

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081320\  
 Data File : BG046256.D  
 Acq On : 13 Aug 2020 19:42  
 Operator : CG/JU  
 Sample : L3629-14  
 Misc :  
 ALS Vial : 10 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 BFM61

## 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\_G\METHODS\SOM-EPA-BG081320PAH.M  
 Title : SVOA CALIBRATION

Signal : TIC

| peak # | R.T. min | first scan | max scan | last scan | PK TY | peak height | corr. area | corr. % max. | % of total |
|--------|----------|------------|----------|-----------|-------|-------------|------------|--------------|------------|
| 1      | 3.170    | 8          | 14       | 43        | rVB   | 1568920     | 3205249    | 100.00%      | 12.075%    |
| 2      | 3.422    | 53         | 57       | 62        | rVB3  | 10090       | 13913      | 0.43%        | 0.052%     |
| 3      | 3.845    | 125        | 129      | 132       | rVB4  | 3834        | 5588       | 0.17%        | 0.021%     |
| 4      | 3.939    | 140        | 145      | 153       | rBV2  | 22315       | 33604      | 1.05%        | 0.127%     |
| 5      | 4.010    | 153        | 157      | 161       | rVB5  | 3289        | 5550       | 0.17%        | 0.021%     |
| 6      | 4.075    | 163        | 168      | 175       | rVV2  | 15419       | 26361      | 0.82%        | 0.099%     |
| 7      | 4.163    | 177        | 183      | 192       | rVB   | 16334       | 27179      | 0.85%        | 0.102%     |
| 8      | 5.073    | 332        | 338      | 348       | rBV2  | 14871       | 26919      | 0.84%        | 0.101%     |
| 9      | 7.118    | 678        | 686      | 696       | rBV   | 204466      | 384904     | 12.01%       | 1.450%     |
| 10     | 7.277    | 705        | 713      | 725       | rBV   | 187642      | 315960     | 9.86%        | 1.190%     |
| 11     | 7.482    | 741        | 748      | 756       | rBV   | 216648      | 387720     | 12.10%       | 1.461%     |
| 12     | 7.565    | 758        | 762      | 764       | rVV2  | 7375        | 11200      | 0.35%        | 0.042%     |
| 13     | 7.582    | 764        | 765      | 772       | rVB2  | 5942        | 8116       | 0.25%        | 0.031%     |
| 14     | 7.952    | 821        | 828      | 835       | rBV   | 325422      | 551873     | 17.22%       | 2.079%     |
| 15     | 8.652    | 938        | 947      | 961       | rBV   | 237516      | 411077     | 12.83%       | 1.549%     |
| 16     | 8.769    | 963        | 967      | 975       | rVB4  | 6063        | 11343      | 0.35%        | 0.043%     |
| 17     | 9.104    | 1015       | 1024     | 1034      | rBV   | 204472      | 377157     | 11.77%       | 1.421%     |
| 18     | 9.827    | 1139       | 1147     | 1156      | rVB   | 173088      | 306446     | 9.56%        | 1.155%     |
| 19     | 10.367   | 1232       | 1239     | 1251      | rBV   | 274949      | 490282     | 15.30%       | 1.847%     |
| 20     | 10.743   | 1293       | 1303     | 1311      | rBV   | 450300      | 825134     | 25.74%       | 3.109%     |
| 21     | 10.872   | 1318       | 1325     | 1339      | rVV   | 200367      | 378500     | 11.81%       | 1.426%     |
| 22     | 12.330   | 1567       | 1573     | 1580      | rBV3  | 4527        | 8611       | 0.27%        | 0.032%     |
| 23     | 13.417   | 1753       | 1758     | 1763      | rBV5  | 4138        | 7049       | 0.22%        | 0.027%     |
| 24     | 13.481   | 1763       | 1769     | 1776      | rVB   | 80466       | 122723     | 3.83%        | 0.462%     |
| 25     | 13.981   | 1845       | 1854     | 1858      | rBV   | 551455      | 805297     | 25.12%       | 3.034%     |
| 26     | 14.028   | 1858       | 1862     | 1868      | rVB   | 195705      | 277429     | 8.66%        | 1.045%     |
| 27     | 14.269   | 1892       | 1903     | 1917      | rBV   | 583209      | 953945     | 29.76%       | 3.594%     |
| 28     | 14.574   | 1948       | 1955     | 1964      | rBV2  | 728591      | 1154974    | 36.03%       | 4.351%     |
| 29     | 14.639   | 1964       | 1966     | 1974      | rVB3  | 6772        | 9236       | 0.29%        | 0.035%     |
| 30     | 14.762   | 1981       | 1987     | 1999      | rBV   | 118592      | 245679     | 7.66%        | 0.926%     |
| 31     | 14.979   | 2019       | 2024     | 2030      | rVB5  | 5569        | 11646      | 0.36%        | 0.044%     |
| 32     | 15.391   | 2087       | 2094     | 2098      | rVV5  | 2917        | 6723       | 0.21%        | 0.025%     |
| 33     | 15.567   | 2117       | 2124     | 2131      | rBV   | 854033      | 1233501    | 38.48%       | 4.647%     |
| 34     | 15.620   | 2131       | 2133     | 2137      | rVV4  | 9561        | 10460      | 0.33%        | 0.039%     |

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Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG081320PAH.M  
 Title : SVOA CALIBRATION

|    |        |      |      |      |      |        |         |        |        |
|----|--------|------|------|------|------|--------|---------|--------|--------|
| 35 | 15.679 | 2137 | 2143 | 2152 | rVV  | 192493 | 304692  | 9.51%  | 1.148% |
| 36 | 15.802 | 2159 | 2164 | 2168 | rBV5 | 9008   | 14518   | 0.45%  | 0.055% |
| 37 | 16.061 | 2202 | 2208 | 2209 | rBV4 | 4912   | 5794    | 0.18%  | 0.022% |
| 38 | 16.296 | 2244 | 2248 | 2251 | rBV5 | 6370   | 9189    | 0.29%  | 0.035% |
| 39 | 16.630 | 2297 | 2305 | 2310 | rVB7 | 3685   | 7821    | 0.24%  | 0.029% |
| 40 | 16.860 | 2337 | 2344 | 2348 | rBV5 | 8438   | 13834   | 0.43%  | 0.052% |
| 41 | 16.901 | 2348 | 2351 | 2353 | rVV2 | 4886   | 5743    | 0.18%  | 0.022% |
| 42 | 16.971 | 2357 | 2363 | 2370 | rVV2 | 11904  | 22109   | 0.69%  | 0.083% |
| 43 | 17.053 | 2372 | 2377 | 2379 | rVV5 | 4974   | 6682    | 0.21%  | 0.025% |
| 44 | 17.100 | 2381 | 2385 | 2388 | rVV5 | 5588   | 9412    | 0.29%  | 0.035% |
| 45 | 17.136 | 2388 | 2391 | 2397 | rVB3 | 6167   | 9663    | 0.30%  | 0.036% |
| 46 | 17.318 | 2414 | 2422 | 2426 | rBV  | 912584 | 1398472 | 43.63% | 5.269% |
| 47 | 17.359 | 2426 | 2429 | 2433 | rVV  | 160127 | 226542  | 7.07%  | 0.853% |
| 48 | 17.412 | 2433 | 2438 | 2451 | rVV2 | 837229 | 1303393 | 40.66% | 4.910% |
| 49 | 17.500 | 2451 | 2453 | 2460 | rVB7 | 7487   | 13388   | 0.42%  | 0.050% |
| 50 | 17.712 | 2481 | 2489 | 2493 | rBV2 | 12391  | 27311   | 0.85%  | 0.103% |
| 51 | 17.835 | 2504 | 2510 | 2514 | rVB6 | 8395   | 13366   | 0.42%  | 0.050% |
| 52 | 17.894 | 2517 | 2520 | 2524 | rBV5 | 6392   | 8167    | 0.25%  | 0.031% |
| 53 | 18.111 | 2554 | 2557 | 2562 | rBV5 | 4151   | 6299    | 0.20%  | 0.024% |
| 54 | 18.217 | 2566 | 2575 | 2578 | rBV2 | 92232  | 180592  | 5.63%  | 0.680% |
| 55 | 18.317 | 2589 | 2592 | 2599 | rVB  | 11147  | 14949   | 0.47%  | 0.056% |
| 56 | 18.387 | 2599 | 2604 | 2607 | rVV2 | 36897  | 66384   | 2.07%  | 0.250% |
| 57 | 18.422 | 2607 | 2610 | 2616 | rVB2 | 26751  | 40321   | 1.26%  | 0.152% |
| 58 | 18.511 | 2622 | 2625 | 2632 | rVB9 | 6305   | 13999   | 0.44%  | 0.053% |
| 59 | 18.693 | 2651 | 2656 | 2658 | rVV  | 27794  | 39063   | 1.22%  | 0.147% |
| 60 | 18.722 | 2658 | 2661 | 2666 | rVB2 | 28332  | 39720   | 1.24%  | 0.150% |
| 61 | 18.828 | 2674 | 2679 | 2682 | rBV7 | 3819   | 6521    | 0.20%  | 0.025% |
| 62 | 18.934 | 2694 | 2697 | 2702 | rVB5 | 11628  | 16046   | 0.50%  | 0.060% |
| 63 | 18.986 | 2702 | 2706 | 2710 | rBV2 | 9797   | 14998   | 0.47%  | 0.057% |
| 64 | 19.022 | 2710 | 2712 | 2716 | rVB5 | 8584   | 9903    | 0.31%  | 0.037% |
| 65 | 19.104 | 2719 | 2726 | 2730 | rBV3 | 39848  | 73513   | 2.29%  | 0.277% |
| 66 | 19.151 | 2730 | 2734 | 2739 | rVV7 | 13786  | 28899   | 0.90%  | 0.109% |
| 67 | 19.192 | 2739 | 2741 | 2745 | rVB4 | 9662   | 11341   | 0.35%  | 0.043% |
| 68 | 19.239 | 2745 | 2749 | 2758 | rVB3 | 29461  | 55026   | 1.72%  | 0.207% |
| 69 | 19.363 | 2761 | 2770 | 2778 | rBV  | 356279 | 541611  | 16.90% | 2.040% |
| 70 | 19.486 | 2788 | 2791 | 2794 | rBV5 | 19556  | 24154   | 0.75%  | 0.091% |
| 71 | 19.556 | 2800 | 2803 | 2805 | rBV4 | 11512  | 13307   | 0.42%  | 0.050% |

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Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG081320PAH.M  
 Title : SVOA CALIBRATION

|     |        |      |      |      |       |         |         |        |        |
|-----|--------|------|------|------|-------|---------|---------|--------|--------|
| 72  | 19.697 | 2821 | 2827 | 2831 | rBV   | 1131621 | 1790231 | 55.85% | 6.744% |
| 73  | 19.803 | 2841 | 2845 | 2850 | rVB2  | 18170   | 29903   | 0.93%  | 0.113% |
| 74  | 19.903 | 2858 | 2862 | 2867 | rBV   | 23773   | 38810   | 1.21%  | 0.146% |
| 75  | 19.968 | 2869 | 2873 | 2878 | rVB2  | 19084   | 30352   | 0.95%  | 0.114% |
| 76  | 20.044 | 2882 | 2886 | 2888 | rBV3  | 34966   | 50181   | 1.57%  | 0.189% |
| 77  | 20.185 | 2906 | 2910 | 2911 | rBV2  | 22435   | 28180   | 0.88%  | 0.106% |
| 78  | 20.214 | 2912 | 2915 | 2920 | rVB   | 48023   | 67950   | 2.12%  | 0.256% |
| 79  | 20.320 | 2929 | 2933 | 2937 | rBV2  | 24429   | 42710   | 1.33%  | 0.161% |
| 80  | 20.373 | 2937 | 2942 | 2947 | rVB3  | 48287   | 83219   | 2.60%  | 0.314% |
| 81  | 20.449 | 2953 | 2955 | 2956 | rBV2  | 12754   | 9393    | 0.29%  | 0.035% |
| 82  | 20.520 | 2962 | 2967 | 2970 | rBV2  | 29540   | 45366   | 1.42%  | 0.171% |
| 83  | 20.561 | 2970 | 2974 | 2978 | rVV4  | 27120   | 43605   | 1.36%  | 0.164% |
| 84  | 20.967 | 3040 | 3043 | 3051 | rVB   | 51235   | 78858   | 2.46%  | 0.297% |
| 85  | 21.137 | 3068 | 3072 | 3073 | rBV2  | 27376   | 35443   | 1.11%  | 0.134% |
| 86  | 21.196 | 3079 | 3082 | 3084 | rBV2  | 30121   | 37567   | 1.17%  | 0.142% |
| 87  | 21.231 | 3084 | 3088 | 3094 | rVV4  | 253704  | 374365  | 11.68% | 1.410% |
| 88  | 21.560 | 3136 | 3144 | 3149 | rBV2  | 1001568 | 1829995 | 57.09% | 6.894% |
| 89  | 21.601 | 3149 | 3151 | 3162 | rVB   | 174865  | 292871  | 9.14%  | 1.103% |
| 90  | 22.371 | 3279 | 3282 | 3289 | rVB2  | 75882   | 134703  | 4.20%  | 0.507% |
| 91  | 23.041 | 3392 | 3396 | 3401 | rVB   | 69880   | 110956  | 3.46%  | 0.418% |
| 92  | 23.693 | 3499 | 3507 | 3515 | rBV2  | 123708  | 426603  | 13.31% | 1.607% |
| 93  | 23.822 | 3523 | 3529 | 3533 | rBV   | 121986  | 235311  | 7.34%  | 0.887% |
| 94  | 23.869 | 3535 | 3537 | 3544 | rVB6  | 49324   | 75983   | 2.37%  | 0.286% |
| 95  | 24.374 | 3620 | 3623 | 3628 | rBV2  | 42839   | 74625   | 2.33%  | 0.281% |
| 96  | 24.457 | 3629 | 3637 | 3644 | rBV   | 505550  | 1258009 | 39.25% | 4.739% |
| 97  | 24.668 | 3663 | 3673 | 3681 | rBV2  | 564169  | 1563380 | 48.78% | 5.890% |
| 98  | 24.733 | 3681 | 3684 | 3694 | rVB5  | 95129   | 209157  | 6.53%  | 0.788% |
| 99  | 25.820 | 3864 | 3869 | 3879 | rVB2  | 90710   | 245409  | 7.66%  | 0.925% |
| 100 | 25.925 | 3883 | 3887 | 3895 | rVV10 | 22325   | 52364   | 1.63%  | 0.197% |

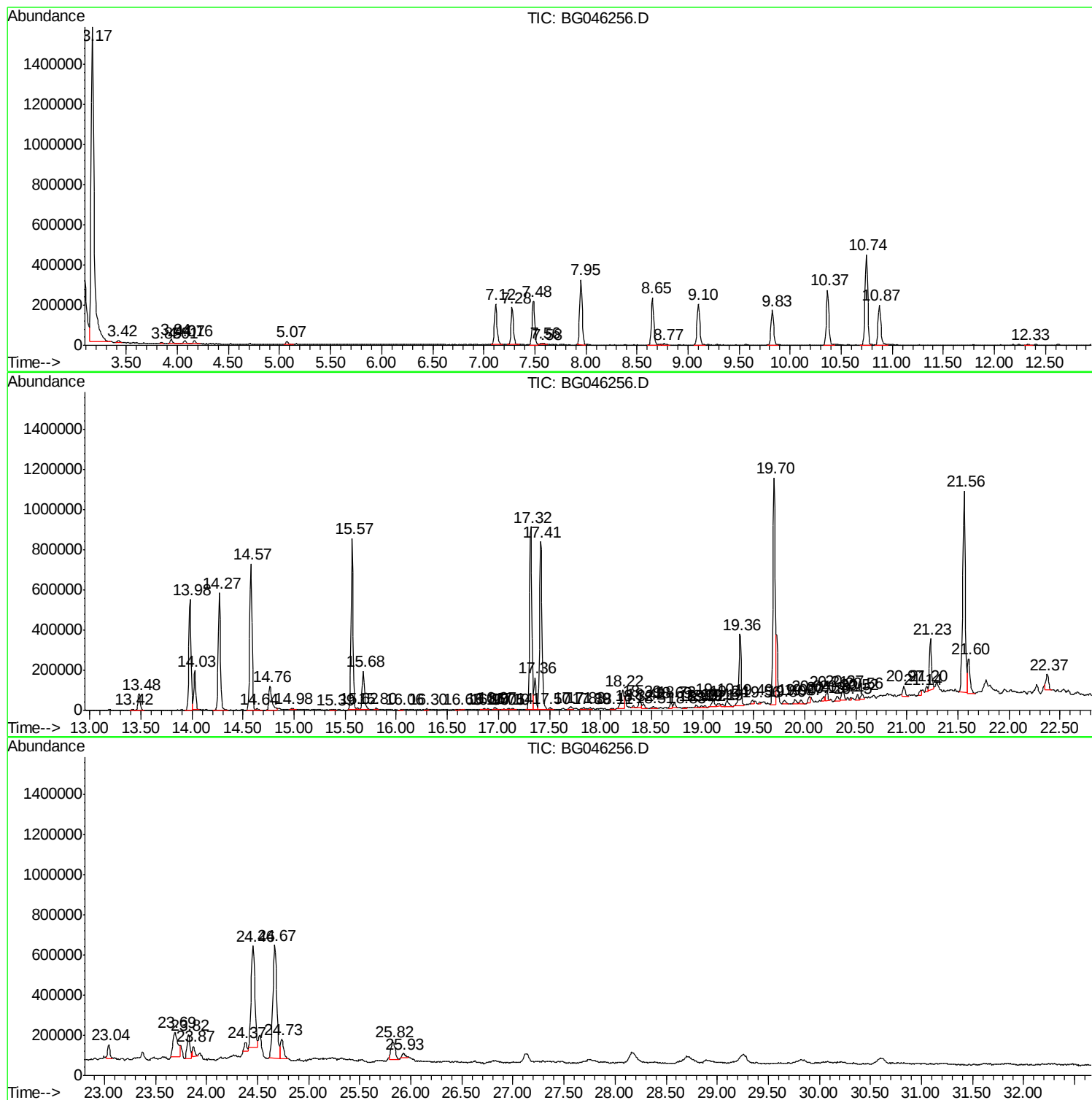
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG081320PAH.M  
Quant Title : SVOA CALIBRATION

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



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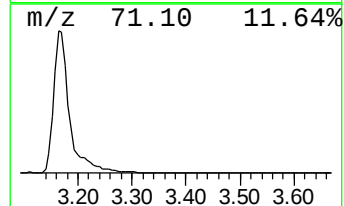
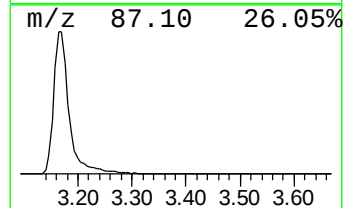
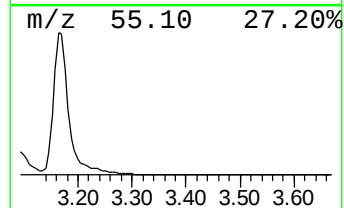
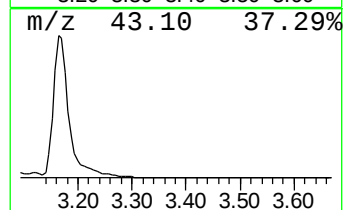
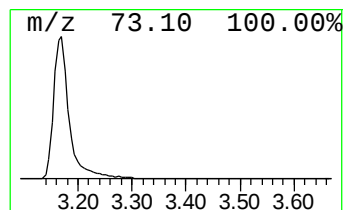
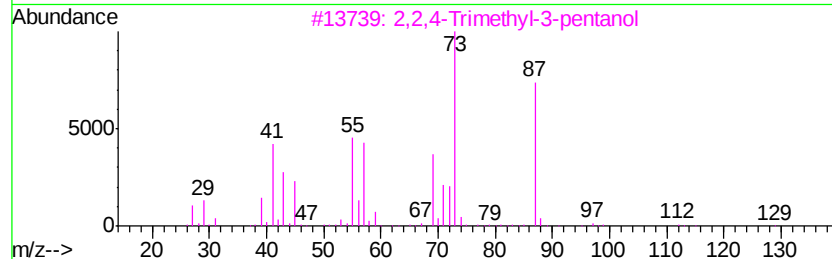
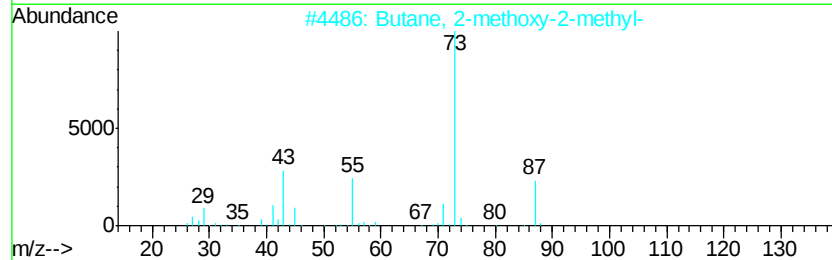
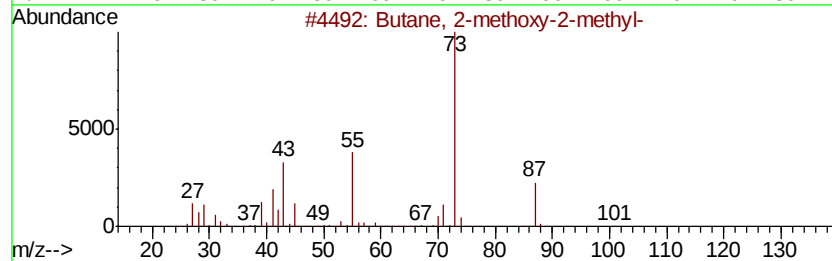
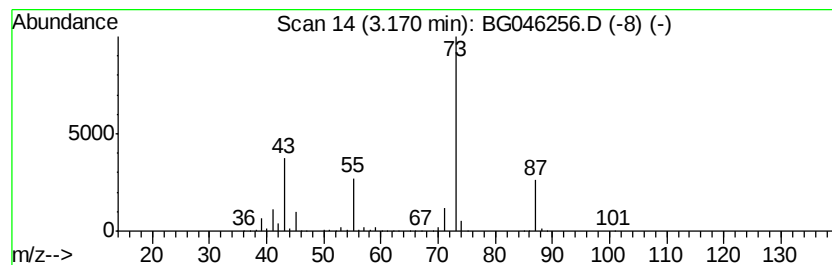
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG081320PAH.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 1

| R.T. | EstConc      | Area    | Relative to ISTD       | R.T. |
|------|--------------|---------|------------------------|------|
| 3.17 | 116.16 ng/ul | 3205250 | 1,4-Dichlorobenzene-d4 | 7.95 |

| Hit# | of | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------------|-----|---------|-------------|------|
| 1    | 5  | Butane, 2-methoxy-2-methyl- | 102 | C6H14O  | 000994-05-8 | 83   |
| 2    |    | Butane, 2-methoxy-2-methyl- | 102 | C6H14O  | 000994-05-8 | 78   |
| 3    |    | 2,2,4-Trimethyl-3-pentanol  | 130 | C8H18O  | 005162-48-1 | 40   |
| 4    |    | 2,2,4-Trimethyl-3-pentanol  | 130 | C8H18O  | 005162-48-1 | 40   |
| 5    |    | 1,3-Dioxolane, 2-methyl-    | 88  | C4H8O2  | 000497-26-7 | 25   |



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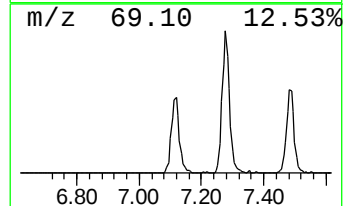
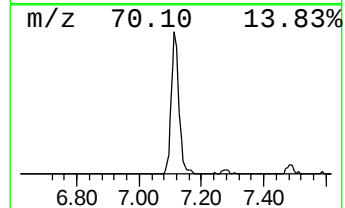
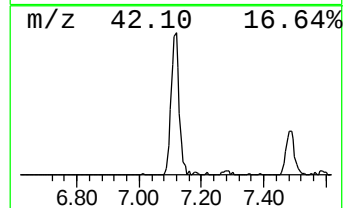
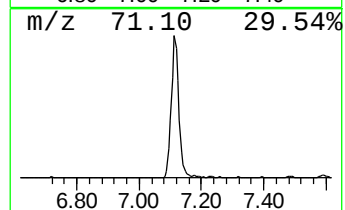
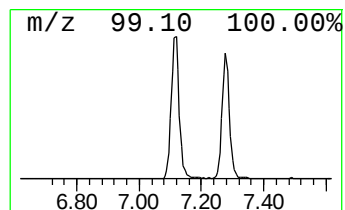
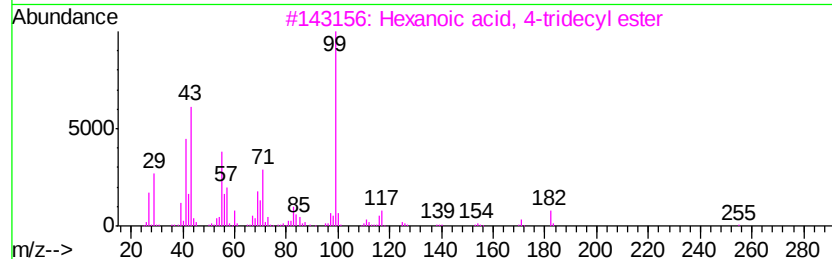
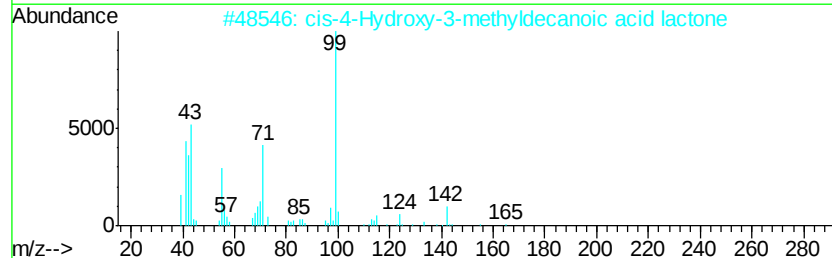
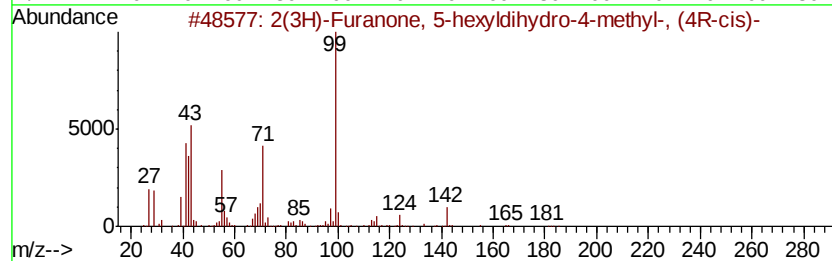
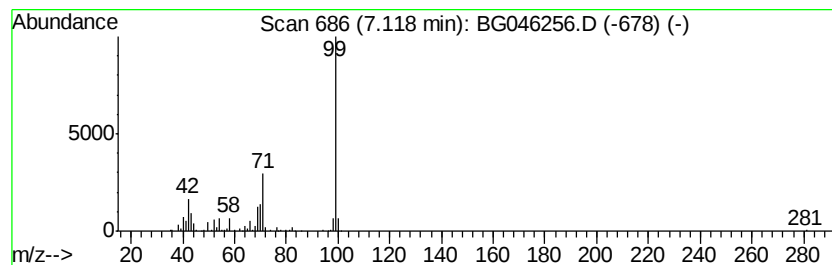
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 2 2(3H)-Furanone, 5-hexyldihy... Concentration Rank 4

| R.T. | EstConc     | Area   | Relative to ISTD       | R.T. |
|------|-------------|--------|------------------------|------|
| 7.12 | 13.95 ng/ul | 384904 | 1,4-Dichlorobenzene-d4 | 7.95 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|-------------------------------------|-----|----------|--------------|------|
| 1    | 5  | 2(3H)-Furanone, 5-hexyldihydro-4... | 184 | C11H20O2 | 121644-12-0  | 64   |
| 2    |    | cis-4-Hydroxy-3-methyldecanoic a... | 184 | C11H20O2 | 147254-32-8  | 64   |
| 3    |    | Hexanoic acid, 4-tridecyl ester     | 298 | C19H38O2 | 1000279-29-3 | 64   |
| 4    |    | Hexanoic acid, 5-tridecyl ester     | 298 | C19H38O2 | 1000279-29-4 | 56   |
| 5    |    | Hexanoic acid, 3-tridecyl ester     | 298 | C19H38O2 | 1000279-29-2 | 56   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081320\  
Data File : BG046256.D  
Acq On : 13 Aug 2020 19:42  
Operator : CG/JU  
Sample : L3629-14  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BFM61

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Quant Title : SVOA CALIBRATION

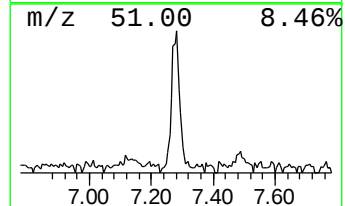
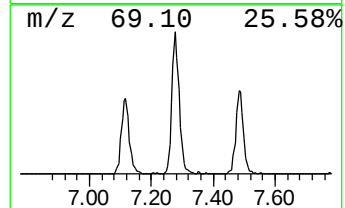
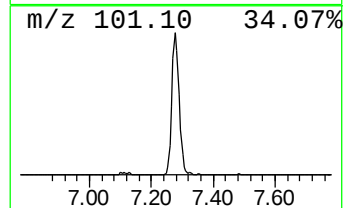
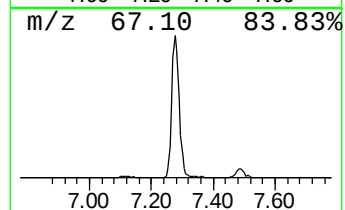
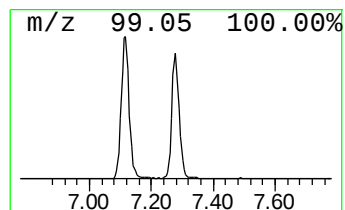
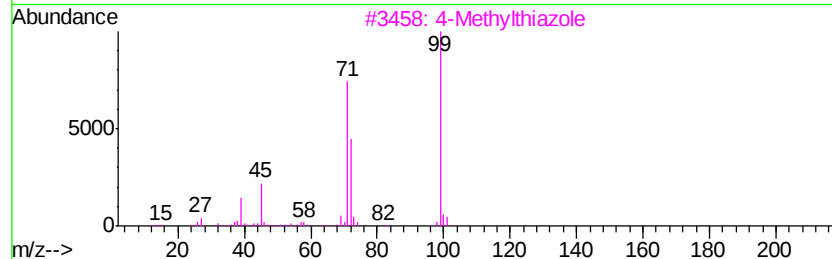
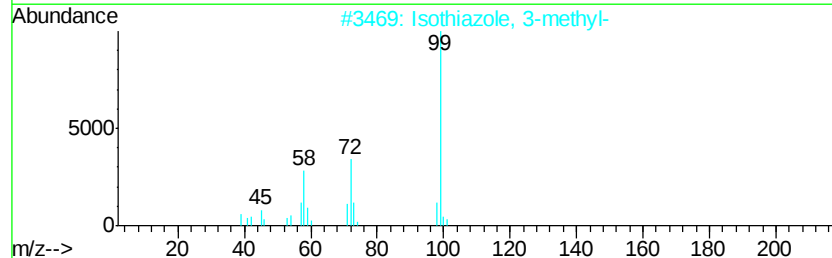
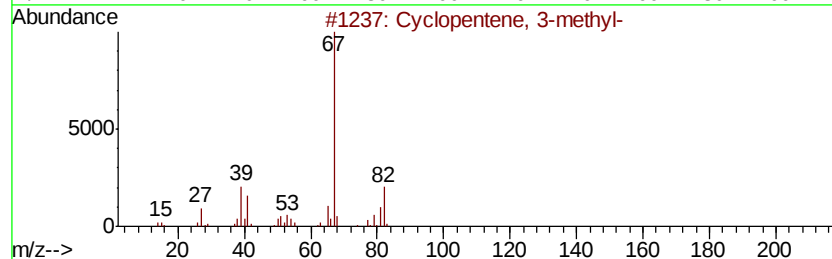
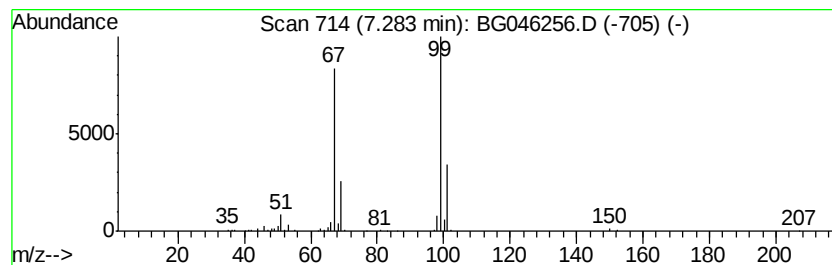
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 3 unknown-01 Concentration Rank 7

| R.T. | EstConc     | Area   | Relative to ISTD       | R.T. |
|------|-------------|--------|------------------------|------|
| 7.28 | 11.45 ng/ul | 315960 | 1,4-Dichlorobenzene-d4 | 7.95 |

| Hit# of | 5 | Tentative ID               | MW | MolForm | CAS#        | Qual |
|---------|---|----------------------------|----|---------|-------------|------|
| 1       |   | Cyclopentene, 3-methyl-    | 82 | C6H10   | 001120-62-3 | 12   |
| 2       |   | Isothiazole, 3-methyl-     | 99 | C4H5NS  | 000693-92-5 | 9    |
| 3       |   | 4-Methylthiazole           | 99 | C4H5NS  | 000693-95-8 | 9    |
| 4       |   | 3,5-Diamino-1,2,4-triazole | 99 | C2H5N5  | 001455-77-2 | 9    |
| 5       |   | Cyclopentene, 1-methyl-    | 82 | C6H10   | 000693-89-0 | 9    |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081320\  
Data File : BG046256.D  
Acq On : 13 Aug 2020 19:42  
Operator : CG/JU  
Sample : L3629-14  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BFM61

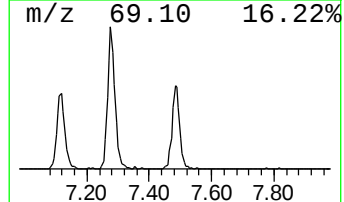
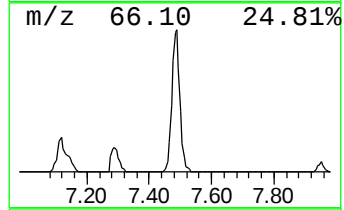
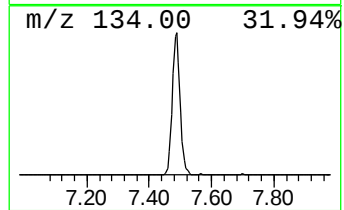
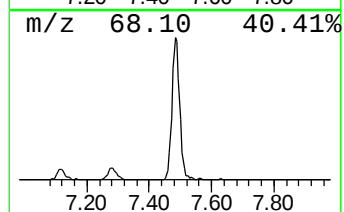
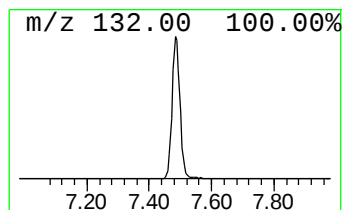
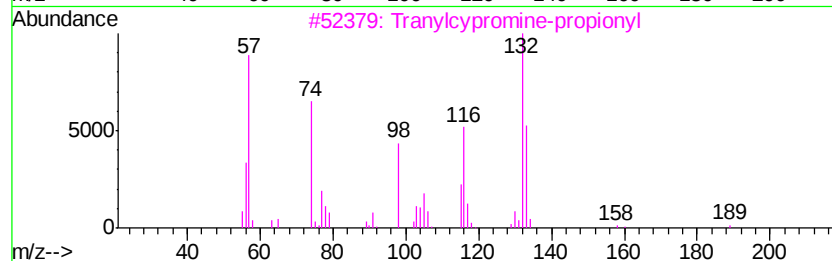
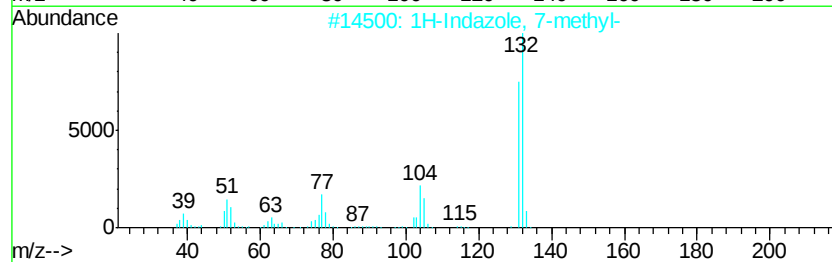
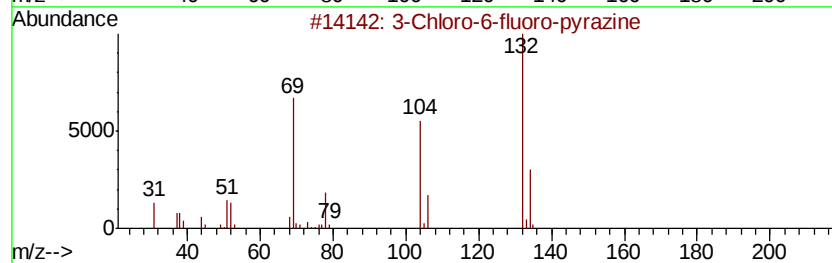
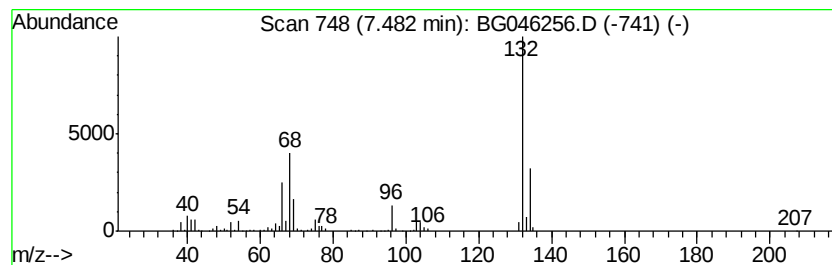
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Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 4 unknown-02 Concentration Rank 3

| R.T. | EstConc     | Area   | Relative to ISTD       | R.T. |
|------|-------------|--------|------------------------|------|
| 7.48 | 14.05 ng/ul | 387720 | 1,4-Dichlorobenzene-d4 | 7.95 |

| Hit# | of | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|-------------------------------------|-----|------------|--------------|------|
| 1    | 5  | 3-Chloro-6-fluoro-pyrazine          | 132 | C4H2ClFN2  | 1000146-10-7 | 27   |
| 2    |    | 1H-Indazole, 7-methyl-              | 132 | C8H8N2     | 1000316-00-7 | 25   |
| 3    |    | Tranvlcypropine-propionyl           | 189 | C12H15NO   | 1000123-86-3 | 12   |
| 4    |    | 3H-Benzo[4,5]imidazo[1,2-c]oxazo... | 218 | C12H14N2O2 | 1000316-99-9 | 10   |
| 5    |    | o-(Acetonylamino)benzoic acid       | 193 | C10H11NO3  | 037144-42-6  | 10   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081320\  
Data File : BG046256.D  
Acq On : 13 Aug 2020 19:42  
Operator : CG/JU  
Sample : L3629-14  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BFM61

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG081320PAH.M  
Quant Title : SVOA CALIBRATION

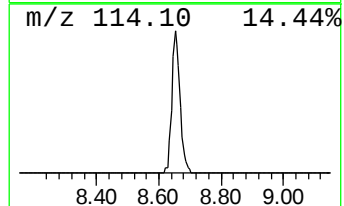
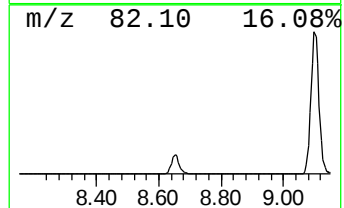
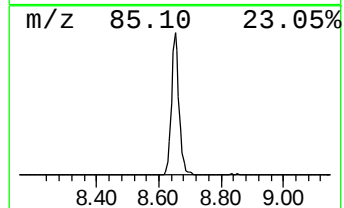
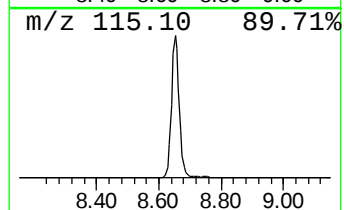
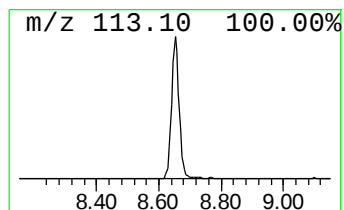
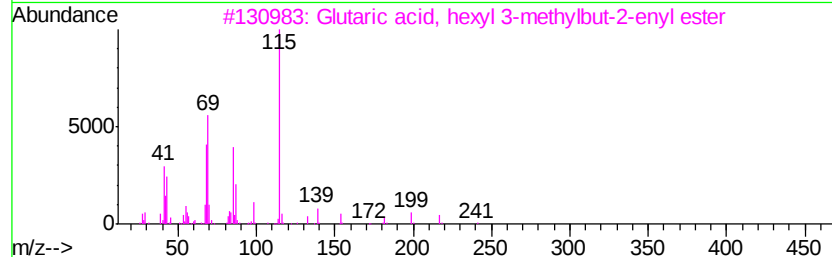
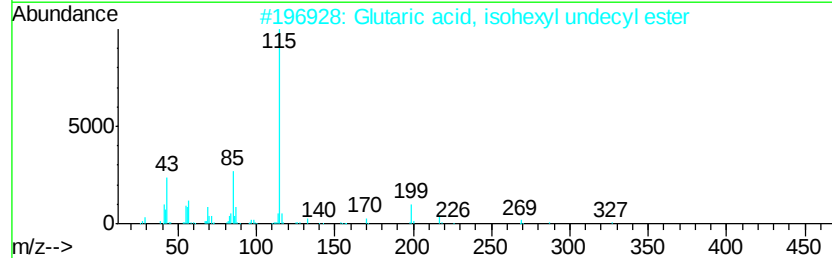
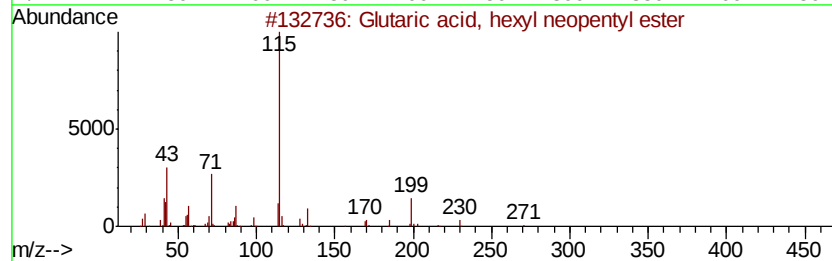
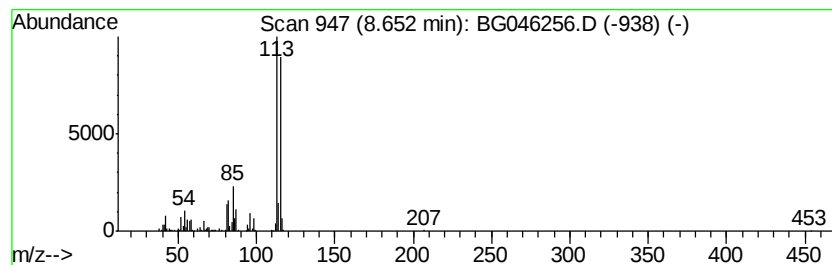
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 5 unknown-03 Concentration Rank 2

| R.T. | EstConc     | Area   | Relative to ISTD       | R.T. |
|------|-------------|--------|------------------------|------|
| 8.65 | 14.90 ng/ul | 411077 | 1,4-Dichlorobenzene-d4 | 7.95 |

| Hit# of | 5                                   | Tentative ID | MW       | MolForm      | CAS# | Qual |
|---------|-------------------------------------|--------------|----------|--------------|------|------|
| 1       | Glutaric acid, hexyl neopentyl e... | 286          | C16H30O4 | 1000358-34-4 | 38   |      |
| 2       | Glutaric acid, isohexyl undecyl ... | 370          | C22H42O4 | 1000358-32-1 | 38   |      |
| 3       | Glutaric acid, hexyl 3-methylbut... | 284          | C16H28O4 | 1000360-09-0 | 38   |      |
| 4       | Glutaric acid, isobutyl neopentv... | 258          | C14H26O4 | 1000358-34-1 | 35   |      |
| 5       | Glutaric acid, isobutyl octyl ester | 300          | C17H32O4 | 1000358-25-6 | 35   |      |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081320\  
Data File : BG046256.D  
Acq On : 13 Aug 2020 19:42  
Operator : CG/JU  
Sample : L3629-14  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BFM61

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Quant Title : SVOA CALIBRATION

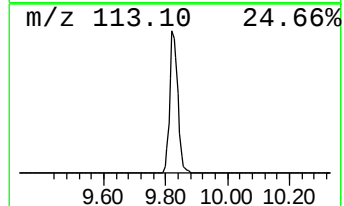
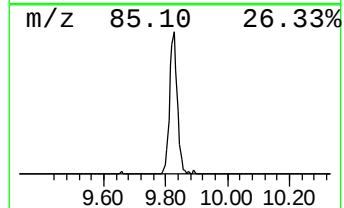
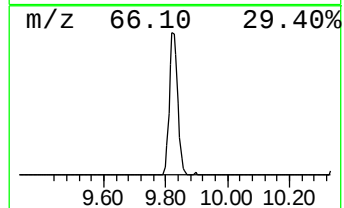
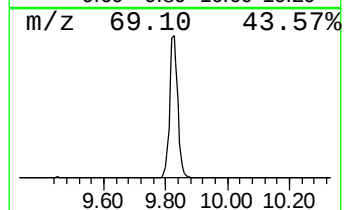
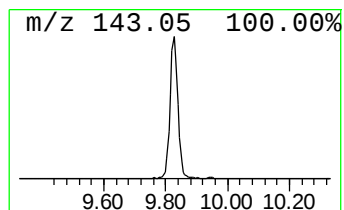
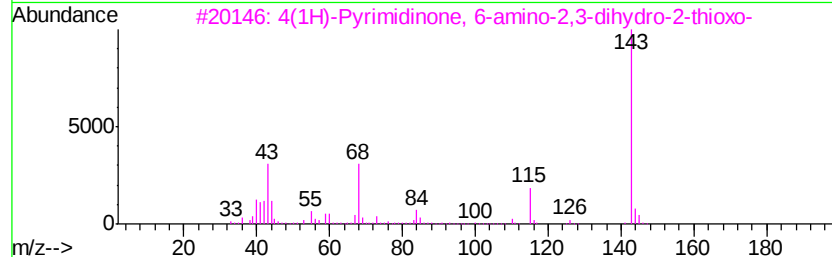
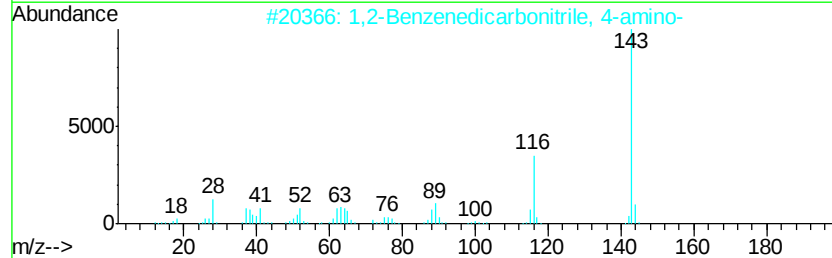
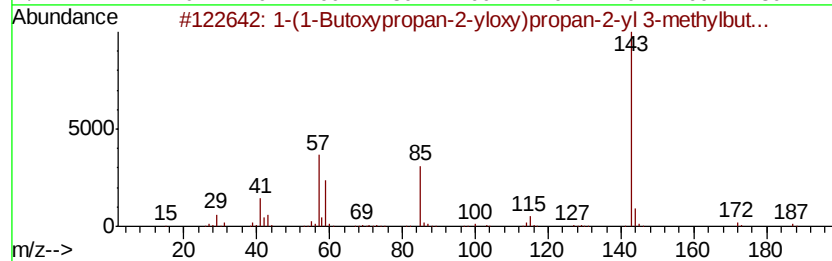
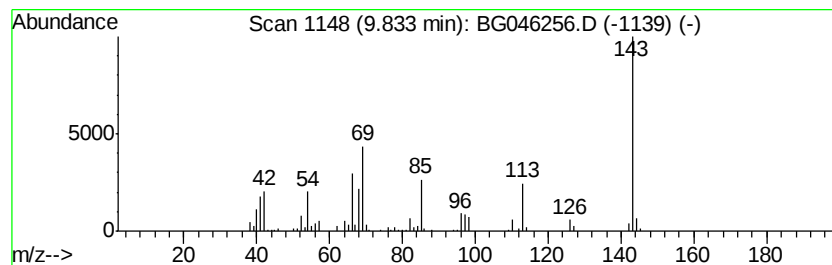
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TIC Integration Parameters: LSCINT.P

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Peak Number 7 unknown-04 Concentration Rank 9

| R.T. | EstConc    | Area   | Relative to ISTD | R.T.  |
|------|------------|--------|------------------|-------|
| 9.83 | 7.43 ng/ul | 306446 | Naphthalene-d8   | 10.75 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|-------------------------------------|-----|----------|--------------|------|
| 1    | 5  | 1-(1-Butoxypropan-2-yloxy)propan... | 274 | C15H30O4 | 1000367-12-9 | 22   |
| 2    |    | 1,2-Benzenedicarbonitrile, 4-amino- | 143 | C8H5N3   | 056765-79-8  | 14   |
| 3    |    | 4(1H)-Pyrimidinone, 6-amino-2,3-... | 143 | C4H5N3OS | 001004-40-6  | 12   |
| 4    |    | 3-Aminophthalonitrile               | 143 | C8H5N3   | 058632-96-5  | 10   |
| 5    |    | 6-Aza-2-thiothymine                 | 143 | C4H5N3OS | 000615-76-9  | 10   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081320\  
Data File : BG046256.D  
Acq On : 13 Aug 2020 19:42  
Operator : CG/JU  
Sample : L3629-14  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BFM61

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG081320PAH.M  
Quant Title : SVOA CALIBRATION

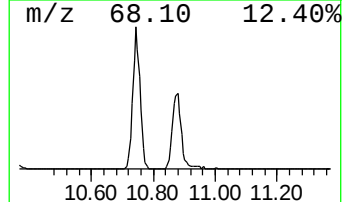
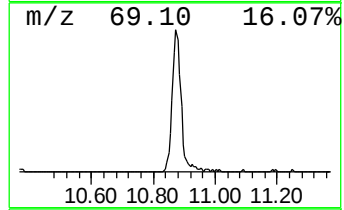
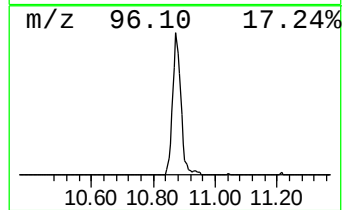
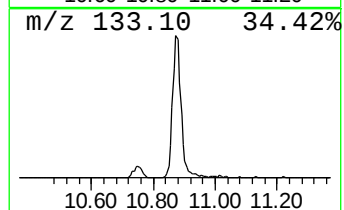
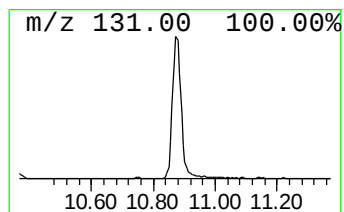
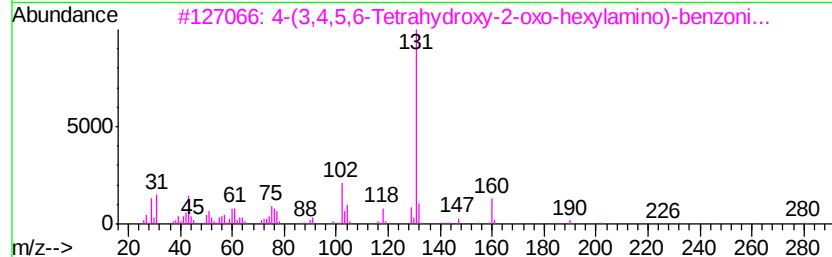
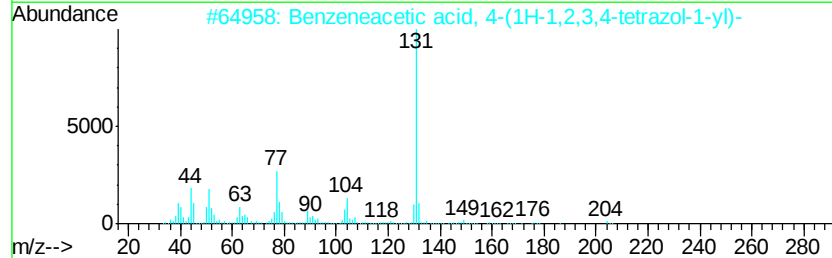
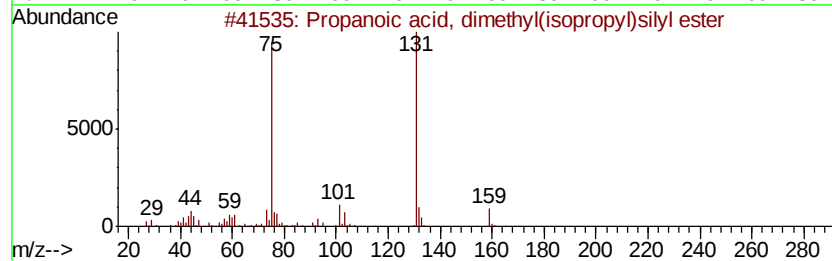
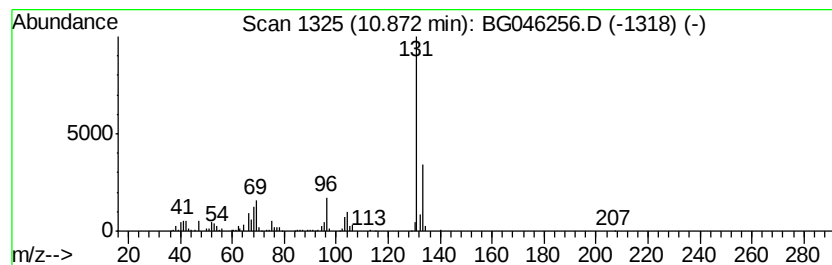
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 8 unknown-05 Concentration Rank 8

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 10.87 | 9.17 ng/ul | 378500 | Napthalene-d8    | 10.75 |

| Hit# | of | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|-------------------------------------|-----|------------|--------------|------|
| 1    | 5  | Propanoic acid, dimethyl(isoprop... | 174 | C8H18O2Si  | 1000279-45-8 | 33   |
| 2    |    | Benzeneacetic acid, 4-(1H-1,2,3,... | 204 | C9H8N4O2   | 1000351-56-5 | 33   |
| 3    |    | 4-(3,4,5,6-Tetrahydroxy-2-oxo-he... | 280 | C13H16N2O5 | 1000189-82-8 | 33   |
| 4    |    | 3-Hydroxy-2-p-tolyl-2-butenenitrile | 173 | C11H11NO   | 1000243-87-8 | 33   |
| 5    |    | Benzenepropanoic acid, 4-(1H-1,2... | 218 | C10H10N4O2 | 1000349-98-4 | 33   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081320\  
Data File : BG046256.D  
Acq On : 13 Aug 2020 19:42  
Operator : CG/JU  
Sample : L3629-14  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BFM61

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG081320PAH.M  
Quant Title : SVOA CALIBRATION

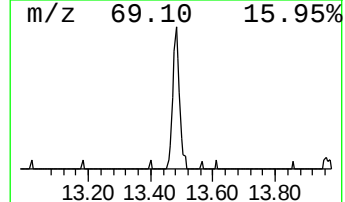
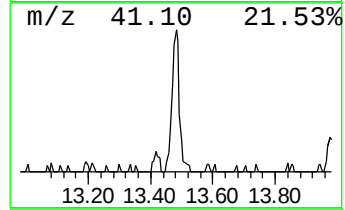
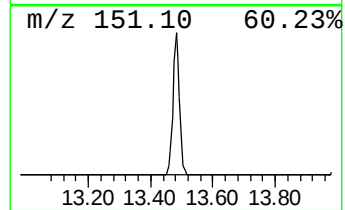
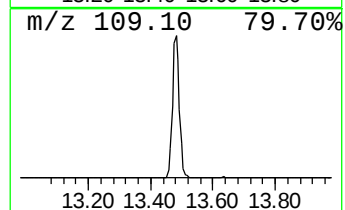
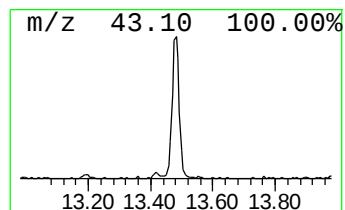
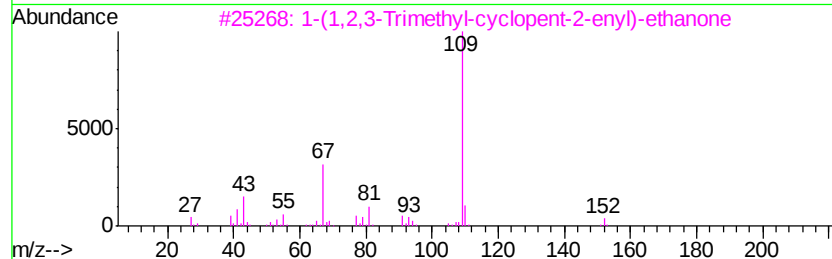
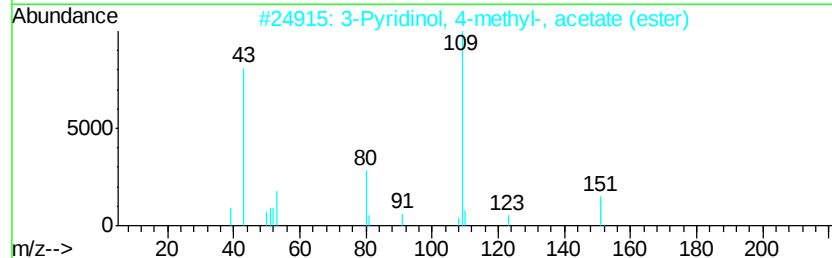
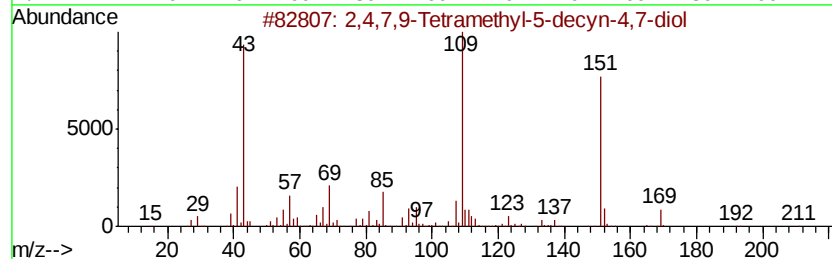
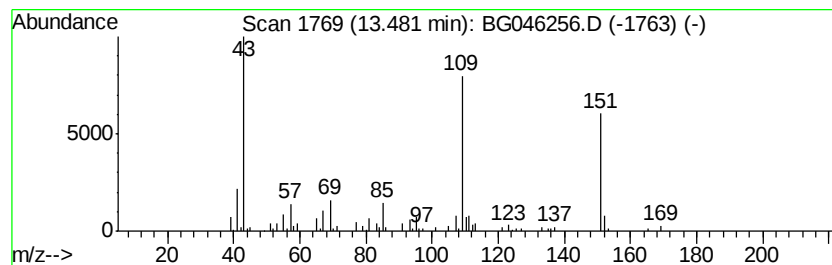
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 9 2,4,7,9-Tetramethyl-5-decyn... Concentration Rank 18

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 13.48 | 2.13 ng/ul | 122723 | Acenaphthene-d10 | 14.57 |

| Hit# | of                                    | 5   | Tentative ID | MW          | MolForm | CAS# | Qual |
|------|---------------------------------------|-----|--------------|-------------|---------|------|------|
| 1    | 2,4,7,9-Tetramethyl-5-decyn-4,7-...   | 226 | C14H26O2     | 000126-86-3 | 91      |      |      |
| 2    | 3-Pyridinol, 4-methyl-, acetate ...   | 151 | C8H9NO2      | 001006-96-8 | 64      |      |      |
| 3    | 1-(1,2,3-Trimethyl-cyclopent-2-en-... | 152 | C10H16O      | 070987-81-4 | 43      |      |      |
| 4    | 3,3,5,5-Tetramethylcyclopentene       | 124 | C9H16        | 038667-10-6 | 38      |      |      |
| 5    | 3,6-Dimethyl-6-hepten-4-yn-3-ol       | 138 | C9H14O       | 003601-67-0 | 38      |      |      |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081320\  
Data File : BG046256.D  
Acq On : 13 Aug 2020 19:42  
Operator : CG/JU  
Sample : L3629-14  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BFM61

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG081320PAH.M  
Quant Title : SVOA CALIBRATION

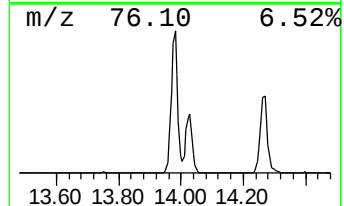
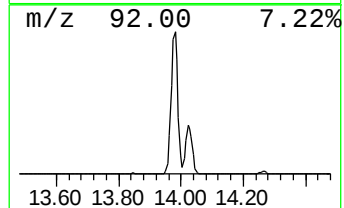
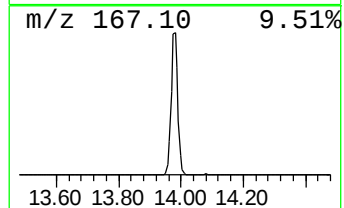
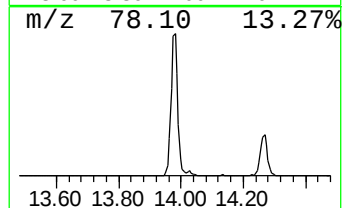
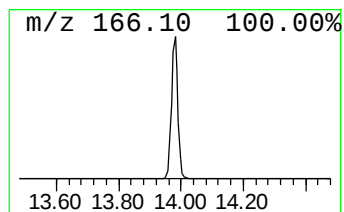
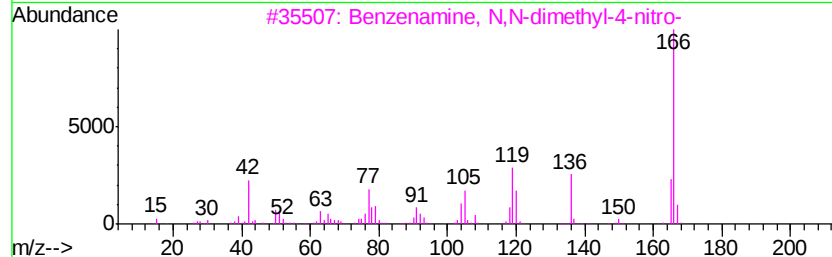
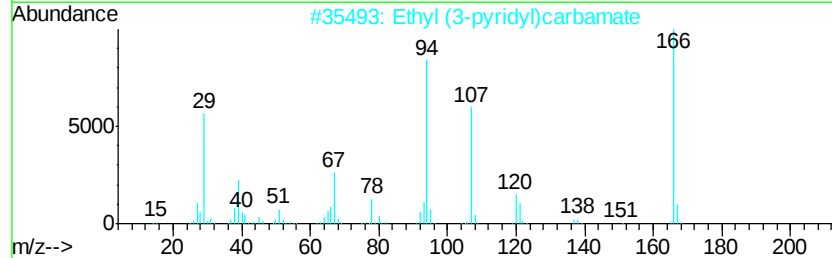
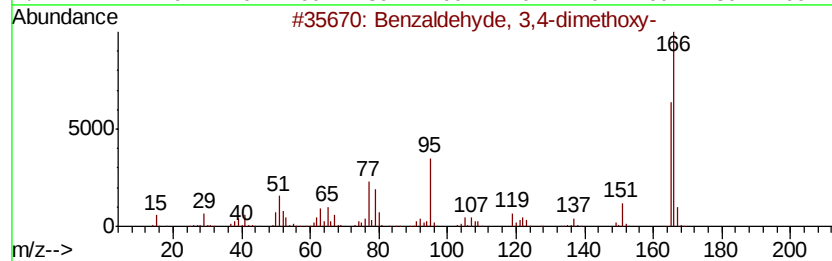
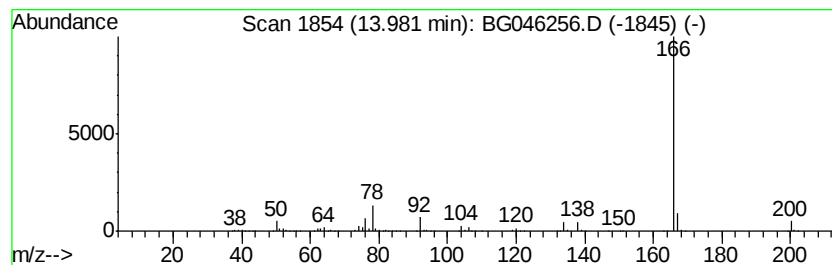
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 10 unknown-06 Concentration Rank 5

| R.T.  | EstConc     | Area   | Relative to ISTD | R.T.  |
|-------|-------------|--------|------------------|-------|
| 13.98 | 13.94 ng/ul | 805297 | Acenaphthene-d10 | 14.57 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#        | Qual |
|------|----|-------------------------------------|-----|-----------|-------------|------|
| 1    | 5  | Benzaldehyde, 3,4-dimethoxy-        | 166 | C9H10O3   | 000120-14-9 | 45   |
| 2    |    | Ethyl (3-pyridyl)carbamate          | 166 | C8H10N2O2 | 006276-11-5 | 39   |
| 3    |    | Benzenamine, N,N-dimethyl-4-nitro-  | 166 | C8H10N2O2 | 000100-23-2 | 38   |
| 4    |    | 6H-Purine-6-thione, 1,7-dihydro-... | 166 | C6H6N4S   | 001126-23-4 | 9    |
| 5    |    | 1,3-Benzodioxole-5-carboxylic acid  | 166 | C8H6O4    | 000094-53-1 | 9    |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081320\  
Data File : BG046256.D  
Acq On : 13 Aug 2020 19:42  
Operator : CG/JU  
Sample : L3629-14  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

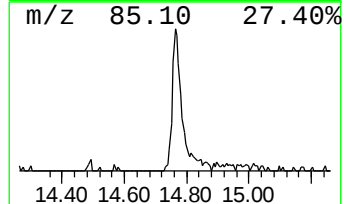
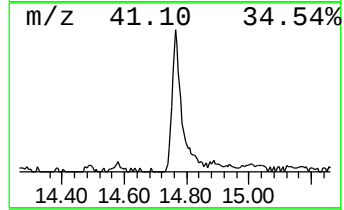
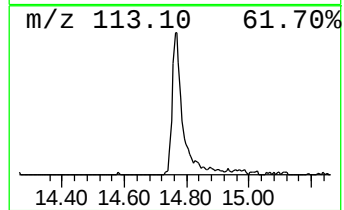
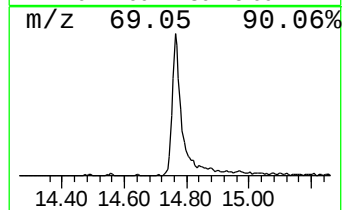
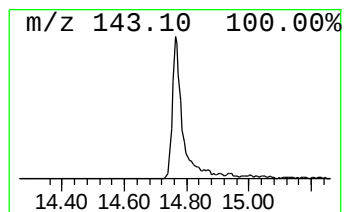
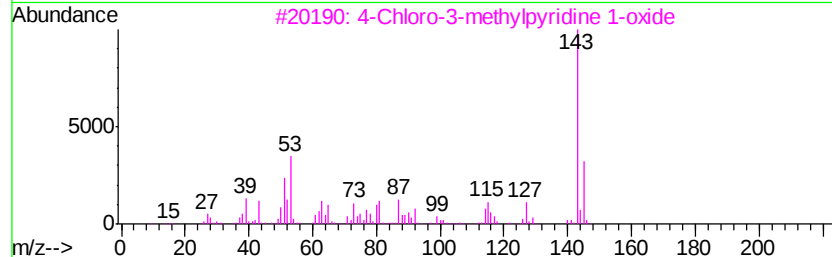
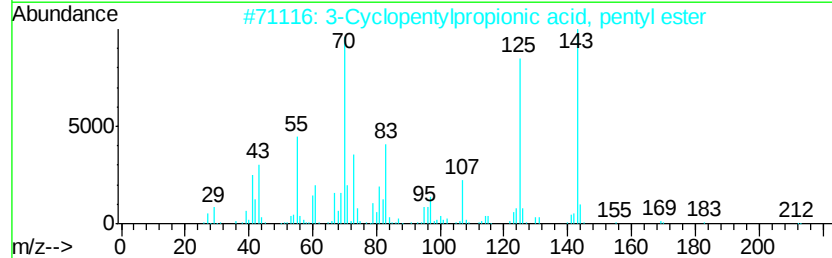
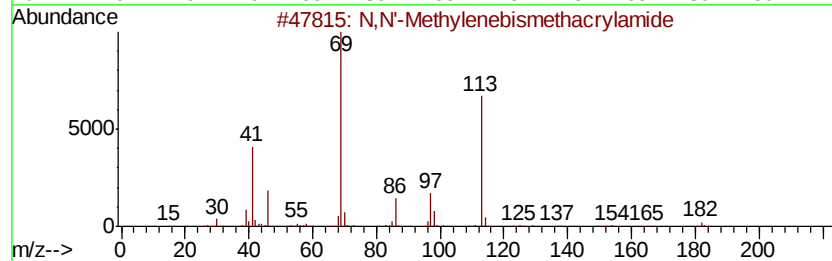
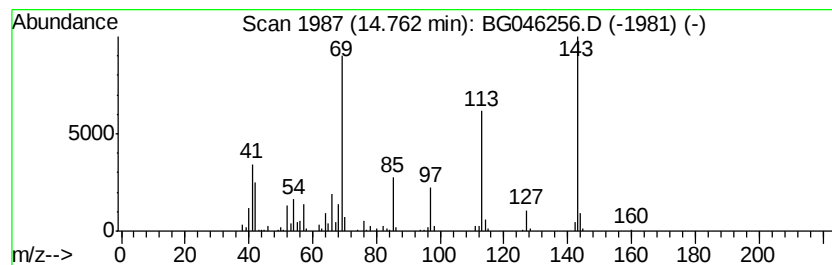
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BNA\_G  
ClientSampleId :  
BFM61

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG081320PAH.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 12 unknown-07 Concentration Rank 12

| R.T.    | EstConc    | Area                                | Relative to ISTD | R.T.            |
|---------|------------|-------------------------------------|------------------|-----------------|
| 14.76   | 4.25 ng/ul | 245679                              | Acenaphthene-d10 | 14.57           |
| Hit# of | 5          | Tentative ID                        | MW MolForm       | CAS# Qual       |
| 1       |            | N,N'-Methylenebismethacrylamide     | 182 C9H14N2O2    | 002359-15-1 38  |
| 2       |            | 3-Cyclopentylpropionic acid, pen... | 212 C13H24O2     | 1000292-33-2 22 |
| 3       |            | 4-Chloro-3-methylpyridine 1-oxide   | 143 C6H6ClNO     | 001073-34-3 22  |
| 4       |            | 4-Amino-2-chloro-6-methylpyrimidine | 143 C5H6ClN3     | 014394-60-6 22  |
| 5       |            | Glutaric acid, ethyl 4-methylhep... | 272 C15H28O4     | 1000377-14-9 22 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081320\  
Data File : BG046256.D  
Acq On : 13 Aug 2020 19:42  
Operator : CG/JU  
Sample : L3629-14  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BFM61

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG081320PAH.M  
Quant Title : SVOA CALIBRATION

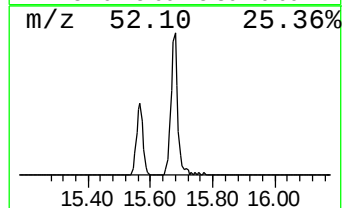
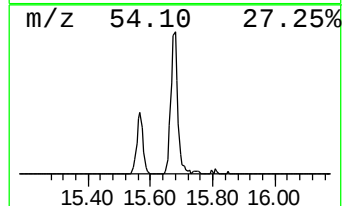
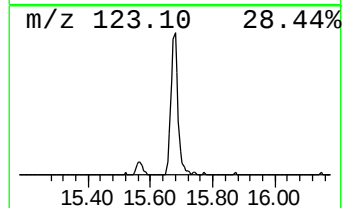
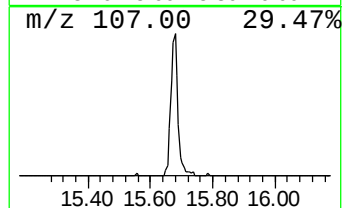
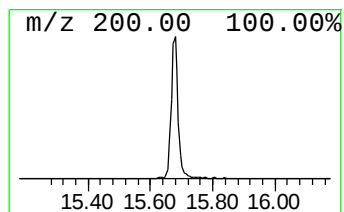
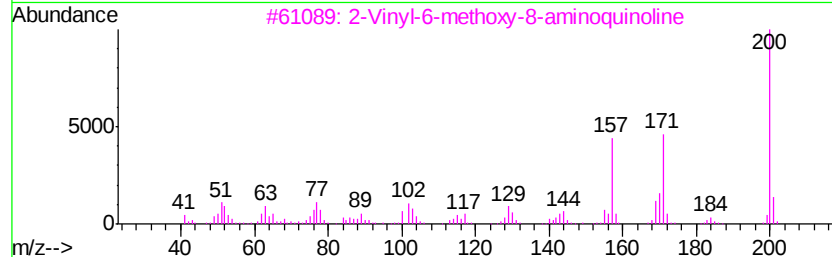
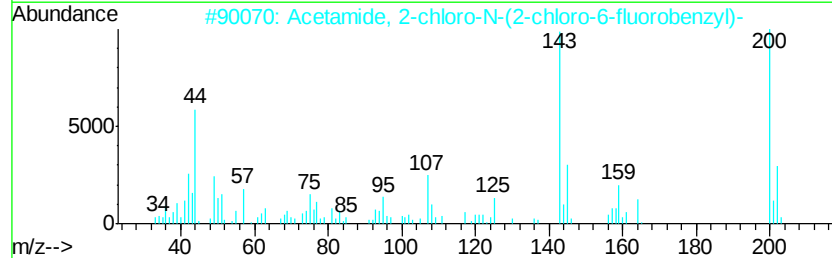
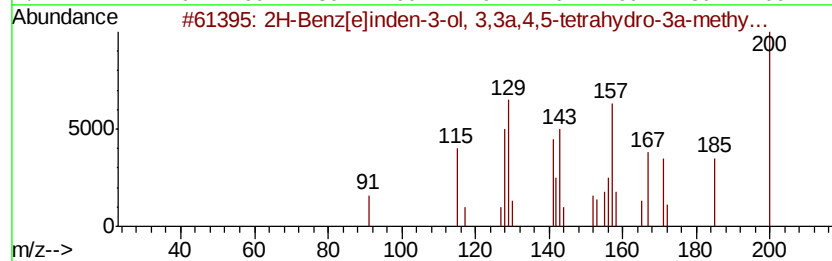
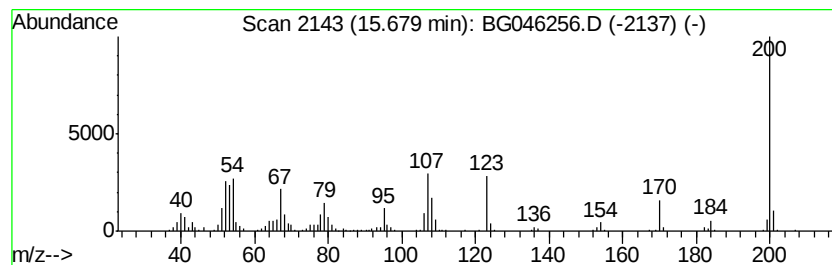
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 13 unknown-08 Concentration Rank 10

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 15.68 | 5.28 ng/ul | 304692 | Acenaphthene-d10 | 14.57 |

| Hit# | of | 5 | Tentative ID                         | MW  | MolForm    | CAS#         | Qual |
|------|----|---|--------------------------------------|-----|------------|--------------|------|
| 1    |    |   | 2H-Benz[e]indene-3-ol, 3,3a,4,5-t... | 200 | C14H16O    | 071805-91-9  | 38   |
| 2    |    |   | Acetamide, 2-chloro-N-(2-chloro-...  | 235 | C9H8Cl2FNO | 1000268-05-5 | 32   |
| 3    |    |   | 2-Vinyl-6-methoxy-8-aminoquinoline   | 200 | C12H12N2O  | 064992-65-0  | 16   |
| 4    |    |   | D-Norleucine, N-neopentylloxycarb... | 455 | C27H53NO4  | 1000344-14-0 | 14   |
| 5    |    |   | Benzenamine, 3-(4-aminophenoxy)-     | 200 | C12H12N2O  | 002657-87-6  | 12   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081320\  
Data File : BG046256.D  
Acq On : 13 Aug 2020 19:42  
Operator : CG/JU  
Sample : L3629-14  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BFM61

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG081320PAH.M  
Quant Title : SVOA CALIBRATION

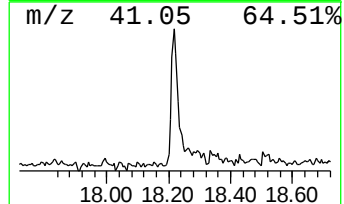
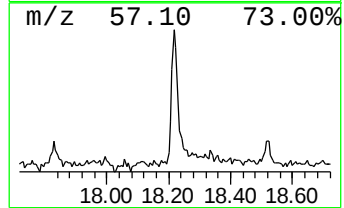
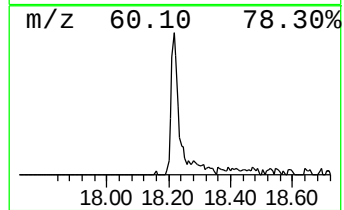
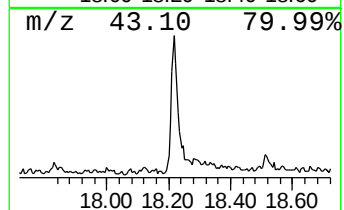
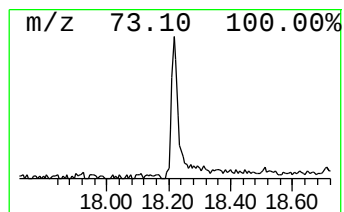
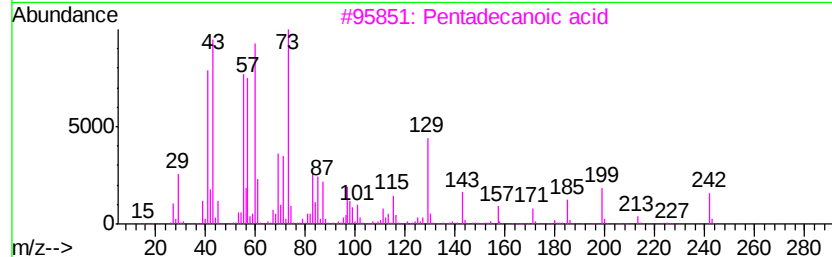
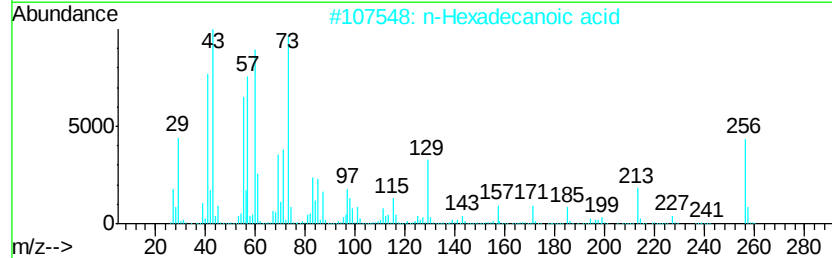
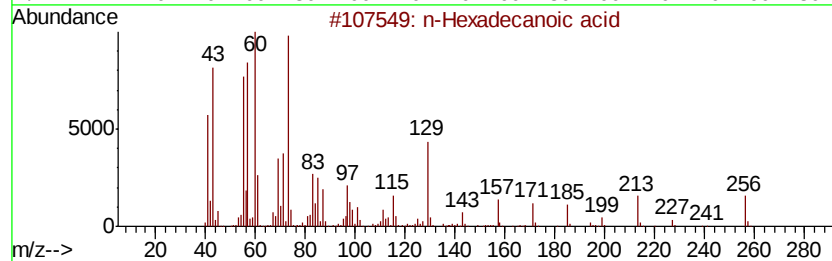
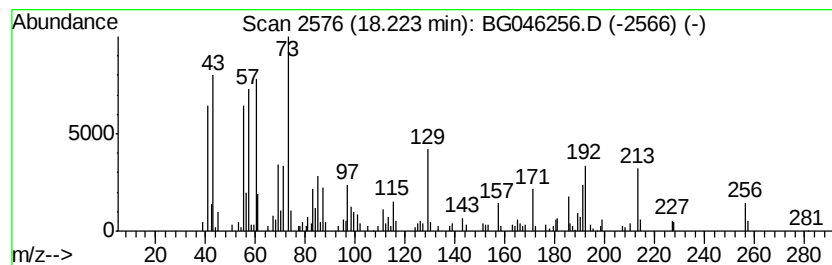
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 14 n-Hexadecanoic acid Concentration Rank 17

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 18.22 | 2.58 ng/ul | 180592 | Phenanthrene-d10 | 17.32 |

| 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    |    |   | Pentadecanoic acid  | 242 | C15H30O2 | 001002-84-2 | 83   |
| 4    |    |   | Dodecanoic acid     | 200 | C12H24O2 | 000143-07-7 | 81   |
| 5    |    |   | Tridecanoic acid    | 214 | C13H26O2 | 000638-53-9 | 70   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081320\  
Data File : BG046256.D  
Acq On : 13 Aug 2020 19:42  
Operator : CG/JU  
Sample : L3629-14  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BFM61

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG081320PAH.M  
Quant Title : SVOA CALIBRATION

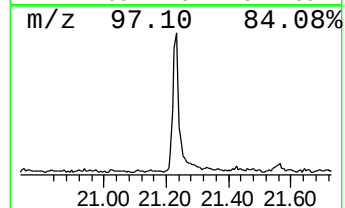
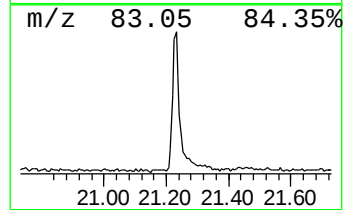
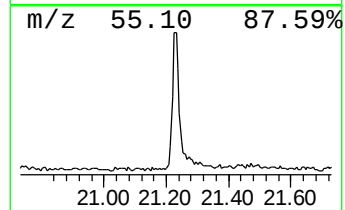
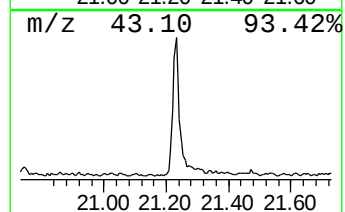
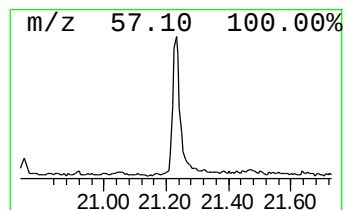
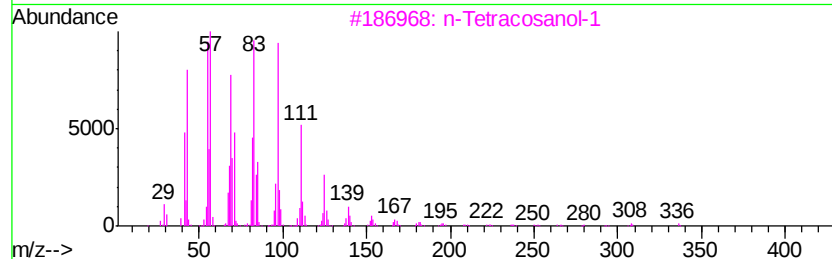
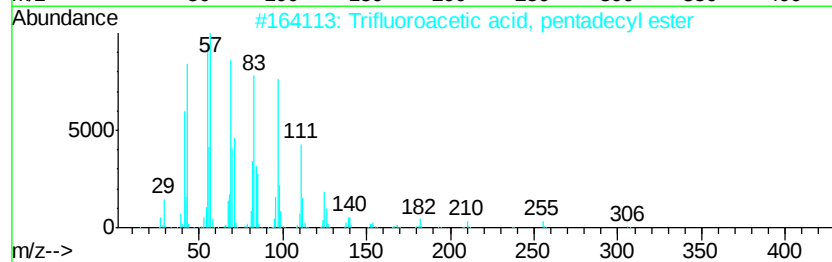
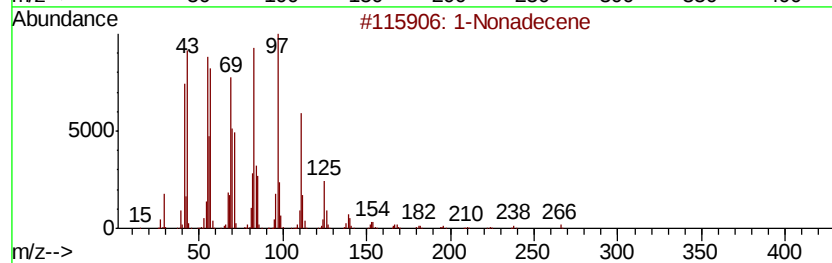
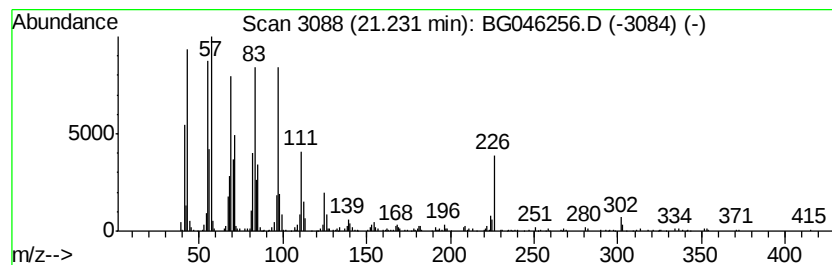
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 15 (DEL) Alkane: Straight-Chai... Concentration Rank 13

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 21.23 | 4.09 ng/ul | 374365 | Chrysene-d12     | 21.56 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm    | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|------------|-------------|------|
| 1    |    |   | 1-Nonadecene                        | 266 | C19H38     | 018435-45-5 | 96   |
| 2    |    |   | Trifluoroacetic acid, pentadecyl... | 324 | C17H31F3O2 | 959010-23-2 | 93   |
| 3    |    |   | n-Tetracosanol-1                    | 354 | C24H50O    | 000506-51-4 | 93   |
| 4    |    |   | Heptafluorobutyric acid, penta...   | 424 | C19H31F7O2 | 959261-23-5 | 93   |
| 5    |    |   | Pentafluoropropionic acid, penta... | 374 | C18H31F5O2 | 959092-08-1 | 93   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081320\  
Data File : BG046256.D  
Acq On : 13 Aug 2020 19:42  
Operator : CG/JU  
Sample : L3629-14  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BFM61

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG081320PAH.M  
Quant Title : SVOA CALIBRATION

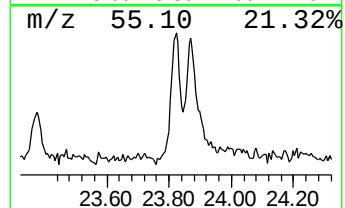
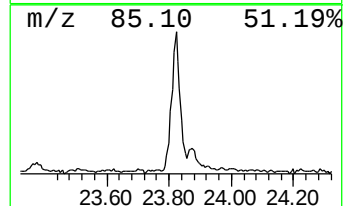
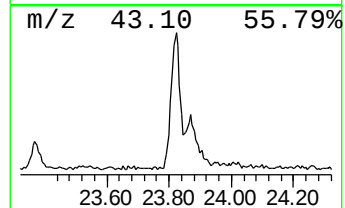
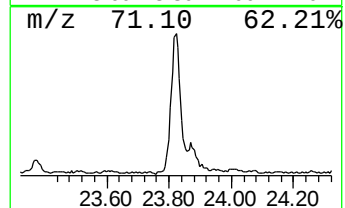
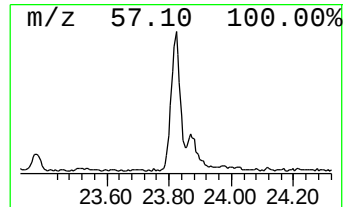
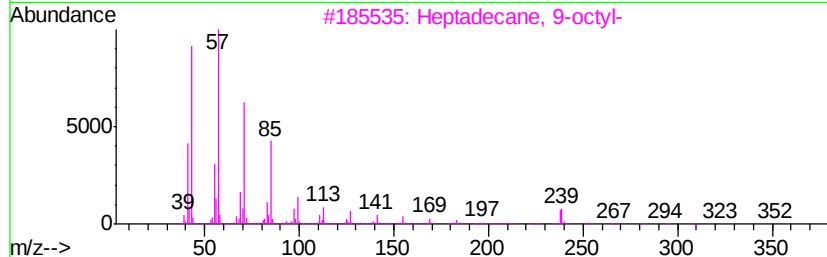
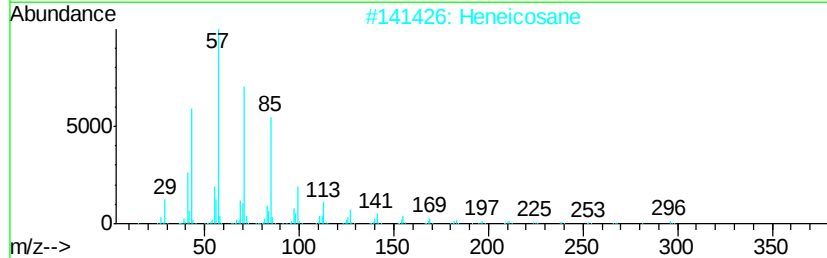
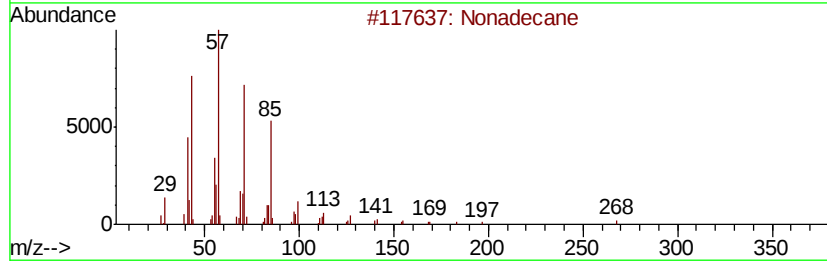
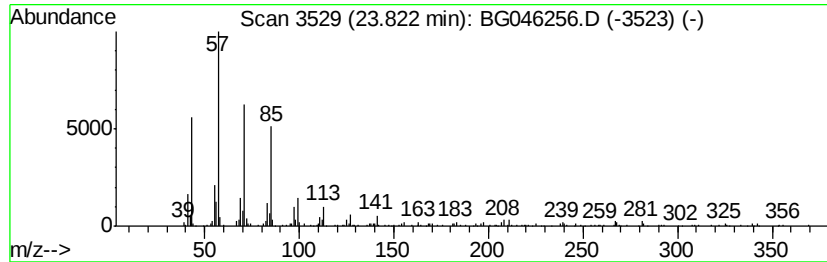
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 16 (DEL) Alkane: Straight-Chai... Concentration Rank 15

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 23.82 | 3.01 ng/ul | 235311 | Perylene-d12     | 24.67 |

| Hit# | of | 5 | Tentative ID          | MW  | MolForm | CAS#         | Qual |
|------|----|---|-----------------------|-----|---------|--------------|------|
| 1    |    |   | Nonadecane            | 268 | C19H40  | 000629-92-5  | 90   |
| 2    |    |   | Heneicosane           | 296 | C21H44  | 000629-94-7  | 87   |
| 3    |    |   | Heptadecane, 9-octyl- | 352 | C25H52  | 007225-64-1  | 87   |
| 4    |    |   | Docosane              | 310 | C22H46  | 000629-97-0  | 87   |
| 5    |    |   | 2-methyloctacosane    | 408 | C29H60  | 1000376-72-8 | 87   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081320\  
Data File : BG046256.D  
Acq On : 13 Aug 2020 19:42  
Operator : CG/JU  
Sample : L3629-14  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BFM61

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG081320PAH.M  
Quant Title : SVOA CALIBRATION

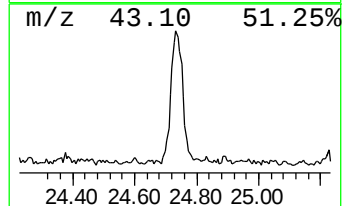
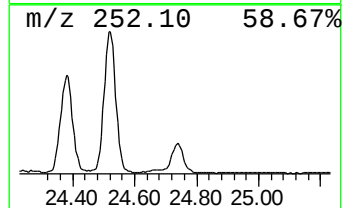
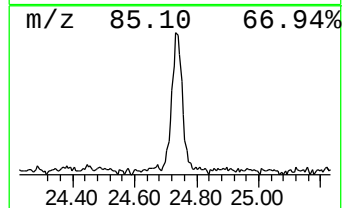
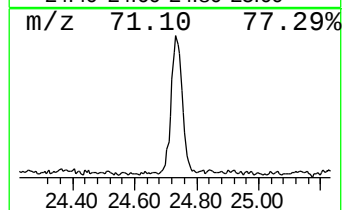
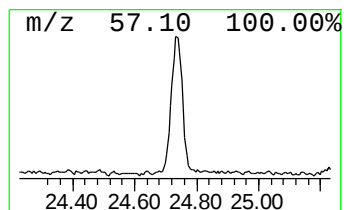
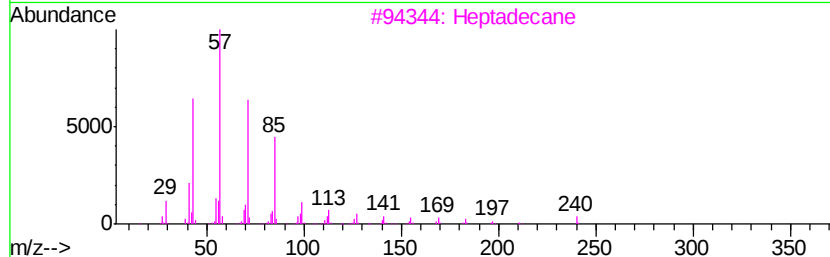
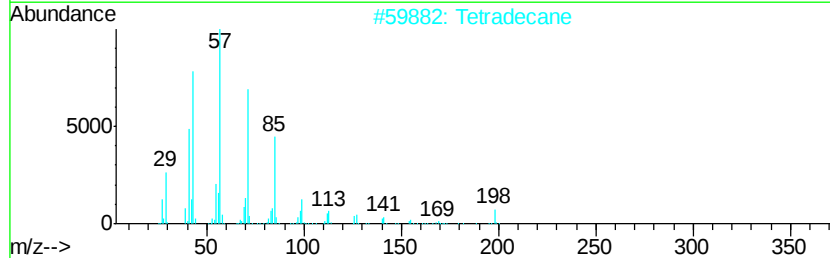
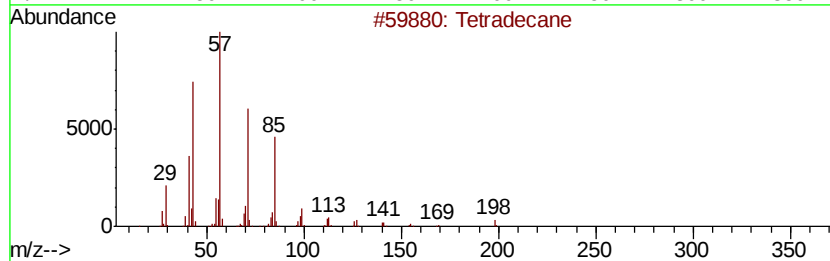
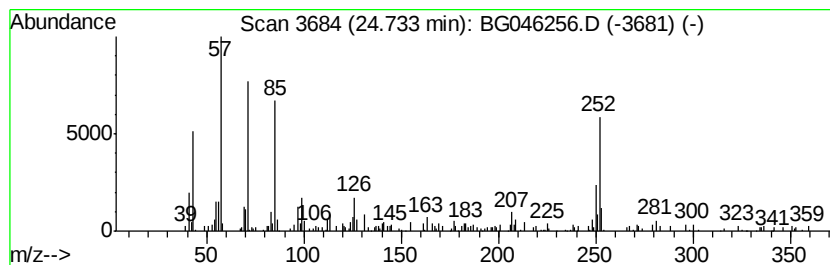
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 17 (DEL) Alkane: Straight-Chai... Concentration Rank 16

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 24.73 | 2.68 ng/ul | 209157 | Perylene-d12     | 24.67 |

| Hit# | of | 5 | Tentative ID             | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------|-----|---------|-------------|------|
| 1    |    |   | Tetradecane              | 198 | C14H30  | 000629-59-4 | 55   |
| 2    |    |   | Tetradecane              | 198 | C14H30  | 000629-59-4 | 49   |
| 3    |    |   | Heptadecane              | 240 | C17H36  | 000629-78-7 | 46   |
| 4    |    |   | Decane, 2,4,6-trimethyl- | 184 | C13H28  | 062108-27-4 | 46   |
| 5    |    |   | Octadecane               | 254 | C18H38  | 000593-45-3 | 46   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081320\  
Data File : BG046256.D  
Acq On : 13 Aug 2020 19:42  
Operator : CG/JU  
Sample : L3629-14  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BFM61

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG081320PAH.M  
Quant Title : SVOA CALIBRATION

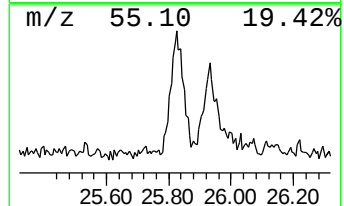
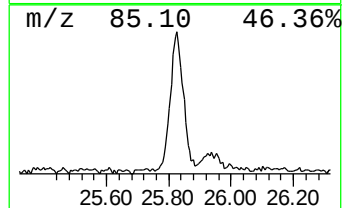
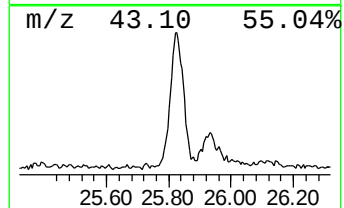
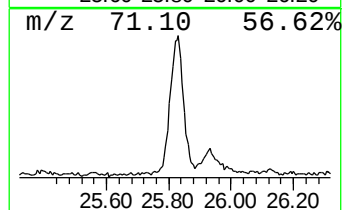
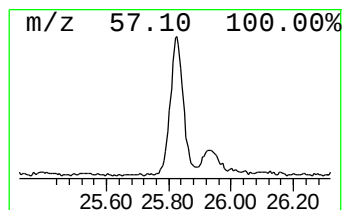
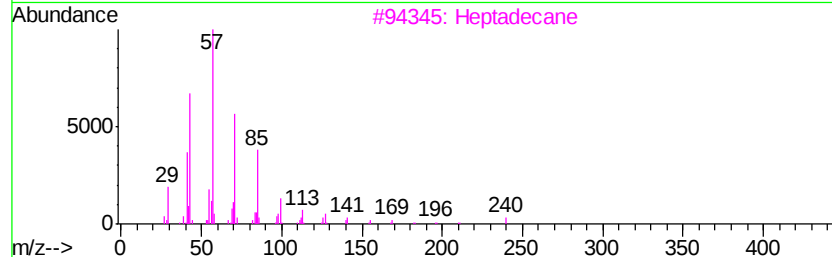
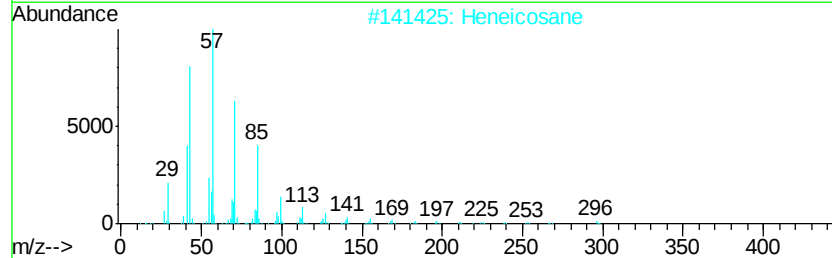
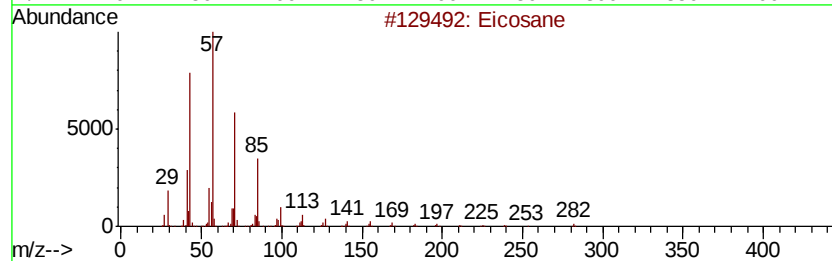
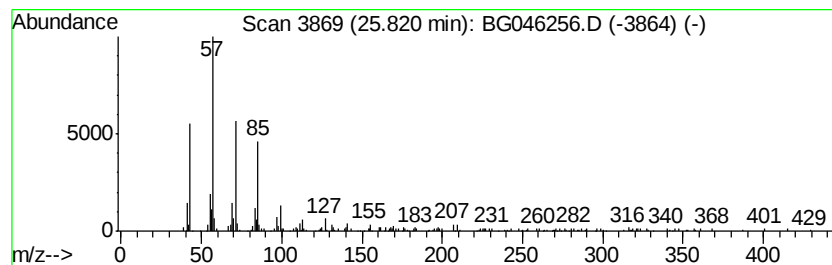
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 18 (DEL) Alkane: Straight-Chai... Concentration Rank 14

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 25.82 | 3.14 ng/ul | 245409 | Perylene-d12     | 24.67 |

| Hit# | of | 5 | Tentative ID       | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------|-----|---------|-------------|------|
| 1    |    |   | Eicosane           | 282 | C20H42  | 000112-95-8 | 92   |
| 2    |    |   | Heneicosane        | 296 | C21H44  | 000629-94-7 | 92   |
| 3    |    |   | Heptadecane        | 240 | C17H36  | 000629-78-7 | 87   |
| 4    |    |   | Tetradecane        | 198 | C14H30  | 000629-59-4 | 87   |
| 5    |    |   | Tridecane, 1-iodo- | 310 | C13H27I | 035599-77-0 | 83   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_G\DATA\BG081320\  
 Data File : BG046256.D  
 Acq On : 13 Aug 2020 19:42  
 Operator : CG/JU  
 Sample : L3629-14  
 Misc :  
 ALS Vial : 10 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 BFM61

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG081320PAH.M  
 Quant Title : SVOA CALIBRATION

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

| TIC Top Hit name     | RT    | EstConc | Units | Response | --Internal Standard-- |       |         |      |
|----------------------|-------|---------|-------|----------|-----------------------|-------|---------|------|
|                      |       |         |       |          | #                     | RT    | Resp    | Conc |
| Butane, 2-methoxy... | 3.17  | 116.2   | ng/ul | 3205250  | 1                     | 7.95  | 551873  | 20.0 |
| 2(3H)-Furanone, 5... | 7.12  | 13.9    | ng/ul | 384904   | 1                     | 7.95  | 551873  | 20.0 |
| unknown-01           | 7.28  | 11.4    | ng/ul | 315960   | 1                     | 7.95  | 551873  | 20.0 |
| unknown-02           | 7.48  | 14.1    | ng/ul | 387720   | 1                     | 7.95  | 551873  | 20.0 |
| unknown-03           | 8.65  | 14.9    | ng/ul | 411077   | 1                     | 7.95  | 551873  | 20.0 |
| unknown-04           | 9.83  | 7.4     | ng/ul | 306446   | 2                     | 10.75 | 825134  | 20.0 |
| unknown-05           | 10.87 | 9.2     | ng/ul | 378500   | 2                     | 10.75 | 825134  | 20.0 |
| 2,4,7,9-Tetrameth... | 13.48 | 2.1     | ng/ul | 122723   | 3                     | 14.57 | 1154970 | 20.0 |
| unknown-06           | 13.98 | 13.9    | ng/ul | 805297   | 3                     | 14.57 | 1154970 | 20.0 |
| unknown-07           | 14.76 | 4.3     | ng/ul | 245679   | 3                     | 14.57 | 1154970 | 20.0 |
| unknown-08           | 15.68 | 5.3     | ng/ul | 304692   | 3                     | 14.57 | 1154970 | 20.0 |
| n-Hexadecanoic acid  | 18.22 | 2.6     | ng/ul | 180592   | 4                     | 17.32 | 1398470 | 20.0 |
| (DEL) Alkane: Str... | 21.23 | 4.1     | ng/ul | 374365   | 5                     | 21.56 | 1830000 | 20.0 |
| (DEL) Alkane: Str... | 23.82 | 3.0     | ng/ul | 235311   | 6                     | 24.67 | 1563380 | 20.0 |
| (DEL) Alkane: Str... | 24.73 | 2.7     | ng/ul | 209157   | 6                     | 24.67 | 1563380 | 20.0 |
| (DEL) Alkane: Str... | 25.82 | 3.1     | ng/ul | 245409   | 6                     | 24.67 | 1563380 | 20.0 |