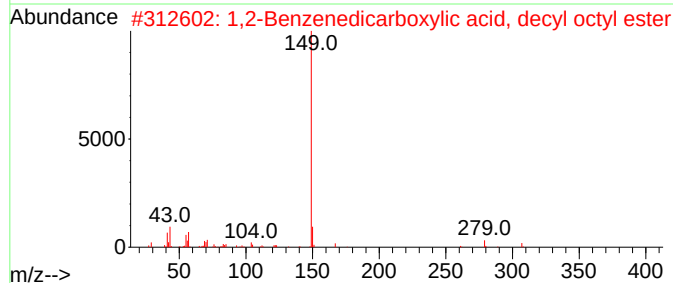
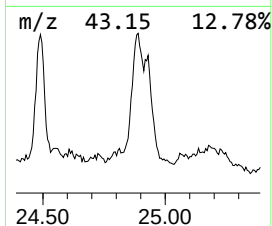
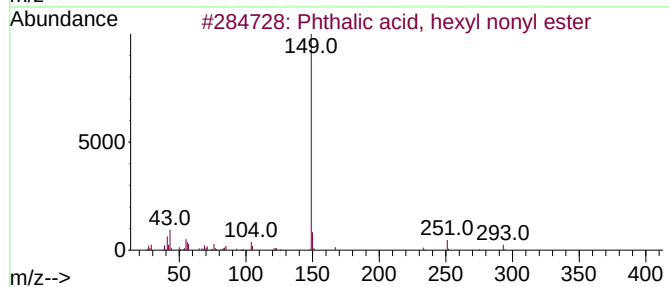
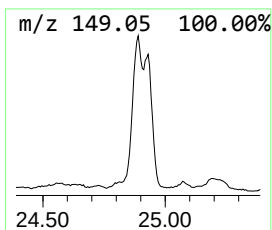
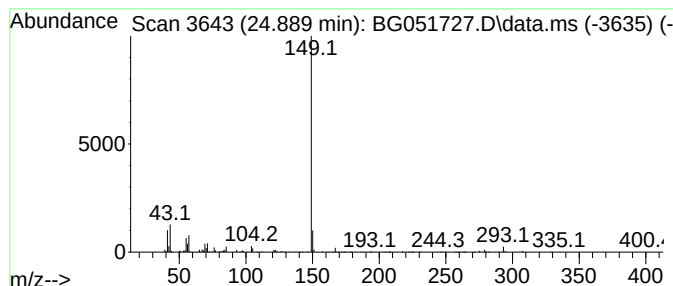
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 31 Phthalic acid, hexyl nonyl ... Concentration Rank 2

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 24.889 | 29.03 ng/ul | 1110420 | Perylene-d12     | 25.482 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Phthalic acid, hexyl nonyl ester    | 376 | C23H36O4 | 1000309-01-1 | 86   |
| 2    |    |   | 1,2-Benzenedicarboxylic acid, de... | 418 | C26H42O4 | 000119-07-3  | 80   |
| 3    |    |   | Phthalic acid, decyl nonyl ester    | 432 | C27H44O4 | 1000308-89-8 | 80   |
| 4    |    |   | Phthalic acid, nonyl pentyl ester   | 362 | C22H34O4 | 1000308-93-1 | 80   |
| 5    |    |   | Phthalic acid, dodecyl pentyl ester | 404 | C25H40O4 | 1000308-93-0 | 72   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122021\  
Data File : BG051727.D  
Acq On : 21 Dec 2021 22:20  
Operator : CG/JU  
Sample : M5121-04DL 10X  
Misc :  
ALS Vial : 54 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL53DL

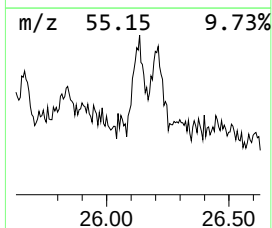
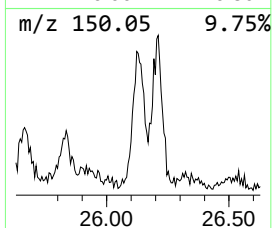
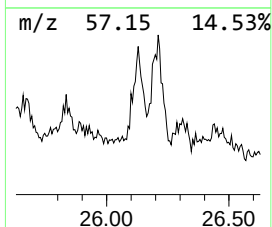
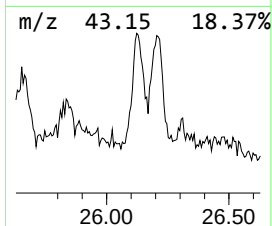
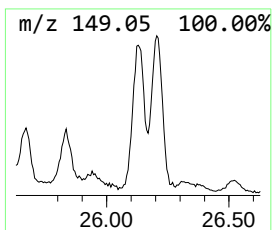
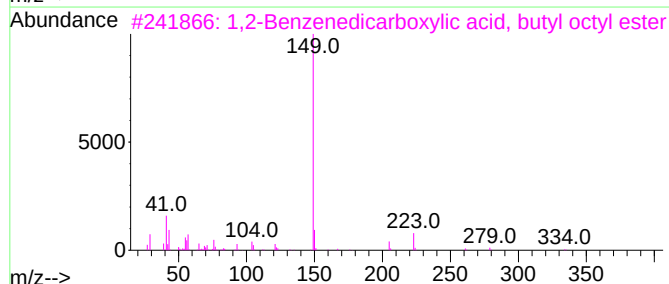
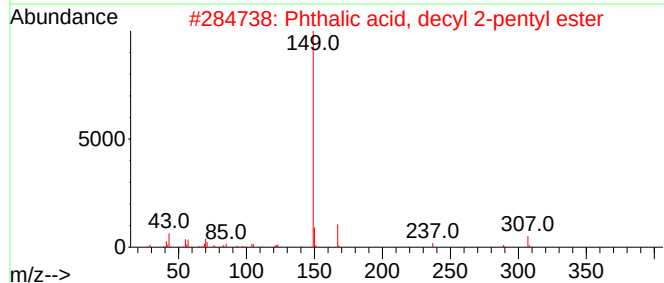
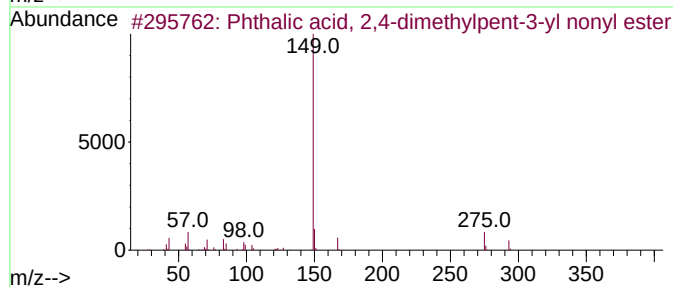
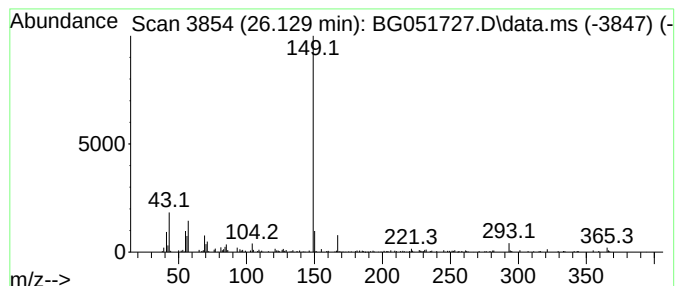
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Quant Title : SVOA CALIBRATION

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

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Peak Number 32 Phthalic acid, 2,4-dimethyl... Concentration Rank 16

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|--------|------------------|--------------|------|
| 26.129    | 8.54 ng/ul                          | 326696 | Perylene-d12     | 25.482       |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#         | Qual |
| 1         | Phthalic acid, 2,4-dimethylpent-... | 390    | C24H38O4         | 1000356-84-8 | 83   |
| 2         | Phthalic acid, decyl 2-pentyl ester | 376    | C23H36O4         | 1000315-48-2 | 78   |
| 3         | 1,2-Benzenedicarboxylic acid, bu... | 334    | C20H30O4         | 000084-78-6  | 74   |
| 4         | Phthalic acid, dec-2-yl isobutyl... | 362    | C22H34O4         | 1000356-93-8 | 72   |
| 5         | Phthalic acid, decyl hex-3-yl ester | 390    | C24H38O4         | 1000356-96-1 | 72   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122021\  
Data File : BG051727.D  
Acq On : 21 Dec 2021 22:20  
Operator : CG/JU  
Sample : M5121-04DL 10X  
Misc :  
ALS Vial : 54 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL53DL

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

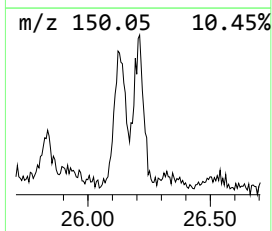
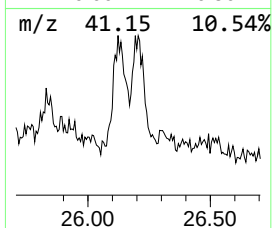
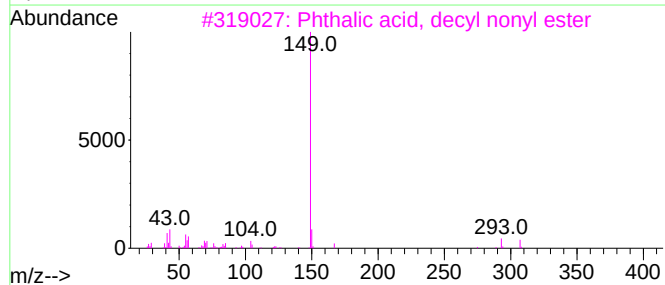
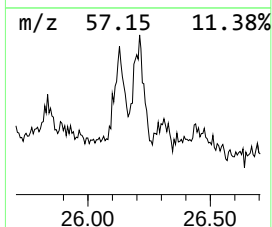
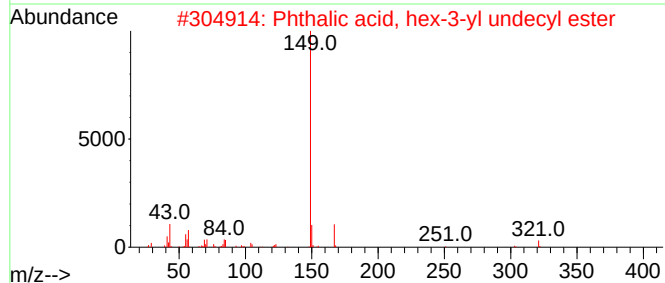
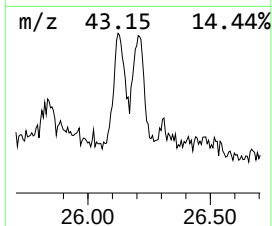
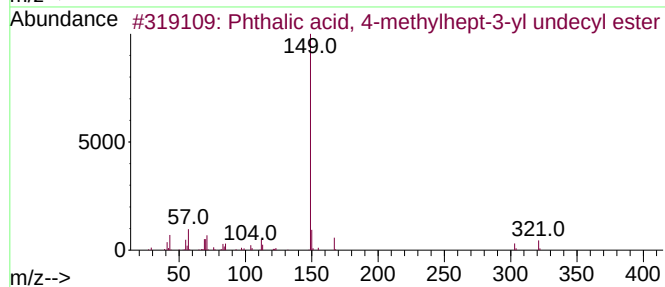
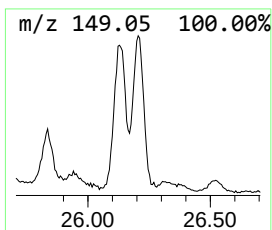
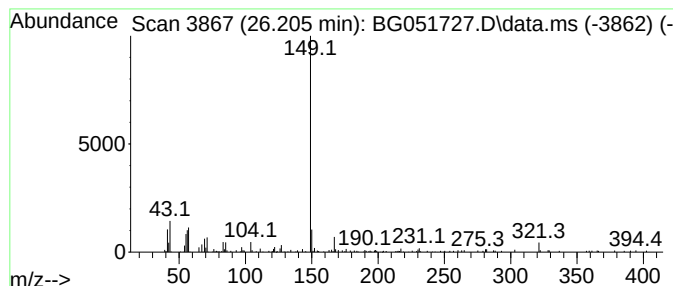
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 33 Phthalic acid, 4-methylhept... Concentration Rank 14

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 26.205 | 9.35 ng/ul | 357580 | Perylene-d12     | 25.482 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Phthalic acid, 4-methylhept-3-yl... | 432 | C27H44O4 | 1000377-94-6 | 80   |
| 2    |    |   | Phthalic acid, hex-3-yl undecyl ... | 404 | C25H40O4 | 1000356-96-2 | 72   |
| 3    |    |   | Phthalic acid, decyl nonyl ester    | 432 | C27H44O4 | 1000308-89-8 | 64   |
| 4    |    |   | Phthalic acid, hexyl dodecyl ester  | 418 | C26H42O4 | 1000308-91-5 | 64   |
| 5    |    |   | Phthalic acid, isobutyl 4-methyl... | 334 | C20H30O4 | 1000377-93-8 | 64   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122021\  
Data File : BG051727.D  
Acq On : 21 Dec 2021 22:20  
Operator : CG/JU  
Sample : M5121-04DL 10X  
Misc :  
ALS Vial : 54 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL53DL

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

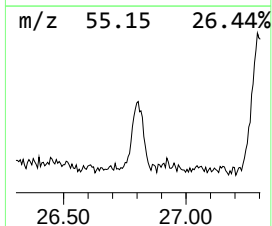
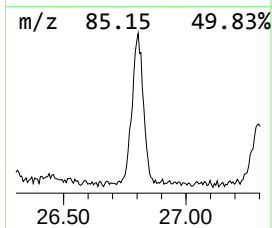
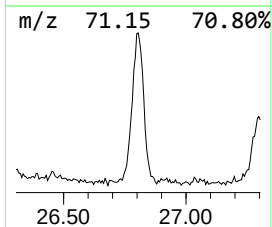
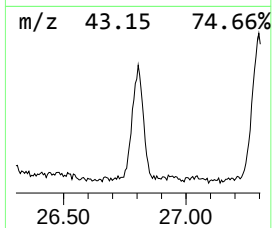
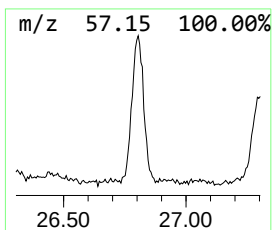
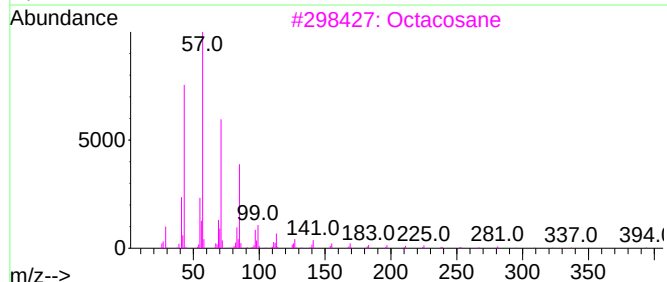
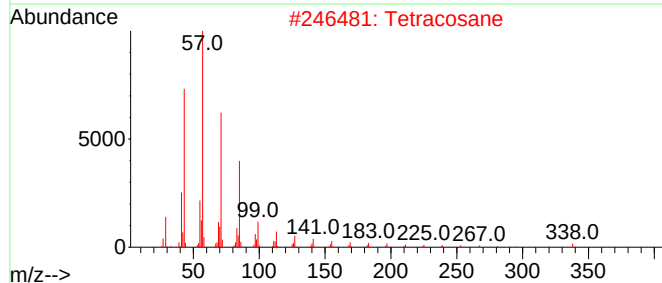
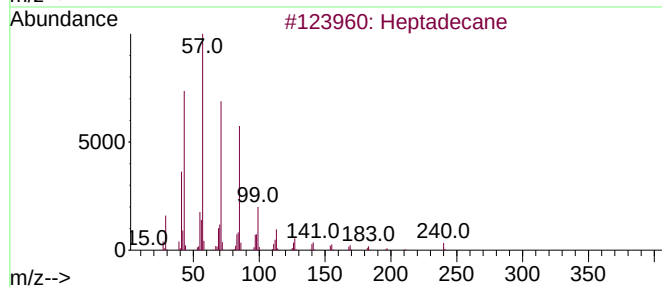
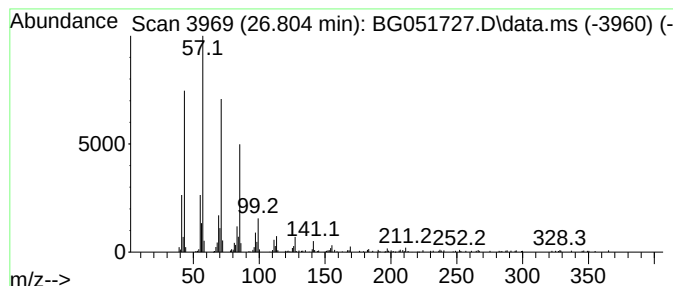
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 34 (DEL) Alkane: Straight-Chai... Concentration Rank 7

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 26.804 | 14.18 ng/ul | 542348 | Perylene-d12     | 25.482 |

| Hit# | of 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|------|--------------|-----|---------|-------------|------|
| 1    |      | Heptadecane  | 240 | C17H36  | 000629-78-7 | 93   |
| 2    |      | Tetracosane  | 338 | C24H50  | 000646-31-1 | 91   |
| 3    |      | Octacosane   | 394 | C28H58  | 000630-02-4 | 91   |
| 4    |      | Tetracosane  | 338 | C24H50  | 000646-31-1 | 91   |
| 5    |      | Hexacosane   | 366 | C26H54  | 000630-01-3 | 91   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122021\  
Data File : BG051727.D  
Acq On : 21 Dec 2021 22:20  
Operator : CG/JU  
Sample : M5121-04DL 10X  
Misc :  
ALS Vial : 54 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL53DL

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

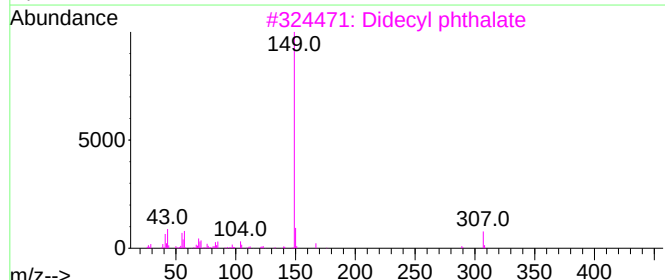
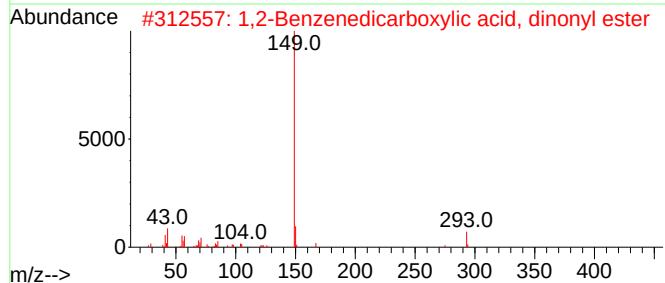
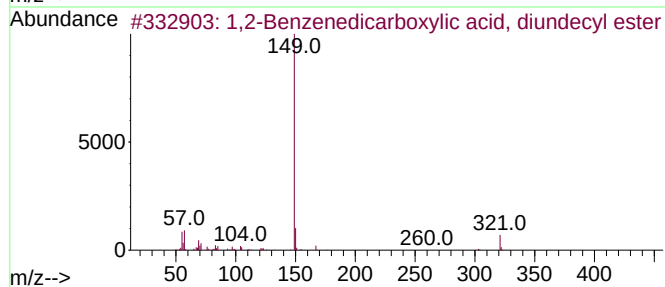
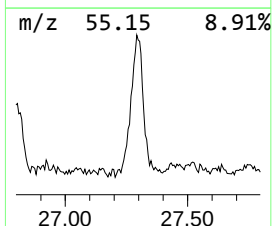
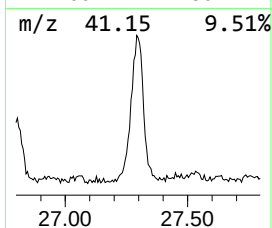
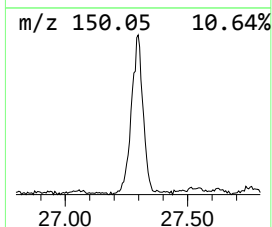
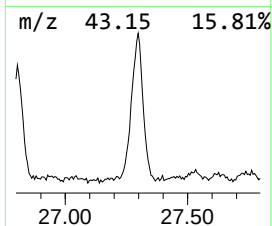
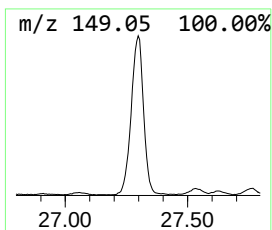
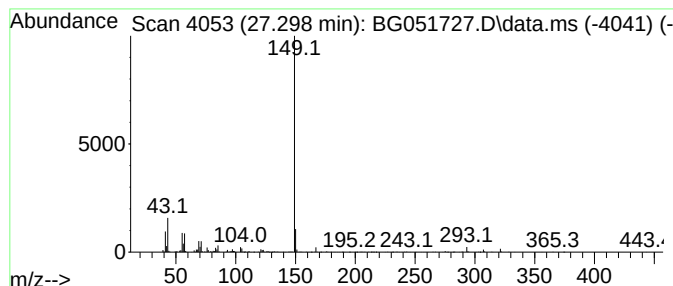
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 35 1,2-Benzenedicarboxylic aci... Concentration Rank 1

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 27.298 | 39.01 ng/ul | 1492150 | Perylene-d12     | 25.482 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | 1,2-Benzenedicarboxylic acid, di... | 474 | C30H5004 | 003648-20-2  | 83   |
| 2    |    |   | 1,2-Benzenedicarboxylic acid, di... | 418 | C26H4204 | 000084-76-4  | 83   |
| 3    |    |   | Didecyl phthalate                   | 446 | C28H4604 | 000084-77-5  | 83   |
| 4    |    |   | Di-n-octyl phthalate                | 390 | C24H3804 | 000117-84-0  | 80   |
| 5    |    |   | Phthalic acid, hept-3-yl undecyl... | 418 | C26H4204 | 1000356-99-5 | 80   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122021\  
Data File : BG051727.D  
Acq On : 21 Dec 2021 22:20  
Operator : CG/JU  
Sample : M5121-04DL 10X  
Misc :  
ALS Vial : 54 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL53DL

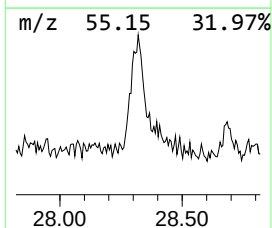
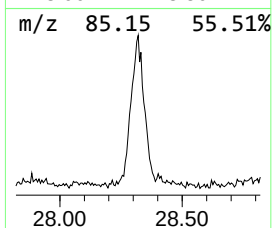
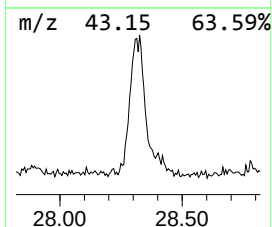
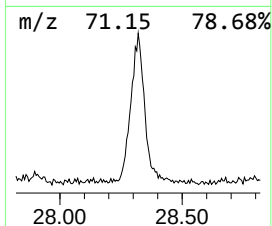
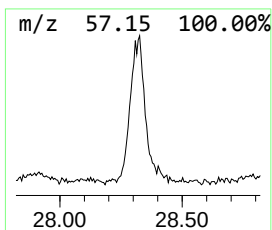
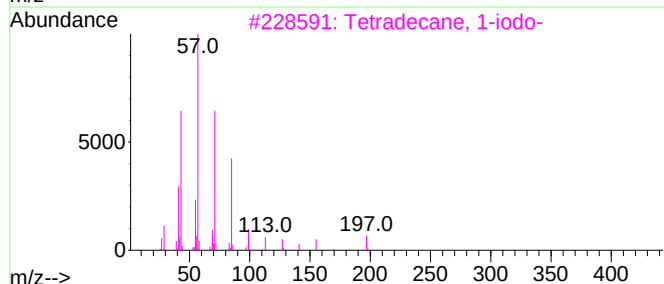
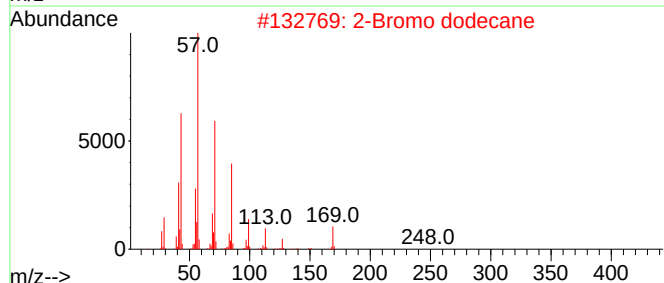
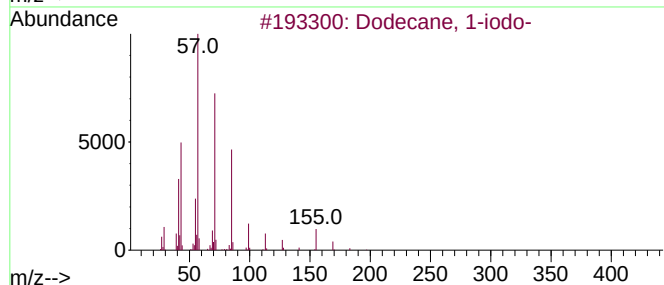
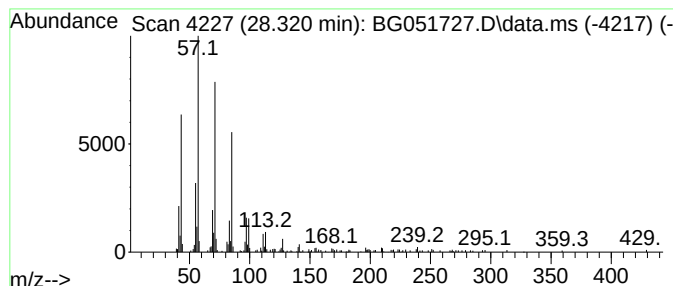
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Quant Title : SVOA CALIBRATION

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

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Peak Number 36 Dodecane, 1-iodo- Concentration Rank 13

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 28.320 | 10.74 ng/ul | 410883 | Perylene-d12     | 25.482 |

| Hit# | of | 5 | Tentative ID         | MW  | MolForm  | CAS#        | Qual |
|------|----|---|----------------------|-----|----------|-------------|------|
| 1    |    |   | Dodecane, 1-iodo-    | 296 | C12H25I  | 004292-19-7 | 86   |
| 2    |    |   | 2-Bromo dodecane     | 248 | C12H25Br | 013187-99-0 | 86   |
| 3    |    |   | Tetradecane, 1-iodo- | 324 | C14H29I  | 019218-94-1 | 86   |
| 4    |    |   | Decane, 1-iodo-      | 268 | C10H21I  | 002050-77-3 | 83   |
| 5    |    |   | Tetradecane          | 198 | C14H30   | 000629-59-4 | 83   |





Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122021\  
Data File : BG051727.D  
Acq On : 21 Dec 2021 22:20  
Operator : CG/JU  
Sample : M5121-04DL 10X  
Misc :  
ALS Vial : 54 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL53DL

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

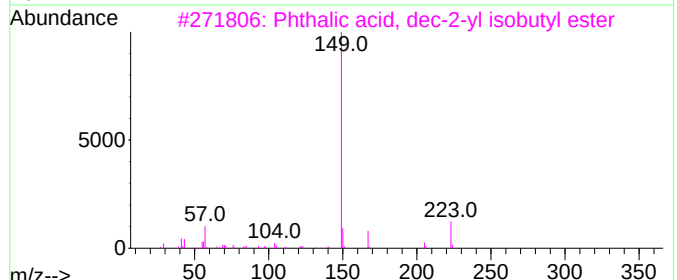
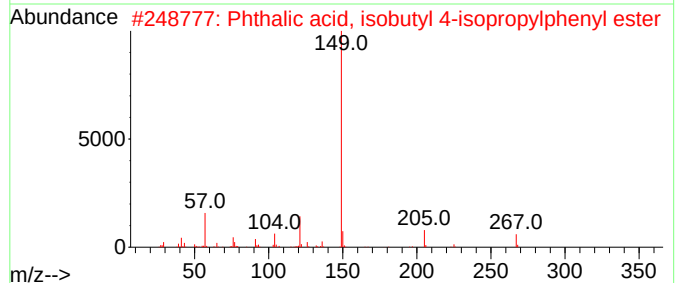
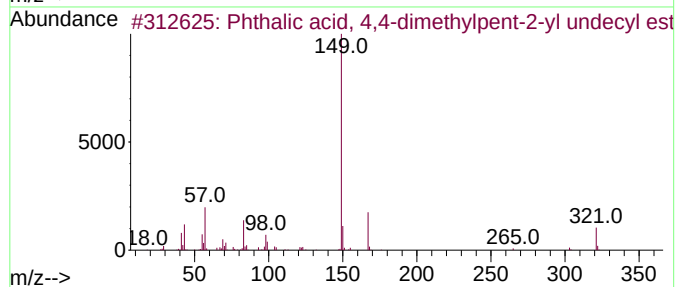
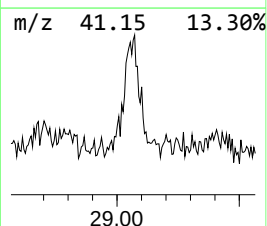
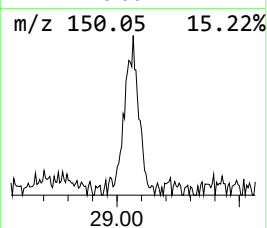
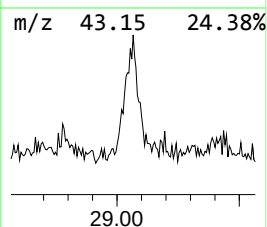
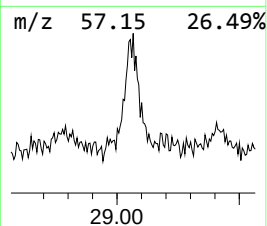
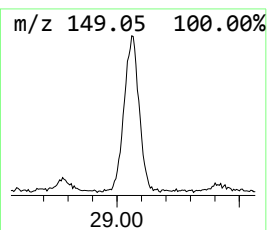
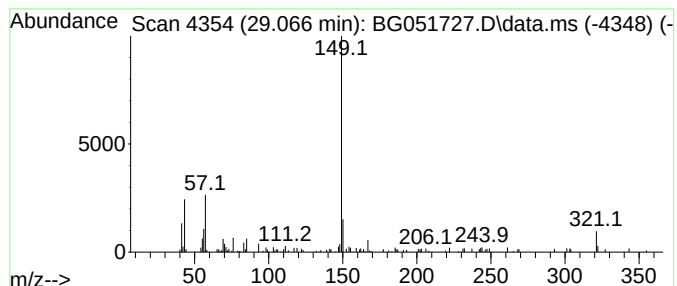
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 37 Phthalic acid, 4,4-dimethyl... Concentration Rank 17

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 29.066 | 7.14 ng/ul | 273104 | Perylene-d12     | 25.482 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Phthalic acid, 4,4-dimethylpent-... | 418 | C26H42O4  | 1000371-13-4 | 78   |
| 2    |    |   | Phthalic acid, isobutyl 4-isopro... | 340 | C21H24O4  | 1000315-22-7 | 59   |
| 3    |    |   | Phthalic acid, dec-2-yl isobutyl... | 362 | C22H34O4  | 1000356-93-8 | 59   |
| 4    |    |   | Phthalic acid, heptyl octyl ester   | 376 | C23H36O4  | 088216-58-4  | 59   |
| 5    |    |   | Phthalic acid, 3-fluorophenyl he... | 484 | C30H41FO4 | 1000356-66-4 | 59   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122021\  
Data File : BG051727.D  
Acq On : 21 Dec 2021 22:20  
Operator : CG/JU  
Sample : M5121-04DL 10X  
Misc :  
ALS Vial : 54 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL53DL

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

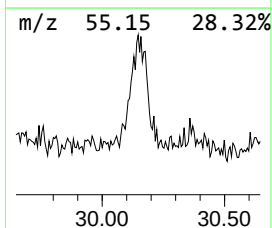
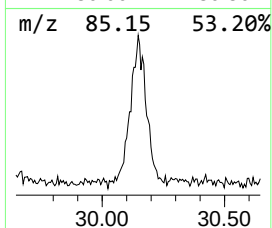
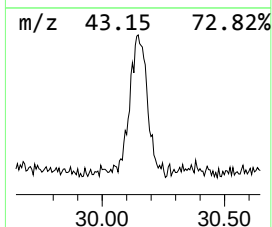
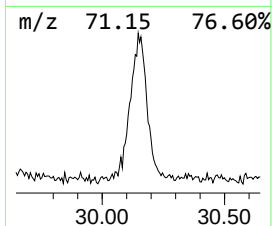
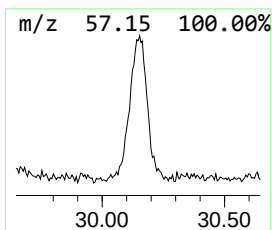
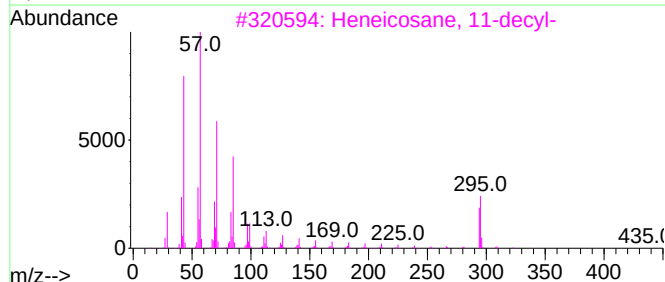
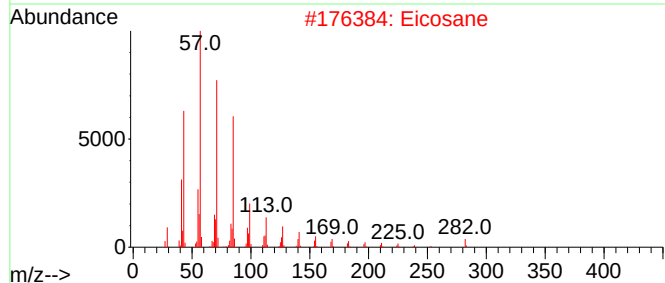
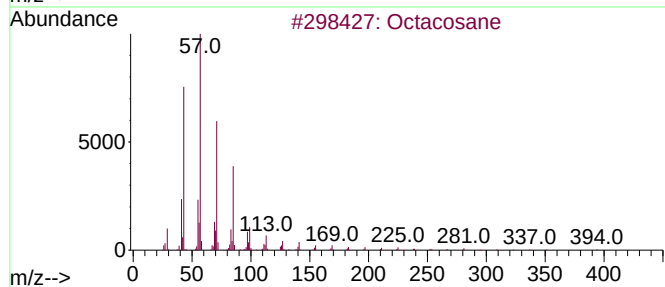
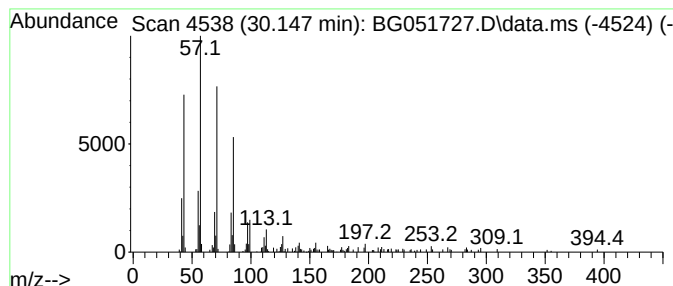
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 38 (DEL) Alkane: Straight-Chai... Concentration Rank 12

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 30.147 | 10.99 ng/ul | 420232 | Perylene-d12     | 25.482 |

| Hit# | of 5 | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|------|------------------------|-----|---------|-------------|------|
| 1    |      | Octacosane             | 394 | C28H58  | 000630-02-4 | 94   |
| 2    |      | Eicosane               | 282 | C20H42  | 000112-95-8 | 94   |
| 3    |      | Heneicosane, 11-decyl- | 437 | C31H64  | 055320-06-4 | 91   |
| 4    |      | Pentacosane            | 352 | C25H52  | 000629-99-2 | 91   |
| 5    |      | Octadecane             | 254 | C18H38  | 000593-45-3 | 91   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122021\  
Data File : BG051727.D  
Acq On : 21 Dec 2021 22:20  
Operator : CG/JU  
Sample : M5121-04DL 10X  
Misc :  
ALS Vial : 54 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL53DL

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

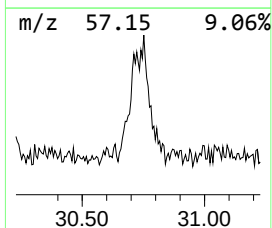
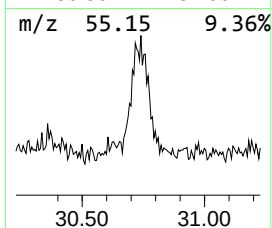
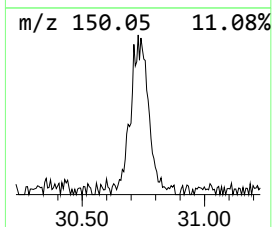
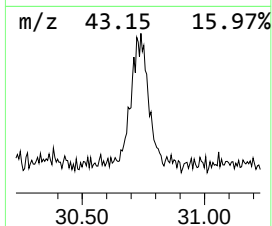
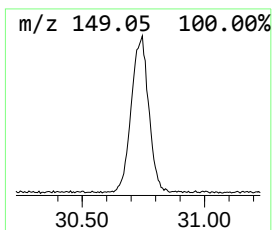
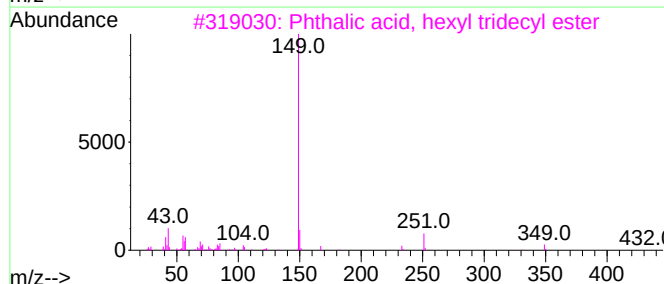
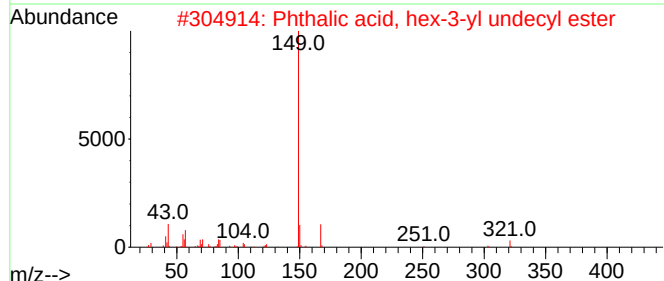
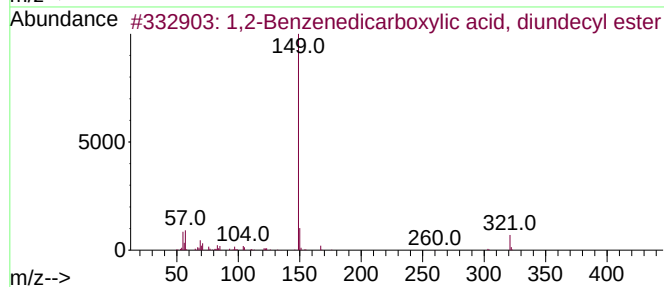
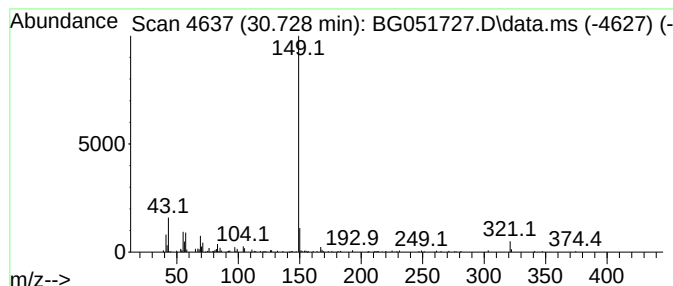
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 39 Phthalic acid, hex-3-yl und... Concentration Rank 21

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 30.728 | 5.47 ng/ul | 209300 | Perylene-d12     | 25.482 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | 1,2-Benzenedicarboxylic acid, di... | 474 | C30H50O4 | 003648-20-2  | 87   |
| 2    |    |   | Phthalic acid, hex-3-yl undecyl ... | 404 | C25H40O4 | 1000356-96-2 | 86   |
| 3    |    |   | Phthalic acid, hexyl tridecyl ester | 432 | C27H44O4 | 1000308-99-1 | 72   |
| 4    |    |   | Phthalic acid, 3-methylbutyl non... | 502 | C32H54O4 | 1010356-82-0 | 72   |
| 5    |    |   | Phthalic acid, dodecyl octyl ester  | 446 | C28H46O4 | 1010308-91-7 | 72   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122021\  
Data File : BG051727.D  
Acq On : 21 Dec 2021 22:20  
Operator : CG/JU  
Sample : M5121-04DL 10X  
Misc :  
ALS Vial : 54 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGL53DL

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

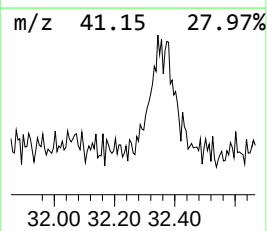
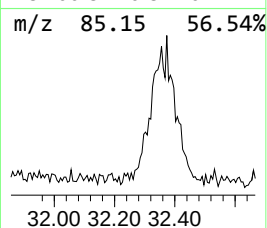
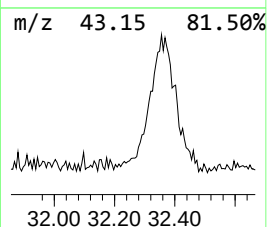
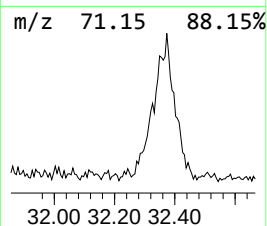
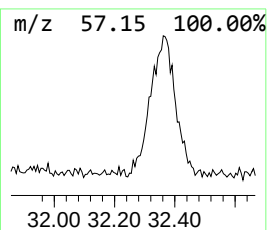
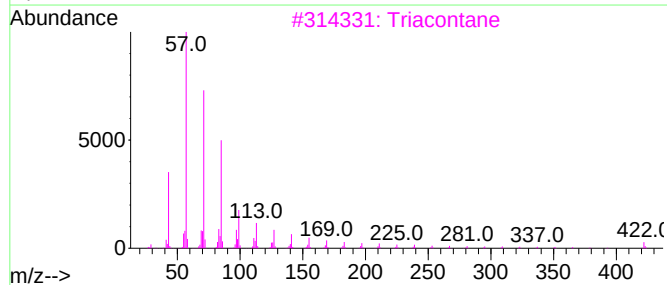
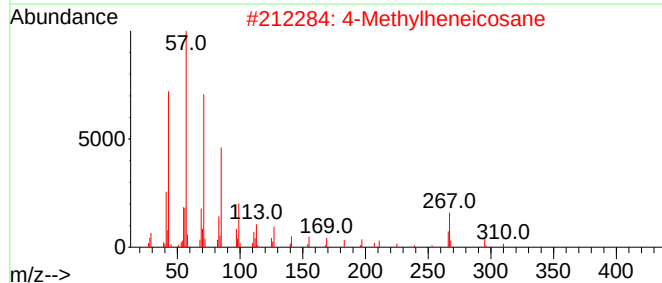
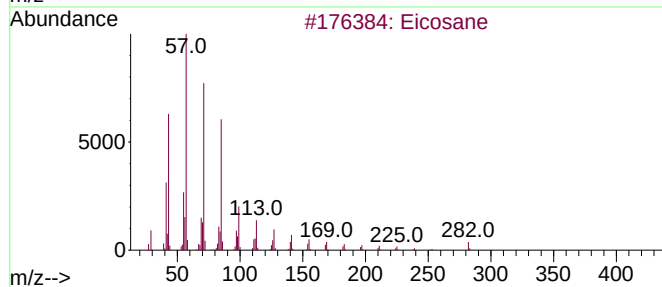
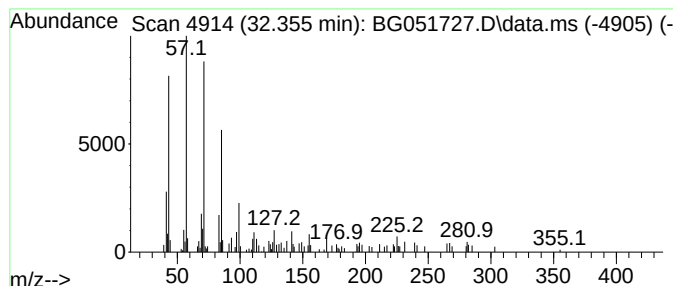
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*

Peak Number 40 (DEL) Alkane: Straight-Chai... Concentration Rank 33

| R.T.   | EstConc    | Area  | Relative to ISTD | R.T.   |
|--------|------------|-------|------------------|--------|
| 32.355 | 2.49 ng/ul | 95324 | Perylene-d12     | 25.482 |

| Hit# | of 5 | Tentative ID           | MW  | MolForm | CAS#         | Qual |
|------|------|------------------------|-----|---------|--------------|------|
| 1    |      | Eicosane               | 282 | C20H42  | 000112-95-8  | 93   |
| 2    |      | 4-Methylheneicosane    | 310 | C22H46  | 025117-29-7  | 86   |
| 3    |      | Triacontane            | 422 | C30H62  | 000638-68-6  | 86   |
| 4    |      | Dotriacontane, 1-iodo- | 576 | C32H65I | 1000406-32-4 | 80   |
| 5    |      | Hexadecane, 1-iodo-    | 352 | C16H33I | 000544-77-4  | 80   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122021\  
 Data File : BG051727.D  
 Acq On : 21 Dec 2021 22:20  
 Operator : CG/JU  
 Sample : M5121-04DL 10X  
 Misc :  
 ALS Vial : 54 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 BGL53DL

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... | 13.722 | 2.9     | ng/ul | 114004   | 3                     | 14.909 | 791628 | 20.0 |
| (DEL) Alkane: S... | 14.785 | 2.3     | ng/ul | 92786    | 3                     | 14.909 | 791628 | 20.0 |
| Tetradecane, 1-... | 15.737 | 3.1     | ng/ul | 121307   | 3                     | 14.909 | 791628 | 20.0 |
| (DEL) Alkane: S... | 16.142 | 2.3     | ng/ul | 92032    | 3                     | 14.909 | 791628 | 20.0 |
| (DEL) Alkane: S... | 16.595 | 3.4     | ng/ul | 156356   | 4                     | 17.658 | 923949 | 20.0 |
| (DEL) Alkane: S... | 16.624 | 4.3     | ng/ul | 199150   | 4                     | 17.658 | 923949 | 20.0 |
| unknown-01         | 16.724 | 2.1     | ng/ul | 96267    | 4                     | 17.658 | 923949 | 20.0 |
| (DEL) Alkane: S... | 17.393 | 2.5     | ng/ul | 114745   | 4                     | 17.658 | 923949 | 20.0 |
| (DEL) Alkane: S... | 17.446 | 2.3     | ng/ul | 106375   | 4                     | 17.658 | 923949 | 20.0 |
| 2-Bromo dodecane   | 18.134 | 3.4     | ng/ul | 156085   | 4                     | 17.658 | 923949 | 20.0 |
| (DEL) Alkane: S... | 18.821 | 2.4     | ng/ul | 108790   | 4                     | 17.658 | 923949 | 20.0 |
| p-Dicyclohexylb... | 19.085 | 3.5     | ng/ul | 161729   | 4                     | 17.658 | 923949 | 20.0 |
| unknown-02         | 19.191 | 3.8     | ng/ul | 175884   | 4                     | 17.658 | 923949 | 20.0 |
| (DEL) Alkane: S... | 19.438 | 3.4     | ng/ul | 159109   | 4                     | 17.658 | 923949 | 20.0 |
| cyclohexane, [1... | 19.626 | 2.4     | ng/ul | 112262   | 4                     | 17.658 | 923949 | 20.0 |
| (DEL) Alkane: S... | 20.542 | 4.8     | ng/ul | 230084   | 5                     | 21.981 | 956749 | 20.0 |
| (DEL) Alkane: S... | 21.047 | 6.1     | ng/ul | 291689   | 5                     | 21.981 | 956749 | 20.0 |
| 1,2-Benzenedica... | 21.312 | 16.2    | ng/ul | 776592   | 5                     | 21.981 | 956749 | 20.0 |
| Phthalic acid, ... | 21.464 | 5.4     | ng/ul | 258369   | 5                     | 21.981 | 956749 | 20.0 |
| (DEL) Alkane: S... | 21.582 | 8.9     | ng/ul | 426324   | 5                     | 21.981 | 956749 | 20.0 |
| Hexacosane         | 22.175 | 16.7    | ng/ul | 800450   | 5                     | 21.981 | 956749 | 20.0 |
| Phthalic acid, ... | 22.398 | 3.0     | ng/ul | 142123   | 5                     | 21.981 | 956749 | 20.0 |
| 1,2-Benzenedica... | 22.463 | 5.9     | ng/ul | 280619   | 5                     | 21.981 | 956749 | 20.0 |
| Phthalic acid, ... | 22.592 | 11.4    | ng/ul | 543758   | 5                     | 21.981 | 956749 | 20.0 |
| 1,2-Benzenedica... | 22.669 | 13.4    | ng/ul | 640205   | 5                     | 21.981 | 956749 | 20.0 |
| (DEL) Alkane: S... | 22.839 | 11.8    | ng/ul | 561872   | 5                     | 21.981 | 956749 | 20.0 |
| (DEL) Alkane: S... | 23.597 | 14.5    | ng/ul | 692102   | 5                     | 21.981 | 956749 | 20.0 |
| 1,2-Benzenedica... | 23.949 | 7.0     | ng/ul | 265910   | 6                     | 25.482 | 764983 | 20.0 |
| 1,2-Benzenedica... | 24.096 | 24.1    | ng/ul | 920892   | 6                     | 25.482 | 764983 | 20.0 |
| (DEL) Alkane: S... | 24.490 | 12.4    | ng/ul | 473825   | 6                     | 25.482 | 764983 | 20.0 |
| Phthalic acid, ... | 24.889 | 29.0    | ng/ul | 1110420  | 6                     | 25.482 | 764983 | 20.0 |
| Phthalic acid, ... | 26.129 | 8.5     | ng/ul | 326696   | 6                     | 25.482 | 764983 | 20.0 |
| Phthalic acid, ... | 26.205 | 9.3     | ng/ul | 357580   | 6                     | 25.482 | 764983 | 20.0 |
| (DEL) Alkane: S... | 26.804 | 14.2    | ng/ul | 542348   | 6                     | 25.482 | 764983 | 20.0 |
| 1,2-Benzenedica... | 27.298 | 39.0    | ng/ul | 1492150  | 6                     | 25.482 | 764983 | 20.0 |
| Dodecane, 1-iodo-  | 28.320 | 10.7    | ng/ul | 410883   | 6                     | 25.482 | 764983 | 20.0 |
| Phthalic acid, ... | 29.066 | 7.1     | ng/ul | 273104   | 6                     | 25.482 | 764983 | 20.0 |
| (DEL) Alkane: S... | 30.147 | 11.0    | ng/ul | 420232   | 6                     | 25.482 | 764983 | 20.0 |
| Phthalic acid, ... | 30.728 | 5.5     | ng/ul | 209300   | 6                     | 25.482 | 764983 | 20.0 |
| (DEL) Alkane: S... | 32.355 | 2.5     | ng/ul | 95324    | 6                     | 25.482 | 764983 | 20.0 |