

Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
 Data File : BN000542.D  
 Acq On : 22 Mar 2018 16:12  
 Operator : JU/SJ  
 Sample : J1905-11DL 10X  
 Misc :  
 ALS Vial : 12 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 DAM46DL

## Integration Parameters: LSCINT.P

Integrator: RTE  
 Smoothing : OFF  
 Sampling : 1  
 Start Thrs: 0.2  
 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: 5

Method : Z:\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN031918.M  
 Title : SVOA CALIBRATION

Signal : TIC

| peak # | R.T. min | first scan | max scan | last scan | PK TY | peak height | corr. area | corr. % max. | % of total |
|--------|----------|------------|----------|-----------|-------|-------------|------------|--------------|------------|
| 1      | 5.408    | 328        | 335      | 346       | rVB   | 1342537     | 2184439    | 2.78%        | 0.467%     |
| 2      | 6.349    | 490        | 495      | 502       | rVB2  | 1976450     | 3204927    | 4.08%        | 0.685%     |
| 3      | 6.861    | 574        | 582      | 587       | rBV   | 1983420     | 3532917    | 4.50%        | 0.755%     |
| 4      | 7.361    | 663        | 667      | 672       | rVB   | 1523468     | 2210345    | 2.81%        | 0.472%     |
| 5      | 7.419    | 672        | 677      | 681       | rBV   | 1601005     | 2393254    | 3.05%        | 0.511%     |
| 6      | 7.472    | 681        | 686      | 691       | rBV   | 1518310     | 2412107    | 3.07%        | 0.515%     |
| 7      | 7.531    | 691        | 696      | 702       | rVB3  | 2642798     | 4328540    | 5.51%        | 0.925%     |
| 8      | 7.813    | 741        | 744      | 752       | rVV   | 1744212     | 2765723    | 3.52%        | 0.591%     |
| 9      | 8.013    | 772        | 778      | 782       | rBV   | 5637987     | 9701634    | 12.34%       | 2.073%     |
| 10     | 8.049    | 782        | 784      | 793       | rVB   | 1959615     | 2863826    | 3.64%        | 0.612%     |
| 11     | 8.366    | 833        | 838      | 843       | rBV   | 1784475     | 2572195    | 3.27%        | 0.550%     |
| 12     | 8.484    | 843        | 858      | 862       | rBV   | 6180531     | 15259678   | 19.42%       | 3.261%     |
| 13     | 8.619    | 877        | 881      | 885       | rVB   | 2452824     | 3332285    | 4.24%        | 0.712%     |
| 14     | 8.860    | 914        | 922      | 927       | rBV   | 3560055     | 6720170    | 8.55%        | 1.436%     |
| 15     | 8.925    | 927        | 933      | 937       | rVB2  | 1059648     | 1834009    | 2.33%        | 0.392%     |
| 16     | 8.984    | 937        | 943      | 949       | rVB2  | 1597726     | 3685814    | 4.69%        | 0.788%     |
| 17     | 9.060    | 949        | 956      | 962       | rBV2  | 3762085     | 7066309    | 8.99%        | 1.510%     |
| 18     | 9.166    | 968        | 974      | 978       | rBV   | 1728894     | 2548052    | 3.24%        | 0.545%     |
| 19     | 9.372    | 1004       | 1009     | 1014      | rBV   | 1005707     | 1912162    | 2.43%        | 0.409%     |
| 20     | 9.431    | 1015       | 1019     | 1024      | rVB   | 1135671     | 1870401    | 2.38%        | 0.400%     |
| 21     | 9.531    | 1031       | 1036     | 1045      | rVB3  | 2039442     | 4126203    | 5.25%        | 0.882%     |
| 22     | 9.643    | 1045       | 1055     | 1060      | rVB   | 6100115     | 11785887   | 15.00%       | 2.519%     |
| 23     | 9.778    | 1066       | 1078     | 1085      | rBV8  | 1219217     | 5369916    | 6.83%        | 1.148%     |
| 24     | 9.872    | 1085       | 1094     | 1098      | rBV4  | 1368526     | 4425025    | 5.63%        | 0.946%     |
| 25     | 10.278   | 1157       | 1163     | 1166      | rBV   | 1661637     | 3054238    | 3.89%        | 0.653%     |
| 26     | 10.360   | 1172       | 1177     | 1183      | rVB4  | 1375263     | 2481495    | 3.16%        | 0.530%     |
| 27     | 10.490   | 1196       | 1199     | 1204      | rVB2  | 1443238     | 2074477    | 2.64%        | 0.443%     |
| 28     | 10.554   | 1205       | 1210     | 1217      | rBV4  | 872148      | 2159336    | 2.75%        | 0.461%     |
| 29     | 10.637   | 1219       | 1224     | 1230      | rVV4  | 1769642     | 3217240    | 4.09%        | 0.688%     |
| 30     | 10.754   | 1237       | 1244     | 1248      | rVV2  | 2265646     | 4817903    | 6.13%        | 1.030%     |
| 31     | 10.996   | 1280       | 1285     | 1291      | rVB   | 1152858     | 1884368    | 2.40%        | 0.403%     |
| 32     | 11.119   | 1292       | 1306     | 1309      | rBV6  | 908408      | 2422092    | 3.08%        | 0.518%     |
| 33     | 11.196   | 1309       | 1319     | 1321      | rBV3  | 3461049     | 8069705    | 10.27%       | 1.725%     |
| 34     | 11.219   | 1321       | 1323     | 1327      | rVB   | 3631748     | 4331646    | 5.51%        | 0.926%     |

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 Stop Thrs : 0

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 Max Peaks: 100  
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >  
 Peak separation: 5

Method : Z:\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN031918.M  
 Title : SVOA CALIBRATION

|    |        |      |      |      |      |         |          |        |        |
|----|--------|------|------|------|------|---------|----------|--------|--------|
| 35 | 11.337 | 1335 | 1343 | 1346 | rBV2 | 2117205 | 3778606  | 4.81%  | 0.808% |
| 36 | 11.372 | 1346 | 1349 | 1355 | rVB4 | 1618033 | 2684929  | 3.42%  | 0.574% |
| 37 | 11.449 | 1355 | 1362 | 1368 | rBV2 | 1109311 | 2077729  | 2.64%  | 0.444% |
| 38 | 11.543 | 1368 | 1378 | 1382 | rBV6 | 934734  | 2557825  | 3.25%  | 0.547% |
| 39 | 11.601 | 1382 | 1388 | 1397 | rVB  | 3693283 | 6196915  | 7.89%  | 1.324% |
| 40 | 11.731 | 1402 | 1410 | 1415 | rVB  | 1610580 | 2700045  | 3.44%  | 0.577% |
| 41 | 11.843 | 1422 | 1429 | 1436 | rVB4 | 974228  | 1973922  | 2.51%  | 0.422% |
| 42 | 11.931 | 1436 | 1444 | 1448 | rBV5 | 1075833 | 2606682  | 3.32%  | 0.557% |
| 43 | 12.119 | 1472 | 1476 | 1481 | rVV3 | 1372653 | 2770539  | 3.53%  | 0.592% |
| 44 | 12.225 | 1489 | 1494 | 1499 | rVV3 | 1648023 | 3217372  | 4.09%  | 0.688% |
| 45 | 12.313 | 1505 | 1509 | 1520 | rVB  | 2539559 | 4170253  | 5.31%  | 0.891% |
| 46 | 12.466 | 1530 | 1535 | 1545 | rVB2 | 2963921 | 4798335  | 6.11%  | 1.025% |
| 47 | 12.613 | 1555 | 1560 | 1564 | rBV  | 3596521 | 6031384  | 7.67%  | 1.289% |
| 48 | 12.660 | 1564 | 1568 | 1574 | rVB  | 2465248 | 3649714  | 4.64%  | 0.780% |
| 49 | 12.748 | 1575 | 1583 | 1585 | rBV2 | 1148642 | 2158999  | 2.75%  | 0.461% |
| 50 | 12.890 | 1604 | 1607 | 1614 | rVB  | 1321855 | 1899143  | 2.42%  | 0.406% |
| 51 | 13.019 | 1625 | 1629 | 1632 | rBV2 | 1268534 | 1876556  | 2.39%  | 0.401% |
| 52 | 13.107 | 1637 | 1644 | 1652 | rVB3 | 666389  | 1759718  | 2.24%  | 0.376% |
| 53 | 13.260 | 1659 | 1670 | 1673 | rBV7 | 932429  | 2843946  | 3.62%  | 0.608% |
| 54 | 13.548 | 1714 | 1719 | 1725 | rBV  | 1550162 | 2675207  | 3.40%  | 0.572% |
| 55 | 13.766 | 1751 | 1756 | 1761 | rVB2 | 1095168 | 1788058  | 2.28%  | 0.382% |
| 56 | 13.831 | 1761 | 1767 | 1777 | rBV  | 4133852 | 7610093  | 9.68%  | 1.626% |
| 57 | 14.072 | 1804 | 1808 | 1821 | rVB3 | 1409097 | 3094210  | 3.94%  | 0.661% |
| 58 | 14.219 | 1822 | 1833 | 1842 | rBV7 | 1437708 | 5423921  | 6.90%  | 1.159% |
| 59 | 14.360 | 1851 | 1857 | 1864 | rBV5 | 1481364 | 4389585  | 5.59%  | 0.938% |
| 60 | 14.478 | 1875 | 1877 | 1881 | rVB3 | 1761756 | 2173950  | 2.77%  | 0.465% |
| 61 | 14.642 | 1902 | 1905 | 1915 | rVB5 | 2238772 | 4456466  | 5.67%  | 0.952% |
| 62 | 14.901 | 1943 | 1949 | 1953 | rVB  | 7096072 | 12058584 | 15.34% | 2.577% |
| 63 | 15.054 | 1968 | 1975 | 1982 | rBV5 | 2022097 | 4697944  | 5.98%  | 1.004% |
| 64 | 15.154 | 1987 | 1992 | 1997 | rBV  | 4314800 | 6433523  | 8.19%  | 1.375% |
| 65 | 15.225 | 2000 | 2004 | 2007 | rBV3 | 1067176 | 1763198  | 2.24%  | 0.377% |
| 66 | 15.325 | 2015 | 2021 | 2028 | rBV2 | 2383943 | 6045703  | 7.69%  | 1.292% |
| 67 | 15.419 | 2034 | 2037 | 2044 | rVB4 | 1161568 | 2230970  | 2.84%  | 0.477% |
| 68 | 15.501 | 2046 | 2051 | 2055 | rVV2 | 1895069 | 2938902  | 3.74%  | 0.628% |
| 69 | 15.560 | 2057 | 2061 | 2070 | rVV8 | 1795693 | 4765014  | 6.06%  | 1.018% |
| 70 | 15.672 | 2070 | 2080 | 2090 | rVV3 | 8463117 | 20519502 | 26.11% | 4.385% |
| 71 | 15.778 | 2092 | 2098 | 2101 | rVV2 | 4512317 | 7615821  | 9.69%  | 1.628% |

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Method : Z:\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN031918.M  
 Title : SVOA CALIBRATION

|     |        |      |      |      |      |          |          |         |         |
|-----|--------|------|------|------|------|----------|----------|---------|---------|
| 72  | 15.831 | 2101 | 2107 | 2114 | rVB2 | 4491496  | 8923942  | 11.36%  | 1.907%  |
| 73  | 16.113 | 2151 | 2155 | 2160 | rVB  | 1262628  | 1887035  | 2.40%   | 0.403%  |
| 74  | 16.231 | 2172 | 2175 | 2179 | rVB  | 1749086  | 2485832  | 3.16%   | 0.531%  |
| 75  | 16.448 | 2209 | 2212 | 2219 | rVB4 | 1120753  | 1746282  | 2.22%   | 0.373%  |
| 76  | 16.619 | 2237 | 2241 | 2248 | rVB2 | 2008489  | 3268246  | 4.16%   | 0.698%  |
| 77  | 16.689 | 2248 | 2253 | 2255 | rBV  | 4323893  | 6666405  | 8.48%   | 1.425%  |
| 78  | 17.048 | 2310 | 2314 | 2321 | rVB6 | 1267938  | 2399890  | 3.05%   | 0.513%  |
| 79  | 17.501 | 2372 | 2391 | 2395 | rBV2 | 13963236 | 78590156 | 100.00% | 16.796% |
| 80  | 17.831 | 2443 | 2447 | 2452 | rVB3 | 1893891  | 3045148  | 3.87%   | 0.651%  |
| 81  | 18.013 | 2473 | 2478 | 2481 | rBV4 | 1026496  | 2193628  | 2.79%   | 0.469%  |
| 82  | 18.207 | 2506 | 2511 | 2520 | rVB  | 3751398  | 5763717  | 7.33%   | 1.232%  |
| 83  | 18.378 | 2537 | 2540 | 2546 | rVB  | 1890180  | 2263775  | 2.88%   | 0.484%  |
| 84  | 18.883 | 2622 | 2626 | 2630 | rVB  | 3171400  | 3782980  | 4.81%   | 0.808%  |
| 85  | 19.501 | 2727 | 2731 | 2733 | rBV  | 2845949  | 3583698  | 4.56%   | 0.766%  |
| 86  | 19.525 | 2733 | 2735 | 2739 | rVB2 | 2372505  | 2631066  | 3.35%   | 0.562%  |
| 87  | 19.654 | 2752 | 2757 | 2761 | rVB2 | 1413347  | 2112355  | 2.69%   | 0.451%  |
| 88  | 20.066 | 2823 | 2827 | 2831 | rVB  | 2207009  | 2629855  | 3.35%   | 0.562%  |
| 89  | 20.589 | 2913 | 2916 | 2921 | rBV2 | 2008994  | 2469165  | 3.14%   | 0.528%  |
| 90  | 21.077 | 2996 | 2999 | 3003 | rBV  | 1808708  | 1908759  | 2.43%   | 0.408%  |
| 91  | 21.542 | 3073 | 3078 | 3085 | rVB  | 3852900  | 4905197  | 6.24%   | 1.048%  |
| 92  | 21.883 | 3131 | 3136 | 3141 | rBV2 | 1690993  | 2554846  | 3.25%   | 0.546%  |
| 93  | 21.983 | 3149 | 3153 | 3160 | rVB  | 1801861  | 2457327  | 3.13%   | 0.525%  |
| 94  | 22.366 | 3214 | 3218 | 3223 | rBV  | 1674742  | 2269106  | 2.89%   | 0.485%  |
| 95  | 22.460 | 3229 | 3234 | 3240 | rVB2 | 2819322  | 4188120  | 5.33%   | 0.895%  |
| 96  | 22.966 | 3316 | 3320 | 3323 | rBV  | 1172139  | 1851113  | 2.36%   | 0.396%  |
| 97  | 23.154 | 3348 | 3352 | 3357 | rVB  | 1037104  | 1898693  | 2.42%   | 0.406%  |
| 98  | 23.307 | 3373 | 3378 | 3384 | rVB  | 1338155  | 2494622  | 3.17%   | 0.533%  |
| 99  | 23.524 | 3411 | 3415 | 3427 | rVV2 | 832185   | 1795263  | 2.28%   | 0.384%  |
| 100 | 24.371 | 3553 | 3559 | 3565 | rVB  | 1096903  | 2101650  | 2.67%   | 0.449%  |

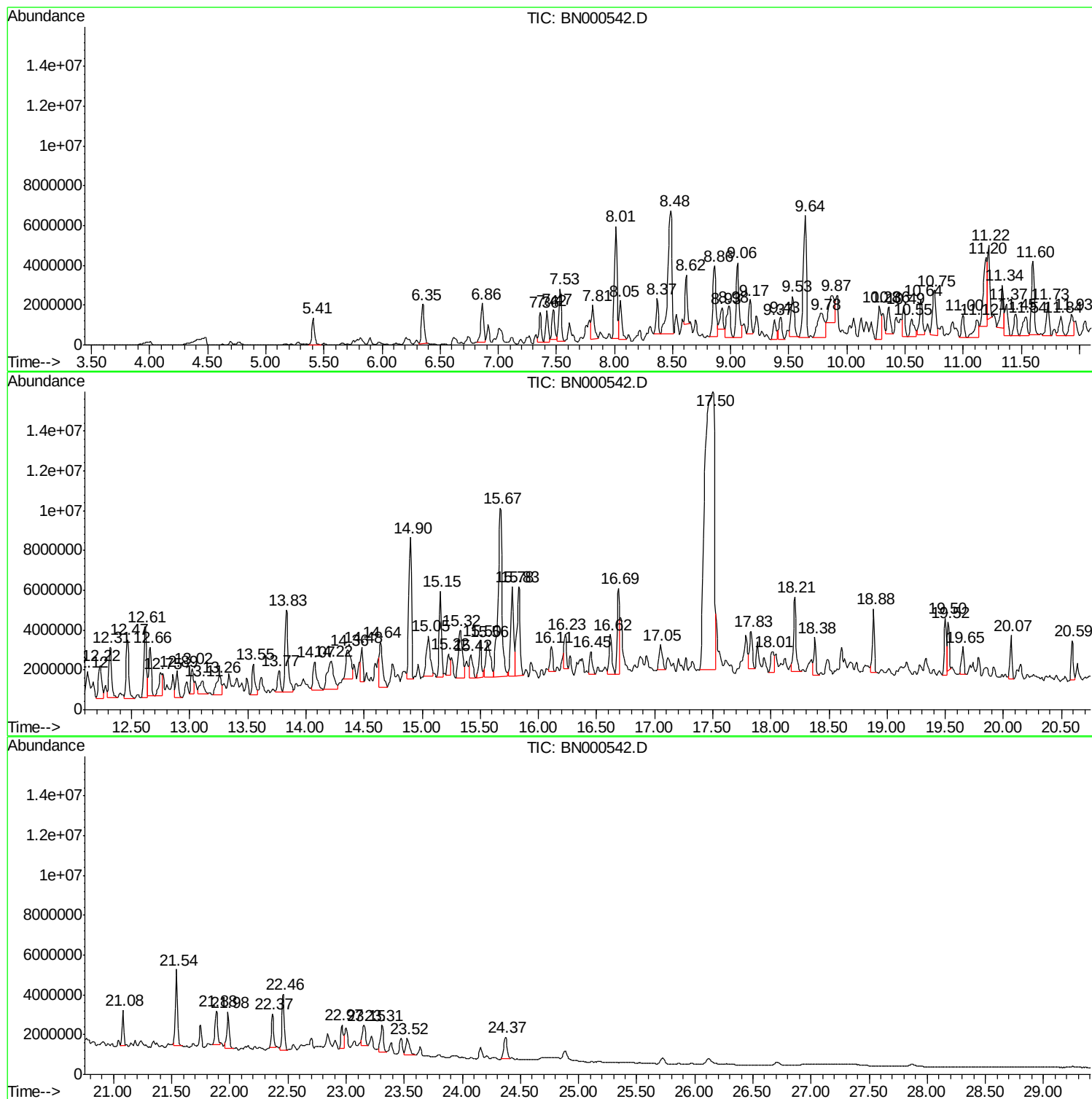
Sum of corrected areas: 467919452

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Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P



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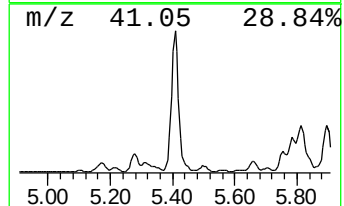
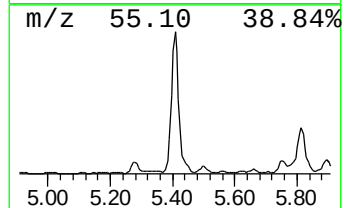
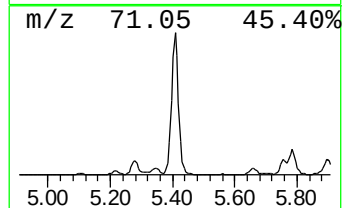
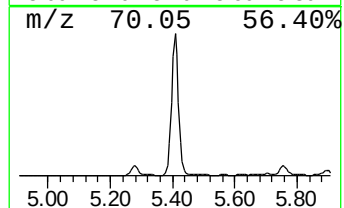
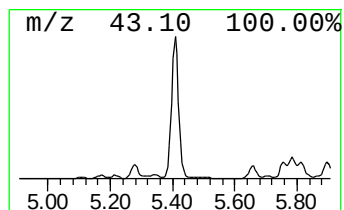
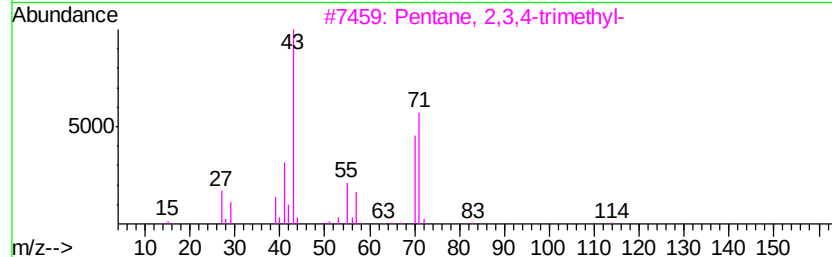
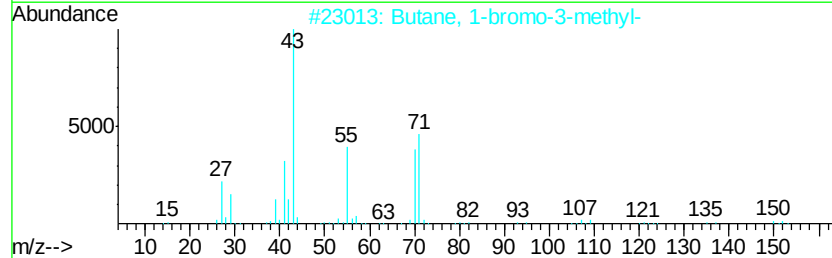
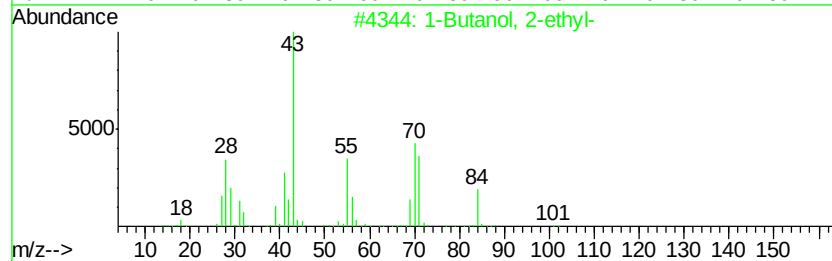
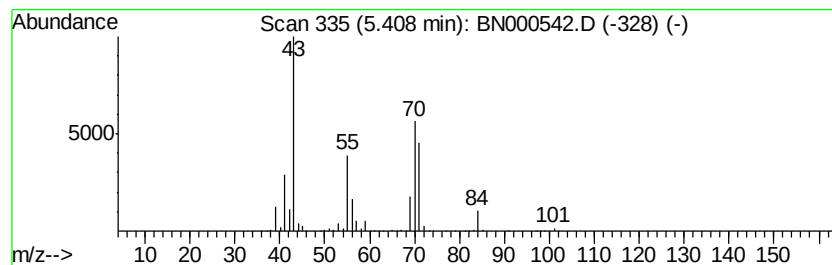
Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 1 1-Butanol, 2-ethyl- Concentration Rank 47

| R.T. | EstConc     | Area    | Relative to ISTD       | R.T. |
|------|-------------|---------|------------------------|------|
| 5.41 | 16.99 ng/ul | 2184440 | 1,4-Dichlorobenzene-d4 | 8.42 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | 1-Butanol, 2-ethyl-                 | 102 | C6H14O  | 000097-95-0 | 83   |
| 2    |    | Butane, 1-bromo-3-methyl-           | 150 | C5H11Br | 000107-82-4 | 64   |
| 3    |    | Pentane, 2,3,4-trimethyl-           | 114 | C8H18   | 000565-75-3 | 59   |
| 4    |    | Pentane, 3-ethyl-                   | 100 | C7H16   | 000617-78-7 | 53   |
| 5    |    | Propanoic acid, 2-methyl-, 2-met... | 158 | C9H18O2 | 002445-69-4 | 53   |



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Quant Title : SVOA CALIBRATION

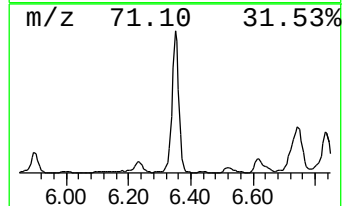
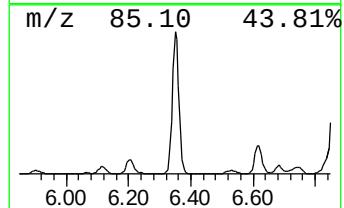
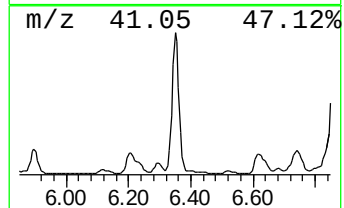
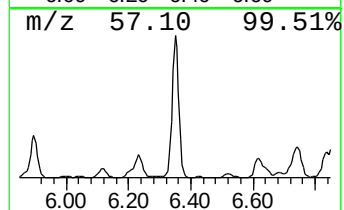
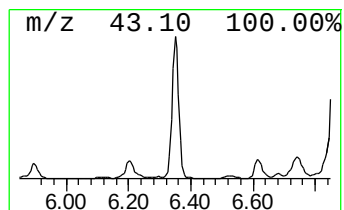
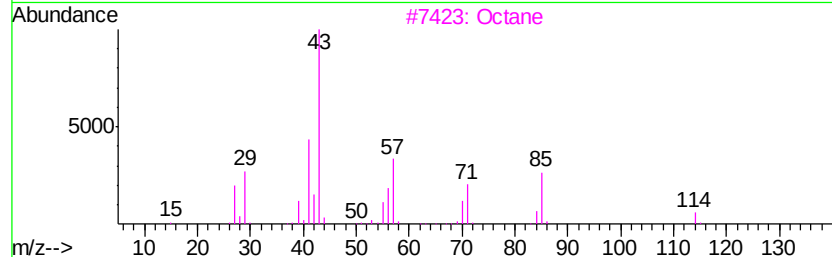
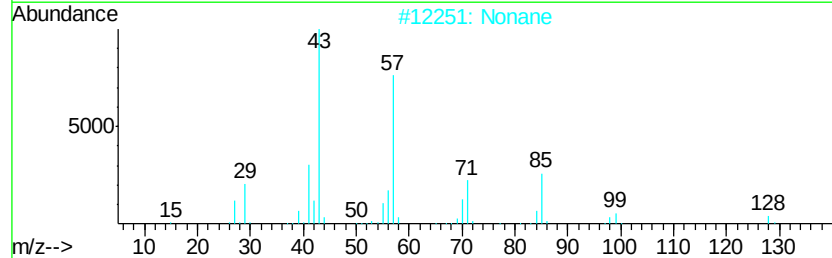
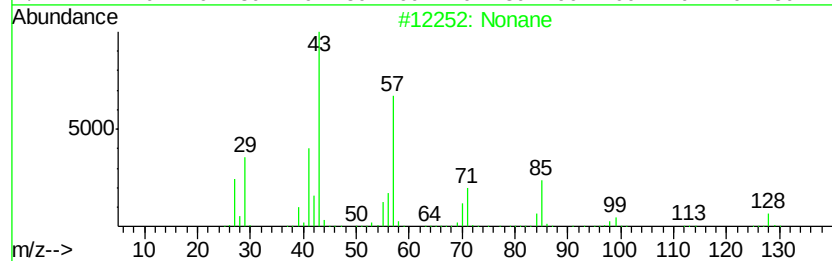
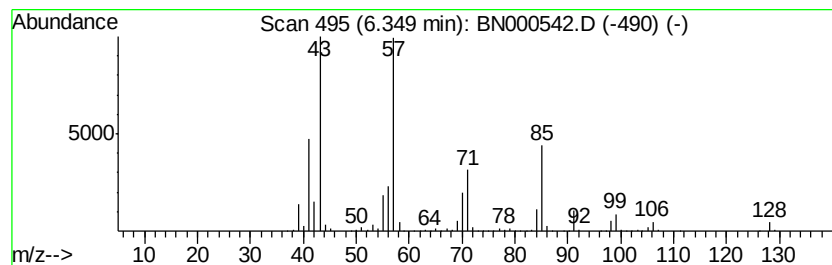
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 21

| R.T. | EstConc     | Area    | Relative to ISTD       | R.T. |
|------|-------------|---------|------------------------|------|
| 6.35 | 24.92 ng/ul | 3204930 | 1,4-Dichlorobenzene-d4 | 8.42 |

| Hit# | of | Tentative ID            | MW  | MolForm | CAS#         | Qual |
|------|----|-------------------------|-----|---------|--------------|------|
| 1    | 5  | Nonane                  | 128 | C9H20   | 000111-84-2  | 95   |
| 2    |    | Nonane                  | 128 | C9H20   | 000111-84-2  | 91   |
| 3    |    | Octane                  | 114 | C8H18   | 000111-65-9  | 72   |
| 4    |    | Octane                  | 114 | C8H18   | 000111-65-9  | 64   |
| 5    |    | 4,4-Dimethylcyclooctene | 138 | C10H18  | 1000113-56-7 | 59   |



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Data File : BN000542.D  
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Operator : JU/SJ  
Sample : J1905-11DL 10X  
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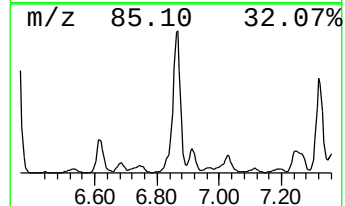
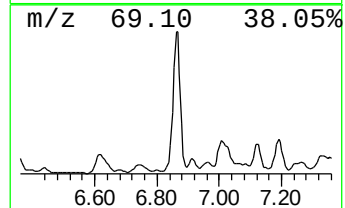
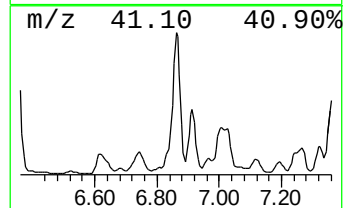
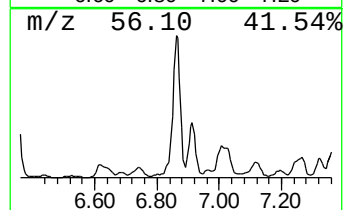
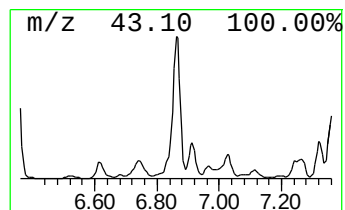
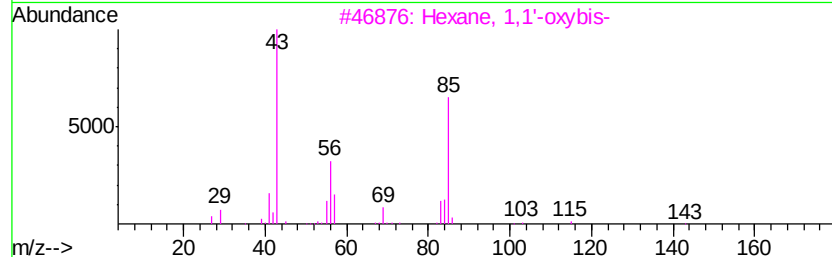
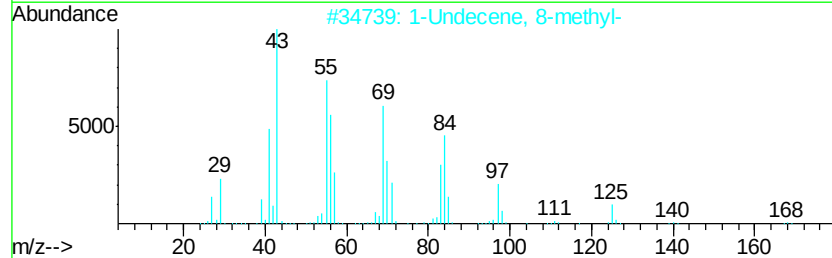
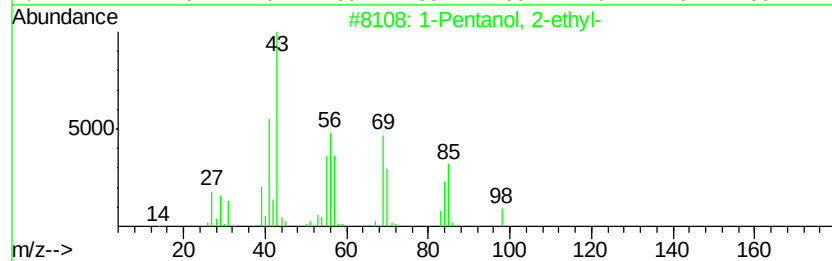
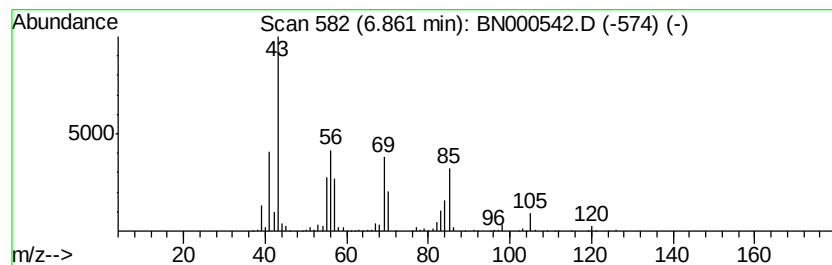
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 3 1-Pentanol, 2-ethyl- Concentration Rank 17

| R.T.    | EstConc               | Area         | Relative to ISTD       | R.T.    |             |      |
|---------|-----------------------|--------------|------------------------|---------|-------------|------|
| 6.86    | 27.47 ng/ul           | 3532920      | 1,4-Dichlorobenzene-d4 | 8.42    |             |      |
| Hit# of | 5                     | Tentative ID | MW                     | MolForm | CAS#        | Qual |
| 1       | 1-Pentanol, 2-ethyl-  |              | 116                    | C7H16O  | 027522-11-8 | 78   |
| 2       | 1-Undecene, 8-methyl- |              | 168                    | C12H24  | 074630-40-3 | 64   |
| 3       | Hexane, 1,1'-oxybis-  |              | 186                    | C12H26O | 000112-58-3 | 59   |
| 4       | Hexane, 1,1'-oxybis-  |              | 186                    | C12H26O | 000112-58-3 | 59   |
| 5       | Hexane, 1-propoxy-    |              | 144                    | C9H20O  | 053685-78-2 | 47   |





Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000542.D  
Acq On : 22 Mar 2018 16:12  
Operator : JU/SJ  
Sample : J1905-11DL 10X  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
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Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

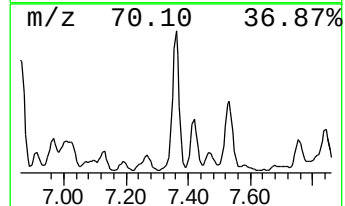
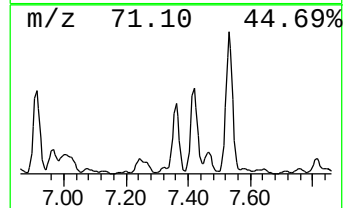
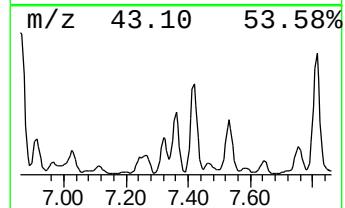
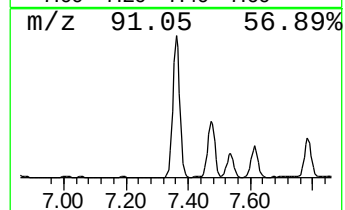
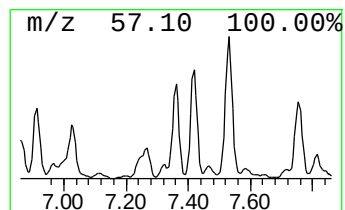
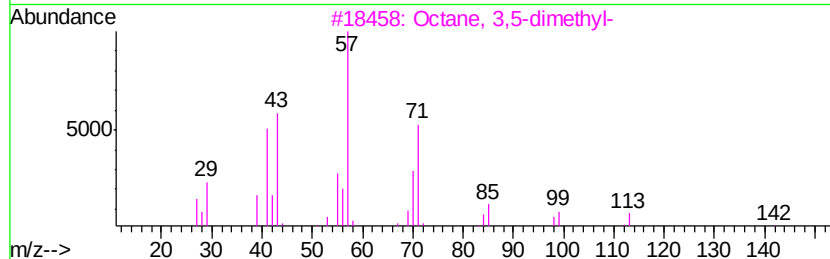
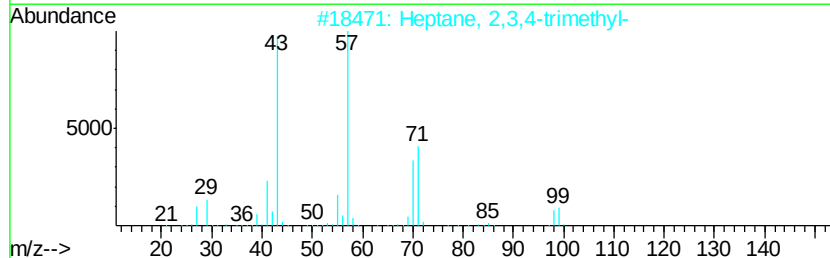
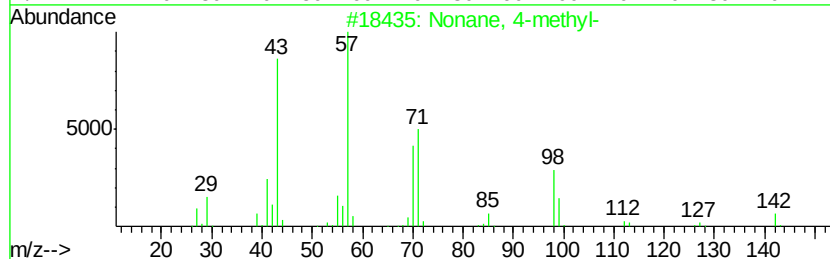
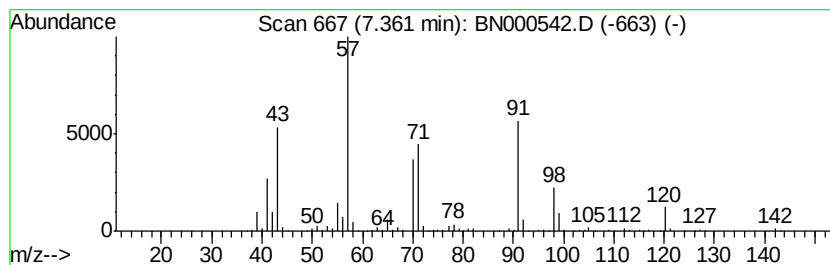
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 45

| R.T. | EstConc     | Area    | Relative to ISTD       | R.T. |
|------|-------------|---------|------------------------|------|
| 7.36 | 17.19 ng/ul | 2210350 | 1,4-Dichlorobenzene-d4 | 8.42 |

| Hit# | of | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------|-----|---------|-------------|------|
| 1    | 5  | Nonane, 4-methyl-             | 142 | C10H22  | 017301-94-9 | 64   |
| 2    |    | Heptane, 2,3,4-trimethyl-     | 142 | C10H22  | 052896-95-4 | 50   |
| 3    |    | Octane, 3,5-dimethyl-         | 142 | C10H22  | 015869-93-9 | 47   |
| 4    |    | Pentane, 2,2,3,3-tetramethyl- | 128 | C9H20   | 007154-79-2 | 47   |
| 5    |    | Undecane, 4,5-dimethyl-       | 184 | C13H28  | 017312-79-7 | 47   |





Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000542.D  
Acq On : 22 Mar 2018 16:12  
Operator : JU/SJ  
Sample : J1905-11DL 10X  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM46DL

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

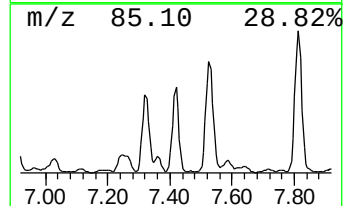
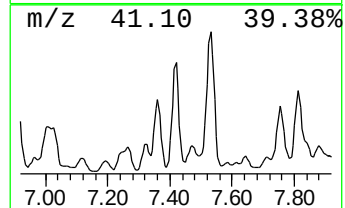
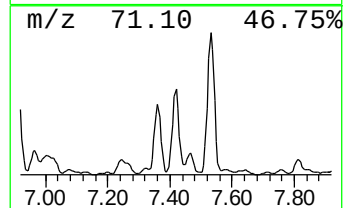
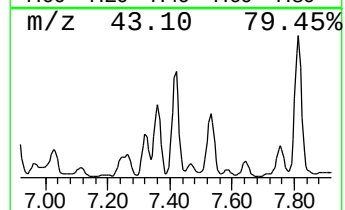
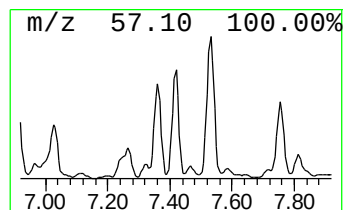
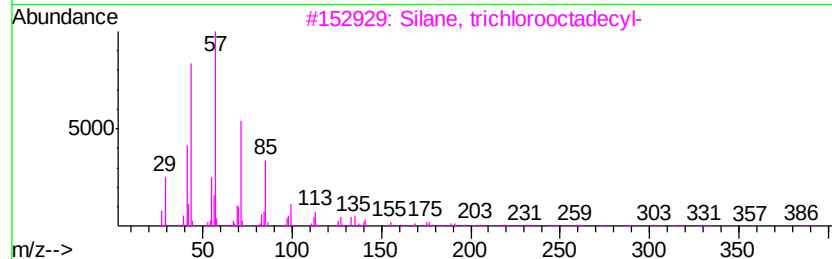
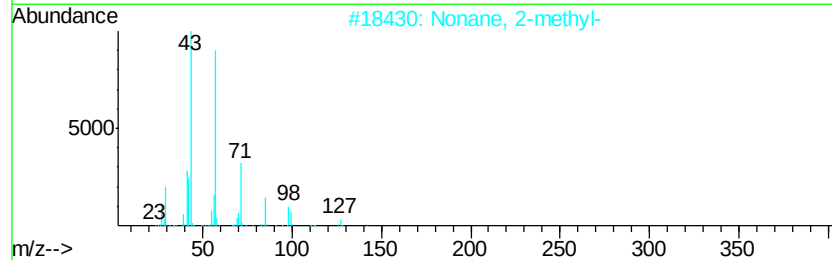
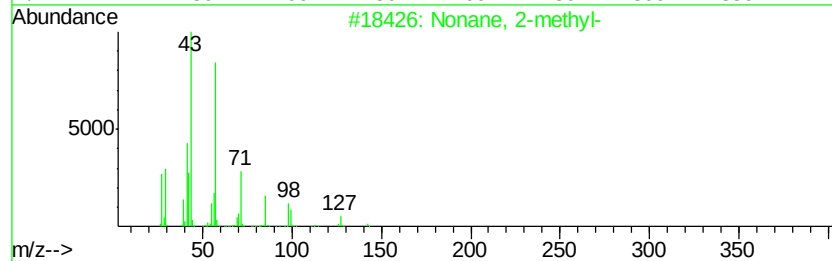
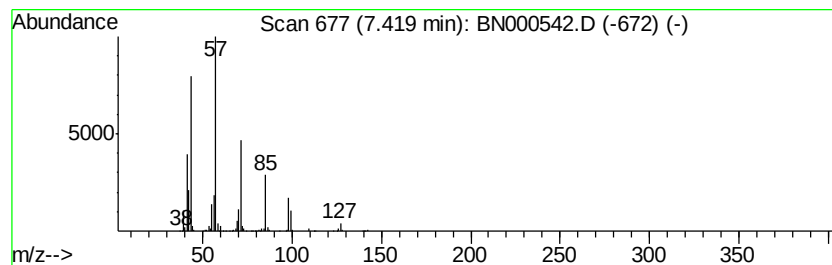
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 41

| R.T. | EstConc     | Area    | Relative to ISTD       | R.T. |
|------|-------------|---------|------------------------|------|
| 7.42 | 18.61 ng/ul | 2393250 | 1,4-Dichlorobenzene-d4 | 8.42 |

| Hit# of | 5 | Tentative ID                | MW  | MolForm     | CAS#        | Qual |
|---------|---|-----------------------------|-----|-------------|-------------|------|
| 1       |   | Nonane, 2-methyl-           | 142 | C10H22      | 000871-83-0 | 86   |
| 2       |   | Nonane, 2-methyl-           | 142 | C10H22      | 000871-83-0 | 81   |
| 3       |   | Silane, trichlorooctadecyl- | 386 | C18H37Cl3Si | 000112-04-9 | 78   |
| 4       |   | Hexadecane                  | 226 | C16H34      | 000544-76-3 | 72   |
| 5       |   | 1-Iodo-2-methylundecane     | 296 | C12H25I     | 073105-67-6 | 72   |



Data Path : Z:\HPCHEM1\BNA\_N\DATA\BN032218\  
Data File : BN000542.D  
Acq On : 22 Mar 2018 16:12  
Operator : JU/SJ  
Sample : J1905-11DL 10X  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

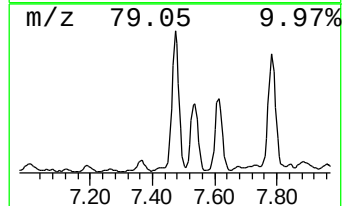
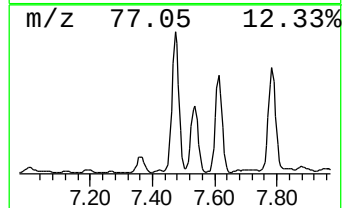
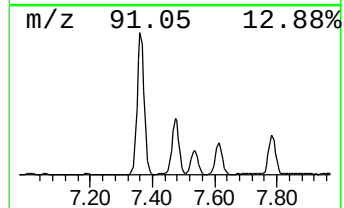
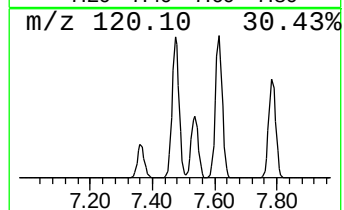
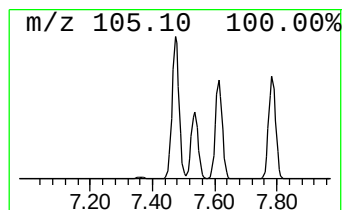
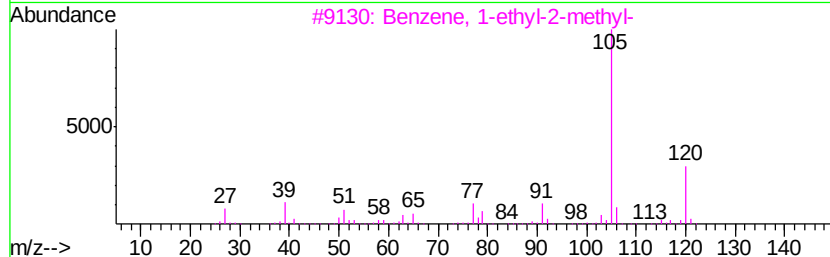
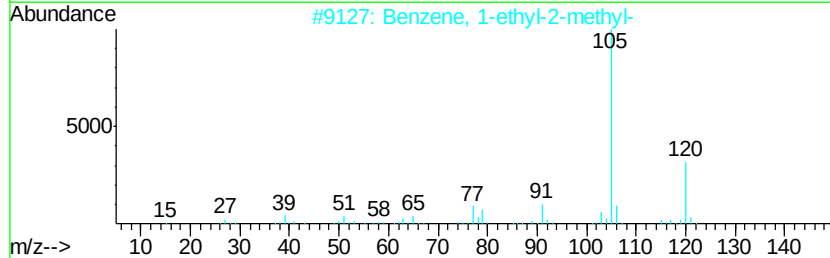
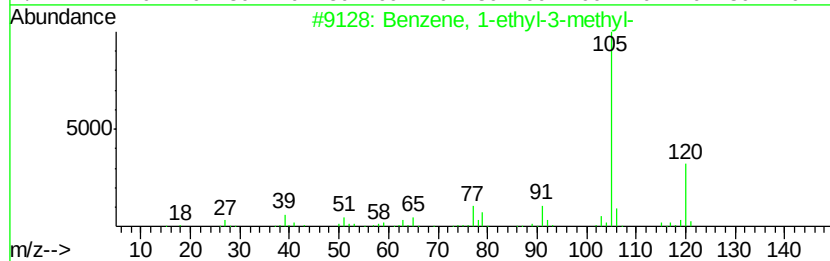
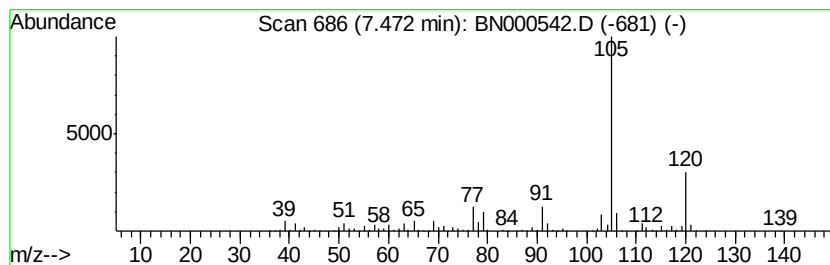
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BNA\_N  
ClientSampled :  
DAM46DL

Quant Method : Z:\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 6 Benzene, 1-ethyl-3-methyl- Concentration Rank 39

| R.T.    | EstConc     | Area                       | Relative to ISTD       | R.T.           |
|---------|-------------|----------------------------|------------------------|----------------|
| 7.47    | 18.76 ng/ul | 2412110                    | 1,4-Dichlorobenzene-d4 | 8.42           |
| Hit# of | 5           | Tentative ID               | MW MolForm             | CAS# Qual      |
| 1       |             | Benzene, 1-ethyl-3-methyl- | 120 C9H12              | 000620-14-4 94 |
| 2       |             | Benzene, 1-ethyl-2-methyl- | 120 C9H12              | 000611-14-3 94 |
| 3       |             | Benzene, 1-ethyl-2-methyl- | 120 C9H12              | 000611-14-3 94 |
| 4       |             | Benzene, 1,3,5-trimethyl-  | 120 C9H12              | 000108-67-8 90 |
| 5       |             | Benzene, 1-ethyl-3-methyl- | 120 C9H12              | 000620-14-4 90 |



Data Path : Z:\HPCHEM1\BNA\_N\DATA\BN032218\  
Data File : BN000542.D  
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Operator : JU/SJ  
Sample : J1905-11DL 10X  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
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Quant Method : Z:\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN031918.M  
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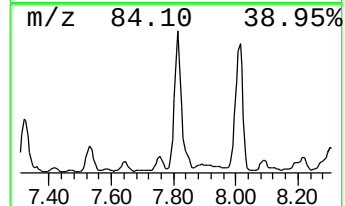
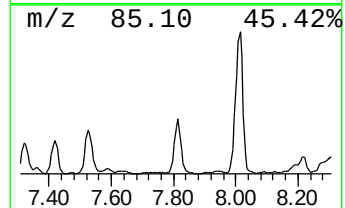
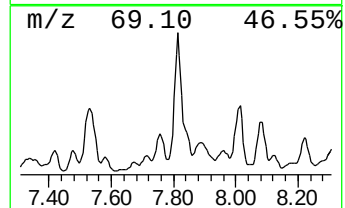
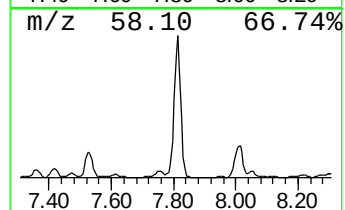
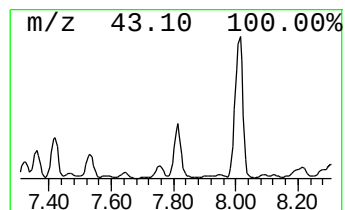
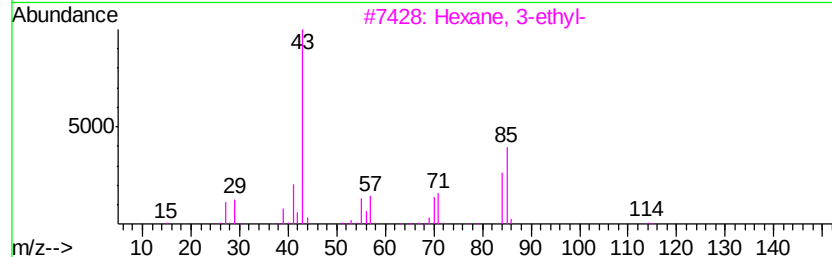
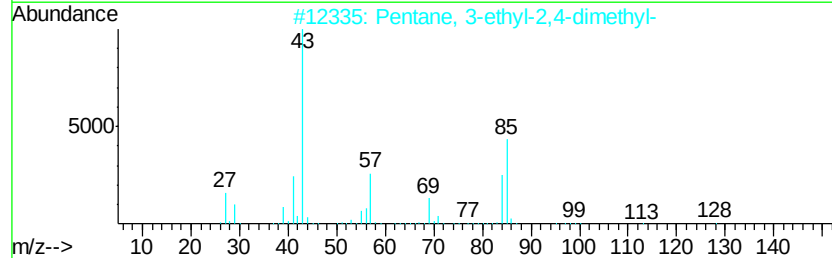
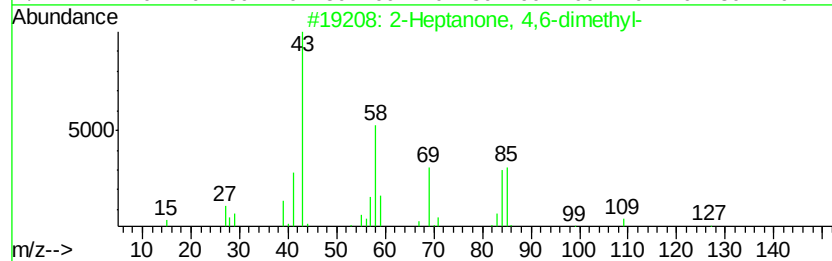
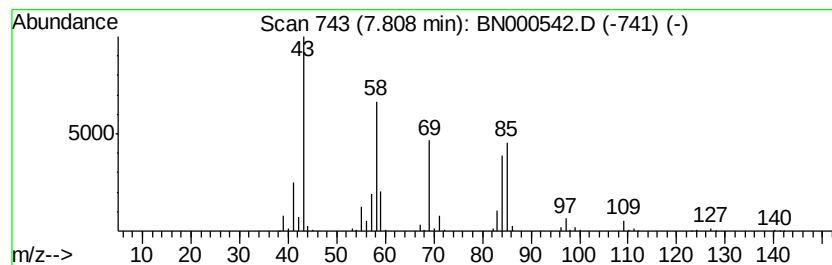
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 7 2-Heptanone, 4,6-dimethyl- Concentration Rank 29

| R.T. | EstConc     | Area    | Relative to ISTD       | R.T. |
|------|-------------|---------|------------------------|------|
| 7.81 | 21.50 ng/ul | 2765720 | 1,4-Dichlorobenzene-d4 | 8.42 |

| Hit# | of | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------|-----|---------|-------------|------|
| 1    | 5  | 2-Heptanone, 4,6-dimethyl-     | 142 | C9H18O  | 019549-80-5 | 90   |
| 2    |    | Pentane, 3-ethyl-2,4-dimethyl- | 128 | C9H20   | 001068-87-7 | 43   |
| 3    |    | Hexane, 3-ethyl-               | 114 | C8H18   | 000619-99-8 | 43   |
| 4    |    | Heptane, 2,3-dimethyl-         | 128 | C9H20   | 003074-71-3 | 37   |
| 5    |    | Heptane, 3,4,5-trimethyl-      | 142 | C10H22  | 020278-89-1 | 35   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000542.D  
Acq On : 22 Mar 2018 16:12  
Operator : JU/SJ  
Sample : J1905-11DL 10X  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampled :  
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Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

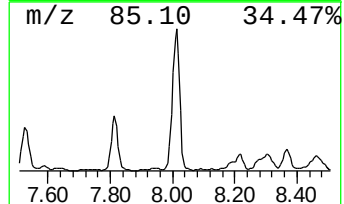
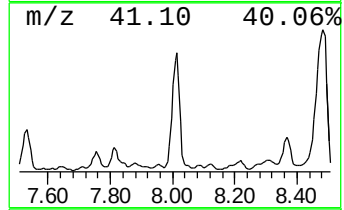
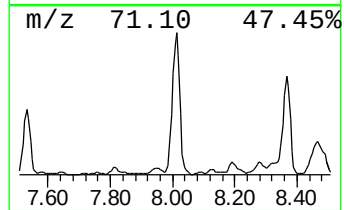
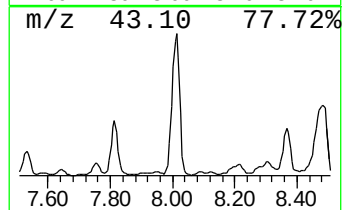
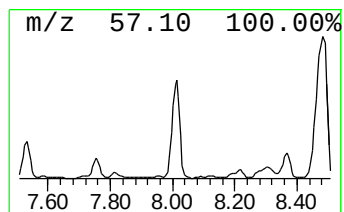
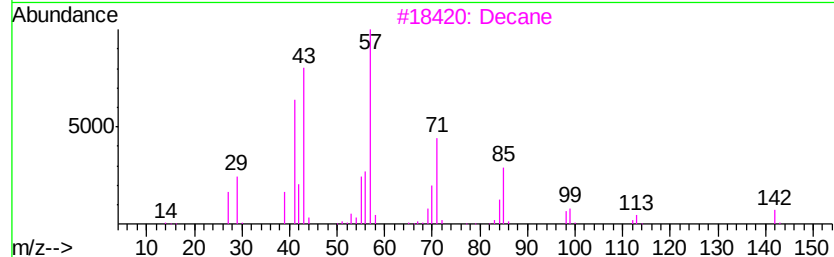
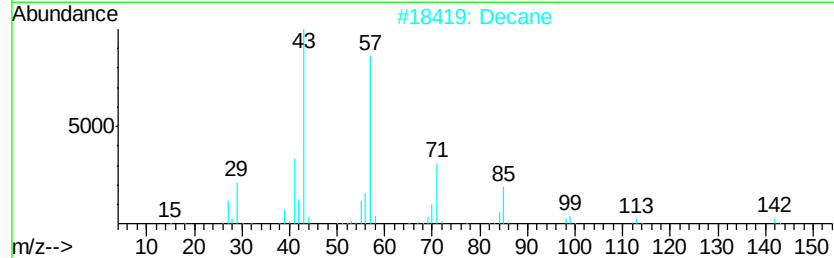
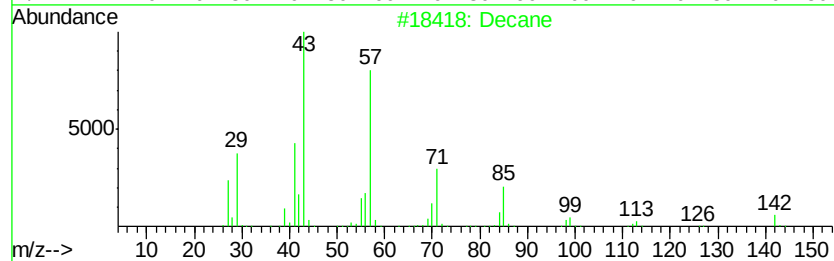
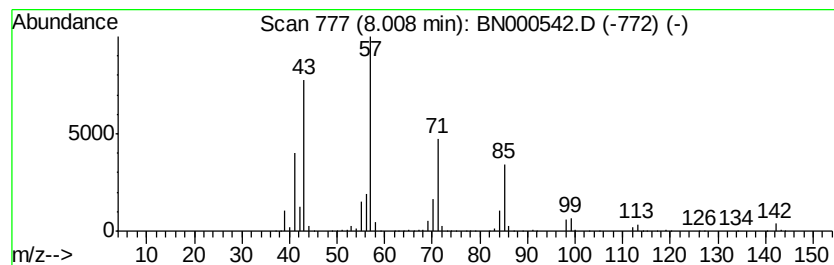
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TIC Integration Parameters: LSCINT.P

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Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 2

| R.T. | EstConc     | Area    | Relative to ISTD       | R.T. |
|------|-------------|---------|------------------------|------|
| 8.01 | 75.43 ng/ul | 9701630 | 1,4-Dichlorobenzene-d4 | 8.42 |

| Hit# | of | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|--------------|-----|---------|-------------|------|
| 1    | 5  | Decane       | 142 | C10H22  | 000124-18-5 | 95   |
| 2    |    | Decane       | 142 | C10H22  | 000124-18-5 | 94   |
| 3    |    | Decane       | 142 | C10H22  | 000124-18-5 | 91   |
| 4    |    | Tridecane    | 184 | C13H28  | 000629-50-5 | 83   |
| 5    |    | Tetradecane  | 198 | C14H30  | 000629-59-4 | 83   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000542.D  
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Operator : JU/SJ  
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Misc :  
ALS Vial : 12 Sample Multiplier: 1

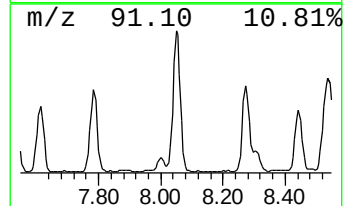
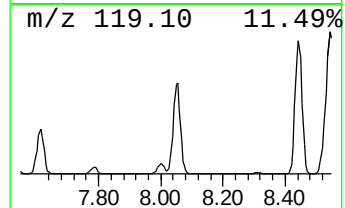
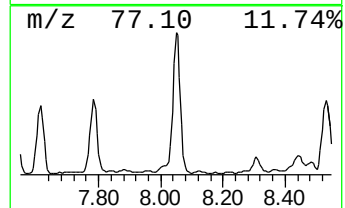
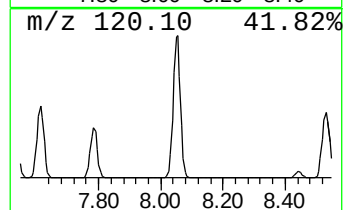
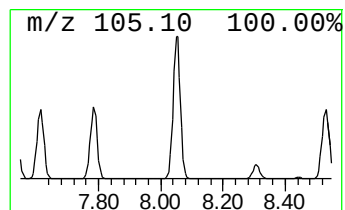
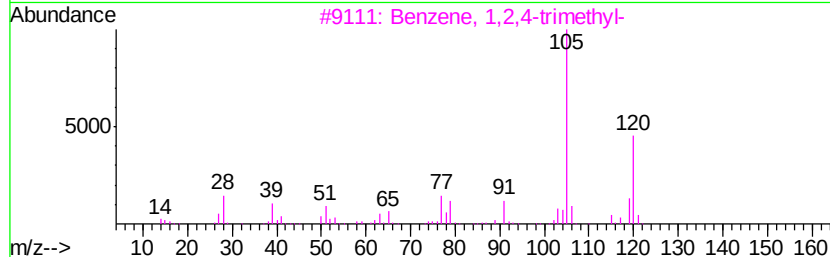
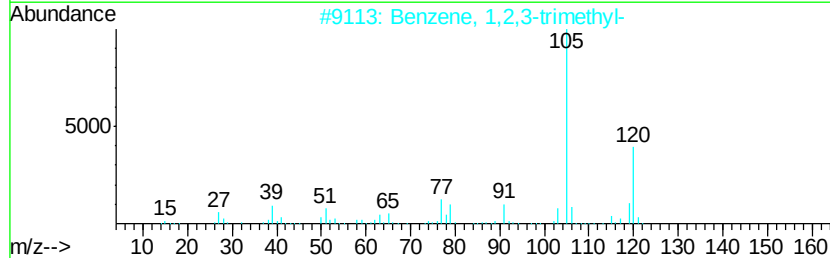
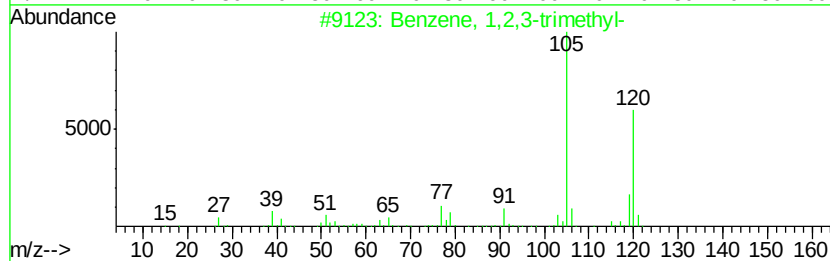
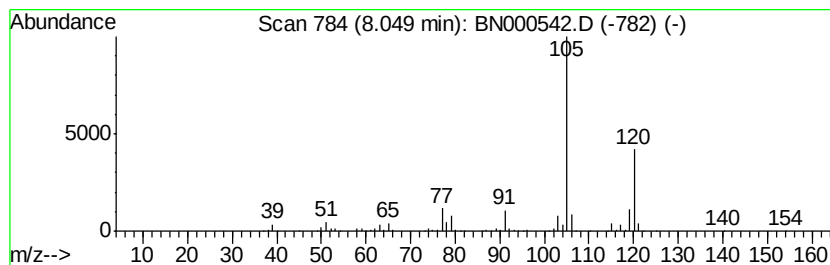
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ClientSampleId :  
DAM46DL

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 9 Benzene, 1,2,3-trimethyl- Concentration Rank 26

| R.T.    | EstConc     | Area                      | Relative to ISTD       | R.T.           |
|---------|-------------|---------------------------|------------------------|----------------|
| 8.05    | 22.27 ng/ul | 2863830                   | 1,4-Dichlorobenzene-d4 | 8.42           |
| Hit# of | 5           | Tentative ID              | MW MolForm             | CAS# Qual      |
| 1       |             | Benzene, 1,2,3-trimethyl- | 120 C9H12              | 000526-73-8 97 |
| 2       |             | Benzene, 1,2,3-trimethyl- | 120 C9H12              | 000526-73-8 95 |
| 3       |             | Benzene, 1,2,4-trimethyl- | 120 C9H12              | 000095-63-6 94 |
| 4       |             | Benzene, 1,2,4-trimethyl- | 120 C9H12              | 000095-63-6 91 |
| 5       |             | Benzene, 1,2,4-trimethyl- | 120 C9H12              | 000095-63-6 91 |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000542.D  
Acq On : 22 Mar 2018 16:12  
Operator : JU/SJ  
Sample : J1905-11DL 10X  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleID :  
DAM46DL

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

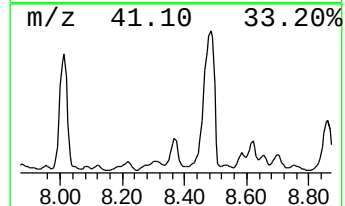
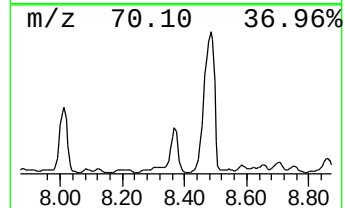
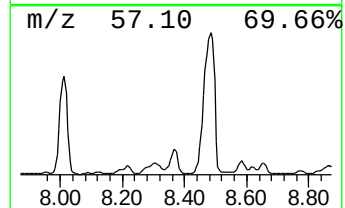
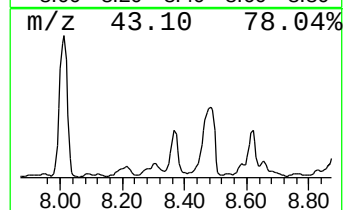
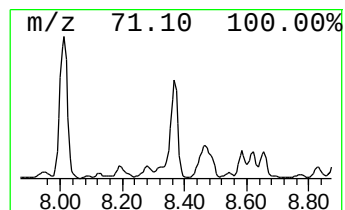
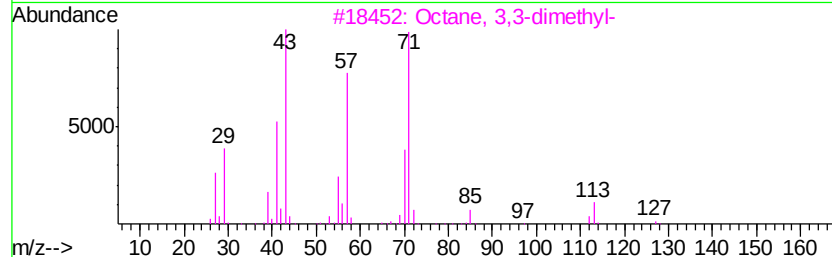
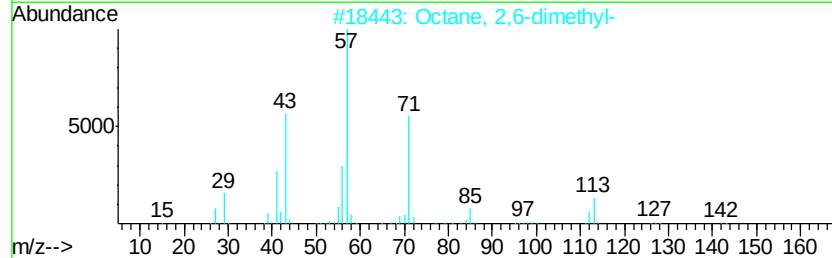
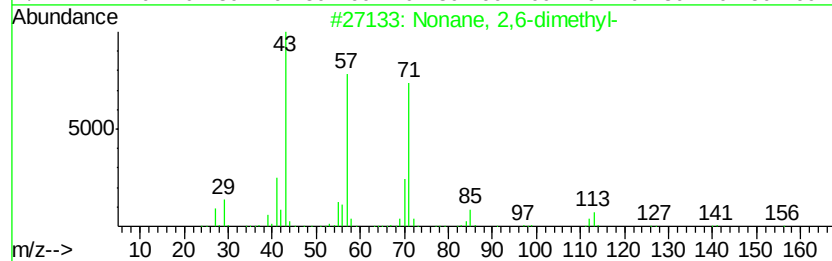
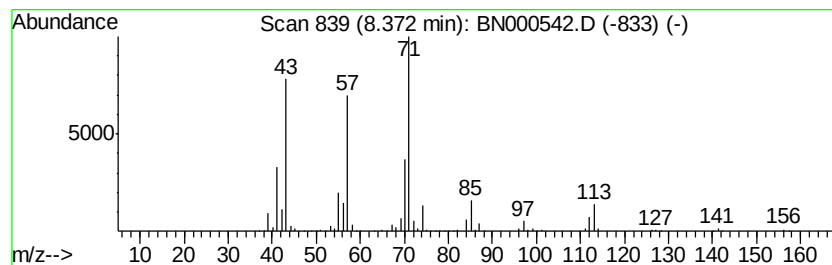
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 34

| R.T. | EstConc     | Area    | Relative to ISTD       | R.T. |
|------|-------------|---------|------------------------|------|
| 8.37 | 20.00 ng/ul | 2572200 | 1,4-Dichlorobenzene-d4 | 8.42 |

| Hit# | of | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------------|-----|---------|-------------|------|
| 1    | 5  | Nonane, 2,6-dimethyl-     | 156 | C11H24  | 017302-28-2 | 87   |
| 2    |    | Octane, 2,6-dimethyl-     | 142 | C10H22  | 002051-30-1 | 72   |
| 3    |    | Octane, 3,3-dimethyl-     | 142 | C10H22  | 004110-44-5 | 64   |
| 4    |    | Octane, 3,3-dimethyl-     | 142 | C10H22  | 004110-44-5 | 64   |
| 5    |    | Heptane, 3,3,5-trimethyl- | 142 | C10H22  | 007154-80-5 | 64   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000542.D  
Acq On : 22 Mar 2018 16:12  
Operator : JU/SJ  
Sample : J1905-11DL 10X  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM46DL

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

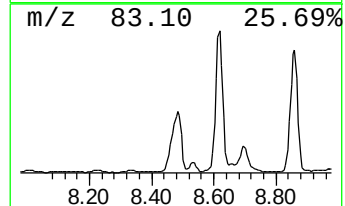
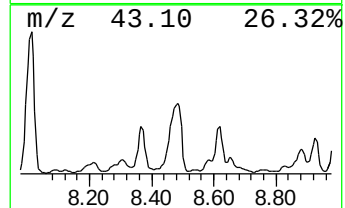
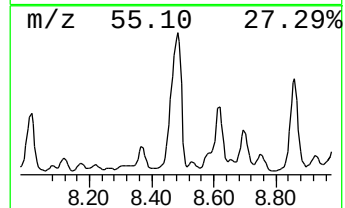
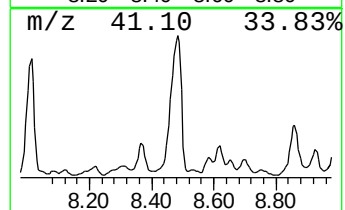
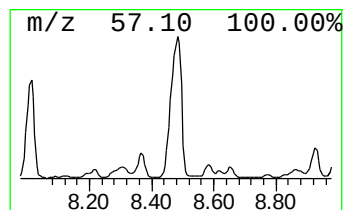
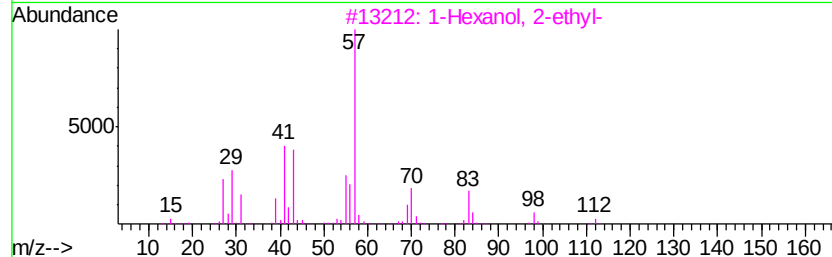
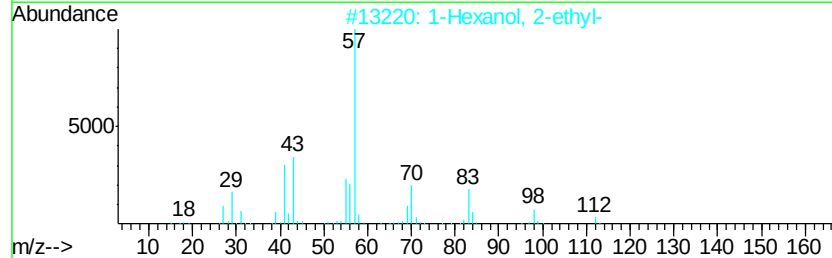
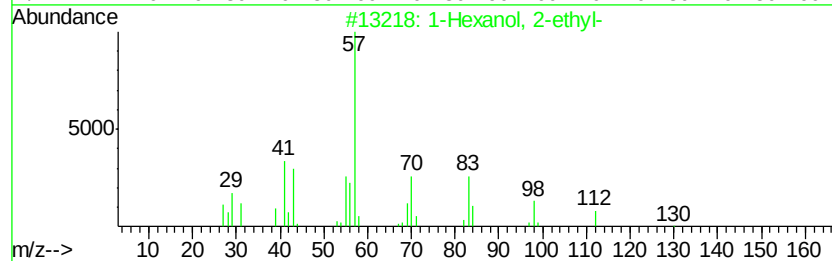
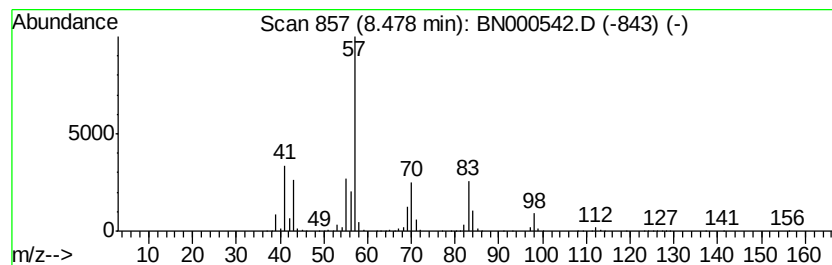
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 11 1-Hexanol, 2-ethyl- Concentration Rank 1

| R.T. | EstConc      | Area     | Relative to ISTD       | R.T. |
|------|--------------|----------|------------------------|------|
| 8.48 | 118.65 ng/ul | 15259700 | 1,4-Dichlorobenzene-d4 | 8.42 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm    | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|------------|-------------|------|
| 1    |    |   | 1-Hexanol, 2-ethyl-                 | 130 | C8H18O     | 000104-76-7 | 83   |
| 2    |    |   | 1-Hexanol, 2-ethyl-                 | 130 | C8H18O     | 000104-76-7 | 78   |
| 3    |    |   | 1-Hexanol, 2-ethyl-                 | 130 | C8H18O     | 000104-76-7 | 64   |
| 4    |    |   | 1-Hexanol, 2-ethyl-                 | 130 | C8H18O     | 000104-76-7 | 64   |
| 5    |    |   | Chloroacetic acid, 2-ethylhexyl ... | 206 | C10H19ClO2 | 005345-58-4 | 59   |





Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000542.D  
Acq On : 22 Mar 2018 16:12  
Operator : JU/SJ  
Sample : J1905-11DL 10X  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

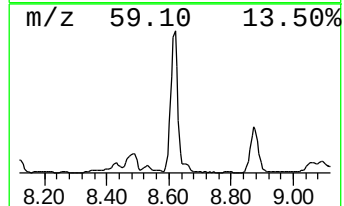
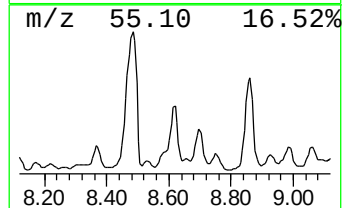
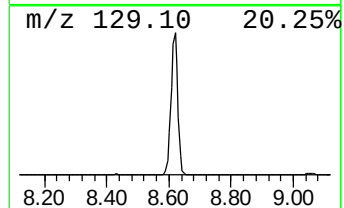
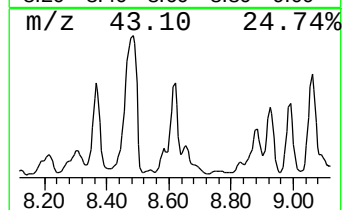
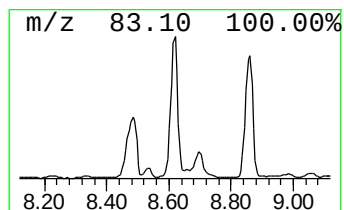
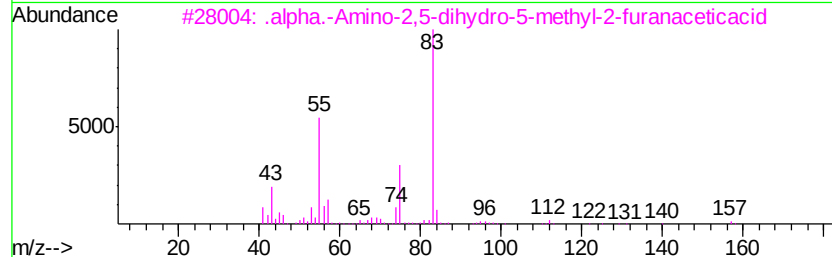
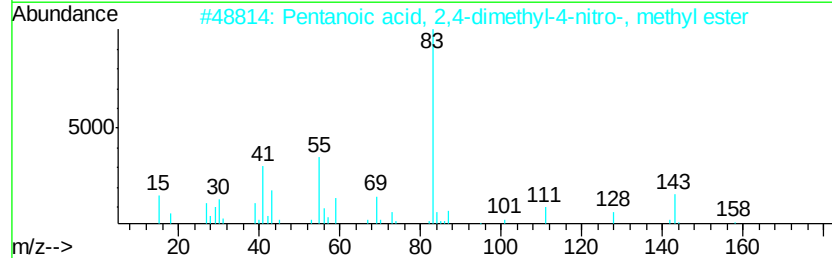
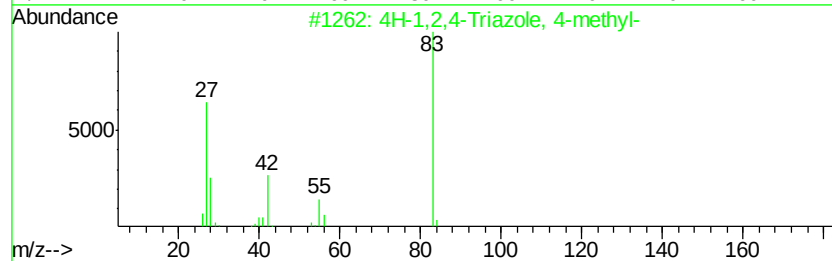
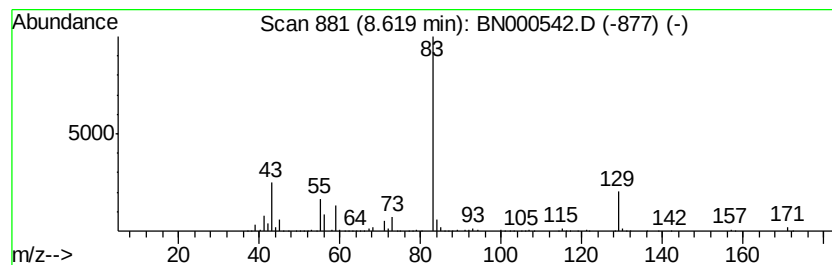
Instrument :  
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ClientSampleId :  
DAM46DL

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 12 unknown-01 Concentration Rank 19

| R.T.    | EstConc                              | Area          | Relative to ISTD       | R.T.      |
|---------|--------------------------------------|---------------|------------------------|-----------|
| 8.62    | 25.91 ng/ul                          | 3332290       | 1,4-Dichlorobenzene-d4 | 8.42      |
| Hit# of | 5                                    | Tentative ID  | MW MolForm             | CAS# Qual |
| 1       | 4H-1,2,4-Triazole, 4-methyl-         | 83 C3H5N3     | 010570-40-8            | 40        |
| 2       | Pentanoic acid, 2,4-dimethyl-4-n...  | 189 C8H15N04  | 005762-40-3            | 38        |
| 3       | .alpha.-Amino-2,5-dihydro-5-meth...  | 157 C7H11N03  | 018455-25-9            | 33        |
| 4       | 1H-Imidazol-2-amine                  | 83 C3H5N3     | 007720-39-0            | 33        |
| 5       | Bis(2-ethylhexyl) hydrogen phosph... | 306 C16H35O3P | 003658-48-8            | 33        |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000542.D  
Acq On : 22 Mar 2018 16:12  
Operator : JU/SJ  
Sample : J1905-11DL 10X  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleID :  
DAM46DL

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

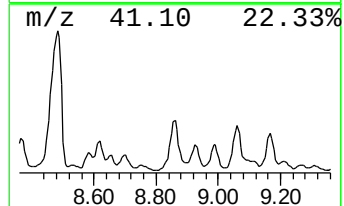
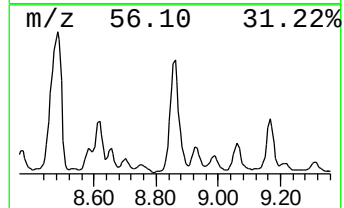
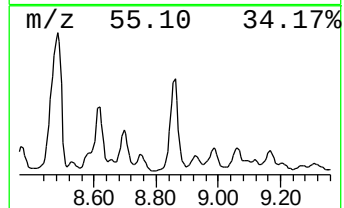
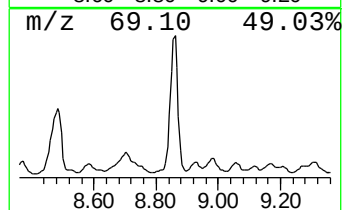
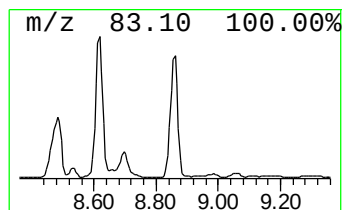
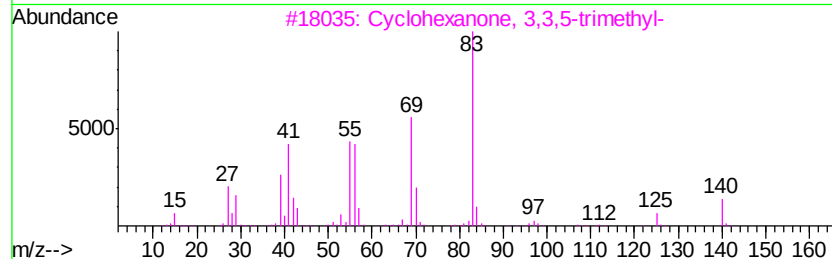
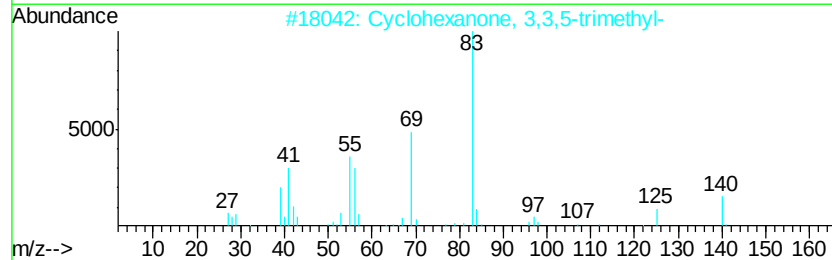
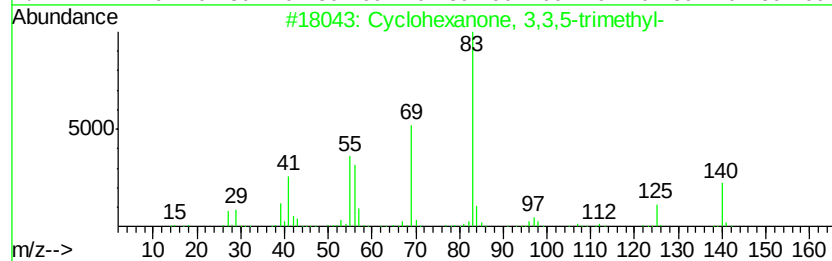
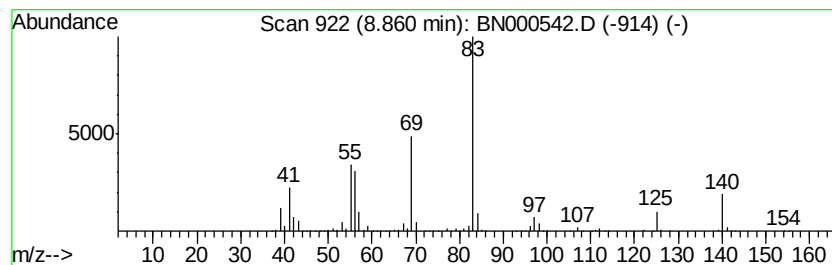
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 13 Cyclohexanone, 3,3,5-trimet... Concentration Rank 4

| R.T. | EstConc     | Area    | Relative to ISTD       | R.T. |
|------|-------------|---------|------------------------|------|
| 8.86 | 52.25 ng/ul | 6720170 | 1,4-Dichlorobenzene-d4 | 8.42 |

| Hit# | of | Tentative ID                    | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------------------|-----|---------|-------------|------|
| 1    | 5  | Cyclohexanone, 3,3,5-trimethyl- | 140 | C9H16O  | 000873-94-9 | 97   |
| 2    |    | Cyclohexanone, 3,3,5-trimethyl- | 140 | C9H16O  | 000873-94-9 | 95   |
| 3    |    | Cyclohexanone, 3,3,5-trimethyl- | 140 | C9H16O  | 000873-94-9 | 91   |
| 4    |    | Cyclohexanone, 3,3,5-trimethyl- | 140 | C9H16O  | 000873-94-9 | 56   |
| 5    |    | 1-Methyl-1H-1,2,4-triazole      | 83  | C3H5N3  | 006086-21-1 | 50   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000542.D  
Acq On : 22 Mar 2018 16:12  
Operator : JU/SJ  
Sample : J1905-11DL 10X  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM46DL

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

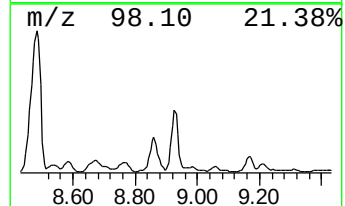
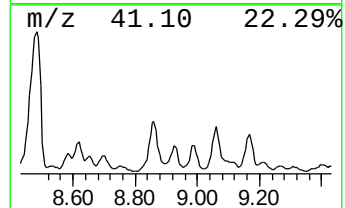
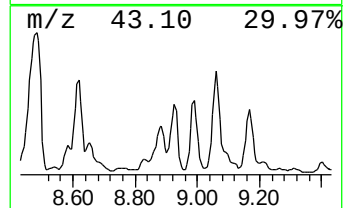
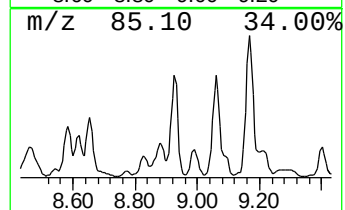
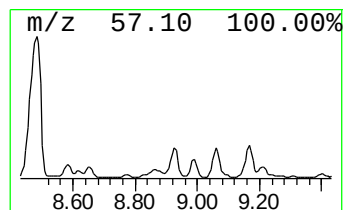
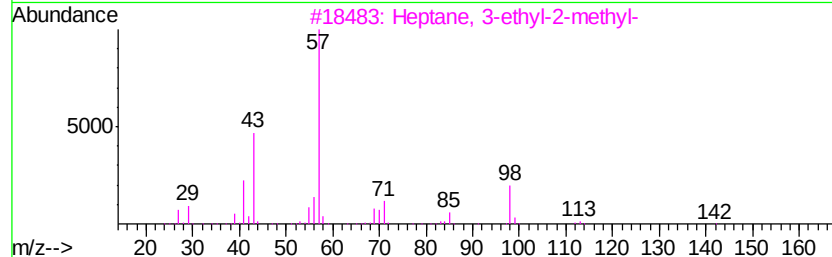
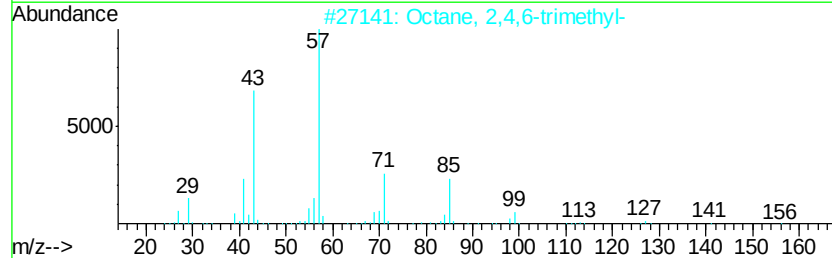
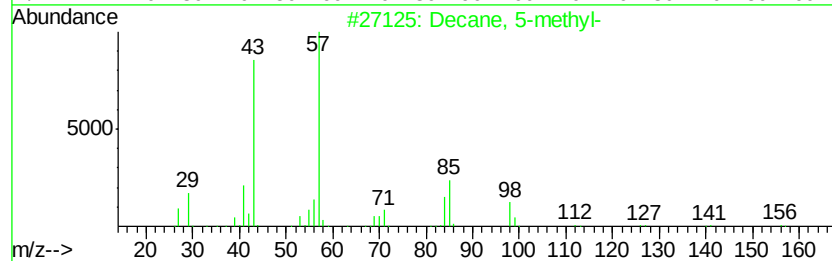
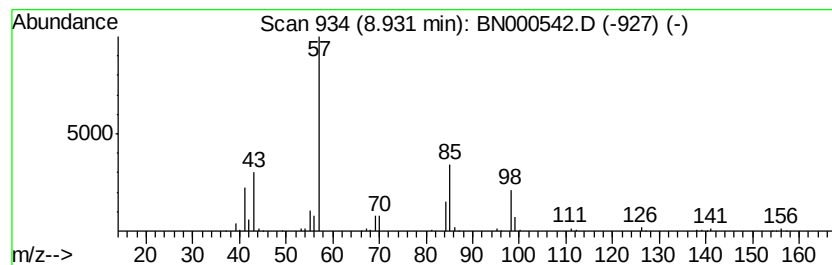
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 14 (DEL) Alkane: Straight-Chai... Concentration Rank 57

| R.T. | EstConc     | Area    | Relative to ISTD       | R.T. |
|------|-------------|---------|------------------------|------|
| 8.93 | 14.26 ng/ul | 1834010 | 1,4-Dichlorobenzene-d4 | 8.42 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Decane, 5-methyl-          | 156 | C11H24  | 013151-35-4 | 64   |
| 2    |    | Octane, 2,4,6-trimethyl-   | 156 | C11H24  | 062016-37-9 | 38   |
| 3    |    | Heptane, 3-ethyl-2-methyl- | 142 | C10H22  | 014676-29-0 | 38   |
| 4    |    | Heptacosane                | 380 | C27H56  | 000593-49-7 | 38   |
| 5    |    | Nonane, 2,5-dimethyl-      | 156 | C11H24  | 017302-27-1 | 38   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000542.D  
Acq On : 22 Mar 2018 16:12  
Operator : JU/SJ  
Sample : J1905-11DL 10X  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM46DL

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

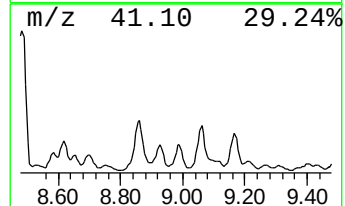
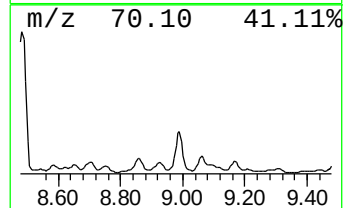
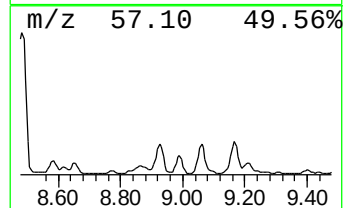
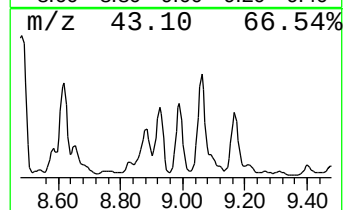
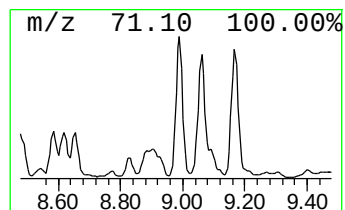
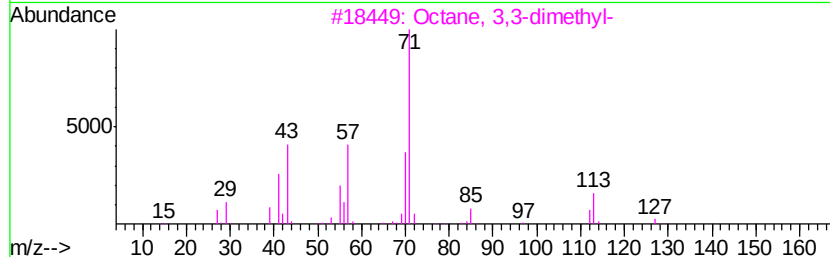
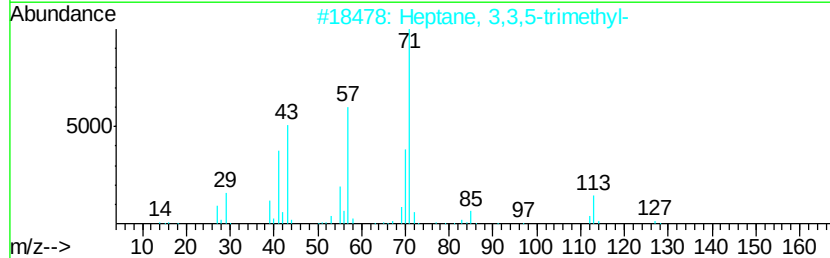
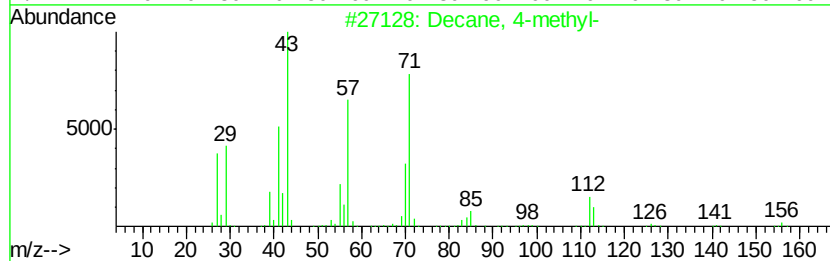
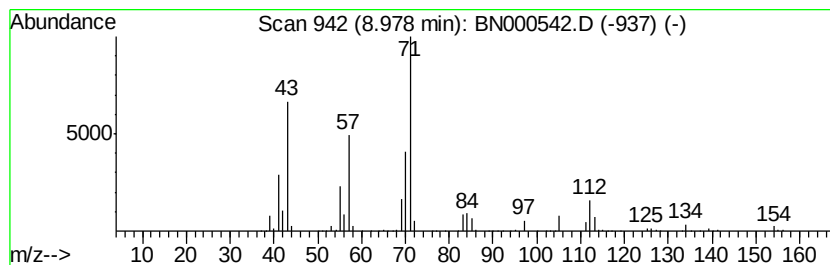
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TIC Integration Parameters: LSCINT.P

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Peak Number 15 (DEL) Alkane: Straight-Chai... Concentration Rank 14

| R.T. | EstConc     | Area    | Relative to ISTD       | R.T. |
|------|-------------|---------|------------------------|------|
| 8.98 | 28.66 ng/ul | 3685810 | 1,4-Dichlorobenzene-d4 | 8.42 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|-------------------------------------|-----|----------|--------------|------|
| 1    | 5  | Decane, 4-methyl-                   | 156 | C11H24   | 002847-72-5  | 83   |
| 2    |    | Heptane, 3,3,5-trimethyl-           | 142 | C10H22   | 007154-80-5  | 78   |
| 3    |    | Octane, 3,3-dimethyl-               | 142 | C10H22   | 004110-44-5  | 72   |
| 4    |    | 1-Hexene, 3,4,5-trimethyl-          | 126 | C9H18    | 056728-10-0  | 64   |
| 5    |    | 1,1-Dimethylpropyl 2-ethylhexanoate | 214 | C13H26O2 | 1000099-99-2 | 64   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000542.D  
Acq On : 22 Mar 2018 16:12  
Operator : JU/SJ  
Sample : J1905-11DL 10X  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleID :  
DAM46DL

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

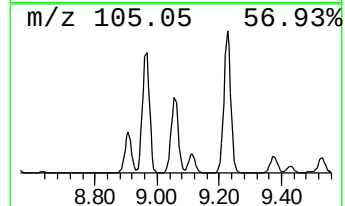
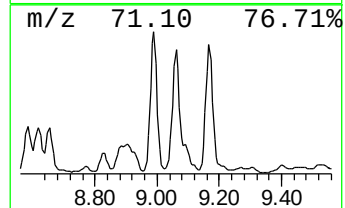
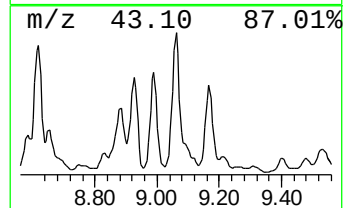
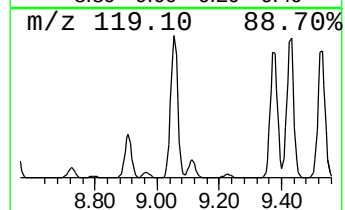
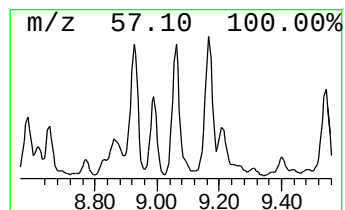
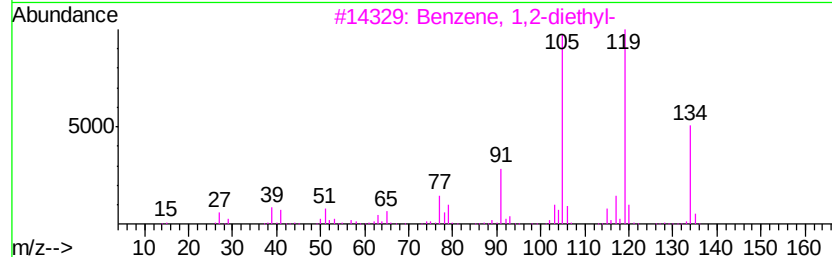
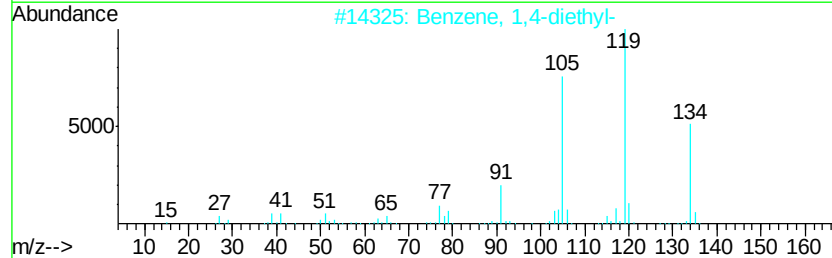
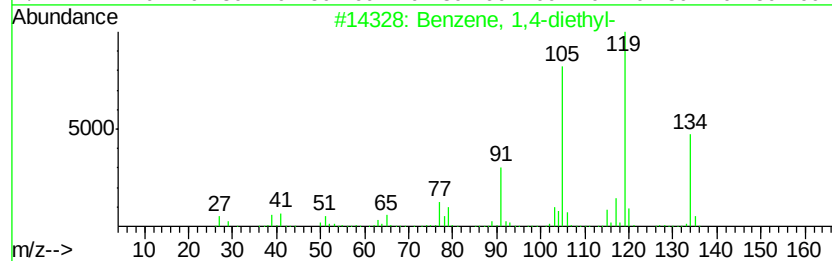
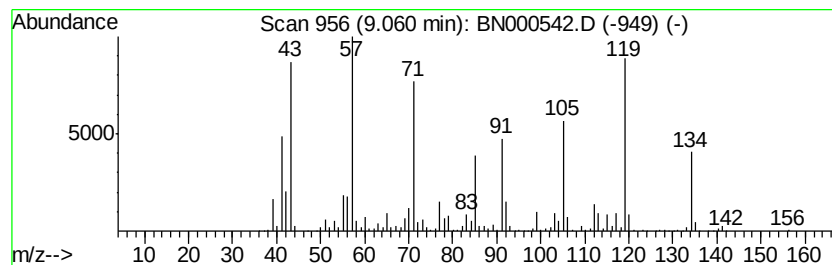
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TIC Integration Parameters: LSCINT.P

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Peak Number 16 Benzene, 1,4-diethyl- Concentration Rank 3

| R.T. | EstConc     | Area    | Relative to ISTD       | R.T. |
|------|-------------|---------|------------------------|------|
| 9.06 | 54.94 ng/ul | 7066310 | 1,4-Dichlorobenzene-d4 | 8.42 |

| Hit# | of | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1,4-diethyl-          | 134 | C10H14  | 000105-05-5 | 94   |
| 2    |    | Benzene, 1,4-diethyl-          | 134 | C10H14  | 000105-05-5 | 83   |
| 3    |    | Benzene, 1,2-diethyl-          | 134 | C10H14  | 000135-01-3 | 60   |
| 4    |    | Benzene, 1-ethyl-2,3-dimethyl- | 134 | C10H14  | 000933-98-2 | 52   |
| 5    |    | Benzene, 1,3-diethyl-          | 134 | C10H14  | 000141-93-5 | 52   |



Data Path : Z:\HPCHEM1\BNA\_N\DATA\BN032218\  
Data File : BN000542.D  
Acq On : 22 Mar 2018 16:12  
Operator : JU/SJ  
Sample : J1905-11DL 10X  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM46DL

Quant Method : Z:\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

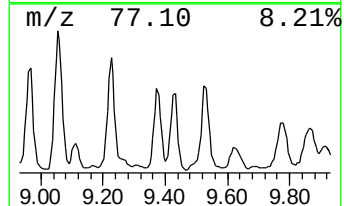
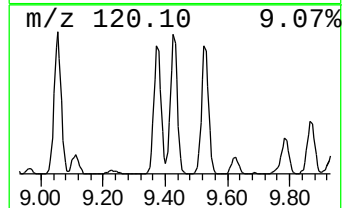
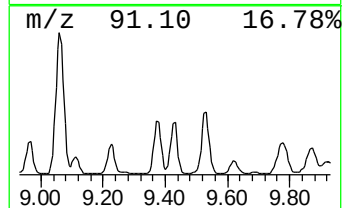
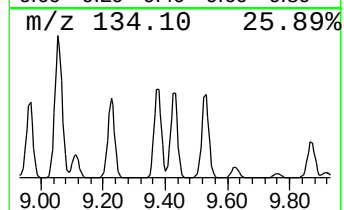
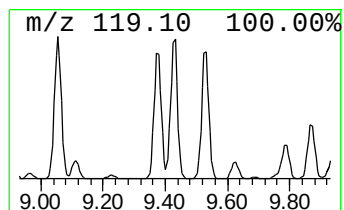
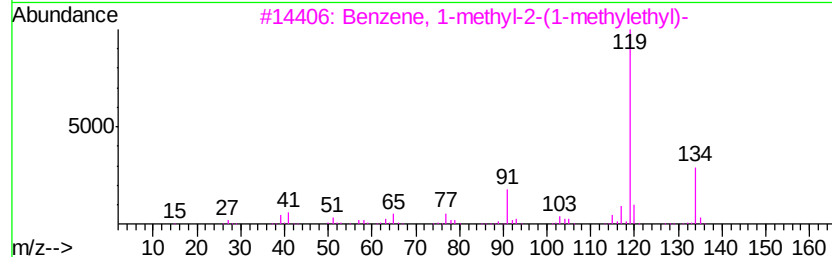
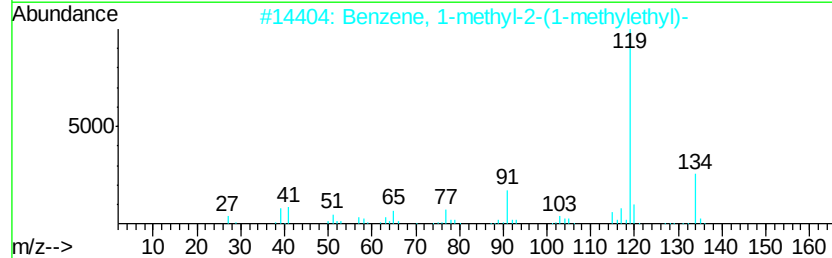
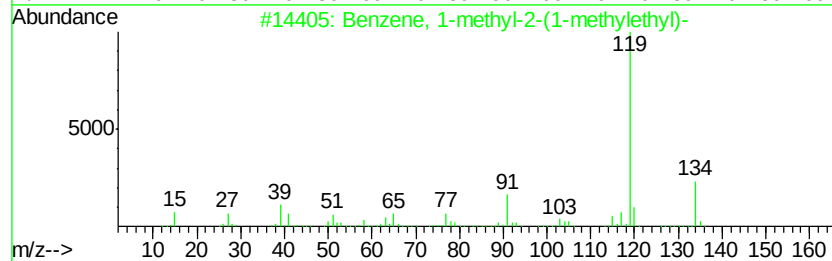
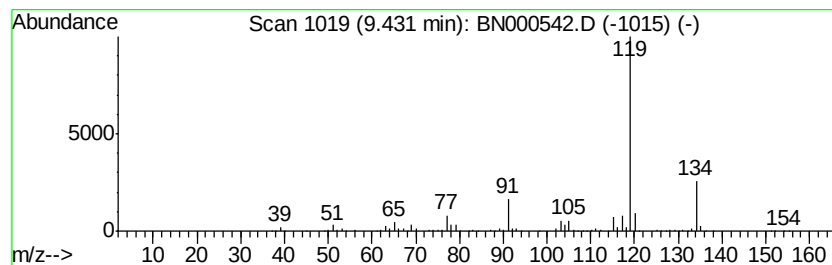
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TIC Integration Parameters: LSCINT.P

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Peak Number 17 Benzene, 1-methyl-2-(1-meth... Concentration Rank 54

| R.T. | EstConc     | Area    | Relative to ISTD       | R.T. |
|------|-------------|---------|------------------------|------|
| 9.43 | 14.54 ng/ul | 1870400 | 1,4-Dichlorobenzene-d4 | 8.42 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1-methyl-2-(1-methyleth... | 134 | C10H14  | 000527-84-4 | 97   |
| 2    |    | Benzene, 1-methyl-2-(1-methyleth... | 134 | C10H14  | 000527-84-4 | 97   |
| 3    |    | Benzene, 1-methyl-2-(1-methyleth... | 134 | C10H14  | 000527-84-4 | 97   |
| 4    |    | Benzene, 2-ethyl-1,3-dimethyl-      | 134 | C10H14  | 002870-04-4 | 95   |
| 5    |    | Benzene, 1-methyl-3-(1-methyleth... | 134 | C10H14  | 000535-77-3 | 95   |



Data Path : Z:\HPCHEM1\BNA\_N\DATA\BN032218\  
Data File : BN000542.D  
Acq On : 22 Mar 2018 16:12  
Operator : JU/SJ  
Sample : J1905-11DL 10X  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampled :  
DAM46DL

Quant Method : Z:\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

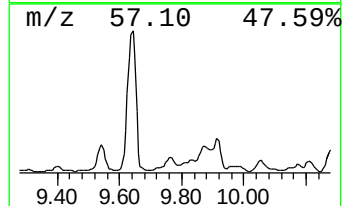
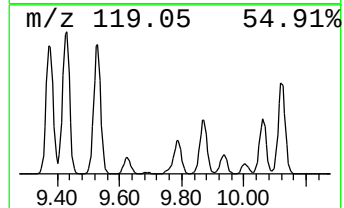
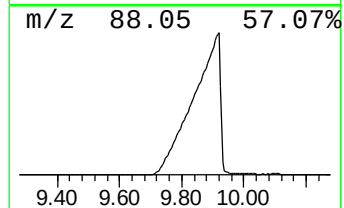
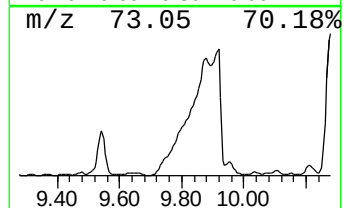
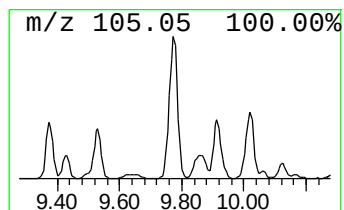
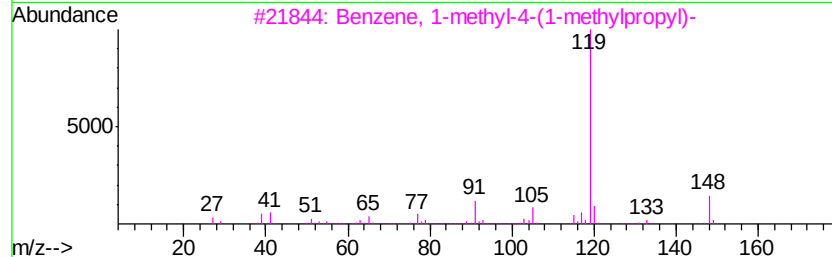
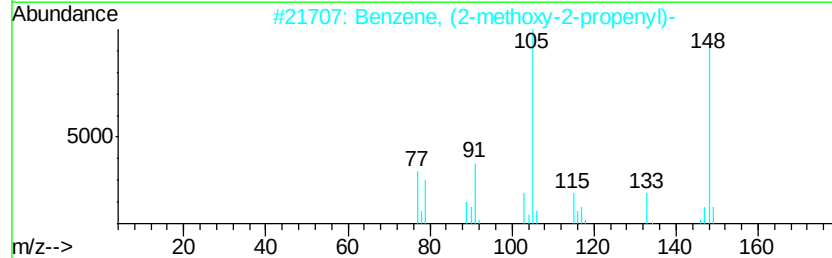
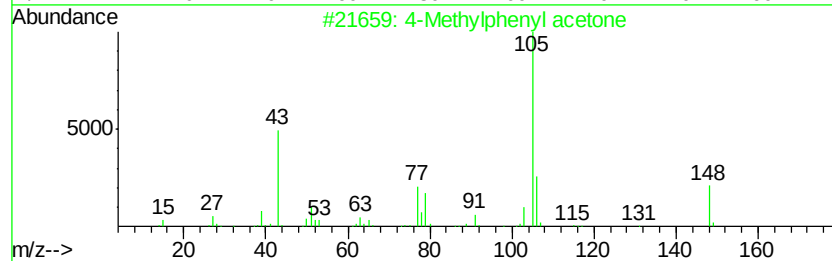
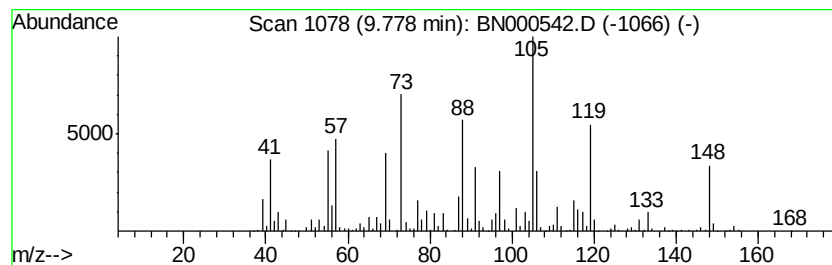
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 19 unknown-02 Concentration Rank 7

| R.T. | EstConc     | Area    | Relative to ISTD       | R.T. |
|------|-------------|---------|------------------------|------|
| 9.78 | 41.75 ng/ul | 5369920 | 1,4-Dichlorobenzene-d4 | 8.42 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | 4-Methylphenyl acetone              | 148 | C10H12O | 002096-86-8 | 35   |
| 2    |    | Benzene, (2-methoxy-2-propenyl)-    | 148 | C10H12O | 026473-60-9 | 30   |
| 3    |    | Benzene, 1-methyl-4-(1-methylpro... | 148 | C11H16  | 001595-16-0 | 27   |
| 4    |    | 2-Ethyl-2-methyl-1,3-dithiolane     | 148 | C6H12S2 | 006008-81-7 | 22   |
| 5    |    | 1,4-Cyclohexadiene, 3-ethenyl-1,... | 134 | C10H14  | 062338-57-2 | 22   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000542.D  
Acq On : 22 Mar 2018 16:12  
Operator : JU/SJ  
Sample : J1905-11DL 10X  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM46DL

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

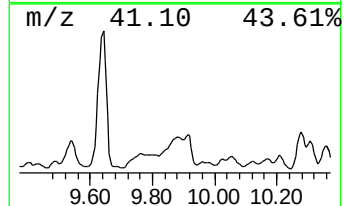
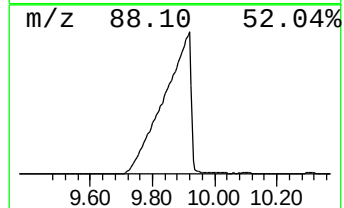
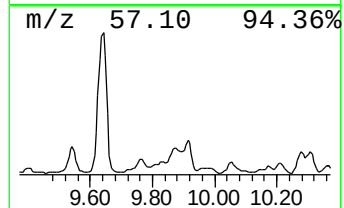
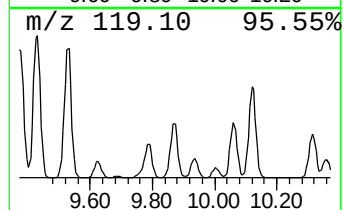
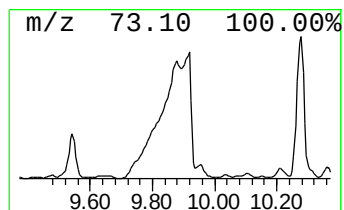
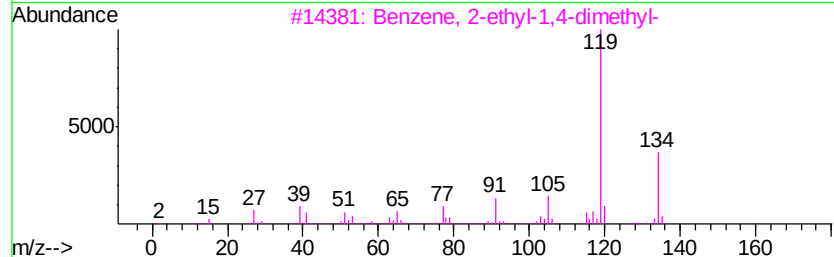
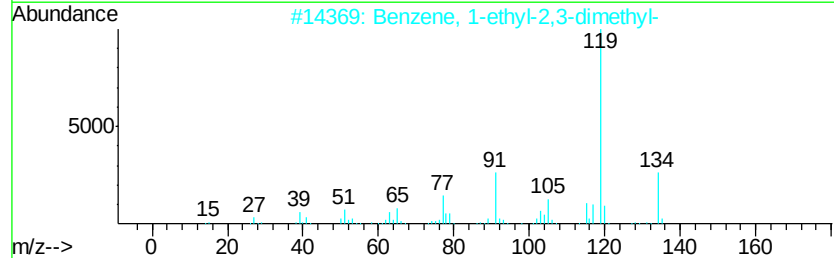
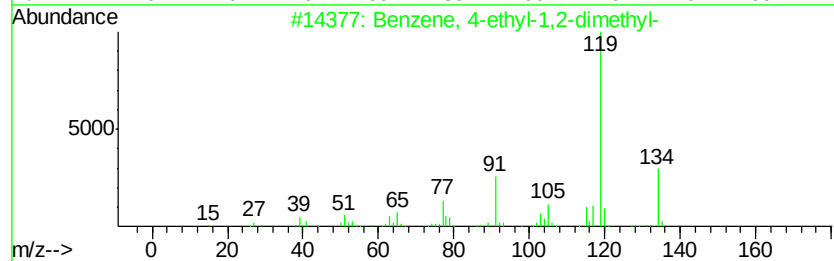
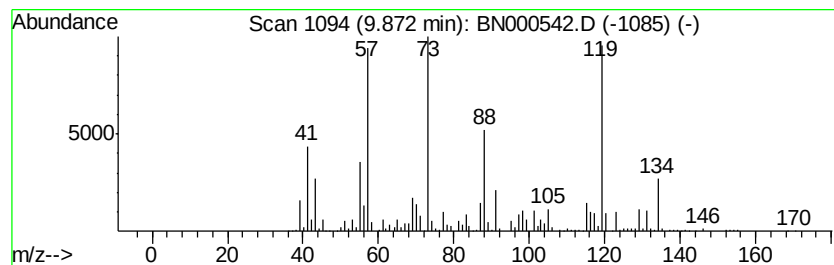
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 20 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 32

| R.T. | EstConc     | Area    | Relative to ISTD | R.T.  |
|------|-------------|---------|------------------|-------|
| 9.87 | 20.43 ng/ul | 4425030 | Napthalene-d8    | 11.26 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 4-ethyl-1,2-dimethyl-      | 134 | C10H14  | 000934-80-5 | 70   |
| 2    |    | Benzene, 1-ethyl-2,3-dimethyl-      | 134 | C10H14  | 000933-98-2 | 70   |
| 3    |    | Benzene, 2-ethyl-1,4-dimethyl-      | 134 | C10H14  | 001758-88-9 | 55   |
| 4    |    | Benzene, 1-methyl-2-(1-methyleth... | 134 | C10H14  | 000527-84-4 | 55   |
| 5    |    | Benzene, 2-ethyl-1,4-dimethyl-      | 134 | C10H14  | 001758-88-9 | 55   |





Data Path : Z:\HPCHEM1\BNA\_N\DATA\BN032218\  
Data File : BN000542.D  
Acq On : 22 Mar 2018 16:12  
Operator : JU/SJ  
Sample : J1905-11DL 10X  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM46DL

Quant Method : Z:\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

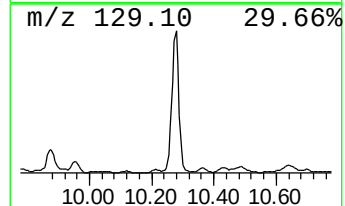
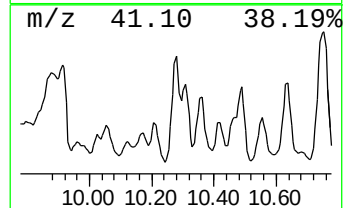
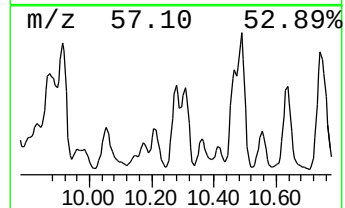
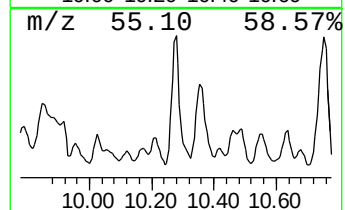
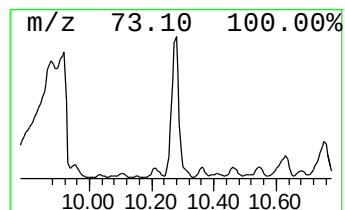
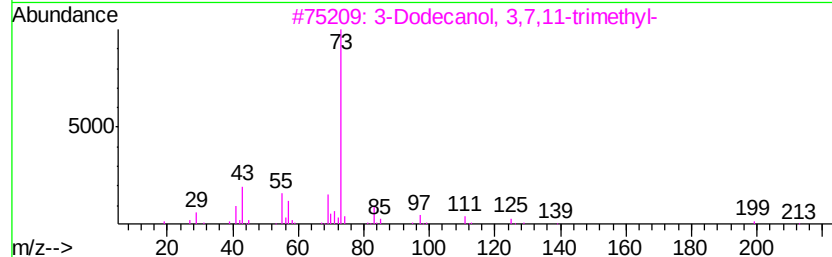
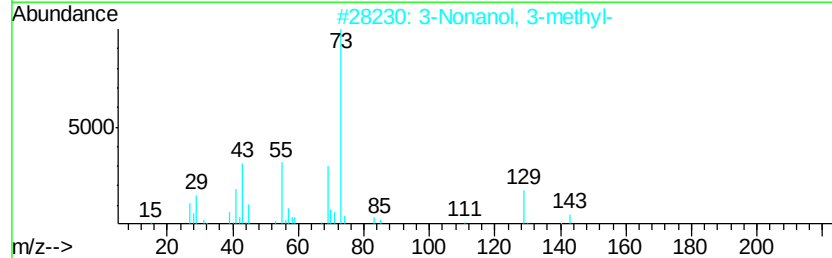
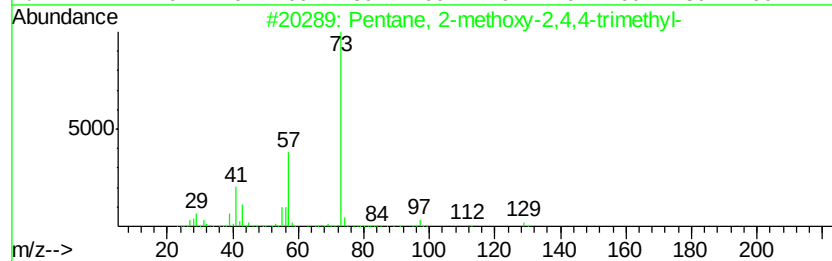
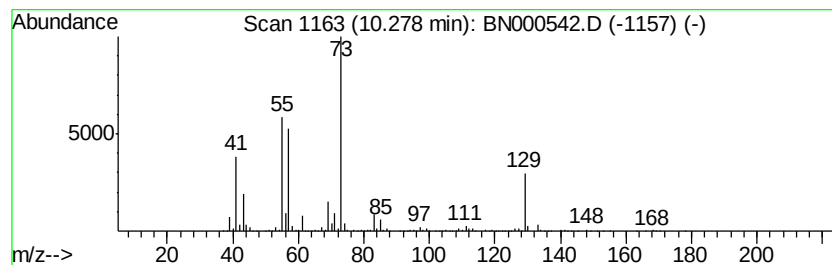
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 21 unknown-03 Concentration Rank 58

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 10.28 | 14.10 ng/ul | 3054240 | Napthalene-d8    | 11.26 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Pentane, 2-methoxy-2,4,4-trimethyl- | 144 | C9H20O  | 062108-41-2 | 32   |
| 2    |    | 3-Nonanol, 3-methyl-                | 158 | C10H22O | 021078-72-8 | 32   |
| 3    |    | 3-Dodecanol, 3,7,11-trimethyl-      | 228 | C15H32O | 007278-65-1 | 16   |
| 4    |    | 1-Heptene, 3-methyl-                | 112 | C8H16   | 004810-09-7 | 14   |
| 5    |    | Cyclopropane, pentamethyl-          | 112 | C8H16   | 014172-83-9 | 12   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000542.D  
Acq On : 22 Mar 2018 16:12  
Operator : JU/SJ  
Sample : J1905-11DL 10X  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

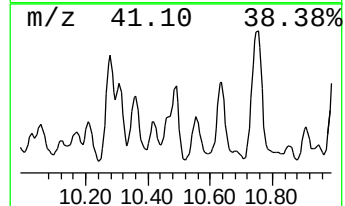
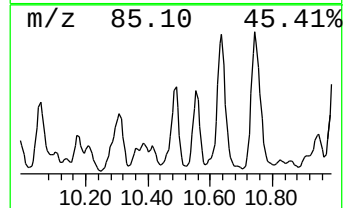
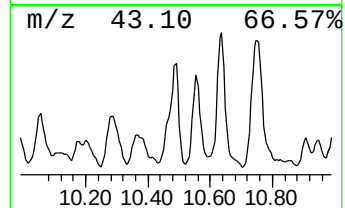
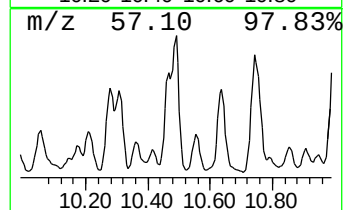
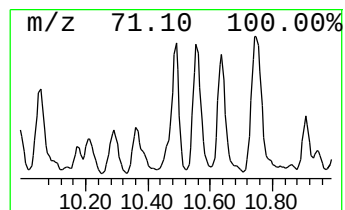
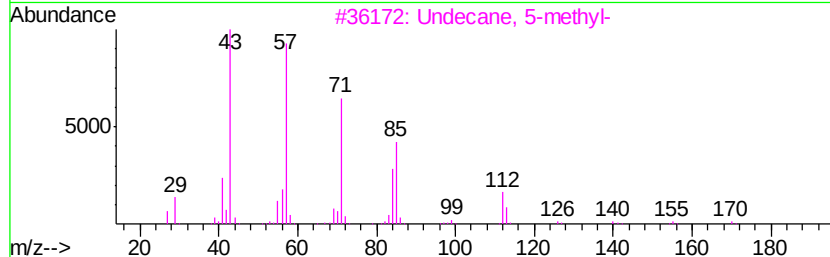
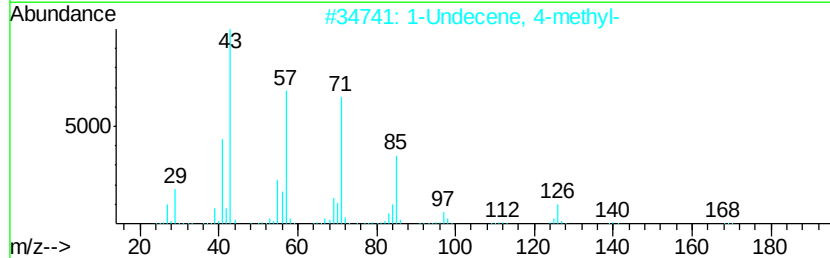
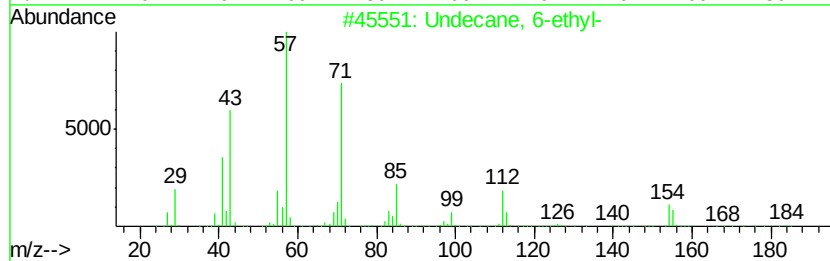
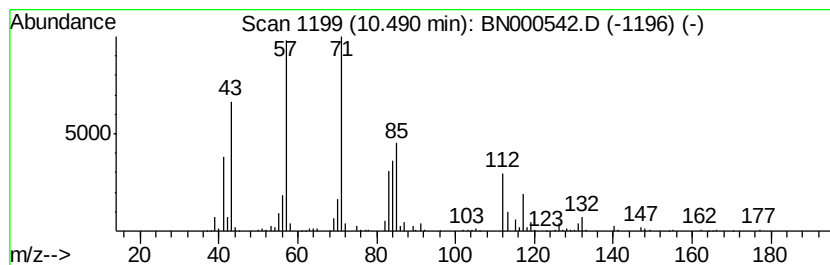
Instrument :  
BNA\_N  
ClientSampleId :  
DAM46DL

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 22 (DEL) Alkane: Straight-Chai... Concentration Rank 73

| R.T.      | EstConc                     | Area    | Relative to ISTD |             | R.T.  |
|-----------|-----------------------------|---------|------------------|-------------|-------|
| 10.49     | 9.58 ng/ul                  | 2074480 | Naphthalene-d8   |             | 11.26 |
| Hit# of 5 | Tentative ID                | MW      | MolForm          | CAS#        | Qual  |
| 1         | Undecane, 6-ethyl-          | 184     | C13H28           | 017312-60-6 | 50    |
| 2         | 1-Undecene, 4-methyl-       | 168     | C12H24           | 074630-39-0 | 50    |
| 3         | Undecane, 5-methyl-         | 170     | C12H26           | 001632-70-8 | 47    |
| 4         | Nonane, 5-(2-methylpropyl)- | 184     | C13H28           | 062185-53-9 | 47    |
| 5         | Dodecane                    | 170     | C12H26           | 000112-40-3 | 43    |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000542.D  
Acq On : 22 Mar 2018 16:12  
Operator : JU/SJ  
Sample : J1905-11DL 10X  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM46DL

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

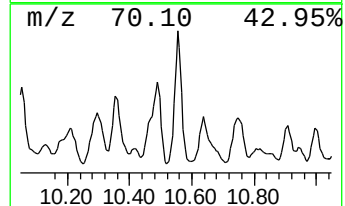
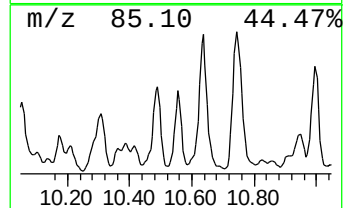
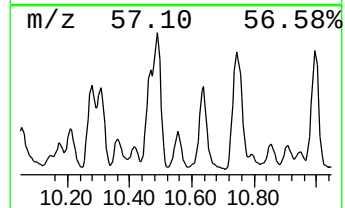
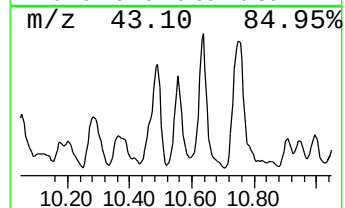
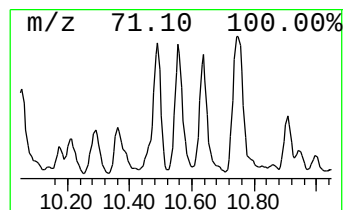
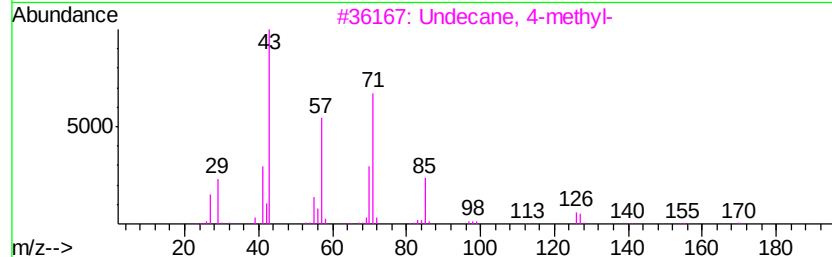
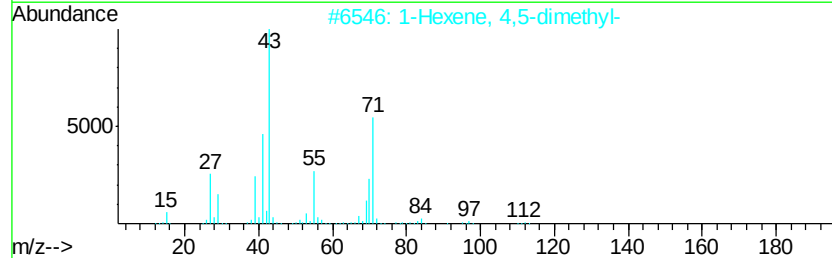
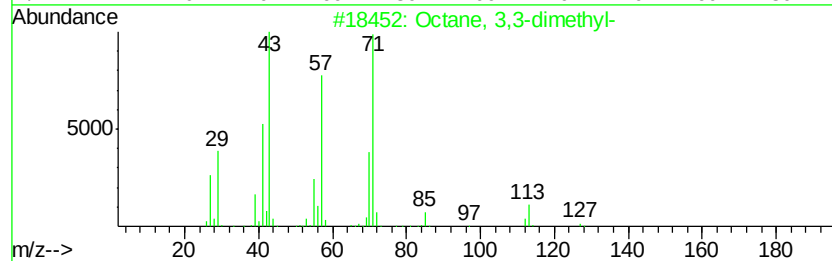
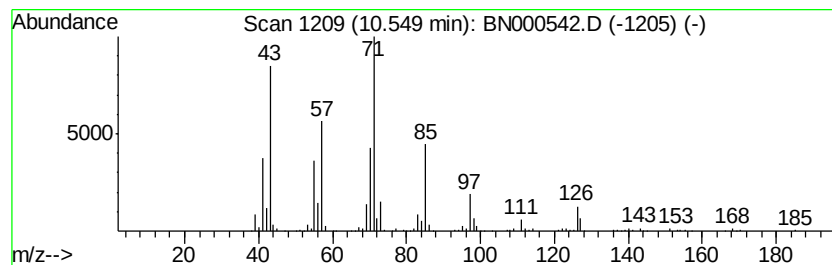
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TIC Integration Parameters: LSCINT.P

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Peak Number 23 (DEL) Alkane: Straight-Chai... Concentration Rank 70

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T.  |
|-------|------------|---------|------------------|-------|
| 10.55 | 9.97 ng/ul | 2159340 | Napthalene-d8    | 11.26 |

| Hit# | of | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------|-----|---------|-------------|------|
| 1    | 5  | Octane, 3,3-dimethyl-   | 142 | C10H22  | 004110-44-5 | 47   |
| 2    |    | 1-Hexene, 4,5-dimethyl- | 112 | C8H16   | 016106-59-5 | 47   |
| 3    |    | Undecane, 4-methyl-     | 170 | C12H26  | 002980-69-0 | 43   |
| 4    |    | Nonane, 5-butyl-        | 184 | C13H28  | 017312-63-9 | 43   |
| 5    |    | Tridecane, 6-methyl-    | 198 | C14H30  | 013287-21-3 | 38   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000542.D  
Acq On : 22 Mar 2018 16:12  
Operator : JU/SJ  
Sample : J1905-11DL 10X  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleID :  
DAM46DL

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

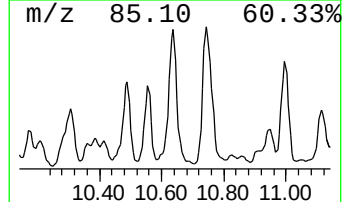
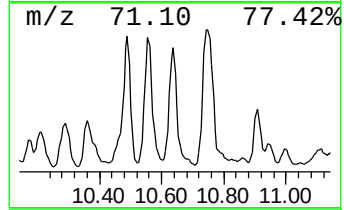
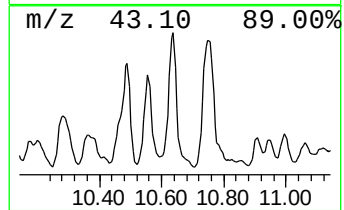
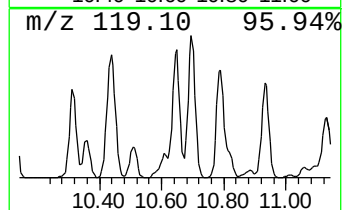
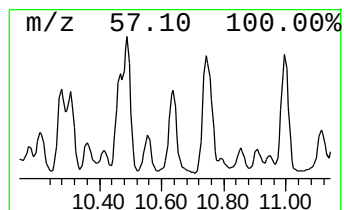
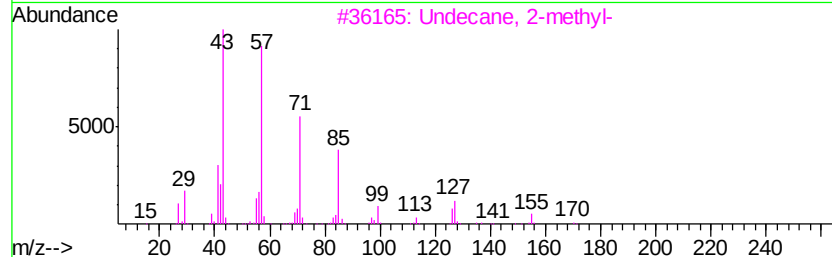
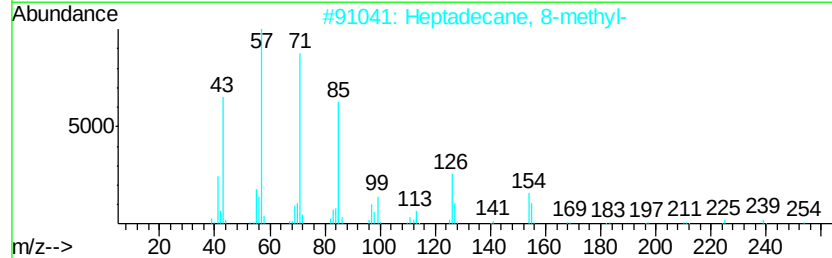
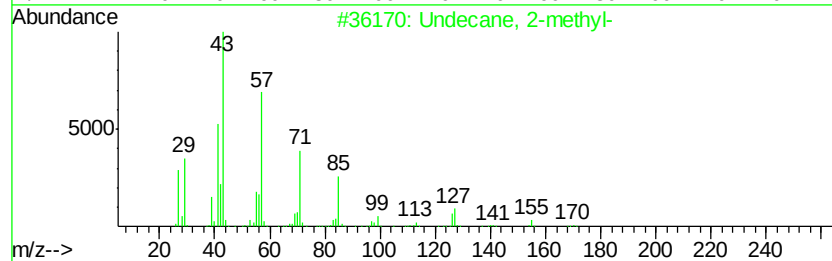
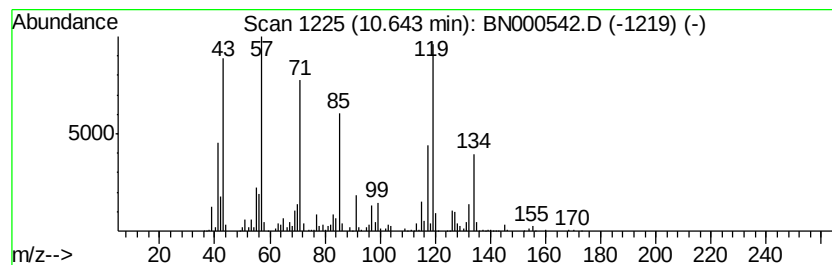
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 24 (DEL) Alkane: Straight-Chai... Concentration Rank 53

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 10.64 | 14.85 ng/ul | 3217240 | Napthalene-d8    | 11.26 |

| Hit# | of | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------|-----|---------|-------------|------|
| 1    | 5  | Undecane, 2-methyl-    | 170 | C12H26  | 007045-71-8 | 47   |
| 2    |    | Heptadecane, 8-methyl- | 254 | C18H38  | 013287-23-5 | 47   |
| 3    |    | Undecane, 2-methyl-    | 170 | C12H26  | 007045-71-8 | 47   |
| 4    |    | Tetracosane            | 338 | C24H50  | 000646-31-1 | 47   |
| 5    |    | 1-Undecene, 4-methyl-  | 168 | C12H24  | 074630-39-0 | 46   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000542.D  
Acq On : 22 Mar 2018 16:12  
Operator : JU/SJ  
Sample : J1905-11DL 10X  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM46DL

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

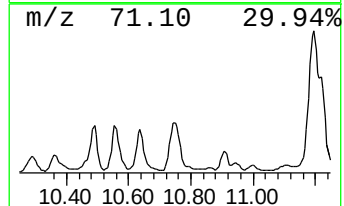
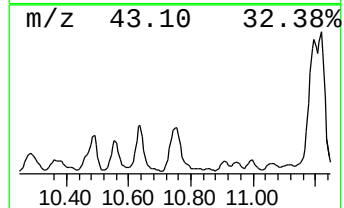
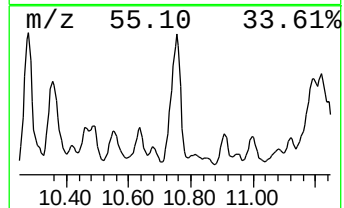
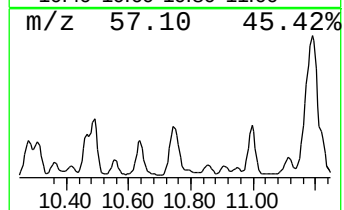
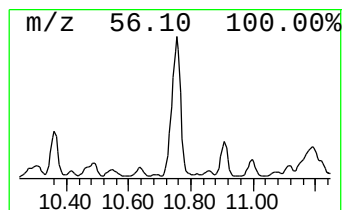
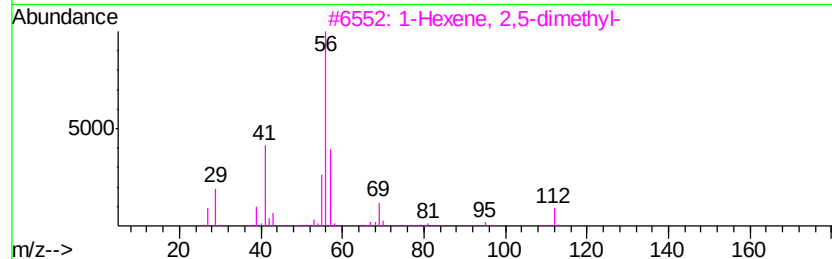
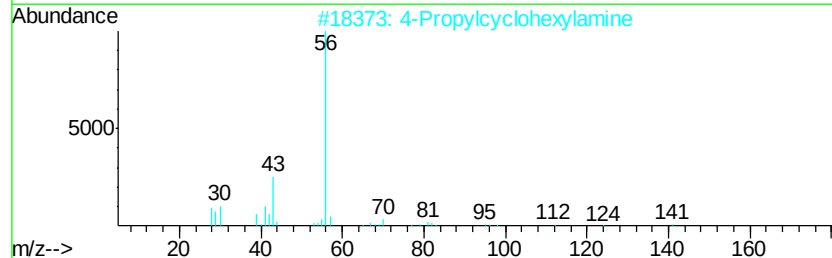
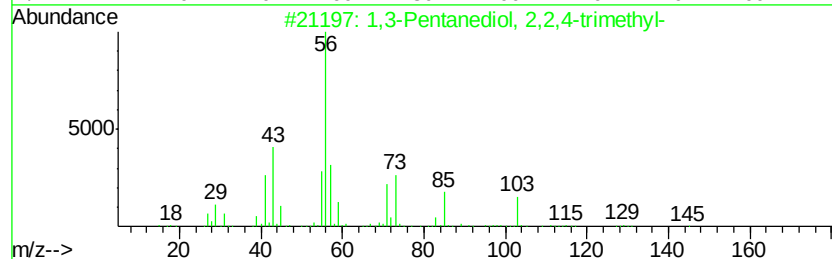
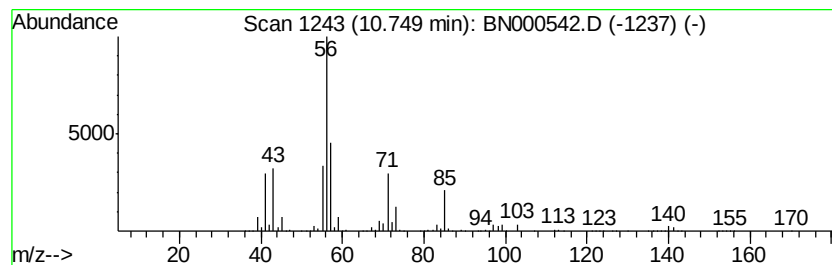
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 25 unknown-04 Concentration Rank 27

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 10.75 | 22.25 ng/ul | 4817900 | Napthalene-d8    | 11.26 |

| Hit# | of | Tentative ID                      | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------------------|-----|---------|-------------|------|
| 1    | 5  | 1,3-Pentenediol, 2,2,4-trimethyl- | 146 | C8H18O2 | 000144-19-4 | 45   |
| 2    |    | 4-Propylcyclohexylamine           | 141 | C9H19N  | 102653-37-2 | 43   |
| 3    |    | 1-Hexene, 2,5-dimethyl-           | 112 | C8H16   | 006975-92-4 | 37   |
| 4    |    | Hexane, 3,4-dimethyl-             | 114 | C8H18   | 000583-48-2 | 33   |
| 5    |    | 1,3-Hexanediol, 2-ethyl-          | 146 | C8H18O2 | 000094-96-2 | 33   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000542.D  
Acq On : 22 Mar 2018 16:12  
Operator : JU/SJ  
Sample : J1905-11DL 10X  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

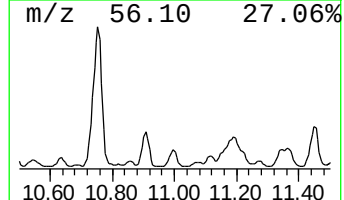
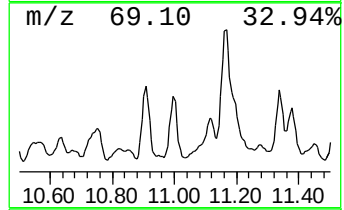
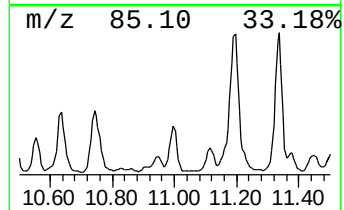
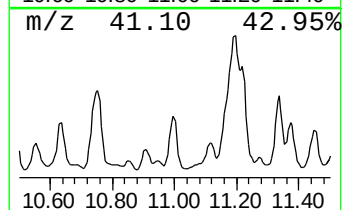
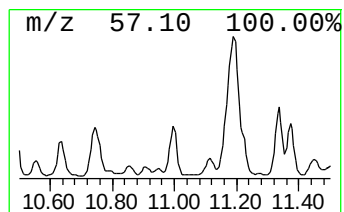
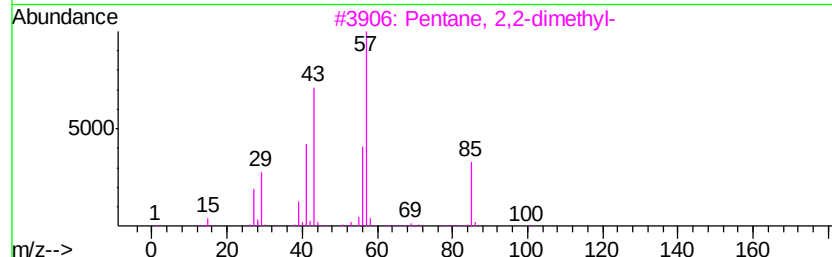
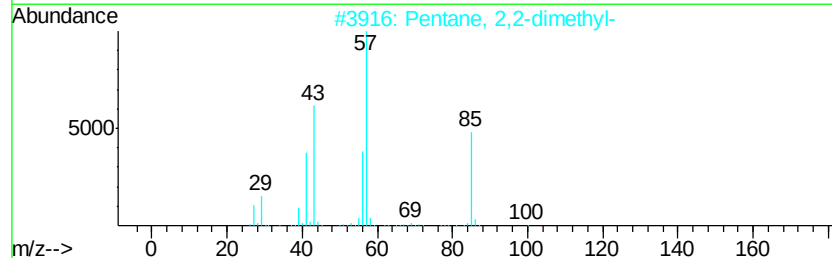
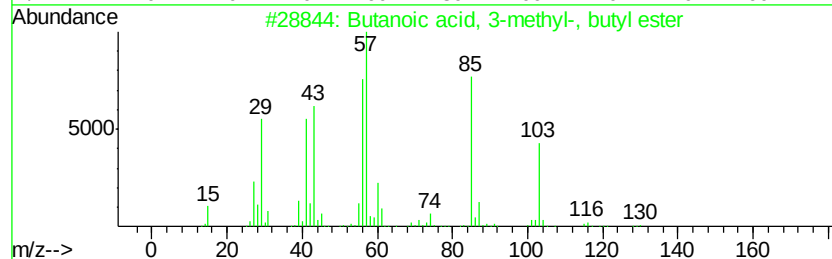
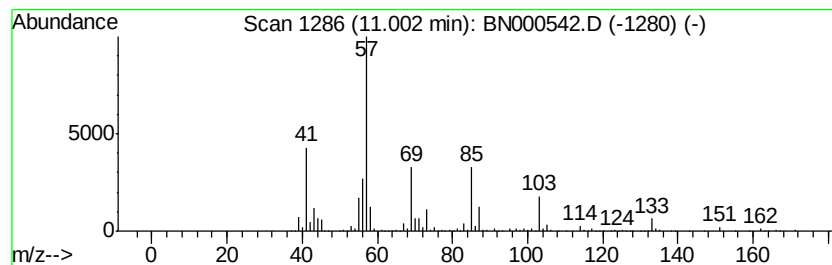
Instrument :  
BNA\_N  
ClientSampleId :  
DAM46DL

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 26 unknown-05 Concentration Rank 77

| R.T.      | EstConc                             | Area    | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------------|---------|------------------|-------------|------|
| 11.00     | 8.70 ng/ul                          | 1884370 | Napthalene-d8    | 11.26       |      |
| Hit# of 5 | Tentative ID                        | MW      | MolForm          | CAS#        | Qual |
| 1         | Butanoic acid, 3-methyl-, butyl ... | 158     | C9H18O2          | 000109-19-3 | 43   |
| 2         | Pentane, 2,2-dimethyl-              | 100     | C7H16            | 000590-35-2 | 37   |
| 3         | Pentane, 2,2-dimethyl-              | 100     | C7H16            | 000590-35-2 | 37   |
| 4         | Pentane, 2,2-dimethyl-              | 100     | C7H16            | 000590-35-2 | 37   |
| 5         | 2-Butanone, 1-chloro-3,3-dimethyl-  | 134     | C6H11ClO         | 013547-70-1 | 35   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000542.D  
Acq On : 22 Mar 2018 16:12  
Operator : JU/SJ  
Sample : J1905-11DL 10X  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

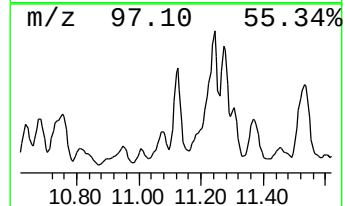
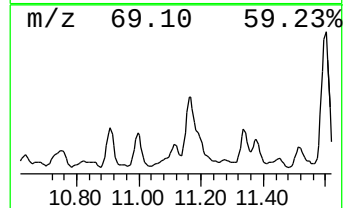
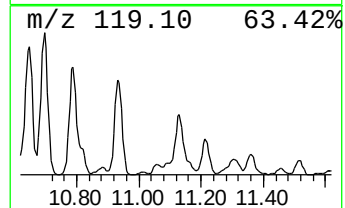
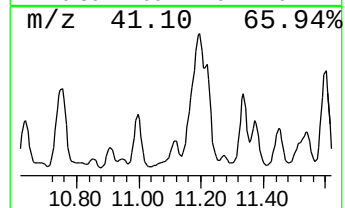
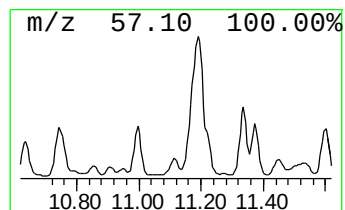
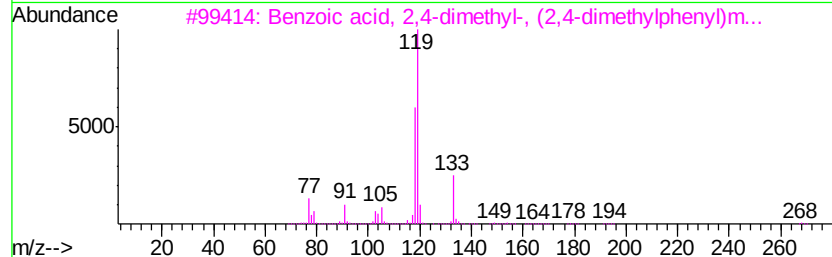
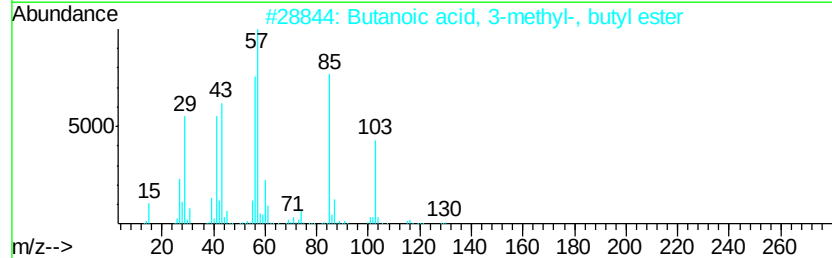
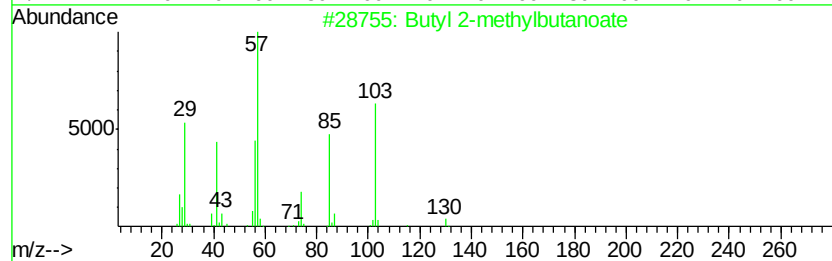
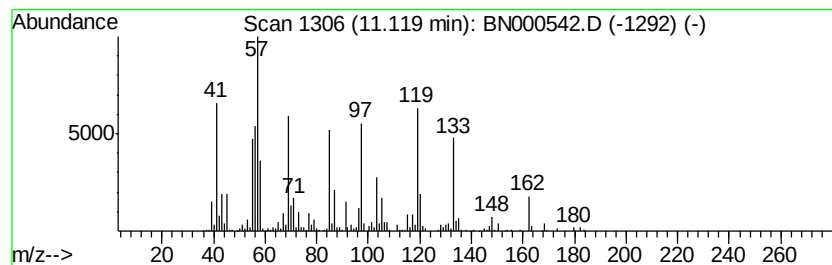
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ClientSampleId :  
DAM46DL

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 27 unknown-06 Concentration Rank 68

| R.T.    | EstConc     | Area                                | Relative to ISTD | R.T.            |
|---------|-------------|-------------------------------------|------------------|-----------------|
| 11.12   | 11.18 ng/ul | 2422090                             | Napthalene-d8    | 11.26           |
| Hit# of | 5           | Tentative ID                        | MW MolForm       | CAS# Qual       |
| 1       |             | Butyl 2-methylbutanoate             | 158 C9H18O2      | 015706-73-7 22  |
| 2       |             | Butanoic acid, 3-methyl-, butyl ... | 158 C9H18O2      | 000109-19-3 14  |
| 3       |             | Benzoic acid, 2,4-dimethyl-, (2,... | 268 C18H20O2     | 055000-43-6 14  |
| 4       |             | Pentanoic acid, butyl ester         | 158 C9H18O2      | 000591-68-4 12  |
| 5       |             | 2-Nitro-1-phenyl-3-(tetrahydro-p... | 281 C14H19NO5    | 1000190-43-3 12 |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000542.D  
Acq On : 22 Mar 2018 16:12  
Operator : JU/SJ  
Sample : J1905-11DL 10X  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleID :  
DAM46DL

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

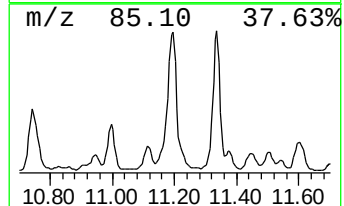
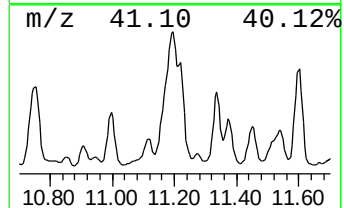
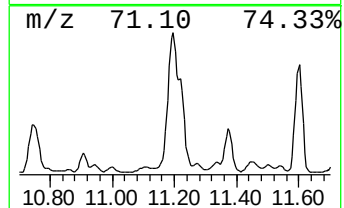
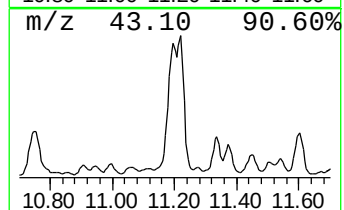
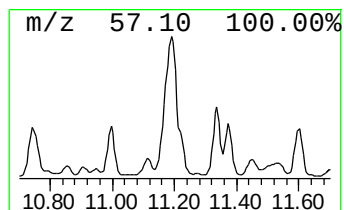
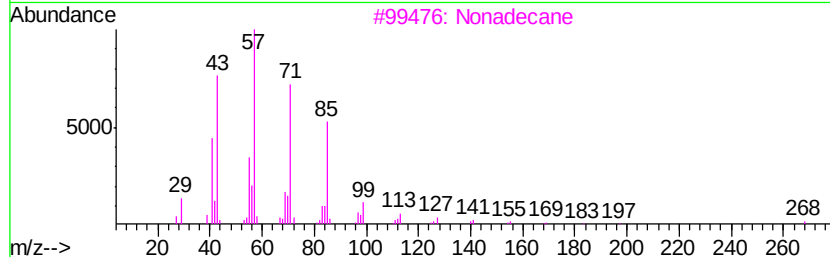
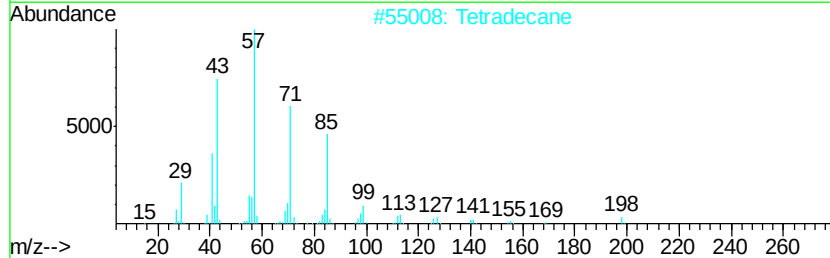
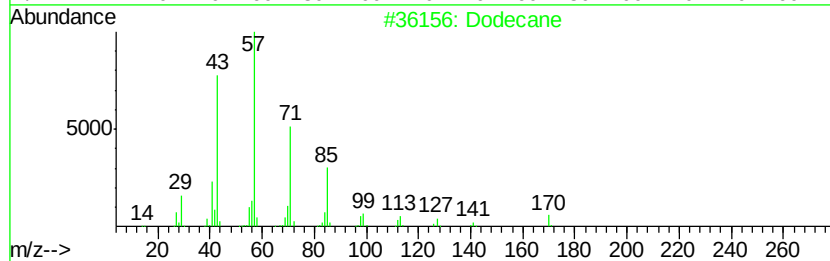
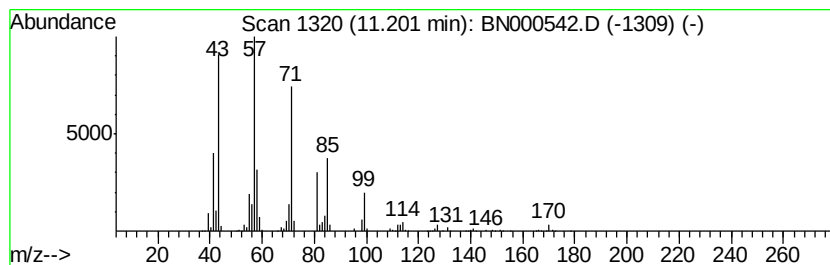
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 28 (DEL) Alkane: Straight-Chai... Concentration Rank 10

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 11.20 | 37.26 ng/ul | 8069710 | Napthalene-d8    | 11.26 |

| Hit# | of | Tentative ID        | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------|-----|---------|-------------|------|
| 1    | 5  | Dodecane            | 170 | C12H26  | 000112-40-3 | 90   |
| 2    |    | Tetradecane         | 198 | C14H30  | 000629-59-4 | 80   |
| 3    |    | Nonadecane          | 268 | C19H40  | 000629-92-5 | 80   |
| 4    |    | Tridecane           | 184 | C13H28  | 000629-50-5 | 80   |
| 5    |    | Undecane, 2-methyl- | 170 | C12H26  | 007045-71-8 | 80   |





Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000542.D  
Acq On : 22 Mar 2018 16:12  
Operator : JU/SJ  
Sample : J1905-11DL 10X  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

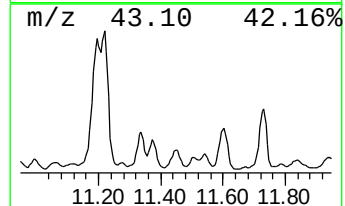
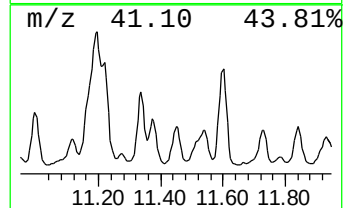
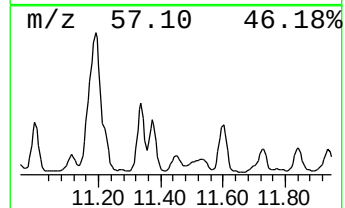
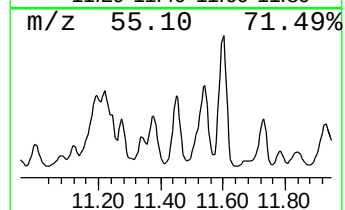
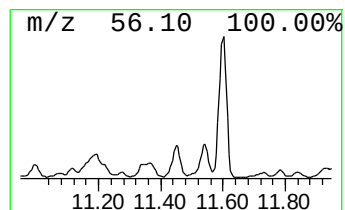
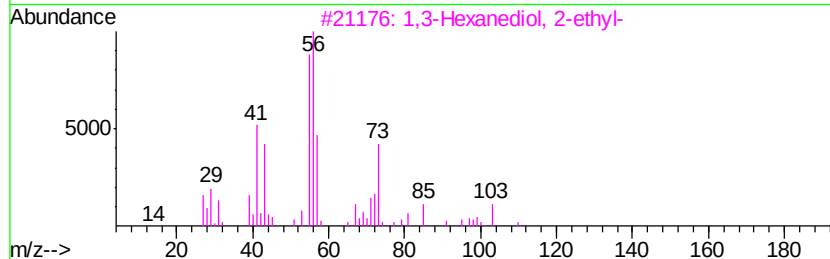
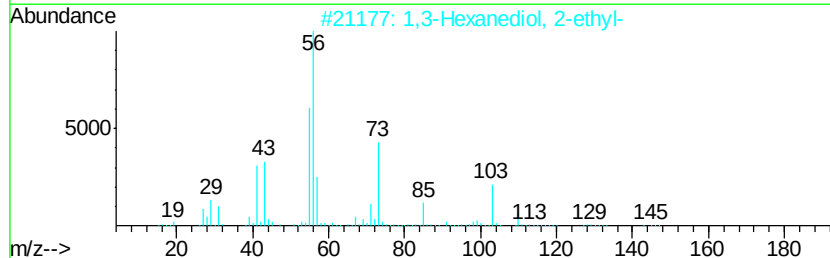
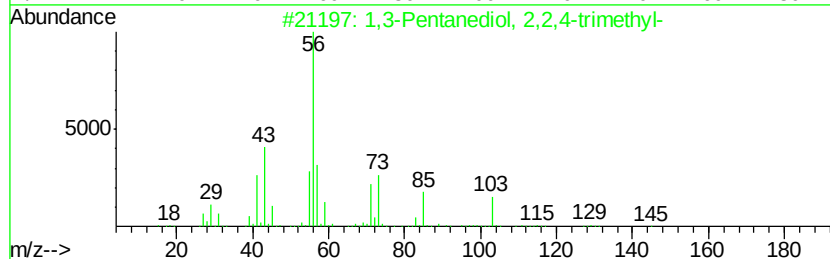
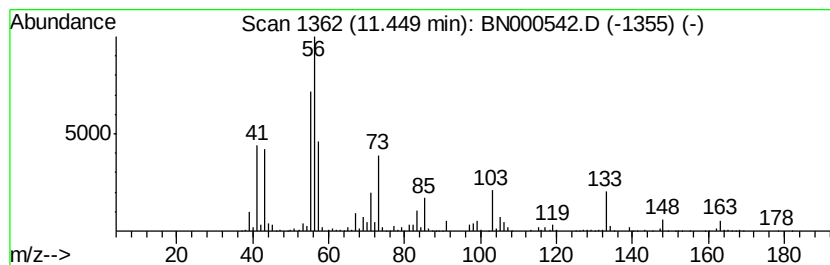
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BNA\_N  
ClientSampleId :  
DAM46DL

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 29 1,3-Pentanediol, 2,2,4-trim... Concentration Rank 72

| R.T.    | EstConc                           | Area         | Relative to ISTD | R.T.      |
|---------|-----------------------------------|--------------|------------------|-----------|
| 11.45   | 9.59 ng/ul                        | 2077730      | Napthalene-d8    | 11.26     |
| Hit# of | 5                                 | Tentative ID | MW MolForm       | CAS# Qual |
| 1       | 1,3-Pentanediol, 2,2,4-trimethyl- | 146 C8H18O2  | 000144-19-4      | 64        |
| 2       | 1,3-Hexanediol, 2-ethyl-          | 146 C8H18O2  | 000094-96-2      | 59        |
| 3       | 1,3-Hexanediol, 2-ethyl-          | 146 C8H18O2  | 000094-96-2      | 53        |
| 4       | Pentanal, 3-methyl-               | 100 C6H12O   | 015877-57-3      | 43        |
| 5       | 1,3-Propanediol, 2,2-dimethyl-    | 104 C5H12O2  | 000126-30-7      | 38        |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000542.D  
Acq On : 22 Mar 2018 16:12  
Operator : JU/SJ  
Sample : J1905-11DL 10X  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM46DL

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

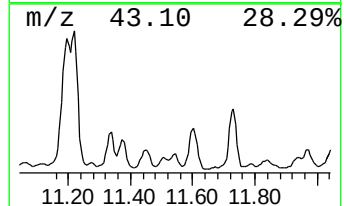
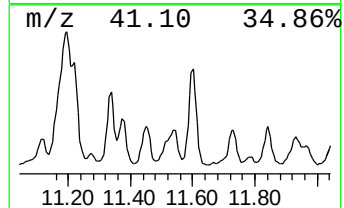
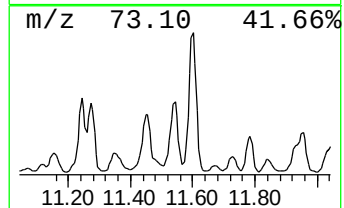
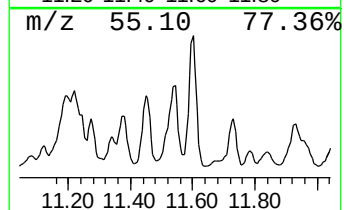
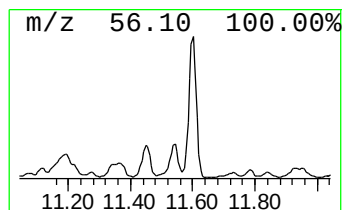
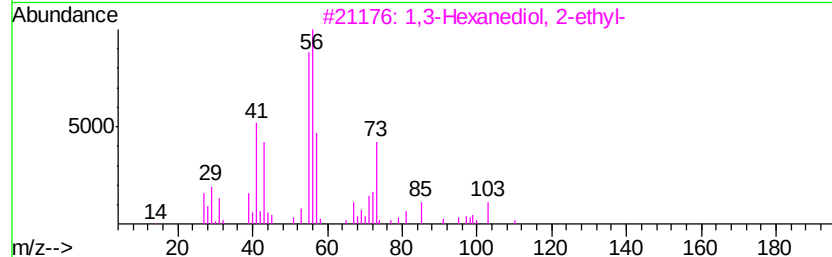
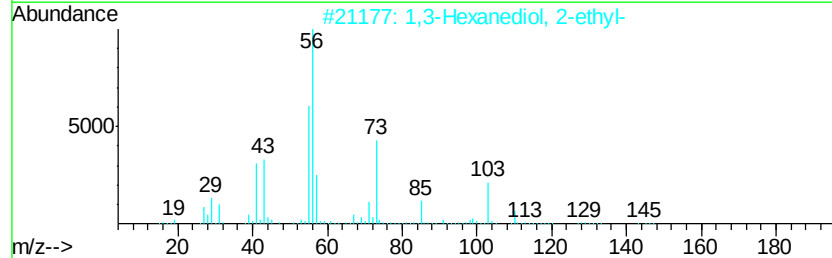
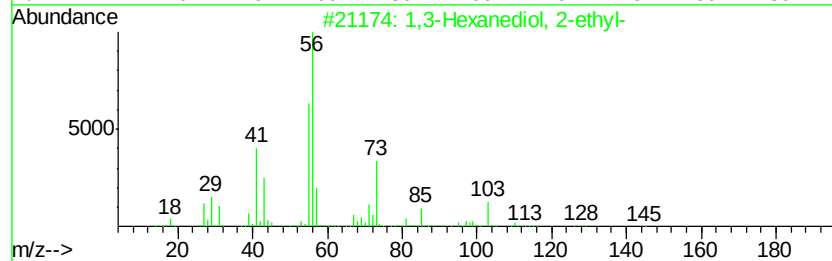
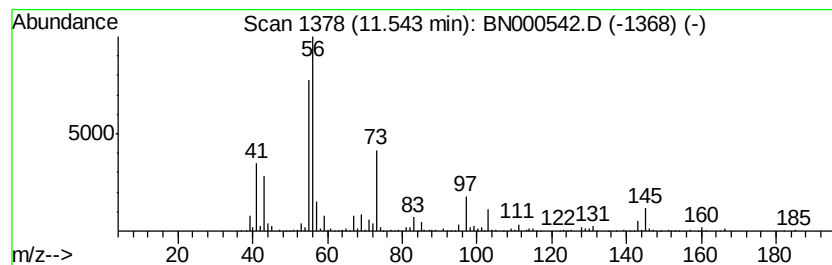
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TIC Integration Parameters: LSCINT.P

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Peak Number 30 1,3-Hexanediol, 2-ethyl- Concentration Rank 66

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 11.54 | 11.81 ng/ul | 2557830 | Napthalene-d8    | 11.26 |

| Hit# | of                               | 5 | Tentative ID | MW  | MolForm  | CAS#        | Qual |
|------|----------------------------------|---|--------------|-----|----------|-------------|------|
| 1    | 1,3-Hexanediol, 2-ethyl-         |   |              | 146 | C8H18O2  | 000094-96-2 | 59   |
| 2    | 1,3-Hexanediol, 2-ethyl-         |   |              | 146 | C8H18O2  | 000094-96-2 | 59   |
| 3    | 1,3-Hexanediol, 2-ethyl-         |   |              | 146 | C8H18O2  | 000094-96-2 | 45   |
| 4    | 3-Buten-1-ol, 2-methyl-          |   |              | 86  | C5H10O   | 004516-90-9 | 43   |
| 5    | 3-Chloro-2,2-dimethyl-1-propanol |   |              | 122 | C5H11ClO | 013401-56-4 | 42   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000542.D  
Acq On : 22 Mar 2018 16:12  
Operator : JU/SJ  
Sample : J1905-11DL 10X  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM46DL

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

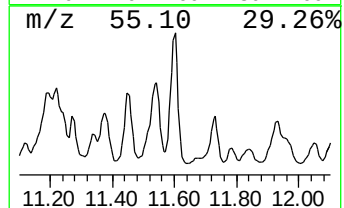
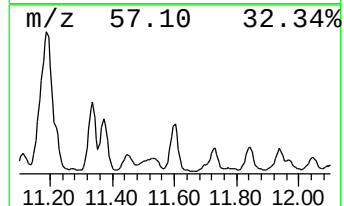
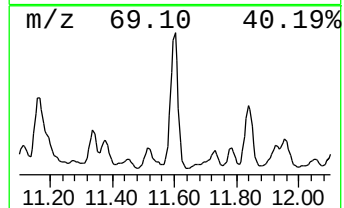
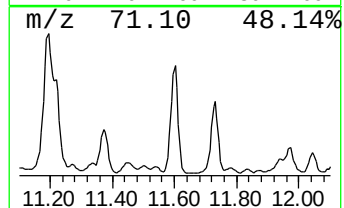
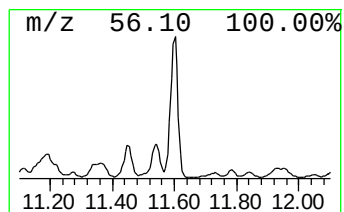
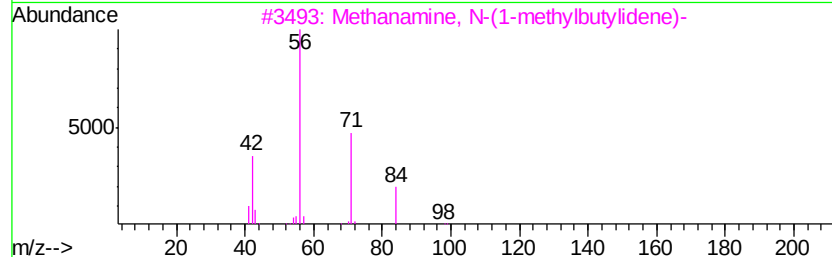
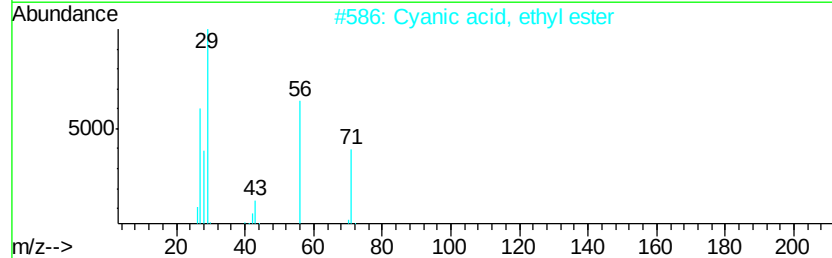
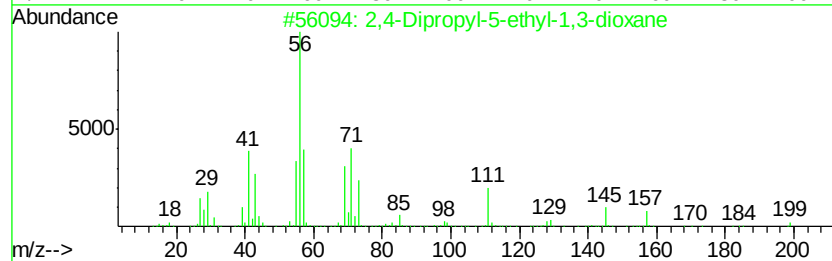
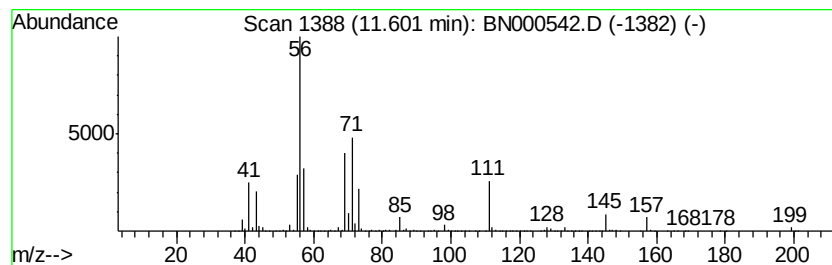
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 31 2,4-Dipropyl-5-ethyl-1,3-di... Concentration Rank 15

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 11.60 | 28.61 ng/ul | 6196920 | Napthalene-d8    | 11.26 |

| Hit# | of | 5 | Tentative ID                         | MW  | MolForm  | CAS#        | Qual |
|------|----|---|--------------------------------------|-----|----------|-------------|------|
| 1    |    |   | 2,4-Dipropyl-5-ethyl-1,3-dioxane     | 200 | C12H24O2 | 006413-83-8 | 59   |
| 2    |    |   | Cyanoic acid, ethyl ester            | 71  | C3H5NO   | 000627-48-5 | 40   |
| 3    |    |   | Methanamine, N-(1-methylbutylidene)- | 99  | C6H13N   | 022431-09-0 | 38   |
| 4    |    |   | Oxirane, 2,3-bis(1-methylethyl)-     | 128 | C8H16O   | 054644-32-5 | 38   |
| 5    |    |   | Cyclohexanamine                      | 99  | C6H13N   | 000108-91-8 | 25   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000542.D  
Acq On : 22 Mar 2018 16:12  
Operator : JU/SJ  
Sample : J1905-11DL 10X  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleID :  
DAM46DL

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

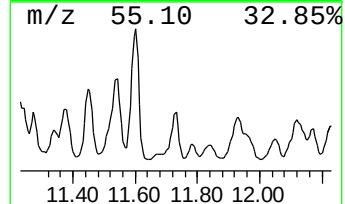
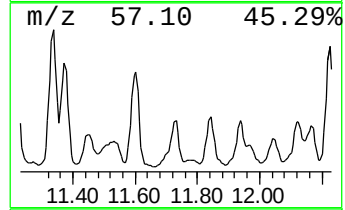
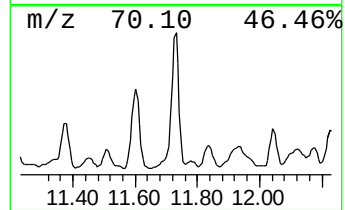
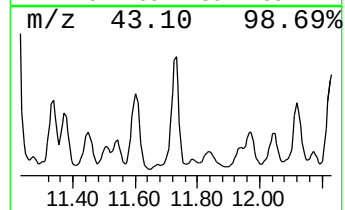
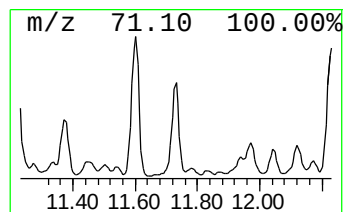
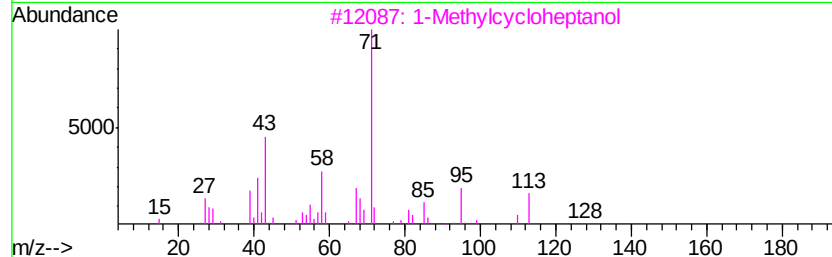
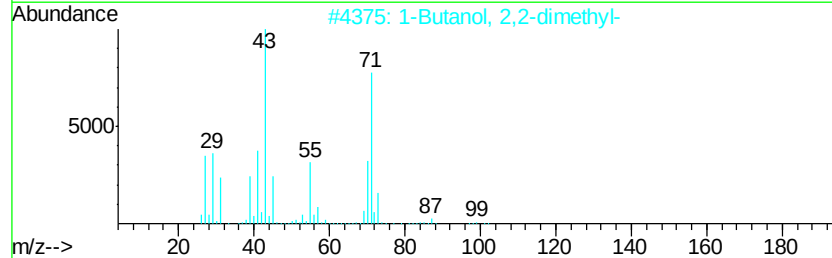
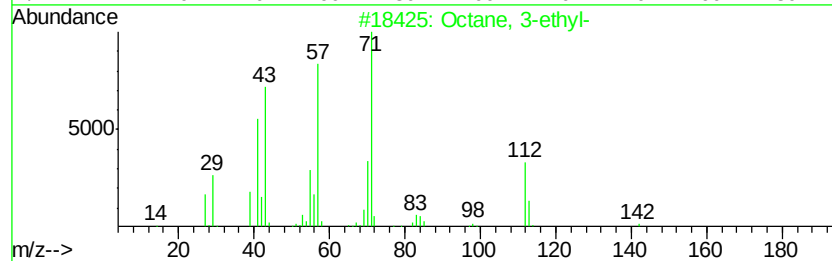
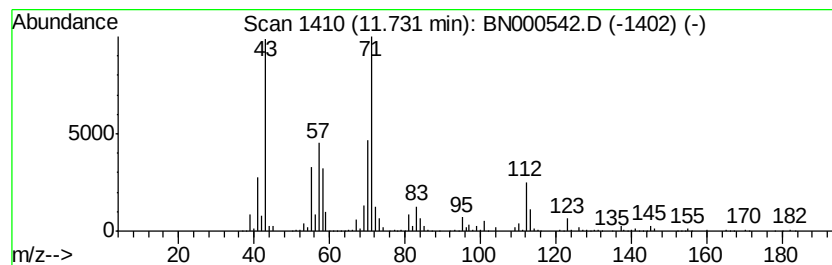
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 32 (DEL) Alkane: Straight-Chai... Concentration Rank 63

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 11.73 | 12.47 ng/ul | 2700050 | Napthalene-d8    | 11.26 |

| Hit# | of | Tentative ID                      | MW  | MolForm  | CAS#        | Qual |
|------|----|-----------------------------------|-----|----------|-------------|------|
| 1    | 5  | Octane, 3-ethyl-                  | 142 | C10H22   | 005881-17-4 | 50   |
| 2    |    | 1-Butanol, 2,2-dimethyl-          | 102 | C6H14O   | 001185-33-7 | 50   |
| 3    |    | 1-Methylcycloheptanol             | 128 | C8H16O   | 003761-94-2 | 50   |
| 4    |    | 2,4,4-Trimethyl-1-hexene          | 126 | C9H18    | 051174-12-0 | 47   |
| 5    |    | n-Butyric acid 2-ethylhexyl ester | 200 | C12H24O2 | 025415-84-3 | 43   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000542.D  
Acq On : 22 Mar 2018 16:12  
Operator : JU/SJ  
Sample : J1905-11DL 10X  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

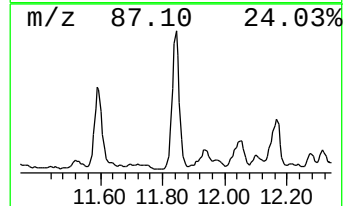
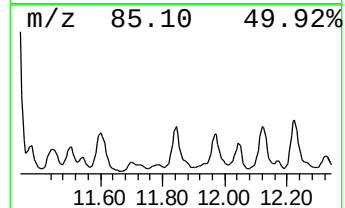
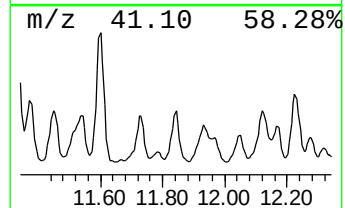
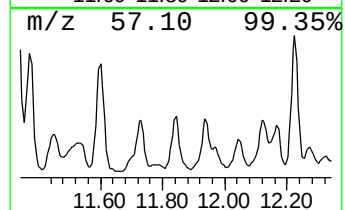
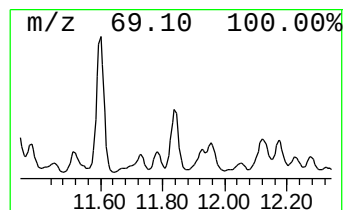
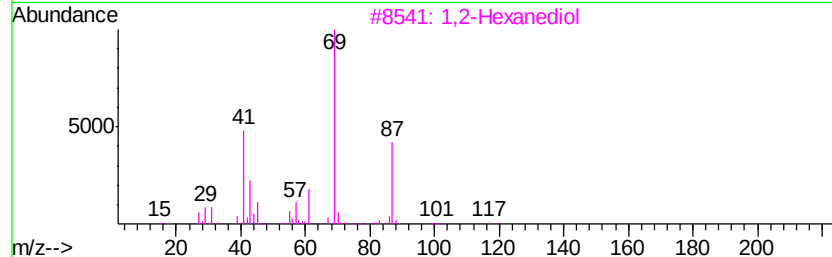
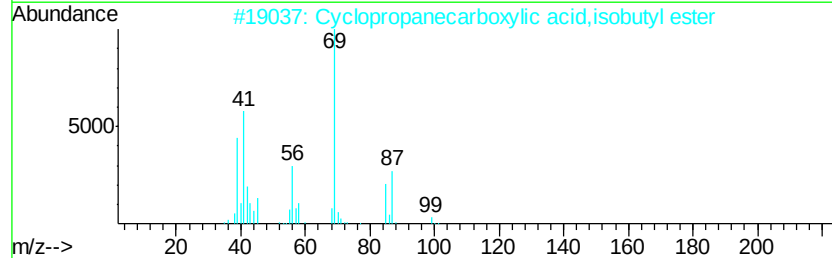
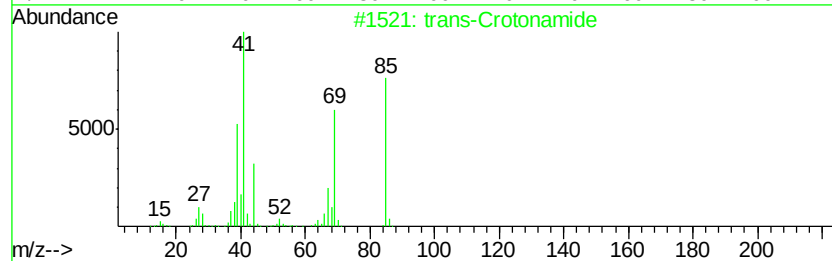
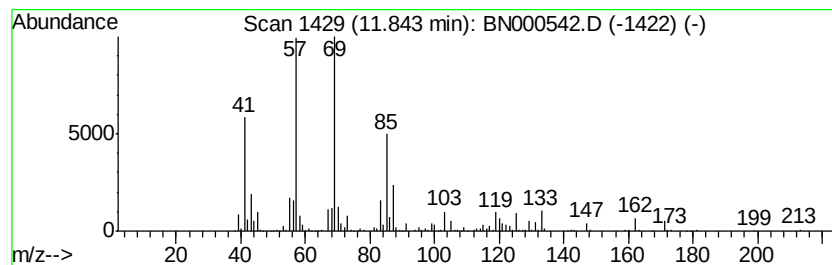
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ClientSampleId :  
DAM46DL

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 33 unknown-07 Concentration Rank 76

| R.T.    | EstConc                             | Area         | Relative to ISTD | R.T.           |
|---------|-------------------------------------|--------------|------------------|----------------|
| 11.84   | 9.11 ng/ul                          | 1973920      | Napthalene-d8    | 11.26          |
| Hit# of | 5                                   | Tentative ID | MW MolForm       | CAS# Qual      |
| 1       | trans-Crotonamide                   | 85           | C4H7NO           | 000625-37-6 43 |
| 2       | Cyclopropanecarboxylic acid,isob... | 142          | C8H14O2          | 064969-79-5 43 |
| 3       | 1,2-Hexanediol                      | 118          | C6H14O2          | 006920-22-5 38 |
| 4       | 5-Tridecanol                        | 200          | C13H28O          | 058783-82-7 35 |
| 5       | 5-Dodecanol                         | 186          | C12H26O          | 010203-33-5 35 |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000542.D  
Acq On : 22 Mar 2018 16:12  
Operator : JU/SJ  
Sample : J1905-11DL 10X  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM46DL

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

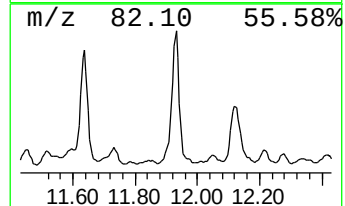
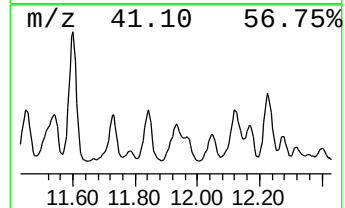
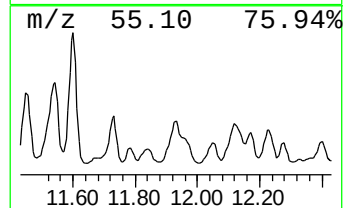
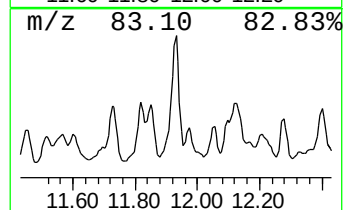
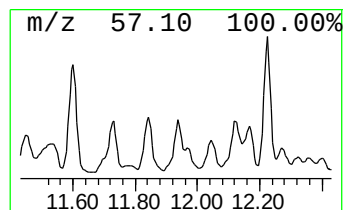
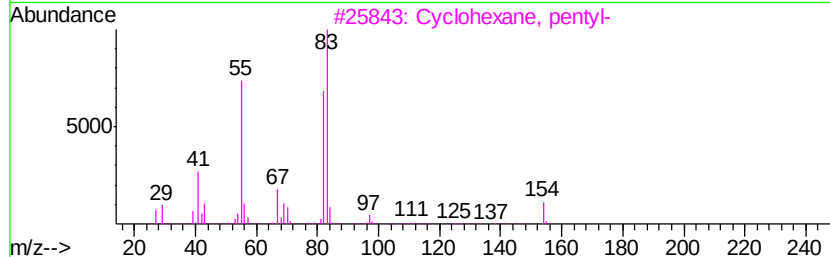
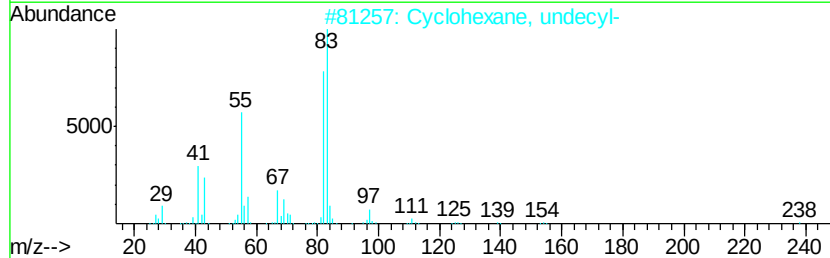
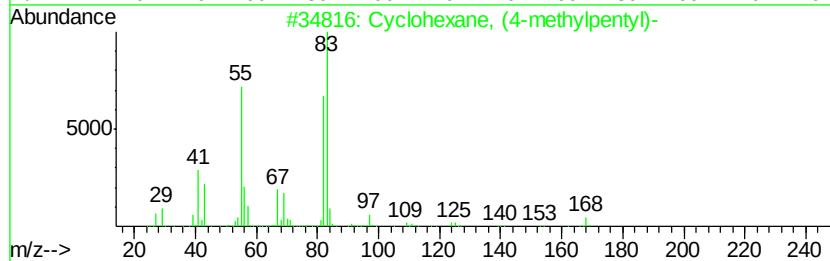
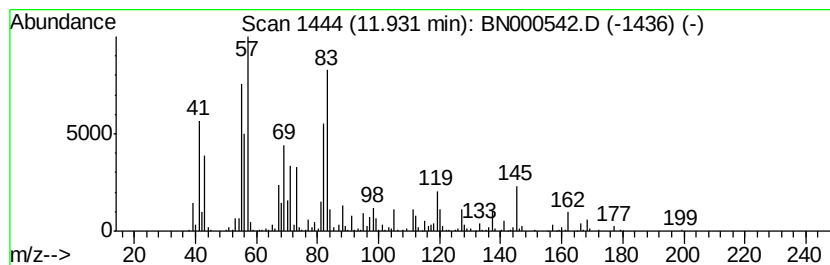
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 34 unknown-08 Concentration Rank 65

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 11.93 | 12.04 ng/ul | 2606680 | Napthalene-d8    | 11.26 |

| Hit# | of | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------|-----|---------|-------------|------|
| 1    | 5  | Cyclohexane, (4-methylpentyl)- | 168 | C12H24  | 061142-20-9 | 43   |
| 2    |    | Cyclohexane, undecyl-          | 238 | C17H34  | 054105-66-7 | 38   |
| 3    |    | Cyclohexane, pentyl-           | 154 | C11H22  | 004292-92-6 | 38   |
| 4    |    | Cyclohexane, hexyl-            | 168 | C12H24  | 004292-75-5 | 38   |
| 5    |    | Cyclohexane, (1-methylethyl)-  | 126 | C9H18   | 000696-29-7 | 38   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000542.D  
Acq On : 22 Mar 2018 16:12  
Operator : JU/SJ  
Sample : J1905-11DL 10X  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

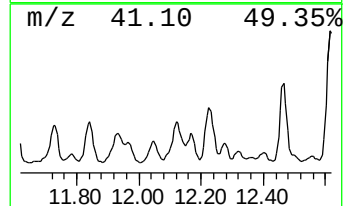
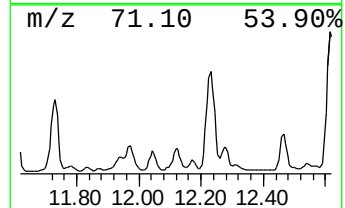
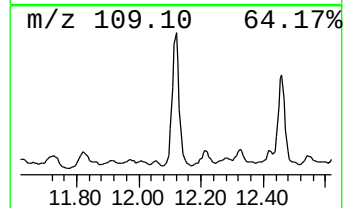
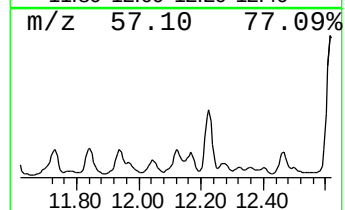
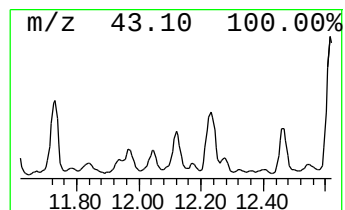
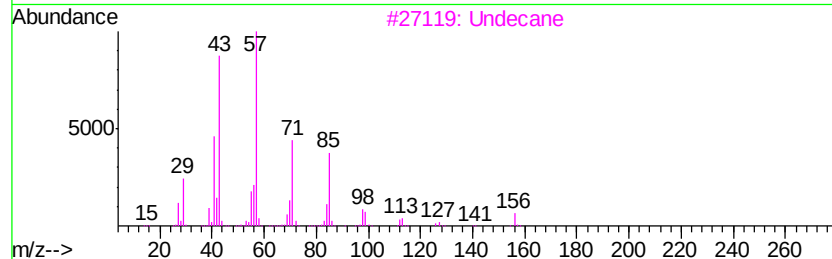
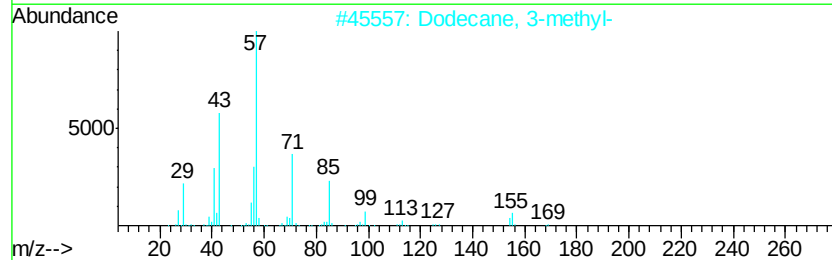
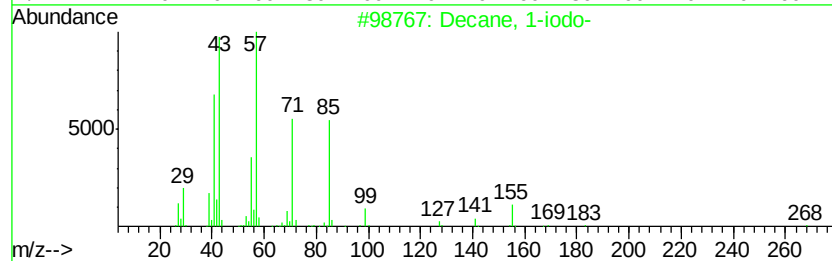
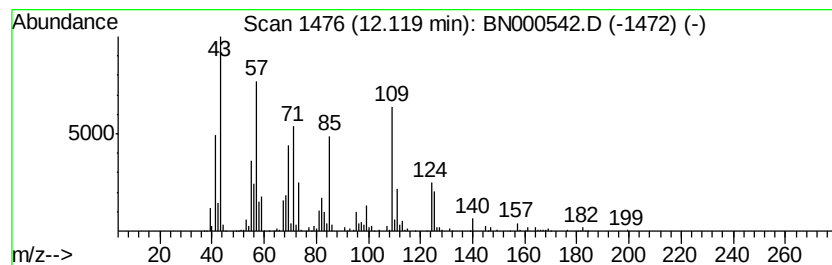
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BNA\_N  
ClientSampleId :  
DAM46DL

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 35 unknown-09 Concentration Rank 61

| R.T.    | EstConc                  | Area         | Relative to ISTD | R.T.      |
|---------|--------------------------|--------------|------------------|-----------|
| 12.12   | 12.79 ng/ul              | 2770540      | Napthalene-d8    | 11.26     |
| Hit# of | 5                        | Tentative ID | MW MolForm       | CAS# Qual |
| 1       | Decane, 1-iodo-          | 268 C10H21I  | 002050-77-3      | 30        |
| 2       | Dodecane, 3-methyl-      | 184 C13H28   | 017312-57-1      | 27        |
| 3       | Undecane                 | 156 C11H24   | 001120-21-4      | 27        |
| 4       | Undecane, 2,4-dimethyl-  | 184 C13H28   | 017312-80-0      | 25        |
| 5       | Hexane, 2,3,4-trimethyl- | 128 C9H20    | 000921-47-1      | 25        |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000542.D  
Acq On : 22 Mar 2018 16:12  
Operator : JU/SJ  
Sample : J1905-11DL 10X  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM46DL

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

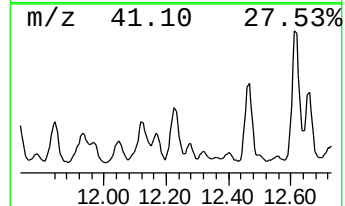
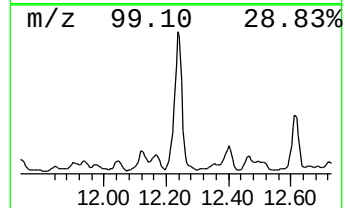
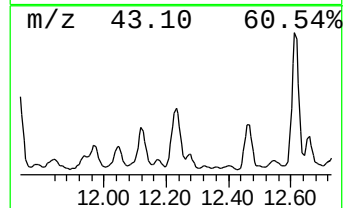
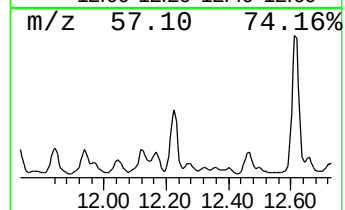
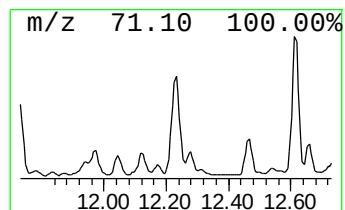
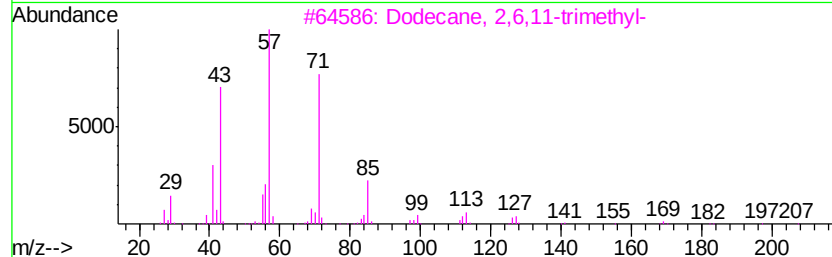
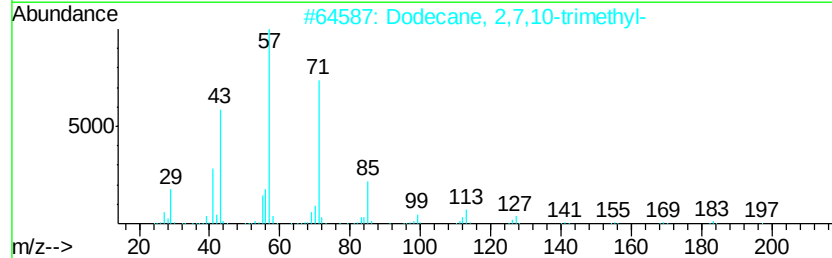
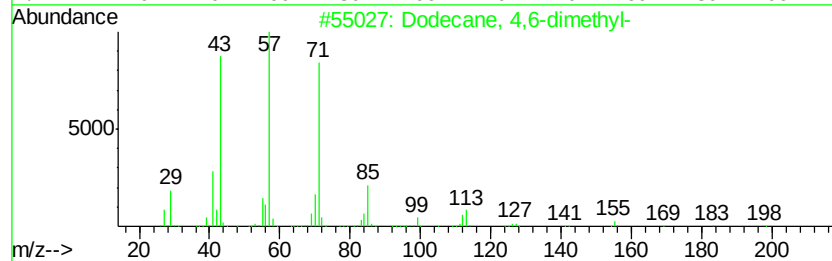
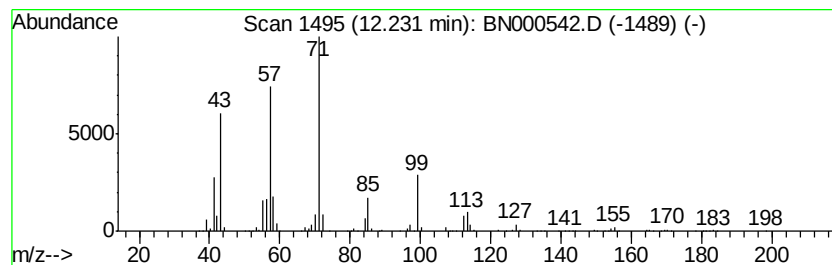
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 36 (DEL) Alkane: Straight-Chai... Concentration Rank 52

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 12.23 | 14.86 ng/ul | 3217370 | Napthalene-d8    | 11.26 |

| Hit# | of | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------------|-----|---------|-------------|------|
| 1    | 5  | Dodecane, 4,6-dimethyl-     | 198 | C14H30  | 061141-72-8 | 90   |
| 2    |    | Dodecane, 2,7,10-trimethyl- | 212 | C15H32  | 074645-98-0 | 83   |
| 3    |    | Dodecane, 2,6,11-trimethyl- | 212 | C15H32  | 031295-56-4 | 83   |
| 4    |    | Nonane, 2,6-dimethyl-       | 156 | C11H24  | 017302-28-2 | 72   |
| 5    |    | Octane, 3,6-dimethyl-       | 142 | C10H22  | 015869-94-0 | 72   |





Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000542.D  
Acq On : 22 Mar 2018 16:12  
Operator : JU/SJ  
Sample : J1905-11DL 10X  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM46DL

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

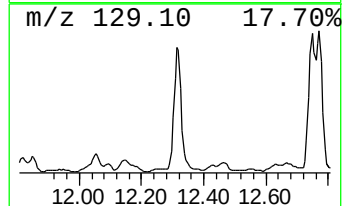
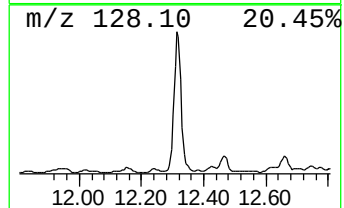
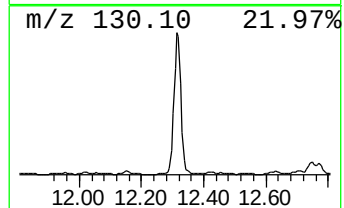
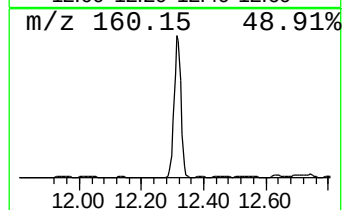
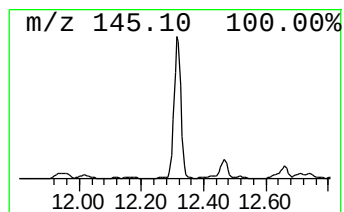
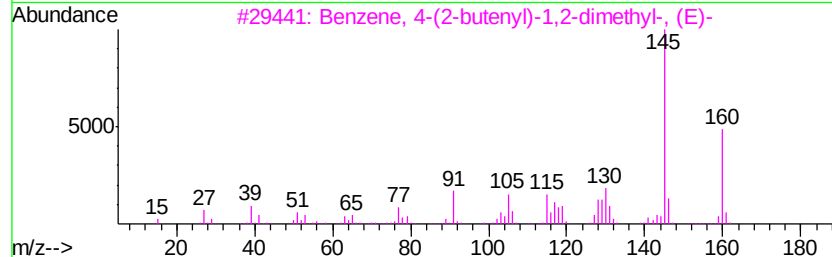
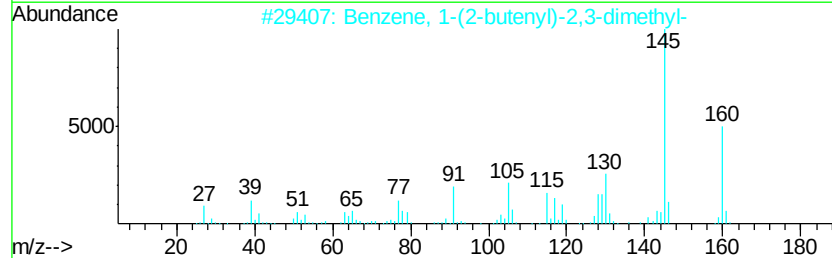
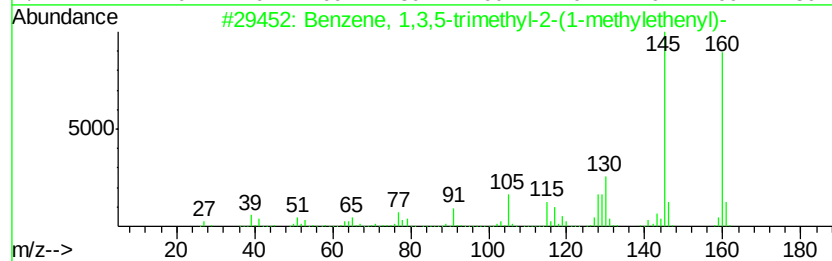
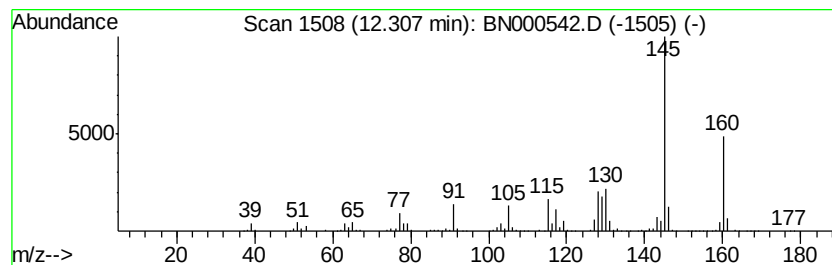
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 37 Benzene, 1,3,5-trimethyl-2-... Concentration Rank 36

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 12.31 | 19.25 ng/ul | 4170250 | Naphthalene-d8   | 11.26 |

| Hit# | of | Tentative ID                                  | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1,3,5-trimethyl-2-(1-methylethenyl)- | 160 | C12H16  | 014679-13-1 | 94   |
| 2    |    | Benzene, 1-(2-butenyl)-2,3-dimethyl-          | 160 | C12H16  | 054340-85-1 | 94   |
| 3    |    | Benzene, 4-(2-butenyl)-1,2-dimethyl-          | 160 | C12H16  | 054340-86-2 | 94   |
| 4    |    | Benzene, 1-(2-butenyl)-2,3-dimethyl-          | 160 | C12H16  | 054340-85-1 | 91   |
| 5    |    | Naphthalene, 1,2,3,4-tetrahydro-              | 160 | C12H16  | 004175-54-6 | 87   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000542.D  
Acq On : 22 Mar 2018 16:12  
Operator : JU/SJ  
Sample : J1905-11DL 10X  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

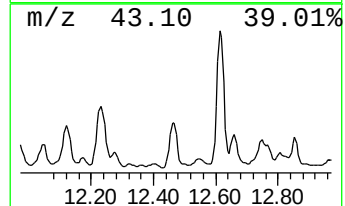
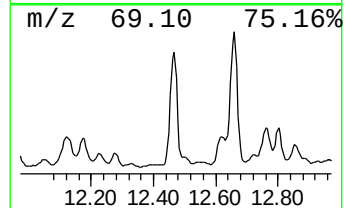
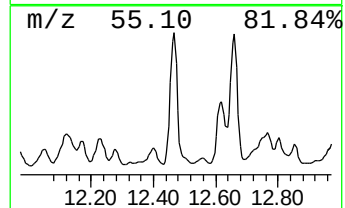
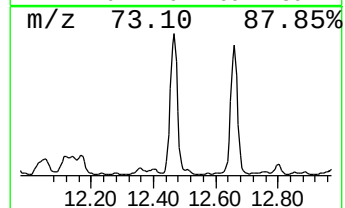
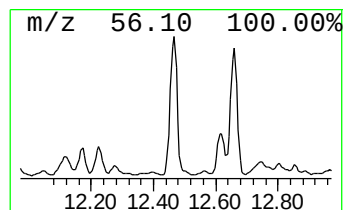
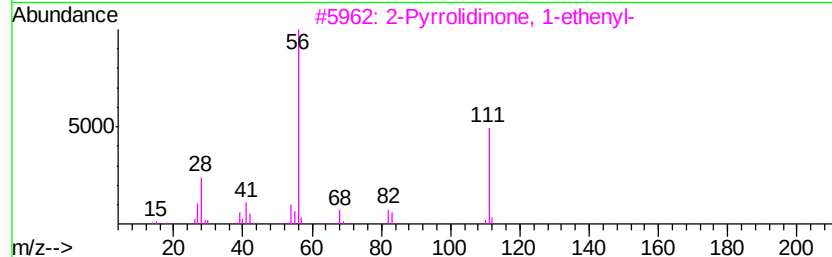
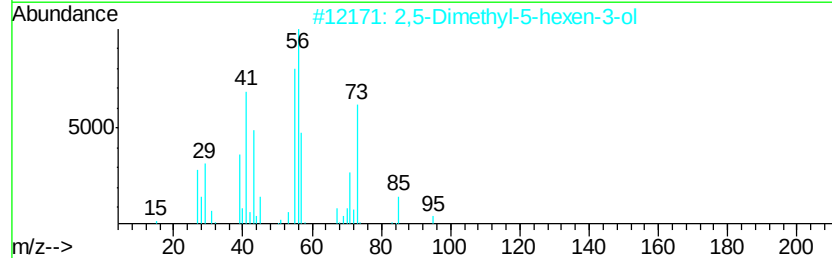
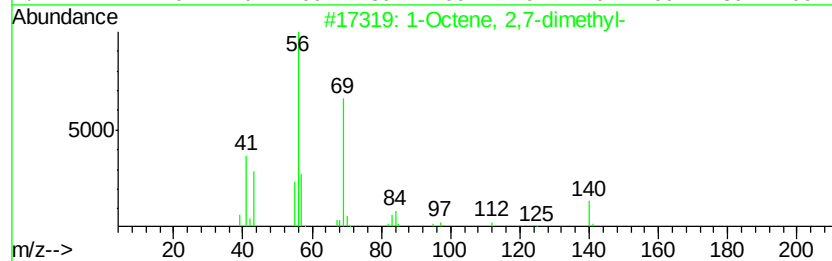
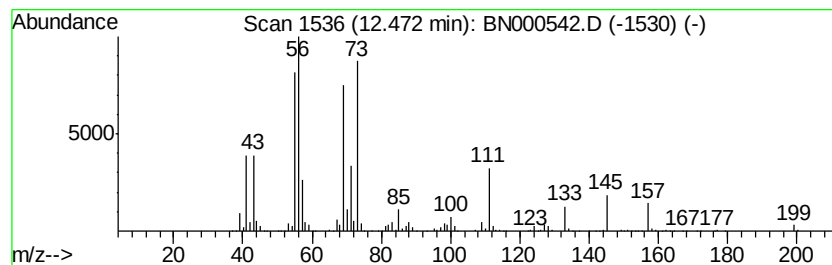
Instrument :  
BNA\_N  
ClientSampleId :  
DAM46DL

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 38 unknown-10 Concentration Rank 28

| R.T.    | EstConc                          | Area         | Relative to ISTD | R.T.      |
|---------|----------------------------------|--------------|------------------|-----------|
| 12.47   | 22.15 ng/ul                      | 4798340      | Napthalene-d8    | 11.26     |
| Hit# of | 5                                | Tentative ID | MW MolForm       | CAS# Qual |
| 1       | 1-Octene, 2,7-dimethyl-          | 140 C10H20   | 033718-03-5      | 47        |
| 2       | 2,5-Dimethyl-5-hexen-3-ol        | 128 C8H16O   | 067760-91-2      | 32        |
| 3       | 2-Pyrrolidinone, 1-ethenvl-      | 111 C6H9NO   | 000088-12-0      | 27        |
| 4       | Decane, 4-methylene-             | 154 C11H22   | 024949-41-5      | 27        |
| 5       | 2,4-Dipropyl-5-ethyl-1,3-dioxane | 200 C12H24O2 | 006413-83-8      | 25        |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000542.D  
Acq On : 22 Mar 2018 16:12  
Operator : JU/SJ  
Sample : J1905-11DL 10X  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM46DL

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

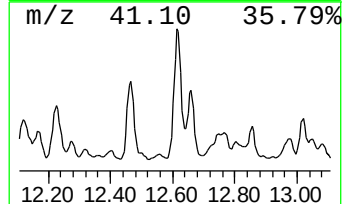
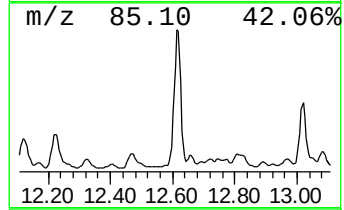
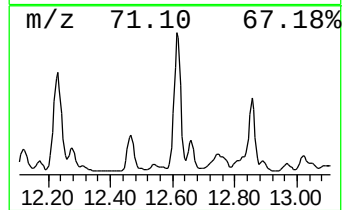
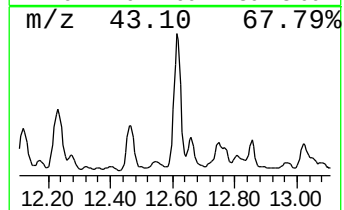
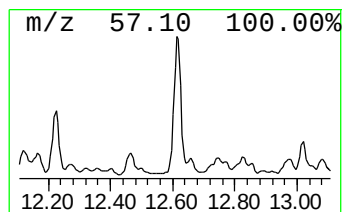
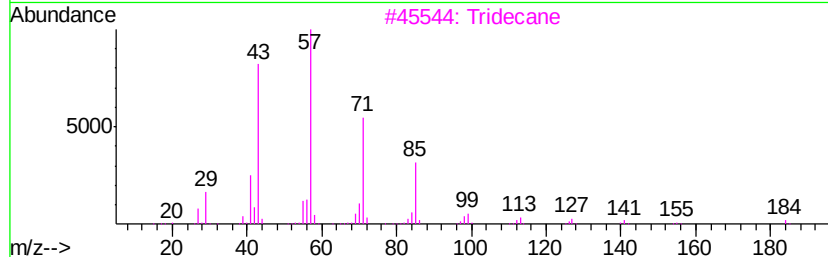
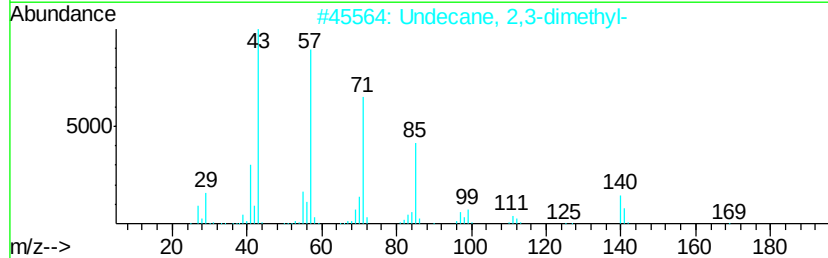
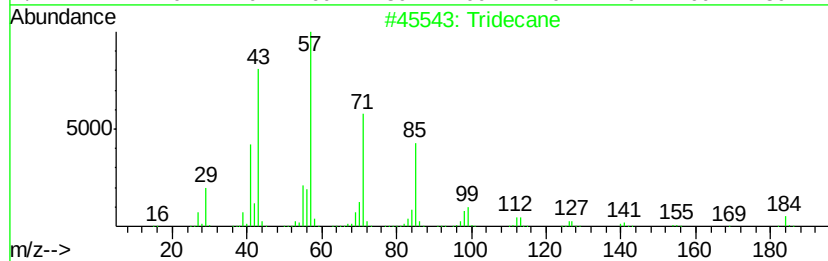
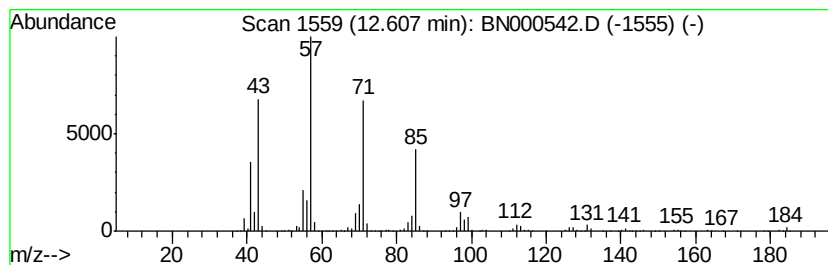
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 39 (DEL) Alkane: Straight-Chai... Concentration Rank 16

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 12.61 | 27.85 ng/ul | 6031380 | Napthalene-d8    | 11.26 |

| Hit# | of | Tentative ID            | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------|-----|----------|-------------|------|
| 1    | 5  | Tridecane               | 184 | C13H28   | 000629-50-5 | 87   |
| 2    |    | Undecane, 2,3-dimethyl- | 184 | C13H28   | 017312-77-5 | 83   |
| 3    |    | Tridecane               | 184 | C13H28   | 000629-50-5 | 81   |
| 4    |    | Tridecane               | 184 | C13H28   | 000629-50-5 | 81   |
| 5    |    | 2-Bromo dodecane        | 248 | C12H25Br | 013187-99-0 | 72   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000542.D  
Acq On : 22 Mar 2018 16:12  
Operator : JU/SJ  
Sample : J1905-11DL 10X  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM46DL

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

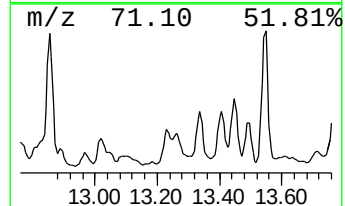
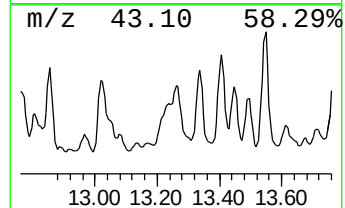
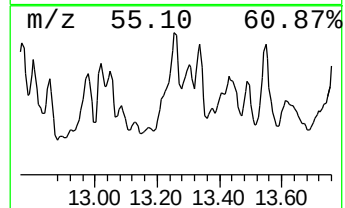
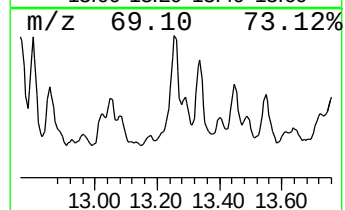
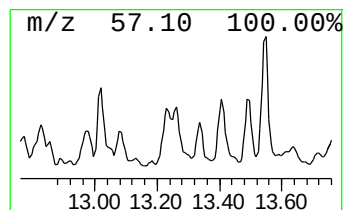
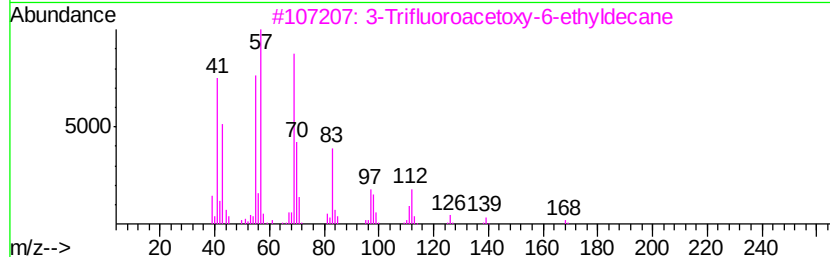
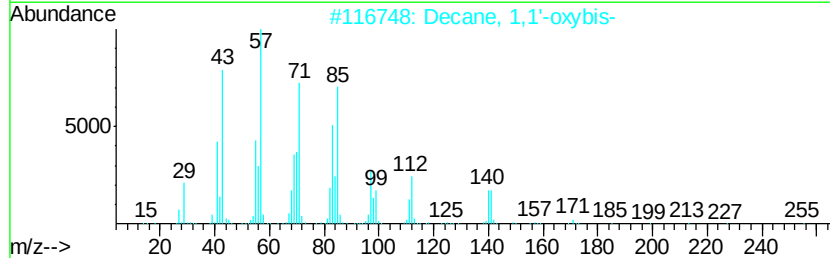
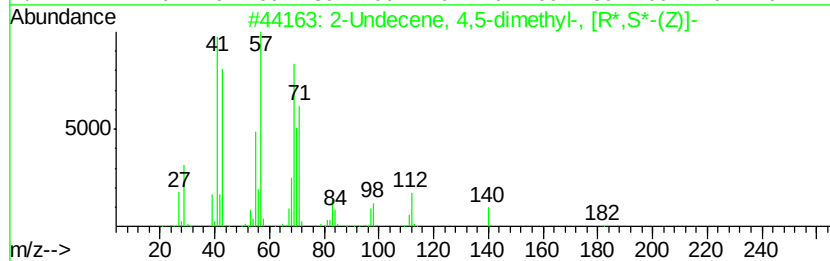
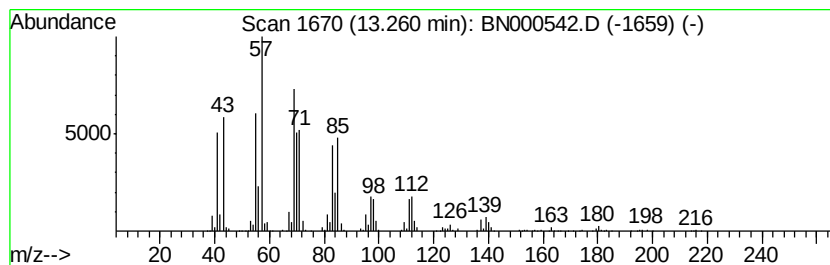
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 44 (DEL) Alkane: Cyclic13.26 Concentration Rank 64

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 13.26 | 12.11 ng/ul | 2843950 | Acenaphthene-d10 | 15.04 |

| Hit# | of | Tentative ID                        | MW  | MolForm    | CAS#        | Qual |
|------|----|-------------------------------------|-----|------------|-------------|------|
| 1    | 5  | 2-Undecene, 4,5-dimethyl-, [R*,S... | 182 | C13H26     | 055170-93-9 | 64   |
| 2    |    | Decane, 1,1'-oxybis-                | 298 | C20H42O    | 002456-28-2 | 50   |
| 3    |    | 3-Trifluoroacetoxy-6-ethyldecane    | 282 | C14H25F3O2 | 116436-59-0 | 43   |
| 4    |    | Cyclooctane, 1,4-dimethyl-, cis-    | 140 | C10H20     | 013151-99-0 | 41   |
| 5    |    | 1-Hexene, 3,3-dimethyl-             | 112 | C8H16      | 003404-77-1 | 38   |



Data Path : Z:\HPCHEM1\BNA\_N\DATA\BN032218\  
Data File : BN000542.D  
Acq On : 22 Mar 2018 16:12  
Operator : JU/SJ  
Sample : J1905-11DL 10X  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM46DL

Quant Method : Z:\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

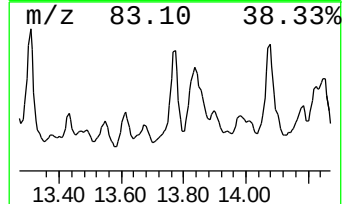
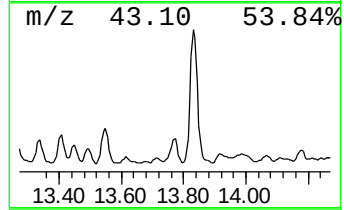
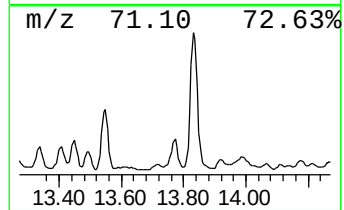
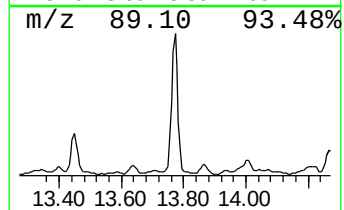
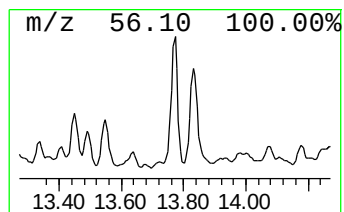
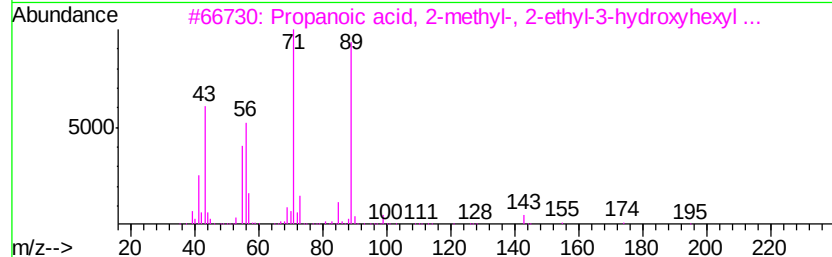
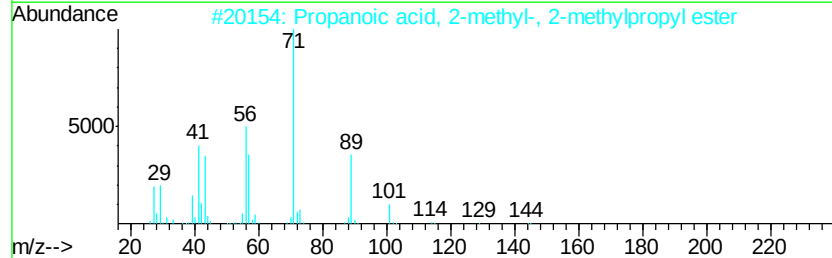
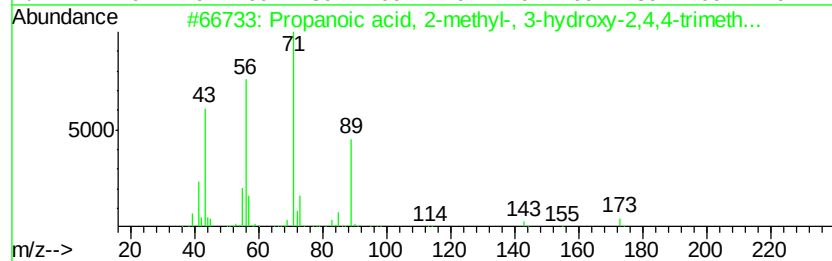
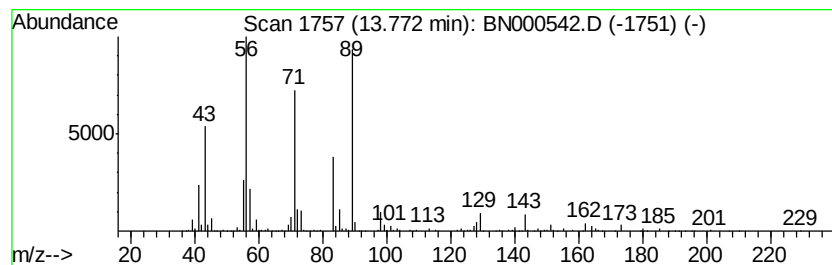
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 45 Propanoic acid, 2-methyl-, ... Concentration Rank 80

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T.  |
|-------|------------|---------|------------------|-------|
| 13.77 | 7.61 ng/ul | 1788060 | Acenaphthene-d10 | 15.04 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | Propanoic acid, 2-methyl-, 3-hyd... | 216 | C12H24O3 | 074367-34-3 | 59   |
| 2    |    | Propanoic acid, 2-methyl-, 2-met... | 144 | C8H16O2  | 000097-85-8 | 59   |
| 3    |    | Propanoic acid, 2-methyl-, 2-eth... | 216 | C12H24O3 | 074367-31-0 | 53   |
| 4    |    | Butyric acid, hexadecyl ester       | 312 | C20H40O2 | 006221-99-4 | 50   |
| 5    |    | Propanoic acid, 2-methyl-, decyl... | 228 | C14H28O2 | 005454-22-8 | 50   |



Data Path : Z:\HPCHEM1\BNA\_N\DATA\BN032218\  
Data File : BN000542.D  
Acq On : 22 Mar 2018 16:12  
Operator : JU/SJ  
Sample : J1905-11DL 10X  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleID :  
DAM46DL

Quant Method : Z:\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

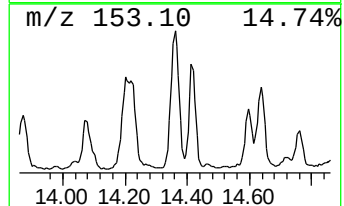
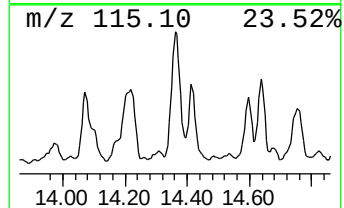
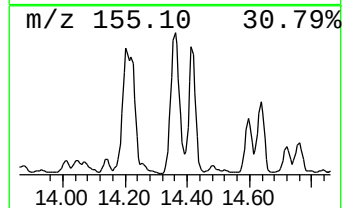
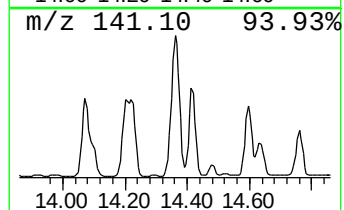
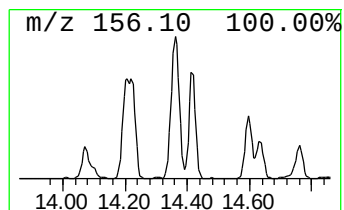
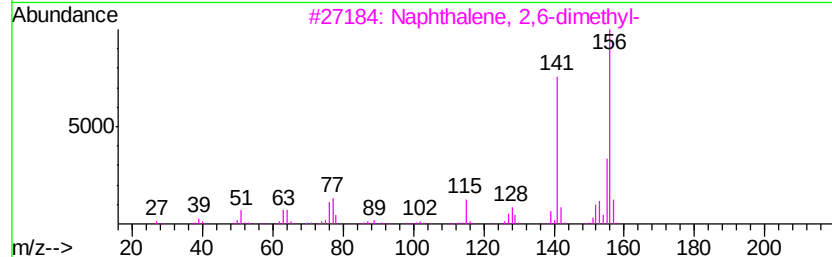
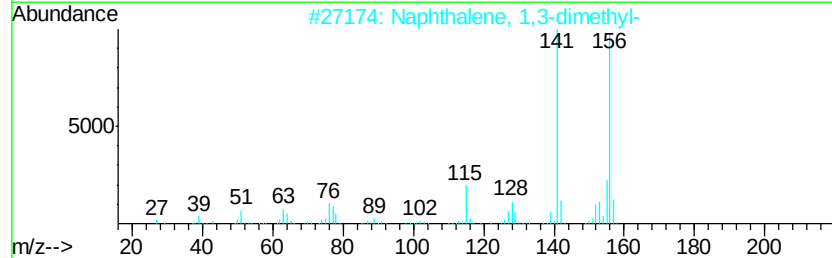
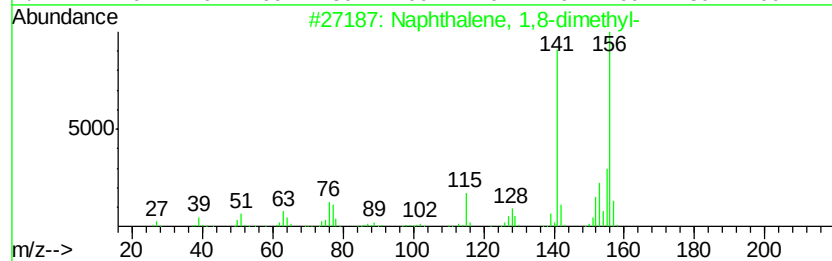
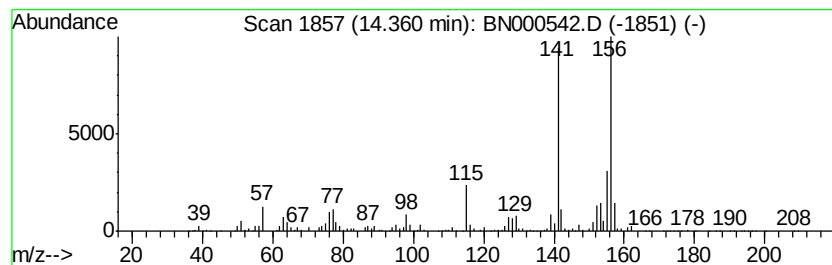
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 48 Naphthalene, 1,8-dimethyl- Concentration Rank 40

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 14.36 | 18.69 ng/ul | 4389590 | Acenaphthene-d10 | 15.04 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 1,8-dimethyl- | 156 | C12H12  | 000569-41-5 | 95   |
| 2    |    | Naphthalene, 1,3-dimethyl- | 156 | C12H12  | 000575-41-7 | 95   |
| 3    |    | Naphthalene, 2,6-dimethyl- | 156 | C12H12  | 000581-42-0 | 94   |
| 4    |    | Naphthalene, 1,3-dimethyl- | 156 | C12H12  | 000575-41-7 | 94   |
| 5    |    | Naphthalene, 1,7-dimethyl- | 156 | C12H12  | 000575-37-1 | 94   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000542.D  
Acq On : 22 Mar 2018 16:12  
Operator : JU/SJ  
Sample : J1905-11DL 10X  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM46DL

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

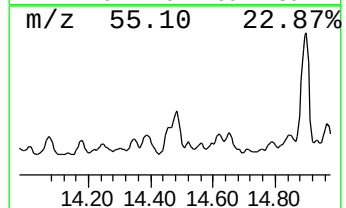
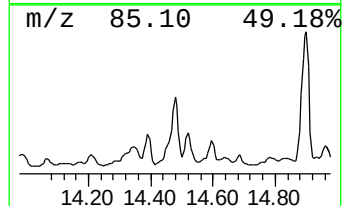
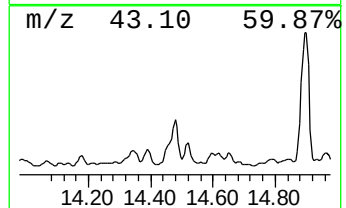
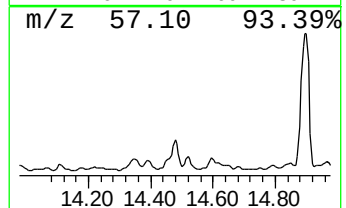
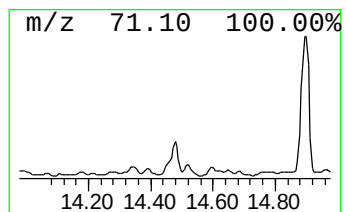
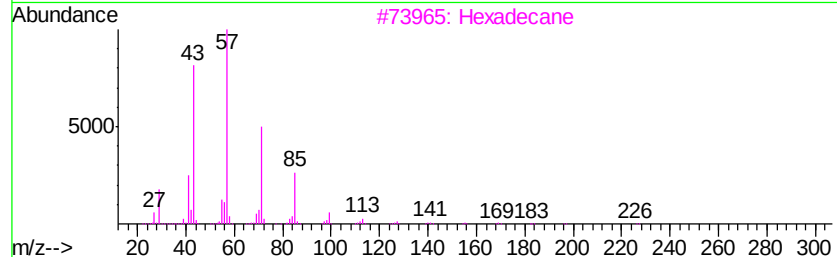
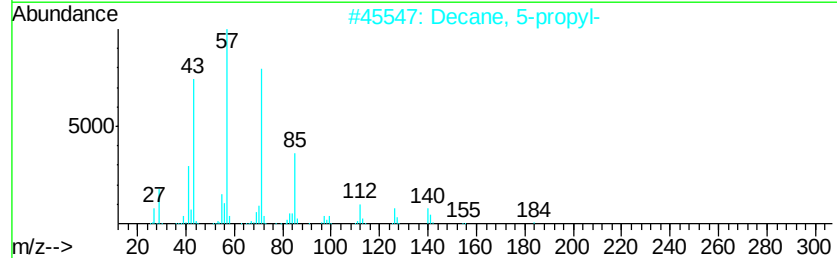
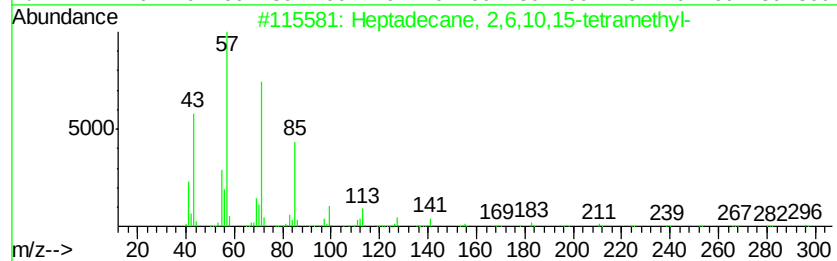
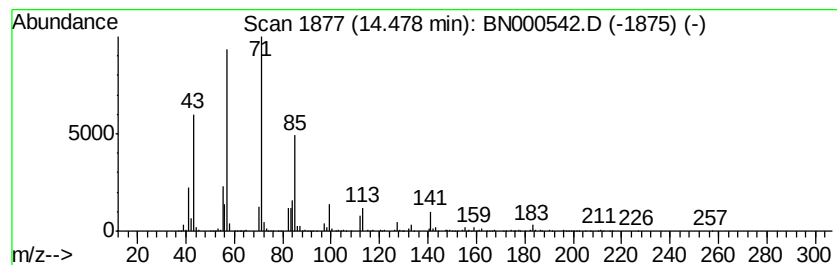
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 49 (DEL) Alkane: Straight-Chai... Concentration Rank 75

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T.  |
|-------|------------|---------|------------------|-------|
| 14.48 | 9.25 ng/ul | 2173950 | Acenaphthene-d10 | 15.04 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Heptadecane, 2,6,10,15-tetramethyl- | 296 | C21H44  | 054833-48-6 | 86   |
| 2    |    |   | Decane, 5-propyl-                   | 184 | C13H28  | 017312-62-8 | 80   |
| 3    |    |   | Hexadecane                          | 226 | C16H34  | 000544-76-3 | 64   |
| 4    |    |   | Tetradecane                         | 198 | C14H30  | 000629-59-4 | 59   |
| 5    |    |   | Undecane, 4,6-dimethyl-             | 184 | C13H28  | 017312-82-2 | 58   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000542.D  
Acq On : 22 Mar 2018 16:12  
Operator : JU/SJ  
Sample : J1905-11DL 10X  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM46DL

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

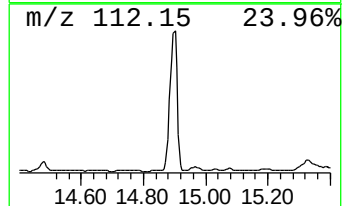
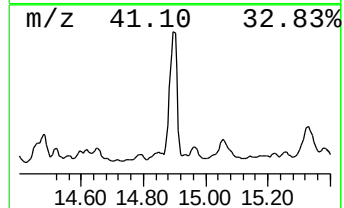
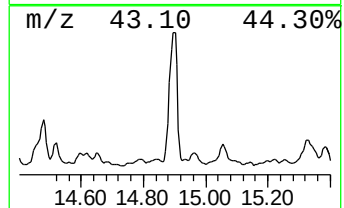
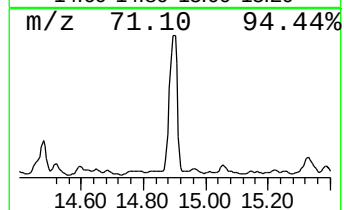
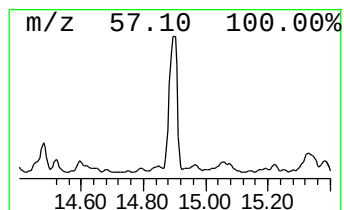
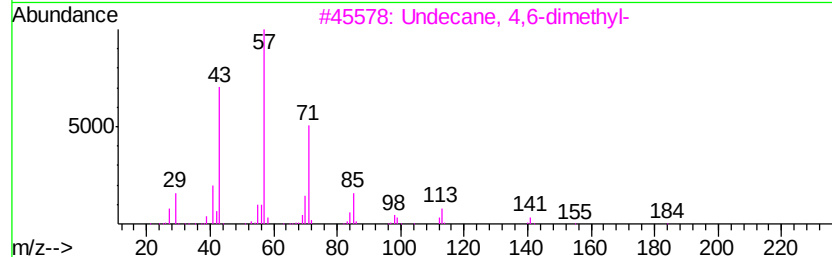
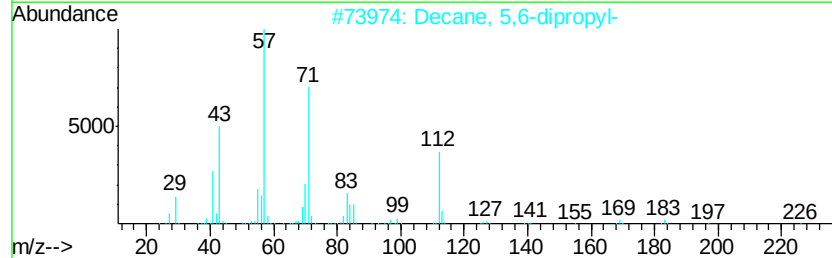
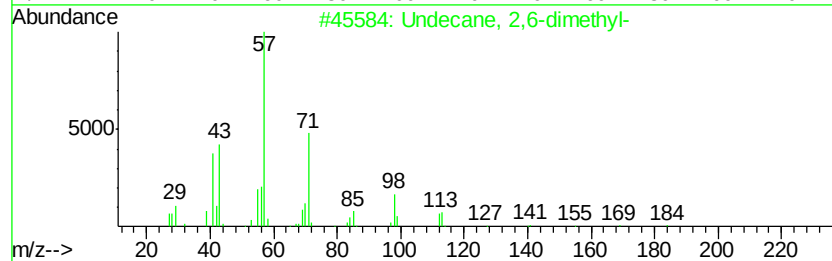
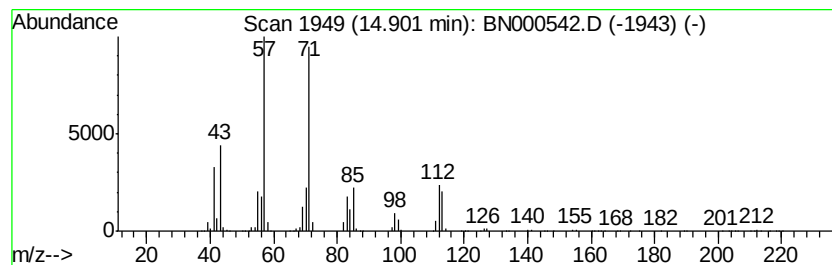
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 51 (DEL) Alkane: Straight-Chai... Concentration Rank 5

| R.T.  | EstConc     | Area     | Relative to ISTD | R.T.  |
|-------|-------------|----------|------------------|-------|
| 14.90 | 51.34 ng/ul | 12058600 | Acenaphthene-d10 | 15.04 |

| Hit# of 5 | Tentative ID             | MW  | MolForm | CAS#        | Qual |
|-----------|--------------------------|-----|---------|-------------|------|
| 1         | Undecane, 2,6-dimethyl-  | 184 | C13H28  | 017301-23-4 | 72   |
| 2         | Decane, 5,6-dipropyl-    | 226 | C16H34  | 119209-20-0 | 64   |
| 3         | Undecane, 4,6-dimethyl-  | 184 | C13H28  | 017312-82-2 | 59   |
| 4         | Octane, 2,3,6-trimethyl- | 156 | C11H24  | 062016-33-5 | 59   |
| 5         | Tridecane, 7-methyl-     | 198 | C14H30  | 026730-14-3 | 53   |





Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000542.D  
Acq On : 22 Mar 2018 16:12  
Operator : JU/SJ  
Sample : J1905-11DL 10X  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM46DL

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

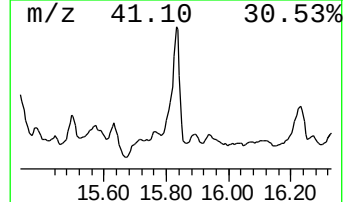
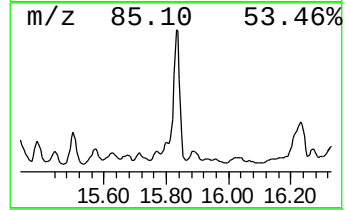
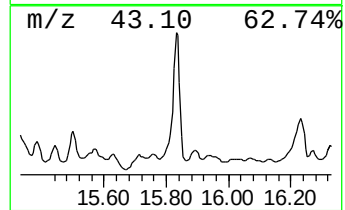
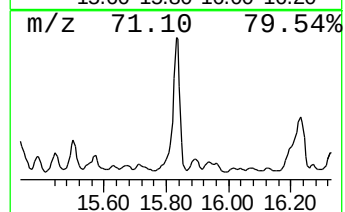
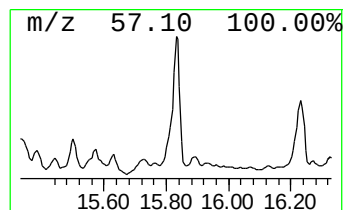
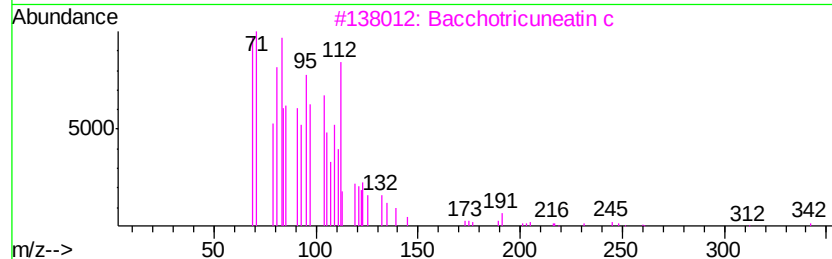
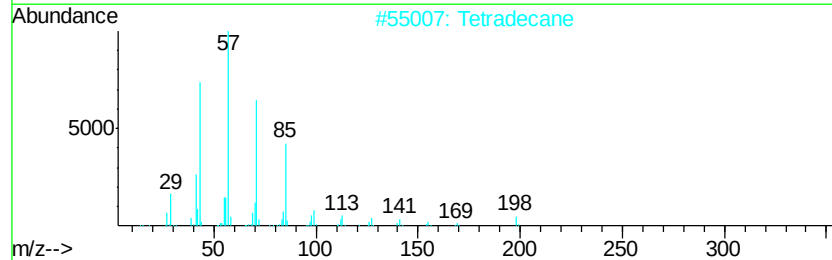
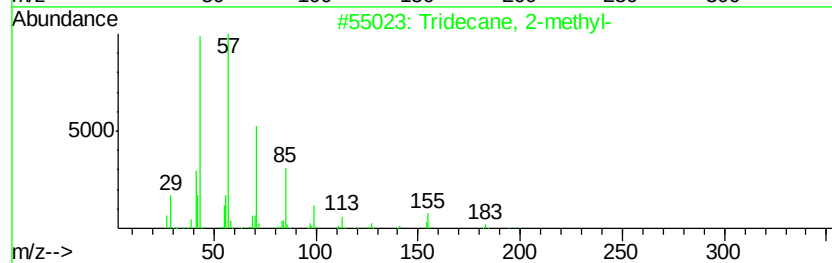
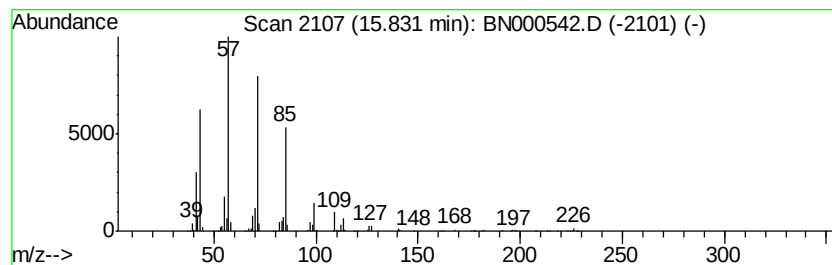
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 58 (DEL) Alkane: Straight-Chai... Concentration Rank 9

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 15.83 | 37.99 ng/ul | 8923940 | Acenaphthene-d10 | 15.04 |

| Hit# | of | Tentative ID             | MW  | MolForm  | CAS#        | Qual |
|------|----|--------------------------|-----|----------|-------------|------|
| 1    | 5  | Tridecane, 2-methyl-     | 198 | C14H30   | 001560-96-9 | 87   |
| 2    |    | Tetradecane              | 198 | C14H30   | 000629-59-4 | 86   |
| 3    |    | Bacchotricuneatin c      | 342 | C20H22O5 | 066563-30-2 | 80   |
| 4    |    | Decane, 2,3,5-trimethyl- | 184 | C13H28   | 062238-11-3 | 78   |
| 5    |    | Heptadecane              | 240 | C17H36   | 000629-78-7 | 78   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000542.D  
Acq On : 22 Mar 2018 16:12  
Operator : JU/SJ  
Sample : J1905-11DL 10X  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM46DL

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

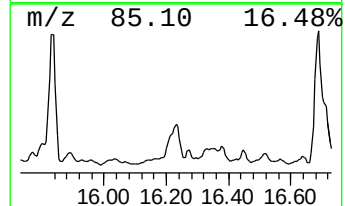
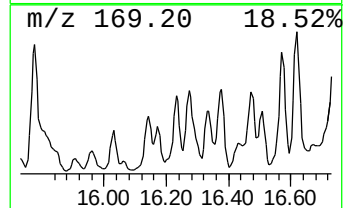
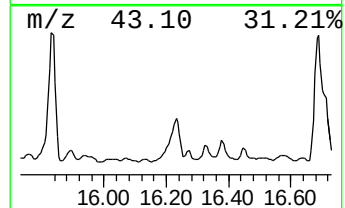
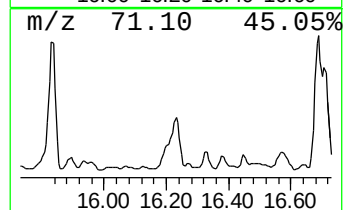
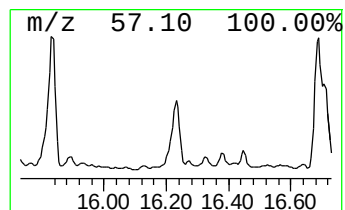
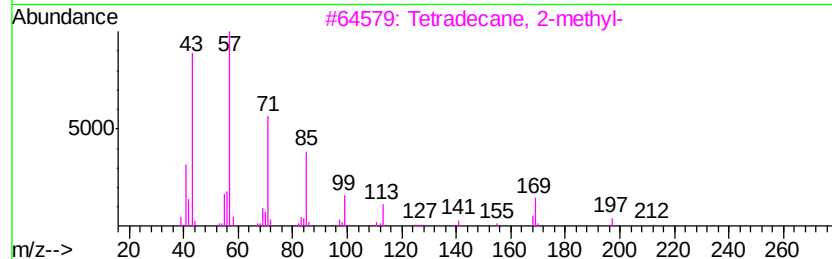
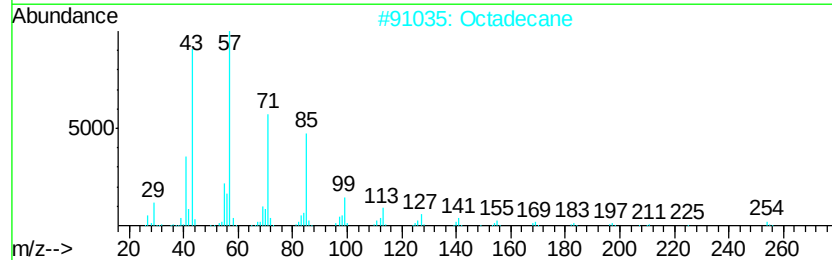
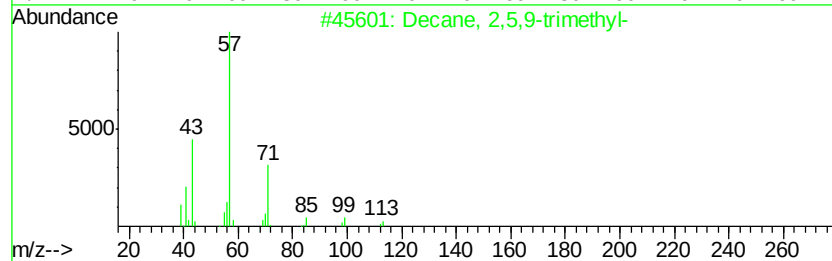
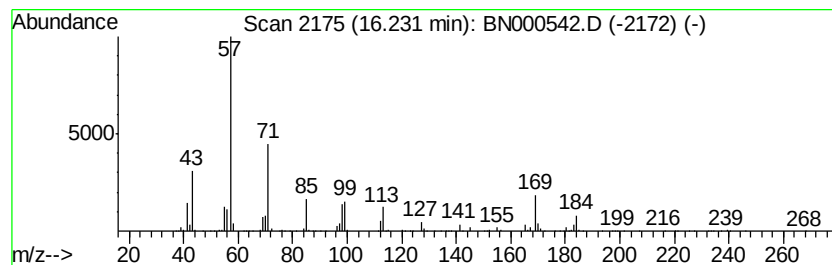
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 59 (DEL) Alkane: Straight-Chai... Concentration Rank 69

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 16.23 | 10.58 ng/ul | 2485830 | Acenaphthene-d10 | 15.04 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Decane, 2,5,9-trimethyl-            | 184 | C13H28  | 062108-22-9 | 59   |
| 2    |    | Octadecane                          | 254 | C18H38  | 000593-45-3 | 53   |
| 3    |    | Tetradecane, 2-methyl-              | 212 | C15H32  | 001560-95-8 | 50   |
| 4    |    | Octane, 2,6-dimethyl-               | 142 | C10H22  | 002051-30-1 | 47   |
| 5    |    | Pentadecane, 2,6,10,14-tetramethyl- | 268 | C19H40  | 001921-70-6 | 47   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000542.D  
Acq On : 22 Mar 2018 16:12  
Operator : JU/SJ  
Sample : J1905-11DL 10X  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM46DL

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

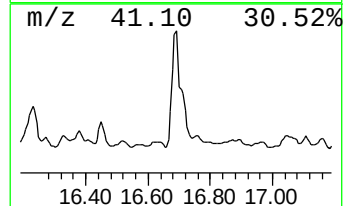
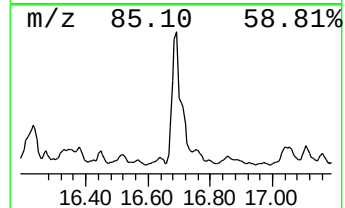
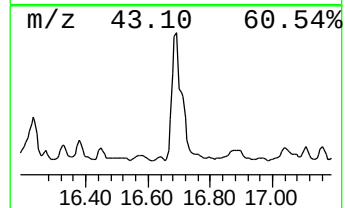
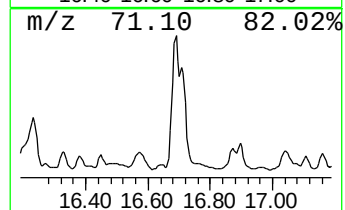
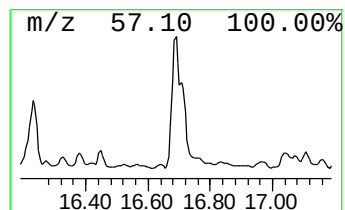
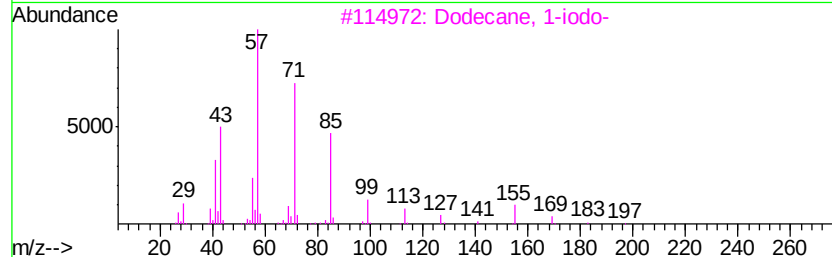
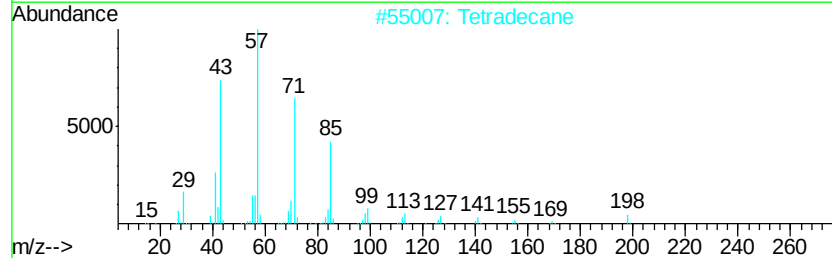
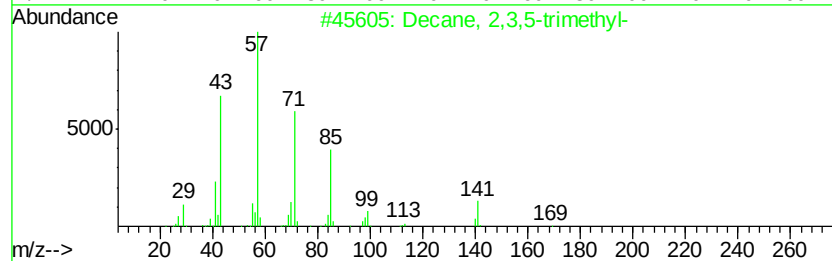
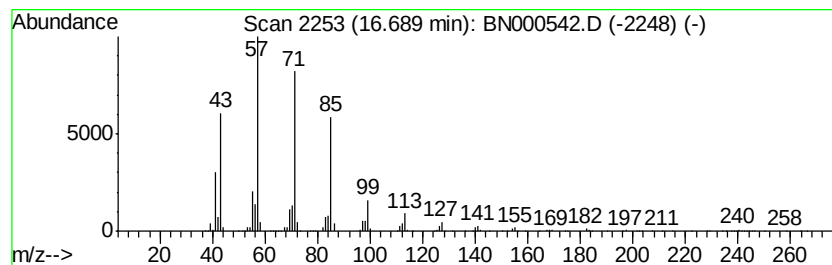
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 62 (DEL) Alkane: Straight-Chai... Concentration Rank 6

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 16.69 | 43.78 ng/ul | 6666410 | Phenanthrene-d10 | 17.78 |

| Hit# | of | Tentative ID             | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------|-----|---------|-------------|------|
| 1    | 5  | Decane, 2,3,5-trimethyl- | 184 | C13H28  | 062238-11-3 | 86   |
| 2    |    | Tetradecane              | 198 | C14H30  | 000629-59-4 | 83   |
| 3    |    | Dodecane, 1-iodo-        | 296 | C12H25I | 004292-19-7 | 78   |
| 4    |    | Hexadecane, 7-methyl-    | 240 | C17H36  | 026730-20-1 | 72   |
| 5    |    | Heptadecane              | 240 | C17H36  | 000629-78-7 | 68   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000542.D  
Acq On : 22 Mar 2018 16:12  
Operator : JU/SJ  
Sample : J1905-11DL 10X  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM46DL

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

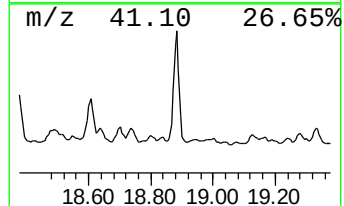
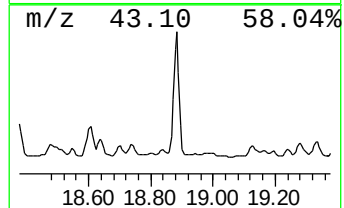
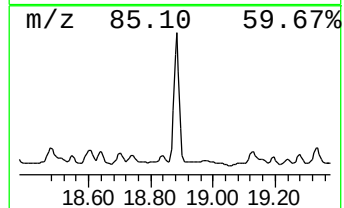
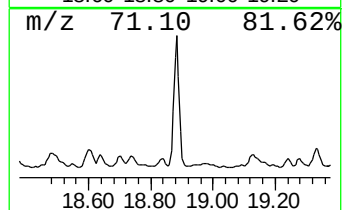
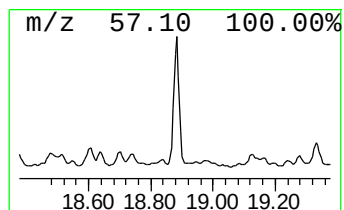
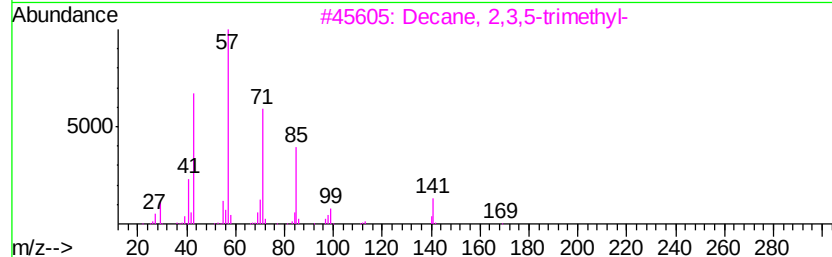
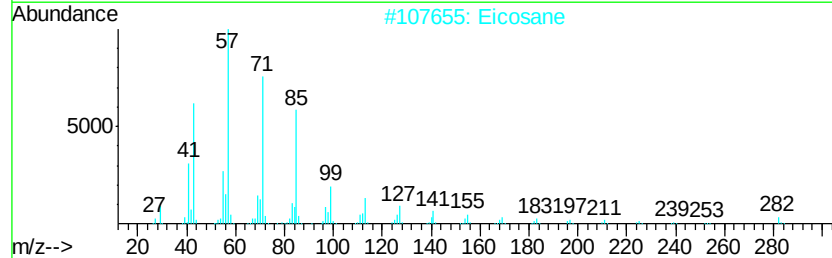
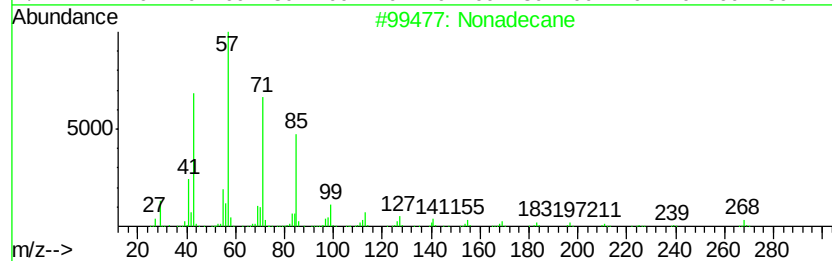
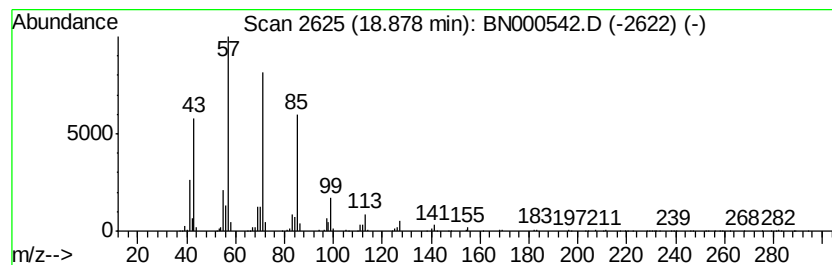
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 66 (DEL) Alkane: Straight-Chai... Concentration Rank 22

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 18.88 | 24.85 ng/ul | 3782980 | Phenanthrene-d10 | 17.78 |

| Hit# | of | Tentative ID             | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------|-----|---------|-------------|------|
| 1    | 5  | Nonadecane               | 268 | C19H40  | 000629-92-5 | 91   |
| 2    |    | Eicosane                 | 282 | C20H42  | 000112-95-8 | 90   |
| 3    |    | Decane, 2,3,5-trimethyl- | 184 | C13H28  | 062238-11-3 | 86   |
| 4    |    | Heptacosane              | 380 | C27H56  | 000593-49-7 | 72   |
| 5    |    | Heptadecane              | 240 | C17H36  | 000629-78-7 | 68   |



Data Path : Z:\HPCHEM1\BNA\_N\DATA\BN032218\  
Data File : BN000542.D  
Acq On : 22 Mar 2018 16:12  
Operator : JU/SJ  
Sample : J1905-11DL 10X  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM46DL

Quant Method : Z:\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

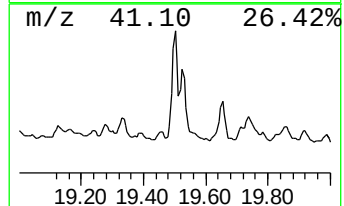
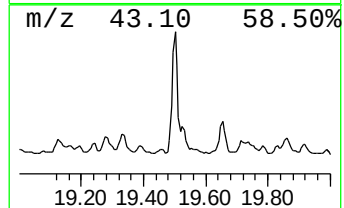
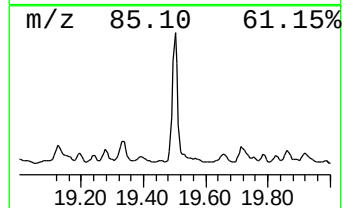
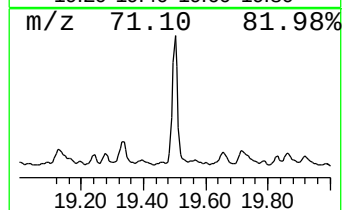
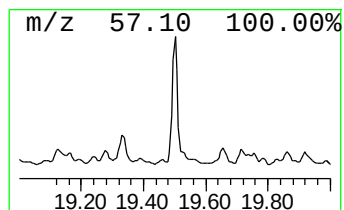
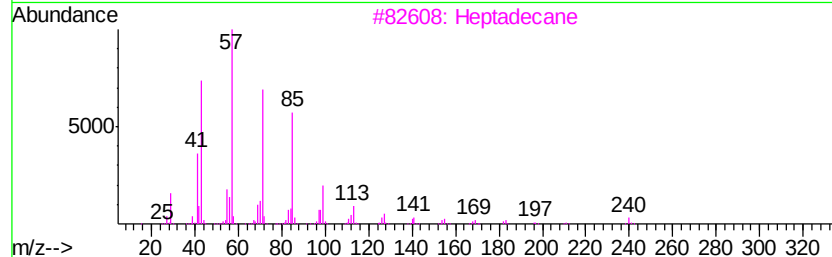
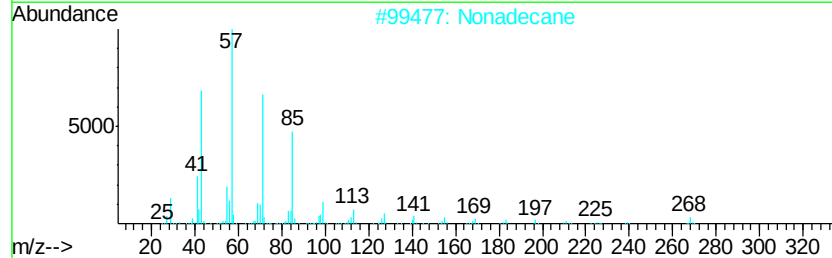
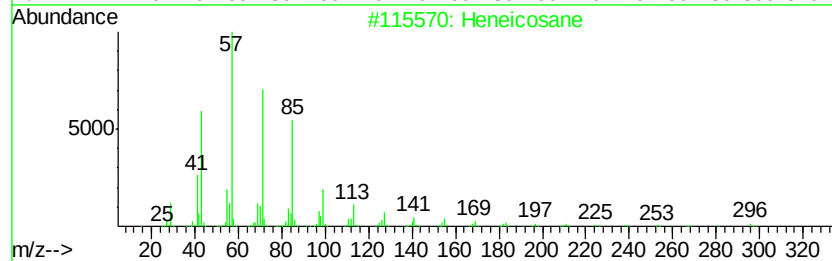
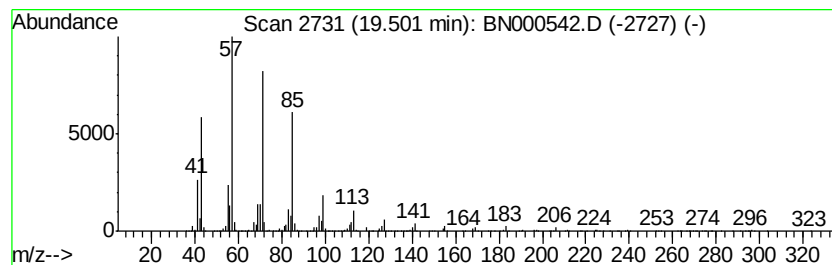
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 67 (DEL) Alkane: Straight-Chai... Concentration Rank 24

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 19.50 | 23.54 ng/ul | 3583700 | Phenanthrene-d10 | 17.78 |

| Hit# | of | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|--------------|-----|---------|-------------|------|
| 1    | 5  | Heneicosane  | 296 | C21H44  | 000629-94-7 | 91   |
| 2    |    | Nonadecane   | 268 | C19H40  | 000629-92-5 | 91   |
| 3    |    | Heptadecane  | 240 | C17H36  | 000629-78-7 | 91   |
| 4    |    | Octadecane   | 254 | C18H38  | 000593-45-3 | 90   |
| 5    |    | Hexadecane   | 226 | C16H34  | 000544-76-3 | 90   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000542.D  
Acq On : 22 Mar 2018 16:12  
Operator : JU/SJ  
Sample : J1905-11DL 10X  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM46DL

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

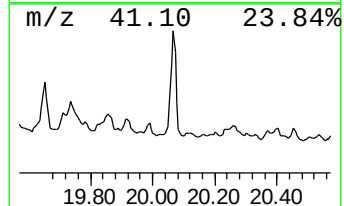
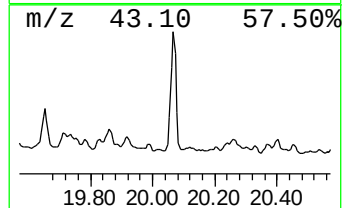
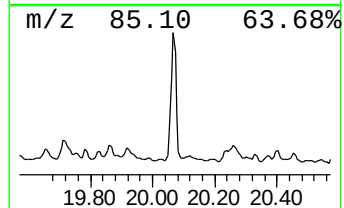
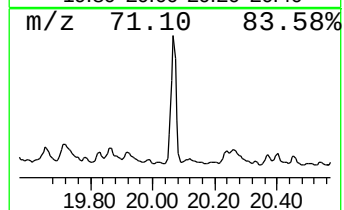
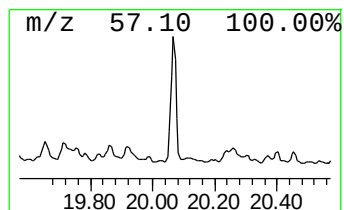
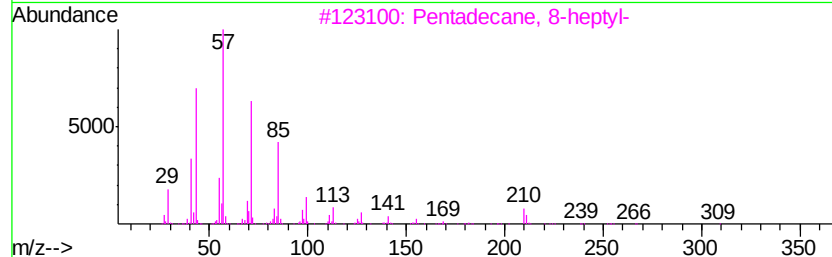
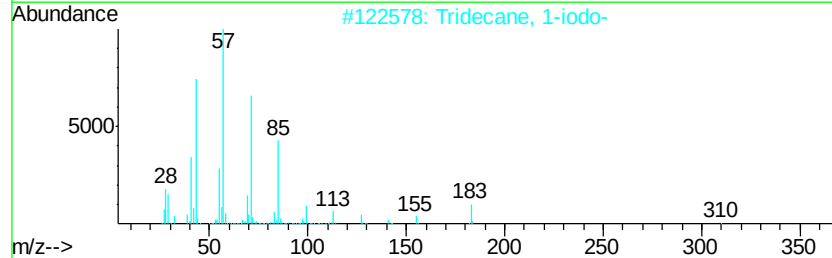
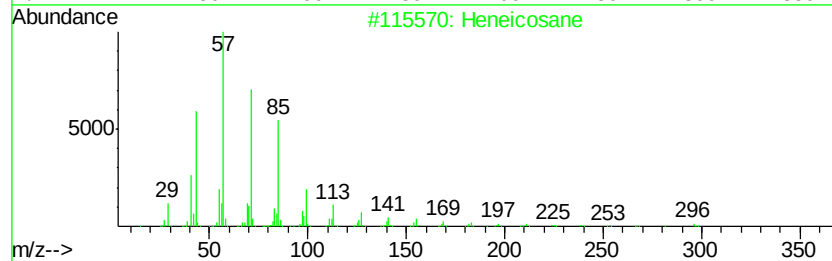
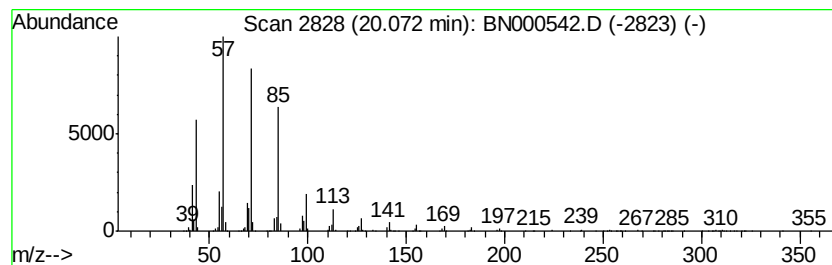
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 70 (DEL) Alkane: Straight-Chai... Concentration Rank 31

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 20.07 | 20.59 ng/ul | 2629860 | Chrysene-d12     | 21.88 |

| Hit# | of | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------|-----|---------|-------------|------|
| 1    | 5  | Heneicosane            | 296 | C21H44  | 000629-94-7 | 91   |
| 2    |    | Tridecane, 1-iodo-     | 310 | C13H27I | 035599-77-0 | 91   |
| 3    |    | Pentadecane, 8-heptyl- | 310 | C22H46  | 071005-15-7 | 91   |
| 4    |    | Heptacosane            | 380 | C27H56  | 000593-49-7 | 90   |
| 5    |    | Hexadecane             | 226 | C16H34  | 000544-76-3 | 90   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000542.D  
Acq On : 22 Mar 2018 16:12  
Operator : JU/SJ  
Sample : J1905-11DL 10X  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM46DL

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

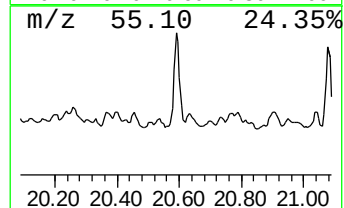
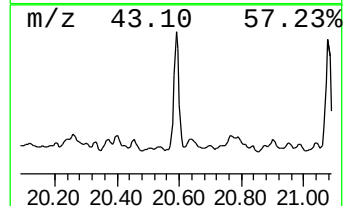
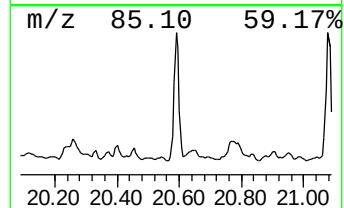
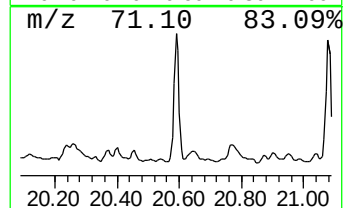
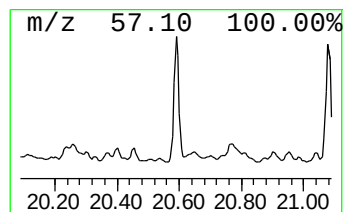
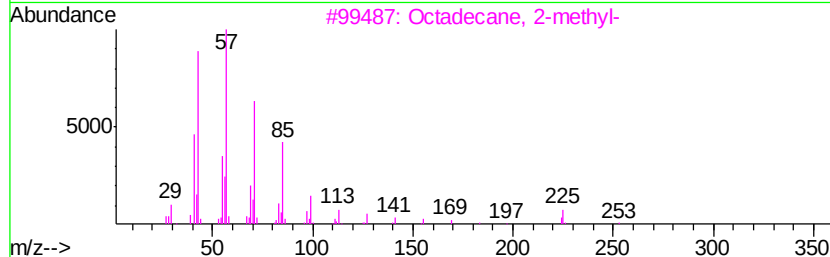
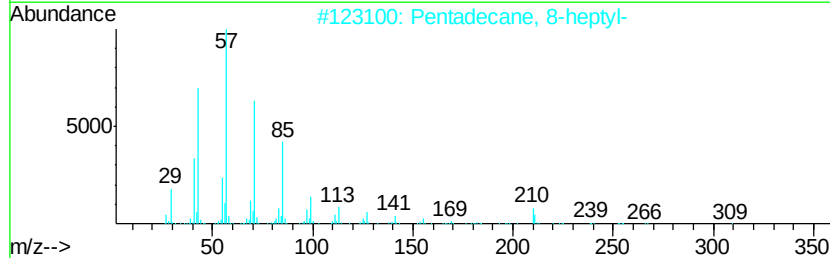
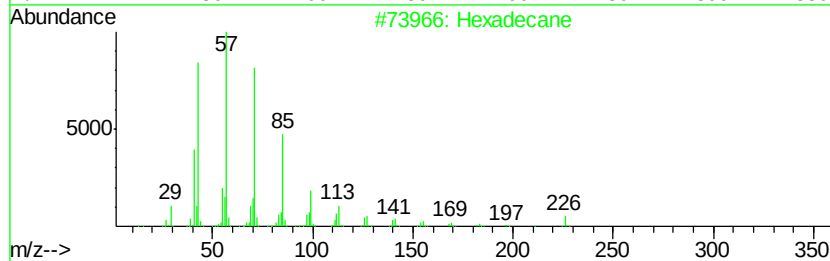
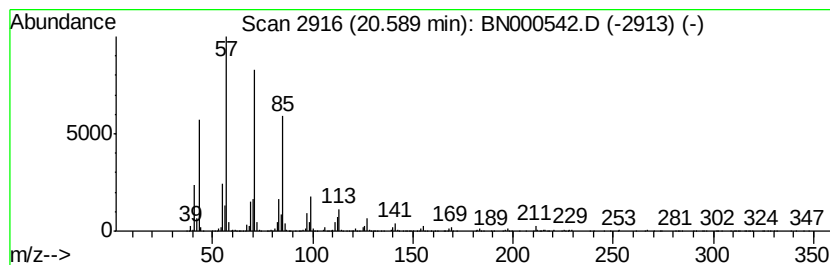
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 71 (DEL) Alkane: Straight-Chai... Concentration Rank 35

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 20.59 | 19.33 ng/ul | 2469170 | Chrysene-d12     | 21.88 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Hexadecane                          | 226 | C16H34  | 000544-76-3 | 91   |
| 2    |    | Pentadecane, 8-heptyl-              | 310 | C22H46  | 071005-15-7 | 90   |
| 3    |    | Octadecane, 2-methyl-               | 268 | C19H40  | 001560-88-9 | 90   |
| 4    |    | Tetracosane                         | 338 | C24H50  | 000646-31-1 | 90   |
| 5    |    | Heptadecane, 2,6,10,15-tetramethyl- | 296 | C21H44  | 054833-48-6 | 90   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000542.D  
Acq On : 22 Mar 2018 16:12  
Operator : JU/SJ  
Sample : J1905-11DL 10X  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM46DL

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

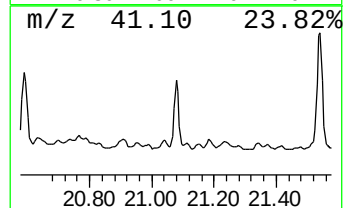
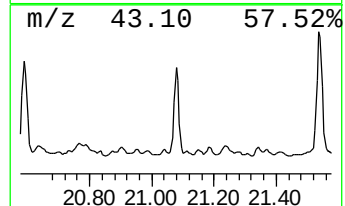
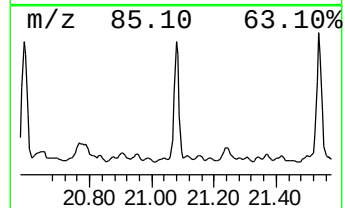
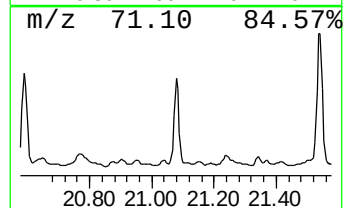
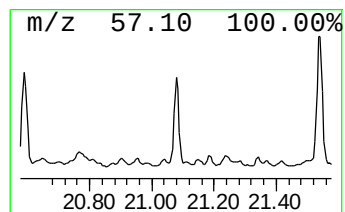
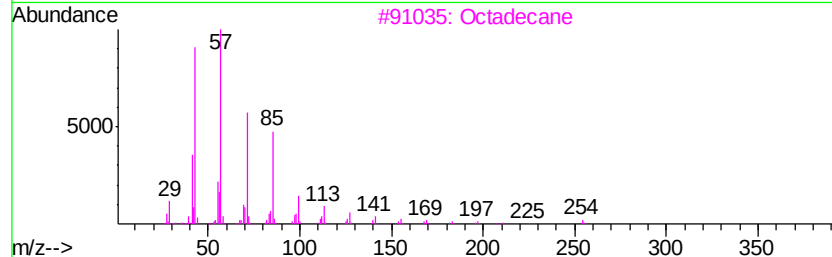
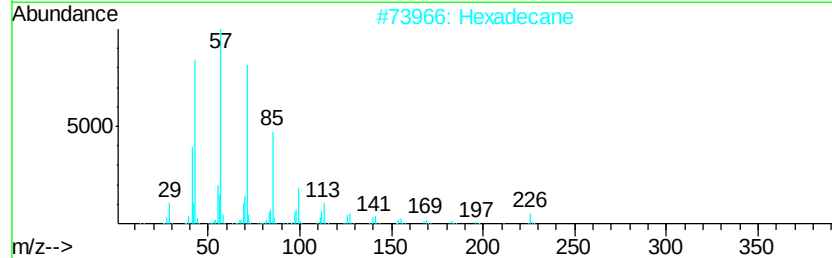
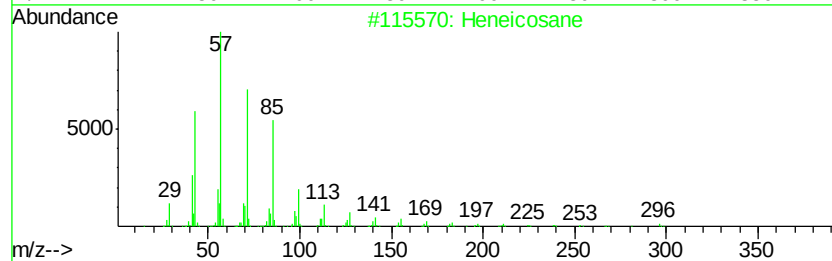
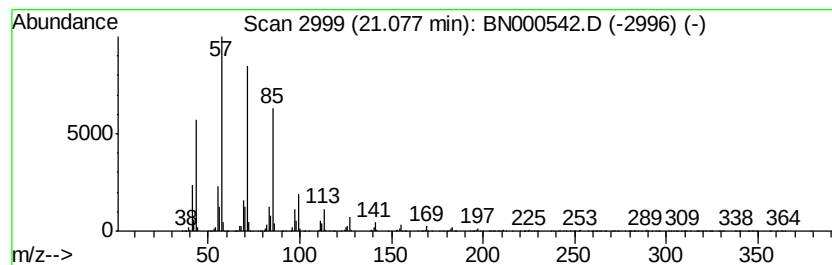
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 72 (DEL) Alkane: Straight-Chai... Concentration Rank 50

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 21.08 | 14.94 ng/ul | 1908760 | Chrysene-d12     | 21.88 |

| Hit# | of | Tentative ID         | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------|-----|---------|-------------|------|
| 1    | 5  | Heneicosane          | 296 | C21H44  | 000629-94-7 | 91   |
| 2    |    | Hexadecane           | 226 | C16H34  | 000544-76-3 | 90   |
| 3    |    | Octadecane           | 254 | C18H38  | 000593-45-3 | 90   |
| 4    |    | Octacosane           | 394 | C28H58  | 000630-02-4 | 87   |
| 5    |    | Eicosane, 10-methyl- | 296 | C21H44  | 054833-23-7 | 87   |





Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000542.D  
Acq On : 22 Mar 2018 16:12  
Operator : JU/SJ  
Sample : J1905-11DL 10X  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM46DL

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

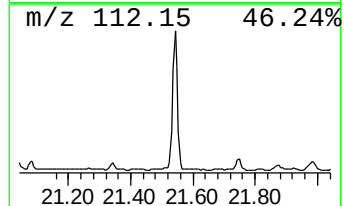
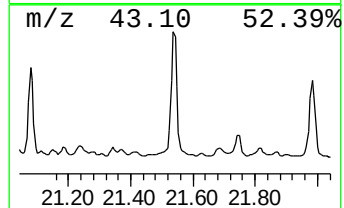
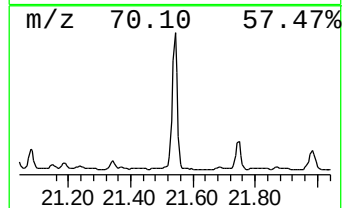
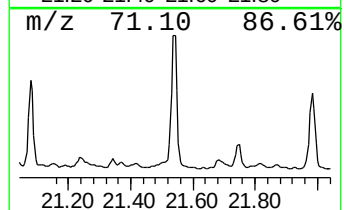
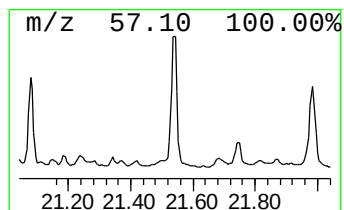
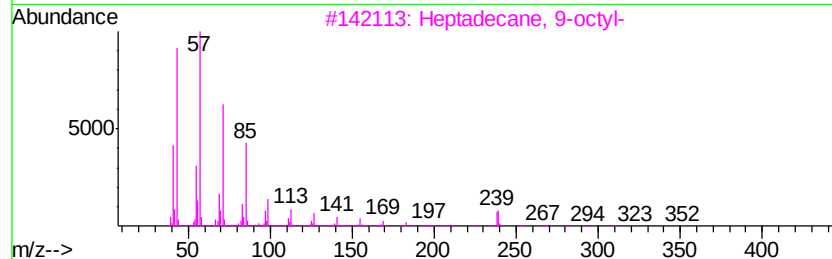
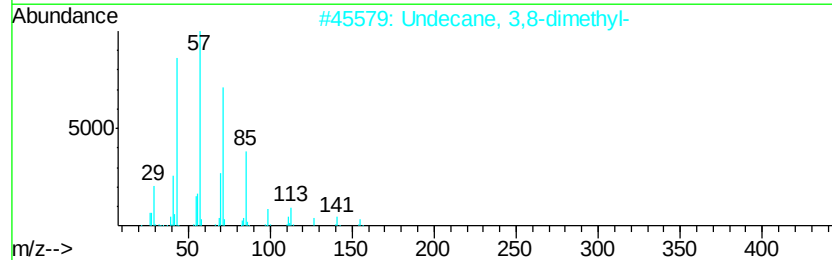
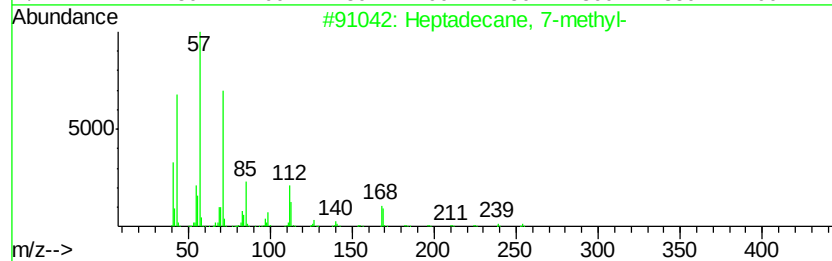
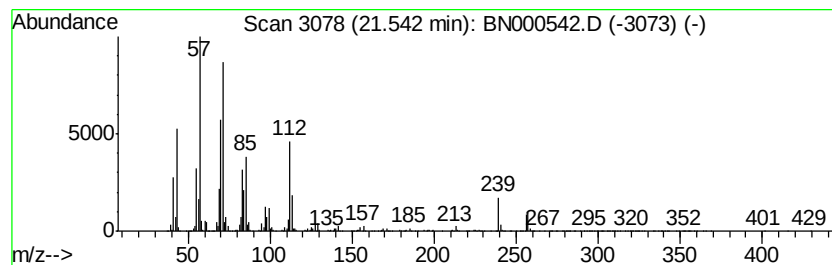
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 73 (DEL) Alkane: Straight-Chai... Concentration Rank 8

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 21.54 | 38.40 ng/ul | 4905200 | Chrysene-d12     | 21.88 |

| Hit# | of | Tentative ID             | MW  | MolForm  | CAS#        | Qual |
|------|----|--------------------------|-----|----------|-------------|------|
| 1    | 5  | Heptadecane, 7-methyl-   | 254 | C18H38   | 020959-33-5 | 58   |
| 2    |    | Undecane, 3,8-dimethyl-  | 184 | C13H28   | 017301-30-3 | 47   |
| 3    |    | Heptadecane, 9-octyl-    | 352 | C25H52   | 007225-64-1 | 41   |
| 4    |    | 2,4,4-Trimethyl-1-hexene | 126 | C9H18    | 051174-12-0 | 38   |
| 5    |    | Hexadecane, 1-chloro-    | 260 | C16H33Cl | 004860-03-1 | 38   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000542.D  
Acq On : 22 Mar 2018 16:12  
Operator : JU/SJ  
Sample : J1905-11DL 10X  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleID :  
DAM46DL

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

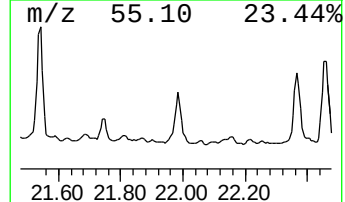
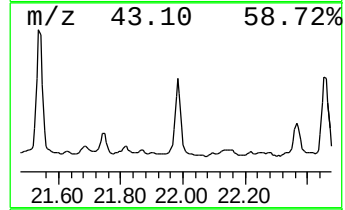
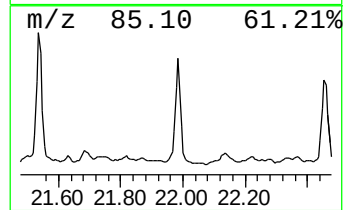
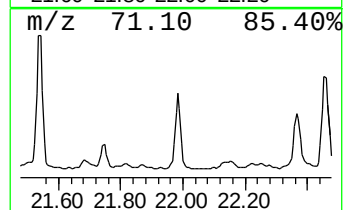
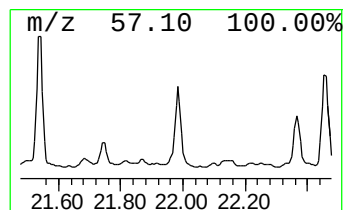
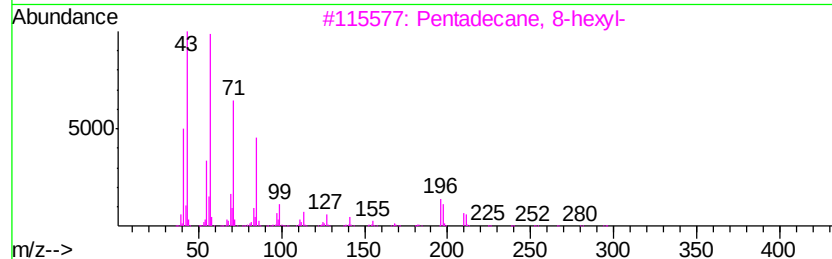
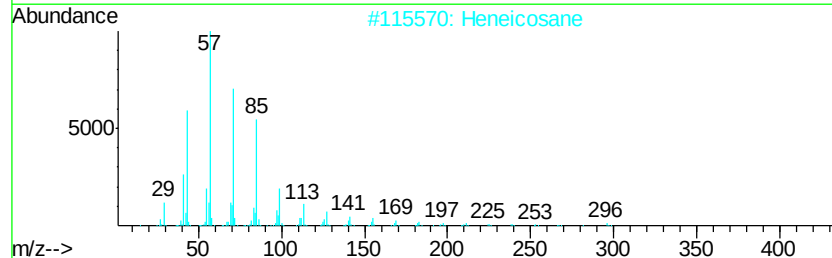
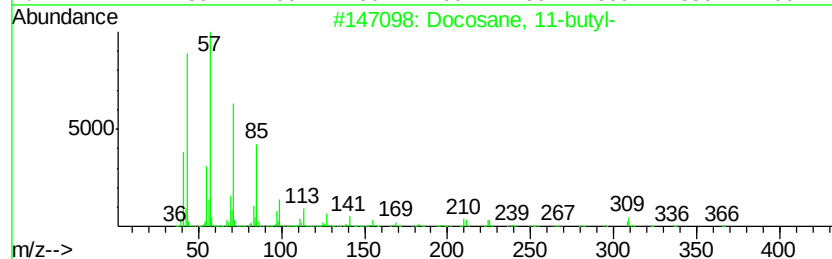
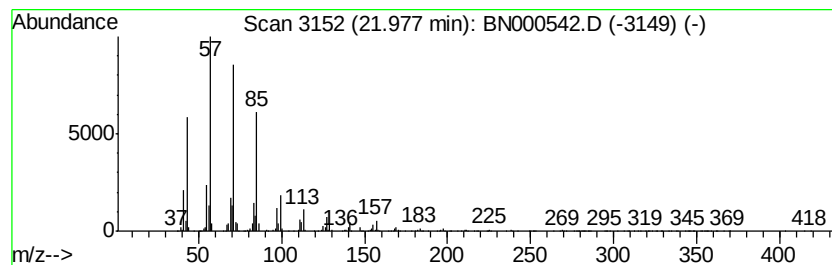
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 74 (DEL) Alkane: Cyclic21.98 Concentration Rank 37

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 21.98 | 19.24 ng/ul | 2457330 | Chrysene-d12     | 21.88 |

| Hit# | of | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------|-----|---------|-------------|------|
| 1    | 5  | Docosane, 11-butyl-    | 366 | C26H54  | 013475-76-8 | 93   |
| 2    |    | Heneicosane            | 296 | C21H44  | 000629-94-7 | 91   |
| 3    |    | Pentadecane, 8-hexyl-  | 296 | C21H44  | 013475-75-7 | 90   |
| 4    |    | Eicosane               | 282 | C20H42  | 000112-95-8 | 90   |
| 5    |    | Pentadecane, 8-heptyl- | 310 | C22H46  | 071005-15-7 | 86   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000542.D  
Acq On : 22 Mar 2018 16:12  
Operator : JU/SJ  
Sample : J1905-11DL 10X  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM46DL

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

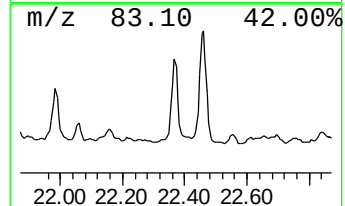
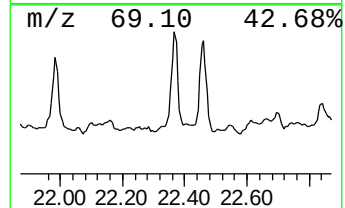
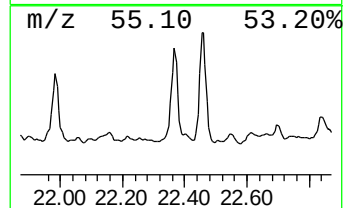
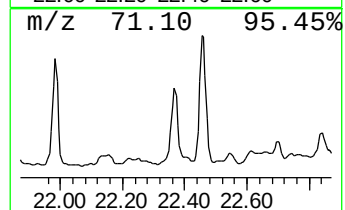
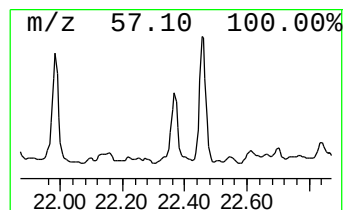
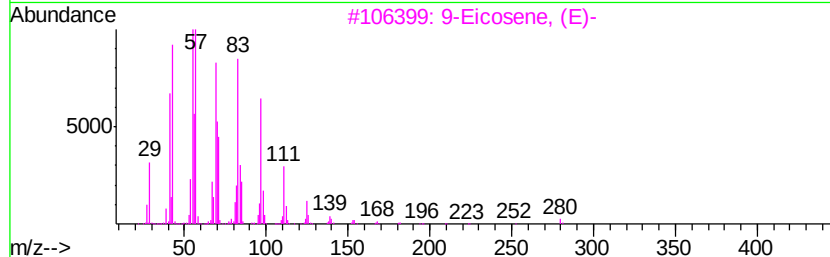
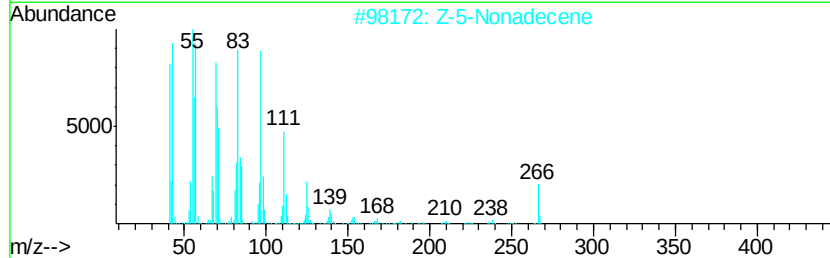
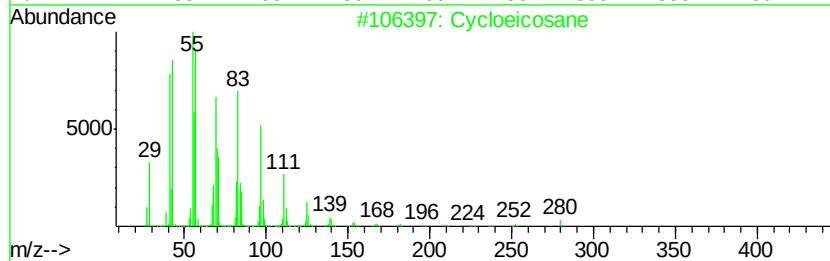
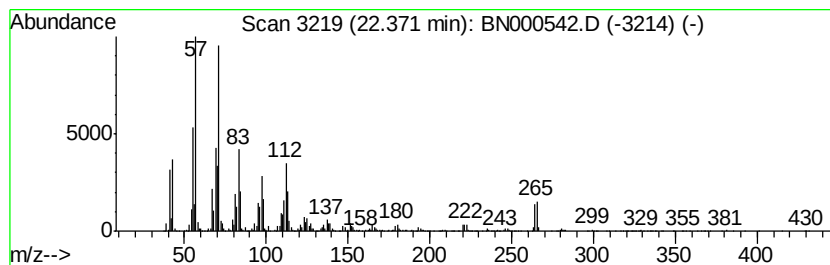
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 75 (DEL) Alkane: Cyclic22.37 Concentration Rank 43

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 22.37 | 17.76 ng/ul | 2269110 | Chrysene-d12     | 21.88 |

| Hit# | of | Tentative ID     | MW  | MolForm | CAS#         | Qual |
|------|----|------------------|-----|---------|--------------|------|
| 1    | 5  | Cycloeicosane    | 280 | C20H40  | 000296-56-0  | 42   |
| 2    |    | Z-5-Nonadecene   | 266 | C19H38  | 1000131-11-8 | 42   |
| 3    |    | 9-Eicosene, (E)- | 280 | C20H40  | 074685-29-3  | 42   |
| 4    |    | 3-Eicosene, (E)- | 280 | C20H40  | 074685-33-9  | 41   |
| 5    |    | 1-Nonadecene     | 266 | C19H38  | 018435-45-5  | 38   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000542.D  
Acq On : 22 Mar 2018 16:12  
Operator : JU/SJ  
Sample : J1905-11DL 10X  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM46DL

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

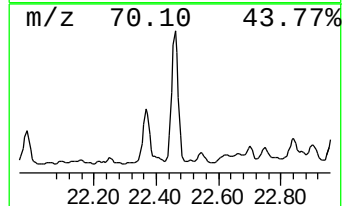
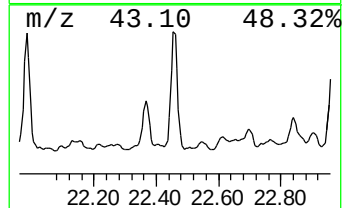
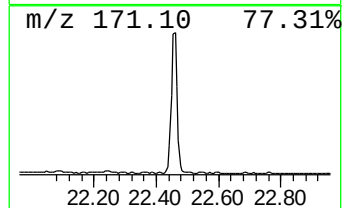
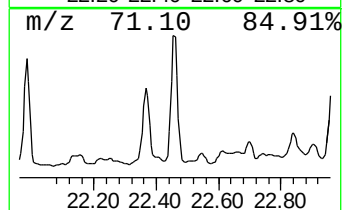
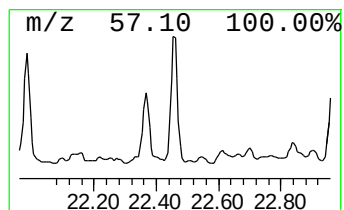
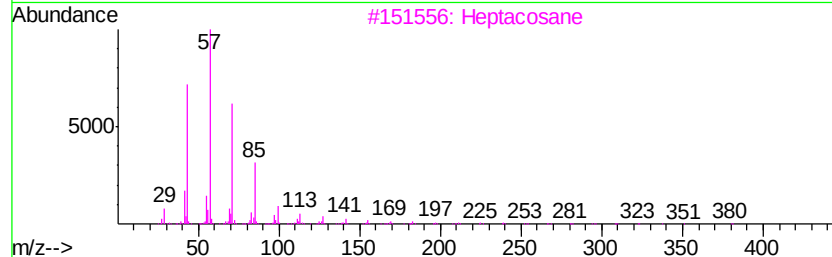
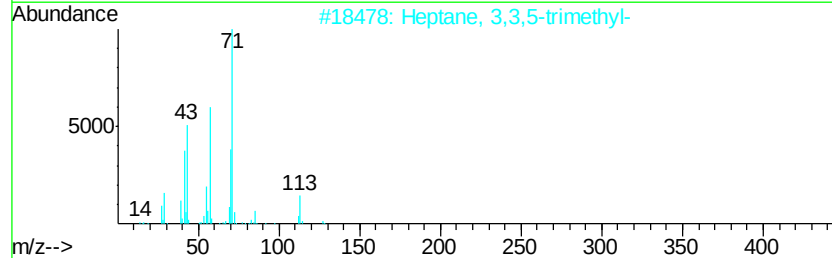
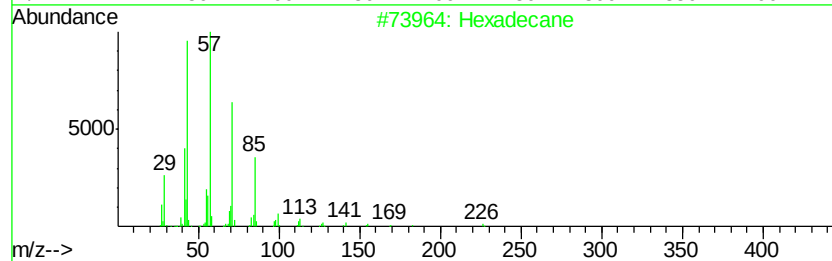
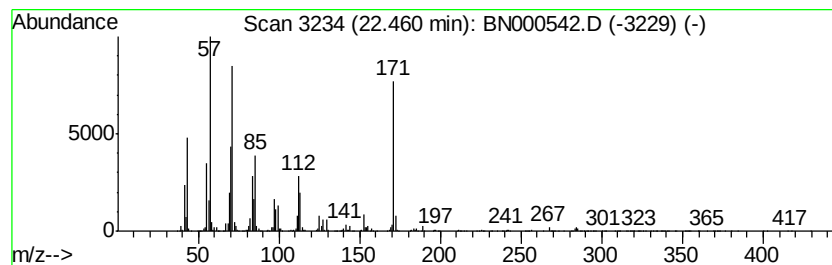
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 76 (DEL) Alkane: Straight-Chai... Concentration Rank 11

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 22.46 | 32.79 ng/ul | 4188120 | Chrysene-d12     | 21.88 |

| Hit# | of | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------------|-----|---------|-------------|------|
| 1    | 5  | Hexadecane                | 226 | C16H34  | 000544-76-3 | 38   |
| 2    |    | Heptane, 3,3,5-trimethyl- | 142 | C10H22  | 007154-80-5 | 35   |
| 3    |    | Heptacosane               | 380 | C27H56  | 000593-49-7 | 35   |
| 4    |    | Pentadecane               | 212 | C15H32  | 000629-62-9 | 35   |
| 5    |    | Hexadecane                | 226 | C16H34  | 000544-76-3 | 35   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000542.D  
Acq On : 22 Mar 2018 16:12  
Operator : JU/SJ  
Sample : J1905-11DL 10X  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM46DL

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

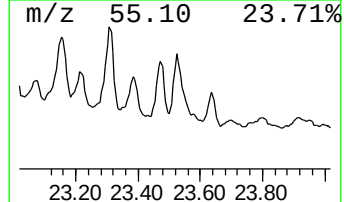
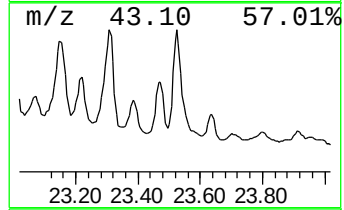
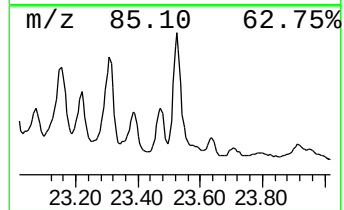
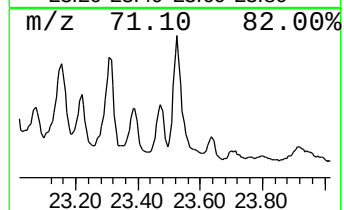
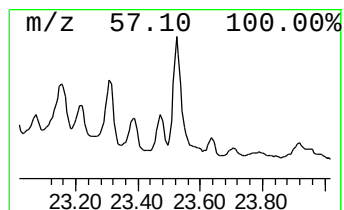
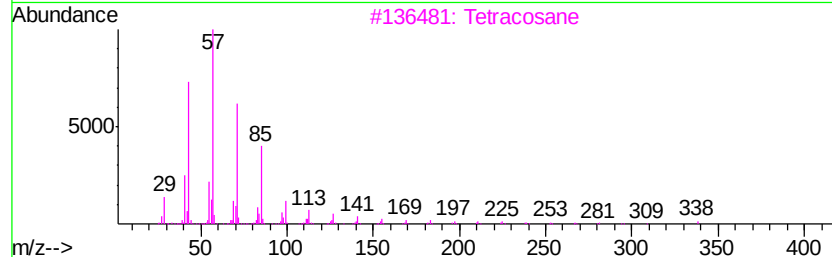
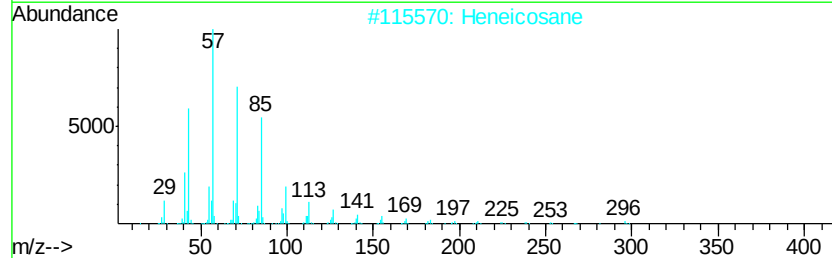
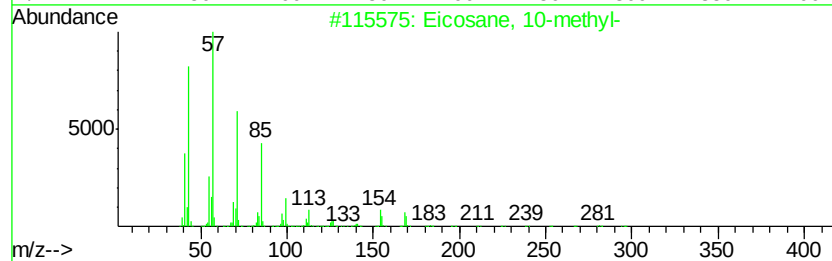
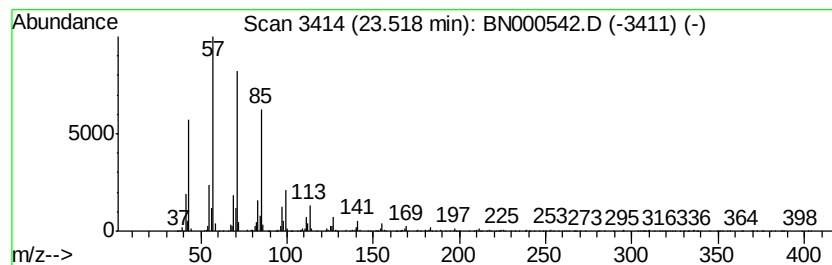
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 80 (DEL) Alkane: Straight-Chai... Concentration Rank 46

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 23.52 | 17.08 ng/ul | 1795260 | Perylene-d12     | 24.37 |

| Hit# | of | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------|-----|---------|-------------|------|
| 1    | 5  | Eicosane, 10-methyl-  | 296 | C21H44  | 054833-23-7 | 96   |
| 2    |    | Heneicosane           | 296 | C21H44  | 000629-94-7 | 96   |
| 3    |    | Tetracosane           | 338 | C24H50  | 000646-31-1 | 93   |
| 4    |    | Pentadecane, 8-hexyl- | 296 | C21H44  | 013475-75-7 | 91   |
| 5    |    | Eicosane              | 282 | C20H42  | 000112-95-8 | 87   |



Data Path : Z:\HPCHEM1\BNA\_N\DATA\BN032218\  
 Data File : BN000542.D  
 Acq On : 22 Mar 2018 16:12  
 Operator : JU/SJ  
 Sample : J1905-11DL 10X  
 Misc :  
 ALS Vial : 12 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 DAM46DL

Quant Method : Z:\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN031918.M  
 Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
 TIC Integration Parameters: LSCINT.P

| TIC Top Hit name     | RT    | EstConc | Units | Response | --Internal Standard-- |       |         |      |
|----------------------|-------|---------|-------|----------|-----------------------|-------|---------|------|
|                      |       |         |       |          | #                     | RT    | Resp    | Conc |
| 1-Butanol, 2-ethyl-  | 5.41  | 17.0    | ng/ul | 2184440  | 1                     | 8.42  | 2572200 | 20.0 |
| (DEL) Alkane: Str... | 6.35  | 24.9    | ng/ul | 3204930  | 1                     | 8.42  | 2572200 | 20.0 |
| 1-Pentanol, 2-ethyl- | 6.86  | 27.5    | ng/ul | 3532920  | 1                     | 8.42  | 2572200 | 20.0 |
| (DEL) Alkane: Str... | 7.36  | 17.2    | ng/ul | 2210350  | 1                     | 8.42  | 2572200 | 20.0 |
| (DEL) Alkane: Str... | 7.42  | 18.6    | ng/ul | 2393250  | 1                     | 8.42  | 2572200 | 20.0 |
| Benzene, 1-ethyl-... | 7.47  | 18.8    | ng/ul | 2412110  | 1                     | 8.42  | 2572200 | 20.0 |
| 2-Heptanone, 4,6-... | 7.81  | 21.5    | ng/ul | 2765720  | 1                     | 8.42  | 2572200 | 20.0 |
| (DEL) Alkane: Str... | 8.01  | 75.4    | ng/ul | 9701630  | 1                     | 8.42  | 2572200 | 20.0 |
| Benzene, 1,2,3-tr... | 8.05  | 22.3    | ng/ul | 2863830  | 1                     | 8.42  | 2572200 | 20.0 |
| (DEL) Alkane: Str... | 8.37  | 20.0    | ng/ul | 2572200  | 1                     | 8.42  | 2572200 | 20.0 |
| 1-Hexanol, 2-ethyl-  | 8.48  | 118.7   | ng/ul | 15259700 | 1                     | 8.42  | 2572200 | 20.0 |
| unknown-01           | 8.62  | 25.9    | ng/ul | 3332290  | 1                     | 8.42  | 2572200 | 20.0 |
| Cyclohexanone, 3,... | 8.86  | 52.3    | ng/ul | 6720170  | 1                     | 8.42  | 2572200 | 20.0 |
| (DEL) Alkane: Str... | 8.93  | 14.3    | ng/ul | 1834010  | 1                     | 8.42  | 2572200 | 20.0 |
| (DEL) Alkane: Str... | 8.98  | 28.7    | ng/ul | 3685810  | 1                     | 8.42  | 2572200 | 20.0 |
| Benzene, 1,4-diet... | 9.06  | 54.9    | ng/ul | 7066310  | 1                     | 8.42  | 2572200 | 20.0 |
| Benzene, 1-methyl... | 9.43  | 14.5    | ng/ul | 1870400  | 1                     | 8.42  | 2572200 | 20.0 |
| unknown-02           | 9.78  | 41.8    | ng/ul | 5369920  | 1                     | 8.42  | 2572200 | 20.0 |
| Benzene, 4-ethyl-... | 9.87  | 20.4    | ng/ul | 4425030  | 2                     | 11.26 | 4331650 | 20.0 |
| unknown-03           | 10.28 | 14.1    | ng/ul | 3054240  | 2                     | 11.26 | 4331650 | 20.0 |
| (DEL) Alkane: Str... | 10.49 | 9.6     | ng/ul | 2074480  | 2                     | 11.26 | 4331650 | 20.0 |
| (DEL) Alkane: Str... | 10.55 | 10.0    | ng/ul | 2159340  | 2                     | 11.26 | 4331650 | 20.0 |
| (DEL) Alkane: Str... | 10.64 | 14.9    | ng/ul | 3217240  | 2                     | 11.26 | 4331650 | 20.0 |
| unknown-04           | 10.75 | 22.3    | ng/ul | 4817900  | 2                     | 11.26 | 4331650 | 20.0 |
| unknown-05           | 11.00 | 8.7     | ng/ul | 1884370  | 2                     | 11.26 | 4331650 | 20.0 |
| unknown-06           | 11.12 | 11.2    | ng/ul | 2422090  | 2                     | 11.26 | 4331650 | 20.0 |
| (DEL) Alkane: Str... | 11.20 | 37.3    | ng/ul | 8069710  | 2                     | 11.26 | 4331650 | 20.0 |
| 1,3-Pentanediol, ... | 11.45 | 9.6     | ng/ul | 2077730  | 2                     | 11.26 | 4331650 | 20.0 |
| 1,3-Hexanediol, 2... | 11.54 | 11.8    | ng/ul | 2557830  | 2                     | 11.26 | 4331650 | 20.0 |
| 2,4-Dipropyl-5-et... | 11.60 | 28.6    | ng/ul | 6196920  | 2                     | 11.26 | 4331650 | 20.0 |
| (DEL) Alkane: Str... | 11.73 | 12.5    | ng/ul | 2700050  | 2                     | 11.26 | 4331650 | 20.0 |
| unknown-07           | 11.84 | 9.1     | ng/ul | 1973920  | 2                     | 11.26 | 4331650 | 20.0 |
| unknown-08           | 11.93 | 12.0    | ng/ul | 2606680  | 2                     | 11.26 | 4331650 | 20.0 |
| unknown-09           | 12.12 | 12.8    | ng/ul | 2770540  | 2                     | 11.26 | 4331650 | 20.0 |
| (DEL) Alkane: Str... | 12.23 | 14.9    | ng/ul | 3217370  | 2                     | 11.26 | 4331650 | 20.0 |
| Benzene, 1,3,5-tr... | 12.31 | 19.3    | ng/ul | 4170250  | 2                     | 11.26 | 4331650 | 20.0 |
| unknown-10           | 12.47 | 22.1    | ng/ul | 4798340  | 2                     | 11.26 | 4331650 | 20.0 |
| (DEL) Alkane: Str... | 12.61 | 27.9    | ng/ul | 6031380  | 2                     | 11.26 | 4331650 | 20.0 |
| (DEL) Alkane: Cyc... | 13.26 | 12.1    | ng/ul | 2843950  | 3                     | 15.04 | 4697940 | 20.0 |
| Propanoic acid, 2... | 13.77 | 7.6     | ng/ul | 1788060  | 3                     | 15.04 | 4697940 | 20.0 |
| Naphthalene, 1,8-... | 14.36 | 18.7    | ng/ul | 4389590  | 3                     | 15.04 | 4697940 | 20.0 |
| (DEL) Alkane: Str... | 14.48 | 9.3     | ng/ul | 2173950  | 3                     | 15.04 | 4697940 | 20.0 |
| (DEL) Alkane: Str... | 14.90 | 51.3    | ng/ul | 12058600 | 3                     | 15.04 | 4697940 | 20.0 |
| (DEL) Alkane: Str... | 15.83 | 38.0    | ng/ul | 8923940  | 3                     | 15.04 | 4697940 | 20.0 |
| (DEL) Alkane: Str... | 16.23 | 10.6    | ng/ul | 2485830  | 3                     | 15.04 | 4697940 | 20.0 |
| (DEL) Alkane: Str... | 16.69 | 43.8    | ng/ul | 6666410  | 4                     | 17.78 | 3045150 | 20.0 |
| (DEL) Alkane: Str... | 18.88 | 24.9    | ng/ul | 3782980  | 4                     | 17.78 | 3045150 | 20.0 |
| (DEL) Alkane: Str... | 19.50 | 23.5    | ng/ul | 3583700  | 4                     | 17.78 | 3045150 | 20.0 |
| (DEL) Alkane: Str... | 20.07 | 20.6    | ng/ul | 2629860  | 5                     | 21.88 | 2554850 | 20.0 |

|       |                |       |      |       |         |   |       |         |      |
|-------|----------------|-------|------|-------|---------|---|-------|---------|------|
| (DEL) | Alkane: Str... | 20.59 | 19.3 | ng/ul | 2469170 | 5 | 21.88 | 2554850 | 20.0 |
| (DEL) | Alkane: Str... | 21.08 | 14.9 | ng/ul | 1908760 | 5 | 21.88 | 2554850 | 20.0 |
| (DEL) | Alkane: Str... | 21.54 | 38.4 | ng/ul | 4905200 | 5 | 21.88 | 2554850 | 20.0 |
| (DEL) | Alkane: Cyc... | 21.98 | 19.2 | ng/ul | 2457330 | 5 | 21.88 | 2554850 | 20.0 |
| (DEL) | Alkane: Cyc... | 22.37 | 17.8 | ng/ul | 2269110 | 5 | 21.88 | 2554850 | 20.0 |
| (DEL) | Alkane: Str... | 22.46 | 32.8 | ng/ul | 4188120 | 5 | 21.88 | 2554850 | 20.0 |
| (DEL) | Alkane: Str... | 23.52 | 17.1 | ng/ul | 1795260 | 6 | 24.37 | 2101650 | 20.0 |

Instrument :  
BNA\_N  
ClientSampleId :  
DAM46DL