

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010719\  
 Data File : BG038973.D  
 Acq On : 7 Jan 2019 19:12  
 Operator : JU/SJ  
 Sample : K1036-05  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 13-E

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-BG121218.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         | 5.602       | 421           | 428         | 441          | rBV      | 362759         | 604736        | 10.85%          | 2.599%        |
| 2         | 7.206       | 693           | 701         | 713          | rBV      | 382118         | 692358        | 12.43%          | 2.976%        |
| 3         | 7.576       | 756           | 764         | 778          | rVB      | 383668         | 679398        | 12.20%          | 2.920%        |
| 4         | 8.046       | 838           | 844         | 851          | rVB      | 120348         | 208999        | 3.75%           | 0.898%        |
| 5         | 8.369       | 891           | 899         | 908          | rVB      | 298117         | 538754        | 9.67%           | 2.316%        |
| 6         | 9.227       | 1034          | 1045        | 1057         | rBV      | 247186         | 488478        | 8.77%           | 2.100%        |
| 7         | 10.784      | 1302          | 1310        | 1316         | rBV      | 62669          | 101199        | 1.82%           | 0.435%        |
| 8         | 10.861      | 1316          | 1323        | 1332         | rVB      | 159871         | 301057        | 5.40%           | 1.294%        |
| 9         | 13.299      | 1731          | 1738        | 1747         | rBV      | 700865         | 1139622       | 20.46%          | 4.898%        |
| 10        | 13.463      | 1759          | 1766        | 1771         | rBV      | 109206         | 155236        | 2.79%           | 0.667%        |
| 11        | 14.680      | 1962          | 1973        | 1981         | rBV4     | 306530         | 577389        | 10.36%          | 2.482%        |
| 12        | 15.150      | 2047          | 2053        | 2059         | rVB      | 415076         | 574726        | 10.32%          | 2.470%        |
| 13        | 15.485      | 2105          | 2110        | 2114         | rBV      | 156125         | 214918        | 3.86%           | 0.924%        |
| 14        | 15.866      | 2169          | 2175        | 2180         | rBV2     | 49306          | 72678         | 1.30%           | 0.312%        |
| 15        | 16.037      | 2200          | 2204        | 2212         | rVB2     | 57837          | 103770        | 1.86%           | 0.446%        |
| 16        | 16.172      | 2220          | 2227        | 2235         | rBV      | 518913         | 848828        | 15.24%          | 3.648%        |
| 17        | 16.272      | 2240          | 2244        | 2251         | rVV      | 41050          | 69979         | 1.26%           | 0.301%        |
| 18        | 16.342      | 2251          | 2256        | 2267         | rVB2     | 38846          | 67572         | 1.21%           | 0.290%        |
| 19        | 17.089      | 2379          | 2383        | 2387         | rVV2     | 55828          | 77562         | 1.39%           | 0.333%        |
| 20        | 17.136      | 2387          | 2391        | 2395         | rVV      | 178896         | 248323        | 4.46%           | 1.067%        |
| 21        | 17.429      | 2435          | 2441        | 2454         | rVB2     | 303311         | 499710        | 8.97%           | 2.148%        |
| 22        | 17.611      | 2465          | 2472        | 2480         | rBV      | 94724          | 140653        | 2.52%           | 0.605%        |
| 23        | 17.876      | 2513          | 2517        | 2523         | rBV      | 42090          | 60005         | 1.08%           | 0.258%        |
| 24        | 18.293      | 2583          | 2588        | 2594         | rBV2     | 187773         | 271245        | 4.87%           | 1.166%        |
| 25        | 18.381      | 2599          | 2603        | 2608         | rVV      | 41576          | 64309         | 1.15%           | 0.276%        |
| 26        | 18.440      | 2608          | 2613        | 2617         | rVV4     | 33312          | 62913         | 1.13%           | 0.270%        |
| 27        | 18.528      | 2624          | 2628        | 2631         | rVV      | 62358          | 92268         | 1.66%           | 0.397%        |
| 28        | 18.563      | 2631          | 2634        | 2638         | rVV      | 161972         | 219853        | 3.95%           | 0.945%        |
| 29        | 18.604      | 2638          | 2641        | 2646         | rVB      | 41255          | 57984         | 1.04%           | 0.249%        |
| 30        | 18.969      | 2696          | 2703        | 2707         | rBV      | 107360         | 143638        | 2.58%           | 0.617%        |
| 31        | 19.556      | 2800          | 2803        | 2811         | rVV6     | 45232          | 108878        | 1.95%           | 0.468%        |
| 32        | 19.674      | 2818          | 2823        | 2827         | rBV4     | 39662          | 67859         | 1.22%           | 0.292%        |
| 33        | 19.744      | 2832          | 2835        | 2837         | rVV      | 73176          | 98704         | 1.77%           | 0.424%        |
| 34        | 19.768      | 2837          | 2839        | 2843         | rVV      | 150576         | 170417        | 3.06%           | 0.732%        |

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 BNA\_G  
 ClientSampleId :  
 13-E

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-BG121218.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |         |         |         |         |
|----|--------|------|------|------|------|---------|---------|---------|---------|
| 35 | 19.815 | 2843 | 2847 | 2852 | rVB  | 41156   | 64527   | 1.16%   | 0.277%  |
| 36 | 20.032 | 2878 | 2884 | 2890 | rVV  | 1170574 | 1548613 | 27.80%  | 6.656%  |
| 37 | 20.109 | 2893 | 2897 | 2903 | rVB  | 118485  | 137894  | 2.48%   | 0.593%  |
| 38 | 20.244 | 2913 | 2920 | 2924 | rBV2 | 47448   | 70392   | 1.26%   | 0.303%  |
| 39 | 20.297 | 2924 | 2929 | 2934 | rBV2 | 54023   | 80581   | 1.45%   | 0.346%  |
| 40 | 20.726 | 2995 | 3002 | 3004 | rVV4 | 35367   | 68301   | 1.23%   | 0.294%  |
| 41 | 20.796 | 3008 | 3014 | 3018 | rVV2 | 132457  | 223686  | 4.02%   | 0.961%  |
| 42 | 20.849 | 3018 | 3023 | 3029 | rVB  | 36796   | 67504   | 1.21%   | 0.290%  |
| 43 | 21.107 | 3063 | 3067 | 3072 | rBV  | 97811   | 117767  | 2.11%   | 0.506%  |
| 44 | 21.295 | 3090 | 3099 | 3107 | rBV4 | 60252   | 153004  | 2.75%   | 0.658%  |
| 45 | 21.401 | 3113 | 3117 | 3124 | rVB2 | 47215   | 77372   | 1.39%   | 0.333%  |
| 46 | 21.554 | 3141 | 3143 | 3148 | rVB2 | 51880   | 65170   | 1.17%   | 0.280%  |
| 47 | 21.619 | 3149 | 3154 | 3160 | rVB  | 156336  | 247660  | 4.45%   | 1.064%  |
| 48 | 21.713 | 3160 | 3170 | 3178 | rBV  | 319893  | 617825  | 11.09%  | 2.655%  |
| 49 | 21.848 | 3178 | 3193 | 3198 | rBV5 | 102490  | 272303  | 4.89%   | 1.170%  |
| 50 | 21.924 | 3204 | 3206 | 3212 | rVB3 | 39954   | 57664   | 1.04%   | 0.248%  |
| 51 | 22.224 | 3251 | 3257 | 3260 | rBV  | 103554  | 210536  | 3.78%   | 0.905%  |
| 52 | 22.394 | 3265 | 3286 | 3313 | rVV3 | 756027  | 5571063 | 100.00% | 23.945% |
| 53 | 22.588 | 3315 | 3319 | 3326 | rVB3 | 38906   | 76161   | 1.37%   | 0.327%  |
| 54 | 23.158 | 3409 | 3416 | 3419 | rBV4 | 147633  | 359457  | 6.45%   | 1.545%  |
| 55 | 23.258 | 3426 | 3433 | 3458 | rVB4 | 375453  | 1678191 | 30.12%  | 7.213%  |
| 56 | 23.646 | 3493 | 3499 | 3504 | rBV2 | 50292   | 103643  | 1.86%   | 0.445%  |
| 57 | 23.963 | 3547 | 3553 | 3558 | rBV2 | 52066   | 111802  | 2.01%   | 0.481%  |
| 58 | 25.009 | 3721 | 3731 | 3740 | rVB2 | 166170  | 464949  | 8.35%   | 1.998%  |
| 59 | 25.608 | 3826 | 3833 | 3836 | rBV2 | 32098   | 74503   | 1.34%   | 0.320%  |
| 60 | 25.679 | 3837 | 3845 | 3859 | rVB2 | 135460  | 433773  | 7.79%   | 1.864%  |
| 61 | 26.055 | 3903 | 3909 | 3919 | rVB5 | 30335   | 95402   | 1.71%   | 0.410%  |
| 62 | 31.049 | 4750 | 4759 | 4761 | rBV4 | 36043   | 101482  | 1.82%   | 0.436%  |
| 63 | 32.136 | 4930 | 4944 | 4959 | rVB4 | 56264   | 290995  | 5.22%   | 1.251%  |

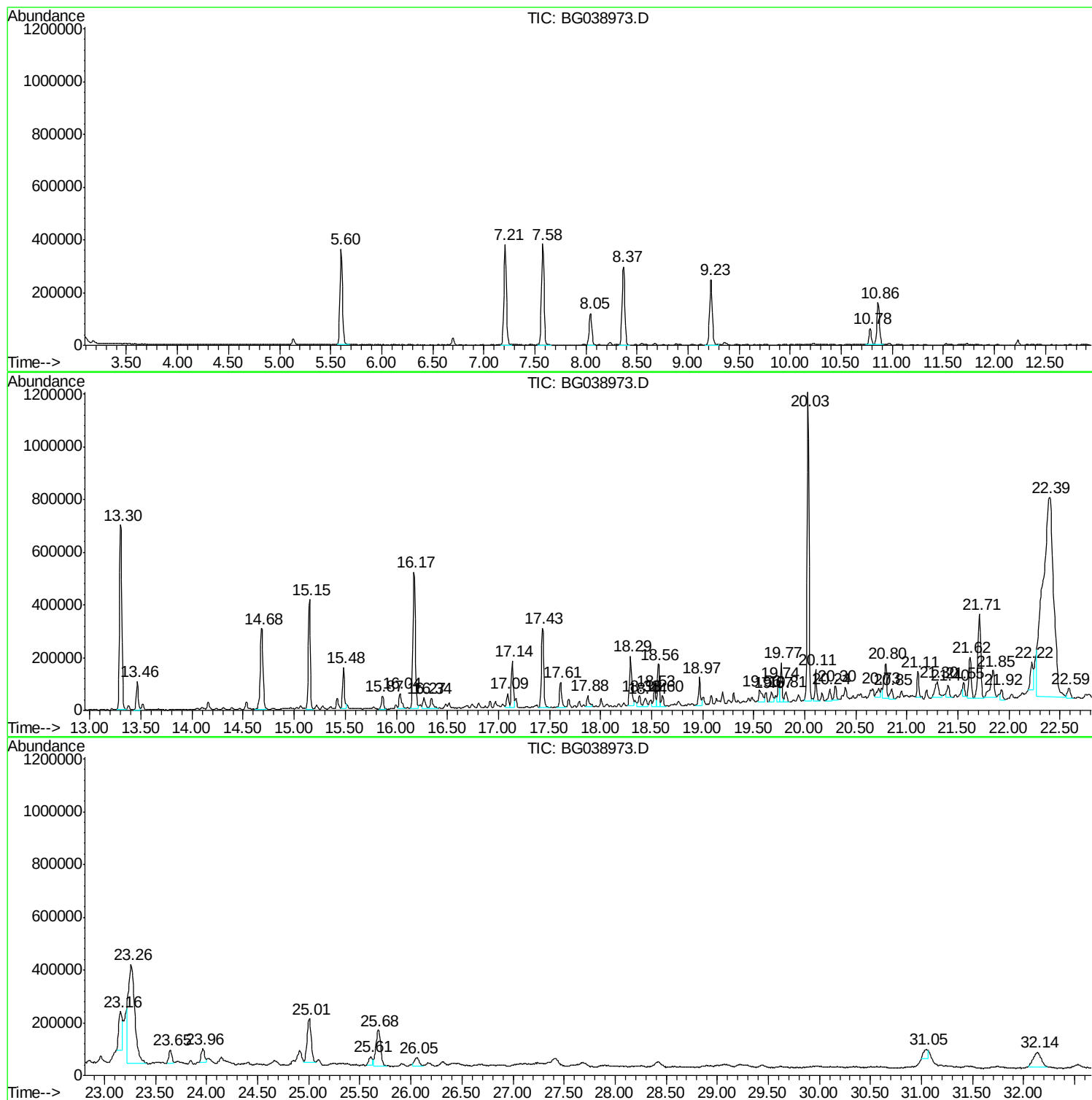
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Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
13-E

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG121218.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\BG010719\  
Data File : BG038973.D  
Acq On : 7 Jan 2019 19:12  
Operator : JU/SJ  
Sample : K1036-05  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
13-E

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG121218.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

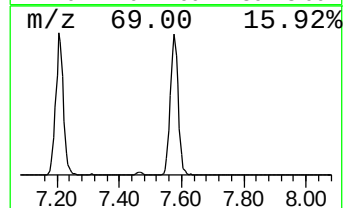
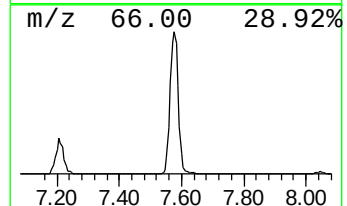
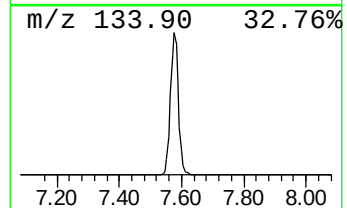
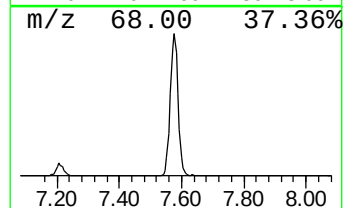
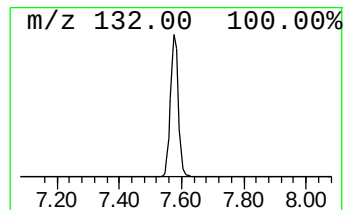
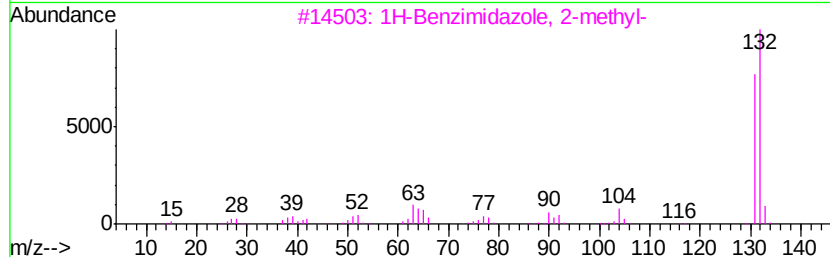
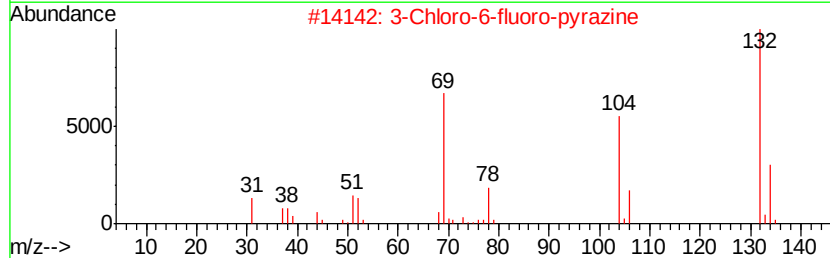
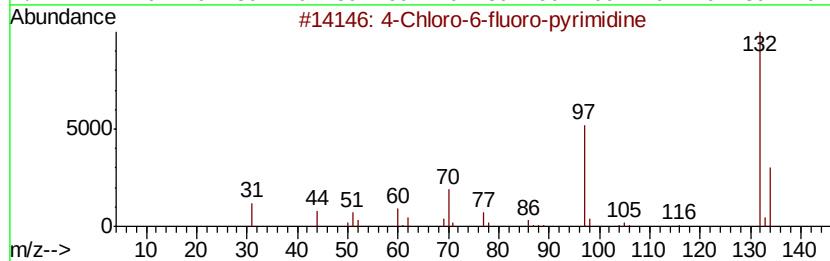
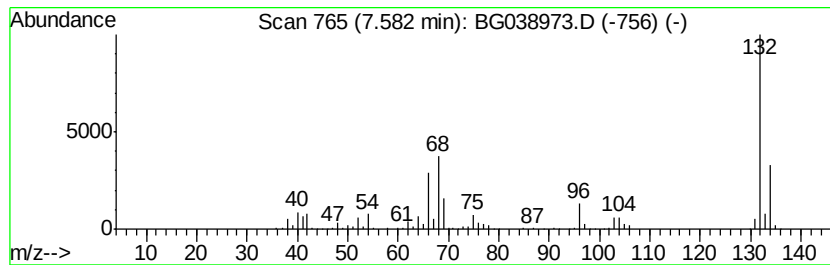
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 1 unknown7.58 Concentration Rank 2

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 7.58 | 65.01 ng | 679398 | 1,4-Dichlorobenzene-d4 | 8.05 |

| Hit# | of | Tentative ID                     | MW  | MolForm   | CAS#         | Qual |
|------|----|----------------------------------|-----|-----------|--------------|------|
| 1    | 5  | 4-Chloro-6-fluoro-pyrimidine     | 132 | C4H2ClFN2 | 051422-01-6  | 27   |
| 2    |    | 3-Chloro-6-fluoro-pyrazine       | 132 | C4H2ClFN2 | 1000146-10-7 | 27   |
| 3    |    | 1H-Benzimidazole, 2-methyl-      | 132 | C8H8N2    | 000615-15-6  | 22   |
| 4    |    | (E)-3-Chloro-2-methyl-2-pentenal | 132 | C6H9ClO   | 031357-76-3  | 12   |
| 5    |    | 5-Aminoindole                    | 132 | C8H8N2    | 005192-03-0  | 12   |



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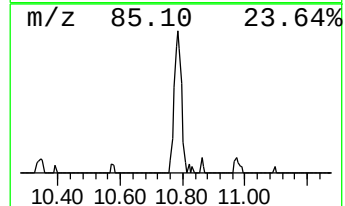
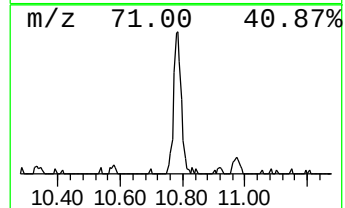
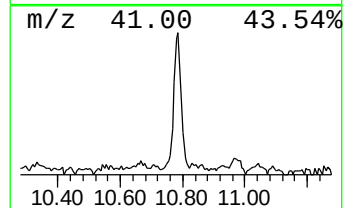
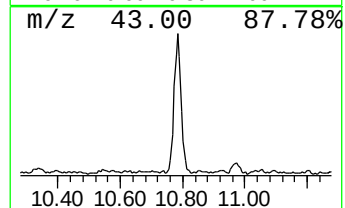
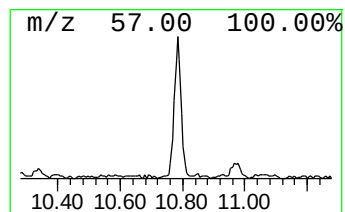
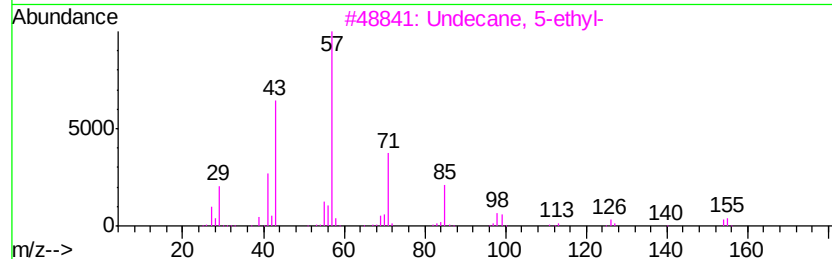
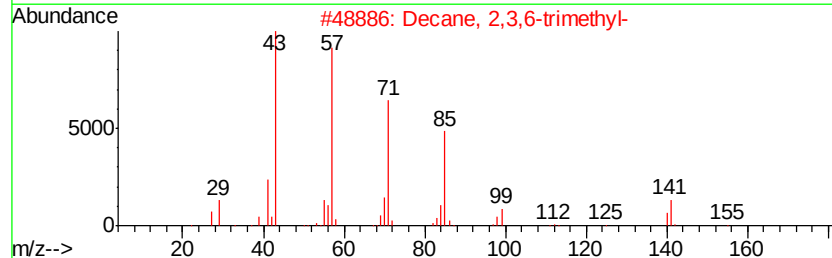
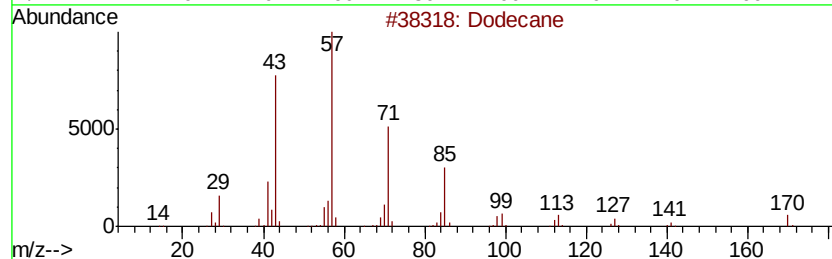
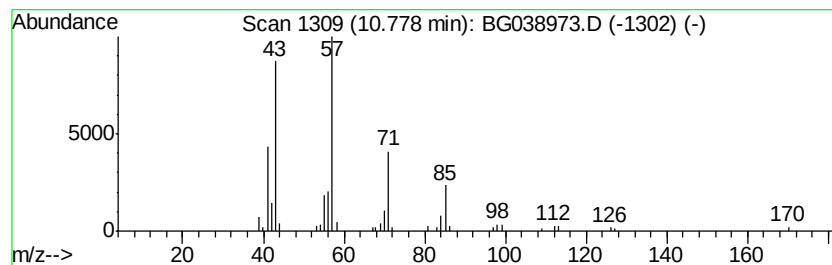
Instrument :  
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG121218.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 3 Dodecane Concentration Rank 17

| R.T.      | EstConc                  | Area   | Relative to ISTD |             | R.T.  |
|-----------|--------------------------|--------|------------------|-------------|-------|
| 10.78     | 6.72 ng                  | 101199 | Naphthalene-d8   |             | 10.86 |
| Hit# of 5 | Tentative ID             | MW     | MolForm          | CAS#        | Qual  |
| 1         | Dodecane                 | 170    | C12H26           | 000112-40-3 | 72    |
| 2         | Decane, 2,3,6-trimethyl- | 184    | C13H28           | 062238-12-4 | 64    |
| 3         | Undecane, 5-ethyl-       | 184    | C13H28           | 017453-94-0 | 59    |
| 4         | Nonane                   | 128    | C9H20            | 000111-84-2 | 59    |
| 5         | 1-Iodoundecane           | 282    | C11H23I          | 004282-44-4 | 59    |



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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG121218.M  
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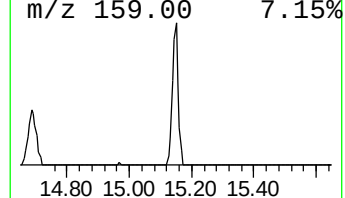
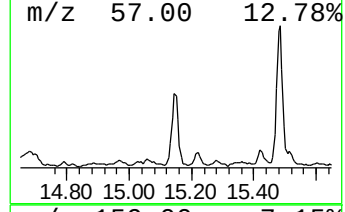
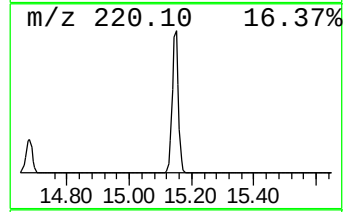
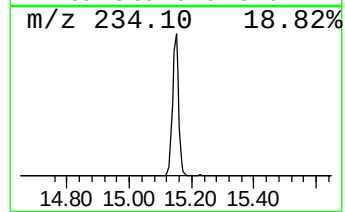
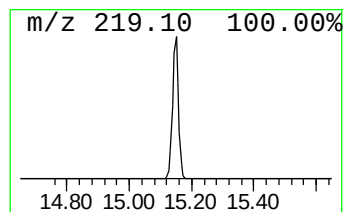
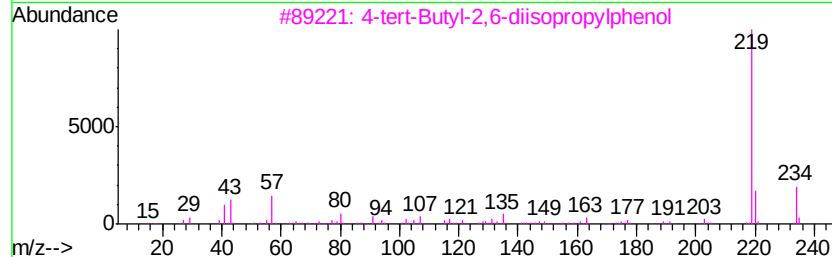
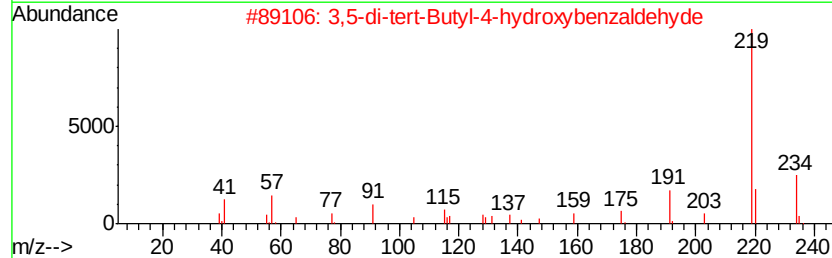
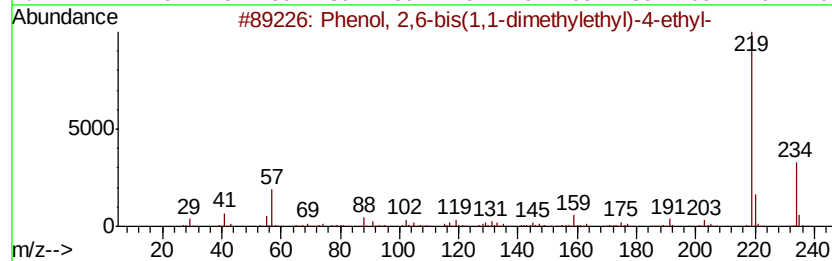
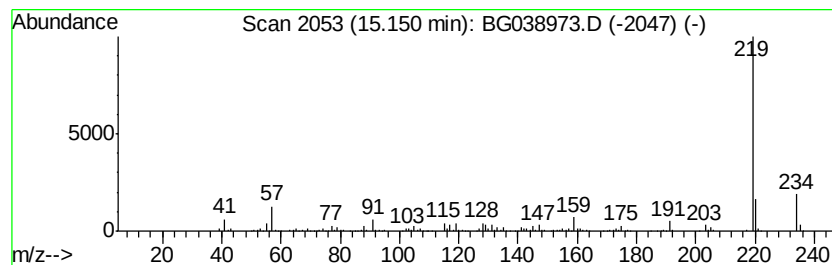
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 4 Phenol, 2,6-bis(1,1-dimethy... Concentration Rank 5

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 15.15 | 19.91 ng | 574726 | Acenaphthene-d10 | 14.69 |

| Hit# of | 5                                   | Tentative ID | MW         | MolForm      | CAS# | Qual |
|---------|-------------------------------------|--------------|------------|--------------|------|------|
| 1       | Phenol, 2,6-bis(1,1-dimethylethy... | 234          | C16H26O    | 004130-42-1  | 91   |      |
| 2       | 3,5-di-tert-Butyl-4-hydroxybenza... | 234          | C15H22O2   | 001620-98-0  | 90   |      |
| 3       | 4-tert-Butyl-2,6-diisopropylphenol  | 234          | C16H26O    | 057354-65-1  | 74   |      |
| 4       | 3,5-Di-tert-butylbenzoic acid       | 234          | C15H22O2   | 016225-26-6  | 72   |      |
| 5       | 5,8-Methano-1H-[1,2,4]triazolo[1... | 395          | C25H21N3O2 | 1000142-91-6 | 64   |      |



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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG121218.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

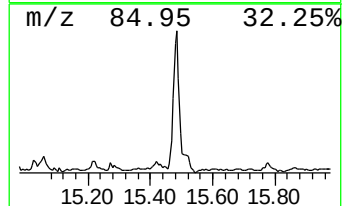
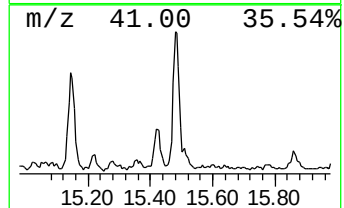
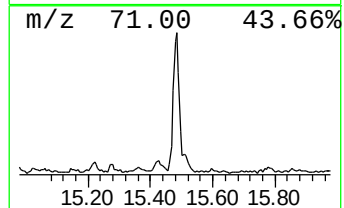
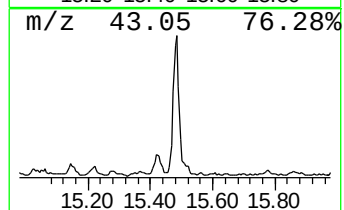
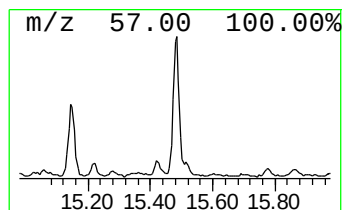
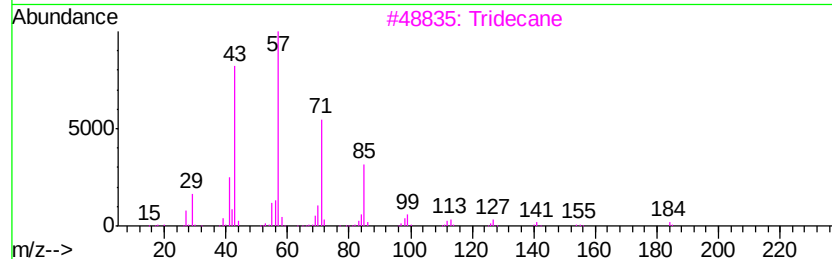
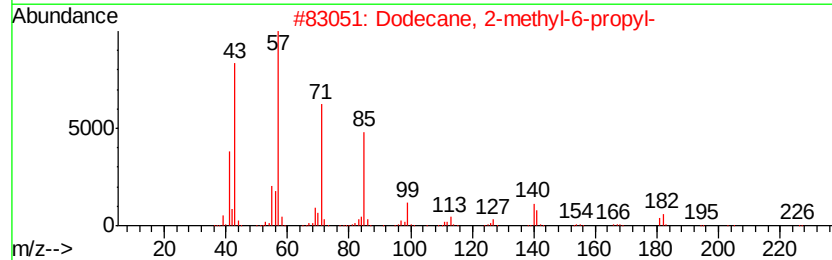
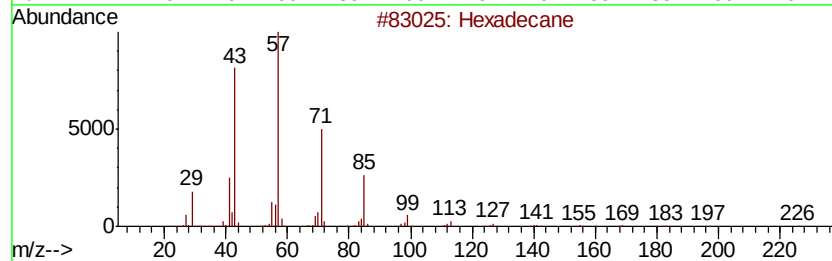
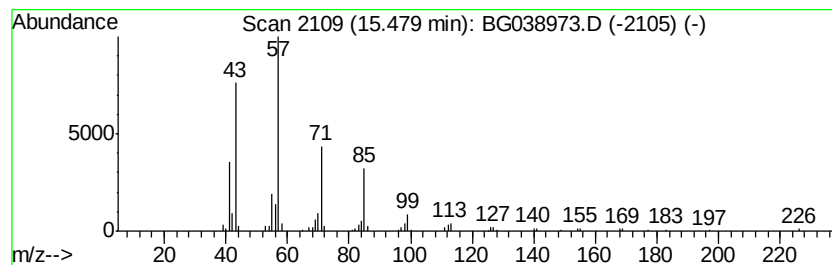
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 TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
 Peak Number 5 Hexadecane Concentration Rank 14

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 15.48 | 7.44 ng | 214918 | Acenaphthene-d10 | 14.69 |

| Hit# of | 5                                   | Tentative ID | MW  | MolForm   | CAS#         | Qual |
|---------|-------------------------------------|--------------|-----|-----------|--------------|------|
| 1       | Hexadecane                          |              | 226 | C16H34    | 000544-76-3  | 97   |
| 2       | Dodecane, 2-methyl-6-propyl-        |              | 226 | C16H34    | 055045-08-4  | 90   |
| 3       | Tridecane                           |              | 184 | C13H28    | 000629-50-5  | 86   |
| 4       | Tetradecane                         |              | 198 | C14H30    | 000629-59-4  | 72   |
| 5       | Sulfurous acid, pentadecyl 2-pro... |              | 334 | C18H38O3S | 1000309-12-6 | 72   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010719\  
Data File : BG038973.D  
Acq On : 7 Jan 2019 19:12  
Operator : JU/SJ  
Sample : K1036-05  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
13-E

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG121218.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

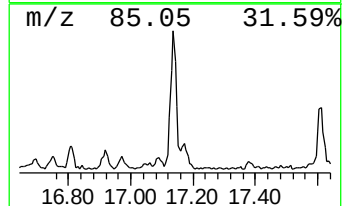
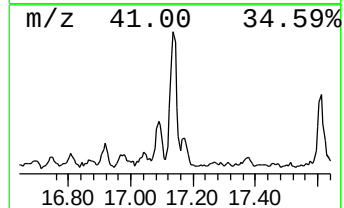
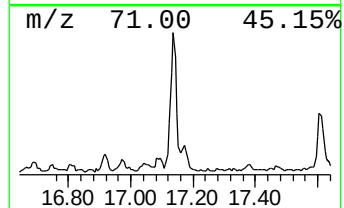
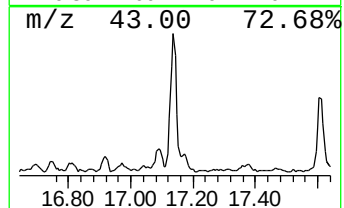
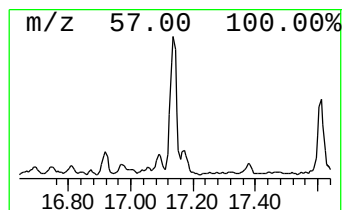
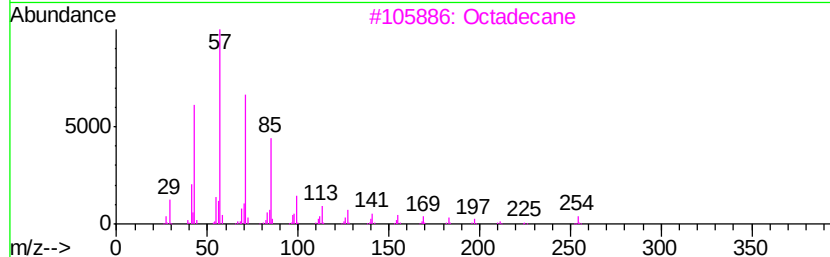
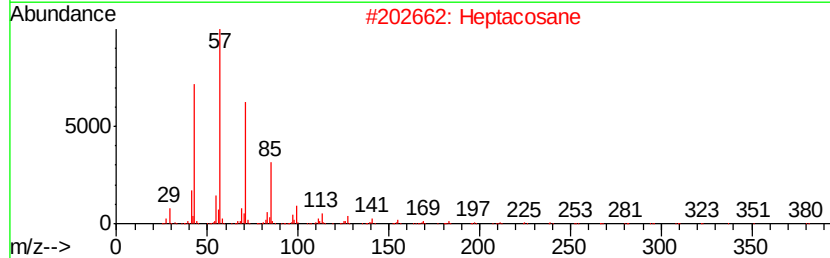
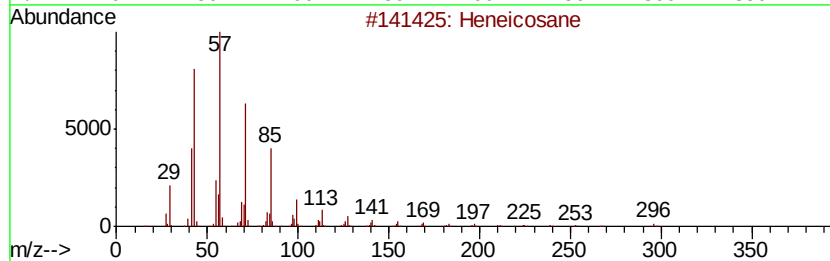
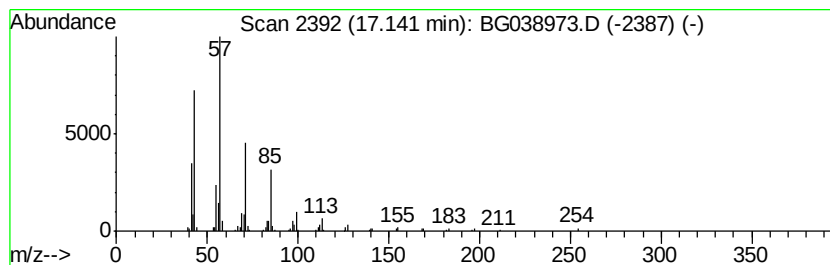
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TIC Integration Parameters: LSCINT.P

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Peak Number 6 Heneicosane Concentration Rank 10

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 17.14 | 9.94 ng | 248323 | Phenanthrene-d10 | 17.43 |

| Hit# of | 5           | Tentative ID | MW  | MolForm | CAS#        | Qual |
|---------|-------------|--------------|-----|---------|-------------|------|
| 1       | Heneicosane |              | 296 | C21H44  | 000629-94-7 | 90   |
| 2       | Heptacosane |              | 380 | C27H56  | 000593-49-7 | 90   |
| 3       | Octadecane  |              | 254 | C18H38  | 000593-45-3 | 87   |
| 4       | Hexadecane  |              | 226 | C16H34  | 000544-76-3 | 87   |
| 5       | Heptadecane |              | 240 | C17H36  | 000629-78-7 | 86   |





Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010719\  
Data File : BG038973.D  
Acq On : 7 Jan 2019 19:12  
Operator : JU/SJ  
Sample : K1036-05  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
13-E

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG121218.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

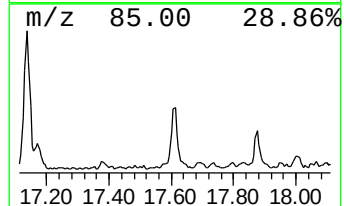
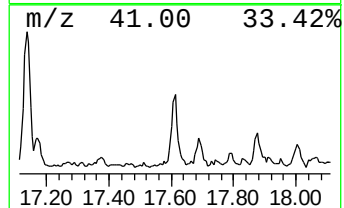
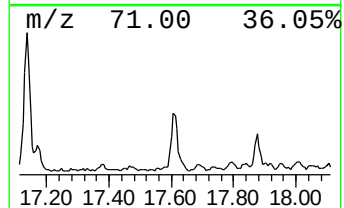
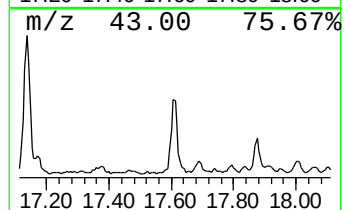
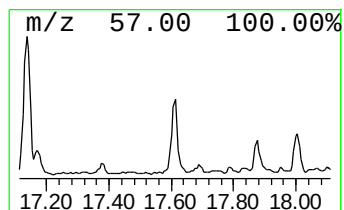
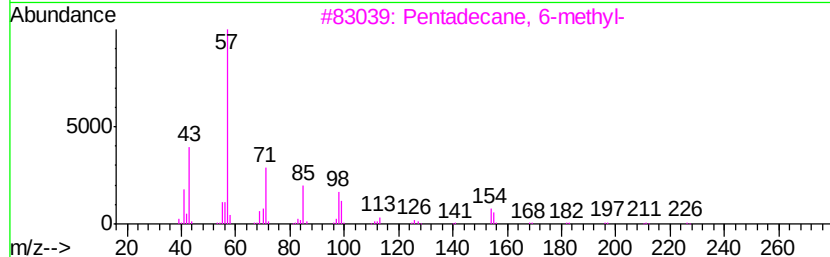
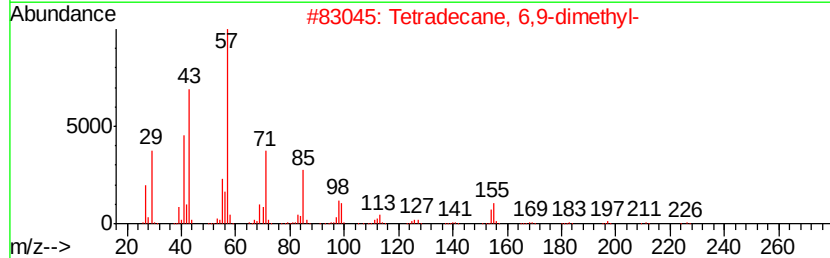
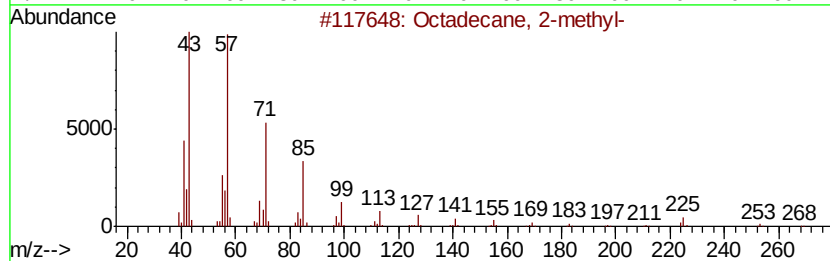
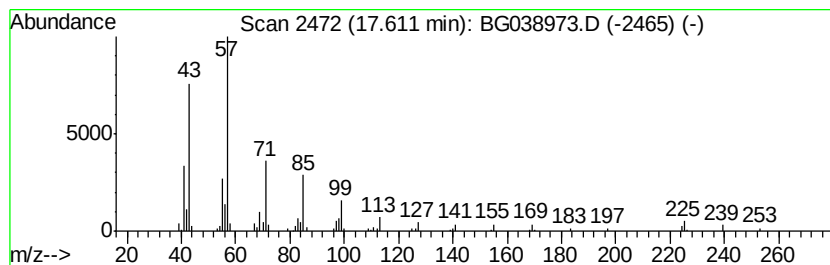
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 7 Octadecane, 2-methyl- Concentration Rank 19

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 17.61 | 5.63 ng | 140653 | Phenanthrene-d10 | 17.43 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Octadecane, 2-methyl-      | 268 | C19H40  | 001560-88-9 | 87   |
| 2    |    |   | Tetradecane, 6,9-dimethyl- | 226 | C16H34  | 055045-13-1 | 81   |
| 3    |    |   | Pentadecane, 6-methyl-     | 226 | C16H34  | 010105-38-1 | 81   |
| 4    |    |   | Hexadecane, 2-methyl-      | 240 | C17H36  | 001560-92-5 | 80   |
| 5    |    |   | Nonadecane, 2-methyl-      | 282 | C20H42  | 001560-86-7 | 72   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010719\  
Data File : BG038973.D  
Acq On : 7 Jan 2019 19:12  
Operator : JU/SJ  
Sample : K1036-05  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
13-E

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG121218.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

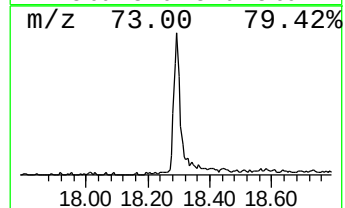
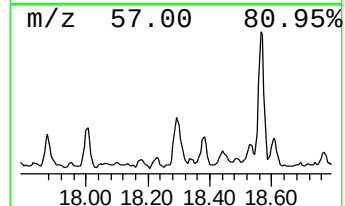
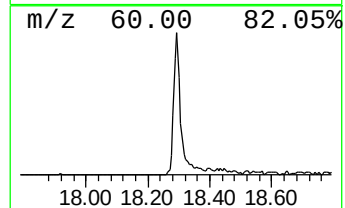
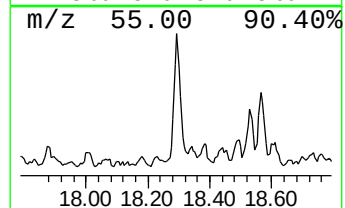
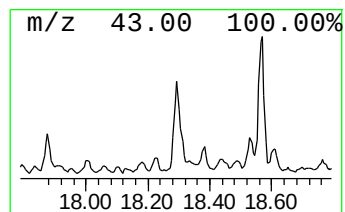
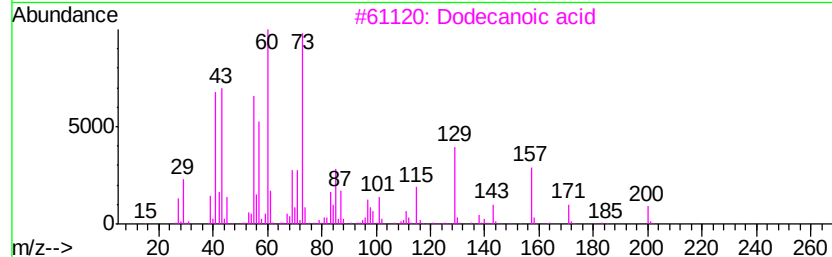
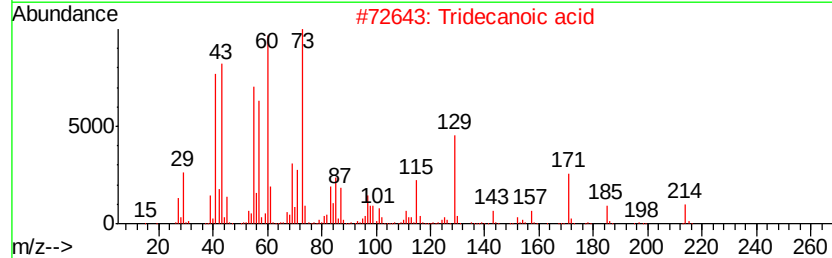
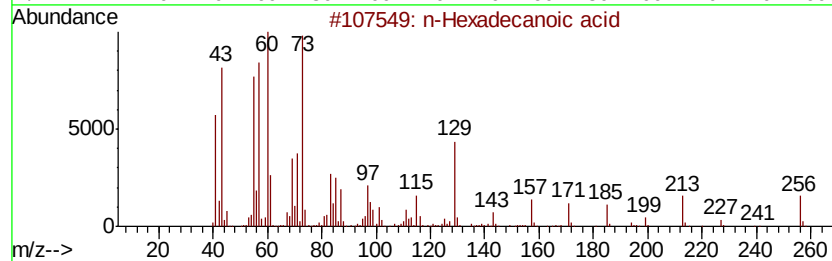
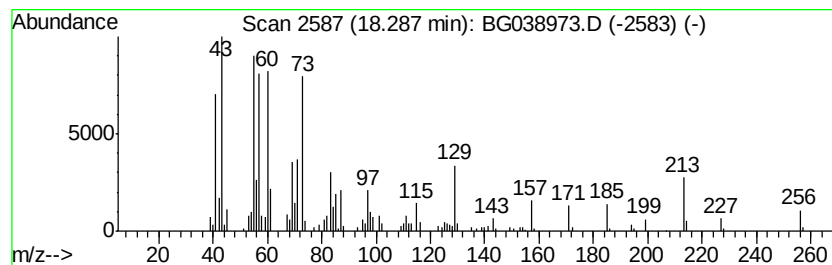
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 8 n-Hexadecanoic acid Concentration Rank 9

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 18.29 | 10.86 ng | 271245 | Phenanthrene-d10 | 17.43 |

| Hit# of | 5 | Tentative ID        | MW  | MolForm  | CAS#        | Qual |
|---------|---|---------------------|-----|----------|-------------|------|
| 1       |   | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 97   |
| 2       |   | Tridecanoic acid    | 214 | C13H26O2 | 000638-53-9 | 93   |
| 3       |   | Dodecanoic acid     | 200 | C12H24O2 | 000143-07-7 | 70   |
| 4       |   | Pentadecanoic acid  | 242 | C15H30O2 | 001002-84-2 | 59   |
| 5       |   | Undecanoic acid     | 186 | C11H22O2 | 000112-37-8 | 55   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010719\  
 Data File : BG038973.D  
 Acq On : 7 Jan 2019 19:12  
 Operator : JU/SJ  
 Sample : K1036-05  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 13-E

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG121218.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

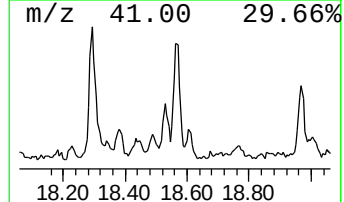
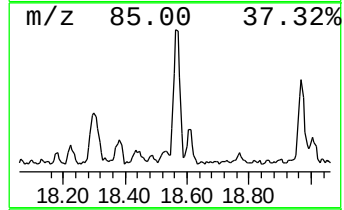
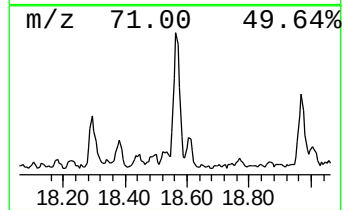
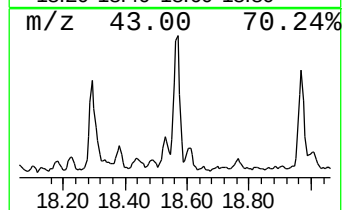
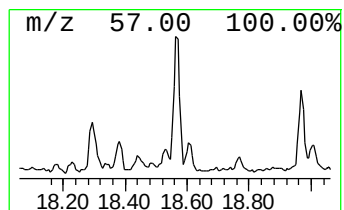
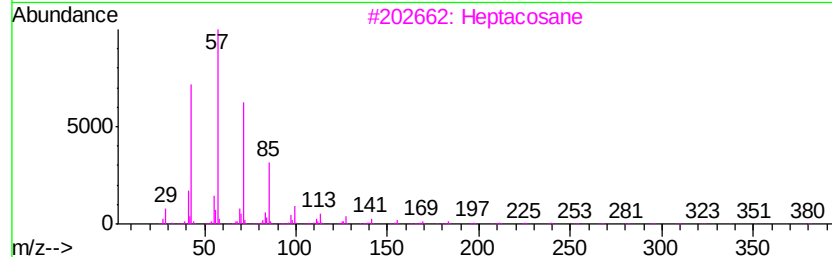
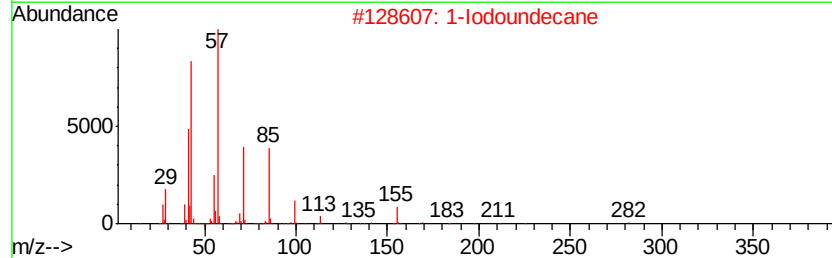
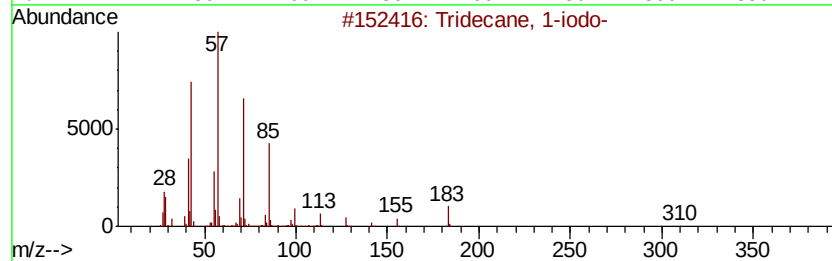
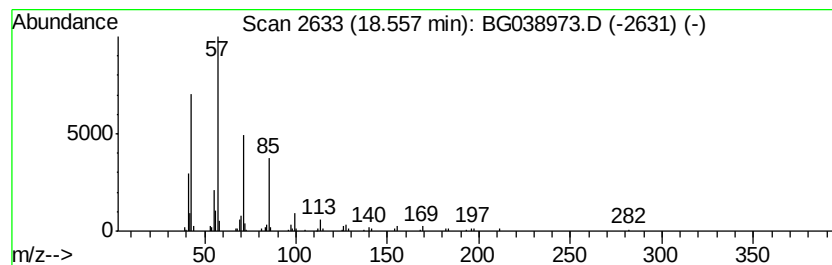
TIC Library : C:\Database\NIST11.L  
 TIC Integration Parameters: LSCINT.P

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 Peak Number 9 Tridecane, 1-iodo- Concentration Rank 12

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 18.56 | 8.80 ng | 219853 | Phenanthrene-d10 | 17.43 |

| Hit# | of | Tentative ID       | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------|-----|---------|-------------|------|
| 1    | 5  | Tridecane, 1-iodo- | 310 | C13H27I | 035599-77-0 | 90   |
| 2    |    | 1-Iodoundecane     | 282 | C11H23I | 004282-44-4 | 87   |
| 3    |    | Heptacosane        | 380 | C27H56  | 000593-49-7 | 87   |
| 4    |    | Tetracosane        | 338 | C24H50  | 000646-31-1 | 86   |
| 5    |    | Hexadecane         | 226 | C16H34  | 000544-76-3 | 86   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010719\  
Data File : BG038973.D  
Acq On : 7 Jan 2019 19:12  
Operator : JU/SJ  
Sample : K1036-05  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
13-E

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG121218.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

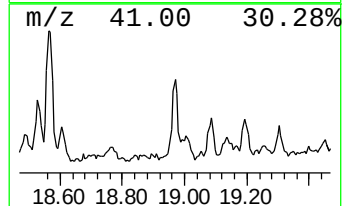
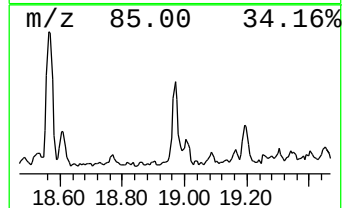
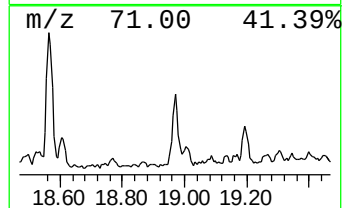
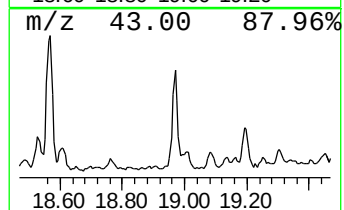
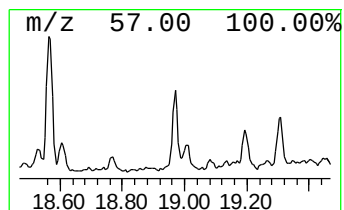
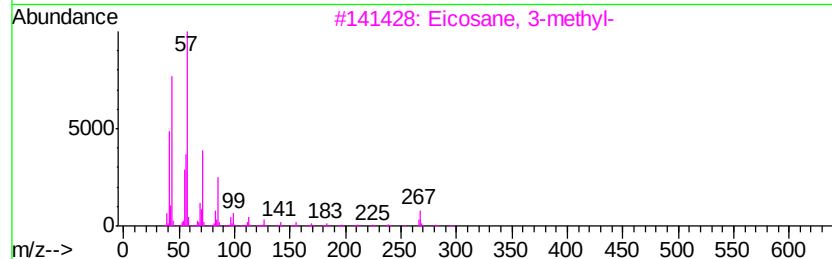
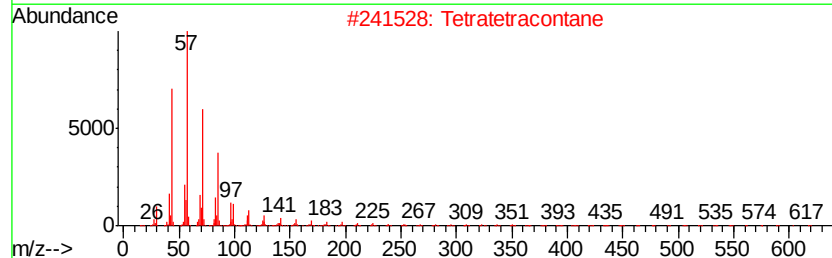
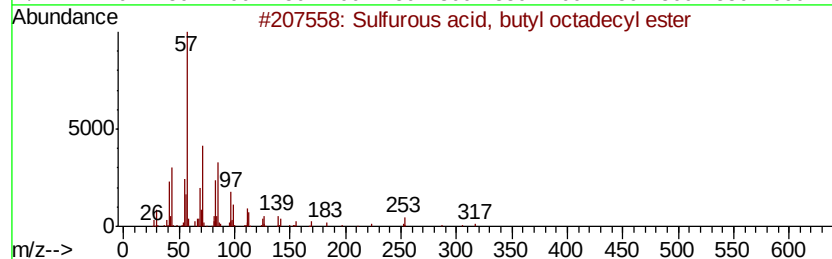
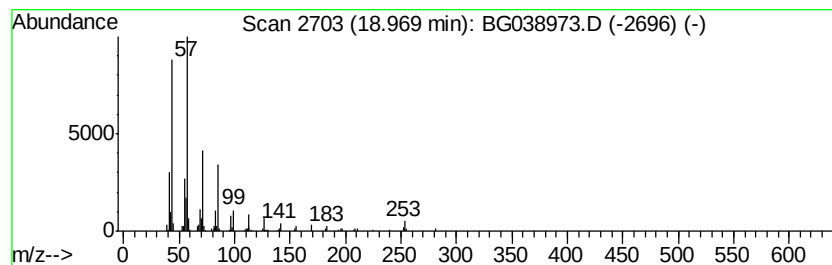
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 10 Sulfurous acid, butyl octadecyl... Concentration Rank 18

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 18.97 | 5.75 ng | 143638 | Phenanthrene-d10 | 17.43 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Sulfurous acid, butyl octadecyl ... | 390 | C22H46O3S | 1000309-18-5 | 86   |
| 2    |    |   | Tetratetracontane                   | 619 | C44H90    | 007098-22-8  | 83   |
| 3    |    |   | Eicosane, 3-methyl-                 | 296 | C21H44    | 006418-46-8  | 80   |
| 4    |    |   | Pentadecane, 2-methyl-              | 226 | C16H34    | 001560-93-6  | 80   |
| 5    |    |   | Heptadecane, 4-propyl-              | 282 | C20H42    | 055044-10-5  | 72   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010719\  
Data File : BG038973.D  
Acq On : 7 Jan 2019 19:12  
Operator : JU/SJ  
Sample : K1036-05  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
13-E

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG121218.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

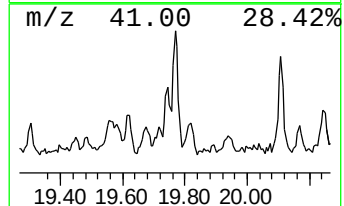
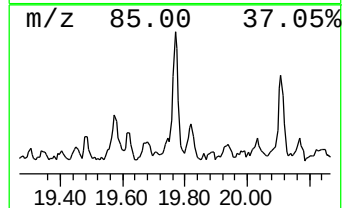
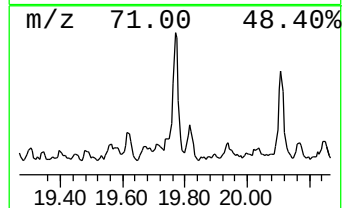
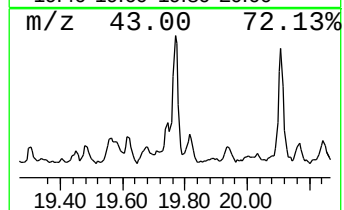
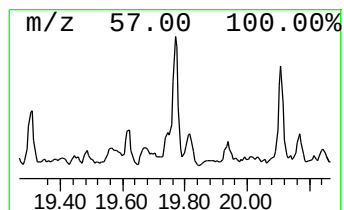
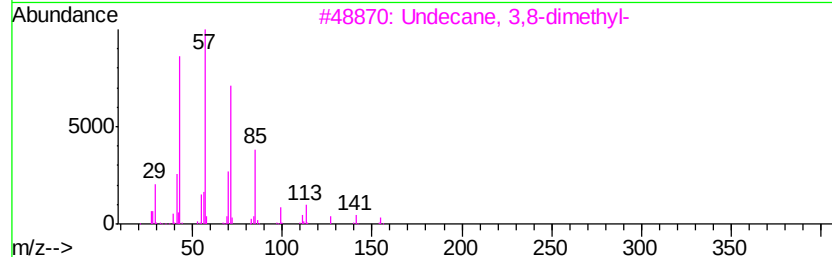
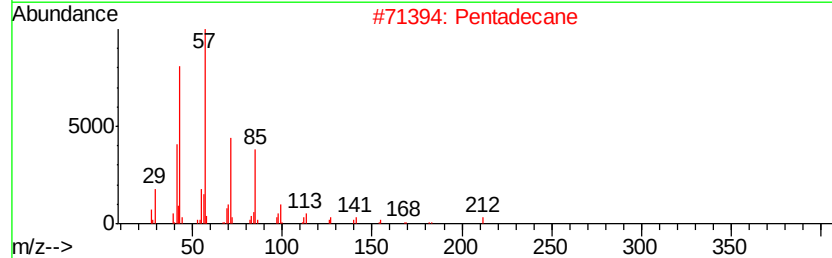
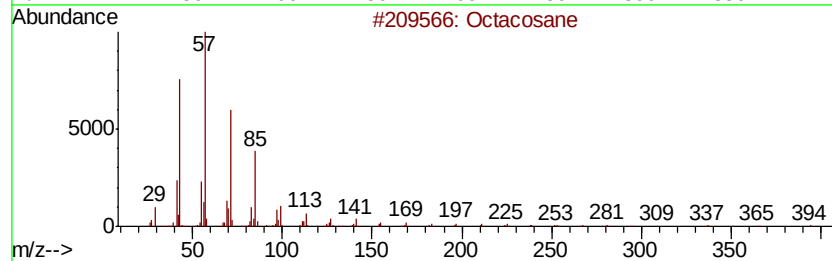
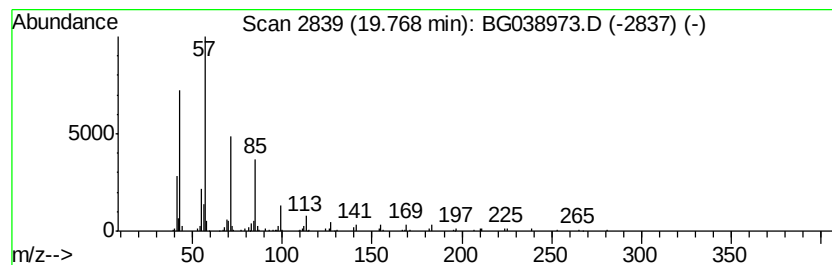
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 11 Octacosane Concentration Rank 20

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 19.77 | 5.52 ng | 170417 | Chrysene-d12     | 21.71 |

| Hit# | of | 5 | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------|-----|---------|-------------|------|
| 1    |    |   | Octacosane              | 394 | C28H58  | 000630-02-4 | 90   |
| 2    |    |   | Pentadecane             | 212 | C15H32  | 000629-62-9 | 86   |
| 3    |    |   | Undecane, 3,8-dimethyl- | 184 | C13H28  | 017301-30-3 | 86   |
| 4    |    |   | Heptadecane             | 240 | C17H36  | 000629-78-7 | 83   |
| 5    |    |   | Octadecane              | 254 | C18H38  | 000593-45-3 | 80   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010719\  
Data File : BG038973.D  
Acq On : 7 Jan 2019 19:12  
Operator : JU/SJ  
Sample : K1036-05  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
13-E

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG121218.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

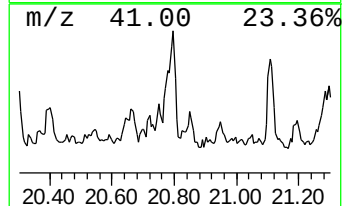
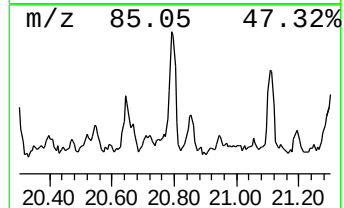
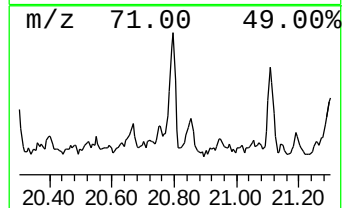
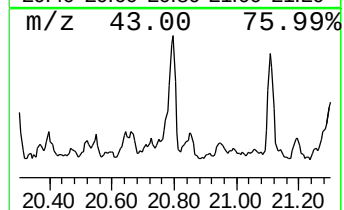
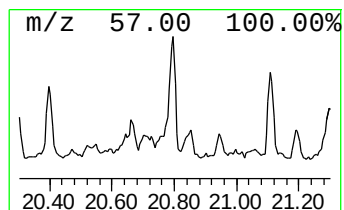
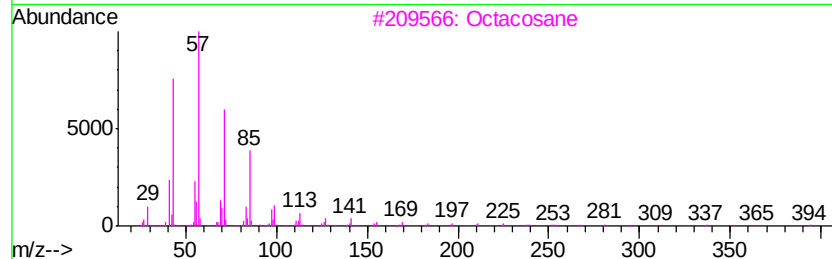
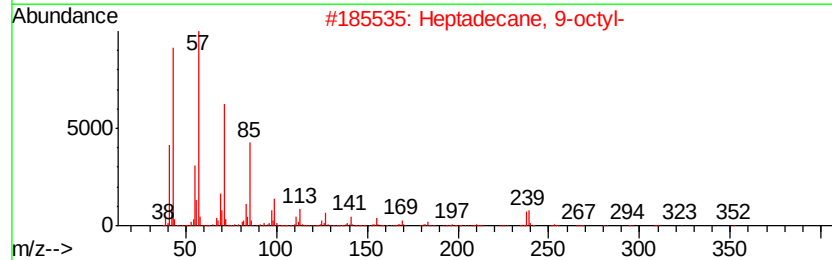
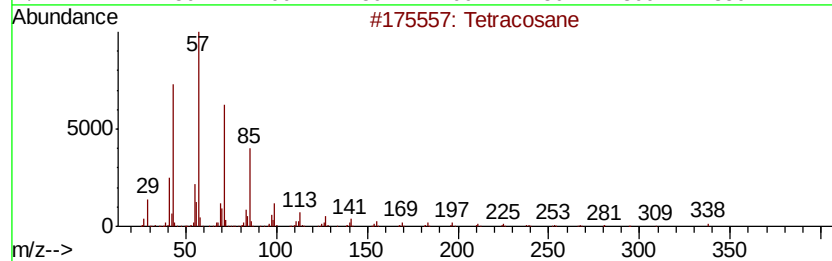
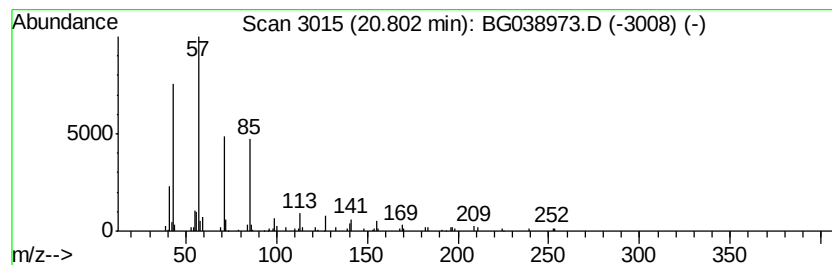
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TIC Integration Parameters: LSCINT.P

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Peak Number 12 Tetracosane Concentration Rank 15

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 20.80 | 7.24 ng | 223686 | Chrysene-d12     | 21.71 |

| Hit# | of | Tentative ID          | MW  | MolForm  | CAS#        | Qual |
|------|----|-----------------------|-----|----------|-------------|------|
| 1    | 5  | Tetracosane           | 338 | C24H50   | 000646-31-1 | 91   |
| 2    |    | Heptadecane, 9-octyl- | 352 | C25H52   | 007225-64-1 | 90   |
| 3    |    | Octacosane            | 394 | C28H58   | 000630-02-4 | 90   |
| 4    |    | 2-Bromo dodecane      | 248 | C12H25Br | 013187-99-0 | 86   |
| 5    |    | Heneicosane           | 296 | C21H44   | 000629-94-7 | 86   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010719\  
Data File : BG038973.D  
Acq On : 7 Jan 2019 19:12  
Operator : JU/SJ  
Sample : K1036-05  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
13-E

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG121218.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

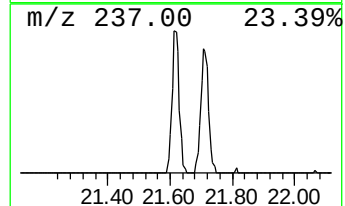
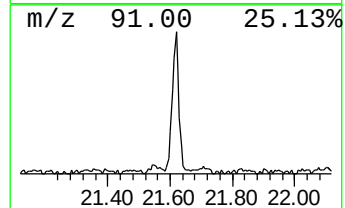
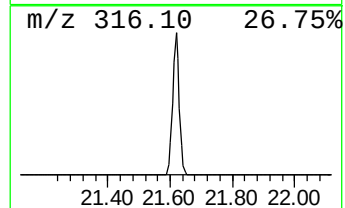
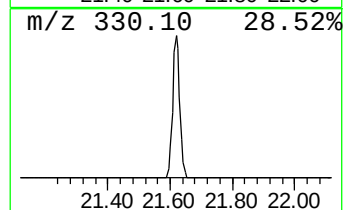
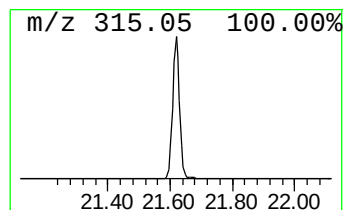
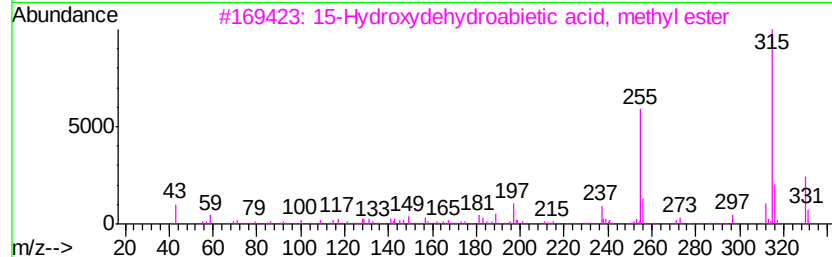
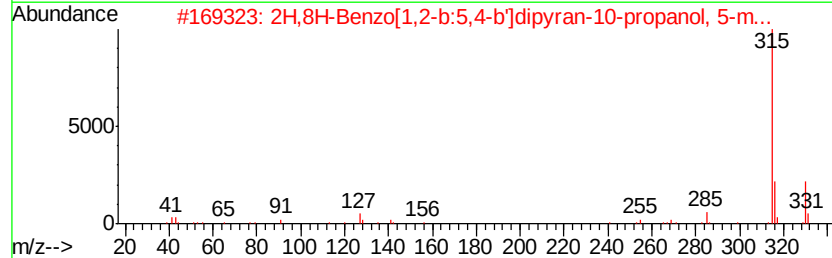
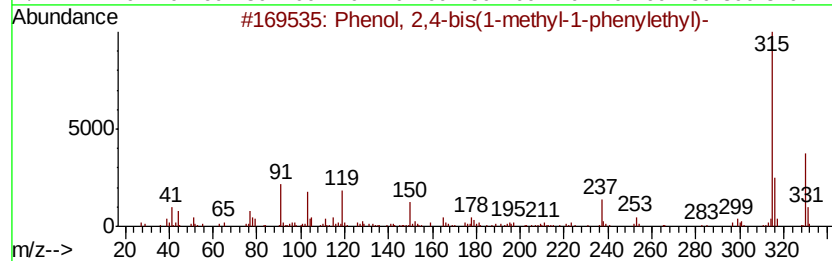
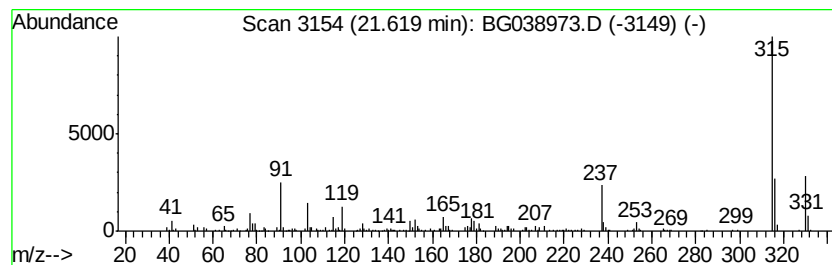
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 13 Phenol, 2,4-bis(1-methyl-1-... Concentration Rank 13

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 21.62 | 8.02 ng | 247660 | Chrysene-d12     | 21.71 |

| Hit# | of | Tentative ID                          | MW  | MolForm    | CAS#         | Qual |
|------|----|---------------------------------------|-----|------------|--------------|------|
| 1    | 5  | Phenol, 2,4-bis(1-methyl-1-phenyl-... | 330 | C24H26O    | 002772-45-4  | 87   |
| 2    |    | 2H,8H-Benzo[1,2-b:5,4-b']dipyrans...  | 330 | C20H26O4   | 026535-37-5  | 58   |
| 3    |    | 15-Hydroxydehydroabiatic acid, m...   | 330 | C21H30O3   | 029461-23-2  | 58   |
| 4    |    | 1H-Dicyclooct[fe,qlisoindole-1,3(...  | 315 | C20H29NO2  | 051624-47-6  | 50   |
| 5    |    | Acridin-9-yl-(4-nitro-phenyl)-amine   | 315 | C19H13N3O2 | 1000317-46-5 | 43   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010719\  
Data File : BG038973.D  
Acq On : 7 Jan 2019 19:12  
Operator : JU/SJ  
Sample : K1036-05  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
13-E

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG121218.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

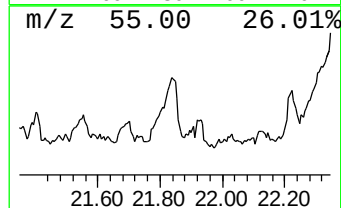
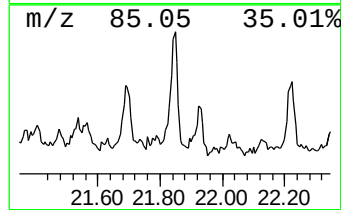
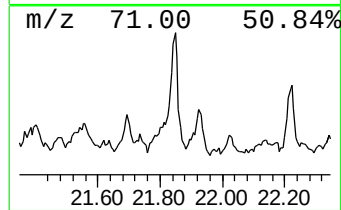
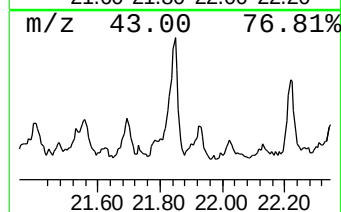
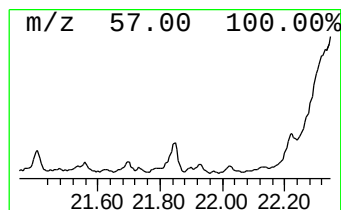
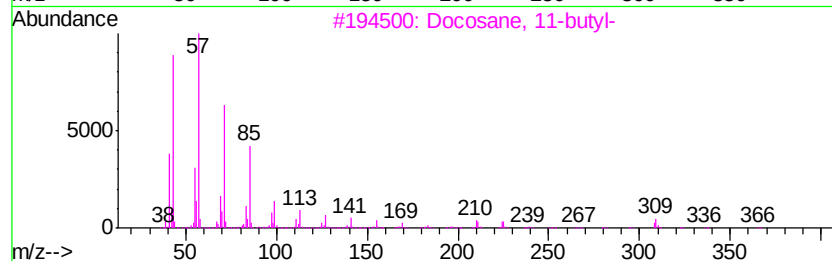
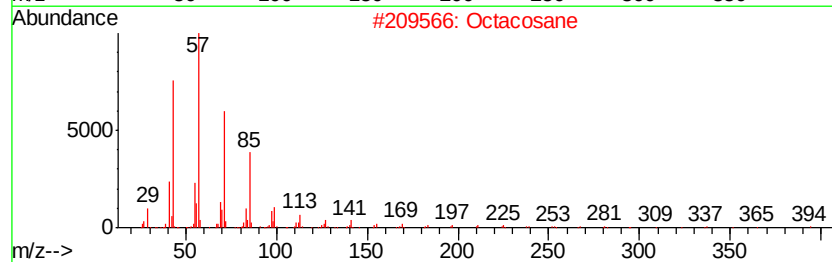
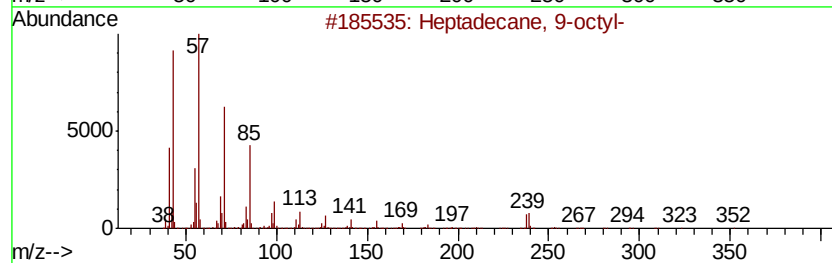
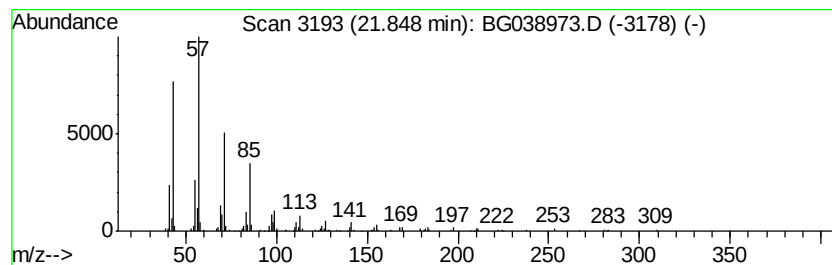
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 14 Heptadecane, 9-octyl- Concentration Rank 11

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 21.85 | 8.81 ng | 272303 | Chrysene-d12     | 21.71 |

| Hit# | of | 5 | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------|-----|---------|-------------|------|
| 1    |    |   | Heptadecane, 9-octyl- | 352 | C25H52  | 007225-64-1 | 91   |
| 2    |    |   | Octacosane            | 394 | C28H58  | 000630-02-4 | 91   |
| 3    |    |   | Docosane, 11-butyl-   | 366 | C26H54  | 013475-76-8 | 91   |
| 4    |    |   | Hentriacontane        | 437 | C31H64  | 000630-04-6 | 91   |
| 5    |    |   | Hexacosane            | 366 | C26H54  | 000630-01-3 | 91   |





Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010719\  
Data File : BG038973.D  
Acq On : 7 Jan 2019 19:12  
Operator : JU/SJ  
Sample : K1036-05  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
13-E

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG121218.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

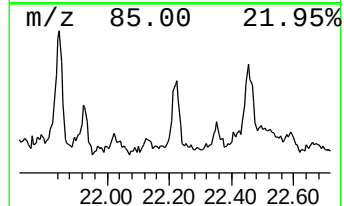
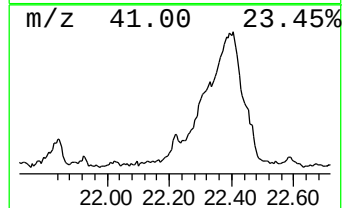
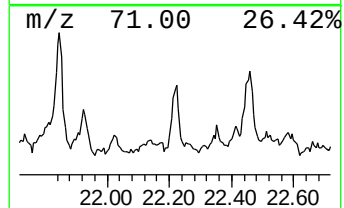
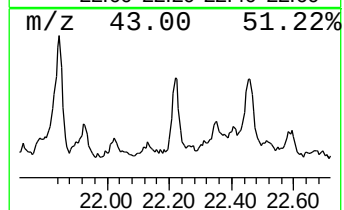
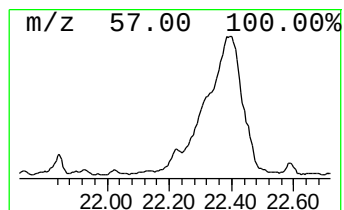
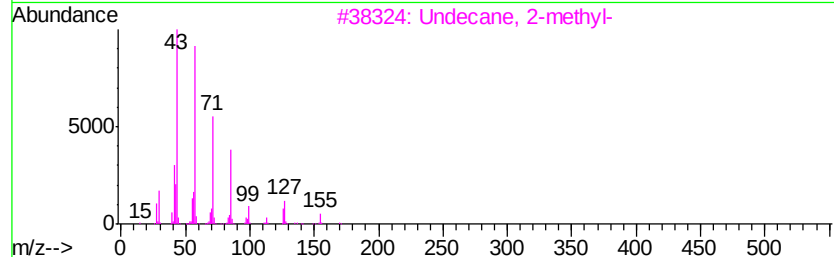
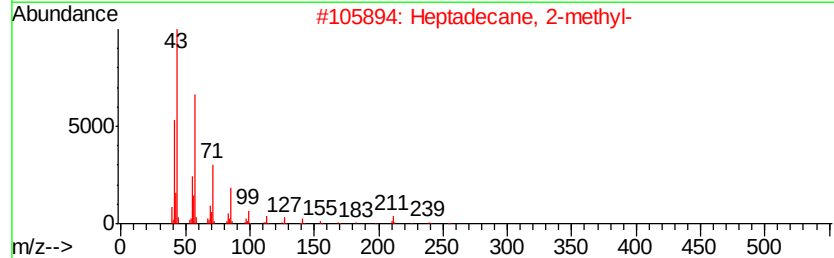
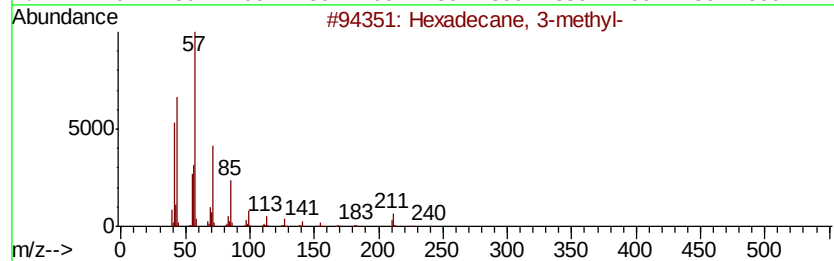
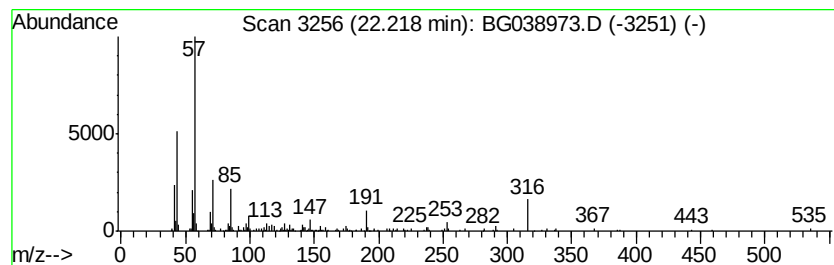
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 15 unknown22.22 Concentration Rank 16

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 22.22 | 6.82 ng | 210536 | Chrysene-d12     | 21.71 |

| Hit# | of | 5 | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------------|-----|---------|-------------|------|
| 1    |    |   | Hexadecane, 3-methyl-  | 240 | C17H36  | 006418-43-5 | 43   |
| 2    |    |   | Heptadecane, 2-methyl- | 254 | C18H38  | 001560-89-0 | 38   |
| 3    |    |   | Undecane, 2-methyl-    | 170 | C12H26  | 007045-71-8 | 38   |
| 4    |    |   | Octacosane             | 394 | C28H58  | 000630-02-4 | 35   |
| 5    |    |   | 1-Iodoundecane         | 282 | C11H23I | 004282-44-4 | 35   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010719\  
Data File : BG038973.D  
Acq On : 7 Jan 2019 19:12  
Operator : JU/SJ  
Sample : K1036-05  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

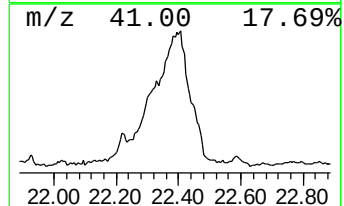
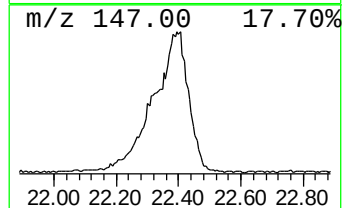
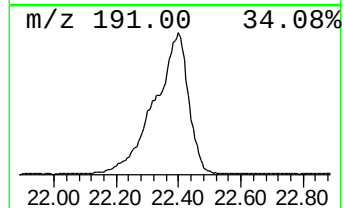
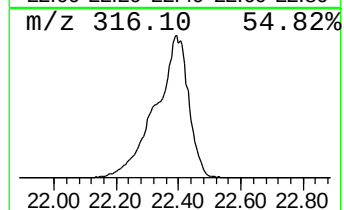
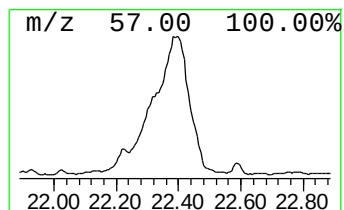
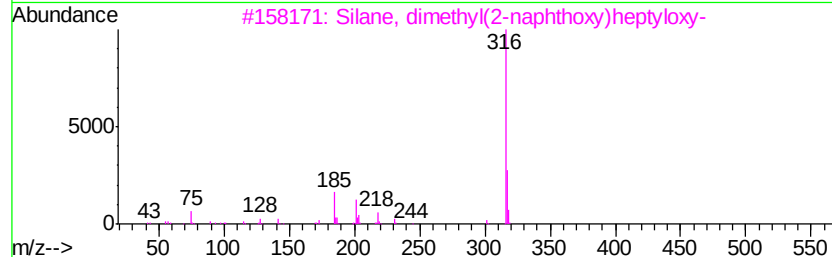
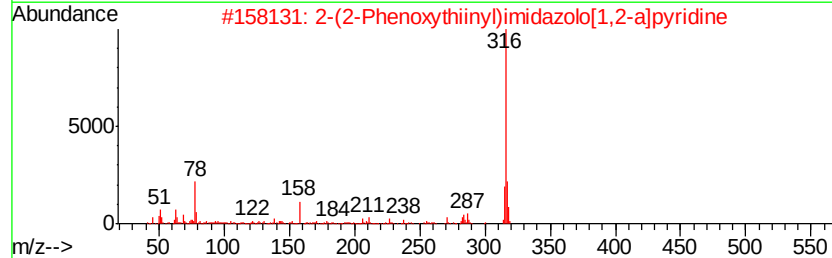
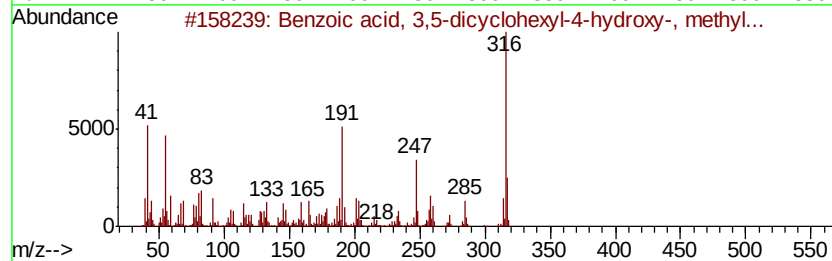
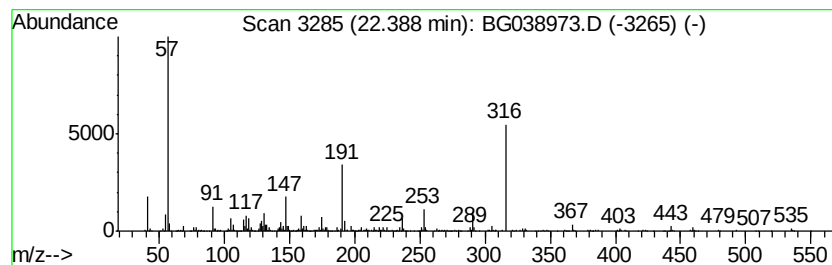
Instrument :  
BNA\_G  
ClientSampleId :  
13-E

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG121218.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 16 unknown22.39 Concentration Rank 1

| R.T.    | EstConc   | Area                                | Relative to ISTD | R.T.       |              |      |
|---------|-----------|-------------------------------------|------------------|------------|--------------|------|
| 22.39   | 180.34 ng | 5571060                             | Chrysene-d12     | 21.71      |              |      |
| Hit# of | 5         | Tentative ID                        | MW               | MolForm    | CAS#         | Qual |
| 1       |           | Benzoic acid, 3,5-dicyclohexyl-4... | 316              | C20H28O3   | 055125-23-0  | 28   |
| 2       |           | 2-(2-Phenoxythiiny)imidazolo[1,...  | 316              | C19H12N2OS | 109534-27-2  | 16   |
| 3       |           | Silane, dimethyl(2-naphthoxy)hep... | 316              | C19H28O2Si | 1000347-21-5 | 16   |
| 4       |           | Imidazo[1,2-b]1,2,4-triazine, 6-... | 316              | C19H16N4O  | 094103-49-8  | 9    |
| 5       |           | 3-(2-Ethyl-piperidin-1-ylmethyl)... | 345              | C22H35NO2  | 1000300-75-1 | 9    |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010719\  
Data File : BG038973.D  
Acq On : 7 Jan 2019 19:12  
Operator : JU/SJ  
Sample : K1036-05  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
13-E

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG121218.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

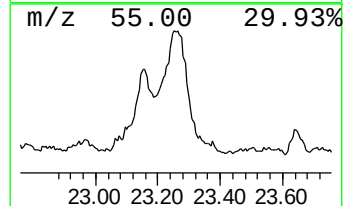
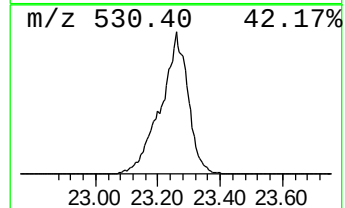
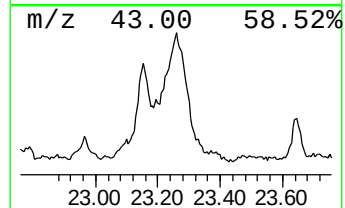
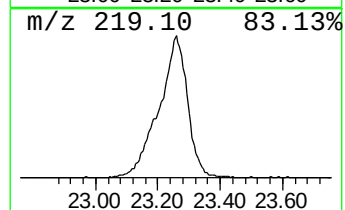
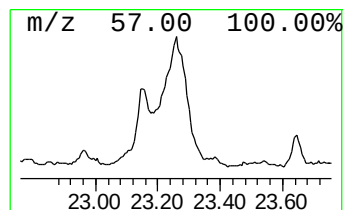
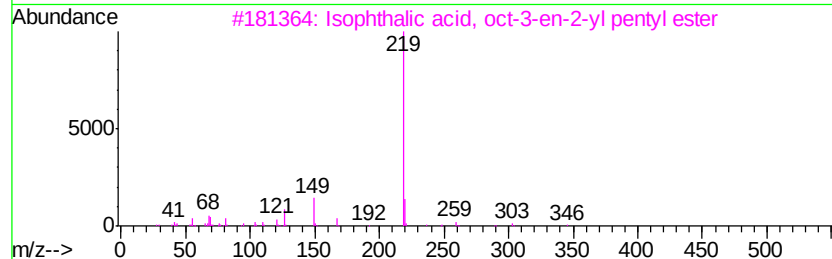
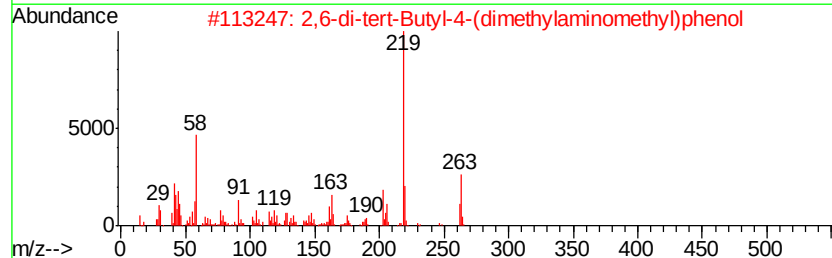
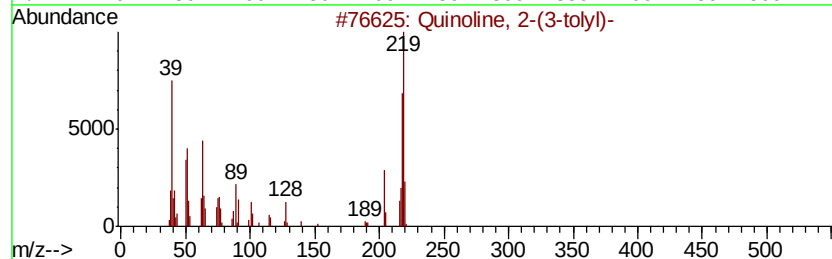
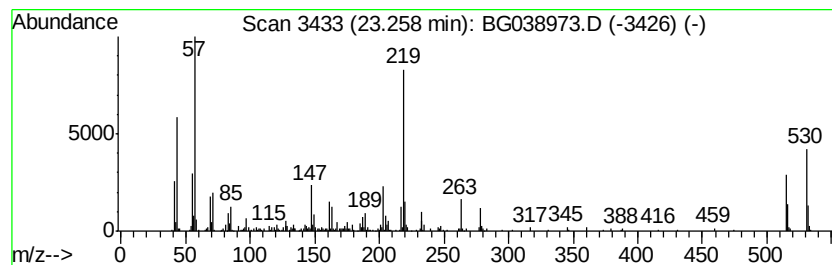
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 18 Quinoline, 2-(3-tolyl)- Concentration Rank 3

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.  |
|-------|----------|---------|------------------|-------|
| 23.26 | 54.33 ng | 1678190 | Chrysene-d12     | 21.71 |

| Hit# | of | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|-------------------------------------|-----|------------|--------------|------|
| 1    | 5  | Quinoline, 2-(3-tolyl)-             | 219 | C16H13N    | 024641-30-3  | 80   |
| 2    |    | 2,6-di-tert-Butyl-4-(dimethylami... | 263 | C17H29N0   | 000088-27-7  | 38   |
| 3    |    | Isophthalic acid, oct-3-en-2-yl ... | 346 | C21H30O4   | 1000343-89-3 | 38   |
| 4    |    | N-(2,6-Dimethyl-phenyl)-2-(4-nit... | 367 | C21H25N3O3 | 1000294-55-6 | 27   |
| 5    |    | Propanoic acid, 2-(1,3-dioxo-2,3... | 249 | C12H11N05  | 1000196-70-9 | 27   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010719\  
Data File : BG038973.D  
Acq On : 7 Jan 2019 19:12  
Operator : JU/SJ  
Sample : K1036-05  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
13-E

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG121218.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

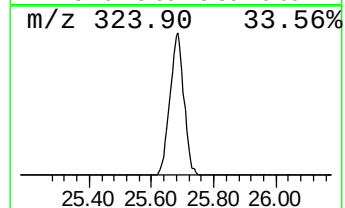
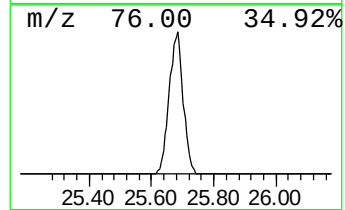
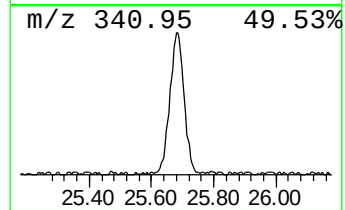
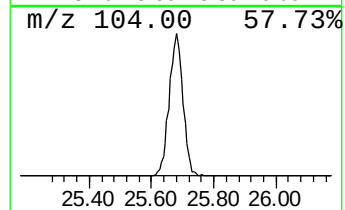
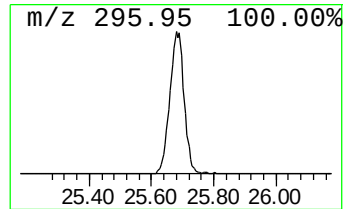
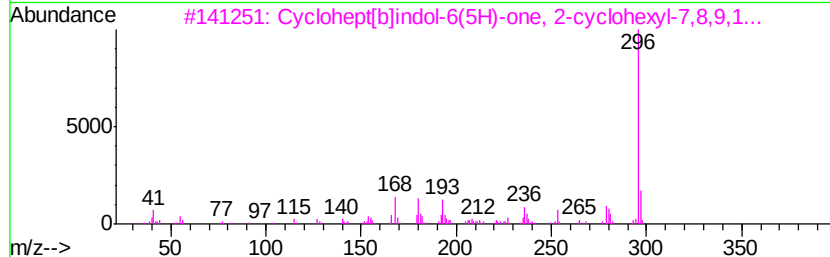
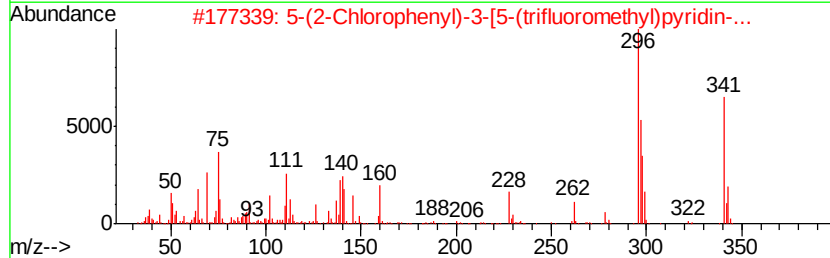
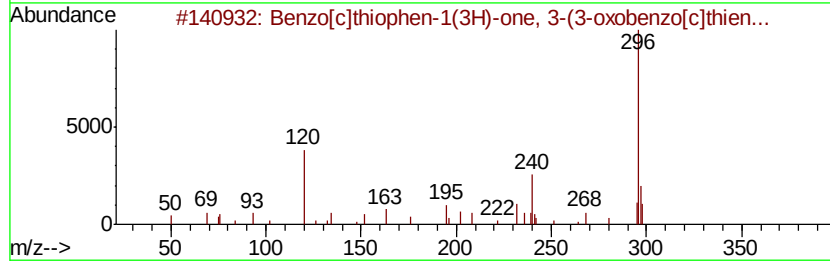
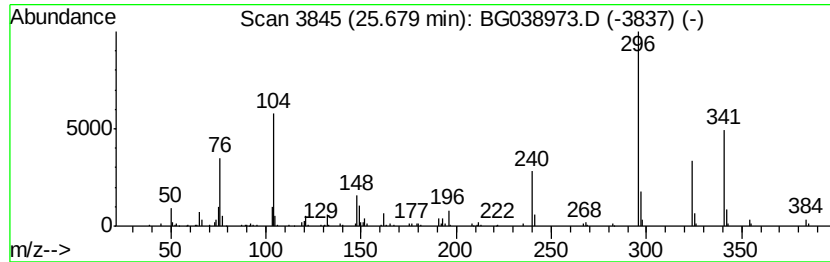
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 19 unknown25.68 Concentration Rank 6

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 25.68 | 18.66 ng | 433773 | Perylene-d12     | 25.00 |

| Hit# | of | Tentative ID                        | MW  | MolForm       | CAS#         | Qual |
|------|----|-------------------------------------|-----|---------------|--------------|------|
| 1    | 5  | Benzo[c]thiophen-1(3H)-one, 3-(3... | 296 | C16H8O2S2     | 023667-32-5  | 35   |
| 2    |    | 5-(2-Chlorophenyl)-3-[5-(trifluo... | 341 | C14H7ClF3N3O2 | 287979-09-3  | 23   |
| 3    |    | Cyclohept[b]indol-6(5H)-one, 2-c... | 296 | C19H24N2O     | 339284-86-5  | 22   |
| 4    |    | 4-(2-Benzofuranyl)-2-chloro-6-(2... | 296 | C16H9ClN2O2   | 131022-75-8  | 22   |
| 5    |    | N-butyl-2,2,3,3,4,4,4-heptafluor... | 387 | C17H20F7NO    | 1000365-85-0 | 22   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010719\  
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Operator : JU/SJ  
Sample : K1036-05  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
13-E

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG121218.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

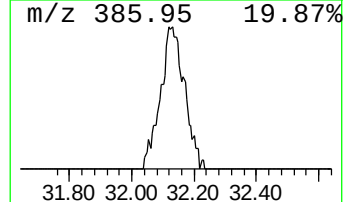
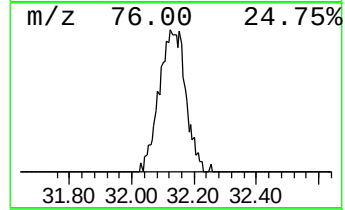
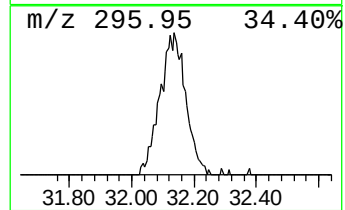
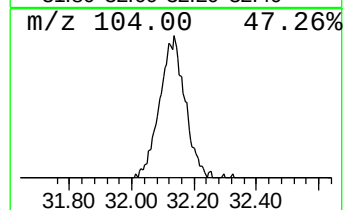
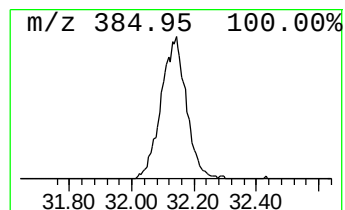
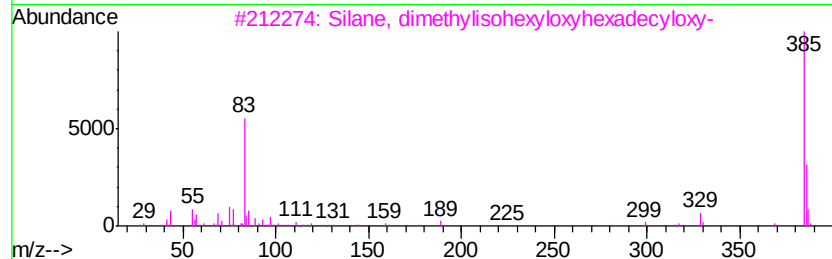
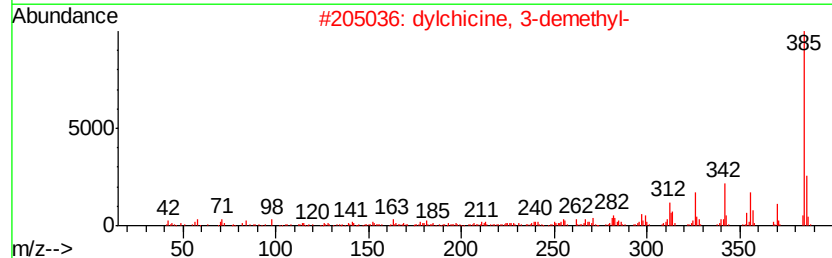
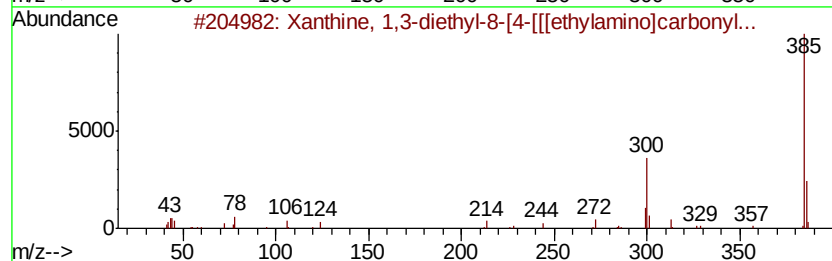
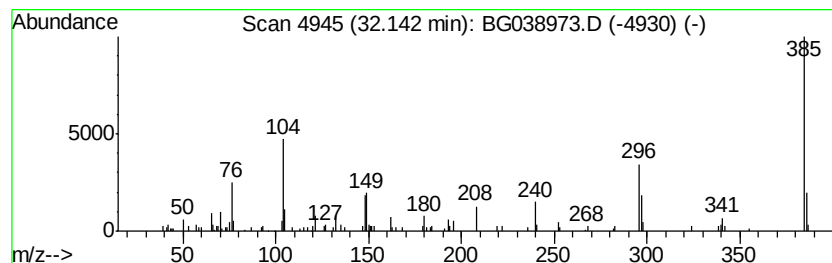
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 20 unknown32.14 Concentration Rank 7

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 32.14 | 12.52 ng | 290995 | Perylene-d12     | 25.00 |

| Hit# | of | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|-------------------------------------|-----|------------|--------------|------|
| 1    | 5  | Xanthine, 1,3-diethyl-8-[4-[[[et... | 385 | C19H23N5O4 | 104576-48-9  | 22   |
| 2    |    | dylchicine, 3-demethyl-             | 385 | C21H23N06  | 007336-34-7  | 16   |
| 3    |    | Silane, dimethylisohexvloxvhexad... | 400 | C24H52O2Si | 1000346-94-5 | 10   |
| 4    |    | Silane, dimethylisobutoxyoctadec... | 400 | C24H52O2Si | 1000346-77-7 | 10   |
| 5    |    | 2,3-Dibenzyl-5,8-dimethoxy-6-am...  | 385 | C24H23N3O2 | 056393-28-3  | 10   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_G\DATA\BG010719\  
 Data File : BG038973.D  
 Acq On : 7 Jan 2019 19:12  
 Operator : JU/SJ  
 Sample : K1036-05  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 13-E

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\8270-BG121218.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 |
| unknown7.58          | 7.58  | 65.0    | ng    | 679398   | 1                     | 8.05  | 208999 | 20.0 |
| Dodecane             | 10.78 | 6.7     | ng    | 101199   | 2                     | 10.86 | 301057 | 20.0 |
| Phenol, 2,6-bis(1... | 15.15 | 19.9    | ng    | 574726   | 3                     | 14.69 | 577389 | 20.0 |
| Hexadecane           | 15.48 | 7.4     | ng    | 214918   | 3                     | 14.69 | 577389 | 20.0 |
| Heneicosane          | 17.14 | 9.9     | ng    | 248323   | 4                     | 17.43 | 499710 | 20.0 |
| Octadecane, 2-met... | 17.61 | 5.6     | ng    | 140653   | 4                     | 17.43 | 499710 | 20.0 |
| n-Hexadecanoic acid  | 18.29 | 10.9    | ng    | 271245   | 4                     | 17.43 | 499710 | 20.0 |
| Tridecane, 1-iodo-   | 18.56 | 8.8     | ng    | 219853   | 4                     | 17.43 | 499710 | 20.0 |
| Sulfurous acid, b... | 18.97 | 5.8     | ng    | 143638   | 4                     | 17.43 | 499710 | 20.0 |
| Octacosane           | 19.77 | 5.5     | ng    | 170417   | 5                     | 21.71 | 617825 | 20.0 |
| Tetracosane          | 20.80 | 7.2     | ng    | 223686   | 5                     | 21.71 | 617825 | 20.0 |
| Phenol, 2,4-bis(1... | 21.62 | 8.0     | ng    | 247660   | 5                     | 21.71 | 617825 | 20.0 |
| Heptadecane, 9-oc... | 21.85 | 8.8     | ng    | 272303   | 5                     | 21.71 | 617825 | 20.0 |
| unknown22.22         | 22.22 | 6.8     | ng    | 210536   | 5                     | 21.71 | 617825 | 20.0 |
| unknown22.39         | 22.39 | 180.3   | ng    | 5571060  | 5                     | 21.71 | 617825 | 20.0 |
| Quinoline, 2-(3-t... | 23.26 | 54.3    | ng    | 1678190  | 5                     | 21.71 | 617825 | 20.0 |
| unknown25.68         | 25.68 | 18.7    | ng    | 433773   | 6                     | 25.00 | 464949 | 20.0 |
| unknown32.14         | 32.14 | 12.5    | ng    | 290995   | 6                     | 25.00 | 464949 | 20.0 |