

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040819\  
 Data File : BM019822.D  
 Acq On : 09 Apr 2019 08:22  
 Operator : JU/SJ  
 Sample : K2188-10  
 Misc :  
 ALS Vial : 35 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 BF654

## 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\_M\METHODS\SOM-EPA-BM032219MA.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         | 4.769       | 298           | 303         | 308          | rBV      | 161448         | 236946        | 9.11%           | 0.540%        |
| 2         | 5.428       | 408           | 415         | 422          | rBV2     | 84750          | 162020        | 6.23%           | 0.369%        |
| 3         | 5.504       | 422           | 428         | 433          | rVV3     | 72206          | 120642        | 4.64%           | 0.275%        |
| 4         | 5.569       | 434           | 439         | 446          | rVB3     | 78446          | 172057        | 6.61%           | 0.392%        |
| 5         | 5.810       | 473           | 480         | 488          | rBV5     | 85784          | 235319        | 9.04%           | 0.536%        |
| 6         | 5.928       | 494           | 500         | 507          | rVB4     | 52513          | 109757        | 4.22%           | 0.250%        |
| 7         | 6.045       | 510           | 520         | 524          | rBV6     | 63375          | 175337        | 6.74%           | 0.399%        |
| 8         | 6.087       | 524           | 527         | 531          | rVV      | 131277         | 172462        | 6.63%           | 0.393%        |
| 9         | 6.175       | 538           | 542         | 557          | rVB4     | 118398         | 327850        | 12.60%          | 0.747%        |
| 10        | 6.345       | 566           | 571         | 577          | rVB3     | 64528          | 106767        | 4.10%           | 0.243%        |
| 11        | 6.522       | 597           | 601         | 606          | rVB3     | 76306          | 115501        | 4.44%           | 0.263%        |
| 12        | 6.634       | 614           | 620         | 622          | rBV3     | 80608          | 154999        | 5.96%           | 0.353%        |
| 13        | 6.669       | 622           | 626         | 634          | rVB3     | 379584         | 645448        | 24.80%          | 1.470%        |
| 14        | 6.757       | 636           | 641         | 653          | rBV2     | 307727         | 713842        | 27.43%          | 1.626%        |
| 15        | 6.851       | 653           | 657         | 662          | rVB3     | 55805          | 98501         | 3.79%           | 0.224%        |
| 16        | 6.916       | 662           | 668         | 676          | rBV2     | 354123         | 778406        | 29.91%          | 1.773%        |
| 17        | 7.016       | 681           | 685         | 693          | rVB2     | 121296         | 200907        | 7.72%           | 0.458%        |
| 18        | 7.098       | 693           | 699         | 704          | rBV      | 428992         | 762181        | 29.29%          | 1.736%        |
| 19        | 7.145       | 704           | 707         | 711          | rVB      | 113442         | 129241        | 4.97%           | 0.294%        |
| 20        | 7.192       | 711           | 715         | 720          | rVB      | 212710         | 305984        | 11.76%          | 0.697%        |
| 21        | 7.340       | 734           | 740         | 747          | rVB4     | 87189          | 172068        | 6.61%           | 0.392%        |
| 22        | 7.439       | 747           | 757         | 762          | rBV4     | 148496         | 379305        | 14.58%          | 0.864%        |
| 23        | 7.492       | 762           | 766         | 772          | rVB      | 253790         | 372709        | 14.32%          | 0.849%        |
| 24        | 7.557       | 772           | 777         | 787          | rVB      | 685436         | 1261335       | 48.47%          | 2.873%        |
| 25        | 7.657       | 787           | 794         | 798          | rBV2     | 184417         | 369055        | 14.18%          | 0.841%        |
| 26        | 7.769       | 806           | 813         | 816          | rBV3     | 91664          | 189494        | 7.28%           | 0.432%        |
| 27        | 7.804       | 816           | 819         | 826          | rVV5     | 104403         | 201806        | 7.75%           | 0.460%        |
| 28        | 7.957       | 837           | 845         | 852          | rVB7     | 50342          | 120429        | 4.63%           | 0.274%        |
| 29        | 8.075       | 852           | 865         | 875          | rVB4     | 349669         | 892147        | 34.28%          | 2.032%        |
| 30        | 8.175       | 875           | 882         | 887          | rBV      | 223442         | 364512        | 14.01%          | 0.830%        |
| 31        | 8.281       | 895           | 900         | 907          | rBV      | 490930         | 836521        | 32.15%          | 1.906%        |
| 32        | 8.345       | 907           | 911         | 916          | rBV2     | 78670          | 166962        | 6.42%           | 0.380%        |
| 33        | 8.486       | 928           | 935         | 940          | rBV3     | 139274         | 286032        | 10.99%          | 0.652%        |
| 34        | 8.539       | 940           | 944         | 947          | rVV      | 101421         | 142613        | 5.48%           | 0.325%        |

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 Stop Thrs : 0

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 Max Peaks: 100  
 Peak Location: TOP

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

Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
 Title : SVOA CALIBRATION

|    |        |      |      |      |      |        |         |        |        |
|----|--------|------|------|------|------|--------|---------|--------|--------|
| 35 | 8.616  | 947  | 957  | 966  | rVV6 | 193579 | 717886  | 27.59% | 1.635% |
| 36 | 8.734  | 969  | 977  | 988  | rVB4 | 616088 | 1515242 | 58.23% | 3.452% |
| 37 | 8.822  | 988  | 992  | 994  | rBV2 | 116487 | 168649  | 6.48%  | 0.384% |
| 38 | 8.869  | 994  | 1000 | 1006 | rVB6 | 173861 | 423394  | 16.27% | 0.964% |
| 39 | 8.934  | 1007 | 1011 | 1013 | rBV3 | 70799  | 99079   | 3.81%  | 0.226% |
| 40 | 8.969  | 1013 | 1017 | 1024 | rVB4 | 155626 | 285998  | 10.99% | 0.651% |
| 41 | 9.039  | 1024 | 1029 | 1034 | rBV7 | 74298  | 175368  | 6.74%  | 0.399% |
| 42 | 9.151  | 1044 | 1048 | 1053 | rBV2 | 154714 | 218025  | 8.38%  | 0.497% |
| 43 | 9.228  | 1055 | 1061 | 1066 | rVV2 | 188391 | 369480  | 14.20% | 0.842% |
| 44 | 9.339  | 1075 | 1080 | 1084 | rBV5 | 103158 | 194001  | 7.45%  | 0.442% |
| 45 | 9.439  | 1092 | 1097 | 1102 | rBV2 | 281162 | 502379  | 19.31% | 1.144% |
| 46 | 9.486  | 1102 | 1105 | 1109 | rVV3 | 174960 | 252844  | 9.72%  | 0.576% |
| 47 | 9.528  | 1109 | 1112 | 1119 | rVB5 | 134073 | 214966  | 8.26%  | 0.490% |
| 48 | 9.651  | 1129 | 1133 | 1137 | rBV3 | 80870  | 107039  | 4.11%  | 0.244% |
| 49 | 9.733  | 1142 | 1147 | 1154 | rVB3 | 254630 | 449732  | 17.28% | 1.024% |
| 50 | 9.792  | 1154 | 1157 | 1161 | rVB2 | 65856  | 96931   | 3.72%  | 0.221% |
| 51 | 9.898  | 1168 | 1175 | 1180 | rBV5 | 142416 | 359589  | 13.82% | 0.819% |
| 52 | 9.980  | 1184 | 1189 | 1195 | rBV  | 357249 | 768910  | 29.55% | 1.751% |
| 53 | 10.092 | 1203 | 1208 | 1211 | rBV6 | 55860  | 97764   | 3.76%  | 0.223% |
| 54 | 10.275 | 1235 | 1239 | 1243 | rBV  | 260568 | 413734  | 15.90% | 0.942% |
| 55 | 10.339 | 1245 | 1250 | 1255 | rVV  | 766069 | 1310542 | 50.36% | 2.985% |
| 56 | 10.380 | 1255 | 1257 | 1262 | rVV2 | 115286 | 160891  | 6.18%  | 0.366% |
| 57 | 10.457 | 1262 | 1270 | 1275 | rVB3 | 208275 | 434620  | 16.70% | 0.990% |
| 58 | 10.516 | 1275 | 1280 | 1290 | rBV4 | 164429 | 458682  | 17.63% | 1.045% |
| 59 | 10.692 | 1305 | 1310 | 1312 | rBV3 | 72666  | 122777  | 4.72%  | 0.280% |
| 60 | 10.786 | 1321 | 1326 | 1332 | rVB  | 216742 | 357567  | 13.74% | 0.814% |
| 61 | 11.216 | 1396 | 1399 | 1407 | rVB8 | 89327  | 180213  | 6.93%  | 0.411% |
| 62 | 11.322 | 1413 | 1417 | 1420 | rBV  | 194551 | 305158  | 11.73% | 0.695% |
| 63 | 11.492 | 1444 | 1446 | 1455 | rVB2 | 83504  | 141585  | 5.44%  | 0.323% |
| 64 | 11.580 | 1457 | 1461 | 1466 | rVB2 | 155574 | 253854  | 9.75%  | 0.578% |
| 65 | 11.733 | 1482 | 1487 | 1492 | rBV2 | 177782 | 288242  | 11.08% | 0.657% |
| 66 | 11.957 | 1523 | 1525 | 1529 | rVB3 | 91207  | 111133  | 4.27%  | 0.253% |
| 67 | 12.016 | 1530 | 1535 | 1538 | rBV  | 112660 | 180797  | 6.95%  | 0.412% |
| 68 | 12.233 | 1569 | 1572 | 1580 | rVB  | 108355 | 191295  | 7.35%  | 0.436% |
| 69 | 12.380 | 1592 | 1597 | 1601 | rBV5 | 79220  | 126786  | 4.87%  | 0.289% |
| 70 | 12.610 | 1632 | 1636 | 1641 | rVB2 | 77079  | 119062  | 4.58%  | 0.271% |
| 71 | 12.710 | 1648 | 1653 | 1658 | rVB2 | 167857 | 255144  | 9.80%  | 0.581% |

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 Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
 Title : SVOA CALIBRATION

|     |        |      |      |      |      |         |         |         |        |
|-----|--------|------|------|------|------|---------|---------|---------|--------|
| 72  | 13.004 | 1701 | 1703 | 1710 | rVB  | 110848  | 148007  | 5.69%   | 0.337% |
| 73  | 13.398 | 1762 | 1770 | 1774 | rBV4 | 101557  | 226570  | 8.71%   | 0.516% |
| 74  | 13.533 | 1788 | 1793 | 1798 | rBV  | 93853   | 173278  | 6.66%   | 0.395% |
| 75  | 13.592 | 1798 | 1803 | 1807 | rBV2 | 106293  | 192207  | 7.39%   | 0.438% |
| 76  | 13.639 | 1807 | 1811 | 1816 | rBV  | 705635  | 1020046 | 39.20%  | 2.324% |
| 77  | 13.916 | 1852 | 1858 | 1869 | rVB  | 790366  | 1308499 | 50.28%  | 2.981% |
| 78  | 14.027 | 1872 | 1877 | 1883 | rVB5 | 67873   | 124985  | 4.80%   | 0.285% |
| 79  | 14.098 | 1885 | 1889 | 1894 | rVB  | 99898   | 133642  | 5.14%   | 0.304% |
| 80  | 14.221 | 1904 | 1910 | 1916 | rVB2 | 1182395 | 1777272 | 68.30%  | 4.048% |
| 81  | 14.510 | 1954 | 1959 | 1969 | rBV4 | 119903  | 343752  | 13.21%  | 0.783% |
| 82  | 14.610 | 1972 | 1976 | 1980 | rVV5 | 63311   | 144441  | 5.55%   | 0.329% |
| 83  | 14.651 | 1981 | 1983 | 1987 | rVV2 | 79324   | 97426   | 3.74%   | 0.222% |
| 84  | 14.698 | 1988 | 1991 | 1999 | rVB4 | 76529   | 161669  | 6.21%   | 0.368% |
| 85  | 15.057 | 2049 | 2052 | 2056 | rVB2 | 91577   | 100790  | 3.87%   | 0.230% |
| 86  | 15.227 | 2075 | 2081 | 2086 | rBV  | 1015888 | 1542256 | 59.26%  | 3.513% |
| 87  | 15.463 | 2117 | 2121 | 2129 | rVB4 | 105784  | 193697  | 7.44%   | 0.441% |
| 88  | 15.945 | 2196 | 2203 | 2210 | rBV2 | 318319  | 613699  | 23.58%  | 1.398% |
| 89  | 16.721 | 2332 | 2335 | 2338 | rVB2 | 100132  | 125128  | 4.81%   | 0.285% |
| 90  | 16.768 | 2338 | 2343 | 2347 | rBV  | 277888  | 423403  | 16.27%  | 0.964% |
| 91  | 16.986 | 2375 | 2380 | 2385 | rBV  | 1400129 | 1953317 | 75.06%  | 4.449% |
| 92  | 17.086 | 2392 | 2397 | 2403 | rVB  | 1040003 | 1467336 | 56.39%  | 3.342% |
| 93  | 17.568 | 2476 | 2479 | 2483 | rVB2 | 209739  | 270411  | 10.39%  | 0.616% |
| 94  | 17.780 | 2513 | 2515 | 2519 | rVB2 | 189404  | 195373  | 7.51%   | 0.445% |
| 95  | 17.880 | 2526 | 2532 | 2535 | rBV2 | 509936  | 822302  | 31.60%  | 1.873% |
| 96  | 18.262 | 2594 | 2597 | 2602 | rVB  | 172155  | 219667  | 8.44%   | 0.500% |
| 97  | 18.792 | 2684 | 2687 | 2694 | rBV5 | 312733  | 651079  | 25.02%  | 1.483% |
| 98  | 19.398 | 2785 | 2790 | 2794 | rBV  | 1452207 | 2167745 | 83.30%  | 4.938% |
| 99  | 19.768 | 2850 | 2853 | 2860 | rVB  | 663115  | 791421  | 30.41%  | 1.803% |
| 100 | 21.203 | 3093 | 3097 | 3103 | rBV  | 2053770 | 2602314 | 100.00% | 5.928% |

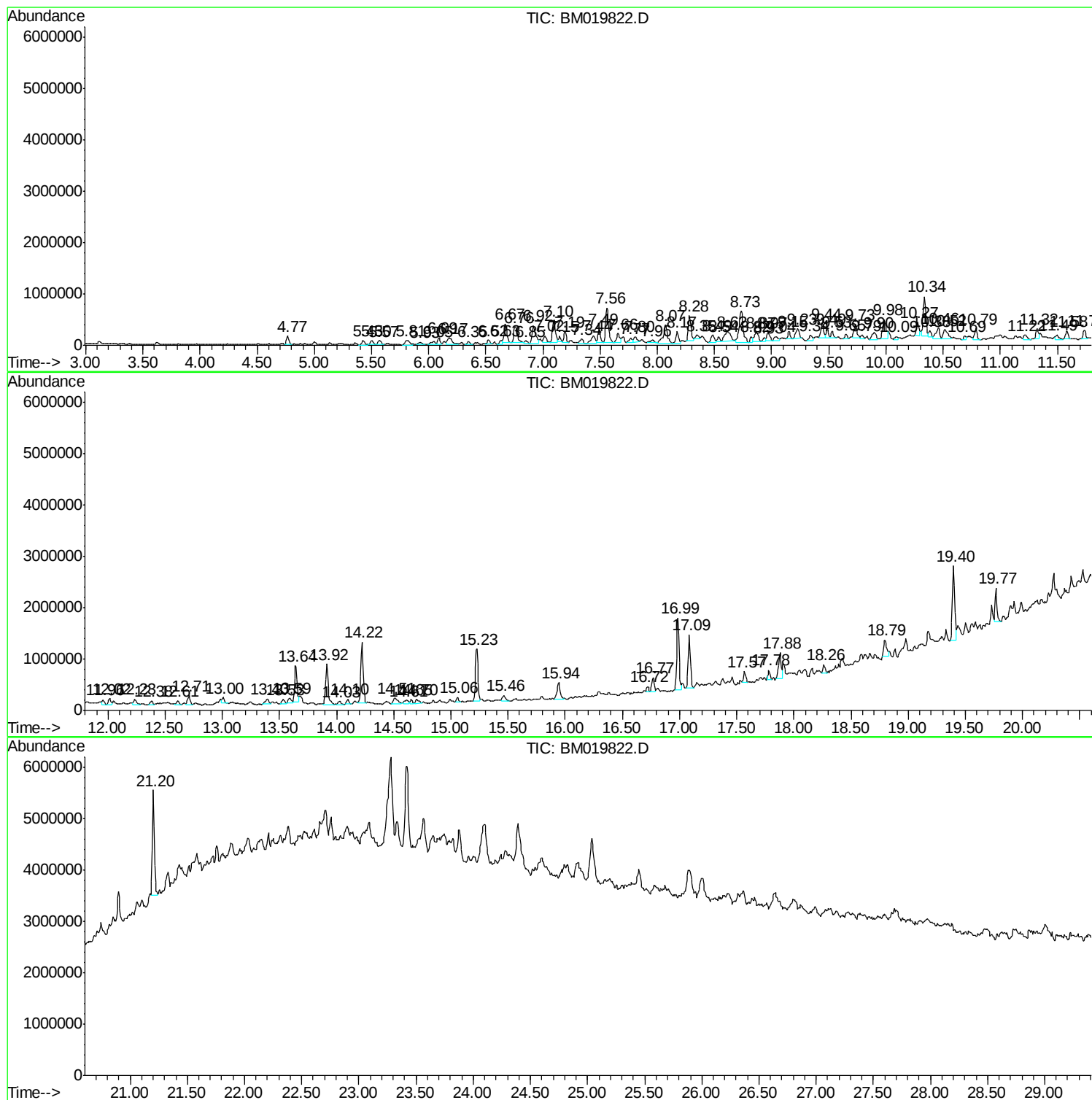
Sum of corrected areas: 43900245

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040819\  
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

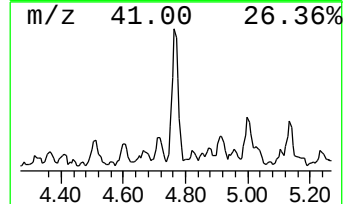
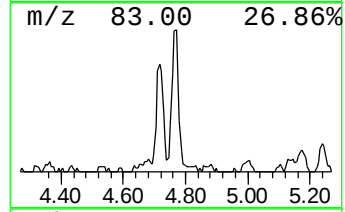
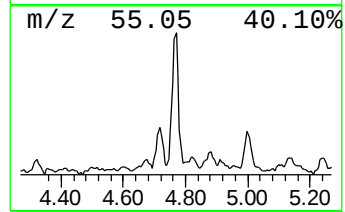
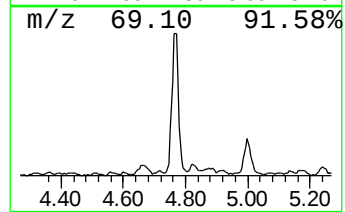
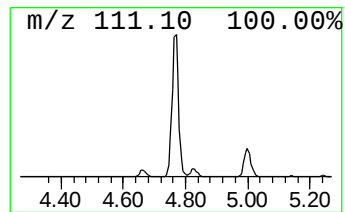
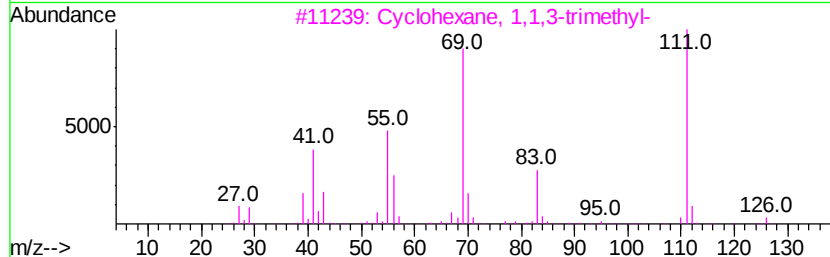
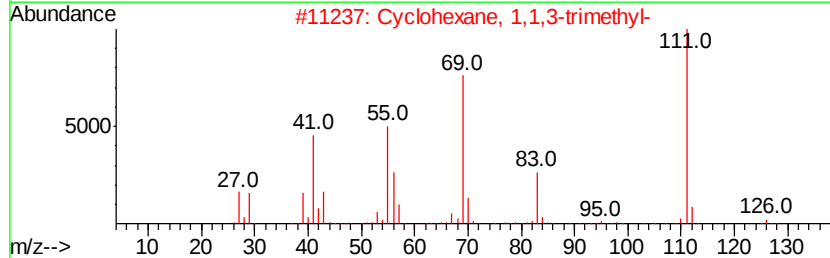
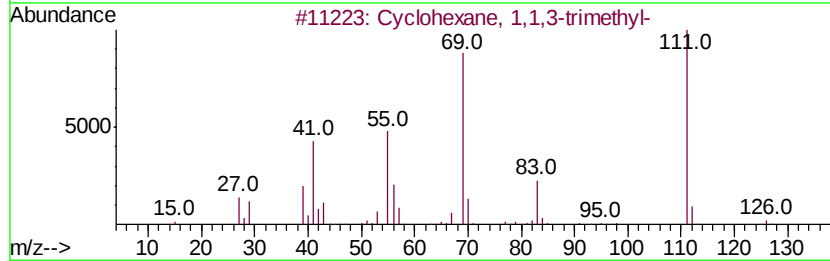
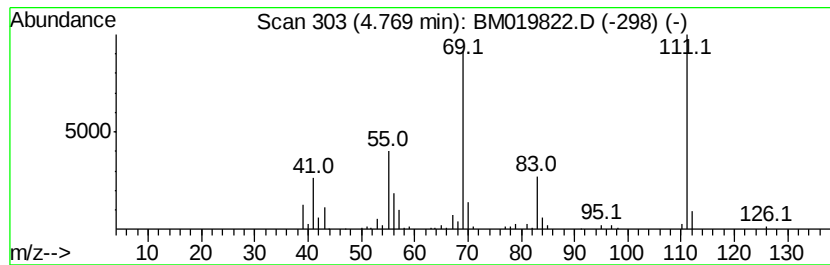
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 1 (DEL) Alkane: Cyclic4.77 Concentration Rank 26

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 4.77 | 3.76 ng/ul | 236946 | 1,4-Dichlorobenzene-d4 | 7.56 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Cyclohexane, 1,1,3-trimethyl-       | 126 | C9H18   | 003073-66-3 | 94   |
| 2    |    |   | Cyclohexane, 1,1,3-trimethyl-       | 126 | C9H18   | 003073-66-3 | 94   |
| 3    |    |   | Cyclohexane, 1,1,3-trimethyl-       | 126 | C9H18   | 003073-66-3 | 90   |
| 4    |    |   | Cyclohexane, 1-ethyl-1,4-dimethy... | 140 | C10H20  | 062238-30-6 | 78   |
| 5    |    |   | Cyclohexane, 1-ethyl-1,4-dimethy... | 140 | C10H20  | 062238-32-8 | 78   |



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Quant Title : SVOA CALIBRATION

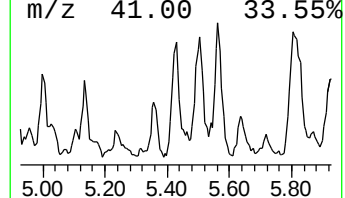
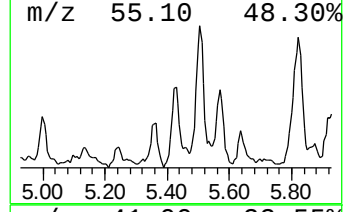
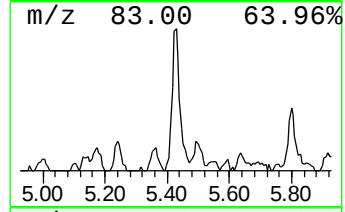
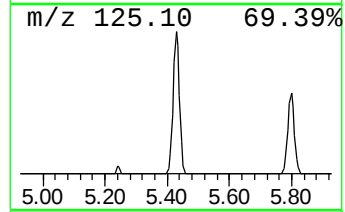
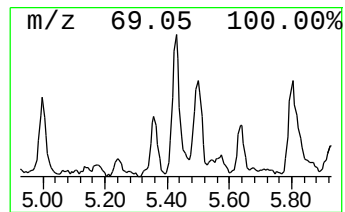
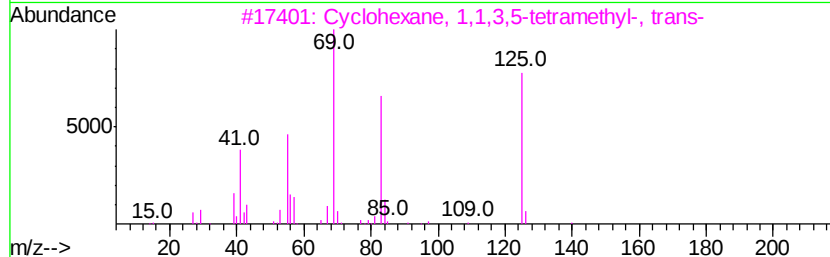
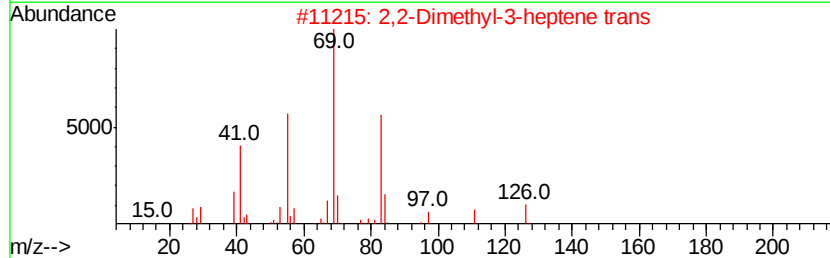
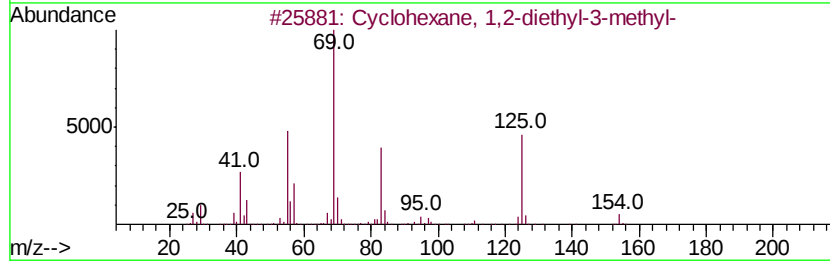
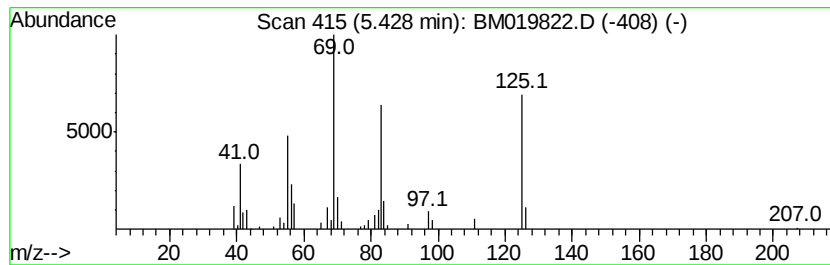
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 2 (DEL) Alkane: Cyclic5.43 Concentration Rank 45

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 5.43 | 2.57 ng/ul | 162020 | 1,4-Dichlorobenzene-d4 | 7.56 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Cyclohexane, 1,2-diethyl-3-methyl-  | 154 | C11H22  | 061141-80-8 | 78   |
| 2    |    | 2,2-Dimethyl-3-heptene trans        | 126 | C9H18   | 019550-75-5 | 64   |
| 3    |    | Cyclohexane, 1,1,3,5-tetramethyl... | 140 | C10H20  | 050876-31-8 | 64   |
| 4    |    | Cyclohexane, 1,1,3,5-tetramethyl... | 140 | C10H20  | 050876-31-8 | 64   |
| 5    |    | Cyclohexane, 1,1,3,5-tetramethyl... | 140 | C10H20  | 050876-32-9 | 64   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040819\  
Data File : BM019822.D  
Acq On : 09 Apr 2019 08:22  
Operator : JU/SJ  
Sample : K2188-10  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF654

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

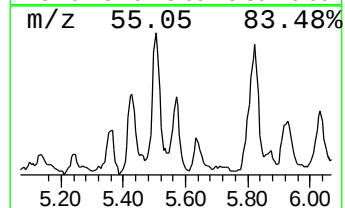
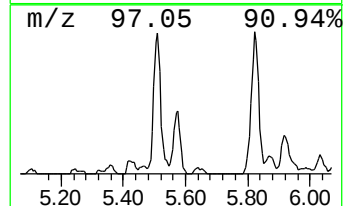
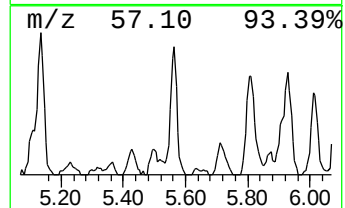
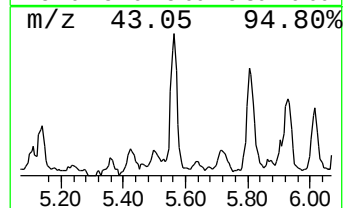
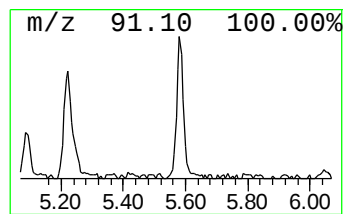
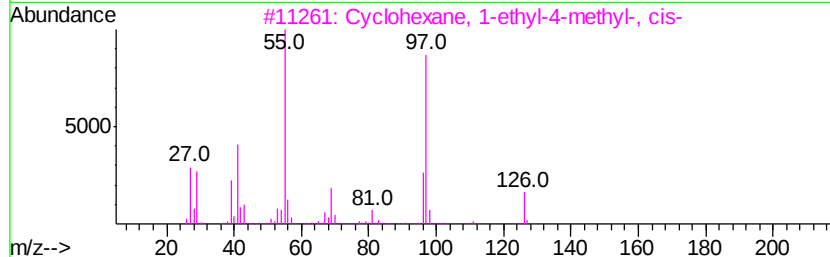
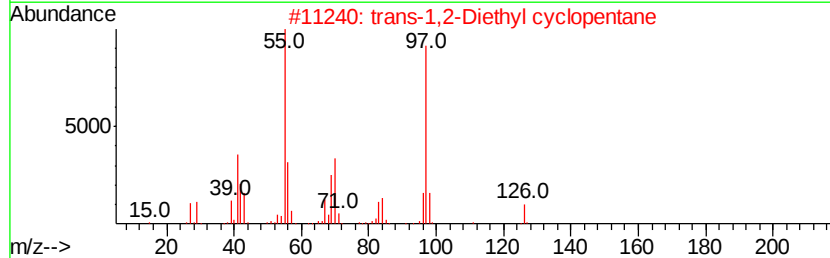
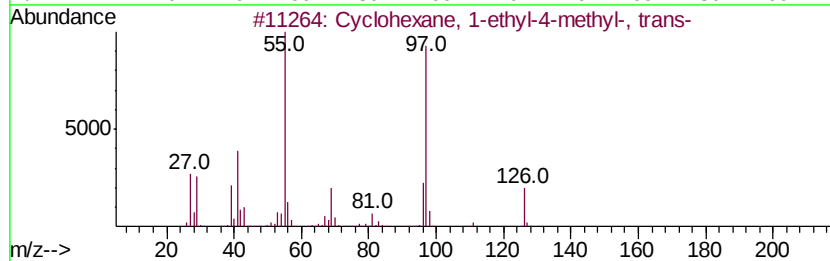
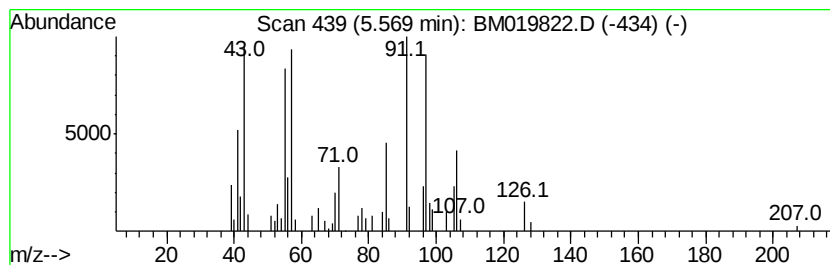
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 3 (DEL) Alkane: Cyclic5.57 Concentration Rank 41

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 5.57 | 2.73 ng/ul | 172057 | 1,4-Dichlorobenzene-d4 | 7.56 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Cyclohexane, 1-ethyl-4-methyl-, ... | 126 | C9H18   | 006236-88-0 | 38   |
| 2    |    | trans-1,2-Diethyl cyclopentane      | 126 | C9H18   | 000932-40-1 | 38   |
| 3    |    | Cyclohexane, 1-ethyl-4-methyl-, ... | 126 | C9H18   | 004926-78-7 | 38   |
| 4    |    | Cyclohexane, 1-ethyl-4-methyl-, ... | 126 | C9H18   | 006236-88-0 | 38   |
| 5    |    | Cyclohexane, 1-ethyl-2-methyl-      | 126 | C9H18   | 003728-54-9 | 38   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040819\  
Data File : BM019822.D  
Acq On : 09 Apr 2019 08:22  
Operator : JU/SJ  
Sample : K2188-10  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF654

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

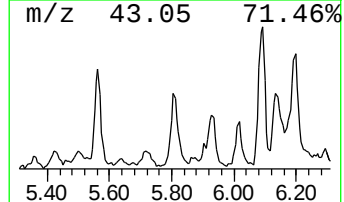
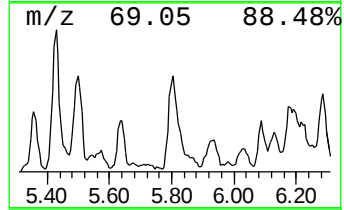
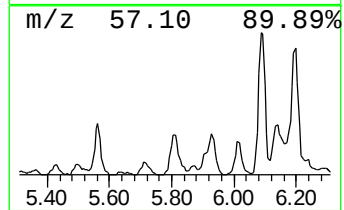
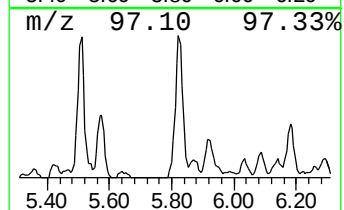
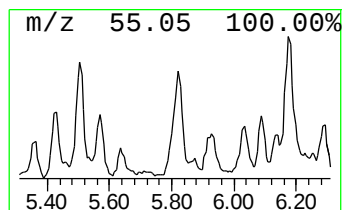
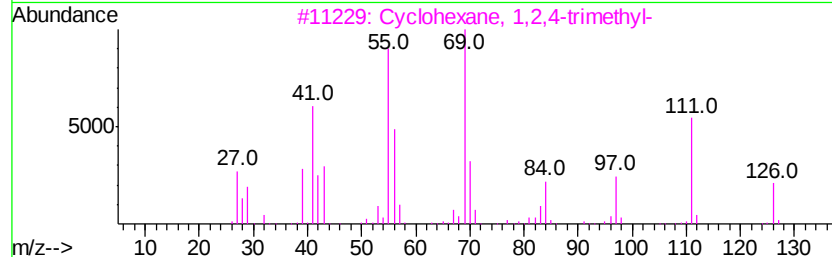
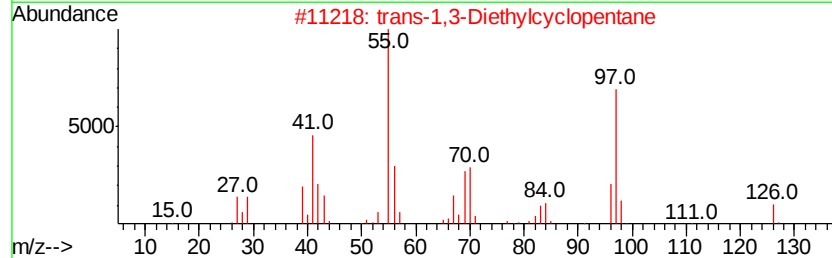
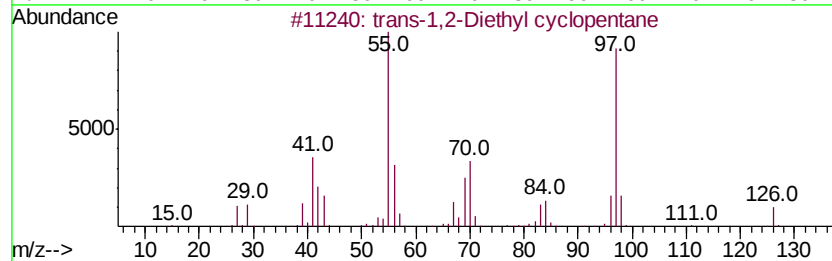
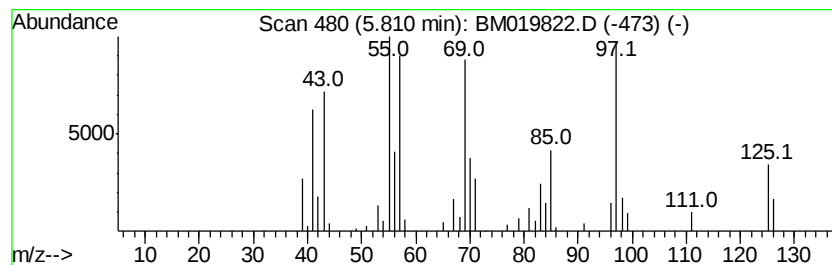
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 4 (DEL) Alkane: Cyclic5.81 Concentration Rank 27

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 5.81 | 3.73 ng/ul | 235319 | 1,4-Dichlorobenzene-d4 | 7.56 |

| Hit# | of                             | 5   | Tentative ID | MW           | MolForm | CAS# | Qual |
|------|--------------------------------|-----|--------------|--------------|---------|------|------|
| 1    | trans-1,2-Diethyl cyclopentane | 126 | C9H18        | 000932-40-1  | 42      |      |      |
| 2    | trans-1,3-Diethylcyclopentane  | 126 | C9H18        | 1000113-87-1 | 38      |      |      |
| 3    | Cyclohexane, 1,2,4-trimethyl-  | 126 | C9H18        | 002234-75-5  | 38      |      |      |
| 4    | 3-Hexene, 2-methyl-, (Z)-      | 98  | C7H14        | 015840-60-5  | 30      |      |      |
| 5    | 1-Ethyl-4-methylcyclohexane    | 126 | C9H18        | 003728-56-1  | 30      |      |      |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040819\  
Data File : BM019822.D  
Acq On : 09 Apr 2019 08:22  
Operator : JU/SJ  
Sample : K2188-10  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF654

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

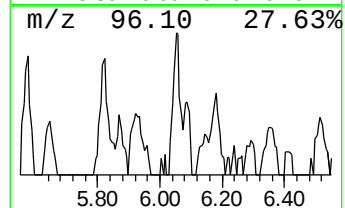
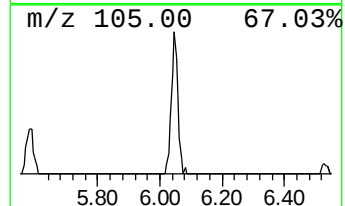
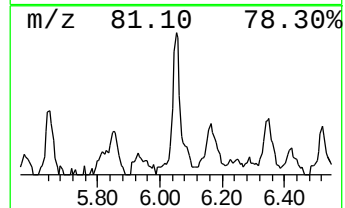
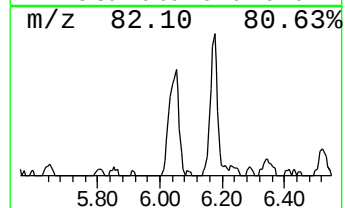
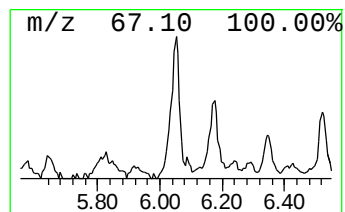
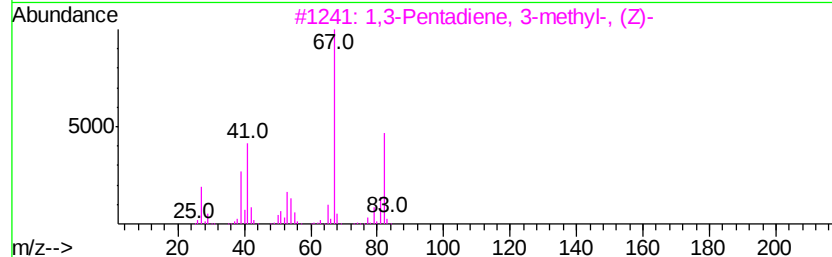
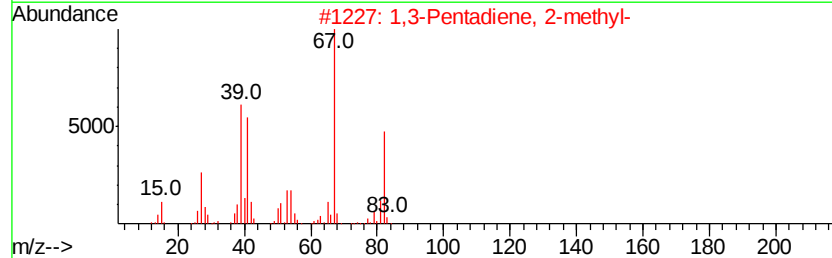
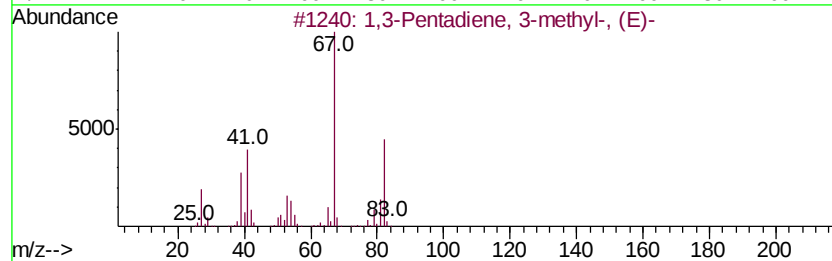
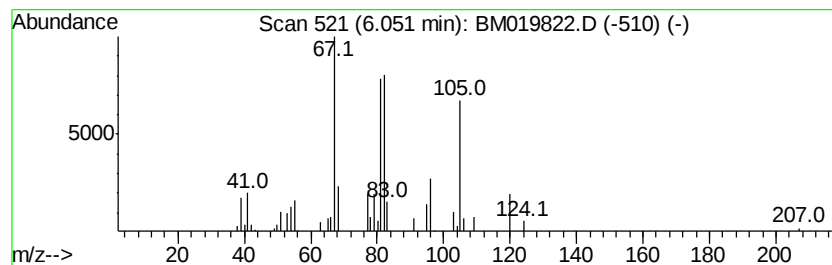
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 5 unknown-01 Concentration Rank 36

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 6.05 | 2.78 ng/ul | 175337 | 1,4-Dichlorobenzene-d4 | 7.56 |

| Hit# | of | Tentative ID                    | MW | MolForm | CAS#        | Qual |
|------|----|---------------------------------|----|---------|-------------|------|
| 1    | 5  | 1,3-Pentadiene, 3-methyl-, (E)- | 82 | C6H10   | 002787-43-1 | 38   |
| 2    |    | 1,3-Pentadiene, 2-methyl-       | 82 | C6H10   | 001118-58-7 | 35   |
| 3    |    | 1,3-Pentadiene, 3-methyl-, (Z)- | 82 | C6H10   | 002787-45-3 | 35   |
| 4    |    | 2,4-Hexadiene, (E,E)-           | 82 | C6H10   | 005194-51-4 | 35   |
| 5    |    | 2,4-Hexadiene                   | 82 | C6H10   | 000592-46-1 | 35   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040819\  
Data File : BM019822.D  
Acq On : 09 Apr 2019 08:22  
Operator : JU/SJ  
Sample : K2188-10  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF654

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

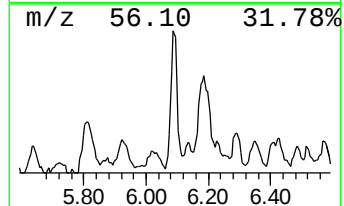
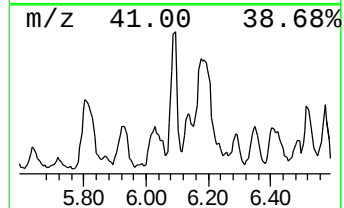
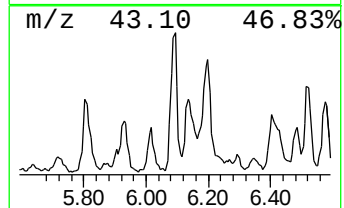
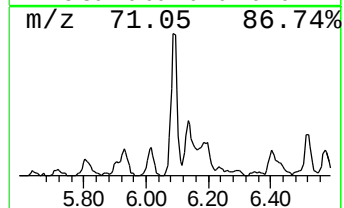
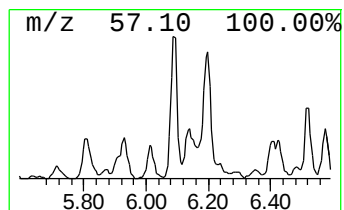
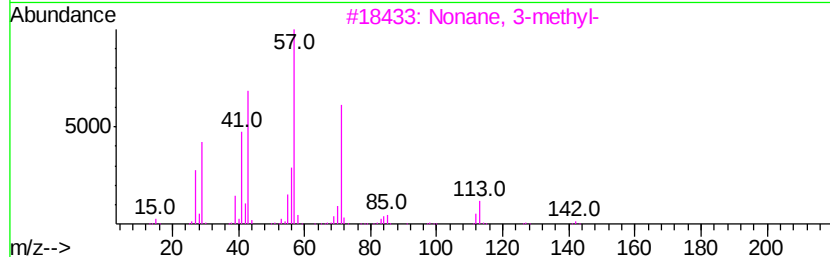
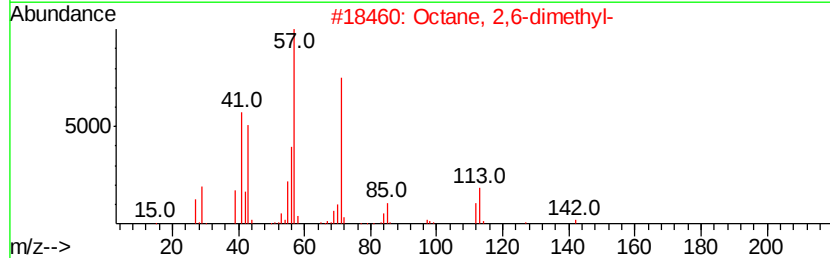
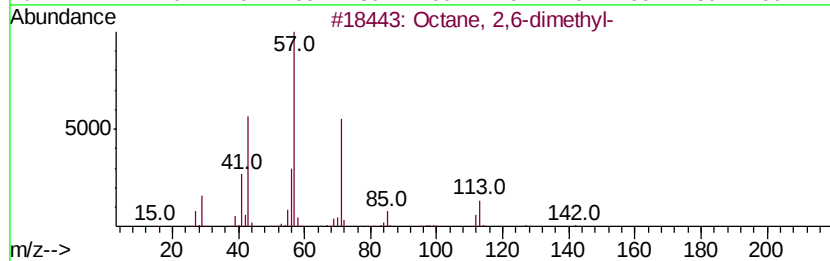
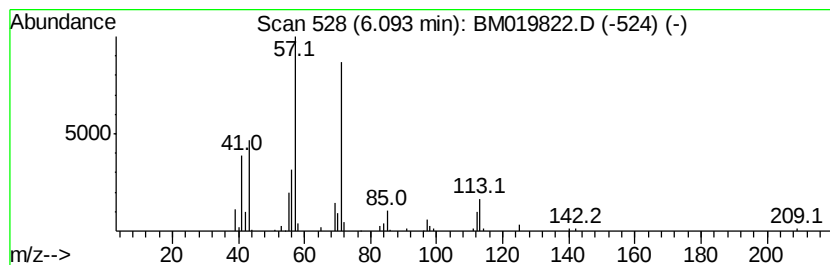
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 39

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 6.09 | 2.73 ng/ul | 172462 | 1,4-Dichlorobenzene-d4 | 7.56 |

| Hit# | of | 5 | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------|-----|---------|-------------|------|
| 1    |    |   | Octane, 2,6-dimethyl- | 142 | C10H22  | 002051-30-1 | 90   |
| 2    |    |   | Octane, 2,6-dimethyl- | 142 | C10H22  | 002051-30-1 | 86   |
| 3    |    |   | Nonane, 3-methyl-     | 142 | C10H22  | 005911-04-6 | 78   |
| 4    |    |   | Octane, 2,6-dimethyl- | 142 | C10H22  | 002051-30-1 | 78   |
| 5    |    |   | Nonane, 3-methyl-     | 142 | C10H22  | 005911-04-6 | 72   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040819\  
Data File : BM019822.D  
Acq On : 09 Apr 2019 08:22  
Operator : JU/SJ  
Sample : K2188-10  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

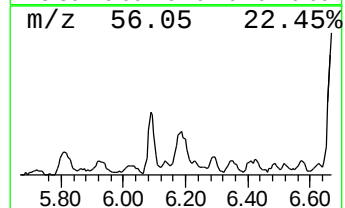
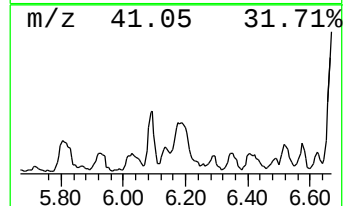
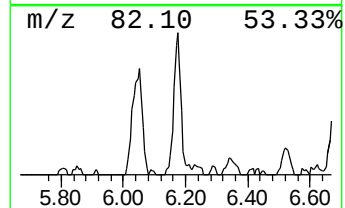
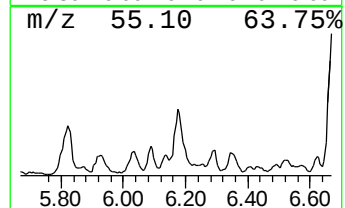
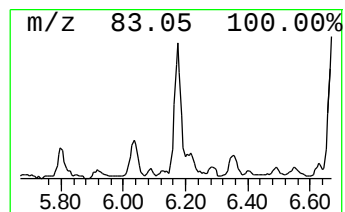
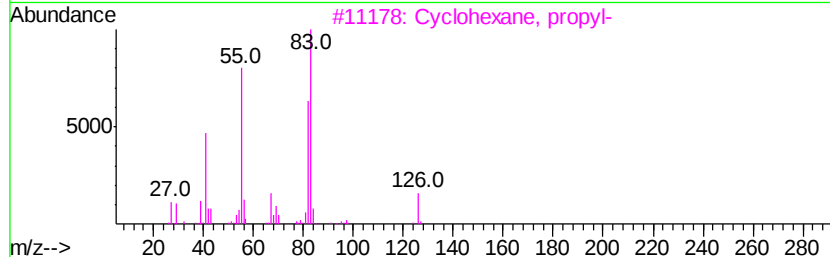
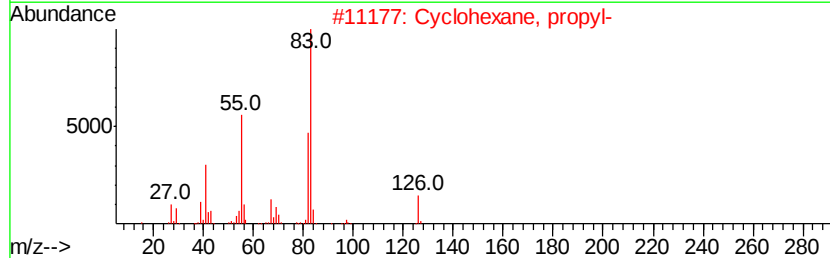
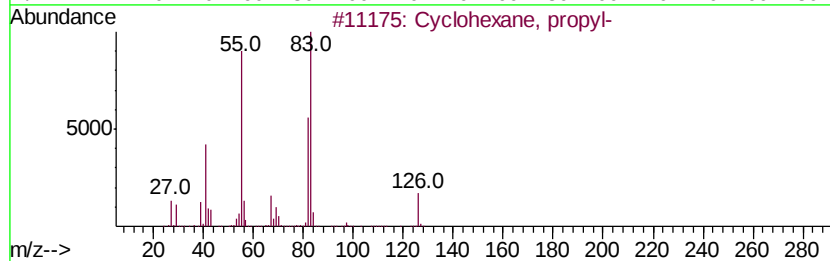
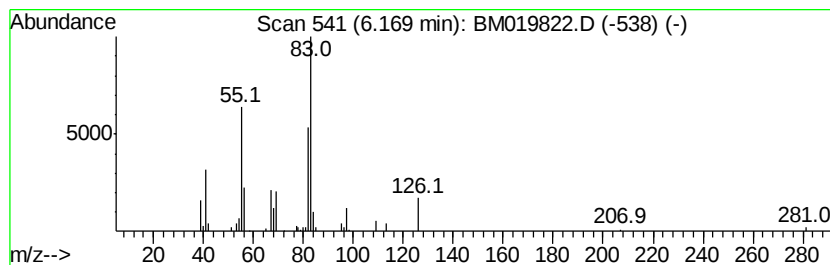
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 7 (DEL) Alkane: Cyclic6.17 Concentration Rank 19

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 6.17 | 5.20 ng/ul | 327850 | 1,4-Dichlorobenzene-d4 | 7.56 |

| Hit# | of | Tentative ID         | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------|-----|---------|-------------|------|
| 1    | 5  | Cyclohexane, propyl- | 126 | C9H18   | 001678-92-8 | 64   |
| 2    |    | Cyclohexane, propyl- | 126 | C9H18   | 001678-92-8 | 64   |
| 3    |    | Cyclohexane, propyl- | 126 | C9H18   | 001678-92-8 | 59   |
| 4    |    | Cyclohexane, propyl- | 126 | C9H18   | 001678-92-8 | 53   |
| 5    |    | Cyclohexane, ethyl-  | 112 | C8H16   | 001678-91-7 | 53   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040819\  
Data File : BM019822.D  
Acq On : 09 Apr 2019 08:22  
Operator : JU/SJ  
Sample : K2188-10  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampled :  
BF654

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

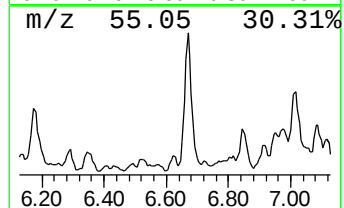
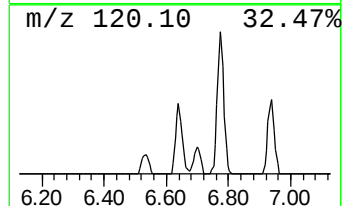
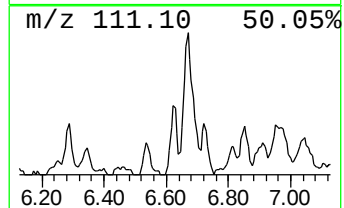
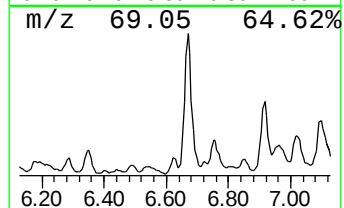
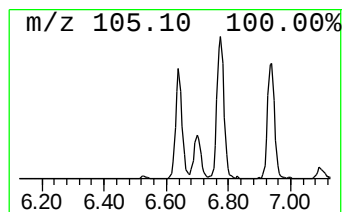
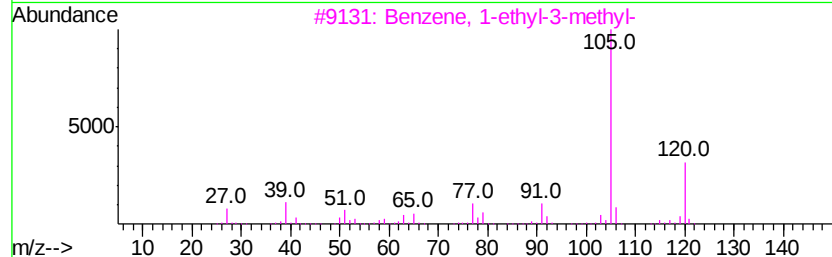
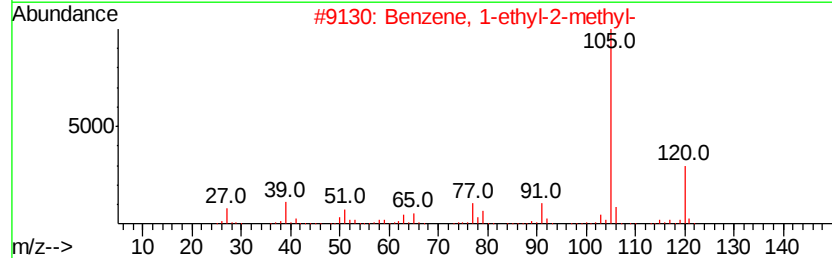
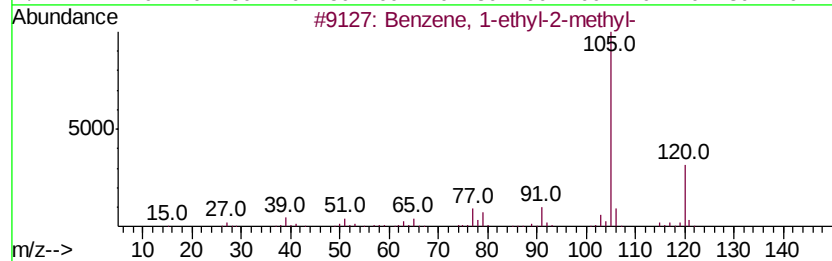
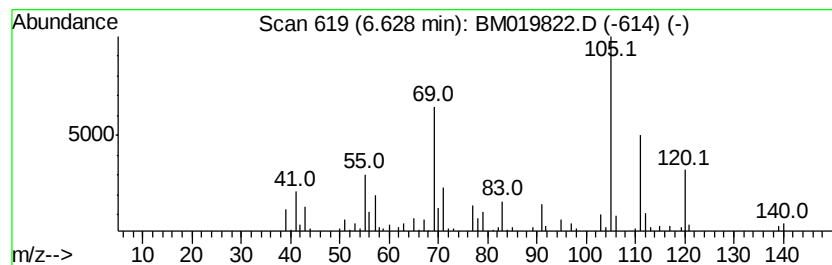
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 8 Benzene, 1-ethyl-2-methyl- Concentration Rank 47

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 6.63 | 2.46 ng/ul | 154999 | 1,4-Dichlorobenzene-d4 | 7.56 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 93   |
| 2    |    |   | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 93   |
| 3    |    |   | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 93   |
| 4    |    |   | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 93   |
| 5    |    |   | Benzene, 1-ethyl-4-methyl- | 120 | C9H12   | 000622-96-8 | 81   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040819\  
Data File : BM019822.D  
Acq On : 09 Apr 2019 08:22  
Operator : JU/SJ  
Sample : K2188-10  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF654

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

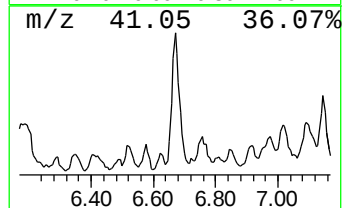
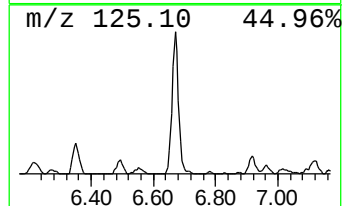
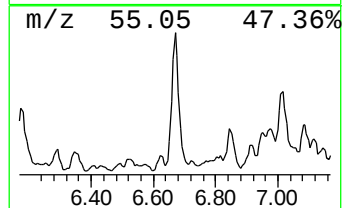
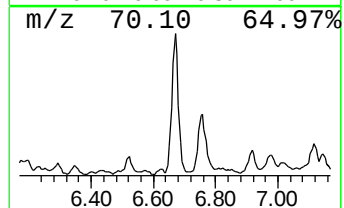
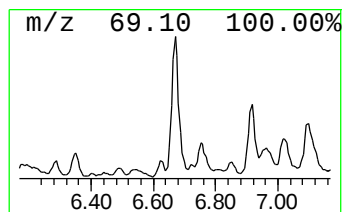
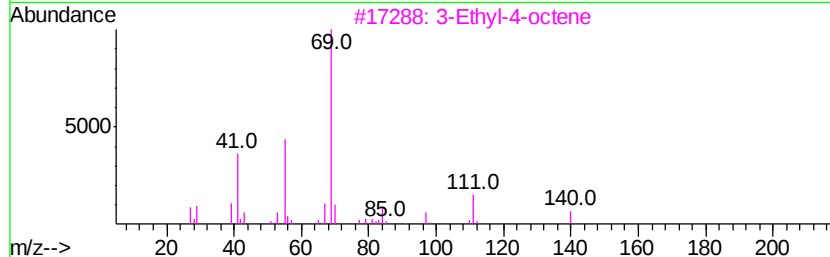
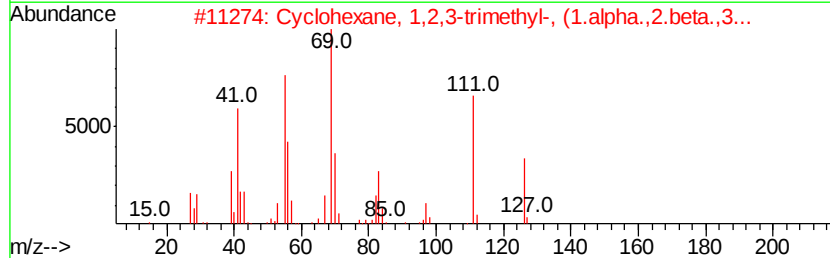
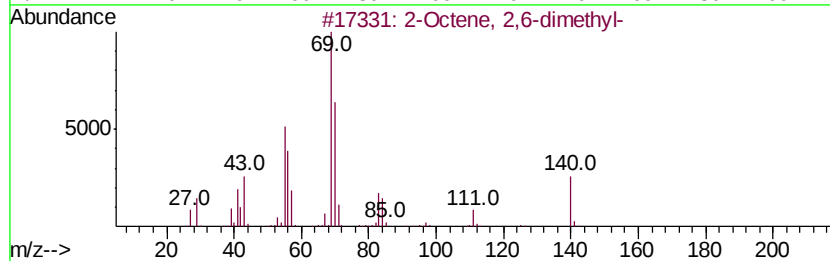
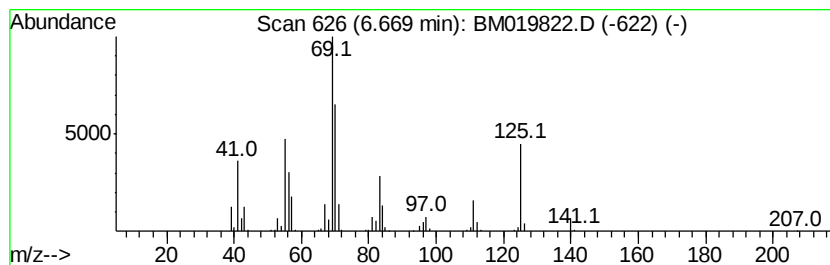
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 9 (DEL) Alkane: Cyclic6.67 Concentration Rank 3

| R.T. | EstConc     | Area   | Relative to ISTD       | R.T. |
|------|-------------|--------|------------------------|------|
| 6.67 | 10.23 ng/ul | 645448 | 1,4-Dichlorobenzene-d4 | 7.56 |

| Hit# | of | Tentative ID                         | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------------|-----|---------|-------------|------|
| 1    | 5  | 2-Octene, 2,6-dimethyl-              | 140 | C10H20  | 004057-42-5 | 62   |
| 2    |    | Cyclohexane, 1,2,3-trimethyl-, (...) | 126 | C9H18   | 001678-81-5 | 50   |
| 3    |    | 3-Ethyl-4-octene                     | 140 | C10H20  | 019781-32-9 | 50   |
| 4    |    | 1-Hexene, 3,3,5-trimethyl-           | 126 | C9H18   | 013427-43-5 | 49   |
| 5    |    | 1-Hexene, 3,3,5-trimethyl-           | 126 | C9H18   | 013427-43-5 | 47   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040819\  
Data File : BM019822.D  
Acq On : 09 Apr 2019 08:22  
Operator : JU/SJ  
Sample : K2188-10  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF654

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

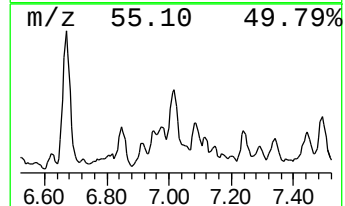
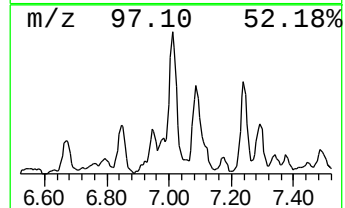
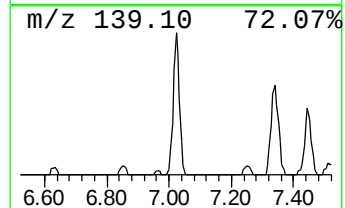
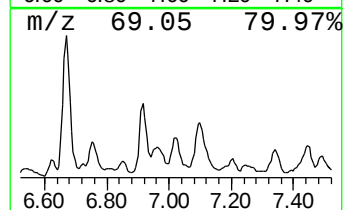
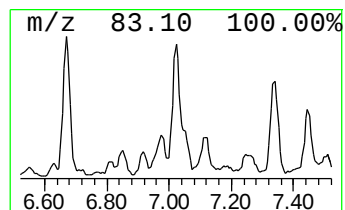
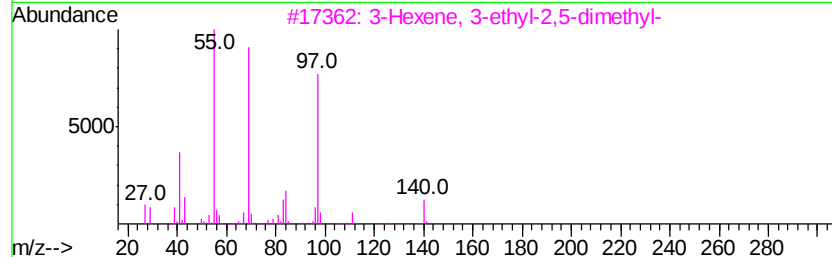
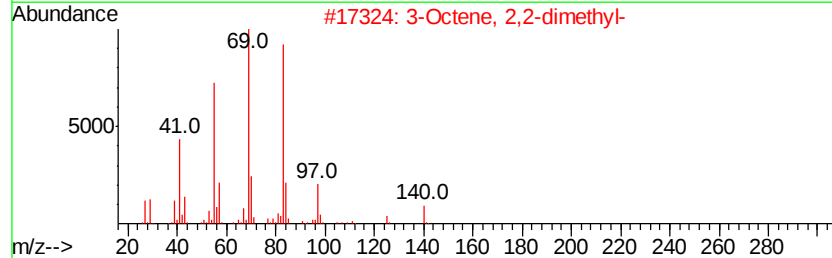
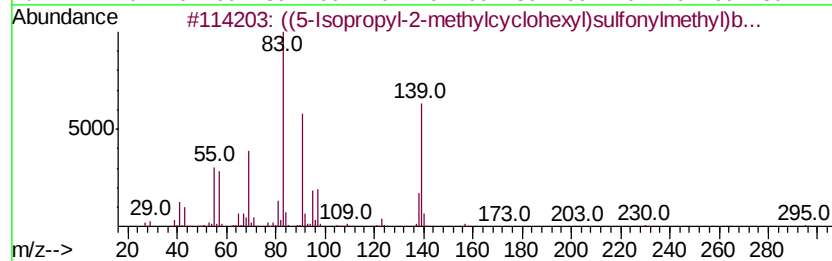
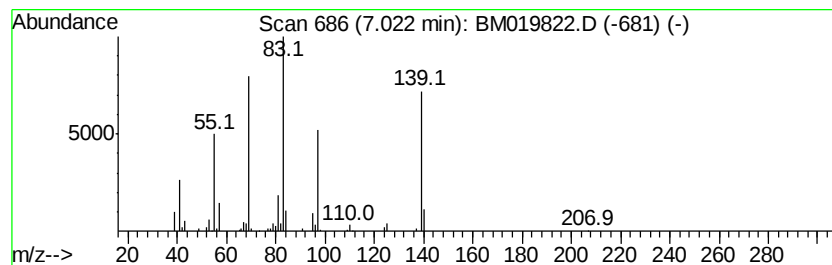
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 10 ((5-Isopropyl-2-methylcyclo... Concentration Rank 31

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 7.02 | 3.19 ng/ul | 200907 | 1,4-Dichlorobenzene-d4 | 7.56 |

| Hit# | of | Tentative ID                         | MW  | MolForm   | CAS#        | Qual |
|------|----|--------------------------------------|-----|-----------|-------------|------|
| 1    | 5  | ((5-Isopropyl-2-methylcyclohexyl...) | 294 | C17H26O2S | 093542-25-7 | 50   |
| 2    |    | 3-Octene, 2,2-dimethyl-              | 140 | C10H20    | 086869-76-3 | 47   |
| 3    |    | 3-Hexene, 3-ethyl-2,5-dimethyl-      | 140 | C10H20    | 062338-08-3 | 38   |
| 4    |    | Cyclohexanol, 5-methyl-2-(1-meth...  | 358 | C20H38O3S | 026510-92-9 | 27   |
| 5    |    | 1-Propanone, 1-cyclohexyl-           | 140 | C9H16O    | 001123-86-0 | 25   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040819\  
Data File : BM019822.D  
Acq On : 09 Apr 2019 08:22  
Operator : JU/SJ  
Sample : K2188-10  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF654

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

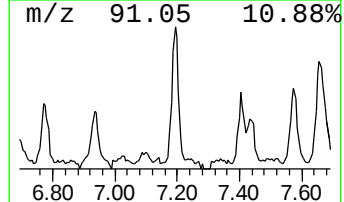
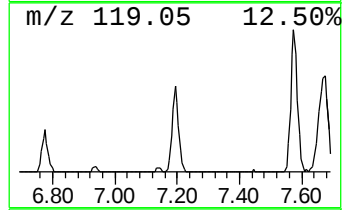
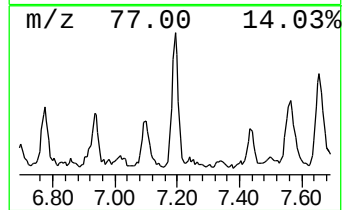
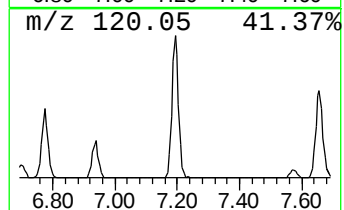
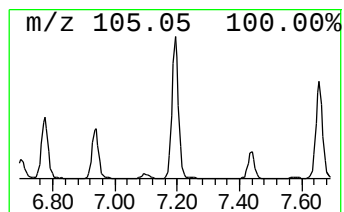
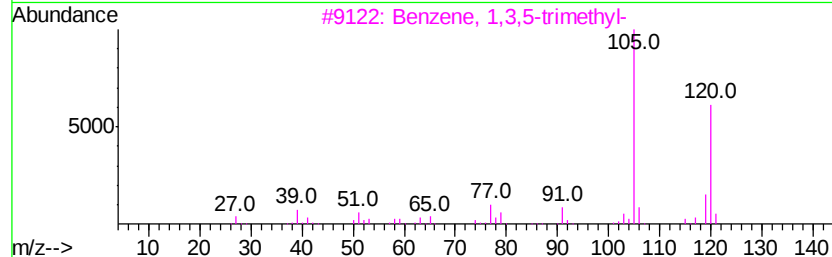
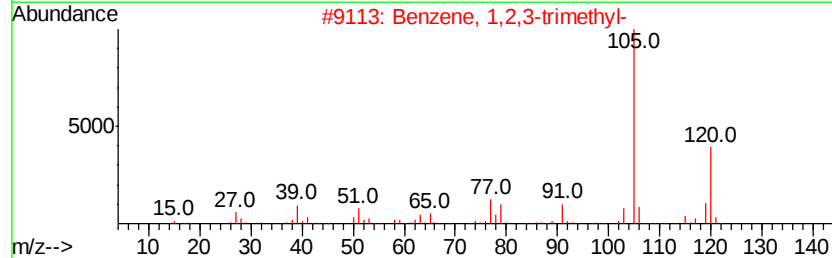
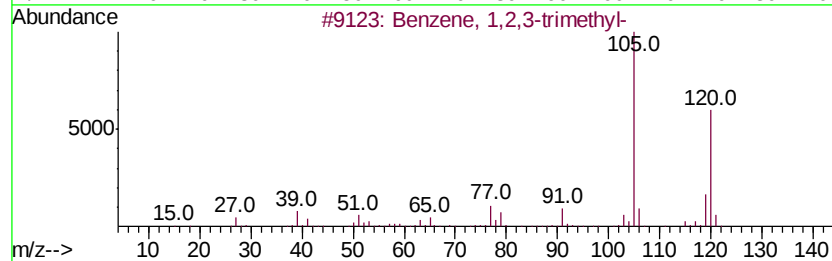
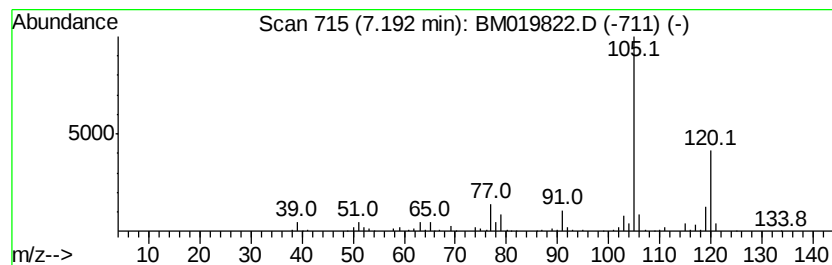
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 11 Benzene, 1,2,3-trimethyl- Concentration Rank 20

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 7.19 | 4.85 ng/ul | 305984 | 1,4-Dichlorobenzene-d4 | 7.56 |

| Hit# | of | 5 | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|------|----|---|---------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1,2,3-trimethyl- | 120 | C9H12   | 000526-73-8 | 95   |
| 2    |    |   | Benzene, 1,2,3-trimethyl- | 120 | C9H12   | 000526-73-8 | 95   |
| 3    |    |   | Benzene, 1,3,5-trimethyl- | 120 | C9H12   | 000108-67-8 | 94   |
| 4    |    |   | Benzene, 1,3,5-trimethyl- | 120 | C9H12   | 000108-67-8 | 91   |
| 5    |    |   | Benzene, 1,2,4-trimethyl- | 120 | C9H12   | 000095-63-6 | 91   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040819\  
Data File : BM019822.D  
Acq On : 09 Apr 2019 08:22  
Operator : JU/SJ  
Sample : K2188-10  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF654

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

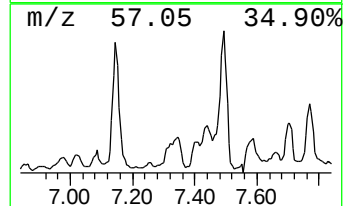
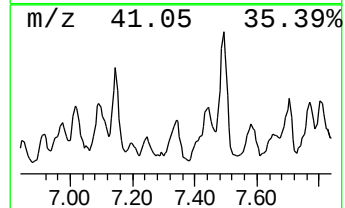
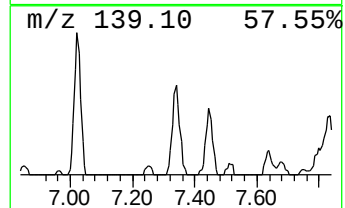
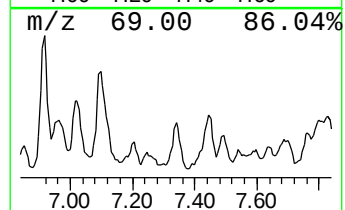
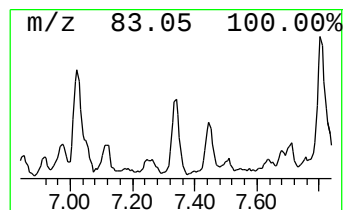
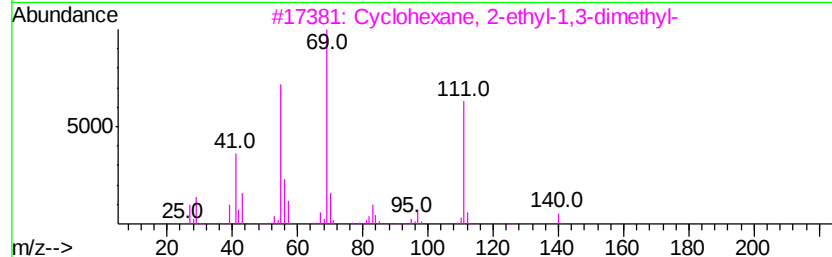
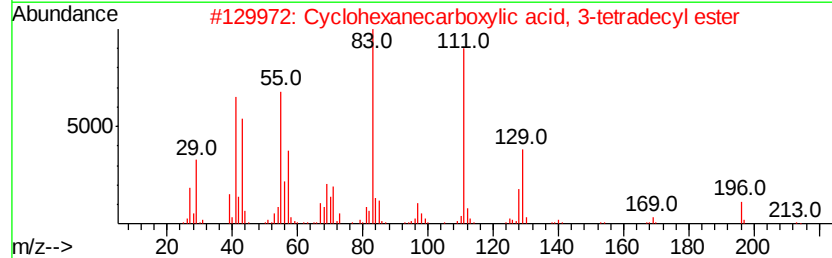
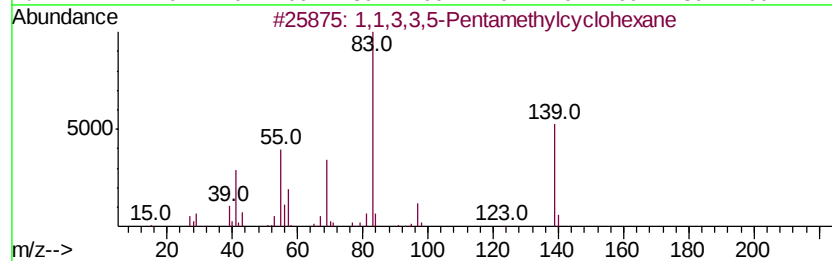
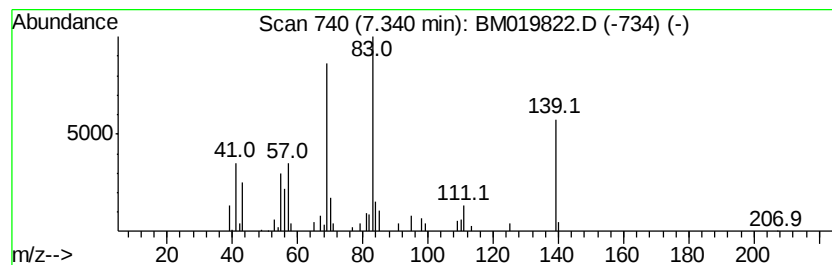
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 12 (DEL) Alkane: Cyclic7.34 Concentration Rank 40

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 7.34 | 2.73 ng/ul | 172068 | 1,4-Dichlorobenzene-d4 | 7.56 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|-------------------------------------|-----|----------|--------------|------|
| 1    | 5  | 1,1,3,3,5-Pentamethylcyclohexane    | 154 | C11H22   | 070810-19-4  | 38   |
| 2    |    | Cyclohexanecarboxylic acid, 3-te... | 324 | C21H40O2 | 1000280-59-7 | 35   |
| 3    |    | Cyclohexane, 2-ethyl-1,3-dimethyl-  | 140 | C10H20   | 007045-67-2  | 35   |
| 4    |    | Oxalic acid, di-(-)-menthyl ester   | 366 | C22H38O4 | 1000158-45-3 | 35   |
| 5    |    | Cyclohexanecarboxylic acid, 3-tr... | 310 | C20H38O2 | 1000280-59-4 | 35   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040819\  
Data File : BM019822.D  
Acq On : 09 Apr 2019 08:22  
Operator : JU/SJ  
Sample : K2188-10  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF654

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

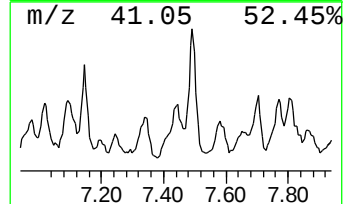
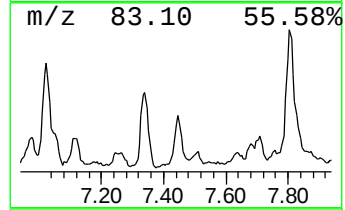
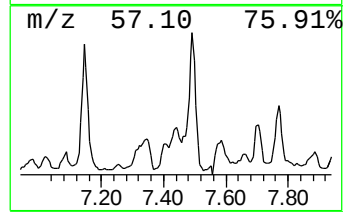
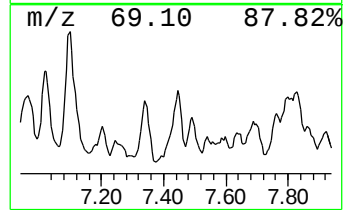
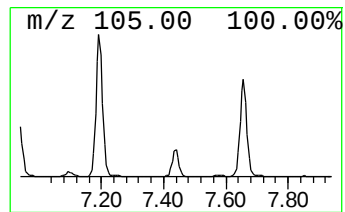
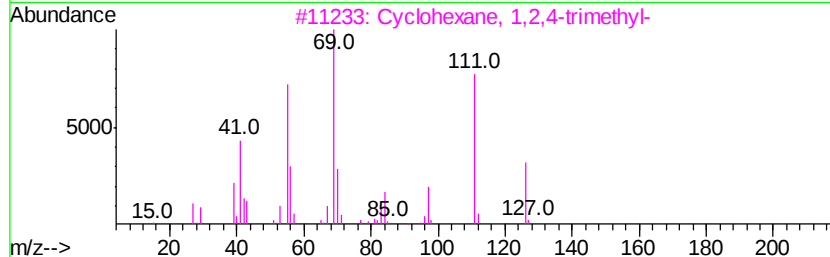
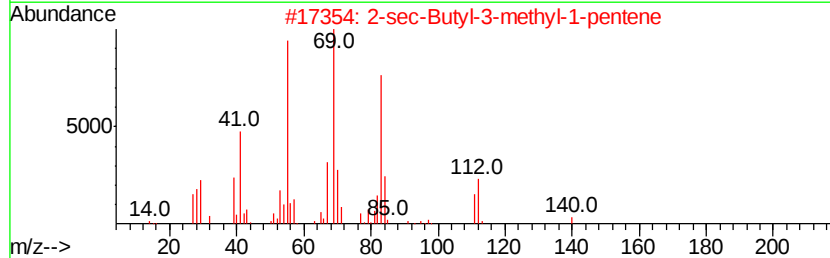
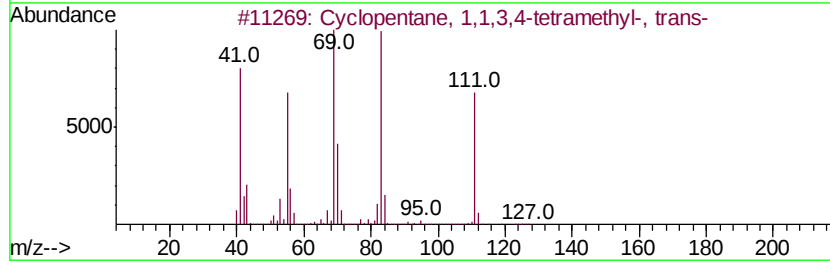
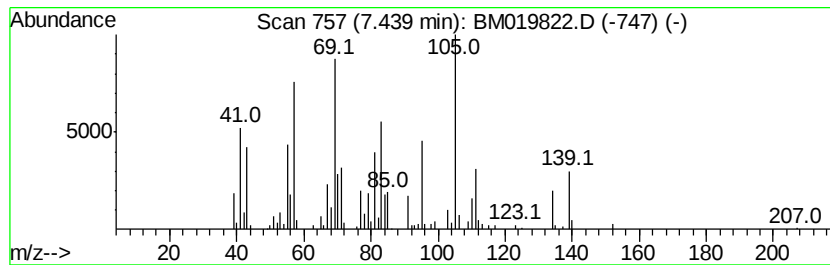
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 13 (DEL) Alkane: Cyclic7.44 Concentration Rank 12

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 7.44 | 6.01 ng/ul | 379305 | 1,4-Dichlorobenzene-d4 | 7.56 |

| Hit# | of | 5 | Tentative ID                         | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Cyclopentane, 1,1,3,4-tetramethyl... | 126 | C9H18   | 020309-77-7 | 38   |
| 2    |    |   | 2-sec-Butyl-3-methyl-1-pentene       | 140 | C10H20  | 075144-24-0 | 35   |
| 3    |    |   | Cyclohexane, 1,2,4-trimethyl-        | 126 | C9H18   | 002234-75-5 | 22   |
| 4    |    |   | 1,1,4-Trimethylcyclohexane           | 126 | C9H18   | 007094-27-1 | 22   |
| 5    |    |   | Cyclohexane, 1,2,4-trimethyl-, (...) | 126 | C9H18   | 007667-60-9 | 22   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040819\  
Data File : BM019822.D  
Acq On : 09 Apr 2019 08:22  
Operator : JU/SJ  
Sample : K2188-10  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF654

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

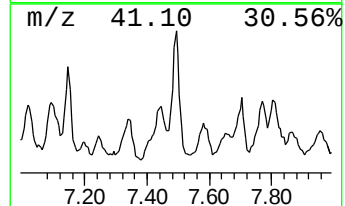
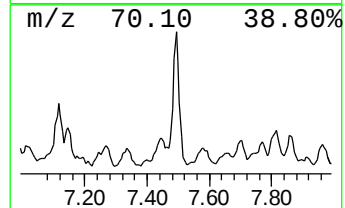
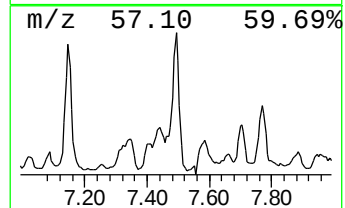
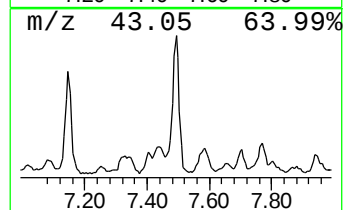
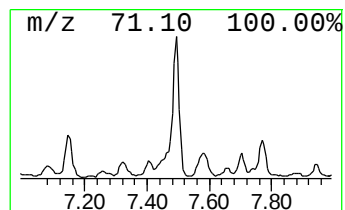
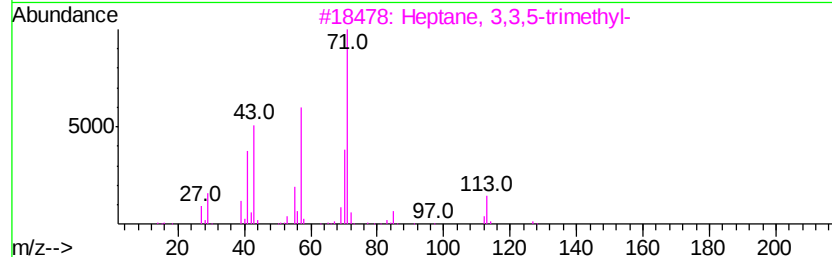
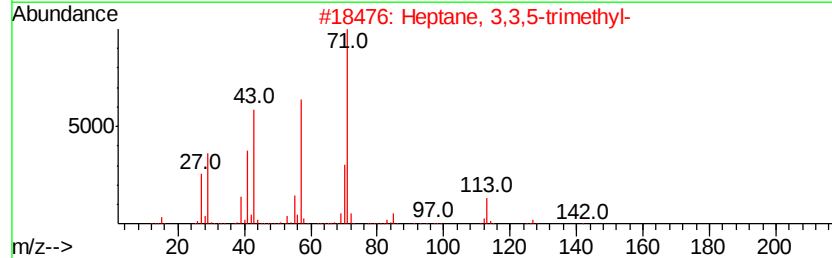
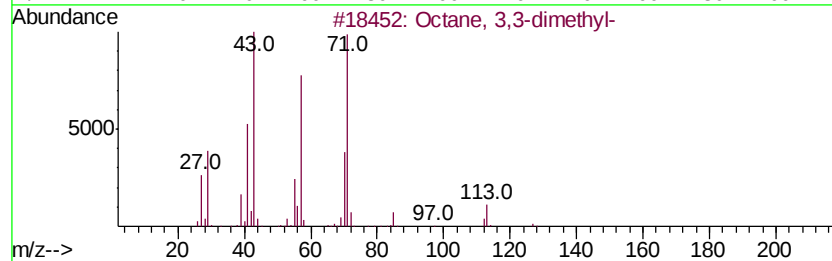
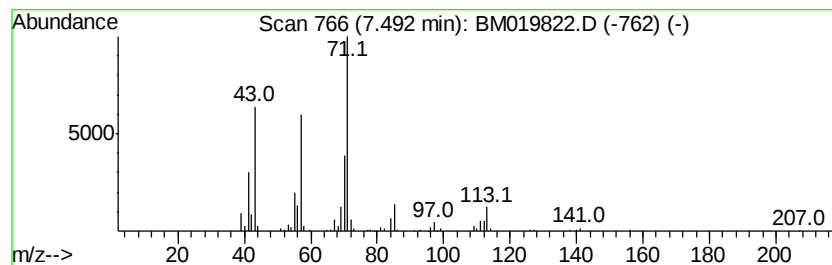
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 14 (DEL) Alkane: Straight-Chai... Concentration Rank 13

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 7.49 | 5.91 ng/ul | 372709 | 1,4-Dichlorobenzene-d4 | 7.56 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Octane, 3,3-dimethyl-      | 142 | C10H22  | 004110-44-5 | 78   |
| 2    |    |   | Heptane, 3,3,5-trimethyl-  | 142 | C10H22  | 007154-80-5 | 72   |
| 3    |    |   | Heptane, 3,3,5-trimethyl-  | 142 | C10H22  | 007154-80-5 | 72   |
| 4    |    |   | 1-Hexene, 3,4,5-trimethyl- | 126 | C9H18   | 056728-10-0 | 64   |
| 5    |    |   | Octane, 3,3-dimethyl-      | 142 | C10H22  | 004110-44-5 | 64   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040819\  
Data File : BM019822.D  
Acq On : 09 Apr 2019 08:22  
Operator : JU/SJ  
Sample : K2188-10  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF654

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

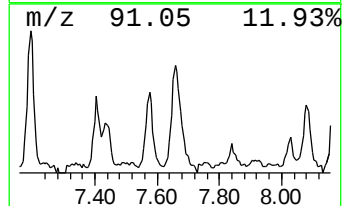
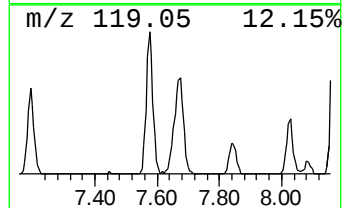
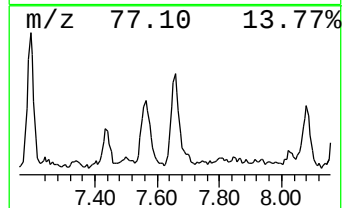
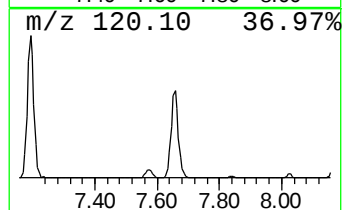
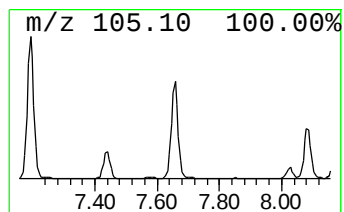
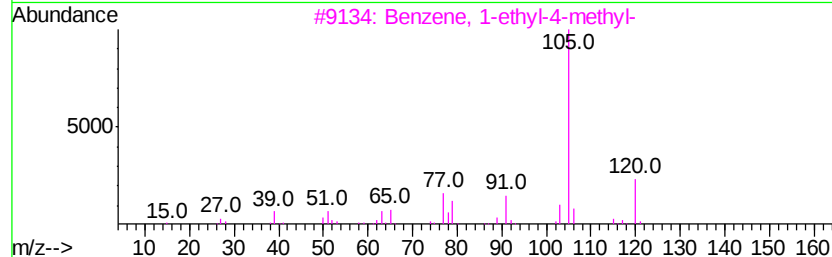
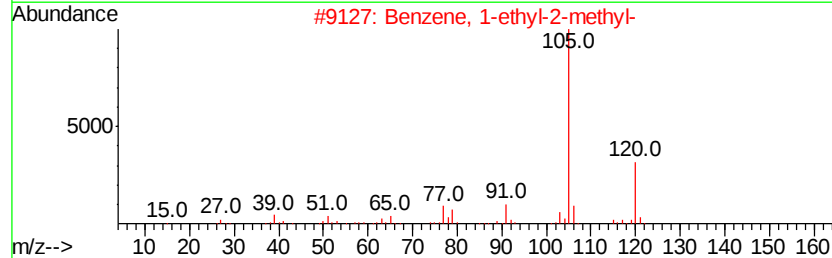
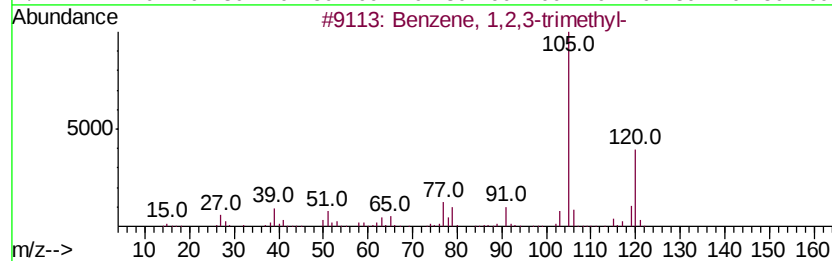
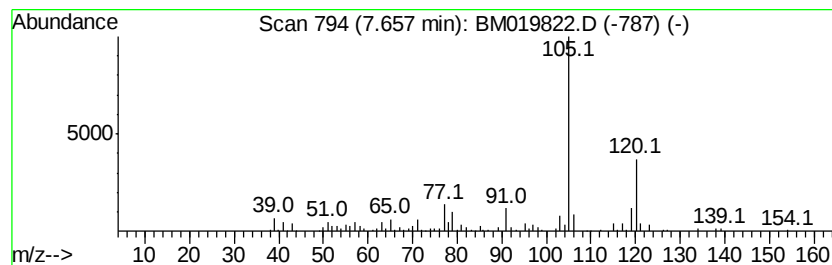
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 15 Benzene, 1,2,4-trimethyl- Concentration Rank 14

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 7.66 | 5.85 ng/ul | 369055 | 1,4-Dichlorobenzene-d4 | 7.56 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1,2,3-trimethyl-  | 120 | C9H12   | 000526-73-8 | 95   |
| 2    |    | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 93   |
| 3    |    | Benzene, 1-ethyl-4-methyl- | 120 | C9H12   | 000622-96-8 | 93   |
| 4    |    | Benzene, 1,2,4-trimethyl-  | 120 | C9H12   | 000095-63-6 | 91   |
| 5    |    | Benzene, 1,2,3-trimethyl-  | 120 | C9H12   | 000526-73-8 | 91   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040819\  
Data File : BM019822.D  
Acq On : 09 Apr 2019 08:22  
Operator : JU/SJ  
Sample : K2188-10  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF654

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

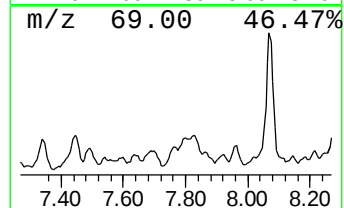
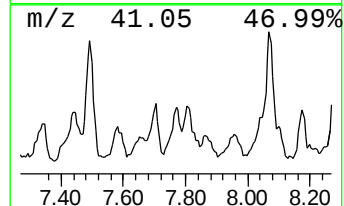
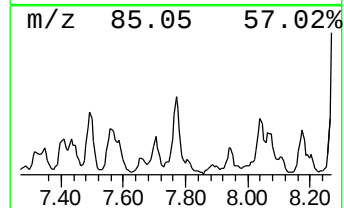
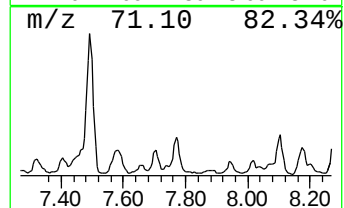
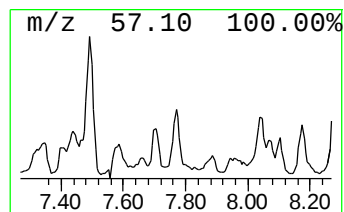
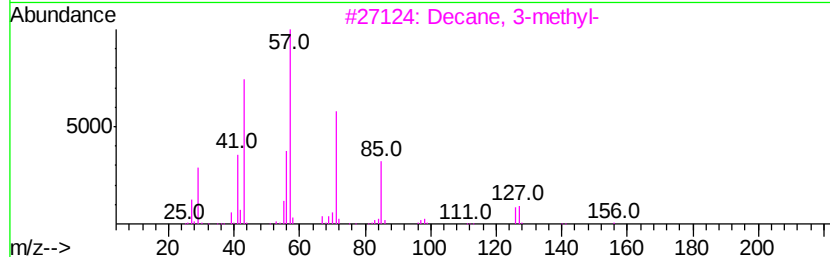
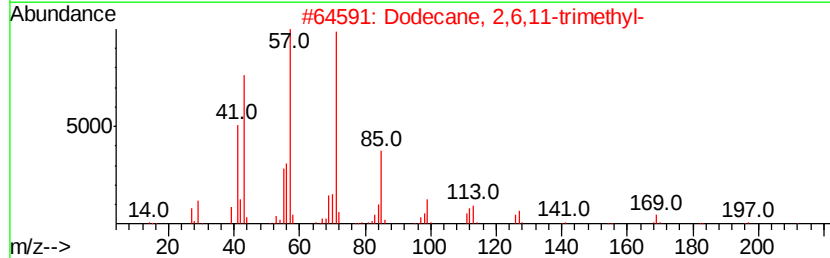
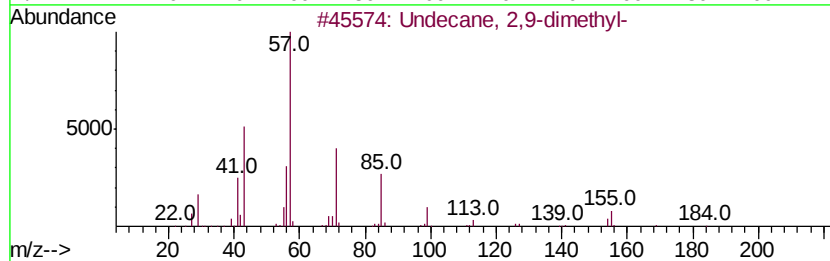
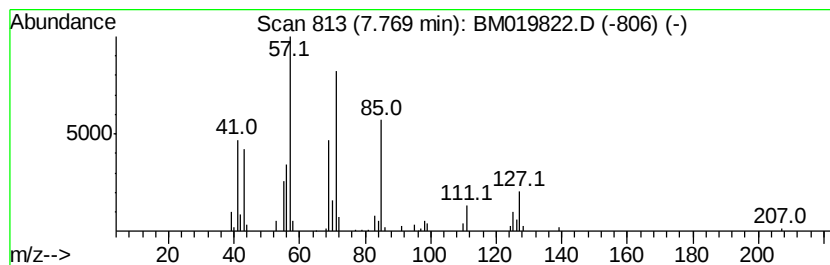
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 16 (DEL) Alkane: Straight-Chai... Concentration Rank 32

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 7.77 | 3.00 ng/ul | 189494 | 1,4-Dichlorobenzene-d4 | 7.56 |

| Hit# | of | 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------------|-----|---------|-------------|------|
| 1    |    |   | Undecane, 2,9-dimethyl-     | 184 | C13H28  | 017301-26-7 | 59   |
| 2    |    |   | Dodecane, 2,6,11-trimethyl- | 212 | C15H32  | 031295-56-4 | 53   |
| 3    |    |   | Decane, 3-methyl-           | 156 | C11H24  | 013151-34-3 | 50   |
| 4    |    |   | Nonane, 2-methyl-5-propyl-  | 184 | C13H28  | 031081-17-1 | 50   |
| 5    |    |   | Nonane, 5-propyl-           | 170 | C12H26  | 000998-35-6 | 43   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040819\  
Data File : BM019822.D  
Acq On : 09 Apr 2019 08:22  
Operator : JU/SJ  
Sample : K2188-10  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF654

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

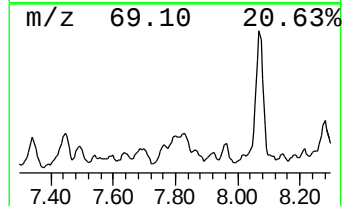
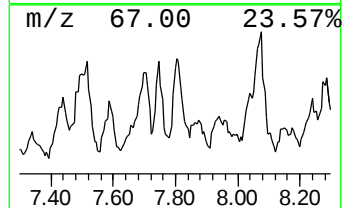
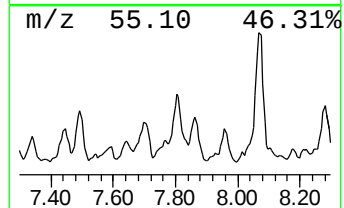
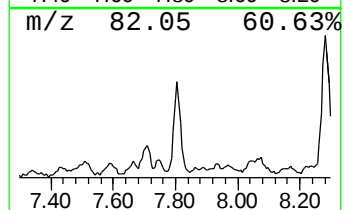
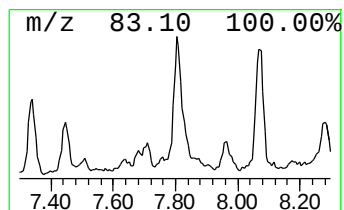
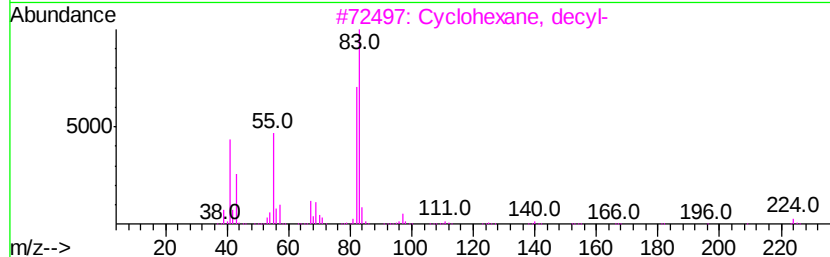
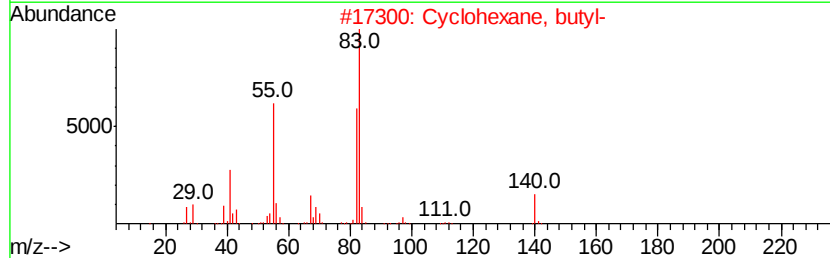
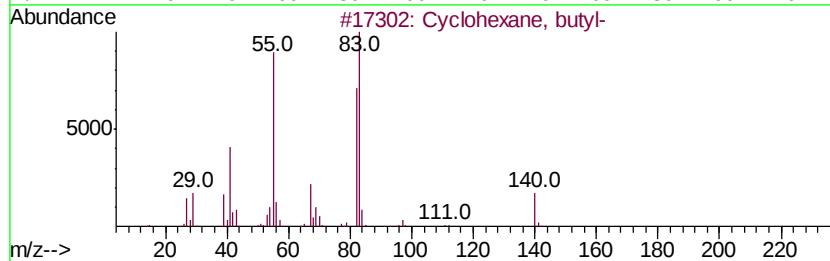
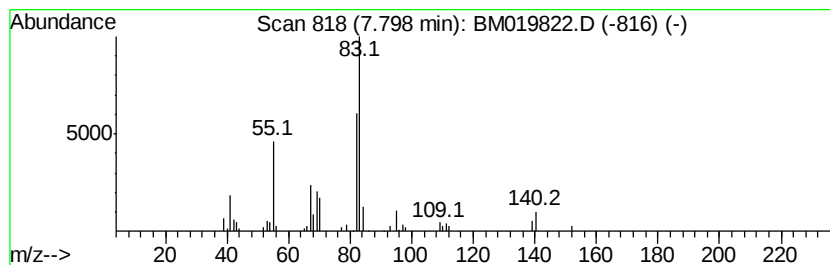
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 17 (DEL) Alkane: Cyclic7.80 Concentration Rank 30

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 7.80 | 3.20 ng/ul | 201806 | 1,4-Dichlorobenzene-d4 | 7.56 |

| Hit# | of | 5 | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------|-----|---------|-------------|------|
| 1    |    |   | Cyclohexane, butyl-           | 140 | C10H20  | 001678-93-9 | 86   |
| 2    |    |   | Cyclohexane, butyl-           | 140 | C10H20  | 001678-93-9 | 72   |
| 3    |    |   | Cyclohexane, decyl-           | 224 | C16H32  | 001795-16-0 | 64   |
| 4    |    |   | Cyclohexane, (1-methylethyl)- | 126 | C9H18   | 000696-29-7 | 64   |
| 5    |    |   | Cyclohexane, 2-propenyl-      | 124 | C9H16   | 002114-42-3 | 64   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040819\  
Data File : BM019822.D  
Acq On : 09 Apr 2019 08:22  
Operator : JU/SJ  
Sample : K2188-10  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF654

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

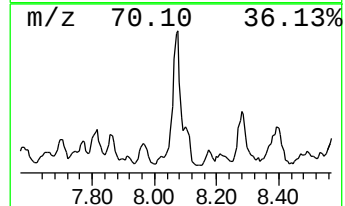
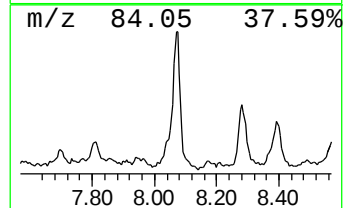
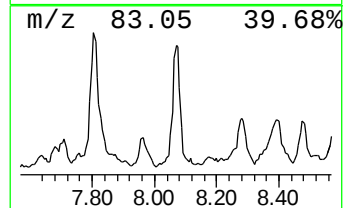
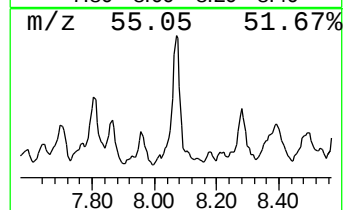
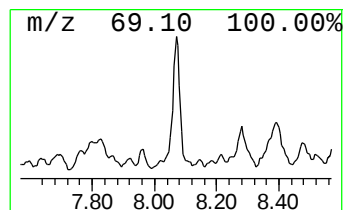
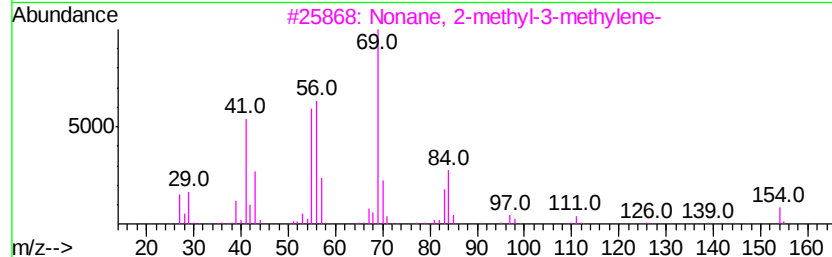
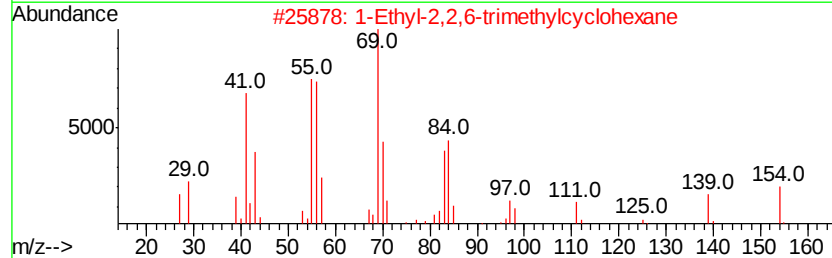
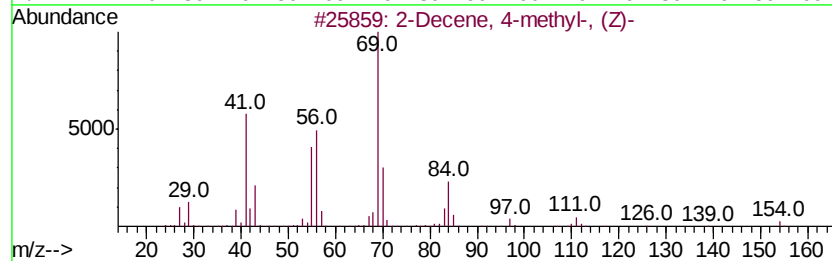
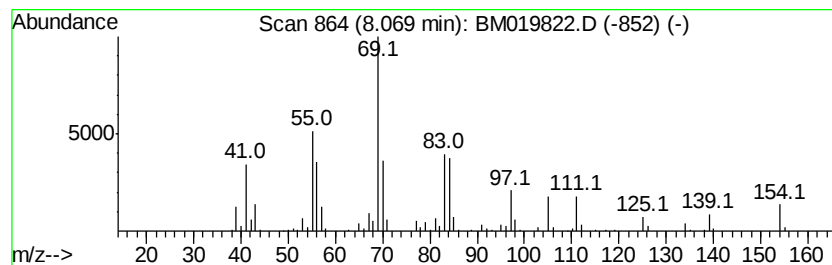
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 18 (DEL) Alkane: Cyclic8.07 Concentration Rank 1

| R.T. | EstConc     | Area   | Relative to ISTD       | R.T. |
|------|-------------|--------|------------------------|------|
| 8.07 | 14.15 ng/ul | 892147 | 1,4-Dichlorobenzene-d4 | 7.56 |

| Hit# | of | Tentative ID                       | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------------------|-----|---------|-------------|------|
| 1    | 5  | 2-Decene, 4-methyl-, (Z)-          | 154 | C11H22  | 074630-30-1 | 64   |
| 2    |    | 1-Ethyl-2,2,6-trimethylcyclohexane | 154 | C11H22  | 071186-27-1 | 58   |
| 3    |    | Nonane, 2-methyl-3-methylene-      | 154 | C11H22  | 055499-08-6 | 53   |
| 4    |    | 1-Hexene, 3,3,5-trimethyl-         | 126 | C9H18   | 013427-43-5 | 47   |
| 5    |    | Cyclohexane, 1,2,3-trimethyl-      | 126 | C9H18   | 001678-97-3 | 47   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040819\  
Data File : BM019822.D  
Acq On : 09 Apr 2019 08:22  
Operator : JU/SJ  
Sample : K2188-10  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF654

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

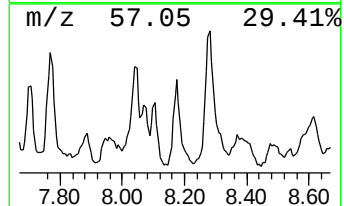
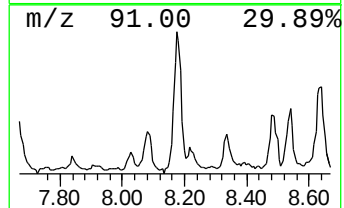
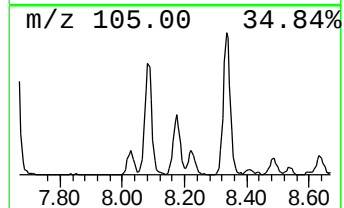
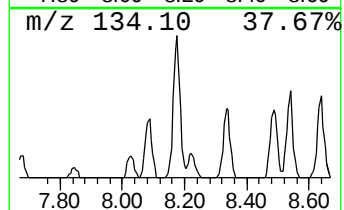
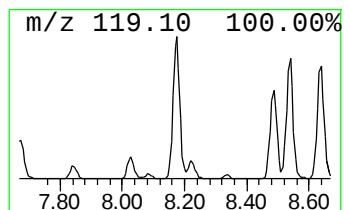
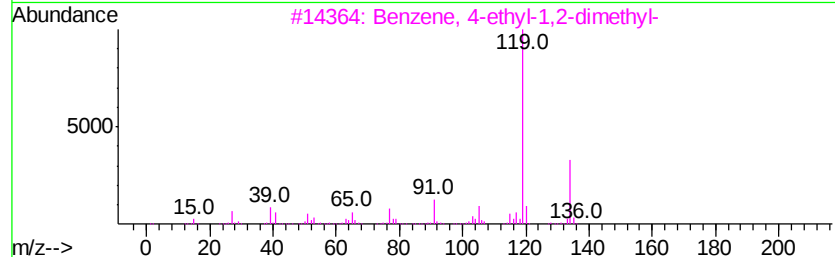
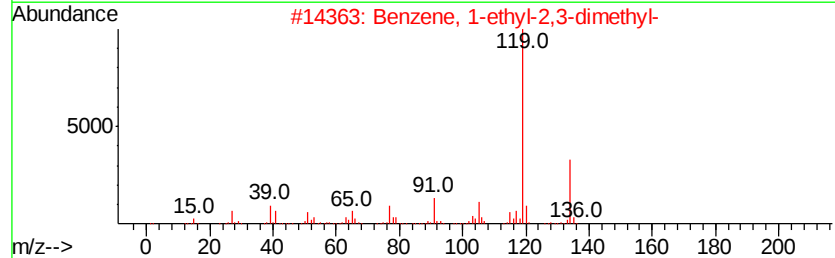
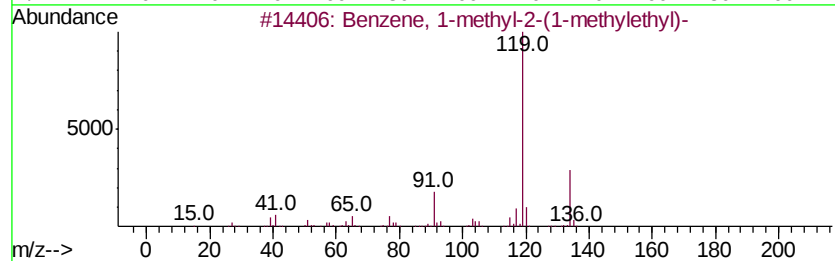
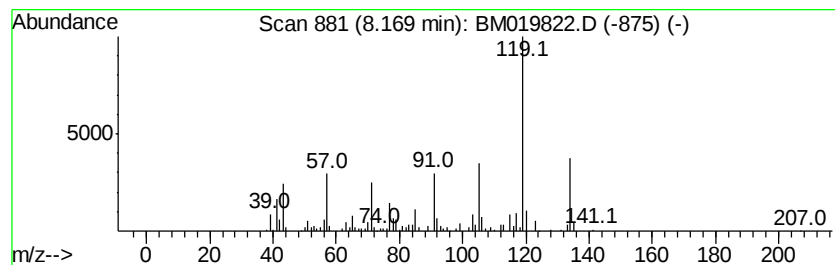
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 19 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 15

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 8.17 | 5.78 ng/ul | 364512 | 1,4-Dichlorobenzene-d4 | 7.56 |

| Hit# | of | Tentative ID                         | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1-methyl-2-(1-methylethyl)- | 134 | C10H14  | 000527-84-4 | 86   |
| 2    |    | Benzene, 1-ethyl-2,3-dimethyl-       | 134 | C10H14  | 000933-98-2 | 76   |
| 3    |    | Benzene, 4-ethyl-1,2-dimethyl-       | 134 | C10H14  | 000934-80-5 | 76   |
| 4    |    | Benzene, 1-ethyl-3,5-dimethyl-       | 134 | C10H14  | 000934-74-7 | 76   |
| 5    |    | Benzene, 1-ethyl-3,5-dimethyl-       | 134 | C10H14  | 000934-74-7 | 76   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040819\  
Data File : BM019822.D  
Acq On : 09 Apr 2019 08:22  
Operator : JU/SJ  
Sample : K2188-10  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampled :  
BF654

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

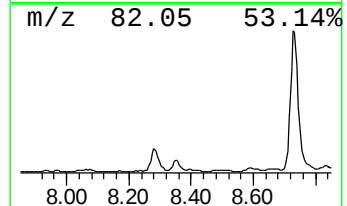
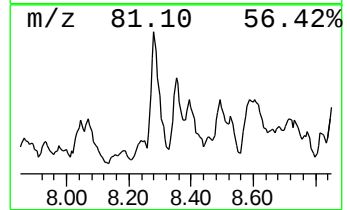
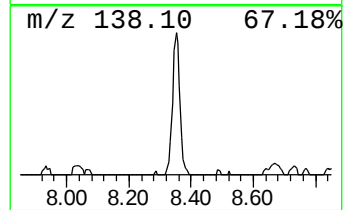
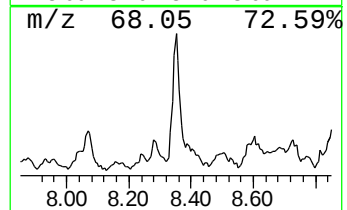
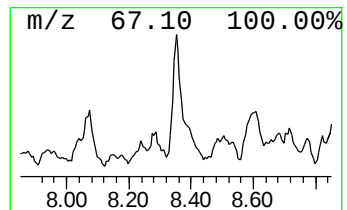
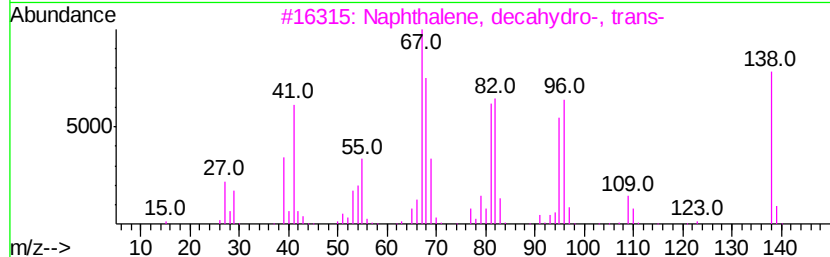
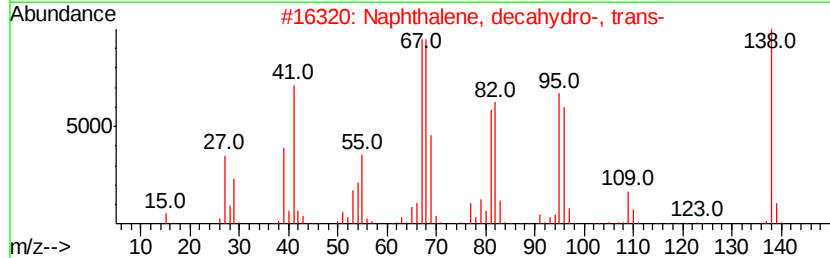
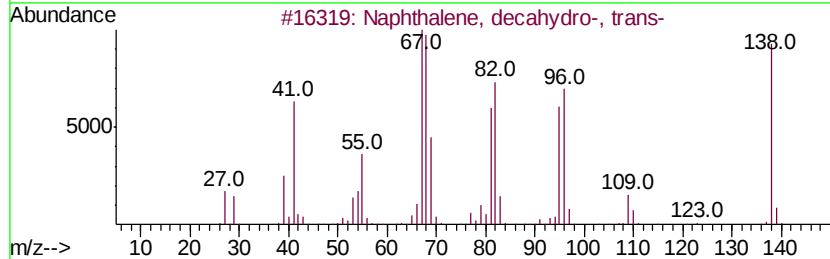
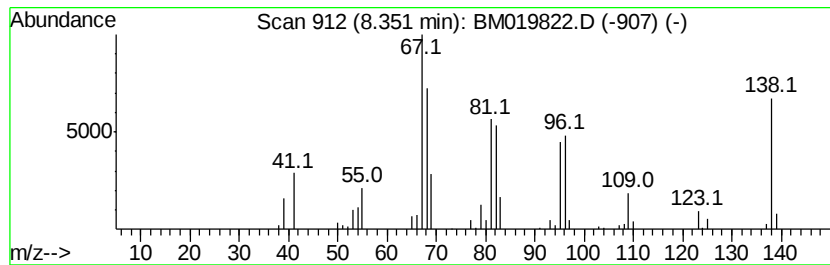
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 20 Naphthalene, decahydro-, tr... Concentration Rank 44

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 8.35 | 2.65 ng/ul | 166962 | 1,4-Dichlorobenzene-d4 | 7.56 |

| Hit# | of | 5 | Tentative ID                    | MW  | MolForm | CAS#        | Qual |
|------|----|---|---------------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, decahydro-, trans- | 138 | C10H18  | 000493-02-7 | 95   |
| 2    |    |   | Naphthalene, decahydro-, trans- | 138 | C10H18  | 000493-02-7 | 94   |
| 3    |    |   | Naphthalene, decahydro-, trans- | 138 | C10H18  | 000493-02-7 | 94   |
| 4    |    |   | Naphthalene, decahydro-         | 138 | C10H18  | 000091-17-8 | 91   |
| 5    |    |   | Naphthalene, decahydro-         | 138 | C10H18  | 000091-17-8 | 70   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040819\  
Data File : BM019822.D  
Acq On : 09 Apr 2019 08:22  
Operator : JU/SJ  
Sample : K2188-10  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF654

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

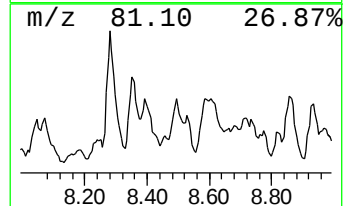
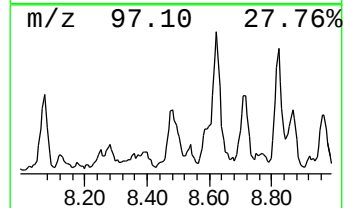
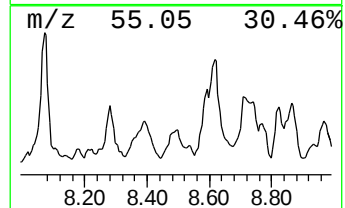
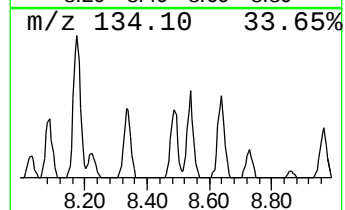
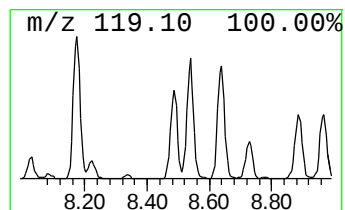
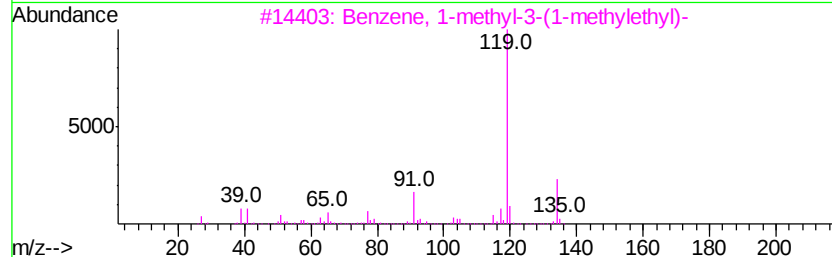
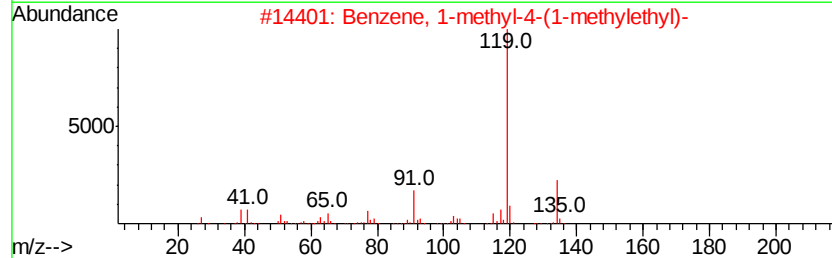
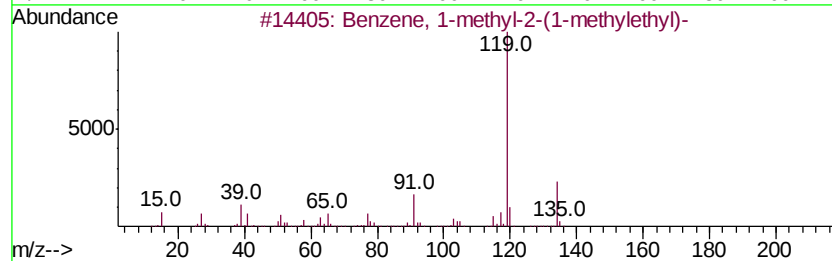
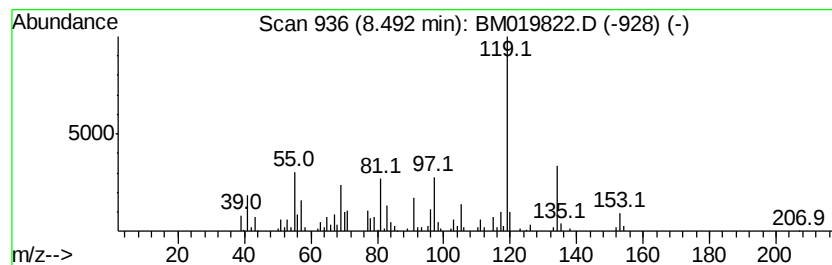
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 21 Benzene, 1-methyl-4-(1-meth... Concentration Rank 21

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 8.49 | 4.54 ng/ul | 286032 | 1,4-Dichlorobenzene-d4 | 7.56 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1-methyl-2-(1-methyleth... | 134 | C10H14  | 000527-84-4 | 93   |
| 2    |    |   | Benzene, 1-methyl-4-(1-methyleth... | 134 | C10H14  | 000099-87-6 | 93   |
| 3    |    |   | Benzene, 1-methyl-3-(1-methyleth... | 134 | C10H14  | 000535-77-3 | 93   |
| 4    |    |   | Benzene, 1-methyl-2-(1-methyleth... | 134 | C10H14  | 000527-84-4 | 93   |
| 5    |    |   | Benzene, 1-methyl-2-(1-methyleth... | 134 | C10H14  | 000527-84-4 | 92   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040819\  
Data File : BM019822.D  
Acq On : 09 Apr 2019 08:22  
Operator : JU/SJ  
Sample : K2188-10  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF654

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

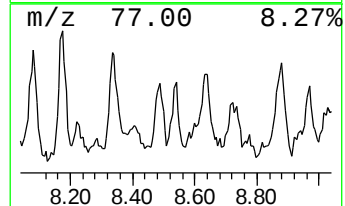
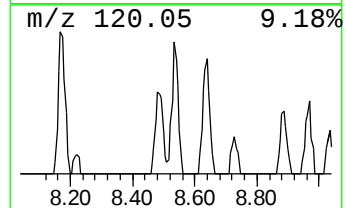
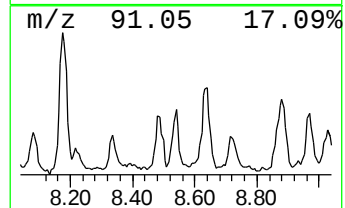
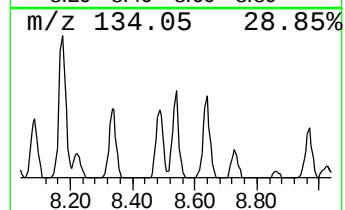
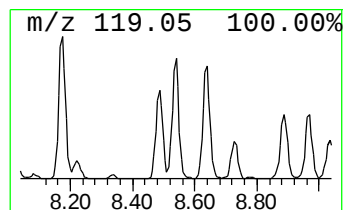
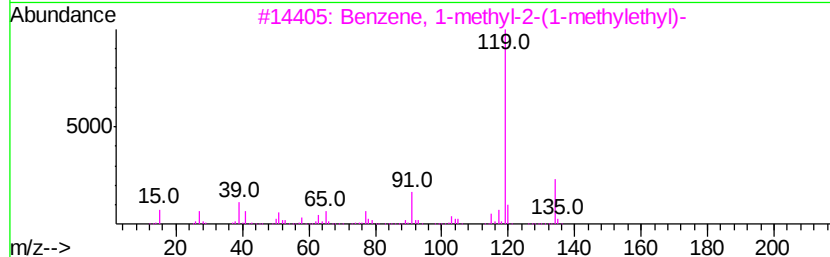
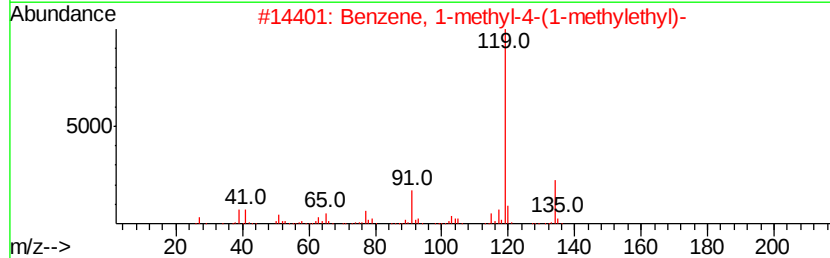
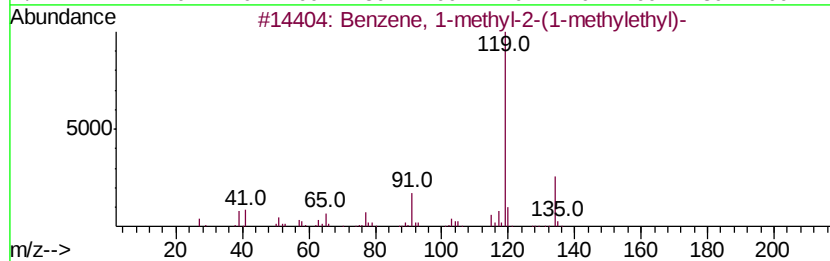
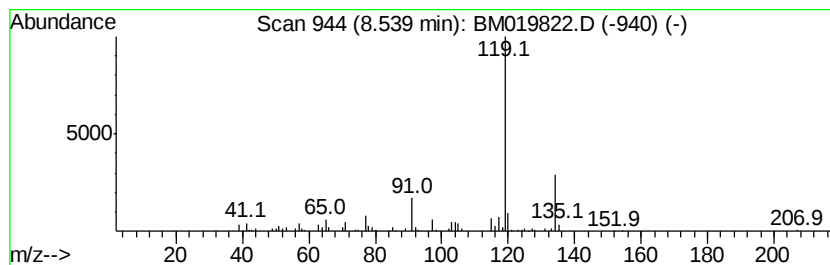
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 22 Benzene, 1-methyl-2-(1-meth... Concentration Rank 48

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 8.54 | 2.26 ng/ul | 142613 | 1,4-Dichlorobenzene-d4 | 7.56 |

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



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040819\  
Data File : BM019822.D  
Acq On : 09 Apr 2019 08:22  
Operator : JU/SJ  
Sample : K2188-10  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampled :  
BF654

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

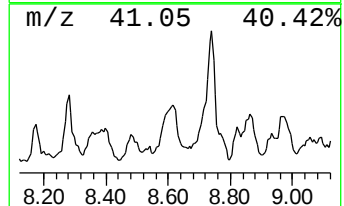
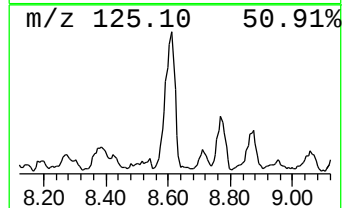
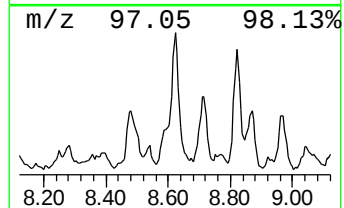
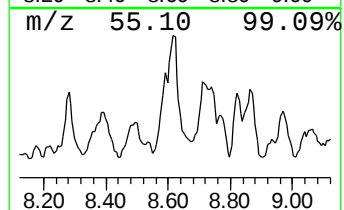
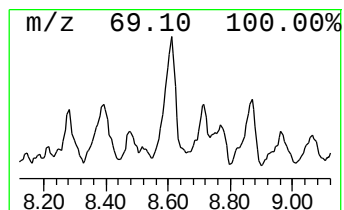
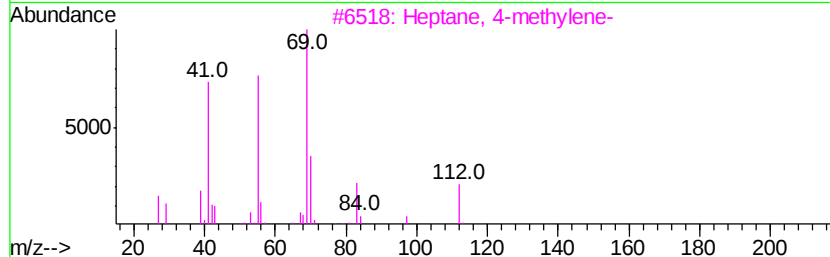
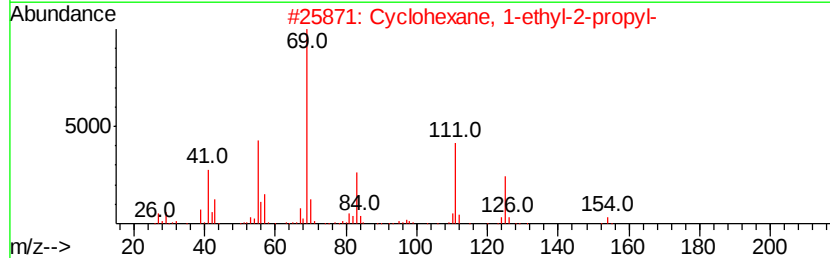
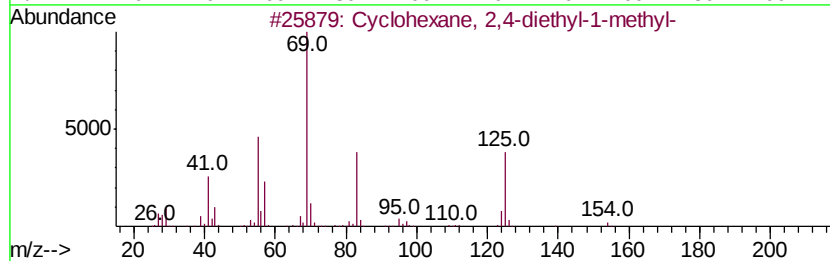
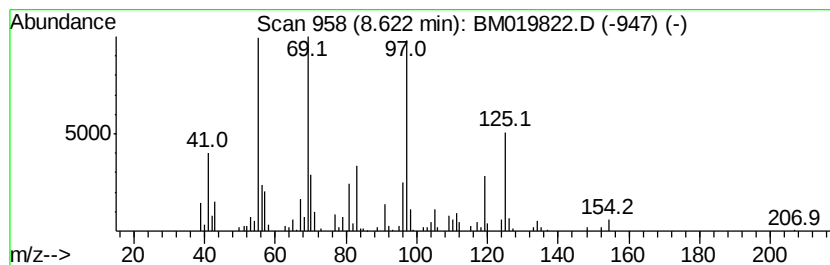
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 23 (DEL) Alkane: Cyclic8.62 Concentration Rank 2

| R.T. | EstConc     | Area   | Relative to ISTD       | R.T. |
|------|-------------|--------|------------------------|------|
| 8.62 | 11.38 ng/ul | 717886 | 1,4-Dichlorobenzene-d4 | 7.56 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Cyclohexane, 2,4-diethyl-1-methyl-  | 154 | C11H22  | 061142-70-9 | 58   |
| 2    |    | Cyclohexane, 1-ethyl-2-propyl-      | 154 | C11H22  | 062238-33-9 | 47   |
| 3    |    | Heptane, 4-methylene-               | 112 | C8H16   | 015918-08-8 | 35   |
| 4    |    | Cyclopentanecarboxylic acid, all... | 154 | C9H14O2 | 178905-33-4 | 35   |
| 5    |    | Cycloheptane, methyl-               | 112 | C8H16   | 004126-78-7 | 27   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040819\  
Data File : BM019822.D  
Acq On : 09 Apr 2019 08:22  
Operator : JU/SJ  
Sample : K2188-10  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampled :  
BF654

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

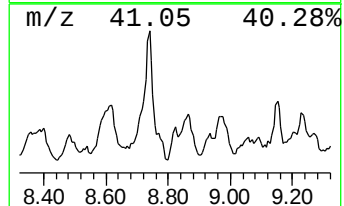
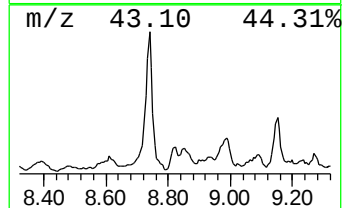
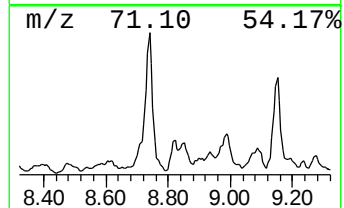
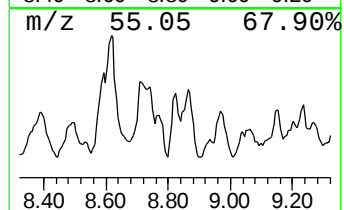
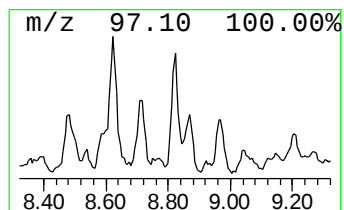
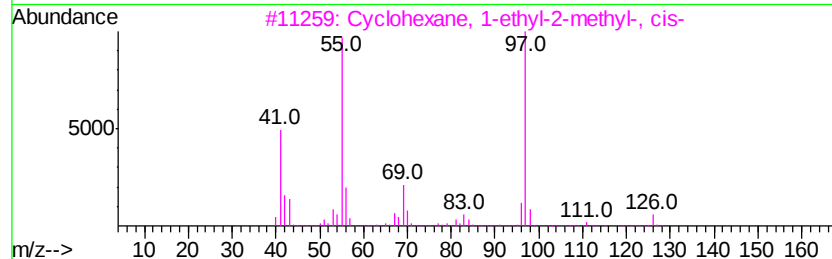
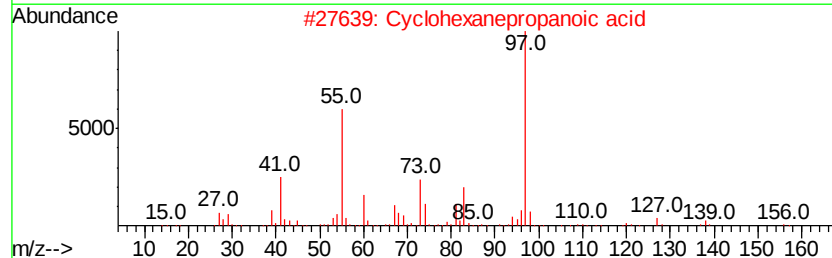
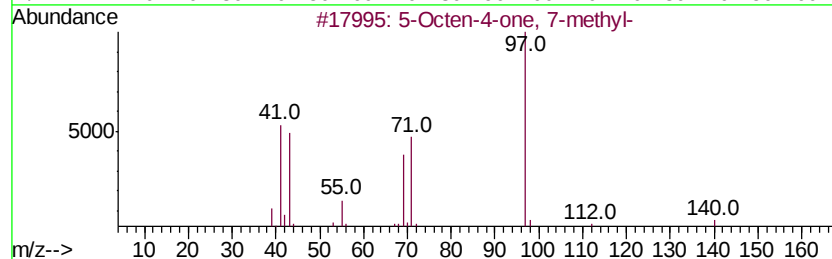
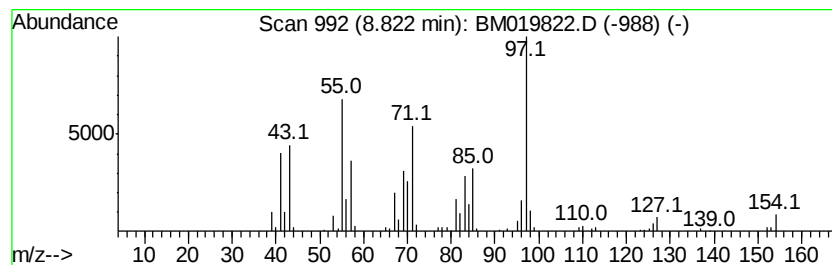
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 24 unknown-02 Concentration Rank 43

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 8.82 | 2.67 ng/ul | 168649 | 1,4-Dichlorobenzene-d4 | 7.56 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|---------|--------------|------|
| 1    |    |   | 5-Octen-4-one, 7-methyl-            | 140 | C9H16O  | 032064-78-1  | 43   |
| 2    |    |   | Cyclohexanepropanoic acid           | 156 | C9H16O2 | 000701-97-3  | 43   |
| 3    |    |   | Cyclohexane, 1-ethyl-2-methyl-, ... | 126 | C9H18   | 004923-77-7  | 43   |
| 4    |    |   | 3-Cyclopent-1-enyl-3-hydroxy-2-m... | 170 | C9H14O3 | 1000191-19-7 | 38   |
| 5    |    |   | Oxazole, 4,5-dimethyl-              | 97  | C5H7NO  | 020662-83-3  | 38   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040819\  
Data File : BM019822.D  
Acq On : 09 Apr 2019 08:22  
Operator : JU/SJ  
Sample : K2188-10  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF654

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

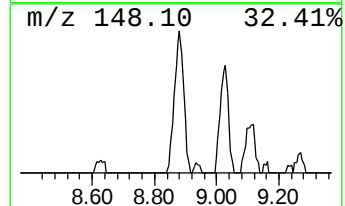
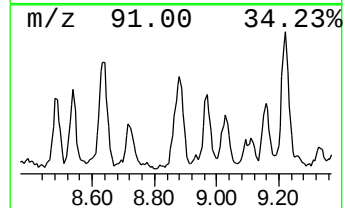
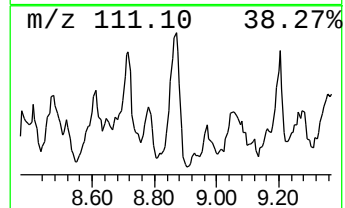
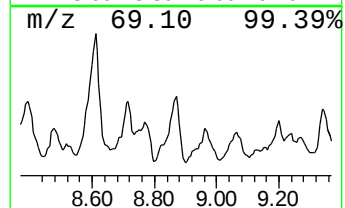
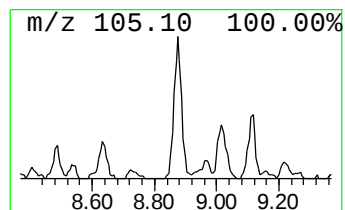
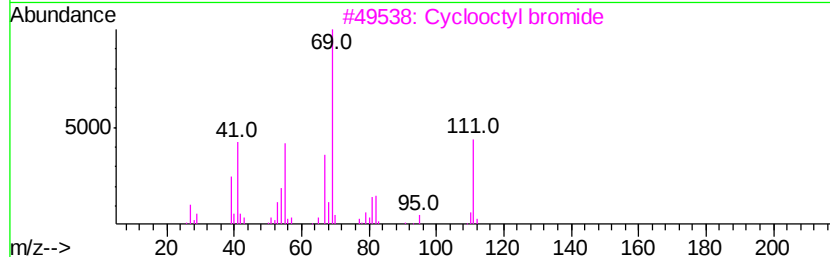
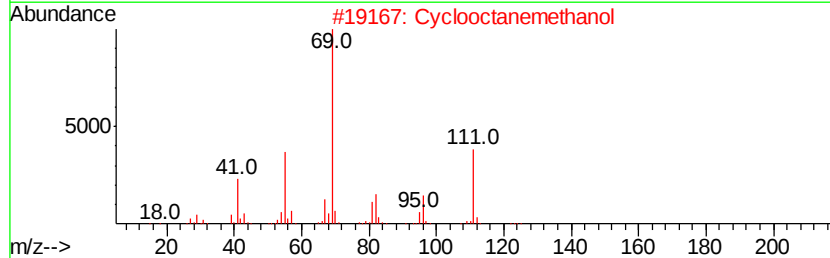
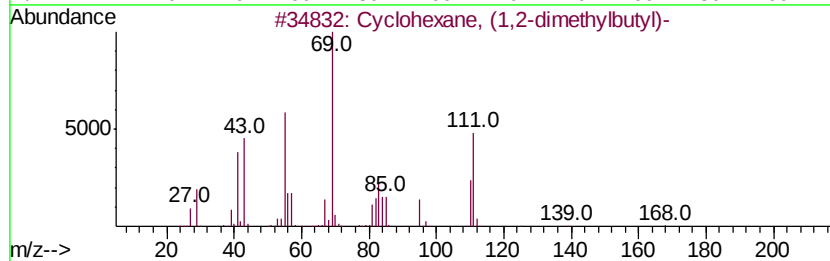
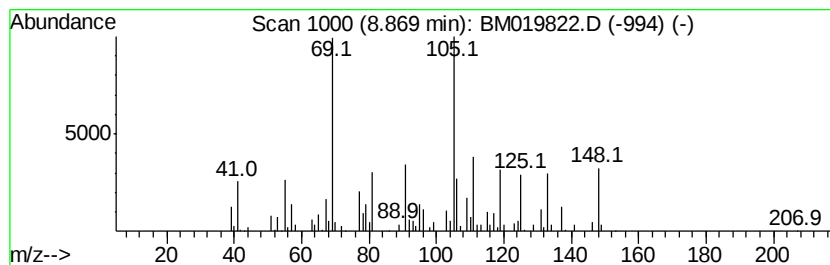
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 25 (DEL) Alkane: Cyclic8.87 Concentration Rank 6

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 8.87 | 6.71 ng/ul | 423394 | 1,4-Dichlorobenzene-d4 | 7.56 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Cyclohexane, (1,2-dimethylbutyl)-   | 168 | C12H24   | 061142-37-8  | 35   |
| 2    |    |   | Cyclooctanemethanol                 | 142 | C9H18O   | 003637-63-6  | 35   |
| 3    |    |   | Cyclooctyl bromide                  | 190 | C8H15Br  | 001556-09-8  | 25   |
| 4    |    |   | Methoxyacetic acid, 2-ethylcyclo... | 200 | C11H20O3 | 1000282-69-7 | 25   |
| 5    |    |   | 1-Butyne, 3-methyl-3-propoxy-       | 126 | C8H14O   | 053907-64-5  | 17   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040819\  
Data File : BM019822.D  
Acq On : 09 Apr 2019 08:22  
Operator : JU/SJ  
Sample : K2188-10  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF654

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

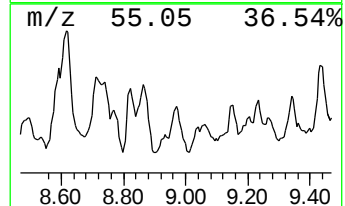
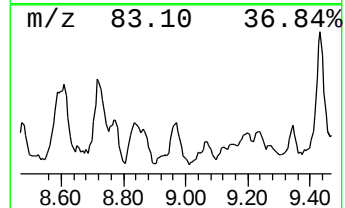
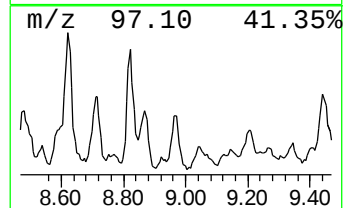
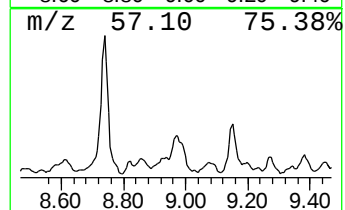
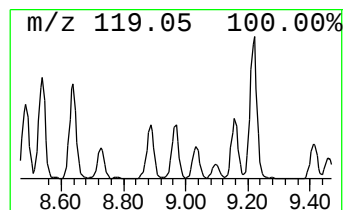
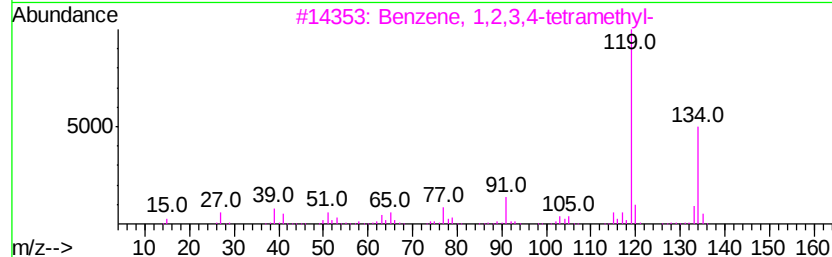
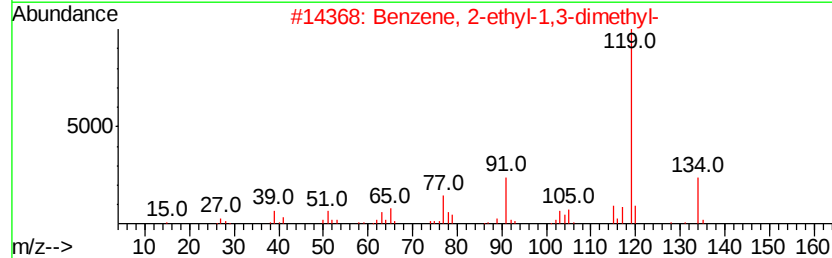
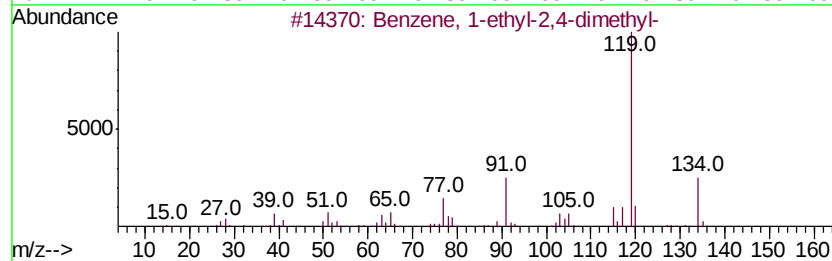
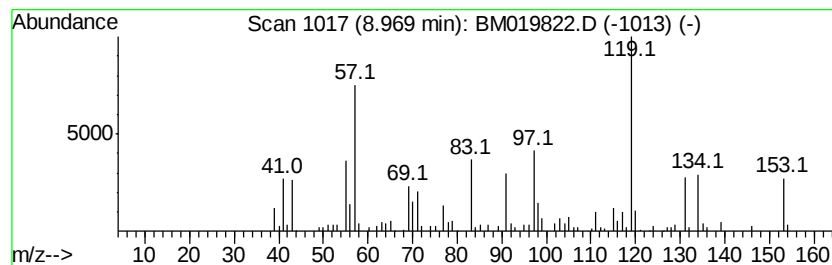
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 26 Benzene, 1-ethyl-2,4-dimethyl- Concentration Rank 23

| R.T. | EstConc    | Area   | Relative to ISTD | R.T.  |
|------|------------|--------|------------------|-------|
| 8.97 | 4.36 ng/ul | 285998 | Naphthalene-d8   | 10.34 |

| Hit# | of | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1-ethyl-2,4-dimethyl- | 134 | C10H14  | 000874-41-9 | 90   |
| 2    |    | Benzene, 2-ethyl-1,3-dimethyl- | 134 | C10H14  | 002870-04-4 | 89   |
| 3    |    | Benzene, 1,2,3,4-tetramethyl-  | 134 | C10H14  | 000488-23-3 | 86   |
| 4    |    | Benzene, 1-ethyl-2,3-dimethyl- | 134 | C10H14  | 000933-98-2 | 64   |
| 5    |    | Benzene, 4-ethyl-1,2-dimethyl- | 134 | C10H14  | 000934-80-5 | 64   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040819\  
Data File : BM019822.D  
Acq On : 09 Apr 2019 08:22  
Operator : JU/SJ  
Sample : K2188-10  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF654

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

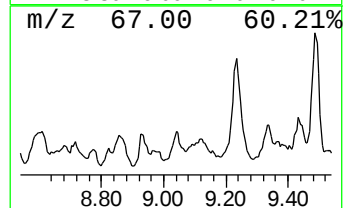
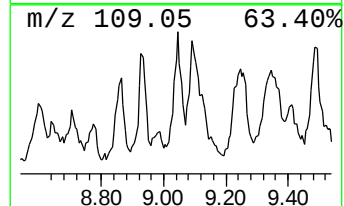
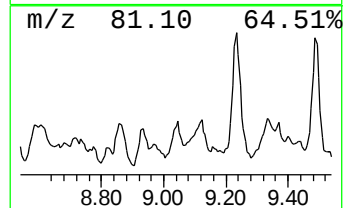
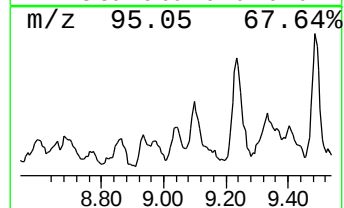
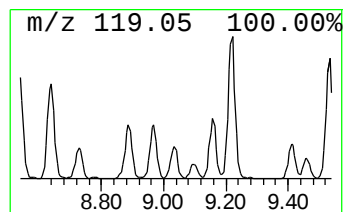
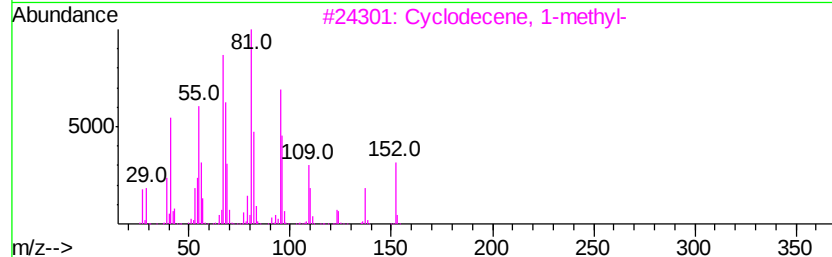
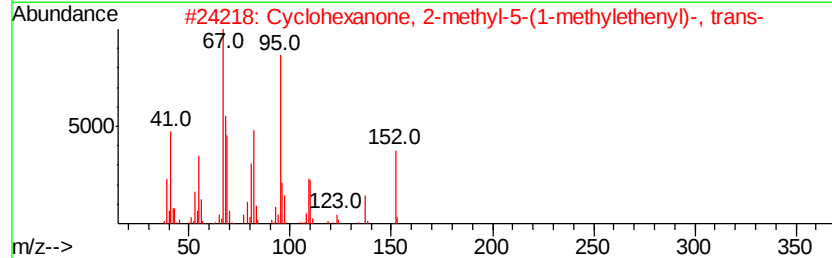
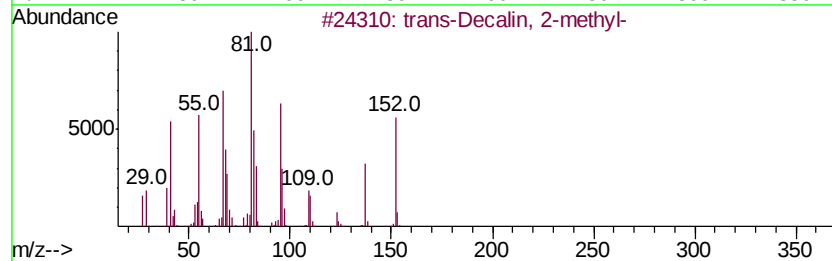
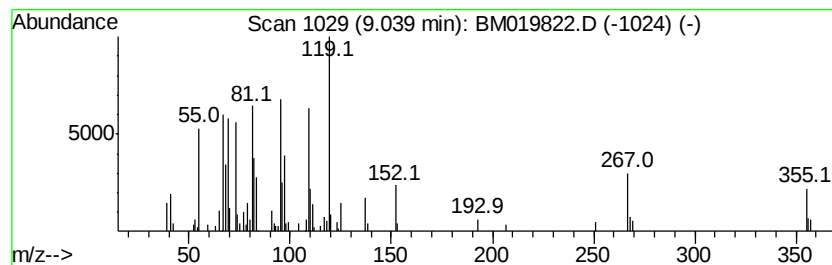
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 27 unknown-03 Concentration Rank 42

| R.T. | EstConc    | Area   | Relative to ISTD | R.T.  |
|------|------------|--------|------------------|-------|
| 9.04 | 2.68 ng/ul | 175368 | Naphthalene-d8   | 10.34 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|-------------------------------------|-----|---------|--------------|------|
| 1    | 5  | trans-Decalin, 2-methyl-            | 152 | C11H20  | 1000152-47-3 | 47   |
| 2    |    | Cyclohexanone, 2-methyl-5-(1-met... | 152 | C10H16O | 005948-04-9  | 42   |
| 3    |    | Cyclodecene, 1-methyl-              | 152 | C11H20  | 066633-38-3  | 42   |
| 4    |    | Cyclohexanone, 2-methyl-5-(1-met... | 152 | C10H16O | 007764-50-3  | 41   |
| 5    |    | Cyclohexanone, 2-methyl-5-(1-met... | 152 | C10H16O | 005948-04-9  | 41   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040819\  
Data File : BM019822.D  
Acq On : 09 Apr 2019 08:22  
Operator : JU/SJ  
Sample : K2188-10  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF654

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

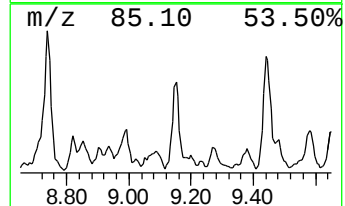
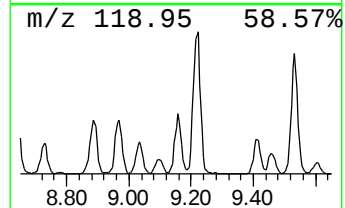
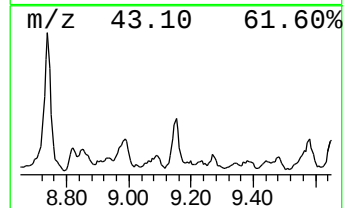
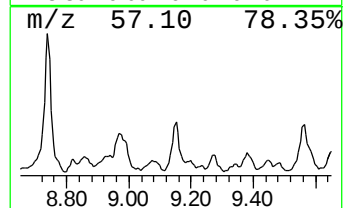
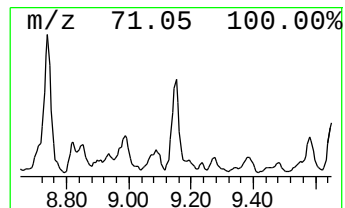
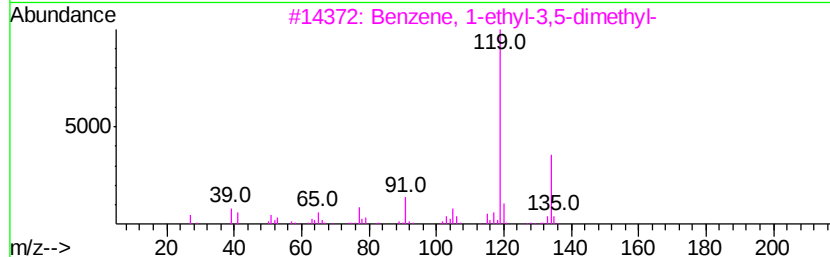
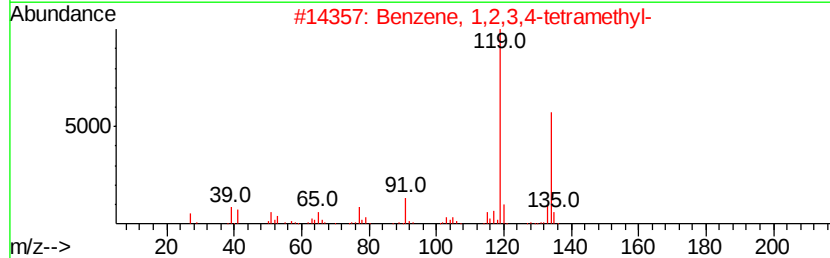
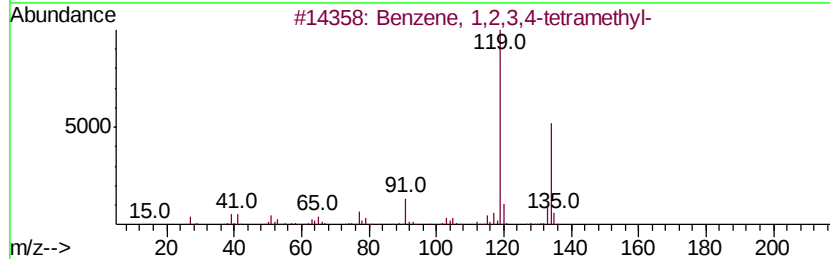
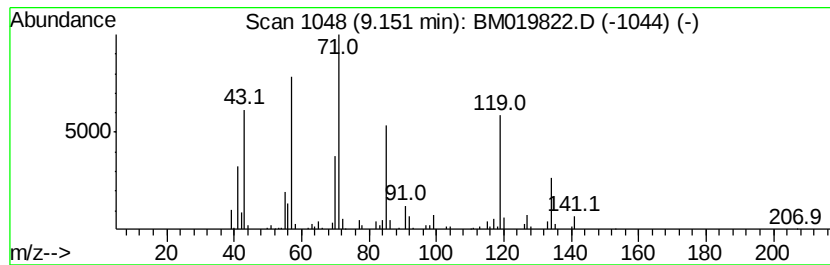
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 28 unknown-04 Concentration Rank 28

| R.T. | EstConc    | Area   | Relative to ISTD | R.T.  |
|------|------------|--------|------------------|-------|
| 9.15 | 3.33 ng/ul | 218025 | Naphthalene-d8   | 10.34 |

| Hit# | of | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1,2,3,4-tetramethyl-  | 134 | C10H14  | 000488-23-3 | 35   |
| 2    |    | Benzene, 1,2,3,4-tetramethyl-  | 134 | C10H14  | 000488-23-3 | 35   |
| 3    |    | Benzene, 1-ethyl-3,5-dimethyl- | 134 | C10H14  | 000934-74-7 | 35   |
| 4    |    | Benzene, 1,2,3,4-tetramethyl-  | 134 | C10H14  | 000488-23-3 | 35   |
| 5    |    | Benzene, 1,2,3,5-tetramethyl-  | 134 | C10H14  | 000527-53-7 | 35   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040819\  
Data File : BM019822.D  
Acq On : 09 Apr 2019 08:22  
Operator : JU/SJ  
Sample : K2188-10  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF654

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

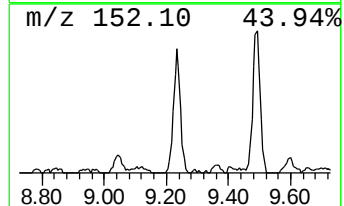
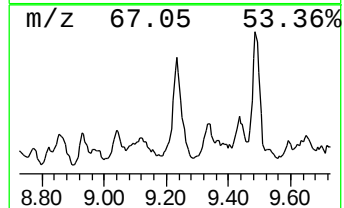
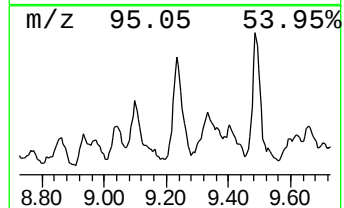
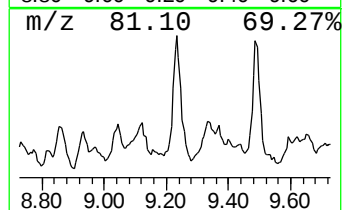
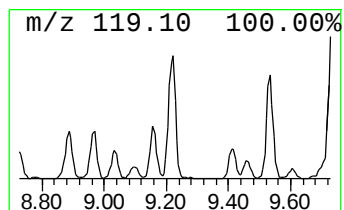
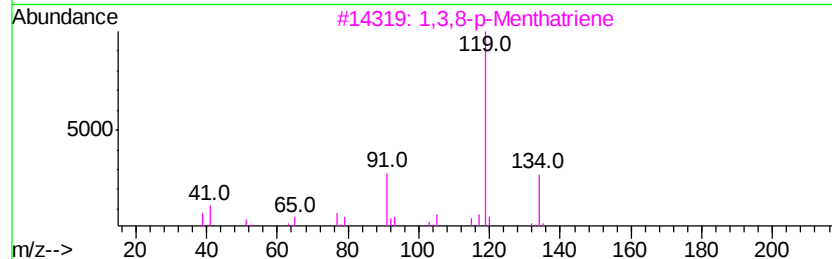
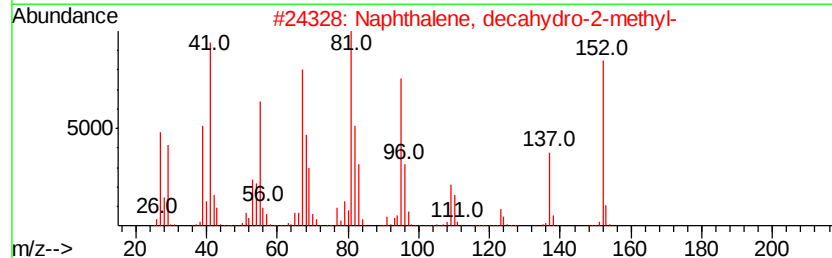
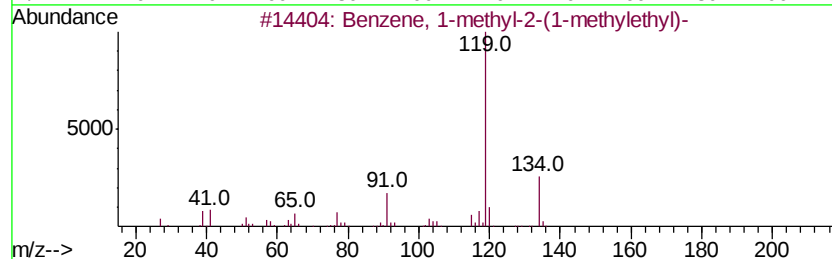
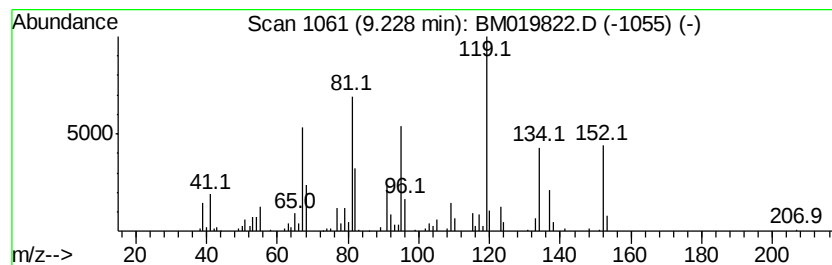
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 29 Naphthalene, decahydro-2-me... Concentration Rank 16

| R.T. | EstConc    | Area   | Relative to ISTD | R.T.  |
|------|------------|--------|------------------|-------|
| 9.23 | 5.64 ng/ul | 369480 | Naphthalene-d8   | 10.34 |

| Hit# | of | Tentative ID                         | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1-methyl-2-(1-methylethyl-) | 134 | C10H14  | 000527-84-4 | 80   |
| 2    |    | Naphthalene, decahydro-2-methyl-     | 152 | C11H20  | 002958-76-1 | 50   |
| 3    |    | 1,3,8-p-Menthatriene                 | 134 | C10H14  | 021195-59-5 | 42   |
| 4    |    | Benzene, 2-ethyl-1,3-dimethyl-       | 134 | C10H14  | 002870-04-4 | 42   |
| 5    |    | Benzene, 1,2,4,5-tetramethyl-        | 134 | C10H14  | 000095-93-2 | 42   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040819\  
Data File : BM019822.D  
Acq On : 09 Apr 2019 08:22  
Operator : JU/SJ  
Sample : K2188-10  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF654

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

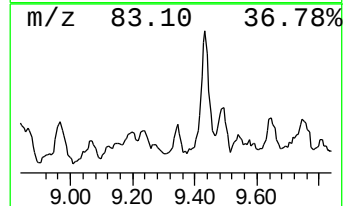
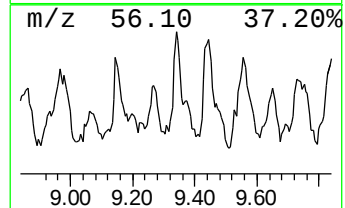
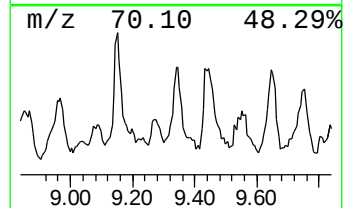
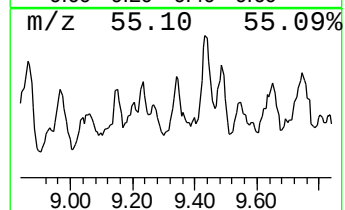
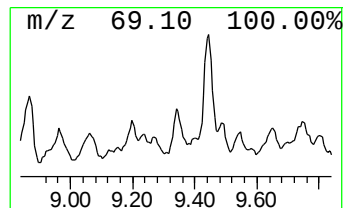
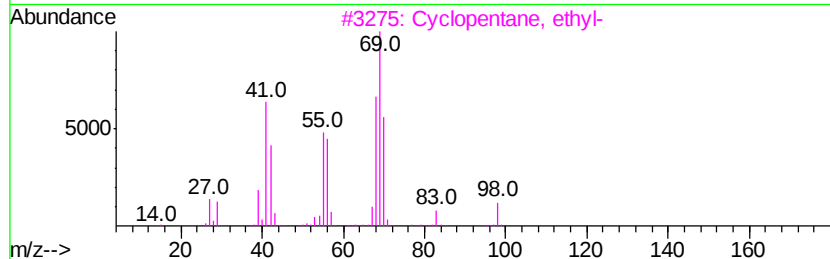
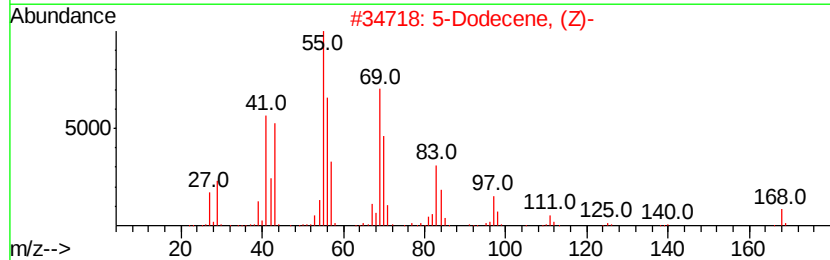
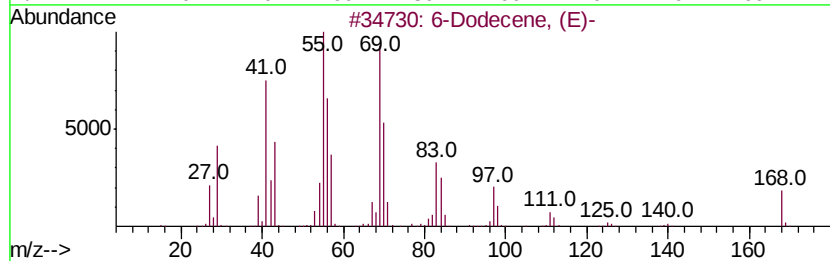
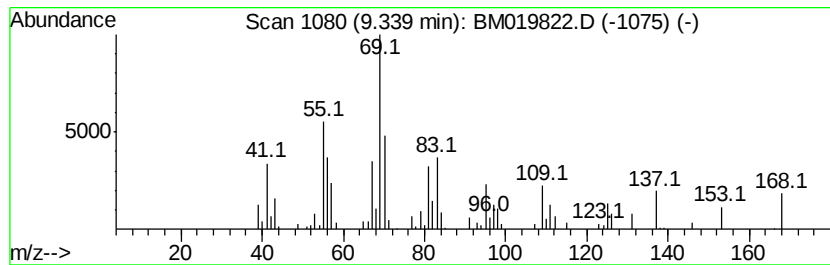
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 30 (DEL) Alkane: Cyclic9.34 Concentration Rank 33

| R.T. | EstConc    | Area   | Relative to ISTD | R.T.  |
|------|------------|--------|------------------|-------|
| 9.34 | 2.96 ng/ul | 194001 | Naphthalene-d8   | 10.34 |

| Hit# | of | 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------------|-----|---------|-------------|------|
| 1    |    |   | 6-Dodecene, (E)-            | 168 | C12H24  | 007206-17-9 | 64   |
| 2    |    |   | 5-Dodecene, (Z)-            | 168 | C12H24  | 007206-28-2 | 47   |
| 3    |    |   | Cyclopentane, ethyl-        | 98  | C7H14   | 001640-89-7 | 46   |
| 4    |    |   | 1-Hexene, 3,3,5-trimethyl-  | 126 | C9H18   | 013427-43-5 | 43   |
| 5    |    |   | 4-Undecene, 4-methyl-, (Z)- | 168 | C12H24  | 074630-57-2 | 43   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040819\  
Data File : BM019822.D  
Acq On : 09 Apr 2019 08:22  
Operator : JU/SJ  
Sample : K2188-10  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF654

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

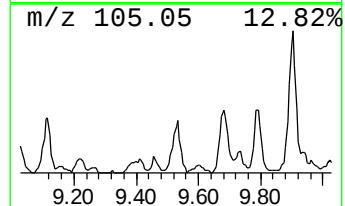
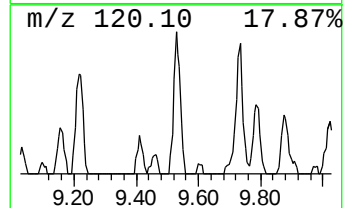
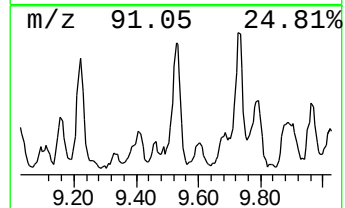
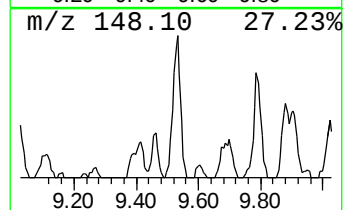
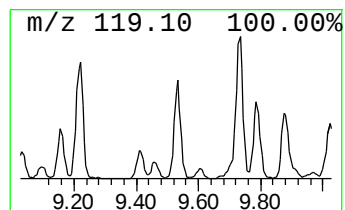
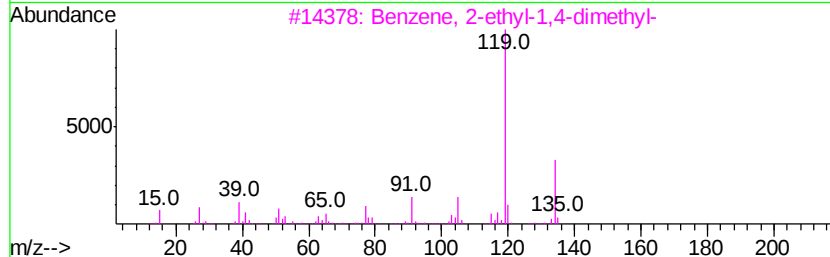
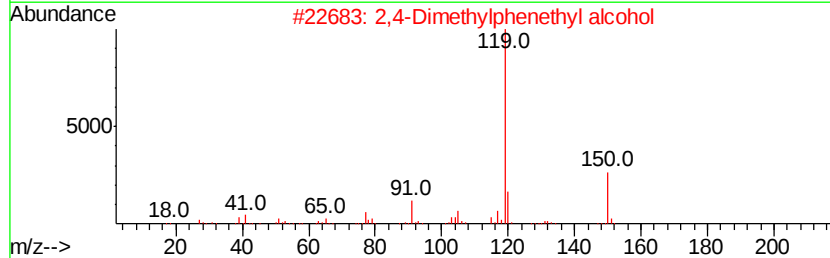
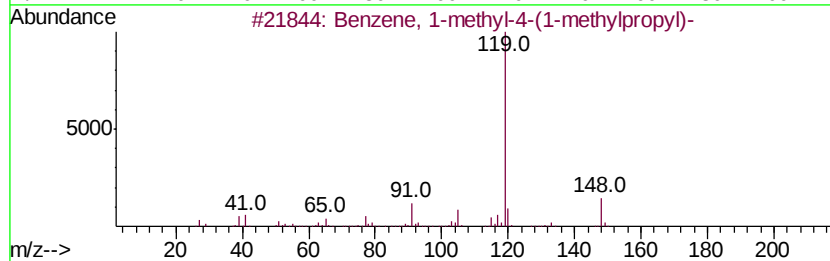
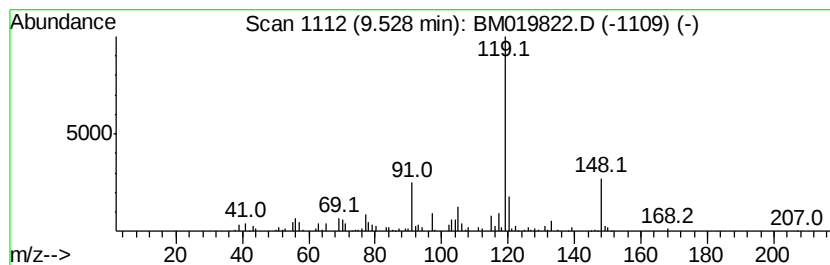
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 31 Benzene, 1-methyl-4-(1-meth... Concentration Rank 29

| R.T. | EstConc    | Area   | Relative to ISTD | R.T.  |
|------|------------|--------|------------------|-------|
| 9.53 | 3.28 ng/ul | 214966 | Napthalene-d8    | 10.34 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1-methyl-4-(1-methylpro... | 148 | C11H16  | 001595-16-0 | 87   |
| 2    |    | 2,4-Dimethylphenethyl alcohol       | 150 | C10H14O | 006597-59-7 | 76   |
| 3    |    | Benzene, 2-ethyl-1,4-dimethyl-      | 134 | C10H14  | 001758-88-9 | 64   |
| 4    |    | Benzene, 1-ethyl-3,5-dimethyl-      | 134 | C10H14  | 000934-74-7 | 64   |
| 5    |    | Benzene, (1,1-dimethylpropyl)-      | 148 | C11H16  | 002049-95-8 | 64   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040819\  
Data File : BM019822.D  
Acq On : 09 Apr 2019 08:22  
Operator : JU/SJ  
Sample : K2188-10  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF654

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

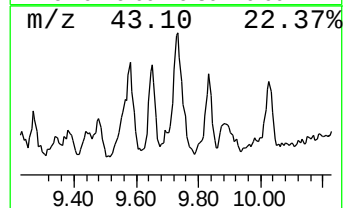
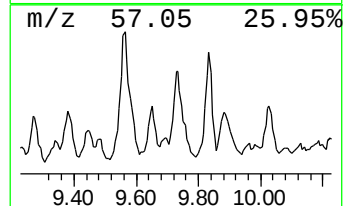
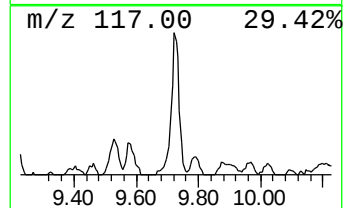
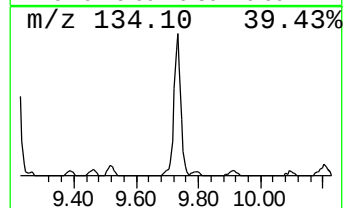
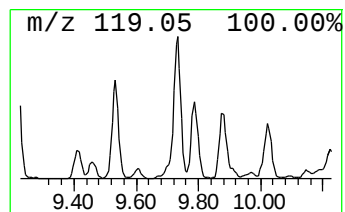
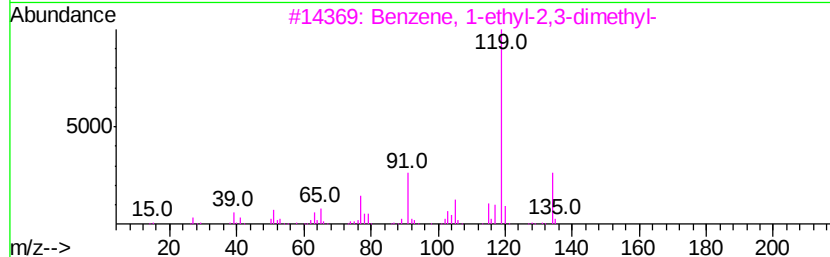
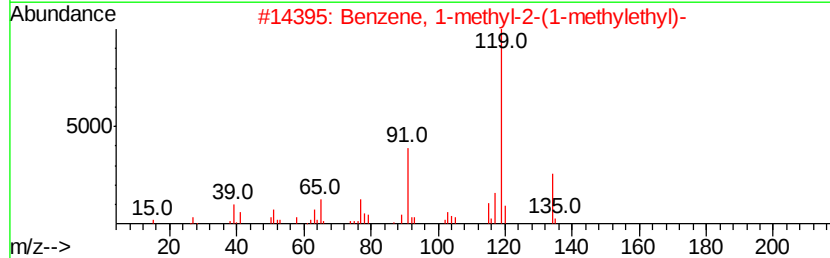
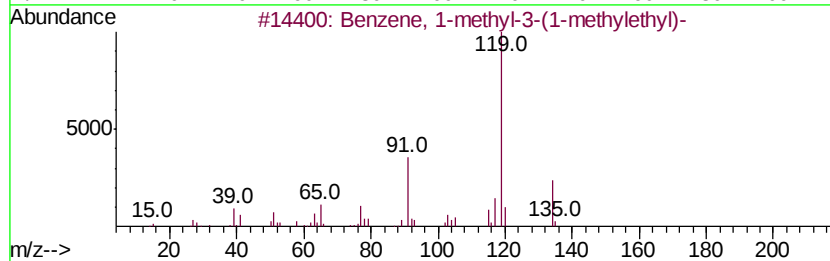
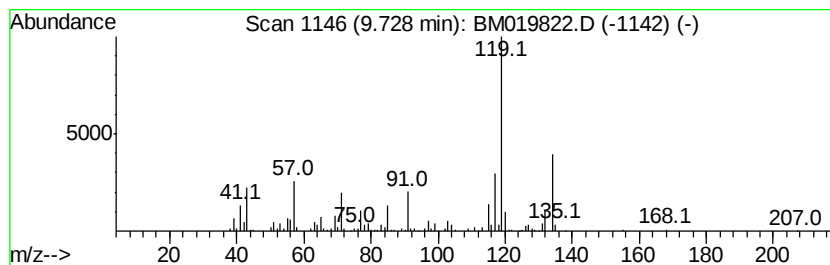
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 32 Benzene, 1-methyl-3-(1-meth... Concentration Rank 5

| R.T. | EstConc    | Area   | Relative to ISTD | R.T.  |
|------|------------|--------|------------------|-------|
| 9.73 | 6.86 ng/ul | 449732 | Naphthalene-d8   | 10.34 |

| Hit# | of | Tentative ID                         | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1-methyl-3-(1-methylethyl)- | 134 | C10H14  | 000535-77-3 | 87   |
| 2    |    | Benzene, 1-methyl-2-(1-methylethyl)- | 134 | C10H14  | 000527-84-4 | 87   |
| 3    |    | Benzene, 1-ethyl-2,3-dimethyl-       | 134 | C10H14  | 000933-98-2 | 81   |
| 4    |    | Benzene, 1-ethyl-3,5-dimethyl-       | 134 | C10H14  | 000934-74-7 | 81   |
| 5    |    | Benzene, 1,2,4,5-tetramethyl-        | 134 | C10H14  | 000095-93-2 | 81   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040819\  
Data File : BM019822.D  
Acq On : 09 Apr 2019 08:22  
Operator : JU/SJ  
Sample : K2188-10  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF654

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

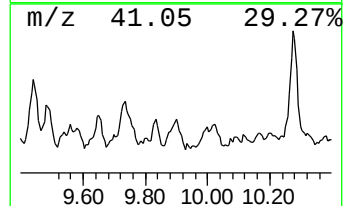
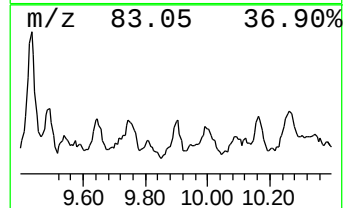
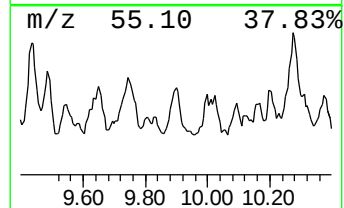
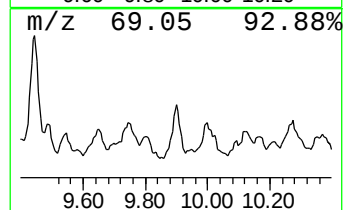
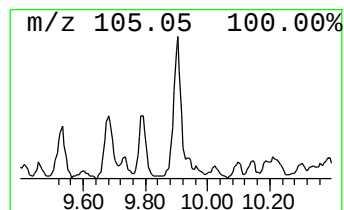
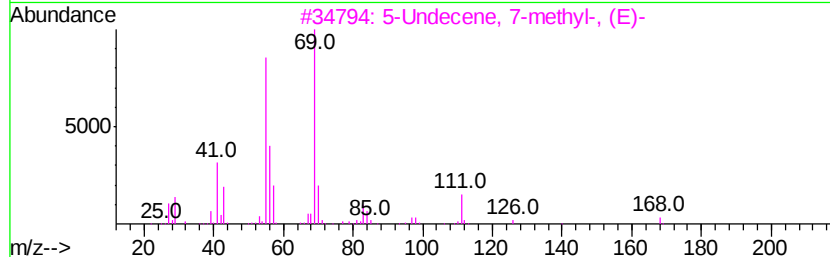
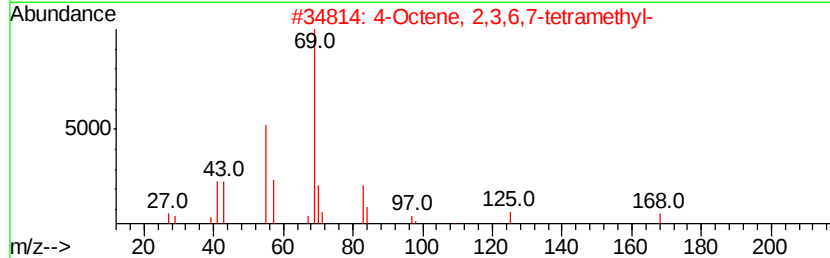
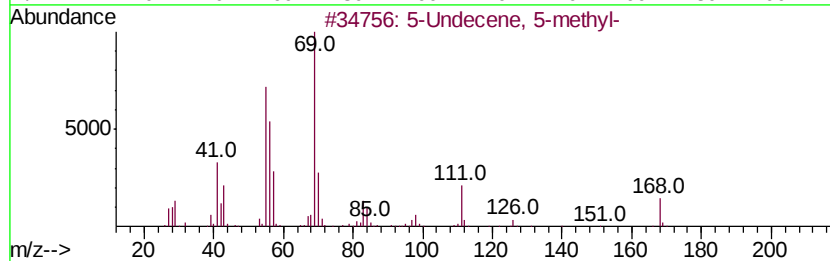
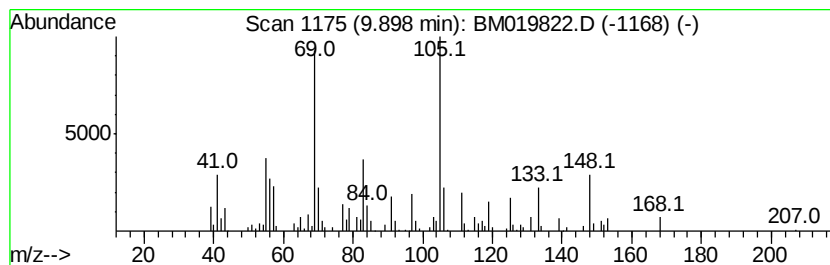
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 33 (DEL) Alkane: Cyclic9.90 Concentration Rank 17

| R.T. | EstConc    | Area   | Relative to ISTD | R.T.  |
|------|------------|--------|------------------|-------|
| 9.90 | 5.49 ng/ul | 359589 | Napthalene-d8    | 10.34 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | 5-Undecene, 5-methyl-               | 168 | C12H24  | 031613-73-7 | 41   |
| 2    |    | 4-Octene, 2,3,6,7-tetramethyl-      | 168 | C12H24  | 063830-66-0 | 38   |
| 3    |    | 5-Undecene, 7-methyl-, (E)-         | 168 | C12H24  | 074630-66-3 | 38   |
| 4    |    | Cyclohexane, 2-propyl-1,1,3-trim... | 168 | C12H24  | 081983-70-2 | 35   |
| 5    |    | 1-Hexene, 3,3,5-trimethyl-          | 126 | C9H18   | 013427-43-5 | 27   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040819\  
Data File : BM019822.D  
Acq On : 09 Apr 2019 08:22  
Operator : JU/SJ  
Sample : K2188-10  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF654

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

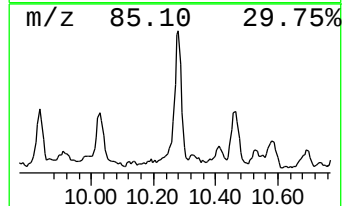
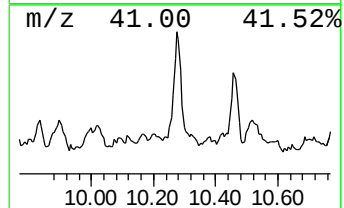
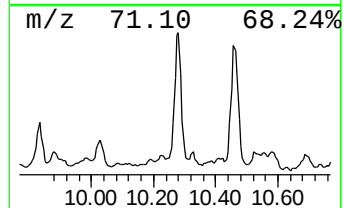
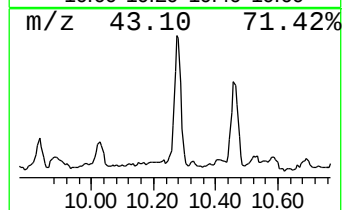
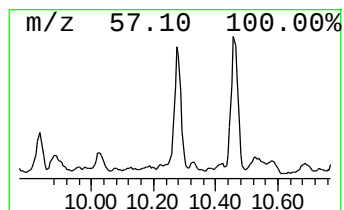
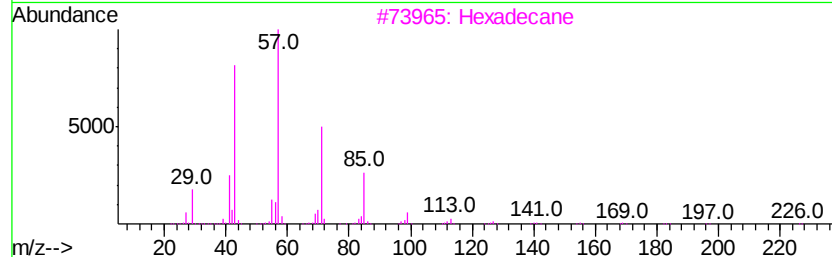
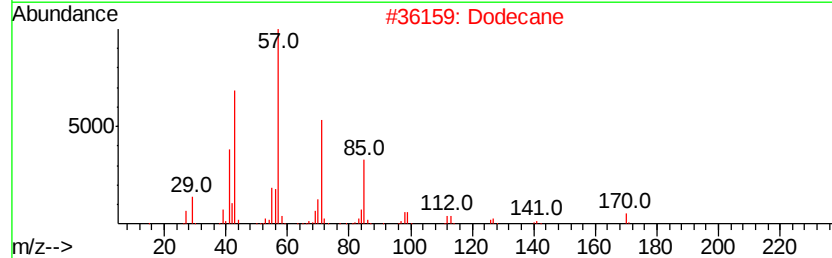
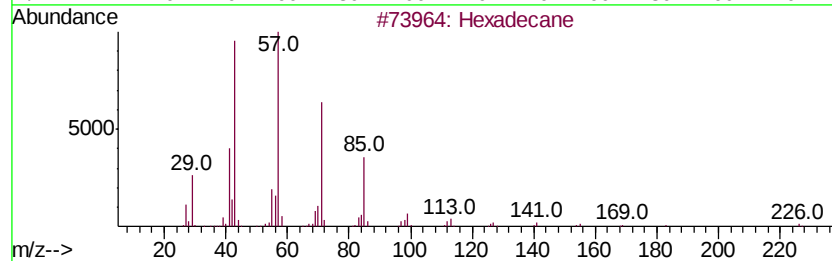
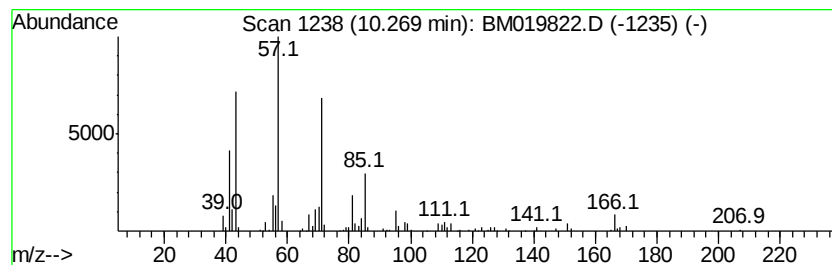
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 34 (DEL) Alkane: Straight-Chai... Concentration Rank 9

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 10.27 | 6.31 ng/ul | 413734 | Napthalene-d8    | 10.34 |

| Hit# | of | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|--------------|-----|---------|-------------|------|
| 1    | 5  | Hexadecane   | 226 | C16H34  | 000544-76-3 | 80   |
| 2    |    | Dodecane     | 170 | C12H26  | 000112-40-3 | 80   |
| 3    |    | Hexadecane   | 226 | C16H34  | 000544-76-3 | 80   |
| 4    |    | Tridecane    | 184 | C13H28  | 000629-50-5 | 80   |
| 5    |    | Tetradecane  | 198 | C14H30  | 000629-59-4 | 72   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040819\  
Data File : BM019822.D  
Acq On : 09 Apr 2019 08:22  
Operator : JU/SJ  
Sample : K2188-10  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF654

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

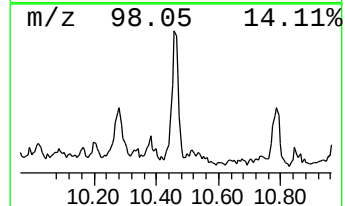
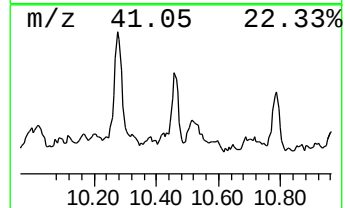
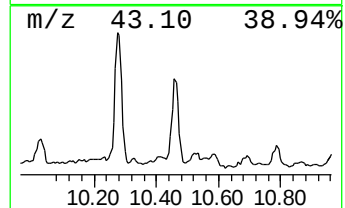
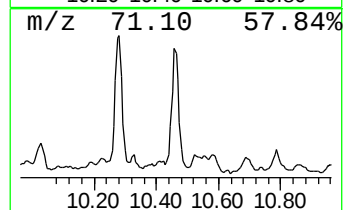
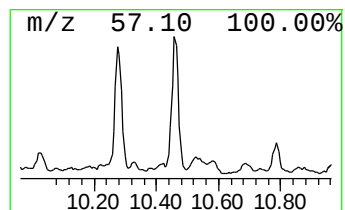
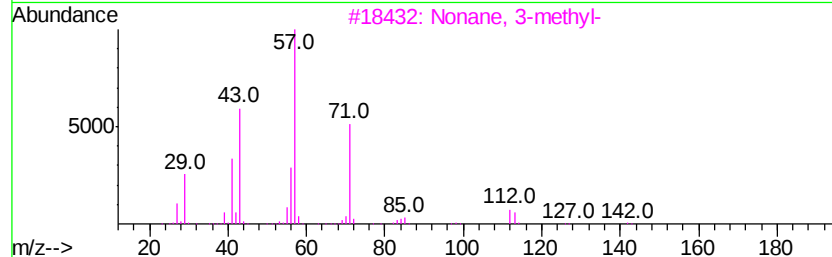
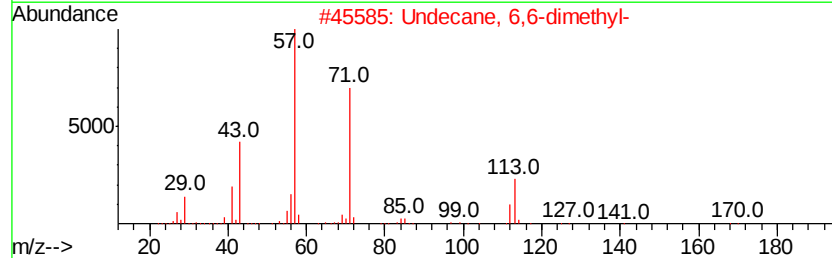
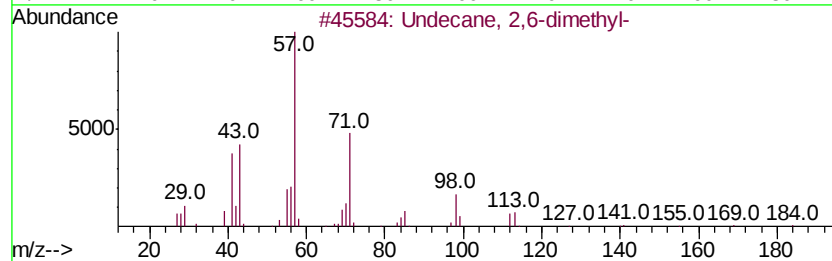
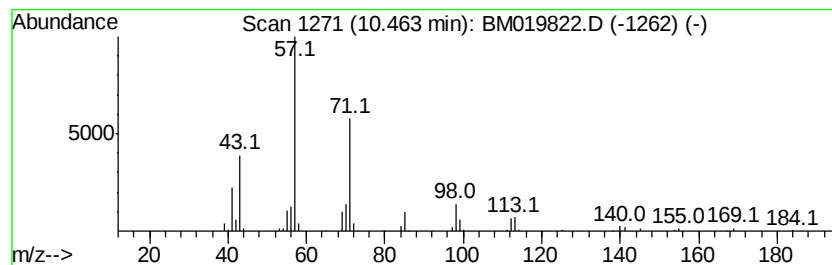
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 35 (DEL) Alkane: Straight-Chai... Concentration Rank 8

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 10.46 | 6.63 ng/ul | 434620 | Napthalene-d8    | 10.34 |

| Hit# | of | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------|-----|---------|-------------|------|
| 1    | 5  | Undecane, 2,6-dimethyl- | 184 | C13H28  | 017301-23-4 | 83   |
| 2    |    | Undecane, 6,6-dimethyl- | 184 | C13H28  | 017312-76-4 | 53   |
| 3    |    | Nonane, 3-methyl-       | 142 | C10H22  | 005911-04-6 | 53   |
| 4    |    | Dodecane, 6-methyl-     | 184 | C13H28  | 006044-71-9 | 53   |
| 5    |    | Undecane, 3,6-dimethyl- | 184 | C13H28  | 017301-28-9 | 53   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040819\  
Data File : BM019822.D  
Acq On : 09 Apr 2019 08:22  
Operator : JU/SJ  
Sample : K2188-10  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF654

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

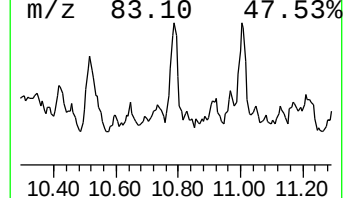
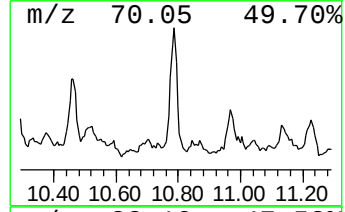
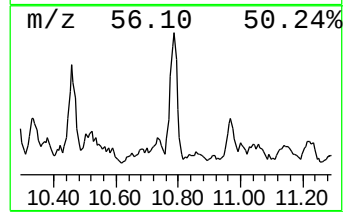
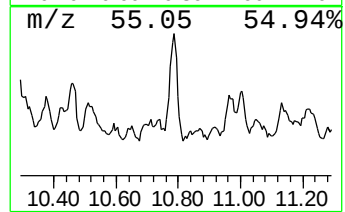
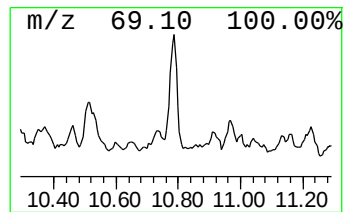
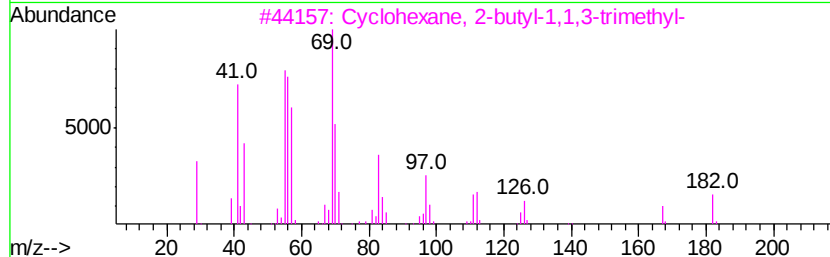
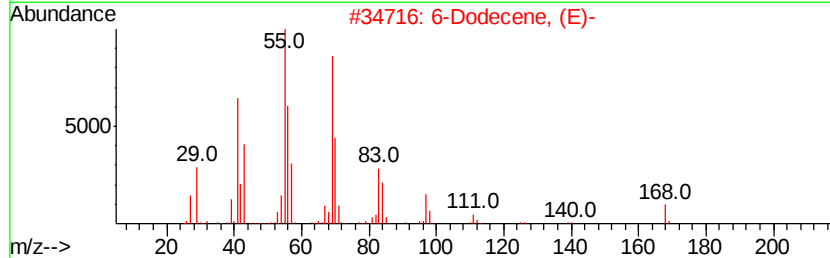
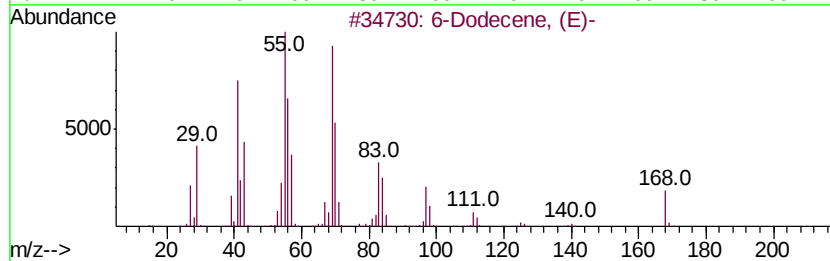
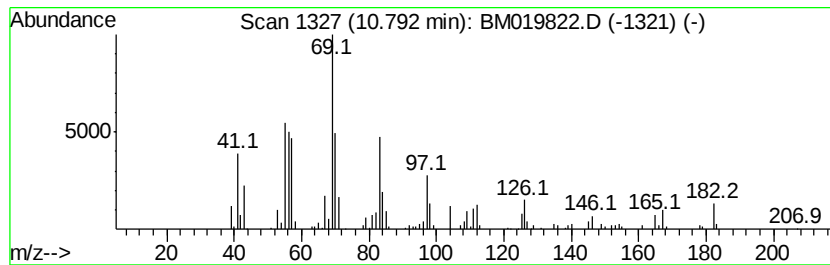
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 36 (DEL) Alkane: Cyclic10.79 Concentration Rank 18

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 10.79 | 5.46 ng/ul | 357567 | Napthalene-d8    | 10.34 |

| Hit# | of                                  | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|-------------------------------------|--------------|-----|---------|-------------|------|
| 1    | 6-Dodecene, (E)-                    |              | 168 | C12H24  | 007206-17-9 | 87   |
| 2    | 6-Dodecene, (E)-                    |              | 168 | C12H24  | 007206-17-9 | 87   |
| 3    | Cyclohexane, 2-butyl-1,1,3-trime... |              | 182 | C13H26  | 054676-39-0 | 81   |
| 4    | Cyclohexane, 2-butyl-1,1,3-trime... |              | 182 | C13H26  | 054676-39-0 | 78   |
| 5    | Cyclohexane, 2-butyl-1,1,3-trime... |              | 182 | C13H26  | 054676-39-0 | 60   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040819\  
Data File : BM019822.D  
Acq On : 09 Apr 2019 08:22  
Operator : JU/SJ  
Sample : K2188-10  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF654

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

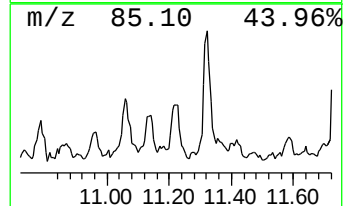
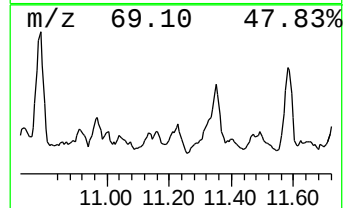
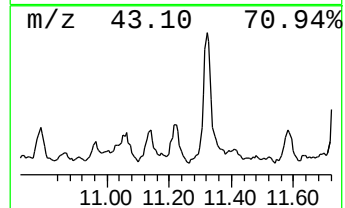
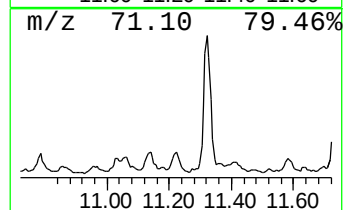
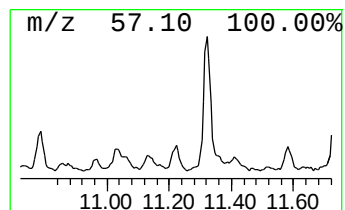
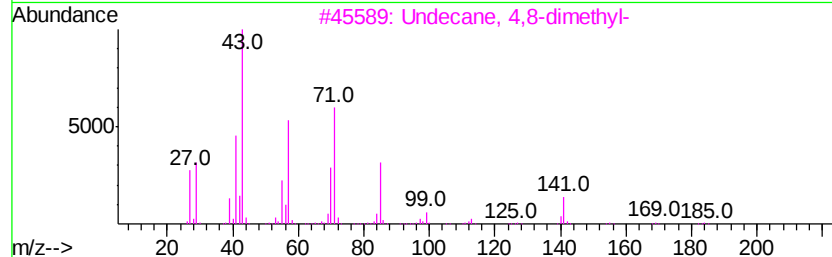
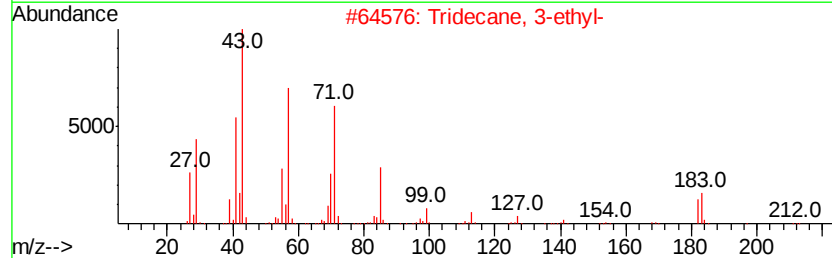
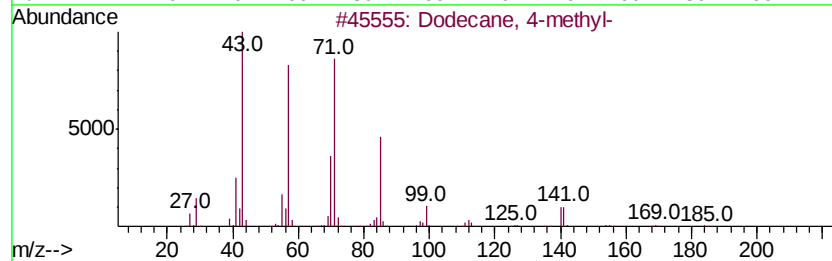
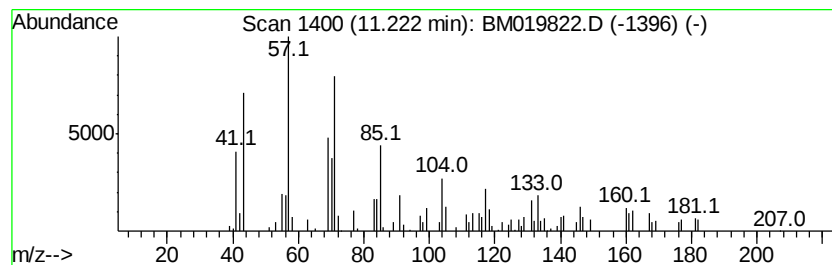
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 37 (DEL) Alkane: Straight-Chai... Concentration Rank 38

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 11.22 | 2.75 ng/ul | 180213 | Napthalene-d8    | 10.34 |

| Hit# | of | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------|-----|---------|-------------|------|
| 1    | 5  | Dodecane, 4-methyl-     | 184 | C13H28  | 006117-97-1 | 43   |
| 2    |    | Tridecane, 3-ethyl-     | 212 | C15H32  | 013286-73-2 | 43   |
| 3    |    | Undecane, 4,8-dimethyl- | 184 | C13H28  | 017301-33-6 | 43   |
| 4    |    | Undecane, 3,8-dimethyl- | 184 | C13H28  | 017301-30-3 | 43   |
| 5    |    | Heptane, 2,4-dimethyl-  | 128 | C9H20   | 002213-23-2 | 43   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040819\  
Data File : BM019822.D  
Acq On : 09 Apr 2019 08:22  
Operator : JU/SJ  
Sample : K2188-10  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF654

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

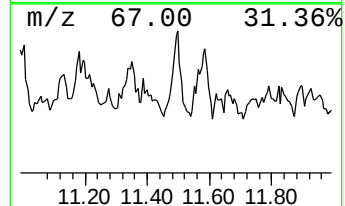
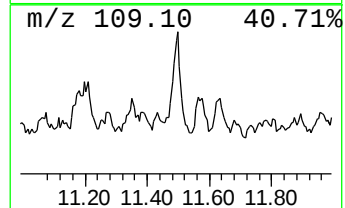
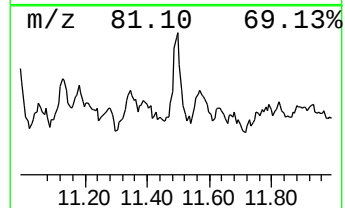
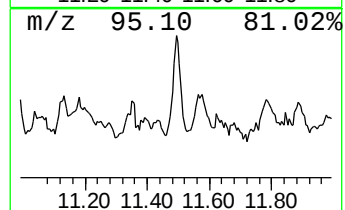
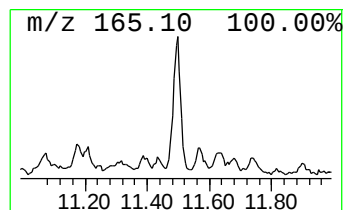
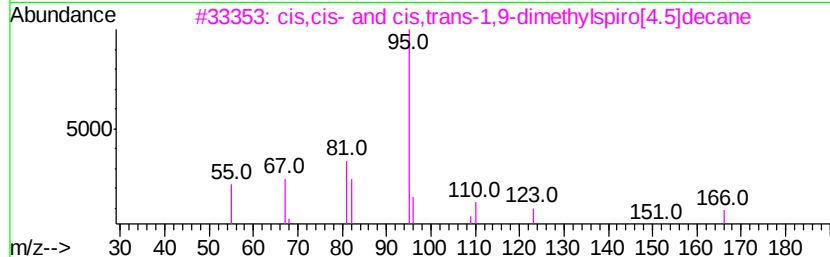
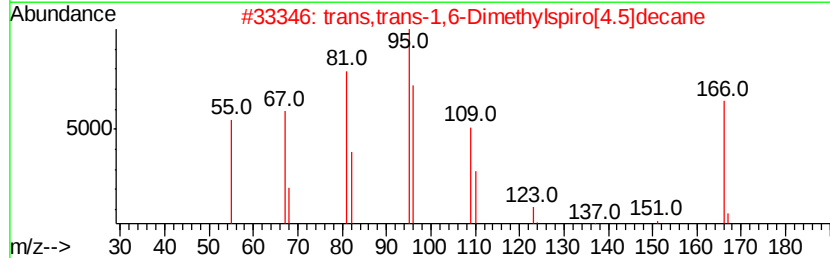
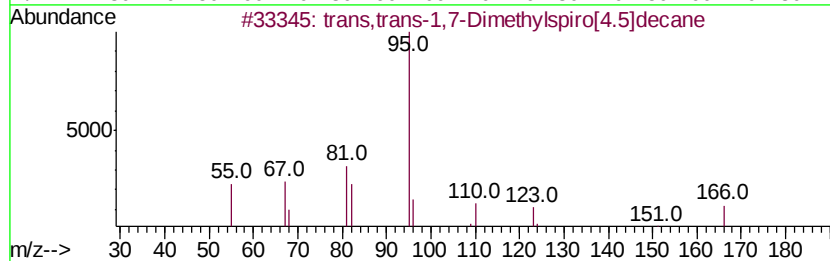
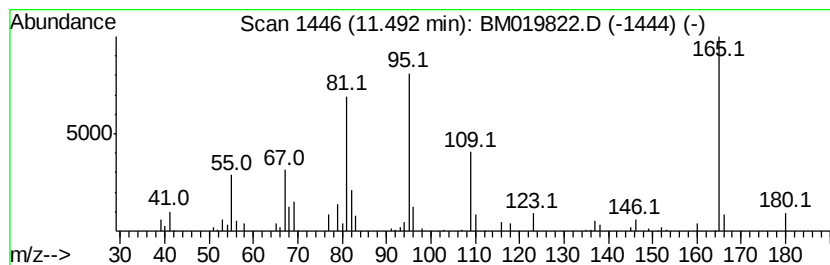
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 38 unknown-05 Concentration Rank 51

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 11.49 | 2.16 ng/ul | 141585 | Naphthalene-d8   | 10.34 |

| Hit# | of                                                    | 5   | Tentative ID | MW           | MolForm | CAS# | Qual |
|------|-------------------------------------------------------|-----|--------------|--------------|---------|------|------|
| 1    | trans,trans-1,7-Dimethylspiro[4.4]undecane            | 166 | C12H22       | 1000111-72-6 | 43      |      |      |
| 2    | trans,trans-1,6-Dimethylspiro[4.4]undecane            | 166 | C12H22       | 1000111-72-1 | 43      |      |      |
| 3    | cis,cis- and cis,trans-1,9-dimethylspiro[4.4]undecane | 166 | C12H22       | 1000111-72-5 | 43      |      |      |
| 4    | cis,trans-1,6-Dimethylspiro[4.5]undecane              | 166 | C12H22       | 1000111-72-3 | 38      |      |      |
| 5    | 1-Propanone, 1-[2-(1,1-dimethylethyl)-2-oxoethyl]-    | 180 | C12H20O      | 019576-21-7  | 38      |      |      |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040819\  
Data File : BM019822.D  
Acq On : 09 Apr 2019 08:22  
Operator : JU/SJ  
Sample : K2188-10  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF654

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

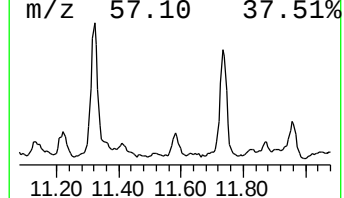
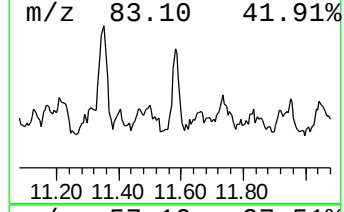
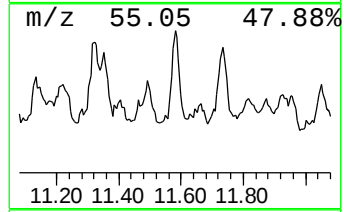
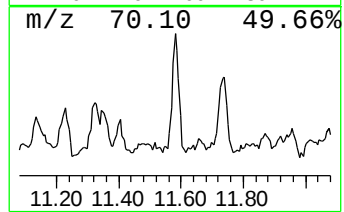
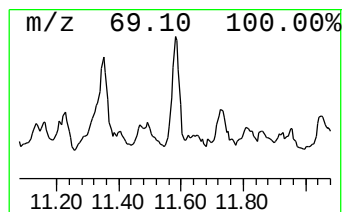
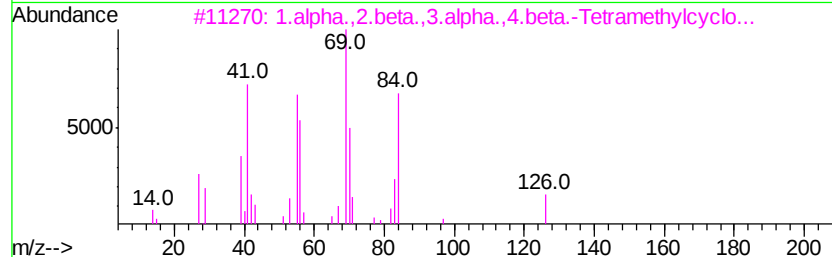
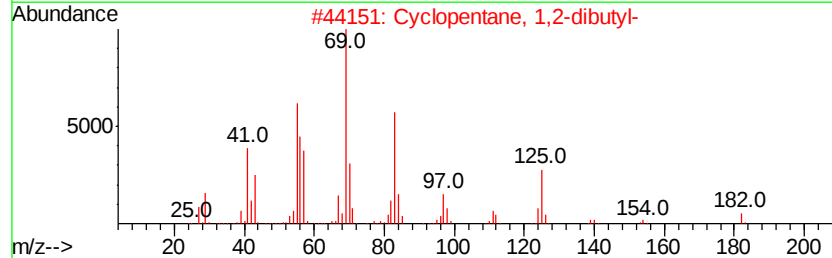
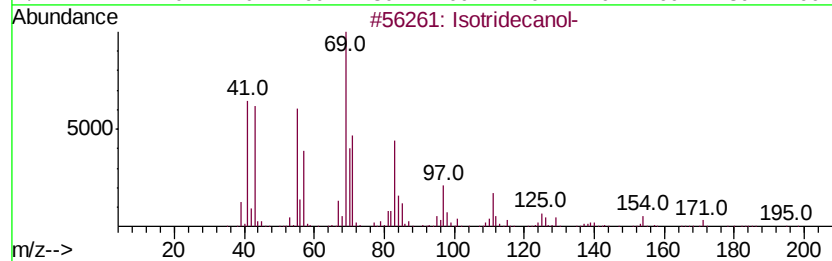
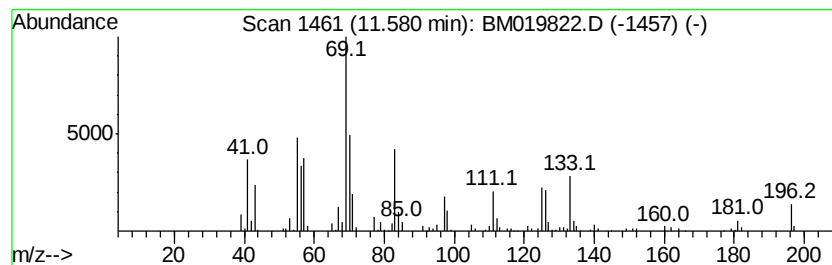
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 39 Isotridecanol- Concentration Rank 25

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 11.58 | 3.87 ng/ul | 253854 | Naphthalene-d8   | 10.34 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Isotridecanol-                      | 200 | C13H28O | 027458-92-0 | 59   |
| 2    |    | Cyclopentane, 1,2-dibutyl-          | 182 | C13H26  | 062199-52-4 | 58   |
| 3    |    | 1.alpha.,2.beta.,3.alpha.,4.beta... | 126 | C9H18   | 002532-67-4 | 47   |
| 4    |    | Cyclohexane, 1,2-diethyl-3-methyl-  | 154 | C11H22  | 061141-80-8 | 47   |
| 5    |    | 2-Methyl-2-octene                   | 126 | C9H18   | 016993-86-5 | 38   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040819\  
Data File : BM019822.D  
Acq On : 09 Apr 2019 08:22  
Operator : JU/SJ  
Sample : K2188-10  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF654

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

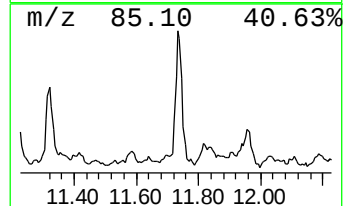
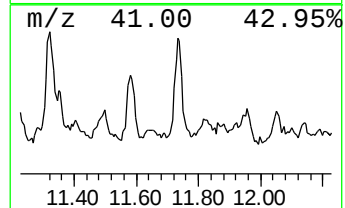
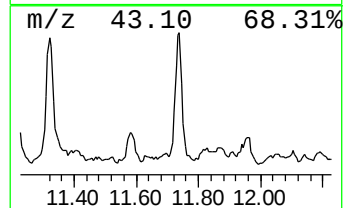
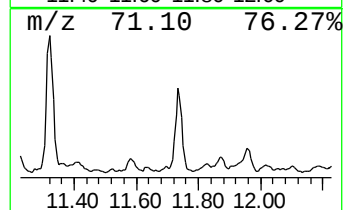
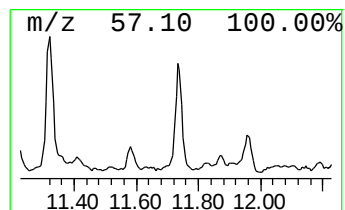
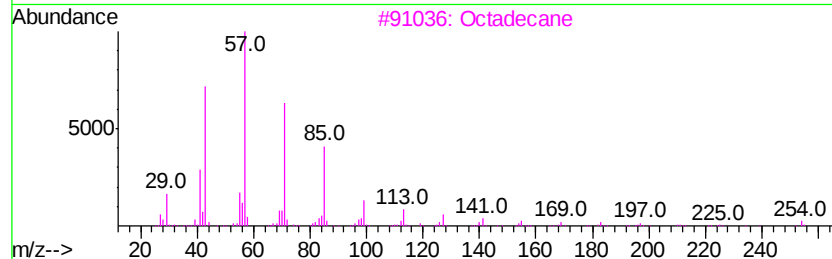
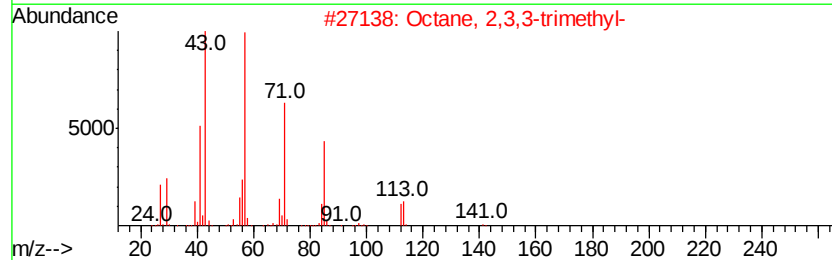
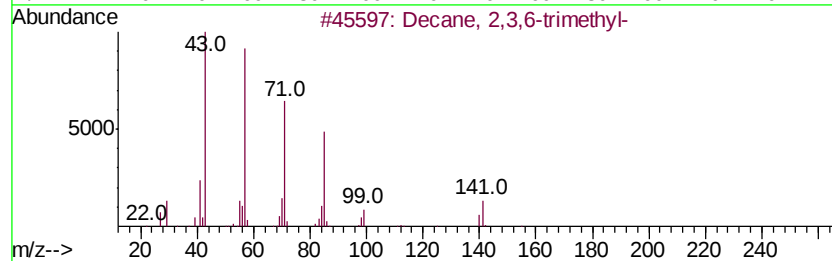
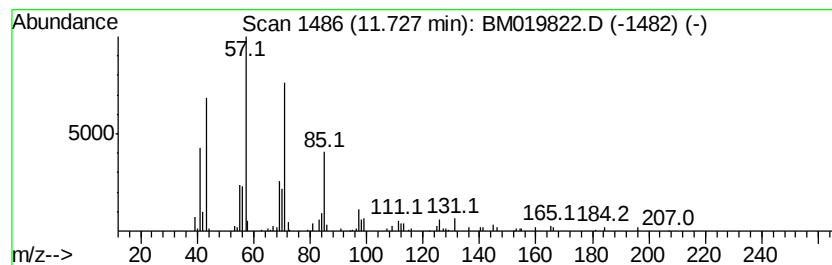
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 40 (DEL) Alkane: Straight-Chai... Concentration Rank 22

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 11.73 | 4.40 ng/ul | 288242 | Napthalene-d8    | 10.34 |

| Hit# | of | 5 | Tentative ID             | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------|-----|---------|-------------|------|
| 1    |    |   | Decane, 2,3,6-trimethyl- | 184 | C13H28  | 062238-12-4 | 72   |
| 2    |    |   | Octane, 2,3,3-trimethyl- | 156 | C11H24  | 062016-30-2 | 72   |
| 3    |    |   | Octadecane               | 254 | C18H38  | 000593-45-3 | 72   |
| 4    |    |   | Undecane, 5,7-dimethyl-  | 184 | C13H28  | 017312-83-3 | 72   |
| 5    |    |   | 3-Ethyl-3-methylheptane  | 142 | C10H22  | 017302-01-1 | 64   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040819\  
Data File : BM019822.D  
Acq On : 09 Apr 2019 08:22  
Operator : JU/SJ  
Sample : K2188-10  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF654

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

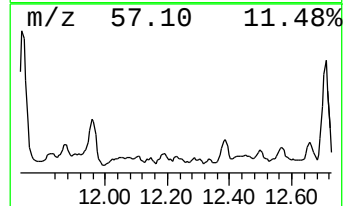
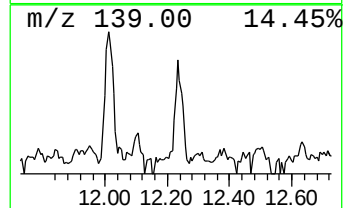
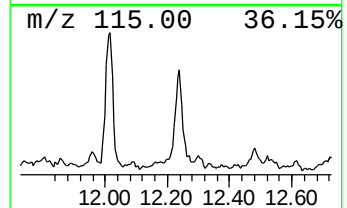
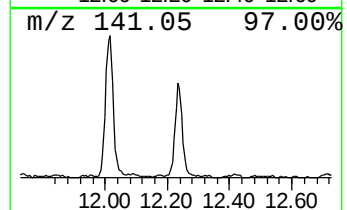
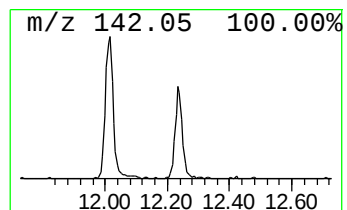
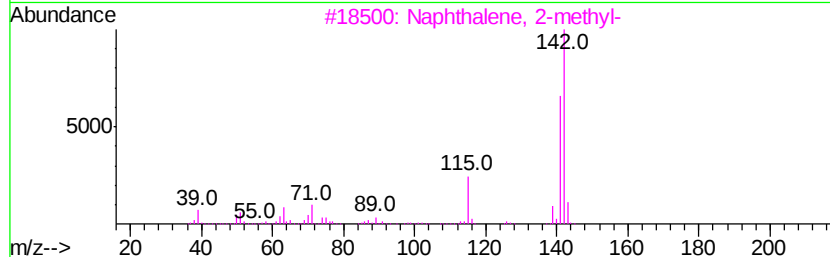
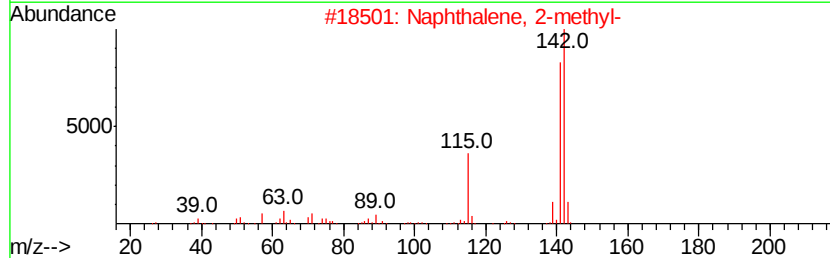
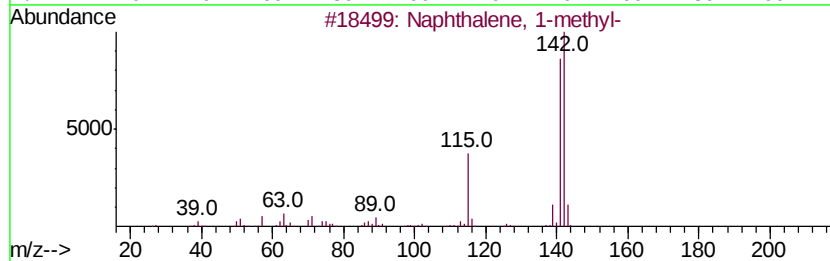
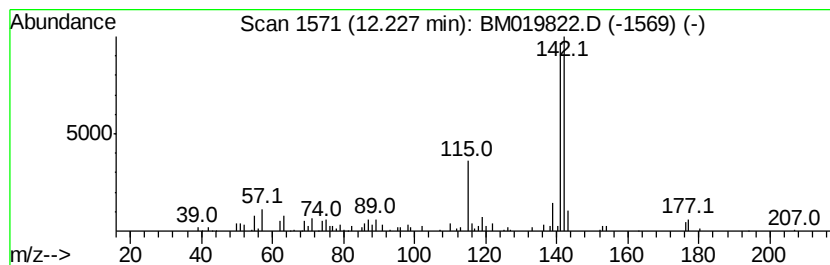
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 41 Naphthalene, 1-methyl- Concentration Rank 34

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 12.23 | 2.92 ng/ul | 191295 | Naphthalene-d8   | 10.34 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 1-methyl-              | 142 | C11H10  | 000090-12-0 | 94   |
| 2    |    | Naphthalene, 2-methyl-              | 142 | C11H10  | 000091-57-6 | 93   |
| 3    |    | Naphthalene, 2-methyl-              | 142 | C11H10  | 000091-57-6 | 93   |
| 4    |    | Naphthalene, 1-methyl-              | 142 | C11H10  | 000090-12-0 | 93   |
| 5    |    | 1,4-Methanonaphthalene, 1,4-dihy... | 142 | C11H10  | 004453-90-1 | 93   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040819\  
Data File : BM019822.D  
Acq On : 09 Apr 2019 08:22  
Operator : JU/SJ  
Sample : K2188-10  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF654

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

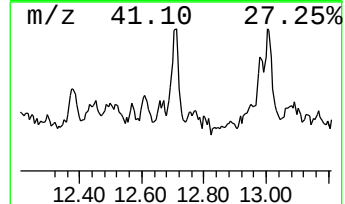
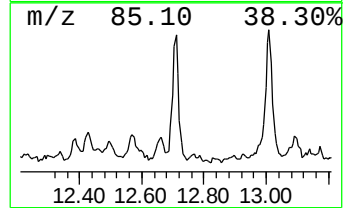
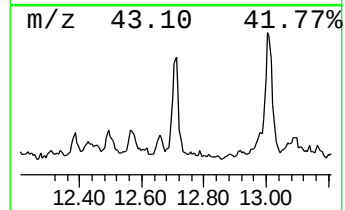
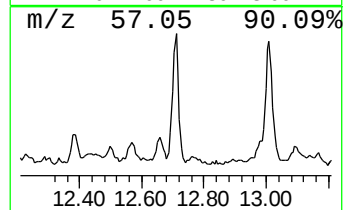
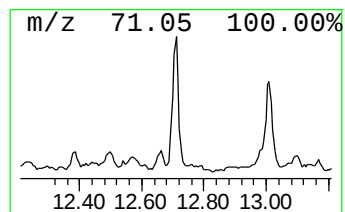
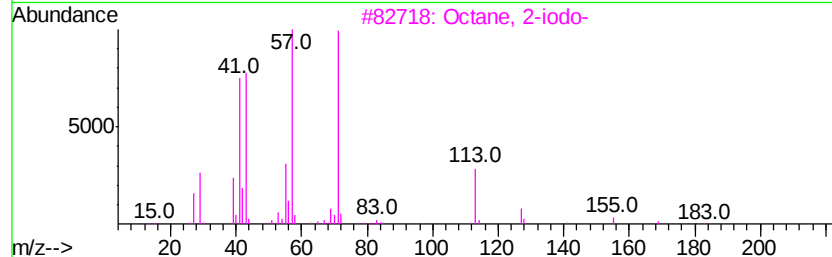
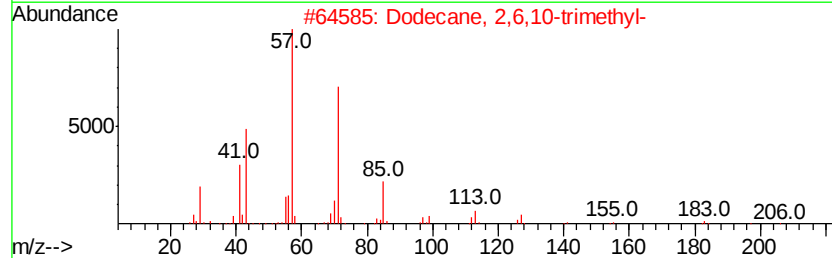
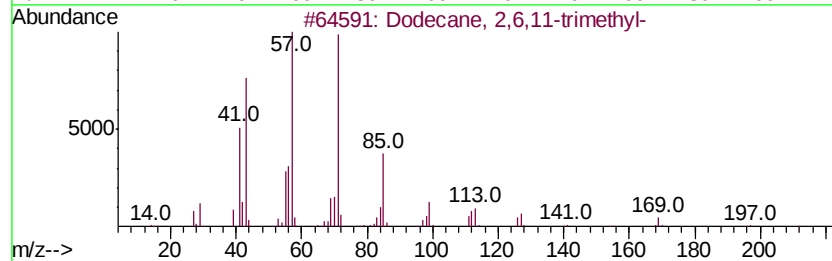
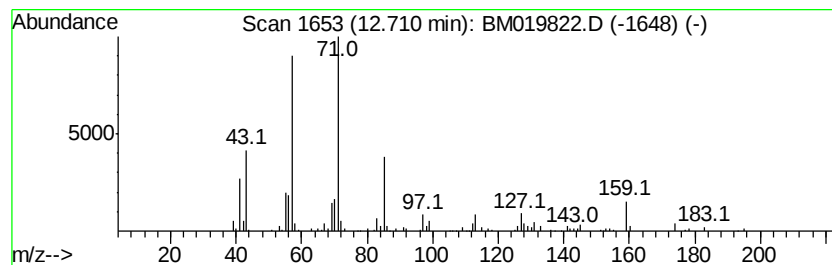
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 42 (DEL) Alkane: Straight-Chai... Concentration Rank 35

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 12.71 | 2.87 ng/ul | 255144 | Acenaphthene-d10 | 14.22 |

| Hit# | of | 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------------|-----|---------|-------------|------|
| 1    |    |   | Dodecane, 2,6,11-trimethyl- | 212 | C15H32  | 031295-56-4 | 64   |
| 2    |    |   | Dodecane, 2,6,10-trimethyl- | 212 | C15H32  | 003891-98-3 | 64   |
| 3    |    |   | Octane, 2-iodo-             | 240 | C8H17I  | 000557-36-8 | 47   |
| 4    |    |   | Octane, 2,6-dimethyl-       | 142 | C10H22  | 002051-30-1 | 47   |
| 5    |    |   | 1-Decene, 4-methyl-         | 154 | C11H22  | 013151-29-6 | 43   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040819\  
Data File : BM019822.D  
Acq On : 09 Apr 2019 08:22  
Operator : JU/SJ  
Sample : K2188-10  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF654

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

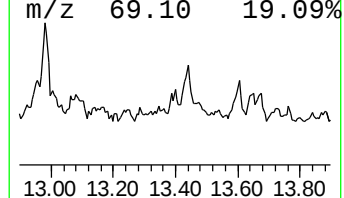
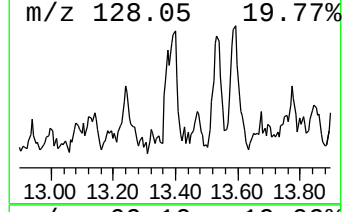
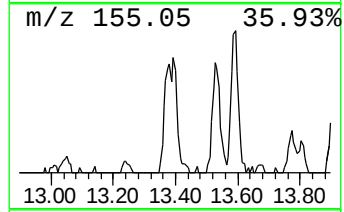
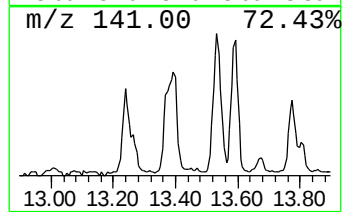
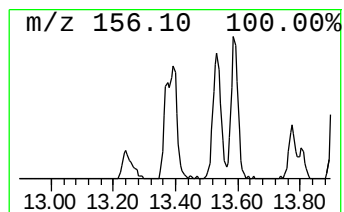
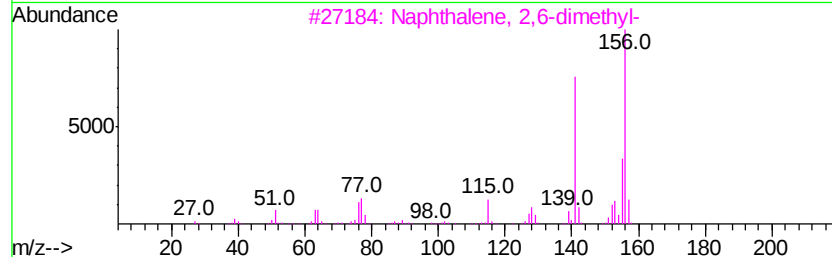
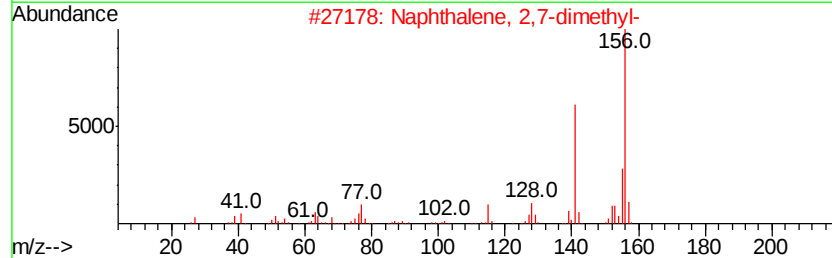
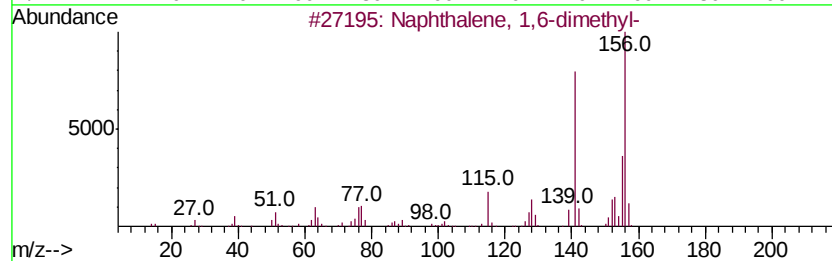
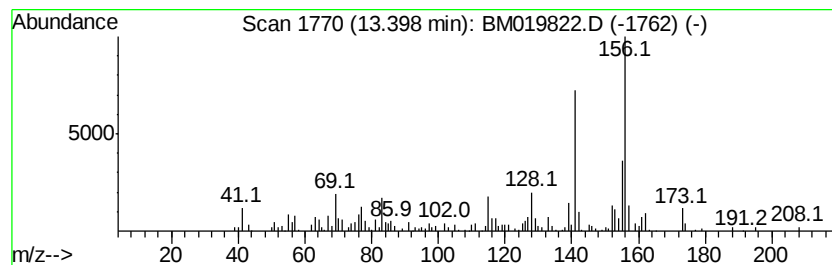
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 43 Naphthalene, 1,6-dimethyl- Concentration Rank 46

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 13.40 | 2.55 ng/ul | 226570 | Acenaphthene-d10 | 14.22 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 1,6-dimethyl- | 156 | C12H12  | 000575-43-9 | 96   |
| 2    |    | Naphthalene, 2,7-dimethyl- | 156 | C12H12  | 000582-16-1 | 95   |
| 3    |    | Naphthalene, 2,6-dimethyl- | 156 | C12H12  | 000581-42-0 | 95   |
| 4    |    | Naphthalene, 2,7-dimethyl- | 156 | C12H12  | 000582-16-1 | 95   |
| 5    |    | Naphthalene, 1,7-dimethyl- | 156 | C12H12  | 000575-37-1 | 94   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040819\  
Data File : BM019822.D  
Acq On : 09 Apr 2019 08:22  
Operator : JU/SJ  
Sample : K2188-10  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF654

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

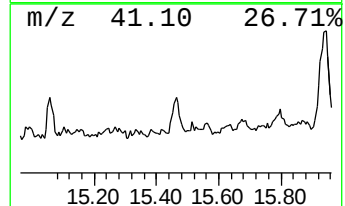
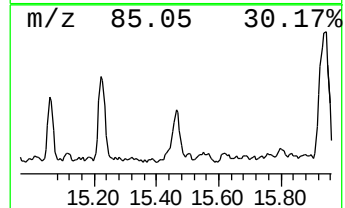
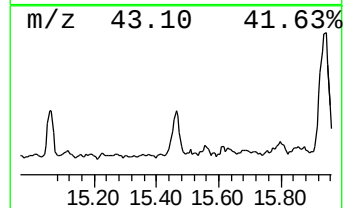
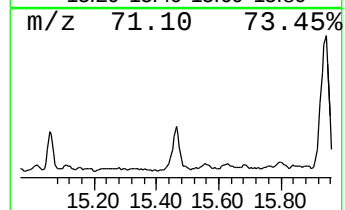
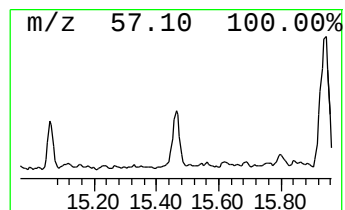
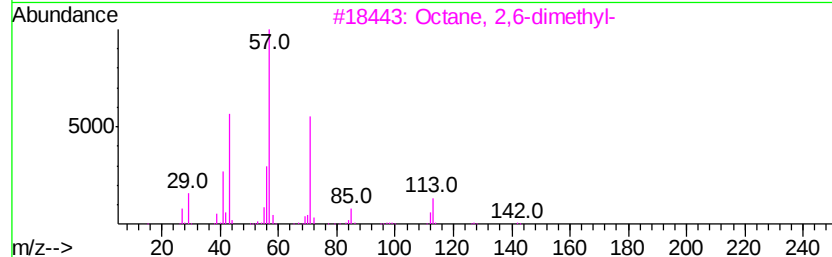
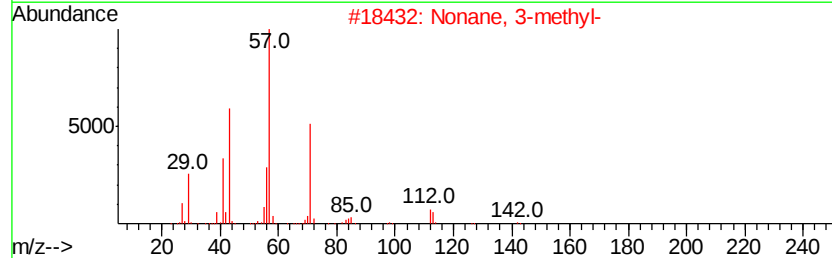
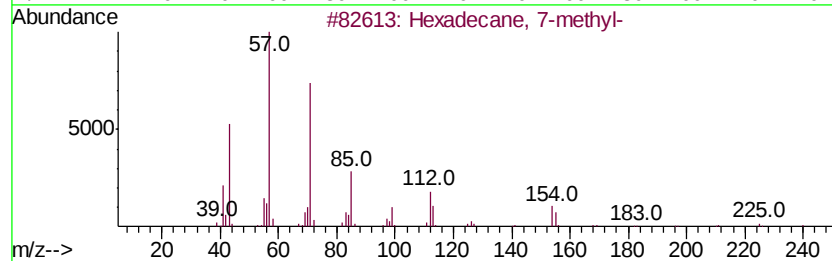
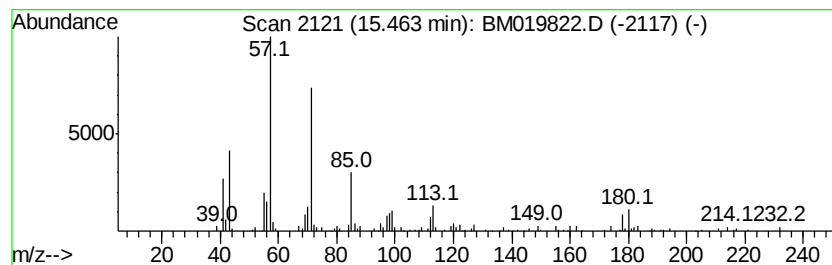
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 44 (DEL) Alkane: Straight-Chai... Concentration Rank 50

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 15.46 | 2.18 ng/ul | 193697 | Acenaphthene-d10 | 14.22 |

| Hit# | of | 5 | Tentative ID             | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------|-----|---------|-------------|------|
| 1    |    |   | Hexadecane, 7-methyl-    | 240 | C17H36  | 026730-20-1 | 64   |
| 2    |    |   | Nonane, 3-methyl-        | 142 | C10H22  | 005911-04-6 | 52   |
| 3    |    |   | Octane, 2,6-dimethyl-    | 142 | C10H22  | 002051-30-1 | 52   |
| 4    |    |   | Undecane, 6,6-dimethyl-  | 184 | C13H28  | 017312-76-4 | 52   |
| 5    |    |   | Decane, 2,3,8-trimethyl- | 184 | C13H28  | 062238-14-6 | 50   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040819\  
Data File : BM019822.D  
Acq On : 09 Apr 2019 08:22  
Operator : JU/SJ  
Sample : K2188-10  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF654

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

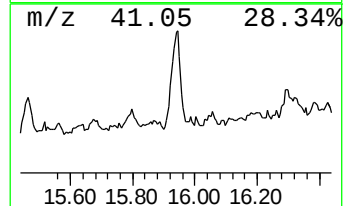
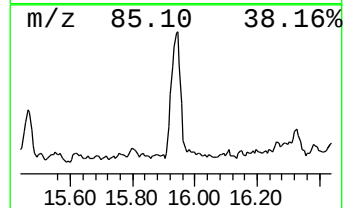
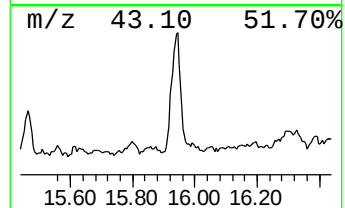
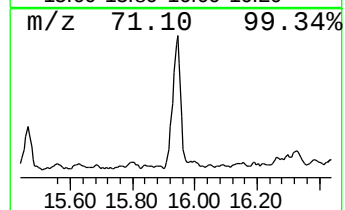
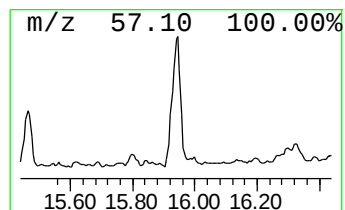
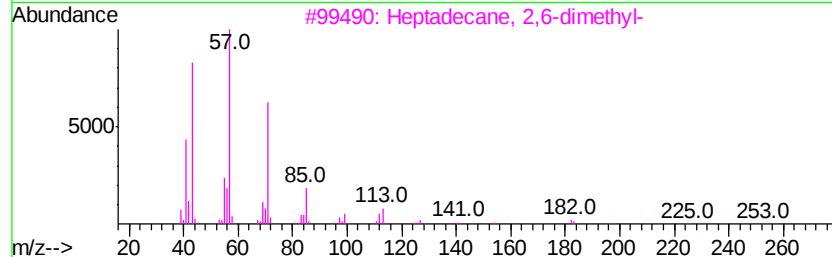
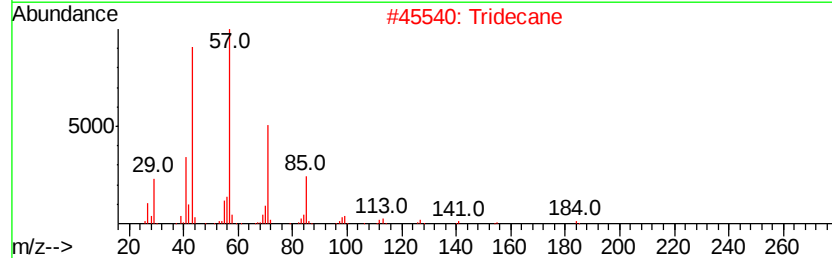
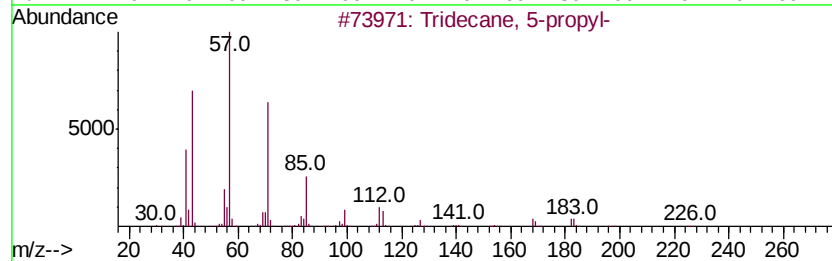
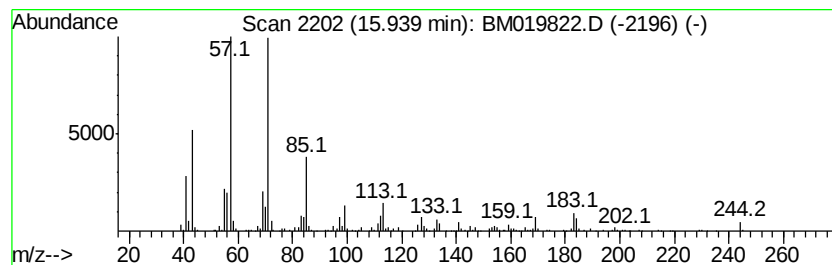
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 45 (DEL) Alkane: Straight-Chai... Concentration Rank 10

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 15.94 | 6.28 ng/ul | 613699 | Phenanthrene-d10 | 16.99 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Tridecane, 5-propyl-                | 226 | C16H34  | 055045-11-9 | 72   |
| 2    |    | Tridecane                           | 184 | C13H28  | 000629-50-5 | 64   |
| 3    |    | Heptadecane, 2,6-dimethyl-          | 268 | C19H40  | 054105-67-8 | 59   |
| 4    |    | Pentadecane, 2,6,10,14-tetramethyl- | 268 | C19H40  | 001921-70-6 | 52   |
| 5    |    | Pentadecane, 2,6,10,14-tetramethyl- | 268 | C19H40  | 001921-70-6 | 52   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040819\  
Data File : BM019822.D  
Acq On : 09 Apr 2019 08:22  
Operator : JU/SJ  
Sample : K2188-10  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF654

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

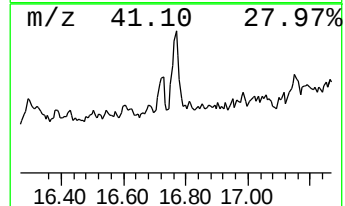
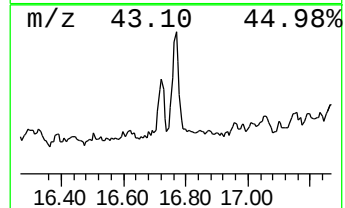
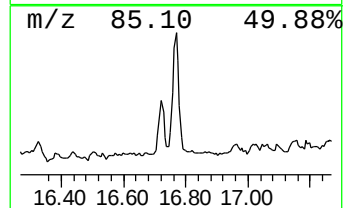
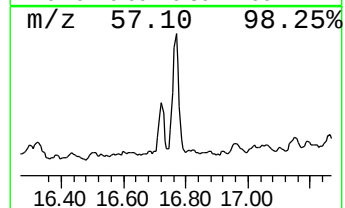
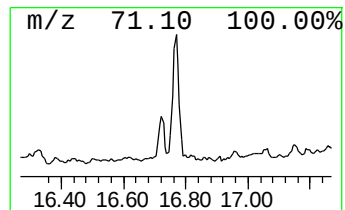
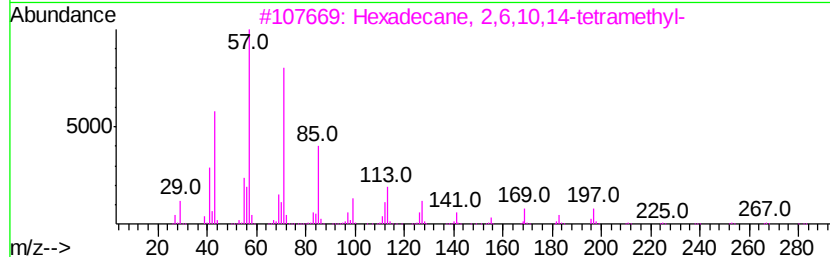
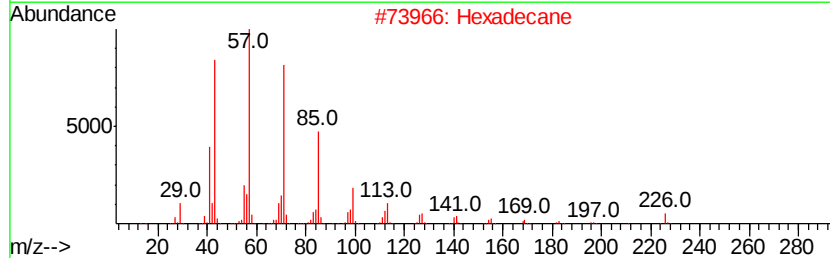
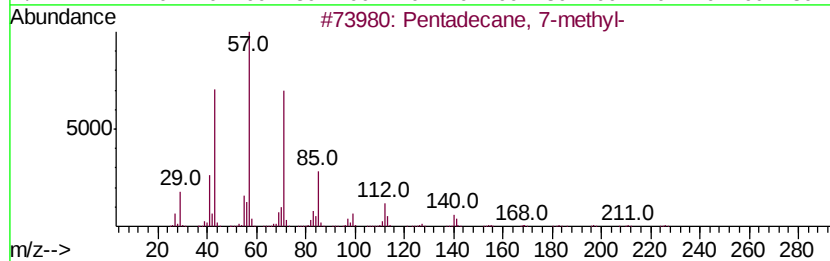
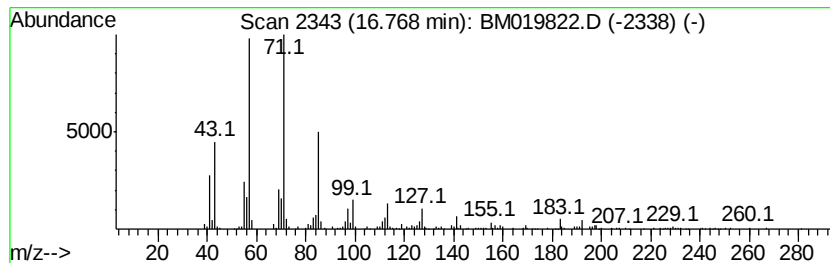
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 46 (DEL) Alkane: Straight-Chai... Concentration Rank 24

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 16.77 | 4.34 ng/ul | 423403 | Phenanthrene-d10 | 16.99 |

| Hit# | of | 5 | Tentative ID                       | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Pentadecane, 7-methyl-             | 226 | C16H34  | 006165-40-8 | 90   |
| 2    |    |   | Hexadecane                         | 226 | C16H34  | 000544-76-3 | 87   |
| 3    |    |   | Hexadecane, 2,6,10,14-tetramethyl- | 282 | C20H42  | 000638-36-8 | 86   |
| 4    |    |   | Octacosane                         | 394 | C28H58  | 000630-02-4 | 80   |
| 5    |    |   | Hexadecane, 7-methyl-              | 240 | C17H36  | 026730-20-1 | 80   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040819\  
Data File : BM019822.D  
Acq On : 09 Apr 2019 08:22  
Operator : JU/SJ  
Sample : K2188-10  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF654

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

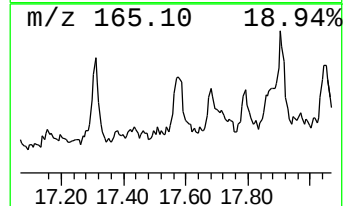
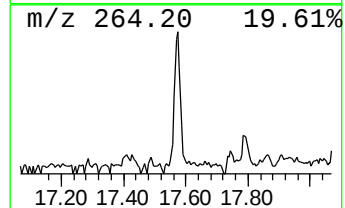
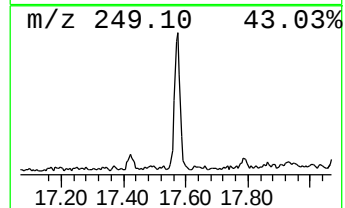
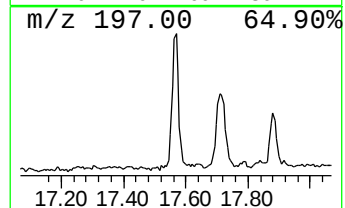
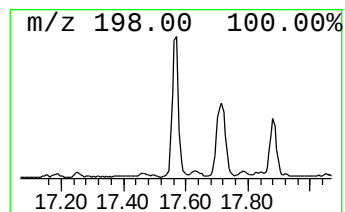
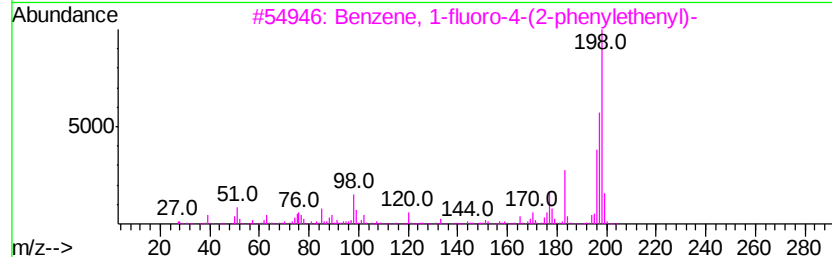
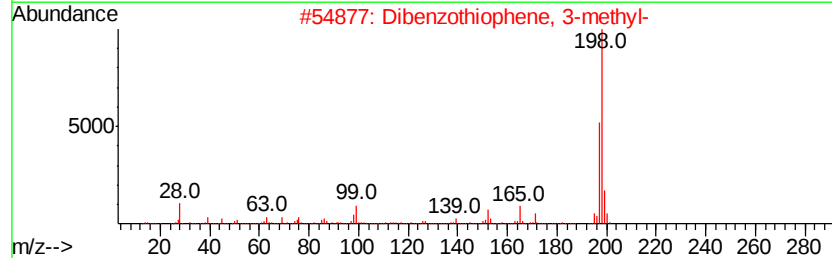
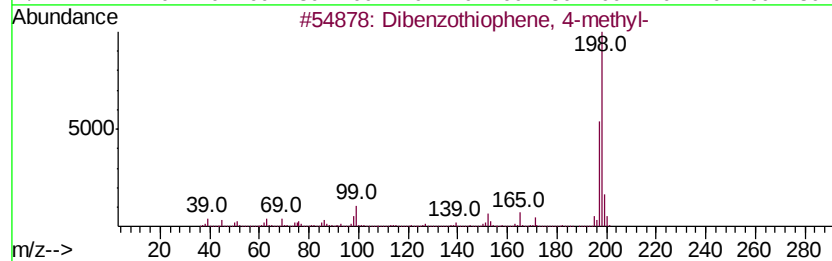
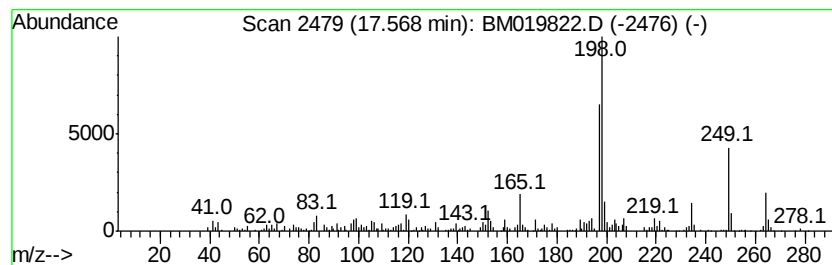
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 47 Dibenzothiophene, 4-methyl- Concentration Rank 37

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 17.57 | 2.77 ng/ul | 270411 | Phenanthrene-d10 | 16.99 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | Dibenzothiophene, 4-methyl-         | 198 | C13H10S  | 007372-88-5 | 55   |
| 2    |    | Dibenzothiophene, 3-methyl-         | 198 | C13H10S  | 016587-52-3 | 55   |
| 3    |    | Benzene, 1-fluoro-4-(2-phenyleth... | 198 | C14H11F  | 017404-69-2 | 47   |
| 4    |    | Benzenamine, 4,4'-methylenebis-     | 198 | C13H14N2 | 000101-77-9 | 46   |
| 5    |    | 9-Acridinamine, 1,2,3,4-tetrahydro- | 198 | C13H14N2 | 000321-64-2 | 43   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040819\  
Data File : BM019822.D  
Acq On : 09 Apr 2019 08:22  
Operator : JU/SJ  
Sample : K2188-10  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF654

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

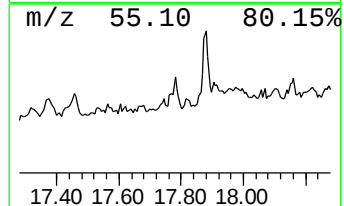
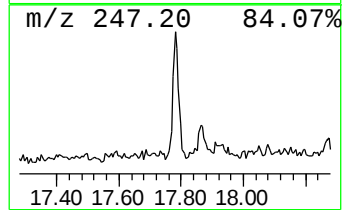
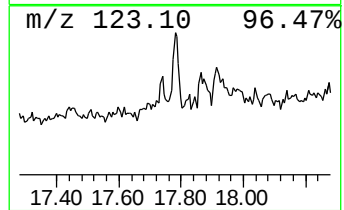
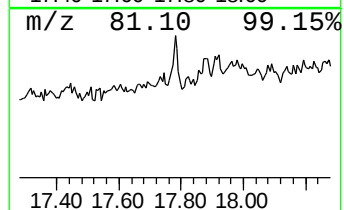
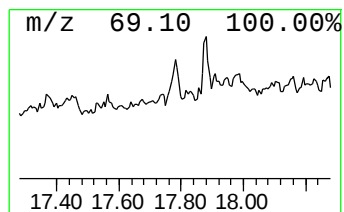
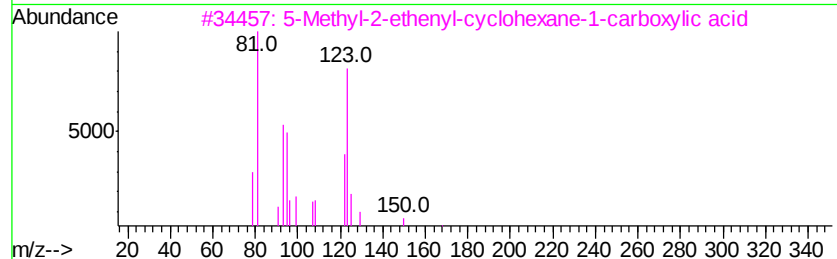
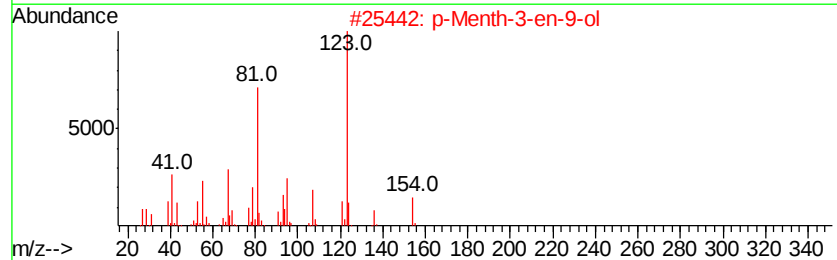
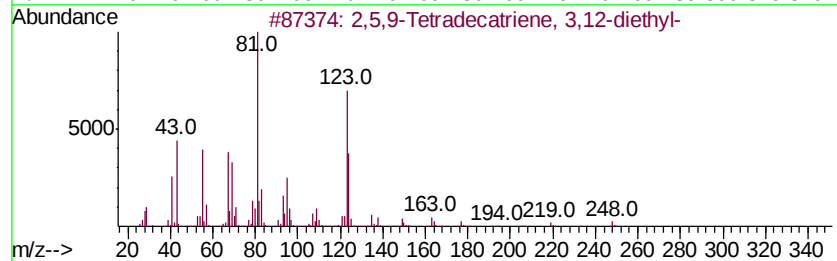
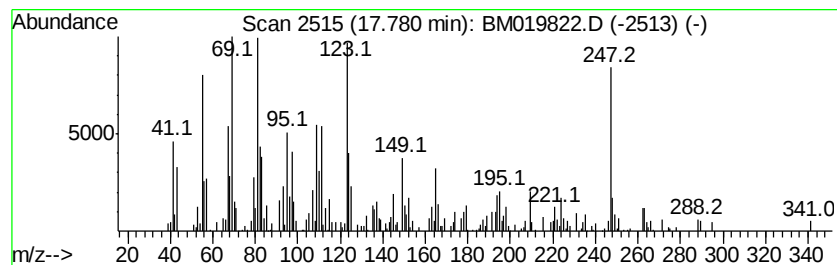
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 48 unknown-06 Concentration Rank 52

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 17.78 | 2.00 ng/ul | 195373 | Phenanthrene-d10 | 16.99 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|-------------------------------------|-----|----------|--------------|------|
| 1    | 5  | 2,5,9-Tetradecatriene, 3,12-diet... | 248 | C18H32   | 074685-87-3  | 38   |
| 2    |    | p-Menth-3-en-9-ol                   | 154 | C10H18O  | 015714-10-0  | 38   |
| 3    |    | 5-Methyl-2-ethenyl-cyclohexane-1... | 168 | C10H16O2 | 1000144-53-6 | 38   |
| 4    |    | (1aR,3S,3aS,7R,7aR,7bS)-(-)-1a,2... | 222 | C15H26O  | 028329-86-4  | 35   |
| 5    |    | 2-Cyclohexene-1-carboxaldehyde, ... | 152 | C10H16O  | 000432-24-6  | 30   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040819\  
Data File : BM019822.D  
Acq On : 09 Apr 2019 08:22  
Operator : JU/SJ  
Sample : K2188-10  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF654

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

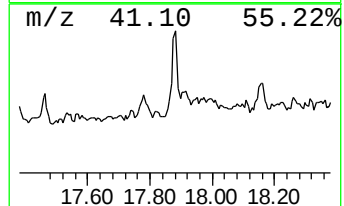
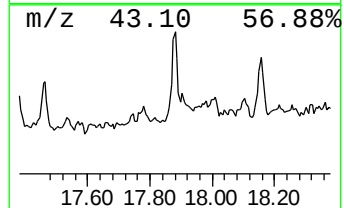
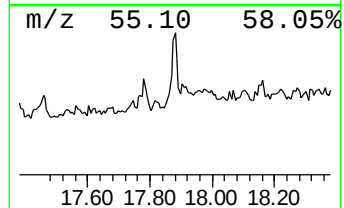
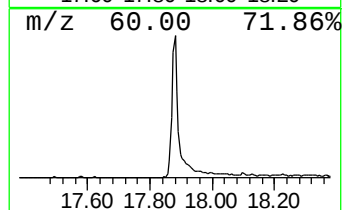
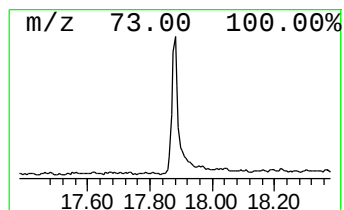
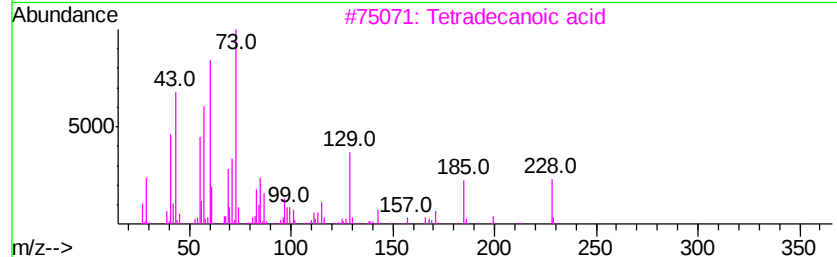
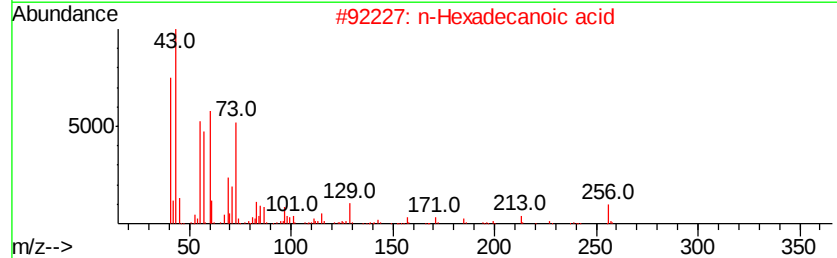
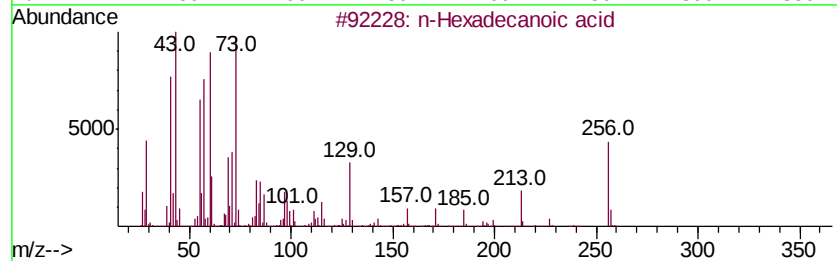
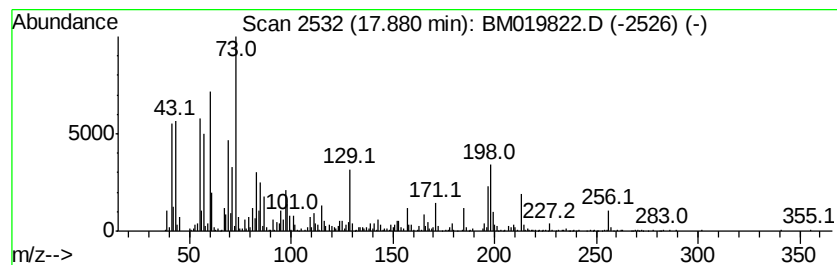
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 49 n-Hexadecanoic acid Concentration Rank 4

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 17.88 | 8.42 ng/ul | 822302 | Phenanthrene-d10 | 16.99 |

| Hit# | of | Tentative ID        | MW  | MolForm  | CAS#        | Qual |
|------|----|---------------------|-----|----------|-------------|------|
| 1    | 5  | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 98   |
| 2    |    | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 94   |
| 3    |    | Tetradecanoic acid  | 228 | C14H28O2 | 000544-63-8 | 91   |
| 4    |    | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 86   |
| 5    |    | Pentadecanoic acid  | 242 | C15H30O2 | 001002-84-2 | 50   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040819\  
Data File : BM019822.D  
Acq On : 09 Apr 2019 08:22  
Operator : JU/SJ  
Sample : K2188-10  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF654

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

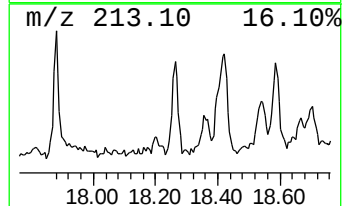
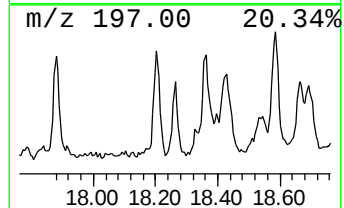
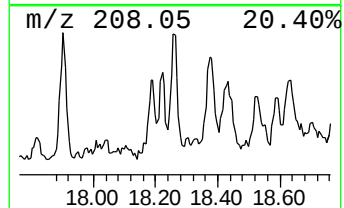
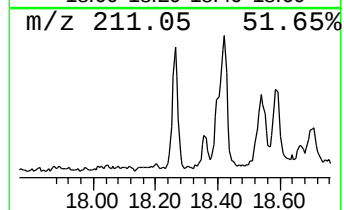
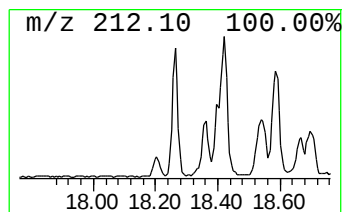
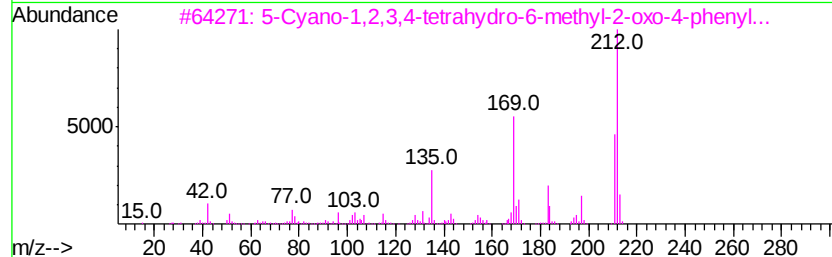
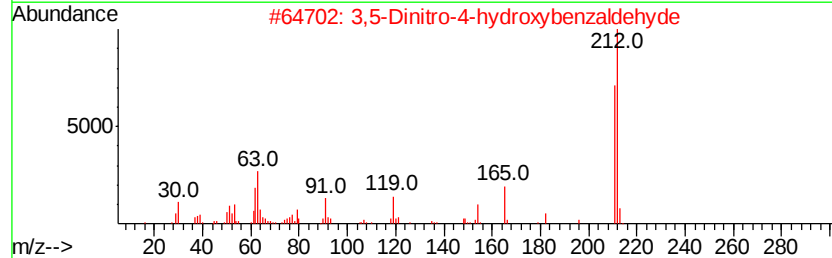
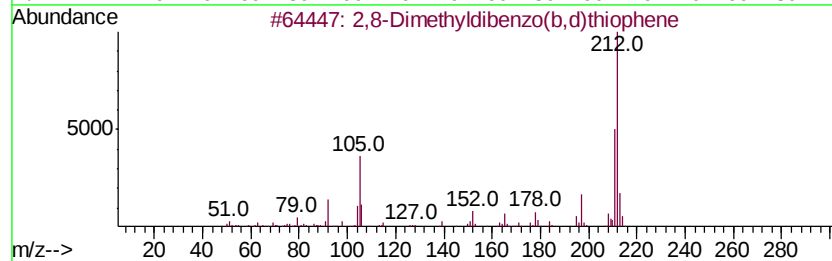
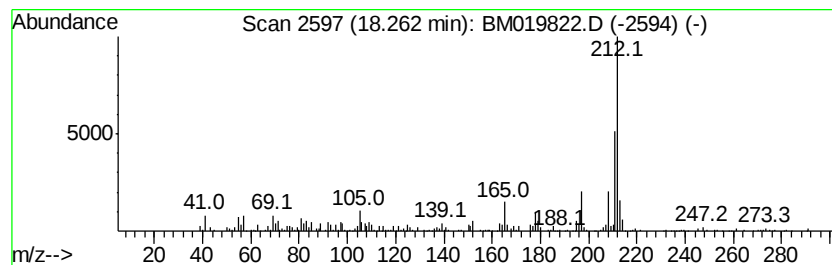
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 50 2,8-Dimethyldibenzo(b,d)thi... Concentration Rank 49

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 18.26 | 2.25 ng/ul | 219667 | Phenanthrene-d10 | 16.99 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#        | Qual |
|------|----|-------------------------------------|-----|-----------|-------------|------|
| 1    | 5  | 2,8-Dimethyldibenzo(b,d)thiophene   | 212 | C14H12S   | 001207-15-4 | 91   |
| 2    |    | 3,5-Dinitro-4-hydroxybenzaldehyde   | 212 | C7H4N2O6  | 052132-61-3 | 50   |
| 3    |    | 5-Cyano-1,2,3,4-tetrahydro-6-met... | 212 | C13H12N2O | 027036-92-6 | 45   |
| 4    |    | Benzo[c]2,7-naphthiridine-4,5(3H... | 212 | C12H8N2O2 | 174565-23-2 | 43   |
| 5    |    | Harmine                             | 212 | C13H12N2O | 000442-51-3 | 43   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040819\  
Data File : BM019822.D  
Acq On : 09 Apr 2019 08:22  
Operator : JU/SJ  
Sample : K2188-10  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF654

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

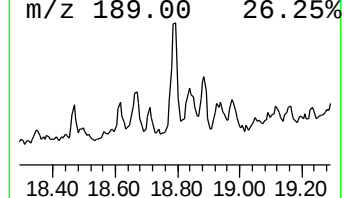
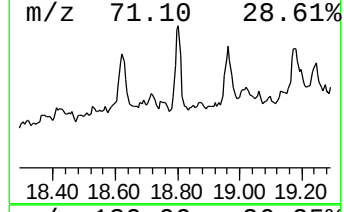
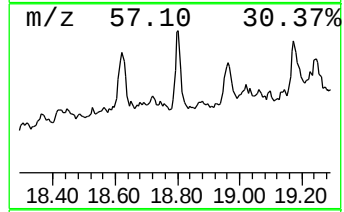
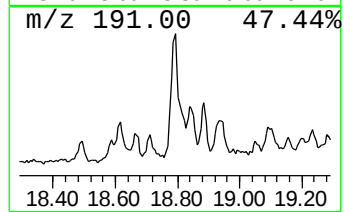
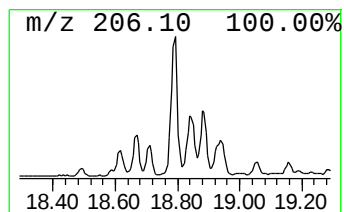
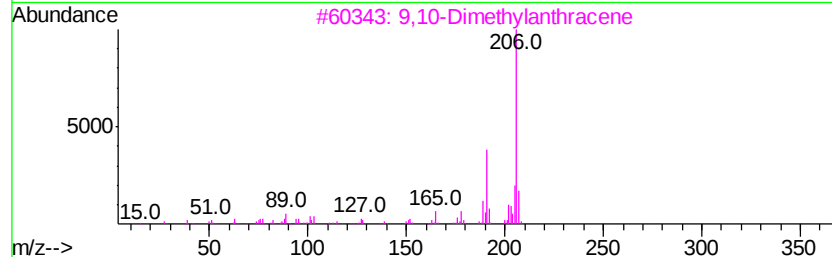
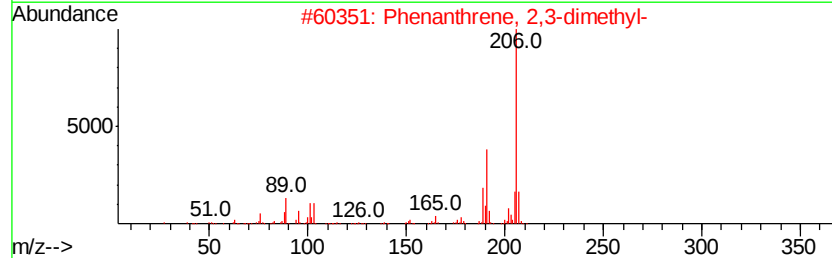
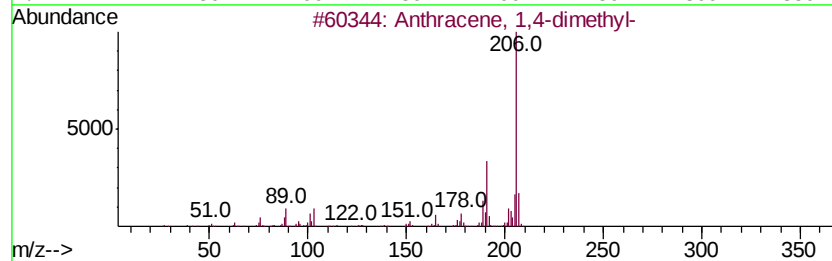
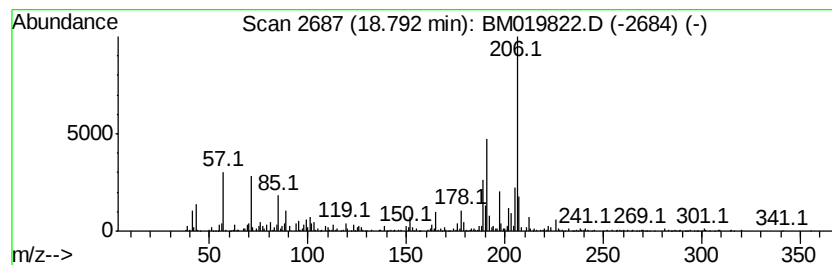
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 51 Anthracene, 1,4-dimethyl- Concentration Rank 7

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 18.79 | 6.67 ng/ul | 651079 | Phenanthrene-d10 | 16.99 |

| Hit# | of | 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------------|-----|---------|-------------|------|
| 1    |    |   | Anthracene, 1,4-dimethyl-   | 206 | C16H14  | 000781-92-0 | 93   |
| 2    |    |   | Phenanthrene, 2,3-dimethyl- | 206 | C16H14  | 003674-65-5 | 90   |
| 3    |    |   | 9,10-Dimethylanthracene     | 206 | C16H14  | 000781-43-1 | 89   |
| 4    |    |   | 9,10-Dimethylanthracene     | 206 | C16H14  | 000781-43-1 | 89   |
| 5    |    |   | di-p-Tolylacetylene         | 206 | C16H14  | 002789-88-0 | 89   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040819\  
Data File : BM019822.D  
Acq On : 09 Apr 2019 08:22  
Operator : JU/SJ  
Sample : K2188-10  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF654

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

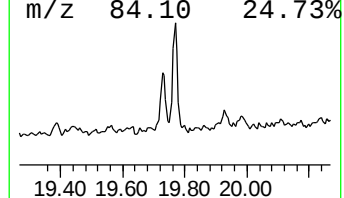
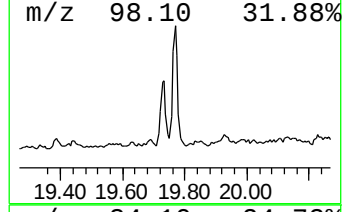
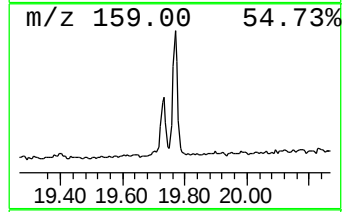
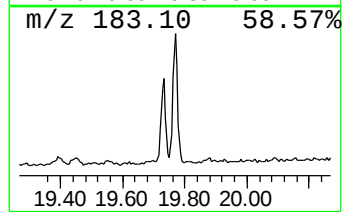
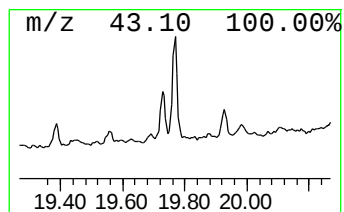
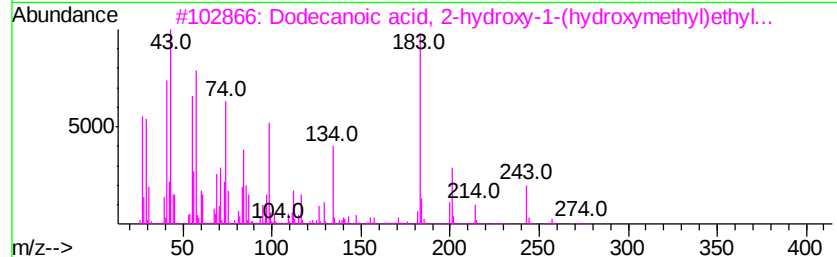
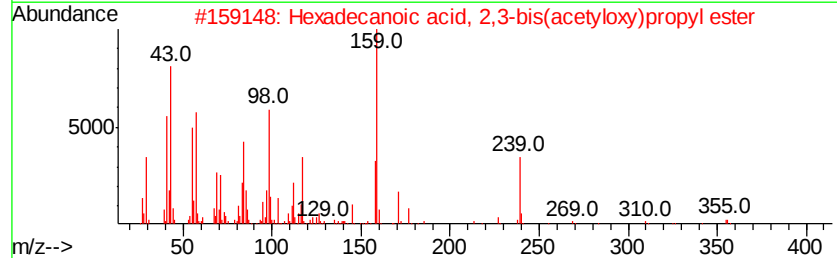
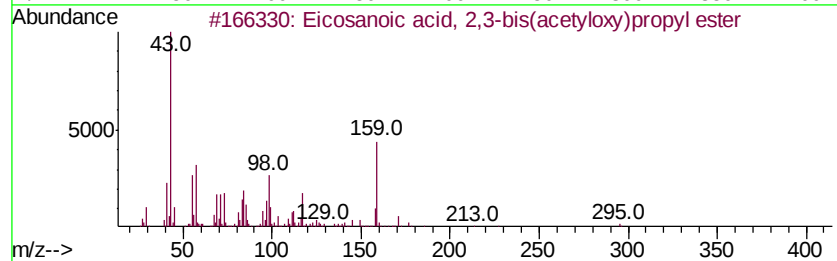
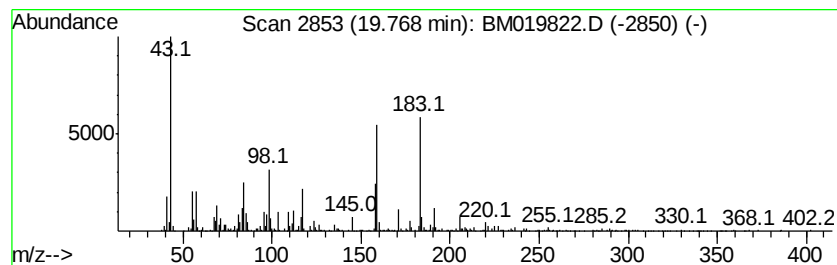
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 52 unknown-07 Concentration Rank 11

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 19.77 | 6.08 ng/ul | 791421 | Chrysene-d12     | 21.20 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|------------|--------------|------|
| 1    |    |   | Eicosanoic acid, 2,3-bis(acetylo... | 470 | C27H50O6   | 055429-67-9  | 49   |
| 2    |    |   | Hexadecanoic acid, 2,3-bis(acety... | 414 | C23H42O6   | 055268-70-7  | 49   |
| 3    |    |   | Dodecanoic acid, 2-hydroxy-1-(hv... | 274 | C15H30O4   | 001678-45-1  | 25   |
| 4    |    |   | Dodecanoyl chloride                 | 218 | C12H23ClO  | 000112-16-3  | 25   |
| 5    |    |   | 2,3,4-Trifluorobenzoic acid, 6-e... | 316 | C17H23F3O2 | 1000282-58-3 | 14   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_M\DATA\BM040819\  
 Data File : BM019822.D  
 Acq On : 09 Apr 2019 08:22  
 Operator : JU/SJ  
 Sample : K2188-10  
 Misc :  
 ALS Vial : 35 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 BF654

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
 Quant Title : SVOA CALIBRATION

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

| TIC Top Hit name     | RT    | EstConc | Units | Response | --Internal Standard-- |       |         |      |
|----------------------|-------|---------|-------|----------|-----------------------|-------|---------|------|
|                      |       |         |       |          | #                     | RT    | Resp    | Conc |
| (DEL) Alkane: Cyc... | 4.77  | 3.8     | ng/ul | 236946   | 1                     | 7.56  | 1261340 | 20.0 |
| (DEL) Alkane: Cyc... | 5.43  | 2.6     | ng/ul | 162020   | 1                     | 7.56  | 1261340 | 20.0 |
| (DEL) Alkane: Cyc... | 5.57  | 2.7     | ng/ul | 172057   | 1                     | 7.56  | 1261340 | 20.0 |
| (DEL) Alkane: Cyc... | 5.81  | 3.7     | ng/ul | 235319   | 1                     | 7.56  | 1261340 | 20.0 |
| unknown-01           | 6.05  | 2.8     | ng/ul | 175337   | 1                     | 7.56  | 1261340 | 20.0 |
| (DEL) Alkane: Str... | 6.09  | 2.7     | ng/ul | 172462   | 1                     | 7.56  | 1261340 | 20.0 |
| (DEL) Alkane: Cyc... | 6.17  | 5.2     | ng/ul | 327850   | 1                     | 7.56  | 1261340 | 20.0 |
| Benzene, 1-ethyl-... | 6.63  | 2.5     | ng/ul | 154999   | 1                     | 7.56  | 1261340 | 20.0 |
| (DEL) Alkane: Cyc... | 6.67  | 10.2    | ng/ul | 645448   | 1                     | 7.56  | 1261340 | 20.0 |
| ((5-Isopropyl-2-m... | 7.02  | 3.2     | ng/ul | 200907   | 1                     | 7.56  | 1261340 | 20.0 |
| Benzene, 1,2,3-tr... | 7.19  | 4.8     | ng/ul | 305984   | 1                     | 7.56  | 1261340 | 20.0 |
| (DEL) Alkane: Cyc... | 7.34  | 2.7     | ng/ul | 172068   | 1                     | 7.56  | 1261340 | 20.0 |
| (DEL) Alkane: Cyc... | 7.44  | 6.0     | ng/ul | 379305   | 1                     | 7.56  | 1261340 | 20.0 |
| (DEL) Alkane: Str... | 7.49  | 5.9     | ng/ul | 372709   | 1                     | 7.56  | 1261340 | 20.0 |
| Benzene, 1,2,4-tr... | 7.66  | 5.8     | ng/ul | 369055   | 1                     | 7.56  | 1261340 | 20.0 |
| (DEL) Alkane: Str... | 7.77  | 3.0     | ng/ul | 189494   | 1                     | 7.56  | 1261340 | 20.0 |
| (DEL) Alkane: Cyc... | 7.80  | 3.2     | ng/ul | 201806   | 1                     | 7.56  | 1261340 | 20.0 |
| (DEL) Alkane: Cyc... | 8.07  | 14.2    | ng/ul | 892147   | 1                     | 7.56  | 1261340 | 20.0 |
| Benzene, 1-ethyl-... | 8.17  | 5.8     | ng/ul | 364512   | 1                     | 7.56  | 1261340 | 20.0 |
| Naphthalene, deca... | 8.35  | 2.6     | ng/ul | 166962   | 1                     | 7.56  | 1261340 | 20.0 |
| Benzene, 1-methyl... | 8.49  | 4.5     | ng/ul | 286032   | 1                     | 7.56  | 1261340 | 20.0 |
| Benzene, 1-methyl... | 8.54  | 2.3     | ng/ul | 142613   | 1                     | 7.56  | 1261340 | 20.0 |
| (DEL) Alkane: Cyc... | 8.62  | 11.4    | ng/ul | 717886   | 1                     | 7.56  | 1261340 | 20.0 |
| unknown-02           | 8.82  | 2.7     | ng/ul | 168649   | 1                     | 7.56  | 1261340 | 20.0 |
| (DEL) Alkane: Cyc... | 8.87  | 6.7     | ng/ul | 423394   | 1                     | 7.56  | 1261340 | 20.0 |
| Benzene, 1-ethyl-... | 8.97  | 4.4     | ng/ul | 285998   | 2                     | 10.34 | 1310540 | 20.0 |
| unknown-03           | 9.04  | 2.7     | ng/ul | 175368   | 2                     | 10.34 | 1310540 | 20.0 |
| unknown-04           | 9.15  | 3.3     | ng/ul | 218025   | 2                     | 10.34 | 1310540 | 20.0 |
| Naphthalene, deca... | 9.23  | 5.6     | ng/ul | 369480   | 2                     | 10.34 | 1310540 | 20.0 |
| (DEL) Alkane: Cyc... | 9.34  | 3.0     | ng/ul | 194001   | 2                     | 10.34 | 1310540 | 20.0 |
| Benzene, 1-methyl... | 9.53  | 3.3     | ng/ul | 214966   | 2                     | 10.34 | 1310540 | 20.0 |
| Benzene, 1-methyl... | 9.73  | 6.9     | ng/ul | 449732   | 2                     | 10.34 | 1310540 | 20.0 |
| (DEL) Alkane: Cyc... | 9.90  | 5.5     | ng/ul | 359589   | 2                     | 10.34 | 1310540 | 20.0 |
| (DEL) Alkane: Str... | 10.27 | 6.3     | ng/ul | 413734   | 2                     | 10.34 | 1310540 | 20.0 |
| (DEL) Alkane: Str... | 10.46 | 6.6     | ng/ul | 434620   | 2                     | 10.34 | 1310540 | 20.0 |
| (DEL) Alkane: Cyc... | 10.79 | 5.5     | ng/ul | 357567   | 2                     | 10.34 | 1310540 | 20.0 |
| (DEL) Alkane: Str... | 11.22 | 2.8     | ng/ul | 180213   | 2                     | 10.34 | 1310540 | 20.0 |
| unknown-05           | 11.49 | 2.2     | ng/ul | 141585   | 2                     | 10.34 | 1310540 | 20.0 |
| Isotridecanol-       | 11.58 | 3.9     | ng/ul | 253854   | 2                     | 10.34 | 1310540 | 20.0 |
| (DEL) Alkane: Str... | 11.73 | 4.4     | ng/ul | 288242   | 2                     | 10.34 | 1310540 | 20.0 |
| Naphthalene, 1-me... | 12.23 | 2.9     | ng/ul | 191295   | 2                     | 10.34 | 1310540 | 20.0 |
| (DEL) Alkane: Str... | 12.71 | 2.9     | ng/ul | 255144   | 3                     | 14.22 | 1777270 | 20.0 |
| Naphthalene, 1,6-... | 13.40 | 2.5     | ng/ul | 226570   | 3                     | 14.22 | 1777270 | 20.0 |
| (DEL) Alkane: Str... | 15.46 | 2.2     | ng/ul | 193697   | 3                     | 14.22 | 1777270 | 20.0 |
| (DEL) Alkane: Str... | 15.94 | 6.3     | ng/ul | 613699   | 4                     | 16.99 | 1953320 | 20.0 |
| (DEL) Alkane: Str... | 16.77 | 4.3     | ng/ul | 423403   | 4                     | 16.99 | 1953320 | 20.0 |
| Dibenzothiophene,... | 17.57 | 2.8     | ng/ul | 270411   | 4                     | 16.99 | 1953320 | 20.0 |
| unknown-06           | 17.78 | 2.0     | ng/ul | 195373   | 4                     | 16.99 | 1953320 | 20.0 |
| n-Hexadecanoic acid  | 17.88 | 8.4     | ng/ul | 822302   | 4                     | 16.99 | 1953320 | 20.0 |

|                      |       |           |        |   |       |         |      |
|----------------------|-------|-----------|--------|---|-------|---------|------|
| 2,8-Dimethyldiben... | 18.26 | 2.3 ng/ul | 219667 | 4 | 16.99 | 1953320 | 20.0 |
| Anthracene, 1,4-d... | 18.79 | 6.7 ng/ul | 651079 | 4 | 16.99 | 1953320 | 20.0 |
| unknown-07           | 19.77 | 6.1 ng/ul | 791421 | 5 | 21.20 | 2602310 | 20.0 |

Instrument :

BNA\_M

ClientSampleId :

BF654