

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022420\  
 Data File : BN009891.D  
 Acq On : 25 Feb 2020 09:04  
 Operator : CG/JU  
 Sample : L1424-22DL 10X  
 Misc :  
 ALS Vial : 29 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 C5B30DL

## 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:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN021220MA.M  
 Title : SVOA CALIBRATION

Signal : TIC

| peak<br># | R.T.<br>min | first<br>scan | max<br>scan | last<br>scan | PK<br>TY | peak<br>height | corr.<br>area | corr.<br>% max. | % of<br>total |
|-----------|-------------|---------------|-------------|--------------|----------|----------------|---------------|-----------------|---------------|
| 1         | 5.452       | 413           | 419         | 428          | rVB      | 142994         | 227804        | 1.46%           | 0.342%        |
| 2         | 6.499       | 593           | 597         | 601          | rVV      | 95158          | 145534        | 0.93%           | 0.219%        |
| 3         | 6.558       | 601           | 607         | 613          | rVV3     | 85732          | 197635        | 1.26%           | 0.297%        |
| 4         | 6.628       | 613           | 619         | 625          | rVB      | 87628          | 135521        | 0.87%           | 0.204%        |
| 5         | 6.787       | 642           | 646         | 654          | rVV      | 170694         | 277999        | 1.78%           | 0.418%        |
| 6         | 7.040       | 682           | 689         | 698          | rBV      | 719442         | 1170493       | 7.49%           | 1.758%        |
| 7         | 7.387       | 739           | 748         | 754          | rBV2     | 471712         | 836694        | 5.35%           | 1.257%        |
| 8         | 7.499       | 761           | 767         | 776          | rVB      | 415603         | 662932        | 4.24%           | 0.996%        |
| 9         | 7.663       | 789           | 795         | 803          | rVB5     | 56153          | 132813        | 0.85%           | 0.199%        |
| 10        | 7.746       | 803           | 809         | 817          | rVB      | 110051         | 203825        | 1.30%           | 0.306%        |
| 11        | 7.922       | 834           | 839         | 845          | rVB4     | 117967         | 210210        | 1.35%           | 0.316%        |
| 12        | 8.016       | 851           | 855         | 859          | rBV2     | 131277         | 213680        | 1.37%           | 0.321%        |
| 13        | 8.181       | 880           | 883         | 889          | rVB2     | 81020          | 132473        | 0.85%           | 0.199%        |
| 14        | 8.328       | 902           | 908         | 912          | rBV      | 108384         | 170956        | 1.09%           | 0.257%        |
| 15        | 8.375       | 912           | 916         | 923          | rVV      | 127440         | 197082        | 1.26%           | 0.296%        |
| 16        | 8.475       | 923           | 933         | 939          | rVV      | 247538         | 438208        | 2.80%           | 0.658%        |
| 17        | 8.546       | 939           | 945         | 951          | rVV3     | 139543         | 313302        | 2.00%           | 0.471%        |
| 18        | 8.604       | 951           | 955         | 963          | rVB      | 229793         | 335911        | 2.15%           | 0.505%        |
| 19        | 8.716       | 963           | 974         | 980          | rVB7     | 62197          | 163258        | 1.04%           | 0.245%        |
| 20        | 8.799       | 980           | 988         | 994          | rBV2     | 91816          | 173405        | 1.11%           | 0.260%        |
| 21        | 8.993       | 1016          | 1021        | 1026         | rVV      | 137660         | 237154        | 1.52%           | 0.356%        |
| 22        | 9.046       | 1026          | 1030        | 1038         | rVB      | 210275         | 349194        | 2.23%           | 0.524%        |
| 23        | 9.546       | 1104          | 1115        | 1120         | rBV3     | 220936         | 498071        | 3.19%           | 0.748%        |
| 24        | 9.769       | 1150          | 1153        | 1162         | rVV3     | 94113          | 217380        | 1.39%           | 0.326%        |
| 25        | 10.057      | 1197          | 1202        | 1206         | rVV5     | 81000          | 171957        | 1.10%           | 0.258%        |
| 26        | 10.134      | 1208          | 1215        | 1219         | rVV      | 793664         | 1411611       | 9.03%           | 2.120%        |
| 27        | 10.181      | 1219          | 1223        | 1232         | rVV      | 1109792        | 1885380       | 12.06%          | 2.832%        |
| 28        | 11.046      | 1365          | 1370        | 1379         | rVB3     | 69081          | 148984        | 0.95%           | 0.224%        |
| 29        | 11.175      | 1385          | 1392        | 1397         | rBV3     | 93204          | 181151        | 1.16%           | 0.272%        |
| 30        | 11.587      | 1454          | 1462        | 1467         | rVV      | 227867         | 367830        | 2.35%           | 0.552%        |
| 31        | 11.804      | 1488          | 1499        | 1508         | rVB      | 746533         | 1262117       | 8.08%           | 1.896%        |
| 32        | 11.993      | 1521          | 1531        | 1532         | rBV2     | 152077         | 256808        | 1.64%           | 0.386%        |
| 33        | 12.028      | 1532          | 1537        | 1543         | rVB      | 1284254        | 1978318       | 12.66%          | 2.971%        |
| 34        | 12.428      | 1599          | 1605        | 1611         | rBV3     | 77855          | 165623        | 1.06%           | 0.249%        |

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## Integration Parameters: LSCINT.P

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 Sampling : 1  
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 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:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN021220MA.M  
 Title : SVOA CALIBRATION

|    |        |      |      |      |      |        |         |        |        |
|----|--------|------|------|------|------|--------|---------|--------|--------|
| 35 | 12.522 | 1616 | 1621 | 1625 | rVV4 | 123727 | 212199  | 1.36%  | 0.319% |
| 36 | 12.745 | 1654 | 1659 | 1670 | rBV  | 505843 | 870890  | 5.57%  | 1.308% |
| 37 | 12.863 | 1671 | 1679 | 1686 | rBV2 | 544557 | 907719  | 5.81%  | 1.363% |
| 38 | 13.045 | 1704 | 1710 | 1712 | rBV  | 245356 | 378407  | 2.42%  | 0.568% |
| 39 | 13.181 | 1727 | 1733 | 1743 | rBV2 | 582224 | 1522829 | 9.74%  | 2.287% |
| 40 | 13.339 | 1754 | 1760 | 1766 | rVV  | 975127 | 1798038 | 11.51% | 2.701% |
| 41 | 13.398 | 1766 | 1770 | 1774 | rVV  | 652995 | 1001037 | 6.41%  | 1.504% |
| 42 | 13.445 | 1774 | 1778 | 1784 | rVV4 | 197994 | 404683  | 2.59%  | 0.608% |
| 43 | 13.534 | 1786 | 1793 | 1796 | rVV2 | 183369 | 339104  | 2.17%  | 0.509% |
| 44 | 13.581 | 1796 | 1801 | 1805 | rVV2 | 524904 | 855925  | 5.48%  | 1.286% |
| 45 | 13.610 | 1805 | 1806 | 1811 | rVV2 | 168480 | 176626  | 1.13%  | 0.265% |
| 46 | 13.710 | 1819 | 1823 | 1826 | rBV  | 146527 | 216060  | 1.38%  | 0.325% |
| 47 | 13.745 | 1826 | 1829 | 1835 | rVB  | 315511 | 444148  | 2.84%  | 0.667% |
| 48 | 13.957 | 1860 | 1865 | 1870 | rBV  | 765994 | 974718  | 6.24%  | 1.464% |
| 49 | 14.022 | 1871 | 1876 | 1883 | rVV2 | 958281 | 1610657 | 10.31% | 2.419% |
| 50 | 14.087 | 1883 | 1887 | 1890 | rVV3 | 171630 | 261814  | 1.68%  | 0.393% |
| 51 | 14.122 | 1890 | 1893 | 1898 | rVV  | 130797 | 182238  | 1.17%  | 0.274% |
| 52 | 14.245 | 1908 | 1914 | 1922 | rBV  | 251801 | 598233  | 3.83%  | 0.899% |
| 53 | 14.439 | 1938 | 1947 | 1954 | rVB3 | 375629 | 762907  | 4.88%  | 1.146% |
| 54 | 14.516 | 1954 | 1960 | 1964 | rBV  | 333794 | 472119  | 3.02%  | 0.709% |
| 55 | 14.586 | 1967 | 1972 | 1978 | rVB3 | 389950 | 583328  | 3.73%  | 0.876% |
| 56 | 14.669 | 1978 | 1986 | 1990 | rBV3 | 545346 | 1125213 | 7.20%  | 1.690% |
| 57 | 14.716 | 1990 | 1994 | 1998 | rVB  | 273915 | 398243  | 2.55%  | 0.598% |
| 58 | 14.828 | 2006 | 2013 | 2017 | rBV3 | 217639 | 460871  | 2.95%  | 0.692% |
| 59 | 14.863 | 2017 | 2019 | 2025 | rVB2 | 116262 | 147146  | 0.94%  | 0.221% |
| 60 | 14.922 | 2025 | 2029 | 2033 | rBV  | 743646 | 843188  | 5.40%  | 1.266% |
| 61 | 14.992 | 2036 | 2041 | 2043 | rBV3 | 84498  | 138738  | 0.89%  | 0.208% |
| 62 | 15.028 | 2043 | 2047 | 2052 | rVB2 | 243045 | 368472  | 2.36%  | 0.553% |
| 63 | 15.086 | 2052 | 2057 | 2062 | rVB3 | 223259 | 353001  | 2.26%  | 0.530% |
| 64 | 15.139 | 2062 | 2066 | 2070 | rBV3 | 126335 | 211364  | 1.35%  | 0.317% |
| 65 | 15.198 | 2070 | 2076 | 2082 | rBV2 | 841327 | 1424276 | 9.11%  | 2.139% |
| 66 | 15.245 | 2082 | 2084 | 2091 | rVB5 | 115326 | 177638  | 1.14%  | 0.267% |
| 67 | 15.328 | 2091 | 2098 | 2103 | rBV3 | 254458 | 523872  | 3.35%  | 0.787% |
| 68 | 15.386 | 2103 | 2108 | 2111 | rBV4 | 127793 | 238649  | 1.53%  | 0.358% |
| 69 | 15.481 | 2118 | 2124 | 2127 | rBV4 | 108123 | 199703  | 1.28%  | 0.300% |
| 70 | 15.545 | 2133 | 2135 | 2141 | rVB4 | 156240 | 228404  | 1.46%  | 0.343% |
| 71 | 15.616 | 2141 | 2147 | 2151 | rBV4 | 97992  | 201481  | 1.29%  | 0.303% |

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 C5B30DL

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 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:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN021220MA.M  
 Title : SVOA CALIBRATION

|     |        |      |      |      |      |          |          |         |         |
|-----|--------|------|------|------|------|----------|----------|---------|---------|
| 72  | 15.786 | 2170 | 2176 | 2179 | rBV2 | 840567   | 1209577  | 7.74%   | 1.817%  |
| 73  | 16.092 | 2222 | 2228 | 2232 | rBV5 | 106033   | 244613   | 1.57%   | 0.367%  |
| 74  | 16.145 | 2232 | 2237 | 2242 | rVV3 | 253367   | 388427   | 2.49%   | 0.583%  |
| 75  | 16.298 | 2260 | 2263 | 2266 | rBV3 | 111256   | 138395   | 0.89%   | 0.208%  |
| 76  | 16.457 | 2283 | 2290 | 2297 | rBV  | 11021915 | 15626986 | 100.00% | 23.471% |
| 77  | 16.581 | 2306 | 2311 | 2314 | rBV  | 643253   | 770605   | 4.93%   | 1.157%  |
| 78  | 16.633 | 2317 | 2320 | 2327 | rVB2 | 171748   | 263308   | 1.68%   | 0.395%  |
| 79  | 16.786 | 2341 | 2346 | 2350 | rBV  | 1050842  | 1370633  | 8.77%   | 2.059%  |
| 80  | 16.828 | 2350 | 2353 | 2358 | rVB  | 455580   | 567365   | 3.63%   | 0.852%  |
| 81  | 16.886 | 2358 | 2363 | 2367 | rBV2 | 256958   | 393209   | 2.52%   | 0.591%  |
| 82  | 16.992 | 2376 | 2381 | 2386 | rBV5 | 91543    | 170518   | 1.09%   | 0.256%  |
| 83  | 17.316 | 2432 | 2436 | 2440 | rVB  | 494100   | 538477   | 3.45%   | 0.809%  |
| 84  | 17.663 | 2491 | 2495 | 2499 | rVV  | 235480   | 334637   | 2.14%   | 0.503%  |
| 85  | 17.710 | 2499 | 2503 | 2509 | rVB  | 296547   | 399965   | 2.56%   | 0.601%  |
| 86  | 17.845 | 2523 | 2526 | 2531 | rBV  | 162344   | 269542   | 1.72%   | 0.405%  |
| 87  | 17.892 | 2531 | 2534 | 2544 | rVB  | 167025   | 245419   | 1.57%   | 0.369%  |
| 88  | 18.010 | 2549 | 2554 | 2559 | rVB  | 401676   | 485639   | 3.11%   | 0.729%  |
| 89  | 18.592 | 2649 | 2653 | 2658 | rBV  | 177338   | 266365   | 1.70%   | 0.400%  |
| 90  | 18.657 | 2658 | 2664 | 2668 | rVV2 | 277828   | 364657   | 2.33%   | 0.548%  |
| 91  | 19.192 | 2751 | 2755 | 2759 | rBV  | 242082   | 342155   | 2.19%   | 0.514%  |
| 92  | 19.245 | 2761 | 2764 | 2771 | rVV2 | 184749   | 252725   | 1.62%   | 0.380%  |
| 93  | 21.004 | 3058 | 3063 | 3068 | rBV  | 1363069  | 1638563  | 10.49%  | 2.461%  |
| 94  | 22.039 | 3236 | 3239 | 3244 | rVB2 | 133501   | 177041   | 1.13%   | 0.266%  |
| 95  | 22.504 | 3315 | 3318 | 3323 | rVB  | 154579   | 201004   | 1.29%   | 0.302%  |
| 96  | 22.974 | 3393 | 3398 | 3402 | rBV  | 166350   | 287085   | 1.84%   | 0.431%  |
| 97  | 23.021 | 3402 | 3406 | 3413 | rVB  | 197908   | 306286   | 1.96%   | 0.460%  |
| 98  | 23.104 | 3413 | 3420 | 3429 | rBV  | 923740   | 1646936  | 10.54%  | 2.474%  |
| 99  | 23.604 | 3501 | 3505 | 3513 | rVB  | 191750   | 317857   | 2.03%   | 0.477%  |
| 100 | 24.268 | 3614 | 3618 | 3627 | rVB  | 142988   | 275840   | 1.77%   | 0.414%  |

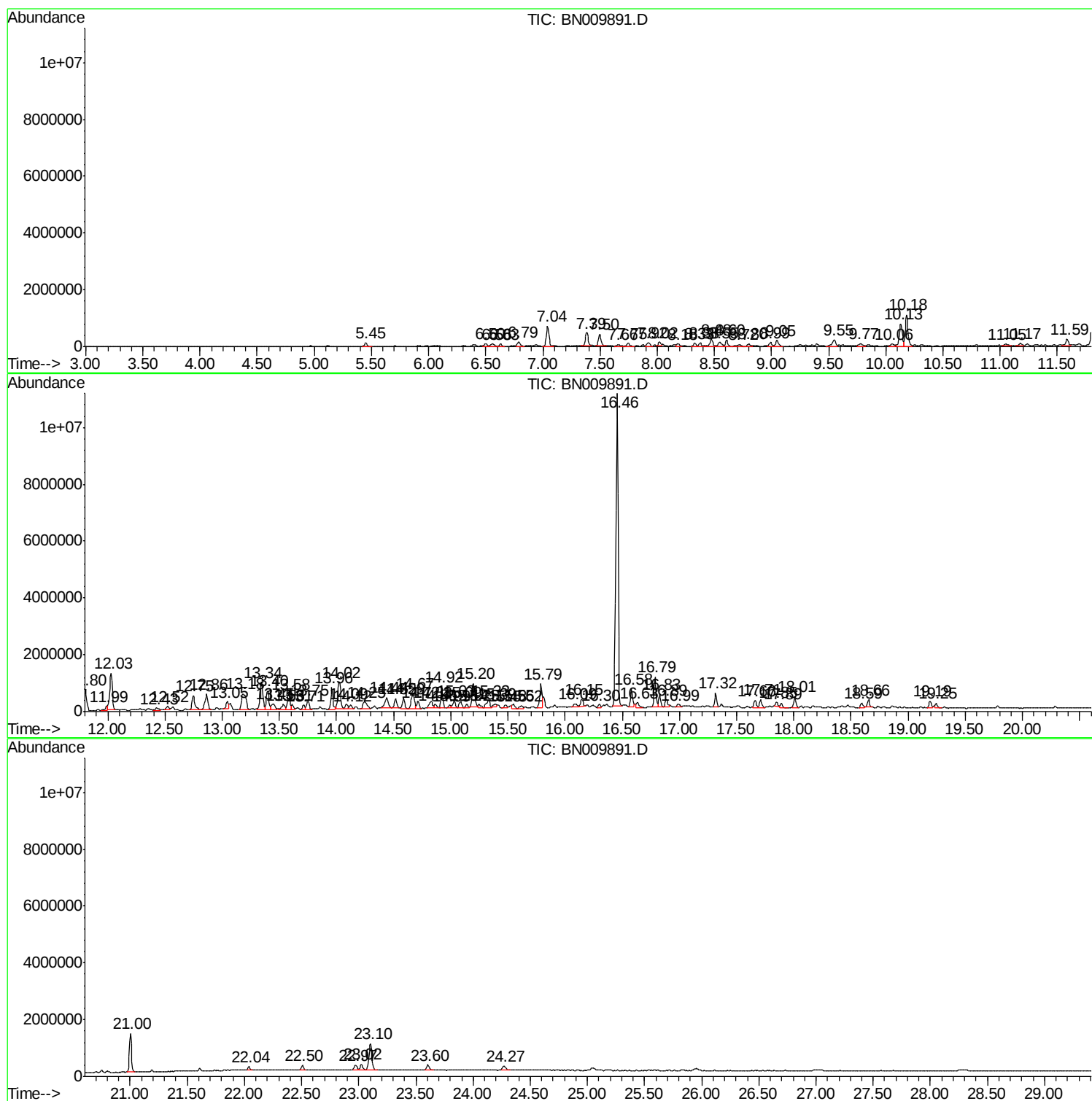
Sum of corrected areas: 66579080

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ALS Vial : 29 Sample Multiplier: 1

Instrument :  
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN021220MA.M  
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_N\DATA\BN022420\  
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C5B30DL

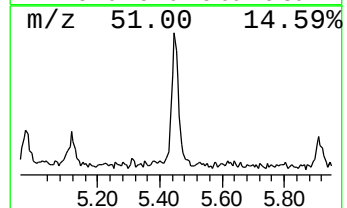
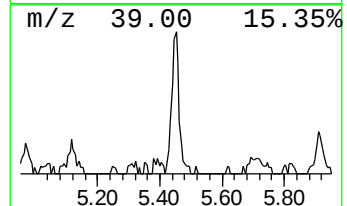
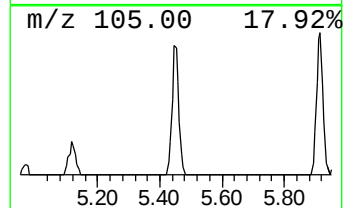
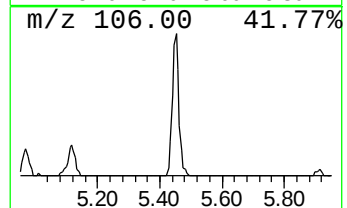
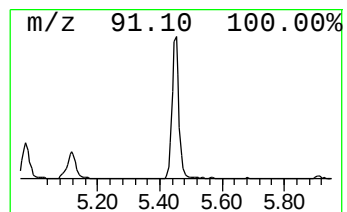
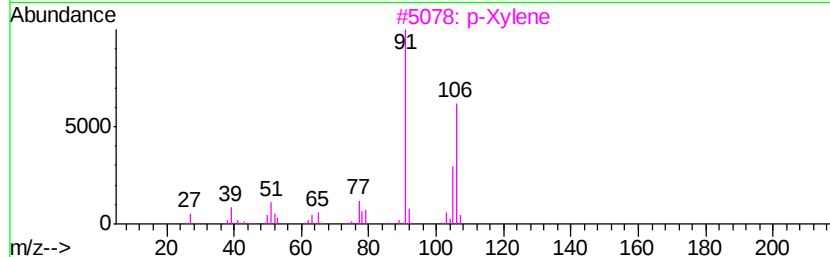
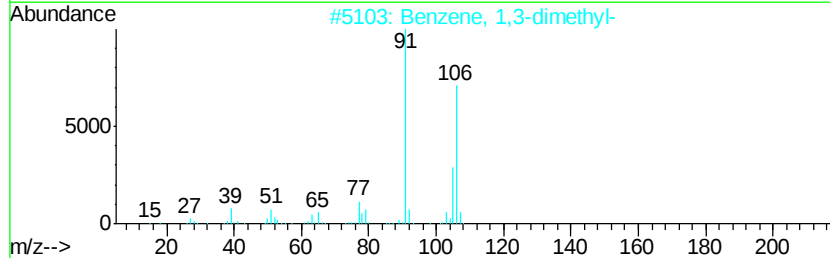
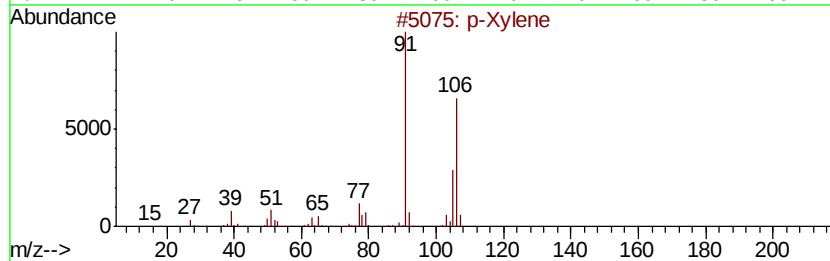
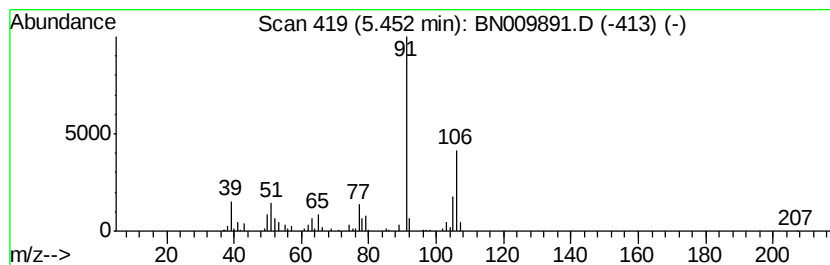
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN021220MA.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 1 p-Xylene Concentration Rank 25

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 5.45 | 5.45 ng/ul | 227804 | 1,4-Dichlorobenzene-d4 | 7.39 |

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



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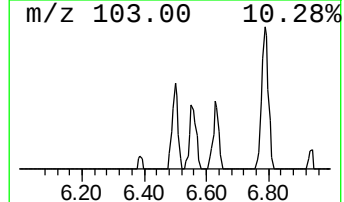
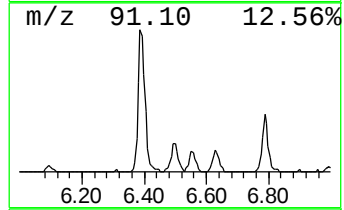
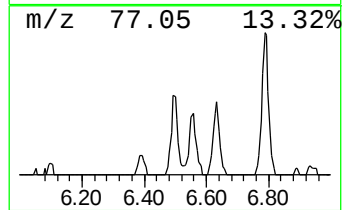
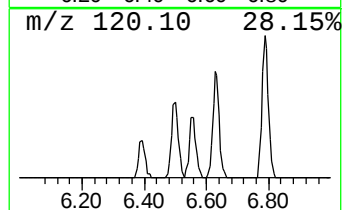
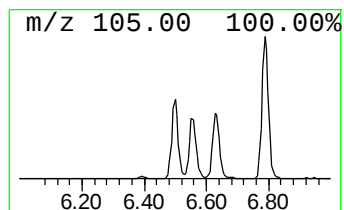
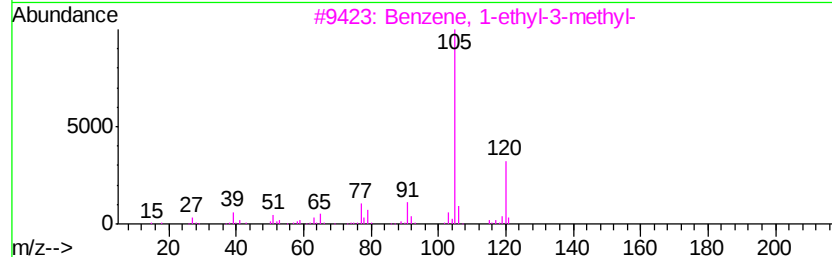
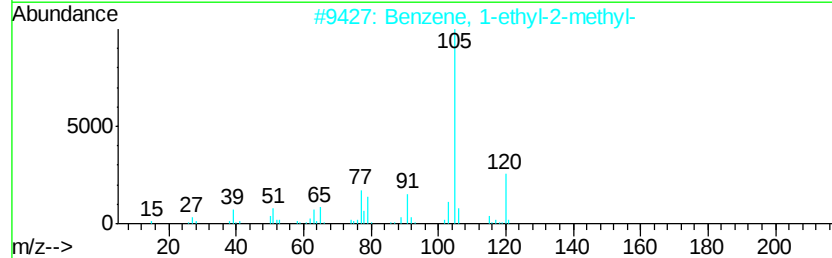
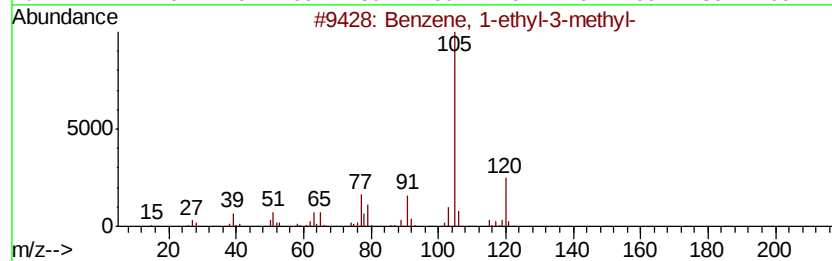
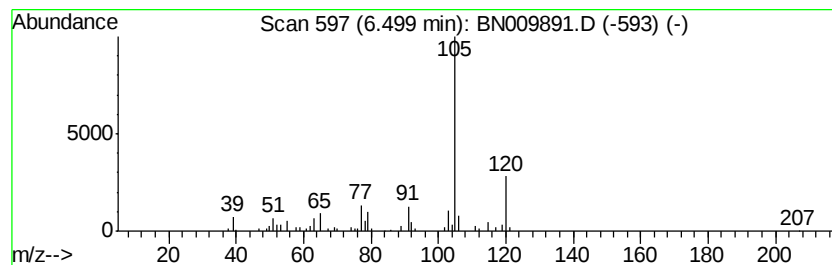
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN021220MA.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 2 Benzene, 1-ethyl-3-methyl- Concentration Rank 45

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 6.50 | 3.48 ng/ul | 145534 | 1,4-Dichlorobenzene-d4 | 7.39 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 95   |
| 2    |    | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 94   |
| 3    |    | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 94   |
| 4    |    | Mesitylene                 | 120 | C9H12   | 000108-67-8 | 91   |
| 5    |    | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 91   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022420\  
Data File : BN009891.D  
Acq On : 25 Feb 2020 09:04  
Operator : CG/JU  
Sample : L1424-22DL 10X  
Misc :  
ALS Vial : 29 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
C5B30DL

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Quant Title : SVOA CALIBRATION

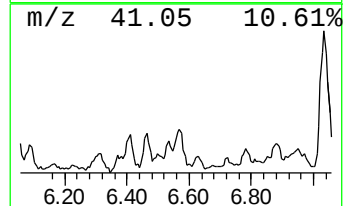
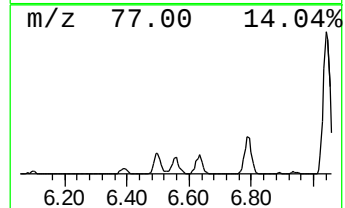
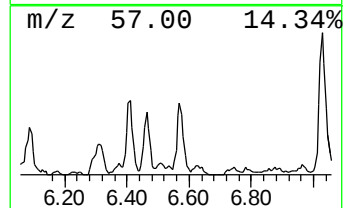
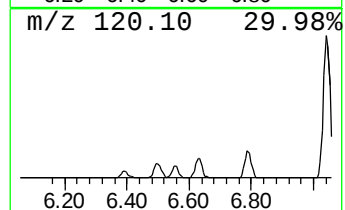
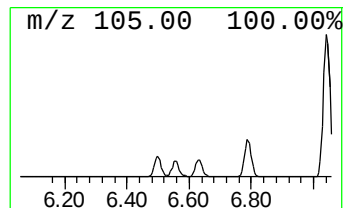
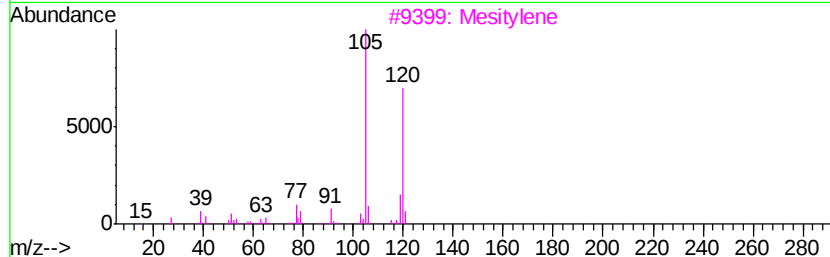
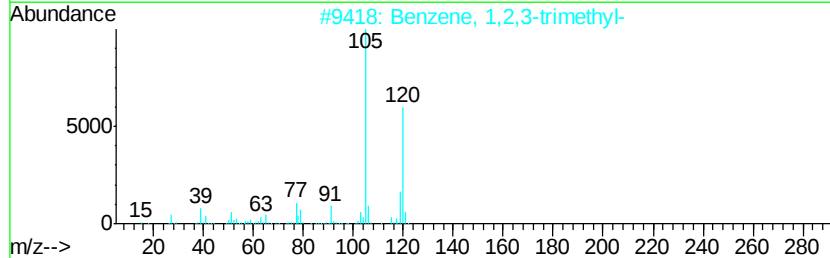
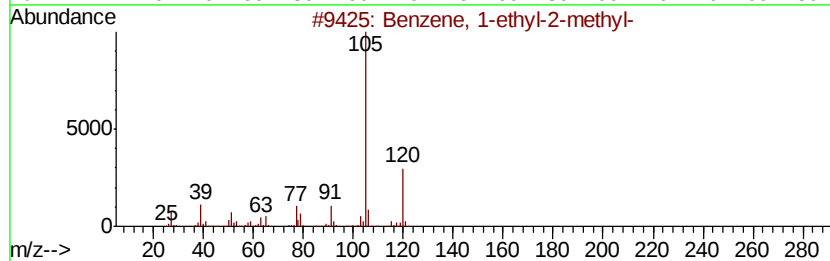
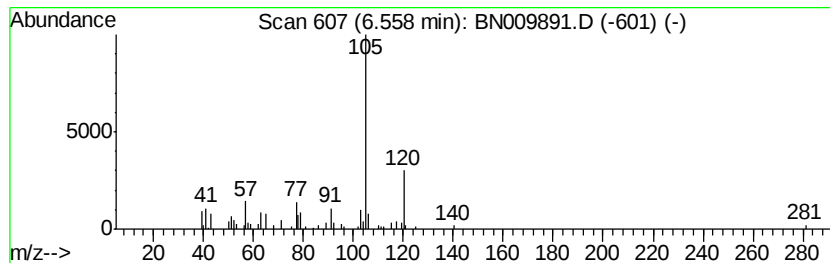
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 3 Benzene, 1-ethyl-2-methyl- Concentration Rank 33

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 6.56 | 4.72 ng/ul | 197635 | 1,4-Dichlorobenzene-d4 | 7.39 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 93   |
| 2    |    | Benzene, 1,2,3-trimethyl-  | 120 | C9H12   | 000526-73-8 | 87   |
| 3    |    | Mesitylene                 | 120 | C9H12   | 000108-67-8 | 87   |
| 4    |    | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 87   |
| 5    |    | Mesitylene                 | 120 | C9H12   | 000108-67-8 | 87   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_N\DATA\BN022420\  
Data File : BN009891.D  
Acq On : 25 Feb 2020 09:04  
Operator : CG/JU  
Sample : L1424-22DL 10X  
Misc :  
ALS Vial : 29 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
C5B30DL

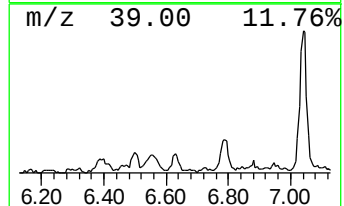
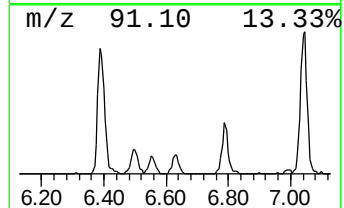
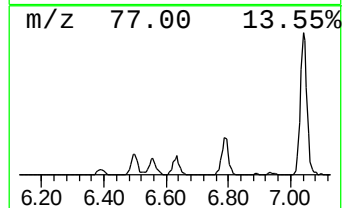
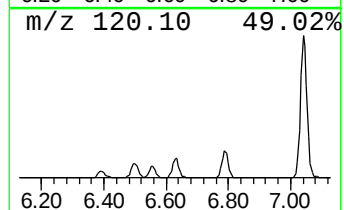
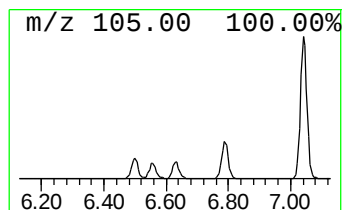
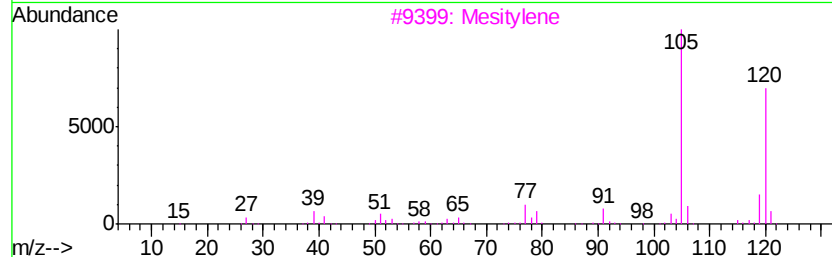
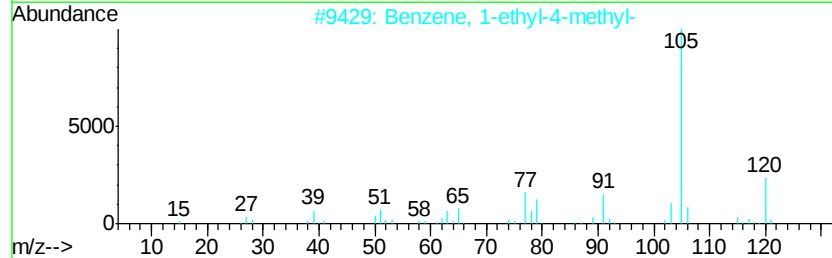
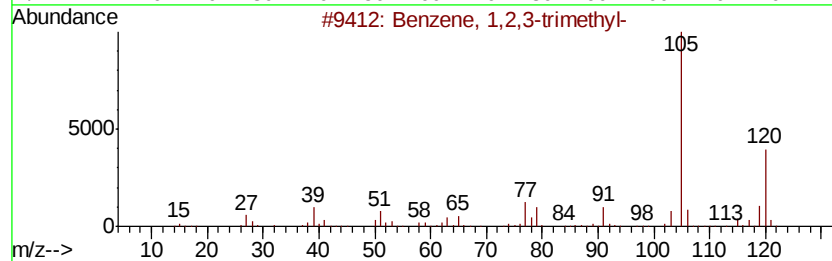
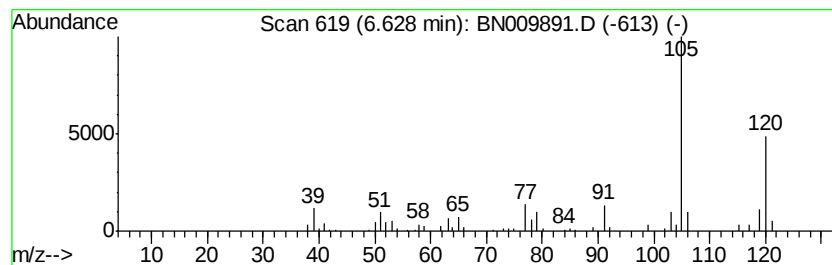
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Quant Title : SVOA CALIBRATION

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

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Peak Number 4 Benzene, 1,2,3-trimethyl- Concentration Rank 49

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 6.63 | 3.24 ng/ul | 135521 | 1,4-Dichlorobenzene-d4 | 7.39 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1,2,3-trimethyl-  | 120 | C9H12   | 000526-73-8 | 95   |
| 2    |    |   | Benzene, 1-ethyl-4-methyl- | 120 | C9H12   | 000622-96-8 | 95   |
| 3    |    |   | Mesitylene                 | 120 | C9H12   | 000108-67-8 | 94   |
| 4    |    |   | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 94   |
| 5    |    |   | Mesitylene                 | 120 | C9H12   | 000108-67-8 | 94   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022420\  
Data File : BN009891.D  
Acq On : 25 Feb 2020 09:04  
Operator : CG/JU  
Sample : L1424-22DL 10X  
Misc :  
ALS Vial : 29 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
C5B30DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN021220MA.M  
Quant Title : SVOA CALIBRATION

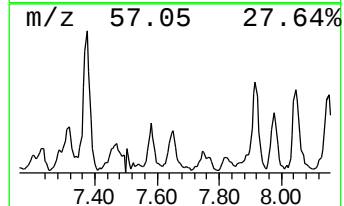
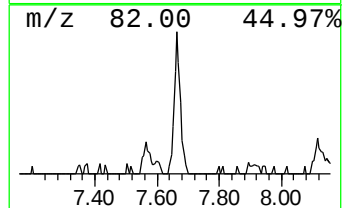
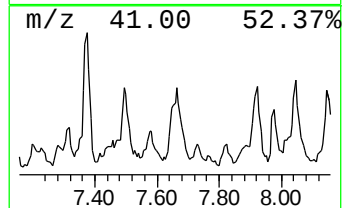
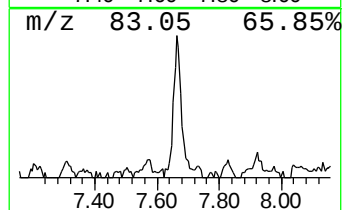
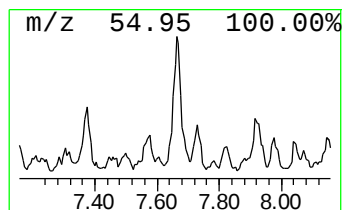
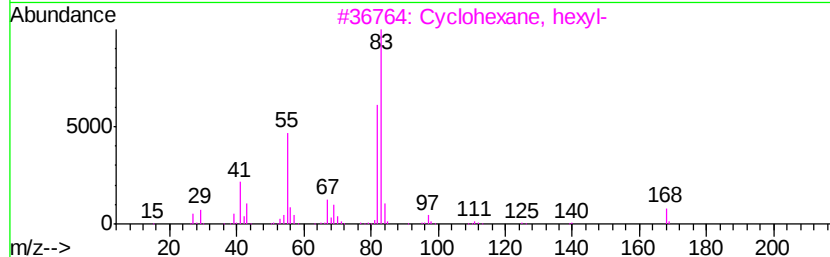
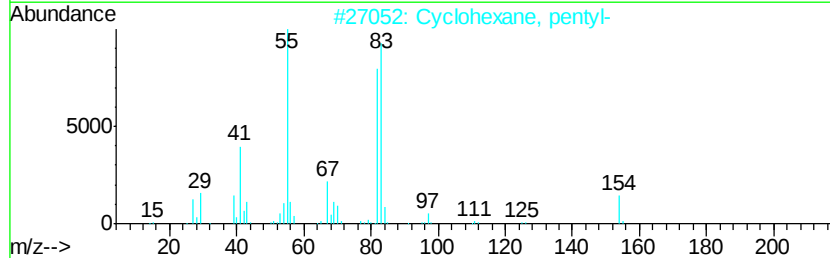
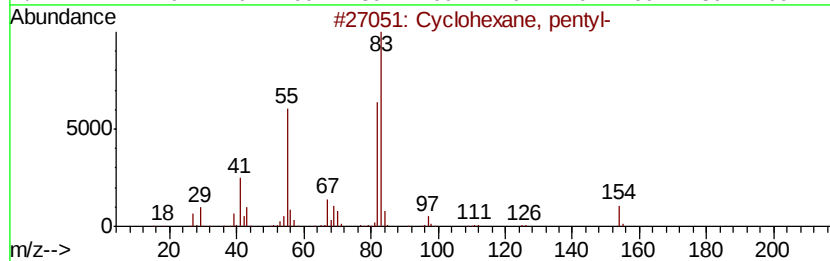
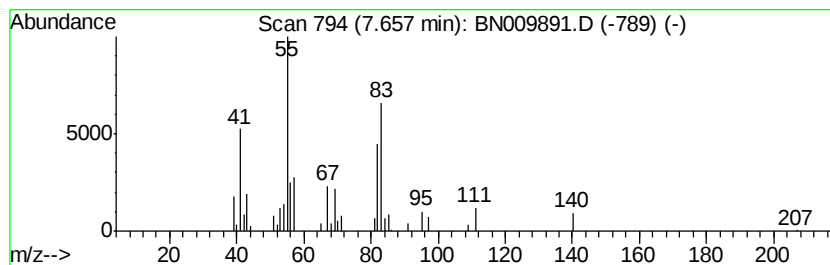
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 7 (DEL) Alkane: Cyclic7.66 Concentration Rank 50

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 7.66 | 3.17 ng/ul | 132813 | 1,4-Dichlorobenzene-d4 | 7.39 |

| Hit# | of | 5 | Tentative ID         | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------|-----|---------|-------------|------|
| 1    |    |   | Cyclohexane, pentyl- | 154 | C11H22  | 004292-92-6 | 72   |
| 2    |    |   | Cyclohexane, pentyl- | 154 | C11H22  | 004292-92-6 | 72   |
| 3    |    |   | Cyclohexane, hexyl-  | 168 | C12H24  | 004292-75-5 | 64   |
| 4    |    |   | Cyclohexane, pentyl- | 154 | C11H22  | 004292-92-6 | 64   |
| 5    |    |   | Cyclohexane, butyl-  | 140 | C10H20  | 001678-93-9 | 59   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022420\  
Data File : BN009891.D  
Acq On : 25 Feb 2020 09:04  
Operator : CG/JU  
Sample : L1424-22DL 10X  
Misc :  
ALS Vial : 29 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
C5B30DL

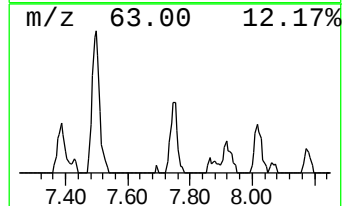
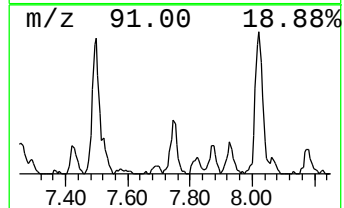
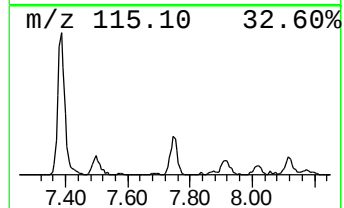
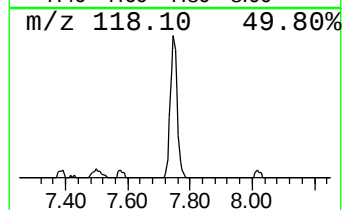
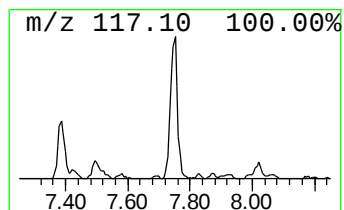
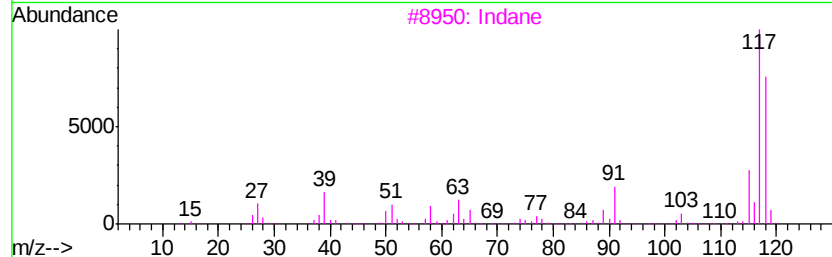
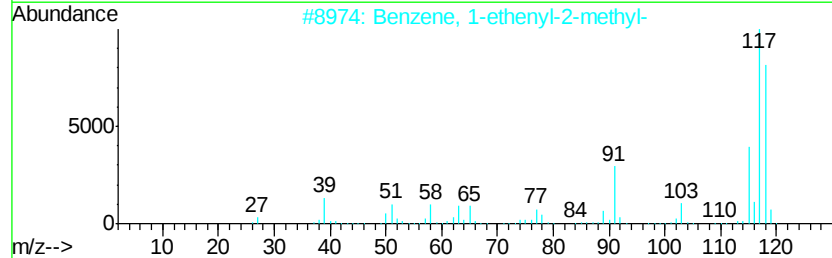
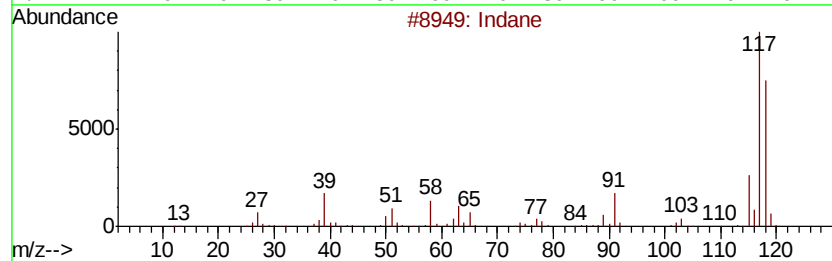
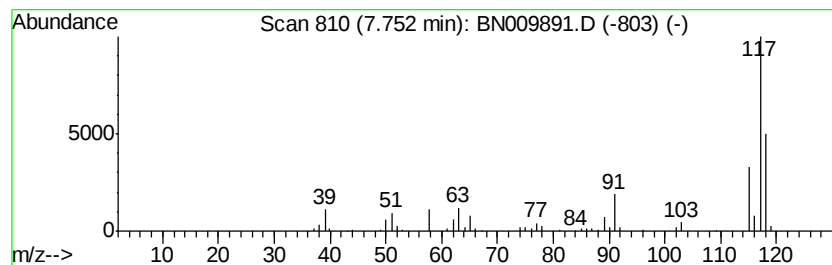
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Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 8 Indane Concentration Rank 32

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 7.75 | 4.87 ng/ul | 203825 | 1,4-Dichlorobenzene-d4 | 7.39 |

| Hit# | of | Tentative ID                 | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------------|-----|---------|-------------|------|
| 1    | 5  | Indane                       | 118 | C9H10   | 000496-11-7 | 91   |
| 2    |    | Benzene, 1-ethenyl-2-methyl- | 118 | C9H10   | 000611-15-4 | 87   |
| 3    |    | Indane                       | 118 | C9H10   | 000496-11-7 | 87   |
| 4    |    | Indane                       | 118 | C9H10   | 000496-11-7 | 87   |
| 5    |    | Benzene, 2-propenyl-         | 118 | C9H10   | 000300-57-2 | 76   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022420\  
Data File : BN009891.D  
Acq On : 25 Feb 2020 09:04  
Operator : CG/JU  
Sample : L1424-22DL 10X  
Misc :  
ALS Vial : 29 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
C5B30DL

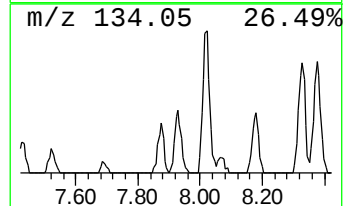
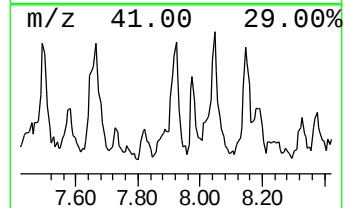
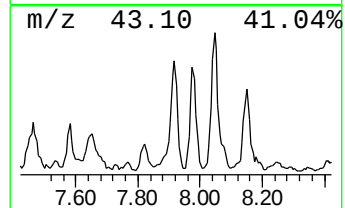
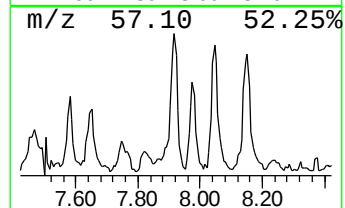
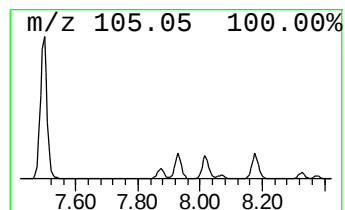
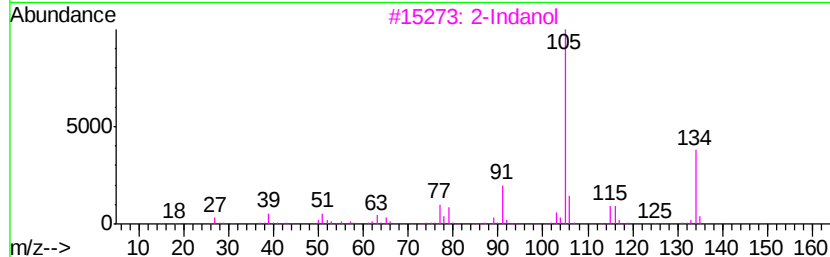
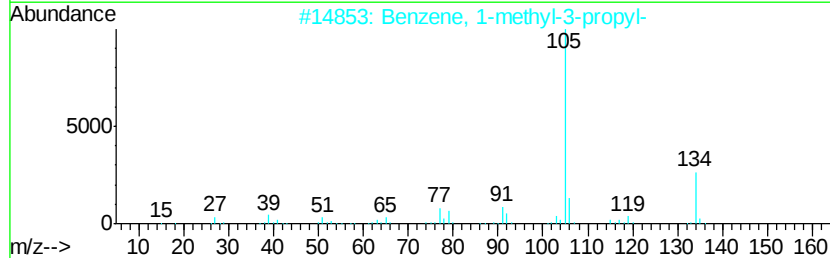
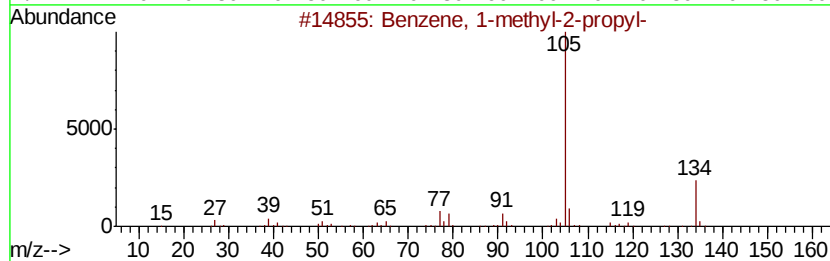
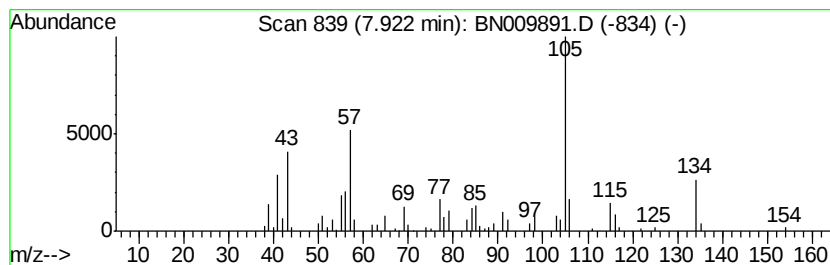
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Quant Title : SVOA CALIBRATION

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

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Peak Number 9 Benzene, 1-methyl-2-propyl- Concentration Rank 29

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 7.92 | 5.02 ng/ul | 210210 | 1,4-Dichlorobenzene-d4 | 7.39 |

| Hit# | of | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1-methyl-2-propyl- | 134 | C10H14  | 001074-17-5 | 50   |
| 2    |    | Benzene, 1-methyl-3-propyl- | 134 | C10H14  | 001074-43-7 | 50   |
| 3    |    | 2-Indanol                   | 134 | C9H10O  | 004254-29-9 | 50   |
| 4    |    | Benzene, 1-methyl-3-propyl- | 134 | C10H14  | 001074-43-7 | 46   |
| 5    |    | 2-Tolyloxirane              | 134 | C9H10O  | 002783-26-8 | 45   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022420\  
Data File : BN009891.D  
Acq On : 25 Feb 2020 09:04  
Operator : CG/JU  
Sample : L1424-22DL 10X  
Misc :  
ALS Vial : 29 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleID :  
C5B30DL

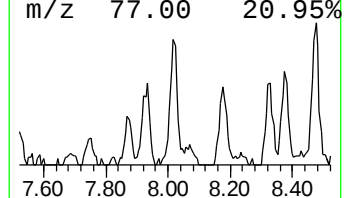
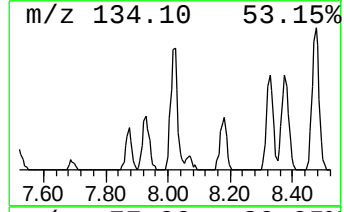
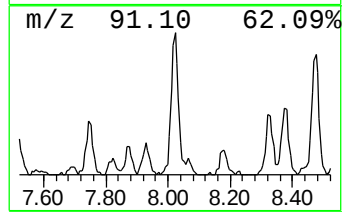
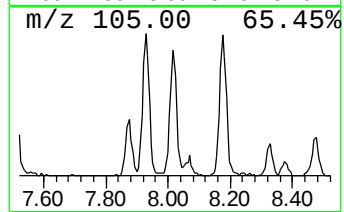
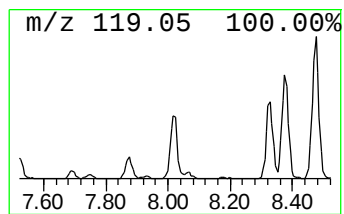
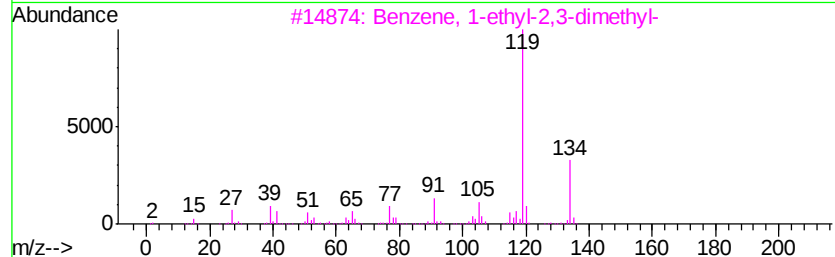
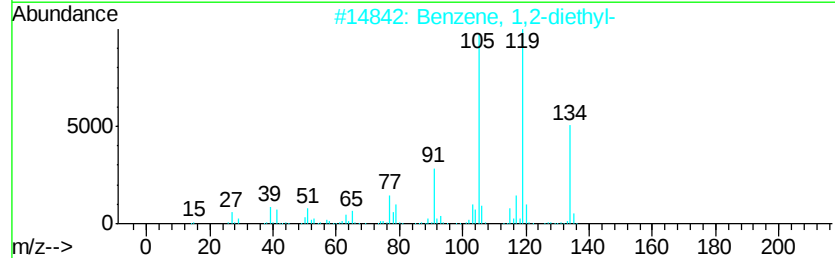
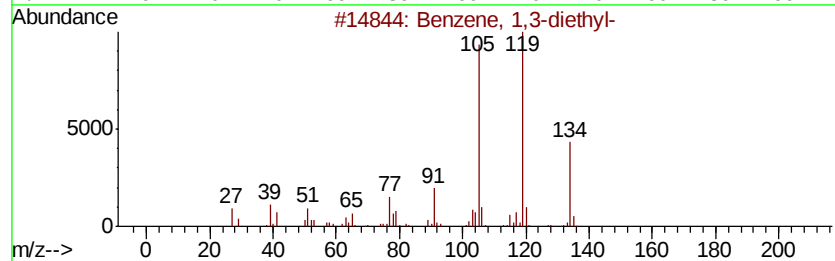
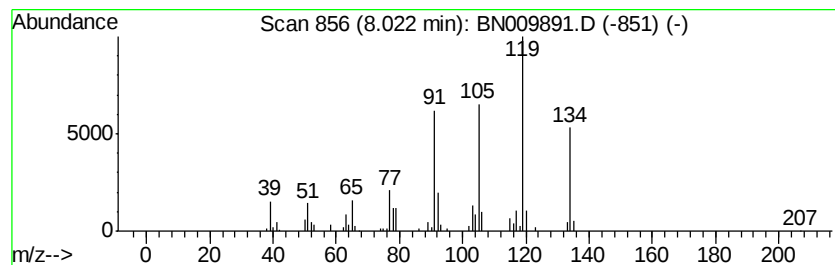
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Quant Title : SVOA CALIBRATION

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

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Peak Number 10 Benzene, 1,3-diethyl- Concentration Rank 28

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 8.02 | 5.11 ng/ul | 213680 | 1,4-Dichlorobenzene-d4 | 7.39 |

| Hit# | of | 5 | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1,3-diethyl-          | 134 | C10H14  | 000141-93-5 | 94   |
| 2    |    |   | Benzene, 1,2-diethyl-          | 134 | C10H14  | 000135-01-3 | 93   |
| 3    |    |   | Benzene, 1-ethyl-2,3-dimethyl- | 134 | C10H14  | 000933-98-2 | 91   |
| 4    |    |   | Benzene, 1,4-diethyl-          | 134 | C10H14  | 000105-05-5 | 90   |
| 5    |    |   | Benzene, 1,2,3,4-tetramethyl-  | 134 | C10H14  | 000488-23-3 | 90   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022420\  
Data File : BN009891.D  
Acq On : 25 Feb 2020 09:04  
Operator : CG/JU  
Sample : L1424-22DL 10X  
Misc :  
ALS Vial : 29 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleID :  
C5B30DL

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Quant Title : SVOA CALIBRATION

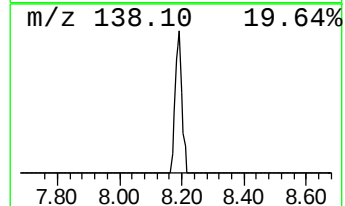
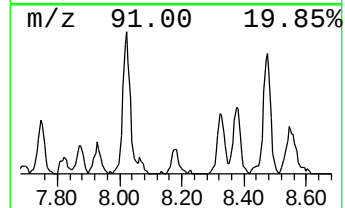
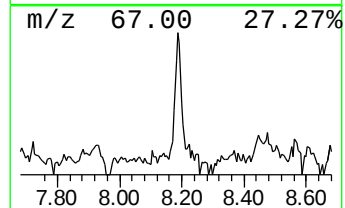
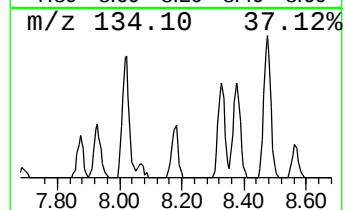
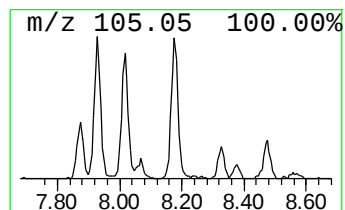
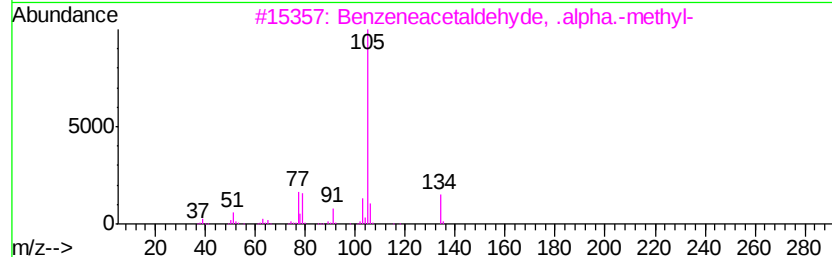
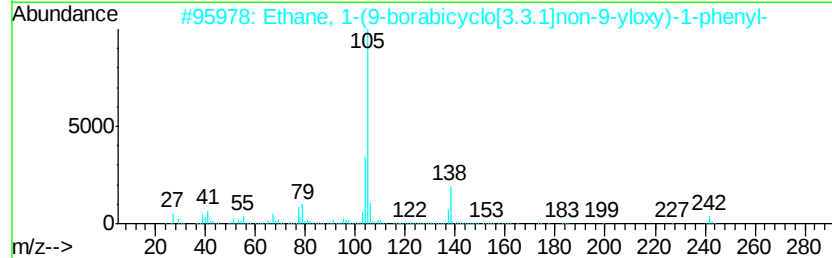
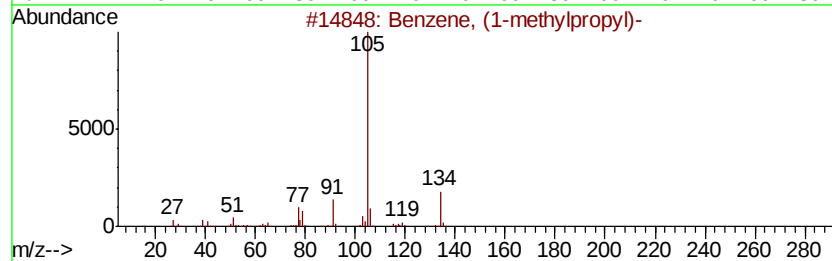
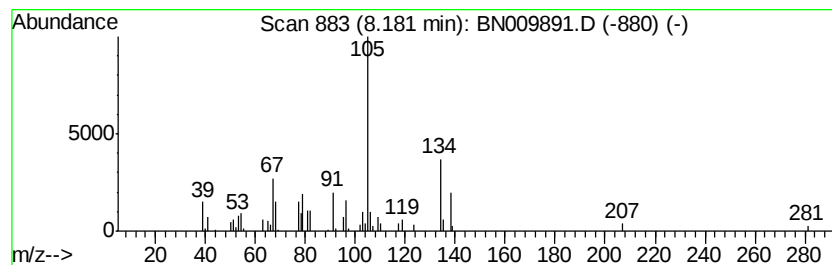
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 11 Benzene, (1-methylpropyl)- Concentration Rank 51

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 8.18 | 3.17 ng/ul | 132473 | 1,4-Dichlorobenzene-d4 | 7.39 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Benzene, (1-methylpropyl)-          | 134 | C10H14   | 000135-98-8  | 50   |
| 2    |    |   | Ethane, 1-(9-borabicyclo[3.3.1]n... | 242 | C16H23BO | 1000160-80-6 | 43   |
| 3    |    |   | Benzeneacetaldehyde, .alpha.-met... | 134 | C9H10O   | 000093-53-8  | 43   |
| 4    |    |   | Oxirane, 2-methyl-2-phenyl-         | 134 | C9H10O   | 002085-88-3  | 43   |
| 5    |    |   | 2-Indanol                           | 134 | C9H10O   | 004254-29-9  | 38   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022420\  
Data File : BN009891.D  
Acq On : 25 Feb 2020 09:04  
Operator : CG/JU  
Sample : L1424-22DL 10X  
Misc :  
ALS Vial : 29 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
C5B30DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN021220MA.M  
Quant Title : SVOA CALIBRATION

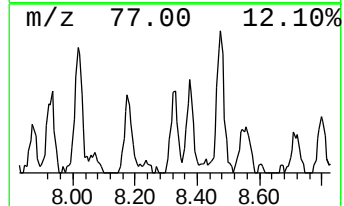
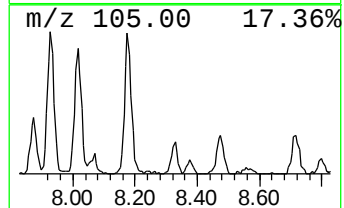
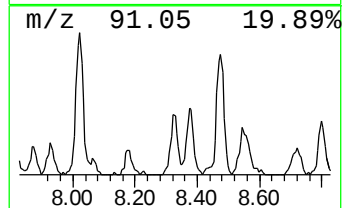
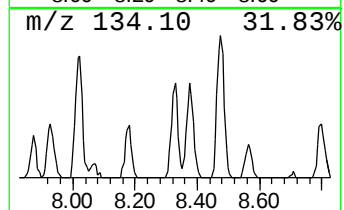
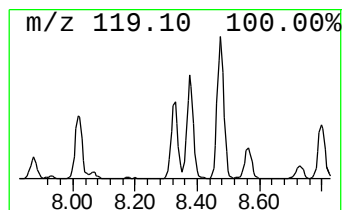
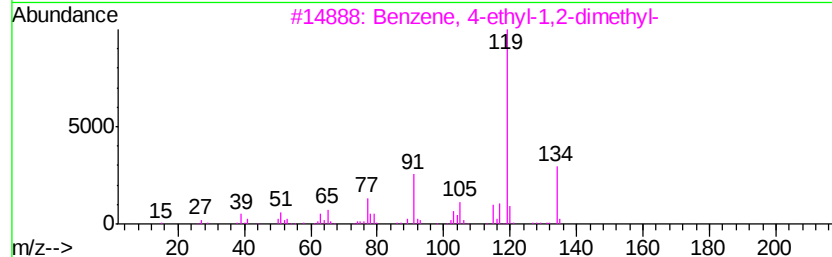
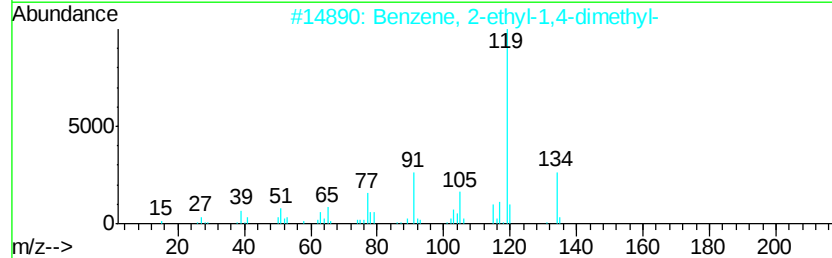
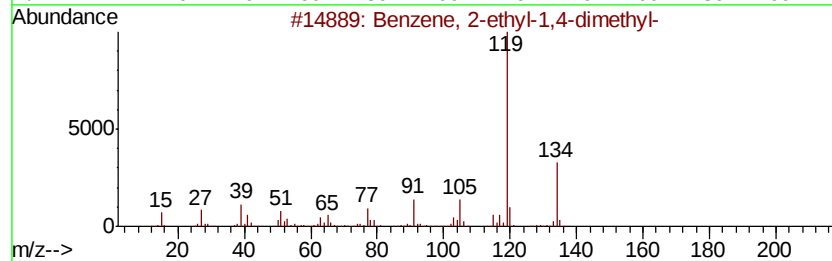
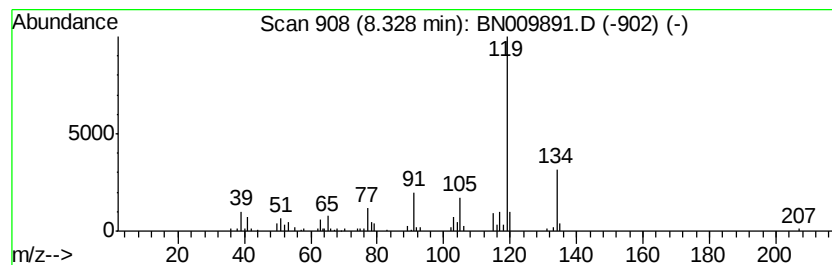
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 12 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 37

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 8.33 | 4.09 ng/ul | 170956 | 1,4-Dichlorobenzene-d4 | 7.39 |

| Hit# | of | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 2-ethyl-1,4-dimethyl- | 134 | C10H14  | 001758-88-9 | 97   |
| 2    |    | Benzene, 2-ethyl-1,4-dimethyl- | 134 | C10H14  | 001758-88-9 | 96   |
| 3    |    | Benzene, 4-ethyl-1,2-dimethyl- | 134 | C10H14  | 000934-80-5 | 96   |
| 4    |    | Benzene, 1-ethyl-2,3-dimethyl- | 134 | C10H14  | 000933-98-2 | 95   |
| 5    |    | o-Cymene                       | 134 | C10H14  | 000527-84-4 | 95   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022420\  
Data File : BN009891.D  
Acq On : 25 Feb 2020 09:04  
Operator : CG/JU  
Sample : L1424-22DL 10X  
Misc :  
ALS Vial : 29 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
C5B30DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN021220MA.M  
Quant Title : SVOA CALIBRATION

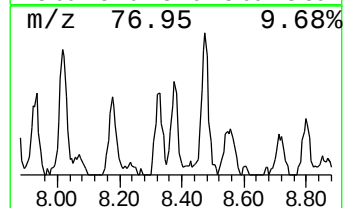
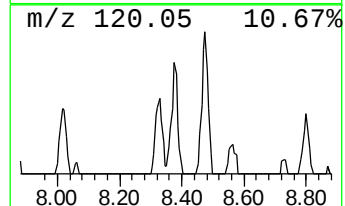
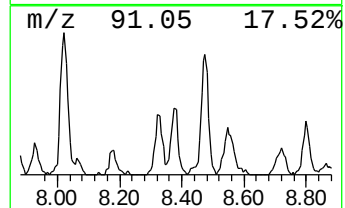
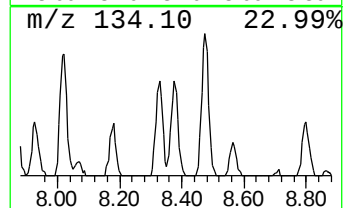
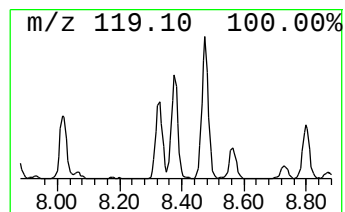
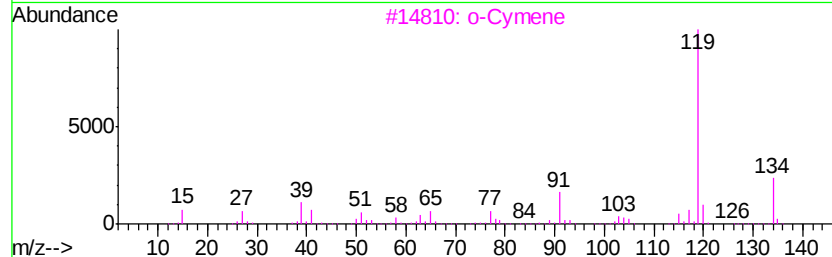
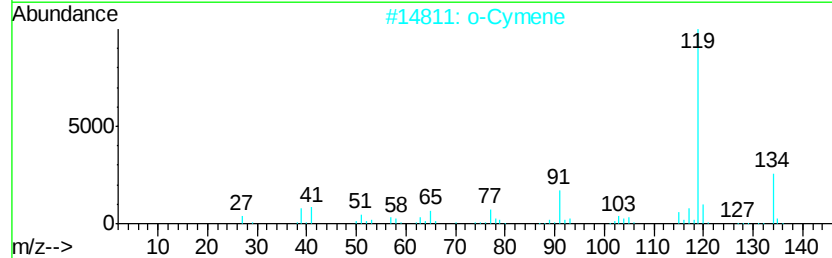
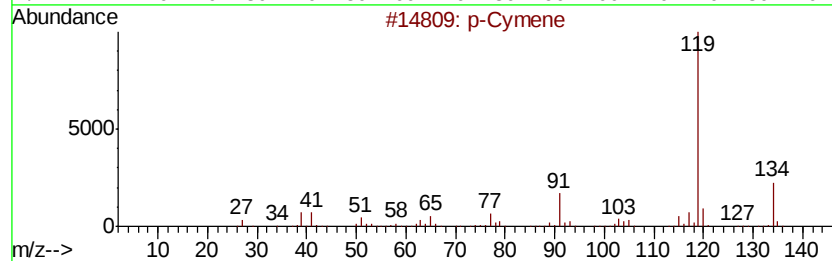
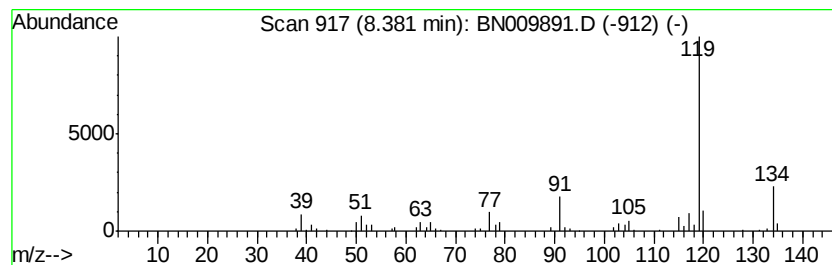
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 13 p-Cymene Concentration Rank 34

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 8.38 | 4.71 ng/ul | 197082 | 1,4-Dichlorobenzene-d4 | 7.39 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | p-Cymene                            | 134 | C10H14  | 000099-87-6 | 97   |
| 2    |    | o-Cymene                            | 134 | C10H14  | 000527-84-4 | 97   |
| 3    |    | o-Cymene                            | 134 | C10H14  | 000527-84-4 | 95   |
| 4    |    | Benzene, 1-methyl-3-(1-methyleth... | 134 | C10H14  | 000535-77-3 | 95   |
| 5    |    | o-Cymene                            | 134 | C10H14  | 000527-84-4 | 94   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022420\  
Data File : BN009891.D  
Acq On : 25 Feb 2020 09:04  
Operator : CG/JU  
Sample : L1424-22DL 10X  
Misc :  
ALS Vial : 29 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
C5B30DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN021220MA.M  
Quant Title : SVOA CALIBRATION

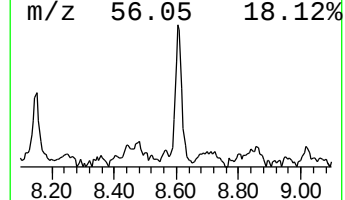
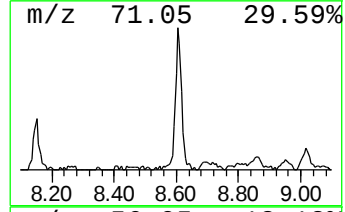
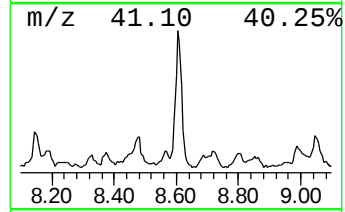
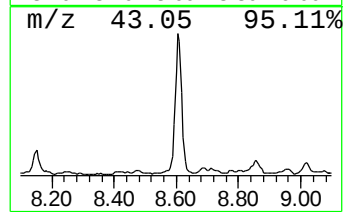
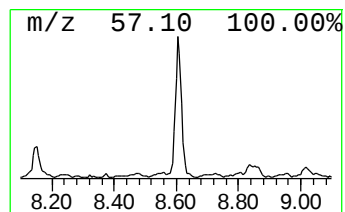
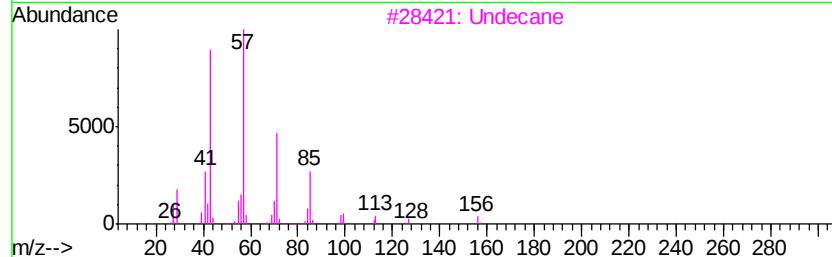
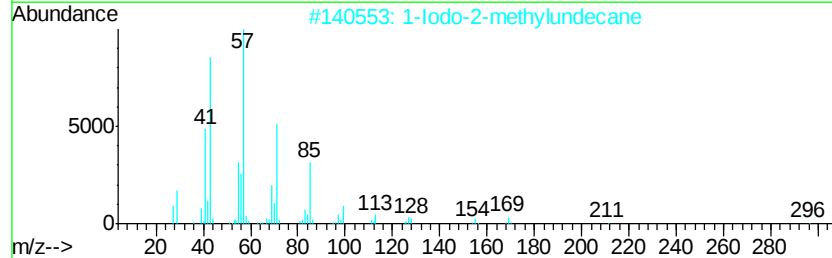
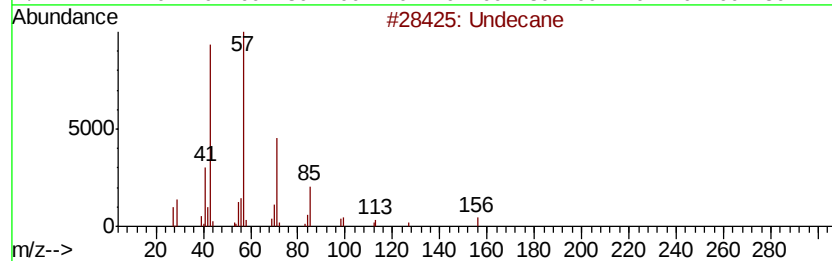
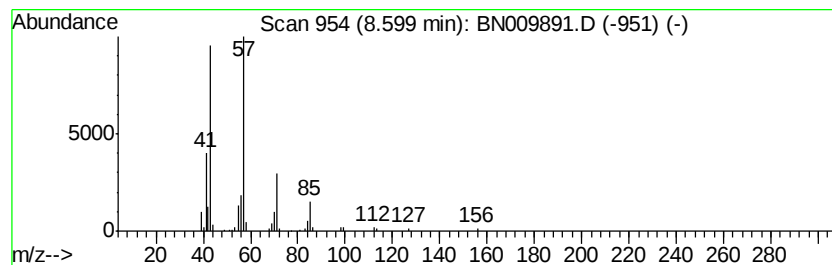
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 8.60 | 8.03 ng/ul | 335911 | 1,4-Dichlorobenzene-d4 | 7.39 |

| Hit# | of | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------------|-----|---------|-------------|------|
| 1    | 5  | Undecane                  | 156 | C11H24  | 001120-21-4 | 83   |
| 2    |    | 1-Iodo-2-methylundecane   | 296 | C12H25I | 073105-67-6 | 78   |
| 3    |    | Undecane                  | 156 | C11H24  | 001120-21-4 | 78   |
| 4    |    | Decane, 6-ethyl-2-methyl- | 184 | C13H28  | 062108-21-8 | 72   |
| 5    |    | Tetradecane               | 198 | C14H30  | 000629-59-4 | 64   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_N\DATA\BN022420\  
Data File : BN009891.D  
Acq On : 25 Feb 2020 09:04  
Operator : CG/JU  
Sample : L1424-22DL 10X  
Misc :  
ALS Vial : 29 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
C5B30DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN021220MA.M  
Quant Title : SVOA CALIBRATION

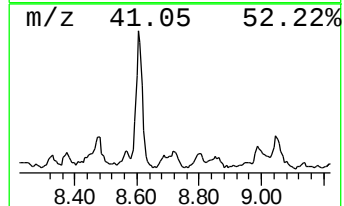
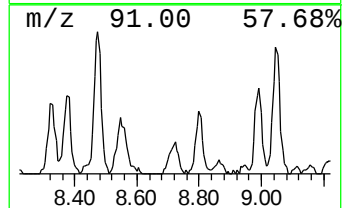
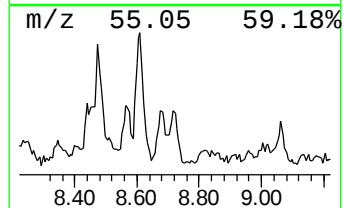
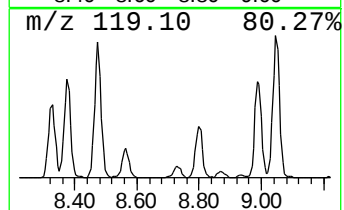
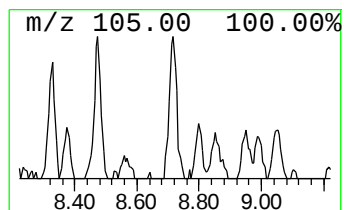
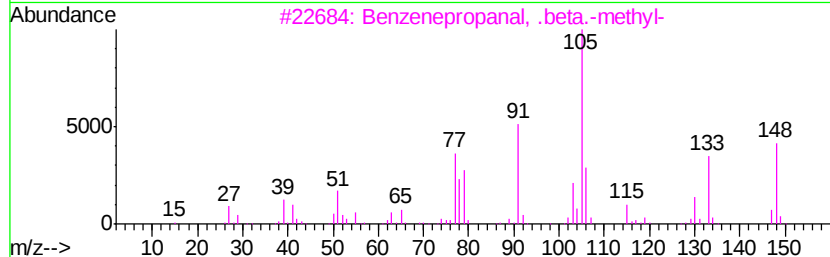
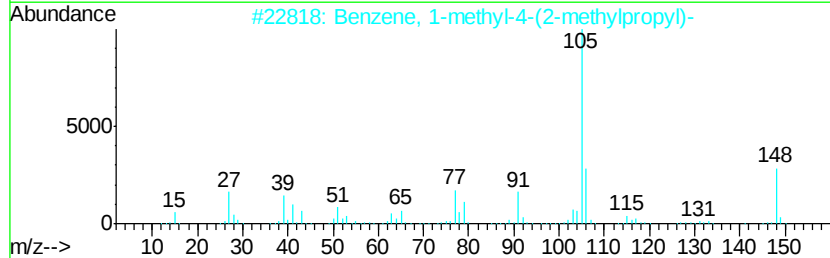
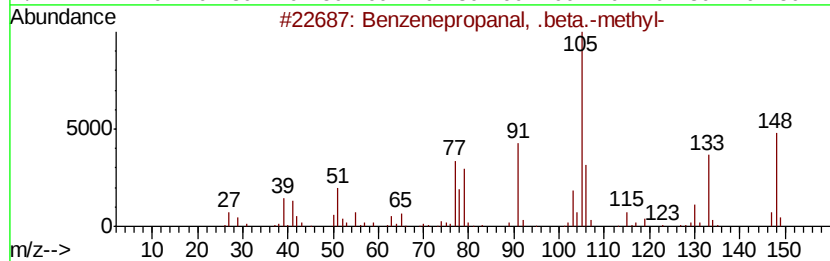
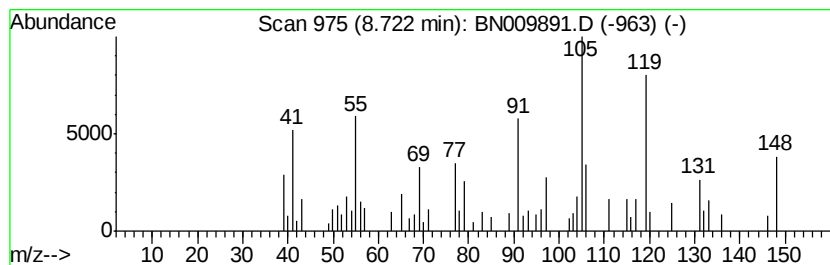
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 16 unknown-01 Concentration Rank 39

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 8.72 | 3.90 ng/ul | 163258 | 1,4-Dichlorobenzene-d4 | 7.39 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | Benzenepropanal, .beta.-methyl-     | 148 | C10H12O  | 016251-77-7 | 45   |
| 2    |    | Benzene, 1-methyl-4-(2-methylpro... | 148 | C11H16   | 005161-04-6 | 35   |
| 3    |    | Benzenepropanal, .beta.-methyl-     | 148 | C10H12O  | 016251-77-7 | 30   |
| 4    |    | Tricyclo[4.2.1.1(2,5)]deca-3,7-d... | 176 | C11H12O2 | 070220-88-1 | 27   |
| 5    |    | Benzene, 1-methyl-4-butyl           | 148 | C11H16   | 001595-05-7 | 27   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_N\DATA\BN022420\  
Data File : BN009891.D  
Acq On : 25 Feb 2020 09:04  
Operator : CG/JU  
Sample : L1424-22DL 10X  
Misc :  
ALS Vial : 29 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
C5B30DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN021220MA.M  
Quant Title : SVOA CALIBRATION

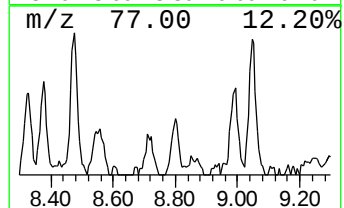
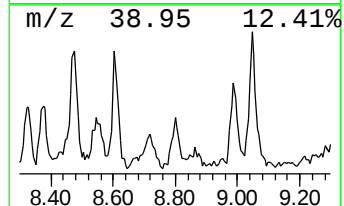
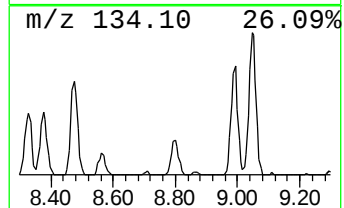
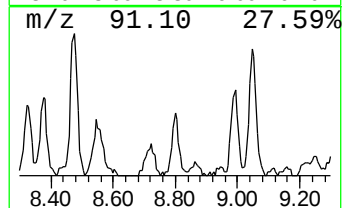
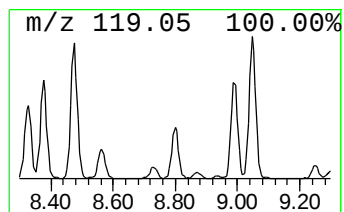
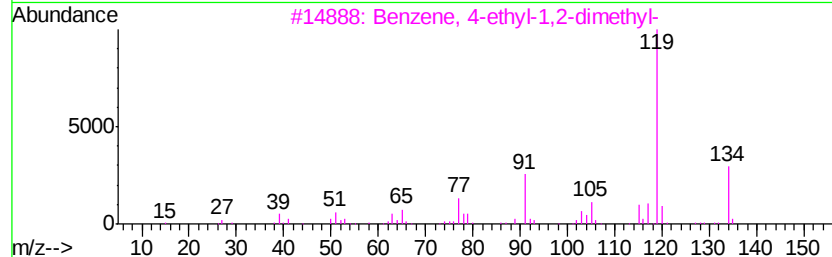
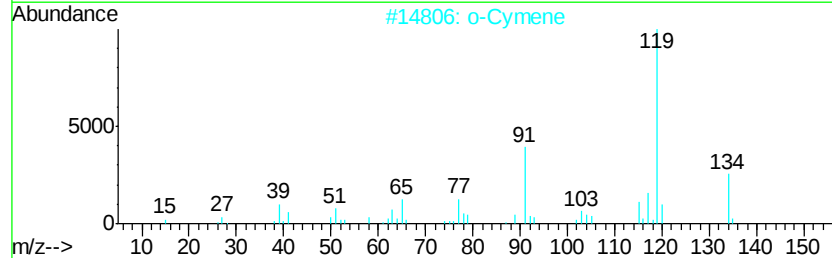
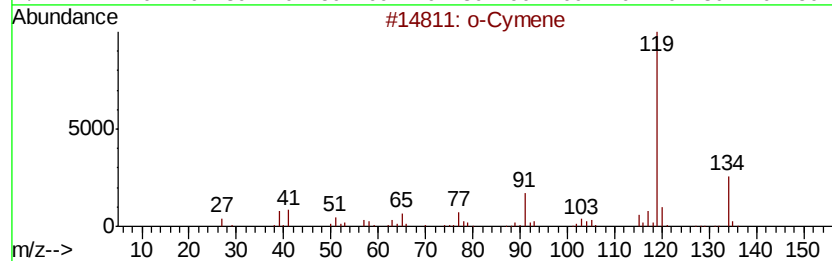
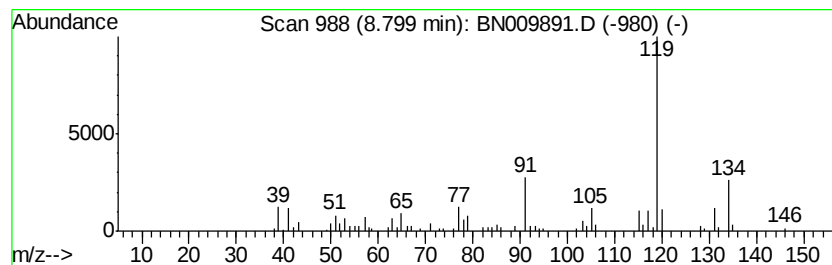
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 17 o-Cymene Concentration Rank 59

| R.T. | EstConc    | Area   | Relative to ISTD | R.T.  |
|------|------------|--------|------------------|-------|
| 8.80 | 2.46 ng/ul | 173405 | Napthalene-d8    | 10.13 |

| Hit# | of | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------|-----|---------|-------------|------|
| 1    | 5  | o-Cymene                       | 134 | C10H14  | 000527-84-4 | 94   |
| 2    |    | o-Cymene                       | 134 | C10H14  | 000527-84-4 | 93   |
| 3    |    | Benzene, 4-ethyl-1,2-dimethyl- | 134 | C10H14  | 000934-80-5 | 93   |
| 4    |    | p-Cymene                       | 134 | C10H14  | 000099-87-6 | 93   |
| 5    |    | Benzene, 1-ethyl-2,3-dimethyl- | 134 | C10H14  | 000933-98-2 | 93   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_N\DATA\BN022420\  
Data File : BN009891.D  
Acq On : 25 Feb 2020 09:04  
Operator : CG/JU  
Sample : L1424-22DL 10X  
Misc :  
ALS Vial : 29 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
C5B30DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN021220MA.M  
Quant Title : SVOA CALIBRATION

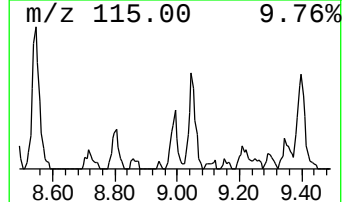
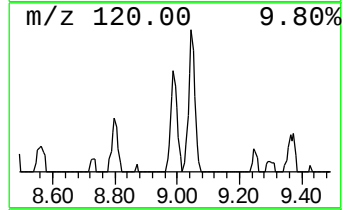
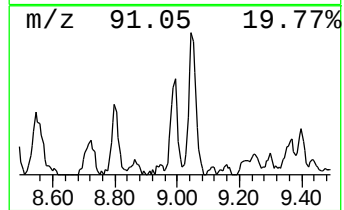
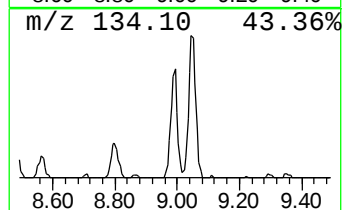
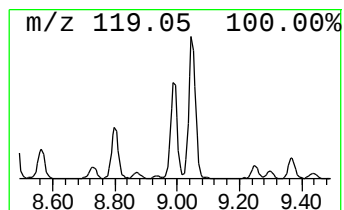
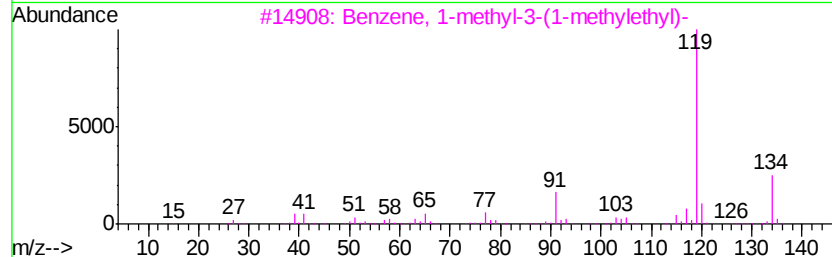
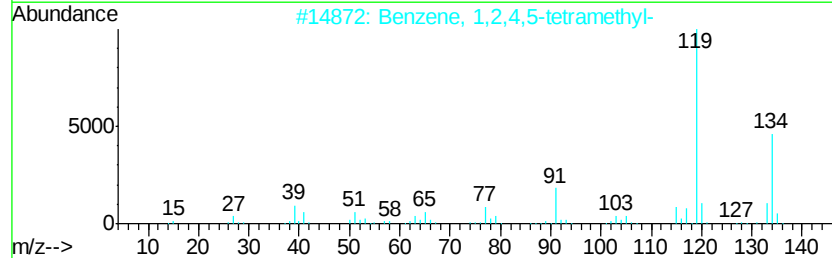
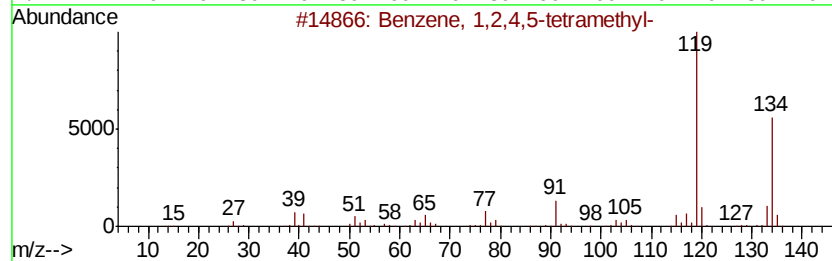
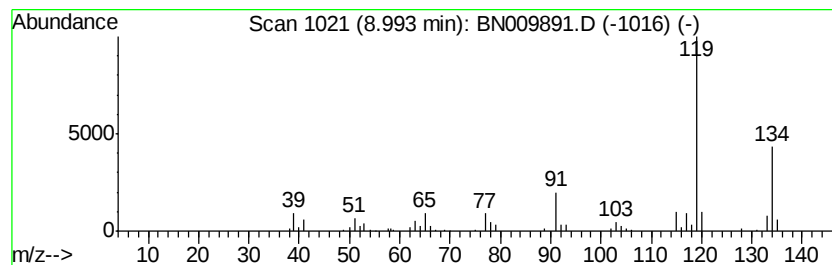
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TIC Integration Parameters: LSCINT.P

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Peak Number 18 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 46

| R.T. | EstConc    | Area   | Relative to ISTD | R.T.  |
|------|------------|--------|------------------|-------|
| 8.99 | 3.36 ng/ul | 237154 | Naphthalene-d8   | 10.13 |

| Hit# | of | Tentative ID                         | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1,2,4,5-tetramethyl-        | 134 | C10H14  | 000095-93-2 | 95   |
| 2    |    | Benzene, 1,2,4,5-tetramethyl-        | 134 | C10H14  | 000095-93-2 | 94   |
| 3    |    | Benzene, 1-methyl-3-(1-methylethyl)- | 134 | C10H14  | 000535-77-3 | 91   |
| 4    |    | o-Cymene                             | 134 | C10H14  | 000527-84-4 | 91   |
| 5    |    | 1,3-Cyclopentadiene, 1,2,3,4-tet...  | 134 | C10H14  | 076089-59-3 | 91   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022420\  
Data File : BN009891.D  
Acq On : 25 Feb 2020 09:04  
Operator : CG/JU  
Sample : L1424-22DL 10X  
Misc :  
ALS Vial : 29 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
C5B30DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN021220MA.M  
Quant Title : SVOA CALIBRATION

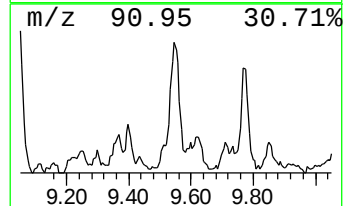
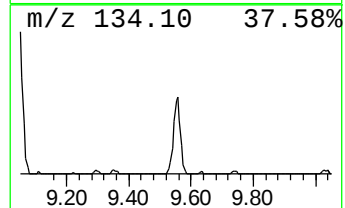
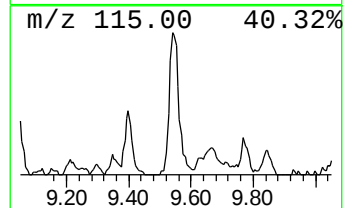
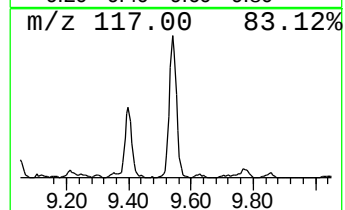
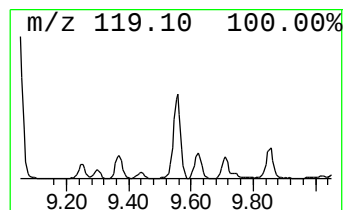
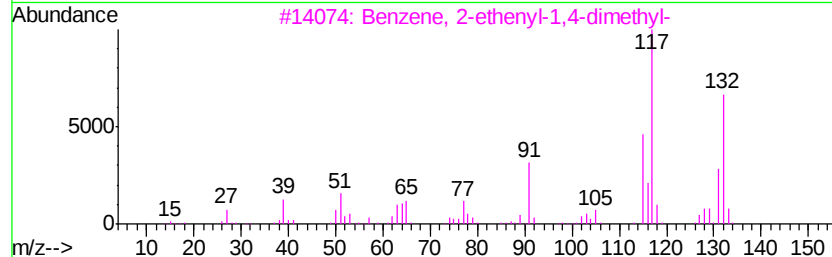
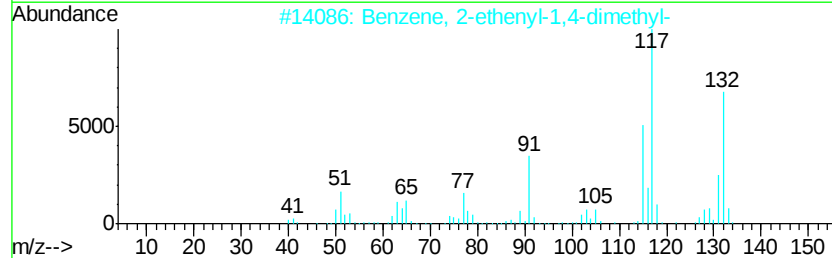
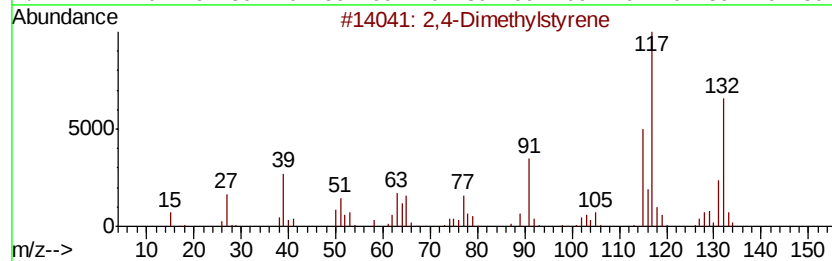
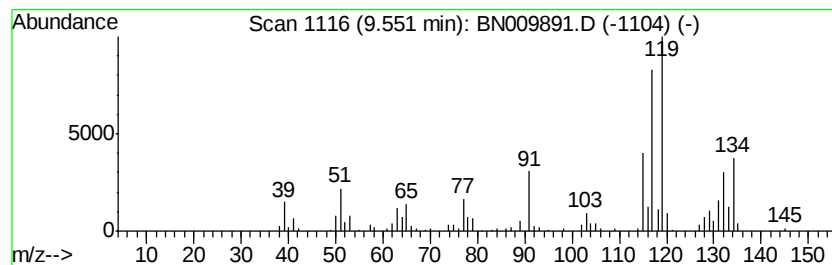
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 20 2,4-Dimethylstyrene Concentration Rank 19

| R.T. | EstConc    | Area   | Relative to ISTD | R.T.  |
|------|------------|--------|------------------|-------|
| 9.55 | 7.06 ng/ul | 498071 | Naphthalene-d8   | 10.13 |

| Hit# | of | Tentative ID                      | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------------------|-----|---------|-------------|------|
| 1    | 5  | 2,4-Dimethylstvrene               | 132 | C10H12  | 002234-20-0 | 91   |
| 2    |    | Benzene, 2-ethenyl-1,4-dimethyl-  | 132 | C10H12  | 002039-89-6 | 90   |
| 3    |    | Benzene, 2-ethenyl-1,4-dimethyl-  | 132 | C10H12  | 002039-89-6 | 90   |
| 4    |    | Benzene, 2-butenyl-               | 132 | C10H12  | 001560-06-1 | 86   |
| 5    |    | Benzene, 1-methyl-4-(2-propenyl)- | 132 | C10H12  | 003333-13-9 | 83   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_N\DATA\BN022420\  
Data File : BN009891.D  
Acq On : 25 Feb 2020 09:04  
Operator : CG/JU  
Sample : L1424-22DL 10X  
Misc :  
ALS Vial : 29 Sample Multiplier: 1

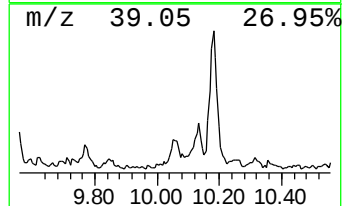
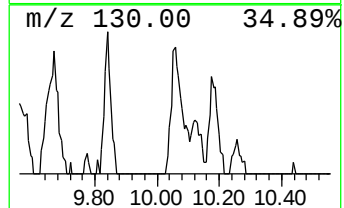
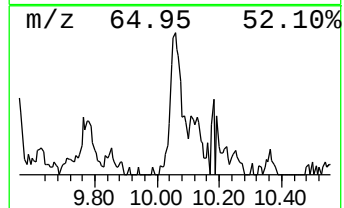
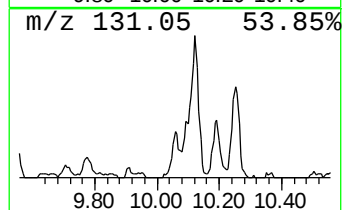
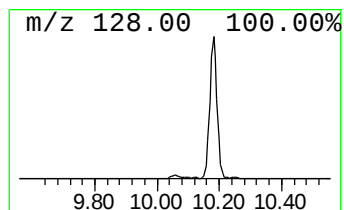
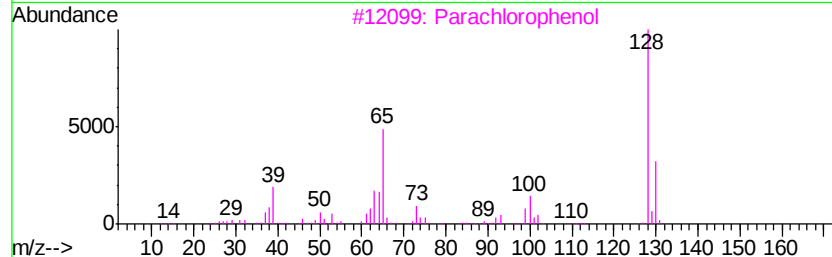
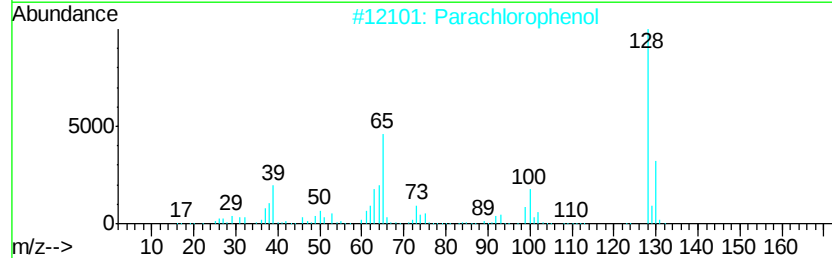
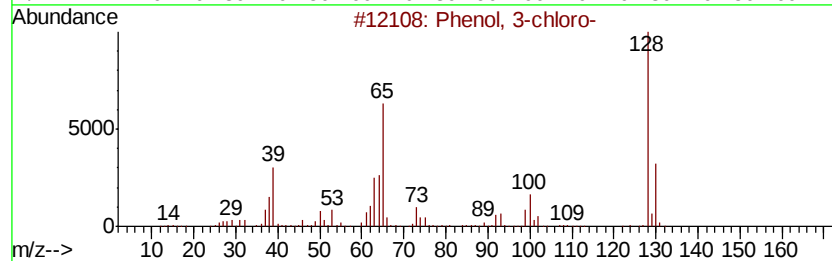
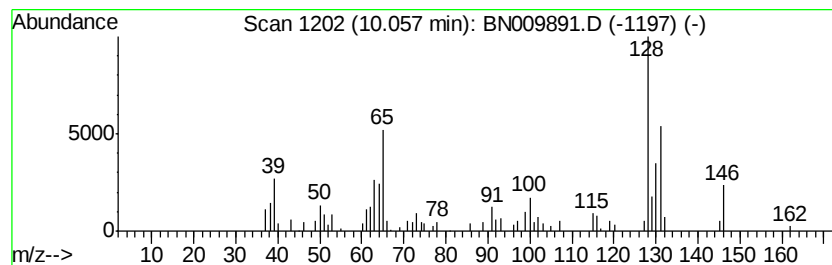
Instrument :  
BNA\_N  
ClientSampleId :  
C5B30DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN021220MA.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 21 Phenol, 3-chloro- Concentration Rank 61

| R.T.    | EstConc           | Area         | Relative to ISTD | R.T.      |
|---------|-------------------|--------------|------------------|-----------|
| 10.06   | 2.44 ng/ul        | 171957       | Naphthalene-d8   | 10.13     |
| Hit# of | 5                 | Tentative ID | MW MolForm       | CAS# Qual |
| 1       | Phenol, 3-chloro- | 128 C6H5ClO  | 000108-43-0      | 97        |
| 2       | Parachlorophenol  | 128 C6H5ClO  | 000106-48-9      | 96        |
| 3       | Parachlorophenol  | 128 C6H5ClO  | 000106-48-9      | 95        |
| 4       | Phenol, 3-chloro- | 128 C6H5ClO  | 000108-43-0      | 95        |
| 5       | Parachlorophenol  | 128 C6H5ClO  | 000106-48-9      | 95        |



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022420\  
Data File : BN009891.D  
Acq On : 25 Feb 2020 09:04  
Operator : CG/JU  
Sample : L1424-22DL 10X  
Misc :  
ALS Vial : 29 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
C5B30DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN021220MA.M  
Quant Title : SVOA CALIBRATION

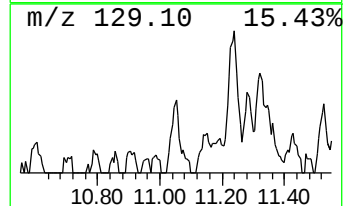
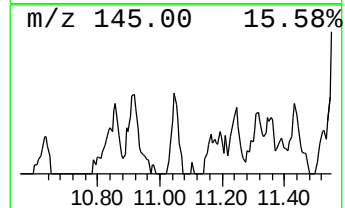
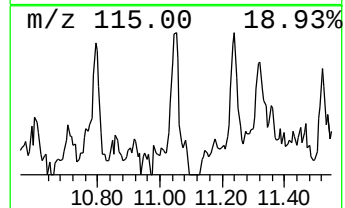
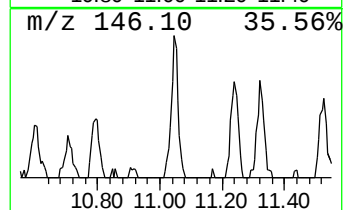
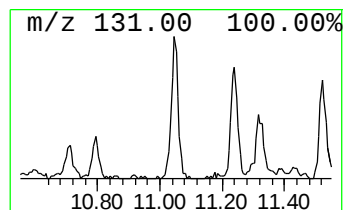
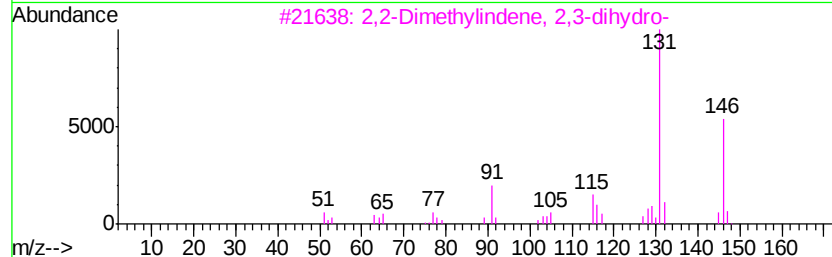
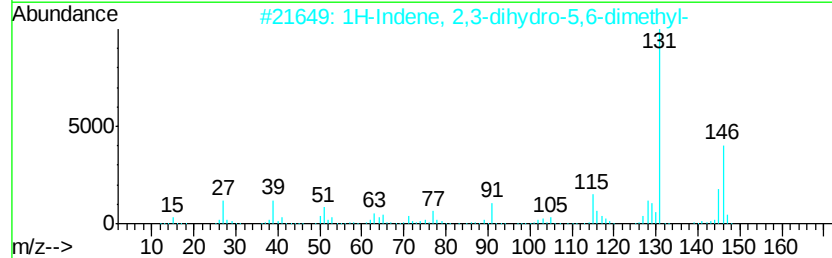
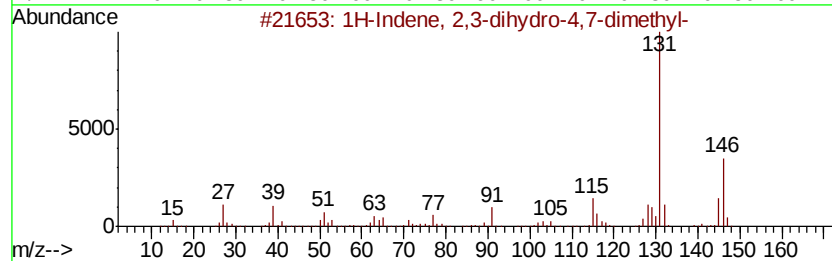
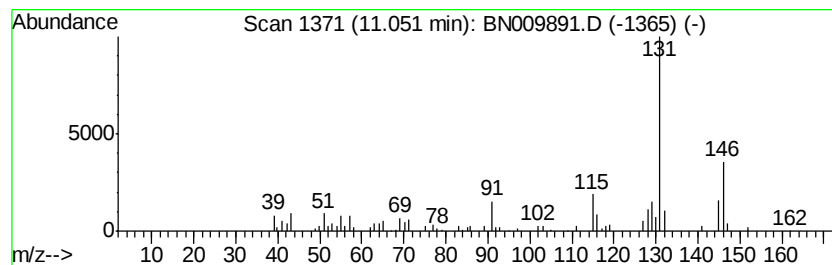
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 22 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 64

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 11.05 | 2.11 ng/ul | 148984 | Napthalene-d8    | 10.13 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | 1H-Indene, 2,3-dihydro-4,7-dimet... | 146 | C11H14  | 006682-71-9 | 94   |
| 2    |    |   | 1H-Indene, 2,3-dihydro-5,6-dimet... | 146 | C11H14  | 001075-22-5 | 91   |
| 3    |    |   | 2,2-Dimethylindene, 2,3-dihydro-    | 146 | C11H14  | 020836-11-7 | 91   |
| 4    |    |   | Benzene, (1-methyl-1-butenyl)-      | 146 | C11H14  | 053172-84-2 | 91   |
| 5    |    |   | 1H-Indene, 2,3-dihydro-1,2-dimet... | 146 | C11H14  | 017057-82-8 | 91   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_N\DATA\BN022420\  
Data File : BN009891.D  
Acq On : 25 Feb 2020 09:04  
Operator : CG/JU  
Sample : L1424-22DL 10X  
Misc :  
ALS Vial : 29 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
C5B30DL

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Quant Title : SVOA CALIBRATION

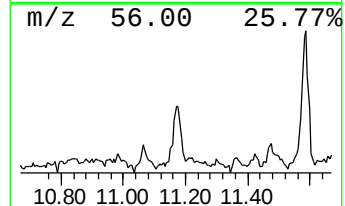
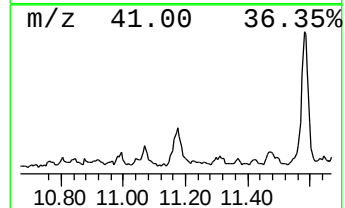
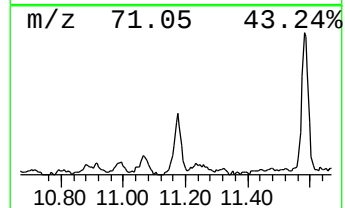
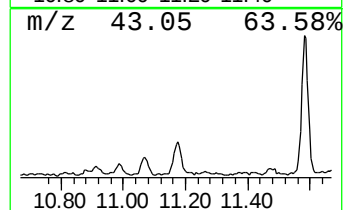
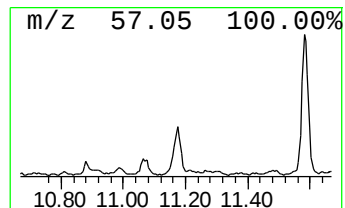
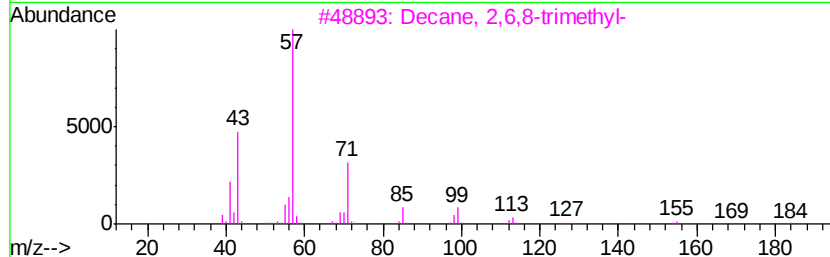
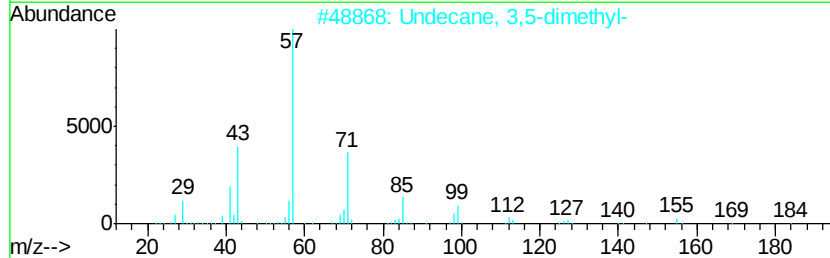
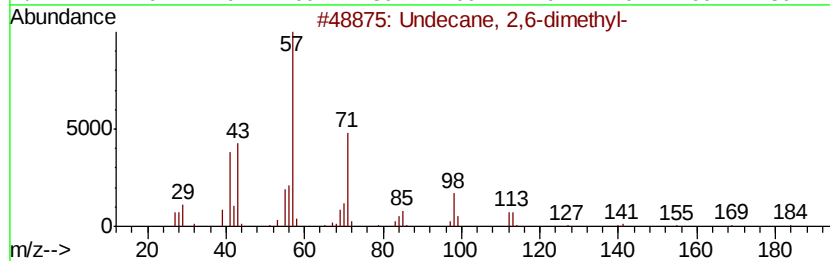
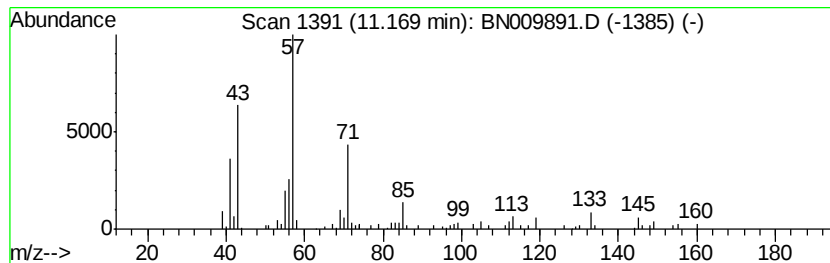
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 23 (DEL) Alkane: Straight-Chai... Concentration Rank 57

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 11.17 | 2.57 ng/ul | 181151 | Naphthalene-d8   | 10.13 |

| Hit# | of | Tentative ID             | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------|-----|---------|-------------|------|
| 1    | 5  | Undecane, 2,6-dimethyl-  | 184 | C13H28  | 017301-23-4 | 53   |
| 2    |    | Undecane, 3,5-dimethyl-  | 184 | C13H28  | 017312-81-1 | 50   |
| 3    |    | Decane, 2,6,8-trimethyl- | 184 | C13H28  | 062108-26-3 | 50   |
| 4    |    | Dodecane, 3-methyl-      | 184 | C13H28  | 017312-57-1 | 50   |
| 5    |    | Butane, 2,2-dimethyl-    | 86  | C6H14   | 000075-83-2 | 47   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022420\  
Data File : BN009891.D  
Acq On : 25 Feb 2020 09:04  
Operator : CG/JU  
Sample : L1424-22DL 10X  
Misc :  
ALS Vial : 29 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
C5B30DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN021220MA.M  
Quant Title : SVOA CALIBRATION

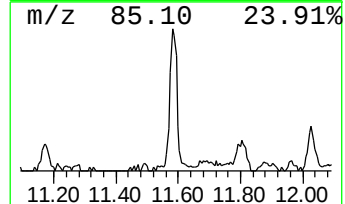
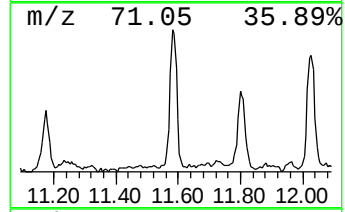
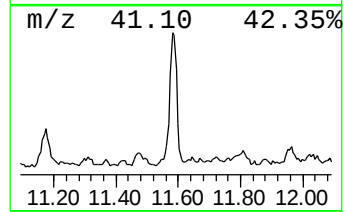
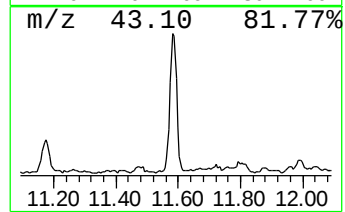
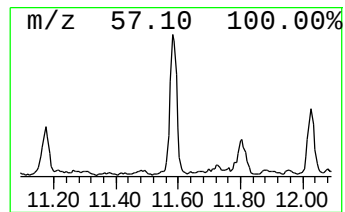
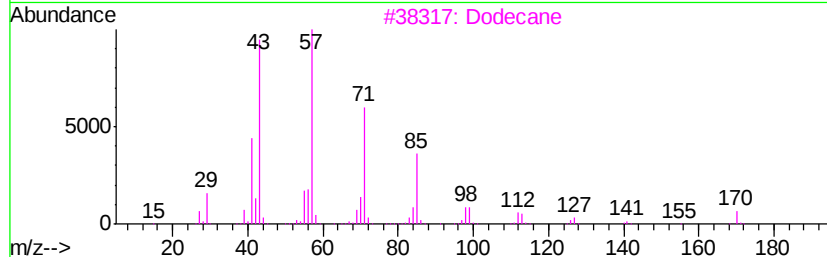
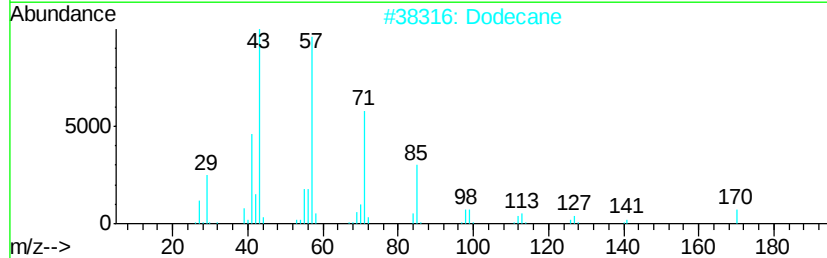
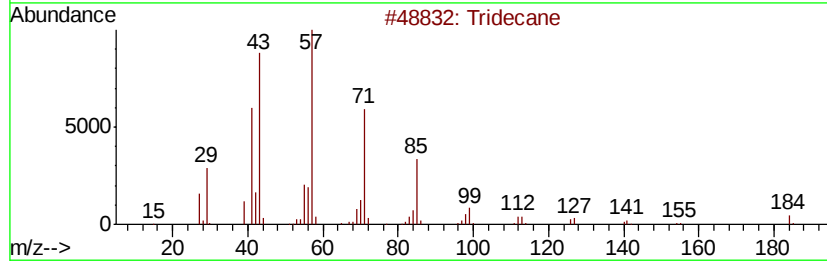
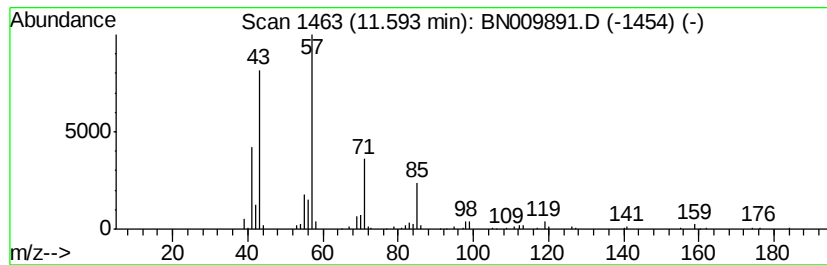
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 24 (DEL) Alkane: Straight-Chai... Concentration Rank 27

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 11.59 | 5.21 ng/ul | 367830 | Napthalene-d8    | 10.13 |

| Hit# | of | 5 | Tentative ID   | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------|-----|---------|-------------|------|
| 1    |    |   | Tridecane      | 184 | C13H28  | 000629-50-5 | 90   |
| 2    |    |   | Dodecane       | 170 | C12H26  | 000112-40-3 | 86   |
| 3    |    |   | Dodecane       | 170 | C12H26  | 000112-40-3 | 78   |
| 4    |    |   | Hexadecane     | 226 | C16H34  | 000544-76-3 | 72   |
| 5    |    |   | 1-Iodoundecane | 282 | C11H23I | 004282-44-4 | 72   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022420\  
Data File : BN009891.D  
Acq On : 25 Feb 2020 09:04  
Operator : CG/JU  
Sample : L1424-22DL 10X  
Misc :  
ALS Vial : 29 Sample Multiplier: 1

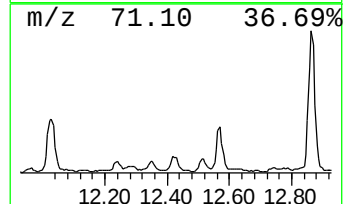
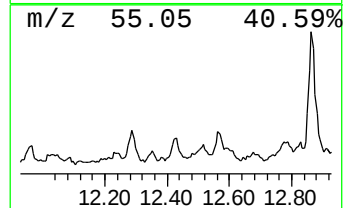
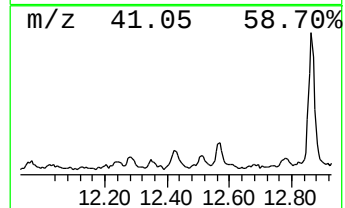
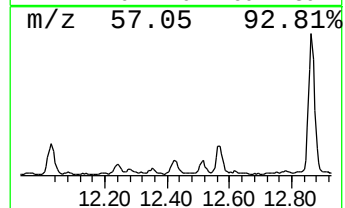
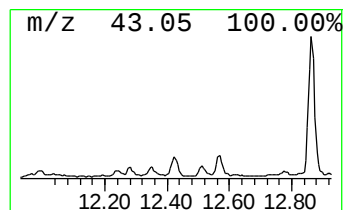
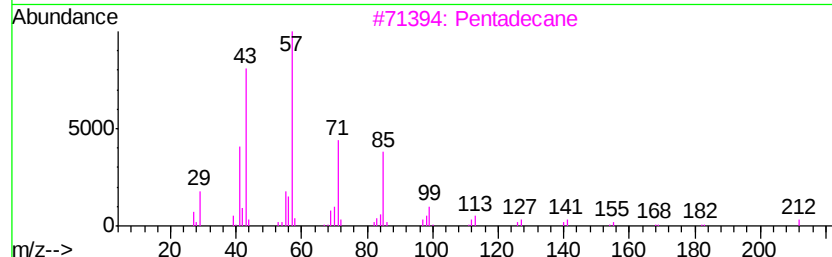
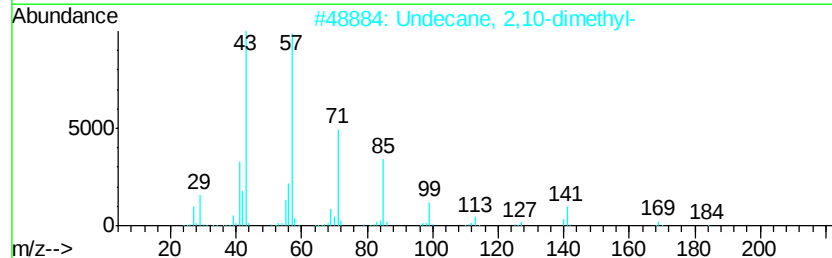
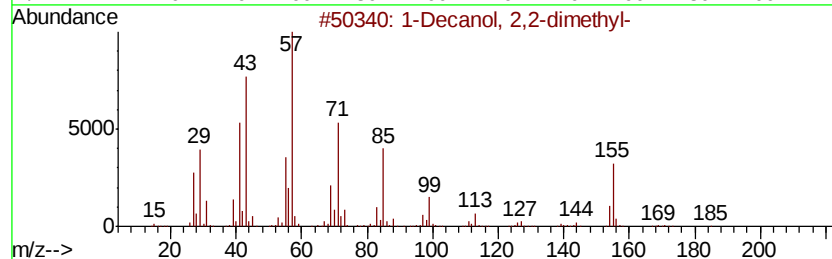
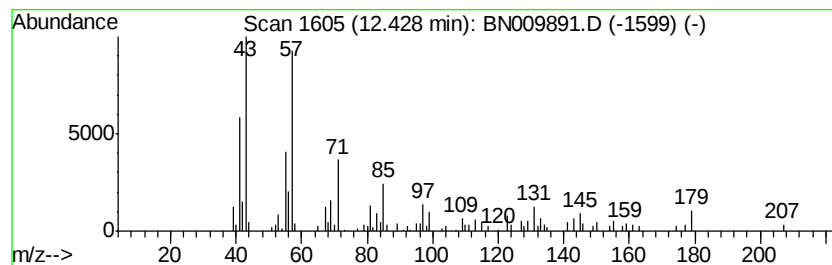
Instrument :  
BNA\_N  
ClientSampleId :  
C5B30DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN021220MA.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 25 1-Decanol, 2,2-dimethyl- Concentration Rank 65

| R.T.    | EstConc                             | Area         | Relative to ISTD | R.T.      |              |      |
|---------|-------------------------------------|--------------|------------------|-----------|--------------|------|
| 12.43   | 2.06 ng/ul                          | 165623       | Acenaphthene-d10 | 14.02     |              |      |
| Hit# of | 5                                   | Tentative ID | MW               | MolForm   | CAS#         | Qual |
| 1       | 1-Decanol, 2,2-dimethyl-            |              | 186              | C12H26O   | 002370-15-2  | 50   |
| 2       | Undecane, 2,10-dimethyl-            |              | 184              | C13H28    | 017301-27-8  | 50   |
| 3       | Pentadecane                         |              | 212              | C15H32    | 000629-62-9  | 50   |
| 4       | Sulfurous acid, butyl heptadecyl... |              | 376              | C21H44O3S | 1000309-18-4 | 50   |
| 5       | 1-Iodo-2-methylundecane             |              | 296              | C12H25I   | 073105-67-6  | 50   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022420\  
Data File : BN009891.D  
Acq On : 25 Feb 2020 09:04  
Operator : CG/JU  
Sample : L1424-22DL 10X  
Misc :  
ALS Vial : 29 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
C5B30DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN021220MA.M  
Quant Title : SVOA CALIBRATION

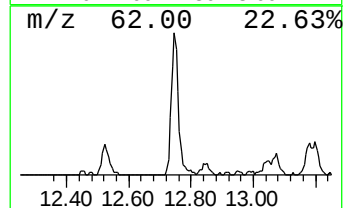
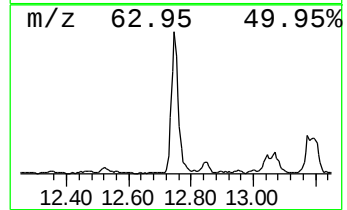
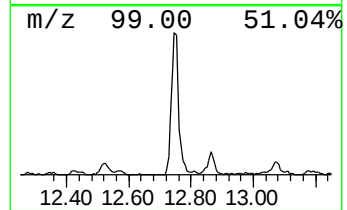
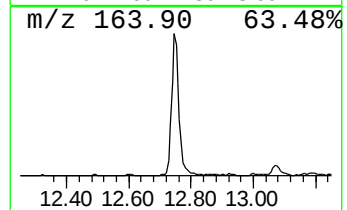
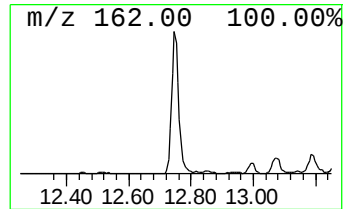
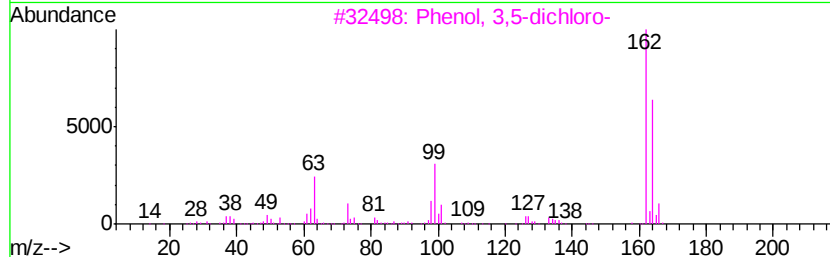
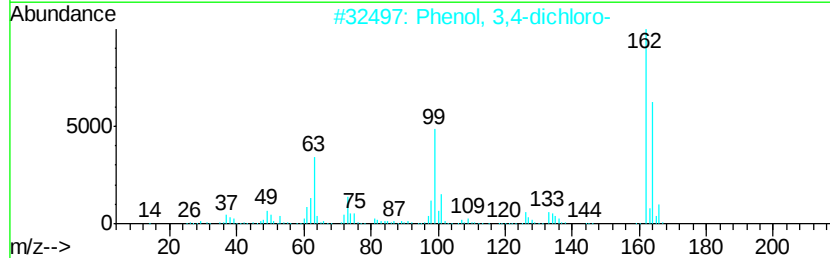
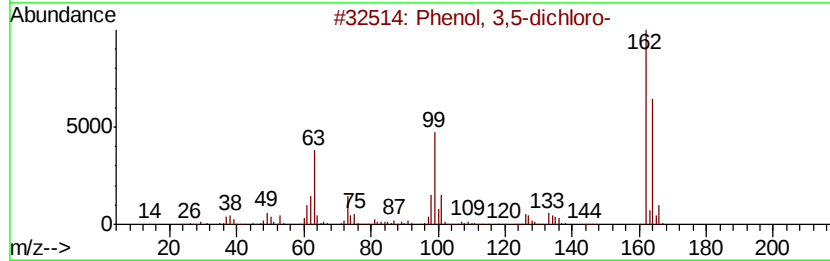
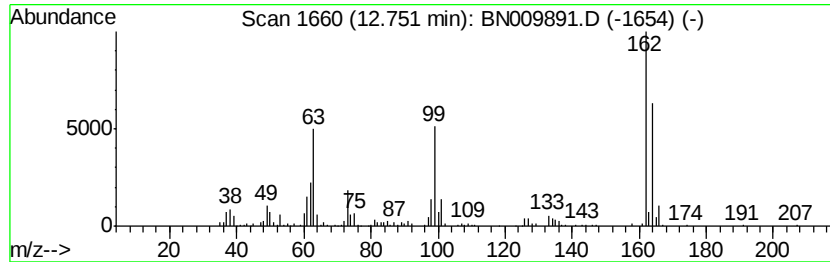
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 26 Phenol, 3,5-dichloro- Concentration Rank 9

| R.T.  | EstConc     | Area   | Relative to ISTD | R.T.  |
|-------|-------------|--------|------------------|-------|
| 12.75 | 10.81 ng/ul | 870890 | Acenaphthene-d10 | 14.02 |

| Hit# | of                    | 5 | Tentative ID | MW  | MolForm  | CAS#        | Qual |
|------|-----------------------|---|--------------|-----|----------|-------------|------|
| 1    | Phenol, 3,5-dichloro- |   |              | 162 | C6H4Cl2O | 000591-35-5 | 97   |
| 2    | Phenol, 3,4-dichloro- |   |              | 162 | C6H4Cl2O | 000095-77-2 | 97   |
| 3    | Phenol, 3,5-dichloro- |   |              | 162 | C6H4Cl2O | 000591-35-5 | 96   |
| 4    | Phenol, 3,4-dichloro- |   |              | 162 | C6H4Cl2O | 000095-77-2 | 95   |
| 5    | Phenol, 3,4-dichloro- |   |              | 162 | C6H4Cl2O | 000095-77-2 | 94   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022420\  
Data File : BN009891.D  
Acq On : 25 Feb 2020 09:04  
Operator : CG/JU  
Sample : L1424-22DL 10X  
Misc :  
ALS Vial : 29 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
C5B30DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN021220MA.M  
Quant Title : SVOA CALIBRATION

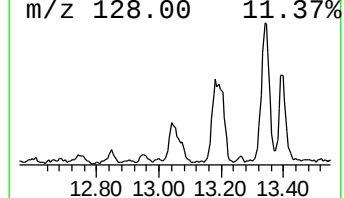
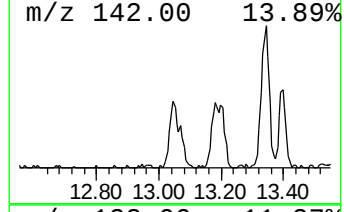
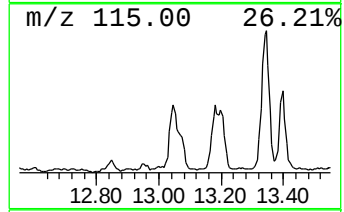
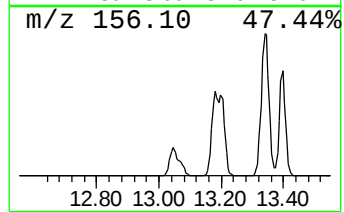
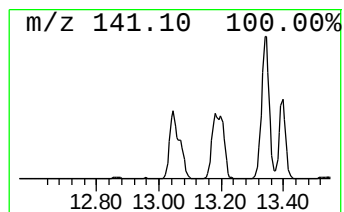
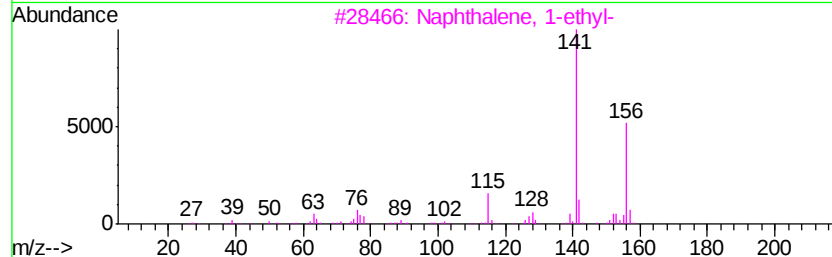
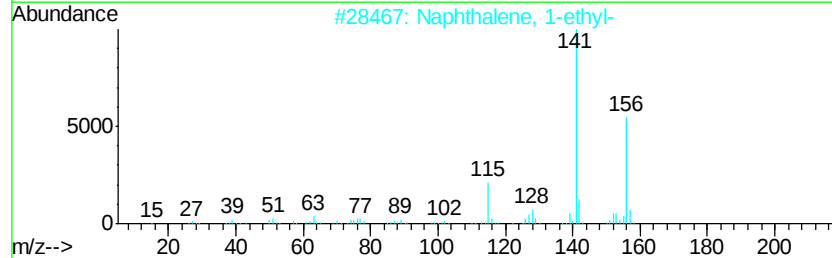
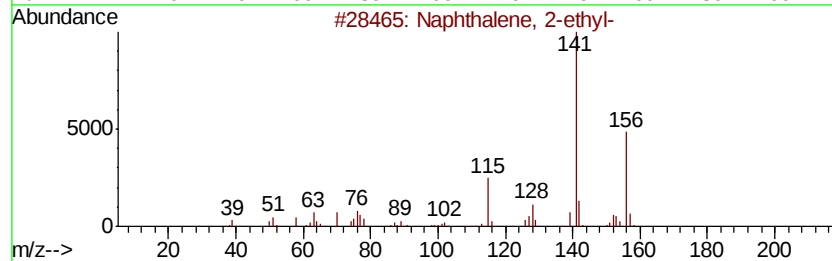
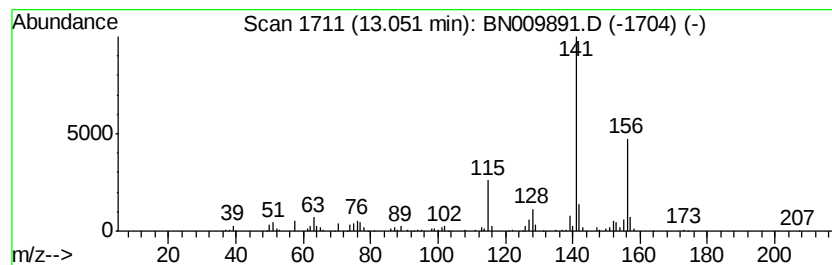
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 27 Naphthalene, 2-ethyl- Concentration Rank 35

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 13.05 | 4.70 ng/ul | 378407 | Acenaphthene-d10 | 14.02 |

| Hit# | of | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 2-ethyl- | 156 | C12H12  | 000939-27-5 | 96   |
| 2    |    | Naphthalene, 1-ethyl- | 156 | C12H12  | 001127-76-0 | 96   |
| 3    |    | Naphthalene, 1-ethyl- | 156 | C12H12  | 001127-76-0 | 95   |
| 4    |    | Naphthalene, 2-ethyl- | 156 | C12H12  | 000939-27-5 | 95   |
| 5    |    | Naphthalene, 2-ethyl- | 156 | C12H12  | 000939-27-5 | 95   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022420\  
Data File : BN009891.D  
Acq On : 25 Feb 2020 09:04  
Operator : CG/JU  
Sample : L1424-22DL 10X  
Misc :  
ALS Vial : 29 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
C5B30DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN021220MA.M  
Quant Title : SVOA CALIBRATION

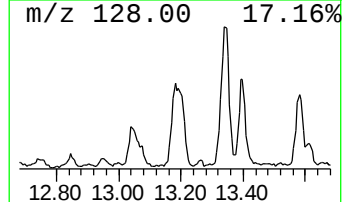
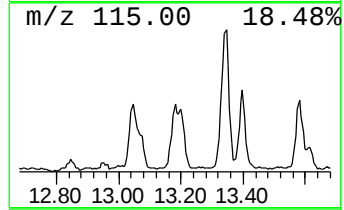
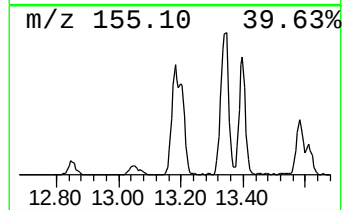
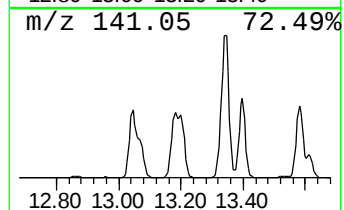
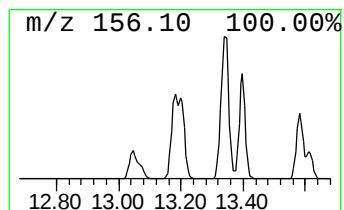
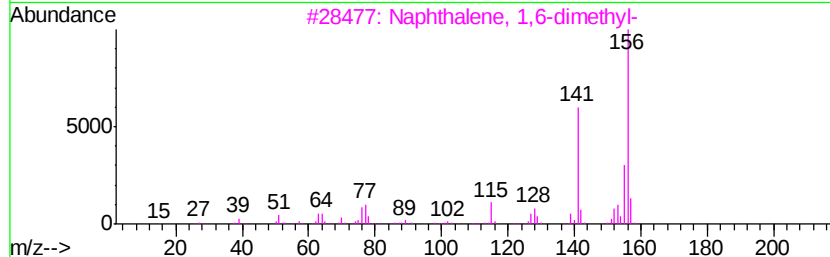
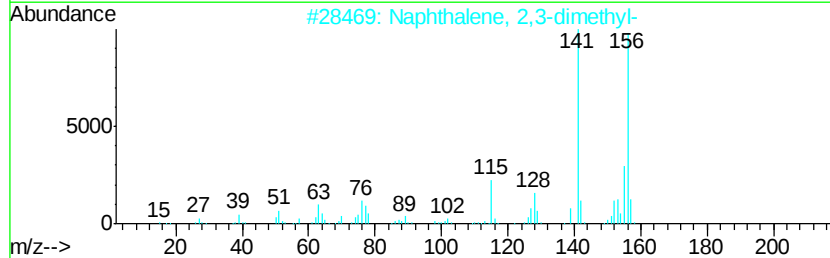
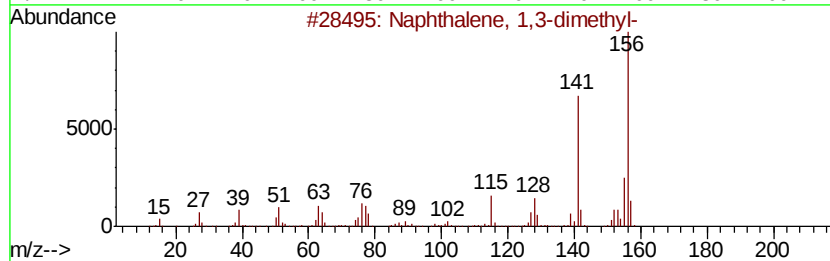
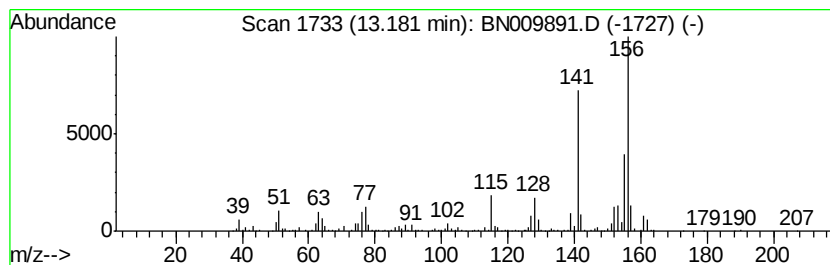
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 28 Naphthalene, 1,3-dimethyl- Concentration Rank 3

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 13.18 | 18.91 ng/ul | 1522830 | Acenaphthene-d10 | 14.02 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 1,3-dimethyl- | 156 | C12H12  | 000575-41-7 | 96   |
| 2    |    | Naphthalene, 2,3-dimethyl- | 156 | C12H12  | 000581-40-8 | 96   |
| 3    |    | Naphthalene, 1,6-dimethyl- | 156 | C12H12  | 000575-43-9 | 96   |
| 4    |    | Naphthalene, 1,3-dimethyl- | 156 | C12H12  | 000575-41-7 | 96   |
| 5    |    | Naphthalene, 1,6-dimethyl- | 156 | C12H12  | 000575-43-9 | 96   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022420\  
Data File : BN009891.D  
Acq On : 25 Feb 2020 09:04  
Operator : CG/JU  
Sample : L1424-22DL 10X  
Misc :  
ALS Vial : 29 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
C5B30DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN021220MA.M  
Quant Title : SVOA CALIBRATION

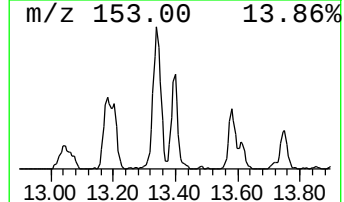
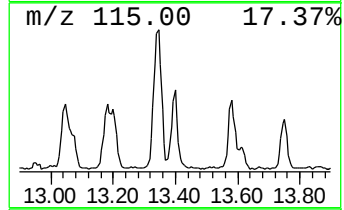
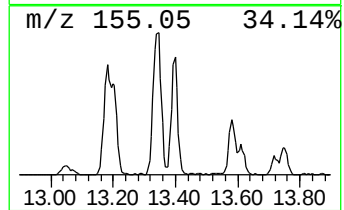
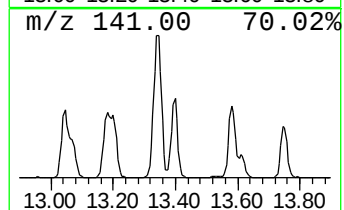
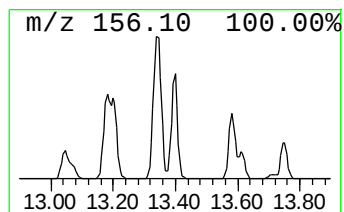
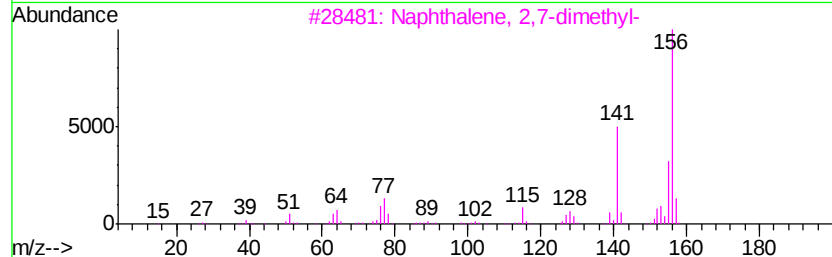
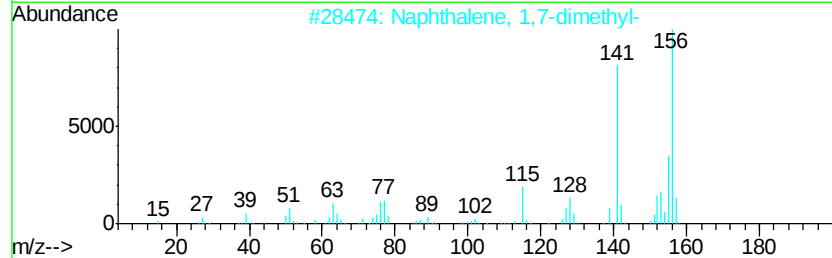
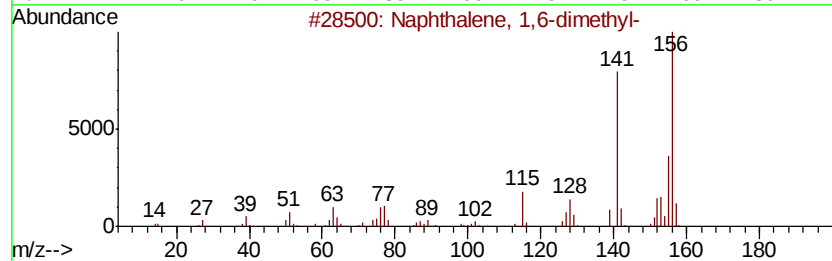
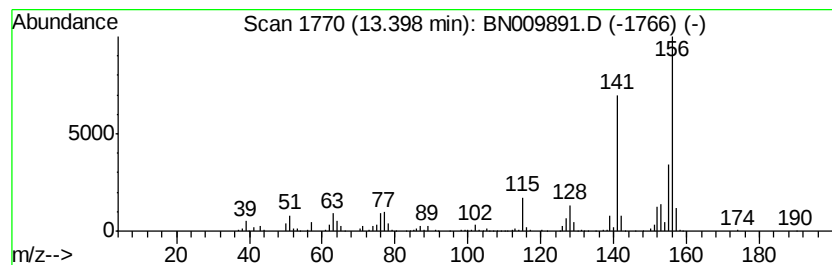
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 30 Naphthalene, 1,6-dimethyl- Concentration Rank 6

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 13.40 | 12.43 ng/ul | 1001040 | Acenaphthene-d10 | 14.02 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 1,6-dimethyl- | 156 | C12H12  | 000575-43-9 | 98   |
| 2    |    | Naphthalene, 1,7-dimethyl- | 156 | C12H12  | 000575-37-1 | 98   |
| 3    |    | Naphthalene, 2,7-dimethyl- | 156 | C12H12  | 000582-16-1 | 97   |
| 4    |    | Naphthalene, 2,6-dimethyl- | 156 | C12H12  | 000581-42-0 | 97   |
| 5    |    | Naphthalene, 1,5-dimethyl- | 156 | C12H12  | 000571-61-9 | 97   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022420\  
Data File : BN009891.D  
Acq On : 25 Feb 2020 09:04  
Operator : CG/JU  
Sample : L1424-22DL 10X  
Misc :  
ALS Vial : 29 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
C5B30DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN021220MA.M  
Quant Title : SVOA CALIBRATION

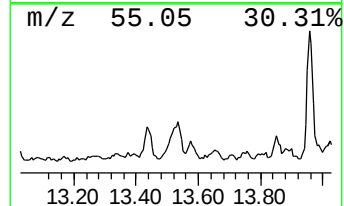
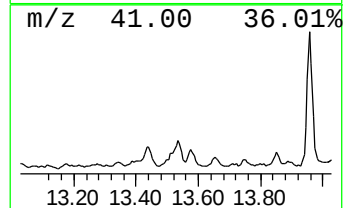
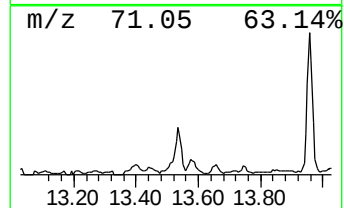
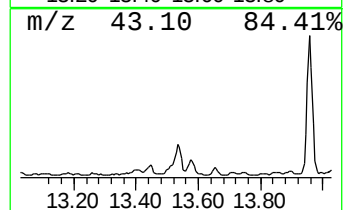
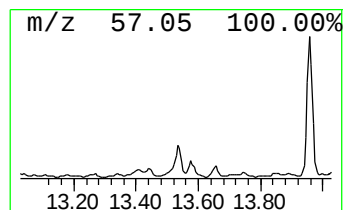
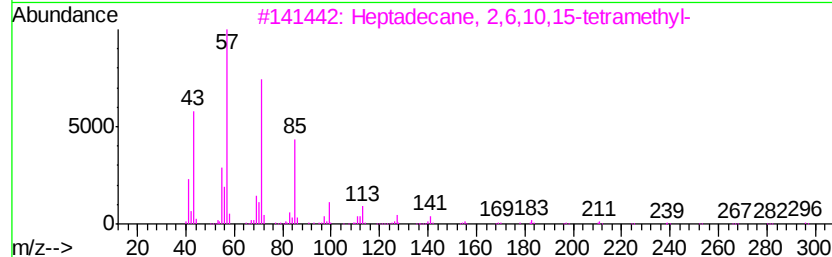
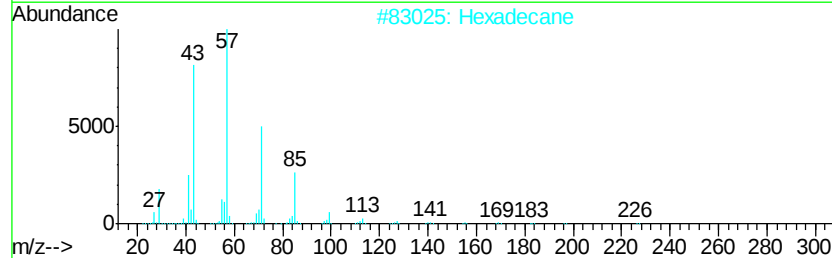
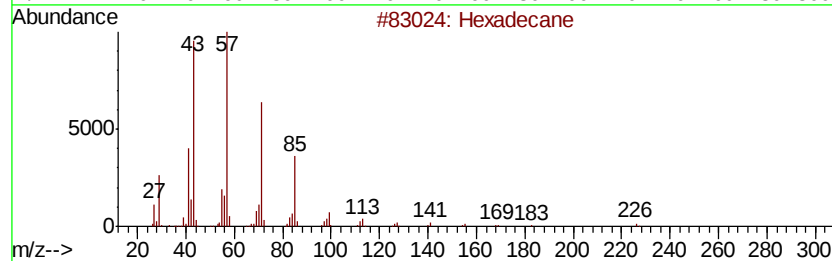
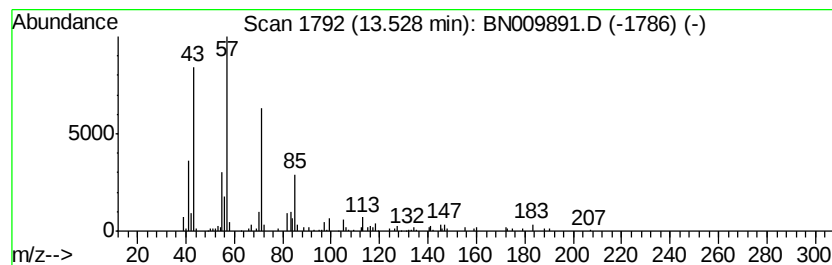
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 31 (DEL) Alkane: Straight-Chai... Concentration Rank 36

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 13.53 | 4.21 ng/ul | 339104 | Acenaphthene-d10 | 14.02 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Hexadecane                          | 226 | C16H34  | 000544-76-3 | 86   |
| 2    |    |   | Hexadecane                          | 226 | C16H34  | 000544-76-3 | 86   |
| 3    |    |   | Heptadecane, 2,6,10,15-tetramethyl- | 296 | C21H44  | 054833-48-6 | 86   |
| 4    |    |   | Heptadecane, 2,6,10,14-tetramethyl- | 296 | C21H44  | 018344-37-1 | 80   |
| 5    |    |   | Tridecane                           | 184 | C13H28  | 000629-50-5 | 78   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_N\DATA\BN022420\  
Data File : BN009891.D  
Acq On : 25 Feb 2020 09:04  
Operator : CG/JU  
Sample : L1424-22DL 10X  
Misc :  
ALS Vial : 29 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
C5B30DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN021220MA.M  
Quant Title : SVOA CALIBRATION

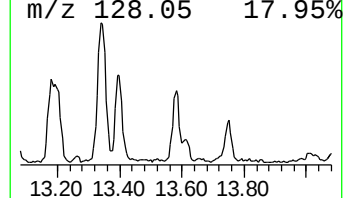
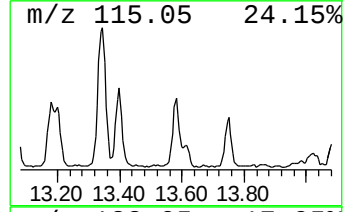
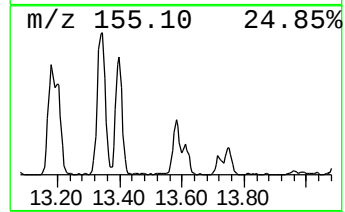
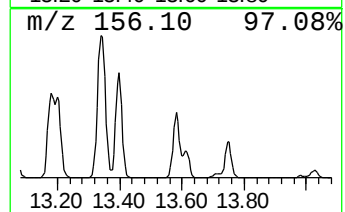
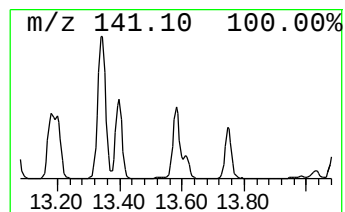
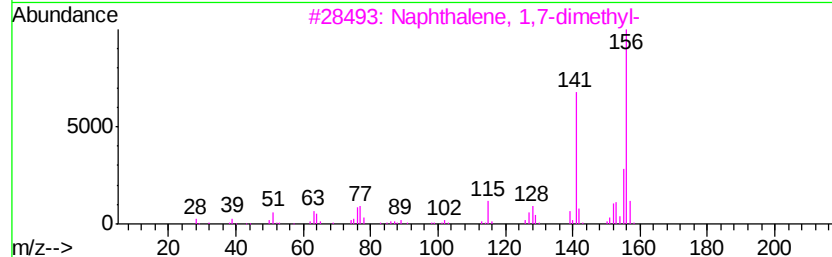
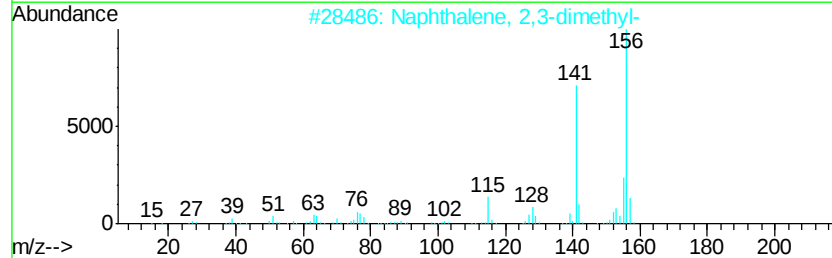
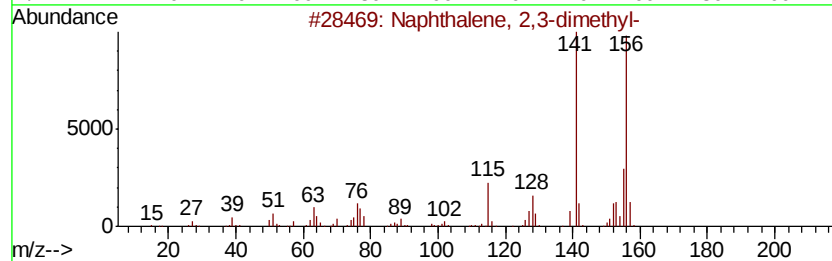
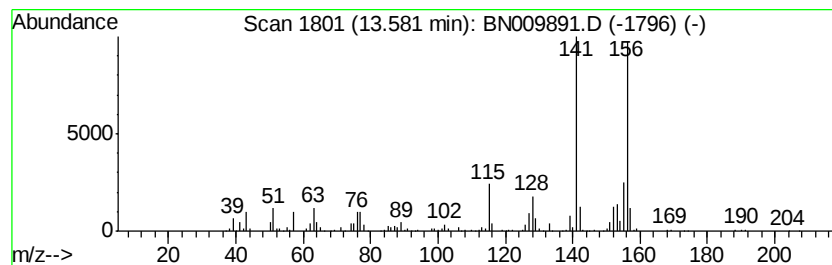
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 32 Naphthalene, 2,3-dimethyl- Concentration Rank 10

| R.T.  | EstConc     | Area   | Relative to ISTD | R.T.  |
|-------|-------------|--------|------------------|-------|
| 13.58 | 10.63 ng/ul | 855925 | Acenaphthene-d10 | 14.02 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 2,3-dimethyl- | 156 | C12H12  | 000581-40-8 | 98   |
| 2    |    | Naphthalene, 2,3-dimethyl- | 156 | C12H12  | 000581-40-8 | 97   |
| 3    |    | Naphthalene, 1,7-dimethyl- | 156 | C12H12  | 000575-37-1 | 97   |
| 4    |    | Naphthalene, 2,3-dimethyl- | 156 | C12H12  | 000581-40-8 | 97   |
| 5    |    | Naphthalene, 1,3-dimethyl- | 156 | C12H12  | 000575-41-7 | 97   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_N\DATA\BN022420\  
Data File : BN009891.D  
Acq On : 25 Feb 2020 09:04  
Operator : CG/JU  
Sample : L1424-22DL 10X  
Misc :  
ALS Vial : 29 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
C5B30DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN021220MA.M  
Quant Title : SVOA CALIBRATION

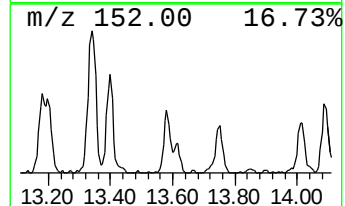
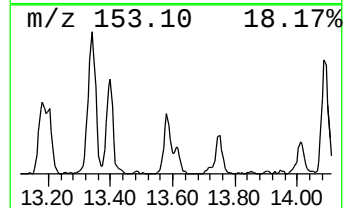
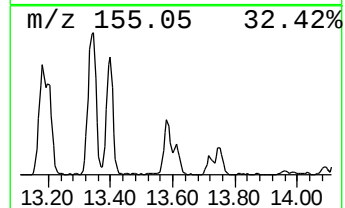
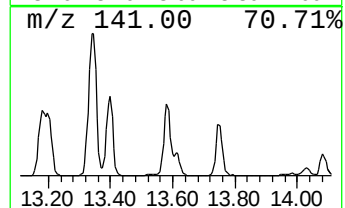
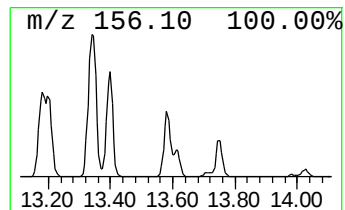
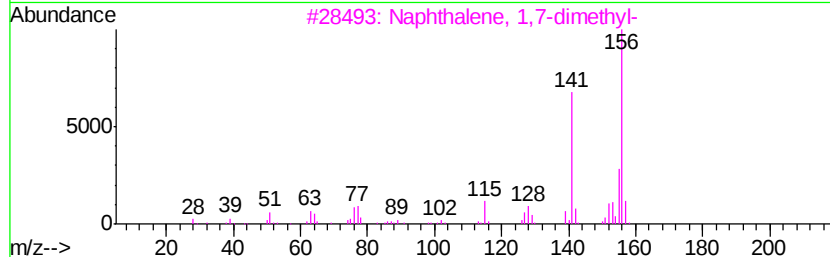
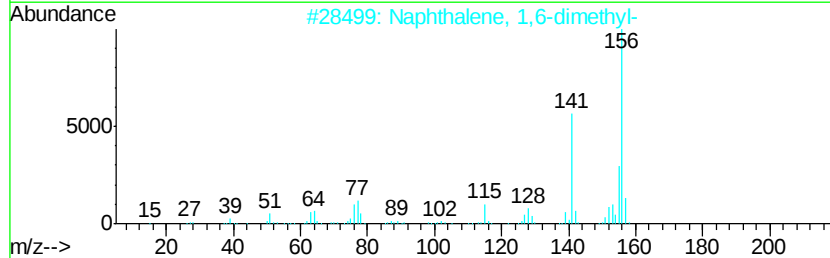
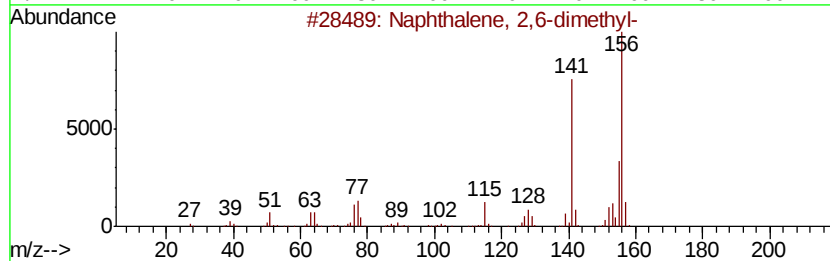
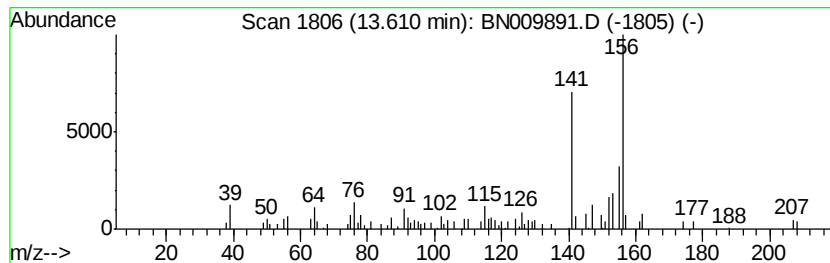
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 33 Naphthalene, 2,6-dimethyl- Concentration Rank 62

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 13.61 | 2.19 ng/ul | 176626 | Acenaphthene-d10 | 14.02 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 2,6-dimethyl- | 156 | C12H12  | 000581-42-0 | 83   |
| 2    |    | Naphthalene, 1,6-dimethyl- | 156 | C12H12  | 000575-43-9 | 80   |
| 3    |    | Naphthalene, 1,7-dimethyl- | 156 | C12H12  | 000575-37-1 | 80   |
| 4    |    | Naphthalene, 2,7-dimethyl- | 156 | C12H12  | 000582-16-1 | 80   |
| 5    |    | Naphthalene, 2,7-dimethyl- | 156 | C12H12  | 000582-16-1 | 72   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022420\  
Data File : BN009891.D  
Acq On : 25 Feb 2020 09:04  
Operator : CG/JU  
Sample : L1424-22DL 10X  
Misc :  
ALS Vial : 29 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
C5B30DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN021220MA.M  
Quant Title : SVOA CALIBRATION

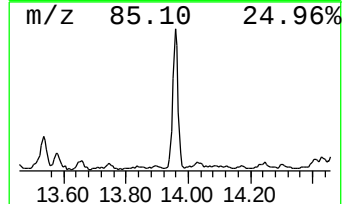
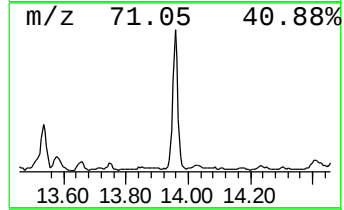
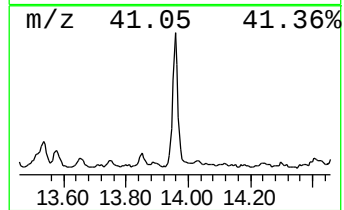
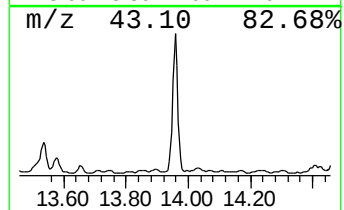
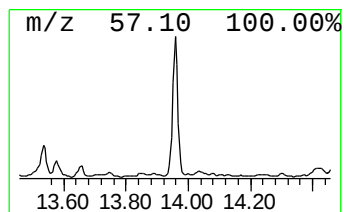
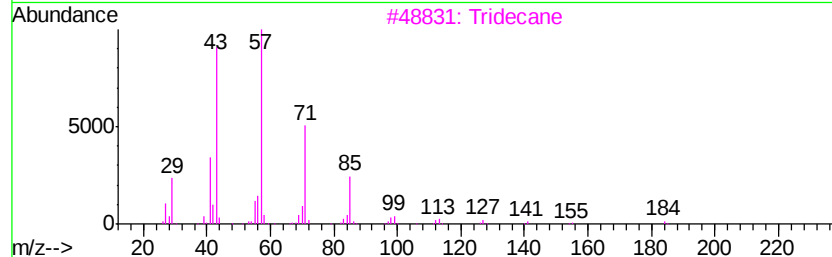
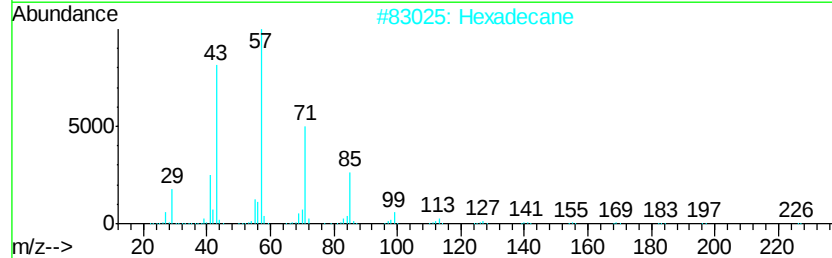
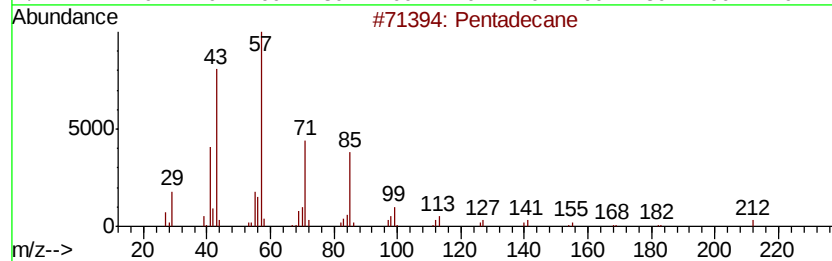
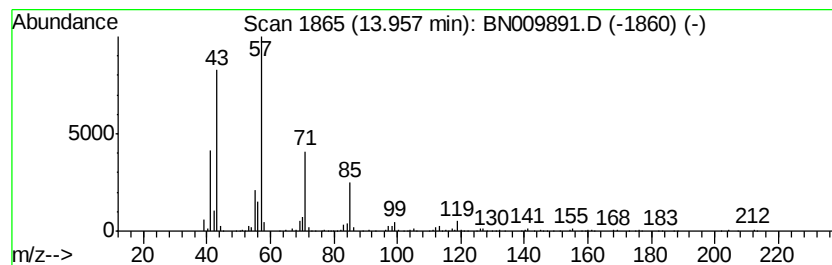
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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

| R.T.  | EstConc     | Area   | Relative to ISTD | R.T.  |
|-------|-------------|--------|------------------|-------|
| 13.96 | 12.10 ng/ul | 974718 | Acenaphthene-d10 | 14.02 |

| Hit# | of | Tentative ID         | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------|-----|---------|-------------|------|
| 1    | 5  | Pentadecane          | 212 | C15H32  | 000629-62-9 | 93   |
| 2    |    | Hexadecane           | 226 | C16H34  | 000544-76-3 | 91   |
| 3    |    | Tridecane            | 184 | C13H28  | 000629-50-5 | 90   |
| 4    |    | Tetradecane, 1-iodo- | 324 | C14H29I | 019218-94-1 | 80   |
| 5    |    | Tridecane, 7-propyl- | 226 | C16H34  | 055045-09-5 | 72   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_N\DATA\BN022420\  
Data File : BN009891.D  
Acq On : 25 Feb 2020 09:04  
Operator : CG/JU  
Sample : L1424-22DL 10X  
Misc :  
ALS Vial : 29 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
C5B30DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN021220MA.M  
Quant Title : SVOA CALIBRATION

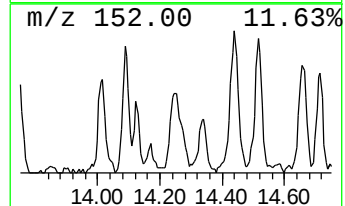
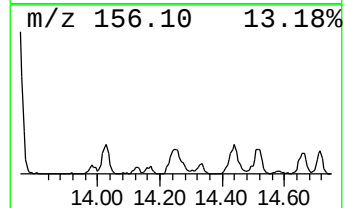
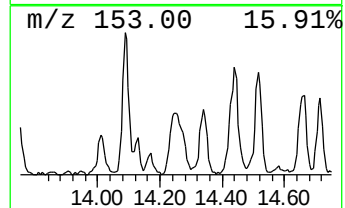
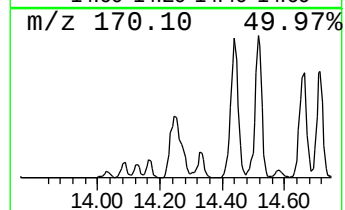
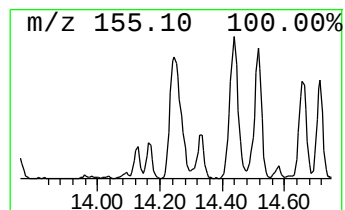
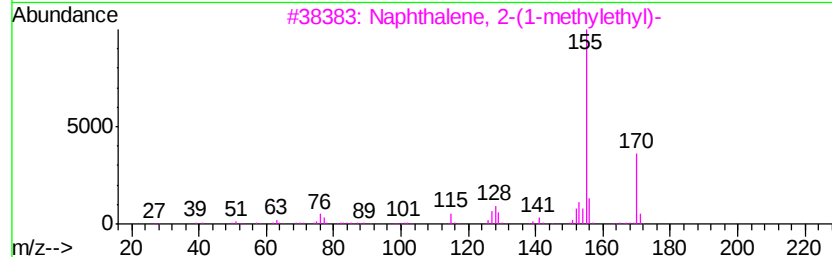
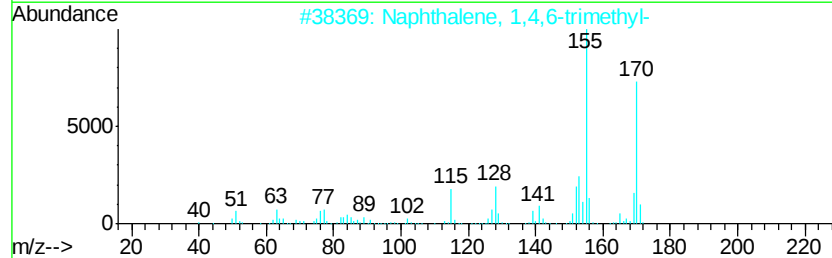
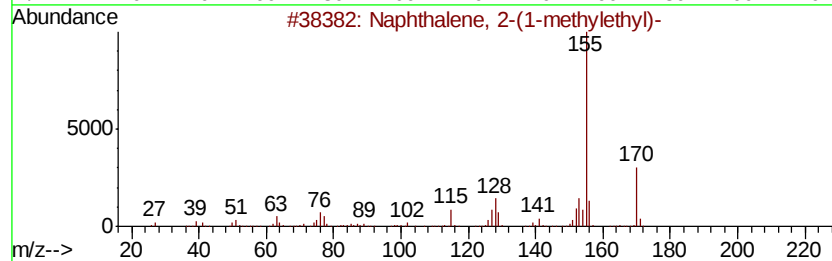
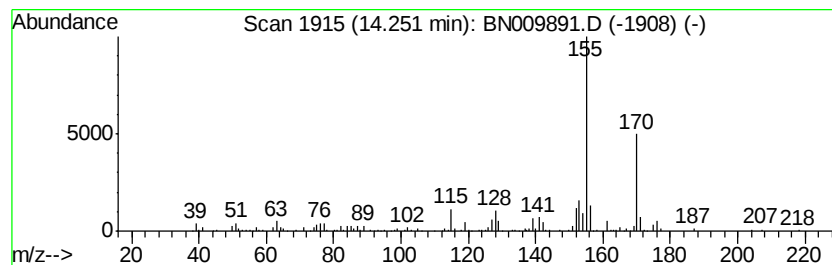
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 35 Naphthalene, 2-(1-methyleth... Concentration Rank 16

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 14.25 | 7.43 ng/ul | 598233 | Acenaphthene-d10 | 14.02 |

| Hit# | of | 5 | Tentative ID                    | MW  | MolForm | CAS#         | Qual |
|------|----|---|---------------------------------|-----|---------|--------------|------|
| 1    |    |   | Naphthalene, 2-(1-methylethyl)- | 170 | C13H14  | 002027-17-0  | 94   |
| 2    |    |   | Naphthalene, 1,4,6-trimethyl-   | 170 | C13H14  | 002131-42-2  | 92   |
| 3    |    |   | Naphthalene, 2-(1-methylethyl)- | 170 | C13H14  | 002027-17-0  | 91   |
| 4    |    |   | 3-(2-Methyl-propenyl)-1H-indene | 170 | C13H14  | 1000187-78-5 | 76   |
| 5    |    |   | Naphthalene, 1,6,7-trimethyl-   | 170 | C13H14  | 002245-38-7  | 70   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022420\  
Data File : BN009891.D  
Acq On : 25 Feb 2020 09:04  
Operator : CG/JU  
Sample : L1424-22DL 10X  
Misc :  
ALS Vial : 29 Sample Multiplier: 1

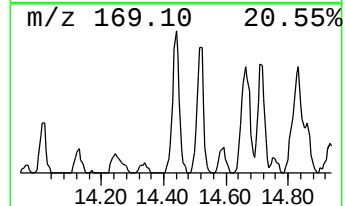
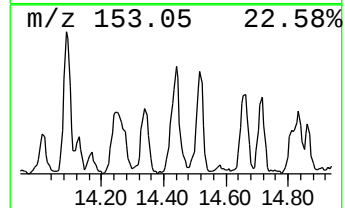
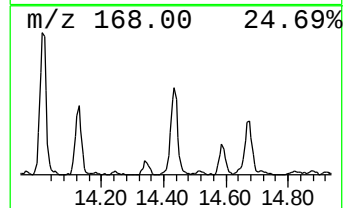
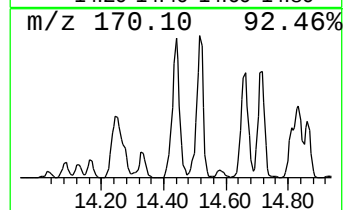
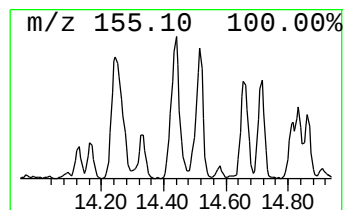
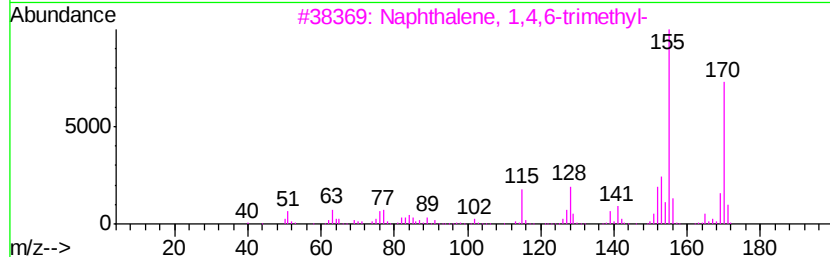
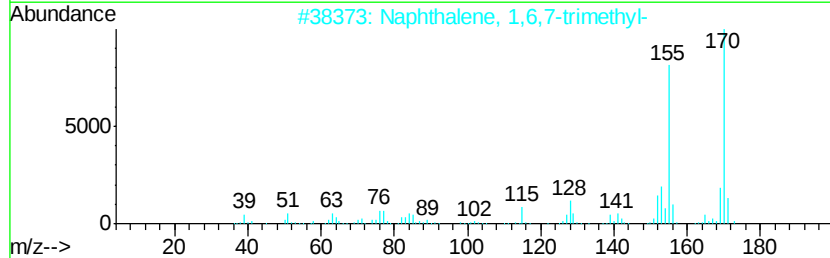
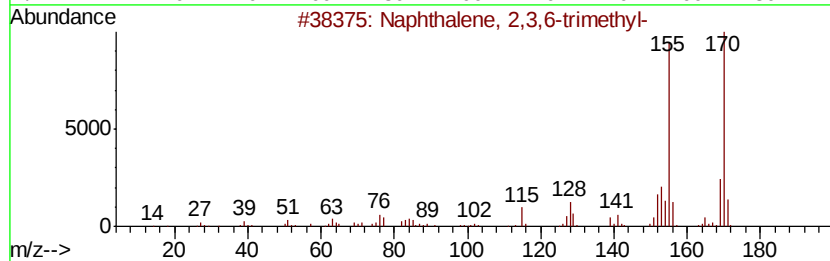
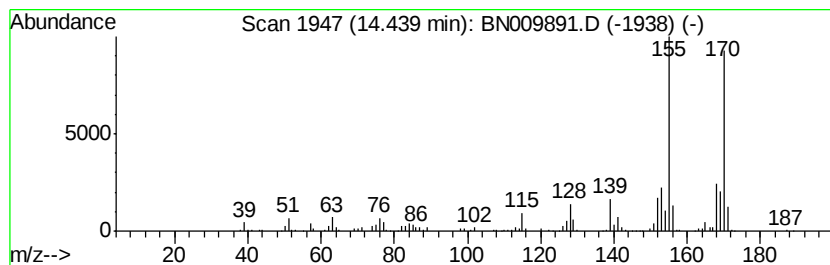
Instrument :  
BNA\_N  
ClientSampleId :  
C5B30DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN021220MA.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 36 Naphthalene, 2,3,6-trimethyl- Concentration Rank 13

| R.T.      | EstConc                       | Area   | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------|--------|------------------|-------------|------|
| 14.44     | 9.47 ng/ul                    | 762907 | Acenaphthene-d10 | 14.02       |      |
| Hit# of 5 | Tentative ID                  | MW     | MolForm          | CAS#        | Qual |
| 1         | Naphthalene, 2,3,6-trimethyl- | 170    | C13H14           | 000829-26-5 | 98   |
| 2         | Naphthalene, 1,6,7-trimethyl- | 170    | C13H14           | 002245-38-7 | 96   |
| 3         | Naphthalene, 1,4,6-trimethyl- | 170    | C13H14           | 002131-42-2 | 95   |
| 4         | Naphthalene, 2,3,6-trimethyl- | 170    | C13H14           | 000829-26-5 | 95   |
| 5         | Naphthalene, 2,3,6-trimethyl- | 170    | C13H14           | 000829-26-5 | 94   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022420\  
Data File : BN009891.D  
Acq On : 25 Feb 2020 09:04  
Operator : CG/JU  
Sample : L1424-22DL 10X  
Misc :  
ALS Vial : 29 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
C5B30DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN021220MA.M  
Quant Title : SVOA CALIBRATION

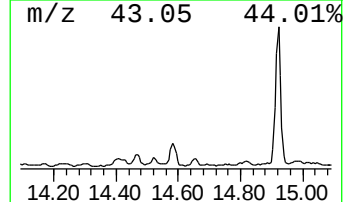
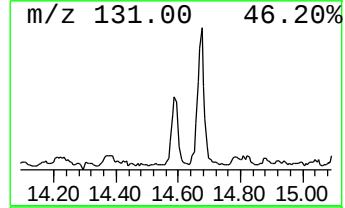
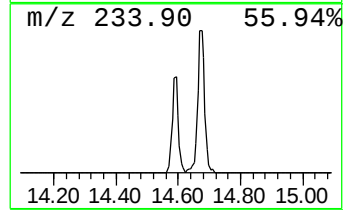
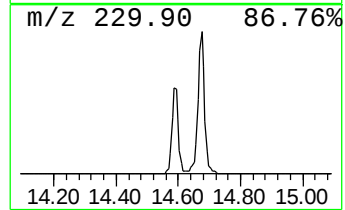
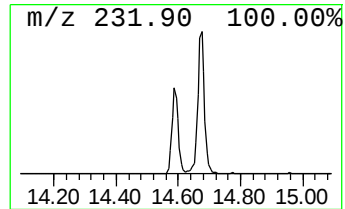
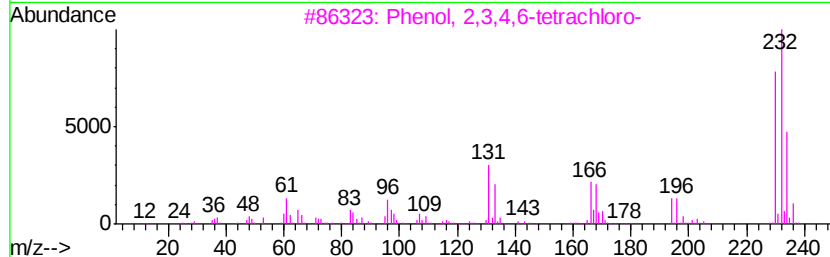
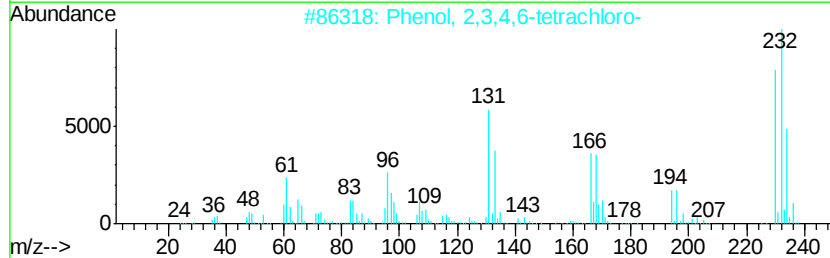
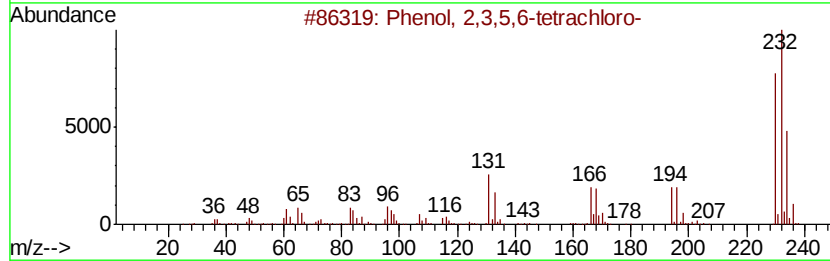
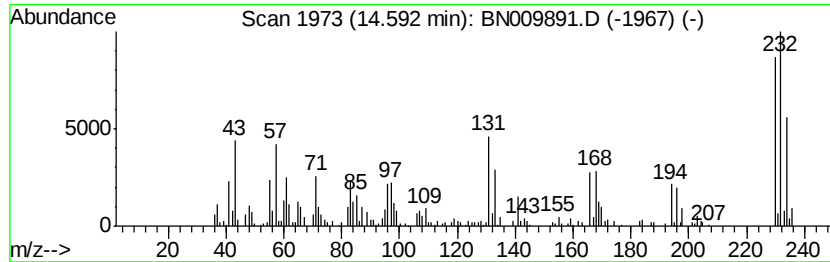
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 38 Phenol, 2,3,5,6-tetrachloro- Concentration Rank 17

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 14.59 | 7.24 ng/ul | 583328 | Acenaphthene-d10 | 14.02 |

| Hit# | of | 5 | Tentative ID                 | MW  | MolForm  | CAS#        | Qual |
|------|----|---|------------------------------|-----|----------|-------------|------|
| 1    |    |   | Phenol, 2,3,5,6-tetrachloro- | 230 | C6H2Cl4O | 000935-95-5 | 99   |
| 2    |    |   | Phenol, 2,3,4,6-tetrachloro- | 230 | C6H2Cl4O | 000058-90-2 | 98   |
| 3    |    |   | Phenol, 2,3,4,6-tetrachloro- | 230 | C6H2Cl4O | 000058-90-2 | 98   |
| 4    |    |   | Phenol, 2,3,4,6-tetrachloro- | 230 | C6H2Cl4O | 000058-90-2 | 96   |
| 5    |    |   | Phenol, 2,3,4,5-tetrachloro- | 230 | C6H2Cl4O | 004901-51-3 | 95   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022420\  
Data File : BN009891.D  
Acq On : 25 Feb 2020 09:04  
Operator : CG/JU  
Sample : L1424-22DL 10X  
Misc :  
ALS Vial : 29 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
C5B30DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN021220MA.M  
Quant Title : SVOA CALIBRATION

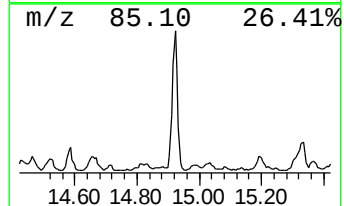
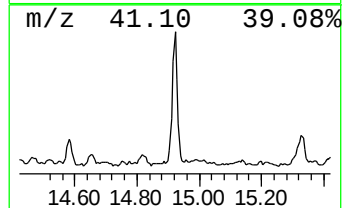
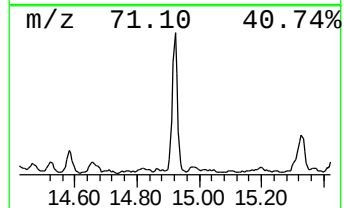
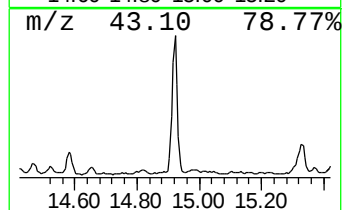
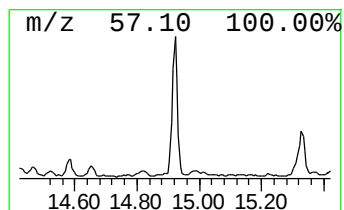
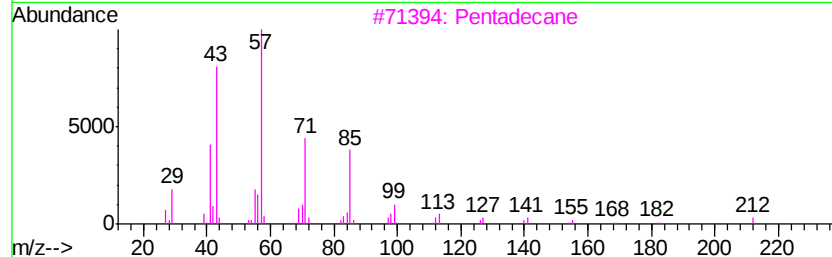
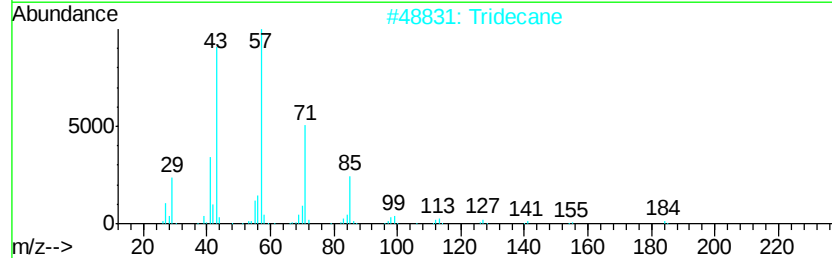
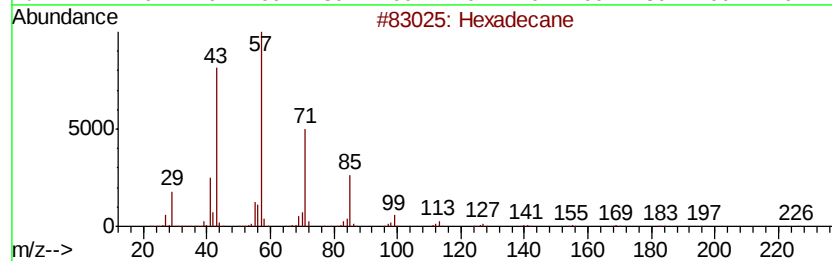
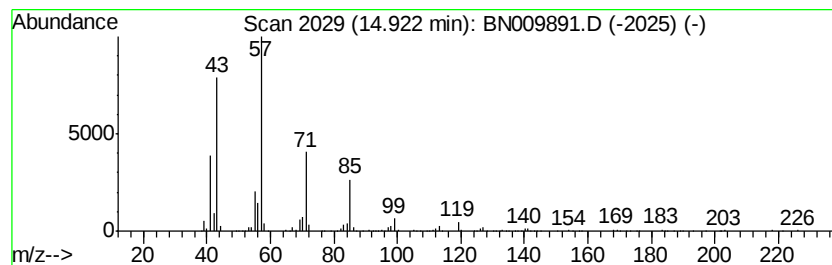
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 40 (DEL) Alkane: Straight-Chai... Concentration Rank 12

| R.T.  | EstConc     | Area   | Relative to ISTD | R.T.  |
|-------|-------------|--------|------------------|-------|
| 14.92 | 10.47 ng/ul | 843188 | Acenaphthene-d10 | 14.02 |

| Hit# | of | Tentative ID         | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------|-----|---------|-------------|------|
| 1    | 5  | Hexadecane           | 226 | C16H34  | 000544-76-3 | 91   |
| 2    |    | Tridecane            | 184 | C13H28  | 000629-50-5 | 90   |
| 3    |    | Pentadecane          | 212 | C15H32  | 000629-62-9 | 86   |
| 4    |    | Tridecane            | 184 | C13H28  | 000629-50-5 | 83   |
| 5    |    | Tetradecane, 1-iodo- | 324 | C14H29I | 019218-94-1 | 80   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022420\  
Data File : BN009891.D  
Acq On : 25 Feb 2020 09:04  
Operator : CG/JU  
Sample : L1424-22DL 10X  
Misc :  
ALS Vial : 29 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
C5B30DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN021220MA.M  
Quant Title : SVOA CALIBRATION

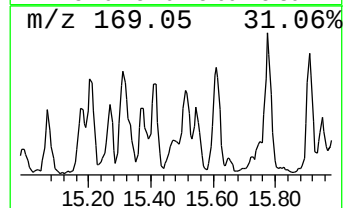
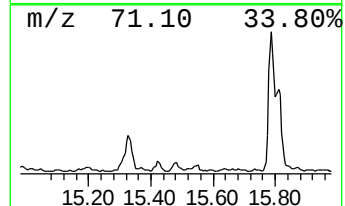
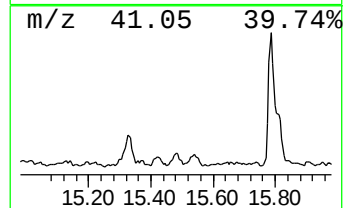
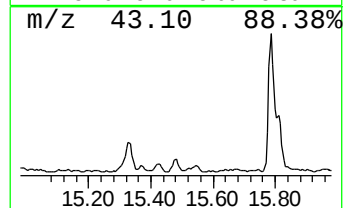
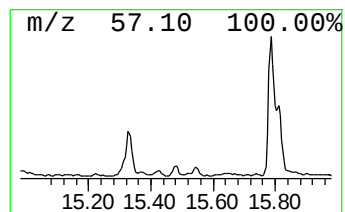
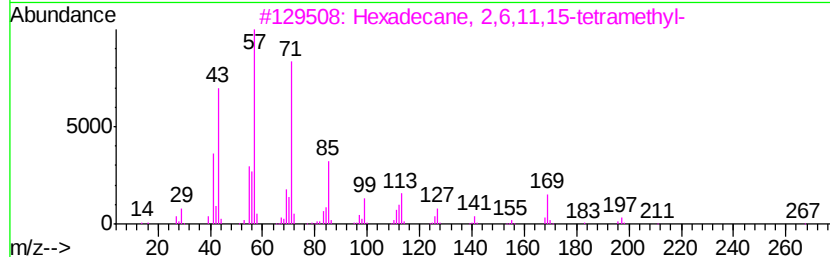
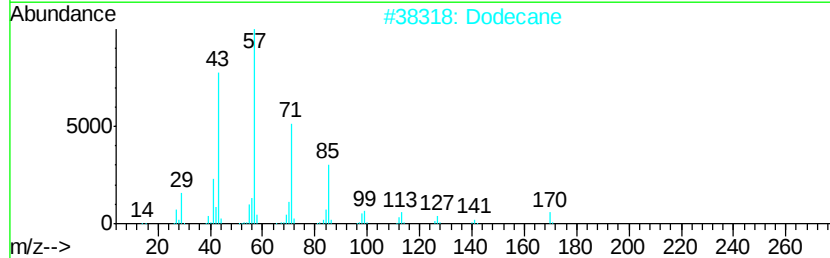
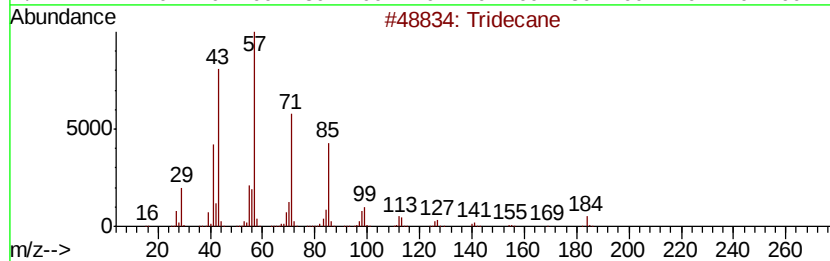
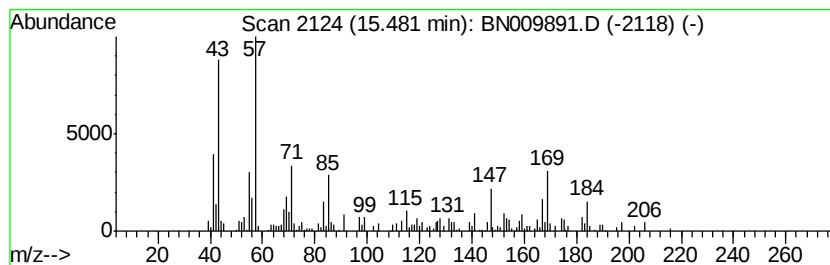
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 44 (DEL) Alkane: Straight-Chai... Concentration Rank 55

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 15.48 | 2.91 ng/ul | 199703 | Phenanthrene-d10 | 16.79 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|-------------------------------------|-----|----------|--------------|------|
| 1    | 5  | Tridecane                           | 184 | C13H28   | 000629-50-5  | 38   |
| 2    |    | Dodecane                            | 170 | C12H26   | 000112-40-3  | 30   |
| 3    |    | Hexadecane, 2,6,11,15-tetramethyl-  | 282 | C20H42   | 000504-44-9  | 27   |
| 4    |    | Pentadecane, 2-methyl-              | 226 | C16H34   | 001560-93-6  | 22   |
| 5    |    | Methoxyacetic acid, 2-tetradecyl... | 286 | C17H34O3 | 1000282-04-8 | 18   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022420\  
Data File : BN009891.D  
Acq On : 25 Feb 2020 09:04  
Operator : CG/JU  
Sample : L1424-22DL 10X  
Misc :  
ALS Vial : 29 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
C5B30DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN021220MA.M  
Quant Title : SVOA CALIBRATION

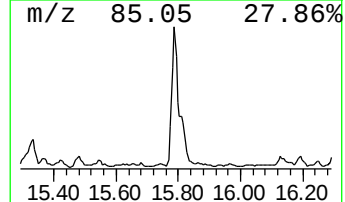
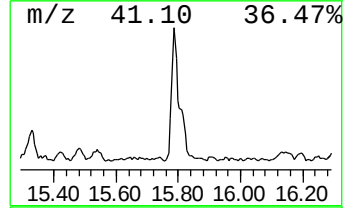
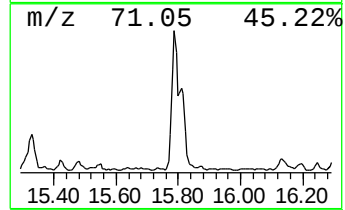
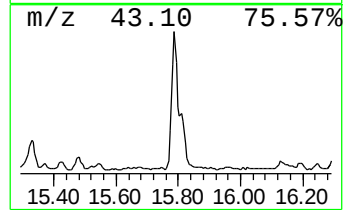
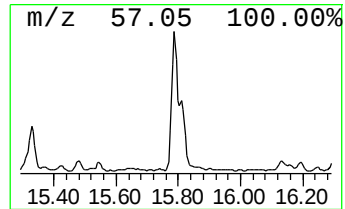
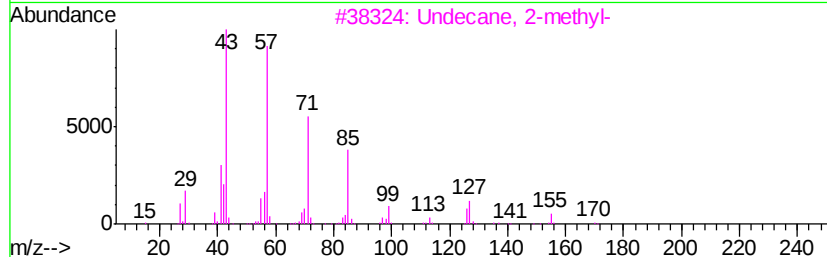
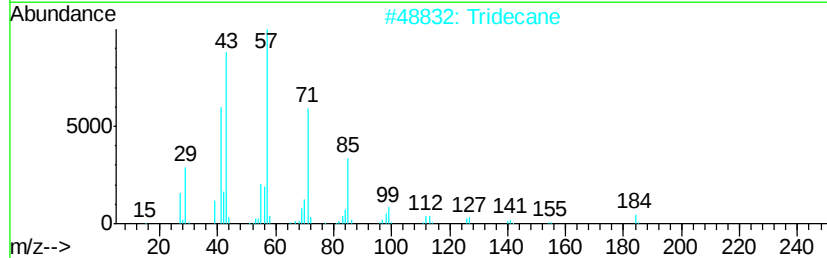
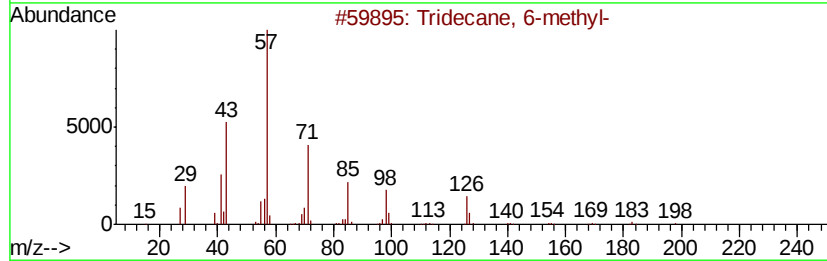
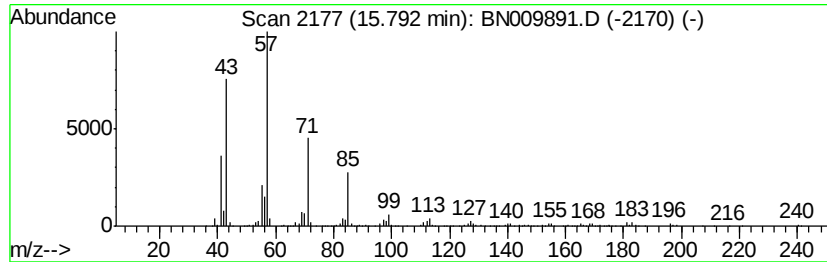
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 47 (DEL) Alkane: Straight-Chai... Concentration Rank 4

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 15.79 | 17.65 ng/ul | 1209580 | Phenanthrene-d10 | 16.79 |

| Hit# | of | Tentative ID         | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------|-----|---------|-------------|------|
| 1    | 5  | Tridecane, 6-methyl- | 198 | C14H30  | 013287-21-3 | 91   |
| 2    |    | Tridecane            | 184 | C13H28  | 000629-50-5 | 86   |
| 3    |    | Undecane, 2-methyl-  | 170 | C12H26  | 007045-71-8 | 86   |
| 4    |    | Octadecane           | 254 | C18H38  | 000593-45-3 | 86   |
| 5    |    | Hexadecane           | 226 | C16H34  | 000544-76-3 | 86   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022420\  
Data File : BN009891.D  
Acq On : 25 Feb 2020 09:04  
Operator : CG/JU  
Sample : L1424-22DL 10X  
Misc :  
ALS Vial : 29 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
C5B30DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN021220MA.M  
Quant Title : SVOA CALIBRATION

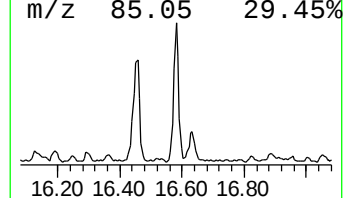
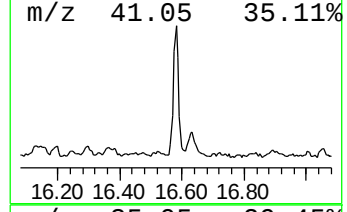
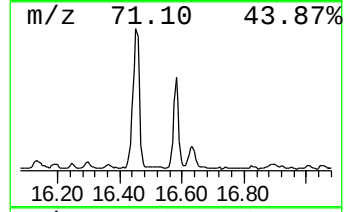
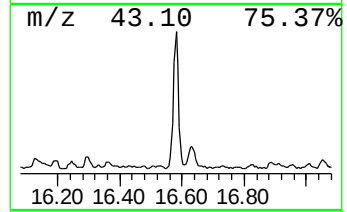
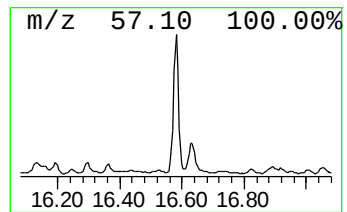
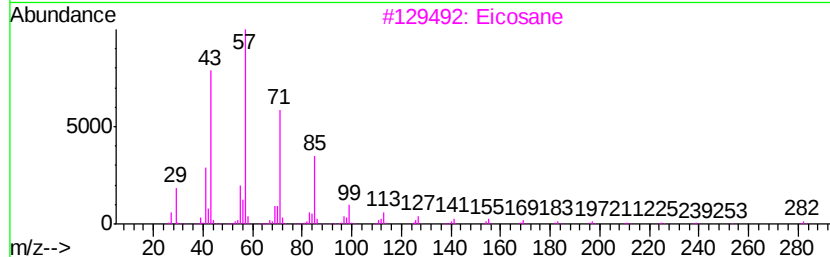
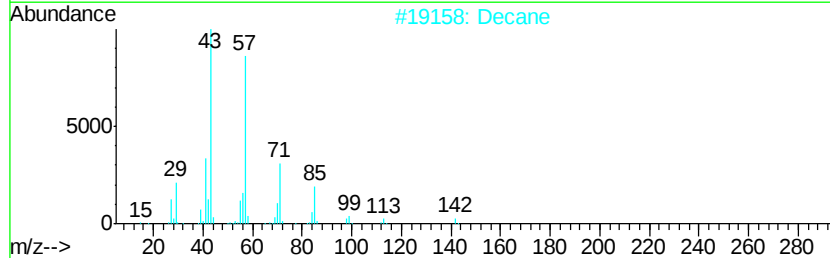
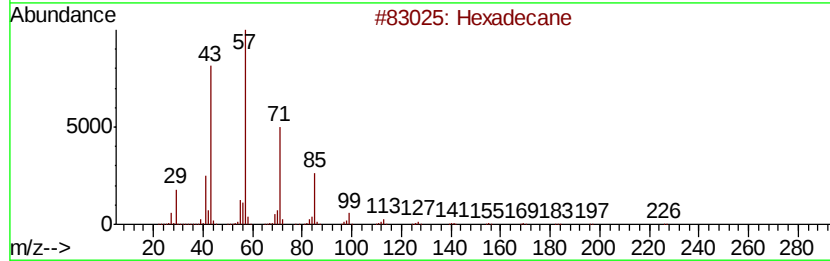
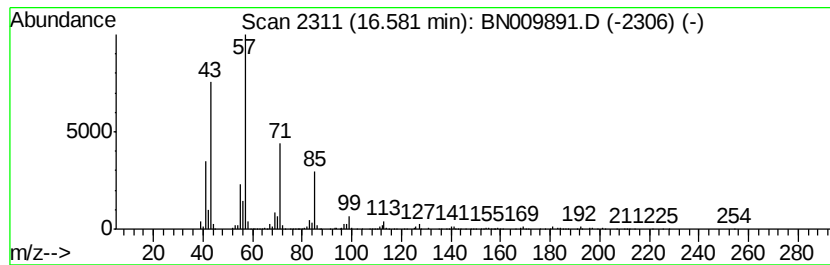
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 51 (DEL) Alkane: Straight-Chai... Concentration Rank 8

| R.T.  | EstConc     | Area   | Relative to ISTD | R.T.  |
|-------|-------------|--------|------------------|-------|
| 16.58 | 11.24 ng/ul | 770605 | Phenanthrene-d10 | 16.79 |

| Hit# | of | 5 | Tentative ID                         | MW  | MolForm   | CAS#         | Qual |
|------|----|---|--------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Hexadecane                           | 226 | C16H34    | 000544-76-3  | 91   |
| 2    |    |   | Decane                               | 142 | C10H22    | 000124-18-5  | 90   |
| 3    |    |   | Eicosane                             | 282 | C20H42    | 000112-95-8  | 86   |
| 4    |    |   | Sulfurous acid, 2-propyl tetradec... | 320 | C17H36O3S | 1000309-12-5 | 86   |
| 5    |    |   | Tetradecane, 1-iodo-                 | 324 | C14H29I   | 019218-94-1  | 86   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_N\DATA\BN022420\  
Data File : BN009891.D  
Acq On : 25 Feb 2020 09:04  
Operator : CG/JU  
Sample : L1424-22DL 10X  
Misc :  
ALS Vial : 29 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
C5B30DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN021220MA.M  
Quant Title : SVOA CALIBRATION

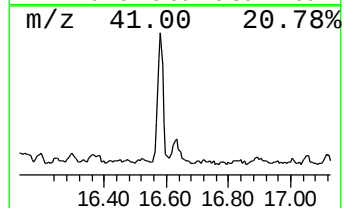
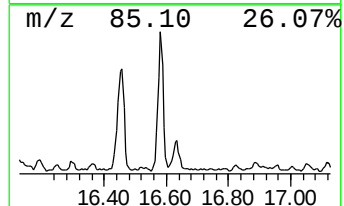
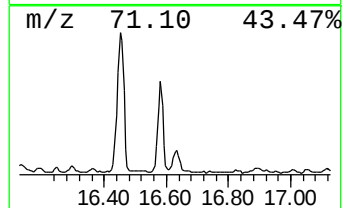
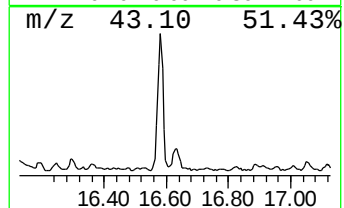
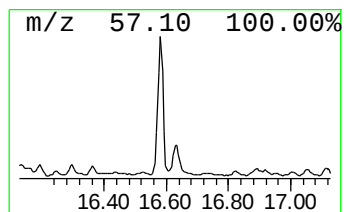
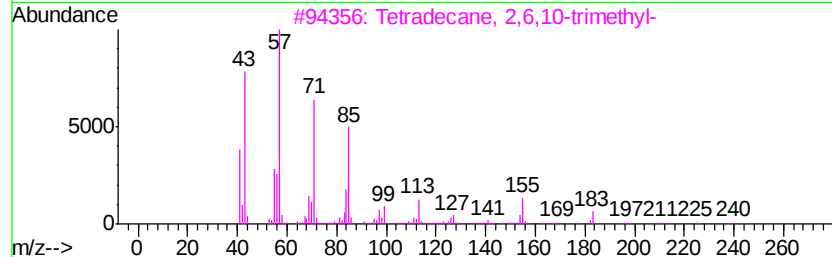
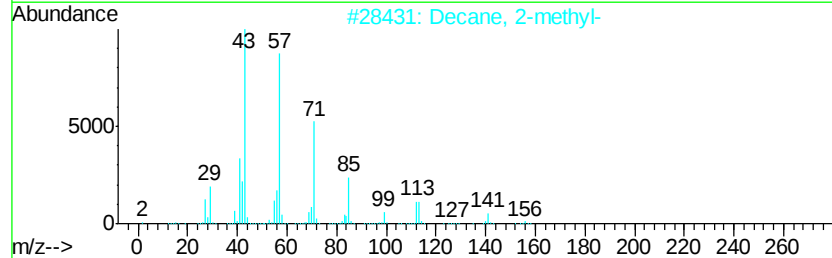
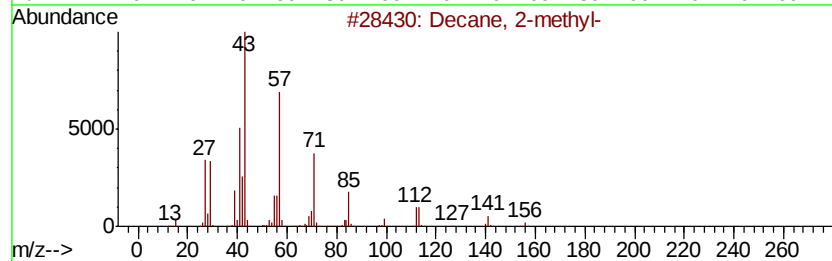
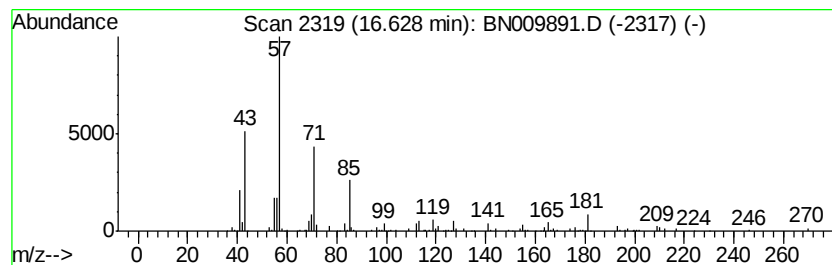
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 52 (DEL) Alkane: Straight-Chai... Concentration Rank 42

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 16.63 | 3.84 ng/ul | 263308 | Phenanthrene-d10 | 16.79 |

| Hit# | of | 5 | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------|-----|---------|-------------|------|
| 1    |    |   | Decane, 2-methyl-              | 156 | C11H24  | 006975-98-0 | 64   |
| 2    |    |   | Decane, 2-methyl-              | 156 | C11H24  | 006975-98-0 | 64   |
| 3    |    |   | Tetradecane, 2,6,10-trimethyl- | 240 | C17H36  | 014905-56-7 | 59   |
| 4    |    |   | Undecane, 5-methyl-            | 170 | C12H26  | 001632-70-8 | 53   |
| 5    |    |   | Octane, 3-ethyl-2,7-dimethyl-  | 170 | C12H26  | 062183-55-5 | 53   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022420\  
Data File : BN009891.D  
Acq On : 25 Feb 2020 09:04  
Operator : CG/JU  
Sample : L1424-22DL 10X  
Misc :  
ALS Vial : 29 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
C5B30DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN021220MA.M  
Quant Title : SVOA CALIBRATION

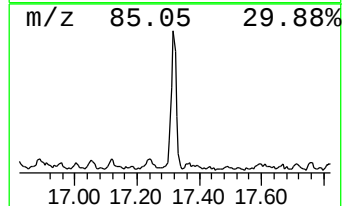
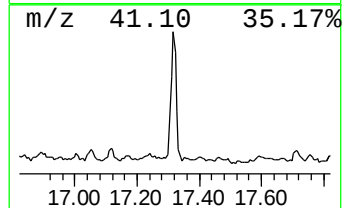
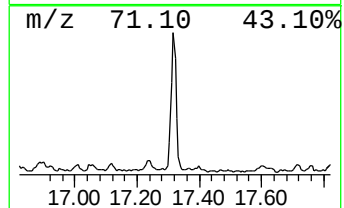
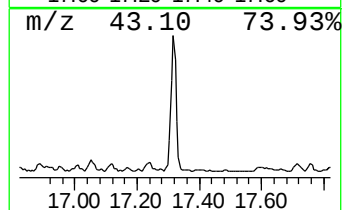
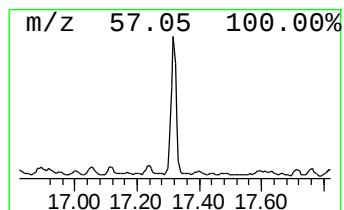
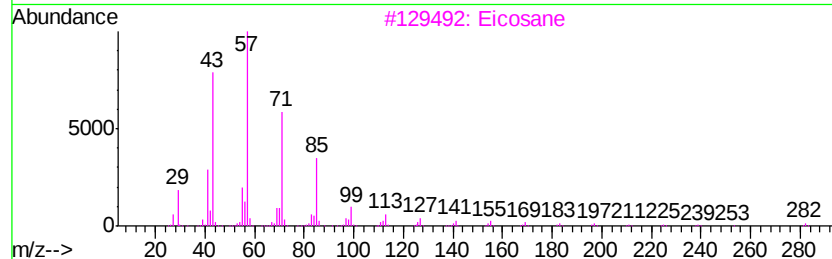
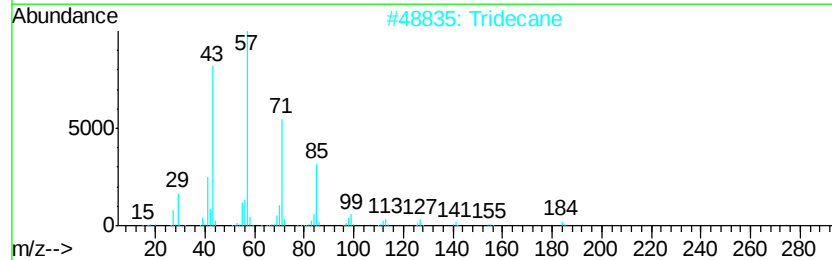
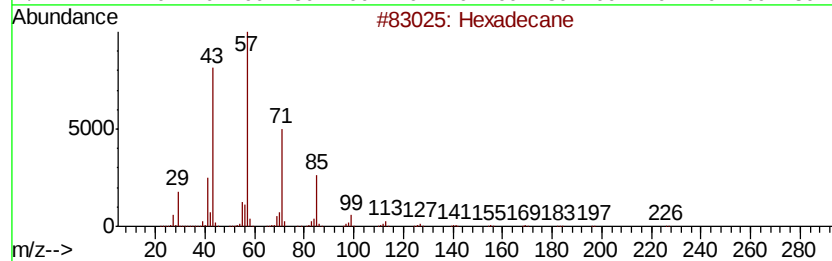
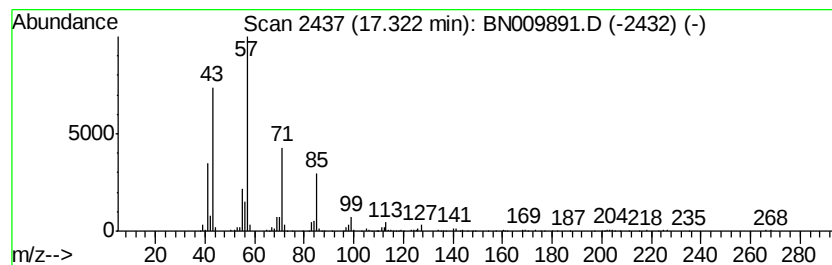
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 54 (DEL) Alkane: Straight-Chai... Concentration Rank 15

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 17.32 | 7.86 ng/ul | 538477 | Phenanthrene-d10 | 16.79 |

| Hit# | of | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------|-----|---------|-------------|------|
| 1    |    |   | Hexadecane   | 226 | C16H34  | 000544-76-3 | 97   |
| 2    |    |   | Tridecane    | 184 | C13H28  | 000629-50-5 | 90   |
| 3    |    |   | Eicosane     | 282 | C20H42  | 000112-95-8 | 86   |
| 4    |    |   | Heptadecane  | 240 | C17H36  | 000629-78-7 | 86   |
| 5    |    |   | Dodecane     | 170 | C12H26  | 000112-40-3 | 83   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022420\  
Data File : BN009891.D  
Acq On : 25 Feb 2020 09:04  
Operator : CG/JU  
Sample : L1424-22DL 10X  
Misc :  
ALS Vial : 29 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
C5B30DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN021220MA.M  
Quant Title : SVOA CALIBRATION

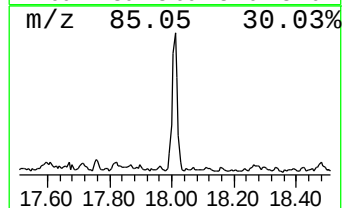
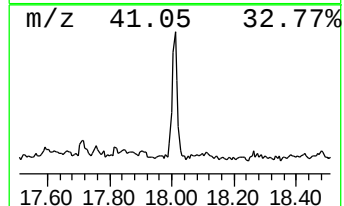
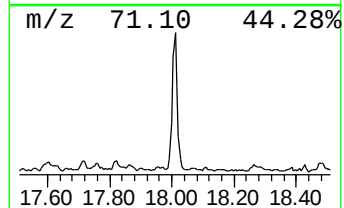
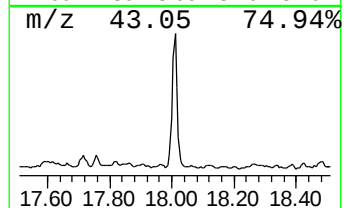
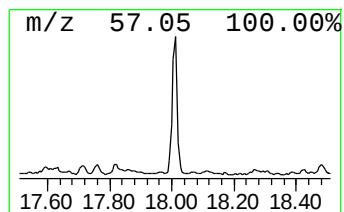
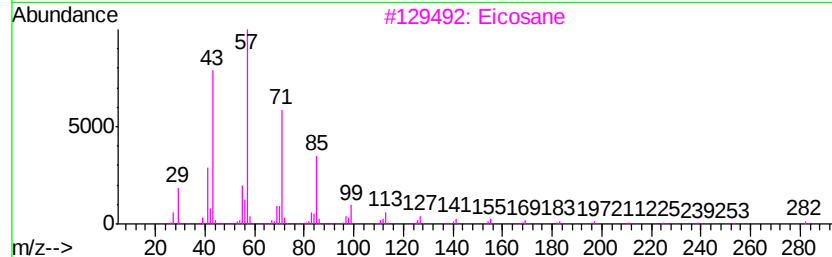
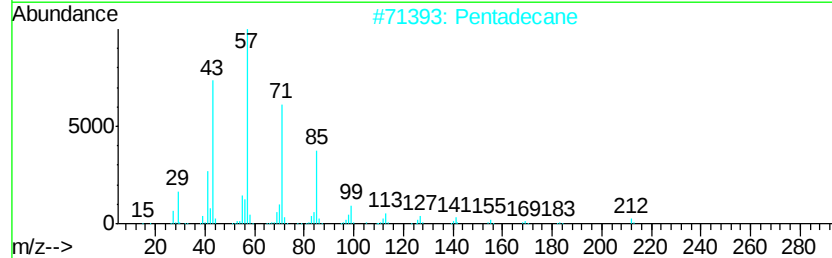
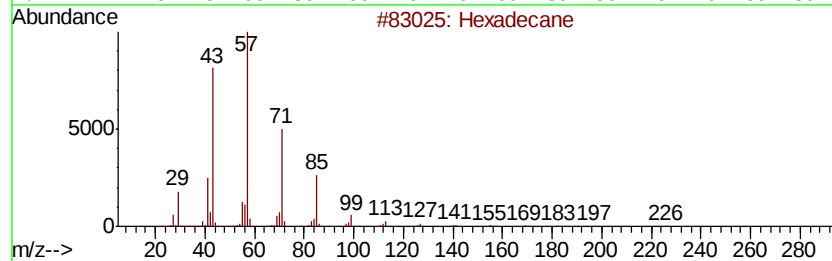
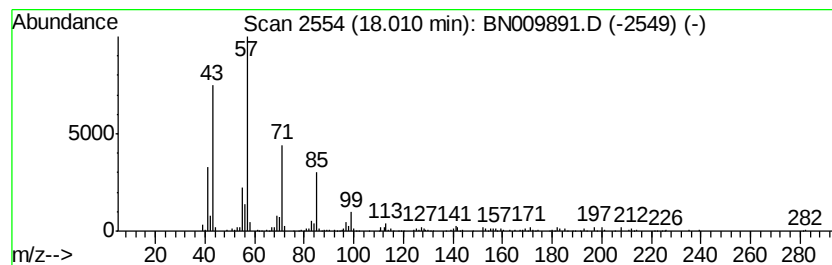
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 59 (DEL) Alkane: Straight-Chai... Concentration Rank 18

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 18.01 | 7.09 ng/ul | 485639 | Phenanthrene-d10 | 16.79 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Hexadecane                 | 226 | C16H34  | 000544-76-3 | 96   |
| 2    |    |   | Pentadecane                | 212 | C15H32  | 000629-62-9 | 95   |
| 3    |    |   | Eicosane                   | 282 | C20H42  | 000112-95-8 | 94   |
| 4    |    |   | Tetradecane, 6,9-dimethyl- | 226 | C16H34  | 055045-13-1 | 94   |
| 5    |    |   | Tridecane, 2-methyl-       | 198 | C14H30  | 001560-96-9 | 87   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022420\  
Data File : BN009891.D  
Acq On : 25 Feb 2020 09:04  
Operator : CG/JU  
Sample : L1424-22DL 10X  
Misc :  
ALS Vial : 29 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
C5B30DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN021220MA.M  
Quant Title : SVOA CALIBRATION

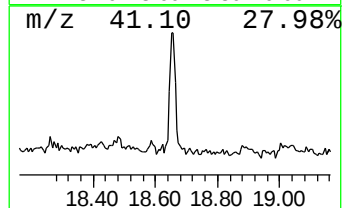
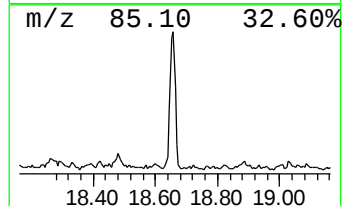
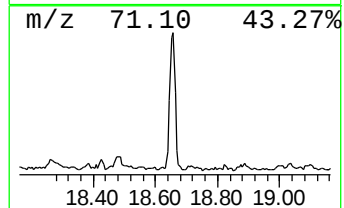
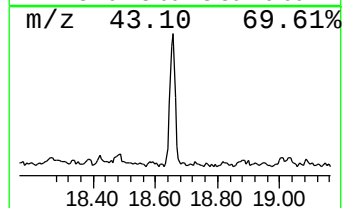
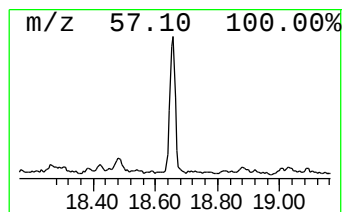
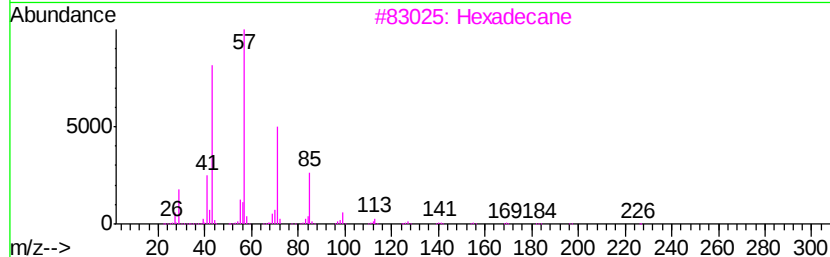
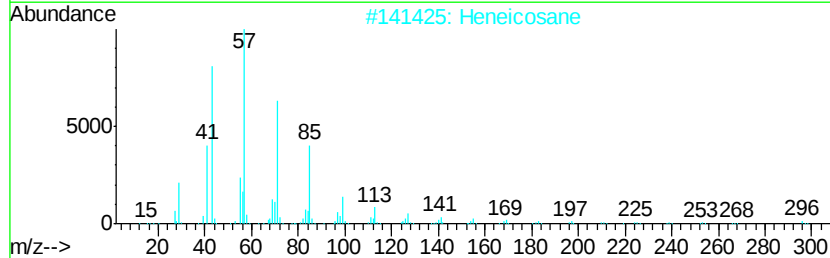
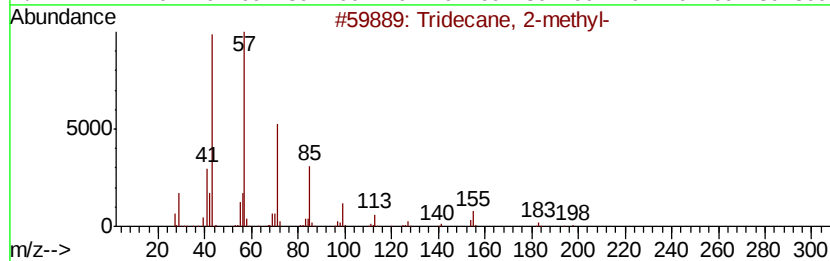
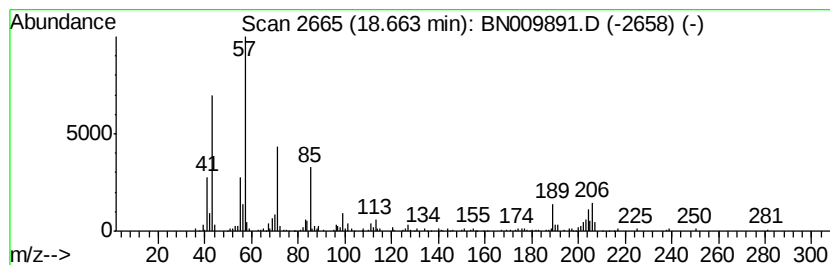
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 61 (DEL) Alkane: Straight-Chai... Concentration Rank 26

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 18.66 | 5.32 ng/ul | 364657 | Phenanthrene-d10 | 16.79 |

| Hit# | of | 5 | Tentative ID         | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------|-----|---------|-------------|------|
| 1    |    |   | Tridecane, 2-methyl- | 198 | C14H30  | 001560-96-9 | 64   |
| 2    |    |   | Heneicosane          | 296 | C21H44  | 000629-94-7 | 52   |
| 3    |    |   | Hexadecane           | 226 | C16H34  | 000544-76-3 | 52   |
| 4    |    |   | Eicosane, 10-methyl- | 296 | C21H44  | 054833-23-7 | 52   |
| 5    |    |   | Tridecane            | 184 | C13H28  | 000629-50-5 | 50   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022420\  
Data File : BN009891.D  
Acq On : 25 Feb 2020 09:04  
Operator : CG/JU  
Sample : L1424-22DL 10X  
Misc :  
ALS Vial : 29 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
C5B30DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN021220MA.M  
Quant Title : SVOA CALIBRATION

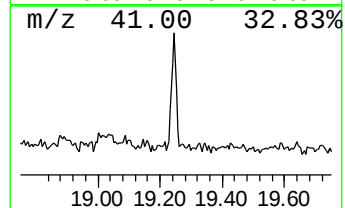
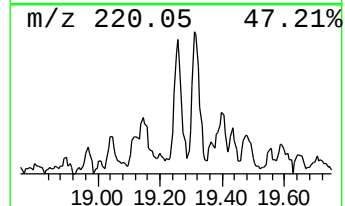
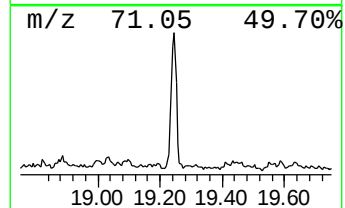
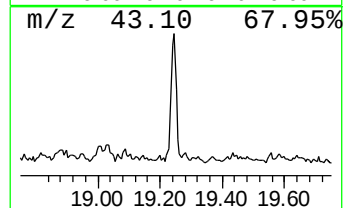
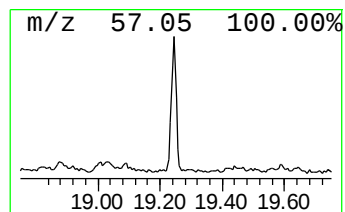
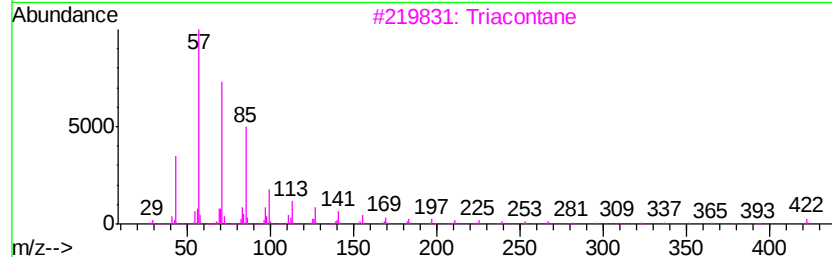
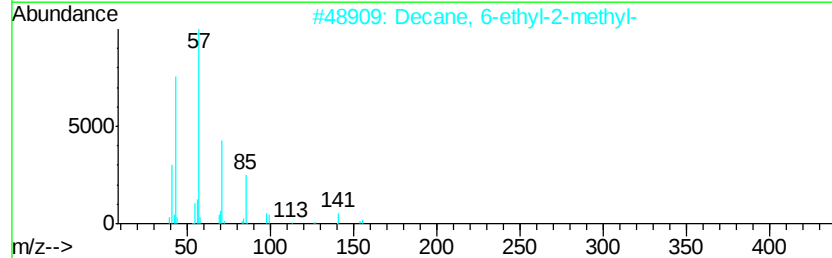
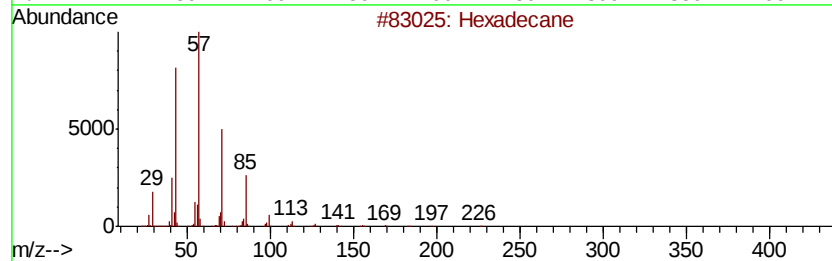
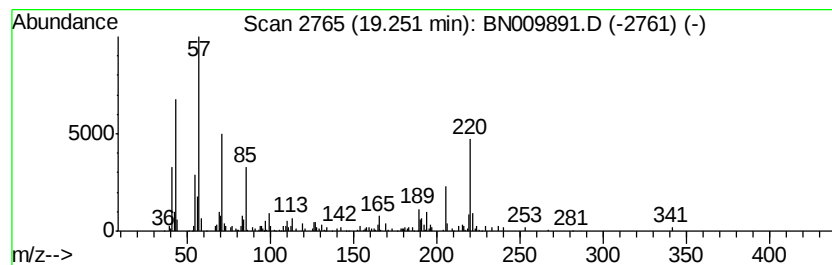
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 62 (DEL) Alkane: Straight-Chai... Concentration Rank 52

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 19.25 | 3.08 ng/ul | 252725 | Chrysene-d12     | 21.00 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Hexadecane                          | 226 | C16H34    | 000544-76-3  | 68   |
| 2    |    |   | Decane, 6-ethyl-2-methyl-           | 184 | C13H28    | 062108-21-8  | 64   |
| 3    |    |   | triacontane                         | 422 | C30H62    | 000638-68-6  | 59   |
| 4    |    |   | Sulfurous acid, 2-ethylhexyl hex... | 278 | C14H30O3S | 1000309-20-2 | 59   |
| 5    |    |   | Undecane, 3-methyl-                 | 170 | C12H26    | 001002-43-3  | 59   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022420\  
Data File : BN009891.D  
Acq On : 25 Feb 2020 09:04  
Operator : CG/JU  
Sample : L1424-22DL 10X  
Misc :  
ALS Vial : 29 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
C5B30DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN021220MA.M  
Quant Title : SVOA CALIBRATION

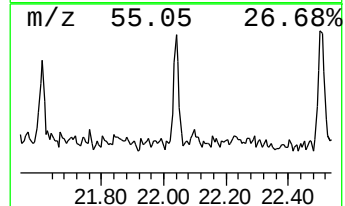
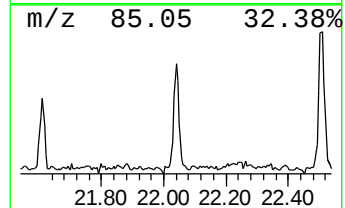
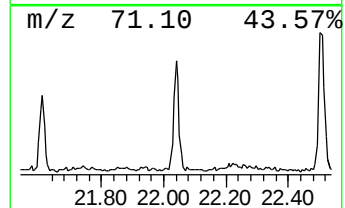
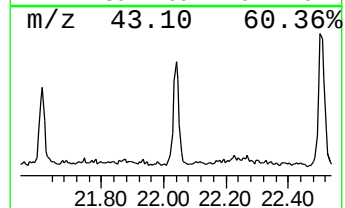
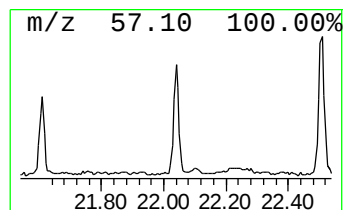
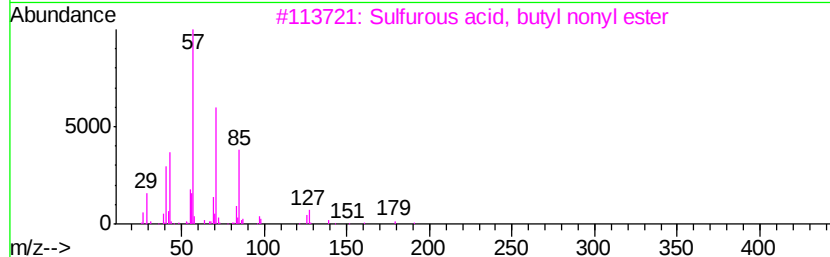
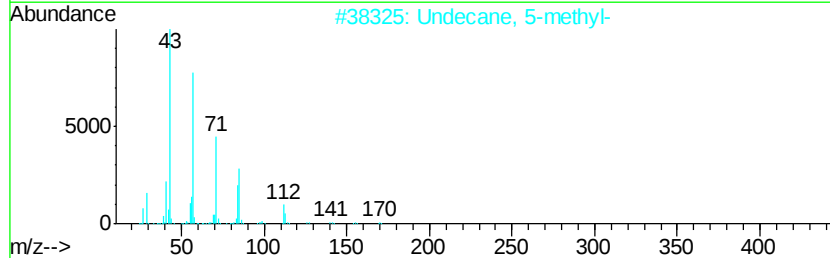
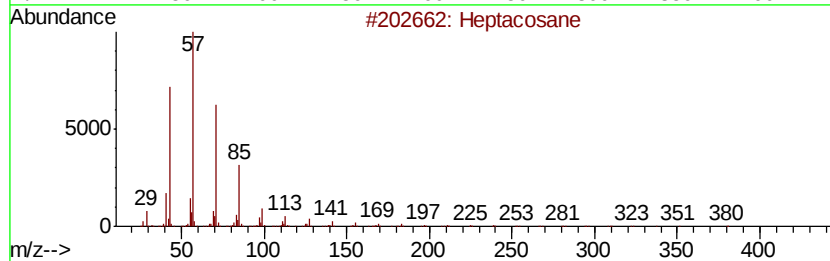
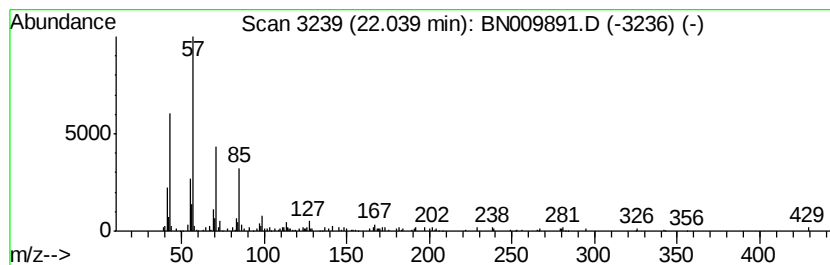
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 22.04 | 2.16 ng/ul | 177041 | Chrysene-d12     | 21.00 |

| Hit# | of | Tentative ID                      | MW  | MolForm   | CAS#         | Qual |
|------|----|-----------------------------------|-----|-----------|--------------|------|
| 1    | 5  | Heptacosane                       | 380 | C27H56    | 000593-49-7  | 74   |
| 2    |    | Undecane, 5-methyl-               | 170 | C12H26    | 001632-70-8  | 72   |
| 3    |    | Sulfurous acid, butyl nonyl ester | 264 | C13H28O3S | 1000309-17-6 | 72   |
| 4    |    | Tetracosane                       | 338 | C24H50    | 000646-31-1  | 72   |
| 5    |    | Tridecane                         | 184 | C13H28    | 000629-50-5  | 70   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022420\  
Data File : BN009891.D  
Acq On : 25 Feb 2020 09:04  
Operator : CG/JU  
Sample : L1424-22DL 10X  
Misc :  
ALS Vial : 29 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
C5B30DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN021220MA.M  
Quant Title : SVOA CALIBRATION

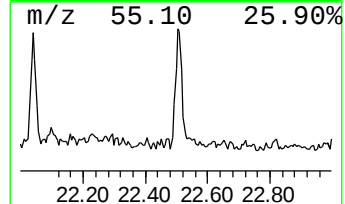
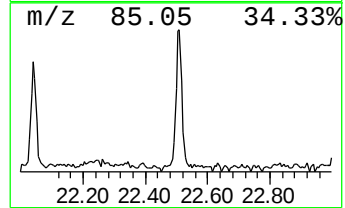
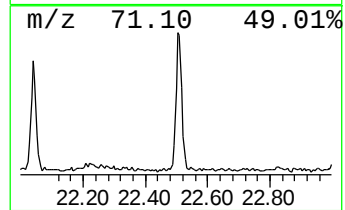
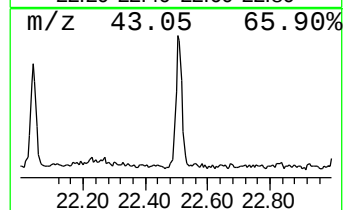
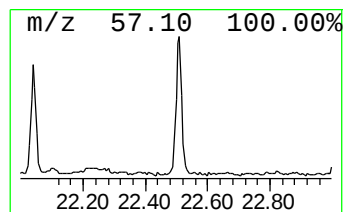
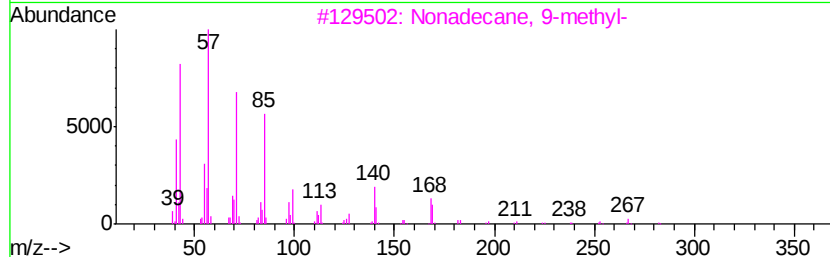
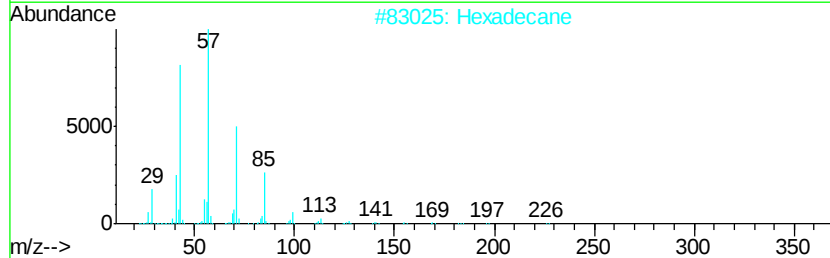
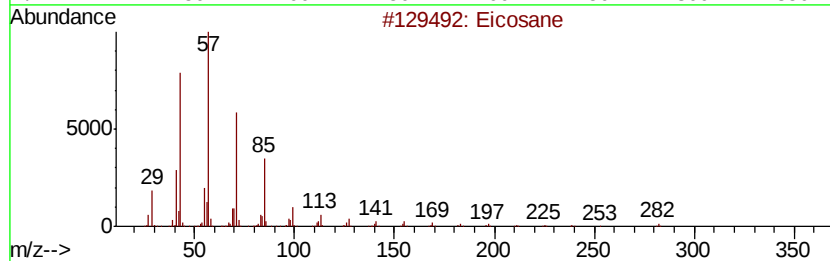
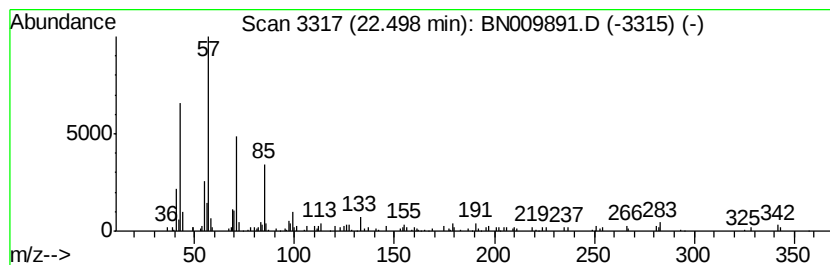
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 64 (DEL) Alkane: Straight-Chai... Concentration Rank 60

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 22.50 | 2.44 ng/ul | 201004 | Perylene-d12     | 23.10 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Eicosane                   | 282 | C20H42  | 000112-95-8 | 98   |
| 2    |    |   | Hexadecane                 | 226 | C16H34  | 000544-76-3 | 94   |
| 3    |    |   | Nonadecane, 9-methyl-      | 282 | C20H42  | 013287-24-6 | 91   |
| 4    |    |   | Octacosane                 | 394 | C28H58  | 000630-02-4 | 81   |
| 5    |    |   | Tetradecane, 6,9-dimethyl- | 226 | C16H34  | 055045-13-1 | 81   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022420\  
Data File : BN009891.D  
Acq On : 25 Feb 2020 09:04  
Operator : CG/JU  
Sample : L1424-22DL 10X  
Misc :  
ALS Vial : 29 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
C5B30DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN021220MA.M  
Quant Title : SVOA CALIBRATION

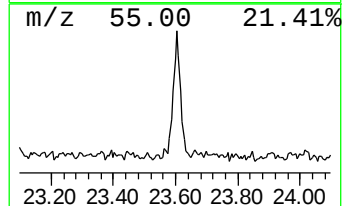
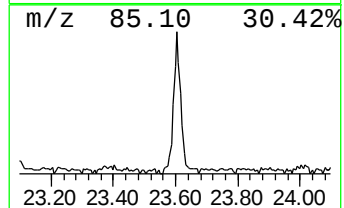
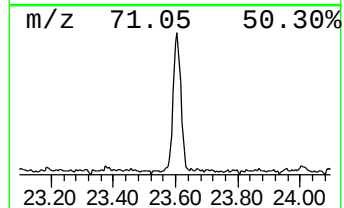
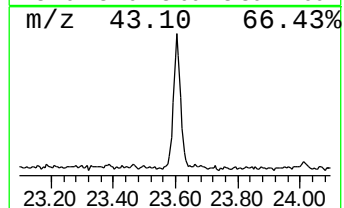
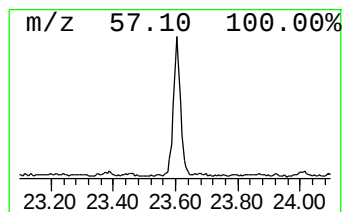
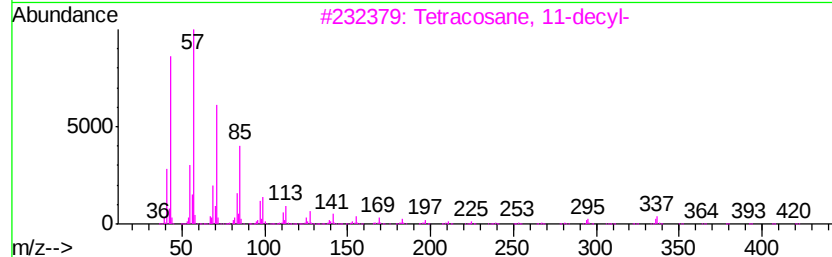
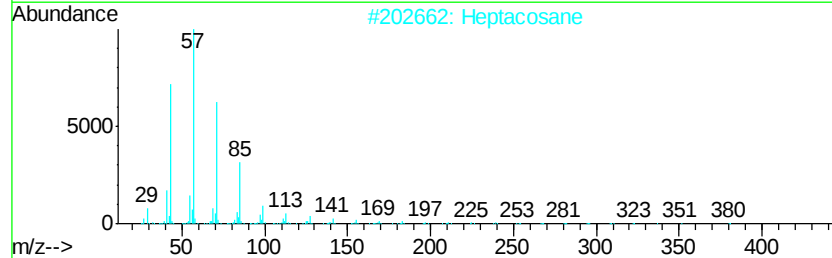
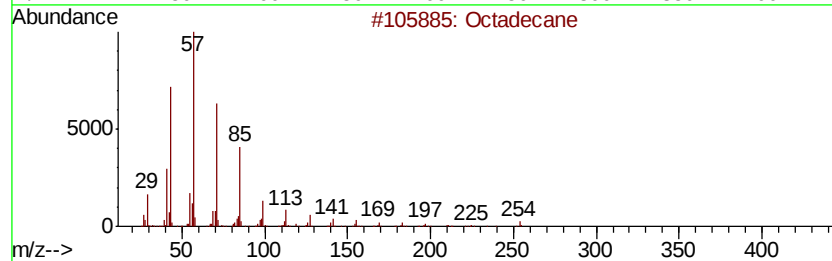
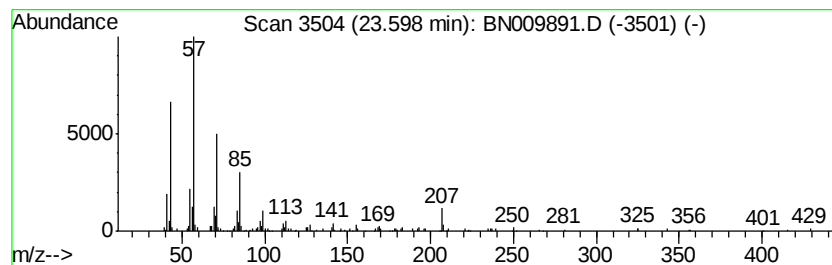
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 65 (DEL) Alkane: Straight-Chai... Concentration Rank 41

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 23.60 | 3.86 ng/ul | 317857 | Perylene-d12     | 23.10 |

| Hit# | of | 5 | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------------|-----|---------|-------------|------|
| 1    |    |   | Octadecane             | 254 | C18H38  | 000593-45-3 | 72   |
| 2    |    |   | Heptacosane            | 380 | C27H56  | 000593-49-7 | 72   |
| 3    |    |   | Tetracosane, 11-decyl- | 479 | C34H70  | 055429-84-0 | 62   |
| 4    |    |   | Heptadecane, 9-octyl-  | 352 | C25H52  | 007225-64-1 | 62   |
| 5    |    |   | Octacosane             | 394 | C28H58  | 000630-02-4 | 62   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022420\  
Data File : BN009891.D  
Acq On : 25 Feb 2020 09:04  
Operator : CG/JU  
Sample : L1424-22DL 10X  
Misc :  
ALS Vial : 29 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
C5B30DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN021220MA.M  
Quant Title : SVOA CALIBRATION

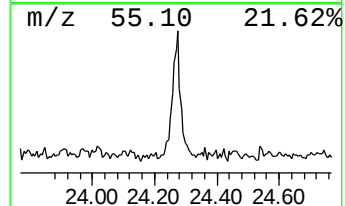
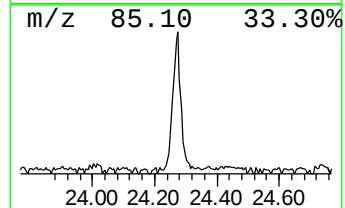
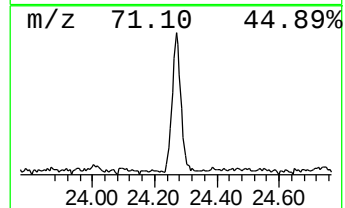
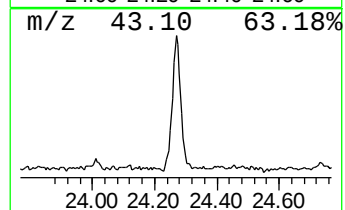
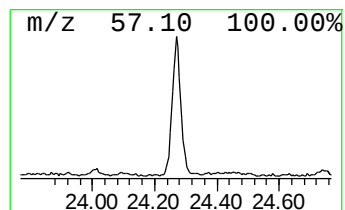
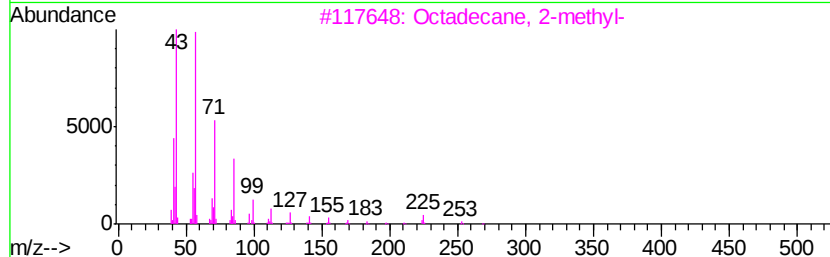
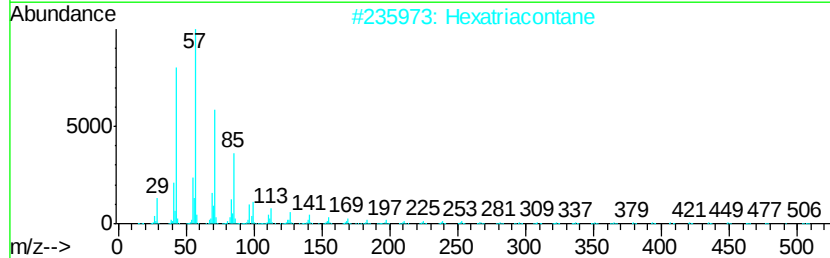
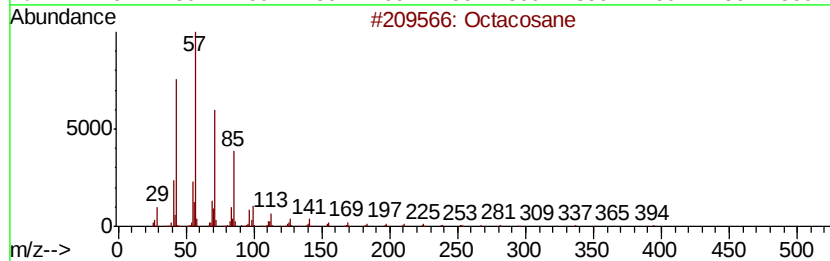
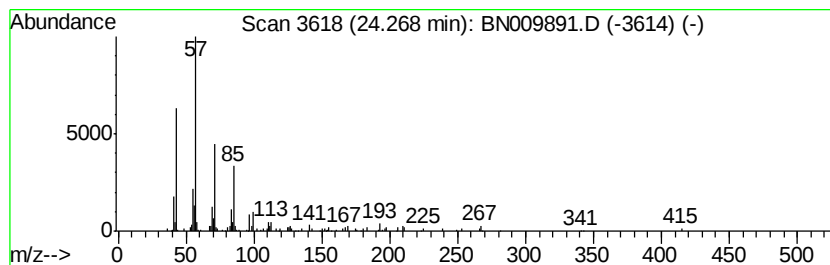
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 66 (DEL) Alkane: Straight-Chai... Concentration Rank 47

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 24.27 | 3.35 ng/ul | 275840 | Perylene-d12     | 23.10 |

| Hit# | of | 5 | Tentative ID                 | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------------------|-----|---------|-------------|------|
| 1    |    |   | Octacosane                   | 394 | C28H58  | 000630-02-4 | 76   |
| 2    |    |   | Hexatriacontane              | 507 | C36H74  | 000630-06-8 | 74   |
| 3    |    |   | Octadecane, 2-methyl-        | 268 | C19H40  | 001560-88-9 | 74   |
| 4    |    |   | Dodecane, 2-methyl-6-propyl- | 226 | C16H34  | 055045-08-4 | 72   |
| 5    |    |   | Tetratetracontane            | 619 | C44H90  | 007098-22-8 | 72   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_N\DATA\BN022420\  
 Data File : BN009891.D  
 Acq On : 25 Feb 2020 09:04  
 Operator : CG/JU  
 Sample : L1424-22DL 10X  
 Misc :  
 ALS Vial : 29 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 C5B30DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN021220MA.M  
 Quant Title : SVOA CALIBRATION

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

| TIC Top Hit name     | RT    | EstConc | Units | Response | --Internal Standard-- |       |         |      |
|----------------------|-------|---------|-------|----------|-----------------------|-------|---------|------|
|                      |       |         |       |          | #                     | RT    | Resp    | Conc |
| p-Xylene             | 5.45  | 5.5     | ng/ul | 227804   | 1                     | 7.39  | 836694  | 20.0 |
| Benzene, 1-ethyl-... | 6.50  | 3.5     | ng/ul | 145534   | 1                     | 7.39  | 836694  | 20.0 |
| Benzene, 1-ethyl-... | 6.56  | 4.7     | ng/ul | 197635   | 1                     | 7.39  | 836694  | 20.0 |
| Benzene, 1,2,3-tr... | 6.63  | 3.2     | ng/ul | 135521   | 1                     | 7.39  | 836694  | 20.0 |
| (DEL) Alkane: Cyc... | 7.66  | 3.2     | ng/ul | 132813   | 1                     | 7.39  | 836694  | 20.0 |
| Indane               | 7.75  | 4.9     | ng/ul | 203825   | 1                     | 7.39  | 836694  | 20.0 |
| Benzene, 1-methyl... | 7.92  | 5.0     | ng/ul | 210210   | 1                     | 7.39  | 836694  | 20.0 |
| Benzene, 1,3-diet... | 8.02  | 5.1     | ng/ul | 213680   | 1                     | 7.39  | 836694  | 20.0 |
| Benzene, (1-methy... | 8.18  | 3.2     | ng/ul | 132473   | 1                     | 7.39  | 836694  | 20.0 |
| Benzene, 2-ethyl-... | 8.33  | 4.1     | ng/ul | 170956   | 1                     | 7.39  | 836694  | 20.0 |
| p-Cymene             | 8.38  | 4.7     | ng/ul | 197082   | 1                     | 7.39  | 836694  | 20.0 |
| (DEL) Alkane: Str... | 8.60  | 8.0     | ng/ul | 335911   | 1                     | 7.39  | 836694  | 20.0 |
| unknown-01           | 8.72  | 3.9     | ng/ul | 163258   | 1                     | 7.39  | 836694  | 20.0 |
| o-Cymene             | 8.80  | 2.5     | ng/ul | 173405   | 2                     | 10.13 | 1411610 | 20.0 |
| Benzene, 1,2,4,5-... | 8.99  | 3.4     | ng/ul | 237154   | 2                     | 10.13 | 1411610 | 20.0 |
| 2,4-Dimethylstyrene  | 9.55  | 7.1     | ng/ul | 498071   | 2                     | 10.13 | 1411610 | 20.0 |
| Phenol, 3-chloro-    | 10.06 | 2.4     | ng/ul | 171957   | 2                     | 10.13 | 1411610 | 20.0 |
| 1H-Indene, 2,3-di... | 11.05 | 2.1     | ng/ul | 148984   | 2                     | 10.13 | 1411610 | 20.0 |
| (DEL) Alkane: Str... | 11.17 | 2.6     | ng/ul | 181151   | 2                     | 10.13 | 1411610 | 20.0 |
| (DEL) Alkane: Str... | 11.59 | 5.2     | ng/ul | 367830   | 2                     | 10.13 | 1411610 | 20.0 |
| 1-Decanol, 2,2-di... | 12.43 | 2.1     | ng/ul | 165623   | 3                     | 14.02 | 1610660 | 20.0 |
| Phenol, 3,5-dichl... | 12.75 | 10.8    | ng/ul | 870890   | 3                     | 14.02 | 1610660 | 20.0 |
| Naphthalene, 2-et... | 13.05 | 4.7     | ng/ul | 378407   | 3                     | 14.02 | 1610660 | 20.0 |
| Naphthalene, 1,3-... | 13.18 | 18.9    | ng/ul | 1522830  | 3                     | 14.02 | 1610660 | 20.0 |
| Naphthalene, 1,6-... | 13.40 | 12.4    | ng/ul | 1001040  | 3                     | 14.02 | 1610660 | 20.0 |
| (DEL) Alkane: Str... | 13.53 | 4.2     | ng/ul | 339104   | 3                     | 14.02 | 1610660 | 20.0 |
| Naphthalene, 2,3-... | 13.58 | 10.6    | ng/ul | 855925   | 3                     | 14.02 | 1610660 | 20.0 |
| Naphthalene, 2,6-... | 13.61 | 2.2     | ng/ul | 176626   | 3                     | 14.02 | 1610660 | 20.0 |
| (DEL) Alkane: Str... | 13.96 | 12.1    | ng/ul | 974718   | 3                     | 14.02 | 1610660 | 20.0 |
| Naphthalene, 2-(1... | 14.25 | 7.4     | ng/ul | 598233   | 3                     | 14.02 | 1610660 | 20.0 |
| Naphthalene, 2,3,... | 14.44 | 9.5     | ng/ul | 762907   | 3                     | 14.02 | 1610660 | 20.0 |
| Phenol, 2,3,5,6-t... | 14.59 | 7.2     | ng/ul | 583328   | 3                     | 14.02 | 1610660 | 20.0 |
| (DEL) Alkane: Str... | 14.92 | 10.5    | ng/ul | 843188   | 3                     | 14.02 | 1610660 | 20.0 |
| (DEL) Alkane: Str... | 15.48 | 2.9     | ng/ul | 199703   | 4                     | 16.79 | 1370630 | 20.0 |
| (DEL) Alkane: Str... | 15.79 | 17.6    | ng/ul | 1209580  | 4                     | 16.79 | 1370630 | 20.0 |
| (DEL) Alkane: Str... | 16.58 | 11.2    | ng/ul | 770605   | 4                     | 16.79 | 1370630 | 20.0 |
| (DEL) Alkane: Str... | 16.63 | 3.8     | ng/ul | 263308   | 4                     | 16.79 | 1370630 | 20.0 |
| (DEL) Alkane: Str... | 17.32 | 7.9     | ng/ul | 538477   | 4                     | 16.79 | 1370630 | 20.0 |
| (DEL) Alkane: Str... | 18.01 | 7.1     | ng/ul | 485639   | 4                     | 16.79 | 1370630 | 20.0 |
| (DEL) Alkane: Str... | 18.66 | 5.3     | ng/ul | 364657   | 4                     | 16.79 | 1370630 | 20.0 |
| (DEL) Alkane: Str... | 19.25 | 3.1     | ng/ul | 252725   | 5                     | 21.00 | 1638560 | 20.0 |
| (DEL) Alkane: Str... | 22.04 | 2.2     | ng/ul | 177041   | 5                     | 21.00 | 1638560 | 20.0 |
| (DEL) Alkane: Str... | 22.50 | 2.4     | ng/ul | 201004   | 6                     | 23.10 | 1646940 | 20.0 |
| (DEL) Alkane: Str... | 23.60 | 3.9     | ng/ul | 317857   | 6                     | 23.10 | 1646940 | 20.0 |
| (DEL) Alkane: Str... | 24.27 | 3.4     | ng/ul | 275840   | 6                     | 23.10 | 1646940 | 20.0 |