

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011522\  
 Data File : BG052012.D  
 Acq On : 16 Jan 2022 1:50  
 Operator : CG/JU  
 Sample : N1132-04 10X  
 Misc :  
 ALS Vial : 23 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 BGLT5

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-BG010622.M  
 Title : SVOA CALIBRATION

Signal : TIC: BG052012.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         | 5.849       | 395           | 402         | 412          | rBV      | 49642          | 110594        | 0.40%           | 0.127%        |
| 2         | 7.382       | 658           | 663         | 670          | rVV3     | 66285          | 121795        | 0.44%           | 0.140%        |
| 3         | 7.458       | 670           | 676         | 683          | rVB      | 59537          | 105263        | 0.38%           | 0.121%        |
| 4         | 7.858       | 739           | 744         | 747          | rBV      | 178012         | 299992        | 1.09%           | 0.344%        |
| 5         | 7.887       | 747           | 749         | 758          | rVB      | 169067         | 235725        | 0.86%           | 0.271%        |
| 6         | 8.245       | 800           | 810         | 815          | rVV3     | 123733         | 321608        | 1.17%           | 0.369%        |
| 7         | 8.369       | 826           | 831         | 837          | rVB3     | 50552          | 95176         | 0.35%           | 0.109%        |
| 8         | 8.798       | 896           | 904         | 908          | rVV4     | 94174          | 228173        | 0.83%           | 0.262%        |
| 9         | 8.903       | 915           | 922         | 935          | rVB6     | 110566         | 289586        | 1.05%           | 0.332%        |
| 10        | 9.015       | 935           | 941         | 945          | rBV2     | 73377          | 118212        | 0.43%           | 0.136%        |
| 11        | 9.367       | 991           | 1001        | 1007         | rBV2     | 84175          | 190744        | 0.69%           | 0.219%        |
| 12        | 9.485       | 1013          | 1021        | 1026         | rVB      | 494361         | 811878        | 2.95%           | 0.932%        |
| 13        | 9.738       | 1053          | 1064        | 1073         | rVB7     | 45159          | 150743        | 0.55%           | 0.173%        |
| 14        | 9.902       | 1086          | 1092        | 1096         | rBV2     | 64027          | 114301        | 0.42%           | 0.131%        |
| 15        | 10.266      | 1147          | 1154        | 1159         | rVV7     | 53903          | 115080        | 0.42%           | 0.132%        |
| 16        | 10.337      | 1159          | 1166        | 1173         | rVB4     | 59441          | 156523        | 0.57%           | 0.180%        |
| 17        | 10.413      | 1173          | 1179        | 1184         | rBV2     | 63062          | 114240        | 0.42%           | 0.131%        |
| 18        | 10.489      | 1185          | 1192        | 1197         | rBV4     | 121536         | 234172        | 0.85%           | 0.269%        |
| 19        | 10.595      | 1205          | 1210        | 1215         | rBV      | 57733          | 94415         | 0.34%           | 0.108%        |
| 20        | 11.042      | 1281          | 1286        | 1291         | rVV      | 503870         | 846270        | 3.07%           | 0.971%        |
| 21        | 11.089      | 1291          | 1294        | 1297         | rVV      | 158554         | 263428        | 0.96%           | 0.302%        |
| 22        | 11.136      | 1297          | 1302        | 1309         | rVB      | 441993         | 839527        | 3.05%           | 0.964%        |
| 23        | 11.236      | 1314          | 1319        | 1324         | rVB      | 168194         | 279982        | 1.02%           | 0.321%        |
| 24        | 11.788      | 1408          | 1413        | 1418         | rBV5     | 74952          | 151965        | 0.55%           | 0.174%        |
| 25        | 11.911      | 1429          | 1434        | 1440         | rVB4     | 87537          | 160042        | 0.58%           | 0.184%        |
| 26        | 11.987      | 1442          | 1447        | 1455         | rVB4     | 120659         | 222900        | 0.81%           | 0.256%        |
| 27        | 12.093      | 1458          | 1465        | 1469         | rBV      | 270549         | 453939        | 1.65%           | 0.521%        |
| 28        | 12.181      | 1477          | 1480        | 1486         | rVB4     | 57346          | 94434         | 0.34%           | 0.108%        |
| 29        | 12.481      | 1523          | 1531        | 1541         | rBV      | 828312         | 1288697       | 4.68%           | 1.479%        |
| 30        | 12.616      | 1551          | 1554        | 1560         | rVV4     | 62497          | 138561        | 0.50%           | 0.159%        |
| 31        | 12.698      | 1562          | 1568        | 1570         | rVV2     | 103629         | 214162        | 0.78%           | 0.246%        |
| 32        | 12.733      | 1570          | 1574        | 1581         | rVB      | 316651         | 537197        | 1.95%           | 0.617%        |
| 33        | 12.951      | 1606          | 1611        | 1618         | rVB      | 206813         | 366203        | 1.33%           | 0.420%        |
| 34        | 13.104      | 1632          | 1637        | 1641         | rBV2     | 111451         | 195742        | 0.71%           | 0.225%        |
| 35        | 13.145      | 1641          | 1644        | 1646         | rVV3     | 75204          | 114065        | 0.41%           | 0.131%        |
| 36        | 13.174      | 1646          | 1649        | 1652         | rVV      | 91549          | 141997        | 0.52%           | 0.163%        |

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Integrator: RTE

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Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010622.M  
 Title : SVOA CALIBRATION

|    |        |      |      |      |      |         |         |        |        |
|----|--------|------|------|------|------|---------|---------|--------|--------|
| 37 | 13.280 | 1663 | 1667 | 1672 | rVB2 | 157158  | 220378  | 0.80%  | 0.253% |
| 38 | 13.362 | 1678 | 1681 | 1686 | rVB  | 103156  | 140043  | 0.51%  | 0.161% |
| 39 | 13.421 | 1686 | 1691 | 1698 | rVB  | 381101  | 536931  | 1.95%  | 0.616% |
| 40 | 13.709 | 1733 | 1740 | 1748 | rBV  | 1896486 | 2871130 | 10.43% | 3.295% |
| 41 | 13.926 | 1772 | 1777 | 1778 | rBV2 | 93230   | 127149  | 0.46%  | 0.146% |
| 42 | 14.055 | 1794 | 1799 | 1810 | rVB3 | 202603  | 615583  | 2.24%  | 0.707% |
| 43 | 14.220 | 1821 | 1827 | 1831 | rBV4 | 362527  | 653619  | 2.37%  | 0.750% |
| 44 | 14.267 | 1831 | 1835 | 1841 | rVB2 | 338507  | 522369  | 1.90%  | 0.600% |
| 45 | 14.331 | 1841 | 1846 | 1847 | rBV2 | 274679  | 395229  | 1.44%  | 0.454% |
| 46 | 14.361 | 1847 | 1851 | 1855 | rVB  | 579925  | 737238  | 2.68%  | 0.846% |
| 47 | 14.396 | 1855 | 1857 | 1862 | rVB3 | 194481  | 226284  | 0.82%  | 0.260% |
| 48 | 14.725 | 1909 | 1913 | 1916 | rBV4 | 166113  | 231878  | 0.84%  | 0.266% |
| 49 | 14.772 | 1916 | 1921 | 1926 | rVV  | 1950255 | 2632799 | 9.56%  | 3.022% |
| 50 | 14.854 | 1930 | 1935 | 1938 | rBV4 | 210891  | 423218  | 1.54%  | 0.486% |
| 51 | 14.889 | 1938 | 1941 | 1947 | rVB  | 305097  | 487105  | 1.77%  | 0.559% |
| 52 | 15.101 | 1972 | 1977 | 1984 | rVB3 | 168857  | 378398  | 1.37%  | 0.434% |
| 53 | 15.207 | 1993 | 1995 | 1998 | rBV4 | 90517   | 125718  | 0.46%  | 0.144% |
| 54 | 15.277 | 2002 | 2007 | 2012 | rVV4 | 323197  | 608374  | 2.21%  | 0.698% |
| 55 | 15.324 | 2013 | 2015 | 2018 | rVV  | 88880   | 102802  | 0.37%  | 0.118% |
| 56 | 15.383 | 2018 | 2025 | 2033 | rVB5 | 302661  | 857474  | 3.12%  | 0.984% |
| 57 | 15.453 | 2035 | 2037 | 2042 | rVB3 | 146725  | 195407  | 0.71%  | 0.224% |
| 58 | 15.506 | 2042 | 2046 | 2051 | rBV3 | 169856  | 298591  | 1.08%  | 0.343% |
| 59 | 15.559 | 2051 | 2055 | 2059 | rBV  | 182743  | 256676  | 0.93%  | 0.295% |
| 60 | 15.682 | 2069 | 2076 | 2078 | rBV6 | 264357  | 503506  | 1.83%  | 0.578% |
| 61 | 15.724 | 2078 | 2083 | 2087 | rVB  | 1903403 | 2511470 | 9.12%  | 2.883% |
| 62 | 15.823 | 2096 | 2100 | 2108 | rVB3 | 252826  | 497294  | 1.81%  | 0.571% |
| 63 | 16.064 | 2139 | 2141 | 2144 | rVB  | 170679  | 177856  | 0.65%  | 0.204% |
| 64 | 16.129 | 2145 | 2152 | 2156 | rBV  | 655753  | 1280314 | 4.65%  | 1.470% |
| 65 | 16.217 | 2165 | 2167 | 2174 | rVB3 | 179021  | 229002  | 0.83%  | 0.263% |
| 66 | 16.340 | 2185 | 2188 | 2192 | rBV3 | 266982  | 328547  | 1.19%  | 0.377% |
| 67 | 16.376 | 2192 | 2194 | 2198 | rVB2 | 203344  | 213329  | 0.77%  | 0.245% |
| 68 | 16.587 | 2225 | 2230 | 2232 | rBV  | 1772979 | 2525973 | 9.18%  | 2.899% |
| 69 | 16.611 | 2232 | 2234 | 2239 | rVB  | 1358397 | 1592238 | 5.78%  | 1.828% |
| 70 | 16.705 | 2246 | 2250 | 2254 | rBV  | 482436  | 711256  | 2.58%  | 0.816% |
| 71 | 16.863 | 2274 | 2277 | 2280 | rVB3 | 235664  | 284619  | 1.03%  | 0.327% |
| 72 | 17.092 | 2313 | 2316 | 2320 | rVB2 | 258687  | 295882  | 1.07%  | 0.340% |
| 73 | 17.380 | 2360 | 2365 | 2369 | rBV  | 1417157 | 1724821 | 6.27%  | 1.980% |
| 74 | 17.433 | 2370 | 2374 | 2378 | rVB  | 603256  | 807992  | 2.94%  | 0.927% |
| 75 | 17.527 | 2385 | 2390 | 2400 | rVB  | 347662  | 710072  | 2.58%  | 0.815% |
| 76 | 18.120 | 2487 | 2491 | 2497 | rVB  | 1294112 | 1657038 | 6.02%  | 1.902% |

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 Title : SVOA CALIBRATION

|     |        |      |      |      |      |          |          |         |         |
|-----|--------|------|------|------|------|----------|----------|---------|---------|
| 77  | 18.526 | 2555 | 2560 | 2563 | rBV  | 1013474  | 1424065  | 5.17%   | 1.635%  |
| 78  | 18.808 | 2604 | 2608 | 2613 | rVB  | 851139   | 990292   | 3.60%   | 1.137%  |
| 79  | 19.078 | 2651 | 2654 | 2660 | rVB4 | 374024   | 483538   | 1.76%   | 0.555%  |
| 80  | 19.424 | 2709 | 2713 | 2718 | rVB2 | 768410   | 1063666  | 3.86%   | 1.221%  |
| 81  | 19.612 | 2742 | 2745 | 2749 | rBV  | 454147   | 507462   | 1.84%   | 0.582%  |
| 82  | 19.783 | 2771 | 2774 | 2779 | rVB  | 372473   | 456472   | 1.66%   | 0.524%  |
| 83  | 19.894 | 2790 | 2793 | 2797 | rVB  | 297835   | 328442   | 1.19%   | 0.377%  |
| 84  | 20.000 | 2807 | 2811 | 2817 | rBV3 | 521580   | 664960   | 2.42%   | 0.763%  |
| 85  | 20.300 | 2859 | 2862 | 2865 | rBV2 | 195882   | 268103   | 0.97%   | 0.308%  |
| 86  | 20.529 | 2897 | 2901 | 2905 | rBV2 | 740348   | 852259   | 3.10%   | 0.978%  |
| 87  | 20.922 | 2964 | 2968 | 2974 | rVB  | 1442509  | 1843877  | 6.70%   | 2.116%  |
| 88  | 21.034 | 2983 | 2987 | 2991 | rVB  | 644642   | 717103   | 2.61%   | 0.823%  |
| 89  | 21.281 | 3026 | 3029 | 3035 | rVB  | 685972   | 928836   | 3.37%   | 1.066%  |
| 90  | 21.569 | 3074 | 3078 | 3083 | rVB  | 709855   | 939252   | 3.41%   | 1.078%  |
| 91  | 21.862 | 3117 | 3128 | 3137 | rVB  | 13264710 | 27527025 | 100.00% | 31.595% |
| 92  | 22.156 | 3170 | 3178 | 3183 | rBV2 | 773116   | 1583689  | 5.75%   | 1.818%  |
| 93  | 22.608 | 3252 | 3255 | 3259 | rVV  | 236979   | 343707   | 1.25%   | 0.394%  |
| 94  | 22.649 | 3259 | 3262 | 3272 | rVB  | 331901   | 789129   | 2.87%   | 0.906%  |
| 95  | 22.814 | 3285 | 3290 | 3296 | rVB  | 603577   | 1010891  | 3.67%   | 1.160%  |
| 96  | 23.037 | 3323 | 3328 | 3332 | rBV2 | 360820   | 745720   | 2.71%   | 0.856%  |
| 97  | 23.143 | 3340 | 3346 | 3358 | rVB  | 667445   | 1571719  | 5.71%   | 1.804%  |
| 98  | 23.572 | 3413 | 3419 | 3425 | rVB  | 434087   | 847772   | 3.08%   | 0.973%  |
| 99  | 24.071 | 3500 | 3504 | 3513 | rVB  | 292674   | 704644   | 2.56%   | 0.809%  |
| 100 | 24.459 | 3565 | 3570 | 3577 | rVB  | 296050   | 619651   | 2.25%   | 0.711%  |

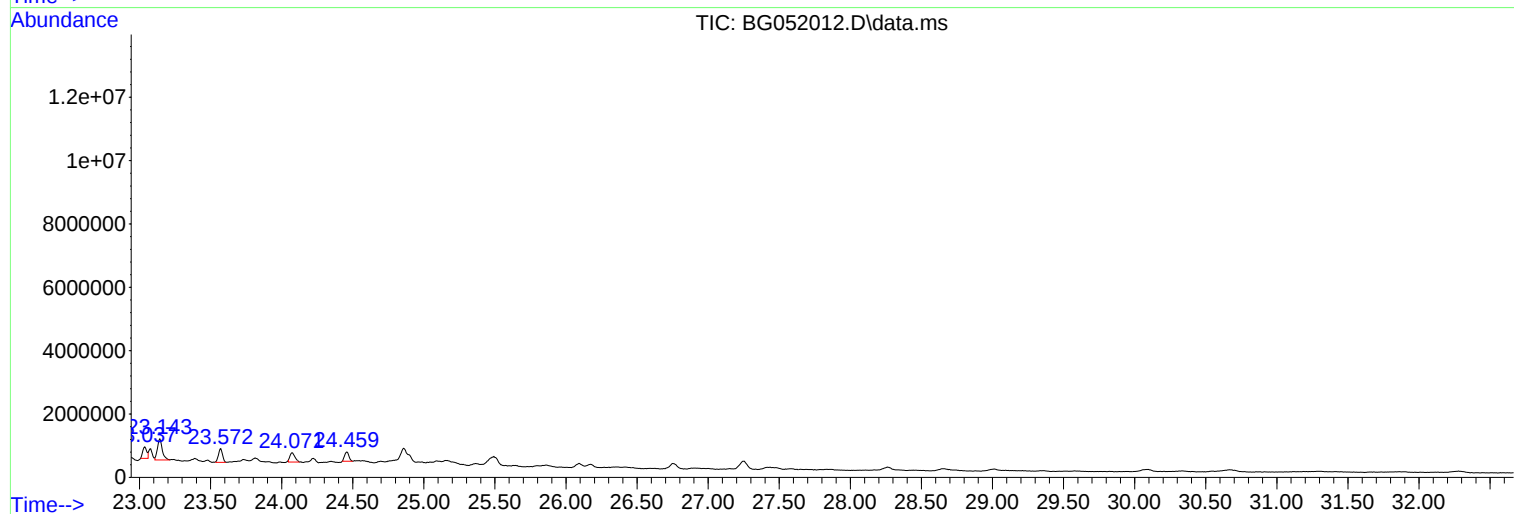
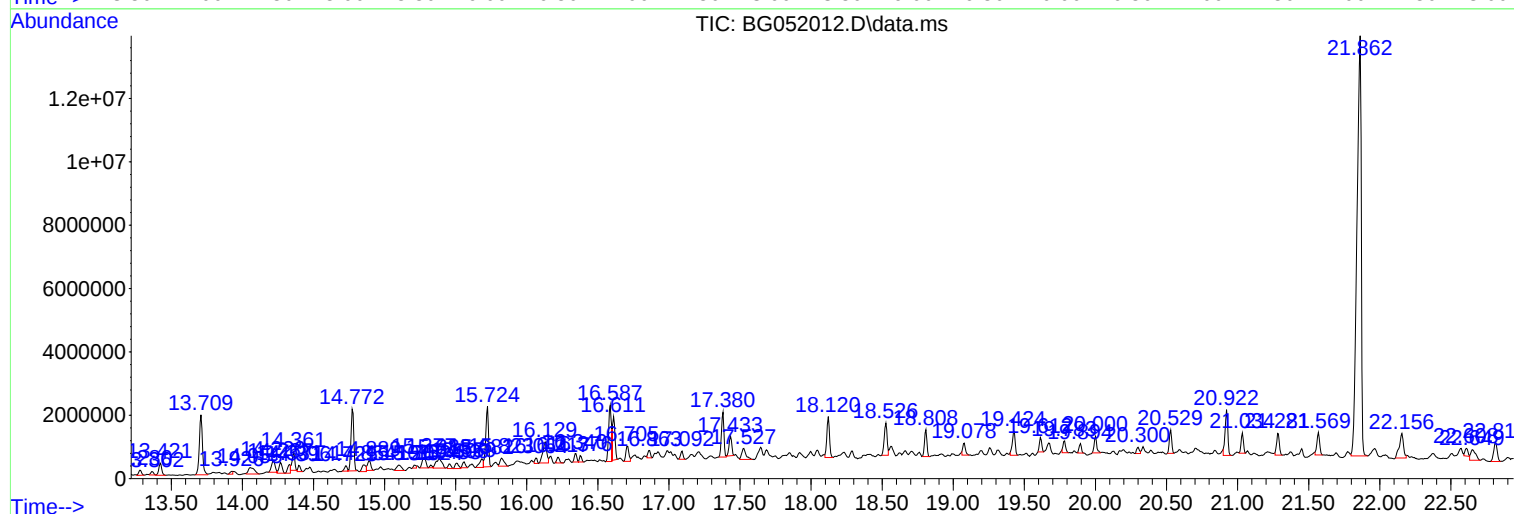
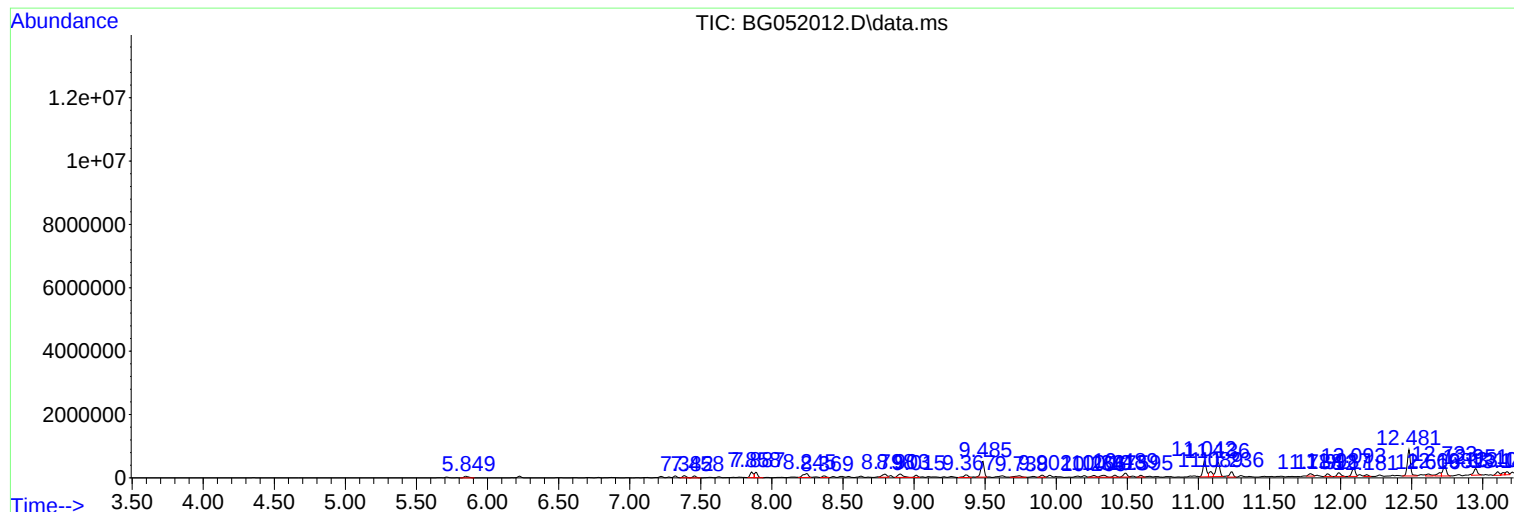
Sum of corrected areas: 87125207

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Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010622.M  
Quant Title : SVOA CALIBRATION

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



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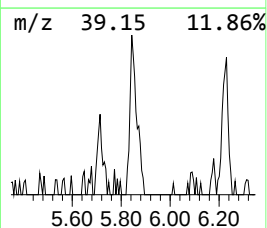
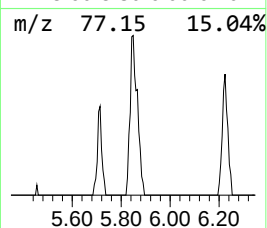
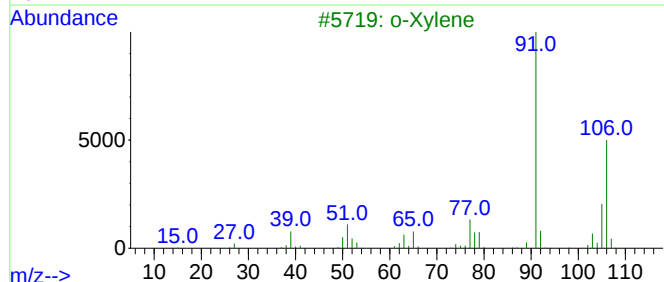
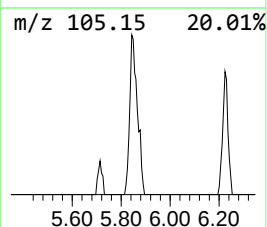
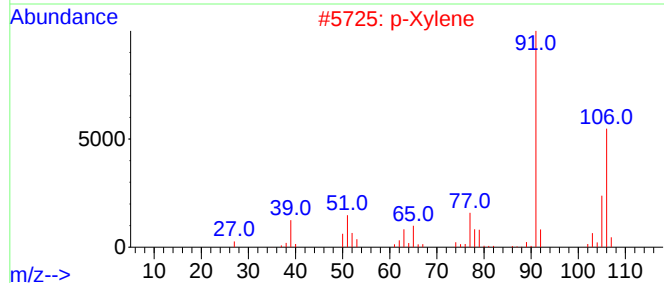
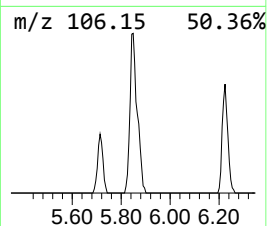
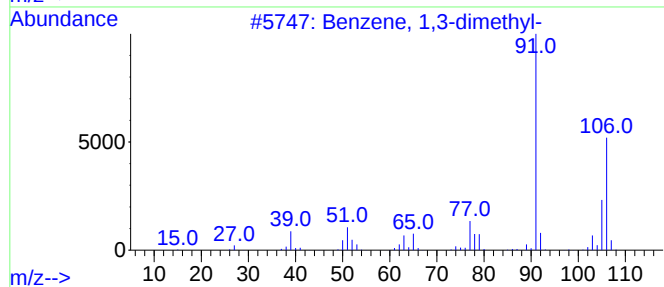
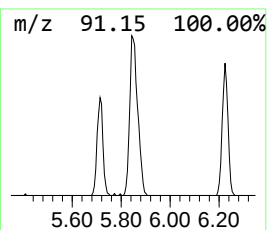
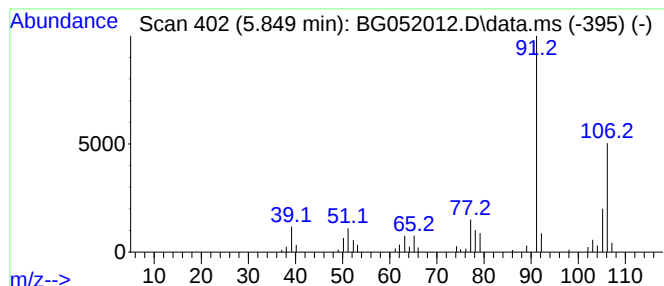
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010622.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 1 Benzene, 1,3-dimethyl- Concentration Rank 59

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 5.849 | 6.88 ng/ul | 110594 | 1,4-Dichlorobenzene-d4 | 8.245 |

| Hit# | of | 5 | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1,3-dimethyl- | 106 | C8H10   | 000108-38-3 | 97   |
| 2    |    |   | p-Xylene               | 106 | C8H10   | 000106-42-3 | 97   |
| 3    |    |   | o-Xylene               | 106 | C8H10   | 000095-47-6 | 97   |
| 4    |    |   | Benzene, 1,3-dimethyl- | 106 | C8H10   | 000108-38-3 | 95   |
| 5    |    |   | o-Xylene               | 106 | C8H10   | 000095-47-6 | 95   |



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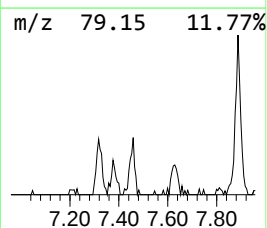
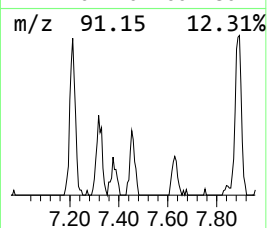
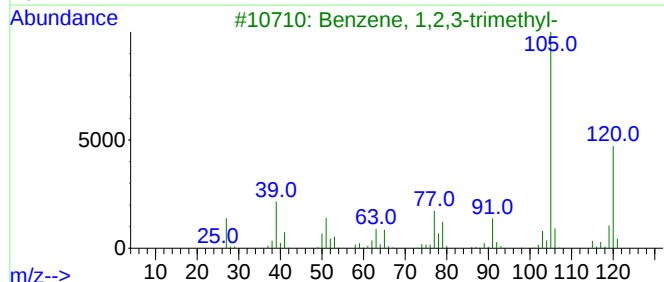
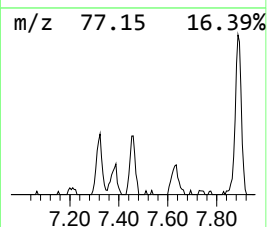
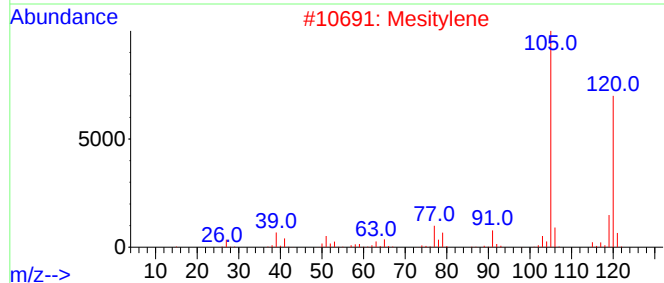
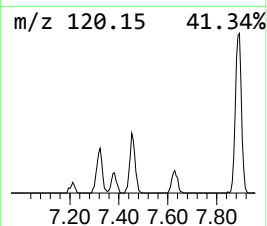
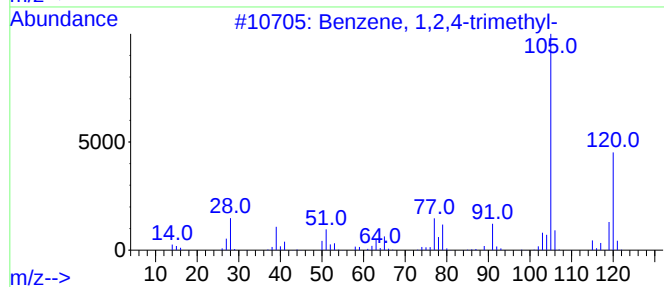
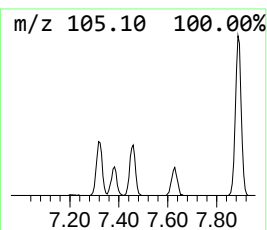
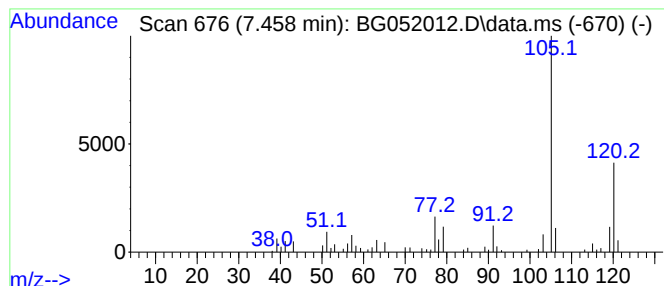
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010622.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 2 Benzene, 1,2,4-trimethyl- Concentration Rank 60

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 7.458 | 6.55 ng/ul | 105263 | 1,4-Dichlorobenzene-d4 | 8.245 |

| Hit# | of | 5 | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|------|----|---|---------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1,2,4-trimethyl- | 120 | C9H12   | 000095-63-6 | 94   |
| 2    |    |   | Mesitylene                | 120 | C9H12   | 000108-67-8 | 93   |
| 3    |    |   | Benzene, 1,2,3-trimethyl- | 120 | C9H12   | 000526-73-8 | 91   |
| 4    |    |   | Benzene, 1,2,4-trimethyl- | 120 | C9H12   | 000095-63-6 | 91   |
| 5    |    |   | Benzene, 1,2,4-trimethyl- | 120 | C9H12   | 000095-63-6 | 90   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011522\  
Data File : BG052012.D  
Acq On : 16 Jan 2022 1:50  
Operator : CG/JU  
Sample : N1132-04 10X  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLT5

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010622.M  
Quant Title : SVOA CALIBRATION

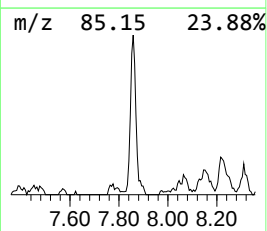
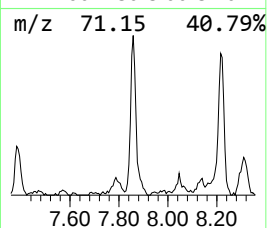
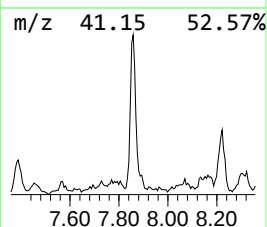
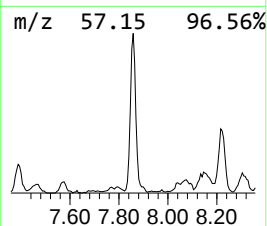
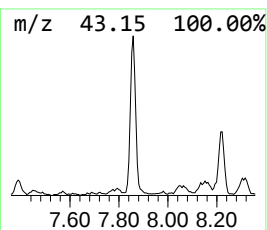
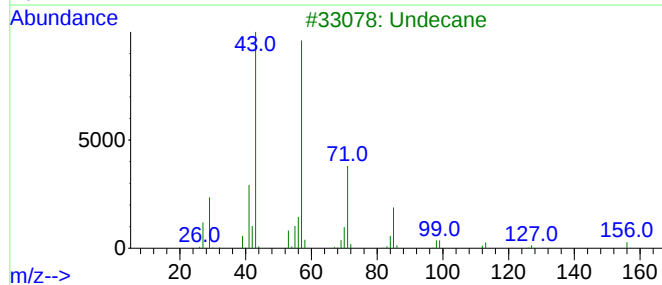
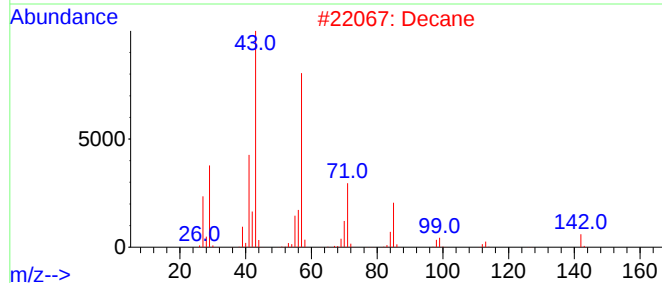
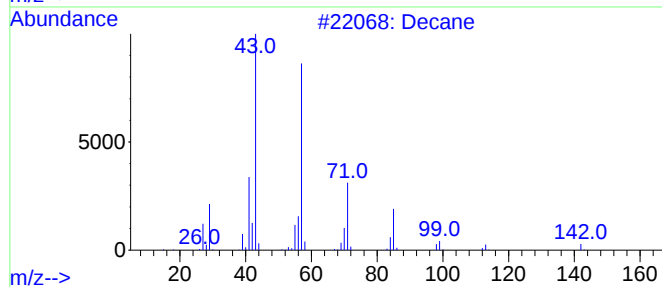
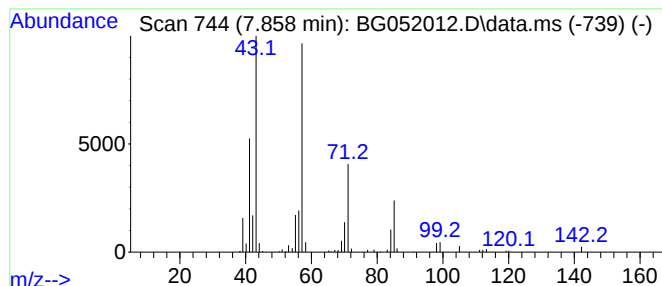
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TIC Integration Parameters: LSCINT.P

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Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 26

| R.T.  | EstConc     | Area   | Relative to ISTD       | R.T.  |
|-------|-------------|--------|------------------------|-------|
| 7.858 | 18.66 ng/ul | 299992 | 1,4-Dichlorobenzene-d4 | 8.245 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Decane                              | 142 | C10H22   | 000124-18-5  | 91   |
| 2    |    |   | Decane                              | 142 | C10H22   | 000124-18-5  | 83   |
| 3    |    |   | Undecane                            | 156 | C11H24   | 001120-21-4  | 83   |
| 4    |    |   | Nonane                              | 128 | C9H20    | 000111-84-2  | 80   |
| 5    |    |   | Carbonic acid, nonyl prop-1-en-2... | 228 | C13H24O3 | 1000382-53-9 | 78   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011522\  
Data File : BG052012.D  
Acq On : 16 Jan 2022 1:50  
Operator : CG/JU  
Sample : N1132-04 10X  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLT5

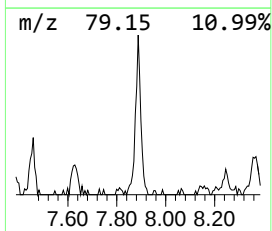
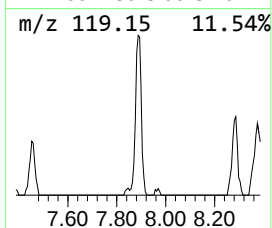
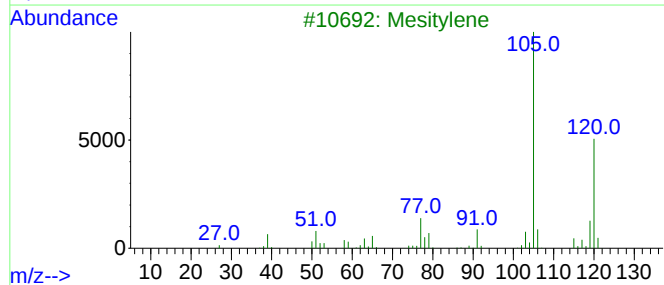
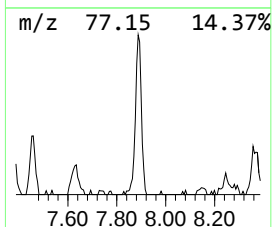
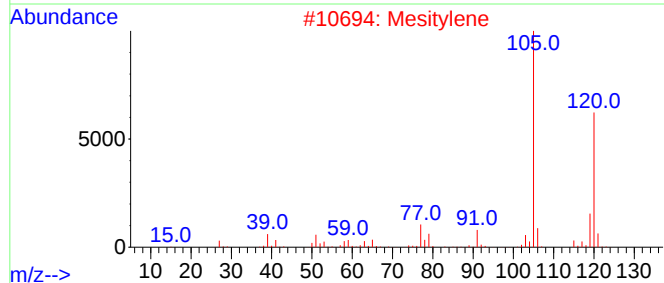
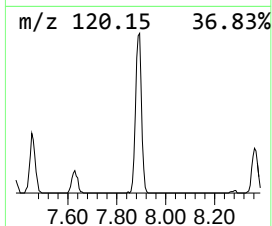
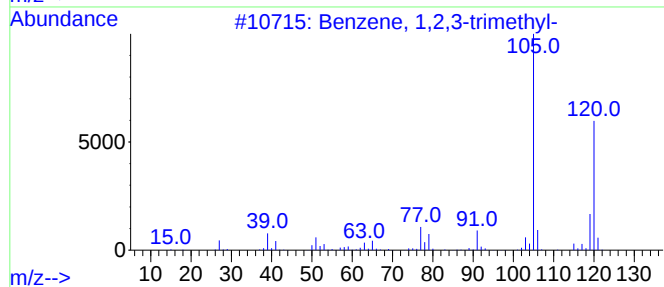
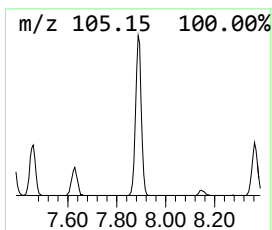
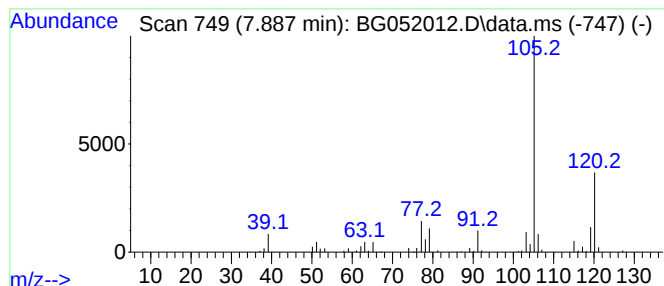
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Quant Title : SVOA CALIBRATION

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

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Peak Number 4 Benzene, 1,2,3-trimethyl- Concentration Rank 31

| R.T.  | EstConc     | Area   | Relative to ISTD       | R.T.  |
|-------|-------------|--------|------------------------|-------|
| 7.887 | 14.66 ng/ul | 235725 | 1,4-Dichlorobenzene-d4 | 8.245 |

| Hit# | of | 5 | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|------|----|---|---------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1,2,3-trimethyl- | 120 | C9H12   | 000526-73-8 | 91   |
| 2    |    |   | Mesitylene                | 120 | C9H12   | 000108-67-8 | 91   |
| 3    |    |   | Mesitylene                | 120 | C9H12   | 000108-67-8 | 91   |
| 4    |    |   | Mesitylene                | 120 | C9H12   | 000108-67-8 | 91   |
| 5    |    |   | Benzene, 1,2,4-trimethyl- | 120 | C9H12   | 000095-63-6 | 91   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011522\  
Data File : BG052012.D  
Acq On : 16 Jan 2022 1:50  
Operator : CG/JU  
Sample : N1132-04 10X  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLT5

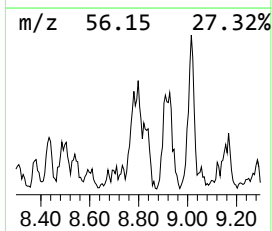
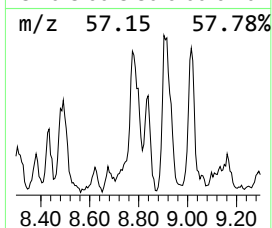
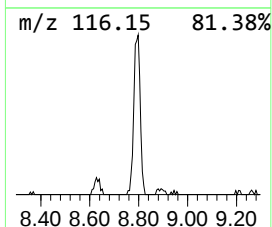
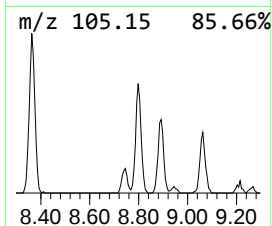
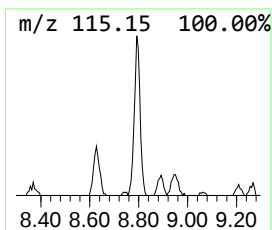
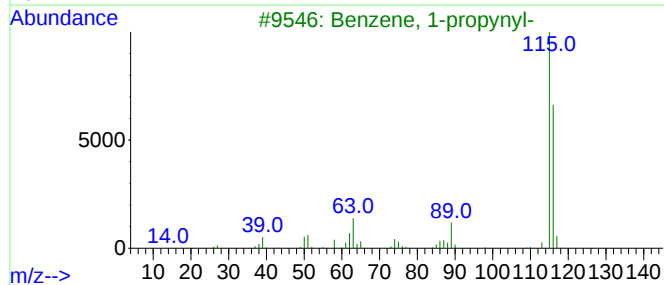
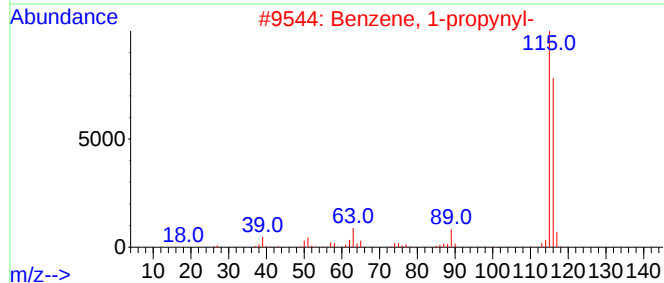
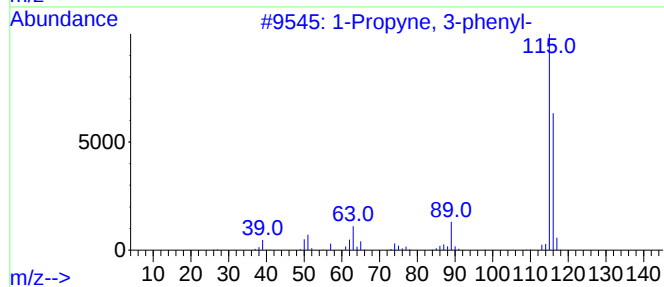
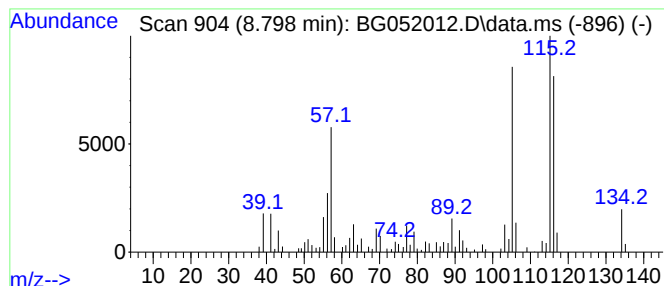
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Quant Title : SVOA CALIBRATION

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

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Peak Number 6 1-Propyne, 3-phenyl- Concentration Rank 33

| R.T.  | EstConc     | Area   | Relative to ISTD       | R.T.  |
|-------|-------------|--------|------------------------|-------|
| 8.798 | 14.19 ng/ul | 228173 | 1,4-Dichlorobenzene-d4 | 8.245 |

| Hit# | of | 5 | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------|-----|---------|-------------|------|
| 1    |    |   | 1-Propyne, 3-phenyl-    | 116 | C9H8    | 010147-11-2 | 93   |
| 2    |    |   | Benzene, 1-propynyl-    | 116 | C9H8    | 000673-32-5 | 92   |
| 3    |    |   | Benzene, 1-propynyl-    | 116 | C9H8    | 000673-32-5 | 83   |
| 4    |    |   | Benzene, 1-propynyl-    | 116 | C9H8    | 000673-32-5 | 70   |
| 5    |    |   | 3-Methylphenylacetylene | 116 | C9H8    | 000766-82-5 | 70   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011522\  
Data File : BG052012.D  
Acq On : 16 Jan 2022 1:50  
Operator : CG/JU  
Sample : N1132-04 10X  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLT5

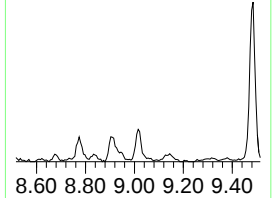
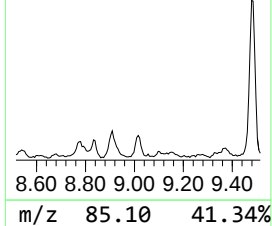
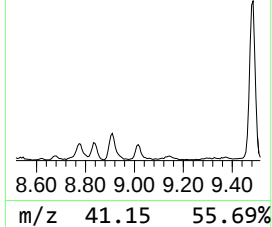
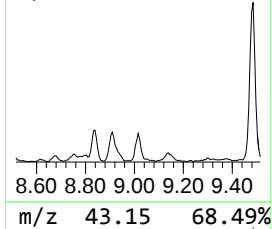
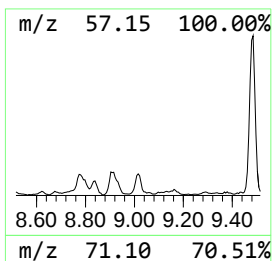
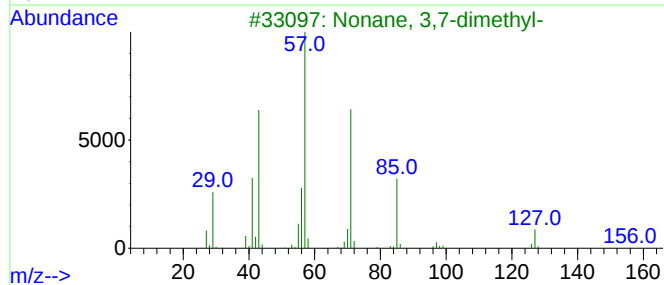
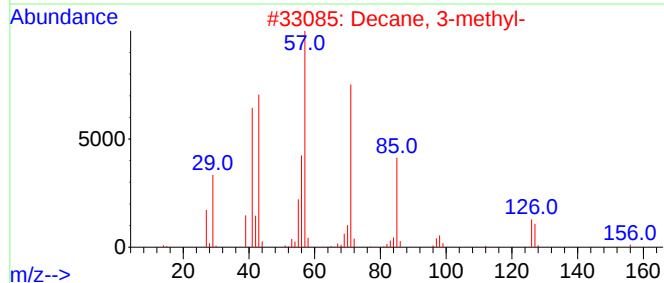
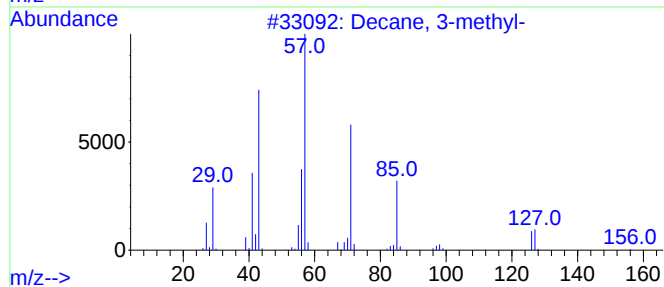
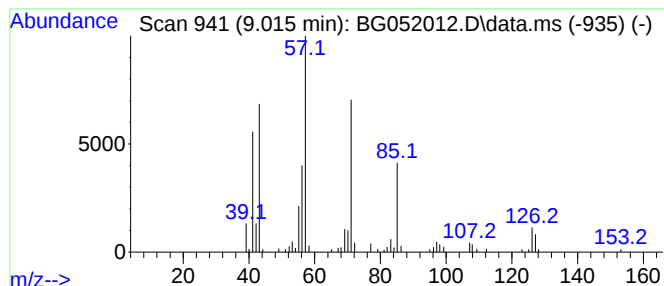
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Quant Title : SVOA CALIBRATION

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

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Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 55

| R.T.      | EstConc               | Area   | Relative to ISTD       |         |             | R.T.  |
|-----------|-----------------------|--------|------------------------|---------|-------------|-------|
| 9.015     | 7.35 ng/ul            | 118212 | 1,4-Dichlorobenzene-d4 |         |             | 8.245 |
| Hit# of 5 | Tentative ID          |        | MW                     | MolForm | CAS#        | Qual  |
| 1         | Decane, 3-methyl-     |        | 156                    | C11H24  | 013151-34-3 | 86    |
| 2         | Decane, 3-methyl-     |        | 156                    | C11H24  | 013151-34-3 | 78    |
| 3         | Nonane, 3,7-dimethyl- |        | 156                    | C11H24  | 017302-32-8 | 59    |
| 4         | Octadecane            |        | 254                    | C18H38  | 000593-45-3 | 53    |
| 5         | 2,6-Dimethyldecane    |        | 170                    | C12H26  | 013150-81-7 | 50    |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011522\  
Data File : BG052012.D  
Acq On : 16 Jan 2022 1:50  
Operator : CG/JU  
Sample : N1132-04 10X  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLT5

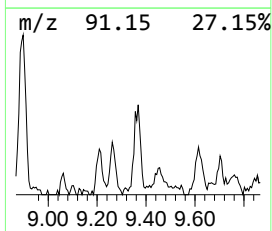
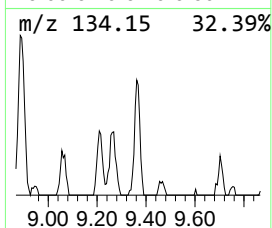
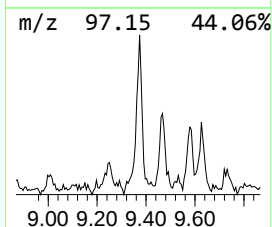
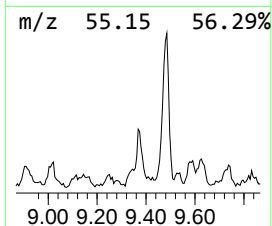
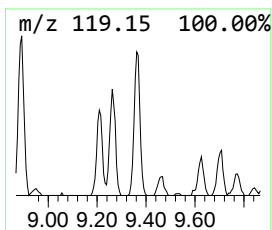
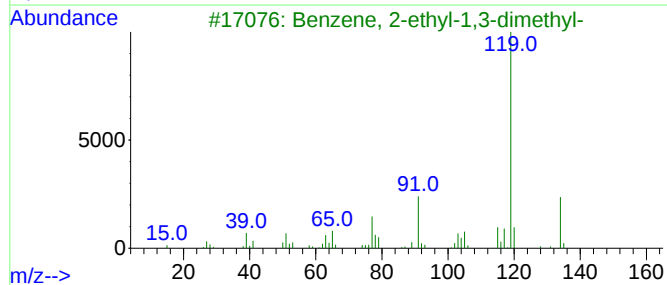
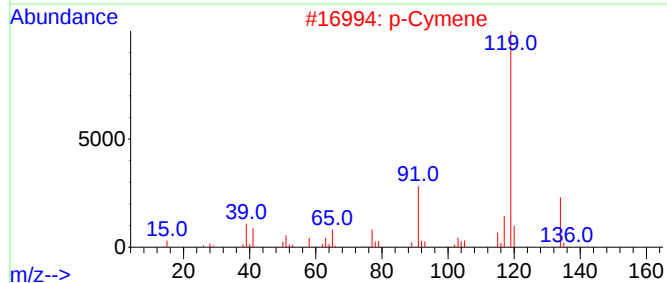
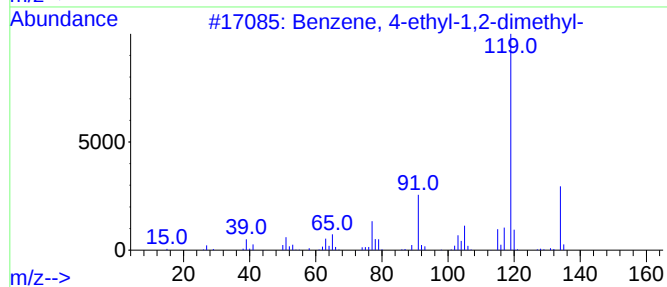
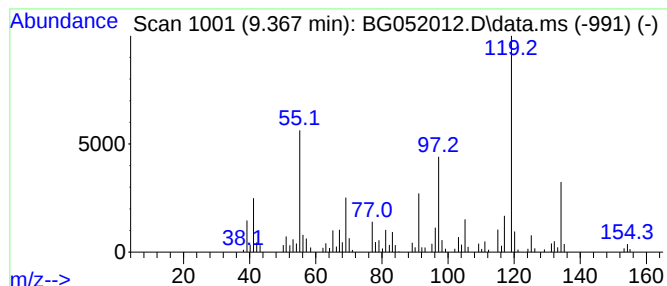
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Quant Title : SVOA CALIBRATION

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

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Peak Number 8 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 39

| R.T.  | EstConc     | Area   | Relative to ISTD       | R.T.  |
|-------|-------------|--------|------------------------|-------|
| 9.367 | 11.86 ng/ul | 190744 | 1,4-Dichlorobenzene-d4 | 8.245 |

| Hit# | of | 5 | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 4-ethyl-1,2-dimethyl- | 134 | C10H14  | 000934-80-5 | 96   |
| 2    |    |   | p-Cymene                       | 134 | C10H14  | 000099-87-6 | 92   |
| 3    |    |   | Benzene, 2-ethyl-1,3-dimethyl- | 134 | C10H14  | 002870-04-4 | 83   |
| 4    |    |   | p-Cymene                       | 134 | C10H14  | 000099-87-6 | 83   |
| 5    |    |   | o-Cymene                       | 134 | C10H14  | 000527-84-4 | 83   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011522\  
Data File : BG052012.D  
Acq On : 16 Jan 2022 1:50  
Operator : CG/JU  
Sample : N1132-04 10X  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLT5

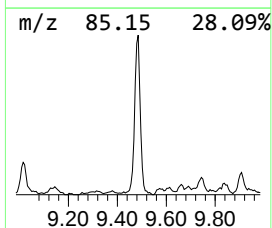
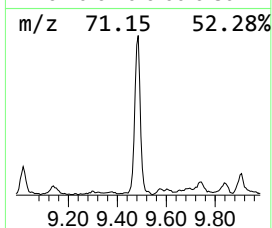
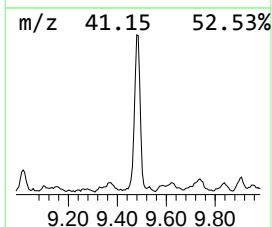
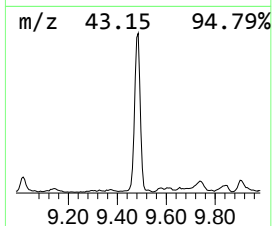
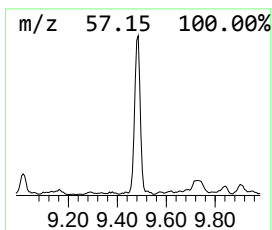
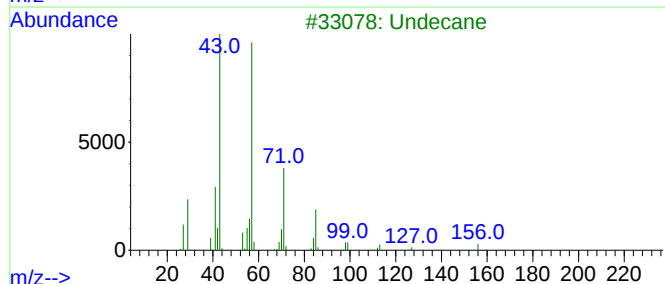
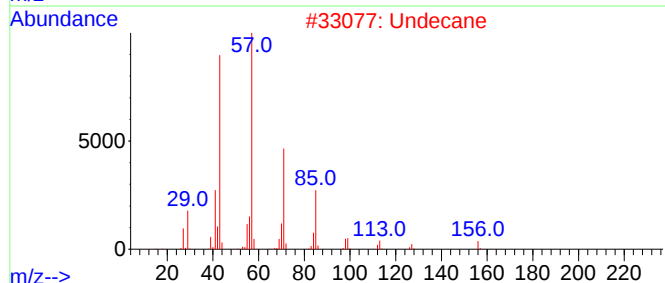
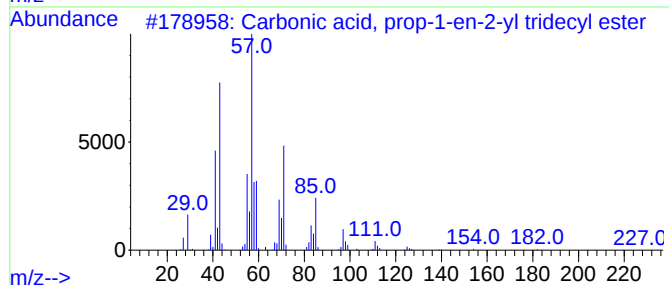
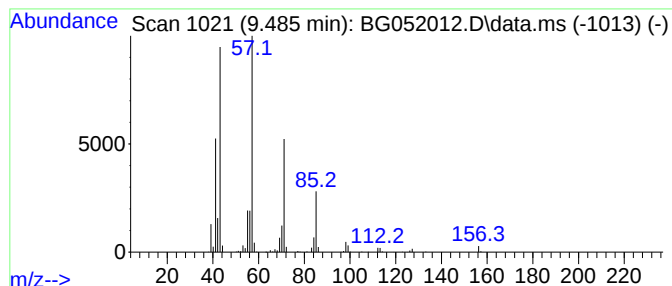
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Quant Title : SVOA CALIBRATION

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

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Peak Number 9 Carbonic acid, prop-1-en-2-... Concentration Rank 6

| R.T.  | EstConc     | Area   | Relative to ISTD       | R.T.  |
|-------|-------------|--------|------------------------|-------|
| 9.485 | 50.49 ng/ul | 811878 | 1,4-Dichlorobenzene-d4 | 8.245 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Carbonic acid, prop-1-en-2-yl tr... | 284 | C17H32O3 | 1000382-90-8 | 90   |
| 2    |    |   | Undecane                            | 156 | C11H24   | 001120-21-4  | 87   |
| 3    |    |   | Undecane                            | 156 | C11H24   | 001120-21-4  | 87   |
| 4    |    |   | Heptadecane                         | 240 | C17H36   | 000629-78-7  | 83   |
| 5    |    |   | Undecane, 4,7-dimethyl-             | 184 | C13H28   | 017301-32-5  | 64   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011522\  
Data File : BG052012.D  
Acq On : 16 Jan 2022 1:50  
Operator : CG/JU  
Sample : N1132-04 10X  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLT5

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010622.M  
Quant Title : SVOA CALIBRATION

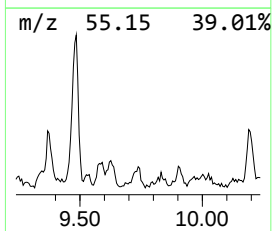
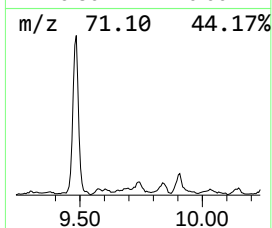
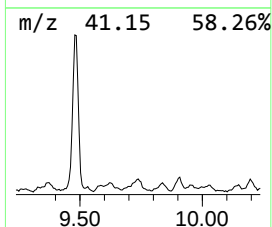
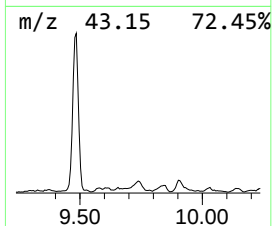
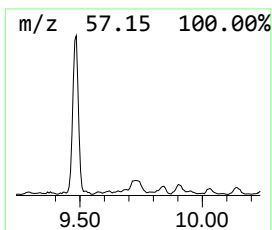
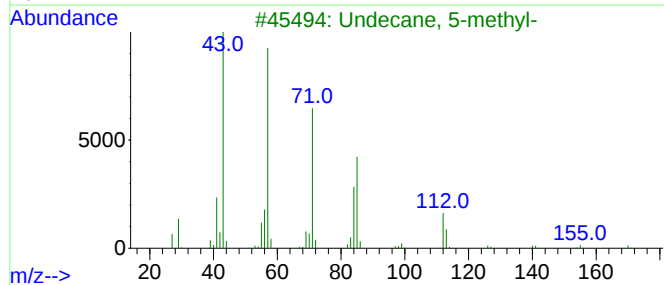
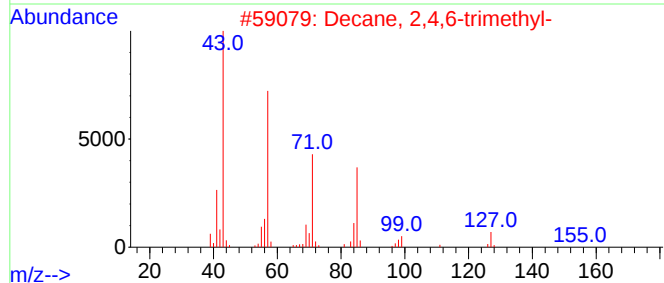
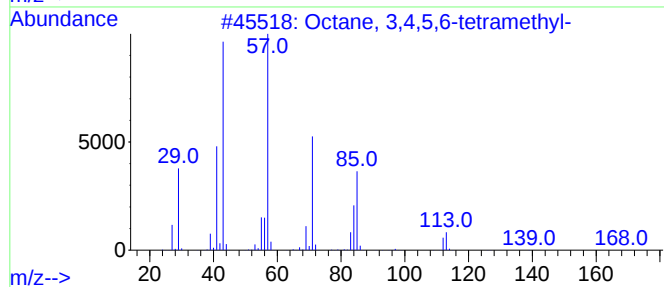
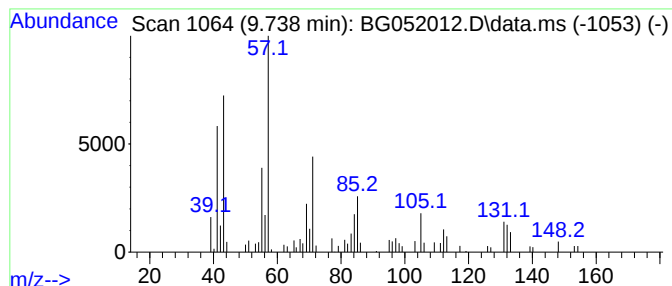
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TIC Integration Parameters: LSCINT.P

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Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 41

| R.T.  | EstConc     | Area   | Relative to ISTD | R.T.   |
|-------|-------------|--------|------------------|--------|
| 9.738 | 11.44 ng/ul | 150743 | Naphthalene-d8   | 11.089 |

| Hit# | of                           | 5 | Tentative ID | MW      | MolForm      | CAS# | Qual |
|------|------------------------------|---|--------------|---------|--------------|------|------|
| 1    | Octane, 3,4,5,6-tetramethyl- |   | 170          | C12H26  | 062185-21-1  | 50   |      |
| 2    | Decane, 2,4,6-trimethyl-     |   | 184          | C13H28  | 062108-27-4  | 47   |      |
| 3    | Undecane, 5-methyl-          |   | 170          | C12H26  | 001632-70-8  | 47   |      |
| 4    | Tridecane, 7-methyl-         |   | 198          | C14H30  | 026730-14-3  | 43   |      |
| 5    | 1-Iodo-2-methylnonane        |   | 268          | C10H21I | 1000101-47-9 | 38   |      |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011522\  
Data File : BG052012.D  
Acq On : 16 Jan 2022 1:50  
Operator : CG/JU  
Sample : N1132-04 10X  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLT5

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010622.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
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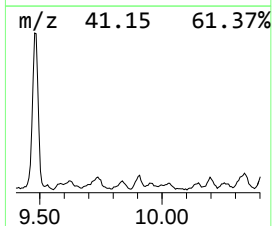
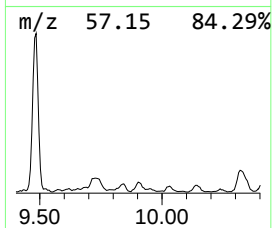
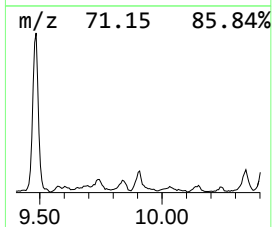
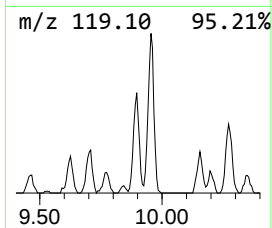
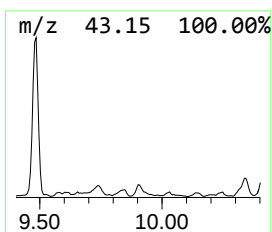
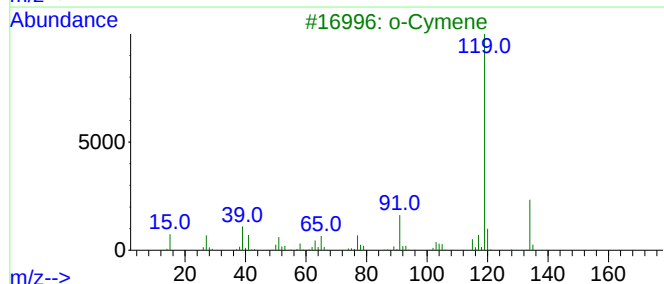
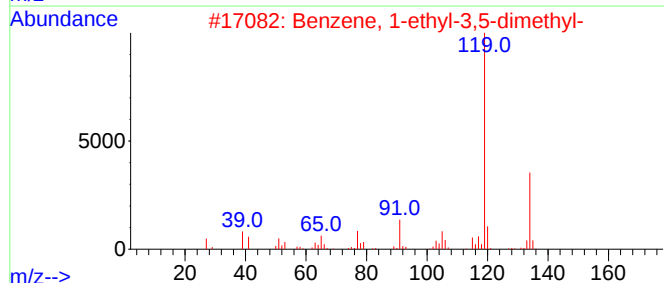
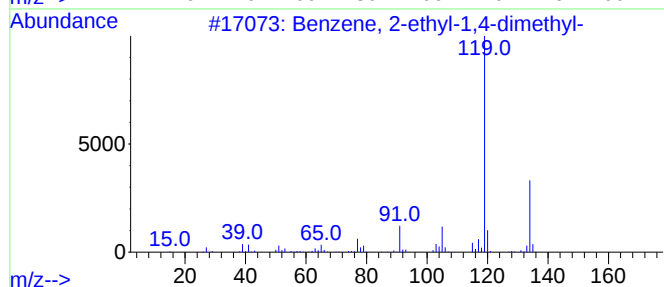
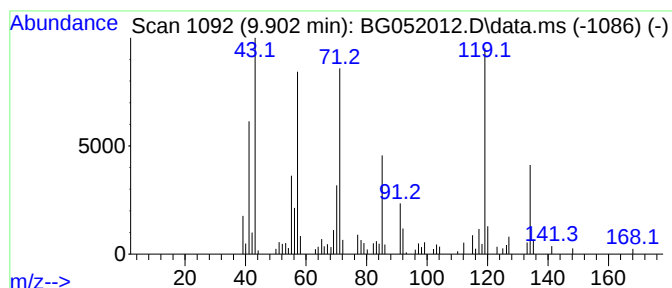
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Peak Number 11 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 49

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.   |
|-------|------------|--------|------------------|--------|
| 9.902 | 8.68 ng/ul | 114301 | Naphthalene-d8   | 11.089 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 2-ethyl-1,4-dimethyl-      | 134 | C10H14  | 001758-88-9 | 42   |
| 2    |    |   | Benzene, 1-ethyl-3,5-dimethyl-      | 134 | C10H14  | 000934-74-7 | 42   |
| 3    |    |   | o-Cymene                            | 134 | C10H14  | 000527-84-4 | 42   |
| 4    |    |   | 1,3-Cyclopentadiene, 1,2,3,4-tet... | 134 | C10H14  | 076089-59-3 | 42   |
| 5    |    |   | Benzene, 2-ethyl-1,4-dimethyl-      | 134 | C10H14  | 001758-88-9 | 42   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011522\  
Data File : BG052012.D  
Acq On : 16 Jan 2022 1:50  
Operator : CG/JU  
Sample : N1132-04 10X  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

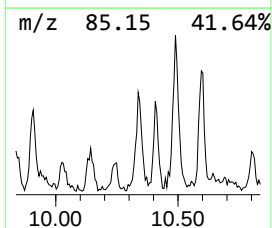
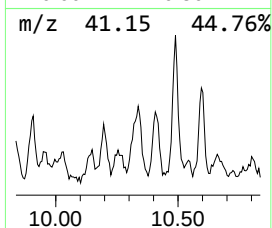
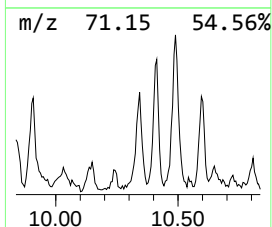
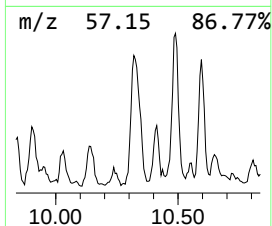
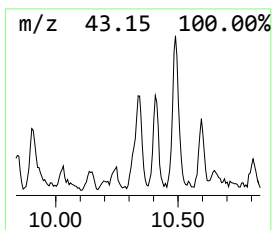
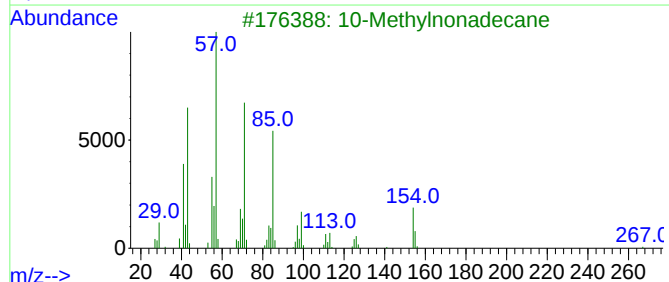
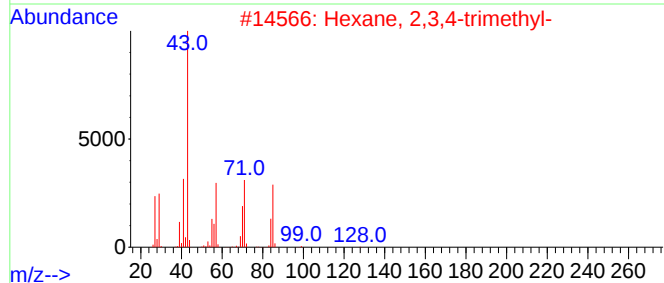
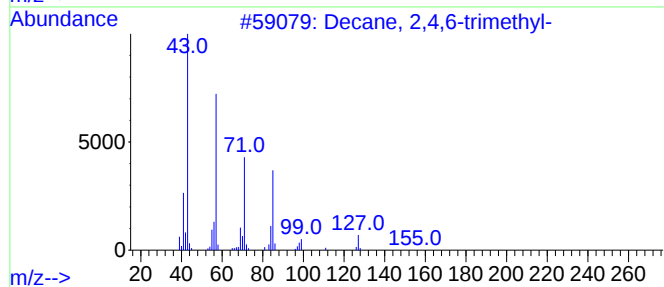
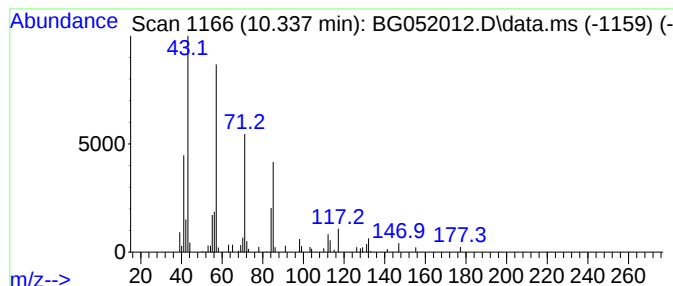
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BNA\_G  
ClientSampleId :  
BGLT5

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010622.M  
Quant Title : SVOA CALIBRATION

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

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

| R.T.      | EstConc                  | Area   | Relative to ISTD |         | R.T.        |      |
|-----------|--------------------------|--------|------------------|---------|-------------|------|
| 10.337    | 11.88 ng/ul              | 156523 | Naphthalene-d8   |         | 11.089      |      |
| Hit# of 5 | Tentative ID             |        | MW               | MolForm | CAS#        | Qual |
| 1         | Decane, 2,4,6-trimethyl- |        | 184              | C13H28  | 062108-27-4 | 53   |
| 2         | Hexane, 2,3,4-trimethyl- |        | 128              | C9H20   | 000921-47-1 | 53   |
| 3         | 10-Methylnonadecane      |        | 282              | C20H42  | 056862-62-5 | 53   |
| 4         | Pentadecane              |        | 212              | C15H32  | 000629-62-9 | 53   |
| 5         | Nonadecane               |        | 268              | C19H40  | 000629-92-5 | 53   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011522\  
Data File : BG052012.D  
Acq On : 16 Jan 2022 1:50  
Operator : CG/JU  
Sample : N1132-04 10X  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLT5

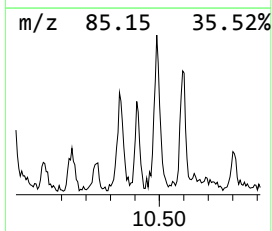
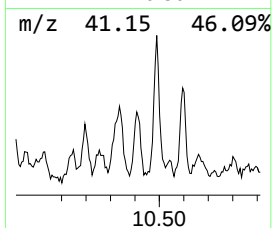
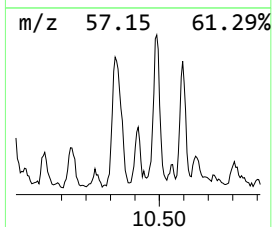
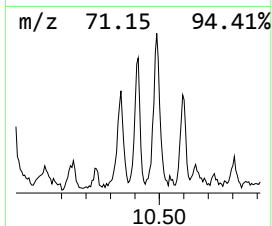
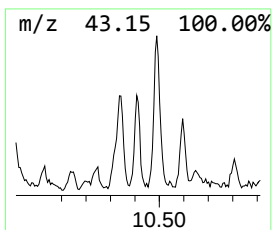
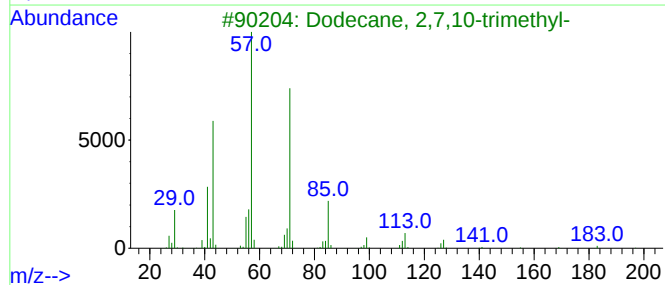
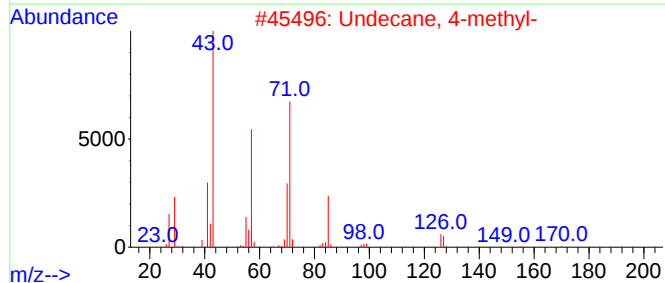
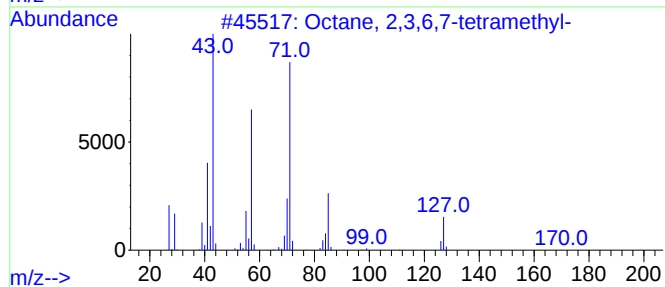
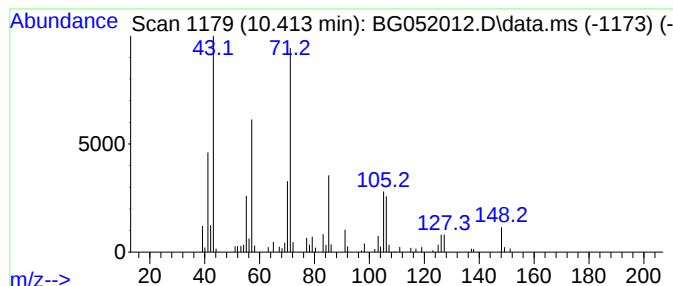
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Quant Title : SVOA CALIBRATION

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

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Peak Number 13 (DEL) Alkane: Straight-Chai... Concentration Rank 50

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 10.413 | 8.67 ng/ul | 114240 | Naphthalene-d8   | 11.089 |

| Hit# | of | 5 | Tentative ID                 | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------------------|-----|---------|-------------|------|
| 1    |    |   | Octane, 2,3,6,7-tetramethyl- | 170 | C12H26  | 052670-34-5 | 50   |
| 2    |    |   | Undecane, 4-methyl-          | 170 | C12H26  | 002980-69-0 | 50   |
| 3    |    |   | Dodecane, 2,7,10-trimethyl-  | 212 | C15H32  | 074645-98-0 | 47   |
| 4    |    |   | 1-Heptanol, 2-propyl-        | 158 | C10H22O | 010042-59-8 | 47   |
| 5    |    |   | Dodecane, 2,6,10-trimethyl-  | 212 | C15H32  | 003891-98-3 | 47   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011522\  
Data File : BG052012.D  
Acq On : 16 Jan 2022 1:50  
Operator : CG/JU  
Sample : N1132-04 10X  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLT5

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010622.M  
Quant Title : SVOA CALIBRATION

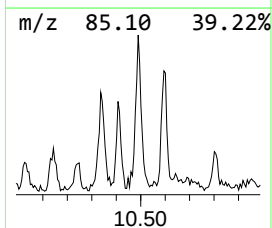
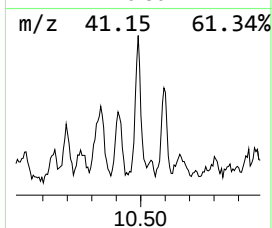
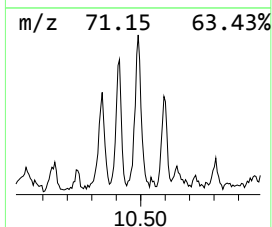
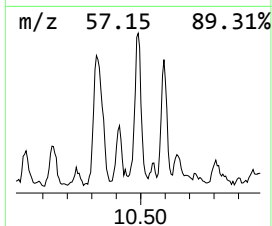
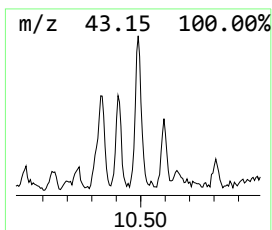
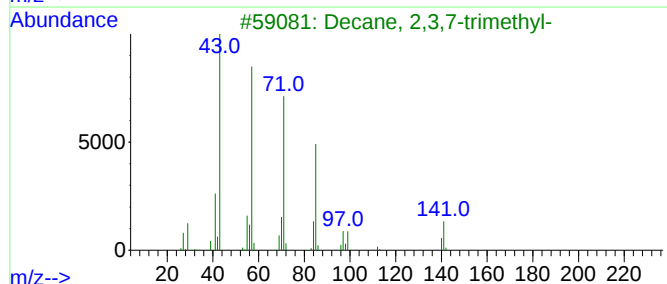
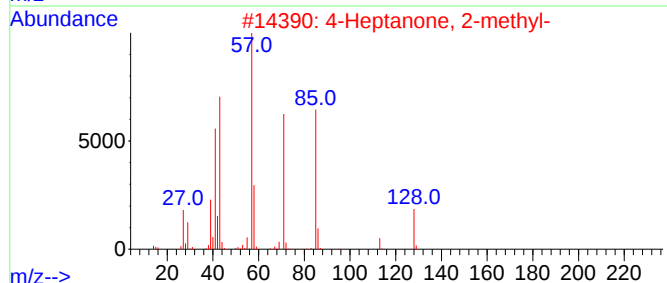
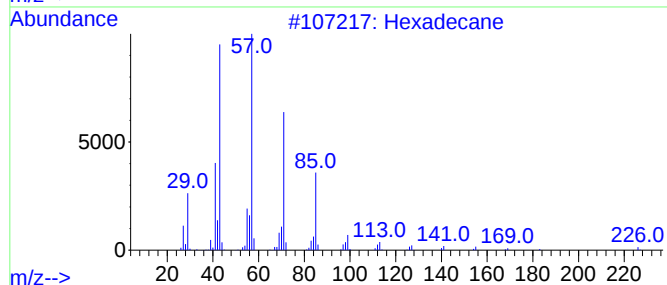
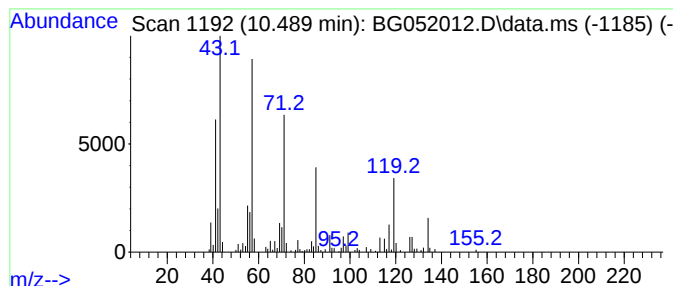
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 10.490 | 17.78 ng/ul | 234172 | Naphthalene-d8   | 11.089 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Hexadecane                          | 226 | C16H34    | 000544-76-3  | 53   |
| 2    |    |   | 4-Heptanone, 2-methyl-              | 128 | C8H16O    | 000626-33-5  | 50   |
| 3    |    |   | Decane, 2,3,7-trimethyl-            | 184 | C13H28    | 062238-13-5  | 47   |
| 4    |    |   | Sulfurous acid, butyl nonyl ester   | 264 | C13H28O3S | 1000309-17-6 | 47   |
| 5    |    |   | Heptadecane, 2,6,10,15-tetramethyl- | 296 | C21H44    | 054833-48-6  | 47   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011522\  
Data File : BG052012.D  
Acq On : 16 Jan 2022 1:50  
Operator : CG/JU  
Sample : N1132-04 10X  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLT5

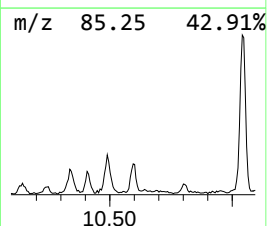
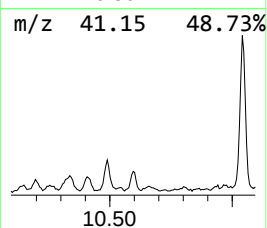
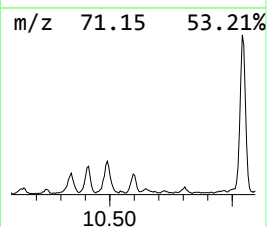
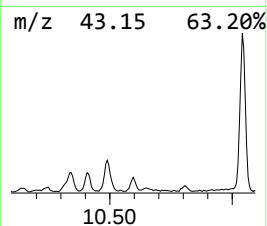
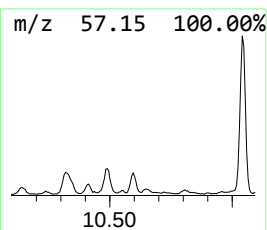
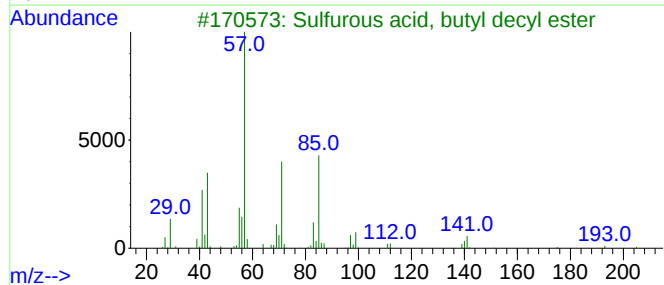
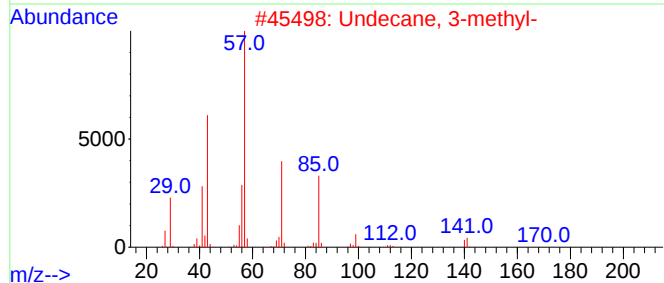
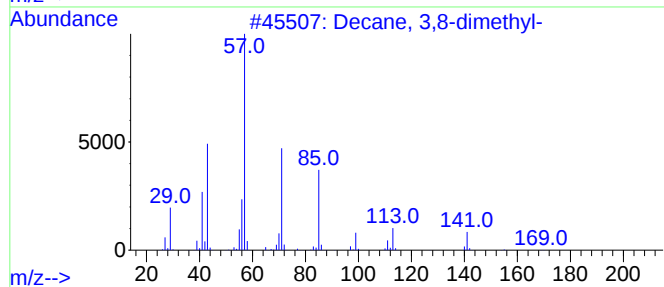
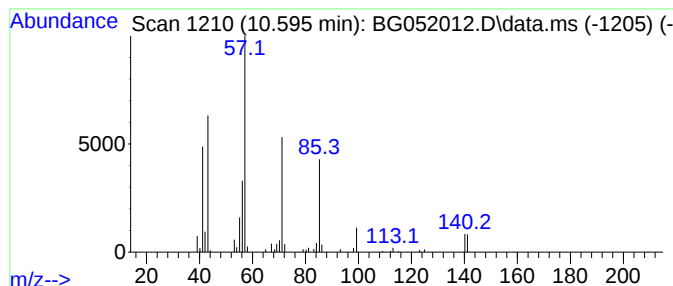
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Quant Title : SVOA CALIBRATION

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

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Peak Number 15 (DEL) Alkane: Straight-Chai... Concentration Rank 58

| R.T.   | EstConc    | Area  | Relative to ISTD | R.T.   |
|--------|------------|-------|------------------|--------|
| 10.595 | 7.17 ng/ul | 94415 | Naphthalene-d8   | 11.089 |

| Hit# | of | 5 | Tentative ID                      | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-----------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Decane, 3,8-dimethyl-             | 170 | C12H26    | 017312-55-9  | 83   |
| 2    |    |   | Undecane, 3-methyl-               | 170 | C12H26    | 001002-43-3  | 83   |
| 3    |    |   | Sulfurous acid, butyl decyl ester | 278 | C14H30O3S | 1000309-17-7 | 78   |
| 4    |    |   | Undecane, 3-methyl-               | 170 | C12H26    | 001002-43-3  | 78   |
| 5    |    |   | Dodecane, 2-methyl-               | 184 | C13H28    | 001560-97-0  | 72   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011522\  
Data File : BG052012.D  
Acq On : 16 Jan 2022 1:50  
Operator : CG/JU  
Sample : N1132-04 10X  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

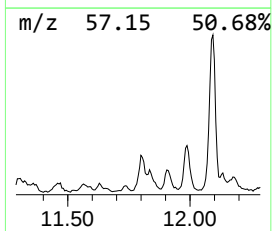
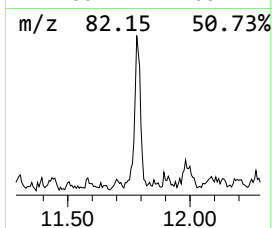
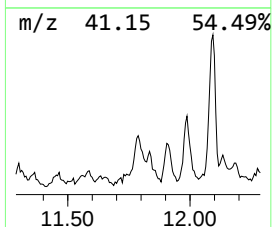
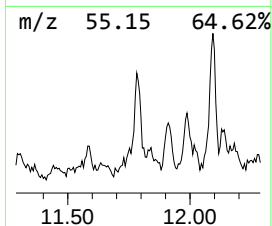
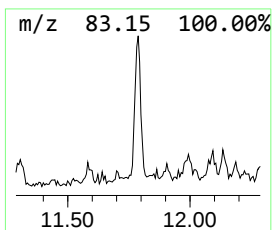
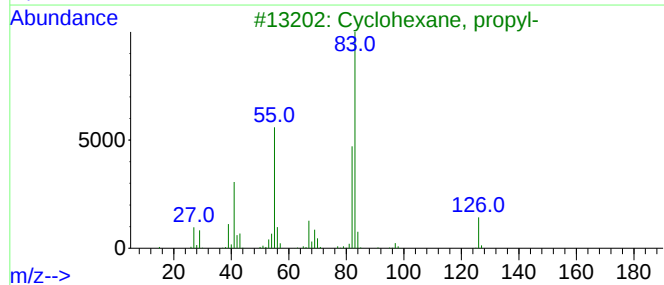
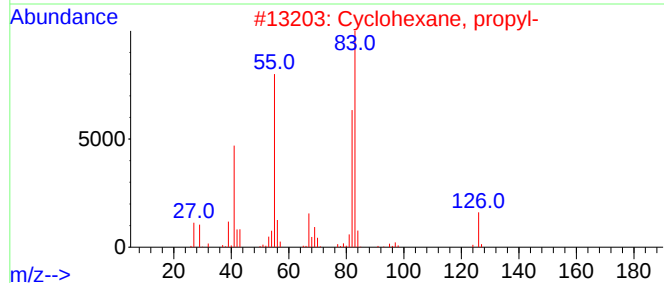
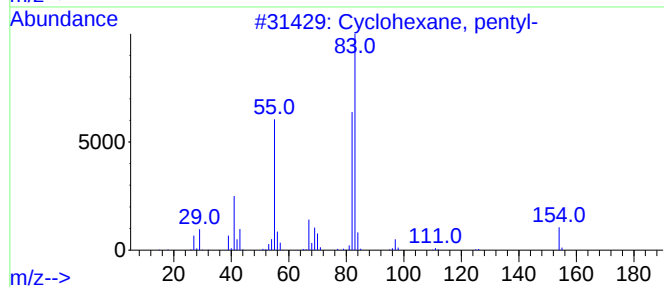
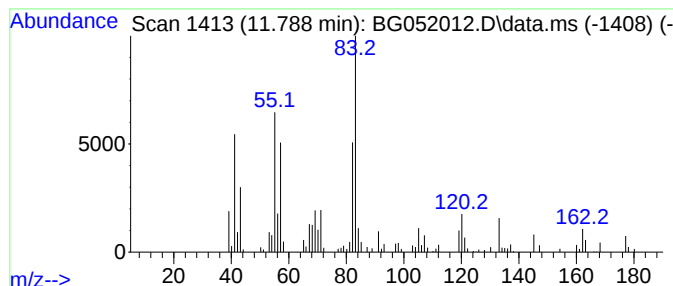
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BNA\_G  
ClientSampleId :  
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Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 16 (DEL) Alkane: Cyclic11.788 Concentration Rank 40

| R.T.      | EstConc              | Area   | Relative to ISTD | R.T.        |      |
|-----------|----------------------|--------|------------------|-------------|------|
| 11.788    | 11.54 ng/ul          | 151965 | Naphthalene-d8   | 11.089      |      |
| Hit# of 5 | Tentative ID         | MW     | MolForm          | CAS#        | Qual |
| 1         | Cyclohexane, pentyl- | 154    | C11H22           | 004292-92-6 | 64   |
| 2         | Cyclohexane, propyl- | 126    | C9H18            | 001678-92-8 | 50   |
| 3         | Cyclohexane, propyl- | 126    | C9H18            | 001678-92-8 | 50   |
| 4         | Cyclohexane, ethyl-  | 112    | C8H16            | 001678-91-7 | 50   |
| 5         | Cyclohexane, octyl-  | 196    | C14H28           | 001795-15-9 | 50   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011522\  
Data File : BG052012.D  
Acq On : 16 Jan 2022 1:50  
Operator : CG/JU  
Sample : N1132-04 10X  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLT5

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010622.M  
Quant Title : SVOA CALIBRATION

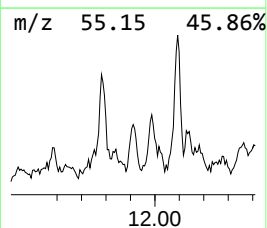
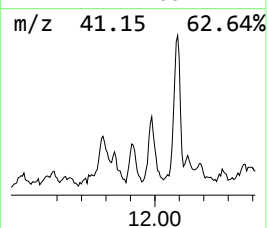
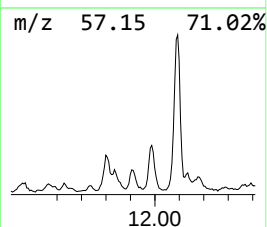
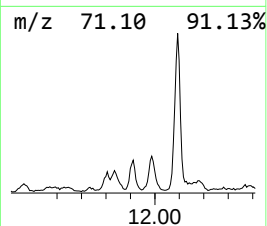
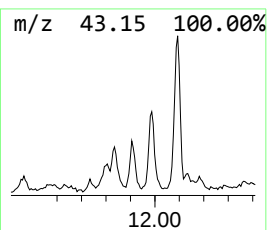
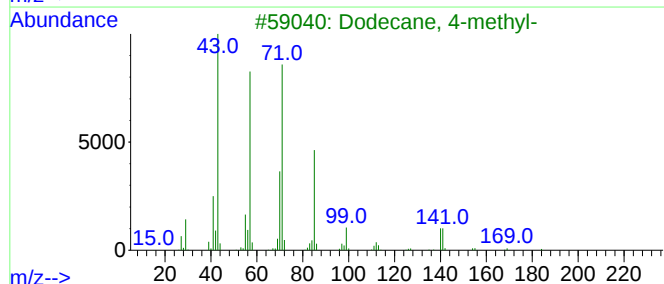
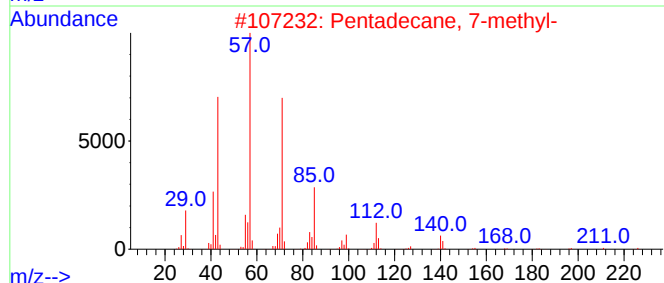
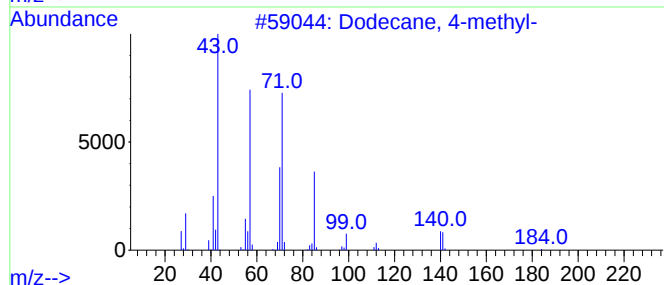
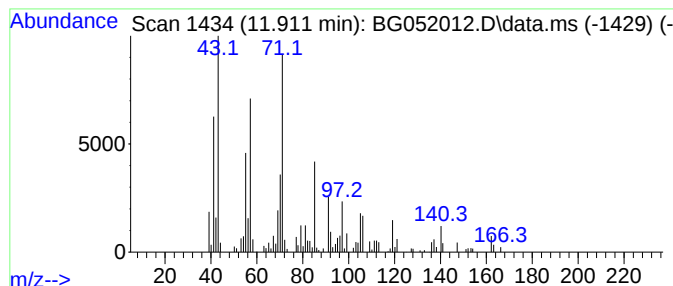
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 17 (DEL) Alkane: Straight-Chai... Concentration Rank 37

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 11.911 | 12.15 ng/ul | 160042 | Naphthalene-d8   | 11.089 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|------|-------------------------------------|-----|-----------|--------------|------|
| 1    |      | Dodecane, 4-methyl-                 | 184 | C13H28    | 006117-97-1  | 58   |
| 2    |      | Pentadecane, 7-methyl-              | 226 | C16H34    | 006165-40-8  | 53   |
| 3    |      | Dodecane, 4-methyl-                 | 184 | C13H28    | 006117-97-1  | 53   |
| 4    |      | Sulfurous acid, hexyl 2-pentyl e... | 236 | C11H24O3S | 1000309-15-6 | 50   |
| 5    |      | Undecane, 3,8-dimethyl-             | 184 | C13H28    | 017301-30-3  | 50   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011522\  
Data File : BG052012.D  
Acq On : 16 Jan 2022 1:50  
Operator : CG/JU  
Sample : N1132-04 10X  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLT5

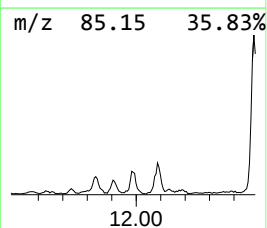
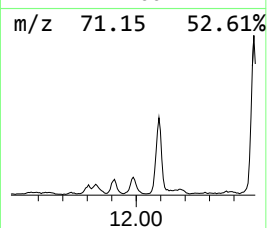
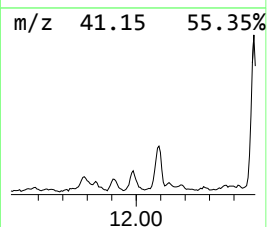
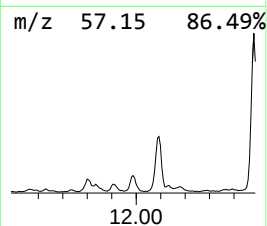
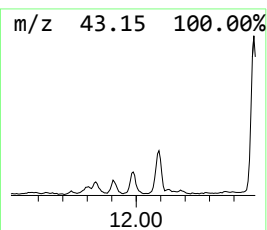
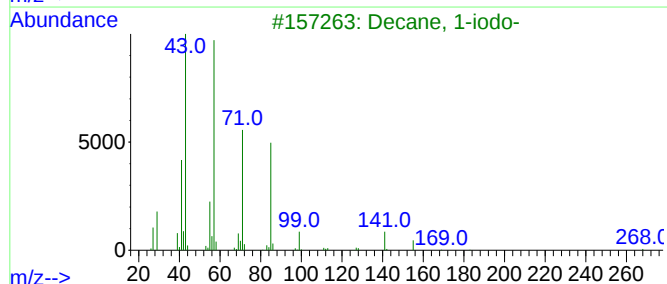
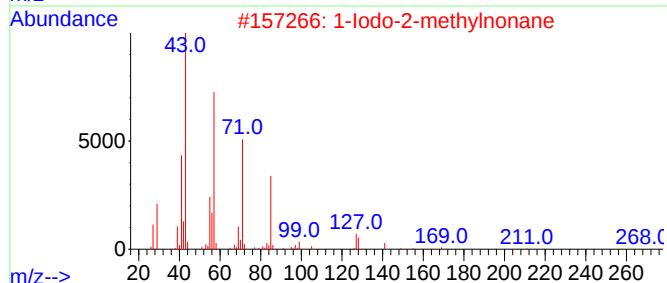
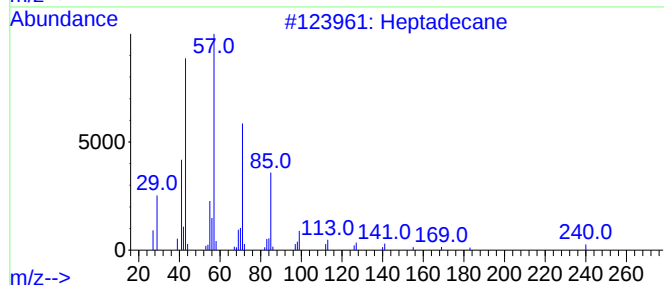
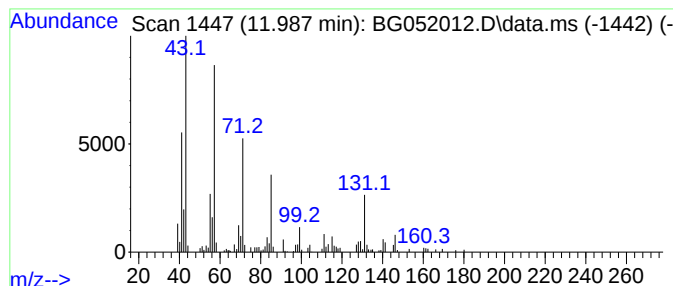
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Quant Title : SVOA CALIBRATION

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

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Peak Number 18 (DEL) Alkane: Straight-Chai... Concentration Rank 28

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 11.987 | 16.92 ng/ul | 222900 | Naphthalene-d8   | 11.089 |

| Hit# | of 5 | Tentative ID          | MW  | MolForm | CAS#         | Qual |
|------|------|-----------------------|-----|---------|--------------|------|
| 1    |      | Heptadecane           | 240 | C17H36  | 000629-78-7  | 52   |
| 2    |      | 1-Iodo-2-methylnonane | 268 | C10H21I | 1000101-47-9 | 52   |
| 3    |      | Decane, 1-iodo-       | 268 | C10H21I | 002050-77-3  | 50   |
| 4    |      | Tetradecane, 1-iodo-  | 324 | C14H29I | 019218-94-1  | 50   |
| 5    |      | Heptacosane           | 380 | C27H56  | 000593-49-7  | 49   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011522\  
Data File : BG052012.D  
Acq On : 16 Jan 2022 1:50  
Operator : CG/JU  
Sample : N1132-04 10X  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLT5

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010622.M  
Quant Title : SVOA CALIBRATION

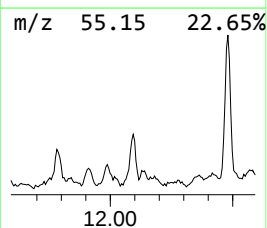
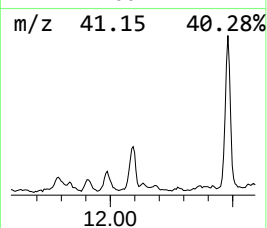
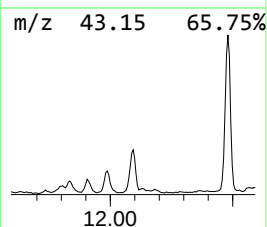
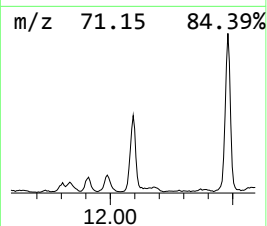
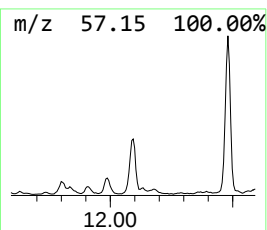
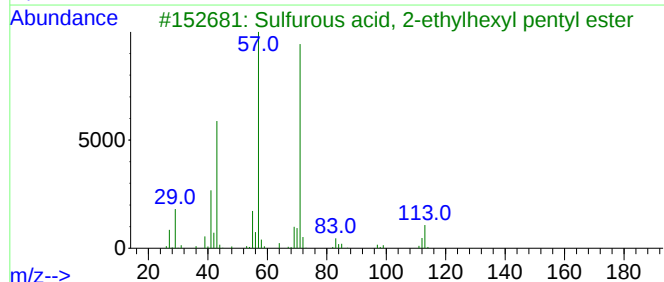
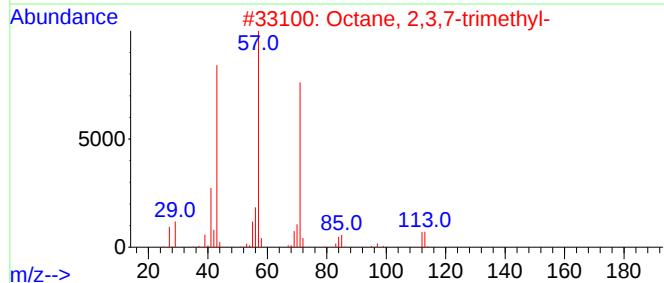
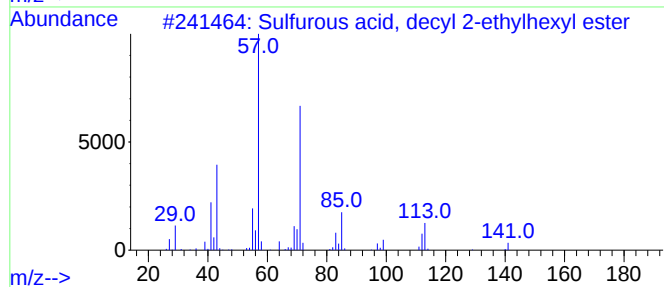
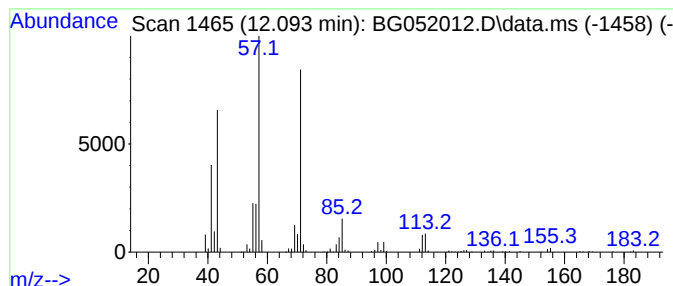
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 19 Sulfurous acid, decyl 2-eth... Concentration Rank 12

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 12.093 | 34.46 ng/ul | 453939 | Naphthalene-d8   | 11.089 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Sulfurous acid, decyl 2-ethylhex... | 334 | C18H38O3S | 1000309-19-3 | 80   |
| 2    |    |   | Octane, 2,3,7-trimethyl-            | 156 | C11H24    | 062016-34-6  | 72   |
| 3    |    |   | Sulfurous acid, 2-ethylhexyl pen... | 264 | C13H28O3S | 1000309-18-9 | 59   |
| 4    |    |   | Decane, 2,6,7-trimethyl-            | 184 | C13H28    | 062108-25-2  | 53   |
| 5    |    |   | Heptadecane, 2,6-dimethyl-          | 268 | C19H40    | 054105-67-8  | 52   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011522\  
Data File : BG052012.D  
Acq On : 16 Jan 2022 1:50  
Operator : CG/JU  
Sample : N1132-04 10X  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLT5

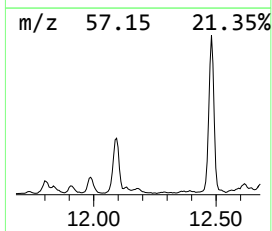
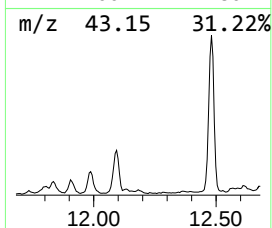
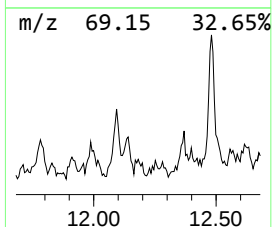
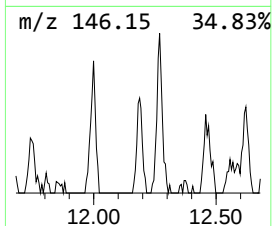
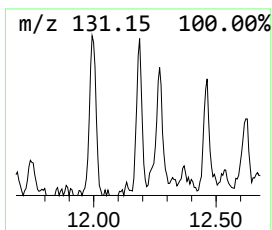
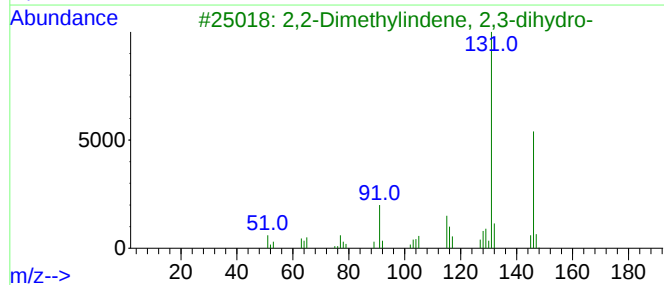
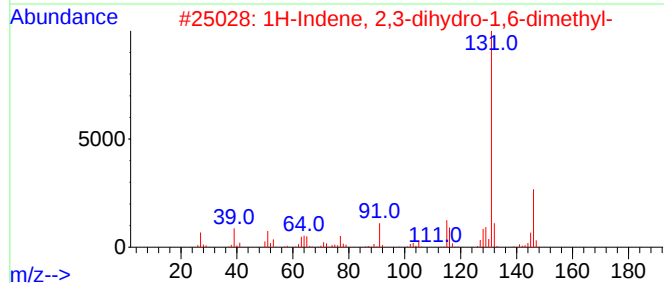
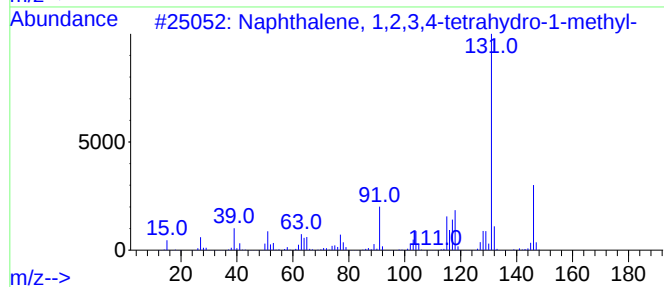
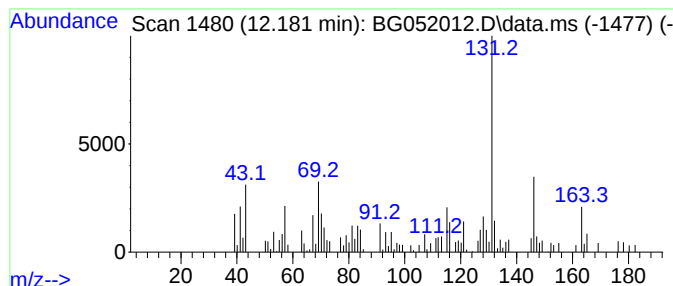
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Quant Title : SVOA CALIBRATION

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

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Peak Number 20 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 57

| R.T.   | EstConc    | Area  | Relative to ISTD | R.T.   |
|--------|------------|-------|------------------|--------|
| 12.181 | 7.17 ng/ul | 94434 | Naphthalene-d8   | 11.089 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, 1,2,3,4-tetrahydro-... | 146 | C11H14  | 001559-81-5 | 81   |
| 2    |    |   | 1H-Indene, 2,3-dihydro-1,6-dimet... | 146 | C11H14  | 017059-48-2 | 76   |
| 3    |    |   | 2,2-Dimethylindene, 2,3-dihydro-    | 146 | C11H14  | 020836-11-7 | 74   |
| 4    |    |   | 1H-Indene, 2,3-dihydro-4,7-dimet... | 146 | C11H14  | 006682-71-9 | 72   |
| 5    |    |   | 1H-Indene, 2,3-dihydro-4,7-dimet... | 146 | C11H14  | 006682-71-9 | 70   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011522\  
Data File : BG052012.D  
Acq On : 16 Jan 2022 1:50  
Operator : CG/JU  
Sample : N1132-04 10X  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLT5

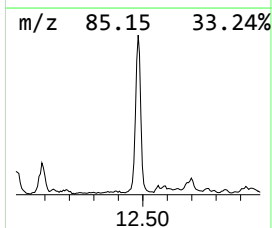
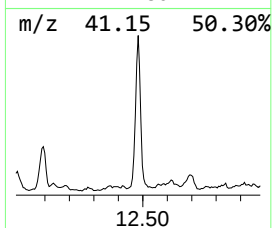
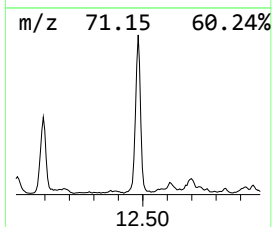
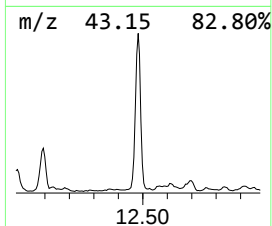
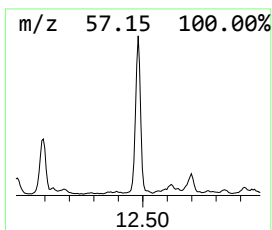
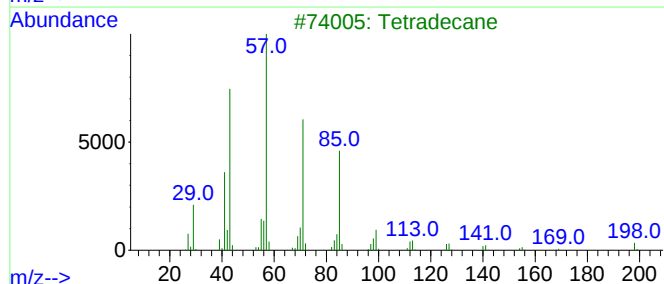
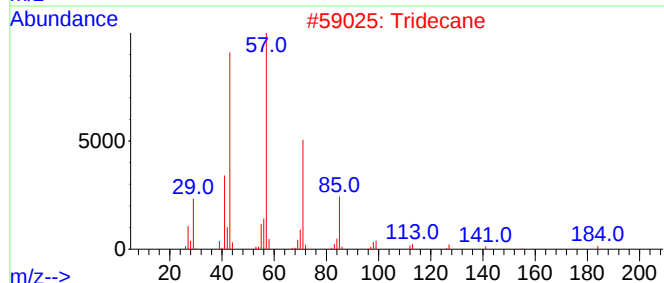
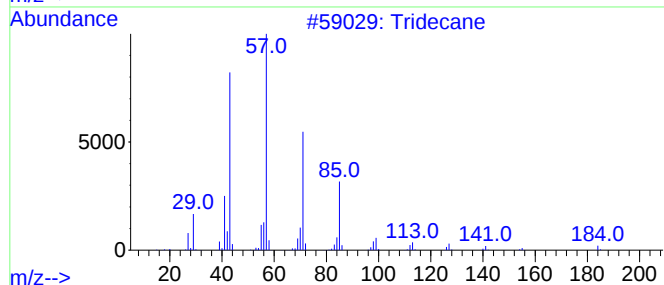
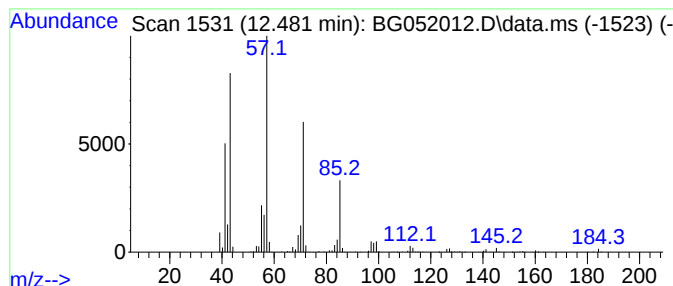
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Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 21 (DEL) Alkane: Straight-Chai... Concentration Rank 3

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 12.481 | 97.84 ng/ul | 1288700 | Naphthalene-d8   | 11.089 |

| Hit# | of 5 | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|------|-------------------------|-----|---------|-------------|------|
| 1    |      | Tridecane               | 184 | C13H28  | 000629-50-5 | 91   |
| 2    |      | Tridecane               | 184 | C13H28  | 000629-50-5 | 90   |
| 3    |      | Tetradecane             | 198 | C14H30  | 000629-59-4 | 86   |
| 4    |      | Hexadecane              | 226 | C16H34  | 000544-76-3 | 86   |
| 5    |      | Undecane, 4,6-dimethyl- | 184 | C13H28  | 017312-82-2 | 80   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011522\  
Data File : BG052012.D  
Acq On : 16 Jan 2022 1:50  
Operator : CG/JU  
Sample : N1132-04 10X  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLT5

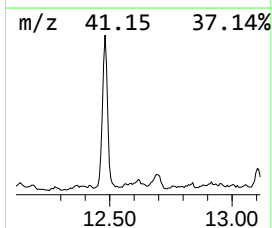
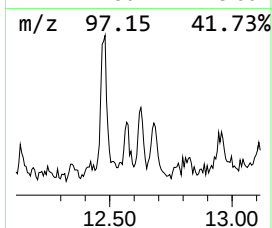
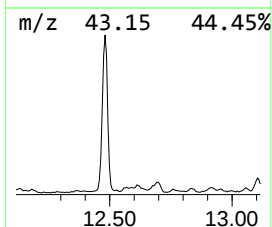
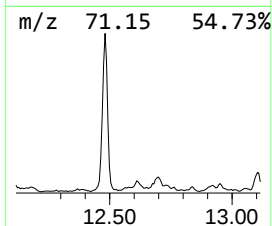
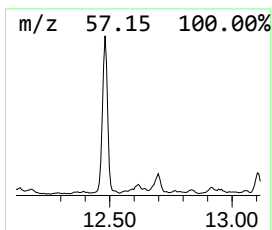
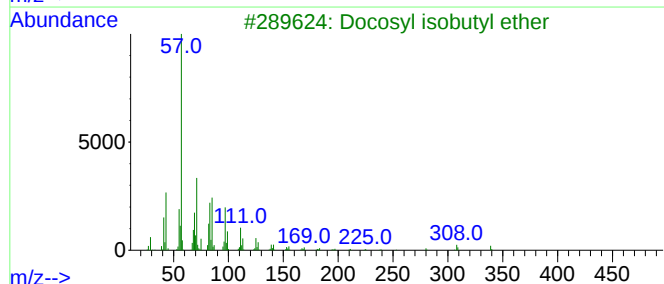
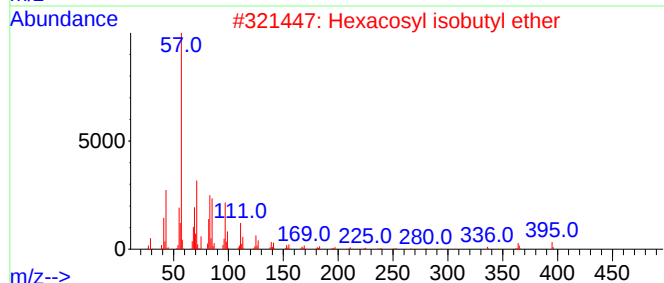
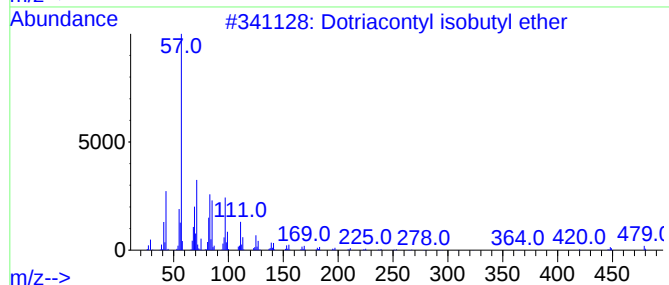
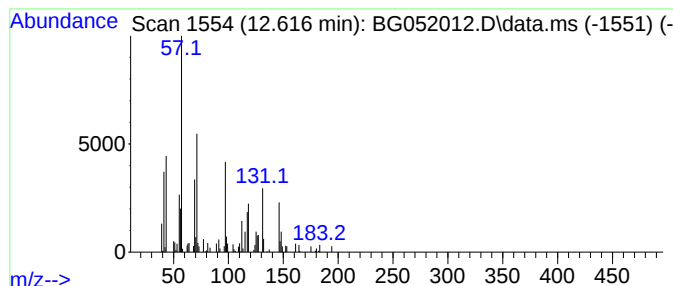
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Quant Title : SVOA CALIBRATION

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

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Peak Number 22 unknown-01 Concentration Rank 43

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 12.616 | 10.52 ng/ul | 138561 | Naphthalene-d8   | 11.089 |

| Hit# | of | 5 | Tentative ID                | MW  | MolForm | CAS#         | Qual |
|------|----|---|-----------------------------|-----|---------|--------------|------|
| 1    |    |   | Dotriacontyl isobutyl ether | 523 | C36H74O | 1010406-33-6 | 25   |
| 2    |    |   | Hexacosyl isobutyl ether    | 438 | C30H62O | 1000406-33-3 | 25   |
| 3    |    |   | Docosyl isobutyl ether      | 382 | C26H54O | 1000406-33-1 | 25   |
| 4    |    |   | Isobutyl octacosyl ether    | 467 | C32H66O | 1000406-33-4 | 25   |
| 5    |    |   | Decane, 2,6,7-trimethyl-    | 184 | C13H28  | 062108-25-2  | 22   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011522\  
Data File : BG052012.D  
Acq On : 16 Jan 2022 1:50  
Operator : CG/JU  
Sample : N1132-04 10X  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLT5

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010622.M  
Quant Title : SVOA CALIBRATION

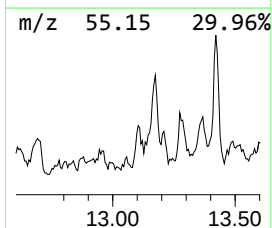
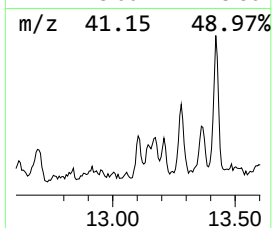
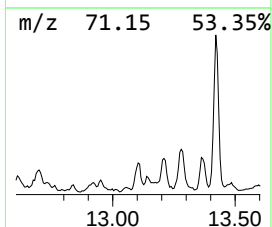
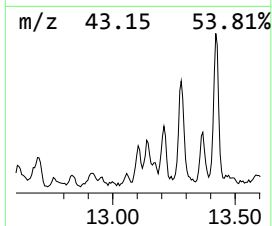
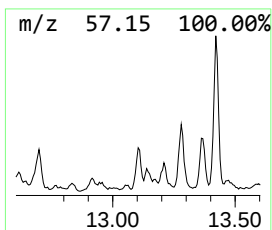
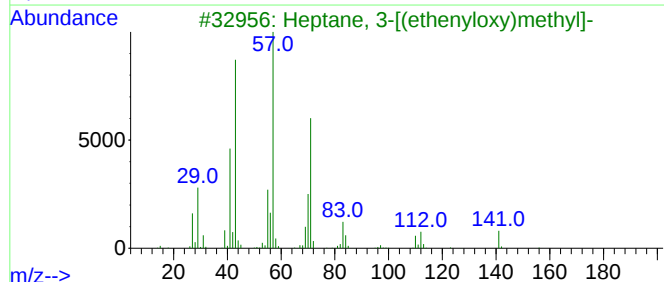
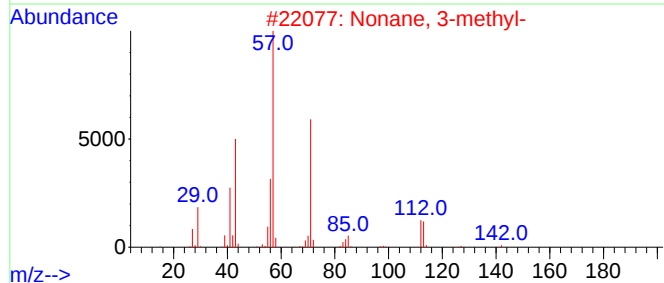
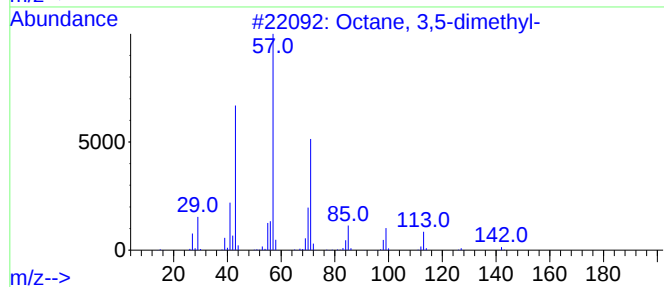
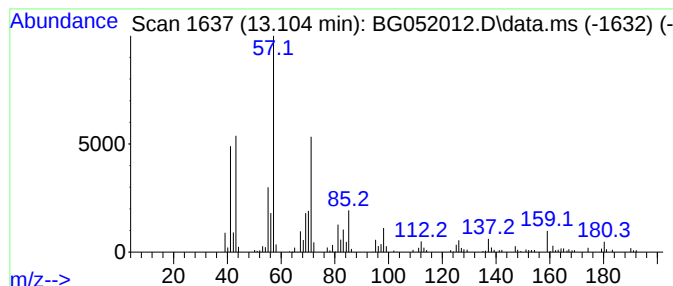
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TIC Integration Parameters: LSCINT.P

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Peak Number 23 (DEL) Alkane: Straight-Chai... Concentration Rank 52

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 13.104 | 8.04 ng/ul | 195742 | Acenaphthene-d10 | 14.895 |

| Hit# | of | 5 | Tentative ID                     | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------------|-----|---------|-------------|------|
| 1    |    |   | Octane, 3,5-dimethyl-            | 142 | C10H22  | 015869-93-9 | 52   |
| 2    |    |   | Nonane, 3-methyl-                | 142 | C10H22  | 005911-04-6 | 43   |
| 3    |    |   | Heptane, 3-[(ethenyloxy)methyl]- | 156 | C10H20O | 000103-44-6 | 43   |
| 4    |    |   | Tridecane                        | 184 | C13H28  | 000629-50-5 | 43   |
| 5    |    |   | Heptane, 2,6-dimethyl-           | 128 | C9H20   | 001072-05-5 | 38   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011522\  
Data File : BG052012.D  
Acq On : 16 Jan 2022 1:50  
Operator : CG/JU  
Sample : N1132-04 10X  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLT5

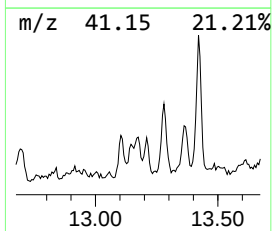
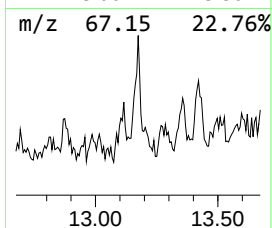
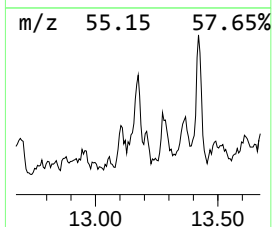
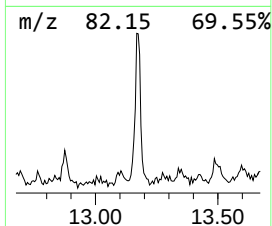
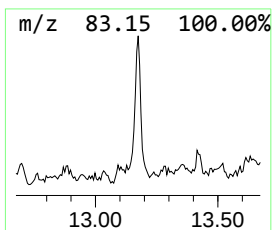
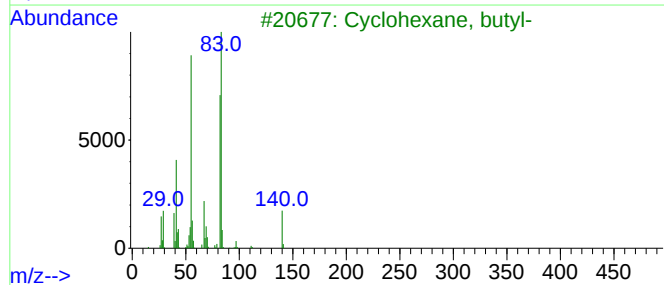
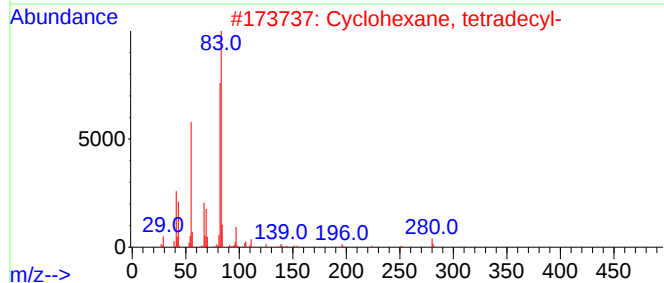
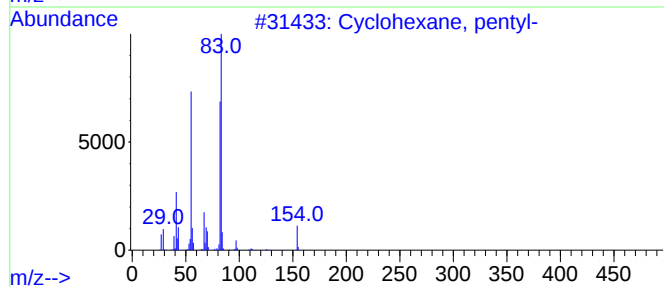
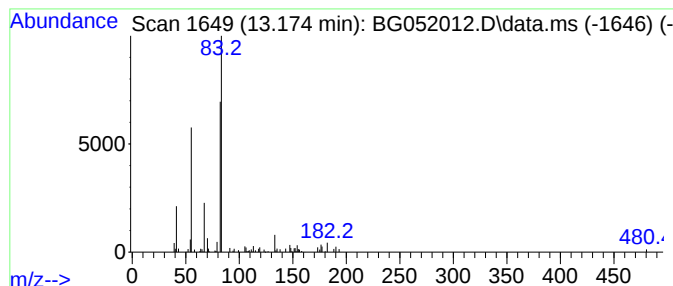
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Quant Title : SVOA CALIBRATION

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

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Peak Number 25 (DEL) Alkane: Cyclic13.174 Concentration Rank 63

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 13.174 | 5.83 ng/ul | 141997 | Acenaphthene-d10 | 14.895 |

| Hit# | of | 5 | Tentative ID             | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------|-----|---------|-------------|------|
| 1    |    |   | Cyclohexane, pentyl-     | 154 | C11H22  | 004292-92-6 | 72   |
| 2    |    |   | Cyclohexane, tetradecyl- | 280 | C20H40  | 001795-18-2 | 72   |
| 3    |    |   | Cyclohexane, butyl-      | 140 | C10H20  | 001678-93-9 | 72   |
| 4    |    |   | Cyclohexane, octyl-      | 196 | C14H28  | 001795-15-9 | 72   |
| 5    |    |   | Cyclohexane, pentyl-     | 154 | C11H22  | 004292-92-6 | 64   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011522\  
Data File : BG052012.D  
Acq On : 16 Jan 2022 1:50  
Operator : CG/JU  
Sample : N1132-04 10X  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLT5

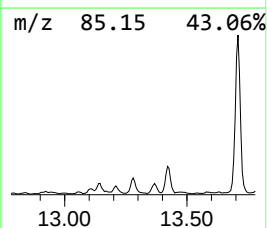
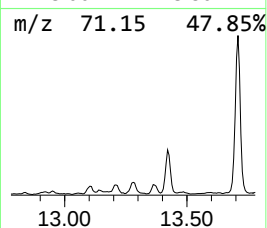
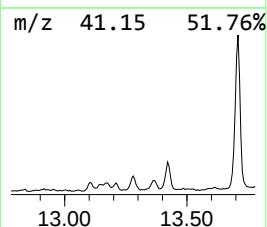
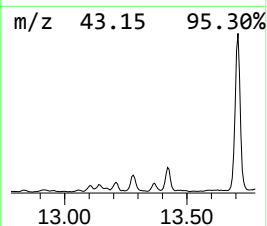
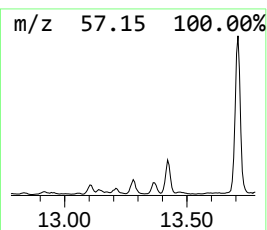
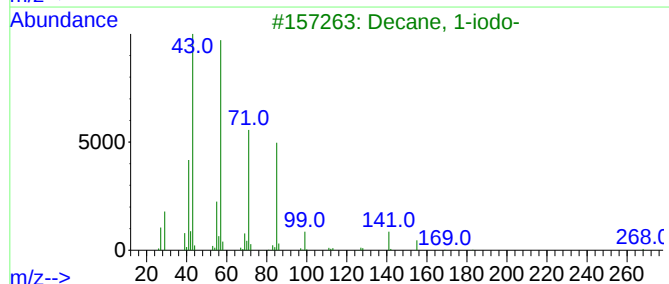
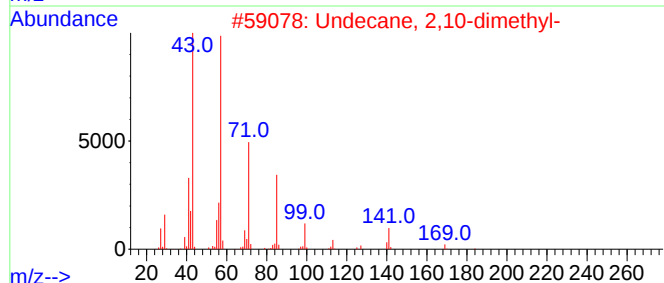
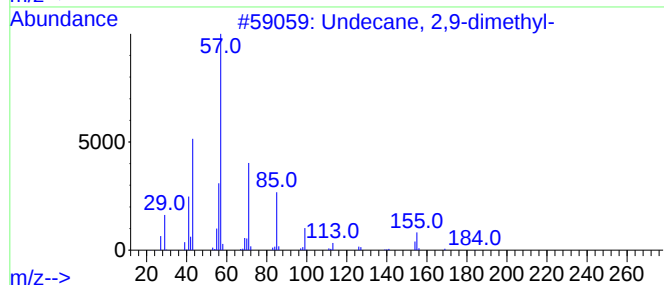
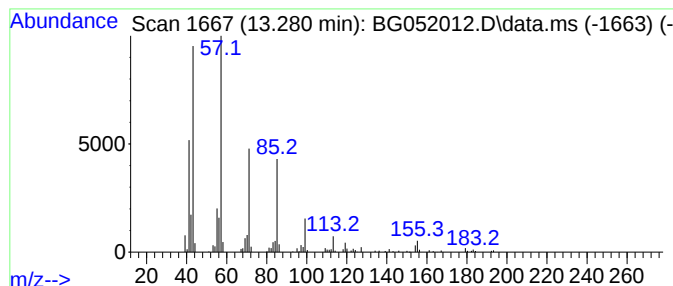
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Quant Title : SVOA CALIBRATION

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

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Peak Number 26 (DEL) Alkane: Straight-Chai... Concentration Rank 48

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 13.280 | 9.05 ng/ul | 220378 | Acenaphthene-d10 | 14.895 |

| Hit# | of | 5 | Tentative ID             | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------|-----|---------|-------------|------|
| 1    |    |   | Undecane, 2,9-dimethyl-  | 184 | C13H28  | 017301-26-7 | 83   |
| 2    |    |   | Undecane, 2,10-dimethyl- | 184 | C13H28  | 017301-27-8 | 81   |
| 3    |    |   | Decane, 1-iodo-          | 268 | C10H21I | 002050-77-3 | 78   |
| 4    |    |   | Dodecane, 2-methyl-      | 184 | C13H28  | 001560-97-0 | 72   |
| 5    |    |   | 4-Methyl-5-nonanone      | 156 | C10H20O | 035900-26-6 | 59   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011522\  
Data File : BG052012.D  
Acq On : 16 Jan 2022 1:50  
Operator : CG/JU  
Sample : N1132-04 10X  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLT5

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010622.M  
Quant Title : SVOA CALIBRATION

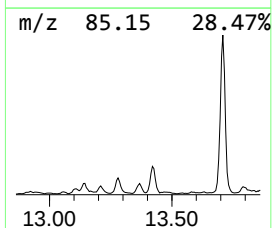
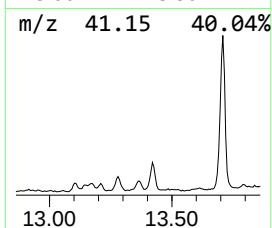
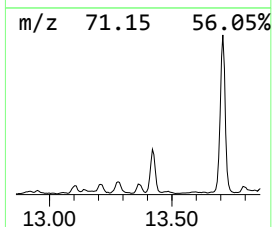
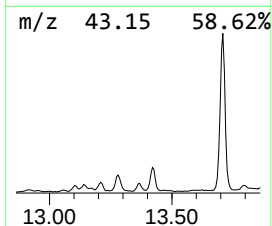
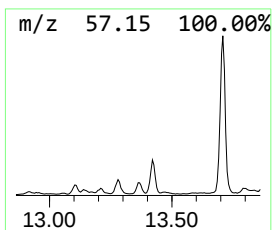
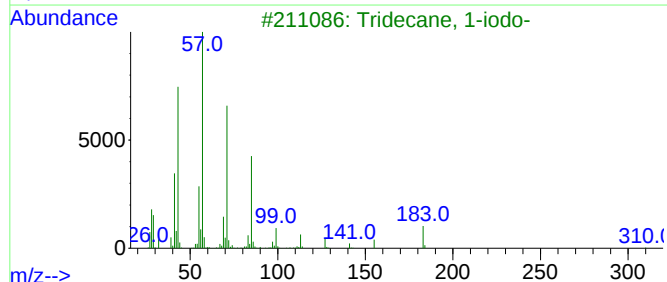
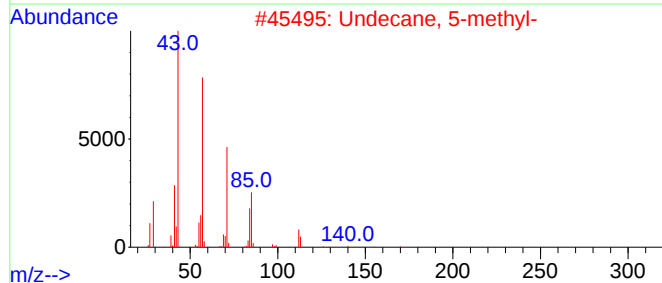
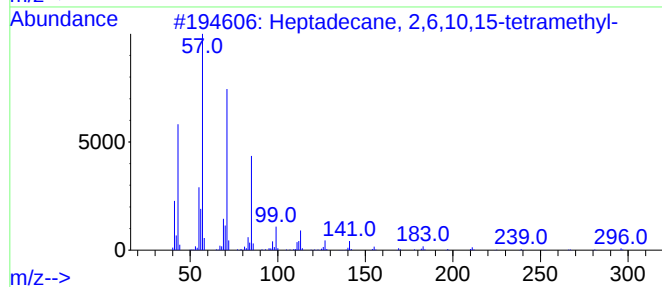
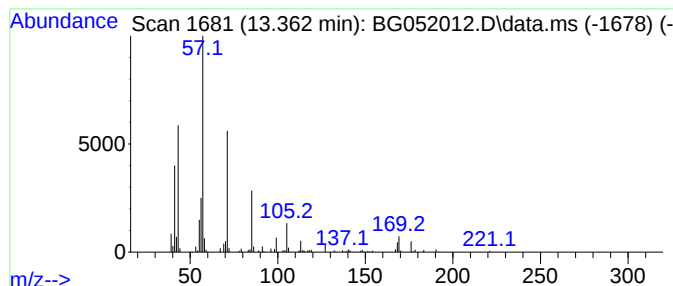
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 27 (DEL) Alkane: Straight-Chai... Concentration Rank 64

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 13.362 | 5.75 ng/ul | 140043 | Acenaphthene-d10 | 14.895 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|------|-------------------------------------|-----|---------|-------------|------|
| 1    |      | Heptadecane, 2,6,10,15-tetramethyl- | 296 | C21H44  | 054833-48-6 | 64   |
| 2    |      | Undecane, 5-methyl-                 | 170 | C12H26  | 001632-70-8 | 58   |
| 3    |      | Tridecane, 1-iodo-                  | 310 | C13H27I | 035599-77-0 | 53   |
| 4    |      | Dodecane, 2-methyl-                 | 184 | C13H28  | 001560-97-0 | 53   |
| 5    |      | Decane                              | 142 | C10H22  | 000124-18-5 | 53   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011522\  
Data File : BG052012.D  
Acq On : 16 Jan 2022 1:50  
Operator : CG/JU  
Sample : N1132-04 10X  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

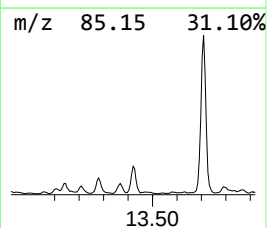
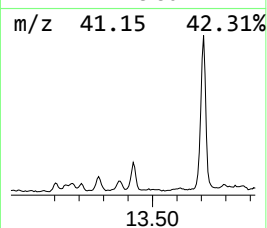
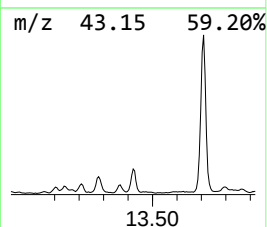
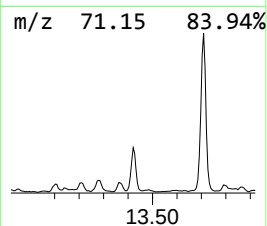
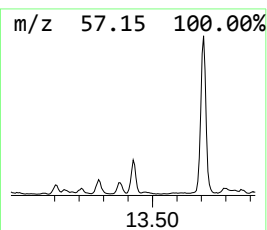
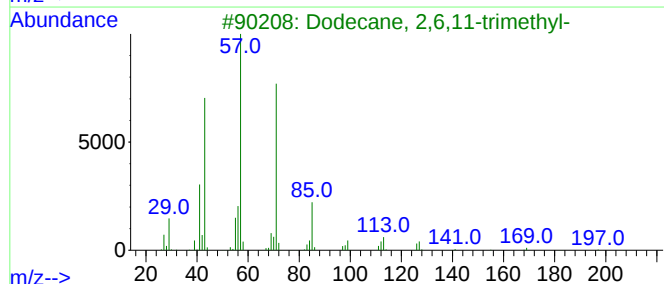
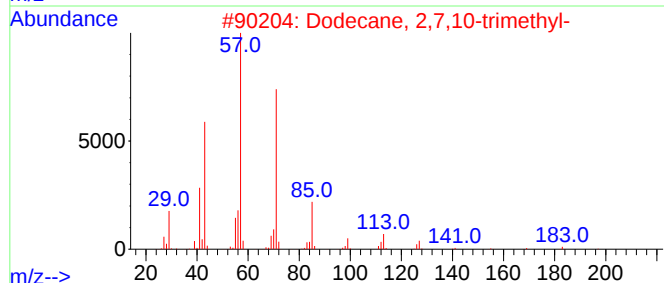
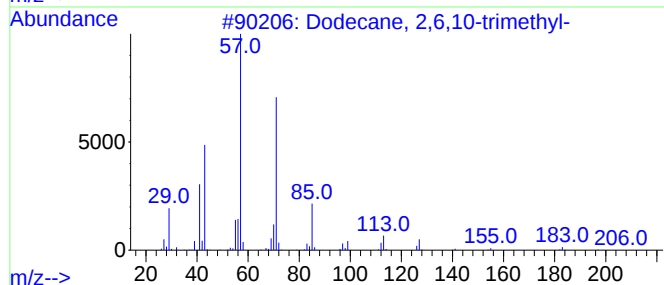
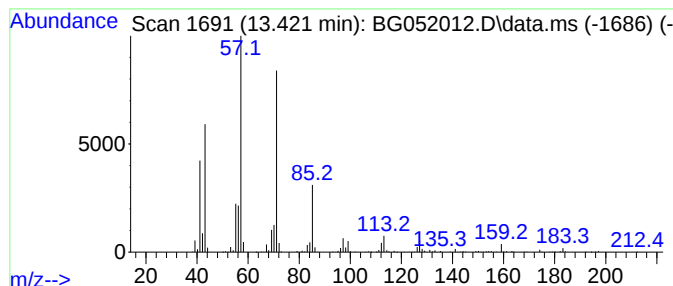
Instrument :  
BNA\_G  
ClientSampleId :  
BGLT5

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010622.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 28 (DEL) Alkane: Straight-Chai... Concentration Rank 20

| R.T.      | EstConc                            | Area   | Relative to ISTD |          |              | R.T.   |
|-----------|------------------------------------|--------|------------------|----------|--------------|--------|
| 13.421    | 22.05 ng/ul                        | 536931 | Acenaphthene-d10 |          |              | 14.895 |
| Hit# of 5 | Tentative ID                       |        | MW               | MolForm  | CAS#         | Qual   |
| 1         | Dodecane, 2,6,10-trimethyl-        |        | 212              | C15H32   | 003891-98-3  | 91     |
| 2         | Dodecane, 2,7,10-trimethyl-        |        | 212              | C15H32   | 074645-98-0  | 91     |
| 3         | Dodecane, 2,6,11-trimethyl-        |        | 212              | C15H32   | 031295-56-4  | 90     |
| 4         | Undecane, 6-ethyl-                 |        | 184              | C13H28   | 017312-60-6  | 64     |
| 5         | Oxalic acid, butyl neopentyl ester |        | 216              | C11H20O4 | 1000309-72-8 | 53     |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011522\  
Data File : BG052012.D  
Acq On : 16 Jan 2022 1:50  
Operator : CG/JU  
Sample : N1132-04 10X  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLT5

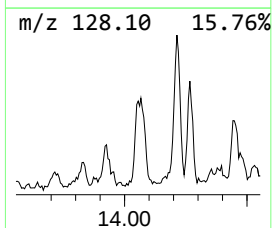
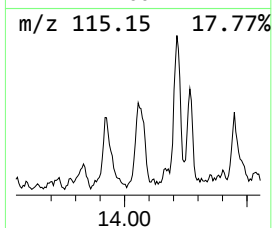
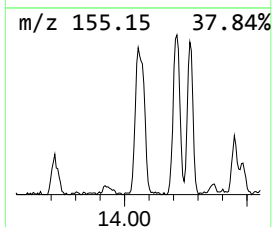
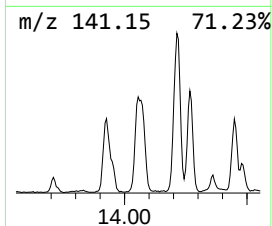
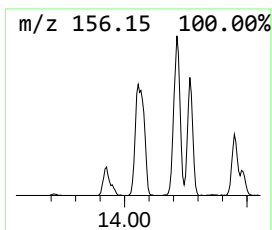
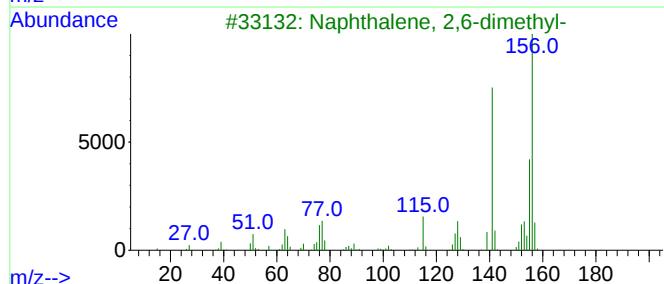
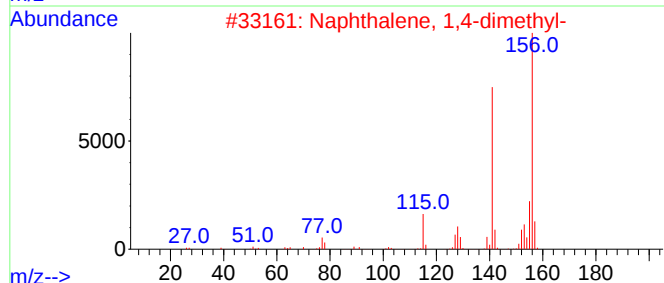
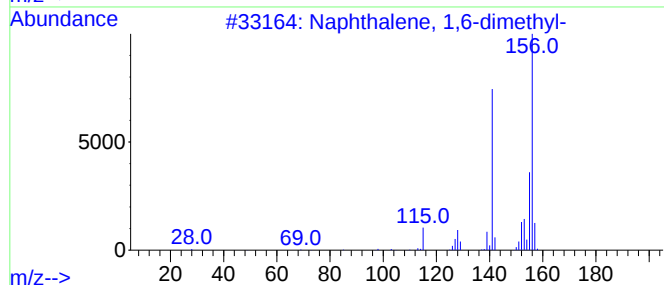
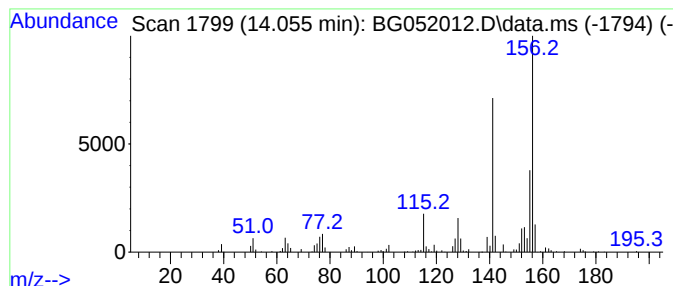
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Quant Title : SVOA CALIBRATION

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

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Peak Number 30 Naphthalene, 1,6-dimethyl- Concentration Rank 17

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 14.055 | 25.28 ng/ul | 615583 | Acenaphthene-d10 | 14.895 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, 1,6-dimethyl- | 156 | C12H12  | 000575-43-9 | 96   |
| 2    |    |   | Naphthalene, 1,4-dimethyl- | 156 | C12H12  | 000571-58-4 | 96   |
| 3    |    |   | Naphthalene, 2,6-dimethyl- | 156 | C12H12  | 000581-42-0 | 95   |
| 4    |    |   | Naphthalene, 1,4-dimethyl- | 156 | C12H12  | 000571-58-4 | 95   |
| 5    |    |   | Naphthalene, 1,7-dimethyl- | 156 | C12H12  | 000575-37-1 | 95   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011522\  
Data File : BG052012.D  
Acq On : 16 Jan 2022 1:50  
Operator : CG/JU  
Sample : N1132-04 10X  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLT5

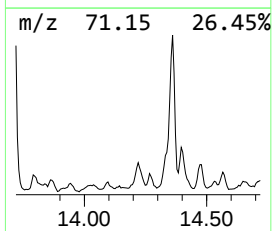
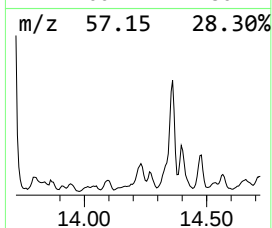
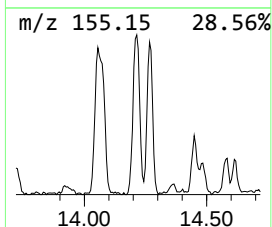
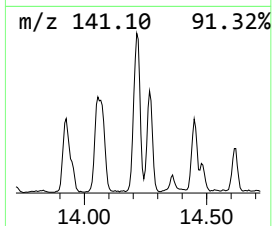
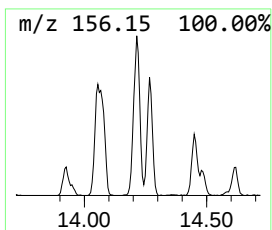
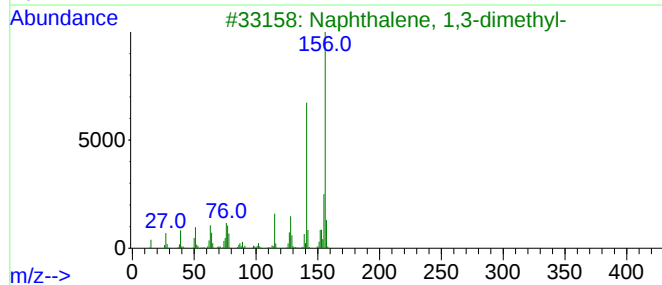
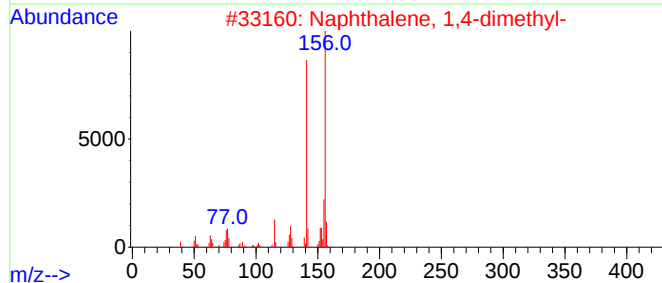
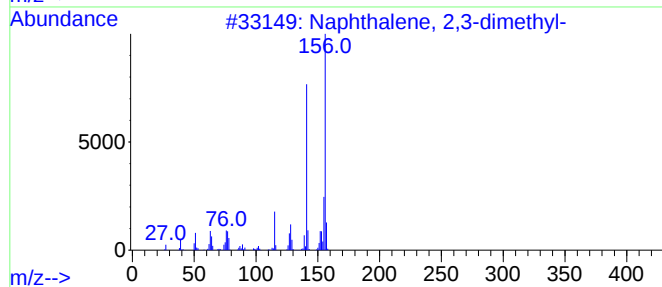
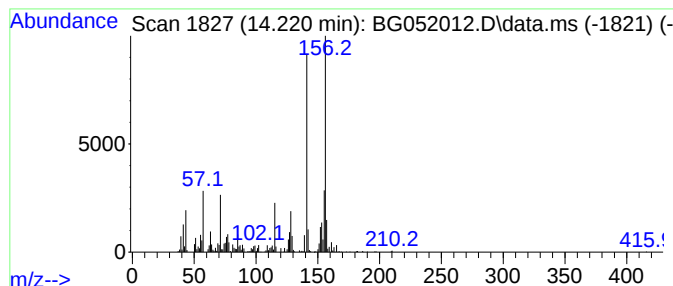
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Quant Title : SVOA CALIBRATION

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

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Peak Number 31 Naphthalene, 2,3-dimethyl- Concentration Rank 16

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 14.220 | 26.84 ng/ul | 653619 | Acenaphthene-d10 | 14.895 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, 2,3-dimethyl- | 156 | C12H12  | 000581-40-8 | 97   |
| 2    |    |   | Naphthalene, 1,4-dimethyl- | 156 | C12H12  | 000571-58-4 | 96   |
| 3    |    |   | Naphthalene, 1,3-dimethyl- | 156 | C12H12  | 000575-41-7 | 96   |
| 4    |    |   | Naphthalene, 1,3-dimethyl- | 156 | C12H12  | 000575-41-7 | 96   |
| 5    |    |   | Naphthalene, 1,4-dimethyl- | 156 | C12H12  | 000571-58-4 | 96   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011522\  
Data File : BG052012.D  
Acq On : 16 Jan 2022 1:50  
Operator : CG/JU  
Sample : N1132-04 10X  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLT5

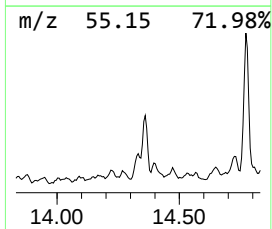
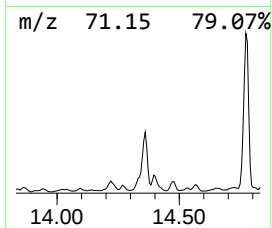
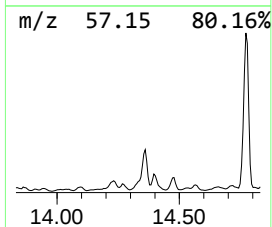
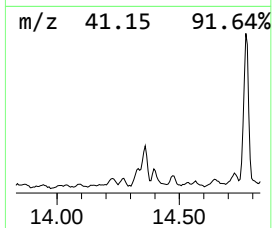
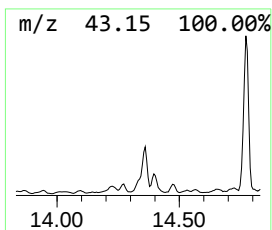
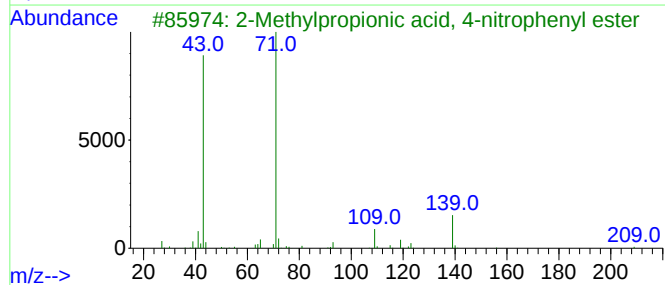
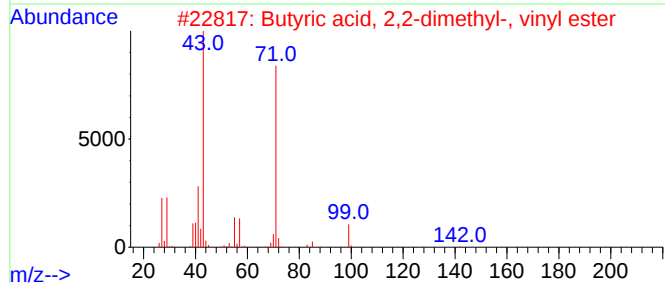
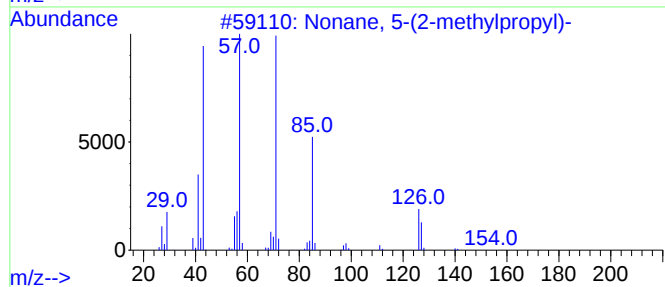
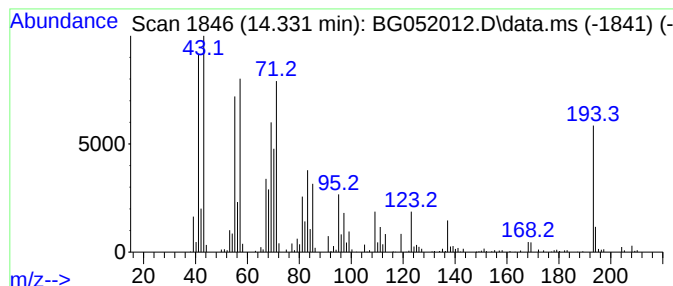
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Quant Title : SVOA CALIBRATION

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

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Peak Number 32 (DEL) Alkane: Straight-Chai... Concentration Rank 29

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 14.331 | 16.23 ng/ul | 395229 | Acenaphthene-d10 | 14.895 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Nonane, 5-(2-methylpropyl)-         | 184 | C13H28    | 062185-53-9  | 27   |
| 2    |    |   | Butyric acid, 2,2-dimethyl-, vin... | 142 | C8H14O2   | 013170-00-8  | 22   |
| 3    |    |   | 2-Methylpropionic acid, 4-nitrop... | 209 | C10H11NO4 | 004195-16-8  | 22   |
| 4    |    |   | Carbonic acid, hexadecyl prop-1-... | 326 | C20H38O3  | 1000382-90-3 | 18   |
| 5    |    |   | Carbonic acid, octadecyl prop-1-... | 354 | C22H42O3  | 1000383-11-5 | 18   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011522\  
Data File : BG052012.D  
Acq On : 16 Jan 2022 1:50  
Operator : CG/JU  
Sample : N1132-04 10X  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLT5

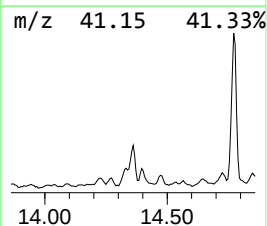
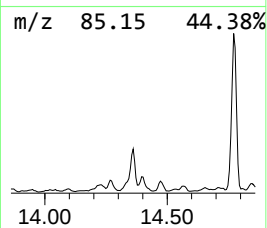
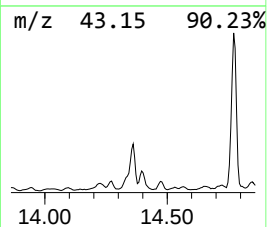
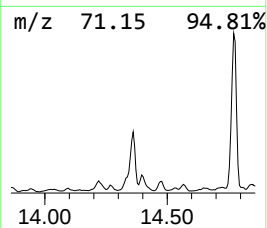
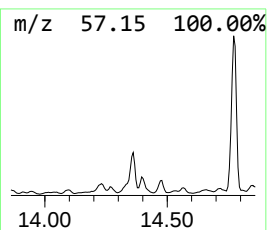
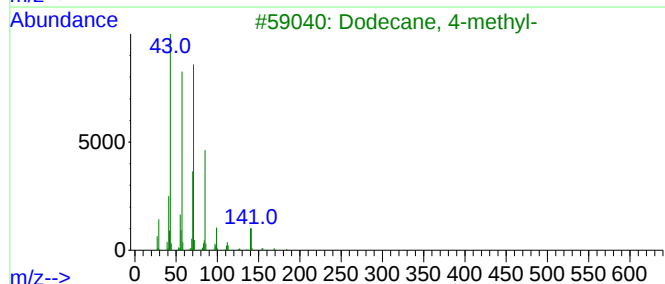
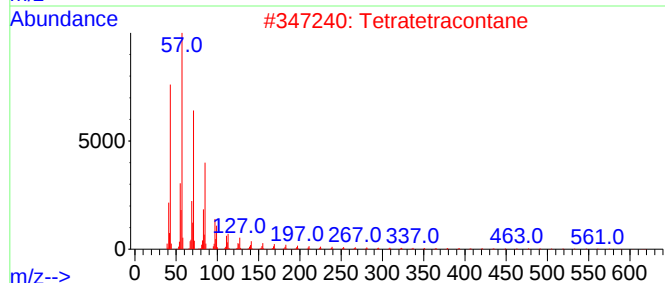
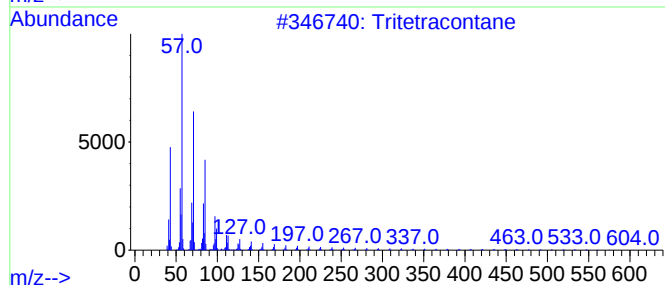
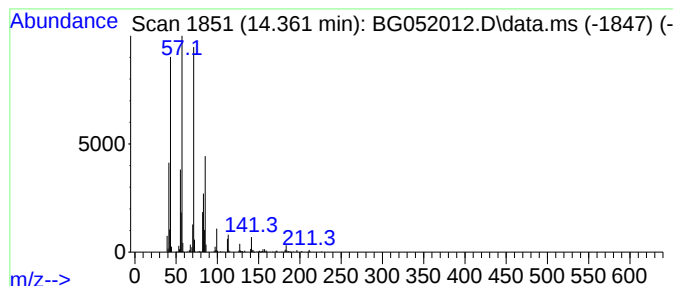
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Quant Title : SVOA CALIBRATION

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

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Peak Number 33 (DEL) Alkane: Straight-Chai... Concentration Rank 13

| R.T.      | EstConc             | Area   | Relative to ISTD |         | R.T.        |      |
|-----------|---------------------|--------|------------------|---------|-------------|------|
| 14.361    | 30.27 ng/ul         | 737238 | Acenaphthene-d10 |         | 14.895      |      |
| Hit# of 5 | Tentative ID        |        | MW               | MolForm | CAS#        | Qual |
| 1         | Tritetracontane     |        | 605              | C43H88  | 007098-21-7 | 90   |
| 2         | Tetratetracontane   |        | 619              | C44H90  | 007098-22-8 | 90   |
| 3         | Dodecane, 4-methyl- |        | 184              | C13H28  | 006117-97-1 | 81   |
| 4         | Decane, 5-propyl-   |        | 184              | C13H28  | 017312-62-8 | 74   |
| 5         | Dodecane, 4-methyl- |        | 184              | C13H28  | 006117-97-1 | 72   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011522\  
Data File : BG052012.D  
Acq On : 16 Jan 2022 1:50  
Operator : CG/JU  
Sample : N1132-04 10X  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLT5

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010622.M  
Quant Title : SVOA CALIBRATION

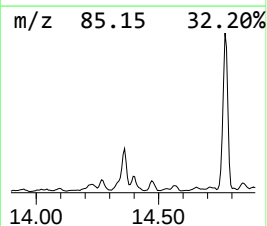
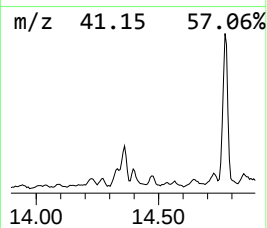
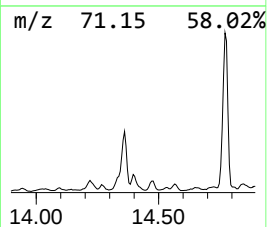
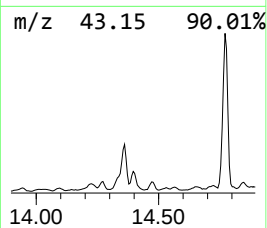
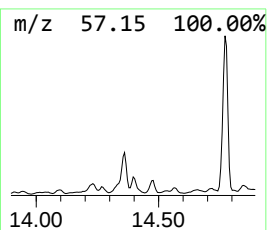
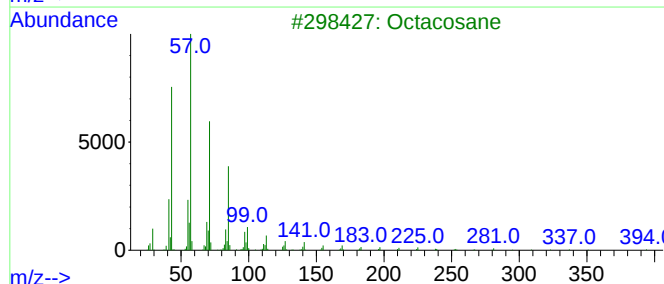
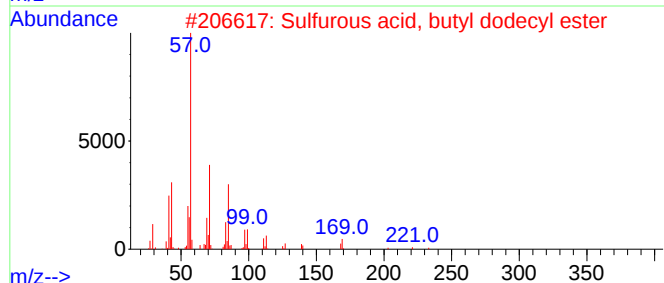
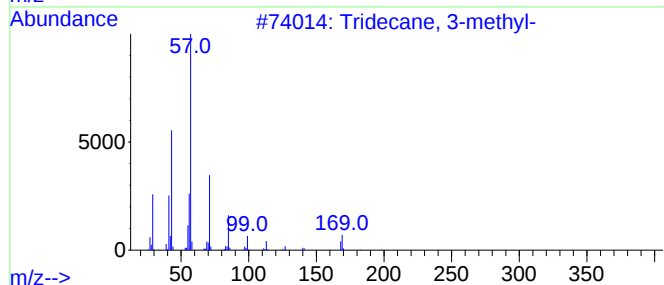
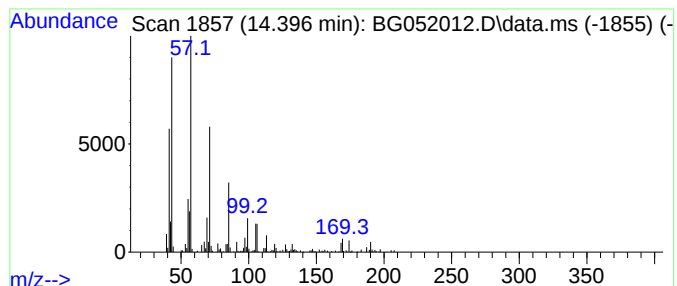
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 34 (DEL) Alkane: Straight-Chai... Concentration Rank 46

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 14.396 | 9.29 ng/ul | 226284 | Acenaphthene-d10 | 14.895 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Tridecane, 3-methyl-                | 198 | C14H30    | 006418-41-3  | 72   |
| 2    |    |   | Sulfurous acid, butyl dodecyl ester | 306 | C16H34O3S | 1000309-17-9 | 72   |
| 3    |    |   | Octacosane                          | 394 | C28H58    | 000630-02-4  | 64   |
| 4    |    |   | Heptacosane                         | 380 | C27H56    | 000593-49-7  | 64   |
| 5    |    |   | Dodecane, 2-methyl-                 | 184 | C13H28    | 001560-97-0  | 64   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011522\  
Data File : BG052012.D  
Acq On : 16 Jan 2022 1:50  
Operator : CG/JU  
Sample : N1132-04 10X  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLT5

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010622.M  
Quant Title : SVOA CALIBRATION

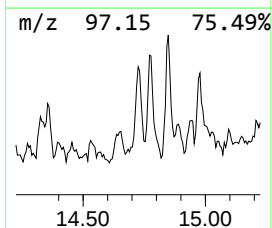
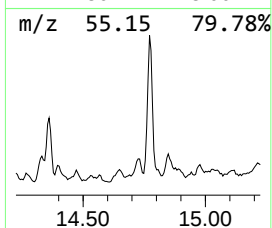
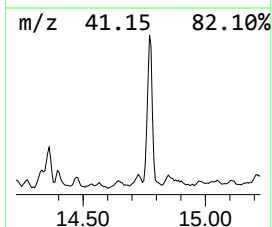
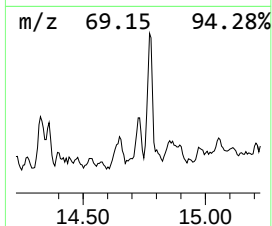
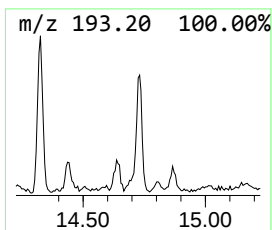
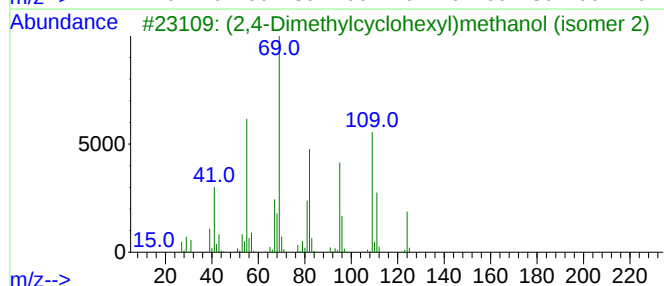
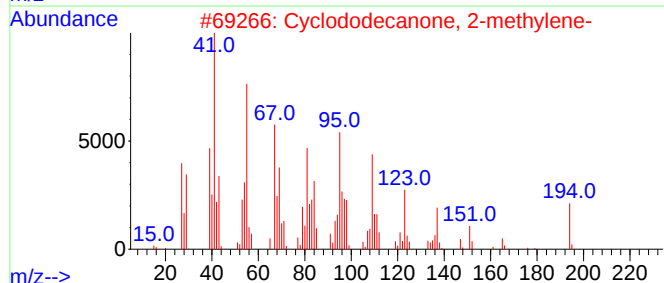
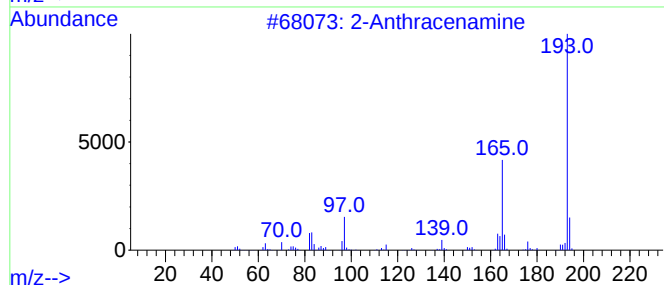
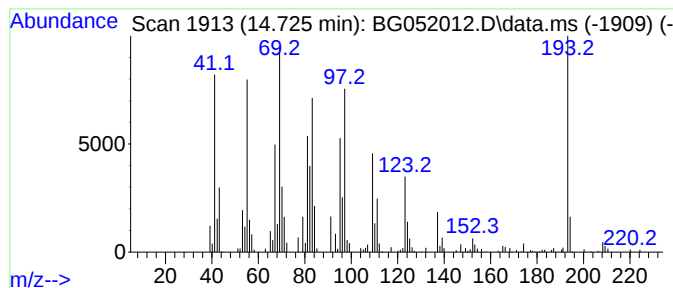
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 35 unknown-02 Concentration Rank 44

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 14.725 | 9.52 ng/ul | 231878 | Acenaphthene-d10 | 14.895 |

| Hit# | of                                  | 5 | Tentative ID | MW  | MolForm | CAS#         | Qual |
|------|-------------------------------------|---|--------------|-----|---------|--------------|------|
| 1    | 2-Anthracenamine                    |   |              | 193 | C14H11N | 000613-13-8  | 38   |
| 2    | Cyclododecanone, 2-methylene-       |   |              | 194 | C13H22O | 003045-76-9  | 38   |
| 3    | (2,4-Dimethylcyclohexyl)methanol... |   |              | 142 | C9H18O  | 1000499-02-7 | 27   |
| 4    | 2-Pentene, 3-ethyl-4,4-dimethyl-    |   |              | 126 | C9H18   | 053907-59-8  | 27   |
| 5    | photocitral B                       |   |              | 152 | C10H16O | 006040-45-5  | 18   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011522\  
Data File : BG052012.D  
Acq On : 16 Jan 2022 1:50  
Operator : CG/JU  
Sample : N1132-04 10X  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLT5

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010622.M  
Quant Title : SVOA CALIBRATION

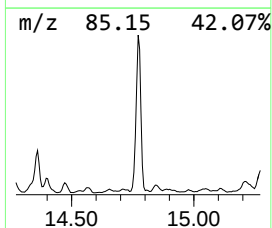
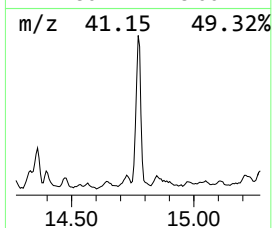
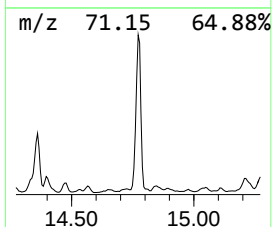
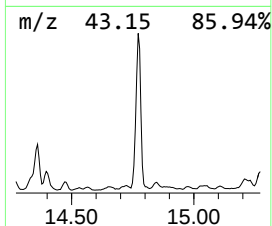
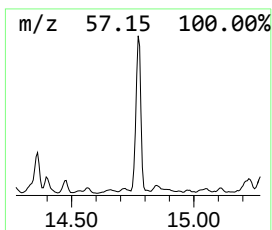
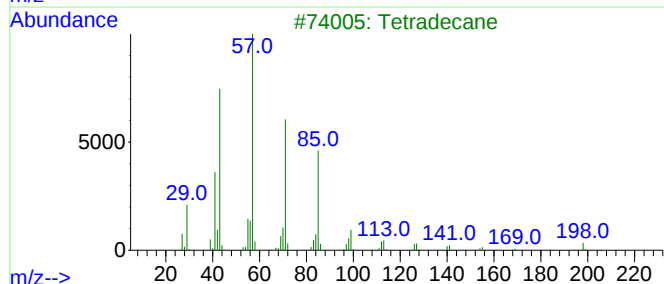
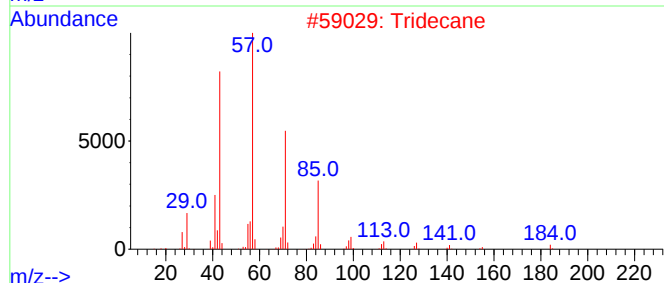
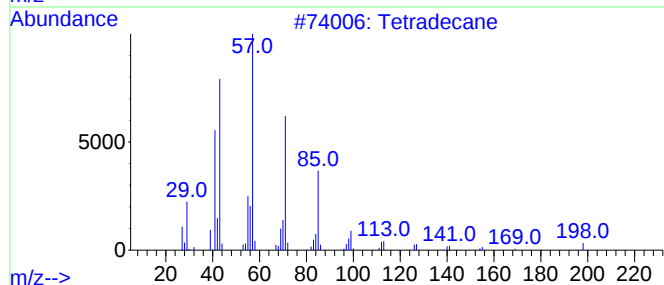
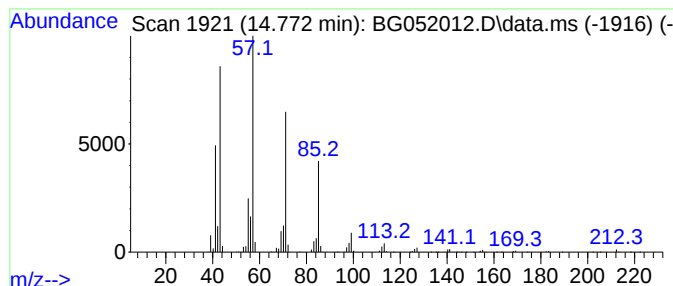
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TIC Integration Parameters: LSCINT.P

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Peak Number 36 (DEL) Alkane: Straight-Chai... Concentration Rank 1

| R.T.   | EstConc      | Area    | Relative to ISTD | R.T.   |
|--------|--------------|---------|------------------|--------|
| 14.772 | 108.10 ng/ul | 2632800 | Acenaphthene-d10 | 14.895 |

| Hit# | of 5 | Tentative ID             | MW  | MolForm | CAS#        | Qual |
|------|------|--------------------------|-----|---------|-------------|------|
| 1    |      | Tetradecane              | 198 | C14H30  | 000629-59-4 | 90   |
| 2    |      | Tridecane                | 184 | C13H28  | 000629-50-5 | 90   |
| 3    |      | Tetradecane              | 198 | C14H30  | 000629-59-4 | 90   |
| 4    |      | Tridecane, 7-propyl-     | 226 | C16H34  | 055045-09-5 | 86   |
| 5    |      | Decane, 2,3,5-trimethyl- | 184 | C13H28  | 062238-11-3 | 83   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011522\  
Data File : BG052012.D  
Acq On : 16 Jan 2022 1:50  
Operator : CG/JU  
Sample : N1132-04 10X  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLT5

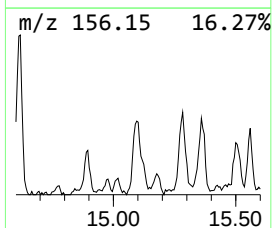
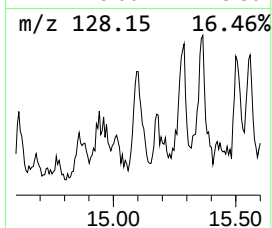
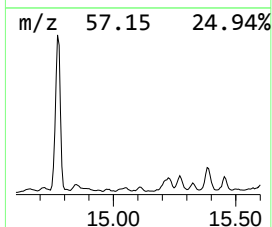
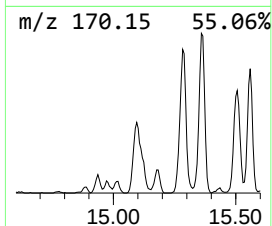
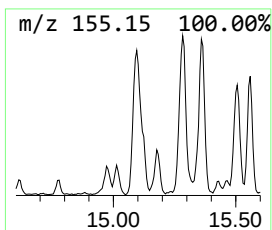
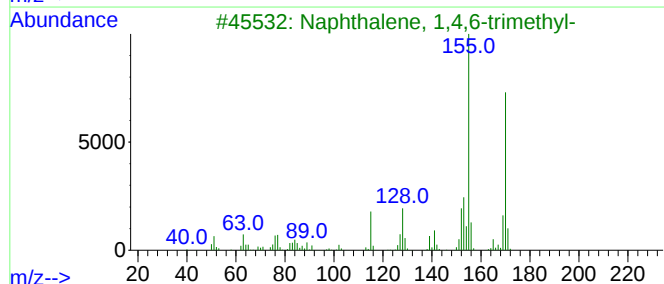
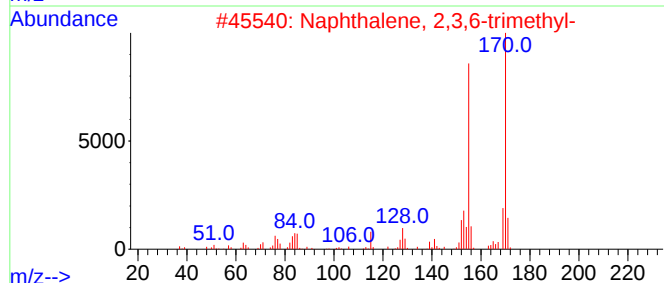
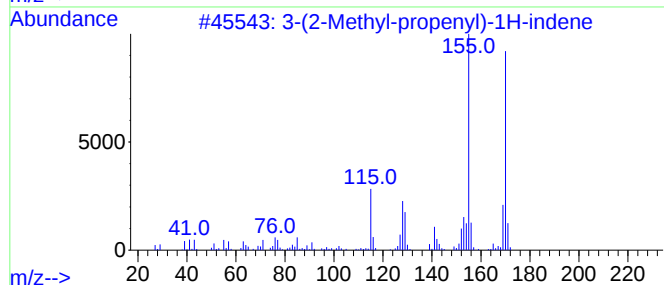
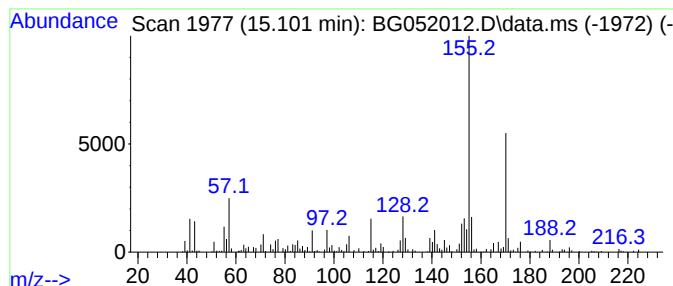
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Quant Title : SVOA CALIBRATION

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

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Peak Number 37 3-(2-Methyl-propenyl)-1H-in... Concentration Rank 30

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 15.101 | 15.54 ng/ul | 378398 | Acenaphthene-d10 | 14.895 |

| Hit# | of 5 | Tentative ID                    | MW  | MolForm | CAS#         | Qual |
|------|------|---------------------------------|-----|---------|--------------|------|
| 1    |      | 3-(2-Methyl-propenyl)-1H-indene | 170 | C13H14  | 1000187-78-5 | 93   |
| 2    |      | Naphthalene, 2,3,6-trimethyl-   | 170 | C13H14  | 000829-26-5  | 93   |
| 3    |      | Naphthalene, 1,4,6-trimethyl-   | 170 | C13H14  | 002131-42-2  | 92   |
| 4    |      | Naphthalene, 1,6,7-trimethyl-   | 170 | C13H14  | 002245-38-7  | 90   |
| 5    |      | Naphthalene, 1,4,6-trimethyl-   | 170 | C13H14  | 002131-42-2  | 90   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011522\  
Data File : BG052012.D  
Acq On : 16 Jan 2022 1:50  
Operator : CG/JU  
Sample : N1132-04 10X  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLT5

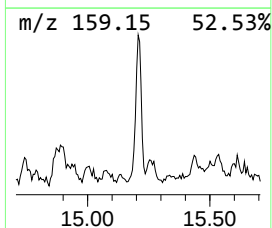
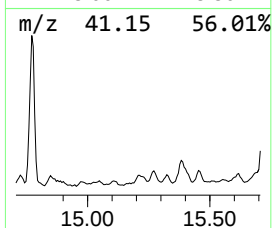
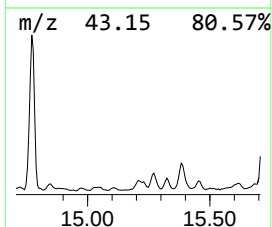
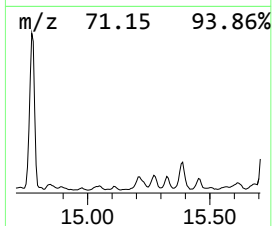
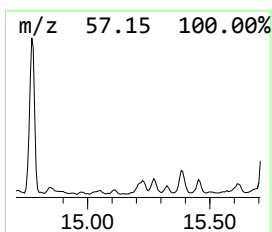
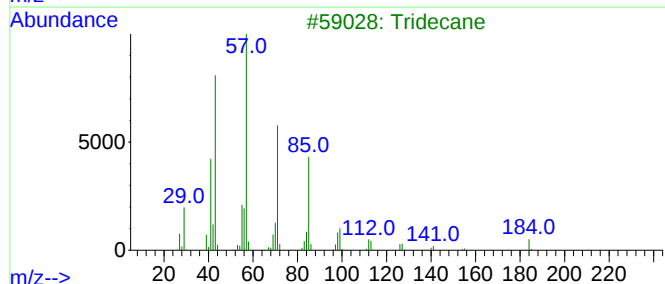
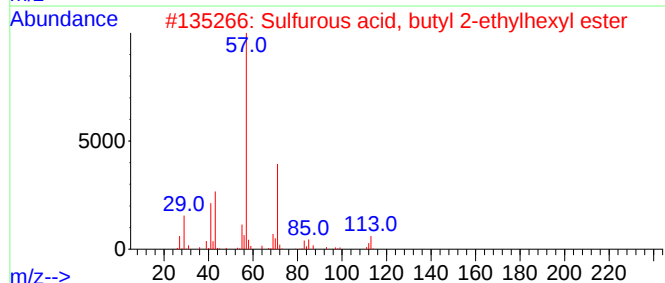
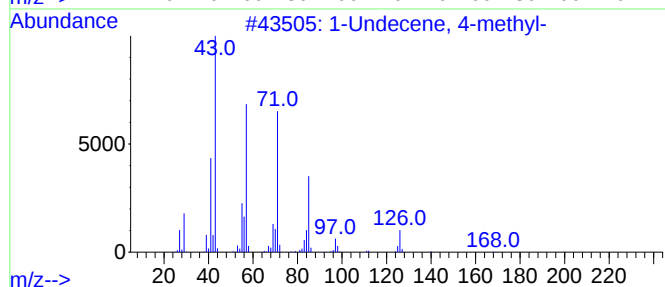
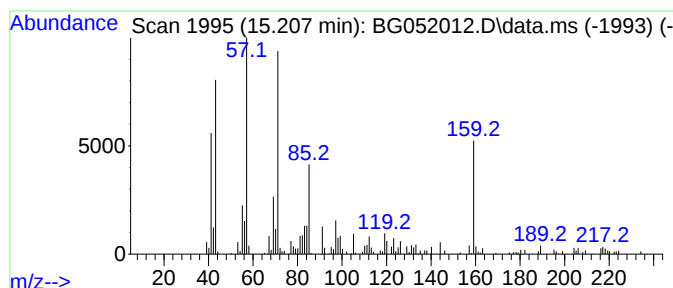
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Quant Title : SVOA CALIBRATION

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

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Peak Number 38 (DEL) Alkane: Straight-Chai... Concentration Rank 66

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|--------|------------------|--------------|------|
| 15.207    | 5.16 ng/ul                          | 125718 | Acenaphthene-d10 | 14.895       |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#         | Qual |
| 1         | 1-Undecene, 4-methyl-               | 168    | C12H24           | 074630-39-0  | 50   |
| 2         | Sulfurous acid, butyl 2-ethylhex... | 250    | C12H26O3S        | 1000309-18-8 | 38   |
| 3         | Tridecane                           | 184    | C13H28           | 000629-50-5  | 35   |
| 4         | Hexadecane                          | 226    | C16H34           | 000544-76-3  | 35   |
| 5         | Tetradecane                         | 198    | C14H30           | 000629-59-4  | 35   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011522\  
Data File : BG052012.D  
Acq On : 16 Jan 2022 1:50  
Operator : CG/JU  
Sample : N1132-04 10X  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLT5

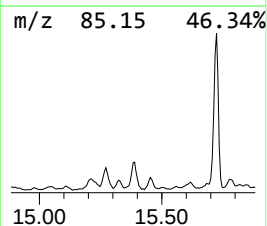
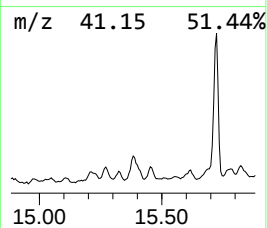
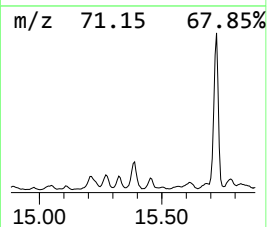
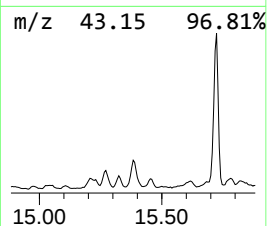
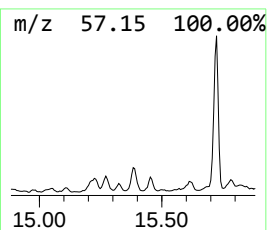
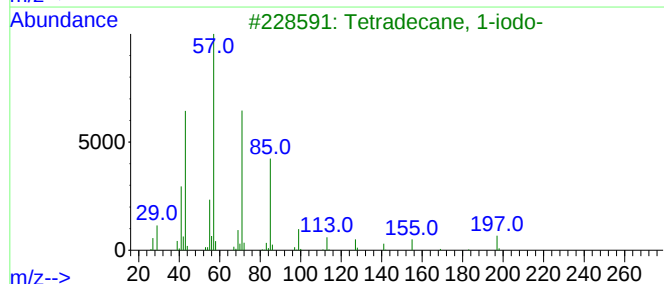
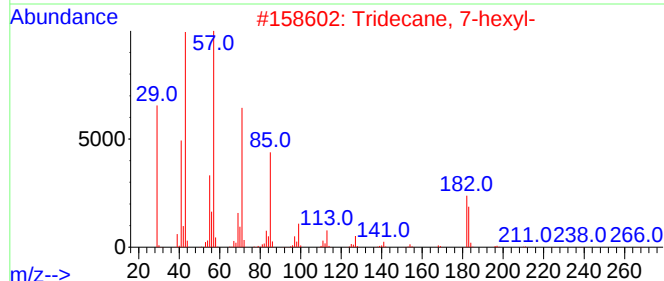
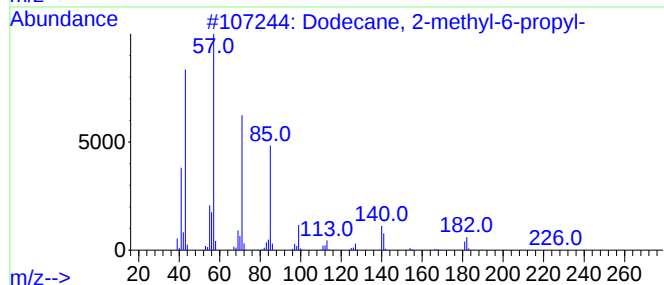
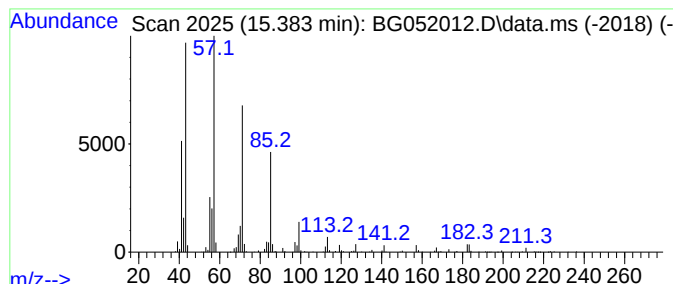
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Quant Title : SVOA CALIBRATION

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

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Peak Number 39 (DEL) Alkane: Straight-Chai... Concentration Rank 11

| R.T.      | EstConc                      | Area   | Relative to ISTD |         | R.T.        |      |
|-----------|------------------------------|--------|------------------|---------|-------------|------|
| 15.383    | 35.21 ng/ul                  | 857474 | Acenaphthene-d10 |         | 14.895      |      |
| Hit# of 5 | Tentative ID                 |        | MW               | MolForm | CAS#        | Qual |
| 1         | Dodecane, 2-methyl-6-propyl- |        | 226              | C16H34  | 055045-08-4 | 78   |
| 2         | Tridecane, 7-hexyl-          |        | 268              | C19H40  | 007225-66-3 | 78   |
| 3         | Tetradecane, 1-iodo-         |        | 324              | C14H29I | 019218-94-1 | 72   |
| 4         | Hexadecane                   |        | 226              | C16H34  | 000544-76-3 | 72   |
| 5         | Nonadecane                   |        | 268              | C19H40  | 000629-92-5 | 64   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011522\  
Data File : BG052012.D  
Acq On : 16 Jan 2022 1:50  
Operator : CG/JU  
Sample : N1132-04 10X  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLT5

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010622.M  
Quant Title : SVOA CALIBRATION

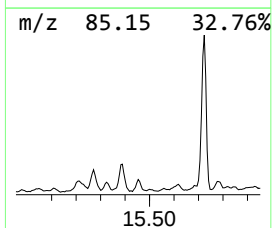
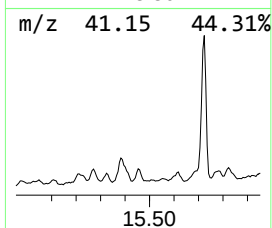
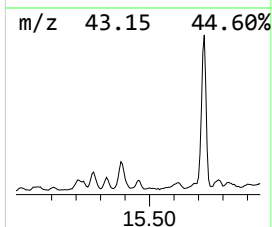
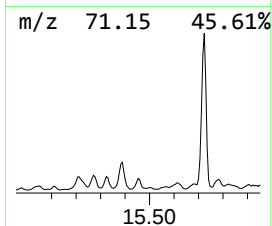
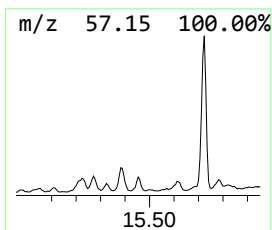
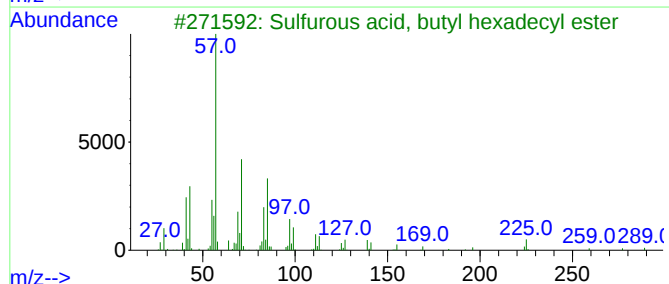
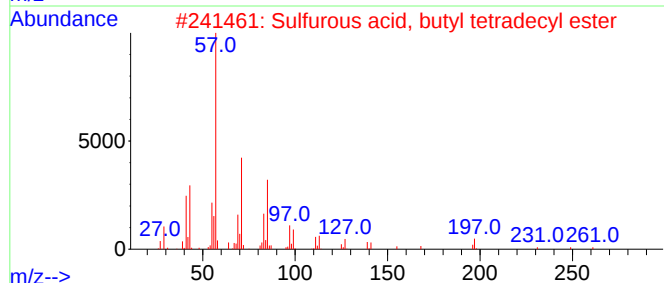
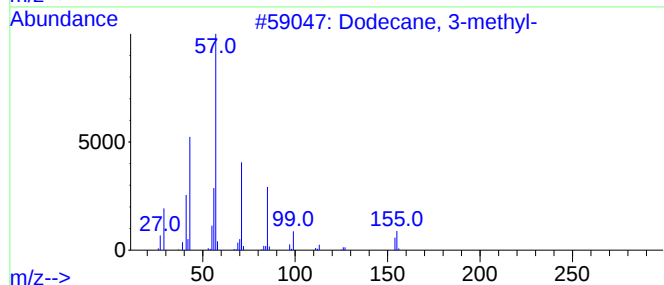
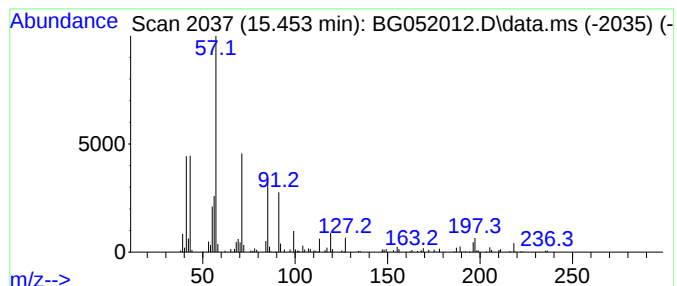
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 40 (DEL) Alkane: Straight-Chai... Concentration Rank 53

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 15.453 | 8.02 ng/ul | 195407 | Acenaphthene-d10 | 14.895 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Dodecane, 3-methyl-                 | 184 | C13H28    | 017312-57-1  | 72   |
| 2    |    |   | Sulfurous acid, butyl tetradecyl... | 334 | C18H38O3S | 1010309-18-1 | 64   |
| 3    |    |   | Sulfurous acid, butyl hexadecyl ... | 362 | C20H42O3S | 1000309-18-3 | 59   |
| 4    |    |   | Octadecane                          | 254 | C18H38    | 000593-45-3  | 59   |
| 5    |    |   | Benzeneacetic acid, 2-tetradecyl... | 332 | C22H36O2  | 1000282-02-4 | 53   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011522\  
Data File : BG052012.D  
Acq On : 16 Jan 2022 1:50  
Operator : CG/JU  
Sample : N1132-04 10X  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLT5

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010622.M  
Quant Title : SVOA CALIBRATION

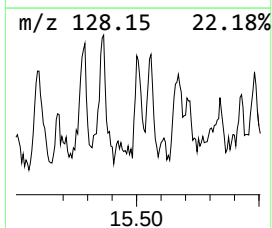
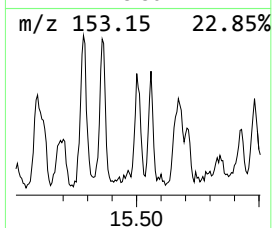
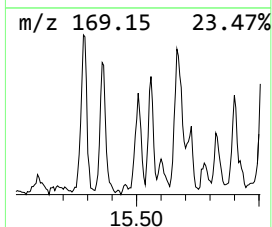
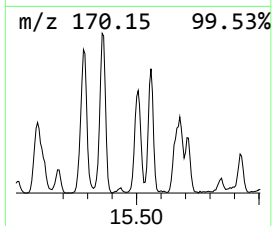
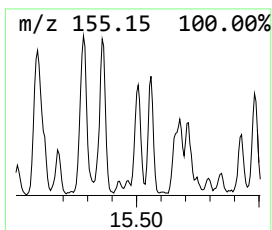
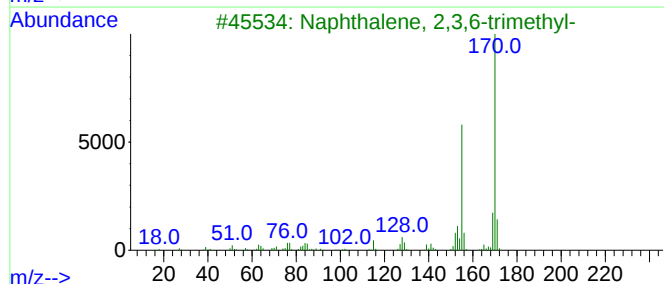
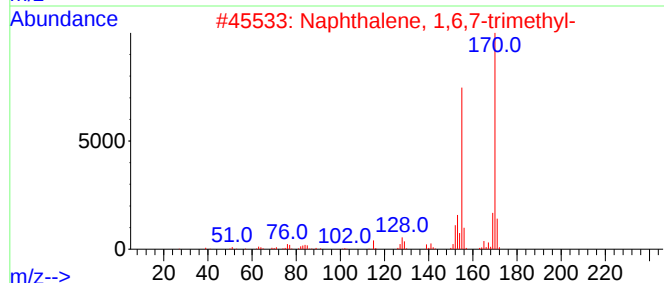
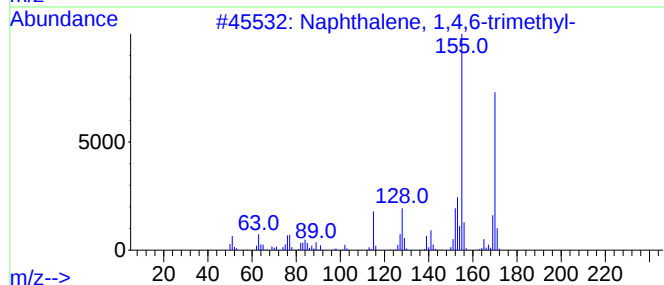
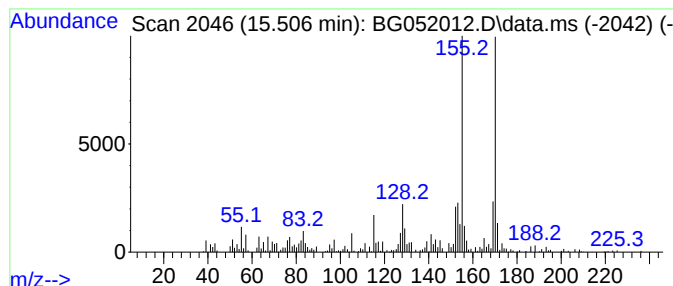
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 41 Naphthalene, 1,4,6-trimethyl- Concentration Rank 36

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 15.506 | 12.26 ng/ul | 298591 | Acenaphthene-d10 | 14.895 |

| Hit# | of | 5 | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, 1,4,6-trimethyl- | 170 | C13H14  | 002131-42-2 | 97   |
| 2    |    |   | Naphthalene, 1,6,7-trimethyl- | 170 | C13H14  | 002245-38-7 | 97   |
| 3    |    |   | Naphthalene, 2,3,6-trimethyl- | 170 | C13H14  | 000829-26-5 | 95   |
| 4    |    |   | Naphthalene, 2,3,6-trimethyl- | 170 | C13H14  | 000829-26-5 | 95   |
| 5    |    |   | Naphthalene, 1,6,7-trimethyl- | 170 | C13H14  | 002245-38-7 | 95   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011522\  
Data File : BG052012.D  
Acq On : 16 Jan 2022 1:50  
Operator : CG/JU  
Sample : N1132-04 10X  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLT5

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010622.M  
Quant Title : SVOA CALIBRATION

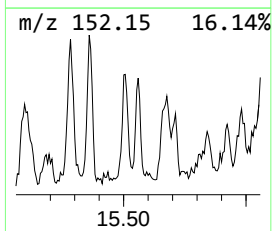
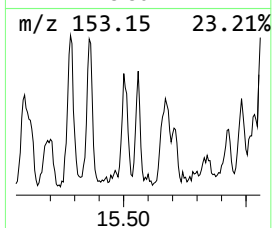
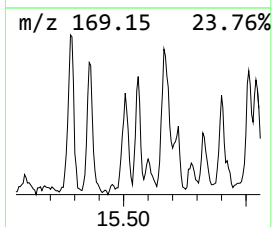
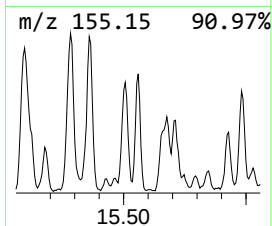
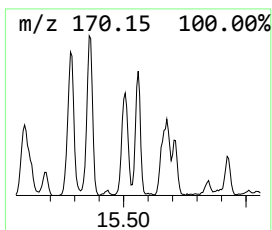
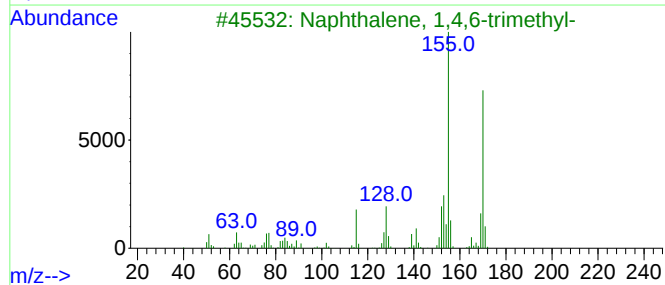
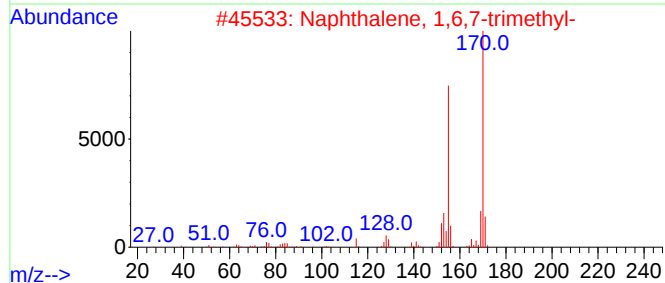
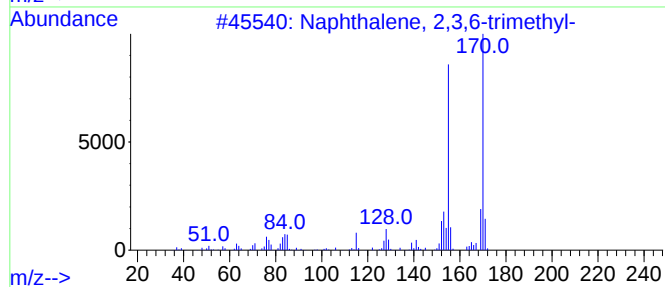
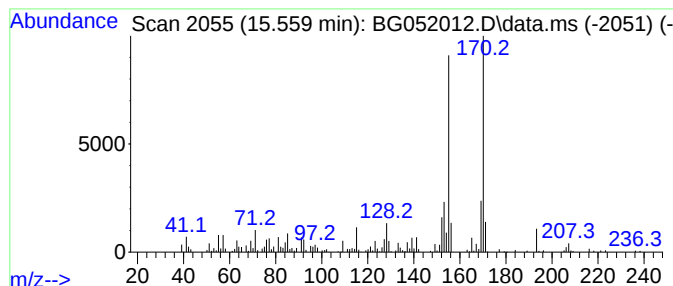
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 42 Naphthalene, 2,3,6-trimethyl- Concentration Rank 42

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 15.559 | 10.54 ng/ul | 256676 | Acenaphthene-d10 | 14.895 |

| Hit# | of | 5 | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, 2,3,6-trimethyl- | 170 | C13H14  | 000829-26-5 | 97   |
| 2    |    |   | Naphthalene, 1,6,7-trimethyl- | 170 | C13H14  | 002245-38-7 | 97   |
| 3    |    |   | Naphthalene, 1,4,6-trimethyl- | 170 | C13H14  | 002131-42-2 | 96   |
| 4    |    |   | Naphthalene, 1,6,7-trimethyl- | 170 | C13H14  | 002245-38-7 | 96   |
| 5    |    |   | Naphthalene, 2,3,6-trimethyl- | 170 | C13H14  | 000829-26-5 | 95   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011522\  
Data File : BG052012.D  
Acq On : 16 Jan 2022 1:50  
Operator : CG/JU  
Sample : N1132-04 10X  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLT5

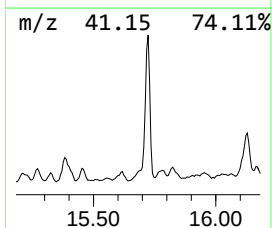
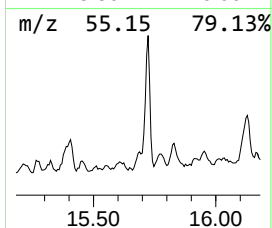
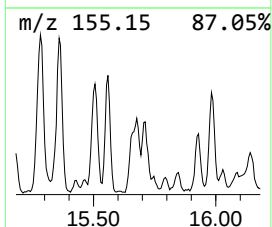
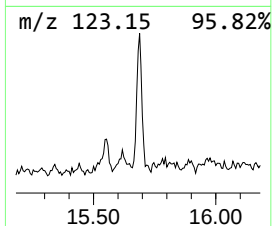
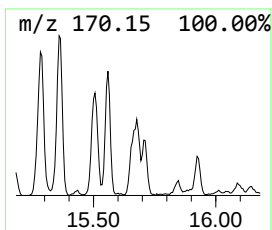
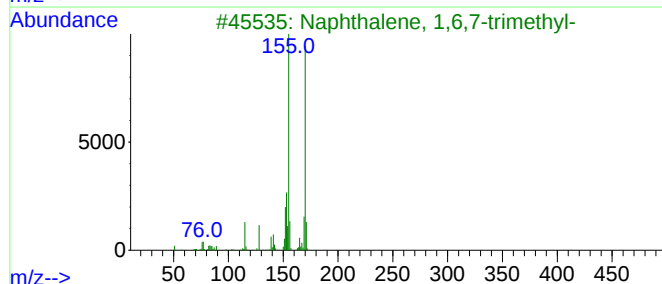
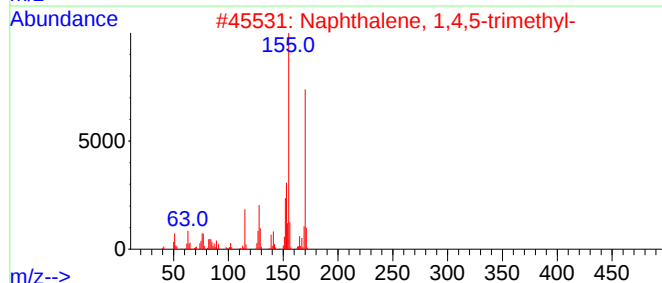
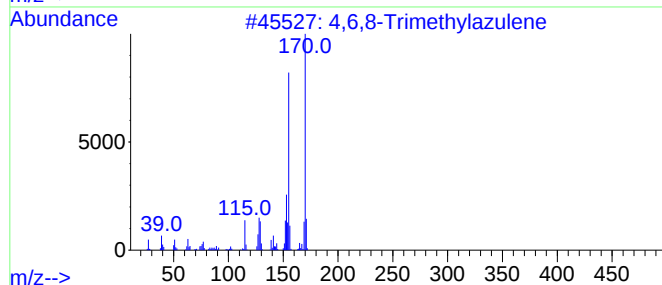
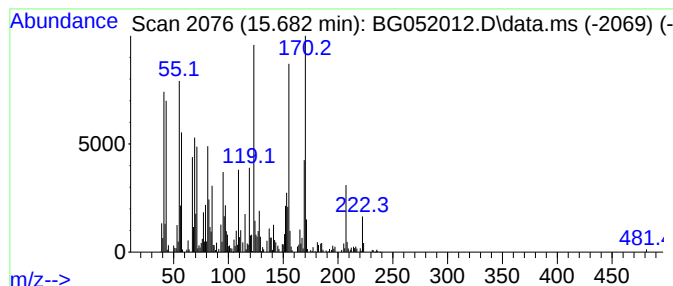
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Quant Title : SVOA CALIBRATION

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

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Peak Number 43 4,6,8-Trimethylazulene Concentration Rank 21

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 15.682 | 20.67 ng/ul | 503506 | Acenaphthene-d10 | 14.895 |

| Hit# | of | 5 | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------|-----|---------|-------------|------|
| 1    |    |   | 4,6,8-Trimethylazulene        | 170 | C13H14  | 000941-81-1 | 60   |
| 2    |    |   | Naphthalene, 1,4,5-trimethyl- | 170 | C13H14  | 002131-41-1 | 52   |
| 3    |    |   | Naphthalene, 1,6,7-trimethyl- | 170 | C13H14  | 002245-38-7 | 46   |
| 4    |    |   | Naphthalene, 1,6,7-trimethyl- | 170 | C13H14  | 002245-38-7 | 46   |
| 5    |    |   | 4,6,8-Trimethylazulene        | 170 | C13H14  | 000941-81-1 | 46   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011522\  
Data File : BG052012.D  
Acq On : 16 Jan 2022 1:50  
Operator : CG/JU  
Sample : N1132-04 10X  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLT5

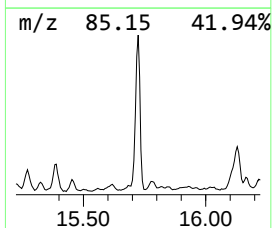
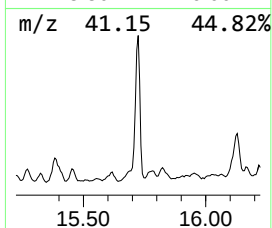
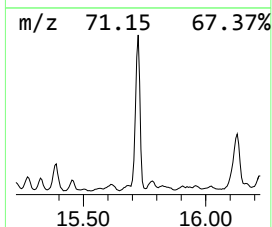
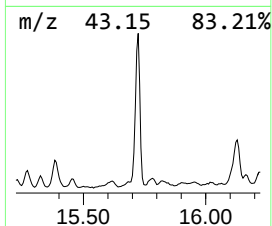
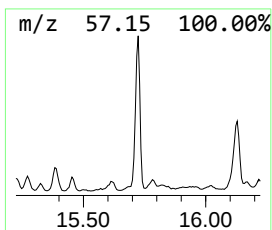
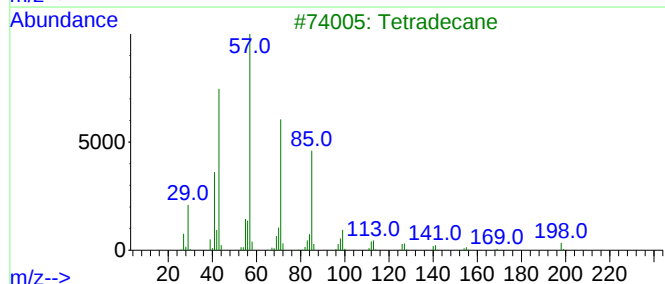
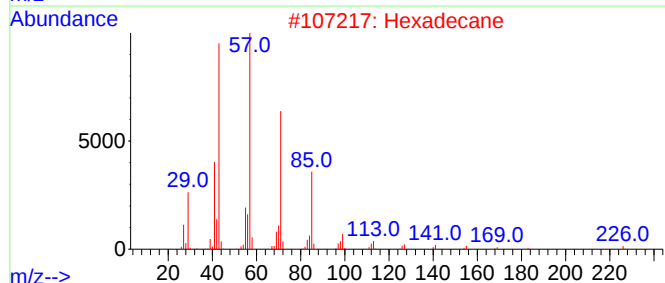
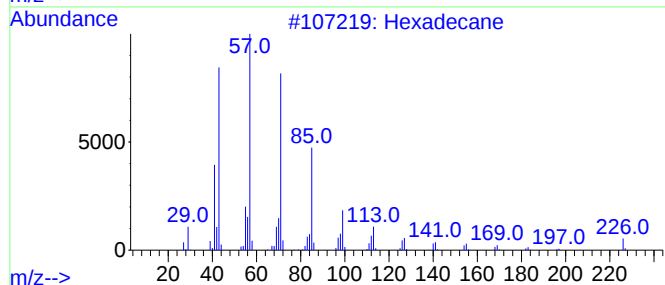
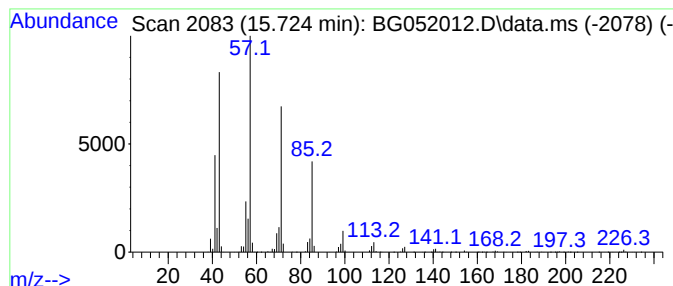
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Quant Title : SVOA CALIBRATION

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

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Peak Number 44 (DEL) Alkane: Straight-Chai... Concentration Rank 2

| R.T.   | EstConc      | Area    | Relative to ISTD | R.T.   |
|--------|--------------|---------|------------------|--------|
| 15.724 | 103.12 ng/ul | 2511470 | Acenaphthene-d10 | 14.895 |

| Hit# | of | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------|-----|---------|-------------|------|
| 1    |    |   | Hexadecane   | 226 | C16H34  | 000544-76-3 | 97   |
| 2    |    |   | Hexadecane   | 226 | C16H34  | 000544-76-3 | 96   |
| 3    |    |   | Tetradecane  | 198 | C14H30  | 000629-59-4 | 90   |
| 4    |    |   | Pentadecane  | 212 | C15H32  | 000629-62-9 | 90   |
| 5    |    |   | Nonadecane   | 268 | C19H40  | 000629-92-5 | 90   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011522\  
Data File : BG052012.D  
Acq On : 16 Jan 2022 1:50  
Operator : CG/JU  
Sample : N1132-04 10X  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLT5

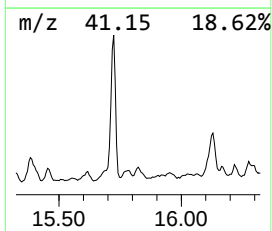
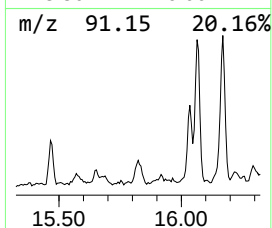
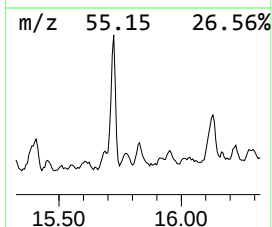
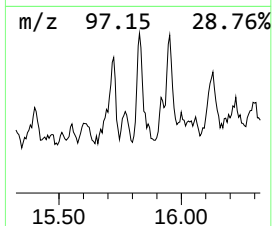
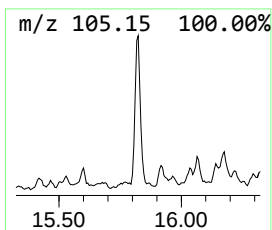
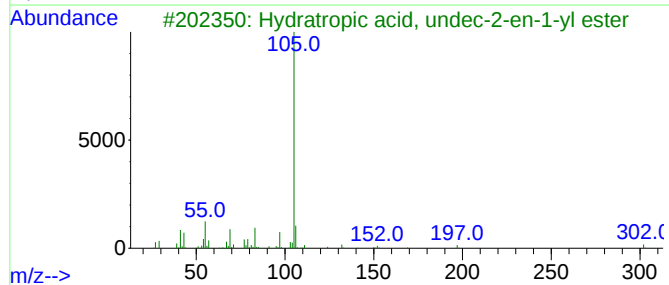
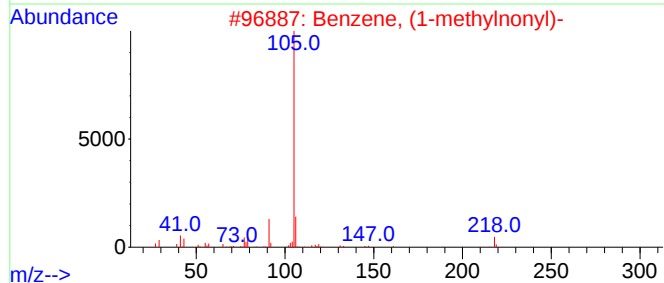
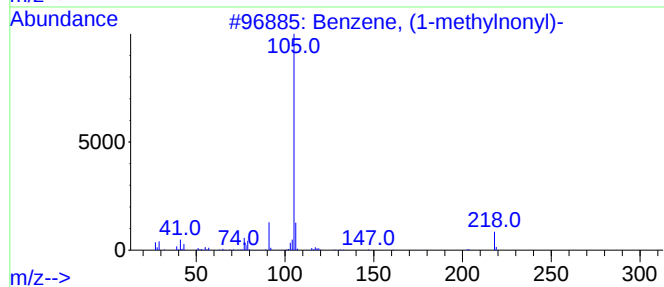
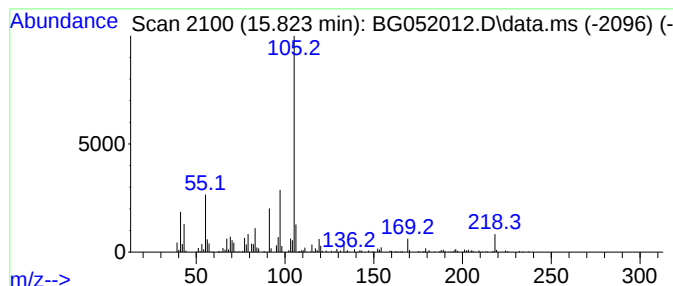
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Quant Title : SVOA CALIBRATION

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

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Peak Number 45 unknown-03 Concentration Rank 22

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 15.823 | 20.42 ng/ul | 497294 | Acenaphthene-d10 | 14.895 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Benzene, (1-methylnonyl)-           | 218 | C16H26   | 004537-13-7  | 46   |
| 2    |    |   | Benzene, (1-methylnonyl)-           | 218 | C16H26   | 004537-13-7  | 43   |
| 3    |    |   | Hydratropic acid, undec-2-en-1-y... | 302 | C20H30O2 | 1000406-96-0 | 43   |
| 4    |    |   | Benzene, (1-methylnonyl)-           | 218 | C16H26   | 004537-13-7  | 43   |
| 5    |    |   | 3-Methylbenzylamine, N,N-dihexyl-   | 289 | C20H35N  | 1010310-40-6 | 43   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011522\  
Data File : BG052012.D  
Acq On : 16 Jan 2022 1:50  
Operator : CG/JU  
Sample : N1132-04 10X  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLT5

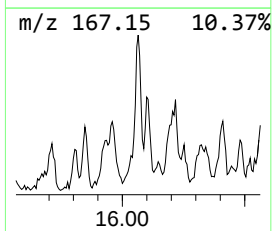
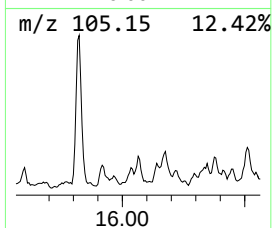
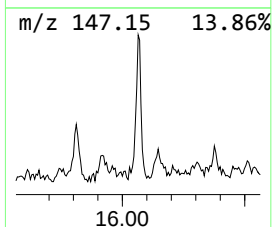
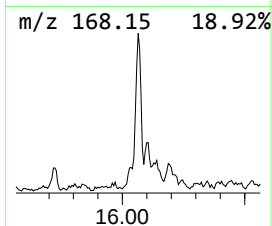
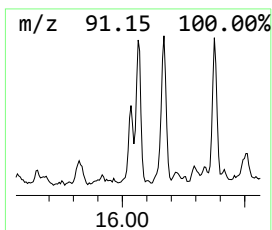
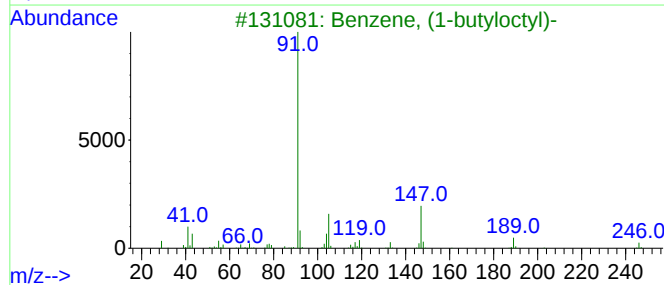
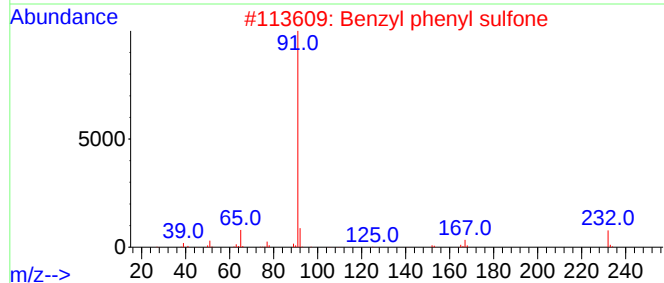
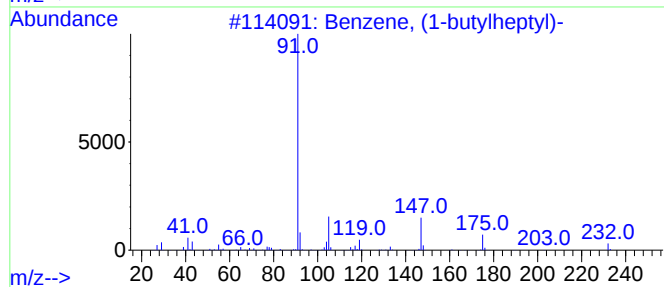
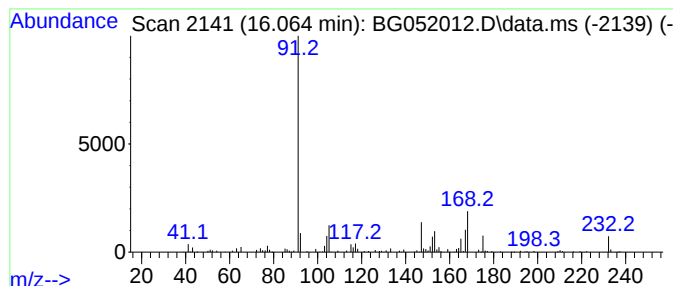
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Quant Title : SVOA CALIBRATION

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

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Peak Number 46 unknown-04 Concentration Rank 56

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 16.064 | 7.30 ng/ul | 177856 | Acenaphthene-d10 | 14.895 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm   | CAS#        | Qual |
|------|----|---|----------------------------|-----|-----------|-------------|------|
| 1    |    |   | Benzene, (1-butylheptyl)-  | 232 | C17H28    | 004537-15-9 | 27   |
| 2    |    |   | Benzyl phenyl sulfone      | 232 | C13H12O2S | 003112-88-7 | 27   |
| 3    |    |   | Benzene, (1-butyloctyl)-   | 246 | C18H30    | 002719-63-3 | 22   |
| 4    |    |   | Benzene, (1-butylhexyl)-   | 218 | C16H26    | 004537-11-5 | 22   |
| 5    |    |   | 4-Benzyloxybenzyl chloride | 232 | C14H13ClO | 000836-42-0 | 14   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011522\  
Data File : BG052012.D  
Acq On : 16 Jan 2022 1:50  
Operator : CG/JU  
Sample : N1132-04 10X  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

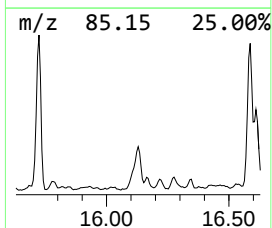
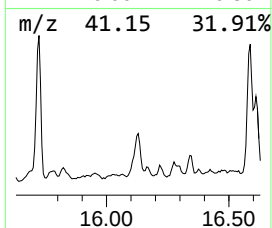
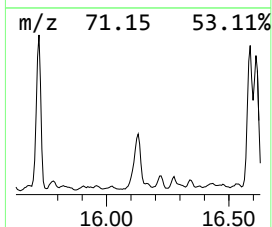
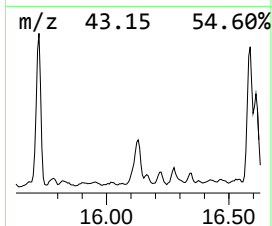
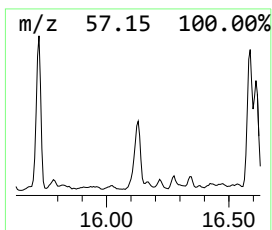
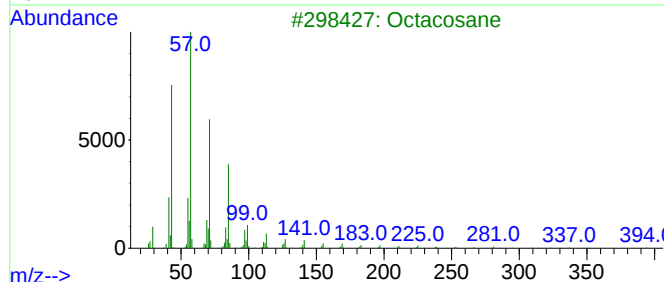
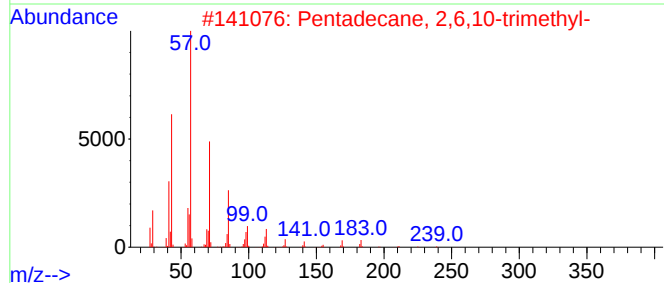
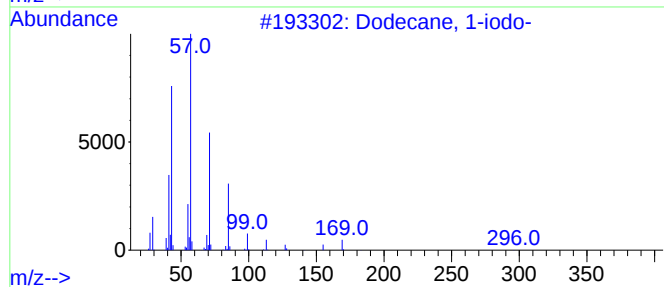
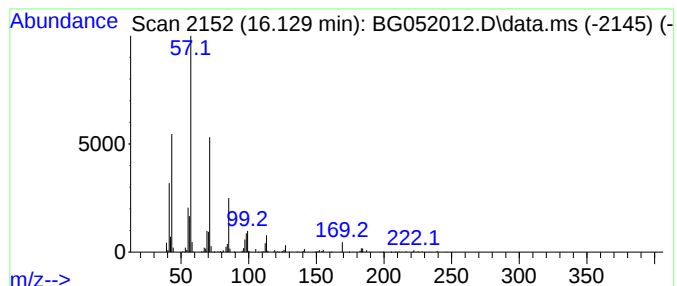
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BNA\_G  
ClientSampleId :  
BGLT5

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010622.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 47 Dodecane, 1-iodo- Concentration Rank 5

| R.T.      | EstConc                        | Area    | Relative to ISTD | R.T.        |      |
|-----------|--------------------------------|---------|------------------|-------------|------|
| 16.129    | 52.57 ng/ul                    | 1280310 | Acenaphthene-d10 | 14.895      |      |
| Hit# of 5 | Tentative ID                   | MW      | MolForm          | CAS#        | Qual |
| 1         | Dodecane, 1-iodo-              | 296     | C12H25I          | 004292-19-7 | 80   |
| 2         | Pentadecane, 2,6,10-trimethyl- | 254     | C18H38           | 003892-00-0 | 78   |
| 3         | Octacosane                     | 394     | C28H58           | 000630-02-4 | 72   |
| 4         | Undecane, 2,6-dimethyl-        | 184     | C13H28           | 017301-23-4 | 64   |
| 5         | Nonadecane                     | 268     | C19H40           | 000629-92-5 | 64   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011522\  
Data File : BG052012.D  
Acq On : 16 Jan 2022 1:50  
Operator : CG/JU  
Sample : N1132-04 10X  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLT5

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010622.M  
Quant Title : SVOA CALIBRATION

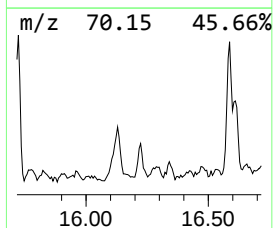
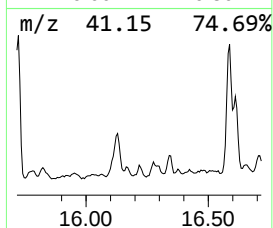
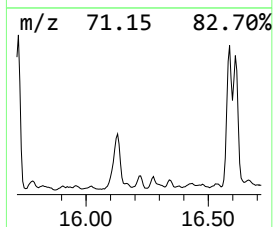
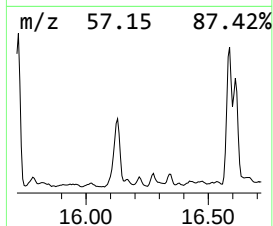
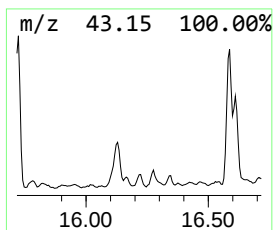
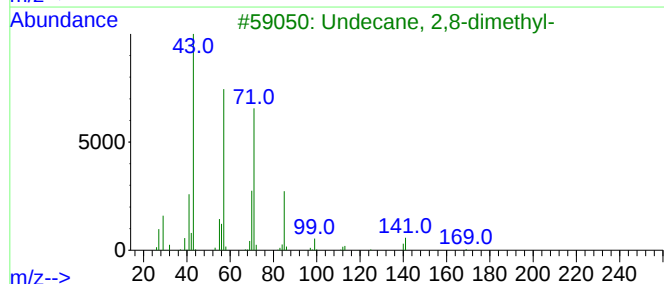
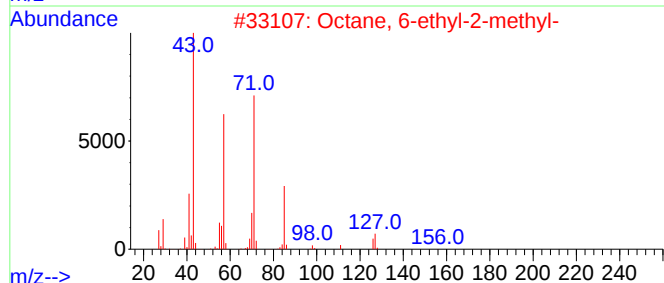
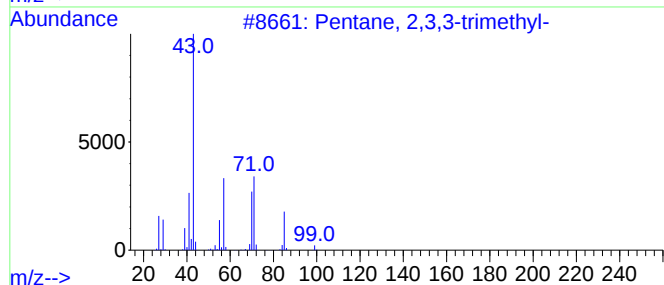
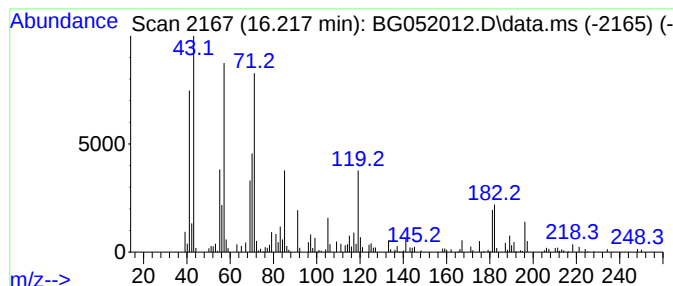
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 48 (DEL) Alkane: Straight-Chai... Concentration Rank 45

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 16.217 | 9.40 ng/ul | 229002 | Acenaphthene-d10 | 14.895 |

| Hit# | of | 5 | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|------|----|---|---------------------------|-----|---------|-------------|------|
| 1    |    |   | Pentane, 2,3,3-trimethyl- | 114 | C8H18   | 000560-21-4 | 47   |
| 2    |    |   | Octane, 6-ethyl-2-methyl- | 156 | C11H24  | 062016-19-7 | 43   |
| 3    |    |   | Undecane, 2,8-dimethyl-   | 184 | C13H28  | 017301-25-6 | 43   |
| 4    |    |   | Hexane, 3,3,4-trimethyl-  | 128 | C9H20   | 016747-31-2 | 43   |
| 5    |    |   | Hexane, 3,3-dimethyl-     | 114 | C8H18   | 000563-16-6 | 43   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011522\  
Data File : BG052012.D  
Acq On : 16 Jan 2022 1:50  
Operator : CG/JU  
Sample : N1132-04 10X  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLT5

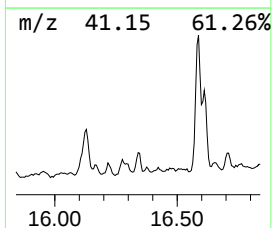
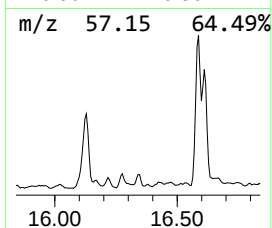
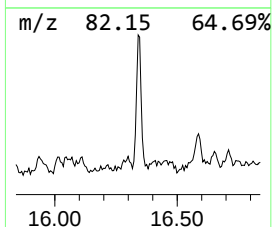
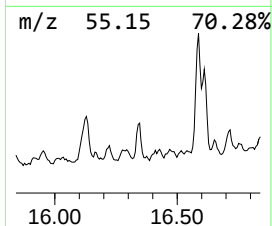
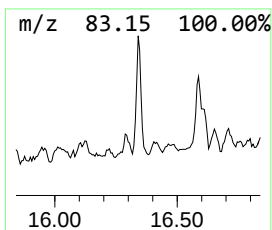
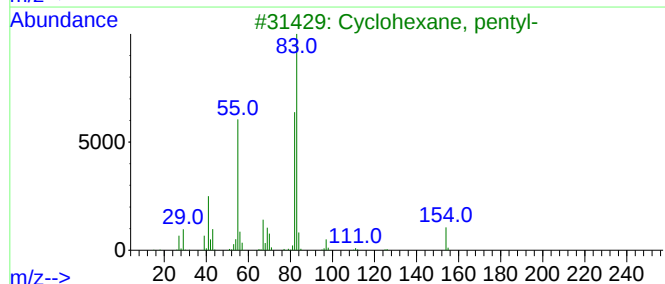
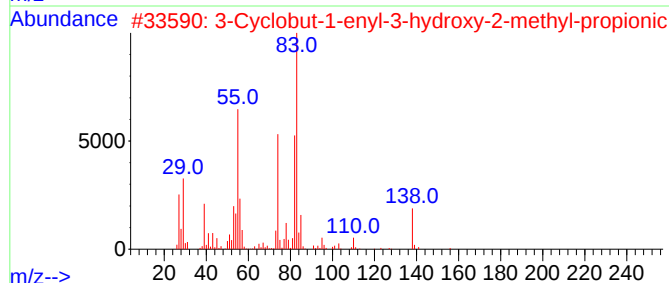
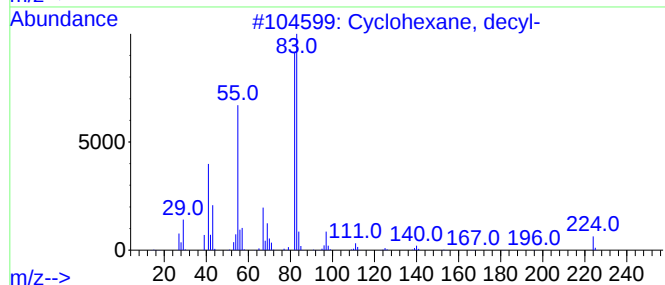
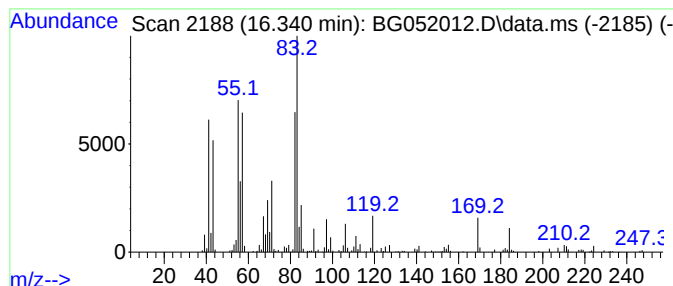
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Quant Title : SVOA CALIBRATION

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

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Peak Number 49 (DEL) Alkane: Cyclic16.340 Concentration Rank 47

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 16.340 | 9.25 ng/ul | 328547 | Phenanthrene-d10 | 17.645 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|---------|--------------|------|
| 1    |    |   | Cyclohexane, decyl-                 | 224 | C16H32  | 001795-16-0  | 53   |
| 2    |    |   | 3-Cyclobut-1-enyl-3-hydroxy-2-me... | 156 | C8H12O3 | 1000186-72-4 | 52   |
| 3    |    |   | Cyclohexane, pentyl-                | 154 | C11H22  | 004292-92-6  | 49   |
| 4    |    |   | Cyclohexane, decyl-                 | 224 | C16H32  | 001795-16-0  | 49   |
| 5    |    |   | Cyclohexane, octyl-                 | 196 | C14H28  | 001795-15-9  | 47   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011522\  
Data File : BG052012.D  
Acq On : 16 Jan 2022 1:50  
Operator : CG/JU  
Sample : N1132-04 10X  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLT5

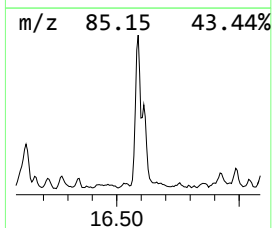
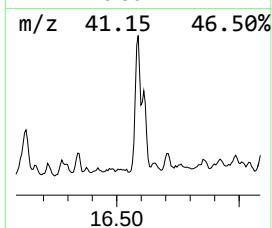
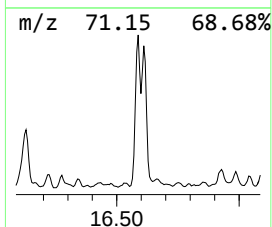
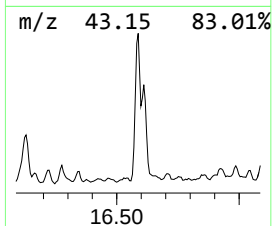
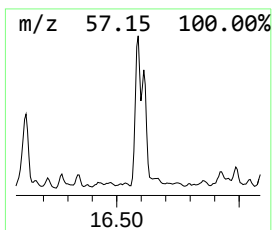
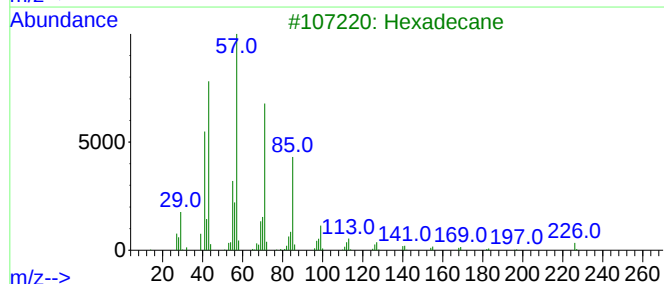
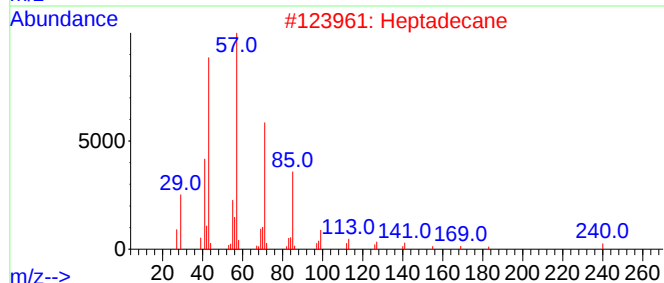
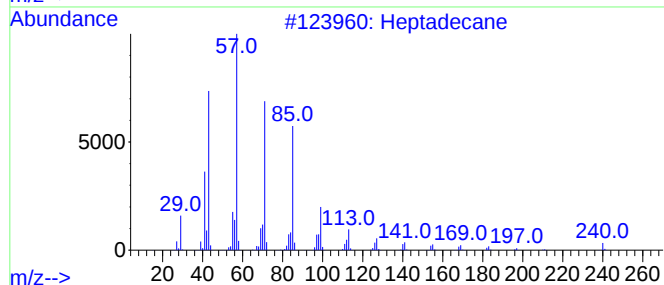
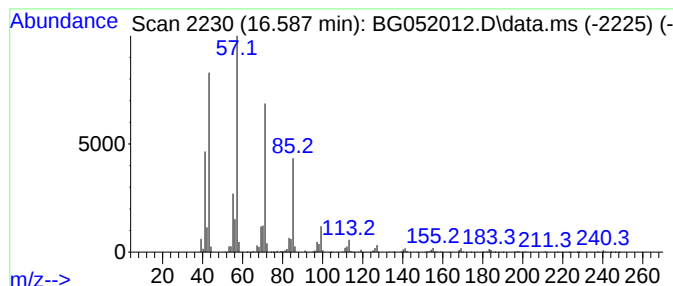
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Quant Title : SVOA CALIBRATION

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

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Peak Number 51 (DEL) Alkane: Straight-Chai... Concentration Rank 4

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 16.587 | 71.15 ng/ul | 2525970 | Phenanthrene-d10 | 17.645 |

| Hit# | of | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------|-----|---------|-------------|------|
| 1    |    |   | Heptadecane  | 240 | C17H36  | 000629-78-7 | 97   |
| 2    |    |   | Heptadecane  | 240 | C17H36  | 000629-78-7 | 93   |
| 3    |    |   | Hexadecane   | 226 | C16H34  | 000544-76-3 | 91   |
| 4    |    |   | Nonadecane   | 268 | C19H40  | 000629-92-5 | 91   |
| 5    |    |   | Hexadecane   | 226 | C16H34  | 000544-76-3 | 91   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011522\  
Data File : BG052012.D  
Acq On : 16 Jan 2022 1:50  
Operator : CG/JU  
Sample : N1132-04 10X  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLT5

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010622.M  
Quant Title : SVOA CALIBRATION

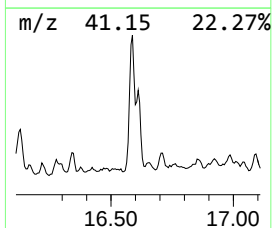
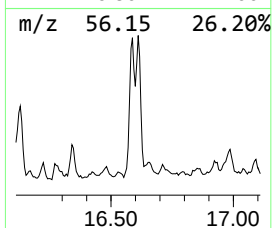
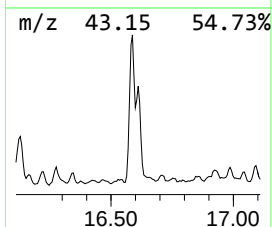
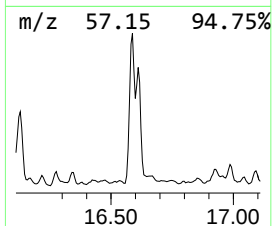
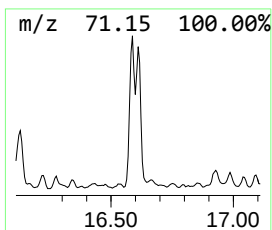
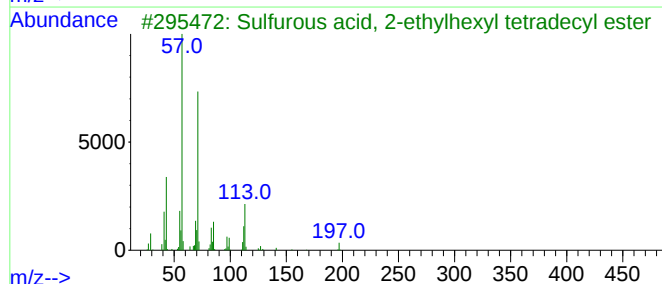
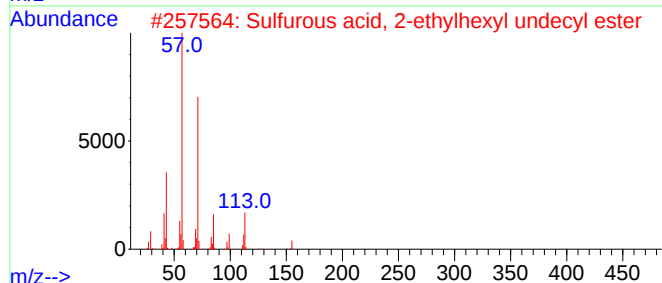
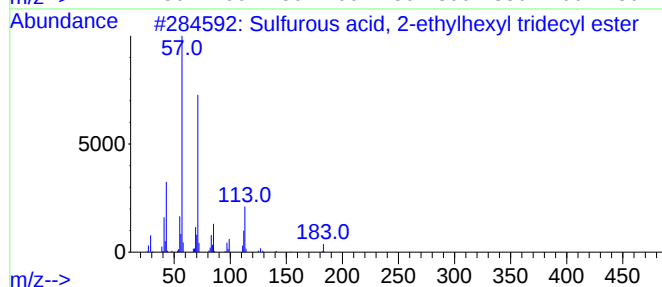
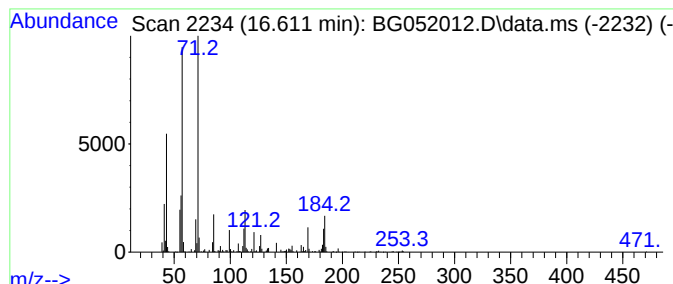
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 52 Sulfurous acid, 2-ethylhexy... Concentration Rank 9

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 16.611 | 44.85 ng/ul | 1592240 | Phenanthrene-d10 | 17.645 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Sulfurous acid, 2-ethylhexyl tri... | 376 | C21H44O3S | 1000309-19-6 | 72   |
| 2    |    |   | Sulfurous acid, 2-ethylhexyl und... | 348 | C19H40O3S | 1000309-19-4 | 64   |
| 3    |    |   | Sulfurous acid, 2-ethylhexyl tet... | 390 | C22H46O3S | 1000309-19-7 | 64   |
| 4    |    |   | Sulfurous acid, 2-ethylhexyl pen... | 264 | C13H28O3S | 1000309-18-9 | 58   |
| 5    |    |   | Octane, 2,6-dimethyl-               | 142 | C10H22    | 002051-30-1  | 53   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011522\  
Data File : BG052012.D  
Acq On : 16 Jan 2022 1:50  
Operator : CG/JU  
Sample : N1132-04 10X  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

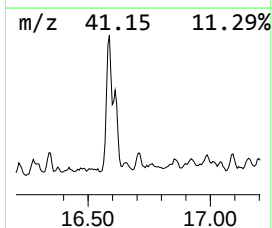
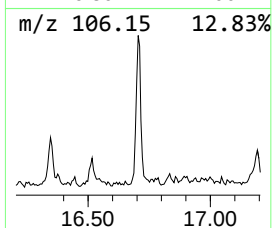
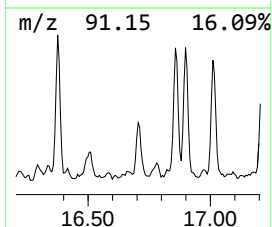
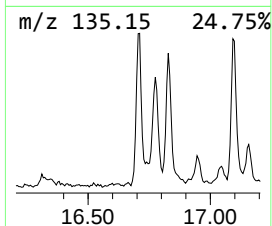
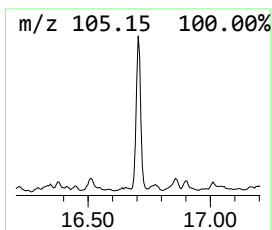
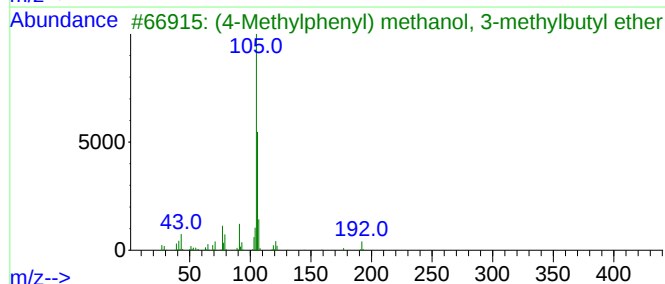
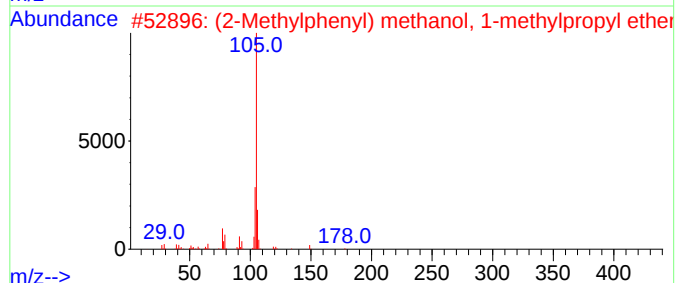
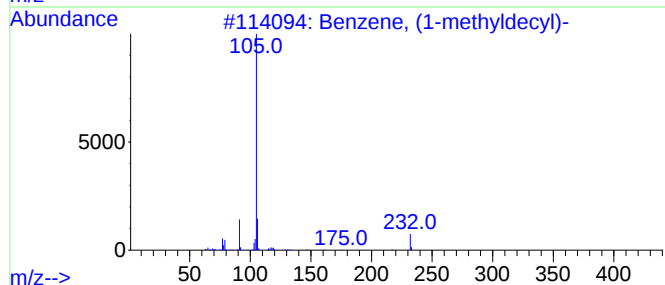
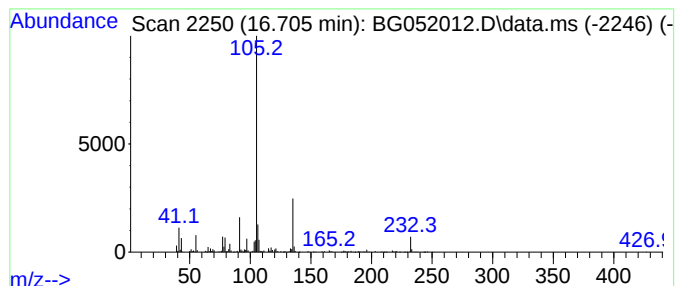
Instrument :  
BNA\_G  
ClientSampleId :  
BGLT5

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010622.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 53 Benzene, (1-methyldecyl)- Concentration Rank 23

| R.T.      | EstConc                             | Area   | Relative to ISTD |             |              | R.T.   |
|-----------|-------------------------------------|--------|------------------|-------------|--------------|--------|
| 16.705    | 20.03 ng/ul                         | 711256 | Phenanthrene-d10 |             |              | 17.645 |
| Hit# of 5 | Tentative ID                        |        | MW               | MolForm     | CAS#         | Qual   |
| 1         | Benzene, (1-methyldecyl)-           |        | 232              | C17H28      | 004536-88-3  | 60     |
| 2         | (2-Methylphenyl) methanol, 1-met... |        | 178              | C12H18O     | 1000374-44-3 | 43     |
| 3         | (4-Methylphenyl) methanol, 3-met... |        | 192              | C13H20O     | 1000374-65-7 | 38     |
| 4         | Succinic acid, 2,4,6-trichloroph... |        | 400              | C18H15Cl3O4 | 1000389-75-5 | 38     |
| 5         | Benzene, (1-methylpropyl)-          |        | 134              | C10H14      | 000135-98-8  | 38     |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011522\  
Data File : BG052012.D  
Acq On : 16 Jan 2022 1:50  
Operator : CG/JU  
Sample : N1132-04 10X  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLT5

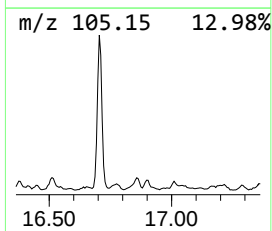
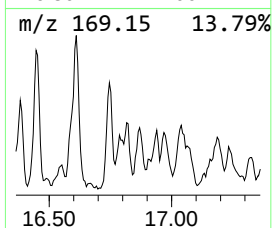
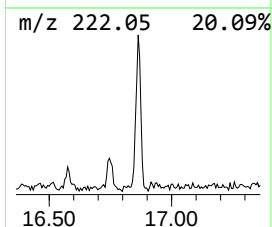
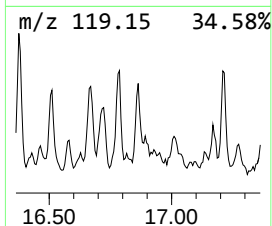
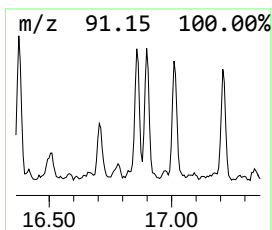
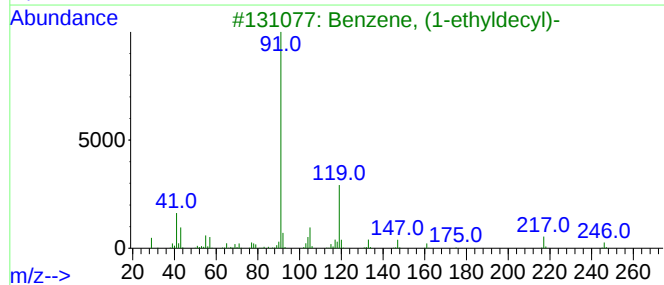
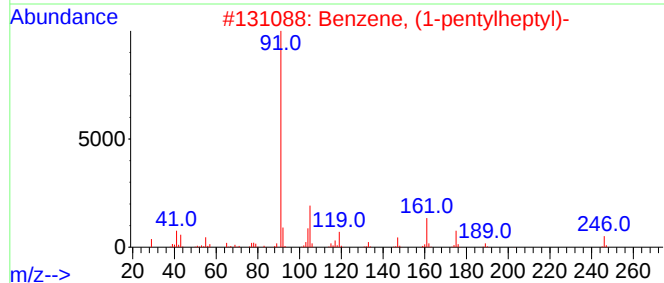
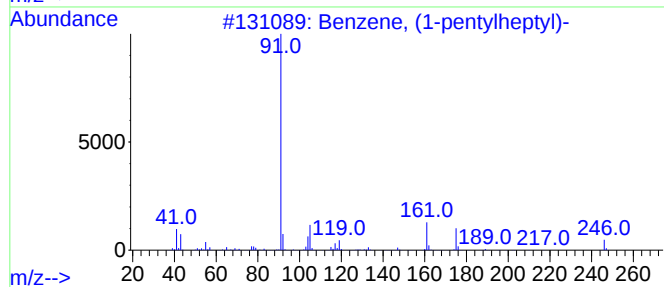
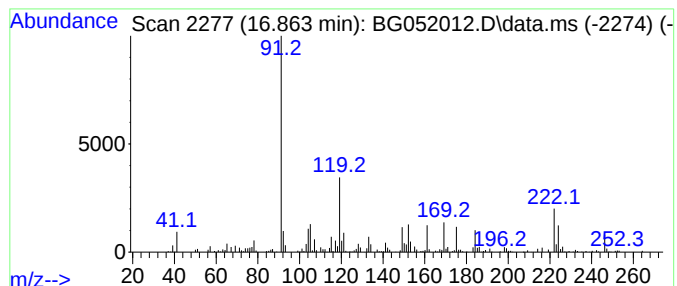
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Quant Title : SVOA CALIBRATION

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

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Peak Number 54 unknown-05 Concentration Rank 54

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 16.863 | 8.02 ng/ul | 284619 | Phenanthrene-d10 | 17.645 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|------------|--------------|------|
| 1    |    |   | Benzene, (1-pentylheptyl)-          | 246 | C18H30     | 002719-62-2  | 11   |
| 2    |    |   | Benzene, (1-pentylheptyl)-          | 246 | C18H30     | 002719-62-2  | 11   |
| 3    |    |   | Benzene, (1-ethyldecyl)-            | 246 | C18H30     | 002400-00-2  | 10   |
| 4    |    |   | Benzene, (1-ethyldecyl)-            | 246 | C18H30     | 002400-00-2  | 10   |
| 5    |    |   | 2-phenylbutyric acid, 2,2,2-trif... | 246 | C12H13F3O2 | 1000376-55-3 | 10   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011522\  
Data File : BG052012.D  
Acq On : 16 Jan 2022 1:50  
Operator : CG/JU  
Sample : N1132-04 10X  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLT5

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010622.M  
Quant Title : SVOA CALIBRATION

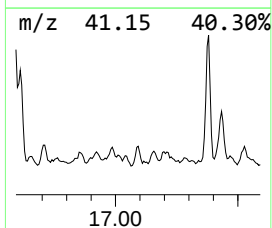
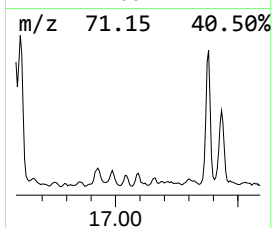
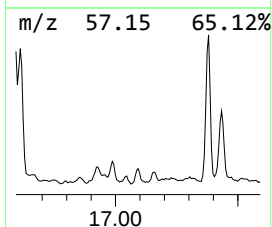
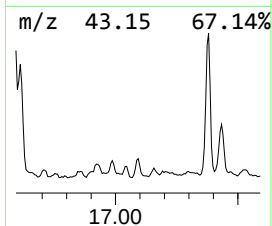
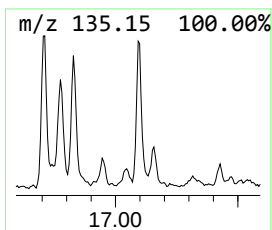
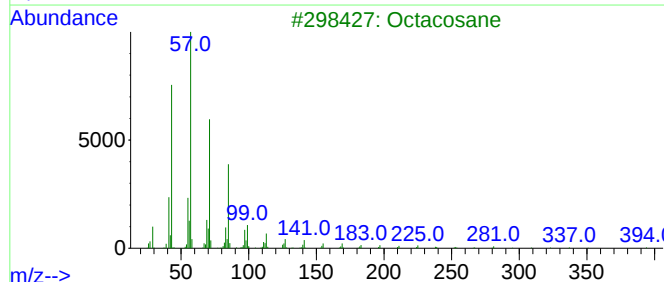
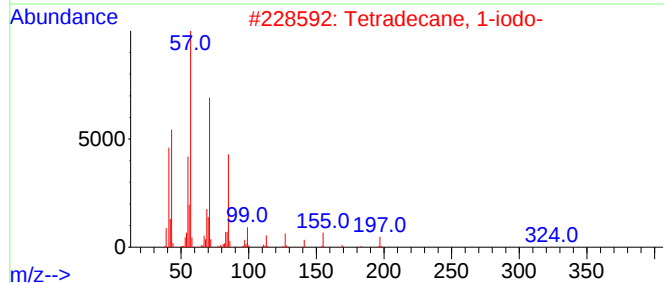
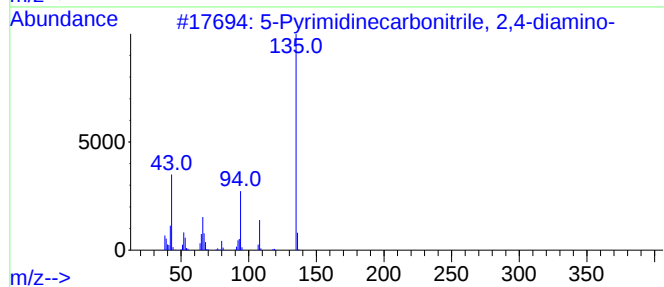
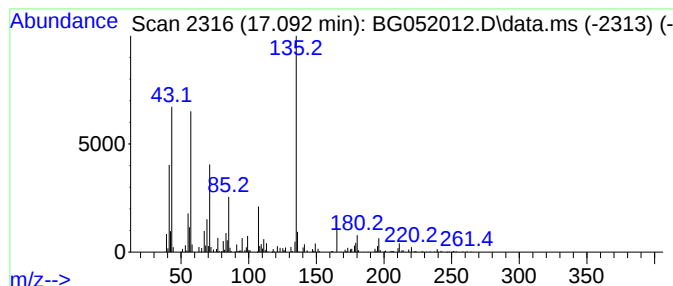
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 55 unknown-06 Concentration Rank 51

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 17.092 | 8.33 ng/ul | 295882 | Phenanthrene-d10 | 17.645 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|---------|--------------|------|
| 1    |    |   | 5-Pyrimidinecarbonitrile, 2,4-di... | 135 | C5H5N5  | 1010338-13-5 | 38   |
| 2    |    |   | Tetradecane, 1-iodo-                | 324 | C14H29I | 019218-94-1  | 18   |
| 3    |    |   | Octacosane                          | 394 | C28H58  | 000630-02-4  | 18   |
| 4    |    |   | Nonadecane                          | 268 | C19H40  | 000629-92-5  | 18   |
| 5    |    |   | Tetratetracontane                   | 619 | C44H90  | 007098-22-8  | 18   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011522\  
Data File : BG052012.D  
Acq On : 16 Jan 2022 1:50  
Operator : CG/JU  
Sample : N1132-04 10X  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

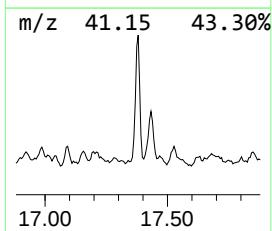
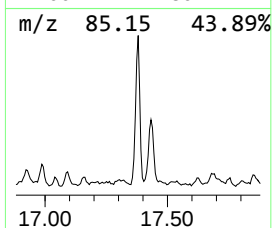
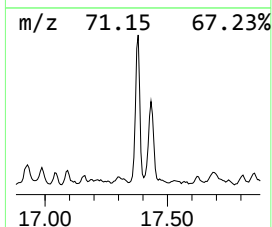
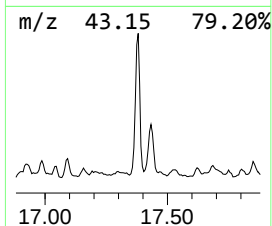
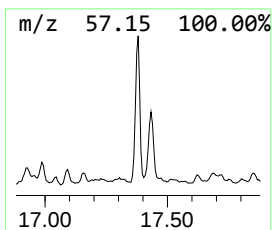
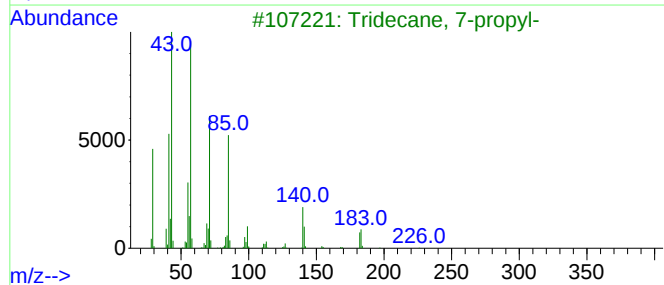
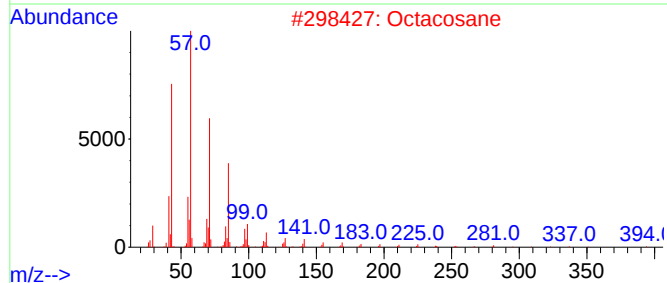
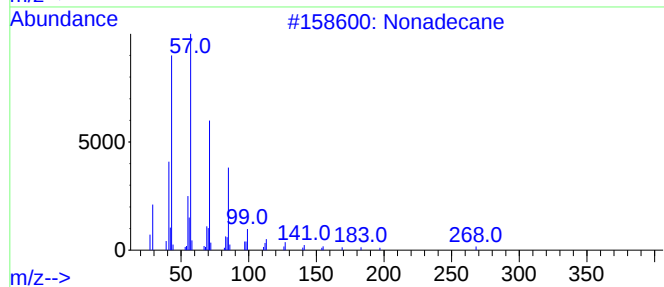
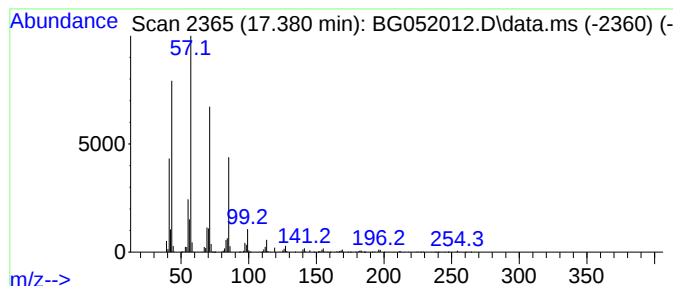
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BNA\_G  
ClientSampleId :  
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Quant Title : SVOA CALIBRATION

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

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Peak Number 56 (DEL) Alkane: Straight-Chai... Concentration Rank 7

| R.T.      | EstConc              | Area    | Relative to ISTD |         | R.T.        |      |
|-----------|----------------------|---------|------------------|---------|-------------|------|
| 17.380    | 48.58 ng/ul          | 1724820 | Phenanthrene-d10 |         | 17.645      |      |
| Hit# of 5 | Tentative ID         |         | MW               | MolForm | CAS#        | Qual |
| 1         | Nonadecane           |         | 268              | C19H40  | 000629-92-5 | 91   |
| 2         | Octacosane           |         | 394              | C28H58  | 000630-02-4 | 90   |
| 3         | Tridecane, 7-propyl- |         | 226              | C16H34  | 055045-09-5 | 90   |
| 4         | Hexadecane, 1-iodo-  |         | 352              | C16H33I | 000544-77-4 | 86   |
| 5         | Nonane, 5-butyl-     |         | 184              | C13H28  | 017312-63-9 | 81   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011522\  
Data File : BG052012.D  
Acq On : 16 Jan 2022 1:50  
Operator : CG/JU  
Sample : N1132-04 10X  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

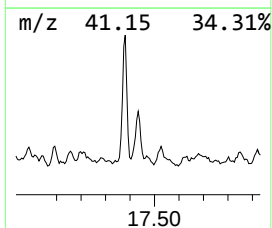
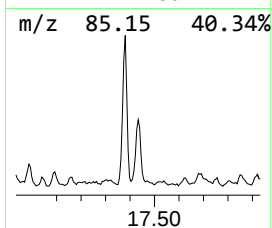
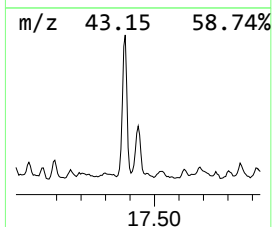
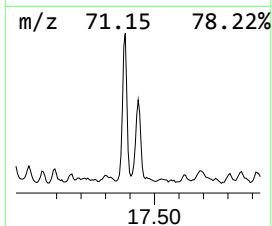
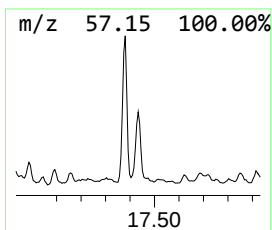
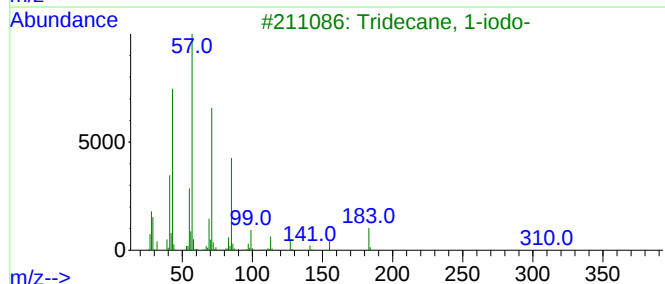
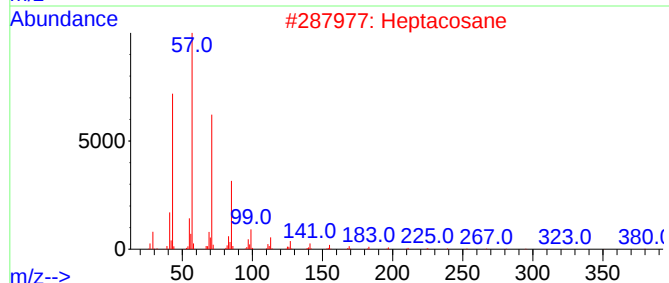
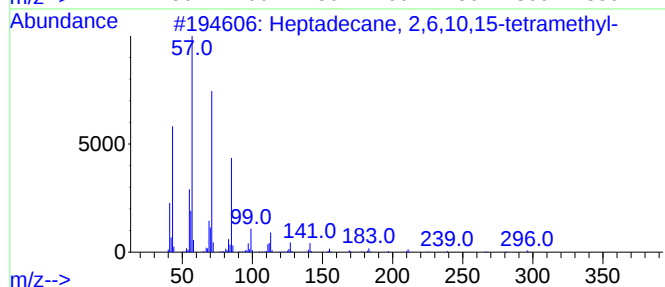
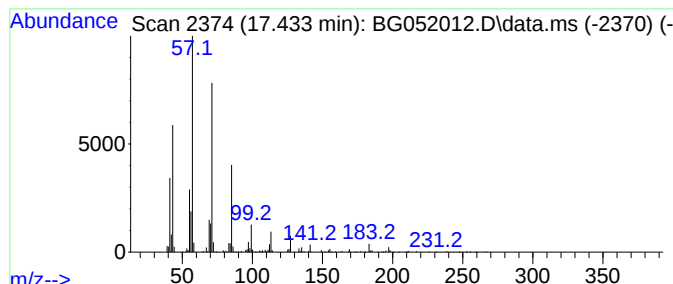
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ClientSampleId :  
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Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010622.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 57 (DEL) Alkane: Straight-Chai... Concentration Rank 18

| R.T.      | EstConc                             | Area   | Relative to ISTD |         | R.T.        |      |
|-----------|-------------------------------------|--------|------------------|---------|-------------|------|
| 17.433    | 22.76 ng/ul                         | 807992 | Phenanthrene-d10 |         | 17.645      |      |
| Hit# of 5 | Tentative ID                        |        | MW               | MolForm | CAS#        | Qual |
| 1         | Heptadecane, 2,6,10,15-tetramethyl- |        | 296              | C21H44  | 054833-48-6 | 91   |
| 2         | Heptacosane                         |        | 380              | C27H56  | 000593-49-7 | 90   |
| 3         | Tridecane, 1-iodo-                  |        | 310              | C13H27I | 035599-77-0 | 90   |
| 4         | Tridecane, 4-methyl-                |        | 198              | C14H30  | 026730-12-1 | 87   |
| 5         | Tetradecane, 1-iodo-                |        | 324              | C14H29I | 019218-94-1 | 86   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011522\  
Data File : BG052012.D  
Acq On : 16 Jan 2022 1:50  
Operator : CG/JU  
Sample : N1132-04 10X  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLT5

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010622.M  
Quant Title : SVOA CALIBRATION

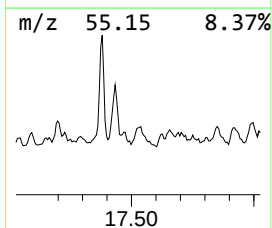
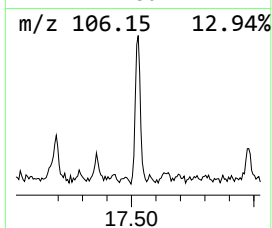
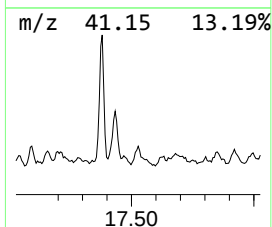
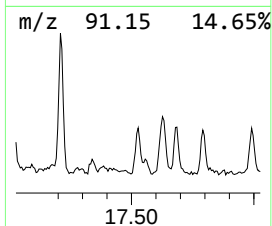
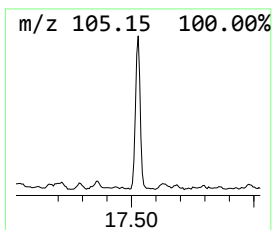
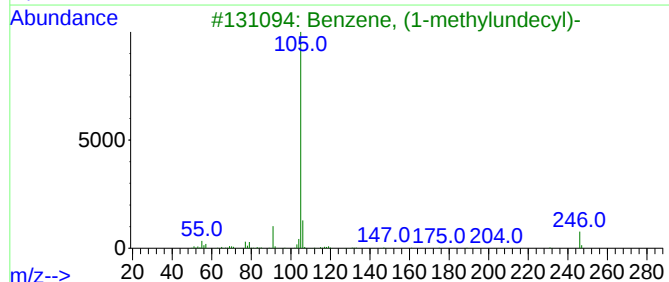
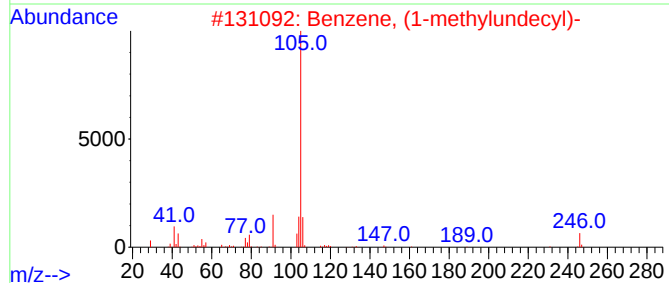
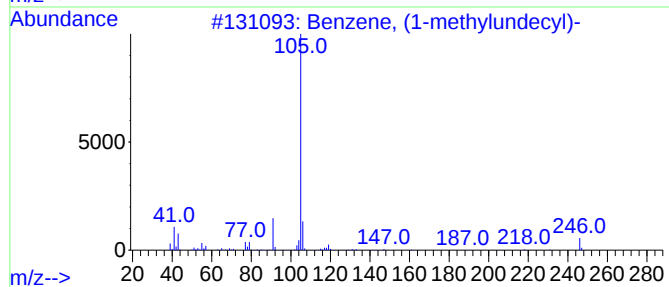
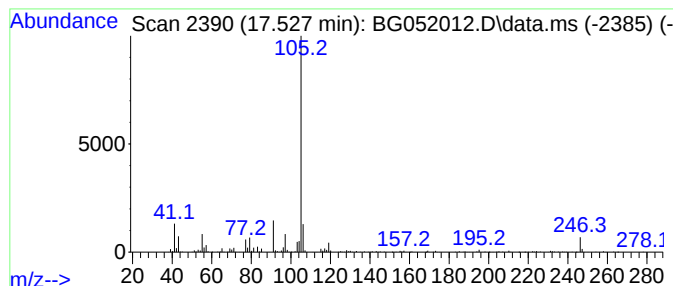
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 58 Benzene, (1-methylundecyl)- Concentration Rank 24

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 17.527 | 20.00 ng/ul | 710072 | Phenanthrene-d10 | 17.645 |

| Hit# | of | 5 | Tentative ID                     | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, (1-methylundecyl)-      | 246 | C18H30  | 002719-61-1 | 91   |
| 2    |    |   | Benzene, (1-methylundecyl)-      | 246 | C18H30  | 002719-61-1 | 89   |
| 3    |    |   | Benzene, (1-methylundecyl)-      | 246 | C18H30  | 002719-61-1 | 64   |
| 4    |    |   | Benzene, (1,3,3-trimethylnonyl)- | 246 | C18H30  | 054986-44-6 | 64   |
| 5    |    |   | Benzene, (1-methyldecyl)-        | 232 | C17H28  | 004536-88-3 | 58   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011522\  
Data File : BG052012.D  
Acq On : 16 Jan 2022 1:50  
Operator : CG/JU  
Sample : N1132-04 10X  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

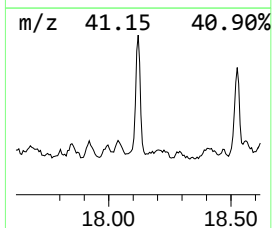
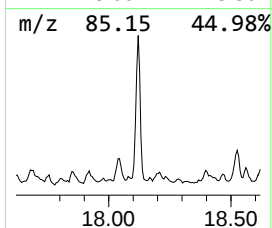
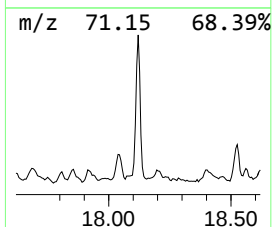
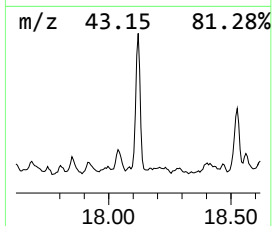
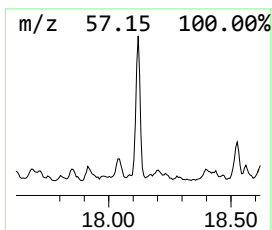
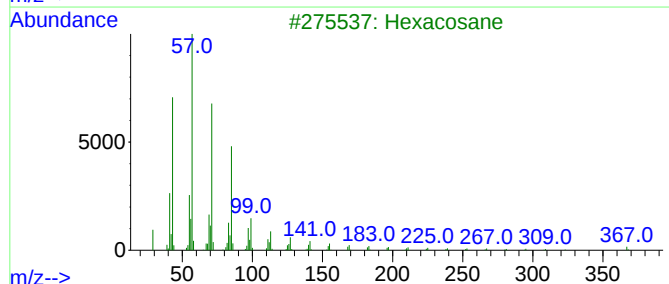
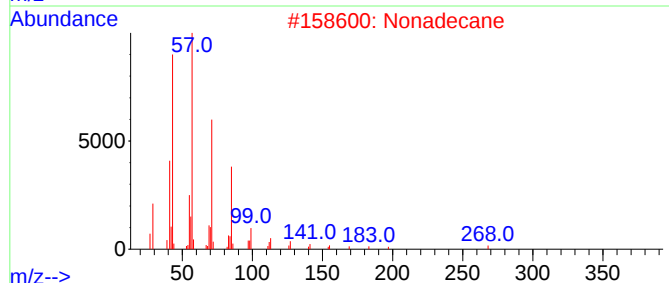
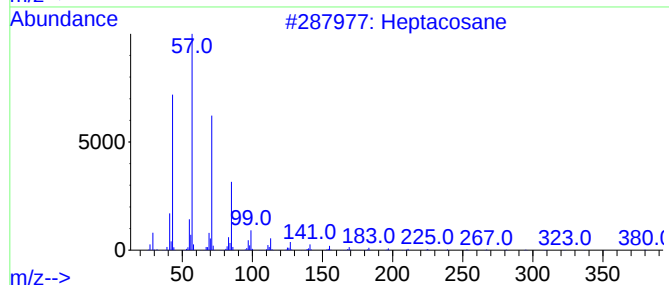
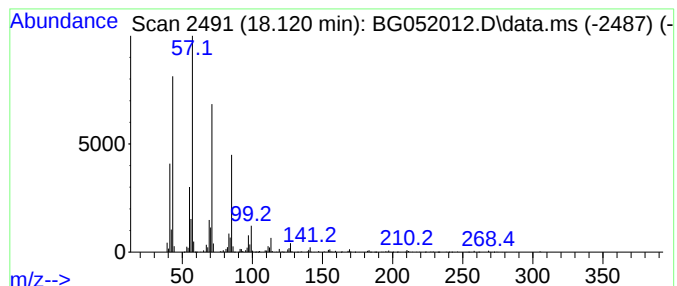
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ClientSampleId :  
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Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010622.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 59 (DEL) Alkane: Straight-Chai... Concentration Rank 8

| R.T.      | EstConc            | Area    | Relative to ISTD |         | R.T.        |      |
|-----------|--------------------|---------|------------------|---------|-------------|------|
| 18.120    | 46.67 ng/ul        | 1657040 | Phenanthrene-d10 |         | 17.645      |      |
| Hit# of 5 | Tentative ID       |         | MW               | MolForm | CAS#        | Qual |
| 1         | Heptacosane        |         | 380              | C27H56  | 000593-49-7 | 91   |
| 2         | Nonadecane         |         | 268              | C19H40  | 000629-92-5 | 91   |
| 3         | Hexacosane         |         | 366              | C26H54  | 000630-01-3 | 91   |
| 4         | Heneicosane        |         | 296              | C21H44  | 000629-94-7 | 91   |
| 5         | Tridecane, 1-iodo- |         | 310              | C13H27I | 035599-77-0 | 90   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011522\  
Data File : BG052012.D  
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Operator : CG/JU  
Sample : N1132-04 10X  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

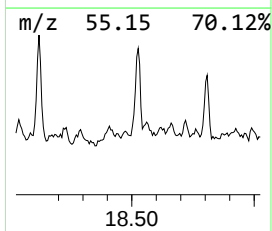
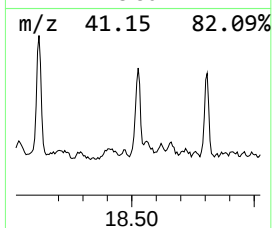
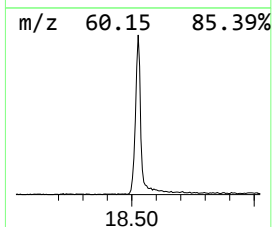
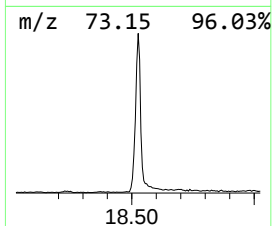
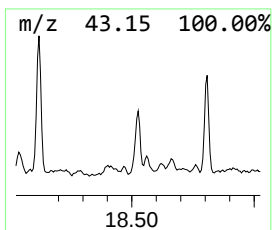
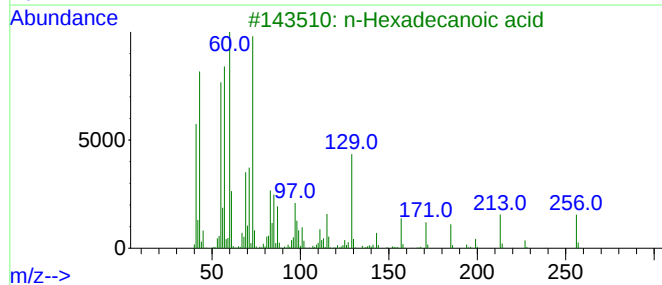
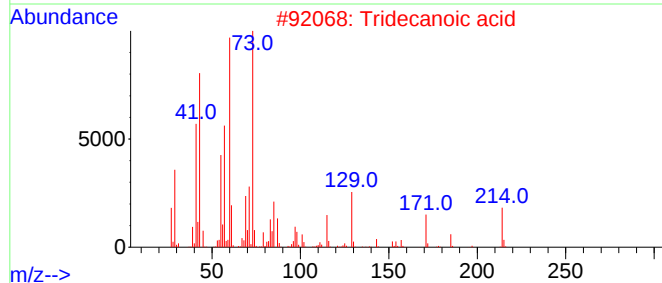
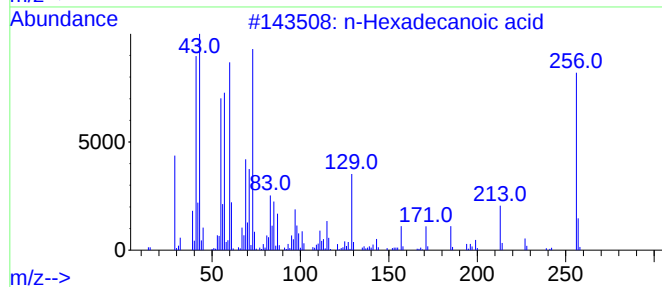
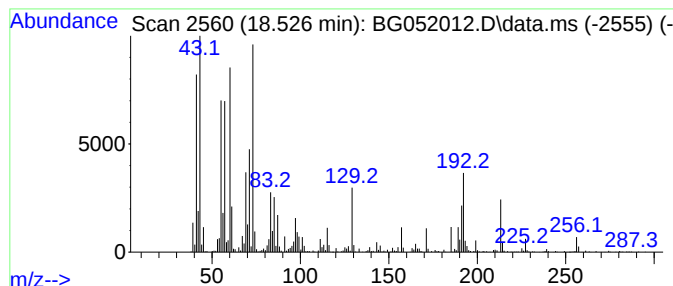
Instrument :  
BNA\_G  
ClientSampleId :  
BGLT5

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010622.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 60 n-Hexadecanoic acid Concentration Rank 10

| R.T.      | EstConc             | Area    | Relative to ISTD |          | R.T.        |      |
|-----------|---------------------|---------|------------------|----------|-------------|------|
| 18.526    | 40.11 ng/ul         | 1424070 | Phenanthrene-d10 |          | 17.645      |      |
| Hit# of 5 | Tentative ID        |         | MW               | MolForm  | CAS#        | Qual |
| 1         | n-Hexadecanoic acid |         | 256              | C16H32O2 | 000057-10-3 | 96   |
| 2         | Tridecanoic acid    |         | 214              | C13H26O2 | 000638-53-9 | 86   |
| 3         | n-Hexadecanoic acid |         | 256              | C16H32O2 | 000057-10-3 | 70   |
| 4         | Pentadecanoic acid  |         | 242              | C15H30O2 | 001002-84-2 | 58   |
| 5         | n-Hexadecanoic acid |         | 256              | C16H32O2 | 000057-10-3 | 58   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011522\  
Data File : BG052012.D  
Acq On : 16 Jan 2022 1:50  
Operator : CG/JU  
Sample : N1132-04 10X  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLT5

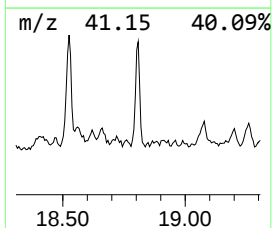
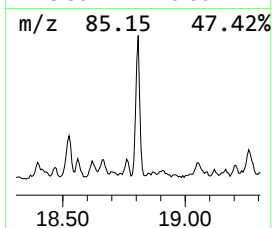
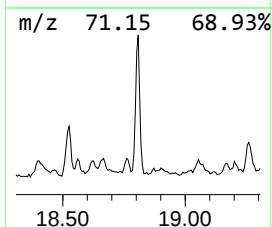
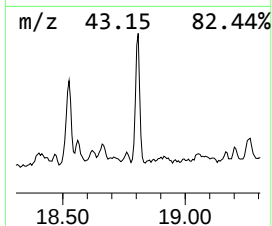
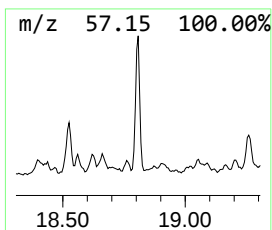
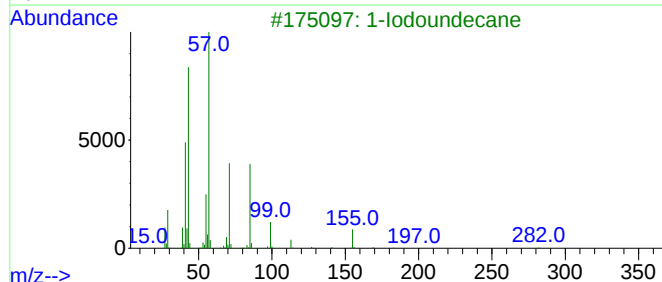
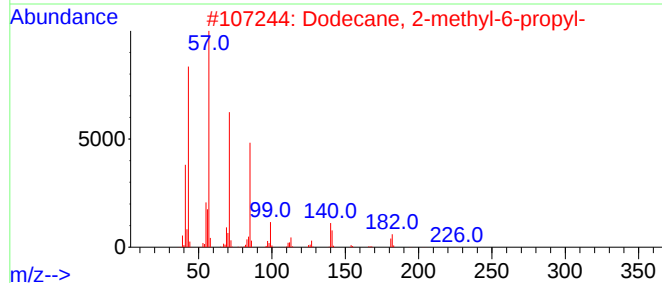
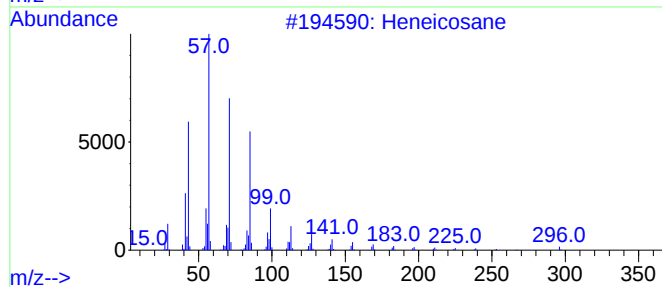
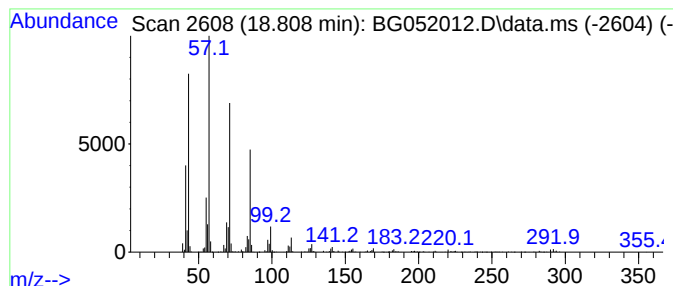
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Quant Title : SVOA CALIBRATION

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

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Peak Number 61 (DEL) Alkane: Straight-Chai... Concentration Rank 15

| R.T.      | EstConc                      | Area   | Relative to ISTD |         | R.T.        |      |
|-----------|------------------------------|--------|------------------|---------|-------------|------|
| 18.808    | 27.89 ng/ul                  | 990292 | Phenanthrene-d10 |         | 17.645      |      |
| Hit# of 5 | Tentative ID                 |        | MW               | MolForm | CAS#        | Qual |
| 1         | Heneicosane                  |        | 296              | C21H44  | 000629-94-7 | 96   |
| 2         | Dodecane, 2-methyl-6-propyl- |        | 226              | C16H34  | 055045-08-4 | 94   |
| 3         | 1-Iodoundecane               |        | 282              | C11H23I | 004282-44-4 | 93   |
| 4         | Nonadecane                   |        | 268              | C19H40  | 000629-92-5 | 91   |
| 5         | Pentadecane                  |        | 212              | C15H32  | 000629-62-9 | 90   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011522\  
Data File : BG052012.D  
Acq On : 16 Jan 2022 1:50  
Operator : CG/JU  
Sample : N1132-04 10X  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLT5

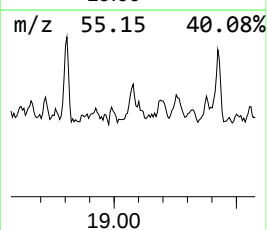
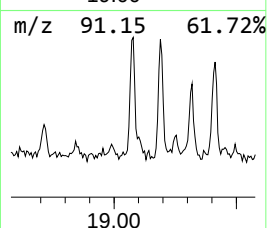
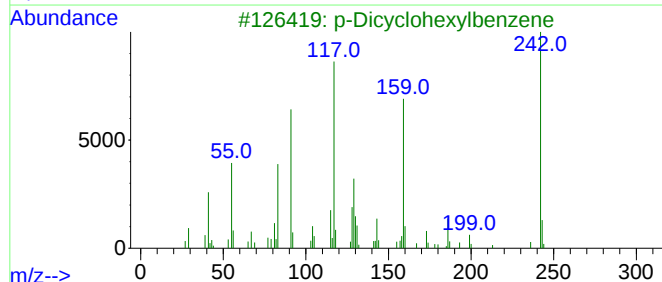
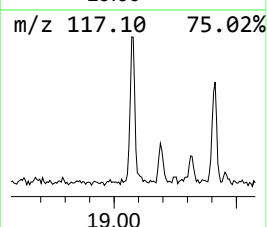
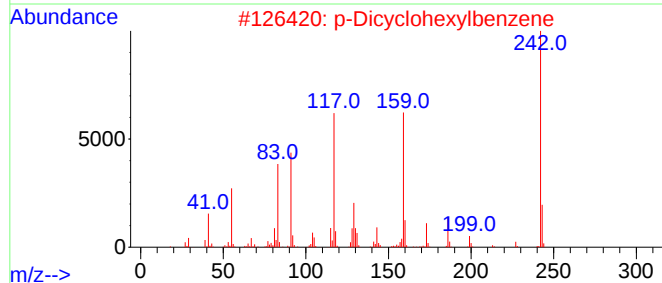
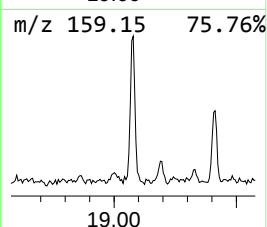
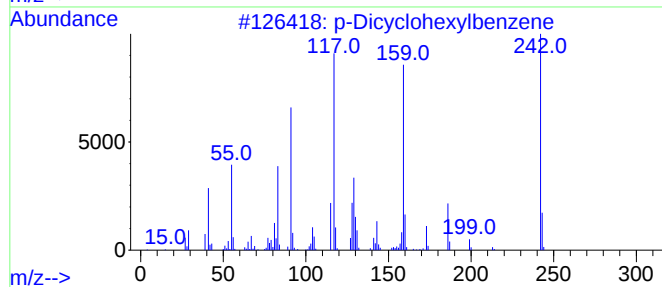
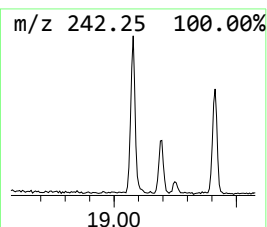
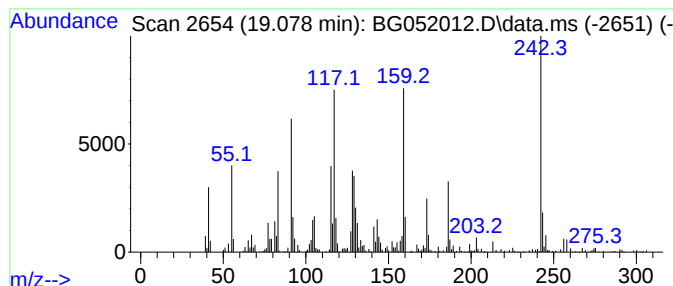
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Quant Title : SVOA CALIBRATION

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

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Peak Number 62 p-Dicyclohexylbenzene Concentration Rank 34

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 19.078 | 13.62 ng/ul | 483538 | Phenanthrene-d10 | 17.645 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | p-Dicyclohexylbenzene               | 242 | C18H26   | 001087-02-1 | 98   |
| 2    |    |   | p-Dicyclohexylbenzene               | 242 | C18H26   | 001087-02-1 | 89   |
| 3    |    |   | p-Dicyclohexylbenzene               | 242 | C18H26   | 001087-02-1 | 89   |
| 4    |    |   | Bicyclohexyl, 4-phenyl-             | 242 | C18H26   | 020273-27-2 | 55   |
| 5    |    |   | 2-(4-(2-Methylallyl)phenyl)propa... | 204 | C13H16O2 | 131734-08-2 | 49   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011522\  
Data File : BG052012.D  
Acq On : 16 Jan 2022 1:50  
Operator : CG/JU  
Sample : N1132-04 10X  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLT5

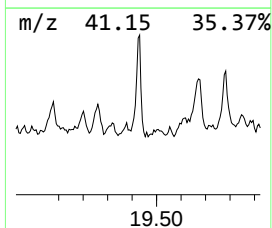
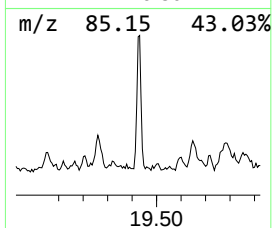
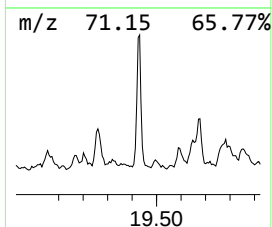
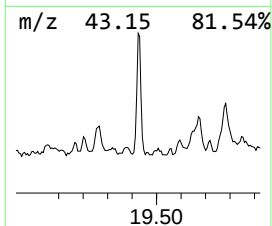
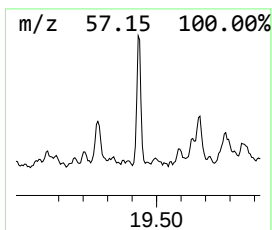
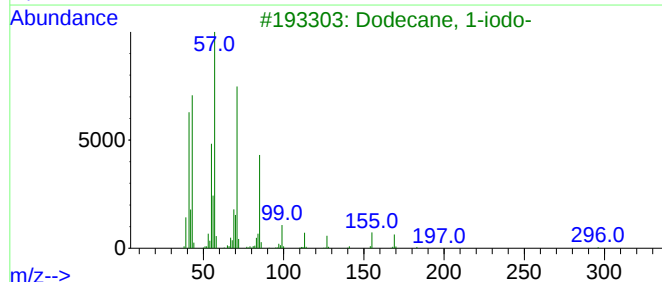
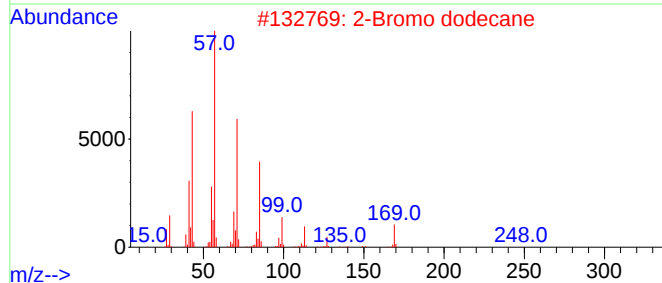
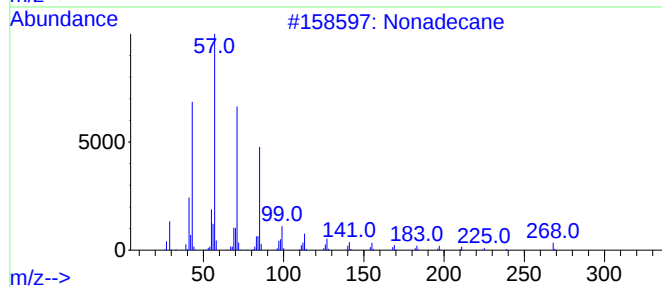
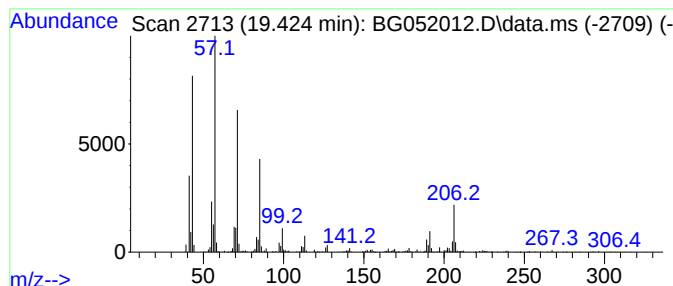
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Quant Title : SVOA CALIBRATION

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

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Peak Number 63 (DEL) Alkane: Straight-Chai... Concentration Rank 14

| R.T.      | EstConc               | Area    | Relative to ISTD |          | R.T.        |      |
|-----------|-----------------------|---------|------------------|----------|-------------|------|
| 19.424    | 29.96 ng/ul           | 1063670 | Phenanthrene-d10 |          | 17.645      |      |
| Hit# of 5 | Tentative ID          |         | MW               | MolForm  | CAS#        | Qual |
| 1         | Nonadecane            |         | 268              | C19H40   | 000629-92-5 | 90   |
| 2         | 2-Bromo dodecane      |         | 248              | C12H25Br | 013187-99-0 | 70   |
| 3         | Dodecane, 1-iodo-     |         | 296              | C12H25I  | 004292-19-7 | 62   |
| 4         | Nonane, 3,7-dimethyl- |         | 156              | C11H24   | 017302-32-8 | 58   |
| 5         | Tridecane, 1-iodo-    |         | 310              | C13H27I  | 035599-77-0 | 58   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011522\  
Data File : BG052012.D  
Acq On : 16 Jan 2022 1:50  
Operator : CG/JU  
Sample : N1132-04 10X  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLT5

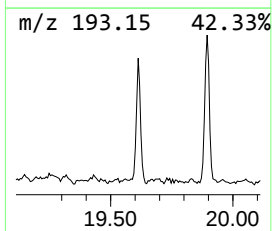
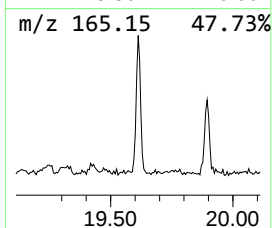
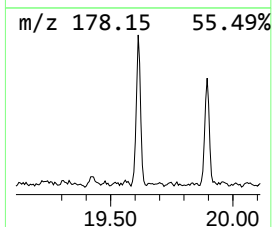
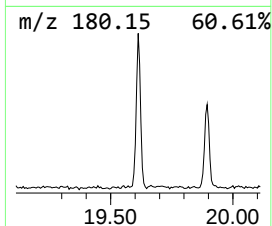
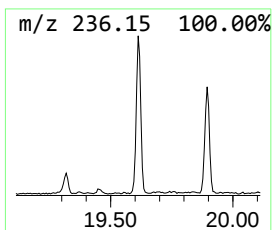
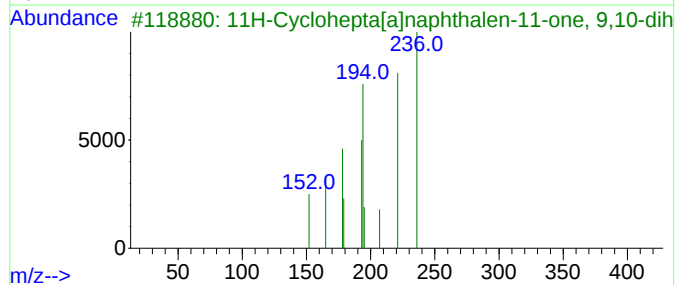
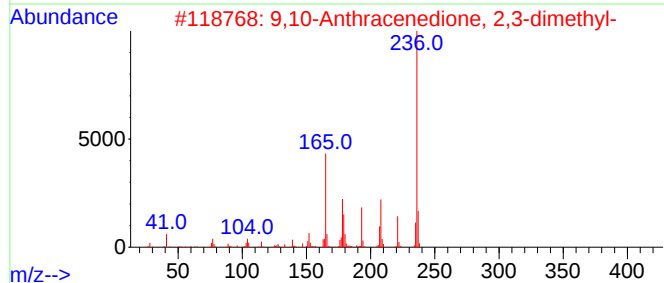
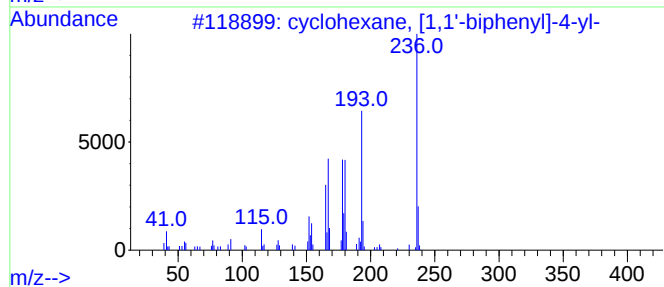
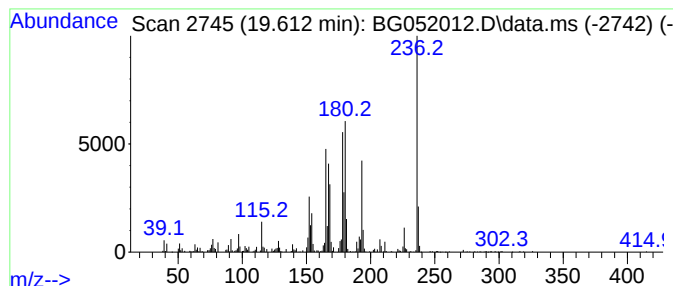
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Quant Title : SVOA CALIBRATION

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

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Peak Number 64 cyclohexane, [1,1'-biphenyl]... Concentration Rank 32

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|--------|------------------|--------------|------|
| 19.612    | 14.29 ng/ul                         | 507462 | Phenanthrene-d10 | 17.645       |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#         | Qual |
| 1         | cyclohexane, [1,1'-biphenyl]-4-yl-  | 236    | C18H20           | 1000401-12-4 | 76   |
| 2         | 9,10-Anthracenedione, 2,3-dimethyl- | 236    | C16H12O2         | 006531-35-7  | 43   |
| 3         | 11H-Cyclohepta[a]naphthalen-11-o... | 236    | C17H16O          | 058111-76-5  | 43   |
| 4         | 9,10-Anthracenedione, 2,7-dimethyl- | 236    | C16H12O2         | 003286-01-9  | 38   |
| 5         | 6-Ethynyl-1,3-dimethylperimidin-... | 236    | C15H12N2O        | 1000443-36-0 | 35   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011522\  
Data File : BG052012.D  
Acq On : 16 Jan 2022 1:50  
Operator : CG/JU  
Sample : N1132-04 10X  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLT5

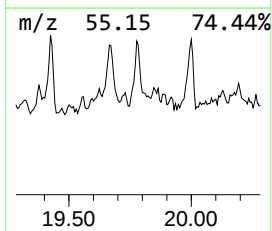
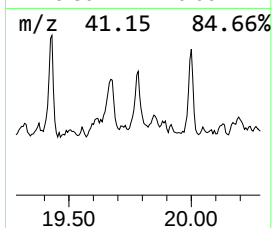
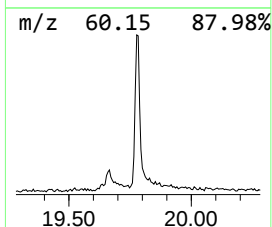
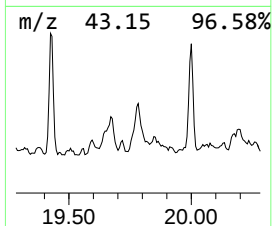
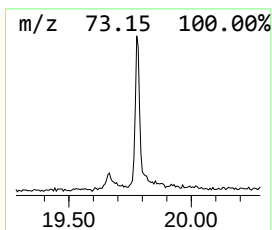
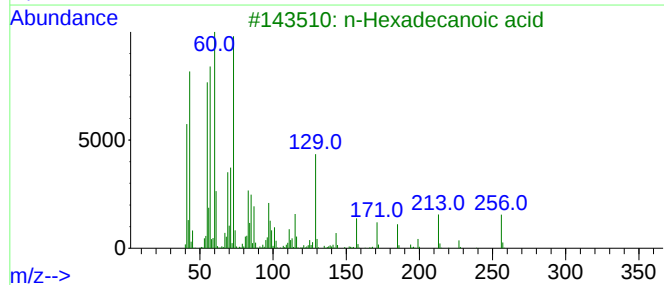
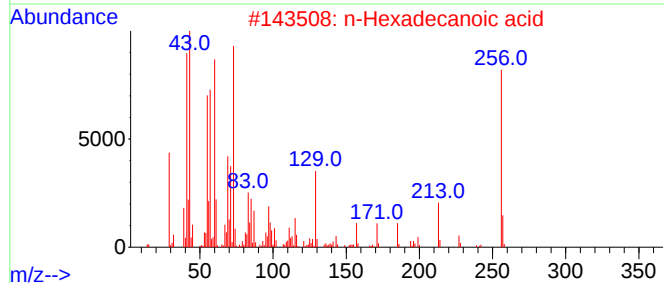
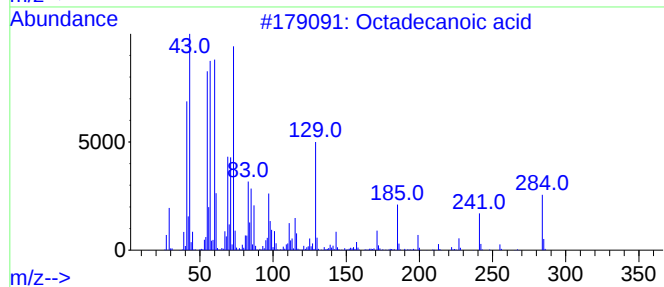
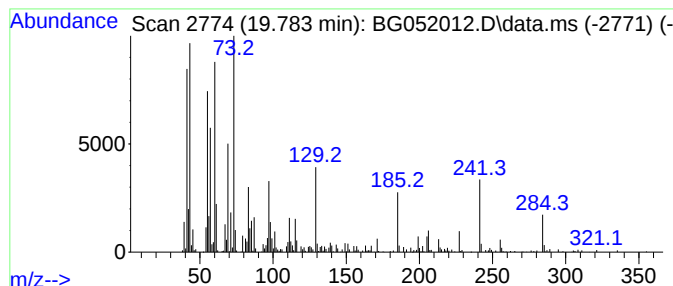
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Quant Title : SVOA CALIBRATION

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

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Peak Number 65 Octadecanoic acid Concentration Rank 35

| R.T.      | EstConc             | Area   | Relative to ISTD | R.T.        |      |
|-----------|---------------------|--------|------------------|-------------|------|
| 19.783    | 12.86 ng/ul         | 456472 | Phenanthrene-d10 | 17.645      |      |
| Hit# of 5 | Tentative ID        | MW     | MolForm          | CAS#        | Qual |
| 1         | Octadecanoic acid   | 284    | C18H36O2         | 000057-11-4 | 78   |
| 2         | n-Hexadecanoic acid | 256    | C16H32O2         | 000057-10-3 | 70   |
| 3         | n-Hexadecanoic acid | 256    | C16H32O2         | 000057-10-3 | 62   |
| 4         | Pentadecanoic acid  | 242    | C15H30O2         | 001002-84-2 | 58   |
| 5         | Pentadecanoic acid  | 242    | C15H30O2         | 001002-84-2 | 58   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011522\  
Data File : BG052012.D  
Acq On : 16 Jan 2022 1:50  
Operator : CG/JU  
Sample : N1132-04 10X  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLT5

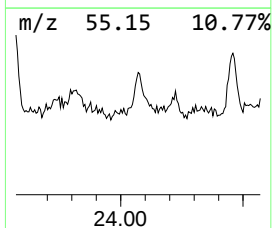
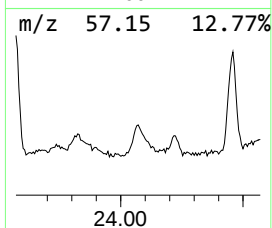
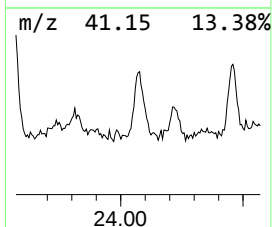
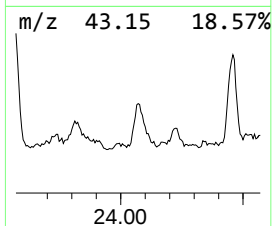
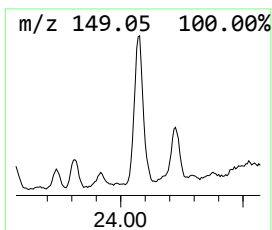
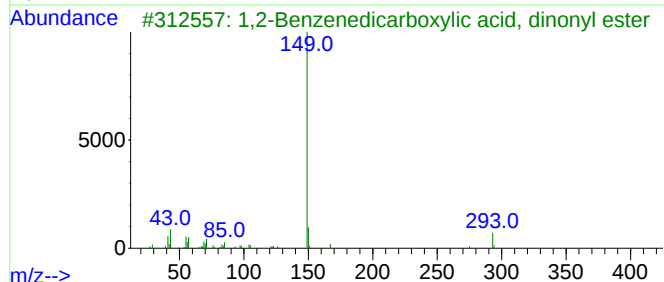
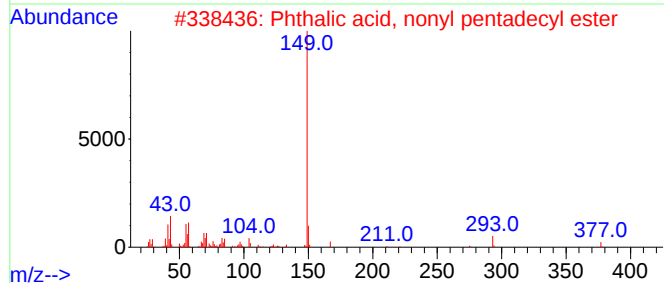
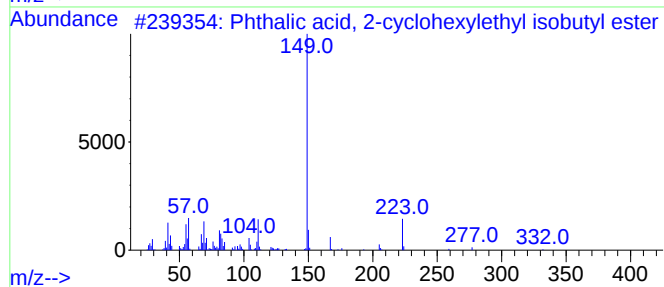
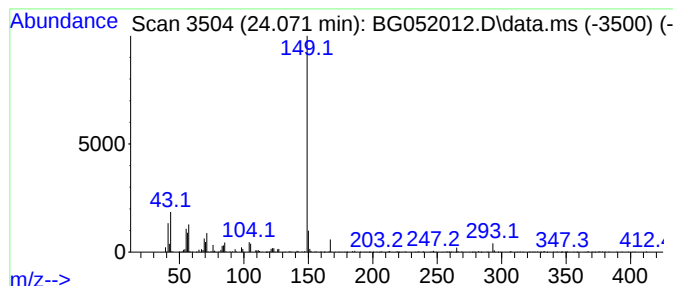
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Quant Title : SVOA CALIBRATION

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

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Peak Number 66 Phthalic acid, 2-cyclohexyl... Concentration Rank 19

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 24.071 | 22.74 ng/ul | 704644 | Perylene-d12     | 25.463 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Phthalic acid, 2-cyclohexylethyl... | 332 | C20H28O4 | 1000309-06-9 | 87   |
| 2    |    |   | Phthalic acid, nonyl pentadecyl ... | 502 | C32H54O4 | 1000308-92-7 | 86   |
| 3    |    |   | 1,2-Benzenedicarboxylic acid, di... | 418 | C26H42O4 | 000084-76-4  | 86   |
| 4    |    |   | Phthalic acid, 2,2-dimethylpent-... | 390 | C24H38O4 | 1000415-53-2 | 86   |
| 5    |    |   | 1,2-Benzenedicarboxylic acid, de... | 418 | C26H42O4 | 000119-07-3  | 80   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011522\  
Data File : BG052012.D  
Acq On : 16 Jan 2022 1:50  
Operator : CG/JU  
Sample : N1132-04 10X  
Misc :  
ALS Vial : 23 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLT5

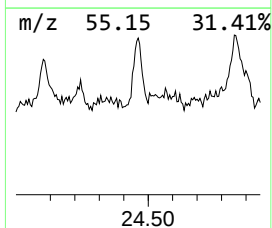
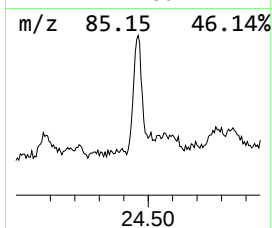
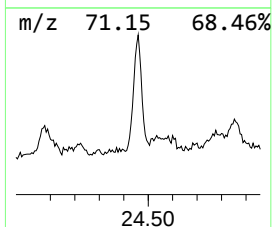
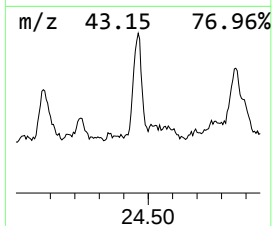
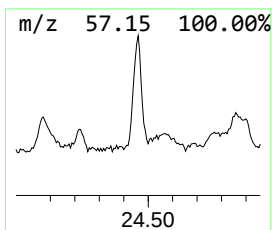
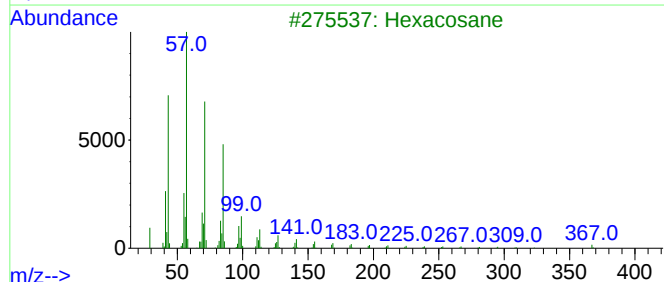
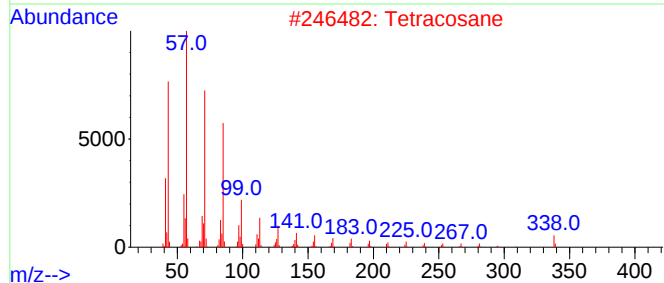
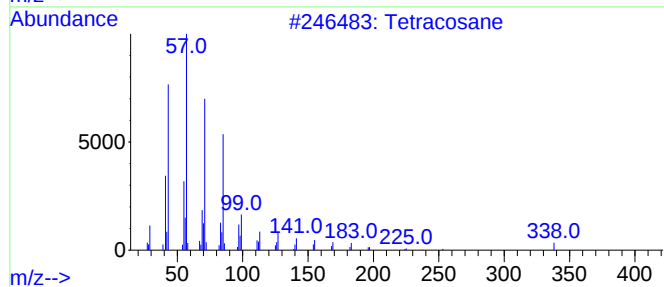
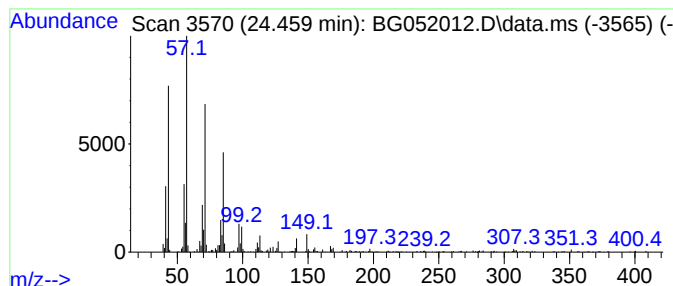
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Quant Title : SVOA CALIBRATION

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

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Peak Number 67 (DEL) Alkane: Straight-Chai... Concentration Rank 25

| R.T.      | EstConc         | Area   | Relative to ISTD | R.T.        |      |
|-----------|-----------------|--------|------------------|-------------|------|
| 24.459    | 20.00 ng/ul     | 619651 | Perylene-d12     | 25.463      |      |
| Hit# of 5 | Tentative ID    | MW     | MolForm          | CAS#        | Qual |
| 1         | Tetracosane     | 338    | C24H50           | 000646-31-1 | 97   |
| 2         | Tetracosane     | 338    | C24H50           | 000646-31-1 | 91   |
| 3         | Hexacosane      | 366    | C26H54           | 000630-01-3 | 87   |
| 4         | Heptadecane     | 240    | C17H36           | 000629-78-7 | 87   |
| 5         | Hexatriacontane | 507    | C36H74           | 000630-06-8 | 87   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011522\  
 Data File : BG052012.D  
 Acq On : 16 Jan 2022 1:50  
 Operator : CG/JU  
 Sample : N1132-04 10X  
 Misc :  
 ALS Vial : 23 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 BGLT5

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010622.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 |
| Benzene, 1,3-di... | 5.849  | 6.9     | ng/ul | 110594   | 1                     | 8.245  | 321608 | 20.0 |
| Benzene, 1,2,4-... | 7.458  | 6.5     | ng/ul | 105263   | 1                     | 8.245  | 321608 | 20.0 |
| (DEL) Alkane: S... | 7.858  | 18.7    | ng/ul | 299992   | 1                     | 8.245  | 321608 | 20.0 |
| Benzene, 1,2,3-... | 7.887  | 14.7    | ng/ul | 235725   | 1                     | 8.245  | 321608 | 20.0 |
| 1-Propyne, 3-ph... | 8.798  | 14.2    | ng/ul | 228173   | 1                     | 8.245  | 321608 | 20.0 |
| (DEL) Alkane: S... | 9.015  | 7.3     | ng/ul | 118212   | 1                     | 8.245  | 321608 | 20.0 |
| Benzene, 4-ethy... | 9.367  | 11.9    | ng/ul | 190744   | 1                     | 8.245  | 321608 | 20.0 |
| Carbonic acid, ... | 9.485  | 50.5    | ng/ul | 811878   | 1                     | 8.245  | 321608 | 20.0 |
| (DEL) Alkane: S... | 9.738  | 11.4    | ng/ul | 150743   | 2                     | 11.089 | 263428 | 20.0 |
| Benzene, 2-ethy... | 9.902  | 8.7     | ng/ul | 114301   | 2                     | 11.089 | 263428 | 20.0 |
| (DEL) Alkane: S... | 10.337 | 11.9    | ng/ul | 156523   | 2                     | 11.089 | 263428 | 20.0 |
| (DEL) Alkane: S... | 10.413 | 8.7     | ng/ul | 114240   | 2                     | 11.089 | 263428 | 20.0 |
| (DEL) Alkane: S... | 10.490 | 17.8    | ng/ul | 234172   | 2                     | 11.089 | 263428 | 20.0 |
| (DEL) Alkane: S... | 10.595 | 7.2     | ng/ul | 94415    | 2                     | 11.089 | 263428 | 20.0 |
| (DEL) Alkane: C... | 11.788 | 11.5    | ng/ul | 151965   | 2                     | 11.089 | 263428 | 20.0 |
| (DEL) Alkane: S... | 11.911 | 12.2    | ng/ul | 160042   | 2                     | 11.089 | 263428 | 20.0 |
| (DEL) Alkane: S... | 11.987 | 16.9    | ng/ul | 222900   | 2                     | 11.089 | 263428 | 20.0 |
| Sulfurous acid,... | 12.093 | 34.5    | ng/ul | 453939   | 2                     | 11.089 | 263428 | 20.0 |
| Naphthalene, 1,... | 12.181 | 7.2     | ng/ul | 94434    | 2                     | 11.089 | 263428 | 20.0 |
| (DEL) Alkane: S... | 12.481 | 97.8    | ng/ul | 1288700  | 2                     | 11.089 | 263428 | 20.0 |
| unknown-01         | 12.616 | 10.5    | ng/ul | 138561   | 2                     | 11.089 | 263428 | 20.0 |
| (DEL) Alkane: S... | 13.104 | 8.0     | ng/ul | 195742   | 3                     | 14.895 | 487105 | 20.0 |
| (DEL) Alkane: C... | 13.174 | 5.8     | ng/ul | 141997   | 3                     | 14.895 | 487105 | 20.0 |
| (DEL) Alkane: S... | 13.280 | 9.1     | ng/ul | 220378   | 3                     | 14.895 | 487105 | 20.0 |
| (DEL) Alkane: S... | 13.362 | 5.8     | ng/ul | 140043   | 3                     | 14.895 | 487105 | 20.0 |
| (DEL) Alkane: S... | 13.421 | 22.1    | ng/ul | 536931   | 3                     | 14.895 | 487105 | 20.0 |
| Naphthalene, 1,... | 14.055 | 25.3    | ng/ul | 615583   | 3                     | 14.895 | 487105 | 20.0 |
| Naphthalene, 2,... | 14.220 | 26.8    | ng/ul | 653619   | 3                     | 14.895 | 487105 | 20.0 |
| (DEL) Alkane: S... | 14.331 | 16.2    | ng/ul | 395229   | 3                     | 14.895 | 487105 | 20.0 |
| (DEL) Alkane: S... | 14.361 | 30.3    | ng/ul | 737238   | 3                     | 14.895 | 487105 | 20.0 |
| (DEL) Alkane: S... | 14.396 | 9.3     | ng/ul | 226284   | 3                     | 14.895 | 487105 | 20.0 |
| unknown-02         | 14.725 | 9.5     | ng/ul | 231878   | 3                     | 14.895 | 487105 | 20.0 |
| (DEL) Alkane: S... | 14.772 | 108.1   | ng/ul | 2632800  | 3                     | 14.895 | 487105 | 20.0 |
| 3-(2-Methyl-pro... | 15.101 | 15.5    | ng/ul | 378398   | 3                     | 14.895 | 487105 | 20.0 |
| (DEL) Alkane: S... | 15.207 | 5.2     | ng/ul | 125718   | 3                     | 14.895 | 487105 | 20.0 |
| (DEL) Alkane: S... | 15.383 | 35.2    | ng/ul | 857474   | 3                     | 14.895 | 487105 | 20.0 |
| (DEL) Alkane: S... | 15.453 | 8.0     | ng/ul | 195407   | 3                     | 14.895 | 487105 | 20.0 |
| Naphthalene, 1,... | 15.506 | 12.3    | ng/ul | 298591   | 3                     | 14.895 | 487105 | 20.0 |
| Naphthalene, 2,... | 15.559 | 10.5    | ng/ul | 256676   | 3                     | 14.895 | 487105 | 20.0 |
| 4,6,8-Trimethyl... | 15.682 | 20.7    | ng/ul | 503506   | 3                     | 14.895 | 487105 | 20.0 |
| (DEL) Alkane: S... | 15.724 | 103.1   | ng/ul | 2511470  | 3                     | 14.895 | 487105 | 20.0 |
| unknown-03         | 15.823 | 20.4    | ng/ul | 497294   | 3                     | 14.895 | 487105 | 20.0 |
| unknown-04         | 16.064 | 7.3     | ng/ul | 177856   | 3                     | 14.895 | 487105 | 20.0 |
| Dodecane, 1-iodo-  | 16.129 | 52.6    | ng/ul | 1280310  | 3                     | 14.895 | 487105 | 20.0 |
| (DEL) Alkane: S... | 16.217 | 9.4     | ng/ul | 229002   | 3                     | 14.895 | 487105 | 20.0 |
| (DEL) Alkane: C... | 16.340 | 9.3     | ng/ul | 328547   | 4                     | 17.645 | 710072 | 20.0 |
| (DEL) Alkane: S... | 16.587 | 71.2    | ng/ul | 2525970  | 4                     | 17.645 | 710072 | 20.0 |
| Sulfurous acid,... | 16.611 | 44.9    | ng/ul | 1592240  | 4                     | 17.645 | 710072 | 20.0 |
| Benzene, (1-met... | 16.705 | 20.0    | ng/ul | 711256   | 4                     | 17.645 | 710072 | 20.0 |
| unknown-05         | 16.863 | 8.0     | ng/ul | 284619   | 4                     | 17.645 | 710072 | 20.0 |
| unknown-06         | 17.092 | 8.3     | ng/ul | 295882   | 4                     | 17.645 | 710072 | 20.0 |
| (DEL) Alkane: S... | 17.380 | 48.6    | ng/ul | 1724820  | 4                     | 17.645 | 710072 | 20.0 |

|                    |        |      |       |         |   |        |        |      |
|--------------------|--------|------|-------|---------|---|--------|--------|------|
| (DEL) Alkane: S... | 17.433 | 22.8 | ng/ul | 807992  | 4 | 17.645 | 710072 | 20.0 |
| Benzene, (1-met... | 17.527 | 20.0 | ng/ul | 710072  | 4 | 17.645 | 710072 | 20.0 |
| (DEL) Alkane: S... | 18.120 | 46.7 | ng/ul | 1657040 | 4 | 17.645 | 710072 | 20.0 |
| n-Hexadecanoic ... | 18.526 | 40.1 | ng/ul | 1424070 | 4 | 17.645 | 710072 | 20.0 |
| (DEL) Alkane: S... | 18.808 | 27.9 | ng/ul | 990292  | 4 | 17.645 | 710072 | 20.0 |
| p-Dicyclohexylb... | 19.078 | 13.6 | ng/ul | 483538  | 4 | 17.645 | 710072 | 20.0 |
| (DEL) Alkane: S... | 19.424 | 30.0 | ng/ul | 1063670 | 4 | 17.645 | 710072 | 20.0 |
| cyclohexane, [1... | 19.612 | 14.3 | ng/ul | 507462  | 4 | 17.645 | 710072 | 20.0 |
| Octadecanoic acid  | 19.783 | 12.9 | ng/ul | 456472  | 4 | 17.645 | 710072 | 20.0 |
| Phthalic acid, ... | 24.071 | 22.7 | ng/ul | 704644  | 6 | 25.463 | 619651 | 20.0 |
| (DEL) Alkane: S... | 24.459 | 20.0 | ng/ul | 619651  | 6 | 25.463 | 619651 | 20.0 |

Instrument :

BNA\_G

ClientSampleId :

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