

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP101719\  
 Data File : BP000831.D  
 Acq On : 17 Oct 2019 10:23  
 Operator : JU  
 Sample : K5191-03  
 Misc :  
 ALS Vial : 4 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 BF6L0

## 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\_P\METHODS\SOM-EPA-BP100919MA.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         | 2.364       | 44            | 47          | 60           | rVB      | 30438          | 44272         | 2.25%           | 0.126%        |
| 2         | 3.752       | 273           | 283         | 291          | rVV      | 73349          | 118309        | 6.01%           | 0.336%        |
| 3         | 5.222       | 529           | 533         | 538          | rBV      | 47693          | 47888         | 2.43%           | 0.136%        |
| 4         | 5.328       | 547           | 551         | 563          | rBV      | 165583         | 219578        | 11.15%          | 0.623%        |
| 5         | 5.593       | 592           | 596         | 601          | rVV      | 162381         | 156917        | 7.97%           | 0.445%        |
| 6         | 6.211       | 692           | 701         | 705          | rBV      | 26217          | 38172         | 1.94%           | 0.108%        |
| 7         | 6.275       | 709           | 712         | 715          | rBV      | 141413         | 110749        | 5.63%           | 0.314%        |
| 8         | 6.305       | 715           | 717         | 721          | rVB      | 55363          | 46510         | 2.36%           | 0.132%        |
| 9         | 6.352       | 721           | 725         | 728          | rVB      | 142551         | 111676        | 5.67%           | 0.317%        |
| 10        | 6.387       | 728           | 731         | 735          | rBV2     | 175174         | 220575        | 11.21%          | 0.626%        |
| 11        | 6.428       | 735           | 738         | 741          | rBV2     | 560285         | 471427        | 23.95%          | 1.337%        |
| 12        | 6.516       | 750           | 753         | 760          | rVB      | 589917         | 530079        | 26.93%          | 1.504%        |
| 13        | 6.575       | 760           | 763         | 775          | rVB      | 436855         | 377231        | 19.16%          | 1.070%        |
| 14        | 6.746       | 789           | 792         | 796          | rBV      | 1216863        | 995119        | 50.55%          | 2.823%        |
| 15        | 6.811       | 800           | 803         | 811          | rVB      | 215450         | 217471        | 11.05%          | 0.617%        |
| 16        | 6.928       | 819           | 823         | 827          | rVB2     | 129195         | 133368        | 6.78%           | 0.378%        |
| 17        | 7.005       | 832           | 836         | 838          | rBV3     | 36851          | 47384         | 2.41%           | 0.134%        |
| 18        | 7.028       | 838           | 840         | 844          | rVB      | 48909          | 41320         | 2.10%           | 0.117%        |
| 19        | 7.069       | 844           | 847         | 850          | rBV      | 107806         | 93743         | 4.76%           | 0.266%        |
| 20        | 7.134       | 855           | 858         | 869          | rBV2     | 486282         | 553825        | 28.13%          | 1.571%        |
| 21        | 7.216       | 869           | 872         | 874          | rBV      | 55454          | 40026         | 2.03%           | 0.114%        |
| 22        | 7.240       | 874           | 876         | 879          | rVB      | 65652          | 47025         | 2.39%           | 0.133%        |
| 23        | 7.287       | 880           | 884         | 885          | rBV      | 132890         | 119386        | 6.06%           | 0.339%        |
| 24        | 7.305       | 885           | 887         | 892          | rVV2     | 684982         | 594974        | 30.22%          | 1.688%        |
| 25        | 7.352       | 892           | 895         | 899          | rVB3     | 45860          | 62320         | 3.17%           | 0.177%        |
| 26        | 7.434       | 906           | 909         | 912          | rBV2     | 53229          | 54194         | 2.75%           | 0.154%        |
| 27        | 7.522       | 921           | 924         | 926          | rBV      | 106108         | 77346         | 3.93%           | 0.219%        |
| 28        | 7.546       | 926           | 928         | 934          | rVB      | 127295         | 110726        | 5.62%           | 0.314%        |
| 29        | 7.634       | 940           | 943         | 946          | rBV      | 554566         | 444375        | 22.57%          | 1.261%        |
| 30        | 7.693       | 950           | 953         | 954          | rBV2     | 106287         | 97721         | 4.96%           | 0.277%        |
| 31        | 7.705       | 954           | 955         | 961          | rVB      | 151111         | 104704        | 5.32%           | 0.297%        |
| 32        | 7.775       | 961           | 967         | 970          | rBV2     | 304936         | 354621        | 18.01%          | 1.006%        |
| 33        | 7.805       | 970           | 972         | 974          | rVV2     | 67916          | 77559         | 3.94%           | 0.220%        |
| 34        | 7.828       | 974           | 976         | 982          | rVV3     | 109588         | 119718        | 6.08%           | 0.340%        |

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 BNA\_P  
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## Integration Parameters: LSCINT.P

Integrator: RTE  
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 Sampling : 1  
 Start Thrs: 0.2  
 Stop Thrs : 0

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

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

Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP100919MA.M  
 Title : SVOA CALIBRATION

|    |        |      |      |      |      |         |         |        |        |
|----|--------|------|------|------|------|---------|---------|--------|--------|
| 35 | 7.887  | 982  | 986  | 993  | rVV  | 762125  | 782477  | 39.75% | 2.220% |
| 36 | 7.999  | 1002 | 1005 | 1006 | rVV  | 47402   | 45937   | 2.33%  | 0.130% |
| 37 | 8.034  | 1007 | 1011 | 1013 | rVV  | 1574826 | 1504142 | 76.41% | 4.267% |
| 38 | 8.052  | 1013 | 1014 | 1017 | rVV  | 792848  | 554378  | 28.16% | 1.573% |
| 39 | 8.081  | 1017 | 1019 | 1021 | rVV  | 61311   | 51965   | 2.64%  | 0.147% |
| 40 | 8.105  | 1021 | 1023 | 1025 | rVV  | 76495   | 59630   | 3.03%  | 0.169% |
| 41 | 8.193  | 1035 | 1038 | 1043 | rBV  | 279580  | 236910  | 12.03% | 0.672% |
| 42 | 8.269  | 1048 | 1051 | 1055 | rBV2 | 32406   | 39815   | 2.02%  | 0.113% |
| 43 | 8.310  | 1055 | 1058 | 1064 | rVB4 | 50535   | 55657   | 2.83%  | 0.158% |
| 44 | 8.393  | 1069 | 1072 | 1075 | rBV  | 77126   | 83089   | 4.22%  | 0.236% |
| 45 | 8.422  | 1075 | 1077 | 1080 | rVB  | 70827   | 54142   | 2.75%  | 0.154% |
| 46 | 8.475  | 1082 | 1086 | 1088 | rBV2 | 43077   | 51565   | 2.62%  | 0.146% |
| 47 | 8.505  | 1088 | 1091 | 1095 | rVV  | 88934   | 95542   | 4.85%  | 0.271% |
| 48 | 8.540  | 1095 | 1097 | 1102 | rVB4 | 40765   | 38992   | 1.98%  | 0.111% |
| 49 | 8.610  | 1105 | 1109 | 1111 | rBV  | 163691  | 148584  | 7.55%  | 0.422% |
| 50 | 8.628  | 1111 | 1112 | 1115 | rVV2 | 50603   | 43832   | 2.23%  | 0.124% |
| 51 | 8.699  | 1122 | 1124 | 1126 | rVV2 | 34509   | 33733   | 1.71%  | 0.096% |
| 52 | 8.752  | 1130 | 1133 | 1136 | rVV  | 608352  | 488674  | 24.82% | 1.386% |
| 53 | 8.846  | 1145 | 1149 | 1153 | rVB  | 347955  | 326234  | 16.57% | 0.925% |
| 54 | 9.152  | 1197 | 1201 | 1206 | rBV7 | 19715   | 38580   | 1.96%  | 0.109% |
| 55 | 9.222  | 1206 | 1213 | 1217 | rBV2 | 142681  | 195029  | 9.91%  | 0.553% |
| 56 | 9.310  | 1226 | 1228 | 1234 | rVB2 | 36362   | 53242   | 2.70%  | 0.151% |
| 57 | 9.381  | 1235 | 1240 | 1243 | rBV  | 76385   | 107709  | 5.47%  | 0.306% |
| 58 | 9.452  | 1249 | 1252 | 1255 | rBV  | 111522  | 114076  | 5.80%  | 0.324% |
| 59 | 9.487  | 1255 | 1258 | 1260 | rVV  | 1296944 | 1163684 | 59.11% | 3.301% |
| 60 | 9.510  | 1260 | 1262 | 1268 | rVV  | 1368141 | 1110557 | 56.42% | 3.150% |
| 61 | 9.569  | 1270 | 1272 | 1274 | rVV  | 52622   | 46304   | 2.35%  | 0.131% |
| 62 | 9.640  | 1281 | 1284 | 1290 | rVB  | 1698670 | 1443034 | 73.31% | 4.094% |
| 63 | 9.799  | 1307 | 1311 | 1314 | rBV  | 1974817 | 1676387 | 85.16% | 4.756% |
| 64 | 9.828  | 1314 | 1316 | 1320 | rVB  | 257678  | 198657  | 10.09% | 0.564% |
| 65 | 9.922  | 1329 | 1332 | 1333 | rBV  | 70835   | 57313   | 2.91%  | 0.163% |
| 66 | 10.040 | 1350 | 1352 | 1355 | rVB  | 46189   | 37499   | 1.90%  | 0.106% |
| 67 | 10.210 | 1377 | 1381 | 1383 | rBV3 | 48390   | 57964   | 2.94%  | 0.164% |
| 68 | 10.322 | 1396 | 1400 | 1403 | rVV  | 1993012 | 1794479 | 91.16% | 5.091% |
| 69 | 10.346 | 1403 | 1404 | 1407 | rVV  | 73818   | 54435   | 2.77%  | 0.154% |
| 70 | 10.393 | 1409 | 1412 | 1417 | rVV  | 361449  | 331045  | 16.82% | 0.939% |
| 71 | 10.434 | 1417 | 1419 | 1422 | rVV3 | 38675   | 46829   | 2.38%  | 0.133% |

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Instrument :  
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ClientSampleId :  
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## 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\_P\METHODS\SOM-EPA-BP100919MA.M  
Title : SVOA CALIBRATION

|     |        |      |      |      |      |         |         |         |        |
|-----|--------|------|------|------|------|---------|---------|---------|--------|
| 72  | 10.463 | 1422 | 1424 | 1428 | rVB3 | 46641   | 39379   | 2.00%   | 0.112% |
| 73  | 10.716 | 1465 | 1467 | 1475 | rVB4 | 60456   | 116852  | 5.94%   | 0.331% |
| 74  | 10.899 | 1495 | 1498 | 1501 | rBV4 | 40101   | 54247   | 2.76%   | 0.154% |
| 75  | 11.299 | 1562 | 1566 | 1568 | rBV  | 1889923 | 1663334 | 84.50%  | 4.719% |
| 76  | 11.316 | 1568 | 1569 | 1572 | rVV  | 126038  | 101373  | 5.15%   | 0.288% |
| 77  | 11.351 | 1572 | 1575 | 1582 | rVB  | 2060496 | 1773759 | 90.11%  | 5.032% |
| 78  | 11.410 | 1583 | 1585 | 1588 | rBV2 | 44032   | 45577   | 2.32%   | 0.129% |
| 79  | 11.804 | 1650 | 1652 | 1655 | rVB2 | 47726   | 35001   | 1.78%   | 0.099% |
| 80  | 11.846 | 1655 | 1659 | 1667 | rBV2 | 678723  | 793541  | 40.31%  | 2.251% |
| 81  | 11.916 | 1668 | 1671 | 1674 | rBV  | 125564  | 122233  | 6.21%   | 0.347% |
| 82  | 12.363 | 1745 | 1747 | 1750 | rBV  | 60215   | 52387   | 2.66%   | 0.149% |
| 83  | 12.522 | 1772 | 1774 | 1777 | rVB  | 138834  | 117038  | 5.95%   | 0.332% |
| 84  | 12.640 | 1791 | 1794 | 1801 | rVB2 | 138001  | 151974  | 7.72%   | 0.431% |
| 85  | 12.740 | 1807 | 1811 | 1816 | rBV  | 2063855 | 1968514 | 100.00% | 5.584% |
| 86  | 12.787 | 1816 | 1819 | 1821 | rBV2 | 73747   | 66057   | 3.36%   | 0.187% |
| 87  | 13.145 | 1878 | 1880 | 1884 | rVB2 | 78925   | 85314   | 4.33%   | 0.242% |
| 88  | 13.493 | 1937 | 1939 | 1942 | rVB2 | 119518  | 102641  | 5.21%   | 0.291% |
| 89  | 13.822 | 1992 | 1995 | 2005 | rVB  | 630636  | 826326  | 41.98%  | 2.344% |
| 90  | 13.975 | 2017 | 2021 | 2024 | rBV  | 1722875 | 1600344 | 81.30%  | 4.540% |
| 91  | 14.151 | 2048 | 2051 | 2054 | rVB  | 196449  | 173998  | 8.84%   | 0.494% |
| 92  | 14.457 | 2100 | 2103 | 2108 | rBV  | 275114  | 303438  | 15.41%  | 0.861% |
| 93  | 14.769 | 2153 | 2156 | 2160 | rBV  | 313699  | 299731  | 15.23%  | 0.850% |
| 94  | 14.845 | 2166 | 2169 | 2172 | rVB  | 120865  | 118180  | 6.00%   | 0.335% |
| 95  | 15.110 | 2210 | 2214 | 2219 | rBV  | 317027  | 352200  | 17.89%  | 0.999% |
| 96  | 15.363 | 2252 | 2257 | 2267 | rVB  | 1456965 | 1701359 | 86.43%  | 4.826% |
| 97  | 15.457 | 2268 | 2273 | 2276 | rVV  | 1275439 | 1612178 | 81.90%  | 4.573% |
| 98  | 15.492 | 2276 | 2279 | 2289 | rVB  | 302165  | 407708  | 20.71%  | 1.157% |
| 99  | 15.934 | 2350 | 2354 | 2359 | rBV  | 208377  | 285974  | 14.53%  | 0.811% |
| 100 | 16.439 | 2436 | 2440 | 2444 | rBV  | 106649  | 177654  | 9.02%   | 0.504% |

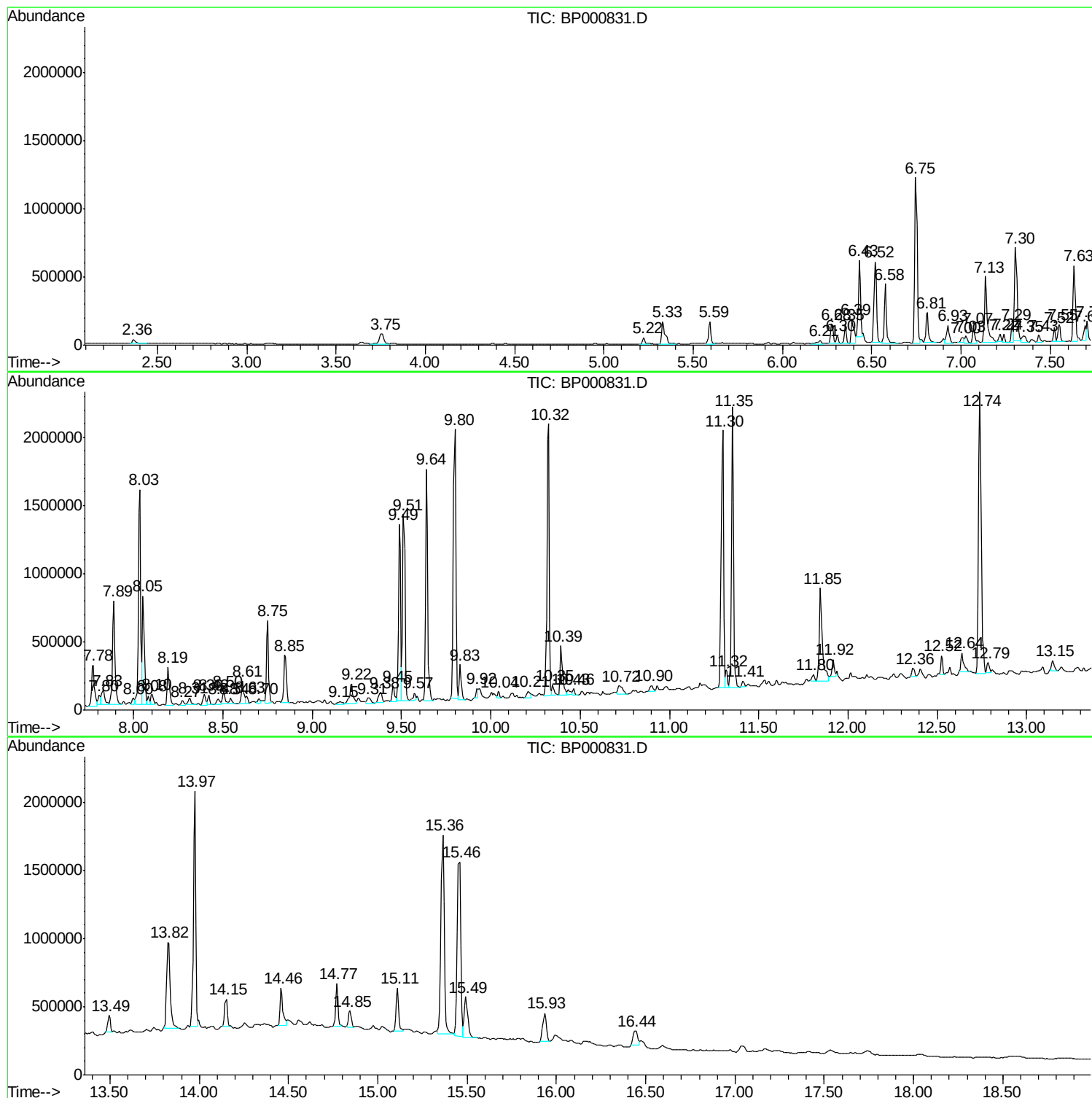
Sum of corrected areas: 35250740

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP100919MA.M  
Quant Title : SVOA CALIBRATION

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



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Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP100919MA.M  
Quant Title : SVOA CALIBRATION

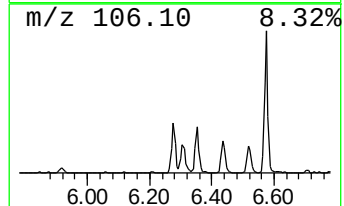
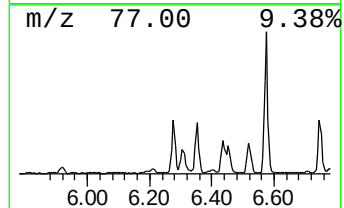
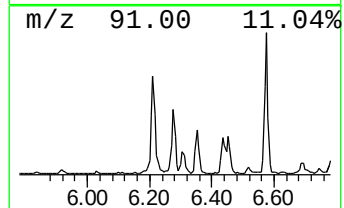
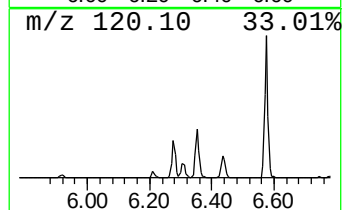
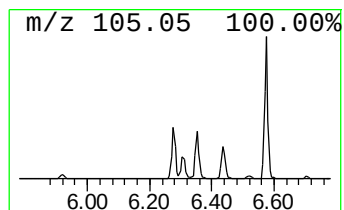
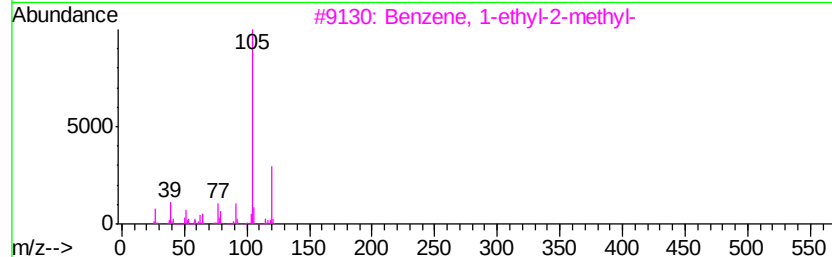
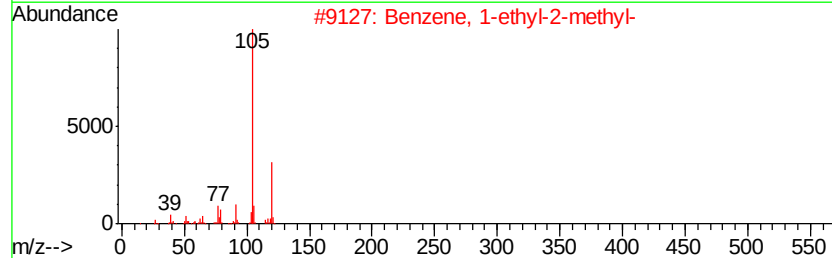
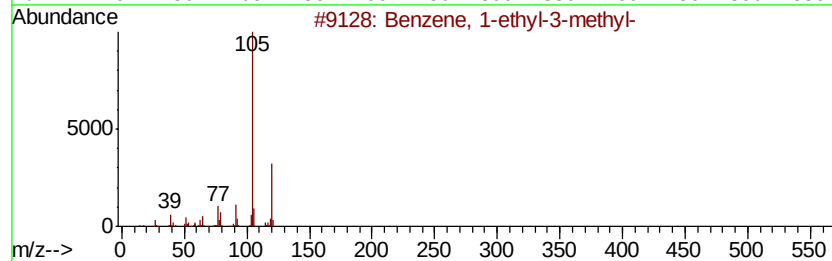
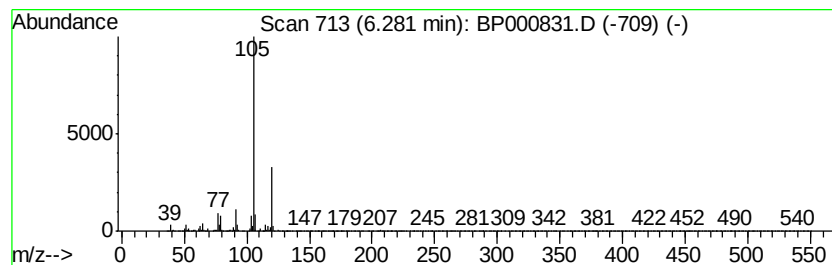
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 4 Benzene, 1-ethyl-3-methyl- Concentration Rank 17

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 6.28 | 2.23 ng/ul | 110749 | 1,4-Dichlorobenzene-d4 | 6.75 |

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



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Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP100919MA.M  
Quant Title : SVOA CALIBRATION

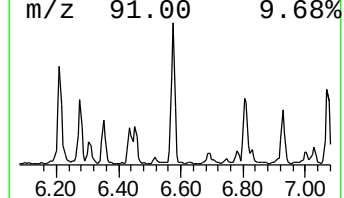
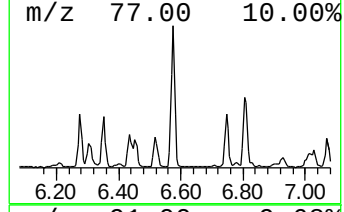
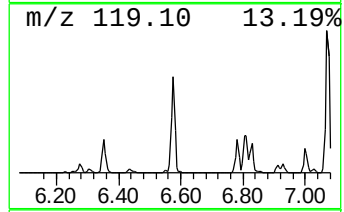
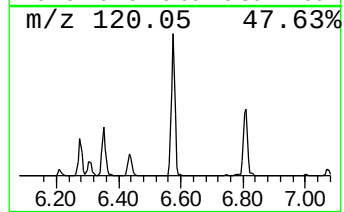
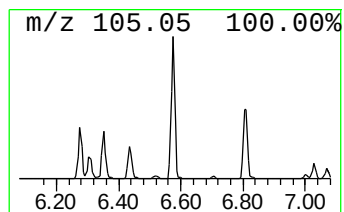
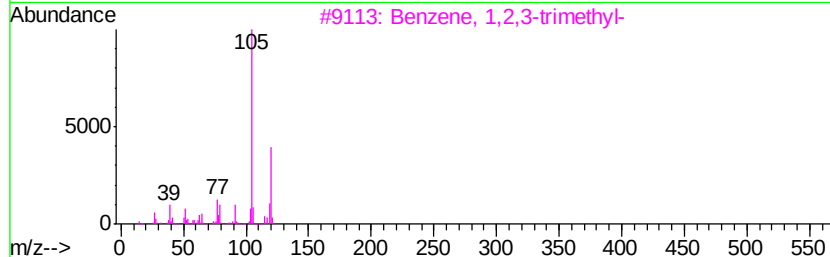
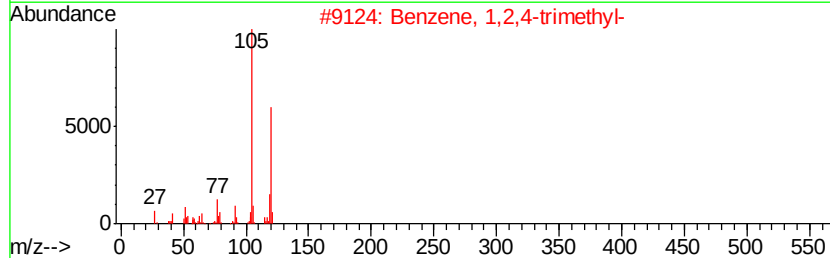
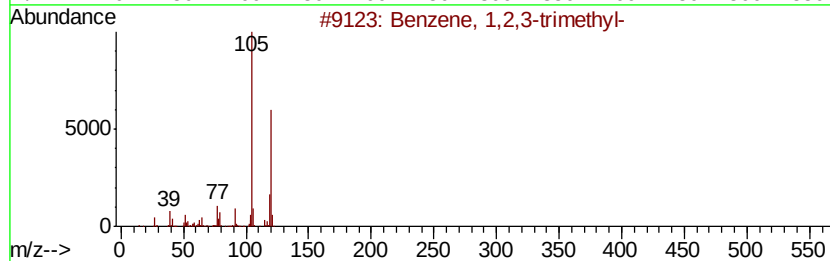
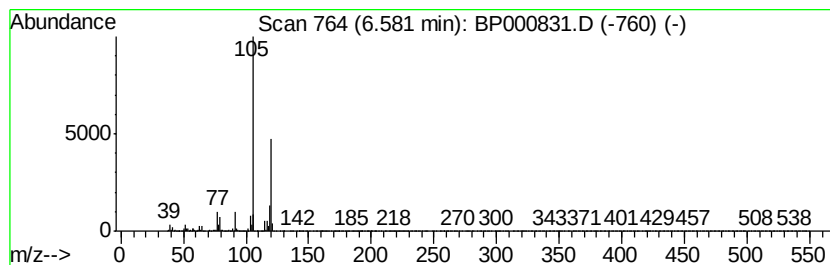
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 5 Benzene, 1,2,3-trimethyl- Concentration Rank 3

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 6.58 | 7.58 ng/ul | 377231 | 1,4-Dichlorobenzene-d4 | 6.75 |

| Hit# | of | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1,2,3-trimethyl- | 120 | C9H12   | 000526-73-8 | 97   |
| 2    |    | Benzene, 1,2,4-trimethyl- | 120 | C9H12   | 000095-63-6 | 95   |
| 3    |    | Benzene, 1,2,3-trimethyl- | 120 | C9H12   | 000526-73-8 | 95   |
| 4    |    | Benzene, 1,3,5-trimethyl- | 120 | C9H12   | 000108-67-8 | 94   |
| 5    |    | Benzene, 1,2,4-trimethyl- | 120 | C9H12   | 000095-63-6 | 94   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP101719\  
Data File : BP000831.D  
Acq On : 17 Oct 2019 10:23  
Operator : JU  
Sample : K5191-03  
Misc :  
ALS Vial : 4 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
BF6L0

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Quant Title : SVOA CALIBRATION

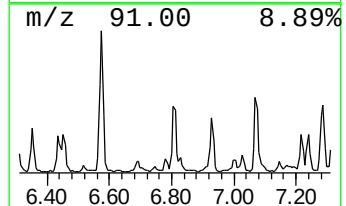
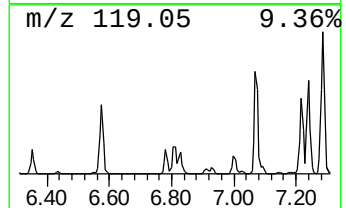
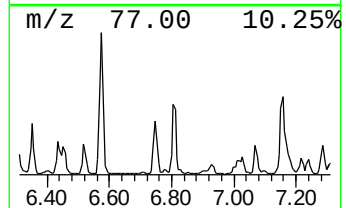
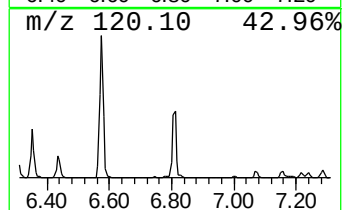
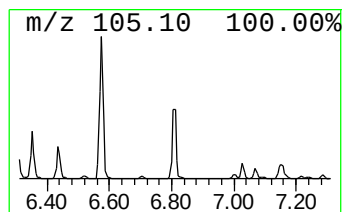
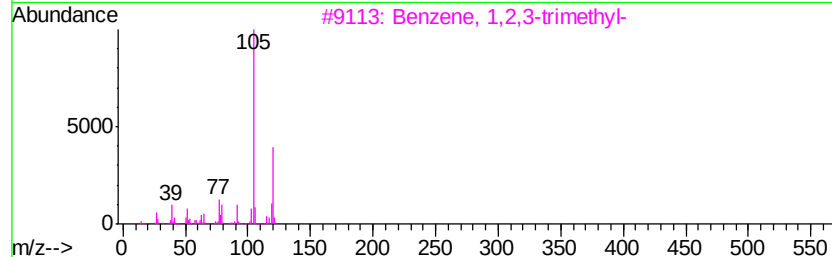
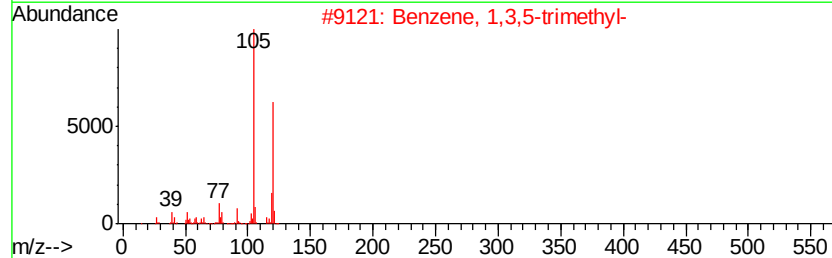
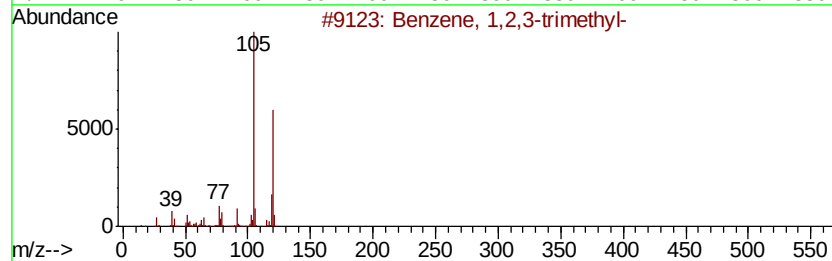
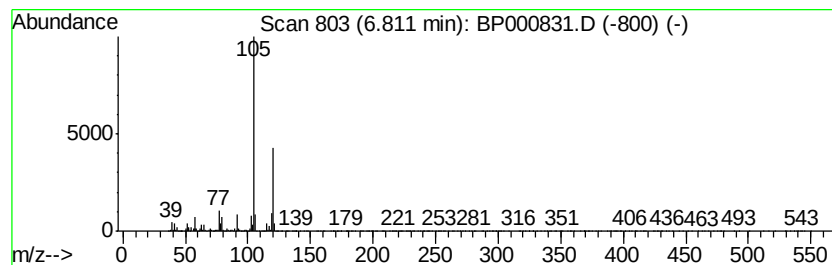
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 6 Benzene, 1,3,5-trimethyl- Concentration Rank 6

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 6.81 | 4.37 ng/ul | 217471 | 1,4-Dichlorobenzene-d4 | 6.75 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1,2,3-trimethyl-  | 120 | C9H12   | 000526-73-8 | 95   |
| 2    |    | Benzene, 1,3,5-trimethyl-  | 120 | C9H12   | 000108-67-8 | 94   |
| 3    |    | Benzene, 1,2,3-trimethyl-  | 120 | C9H12   | 000526-73-8 | 94   |
| 4    |    | Benzene, 1,2,4-trimethyl-  | 120 | C9H12   | 000095-63-6 | 94   |
| 5    |    | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 91   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP101719\  
Data File : BP000831.D  
Acq On : 17 Oct 2019 10:23  
Operator : JU  
Sample : K5191-03  
Misc :  
ALS Vial : 4 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
BF6L0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP100919MA.M  
Quant Title : SVOA CALIBRATION

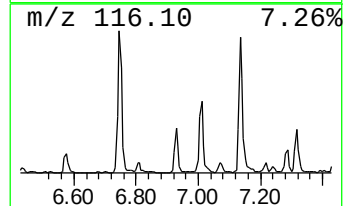
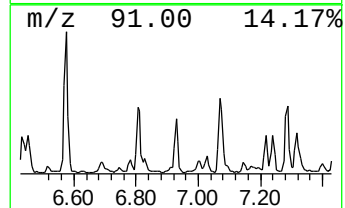
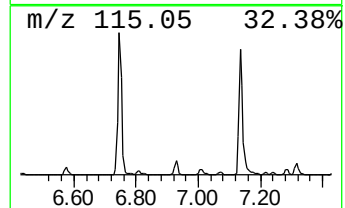
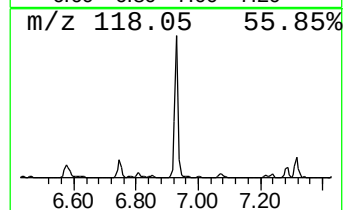
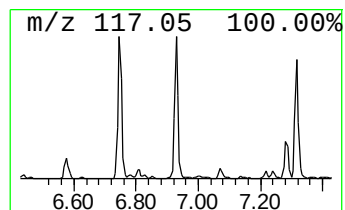
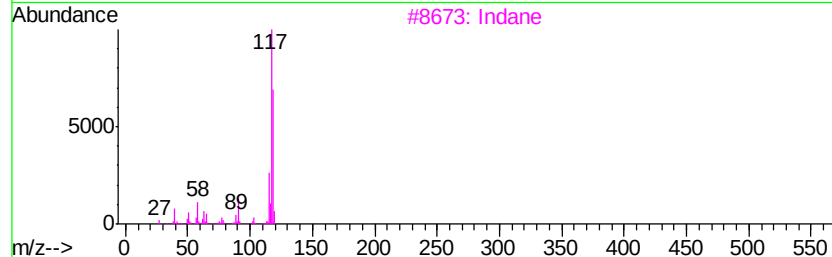
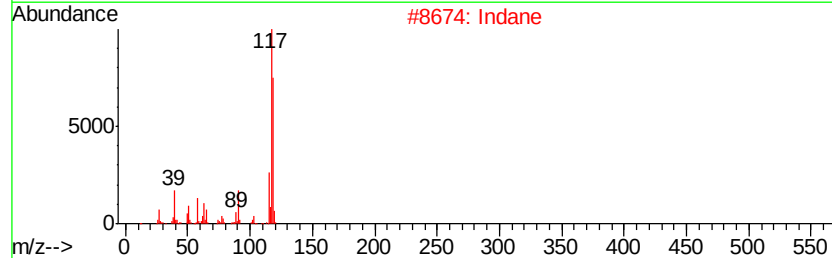
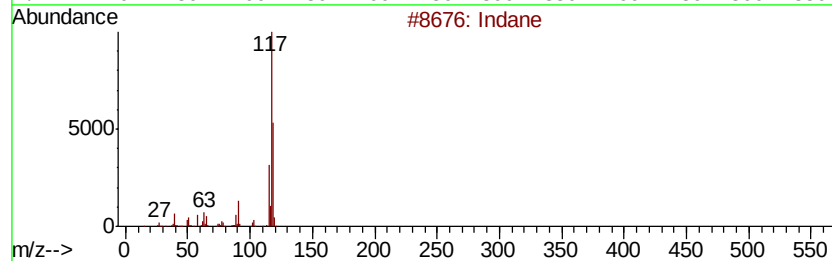
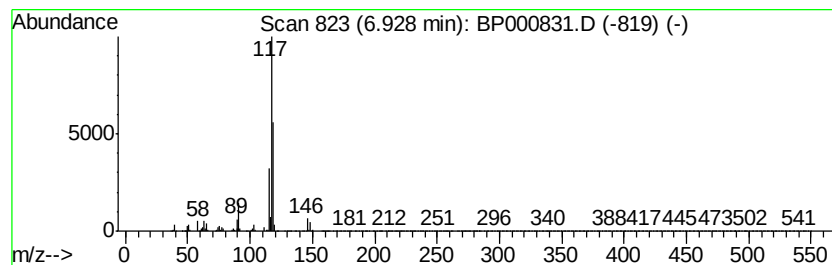
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 7 Indane Concentration Rank 14

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 6.93 | 2.68 ng/ul | 133368 | 1,4-Dichlorobenzene-d4 | 6.75 |

| Hit# of | 5                     | Tentative ID | MW  | MolForm | CAS#        | Qual |
|---------|-----------------------|--------------|-----|---------|-------------|------|
| 1       | Indane                |              | 118 | C9H10   | 000496-11-7 | 93   |
| 2       | Indane                |              | 118 | C9H10   | 000496-11-7 | 74   |
| 3       | Indane                |              | 118 | C9H10   | 000496-11-7 | 72   |
| 4       | Benzene, cyclopropyl- |              | 118 | C9H10   | 000873-49-4 | 59   |
| 5       | Deltacyclene          |              | 118 | C9H10   | 007785-10-6 | 50   |





Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP101719\  
Data File : BP000831.D  
Acq On : 17 Oct 2019 10:23  
Operator : JU  
Sample : K5191-03  
Misc :  
ALS Vial : 4 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
BF6L0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP100919MA.M  
Quant Title : SVOA CALIBRATION

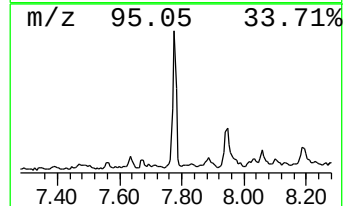
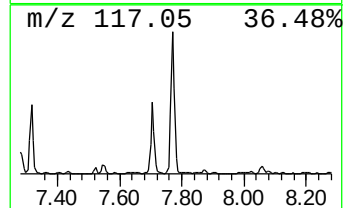
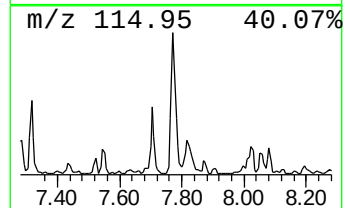
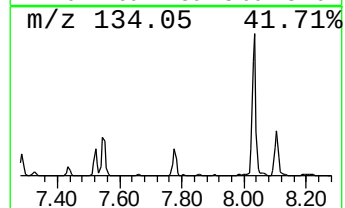
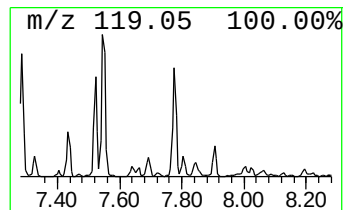
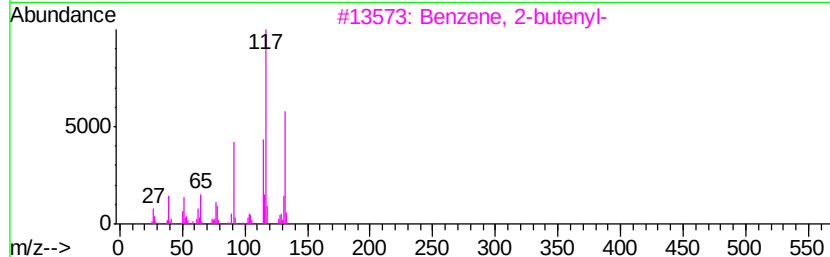
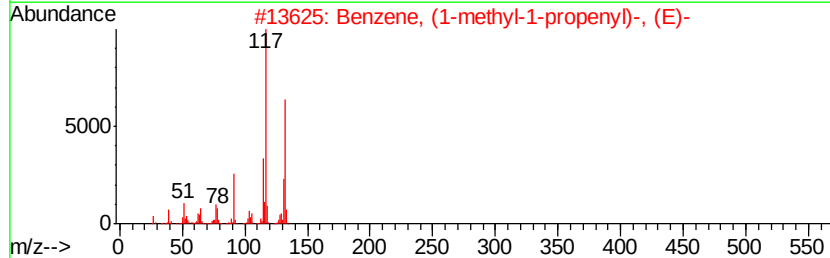
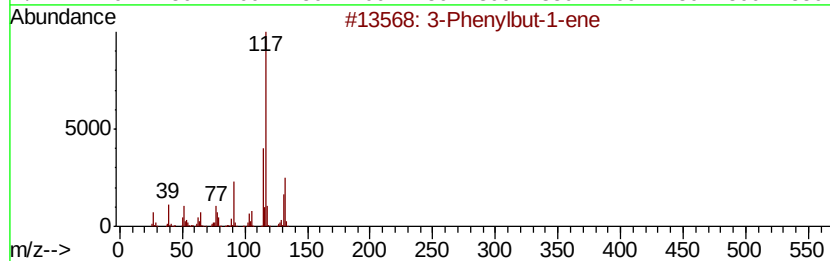
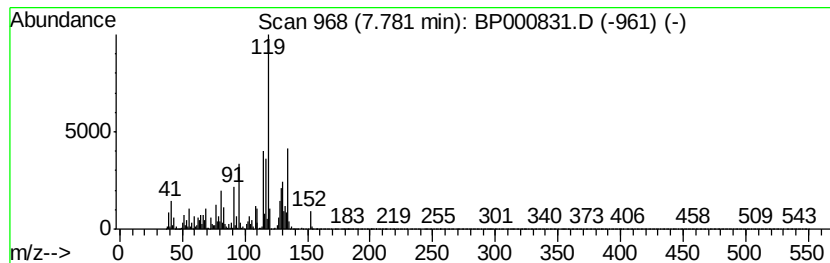
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 8 3-Phenylbut-1-ene Concentration Rank 4

| R.T. | EstConc    | Area   | Relative to ISTD | R.T. |
|------|------------|--------|------------------|------|
| 7.78 | 4.72 ng/ul | 354621 | Napthalene-d8    | 8.03 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | 3-Phenylbut-1-ene                   | 132 | C10H12  | 000934-10-1 | 50   |
| 2    |    | Benzene, (1-methyl-1-propenyl)-,... | 132 | C10H12  | 000768-00-3 | 49   |
| 3    |    | Benzene, 2-butenyl-                 | 132 | C10H12  | 001560-06-1 | 46   |
| 4    |    | Benzene, 1-ethenyl-3-ethyl-         | 132 | C10H12  | 007525-62-4 | 42   |
| 5    |    | Benzene, 1-methyl-2-(1-methyleth... | 134 | C10H14  | 000527-84-4 | 38   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP101719\  
Data File : BP000831.D  
Acq On : 17 Oct 2019 10:23  
Operator : JU  
Sample : K5191-03  
Misc :  
ALS Vial : 4 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP100919MA.M  
Quant Title : SVOA CALIBRATION

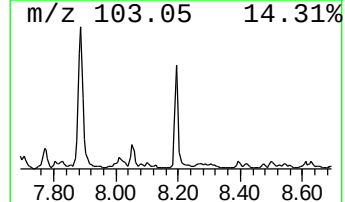
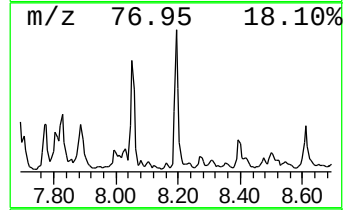
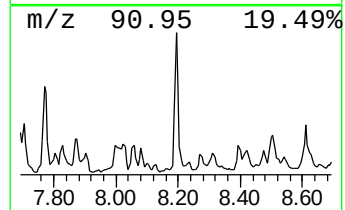
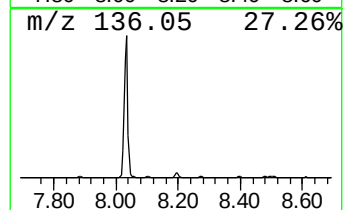
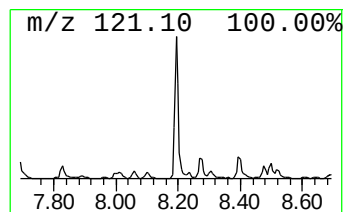
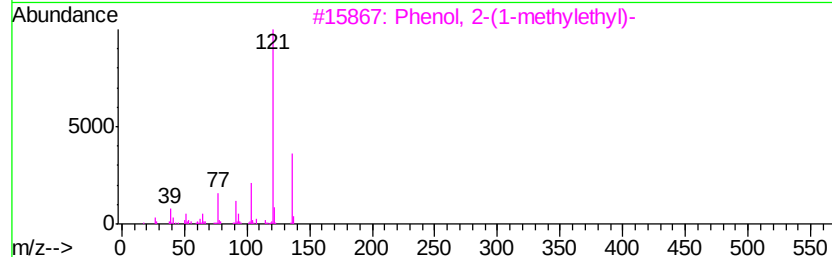
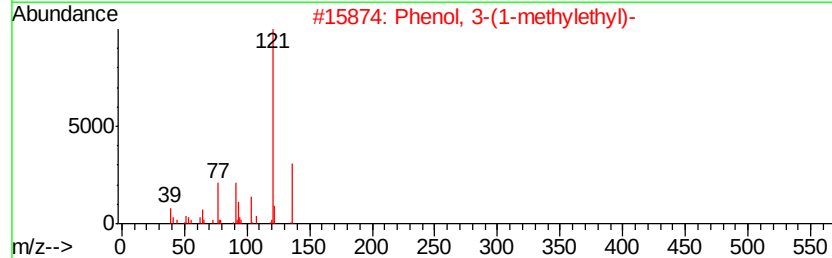
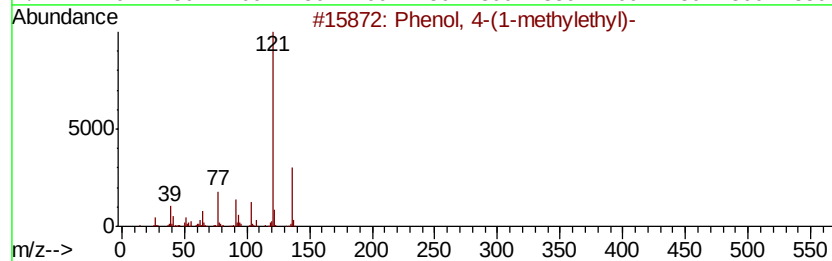
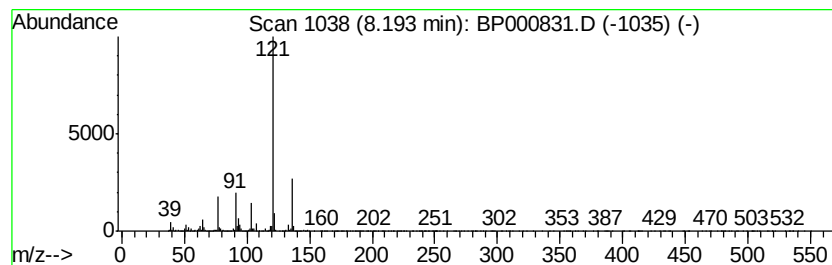
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 9 Phenol, 4-(1-methylethyl)- Concentration Rank 13

| R.T. | EstConc    | Area   | Relative to ISTD | R.T. |
|------|------------|--------|------------------|------|
| 8.19 | 3.15 ng/ul | 236910 | Napthalene-d8    | 8.03 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Phenol, 4-(1-methylethyl)- | 136 | C9H12O  | 000099-89-8 | 95   |
| 2    |    | Phenol, 3-(1-methylethyl)- | 136 | C9H12O  | 000618-45-1 | 91   |
| 3    |    | Phenol, 2-(1-methylethyl)- | 136 | C9H12O  | 000088-69-7 | 91   |
| 4    |    | Phenol, 3-(1-methylethyl)- | 136 | C9H12O  | 000618-45-1 | 91   |
| 5    |    | Phenol, 2-ethyl-5-methyl-  | 136 | C9H12O  | 001687-61-2 | 90   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP101719\  
Data File : BP000831.D  
Acq On : 17 Oct 2019 10:23  
Operator : JU  
Sample : K5191-03  
Misc :  
ALS Vial : 4 Sample Multiplier: 1

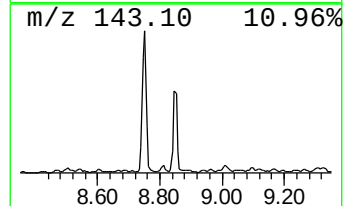
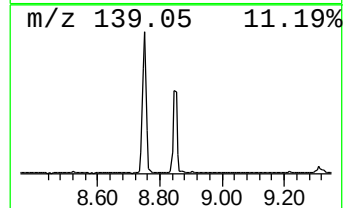
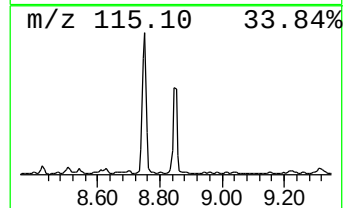
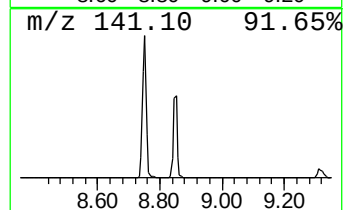
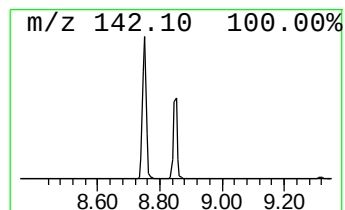
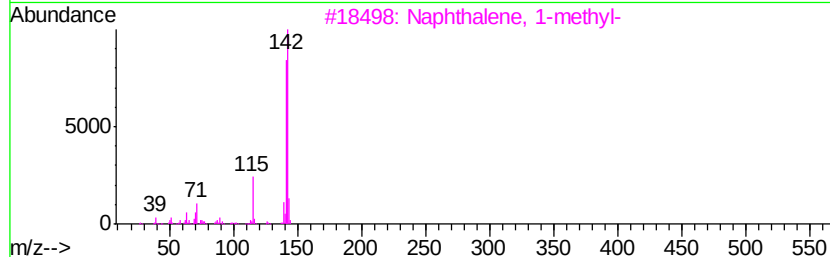
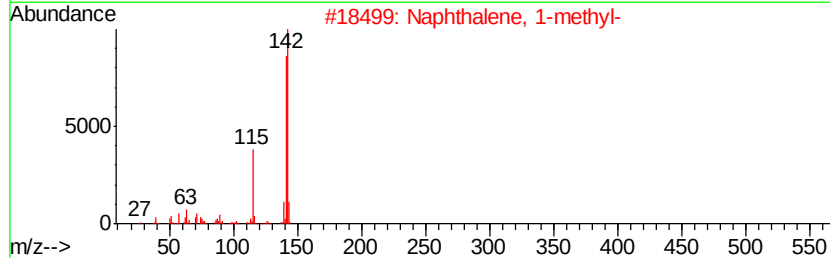
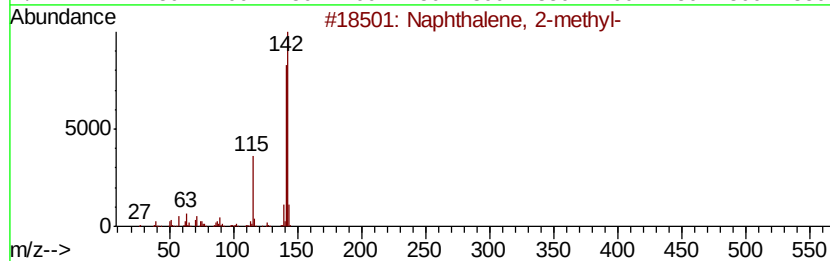
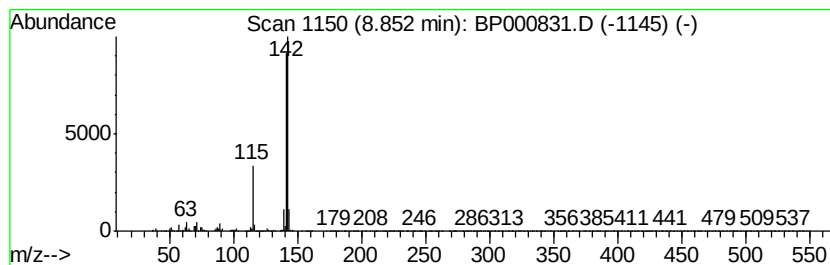
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ClientSampleId :  
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP100919MA.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 10 Naphthalene, 1-methyl- Concentration Rank 8

| R.T.    | EstConc    | Area                   | Relative to ISTD | R.T.           |
|---------|------------|------------------------|------------------|----------------|
| 8.85    | 4.34 ng/ul | 326234                 | Naphthalene-d8   | 8.03           |
| Hit# of | 5          | Tentative ID           | MW MolForm       | CAS# Qual      |
| 1       |            | Naphthalene, 2-methyl- | 142 C11H10       | 000091-57-6 96 |
| 2       |            | Naphthalene, 1-methyl- | 142 C11H10       | 000090-12-0 96 |
| 3       |            | Naphthalene, 1-methyl- | 142 C11H10       | 000090-12-0 94 |
| 4       |            | Naphthalene, 2-methyl- | 142 C11H10       | 000091-57-6 94 |
| 5       |            | Naphthalene, 1-methyl- | 142 C11H10       | 000090-12-0 91 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP101719\  
Data File : BP000831.D  
Acq On : 17 Oct 2019 10:23  
Operator : JU  
Sample : K5191-03  
Misc :  
ALS Vial : 4 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
BF6L0

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Quant Title : SVOA CALIBRATION

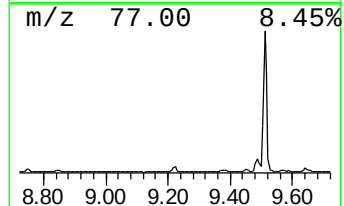
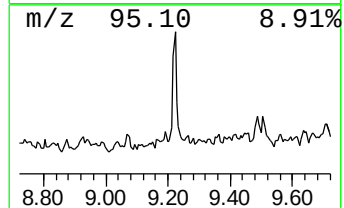
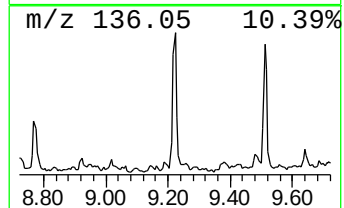
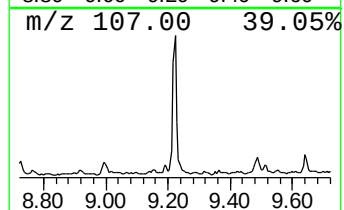
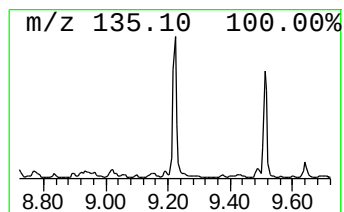
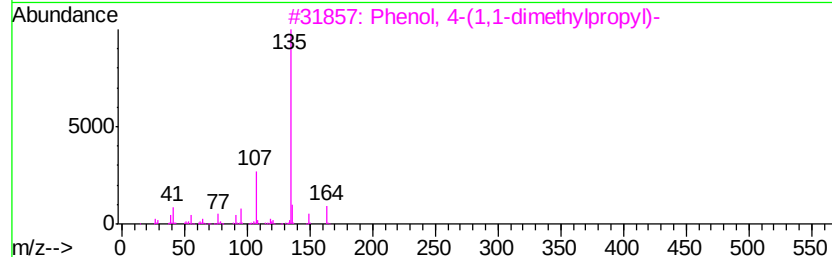
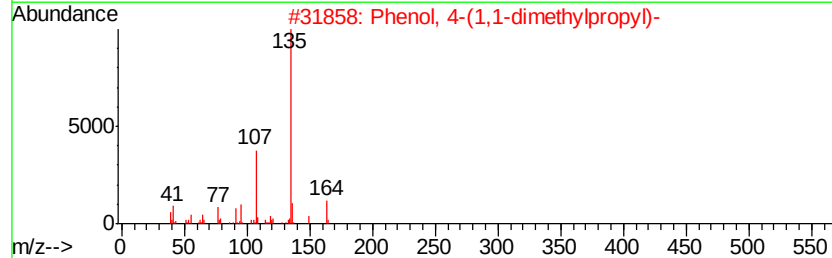
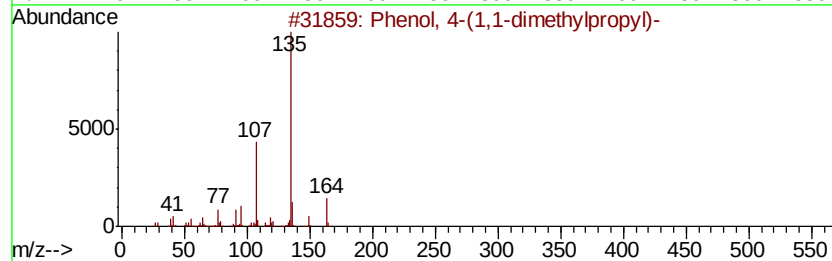
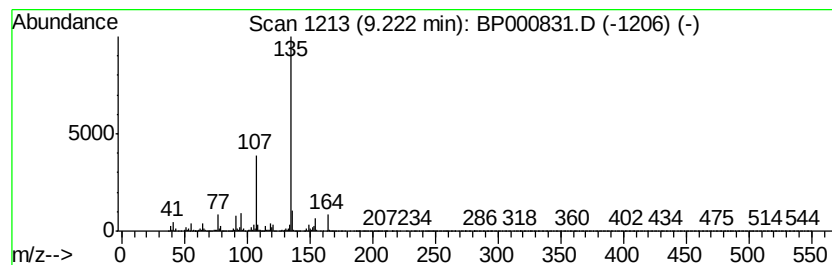
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 11 Phenol, 4-(1,1-dimethylprop... Concentration Rank 16

| R.T. | EstConc    | Area   | Relative to ISTD | R.T. |
|------|------------|--------|------------------|------|
| 9.22 | 2.33 ng/ul | 195029 | Acenaphthene-d10 | 9.80 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | Phenol, 4-(1,1-dimethylpropyl)-     | 164 | C11H16O  | 000080-46-6 | 94   |
| 2    |    |   | Phenol, 4-(1,1-dimethylpropyl)-     | 164 | C11H16O  | 000080-46-6 | 93   |
| 3    |    |   | Phenol, 4-(1,1-dimethylpropyl)-     | 164 | C11H16O  | 000080-46-6 | 81   |
| 4    |    |   | Phenol, m-tert-butyl-               | 150 | C10H14O  | 000585-34-2 | 80   |
| 5    |    |   | Phenol, 4,4'-(1,2-diethyl-1,2-et... | 270 | C18H22O2 | 000084-16-2 | 72   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP101719\  
Data File : BP000831.D  
Acq On : 17 Oct 2019 10:23  
Operator : JU  
Sample : K5191-03  
Misc :  
ALS Vial : 4 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP100919MA.M  
Quant Title : SVOA CALIBRATION

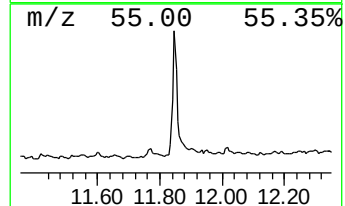
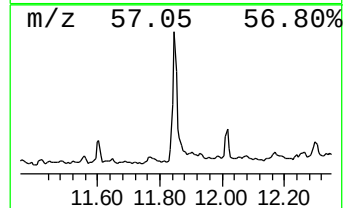
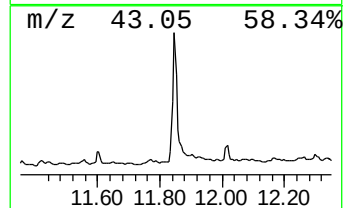
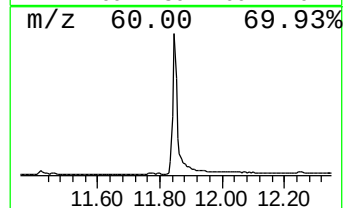
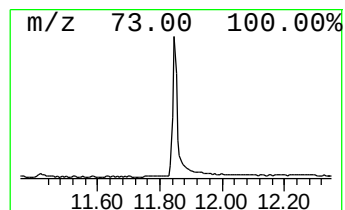
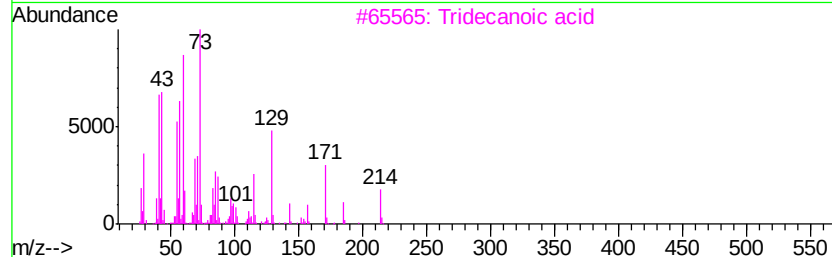
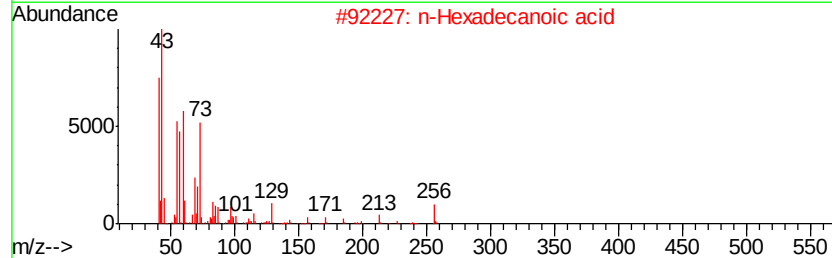
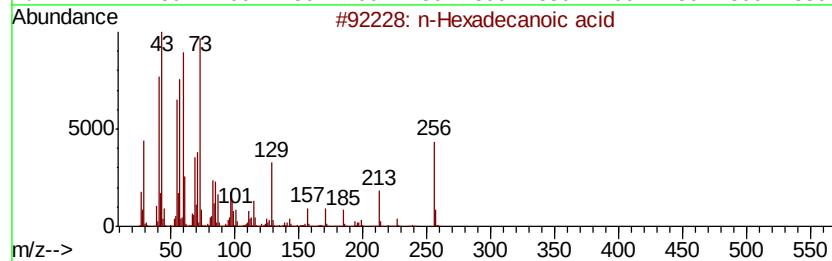
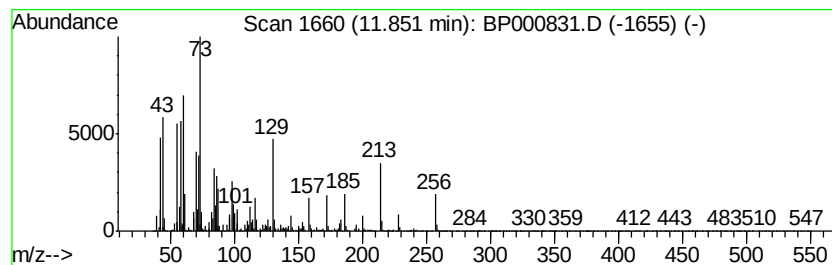
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TIC Integration Parameters: LSCINT.P

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Peak Number 12 n-Hexadecanoic acid Concentration Rank 2

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 11.85 | 9.54 ng/ul | 793541 | Phenanthrene-d10 | 11.30 |

| Hit# | of | 5 | Tentative ID        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|---------------------|-----|----------|-------------|------|
| 1    |    |   | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 98   |
| 2    |    |   | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 98   |
| 3    |    |   | Tridecanoic acid    | 214 | C13H26O2 | 000638-53-9 | 93   |
| 4    |    |   | Tetradecanoic acid  | 228 | C14H28O2 | 000544-63-8 | 87   |
| 5    |    |   | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 86   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP101719\  
Data File : BP000831.D  
Acq On : 17 Oct 2019 10:23  
Operator : JU  
Sample : K5191-03  
Misc :  
ALS Vial : 4 Sample Multiplier: 1

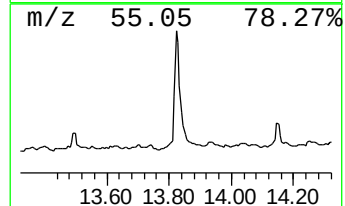
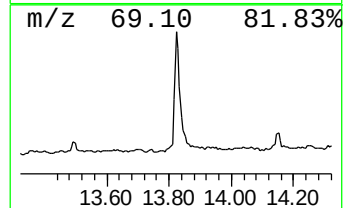
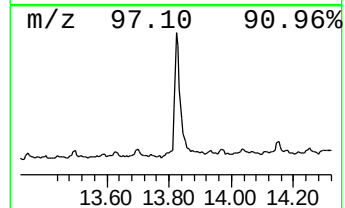
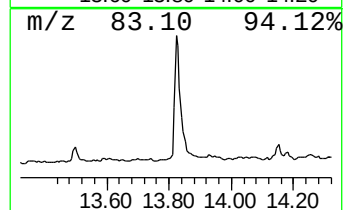
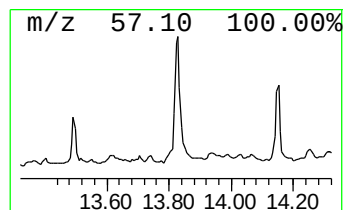
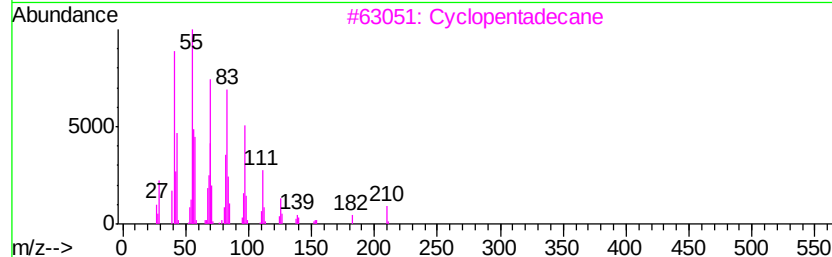
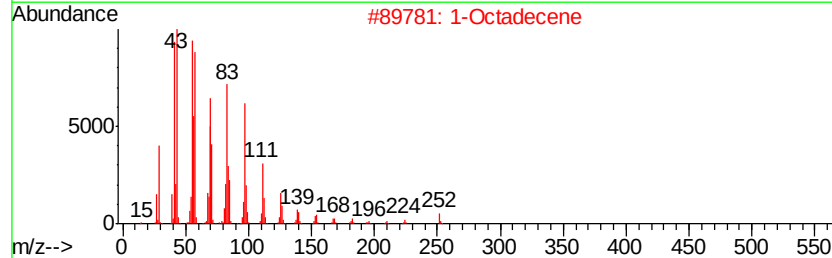
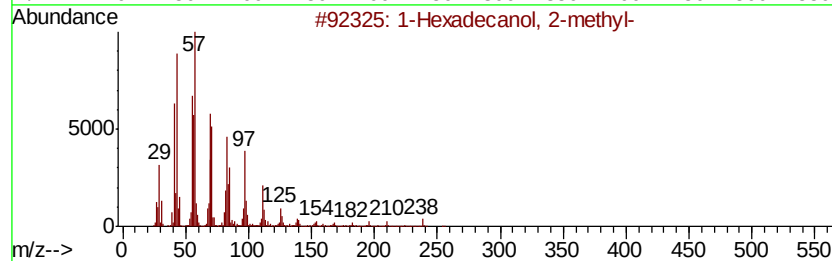
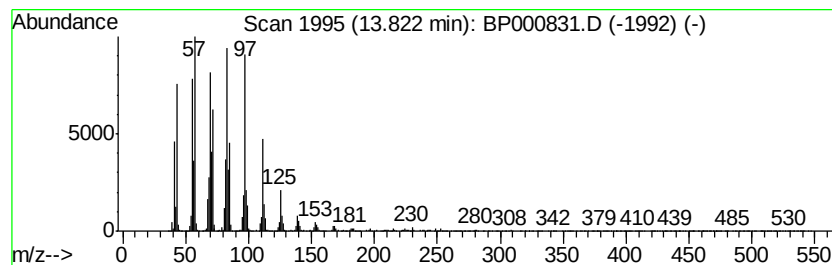
Instrument :  
BNA\_P  
ClientSampleId :  
BF6L0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP100919MA.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 13 1-Hexadecanol, 2-methyl- Concentration Rank 1

| R.T.    | EstConc     | Area                                 | Relative to ISTD | R.T.       |             |      |
|---------|-------------|--------------------------------------|------------------|------------|-------------|------|
| 13.82   | 10.33 ng/ul | 826326                               | Chrysene-d12     | 13.97      |             |      |
| Hit# of | 5           | Tentative ID                         | MW               | MolForm    | CAS#        | Qual |
| 1       |             | 1-Hexadecanol, 2-methyl-             | 256              | C17H36O    | 002490-48-4 | 97   |
| 2       |             | 1-Octadecene                         | 252              | C18H36     | 000112-88-9 | 93   |
| 3       |             | Cyclopentadecane                     | 210              | C15H30     | 000295-48-7 | 93   |
| 4       |             | Trifluoroacetic acid, n-octadecyl... | 366              | C20H37F3O2 | 079392-43-1 | 91   |
| 5       |             | 3-Eicosene, (E)-                     | 280              | C20H40     | 074685-33-9 | 91   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP101719\  
Data File : BP000831.D  
Acq On : 17 Oct 2019 10:23  
Operator : JU  
Sample : K5191-03  
Misc :  
ALS Vial : 4 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
BF6L0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP100919MA.M  
Quant Title : SVOA CALIBRATION

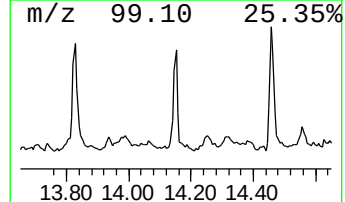
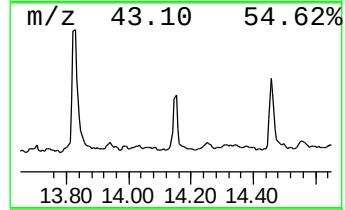
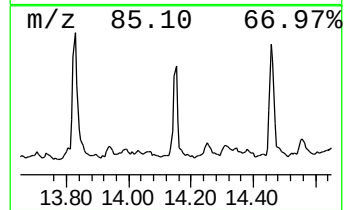
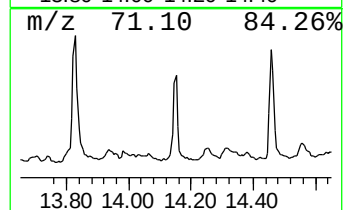
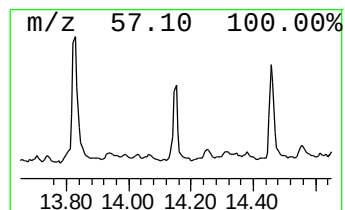
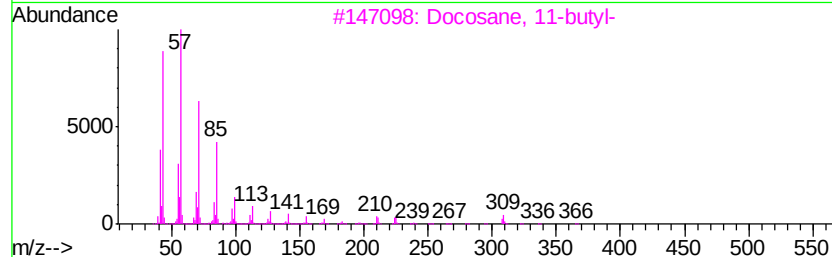
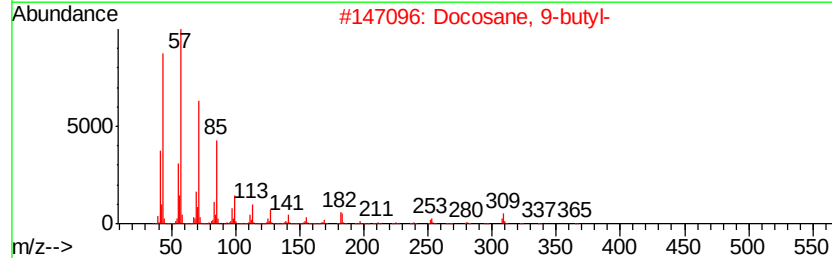
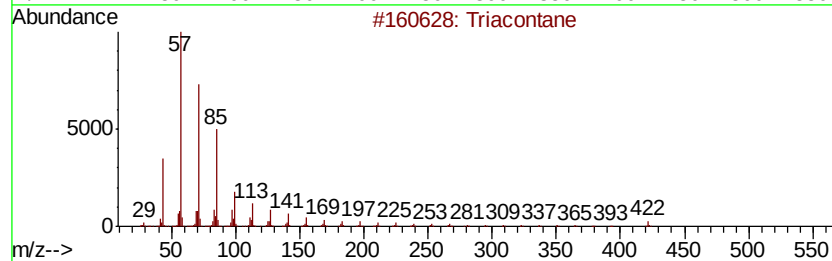
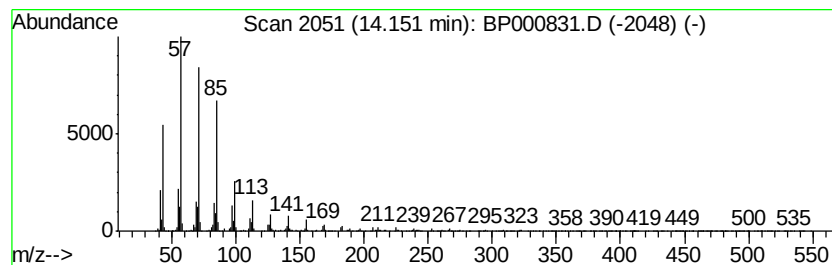
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 14 (DEL) Alkane: Straight-Chai... Concentration Rank 19

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 14.15 | 2.17 ng/ul | 173998 | Chrysene-d12     | 13.97 |

| Hit# | of | Tentative ID        | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------|-----|---------|-------------|------|
| 1    | 5  | Triacontane         | 422 | C30H62  | 000638-68-6 | 91   |
| 2    |    | Docosane, 9-butyl-  | 366 | C26H54  | 055282-14-9 | 91   |
| 3    |    | Docosane, 11-butyl- | 366 | C26H54  | 013475-76-8 | 91   |
| 4    |    | Heneicosane         | 296 | C21H44  | 000629-94-7 | 91   |
| 5    |    | Tetratriacontane    | 479 | C34H70  | 014167-59-0 | 91   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP101719\  
Data File : BP000831.D  
Acq On : 17 Oct 2019 10:23  
Operator : JU  
Sample : K5191-03  
Misc :  
ALS Vial : 4 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP100919MA.M  
Quant Title : SVOA CALIBRATION

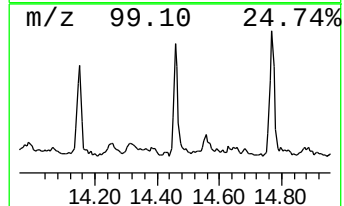
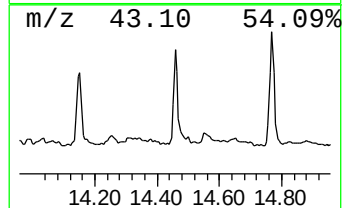
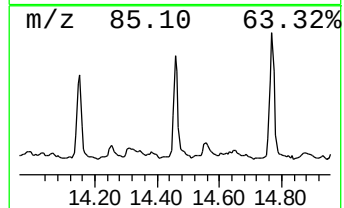
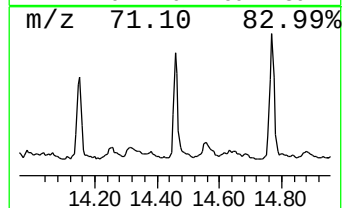
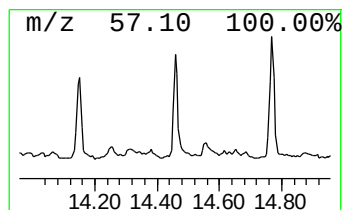
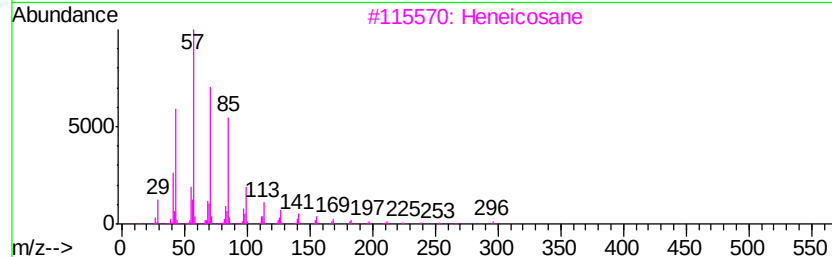
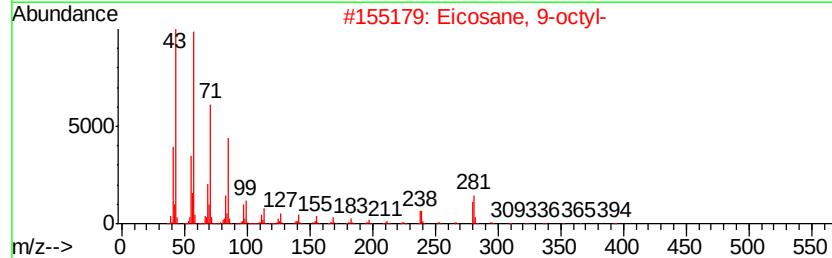
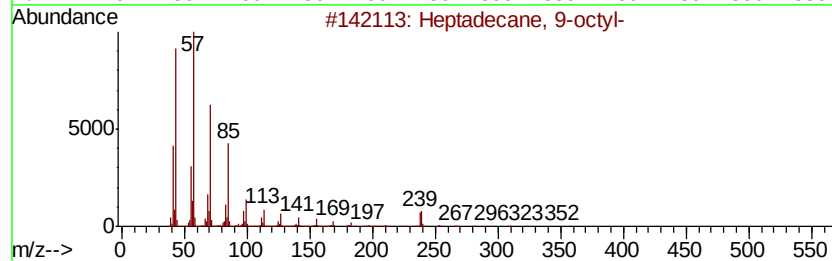
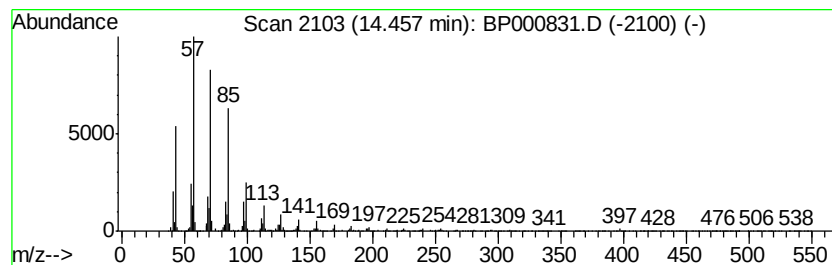
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 15 (DEL) Alkane: Straight-Chai... Concentration Rank 9

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 14.46 | 3.79 ng/ul | 303438 | Chrysene-d12     | 13.97 |

| Hit# | of | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------|-----|---------|-------------|------|
| 1    | 5  | Heptadecane, 9-octyl- | 352 | C25H52  | 007225-64-1 | 93   |
| 2    |    | Eicosane, 9-octyl-    | 394 | C28H58  | 013475-77-9 | 93   |
| 3    |    | Heneicosane           | 296 | C21H44  | 000629-94-7 | 91   |
| 4    |    | Docosane, 7-hexyl-    | 394 | C28H58  | 055373-86-9 | 90   |
| 5    |    | Nonacosane            | 408 | C29H60  | 000630-03-5 | 87   |





Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP101719\  
Data File : BP000831.D  
Acq On : 17 Oct 2019 10:23  
Operator : JU  
Sample : K5191-03  
Misc :  
ALS Vial : 4 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
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Quant Title : SVOA CALIBRATION

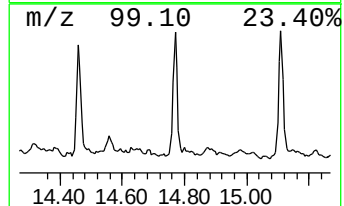
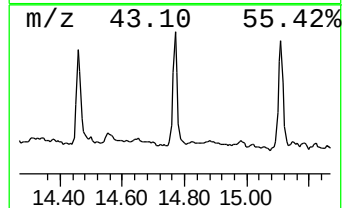
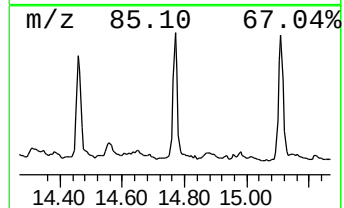
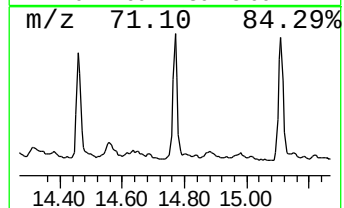
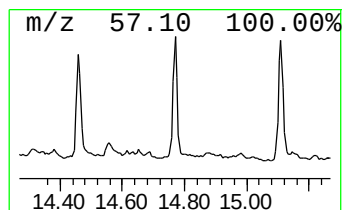
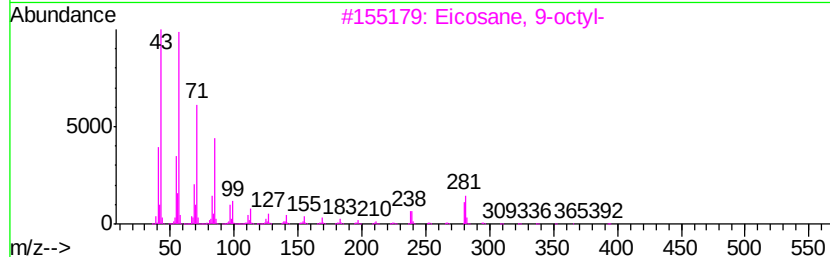
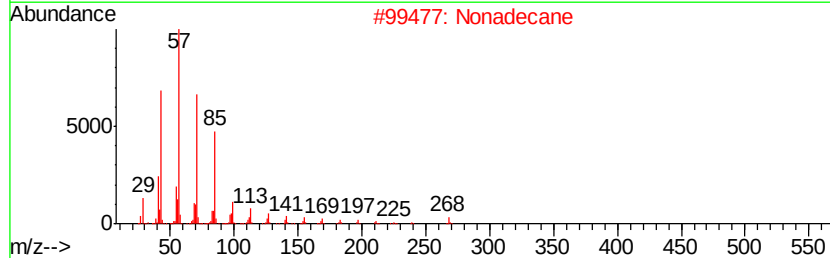
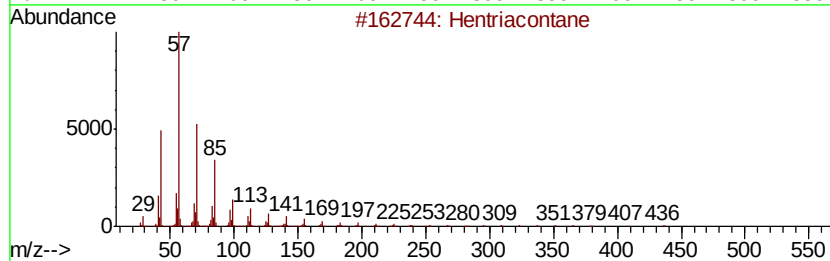
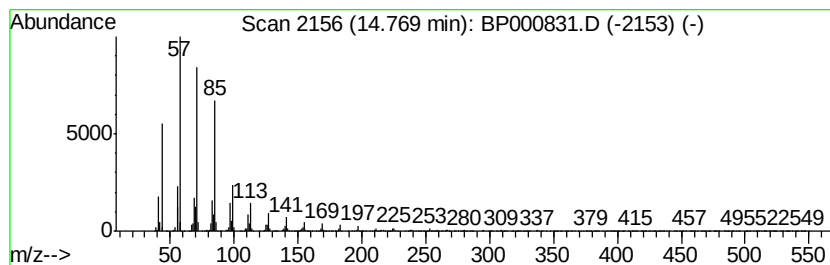
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 16 (DEL) Alkane: Straight-Chain Concentration Rank 10

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 14.77 | 3.72 ng/ul | 299731 | Perylene-d12     | 15.46 |

| Hit# | of | Tentative ID       | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------|-----|---------|-------------|------|
| 1    | 5  | Hentriacontane     | 437 | C31H64  | 000630-04-6 | 93   |
| 2    |    | Nonadecane         | 268 | C19H40  | 000629-92-5 | 93   |
| 3    |    | Eicosane, 9-octyl- | 394 | C28H58  | 013475-77-9 | 93   |
| 4    |    | Heneicosane        | 296 | C21H44  | 000629-94-7 | 91   |
| 5    |    | triacontane        | 422 | C30H62  | 000638-68-6 | 91   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP101719\  
Data File : BP000831.D  
Acq On : 17 Oct 2019 10:23  
Operator : JU  
Sample : K5191-03  
Misc :  
ALS Vial : 4 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
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Quant Title : SVOA CALIBRATION

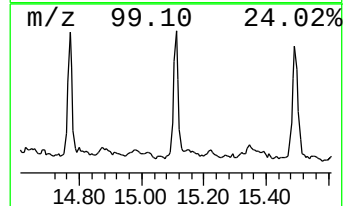
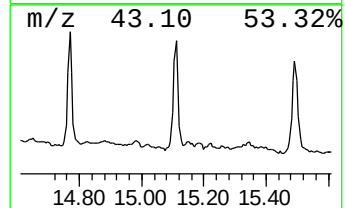
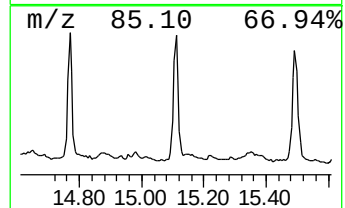
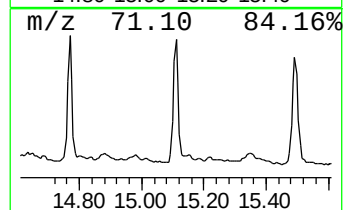
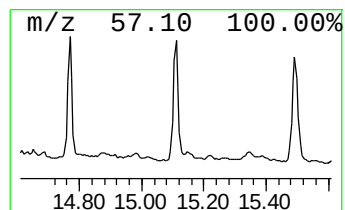
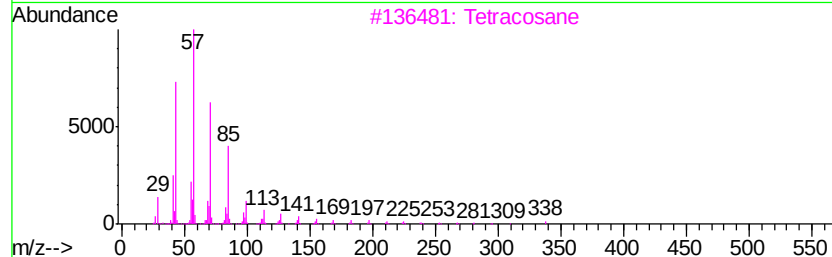
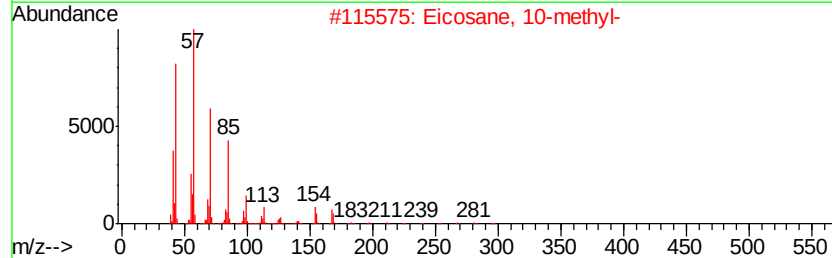
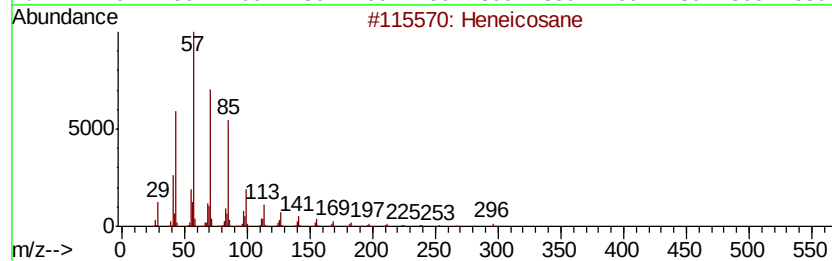
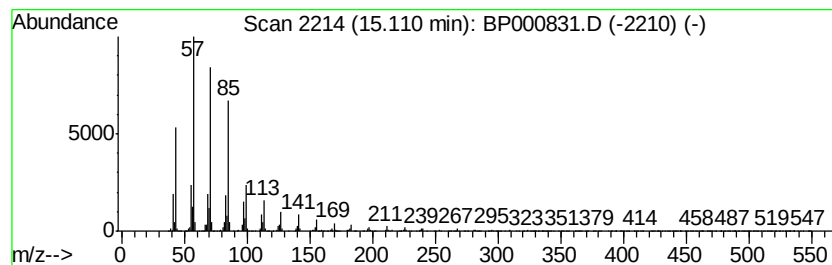
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 17 (DEL) Alkane: Straight-Chai... Concentration Rank 7

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 15.11 | 4.37 ng/ul | 352200 | Perylene-d12     | 15.46 |

| Hit# of | 5                    | Tentative ID | MW  | MolForm | CAS#        | Qual |
|---------|----------------------|--------------|-----|---------|-------------|------|
| 1       | Heneicosane          |              | 296 | C21H44  | 000629-94-7 | 96   |
| 2       | Eicosane, 10-methyl- |              | 296 | C21H44  | 054833-23-7 | 95   |
| 3       | Tetracosane          |              | 338 | C24H50  | 000646-31-1 | 94   |
| 4       | Heptadecane          |              | 240 | C17H36  | 000629-78-7 | 93   |
| 5       | triacontane          |              | 422 | C30H62  | 000638-68-6 | 91   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP101719\  
Data File : BP000831.D  
Acq On : 17 Oct 2019 10:23  
Operator : JU  
Sample : K5191-03  
Misc :  
ALS Vial : 4 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
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Quant Title : SVOA CALIBRATION

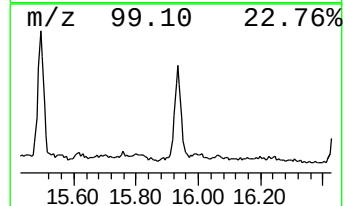
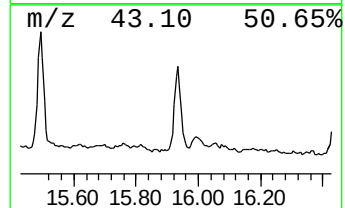
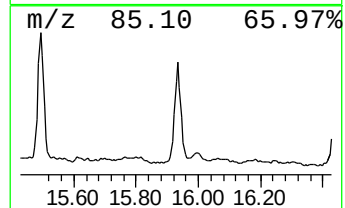
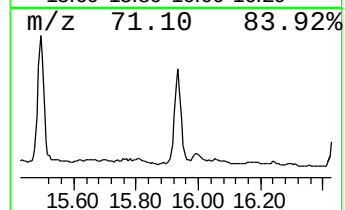
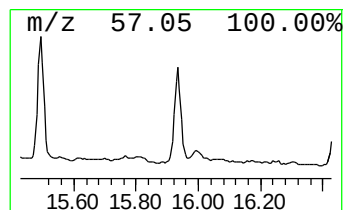
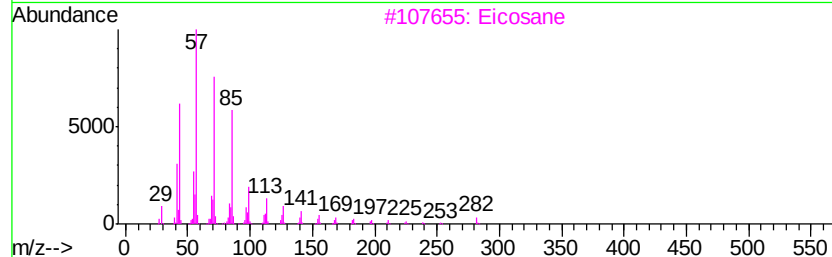
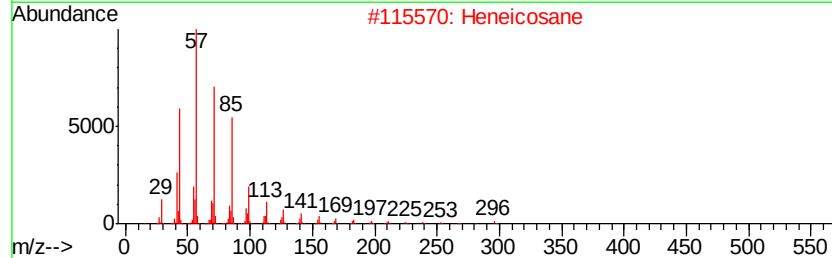
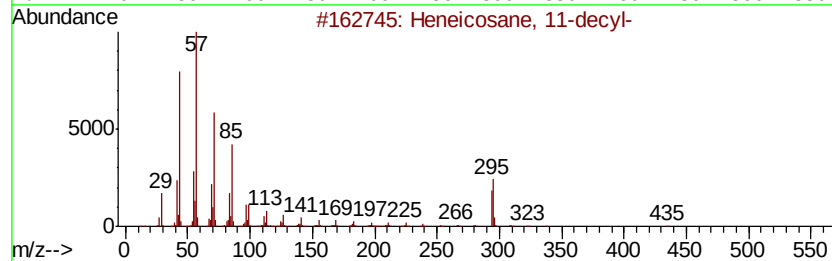
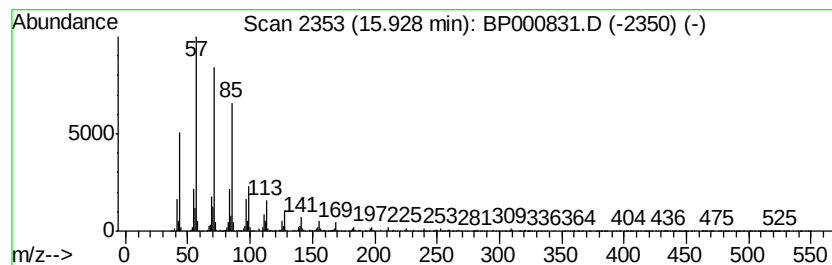
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 18 (DEL) Alkane: Straight-Chai... Concentration Rank 11

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 15.93 | 3.55 ng/ul | 285974 | Perylene-d12     | 15.46 |

| Hit# | of | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------|-----|---------|-------------|------|
| 1    | 5  | Heneicosane, 11-decyl- | 437 | C31H64  | 055320-06-4 | 93   |
| 2    |    | Heneicosane            | 296 | C21H44  | 000629-94-7 | 91   |
| 3    |    | Eicosane               | 282 | C20H42  | 000112-95-8 | 91   |
| 4    |    | Octacosane             | 394 | C28H58  | 000630-02-4 | 91   |
| 5    |    | Eicosane, 7-hexyl-     | 366 | C26H54  | 055333-99-8 | 91   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP101719\  
Data File : BP000831.D  
Acq On : 17 Oct 2019 10:23  
Operator : JU  
Sample : K5191-03  
Misc :  
ALS Vial : 4 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
BF6L0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP100919MA.M  
Quant Title : SVOA CALIBRATION

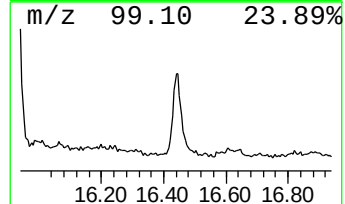
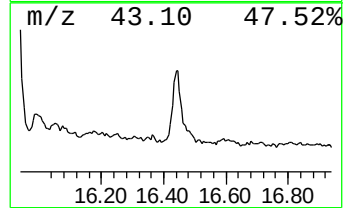
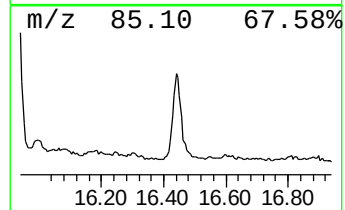
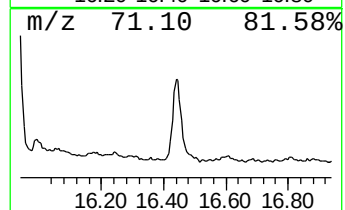
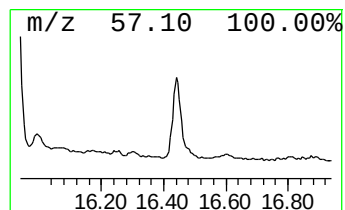
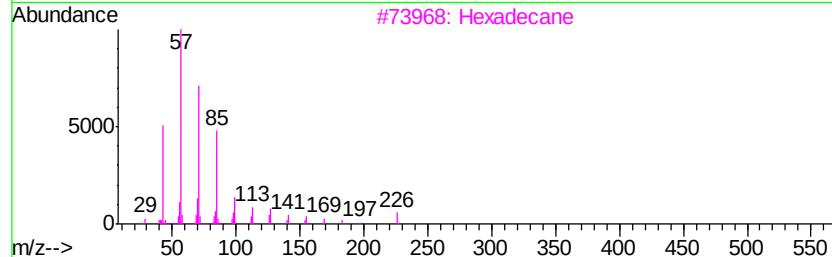
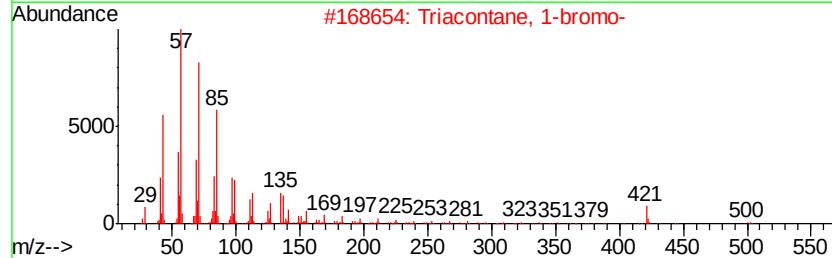
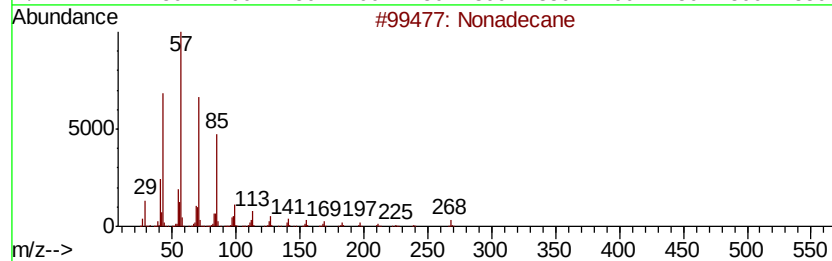
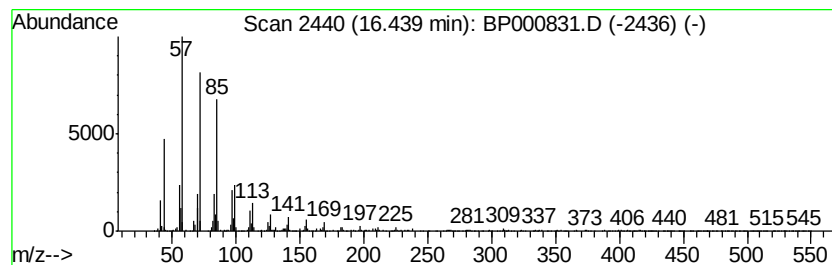
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 19 (DEL) Alkane: Straight-Chai... Concentration Rank 18

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 16.44 | 2.20 ng/ul | 177654 | Perylene-d12     | 15.46 |

| Hit# | of | Tentative ID          | MW  | MolForm  | CAS#        | Qual |
|------|----|-----------------------|-----|----------|-------------|------|
| 1    | 5  | Nonadecane            | 268 | C19H40   | 000629-92-5 | 95   |
| 2    |    | Triacontane, 1-bromo- | 500 | C30H61Br | 004209-22-7 | 90   |
| 3    |    | Hexadecane            | 226 | C16H34   | 000544-76-3 | 86   |
| 4    |    | Hentriacontane        | 437 | C31H64   | 000630-04-6 | 83   |
| 5    |    | Tetratriacontane      | 479 | C34H70   | 014167-59-0 | 83   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_P\DATA\BP101719\  
 Data File : BP000831.D  
 Acq On : 17 Oct 2019 10:23  
 Operator : JU  
 Sample : K5191-03  
 Misc :  
 ALS Vial : 4 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 BF6L0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\SOM-EPA-BP100919MA.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 |
| Benzene, 1-ethyl-... | 6.28  | 2.2     | ng/ul | 110749   | 1                     | 6.75  | 995119  | 20.0 |
| Benzene, 1,2,3-tr... | 6.58  | 7.6     | ng/ul | 377231   | 1                     | 6.75  | 995119  | 20.0 |
| Benzene, 1,3,5-tr... | 6.81  | 4.4     | ng/ul | 217471   | 1                     | 6.75  | 995119  | 20.0 |
| Indane               | 6.93  | 2.7     | ng/ul | 133368   | 1                     | 6.75  | 995119  | 20.0 |
| 3-Phenylbut-1-ene    | 7.78  | 4.7     | ng/ul | 354621   | 2                     | 8.03  | 1504140 | 20.0 |
| Phenol, 4-(1-meth... | 8.19  | 3.1     | ng/ul | 236910   | 2                     | 8.03  | 1504140 | 20.0 |
| Naphthalene, 1-me... | 8.85  | 4.3     | ng/ul | 326234   | 2                     | 8.03  | 1504140 | 20.0 |
| Phenol, 4-(1,1-di... | 9.22  | 2.3     | ng/ul | 195029   | 3                     | 9.80  | 1676390 | 20.0 |
| n-Hexadecanoic acid  | 11.85 | 9.5     | ng/ul | 793541   | 4                     | 11.30 | 1663330 | 20.0 |
| 1-Hexadecanol, 2-... | 13.82 | 10.3    | ng/ul | 826326   | 5                     | 13.97 | 1600340 | 20.0 |
| (DEL) Alkane: Str... | 14.15 | 2.2     | ng/ul | 173998   | 5                     | 13.97 | 1600340 | 20.0 |
| (DEL) Alkane: Str... | 14.46 | 3.8     | ng/ul | 303438   | 5                     | 13.97 | 1600340 | 20.0 |
| (DEL) Alkane: Str... | 14.77 | 3.7     | ng/ul | 299731   | 6                     | 15.46 | 1612180 | 20.0 |
| (DEL) Alkane: Str... | 15.11 | 4.4     | ng/ul | 352200   | 6                     | 15.46 | 1612180 | 20.0 |
| (DEL) Alkane: Str... | 15.93 | 3.5     | ng/ul | 285974   | 6                     | 15.46 | 1612180 | 20.0 |
| (DEL) Alkane: Str... | 16.44 | 2.2     | ng/ul | 177654   | 6                     | 15.46 | 1612180 | 20.0 |