

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021020\  
 Data File : BG044380.D  
 Acq On : 11 Feb 2020 8:27  
 Operator : CG/JU  
 Sample : L1446-01  
 Misc :  
 ALS Vial : 24 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 HP-640

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 3 % 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\8270-BG013020.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

| peak<br># | R.T.<br>min | first<br>scan | max<br>scan | last<br>scan | PK<br>TY | peak<br>height | corr.<br>area | corr.<br>% max. | % of<br>total |
|-----------|-------------|---------------|-------------|--------------|----------|----------------|---------------|-----------------|---------------|
| 1         | 5.488       | 400           | 408         | 416          | rBV      | 385638         | 675382        | 13.31%          | 1.310%        |
| 2         | 7.080       | 671           | 679         | 693          | rVB      | 229528         | 446061        | 8.79%           | 0.865%        |
| 3         | 7.902       | 810           | 819         | 827          | rBV      | 307477         | 546305        | 10.77%          | 1.059%        |
| 4         | 8.226       | 866           | 874         | 890          | rVB      | 1126511        | 2017824       | 39.78%          | 3.913%        |
| 5         | 9.060       | 1008          | 1016        | 1027         | rBV      | 871263         | 1563193       | 30.81%          | 3.032%        |
| 6         | 9.348       | 1041          | 1065        | 1080         | rBV      | 511769         | 2475455       | 48.80%          | 4.801%        |
| 7         | 10.035      | 1175          | 1182        | 1185         | rBV6     | 30705          | 61729         | 1.22%           | 0.120%        |
| 8         | 10.076      | 1185          | 1189        | 1204         | rVB2     | 38331          | 119909        | 2.36%           | 0.233%        |
| 9         | 10.488      | 1252          | 1259        | 1270         | rBV3     | 37219          | 79944         | 1.58%           | 0.155%        |
| 10        | 10.705      | 1288          | 1296        | 1302         | rBV      | 416195         | 767886        | 15.14%          | 1.489%        |
| 11        | 10.764      | 1302          | 1306        | 1312         | rVB      | 149135         | 236459        | 4.66%           | 0.459%        |
| 12        | 11.251      | 1381          | 1389        | 1394         | rBV      | 2544672        | 4403875       | 86.81%          | 8.541%        |
| 13        | 11.316      | 1394          | 1400        | 1405         | rVV      | 2947961        | 4838759       | 95.38%          | 9.384%        |
| 14        | 11.504      | 1425          | 1432        | 1437         | rBV3     | 46218          | 106410        | 2.10%           | 0.206%        |
| 15        | 11.575      | 1439          | 1444        | 1453         | rVV      | 164784         | 320119        | 6.31%           | 0.621%        |
| 16        | 11.663      | 1454          | 1459        | 1465         | rVV4     | 24350          | 63247         | 1.25%           | 0.123%        |
| 17        | 11.745      | 1469          | 1473        | 1482         | rVB3     | 26493          | 55366         | 1.09%           | 0.107%        |
| 18        | 12.526      | 1601          | 1606        | 1611         | rBV      | 101625         | 150448        | 2.97%           | 0.292%        |
| 19        | 12.979      | 1676          | 1683        | 1693         | rBV3     | 50574          | 131723        | 2.60%           | 0.255%        |
| 20        | 13.167      | 1707          | 1715        | 1723         | rBV      | 2286352        | 3627880       | 71.51%          | 7.036%        |
| 21        | 13.514      | 1768          | 1774        | 1779         | rBV3     | 41260          | 73502         | 1.45%           | 0.143%        |
| 22        | 13.989      | 1850          | 1855        | 1860         | rVB      | 48437          | 71619         | 1.41%           | 0.139%        |
| 23        | 14.089      | 1867          | 1872        | 1880         | rVB3     | 50382          | 83899         | 1.65%           | 0.163%        |
| 24        | 14.195      | 1885          | 1890        | 1903         | rBV2     | 108343         | 226993        | 4.47%           | 0.440%        |
| 25        | 14.542      | 1941          | 1949        | 1954         | rBV2     | 703925         | 1237611       | 24.40%          | 2.400%        |
| 26        | 14.595      | 1954          | 1958        | 1978         | rVB2     | 168783         | 431407        | 8.50%           | 0.837%        |
| 27        | 14.777      | 1984          | 1989        | 1994         | rBV      | 124712         | 189331        | 3.73%           | 0.367%        |
| 28        | 15.558      | 2117          | 2122        | 2127         | rBV2     | 42969          | 62479         | 1.23%           | 0.121%        |
| 29        | 16.028      | 2196          | 2202        | 2209         | rBV      | 1219063        | 1872753       | 36.92%          | 3.632%        |
| 30        | 16.087      | 2209          | 2212        | 2222         | rVB2     | 64379          | 111038        | 2.19%           | 0.215%        |
| 31        | 16.181      | 2224          | 2228        | 2234         | rBV3     | 75933          | 117923        | 2.32%           | 0.229%        |
| 32        | 16.322      | 2249          | 2252        | 2256         | rBV      | 43389          | 57659         | 1.14%           | 0.112%        |
| 33        | 16.428      | 2266          | 2270        | 2279         | rVB      | 136845         | 220191        | 4.34%           | 0.427%        |
| 34        | 16.998      | 2362          | 2367        | 2372         | rBV      | 147356         | 207803        | 4.10%           | 0.403%        |

Data Path : Z:\SVOASRV\HPCHEM1\BNA\_G\DATA\BG021020\  
 Data File : BG044380.D  
 Acq On : 11 Feb 2020 8:27  
 Operator : CG/JU  
 Sample : L1446-01  
 Misc :  
 ALS Vial : 24 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 HP-640

Integration Parameters: rteint.p  
 Integrator: RTE  
 Smoothing : ON  
 Sampling : 1  
 Start Thrs: 0.2  
 Stop Thrs : 0

Filtering: 5  
 Min Area: 3 % 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\8270-BG013020.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |         |         |         |        |
|----|--------|------|------|------|------|---------|---------|---------|--------|
| 35 | 17.045 | 2372 | 2375 | 2385 | rVB5 | 51194   | 93203   | 1.84%   | 0.181% |
| 36 | 17.286 | 2410 | 2416 | 2427 | rVB  | 804503  | 1229619 | 24.24%  | 2.385% |
| 37 | 17.579 | 2461 | 2466 | 2472 | rVB2 | 70288   | 134213  | 2.65%   | 0.260% |
| 38 | 17.756 | 2493 | 2496 | 2505 | rVB  | 88254   | 134918  | 2.66%   | 0.262% |
| 39 | 17.979 | 2528 | 2534 | 2540 | rVB5 | 45837   | 75107   | 1.48%   | 0.146% |
| 40 | 18.196 | 2566 | 2571 | 2576 | rBV  | 152986  | 257173  | 5.07%   | 0.499% |
| 41 | 18.255 | 2576 | 2581 | 2587 | rVB  | 399134  | 579065  | 11.41%  | 1.123% |
| 42 | 18.373 | 2597 | 2601 | 2606 | rVB  | 34555   | 54943   | 1.08%   | 0.107% |
| 43 | 18.467 | 2613 | 2617 | 2624 | rBV  | 190709  | 296433  | 5.84%   | 0.575% |
| 44 | 18.925 | 2685 | 2695 | 2702 | rBV3 | 233265  | 396050  | 7.81%   | 0.768% |
| 45 | 19.119 | 2722 | 2728 | 2734 | rVB6 | 42667   | 88818   | 1.75%   | 0.172% |
| 46 | 19.354 | 2759 | 2768 | 2774 | rBV5 | 163982  | 386748  | 7.62%   | 0.750% |
| 47 | 19.530 | 2794 | 2798 | 2806 | rVB  | 172149  | 246062  | 4.85%   | 0.477% |
| 48 | 19.712 | 2822 | 2829 | 2835 | rBV6 | 77641   | 172901  | 3.41%   | 0.335% |
| 49 | 19.906 | 2856 | 2862 | 2874 | rVB  | 3587410 | 5073072 | 100.00% | 9.839% |
| 50 | 20.082 | 2889 | 2892 | 2905 | rVB  | 193498  | 318470  | 6.28%   | 0.618% |
| 51 | 20.447 | 2950 | 2954 | 2963 | rBV  | 309049  | 426497  | 8.41%   | 0.827% |
| 52 | 20.593 | 2973 | 2979 | 2985 | rVB  | 719816  | 917295  | 18.08%  | 1.779% |
| 53 | 20.693 | 2991 | 2996 | 3004 | rVB  | 3825101 | 4552171 | 89.73%  | 8.828% |
| 54 | 20.940 | 3034 | 3038 | 3048 | rBV  | 228193  | 338270  | 6.67%   | 0.656% |
| 55 | 21.105 | 3063 | 3066 | 3070 | rBV3 | 40041   | 51836   | 1.02%   | 0.101% |
| 56 | 21.204 | 3078 | 3083 | 3089 | rBV3 | 286750  | 491444  | 9.69%   | 0.953% |
| 57 | 21.269 | 3090 | 3094 | 3097 | rVV  | 171586  | 235204  | 4.64%   | 0.456% |
| 58 | 21.304 | 3097 | 3100 | 3111 | rVB  | 148004  | 227335  | 4.48%   | 0.441% |
| 59 | 21.440 | 3119 | 3123 | 3124 | rBV  | 100392  | 124180  | 2.45%   | 0.241% |
| 60 | 21.463 | 3124 | 3127 | 3132 | rVV  | 135595  | 212064  | 4.18%   | 0.411% |
| 61 | 21.534 | 3133 | 3139 | 3146 | rVB  | 794964  | 1305971 | 25.74%  | 2.533% |
| 62 | 21.739 | 3170 | 3174 | 3178 | rVB  | 116250  | 152857  | 3.01%   | 0.296% |
| 63 | 21.851 | 3188 | 3193 | 3201 | rBV2 | 256076  | 501254  | 9.88%   | 0.972% |
| 64 | 22.327 | 3269 | 3274 | 3288 | rVB  | 161598  | 304737  | 6.01%   | 0.591% |
| 65 | 22.456 | 3292 | 3296 | 3310 | rVB2 | 137899  | 289860  | 5.71%   | 0.562% |
| 66 | 22.679 | 3329 | 3334 | 3341 | rVB2 | 133605  | 235940  | 4.65%   | 0.458% |
| 67 | 22.938 | 3374 | 3378 | 3382 | rBV3 | 74836   | 133741  | 2.64%   | 0.259% |
| 68 | 22.991 | 3382 | 3387 | 3394 | rVB2 | 206497  | 370546  | 7.30%   | 0.719% |
| 69 | 23.155 | 3410 | 3415 | 3422 | rBV5 | 60718   | 144699  | 2.85%   | 0.281% |
| 70 | 23.702 | 3502 | 3508 | 3512 | rBV2 | 132736  | 283933  | 5.60%   | 0.551% |
| 71 | 23.760 | 3513 | 3518 | 3526 | rVB  | 196540  | 410910  | 8.10%   | 0.797% |

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021020\  
Data File : BG044380.D  
Acq On : 11 Feb 2020 8:27  
Operator : CG/JU  
Sample : L1446-01  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
HP-640

Integration Parameters: rteint.p  
Integrator: RTE  
Smoothing : ON Filtering: 5  
Sampling : 1 Min Area: 3 % of largest Peak  
Start Thrs: 0.2 Max Peaks: 100  
Stop Thrs : 0 Peak Location: TOP

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

Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\8270-BG013020.M  
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |        |         |        |        |
|----|--------|------|------|------|------|--------|---------|--------|--------|
| 72 | 24.612 | 3653 | 3663 | 3684 | rVB3 | 550852 | 1968188 | 38.80% | 3.817% |
| 73 | 25.329 | 3780 | 3785 | 3792 | rVB4 | 28295  | 68237   | 1.35%  | 0.132% |
| 74 | 25.746 | 3849 | 3856 | 3866 | rVB2 | 104310 | 305808  | 6.03%  | 0.593% |
| 75 | 26.516 | 3980 | 3987 | 4001 | rVB5 | 54214  | 201881  | 3.98%  | 0.392% |
| 76 | 27.033 | 4071 | 4075 | 4084 | rVB4 | 46511  | 118985  | 2.35%  | 0.231% |

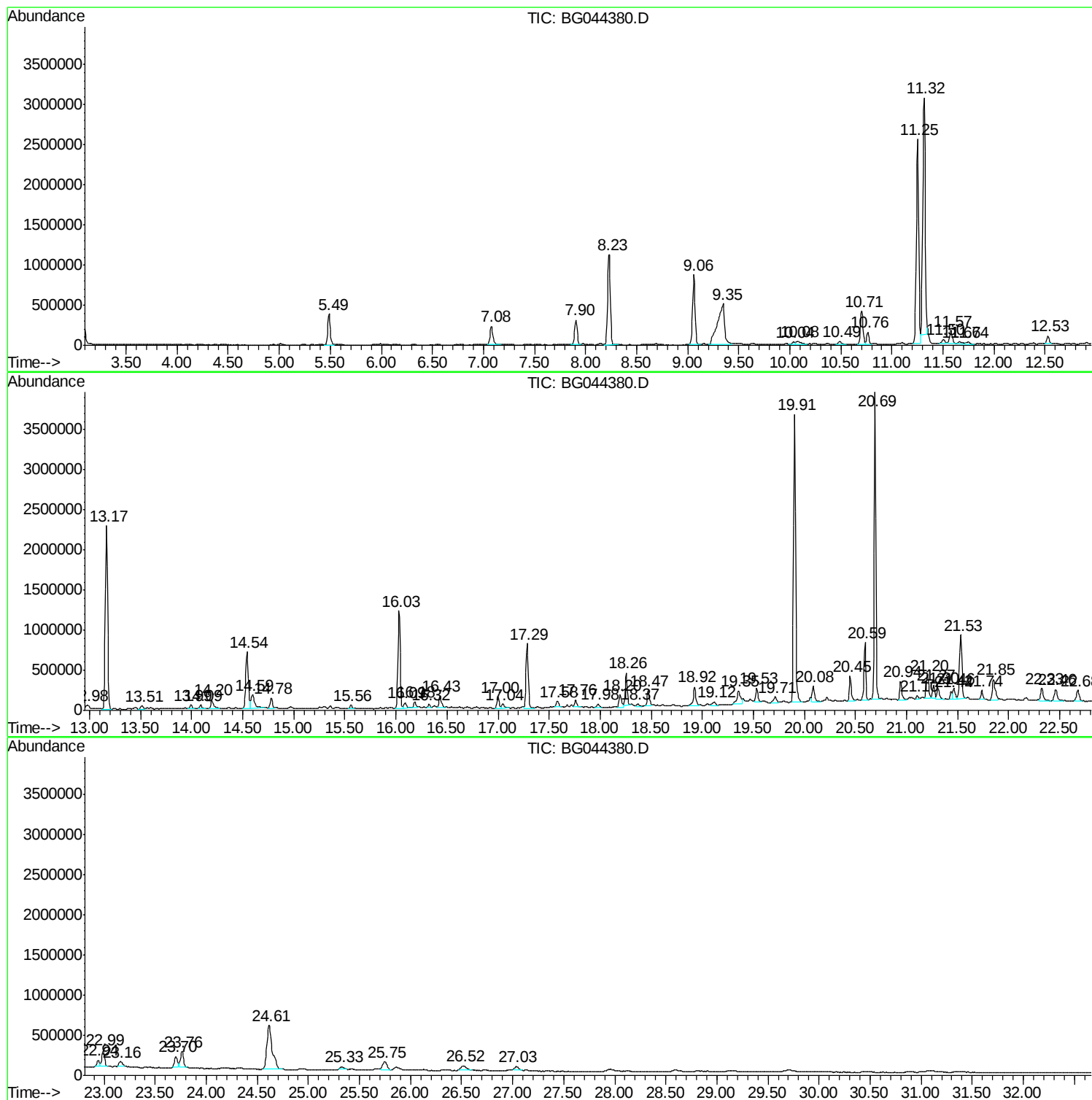
Sum of corrected areas: 51562820

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021020\  
Data File : BG044380.D  
Acq On : 11 Feb 2020 8:27  
Operator : CG/JU  
Sample : L1446-01  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
HP-640

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG013020.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021020\  
Data File : BG044380.D  
Acq On : 11 Feb 2020 8:27  
Operator : CG/JU  
Sample : L1446-01  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
HP-640

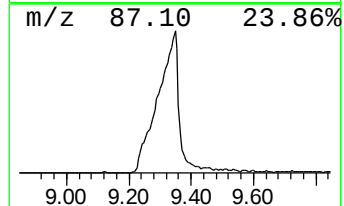
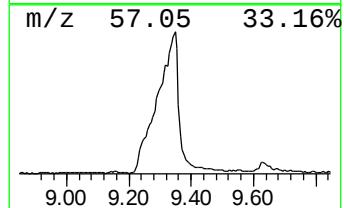
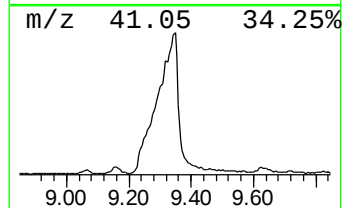
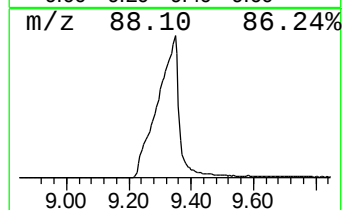
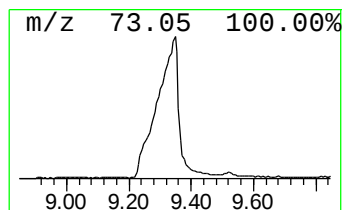
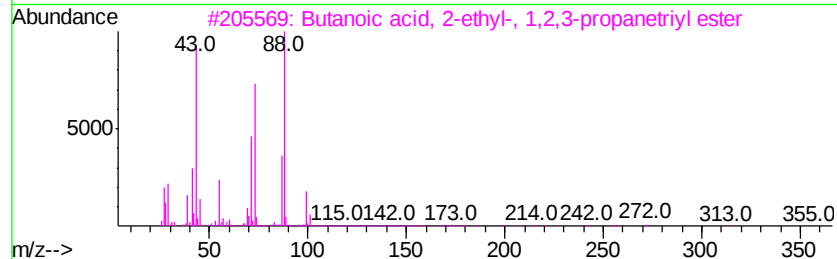
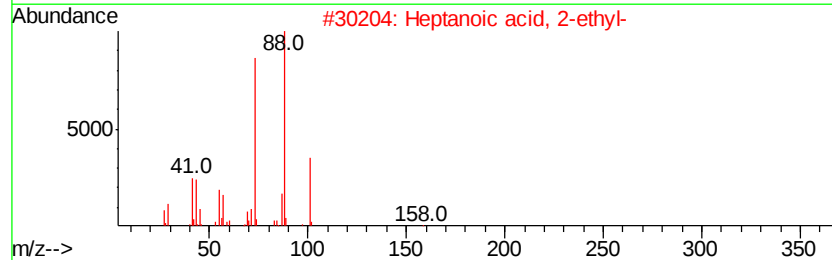
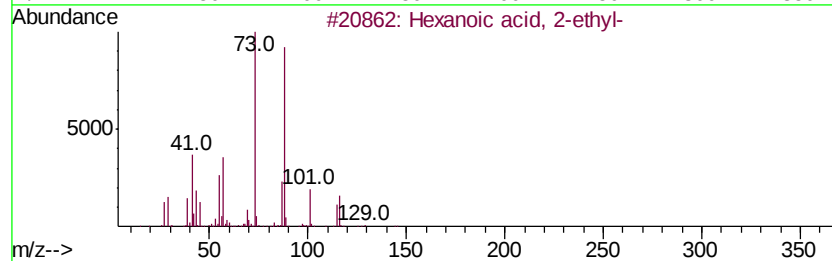
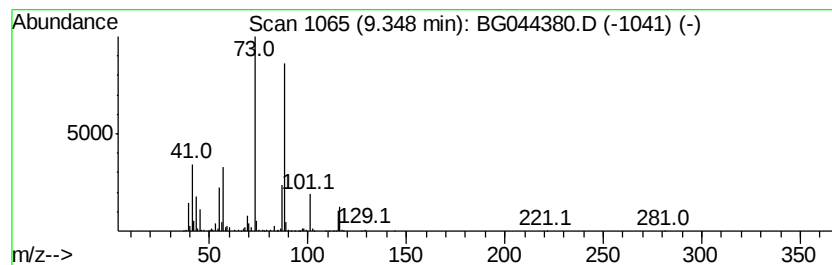
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG013020.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 2 Hexanoic acid, 2-ethyl- Concentration Rank 5

| R.T. | EstConc  | Area    | Relative to ISTD | R.T.  |
|------|----------|---------|------------------|-------|
| 9.35 | 64.47 ng | 2475460 | Naphthalene-d8   | 10.71 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | Hexanoic acid, 2-ethyl-             | 144 | C8H16O2  | 000149-57-5 | 90   |
| 2    |    | Heptanoic acid, 2-ethyl-            | 158 | C9H18O2  | 003274-29-1 | 59   |
| 3    |    | Butanoic acid, 2-ethyl-, 1,2,3-p... | 386 | C21H38O6 | 056554-54-2 | 53   |
| 4    |    | Octanoic acid, 2-methyl-, methyl... | 172 | C10H20O2 | 002177-86-8 | 47   |
| 5    |    | Methyl 2,3,4-tri-O-methyl-6-deox... | 220 | C10H20O5 | 072983-16-5 | 40   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021020\  
Data File : BG044380.D  
Acq On : 11 Feb 2020 8:27  
Operator : CG/JU  
Sample : L1446-01  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
HP-640

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG013020.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

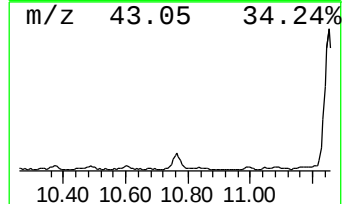
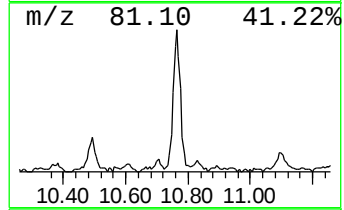
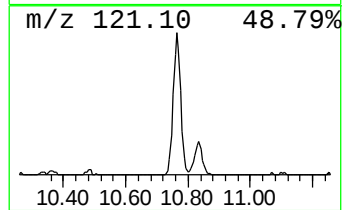
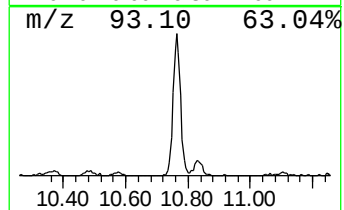
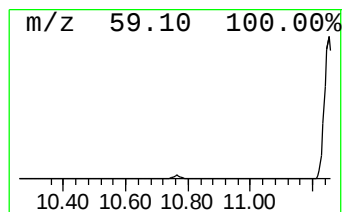
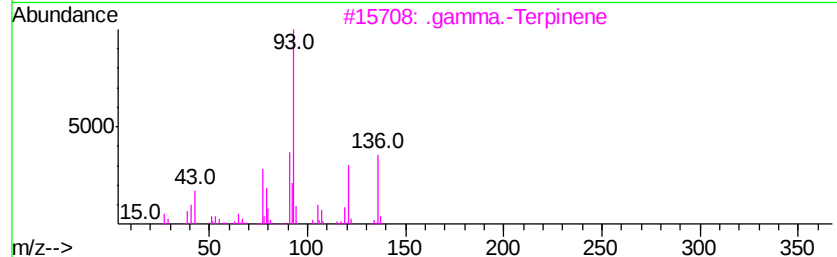
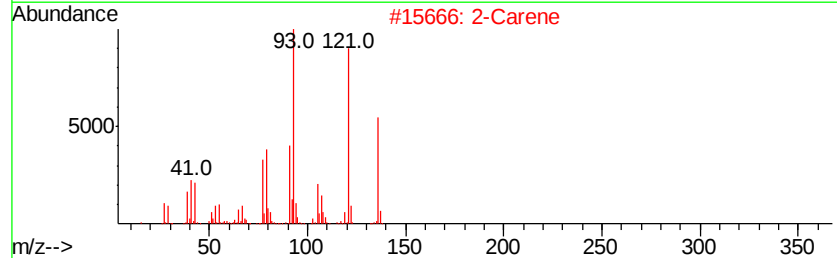
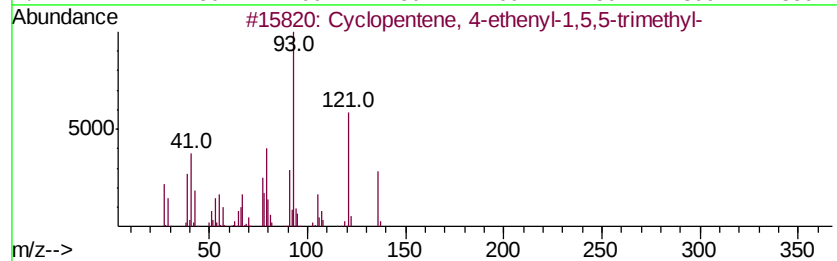
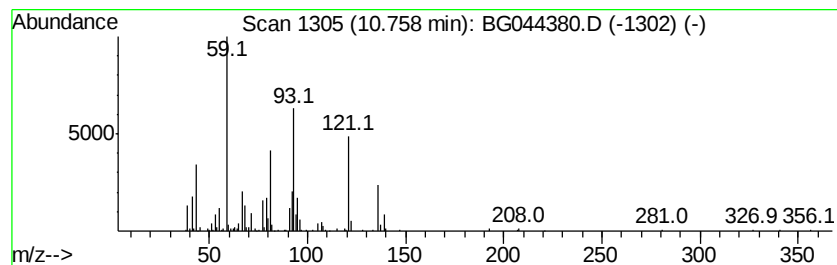
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 3 unknown10.76 Concentration Rank 13

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 10.76 | 6.16 ng | 236459 | Naphthalene-d8   | 10.71 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Cyclopentene, 4-ethenyl-1,5,5-tr... | 136 | C10H16  | 001727-69-1 | 41   |
| 2    |    | 2-Carene                            | 136 | C10H16  | 000554-61-0 | 41   |
| 3    |    | .gamma.-Terpinene                   | 136 | C10H16  | 000099-85-4 | 38   |
| 4    |    | Tricyclo[2.2.1.0(2,6)]heptane, 1... | 136 | C10H16  | 000508-32-7 | 38   |
| 5    |    | Bicyclo[2.2.1]hept-2-ene, 1,7,7-... | 136 | C10H16  | 000464-17-5 | 38   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021020\  
Data File : BG044380.D  
Acq On : 11 Feb 2020 8:27  
Operator : CG/JU  
Sample : L1446-01  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
HP-640

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG013020.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

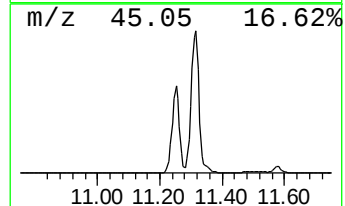
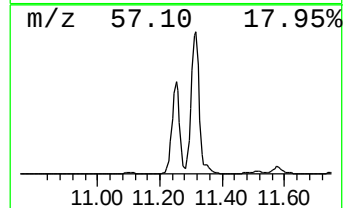
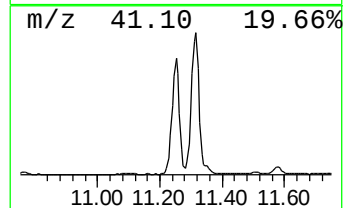
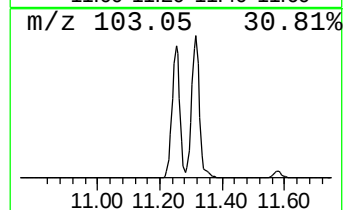
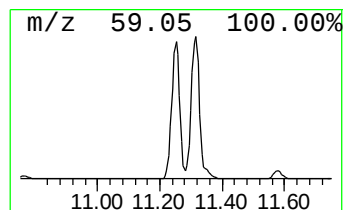
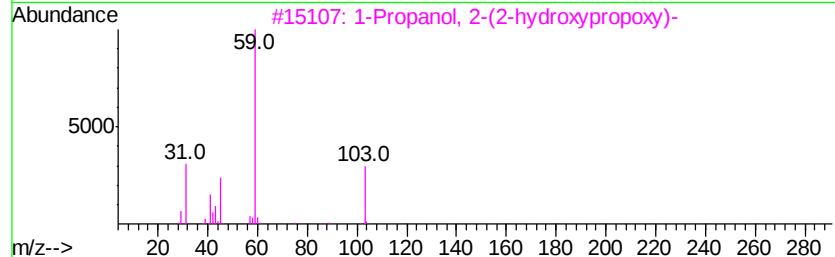
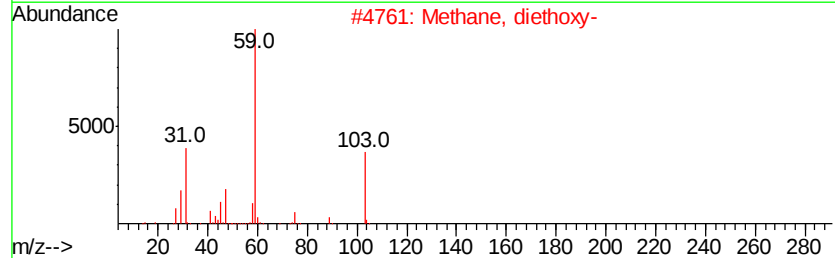
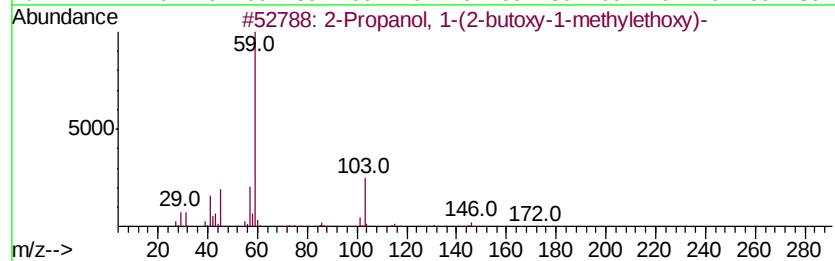
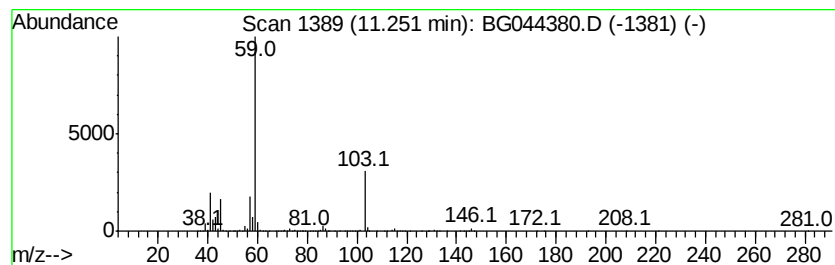
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 4 2-Propanol, 1-(2-butoxy-1-m... Concentration Rank 2

| R.T.  | EstConc   | Area    | Relative to ISTD | R.T.  |
|-------|-----------|---------|------------------|-------|
| 11.25 | 114.70 ng | 4403880 | Naphthalene-d8   | 10.71 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | 2-Propanol, 1-(2-butoxy-1-methyl... | 190 | C10H22O3 | 029911-28-2 | 83   |
| 2    |    | Methane, diethoxy-                  | 104 | C5H12O2  | 000462-95-3 | 80   |
| 3    |    | 1-Propanol, 2-(2-hydroxypropoxy)-   | 134 | C6H14O3  | 000106-62-7 | 72   |
| 4    |    | 1-(1-tert-Butoxypropan-2-yloxy)p... | 190 | C10H22O3 | 132739-31-2 | 72   |
| 5    |    | 1-Propanol, 2,2'-oxybis-            | 134 | C6H14O3  | 000108-61-2 | 64   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021020\  
Data File : BG044380.D  
Acq On : 11 Feb 2020 8:27  
Operator : CG/JU  
Sample : L1446-01  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
HP-640

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG013020.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

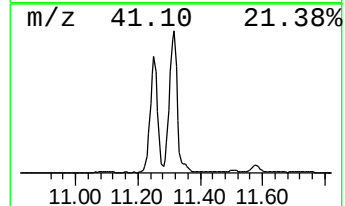
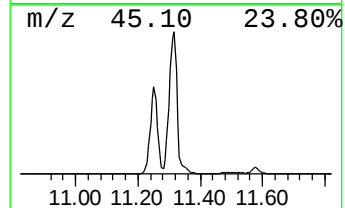
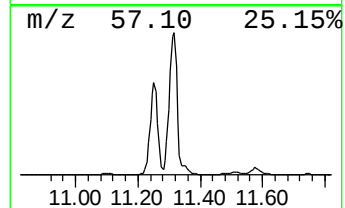
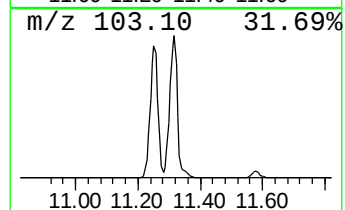
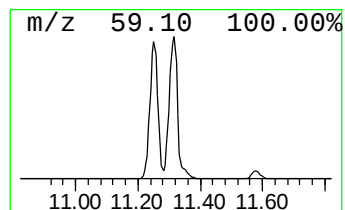
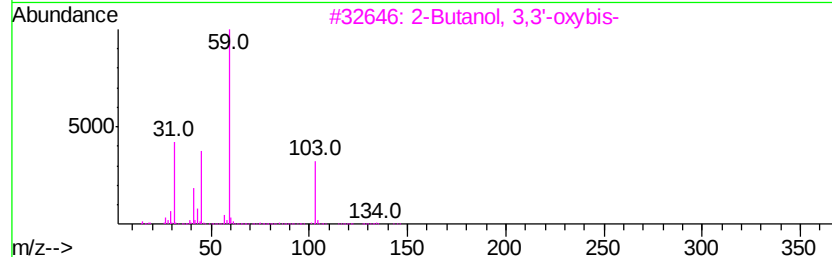
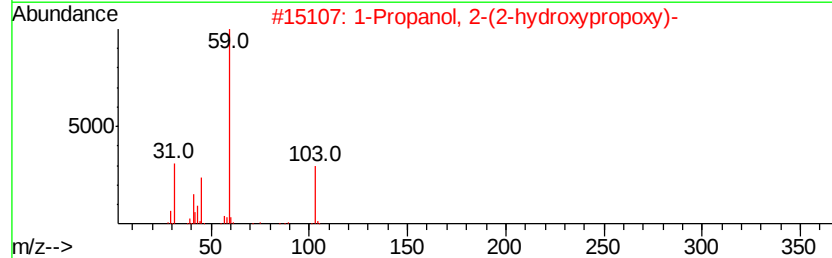
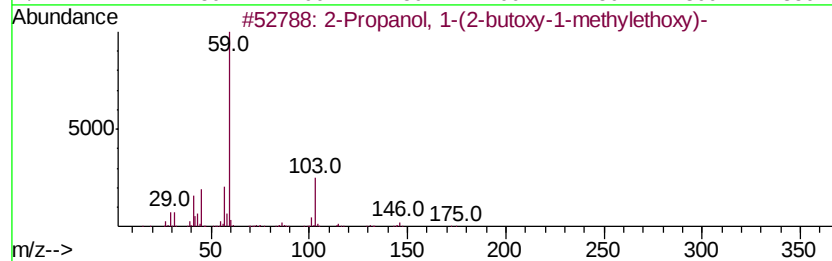
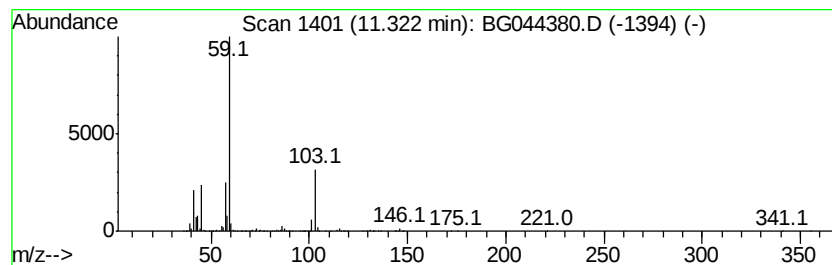
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 5 1-Propanol, 2-(2-hydroxypro... Concentration Rank 1

| R.T.  | EstConc   | Area    | Relative to ISTD | R.T.  |
|-------|-----------|---------|------------------|-------|
| 11.32 | 126.03 ng | 4838760 | Naphthalene-d8   | 10.71 |

| Hit# of | 5                                    | Tentative ID | MW       | MolForm | CAS#        | Qual |
|---------|--------------------------------------|--------------|----------|---------|-------------|------|
| 1       | 2-Propanol, 1-(2-butoxy-1-methyl...  | 190          | C10H22O3 |         | 029911-28-2 | 90   |
| 2       | 1-Propanol, 2-(2-hydroxypropoxy)-    | 134          | C6H14O3  |         | 000106-62-7 | 78   |
| 3       | 2-Butanol, 3,3'-oxybis-              | 162          | C8H18O3  |         | 054305-61-2 | 78   |
| 4       | 2-Propanol, 1-(2-methoxy-1-methyl... | 148          | C7H16O3  |         | 020324-32-7 | 72   |
| 5       | Methane, diethoxy-                   | 104          | C5H12O2  |         | 000462-95-3 | 72   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021020\  
Data File : BG044380.D  
Acq On : 11 Feb 2020 8:27  
Operator : CG/JU  
Sample : L1446-01  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
HP-640

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG013020.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

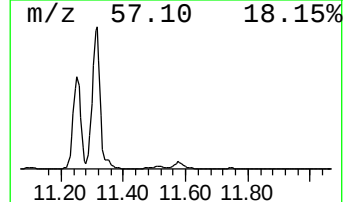
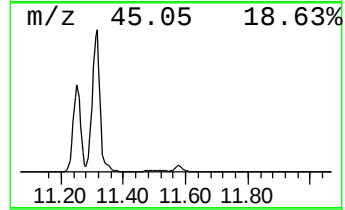
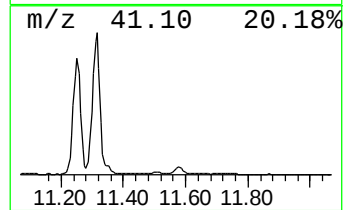
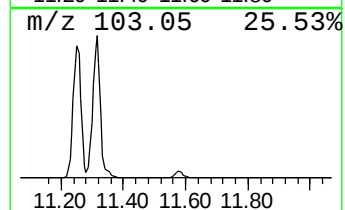
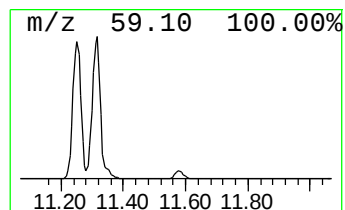
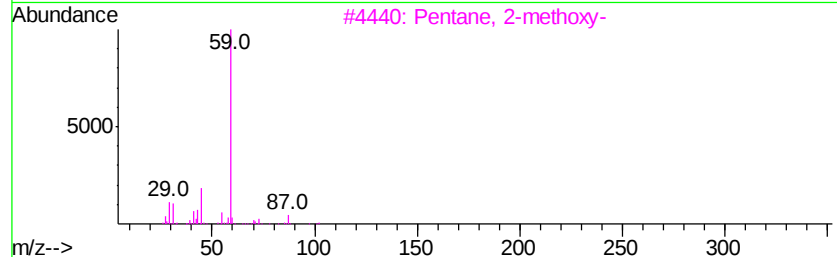
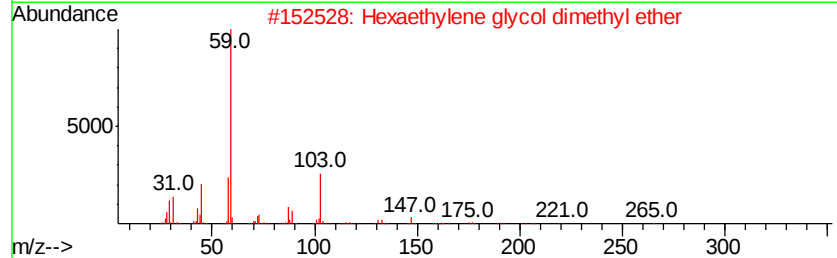
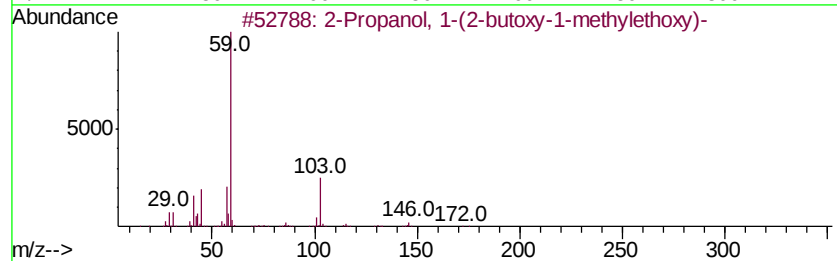
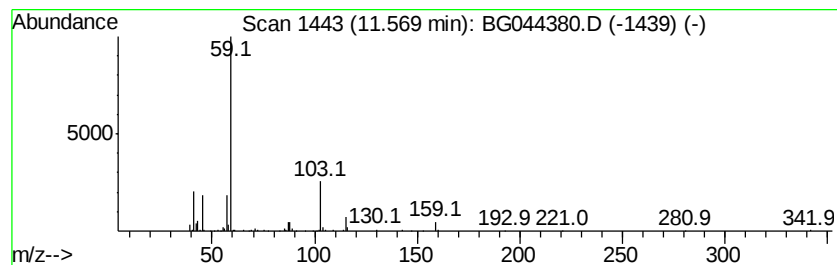
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 6 Hexaethylene glycol dimethy... Concentration Rank 6

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 11.57 | 8.34 ng | 320119 | Naphthalene-d8   | 10.71 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | 2-Propanol, 1-(2-butoxy-1-methyl... | 190 | C10H22O3 | 029911-28-2 | 72   |
| 2    |    | Hexaethylene glycol dimethyl ether  | 310 | C14H30O7 | 001072-40-8 | 64   |
| 3    |    | Pentane, 2-methoxy-                 | 102 | C6H14O   | 006795-88-6 | 58   |
| 4    |    | 2-Propanol, 1-[1-methyl-2-(2-pro... | 174 | C9H18O3  | 055956-25-7 | 56   |
| 5    |    | 1-Propanol, 2-(2-hydroxypropoxy)-   | 134 | C6H14O3  | 000106-62-7 | 50   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021020\  
Data File : BG044380.D  
Acq On : 11 Feb 2020 8:27  
Operator : CG/JU  
Sample : L1446-01  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

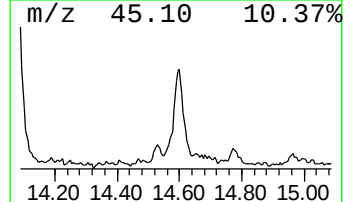
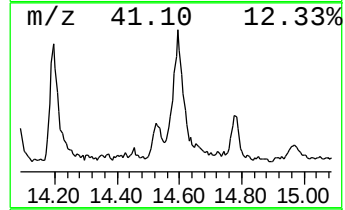
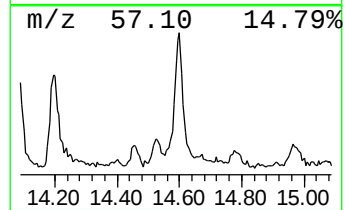
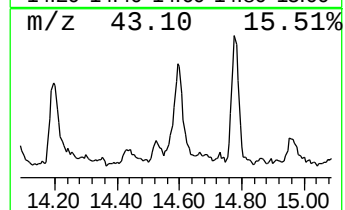
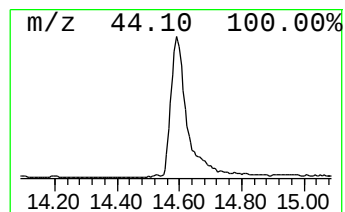
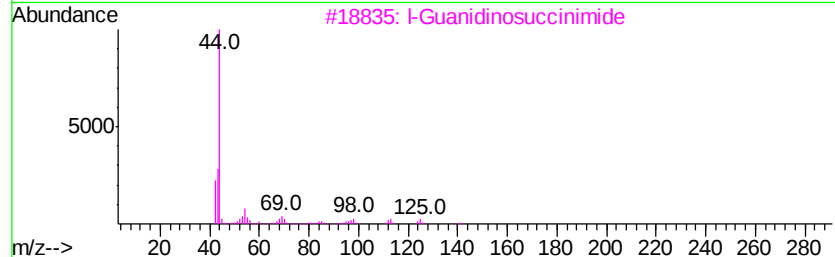
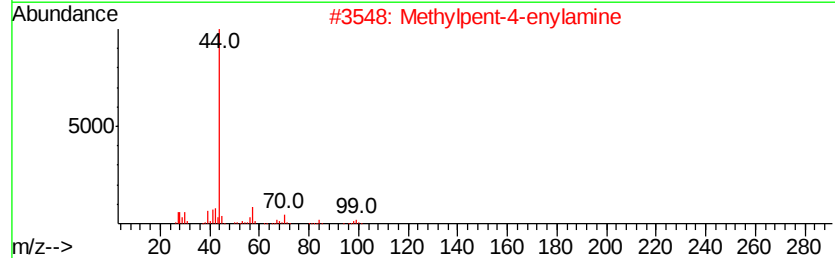
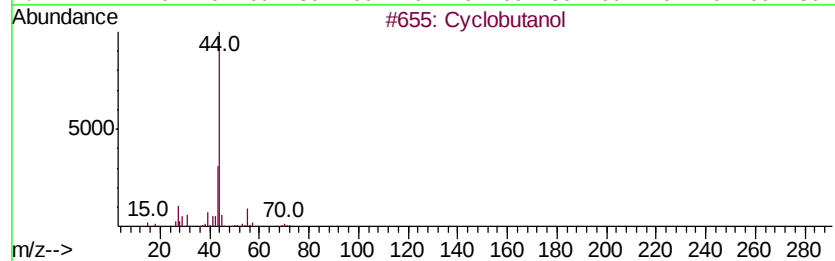
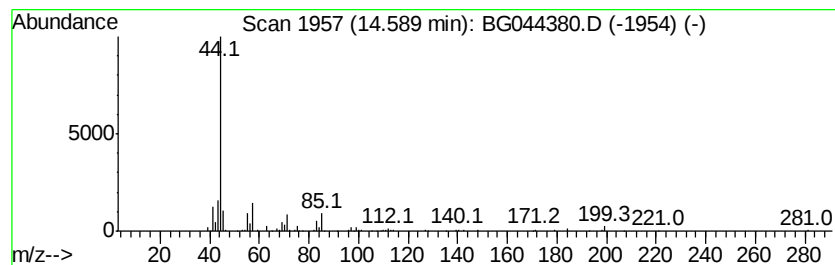
Instrument :  
BNA\_G  
ClientSampleId :  
HP-640

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG013020.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 7 Cyclobutanol Concentration Rank 9

| R.T.    | EstConc | Area                           | Relative to ISTD | R.T.            |
|---------|---------|--------------------------------|------------------|-----------------|
| 14.59   | 6.97 ng | 431407                         | Acenaphthene-d10 | 14.54           |
| Hit# of | 5       | Tentative ID                   | MW MolForm       | CAS# Qual       |
| 1       |         | Cyclobutanol                   | 72 C4H8O         | 002919-23-5 53  |
| 2       |         | Methylpent-4-enylamine         | 99 C6H13N        | 005831-72-1 50  |
| 3       |         | l-Guanidinosuccinimide         | 141 C5H7N3O2     | 1000130-20-8 40 |
| 4       |         | N-Dodecylmethylamine           | 199 C13H29N      | 1000342-26-1 40 |
| 5       |         | Glycine, N-(N-L-alanylglycyl)- | 203 C7H13N3O4    | 003146-40-5 36  |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021020\  
Data File : BG044380.D  
Acq On : 11 Feb 2020 8:27  
Operator : CG/JU  
Sample : L1446-01  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

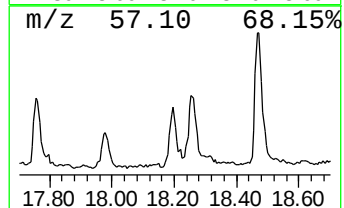
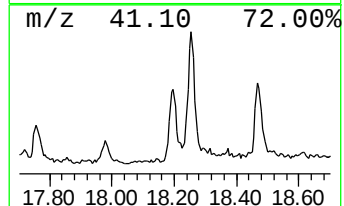
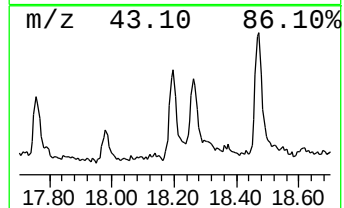
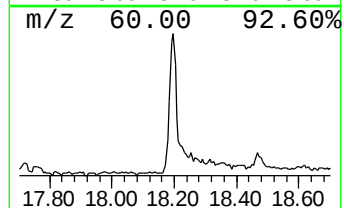
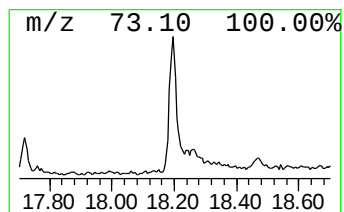
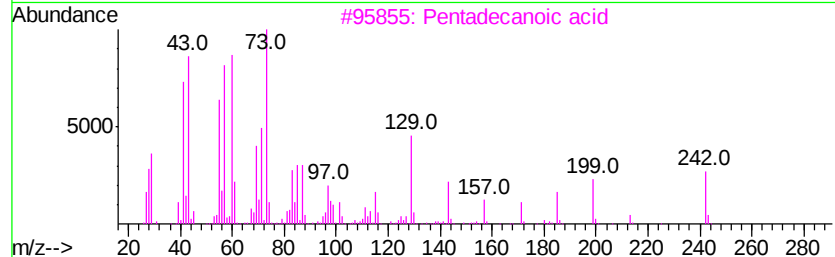
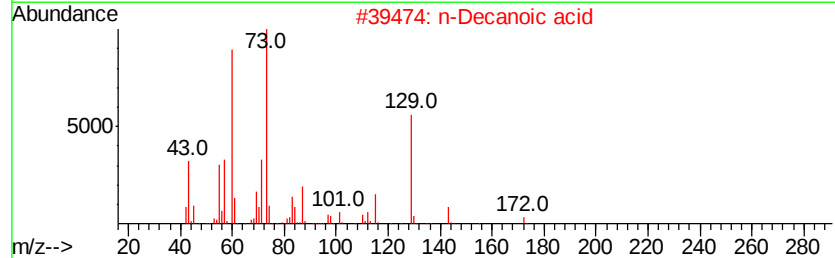
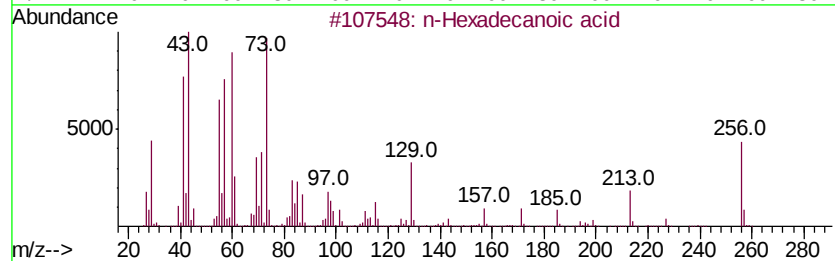
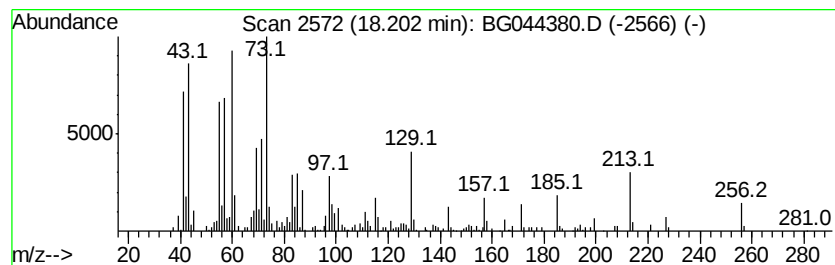
Instrument :  
BNA\_G  
ClientSampleId :  
HP-640

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG013020.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 8 n-Hexadecanoic acid Concentration Rank 20

| R.T.    | EstConc             | Area         | Relative to ISTD |          | R.T.        |      |
|---------|---------------------|--------------|------------------|----------|-------------|------|
| 18.20   | 4.18 ng             | 257173       | Phenanthrene-d10 |          | 17.29       |      |
| Hit# of | 5                   | Tentative ID | MW               | MolForm  | CAS#        | Qual |
| 1       | n-Hexadecanoic acid |              | 256              | C16H32O2 | 000057-10-3 | 98   |
| 2       | n-Decanoic acid     |              | 172              | C10H20O2 | 000334-48-5 | 91   |
| 3       | Pentadecanoic acid  |              | 242              | C15H30O2 | 001002-84-2 | 81   |
| 4       | Tridecanoic acid    |              | 214              | C13H26O2 | 000638-53-9 | 81   |
| 5       | Tetradecanoic acid  |              | 228              | C14H28O2 | 000544-63-8 | 74   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021020\  
Data File : BG044380.D  
Acq On : 11 Feb 2020 8:27  
Operator : CG/JU  
Sample : L1446-01  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
HP-640

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG013020.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

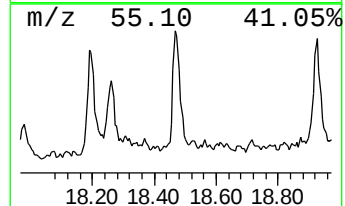
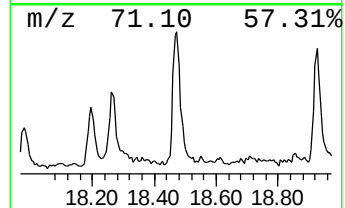
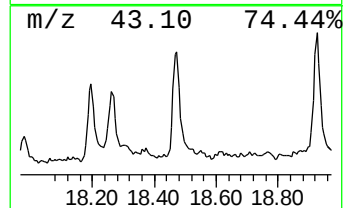
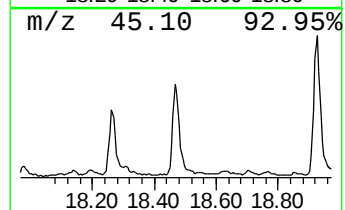
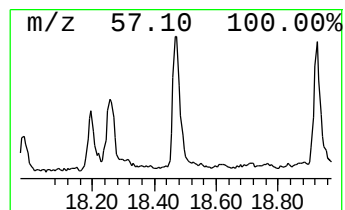
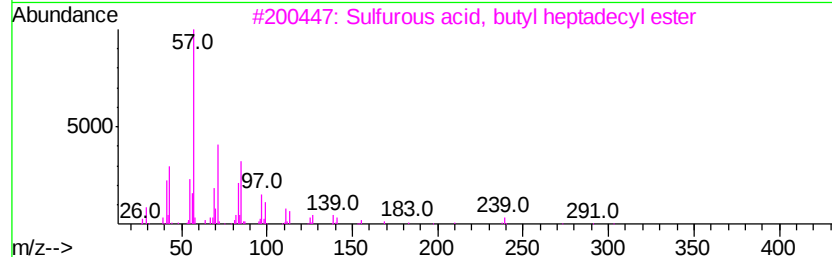
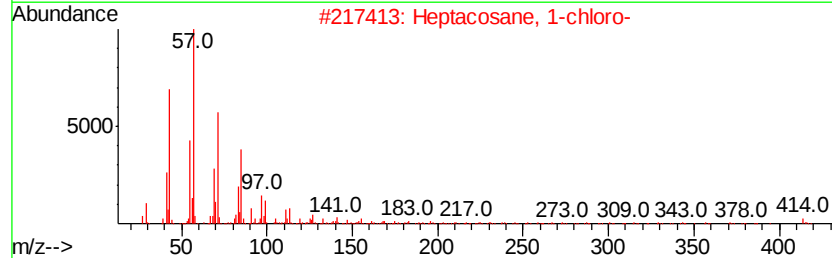
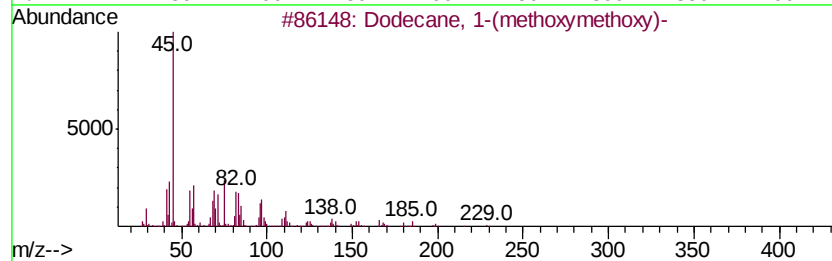
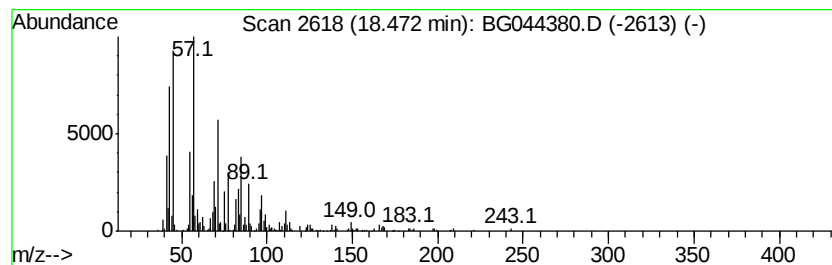
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 9 unknown18.47 Concentration Rank 17

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 18.47 | 4.82 ng | 296433 | Phenanthrene-d10 | 17.29 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | Dodecane, 1-(methoxymethoxy)-       | 230 | C14H30O2  | 034458-41-8  | 43   |
| 2    |    | Heptacosane, 1-chloro-              | 414 | C27H55Cl  | 062016-79-9  | 38   |
| 3    |    | Sulfurous acid, butyl heptadecyl... | 376 | C21H44O3S | 1000309-18-4 | 38   |
| 4    |    | Hexaethylene glycol                 | 282 | C12H26O7  | 002615-15-8  | 27   |
| 5    |    | Oxalic acid, isobutyl pentadecyl... | 356 | C21H40O4  | 1000309-38-0 | 25   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021020\  
Data File : BG044380.D  
Acq On : 11 Feb 2020 8:27  
Operator : CG/JU  
Sample : L1446-01  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
HP-640

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG013020.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

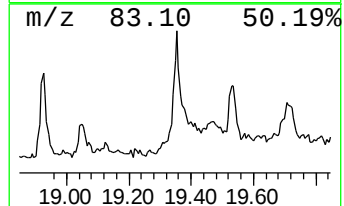
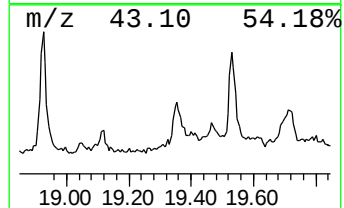
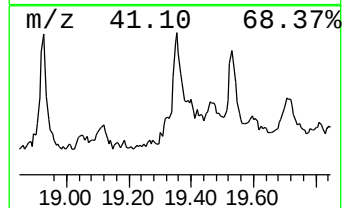
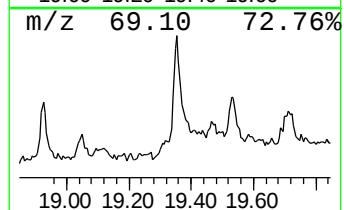
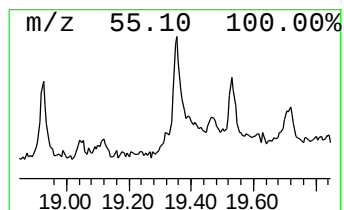
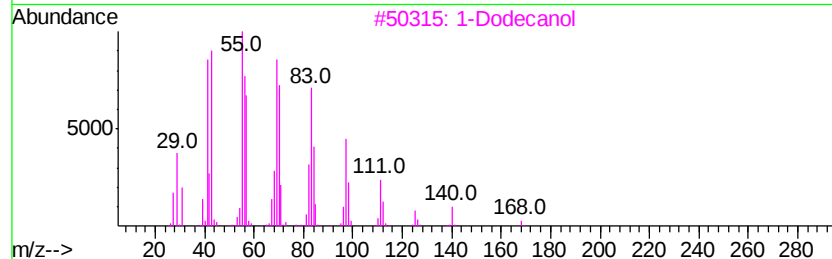
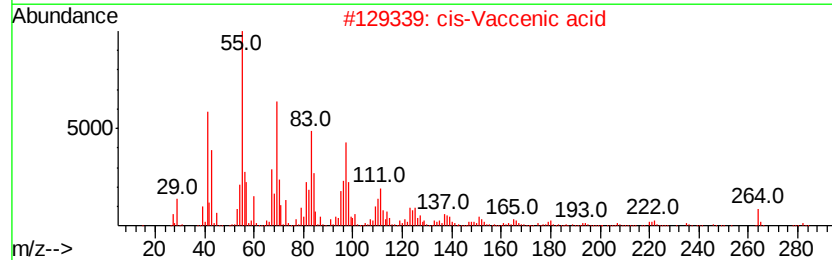
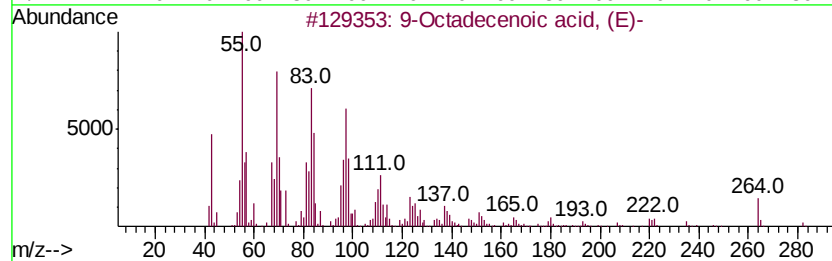
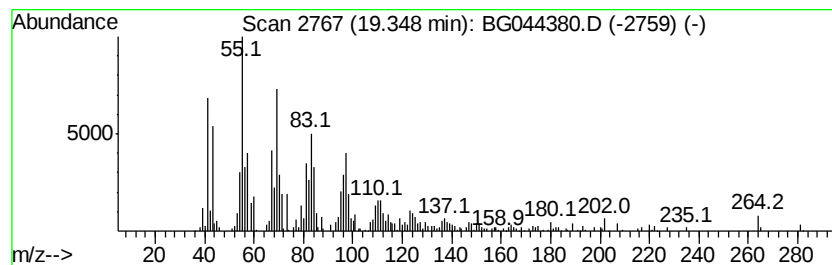
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 11 9-Octadecenoic acid, (E)- Concentration Rank 12

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 19.35 | 6.29 ng | 386748 | Phenanthrene-d10 | 17.29 |

| Hit# | of | Tentative ID              | MW  | MolForm  | CAS#        | Qual |
|------|----|---------------------------|-----|----------|-------------|------|
| 1    | 5  | 9-Octadecenoic acid, (E)- | 282 | C18H34O2 | 000112-79-8 | 93   |
| 2    |    | cis-Vaccenic acid         | 282 | C18H34O2 | 000506-17-2 | 70   |
| 3    |    | 1-Dodecanol               | 186 | C12H26O  | 000112-53-8 | 68   |
| 4    |    | 9-Hexadecenoic acid       | 254 | C16H30O2 | 002091-29-4 | 64   |
| 5    |    | Oleic Acid                | 282 | C18H34O2 | 000112-80-1 | 64   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021020\  
Data File : BG044380.D  
Acq On : 11 Feb 2020 8:27  
Operator : CG/JU  
Sample : L1446-01  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
HP-640

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG013020.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

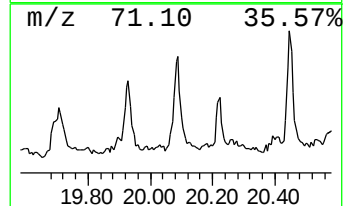
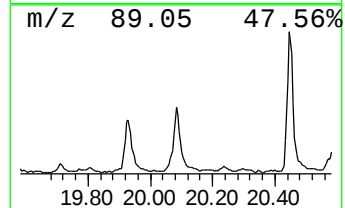
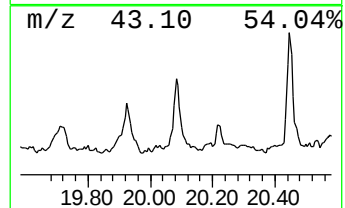
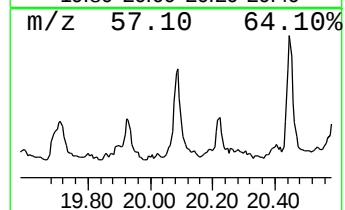
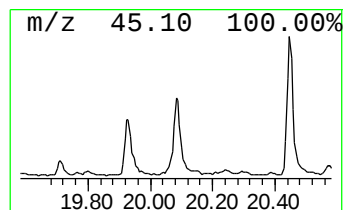
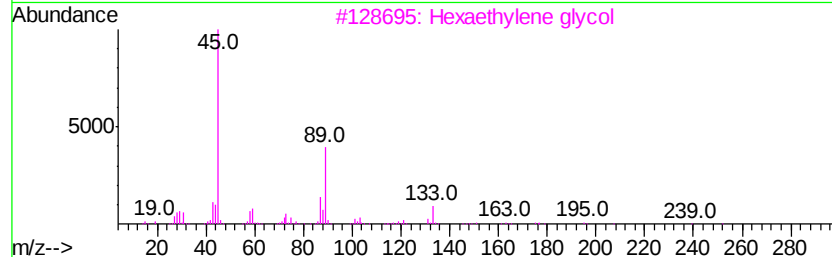
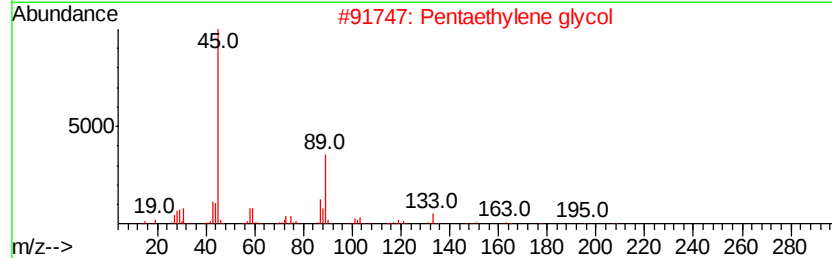
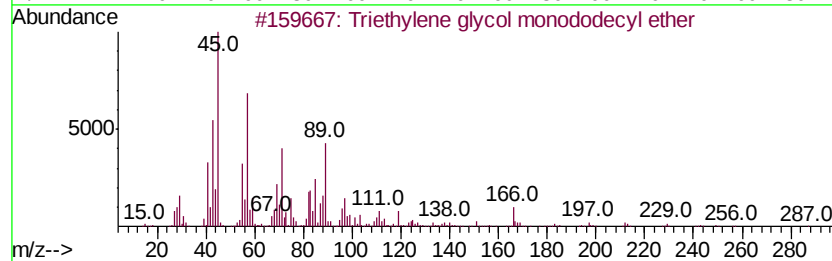
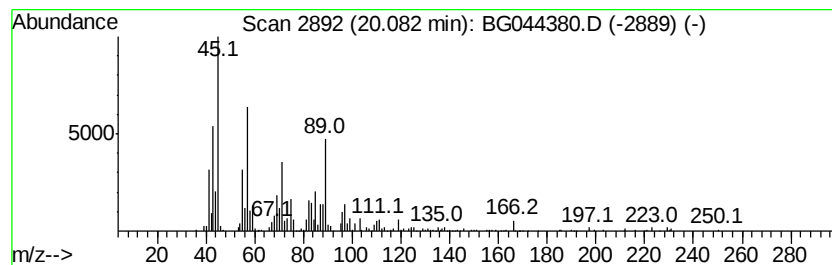
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 12 Triethylene glycol monododecyl ether Concentration Rank 16

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 20.08 | 4.88 ng | 318470 | Chrysene-d12     | 21.53 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | Triethylene glycol monododecyl e... | 318 | C18H38O4 | 003055-94-5 | 91   |
| 2    |    |   | Pentaethylene glycol                | 238 | C10H22O6 | 004792-15-8 | 58   |
| 3    |    |   | Hexaethylene glycol                 | 282 | C12H26O7 | 002615-15-8 | 53   |
| 4    |    |   | Heptaethylene glycol                | 326 | C14H30O8 | 005617-32-3 | 50   |
| 5    |    |   | 2-[2-[2-[2-[2-[2-(2-Hydroxyet...    | 370 | C16H34O9 | 005117-19-1 | 50   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021020\  
Data File : BG044380.D  
Acq On : 11 Feb 2020 8:27  
Operator : CG/JU  
Sample : L1446-01  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
HP-640

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG013020.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

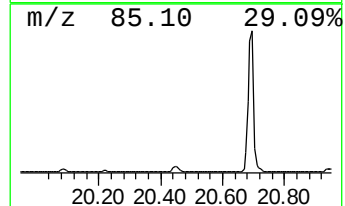
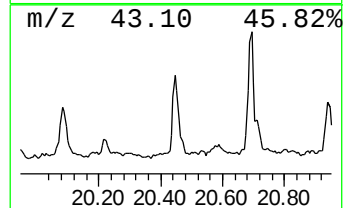
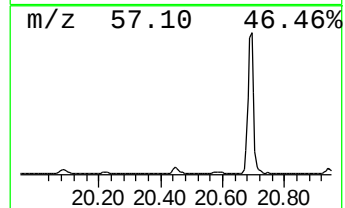
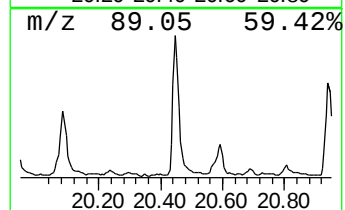
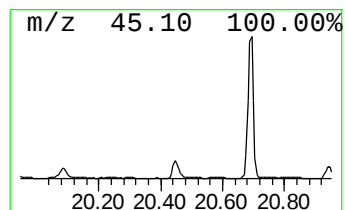
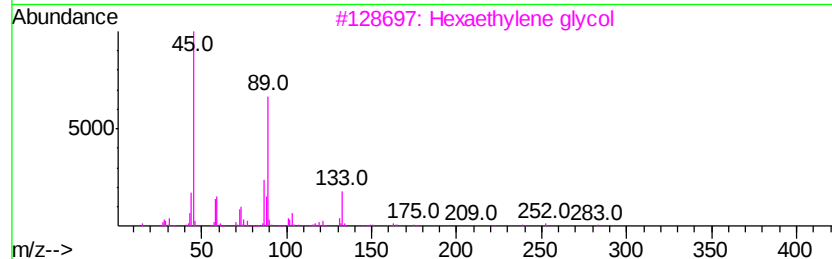
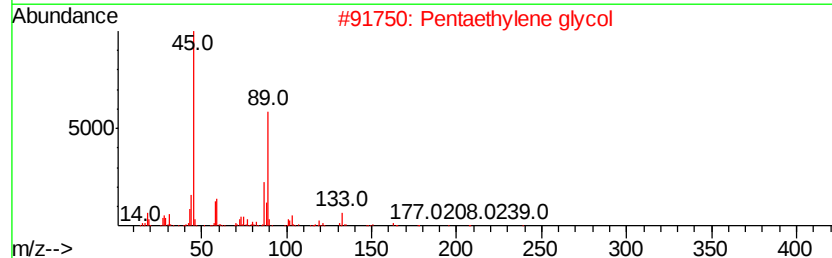
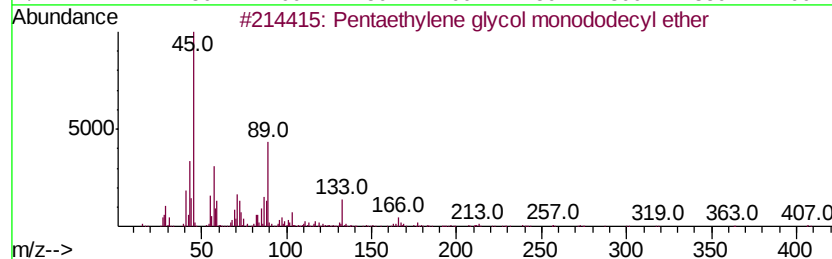
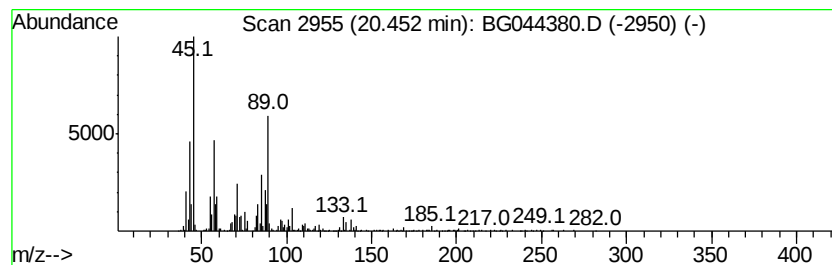
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 13 Pentaethylene glycol monodo... Concentration Rank 10

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 20.45 | 6.53 ng | 426497 | Chrysene-d12     | 21.53 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Pentaethylene glycol monododecyl... | 406 | C22H46O6  | 003055-95-6  | 80   |
| 2    |    |   | Pentaethylene glycol                | 238 | C10H22O6  | 004792-15-8  | 72   |
| 3    |    |   | Hexaethylene glycol                 | 282 | C12H26O7  | 002615-15-8  | 64   |
| 4    |    |   | 2-[2-[2-[2-[2-[2-(2-Hydroxyvet...   | 370 | C16H34O9  | 005117-19-1  | 64   |
| 5    |    |   | 2-[2-[2-[2-[2-[2-[2-(2-Hydrox...    | 414 | C18H38O10 | 1000352-12-5 | 58   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021020\  
Data File : BG044380.D  
Acq On : 11 Feb 2020 8:27  
Operator : CG/JU  
Sample : L1446-01  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
HP-640

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG013020.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

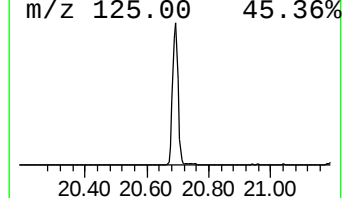
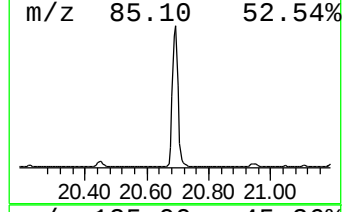
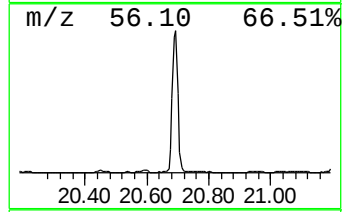
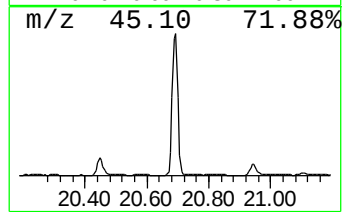
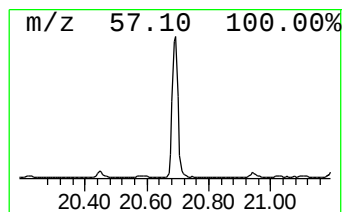
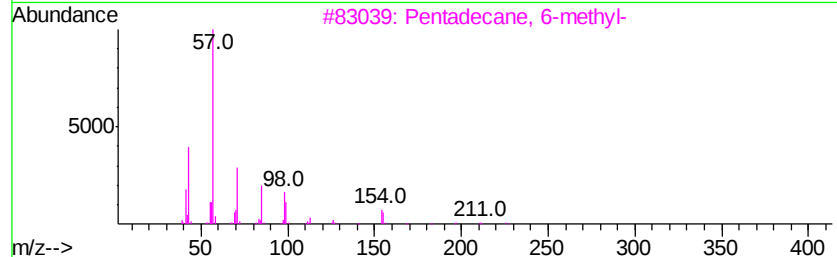
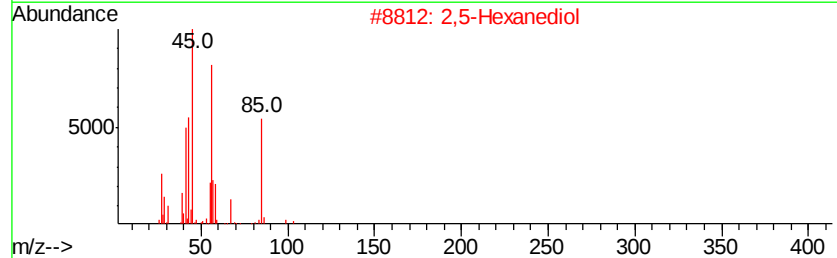
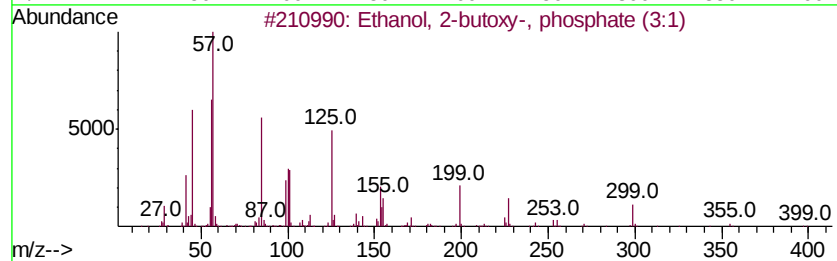
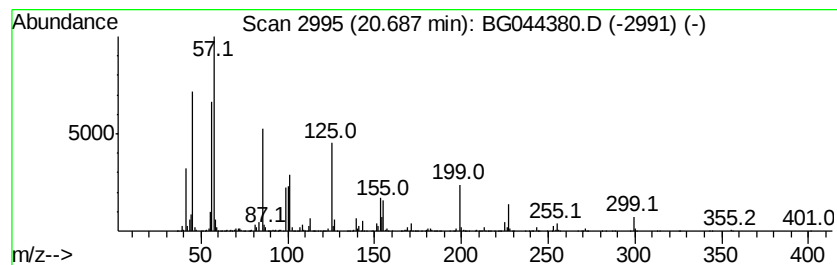
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 14 Ethanol, 2-butoxy-, phospho... Concentration Rank 4

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.  |
|-------|----------|---------|------------------|-------|
| 20.69 | 69.71 ng | 4552170 | Chrysene-d12     | 21.53 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | Ethanol, 2-butoxy-, phosphate (3:1) | 398 | C18H39O7P | 000078-51-3  | 94   |
| 2    |    | 2,5-Hexanediol                      | 118 | C6H14O2   | 002935-44-6  | 35   |
| 3    |    | Pentadecane, 6-methyl-              | 226 | C16H34    | 010105-38-1  | 14   |
| 4    |    | 2-Propanone, ethylhydrazone         | 100 | C5H12N2   | 007422-99-3  | 12   |
| 5    |    | Methoxyacetic acid, 4-tridecyl e... | 272 | C16H32O3  | 1000282-04-7 | 12   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021020\  
Data File : BG044380.D  
Acq On : 11 Feb 2020 8:27  
Operator : CG/JU  
Sample : L1446-01  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
HP-640

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG013020.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

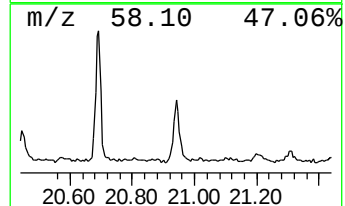
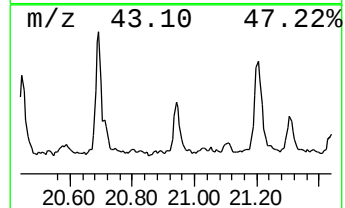
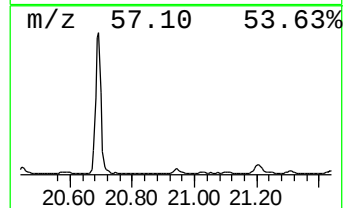
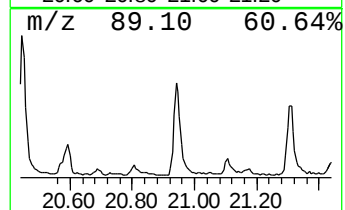
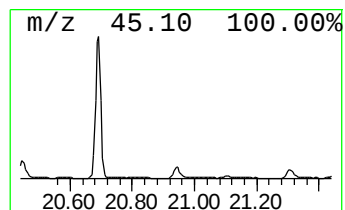
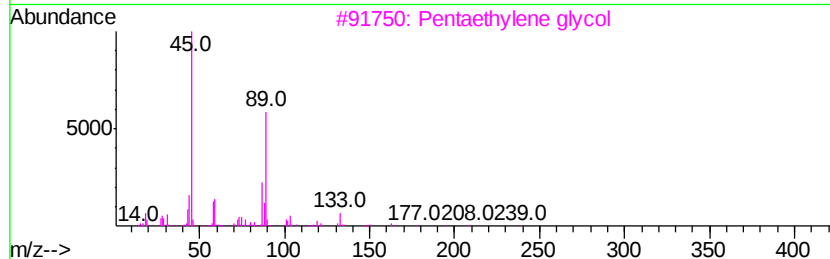
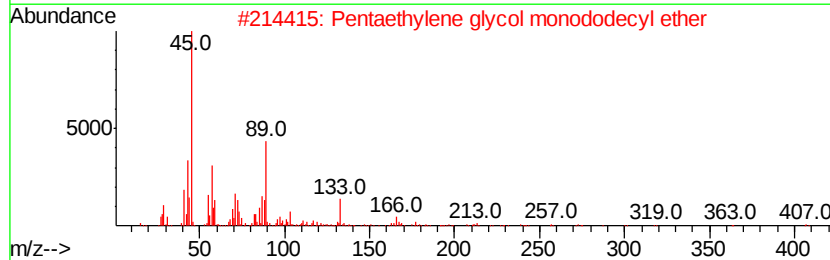
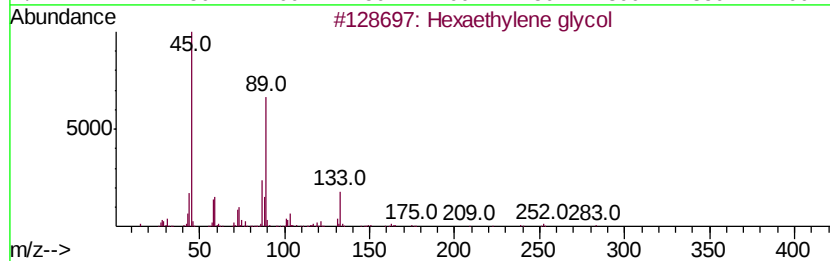
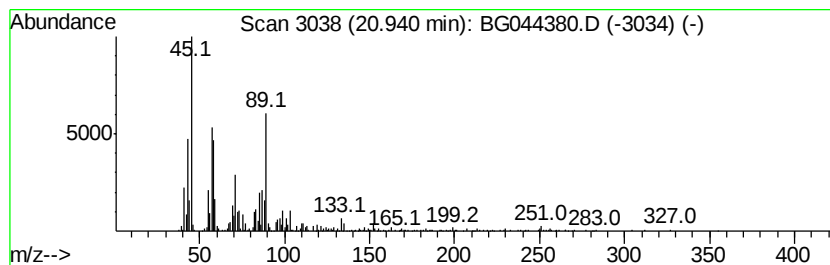
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 15 Hexaethylene glycol Concentration Rank 15

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 20.94 | 5.18 ng | 338270 | Chrysene-d12     | 21.53 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | Hexaethylene glycol                 | 282 | C12H26O7 | 002615-15-8 | 68   |
| 2    |    | Pentaethylene glycol monododecyl... | 406 | C22H46O6 | 003055-95-6 | 64   |
| 3    |    | Pentaethylene glycol                | 238 | C10H22O6 | 004792-15-8 | 62   |
| 4    |    | Heptaethylene glycol                | 326 | C14H30O8 | 005617-32-3 | 58   |
| 5    |    | 2-[2-[2-[2-[2-[2-(2-Hydroxyet...    | 370 | C16H34O9 | 005117-19-1 | 58   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021020\  
Data File : BG044380.D  
Acq On : 11 Feb 2020 8:27  
Operator : CG/JU  
Sample : L1446-01  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
HP-640

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG013020.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

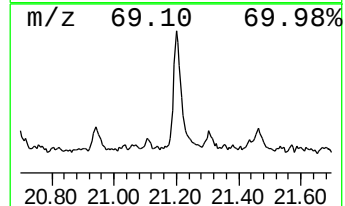
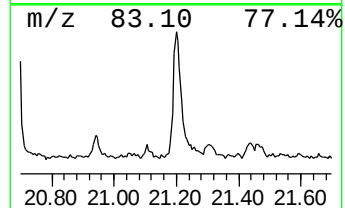
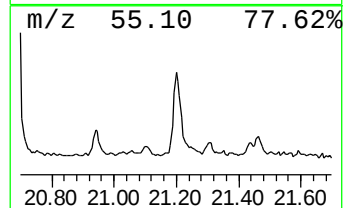
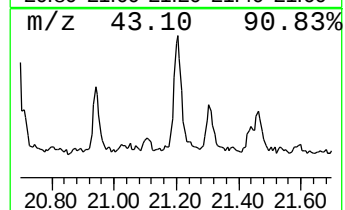
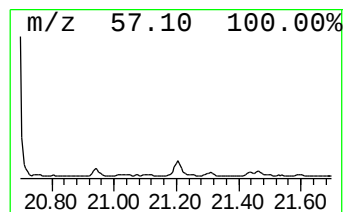
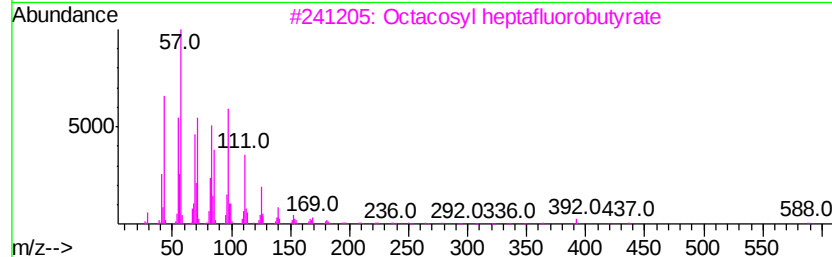
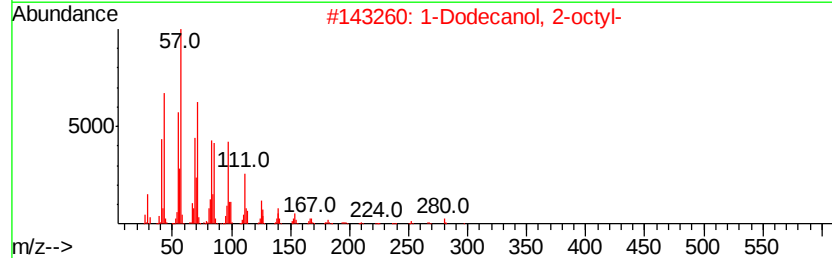
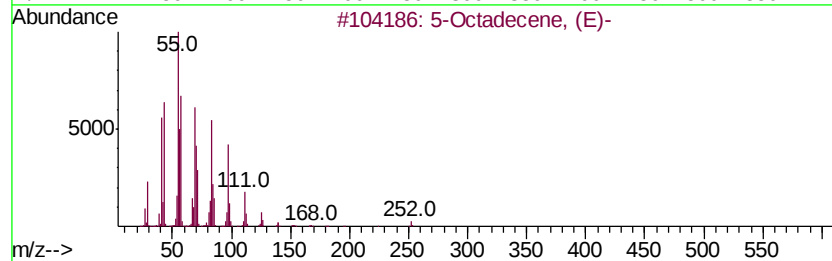
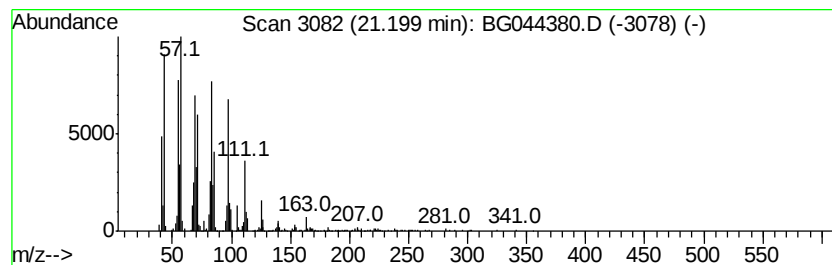
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 16 5-Octadecene, (E)- Concentration Rank 8

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 21.20 | 7.53 ng | 491444 | Chrysene-d12     | 21.53 |

| Hit# | of | 5 | Tentative ID                  | MW  | MolForm     | CAS#         | Qual |
|------|----|---|-------------------------------|-----|-------------|--------------|------|
| 1    |    |   | 5-Octadecene, (E)-            | 252 | C18H36      | 007206-21-5  | 90   |
| 2    |    |   | 1-Dodecanol, 2-octyl-         | 298 | C20H42O     | 005333-42-6  | 87   |
| 3    |    |   | Octacosyl heptafluorobutylate | 606 | C32H57F7O2  | 1000351-83-6 | 74   |
| 4    |    |   | 1-Hexadecanesulfonyl chloride | 324 | C16H33ClO2S | 038775-38-1  | 74   |
| 5    |    |   | Hexacosyl trifluoroacetate    | 478 | C28H53F3O2  | 1000351-75-0 | 74   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021020\  
Data File : BG044380.D  
Acq On : 11 Feb 2020 8:27  
Operator : CG/JU  
Sample : L1446-01  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
HP-640

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG013020.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

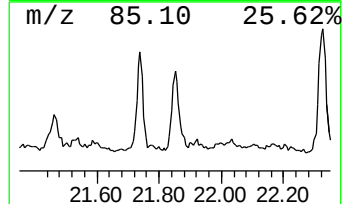
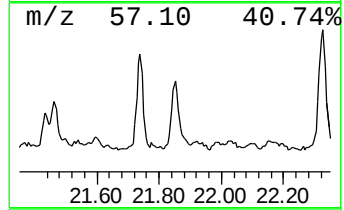
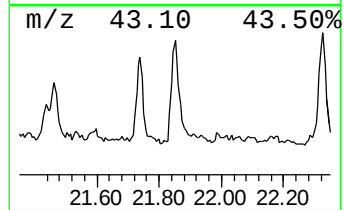
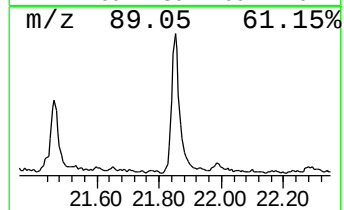
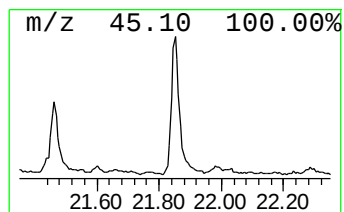
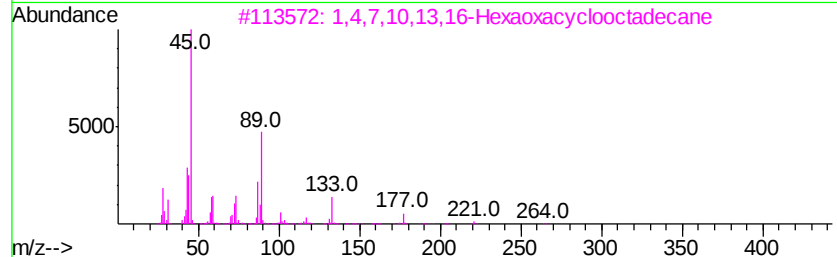
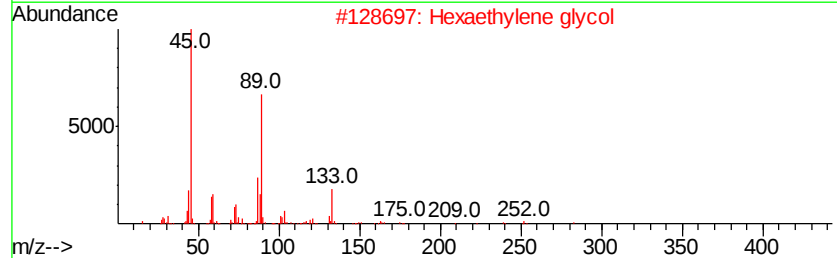
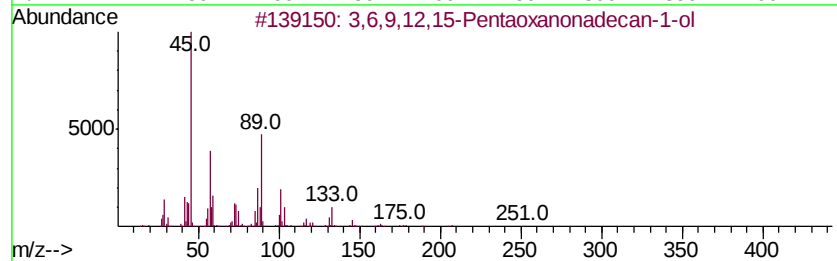
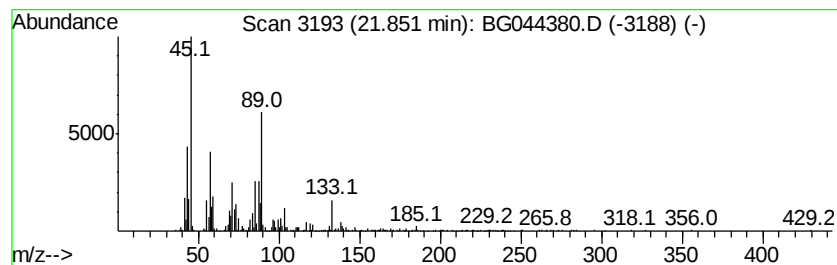
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 17 3,6,9,12,15-Pentaoxanonadec... Concentration Rank 7

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 21.85 | 7.68 ng | 501254 | Chrysene-d12     | 21.53 |

| Hit# | of                                      | 5   | Tentative ID | MW           | MolForm | CAS# | Qual |
|------|-----------------------------------------|-----|--------------|--------------|---------|------|------|
| 1    | 3,6,9,12,15-Pentaoxanonadecan-1-ol      | 294 | C14H30O6     | 001786-94-3  | 83      |      |      |
| 2    | Hexaethylene glycol                     | 282 | C12H26O7     | 002615-15-8  | 74      |      |      |
| 3    | 1,4,7,10,13,16-Hexaoxacyclooctadecane   | 264 | C12H24O6     | 017455-13-9  | 72      |      |      |
| 4    | 2-[2-[2-[2-[2-[2-[2-(2-Hydroxyl)]]]]]]] | 414 | C18H38O10    | 1000352-12-5 | 72      |      |      |
| 5    | 2-[2-[2-[2-[2-[2-[2-(2-Hydroxyl)]]]]]]] | 458 | C20H42O11    | 1000352-12-6 | 68      |      |      |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021020\  
Data File : BG044380.D  
Acq On : 11 Feb 2020 8:27  
Operator : CG/JU  
Sample : L1446-01  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
HP-640

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG013020.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

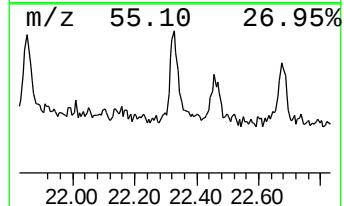
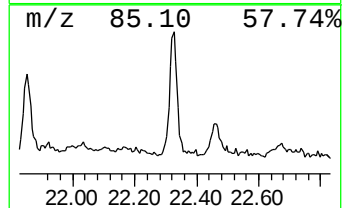
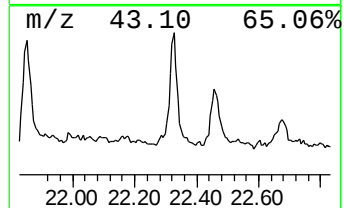
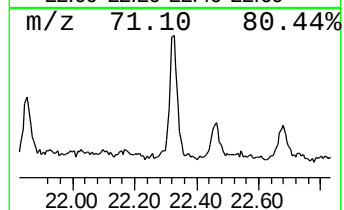
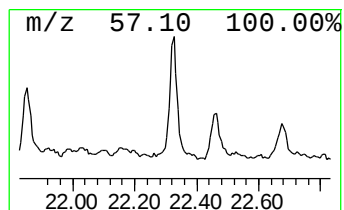
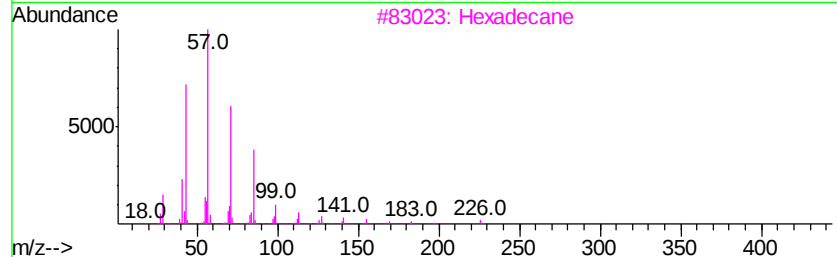
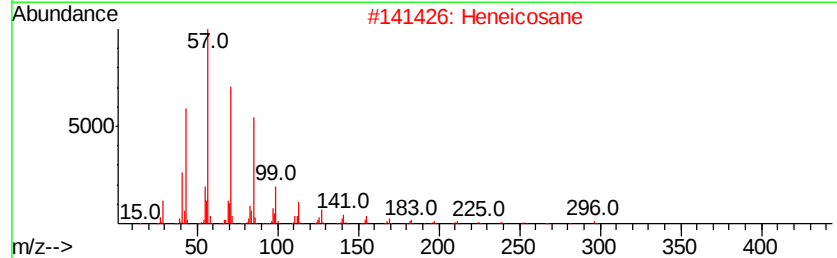
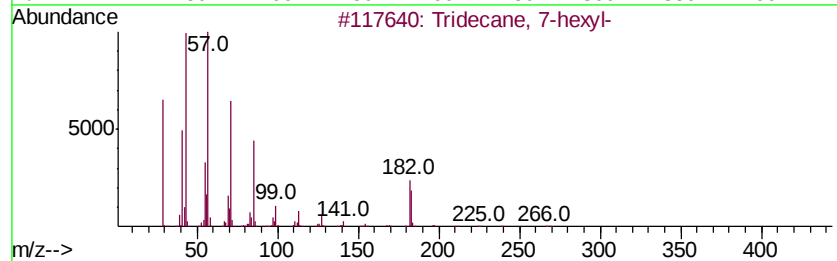
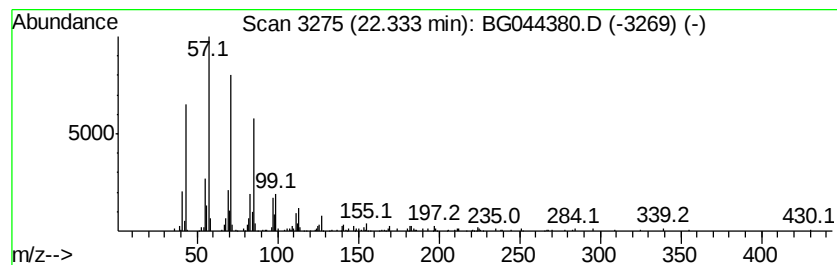
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 18 Tridecane, 7-hexyl- Concentration Rank 18

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 22.33 | 4.67 ng | 304737 | Chrysene-d12     | 21.53 |

| Hit# | of | 5 | Tentative ID         | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------|-----|---------|-------------|------|
| 1    |    |   | Tridecane, 7-hexyl-  | 268 | C19H40  | 007225-66-3 | 94   |
| 2    |    |   | Heneicosane          | 296 | C21H44  | 000629-94-7 | 91   |
| 3    |    |   | Hexadecane           | 226 | C16H34  | 000544-76-3 | 91   |
| 4    |    |   | Tetratetracontane    | 619 | C44H90  | 007098-22-8 | 91   |
| 5    |    |   | Tridecane, 6-propyl- | 226 | C16H34  | 055045-10-8 | 91   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021020\  
Data File : BG044380.D  
Acq On : 11 Feb 2020 8:27  
Operator : CG/JU  
Sample : L1446-01  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

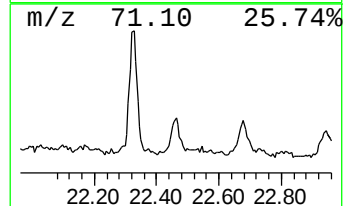
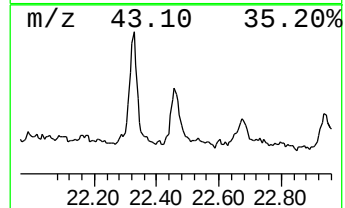
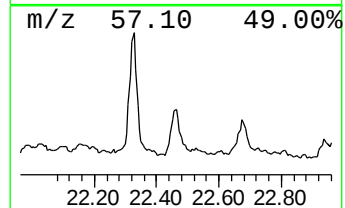
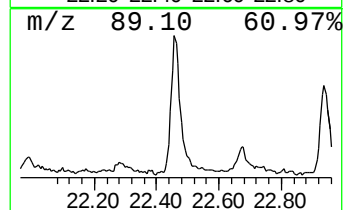
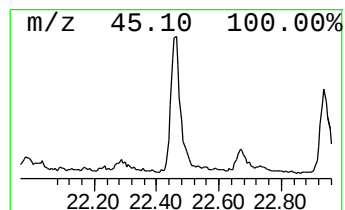
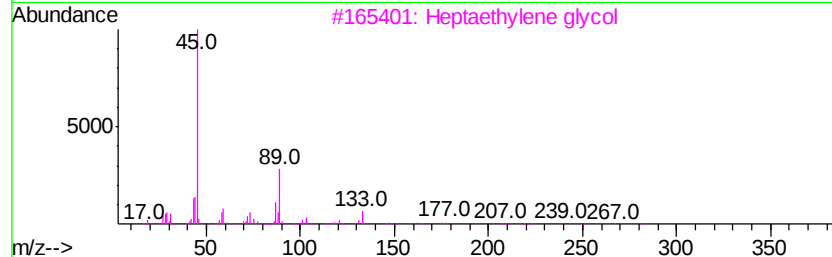
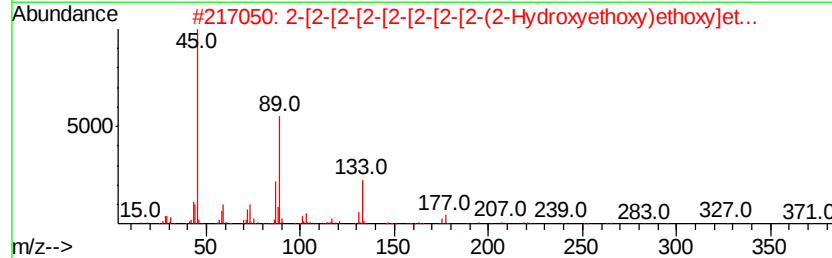
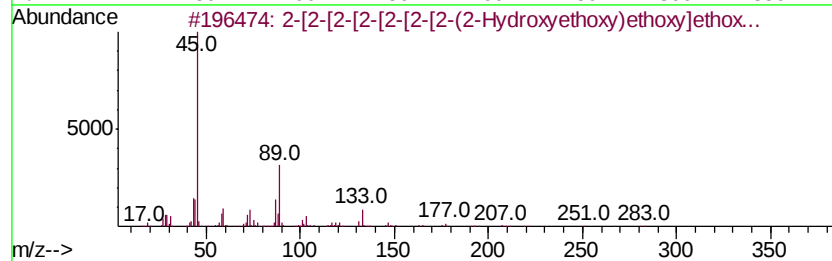
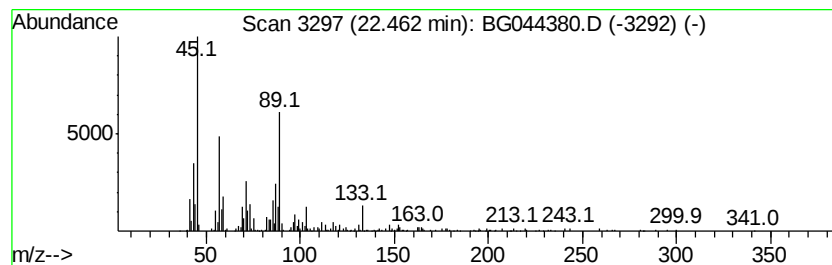
Instrument :  
BNA\_G  
ClientSampleId :  
HP-640

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG013020.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 19 2-[2-[2-[2-[2-[2-(2-Hydr... Concentration Rank 19

| R.T.    | EstConc                            | Area          | Relative to ISTD | R.T.      |
|---------|------------------------------------|---------------|------------------|-----------|
| 22.46   | 4.44 ng                            | 289860        | Chrysene-d12     | 21.53     |
| Hit# of | 5                                  | Tentative ID  | MW MolForm       | CAS# Qual |
| 1       | 2-[2-[2-[2-[2-[2-(2-Hydroxvet...   | 370 C16H3409  | 005117-19-1      | 90        |
| 2       | 2-[2-[2-[2-[2-[2-[2-(2-Hydrox...   | 414 C18H38010 | 1000352-12-5     | 86        |
| 3       | Heptaethylene glycol               | 326 C14H3008  | 005617-32-3      | 86        |
| 4       | 3,6,9,12,15-Pentaoxanonadecan-1-ol | 294 C14H3006  | 001786-94-3      | 86        |
| 5       | 2-[2-[2-[2-[2-[2-[2-[2-[2-(2-...   | 502 C22H46012 | 1000352-13-2     | 83        |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021020\  
Data File : BG044380.D  
Acq On : 11 Feb 2020 8:27  
Operator : CG/JU  
Sample : L1446-01  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

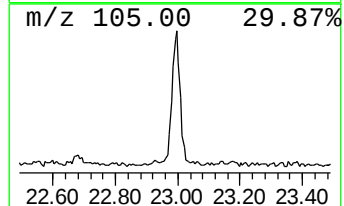
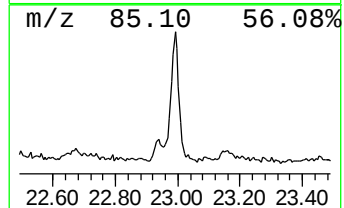
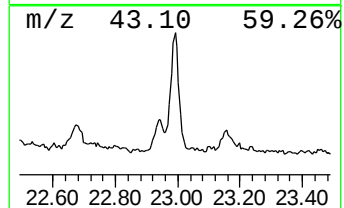
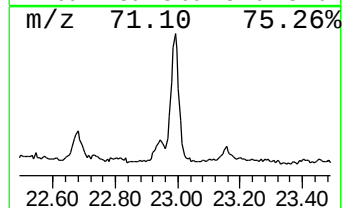
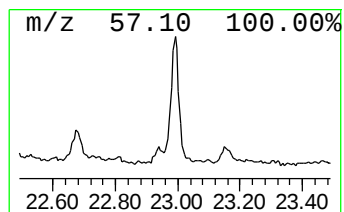
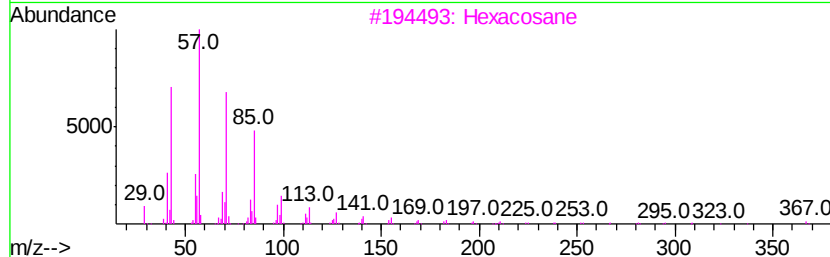
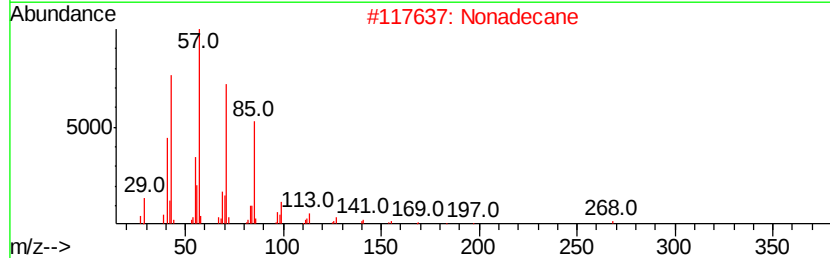
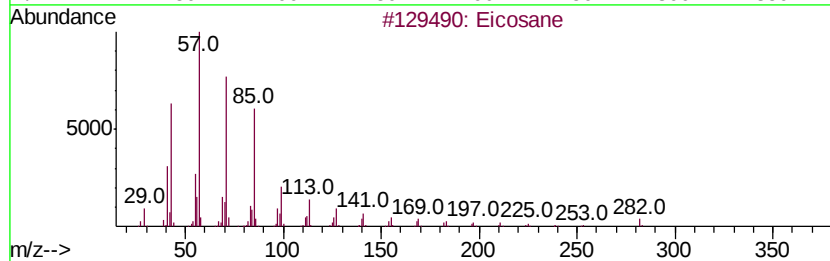
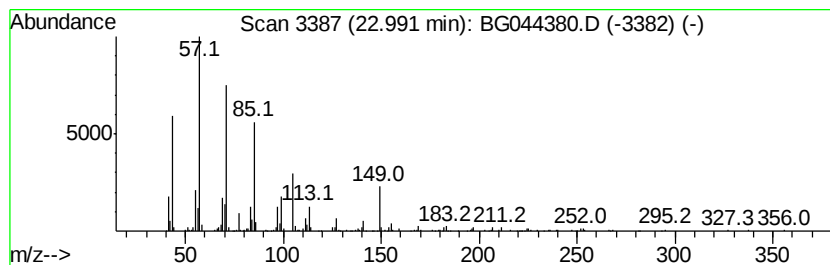
Instrument :  
BNA\_G  
ClientSampleId :  
HP-640

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG013020.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 20 Eicosane Concentration Rank 14

| R.T.    | EstConc     | Area         | Relative to ISTD | R.T.      |
|---------|-------------|--------------|------------------|-----------|
| 22.99   | 5.67 ng     | 370546       | Chrysene-d12     | 21.53     |
| Hit# of | 5           | Tentative ID | MW MolForm       | CAS# Qual |
| 1       | Eicosane    | 282 C20H42   | 000112-95-8      | 90        |
| 2       | Nonadecane  | 268 C19H40   | 000629-92-5      | 81        |
| 3       | Hexacosane  | 366 C26H54   | 000630-01-3      | 81        |
| 4       | Octadecane  | 254 C18H38   | 000593-45-3      | 76        |
| 5       | Tetracosane | 338 C24H50   | 000646-31-1      | 74        |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_G\DATA\BG021020\  
 Data File : BG044380.D  
 Acq On : 11 Feb 2020 8:27  
 Operator : CG/JU  
 Sample : L1446-01  
 Misc :  
 ALS Vial : 24 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 HP-640

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\8270-BG013020.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT 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 |
| Hexanoic acid, 2-... | 9.35  | 64.5    | ng    | 2475460  | 2                     | 10.71 | 767886  | 20.0 |
| unknown10.76         | 10.76 | 6.2     | ng    | 236459   | 2                     | 10.71 | 767886  | 20.0 |
| 2-Propanol, 1-(2-... | 11.25 | 114.7   | ng    | 4403880  | 2                     | 10.71 | 767886  | 20.0 |
| 1-Propanol, 2-(2-... | 11.32 | 126.0   | ng    | 4838760  | 2                     | 10.71 | 767886  | 20.0 |
| Hexaethylene glyc... | 11.57 | 8.3     | ng    | 320119   | 2                     | 10.71 | 767886  | 20.0 |
| Cyclobutanol         | 14.59 | 7.0     | ng    | 431407   | 3                     | 14.54 | 1237610 | 20.0 |
| n-Hexadecanoic acid  | 18.20 | 4.2     | ng    | 257173   | 4                     | 17.29 | 1229620 | 20.0 |
| unknown18.47         | 18.47 | 4.8     | ng    | 296433   | 4                     | 17.29 | 1229620 | 20.0 |
| 9-Octadecenoic ac... | 19.35 | 6.3     | ng    | 386748   | 4                     | 17.29 | 1229620 | 20.0 |
| Triethylene glyco... | 20.08 | 4.9     | ng    | 318470   | 5                     | 21.53 | 1305970 | 20.0 |
| Pentaethylene gly... | 20.45 | 6.5     | ng    | 426497   | 5                     | 21.53 | 1305970 | 20.0 |
| Ethanol, 2-butoxy... | 20.69 | 69.7    | ng    | 4552170  | 5                     | 21.53 | 1305970 | 20.0 |
| Hexaethylene glycol  | 20.94 | 5.2     | ng    | 338270   | 5                     | 21.53 | 1305970 | 20.0 |
| 5-Octadecene, (E)-   | 21.20 | 7.5     | ng    | 491444   | 5                     | 21.53 | 1305970 | 20.0 |
| 3,6,9,12,15-Penta... | 21.85 | 7.7     | ng    | 501254   | 5                     | 21.53 | 1305970 | 20.0 |
| Tridecane, 7-hexyl-  | 22.33 | 4.7     | ng    | 304737   | 5                     | 21.53 | 1305970 | 20.0 |
| 2-[2-[2-[2-[2-...    | 22.46 | 4.4     | ng    | 289860   | 5                     | 21.53 | 1305970 | 20.0 |
| Eicosane             | 22.99 | 5.7     | ng    | 370546   | 5                     | 21.53 | 1305970 | 20.0 |