

Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG121016\  
 Data File : BG025074.D  
 Acq On : 10 Dec 2016 18:37  
 Operator : UM/SJ  
 Sample : H5861-07  
 Misc :  
 ALS Vial : 11 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 DA8D9

## Integration Parameters: LSCINT.P

Integrator: RTE  
 Smoothing : ON  
 Sampling : 1  
 Start Thrs: 0.1  
 Stop Thrs : 0

Filtering: 5  
 Min Area: 0.1 % of largest Peak  
 Max Peaks: 100  
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >  
 Peak separation: 1

Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG113016.M  
 Title : SVOA CALIBRATION

Signal : TIC

| peak<br># | R.T.<br>min | first<br>scan | max<br>scan | last<br>scan | PK<br>TY | peak<br>height | corr.<br>area | corr.<br>% max. | % of<br>total |
|-----------|-------------|---------------|-------------|--------------|----------|----------------|---------------|-----------------|---------------|
| 1         | 3.943       | 185           | 188         | 193          | rVB6     | 7438           | 15074         | 0.44%           | 0.049%        |
| 2         | 4.284       | 245           | 246         | 252          | rBV3     | 6743           | 10825         | 0.32%           | 0.035%        |
| 3         | 4.360       | 255           | 259         | 262          | rBV3     | 7030           | 10385         | 0.30%           | 0.034%        |
| 4         | 4.654       | 304           | 309         | 313          | rVB4     | 9468           | 12773         | 0.37%           | 0.041%        |
| 5         | 4.901       | 345           | 351         | 358          | rBV4     | 56845          | 107007        | 3.12%           | 0.346%        |
| 6         | 5.306       | 411           | 420         | 428          | rVV7     | 15297          | 46350         | 1.35%           | 0.150%        |
| 7         | 5.800       | 500           | 504         | 510          | rBV4     | 5752           | 11980         | 0.35%           | 0.039%        |
| 8         | 6.422       | 608           | 610         | 614          | rVB4     | 8362           | 10427         | 0.30%           | 0.034%        |
| 9         | 7.192       | 735           | 741         | 754          | rVV3     | 48392          | 99301         | 2.90%           | 0.321%        |
| 10        | 7.292       | 754           | 758         | 764          | rVB3     | 18624          | 33696         | 0.98%           | 0.109%        |
| 11        | 7.380       | 766           | 773         | 785          | rBV2     | 93457          | 169029        | 4.93%           | 0.546%        |
| 12        | 7.550       | 795           | 802         | 813          | rVB      | 259599         | 445328        | 13.00%          | 1.438%        |
| 13        | 7.639       | 813           | 817         | 821          | rVB6     | 7557           | 13608         | 0.40%           | 0.044%        |
| 14        | 7.691       | 821           | 826         | 828          | rBV3     | 6246           | 10257         | 0.30%           | 0.033%        |
| 15        | 7.762       | 830           | 838         | 849          | rBV      | 292907         | 523721        | 15.29%          | 1.692%        |
| 16        | 8.238       | 911           | 919         | 928          | rBV      | 261363         | 463039        | 13.52%          | 1.496%        |
| 17        | 8.784       | 1004          | 1012        | 1017         | rBV7     | 15050          | 38752         | 1.13%           | 0.125%        |
| 18        | 8.937       | 1027          | 1038        | 1046         | rBV2     | 242050         | 439942        | 12.84%          | 1.421%        |
| 19        | 9.066       | 1059          | 1060        | 1066         | rBV3     | 7242           | 12647         | 0.37%           | 0.041%        |
| 20        | 9.413       | 1110          | 1119        | 1127         | rBV      | 361170         | 673250        | 19.65%          | 2.175%        |
| 21        | 9.489       | 1127          | 1132        | 1138         | rVV5     | 40787          | 85581         | 2.50%           | 0.276%        |
| 22        | 9.542       | 1138          | 1141        | 1154         | rVB7     | 18083          | 47209         | 1.38%           | 0.152%        |
| 23        | 9.742       | 1170          | 1175        | 1182         | rVB7     | 25149          | 46644         | 1.36%           | 0.151%        |
| 24        | 9.801       | 1182          | 1185        | 1191         | rBV6     | 6560           | 11460         | 0.33%           | 0.037%        |
| 25        | 9.954       | 1209          | 1211        | 1217         | rVB5     | 6882           | 11023         | 0.32%           | 0.036%        |
| 26        | 10.018      | 1218          | 1222        | 1228         | rVB3     | 5865           | 9736          | 0.28%           | 0.031%        |
| 27        | 10.142      | 1228          | 1243        | 1252         | rBV2     | 336237         | 649275        | 18.95%          | 2.097%        |
| 28        | 10.330      | 1270          | 1275        | 1289         | rBV5     | 37990          | 92744         | 2.71%           | 0.300%        |
| 29        | 10.541      | 1307          | 1311        | 1314         | rBV5     | 6348           | 10738         | 0.31%           | 0.035%        |
| 30        | 10.570      | 1314          | 1316        | 1325         | rVB7     | 7076           | 15351         | 0.45%           | 0.050%        |
| 31        | 10.676      | 1325          | 1334        | 1348         | rBV      | 421307         | 858903        | 25.07%          | 2.774%        |
| 32        | 10.806      | 1352          | 1356        | 1364         | rBV4     | 31640          | 68395         | 2.00%           | 0.221%        |
| 33        | 11.064      | 1391          | 1400        | 1410         | rVB      | 364962         | 658759        | 19.23%          | 2.128%        |
| 34        | 11.211      | 1414          | 1425        | 1428         | rBV8     | 8905           | 23458         | 0.68%           | 0.076%        |

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 Misc :  
 ALS Vial : 11 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 DA8D9

## Integration Parameters: LSCINT.P

Integrator: RTE  
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 Sampling : 1  
 Start Thrs: 0.1  
 Stop Thrs : 0

Filtering: 5  
 Min Area: 0.1 % of largest Peak  
 Max Peaks: 100  
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >  
 Peak separation: 1

Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG113016.M  
 Title : SVOA CALIBRATION

|    |        |      |      |      |       |         |         |        |        |
|----|--------|------|------|------|-------|---------|---------|--------|--------|
| 35 | 11.563 | 1478 | 1485 | 1490 | rBV8  | 9080    | 19541   | 0.57%  | 0.063% |
| 36 | 11.616 | 1490 | 1494 | 1501 | rVB7  | 11553   | 24159   | 0.71%  | 0.078% |
| 37 | 11.816 | 1520 | 1528 | 1537 | rBV3  | 86907   | 183863  | 5.37%  | 0.594% |
| 38 | 12.169 | 1581 | 1588 | 1590 | rBV6  | 6632    | 14620   | 0.43%  | 0.047% |
| 39 | 12.233 | 1590 | 1599 | 1605 | rVB4  | 23788   | 47207   | 1.38%  | 0.152% |
| 40 | 12.380 | 1623 | 1624 | 1631 | rVB6  | 6547    | 12132   | 0.35%  | 0.039% |
| 41 | 12.439 | 1631 | 1634 | 1636 | rBV2  | 6864    | 10570   | 0.31%  | 0.034% |
| 42 | 12.509 | 1644 | 1646 | 1652 | rBV5  | 10490   | 15087   | 0.44%  | 0.049% |
| 43 | 12.639 | 1661 | 1668 | 1672 | rBV5  | 10688   | 23221   | 0.68%  | 0.075% |
| 44 | 12.756 | 1683 | 1688 | 1690 | rBV4  | 7052    | 10465   | 0.31%  | 0.034% |
| 45 | 12.932 | 1715 | 1718 | 1720 | rBV4  | 8634    | 10532   | 0.31%  | 0.034% |
| 46 | 13.115 | 1742 | 1749 | 1754 | rBV7  | 17722   | 30210   | 0.88%  | 0.098% |
| 47 | 13.203 | 1758 | 1764 | 1775 | rBV3  | 146288  | 278230  | 8.12%  | 0.899% |
| 48 | 13.303 | 1778 | 1781 | 1786 | rVB5  | 6032    | 10026   | 0.29%  | 0.032% |
| 49 | 13.444 | 1794 | 1805 | 1814 | rBV   | 229401  | 369183  | 10.78% | 1.192% |
| 50 | 13.602 | 1829 | 1832 | 1836 | rVB5  | 12802   | 18358   | 0.54%  | 0.059% |
| 51 | 13.720 | 1846 | 1852 | 1855 | rBV4  | 22165   | 40083   | 1.17%  | 0.129% |
| 52 | 13.778 | 1858 | 1862 | 1875 | rVB5  | 47044   | 101090  | 2.95%  | 0.327% |
| 53 | 13.914 | 1880 | 1885 | 1889 | rBV5  | 7887    | 17430   | 0.51%  | 0.056% |
| 54 | 14.119 | 1917 | 1920 | 1924 | rBV5  | 7715    | 11759   | 0.34%  | 0.038% |
| 55 | 14.249 | 1930 | 1942 | 1954 | rBV   | 1003832 | 1505839 | 43.96% | 4.864% |
| 56 | 14.354 | 1954 | 1960 | 1969 | rVB10 | 21280   | 35921   | 1.05%  | 0.116% |
| 57 | 14.466 | 1969 | 1979 | 1986 | rBV9  | 37909   | 86983   | 2.54%  | 0.281% |
| 58 | 14.554 | 1987 | 1994 | 2004 | rVV   | 939939  | 1515414 | 44.24% | 4.895% |
| 59 | 14.701 | 2018 | 2019 | 2026 | rVB6  | 9286    | 11554   | 0.34%  | 0.037% |
| 60 | 14.854 | 2039 | 2045 | 2055 | rBV   | 613947  | 1017687 | 29.71% | 3.287% |
| 61 | 15.053 | 2072 | 2079 | 2088 | rBV2  | 112655  | 228300  | 6.66%  | 0.737% |
| 62 | 15.224 | 2103 | 2108 | 2118 | rBV   | 139075  | 233201  | 6.81%  | 0.753% |
| 63 | 15.353 | 2126 | 2130 | 2135 | rVB8  | 17048   | 33382   | 0.97%  | 0.108% |
| 64 | 15.424 | 2135 | 2142 | 2149 | rBV8  | 12674   | 32297   | 0.94%  | 0.104% |
| 65 | 15.524 | 2158 | 2159 | 2165 | rBV6  | 7360    | 13215   | 0.39%  | 0.043% |
| 66 | 15.606 | 2165 | 2173 | 2178 | rBV   | 166393  | 252204  | 7.36%  | 0.815% |
| 67 | 15.788 | 2199 | 2204 | 2205 | rVB4  | 9977    | 13894   | 0.41%  | 0.045% |
| 68 | 15.847 | 2205 | 2214 | 2222 | rVB   | 1331839 | 2145161 | 62.63% | 6.929% |
| 69 | 15.958 | 2222 | 2233 | 2248 | rBV   | 622187  | 998113  | 29.14% | 3.224% |
| 70 | 16.076 | 2249 | 2253 | 2258 | rVB4  | 19623   | 32960   | 0.96%  | 0.106% |
| 71 | 16.305 | 2289 | 2292 | 2295 | rVB5  | 8261    | 10091   | 0.29%  | 0.033% |

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Misc :  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
DA8D9

## Integration Parameters: LSCINT.P

Integrator: RTE  
Smoothing : ON  
Sampling : 1  
Start Thrs: 0.1  
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Filtering: 5  
Min Area: 0.1 % of largest Peak  
Max Peaks: 100  
Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >  
Peak separation: 1

Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG113016.M  
Title : SVOA CALIBRATION

|     |        |      |      |      |       |         |         |         |         |
|-----|--------|------|------|------|-------|---------|---------|---------|---------|
| 72  | 16.399 | 2304 | 2308 | 2315 | rBV7  | 18665   | 42182   | 1.23%   | 0.136%  |
| 73  | 16.640 | 2343 | 2349 | 2353 | rVV8  | 11053   | 22335   | 0.65%   | 0.072%  |
| 74  | 16.687 | 2353 | 2357 | 2364 | rVV9  | 24477   | 49286   | 1.44%   | 0.159%  |
| 75  | 16.945 | 2394 | 2401 | 2405 | rBV6  | 36892   | 68721   | 2.01%   | 0.222%  |
| 76  | 17.069 | 2415 | 2422 | 2427 | rVB5  | 22702   | 42974   | 1.25%   | 0.139%  |
| 77  | 17.263 | 2449 | 2455 | 2457 | rVV7  | 7553    | 12415   | 0.36%   | 0.040%  |
| 78  | 17.321 | 2462 | 2465 | 2470 | rVB5  | 26336   | 29336   | 0.86%   | 0.095%  |
| 79  | 17.392 | 2472 | 2477 | 2480 | rBV5  | 8498    | 15930   | 0.47%   | 0.051%  |
| 80  | 17.421 | 2480 | 2482 | 2488 | rVB6  | 9934    | 13307   | 0.39%   | 0.043%  |
| 81  | 17.598 | 2504 | 2512 | 2522 | rBV2  | 942877  | 1459224 | 42.60%  | 4.713%  |
| 82  | 17.697 | 2522 | 2529 | 2538 | rBV2  | 1659781 | 2555372 | 74.60%  | 8.254%  |
| 83  | 17.844 | 2550 | 2554 | 2561 | rVB3  | 54580   | 74494   | 2.17%   | 0.241%  |
| 84  | 18.150 | 2601 | 2606 | 2608 | rBV5  | 7716    | 13471   | 0.39%   | 0.044%  |
| 85  | 18.226 | 2616 | 2619 | 2626 | rVB9  | 15475   | 22922   | 0.67%   | 0.074%  |
| 86  | 18.444 | 2651 | 2656 | 2660 | rBV8  | 22409   | 38977   | 1.14%   | 0.126%  |
| 87  | 18.526 | 2666 | 2670 | 2677 | rBV   | 56251   | 75259   | 2.20%   | 0.243%  |
| 88  | 18.637 | 2685 | 2689 | 2693 | rBV7  | 9025    | 16153   | 0.47%   | 0.052%  |
| 89  | 18.972 | 2742 | 2746 | 2754 | rBV10 | 15756   | 40000   | 1.17%   | 0.129%  |
| 90  | 19.384 | 2811 | 2816 | 2823 | rBV10 | 24127   | 45008   | 1.31%   | 0.145%  |
| 91  | 19.607 | 2849 | 2854 | 2859 | rBV3  | 27132   | 53232   | 1.55%   | 0.172%  |
| 92  | 19.865 | 2893 | 2898 | 2903 | rBV3  | 30865   | 48991   | 1.43%   | 0.158%  |
| 93  | 19.971 | 2908 | 2916 | 2925 | rVB   | 2355747 | 3425376 | 100.00% | 11.064% |
| 94  | 20.183 | 2950 | 2952 | 2956 | rVB4  | 21863   | 25376   | 0.74%   | 0.082%  |
| 95  | 21.887 | 3236 | 3242 | 3251 | rBV   | 1101462 | 2004858 | 58.53%  | 6.476%  |
| 96  | 23.379 | 3493 | 3496 | 3503 | rVB6  | 51660   | 75412   | 2.20%   | 0.244%  |
| 97  | 24.219 | 3634 | 3639 | 3651 | rVB5  | 61515   | 154870  | 4.52%   | 0.500%  |
| 98  | 25.065 | 3771 | 3783 | 3795 | rBV2  | 1082634 | 3293883 | 96.16%  | 10.639% |
| 99  | 25.212 | 3803 | 3808 | 3813 | rBV5  | 71196   | 152150  | 4.44%   | 0.491%  |
| 100 | 25.300 | 3814 | 3823 | 3835 | rVB2  | 629979  | 1937193 | 56.55%  | 6.257%  |

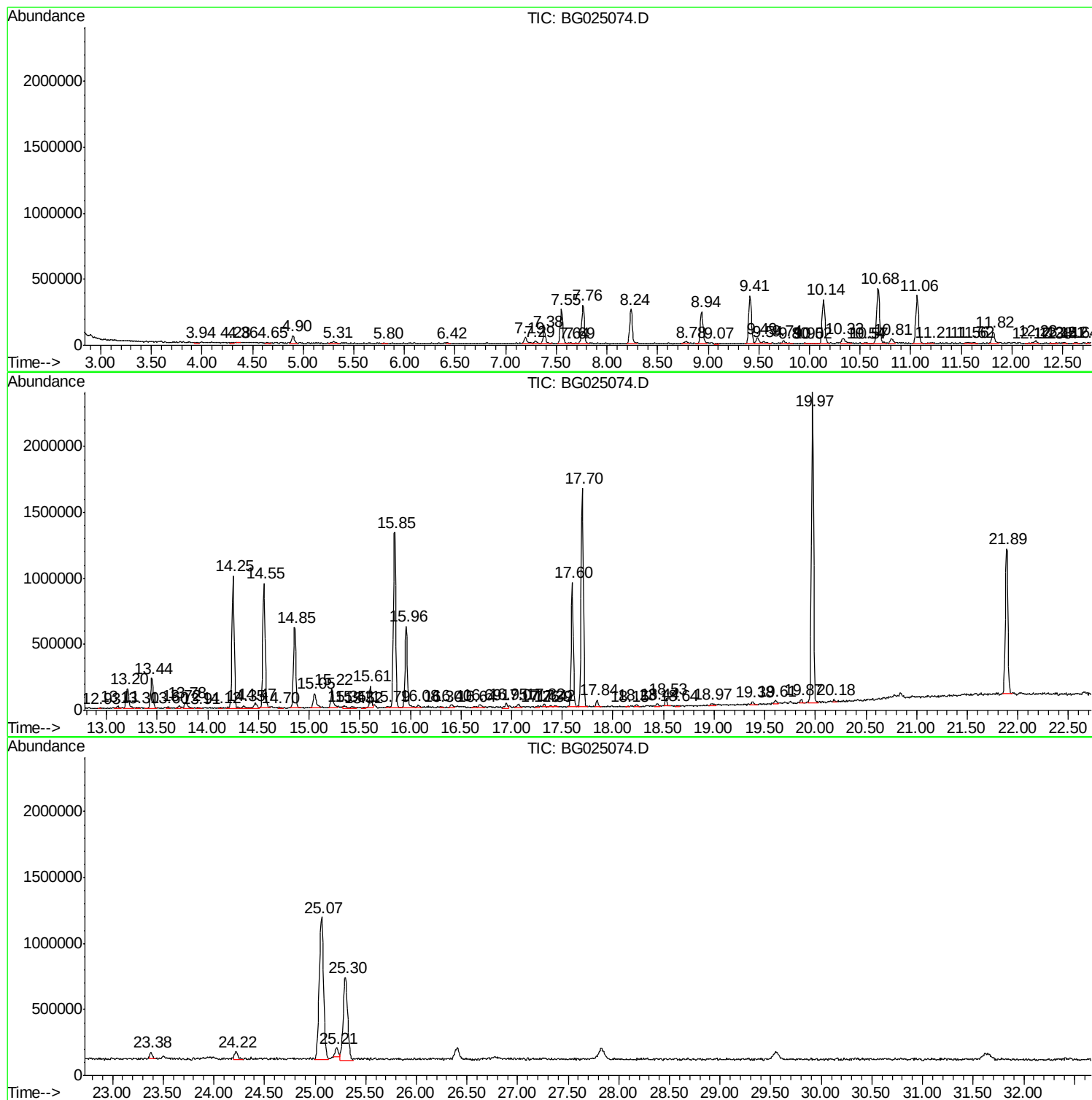
Sum of corrected areas: 30959356

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Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG113016.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P



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Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG113016.M  
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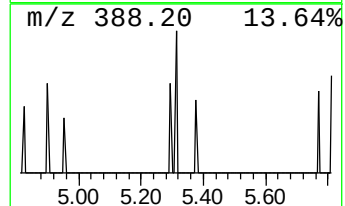
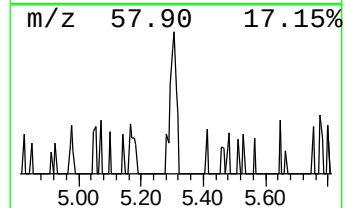
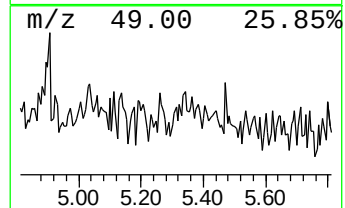
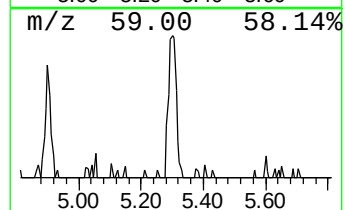
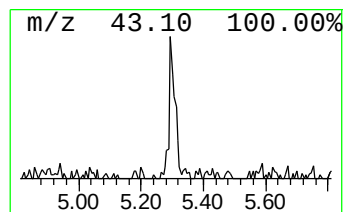
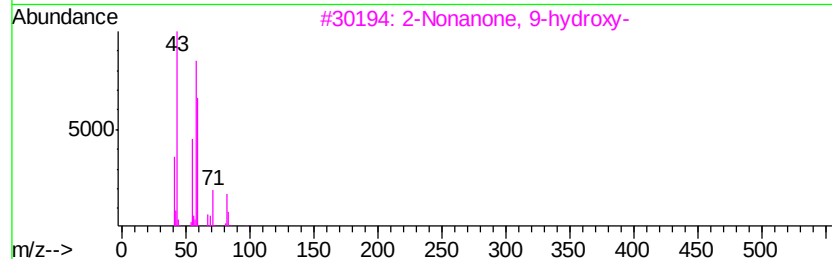
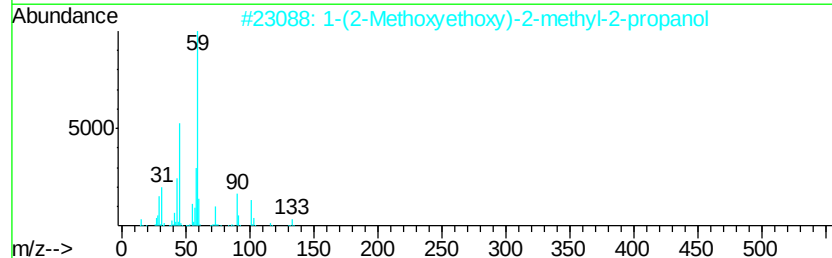
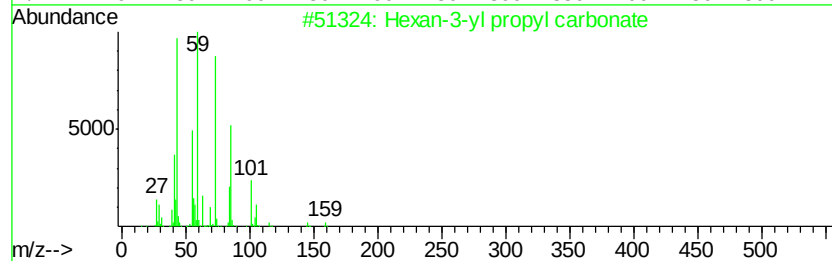
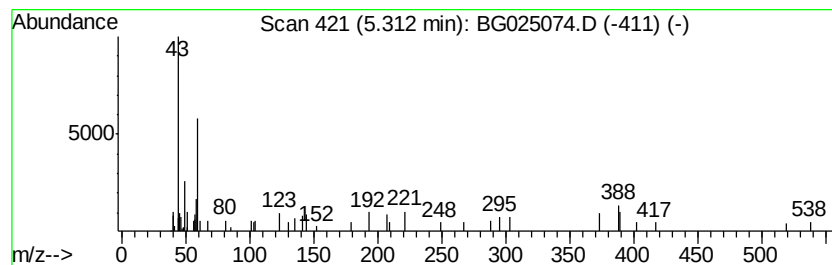
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 2 unknown-01 Concentration Rank 12

| R.T. | EstConc    | Area  | Relative to ISTD       | R.T. |
|------|------------|-------|------------------------|------|
| 5.31 | 2.00 ng/ul | 46350 | 1,4-Dichlorobenzene-d4 | 8.24 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|-------------------------------------|-----|----------|--------------|------|
| 1    | 5  | Hexan-3-yl propyl carbonate         | 188 | C10H20O3 | 1000372-82-0 | 23   |
| 2    |    | 1-(2-Methoxyethoxy)-2-methyl-2-p... | 148 | C7H16O3  | 211321-90-3  | 17   |
| 3    |    | 2-Nonanone, 9-hydroxy-              | 158 | C9H18O2  | 025368-56-3  | 17   |
| 4    |    | Propanedioic acid, bromo-, dimet... | 210 | C5H7BrO4 | 000868-26-8  | 17   |
| 5    |    | Ethanone, 1-cyclopropyl-            | 84  | C5H8O    | 000765-43-5  | 10   |



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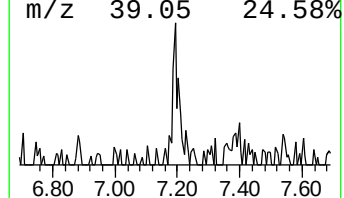
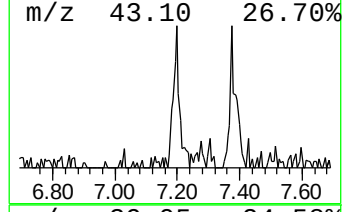
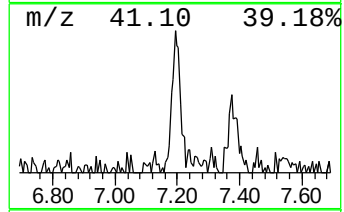
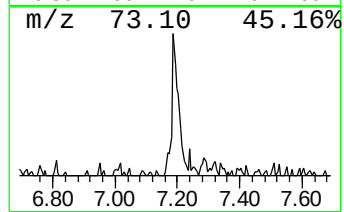
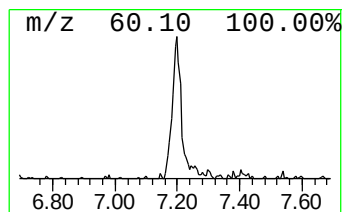
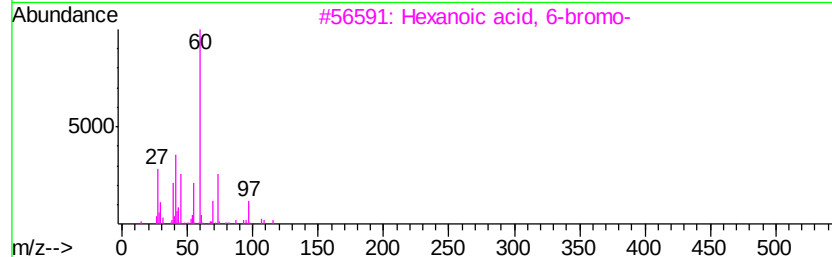
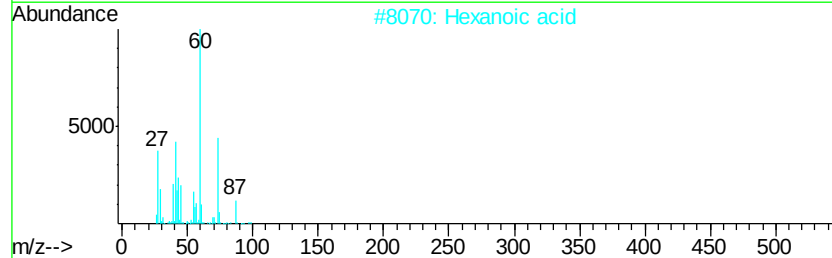
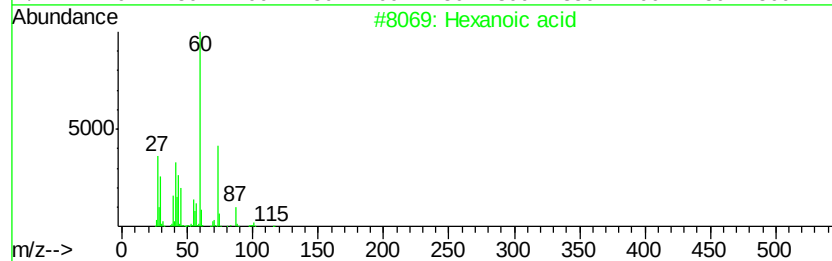
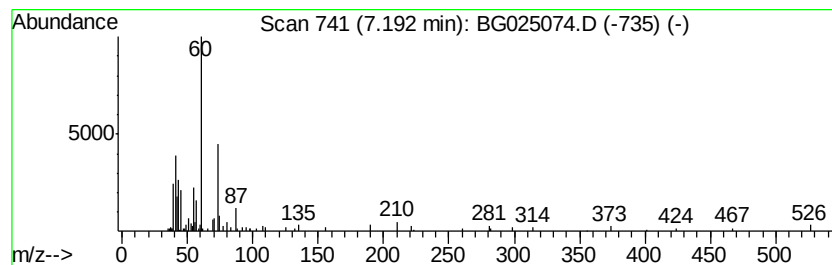
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 3 Hexanoic acid Concentration Rank 7

| R.T. | EstConc    | Area  | Relative to ISTD       | R.T. |
|------|------------|-------|------------------------|------|
| 7.19 | 4.29 ng/ul | 99301 | 1,4-Dichlorobenzene-d4 | 8.24 |

| Hit# | of | Tentative ID            | MW  | MolForm   | CAS#        | Qual |
|------|----|-------------------------|-----|-----------|-------------|------|
| 1    | 5  | Hexanoic acid           | 116 | C6H12O2   | 000142-62-1 | 72   |
| 2    |    | Hexanoic acid           | 116 | C6H12O2   | 000142-62-1 | 72   |
| 3    |    | Hexanoic acid, 6-bromo- | 194 | C6H11BrO2 | 004224-70-8 | 64   |
| 4    |    | Hexanoic acid           | 116 | C6H12O2   | 000142-62-1 | 56   |
| 5    |    | Heptanoic acid          | 130 | C7H14O2   | 000111-14-8 | 53   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG121016\  
Data File : BG025074.D  
Acq On : 10 Dec 2016 18:37  
Operator : UM/SJ  
Sample : H5861-07  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
DA8D9

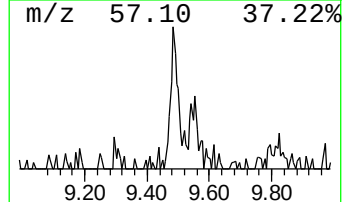
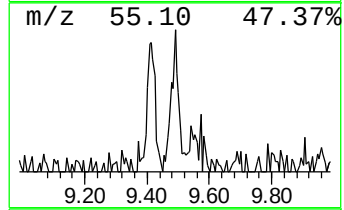
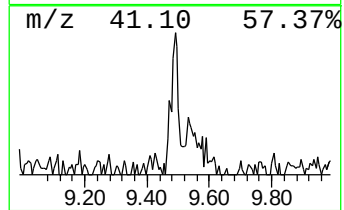
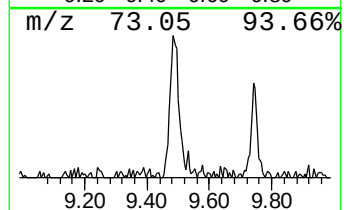
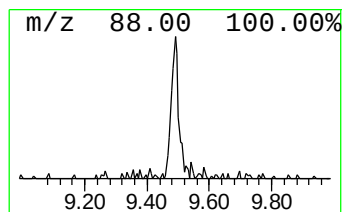
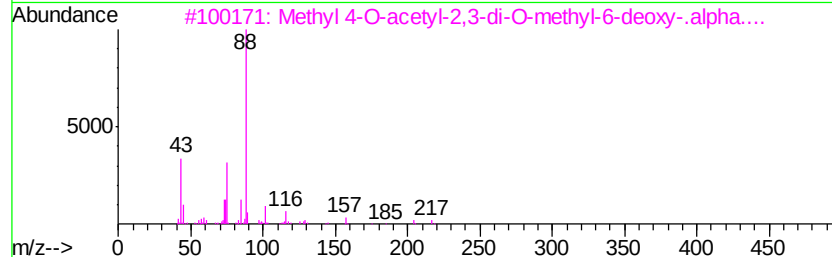
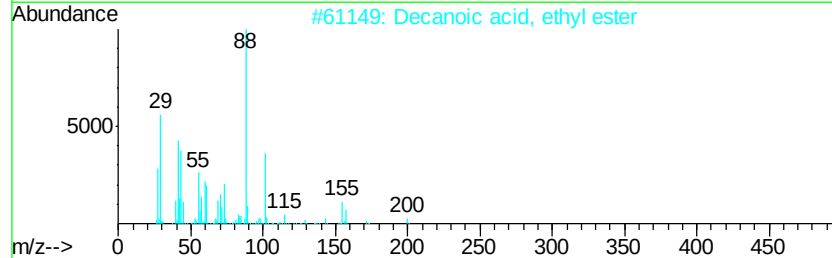
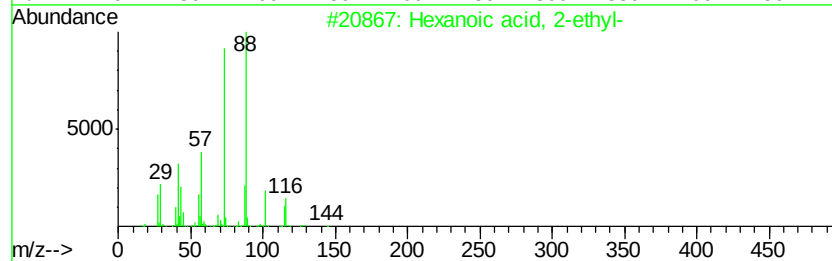
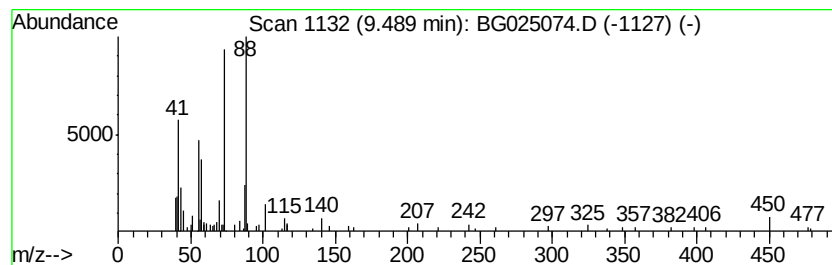
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 4 Hexanoic acid, 2-ethyl- Concentration Rank 8

| R.T. | EstConc    | Area  | Relative to ISTD       | R.T. |
|------|------------|-------|------------------------|------|
| 9.49 | 3.70 ng/ul | 85581 | 1,4-Dichlorobenzene-d4 | 8.24 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | Hexanoic acid, 2-ethyl-             | 144 | C8H16O2  | 000149-57-5 | 72   |
| 2    |    | Decanoic acid, ethyl ester          | 200 | C12H24O2 | 000110-38-3 | 40   |
| 3    |    | Methyl 4-O-acetyl-2,3-di-O-methy... | 248 | C11H20O6 | 072945-56-3 | 40   |
| 4    |    | Heptanoic acid, 2-ethyl-            | 158 | C9H18O2  | 003274-29-1 | 39   |
| 5    |    | Hexanoic acid, 2-ethyl-             | 144 | C8H16O2  | 000149-57-5 | 38   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG121016\  
Data File : BG025074.D  
Acq On : 10 Dec 2016 18:37  
Operator : UM/SJ  
Sample : H5861-07  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
DA8D9

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG113016.M  
Quant Title : SVOA CALIBRATION

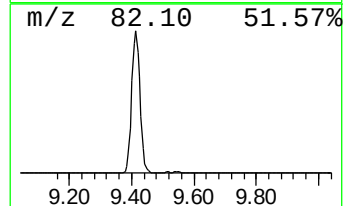
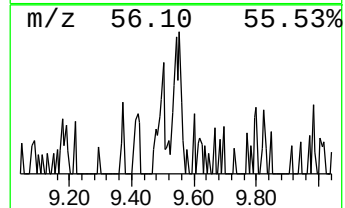
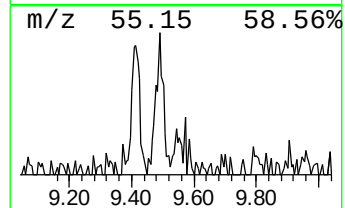
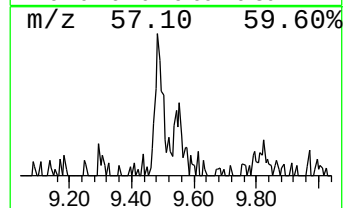
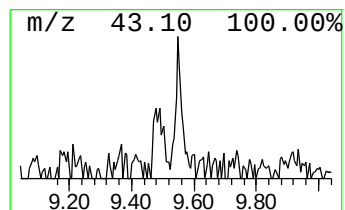
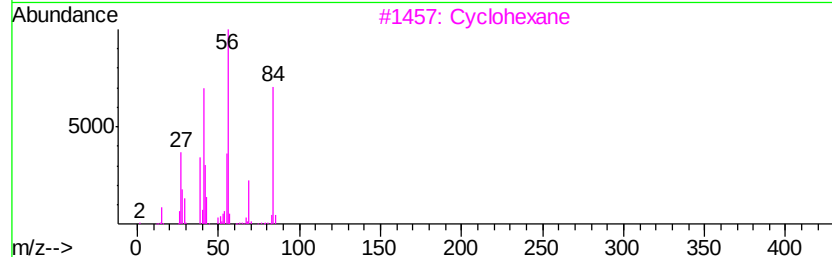
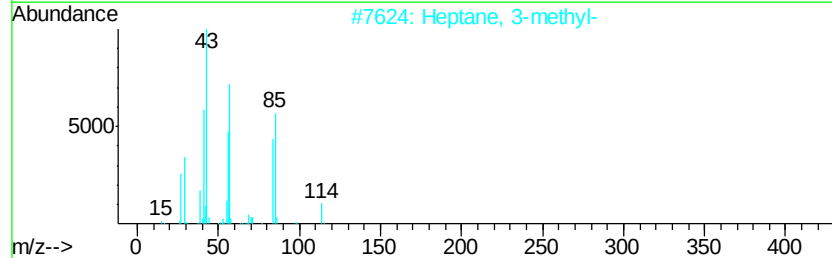
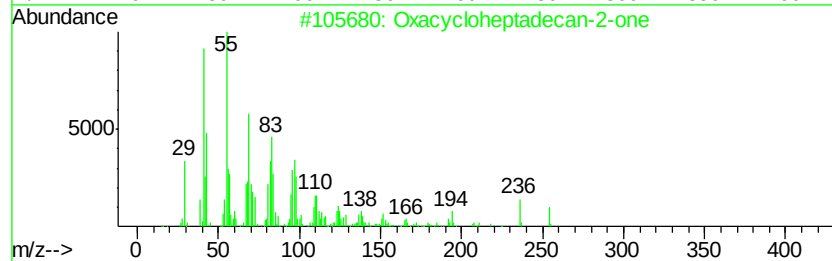
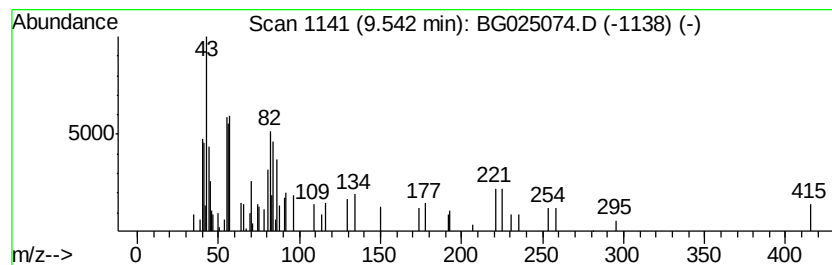
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 5 unknown-02 Concentration Rank 11

| R.T. | EstConc    | Area  | Relative to ISTD       | R.T. |
|------|------------|-------|------------------------|------|
| 9.54 | 2.04 ng/ul | 47209 | 1,4-Dichlorobenzene-d4 | 8.24 |

| Hit# | of | Tentative ID                   | MW  | MolForm  | CAS#        | Qual |
|------|----|--------------------------------|-----|----------|-------------|------|
| 1    | 5  | Oxacycloheptadecan-2-one       | 254 | C16H30O2 | 000109-29-5 | 18   |
| 2    |    | Heptane, 3-methyl-             | 114 | C8H18    | 000589-81-1 | 11   |
| 3    |    | Cyclohexane                    | 84  | C6H12    | 000110-82-7 | 10   |
| 4    |    | 1,3-Cyclopentanediol, cis-     | 102 | C5H10O2  | 016326-97-9 | 10   |
| 5    |    | 1-Pentanol, 2-methyl-, acetate | 144 | C8H16O2  | 007789-99-3 | 9    |





Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG121016\  
Data File : BG025074.D  
Acq On : 10 Dec 2016 18:37  
Operator : UM/SJ  
Sample : H5861-07  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
DA8D9

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG113016.M  
Quant Title : SVOA CALIBRATION

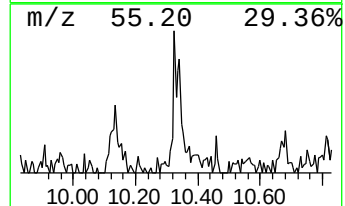
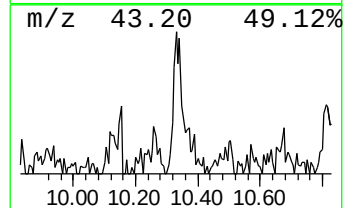
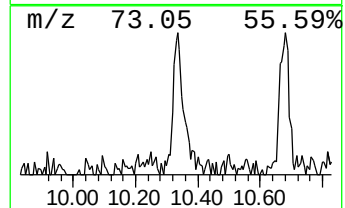
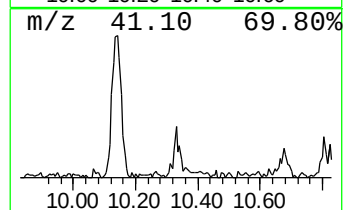
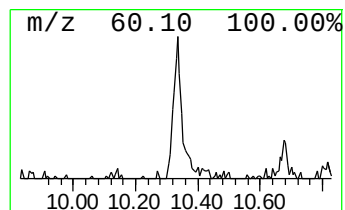
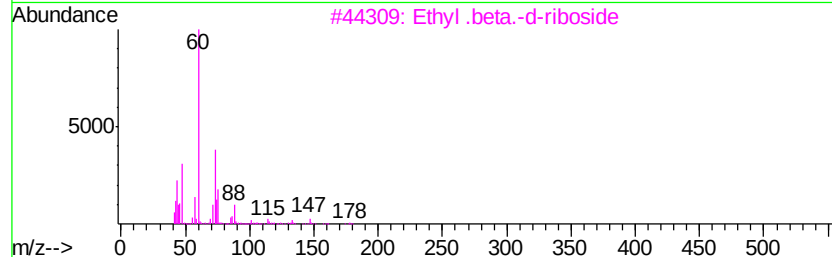
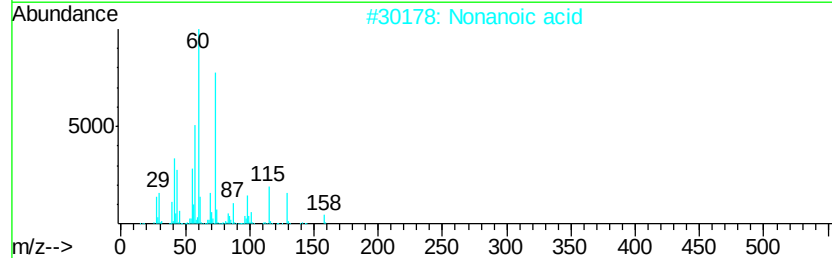
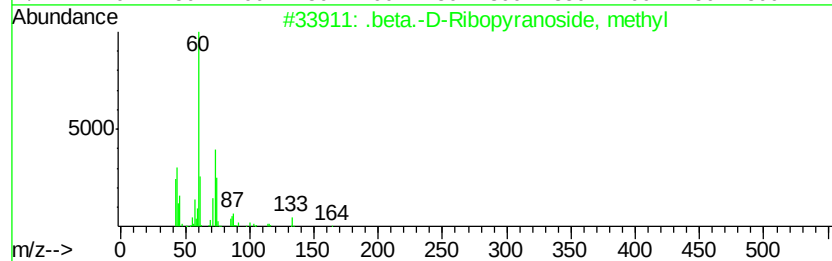
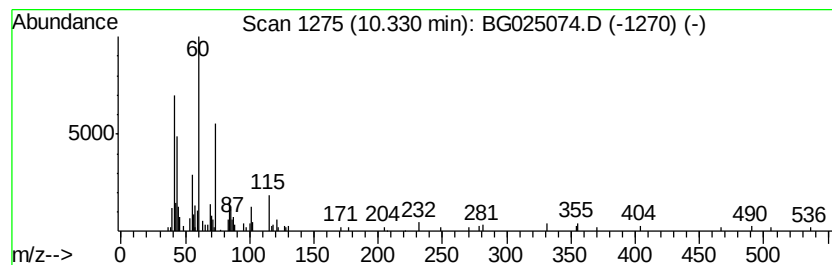
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 6 .beta.-D-Ribopyranoside, me... Concentration Rank 9

| R.T.  | EstConc    | Area  | Relative to ISTD | R.T.  |
|-------|------------|-------|------------------|-------|
| 10.33 | 2.82 ng/ul | 92744 | Napthalene-d8    | 11.06 |

| Hit# of | 5 | Tentative ID                    | MW  | MolForm | CAS#         | Qual |
|---------|---|---------------------------------|-----|---------|--------------|------|
| 1       |   | .beta.-D-Ribopyranoside, methyl | 164 | C6H12O5 | 017289-61-1  | 53   |
| 2       |   | Nonanoic acid                   | 158 | C9H18O2 | 000112-05-0  | 42   |
| 3       |   | Ethyl .beta.-d-ribose           | 178 | C7H14O5 | 1000126-95-4 | 42   |
| 4       |   | Octane, 1-(ethenylthio)-        | 172 | C10H20S | 042779-08-8  | 42   |
| 5       |   | 3-Methyl-hexanoic acid          | 130 | C7H14O2 | 1000365-65-4 | 42   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG121016\  
Data File : BG025074.D  
Acq On : 10 Dec 2016 18:37  
Operator : UM/SJ  
Sample : H5861-07  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
DA8D9

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG113016.M  
Quant Title : SVOA CALIBRATION

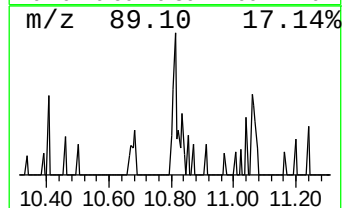
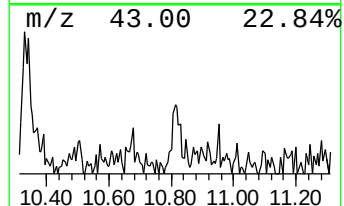
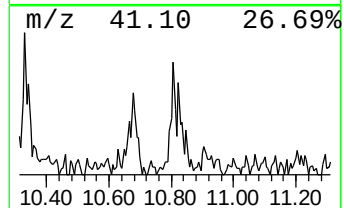
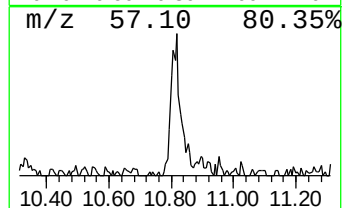
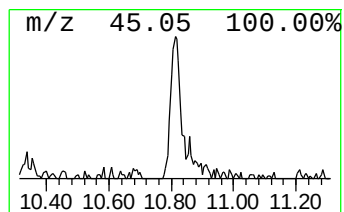
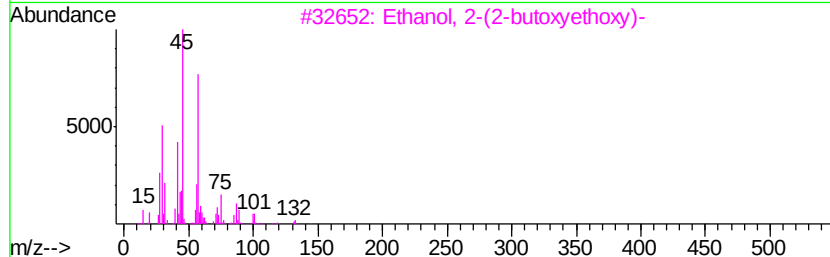
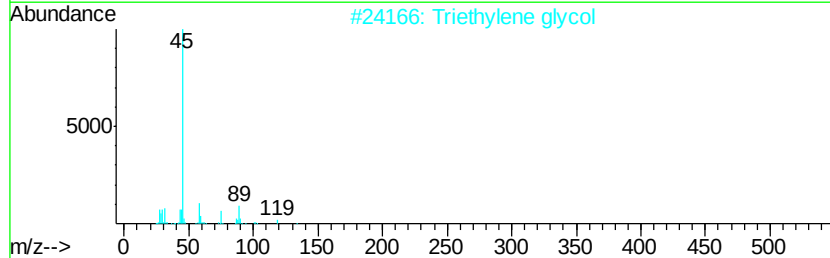
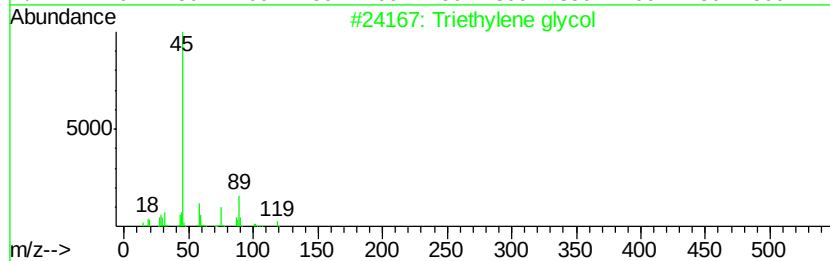
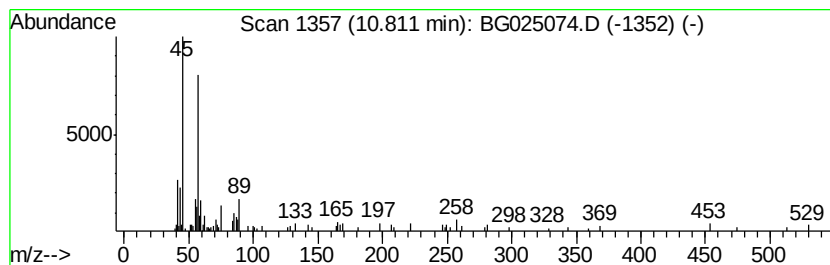
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 7 unknown-03 Concentration Rank 10

| R.T.  | EstConc    | Area  | Relative to ISTD | R.T.  |
|-------|------------|-------|------------------|-------|
| 10.81 | 2.08 ng/ul | 68395 | Naphthalene-d8   | 11.06 |

| Hit# | of | Tentative ID                 | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------------|-----|---------|-------------|------|
| 1    | 5  | Triethylene glycol           | 150 | C6H14O4 | 000112-27-6 | 47   |
| 2    |    | Triethylene glycol           | 150 | C6H14O4 | 000112-27-6 | 47   |
| 3    |    | Ethanol, 2-(2-butoxyethoxy)- | 162 | C8H18O3 | 000112-34-5 | 47   |
| 4    |    | Triethylene glycol           | 150 | C6H14O4 | 000112-27-6 | 47   |
| 5    |    | Ethanol, 2-(2-butoxyethoxy)- | 162 | C8H18O3 | 000112-34-5 | 47   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG121016\  
Data File : BG025074.D  
Acq On : 10 Dec 2016 18:37  
Operator : UM/SJ  
Sample : H5861-07  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
DA8D9

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG113016.M  
Quant Title : SVOA CALIBRATION

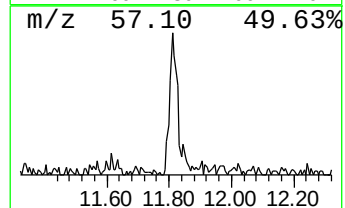
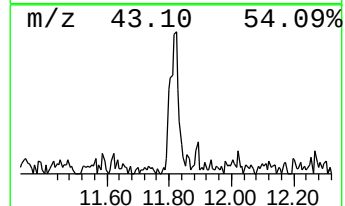
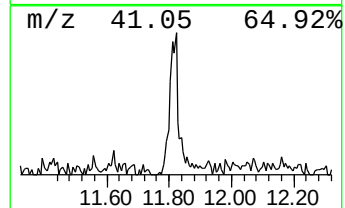
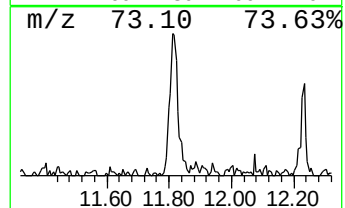
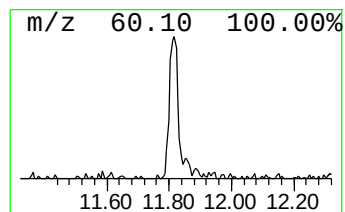
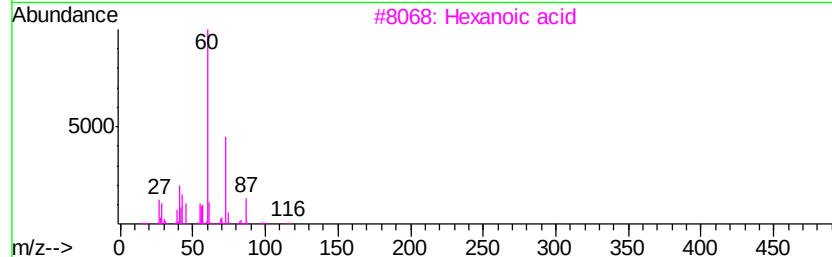
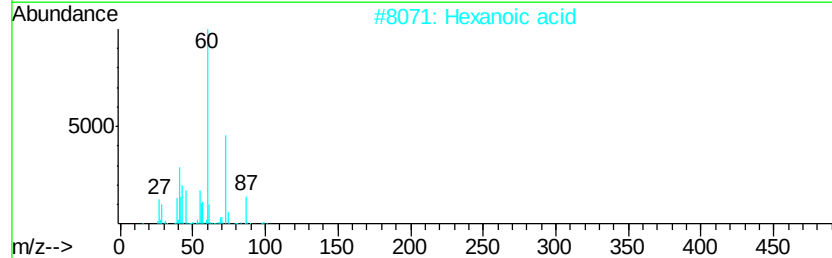
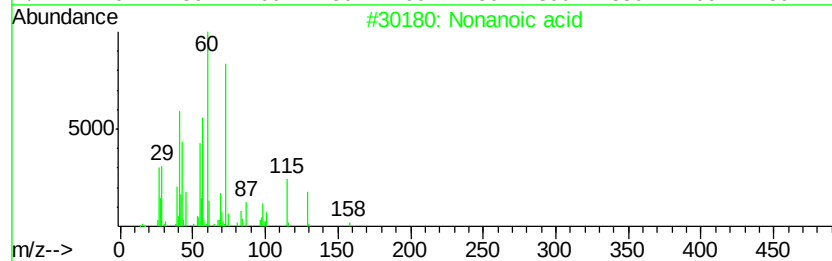
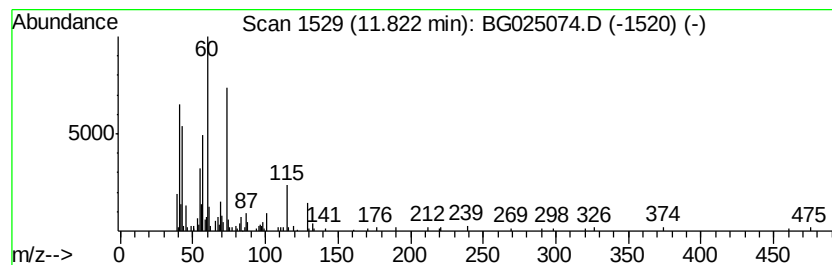
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 8 Nonanoic acid Concentration Rank 2

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 11.82 | 5.58 ng/ul | 183863 | Napthalene-d8    | 11.06 |

| Hit# of | 5               | Tentative ID | MW  | MolForm  | CAS#        | Qual |
|---------|-----------------|--------------|-----|----------|-------------|------|
| 1       | Nonanoic acid   |              | 158 | C9H18O2  | 000112-05-0 | 64   |
| 2       | Hexanoic acid   |              | 116 | C6H12O2  | 000142-62-1 | 50   |
| 3       | Hexanoic acid   |              | 116 | C6H12O2  | 000142-62-1 | 47   |
| 4       | Nonanoic acid   |              | 158 | C9H18O2  | 000112-05-0 | 45   |
| 5       | n-Decanoic acid |              | 172 | C10H20O2 | 000334-48-5 | 45   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG121016\  
Data File : BG025074.D  
Acq On : 10 Dec 2016 18:37  
Operator : UM/SJ  
Sample : H5861-07  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
DA8D9

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG113016.M  
Quant Title : SVOA CALIBRATION

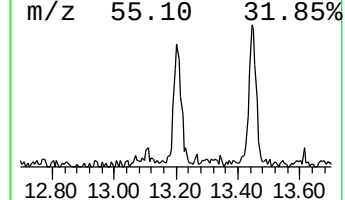
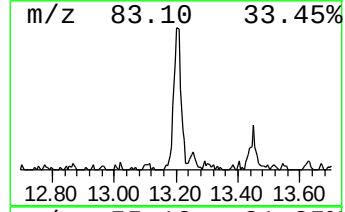
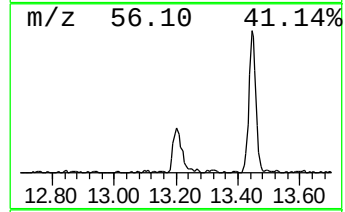
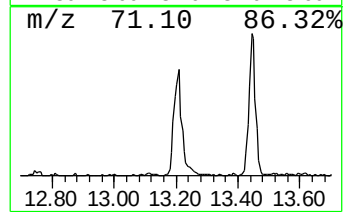
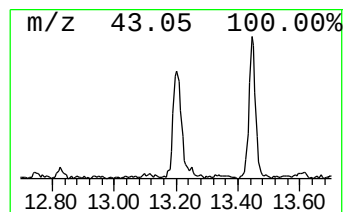
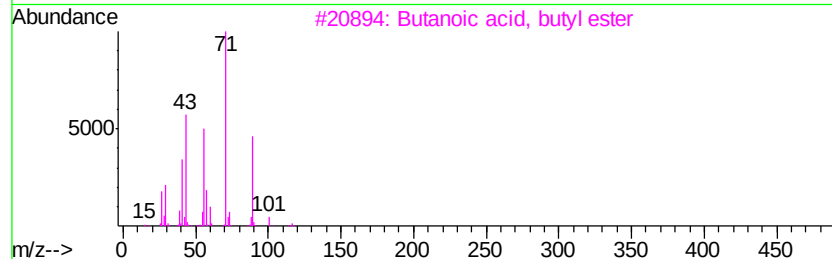
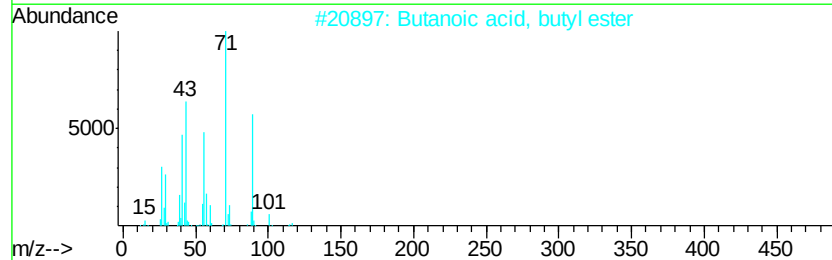
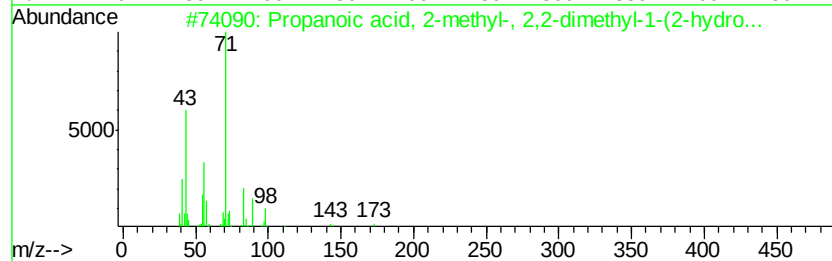
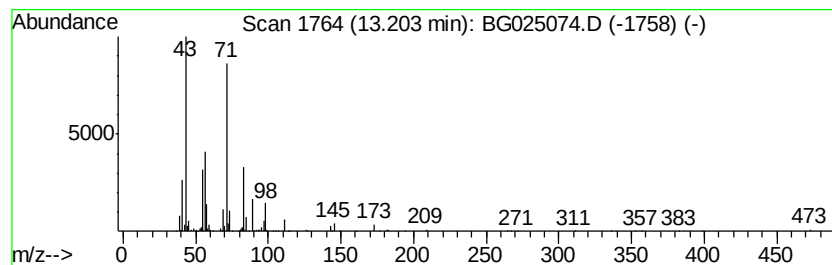
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 9 Propanoic acid, 2-methyl-, ... Concentration Rank 3

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 13.20 | 5.47 ng/ul | 278230 | Acenaphthene-d10 | 14.86 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Propanoic acid, 2-methyl-, 2,2-d... | 216 | C12H24O3 | 074367-33-2  | 78   |
| 2    |    |   | Butanoic acid, butyl ester          | 144 | C8H16O2  | 000109-21-7  | 59   |
| 3    |    |   | Butanoic acid, butyl ester          | 144 | C8H16O2  | 000109-21-7  | 45   |
| 4    |    |   | Propanoic acid, 2-methyl-, 2-met... | 144 | C8H16O2  | 000097-85-8  | 42   |
| 5    |    |   | Butyric acid, 3-pentadecyl ester    | 298 | C19H38O2 | 1000280-57-3 | 38   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG121016\  
Data File : BG025074.D  
Acq On : 10 Dec 2016 18:37  
Operator : UM/SJ  
Sample : H5861-07  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

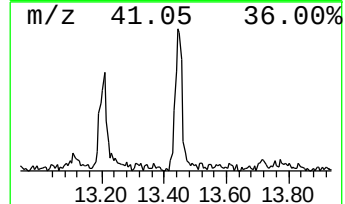
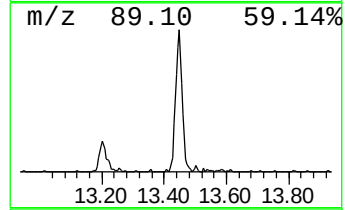
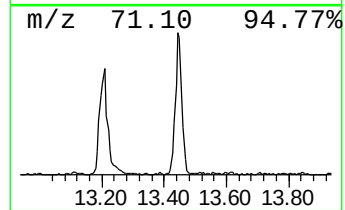
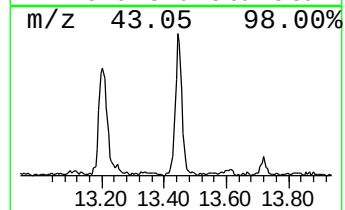
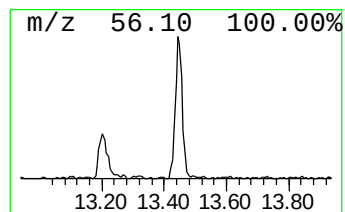
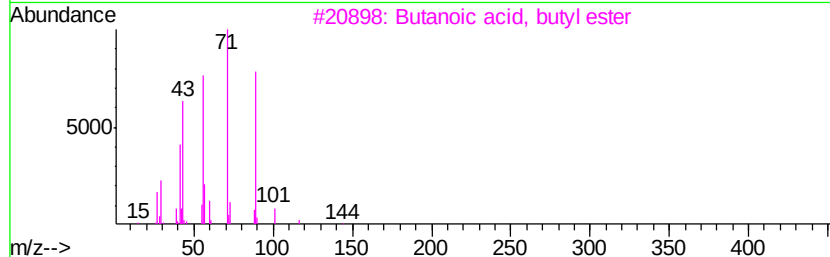
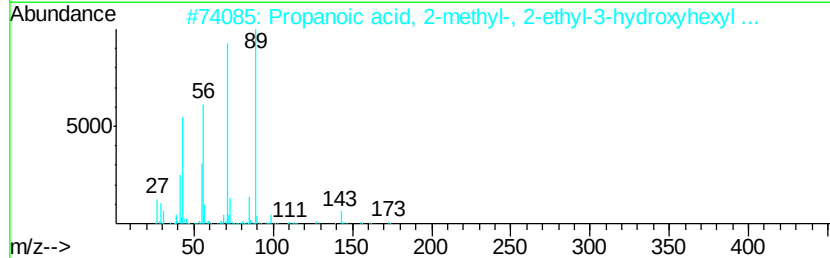
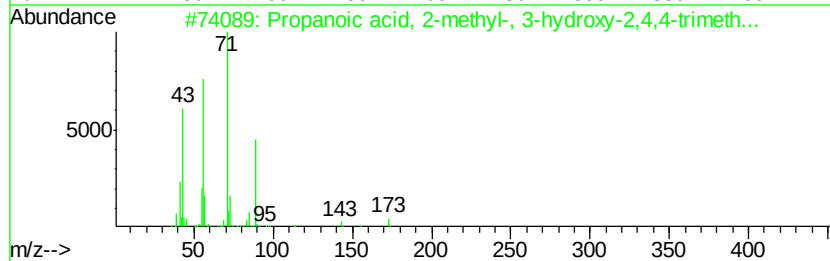
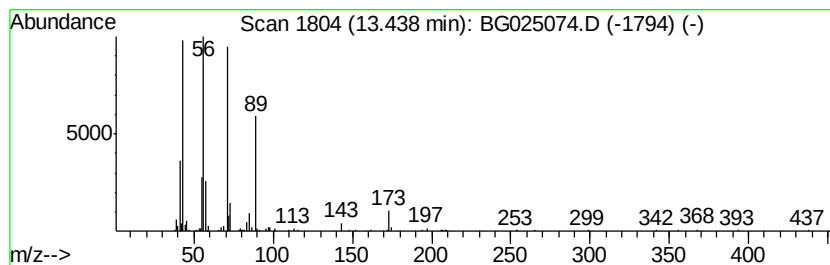
Instrument :  
BNA\_G  
ClientSampleId :  
DA8D9

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG113016.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 10 Propanoic acid, 2-methyl-, ... Concentration Rank 1

| R.T.    | EstConc                             | Area         | Relative to ISTD | R.T.      |
|---------|-------------------------------------|--------------|------------------|-----------|
| 13.44   | 7.26 ng/ul                          | 369183       | Acenaphthene-d10 | 14.86     |
| Hit# of | 5                                   | Tentative ID | MW MolForm       | CAS# Qual |
| 1       | Propanoic acid, 2-methyl-, 3-hyd... | 216 C12H24O3 | 074367-34-3      | 90        |
| 2       | Propanoic acid, 2-methyl-, 2-eth... | 216 C12H24O3 | 074367-31-0      | 78        |
| 3       | Butanoic acid, butyl ester          | 144 C8H16O2  | 000109-21-7      | 56        |
| 4       | Propanoic acid, 2-methyl-, butyl... | 144 C8H16O2  | 000097-87-0      | 53        |
| 5       | Propanoic acid, 2-methyl-, 2-met... | 144 C8H16O2  | 000097-85-8      | 38        |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG121016\  
Data File : BG025074.D  
Acq On : 10 Dec 2016 18:37  
Operator : UM/SJ  
Sample : H5861-07  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
DA8D9

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG113016.M  
Quant Title : SVOA CALIBRATION

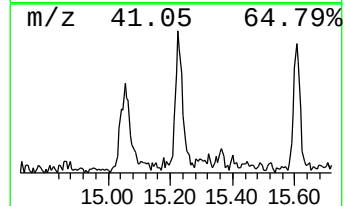
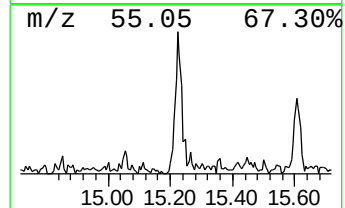
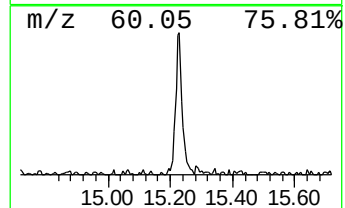
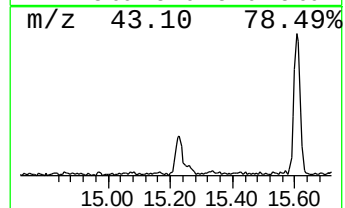
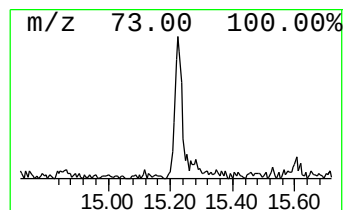
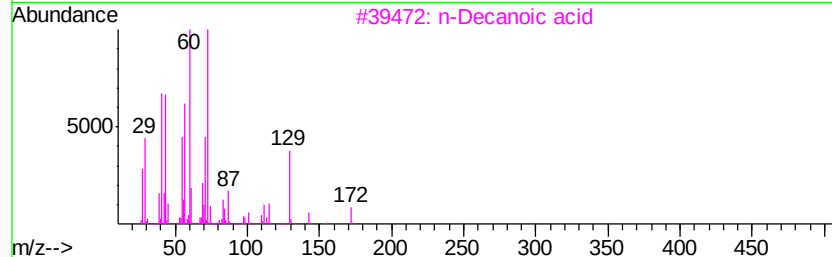
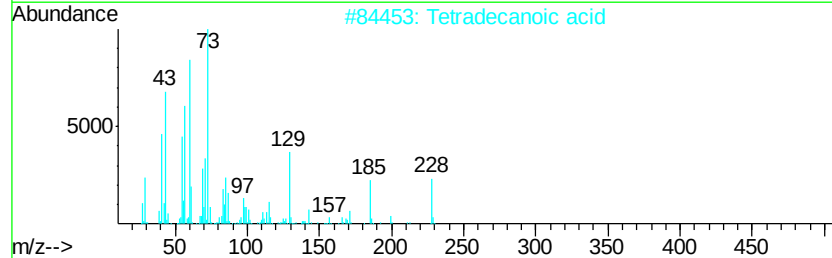
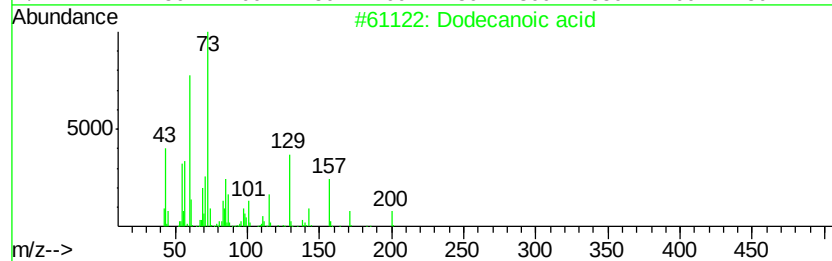
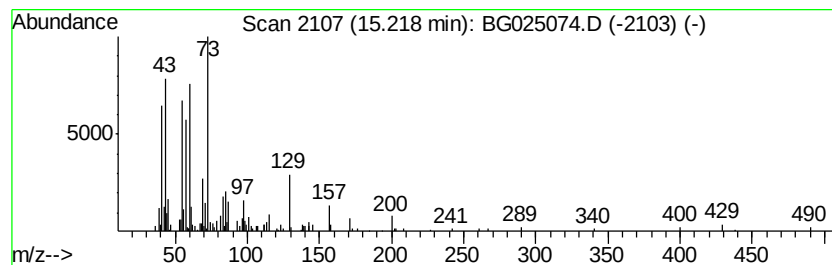
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 11 Dodecanoic acid Concentration Rank 6

| R.T.  | EstConc    | Area   | Relative to ISTD |  | R.T.  |
|-------|------------|--------|------------------|--|-------|
| 15.22 | 4.58 ng/ul | 233201 | Acenaphthene-d10 |  | 14.86 |

| Hit# | of | 5 | Tentative ID        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|---------------------|-----|----------|-------------|------|
| 1    |    |   | Dodecanoic acid     | 200 | C12H24O2 | 000143-07-7 | 60   |
| 2    |    |   | Tetradecanoic acid  | 228 | C14H28O2 | 000544-63-8 | 59   |
| 3    |    |   | n-Decanoic acid     | 172 | C10H20O2 | 000334-48-5 | 58   |
| 4    |    |   | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 56   |
| 5    |    |   | n-Decanoic acid     | 172 | C10H20O2 | 000334-48-5 | 52   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG121016\  
Data File : BG025074.D  
Acq On : 10 Dec 2016 18:37  
Operator : UM/SJ  
Sample : H5861-07  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
DA8D9

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG113016.M  
Quant Title : SVOA CALIBRATION

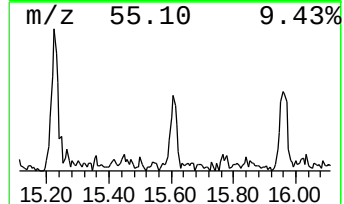
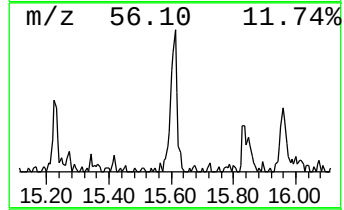
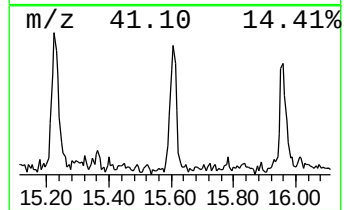
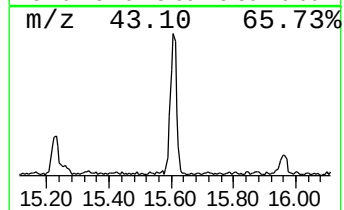
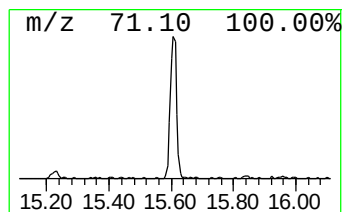
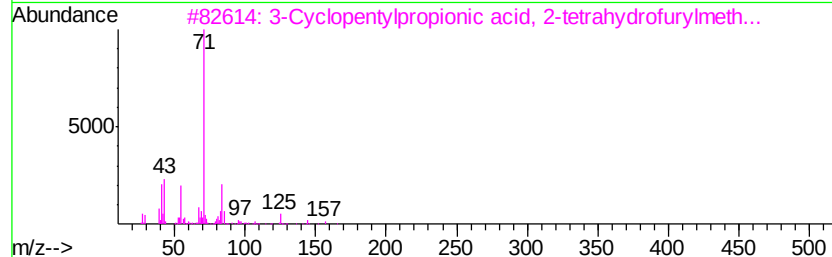
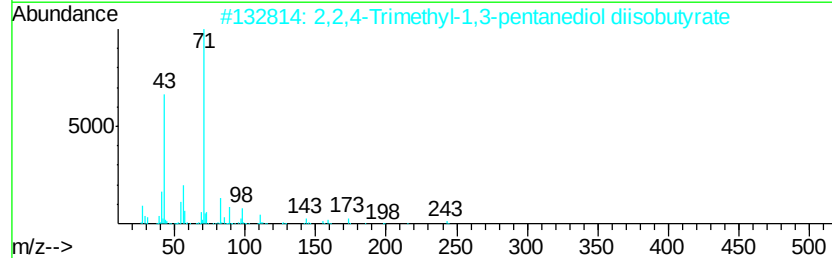
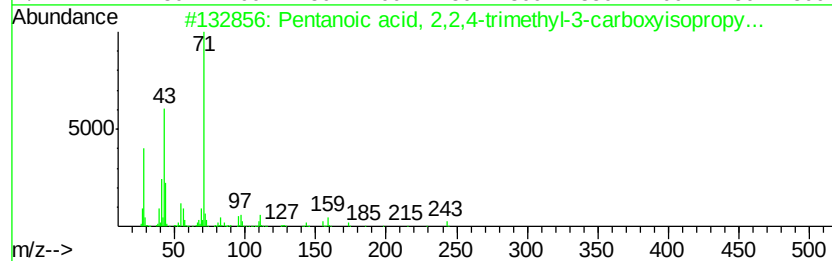
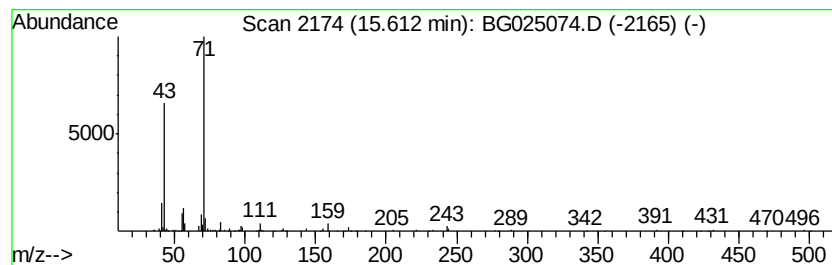
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 12 Pentanoic acid, 2,2,4-trime... Concentration Rank 4

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 15.61 | 4.96 ng/ul | 252204 | Acenaphthene-d10 | 14.86 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Pentanoic acid, 2,2,4-trimethyl-... | 286 | C16H30O4  | 1000140-77-5 | 72   |
| 2    |    |   | 2,2,4-Trimethyl-1,3-pentanediol ... | 286 | C16H30O4  | 006846-50-0  | 64   |
| 3    |    |   | 3-Cyclopentylpropionic acid, 2-t... | 226 | C13H22O3  | 1000293-46-9 | 53   |
| 4    |    |   | Sulfurous acid, octadecyl 2-pent... | 404 | C23H48O3S | 1000309-16-6 | 39   |
| 5    |    |   | Butyric acid, 1-propylpentyl ester  | 200 | C12H24O2  | 020286-46-8  | 36   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG121016\  
Data File : BG025074.D  
Acq On : 10 Dec 2016 18:37  
Operator : UM/SJ  
Sample : H5861-07  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
DA8D9

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG113016.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

| TIC Top Hit name     | RT    | EstConc | Units | Response | --Internal Standard-- |       |         |      |
|----------------------|-------|---------|-------|----------|-----------------------|-------|---------|------|
|                      |       |         |       |          | #                     | RT    | Resp    | Conc |
| unknown-01           | 5.31  | 2.0     | ng/ul | 46350    | 1                     | 8.24  | 463039  | 20.0 |
| Hexanoic acid        | 7.19  | 4.3     | ng/ul | 99301    | 1                     | 8.24  | 463039  | 20.0 |
| Hexanoic acid, 2-... | 9.49  | 3.7     | ng/ul | 85581    | 1                     | 8.24  | 463039  | 20.0 |
| unknown-02           | 9.54  | 2.0     | ng/ul | 47209    | 1                     | 8.24  | 463039  | 20.0 |
| .beta.-D-Ribopyra... | 10.33 | 2.8     | ng/ul | 92744    | 2                     | 11.06 | 658759  | 20.0 |
| unknown-03           | 10.81 | 2.1     | ng/ul | 68395    | 2                     | 11.06 | 658759  | 20.0 |
| Nonanoic acid        | 11.82 | 5.6     | ng/ul | 183863   | 2                     | 11.06 | 658759  | 20.0 |
| Propanoic acid, 2... | 13.20 | 5.5     | ng/ul | 278230   | 3                     | 14.86 | 1017690 | 20.0 |
| Propanoic acid, 2... | 13.44 | 7.3     | ng/ul | 369183   | 3                     | 14.86 | 1017690 | 20.0 |
| Dodecanoic acid      | 15.22 | 4.6     | ng/ul | 233201   | 3                     | 14.86 | 1017690 | 20.0 |
| Pentanoic acid, 2... | 15.61 | 5.0     | ng/ul | 252204   | 3                     | 14.86 | 1017690 | 20.0 |