

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM122718\  
 Data File : BM018402.D  
 Acq On : 27 Dec 2018 18:41  
 Operator : JU/SJ  
 Sample : J6569-03  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 381128115-4

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\_M\METHODS\8270-BM122718.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         | 5.357       | 415           | 420         | 427          | rVB      | 275883         | 417312        | 3.35%           | 0.439%        |
| 2         | 6.128       | 546           | 551         | 553          | rBV2     | 155594         | 251129        | 2.02%           | 0.264%        |
| 3         | 6.622       | 631           | 635         | 641          | rVB2     | 166908         | 283103        | 2.27%           | 0.298%        |
| 4         | 6.934       | 683           | 688         | 697          | rVB2     | 235647         | 374597        | 3.01%           | 0.394%        |
| 5         | 7.781       | 826           | 832         | 838          | rBV      | 1154180        | 1846899       | 14.83%          | 1.945%        |
| 6         | 8.175       | 894           | 899         | 903          | rBV3     | 101236         | 191329        | 1.54%           | 0.201%        |
| 7         | 8.881       | 1014          | 1019        | 1024         | rBV      | 231746         | 353695        | 2.84%           | 0.372%        |
| 8         | 8.945       | 1024          | 1030        | 1036         | rVV      | 802117         | 1330527       | 10.69%          | 1.401%        |
| 9         | 10.575      | 1300          | 1307        | 1317         | rBV      | 1757635        | 2903309       | 23.32%          | 3.057%        |
| 10        | 11.104      | 1392          | 1397        | 1402         | rBV      | 82828          | 125151        | 1.01%           | 0.132%        |
| 11        | 11.163      | 1402          | 1407        | 1415         | rVV      | 90476          | 137289        | 1.10%           | 0.145%        |
| 12        | 11.251      | 1415          | 1422        | 1430         | rVV2     | 172270         | 374901        | 3.01%           | 0.395%        |
| 13        | 11.974      | 1540          | 1545        | 1556         | rVB9     | 55095          | 133008        | 1.07%           | 0.140%        |
| 14        | 13.045      | 1719          | 1727        | 1736         | rVB      | 2818033        | 4026965       | 32.34%          | 4.241%        |
| 15        | 13.316      | 1766          | 1773        | 1780         | rBV      | 690501         | 1065703       | 8.56%           | 1.122%        |
| 16        | 14.068      | 1895          | 1901        | 1913         | rBV      | 1585527        | 2112305       | 16.97%          | 2.224%        |
| 17        | 14.386      | 1949          | 1955        | 1957         | rBV      | 719367         | 1075233       | 8.64%           | 1.132%        |
| 18        | 14.427      | 1957          | 1962        | 1971         | rVB2     | 2777307        | 4303618       | 34.57%          | 4.532%        |
| 19        | 15.427      | 2127          | 2132        | 2143         | rBV      | 221506         | 349982        | 2.81%           | 0.369%        |
| 20        | 15.545      | 2147          | 2152        | 2157         | rVB      | 312541         | 422769        | 3.40%           | 0.445%        |
| 21        | 15.915      | 2209          | 2215        | 2220         | rBV      | 1896209        | 2537062       | 20.38%          | 2.672%        |
| 22        | 15.968      | 2220          | 2224        | 2232         | rVV      | 1628396        | 2042741       | 16.41%          | 2.151%        |
| 23        | 16.051      | 2233          | 2238        | 2249         | rVB9     | 81585          | 193732        | 1.56%           | 0.204%        |
| 24        | 16.198      | 2260          | 2263        | 2268         | rVB      | 114071         | 151874        | 1.22%           | 0.160%        |
| 25        | 16.257      | 2268          | 2273        | 2278         | rBV      | 121764         | 160051        | 1.29%           | 0.169%        |
| 26        | 16.315      | 2278          | 2283        | 2289         | rBV2     | 86379          | 180666        | 1.45%           | 0.190%        |
| 27        | 16.374      | 2289          | 2293        | 2297         | rBV3     | 110565         | 159108        | 1.28%           | 0.168%        |
| 28        | 16.415      | 2297          | 2300        | 2306         | rVB      | 104111         | 131112        | 1.05%           | 0.138%        |
| 29        | 16.521      | 2315          | 2318        | 2322         | rVB      | 116430         | 137317        | 1.10%           | 0.145%        |
| 30        | 16.574      | 2322          | 2327        | 2334         | rVB4     | 248835         | 457026        | 3.67%           | 0.481%        |
| 31        | 16.645      | 2334          | 2339        | 2342         | rBV      | 169887         | 255161        | 2.05%           | 0.269%        |
| 32        | 16.715      | 2347          | 2351        | 2356         | rVB      | 144472         | 189880        | 1.53%           | 0.200%        |
| 33        | 16.933      | 2383          | 2388        | 2400         | rVB      | 631230         | 879420        | 7.06%           | 0.926%        |
| 34        | 17.045      | 2400          | 2407        | 2412         | rBV      | 148996         | 243004        | 1.95%           | 0.256%        |

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 381128115-4

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\_M\METHODS\8270-BM122718.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |       |          |          |         |         |
|----|--------|------|------|------|-------|----------|----------|---------|---------|
| 35 | 17.174 | 2423 | 2429 | 2439 | rVB   | 3673219  | 5054768  | 40.60%  | 5.323%  |
| 36 | 17.480 | 2474 | 2481 | 2487 | rBV   | 1029888  | 1502856  | 12.07%  | 1.583%  |
| 37 | 17.556 | 2487 | 2494 | 2502 | rVV2  | 344278   | 538820   | 4.33%   | 0.567%  |
| 38 | 17.733 | 2517 | 2524 | 2532 | rBV2  | 129394   | 251813   | 2.02%   | 0.265%  |
| 39 | 18.068 | 2575 | 2581 | 2587 | rBV3  | 307361   | 542855   | 4.36%   | 0.572%  |
| 40 | 18.409 | 2634 | 2639 | 2641 | rBV2  | 96392    | 136862   | 1.10%   | 0.144%  |
| 41 | 18.533 | 2656 | 2660 | 2666 | rVB2  | 97810    | 155120   | 1.25%   | 0.163%  |
| 42 | 18.933 | 2723 | 2728 | 2736 | rBV   | 245228   | 376155   | 3.02%   | 0.396%  |
| 43 | 19.162 | 2763 | 2767 | 2771 | rBV   | 97300    | 159584   | 1.28%   | 0.168%  |
| 44 | 19.356 | 2795 | 2800 | 2808 | rBV   | 449295   | 673503   | 5.41%   | 0.709%  |
| 45 | 19.456 | 2813 | 2817 | 2819 | rBV4  | 109739   | 137992   | 1.11%   | 0.145%  |
| 46 | 19.497 | 2820 | 2824 | 2836 | rVV   | 5776525  | 6764673  | 54.33%  | 7.124%  |
| 47 | 19.809 | 2871 | 2877 | 2882 | rBV   | 10291902 | 12450615 | 100.00% | 13.111% |
| 48 | 19.939 | 2895 | 2899 | 2904 | rBV3  | 108460   | 221045   | 1.78%   | 0.233%  |
| 49 | 20.133 | 2922 | 2932 | 2935 | rBV5  | 223246   | 595297   | 4.78%   | 0.627%  |
| 50 | 20.209 | 2941 | 2945 | 2949 | rBV   | 181080   | 221614   | 1.78%   | 0.233%  |
| 51 | 20.274 | 2953 | 2956 | 2961 | rVB   | 144493   | 191890   | 1.54%   | 0.202%  |
| 52 | 20.609 | 3008 | 3013 | 3020 | rBV   | 1031286  | 1389817  | 11.16%  | 1.464%  |
| 53 | 21.086 | 3091 | 3094 | 3100 | rBV3  | 184150   | 231123   | 1.86%   | 0.243%  |
| 54 | 21.174 | 3106 | 3109 | 3113 | rBV   | 97601    | 125593   | 1.01%   | 0.132%  |
| 55 | 21.286 | 3125 | 3128 | 3131 | rBV   | 231909   | 221595   | 1.78%   | 0.233%  |
| 56 | 21.368 | 3137 | 3142 | 3147 | rBV   | 5526374  | 6593597  | 52.96%  | 6.943%  |
| 57 | 21.968 | 3240 | 3244 | 3250 | rBV3  | 250084   | 354153   | 2.84%   | 0.373%  |
| 58 | 22.233 | 3284 | 3289 | 3297 | rBV   | 1965036  | 2525030  | 20.28%  | 2.659%  |
| 59 | 22.444 | 3321 | 3325 | 3330 | rVB   | 270816   | 363542   | 2.92%   | 0.383%  |
| 60 | 22.574 | 3342 | 3347 | 3353 | rBV   | 1050485  | 1402016  | 11.26%  | 1.476%  |
| 61 | 22.815 | 3384 | 3388 | 3393 | rBV6  | 208103   | 352975   | 2.84%   | 0.372%  |
| 62 | 22.968 | 3409 | 3414 | 3424 | rBV   | 431881   | 691271   | 5.55%   | 0.728%  |
| 63 | 23.421 | 3487 | 3491 | 3497 | rVB8  | 191305   | 327211   | 2.63%   | 0.345%  |
| 64 | 23.556 | 3510 | 3514 | 3521 | rVB   | 415150   | 659765   | 5.30%   | 0.695%  |
| 65 | 23.662 | 3524 | 3532 | 3545 | rVB   | 3559609  | 6975743  | 56.03%  | 7.346%  |
| 66 | 24.232 | 3624 | 3629 | 3635 | rVB   | 252218   | 474526   | 3.81%   | 0.500%  |
| 67 | 24.709 | 3702 | 3710 | 3725 | rVB2  | 1611653  | 3854030  | 30.95%  | 4.059%  |
| 68 | 24.985 | 3749 | 3757 | 3773 | rBV3  | 2616892  | 7536690  | 60.53%  | 7.937%  |
| 69 | 26.456 | 4001 | 4007 | 4017 | rVB10 | 206593   | 572021   | 4.59%   | 0.602%  |
| 70 | 26.821 | 4064 | 4069 | 4077 | rVB7  | 212337   | 538151   | 4.32%   | 0.567%  |

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Sample : J6569-03  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
381128115-4

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\_M\METHODS\8270-BM122718.M  
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

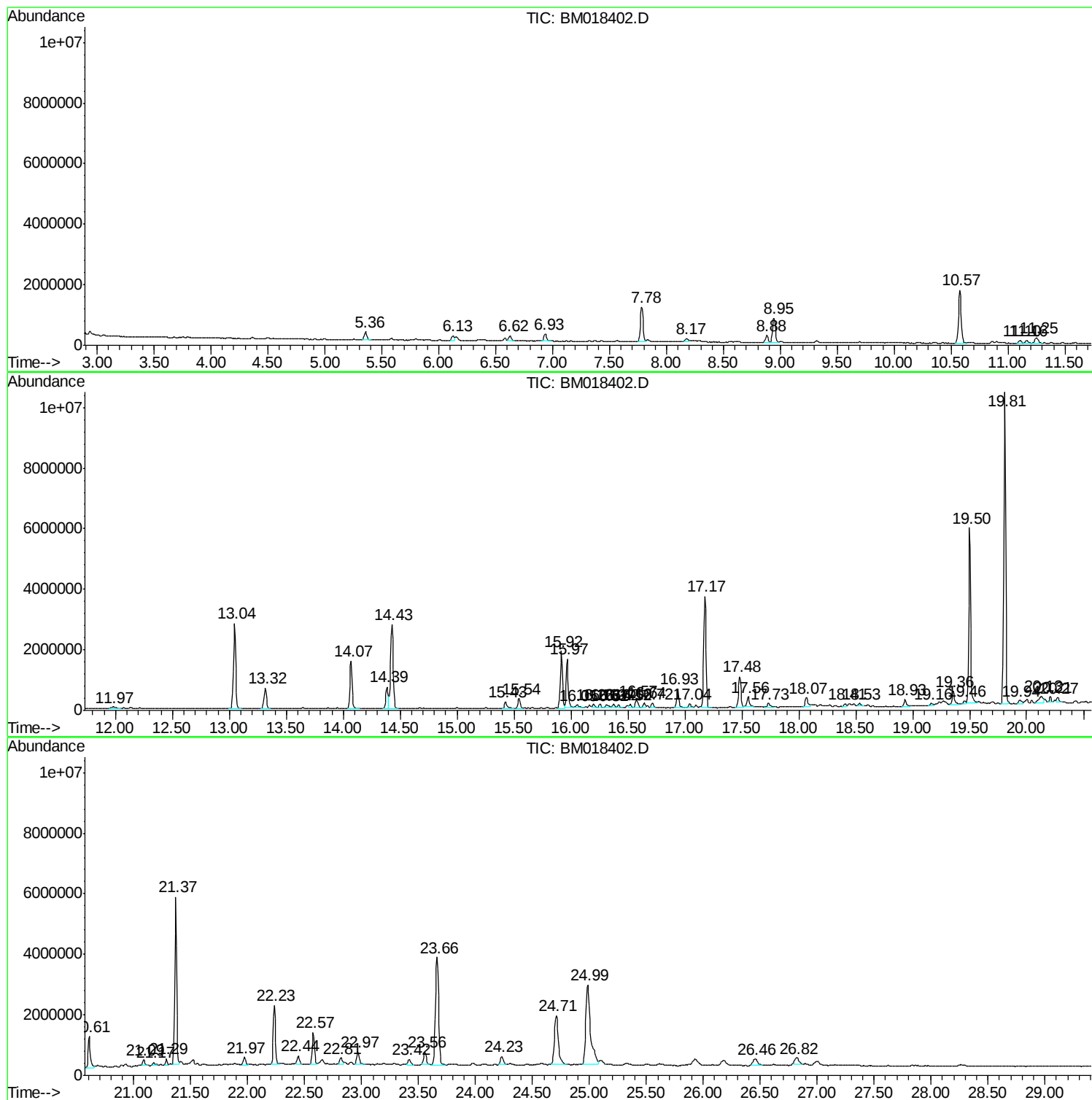
Sum of corrected areas: 94961289

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM122718\  
Data File : BM018402.D  
Acq On : 27 Dec 2018 18:41  
Operator : JU/SJ  
Sample : J6569-03  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_M  
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381128115-4

Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM122718.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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Sample : J6569-03  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
381128115-4

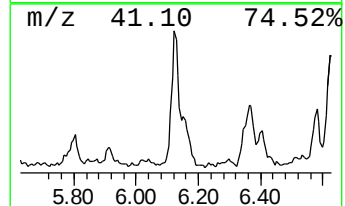
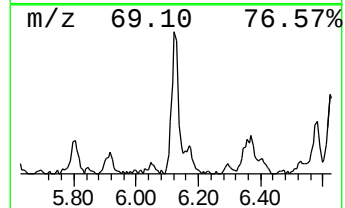
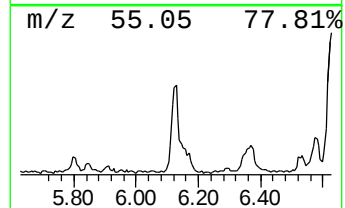
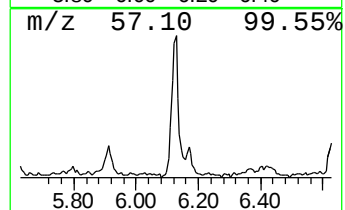
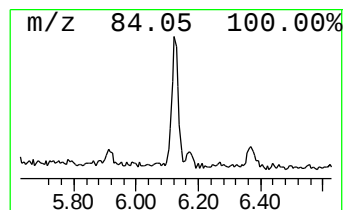
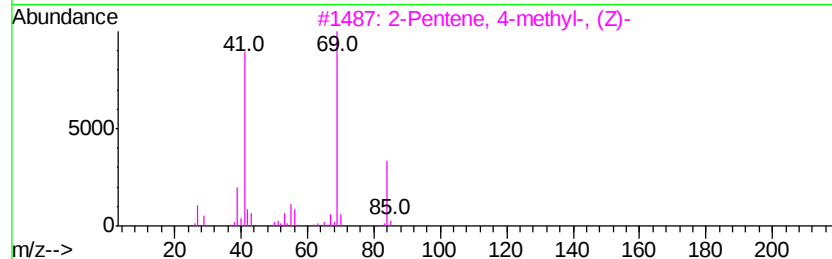
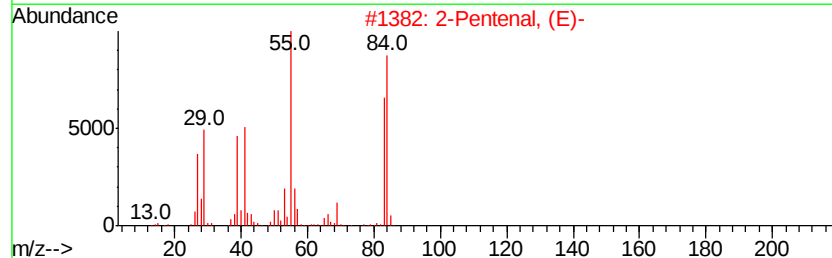
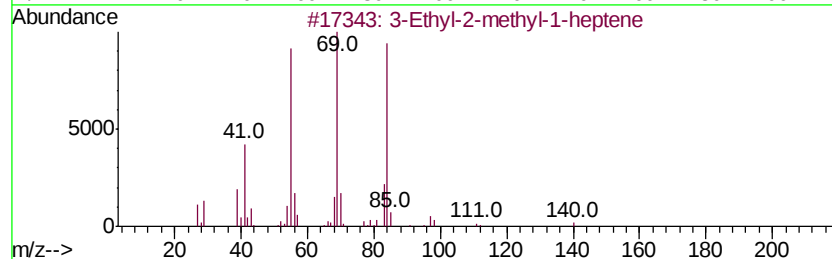
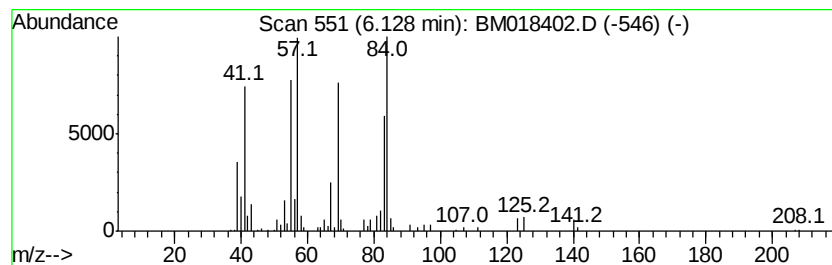
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM122718.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 1 3-Ethyl-2-methyl-1-heptene Concentration Rank 14

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 6.13 | 2.72 ng | 251129 | 1,4-Dichlorobenzene-d4 | 7.78 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|---------|--------------|------|
| 1    |    |   | 3-Ethyl-2-methyl-1-heptene          | 140 | C10H20  | 019780-60-0  | 59   |
| 2    |    |   | 2-Pentenal, (E)-                    | 84  | C5H8O   | 001576-87-0  | 55   |
| 3    |    |   | 2-Pentene, 4-methyl-, (Z)-          | 84  | C6H12   | 000691-38-3  | 47   |
| 4    |    |   | 2-Butene, 2,3-dimethyl-             | 84  | C6H12   | 000563-79-1  | 47   |
| 5    |    |   | Cyclopentane, 1,2,3,4,5-pentamet... | 140 | C10H20  | 1000152-79-7 | 46   |



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Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM122718.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

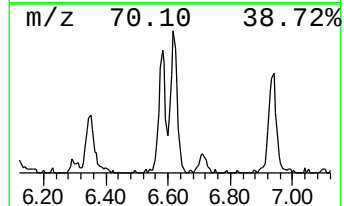
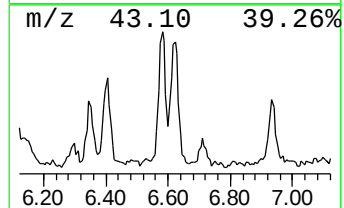
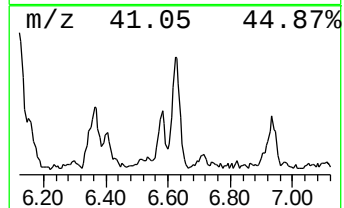
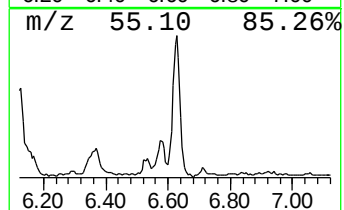
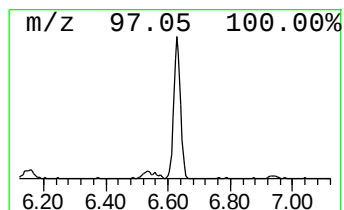
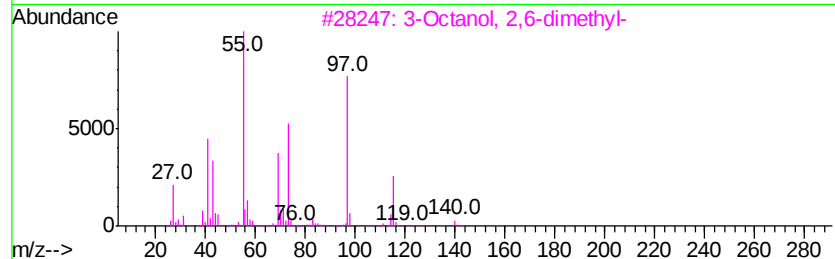
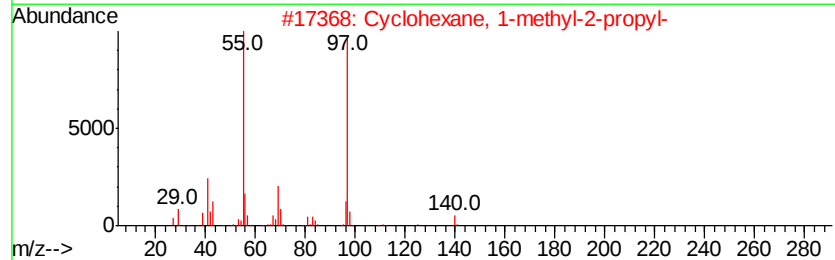
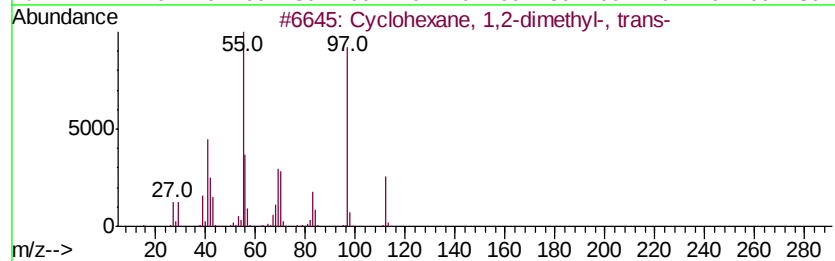
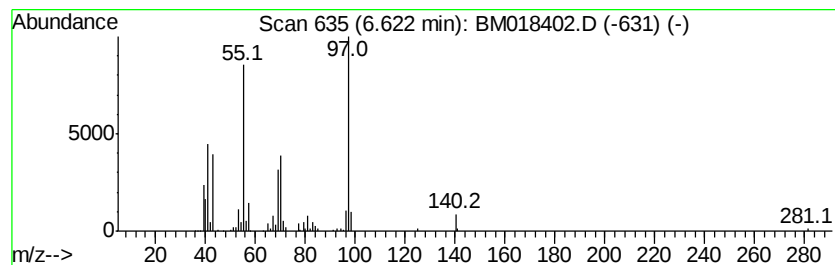
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 2 Cyclohexane, 1,2-dimethyl-,... Concentration Rank 13

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 6.62 | 3.07 ng | 283103 | 1,4-Dichlorobenzene-d4 | 7.78 |

| Hit# | of | Tentative ID                       | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Cyclohexane, 1,2-dimethyl-, trans- | 112 | C8H16   | 006876-23-9 | 72   |
| 2    |    | Cyclohexane, 1-methyl-2-propyl-    | 140 | C10H20  | 004291-79-6 | 53   |
| 3    |    | 3-Octanol, 2,6-dimethyl-           | 158 | C10H22O | 018479-55-5 | 50   |
| 4    |    | 7-Tridecanol                       | 200 | C13H28O | 000927-45-7 | 50   |
| 5    |    | 2-Pentene, 2,3,4-trimethyl-        | 112 | C8H16   | 000565-77-5 | 50   |



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BNA\_M  
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381128115-4

Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM122718.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

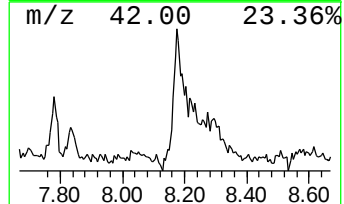
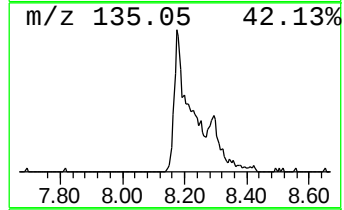
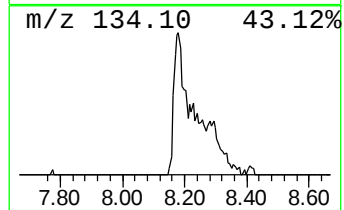
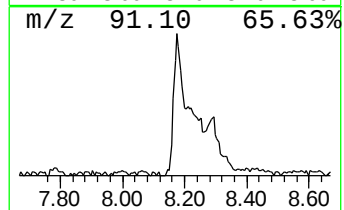
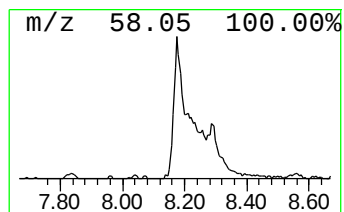
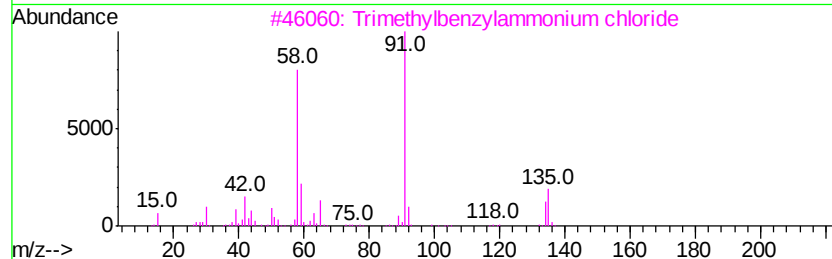
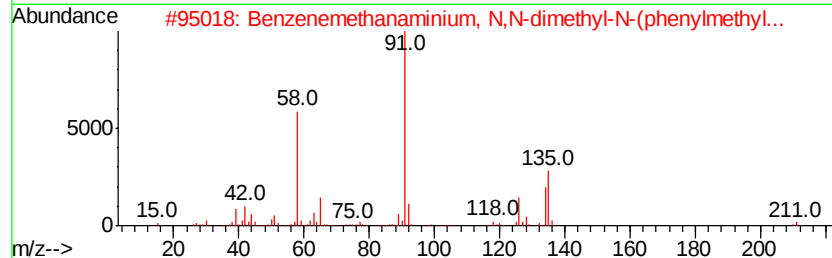
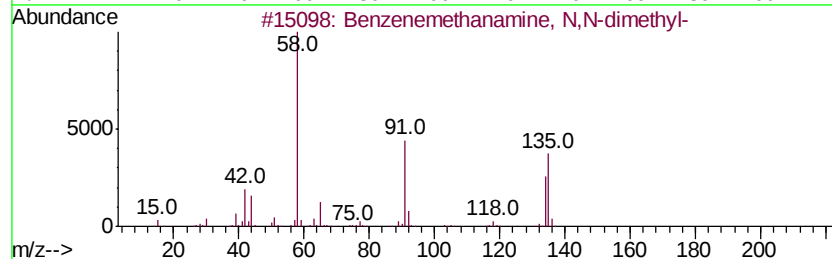
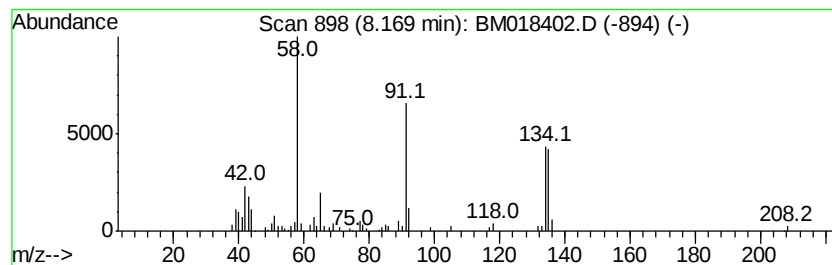
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 3 Benzenemethanamine, N,N-dim... Concentration Rank 18

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 8.17 | 2.07 ng | 191329 | 1,4-Dichlorobenzene-d4 | 7.78 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#        | Qual |
|------|----|-------------------------------------|-----|-----------|-------------|------|
| 1    | 5  | Benzenemethanamine, N,N-dimethyl-   | 135 | C9H13N    | 000103-83-3 | 90   |
| 2    |    | Benzenemethanaminium, N,N-dimeth... | 261 | C16H20ClN | 000100-94-7 | 72   |
| 3    |    | Trimethylbenzylammonium chloride    | 185 | C10H16ClN | 000056-93-9 | 56   |
| 4    |    | N,N-Dimethyl-2-aminoethanol         | 89  | C4H11NO   | 000108-01-0 | 38   |
| 5    |    | 2-Dimethylamino-2'-methoxyacetop... | 193 | C11H15NO2 | 031543-50-7 | 32   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM122718\  
Data File : BM018402.D  
Acq On : 27 Dec 2018 18:41  
Operator : JU/SJ  
Sample : J6569-03  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
381128115-4

Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM122718.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

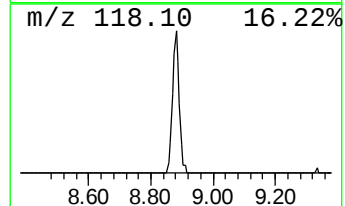
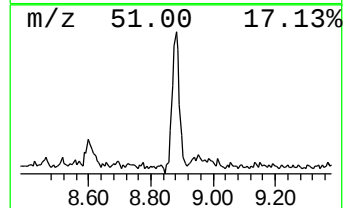
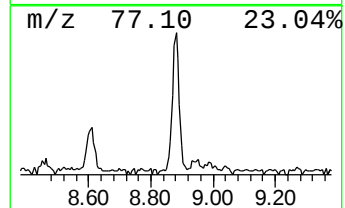
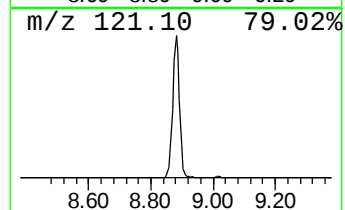
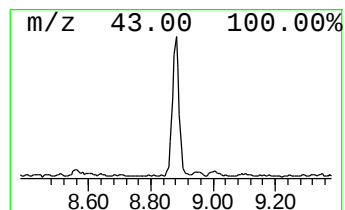
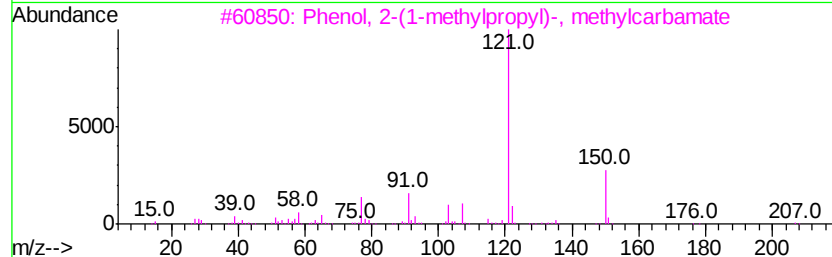
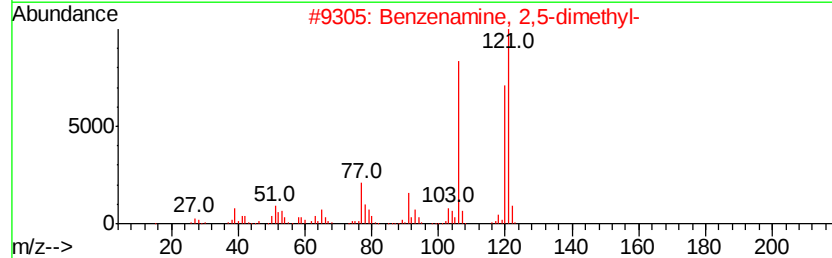
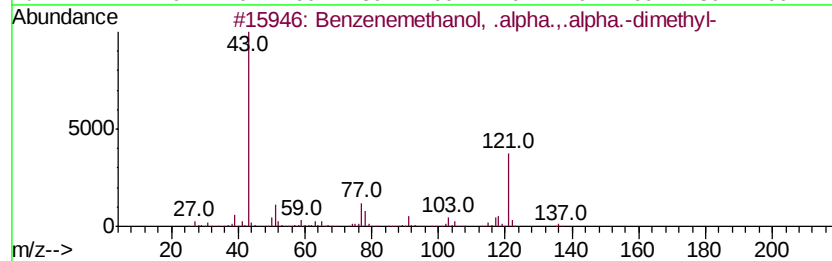
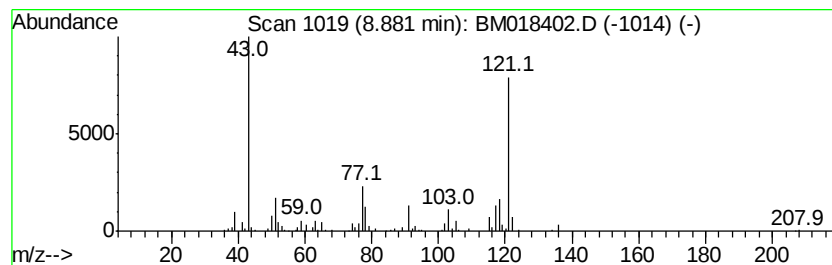
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TIC Integration Parameters: LSCINT.P

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Peak Number 4 Benzenemethanol, .alpha.,.a... Concentration Rank 11

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 8.88 | 3.83 ng | 353695 | 1,4-Dichlorobenzene-d4 | 7.78 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#        | Qual |
|------|----|-------------------------------------|-----|-----------|-------------|------|
| 1    | 5  | Benzenemethanol, .alpha.,.alpha.... | 136 | C9H12O    | 000617-94-7 | 95   |
| 2    |    | Benzenamine, 2,5-dimethyl-          | 121 | C8H11N    | 000095-78-3 | 72   |
| 3    |    | Phenol, 2-(1-methylpropyl)-, met... | 207 | C12H17NO2 | 003766-81-2 | 53   |
| 4    |    | Benzenamine, 3,4-dimethyl-          | 121 | C8H11N    | 000095-64-7 | 53   |
| 5    |    | Benzene, (1-methoxyethyl)-          | 136 | C9H12O    | 004013-34-7 | 53   |



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Acq On : 27 Dec 2018 18:41  
Operator : JU/SJ  
Sample : J6569-03  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
381128115-4

Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM122718.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

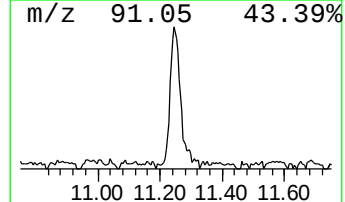
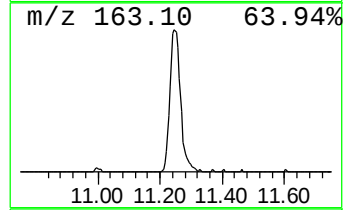
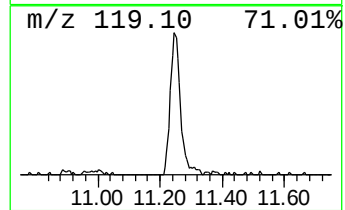
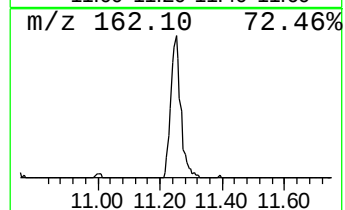
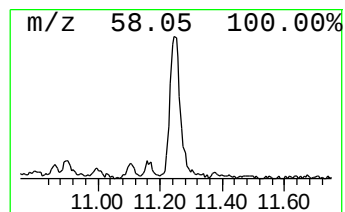
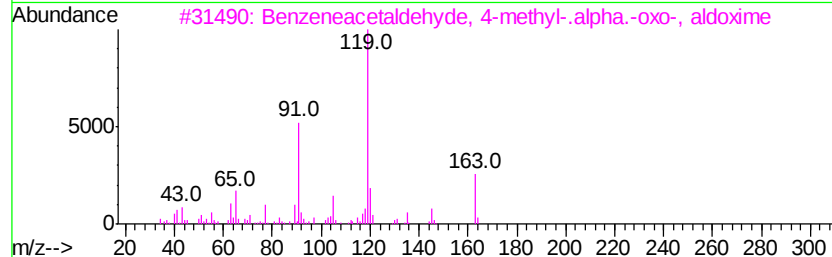
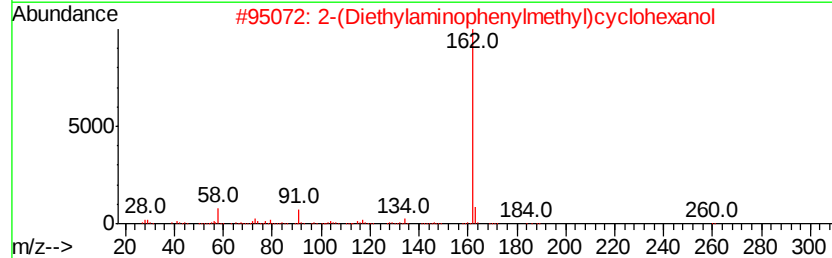
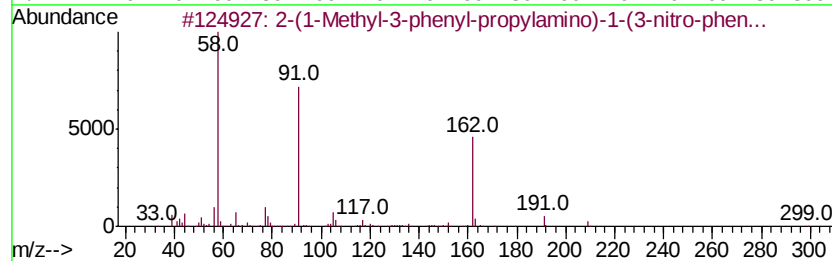
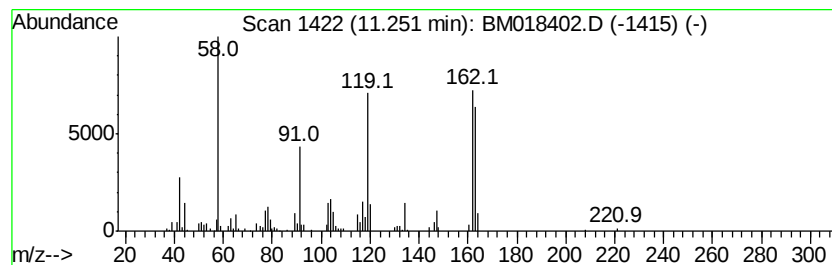
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TIC Integration Parameters: LSCINT.P

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Peak Number 5 unknown11.25 Concentration Rank 15

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 11.25 | 2.58 ng | 374901 | Naphthalene-d8   | 10.57 |

| Hit# | of | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|-------------------------------------|-----|------------|--------------|------|
| 1    | 5  | 2-(1-Methyl-3-phenyl-propylamino... | 314 | C18H22N2O3 | 1000296-61-0 | 38   |
| 2    |    | 2-(Diethylaminophenylmethyl)cycl... | 261 | C17H27NO   | 1000192-67-4 | 27   |
| 3    |    | Benzeneacetaldehyde, 4-methyl-.a... | 163 | C9H9NO2    | 017628-76-1  | 25   |
| 4    |    | Benzonitrile, 2,4-dimethoxy-        | 163 | C9H9NO2    | 004107-65-7  | 22   |
| 5    |    | 2-(2-Propenyl)-m-anisidine          | 163 | C10H13NO   | 194782-58-6  | 18   |



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Sample : J6569-03  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
381128115-4

Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM122718.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

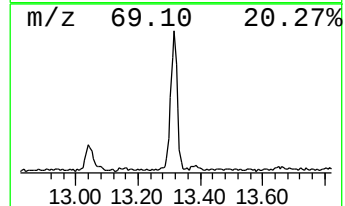
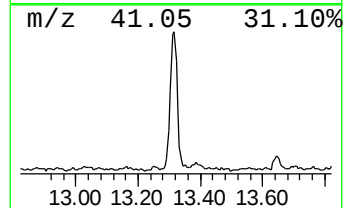
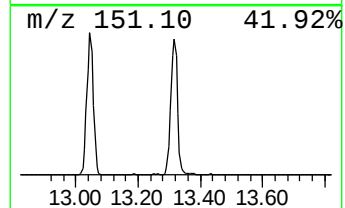
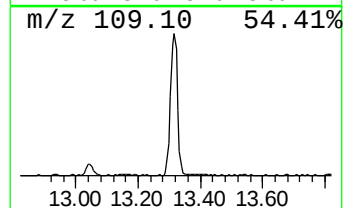
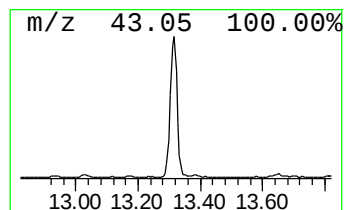
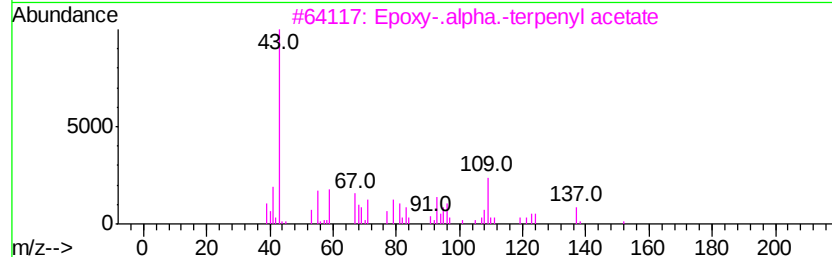
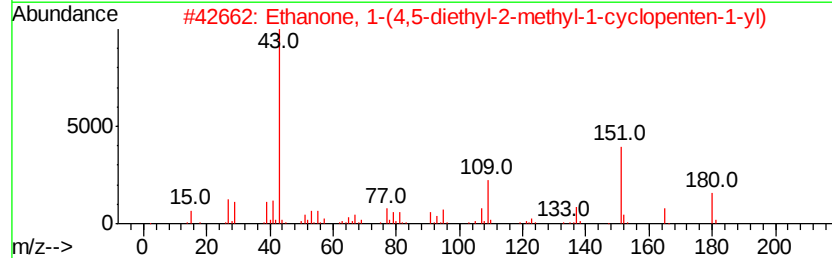
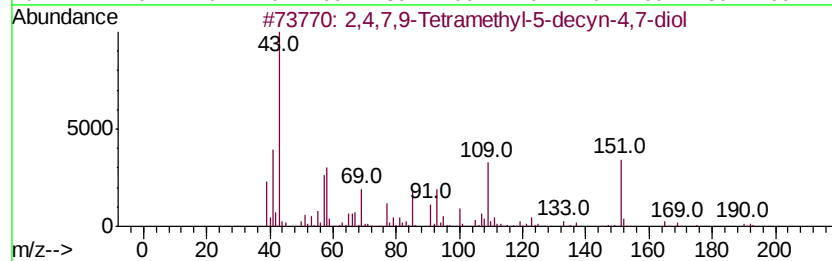
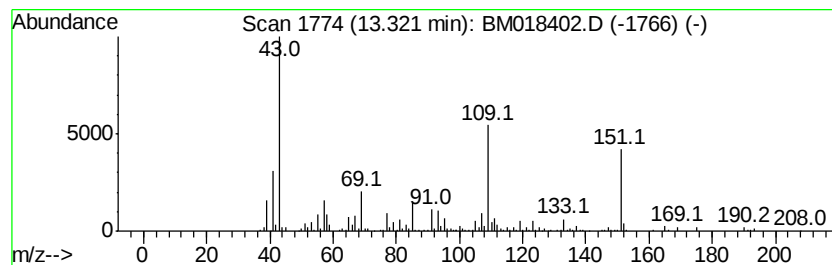
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TIC Integration Parameters: LSCINT.P

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Peak Number 6 unknown13.32 Concentration Rank 8

| R.T.  | EstConc | Area    | Relative to ISTD | R.T.  |
|-------|---------|---------|------------------|-------|
| 13.32 | 4.95 ng | 1065700 | Acenaphthene-d10 | 14.43 |

| Hit# | of                                    | 5   | Tentative ID | MW           | MolForm | CAS# | Qual |
|------|---------------------------------------|-----|--------------|--------------|---------|------|------|
| 1    | 2,4,7,9-Tetramethyl-5-decyn-4,7-...   | 226 | C14H26O2     | 000126-86-3  | 40      |      |      |
| 2    | Ethanone, 1-(4,5-diethyl-2-methyl-... | 180 | C12H20O      | 062338-24-3  | 36      |      |      |
| 3    | Epoxy-.alpha.-terpenyl acetate        | 212 | C12H20O3     | 1000293-00-8 | 25      |      |      |
| 4    | Geranyl acetate, 2,3-epoxy-           | 212 | C12H20O3     | 050727-95-2  | 22      |      |      |
| 5    | 3-Hexyne-2,5-diol, 2,5-dimethyl-      | 142 | C8H14O2      | 000142-30-3  | 12      |      |      |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM122718\  
Data File : BM018402.D  
Acq On : 27 Dec 2018 18:41  
Operator : JU/SJ  
Sample : J6569-03  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

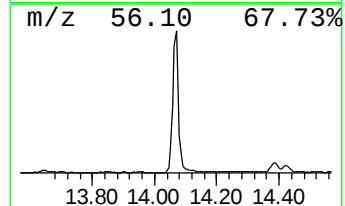
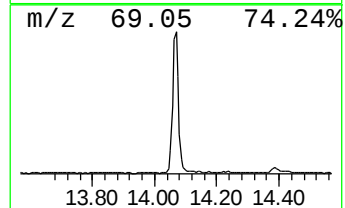
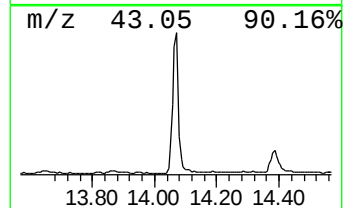
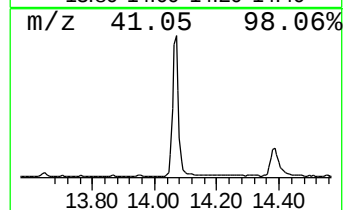
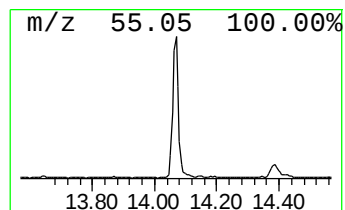
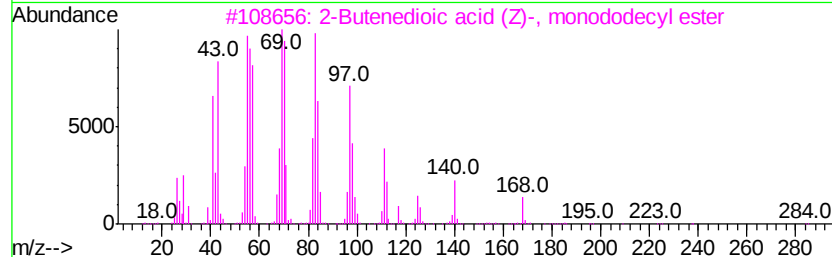
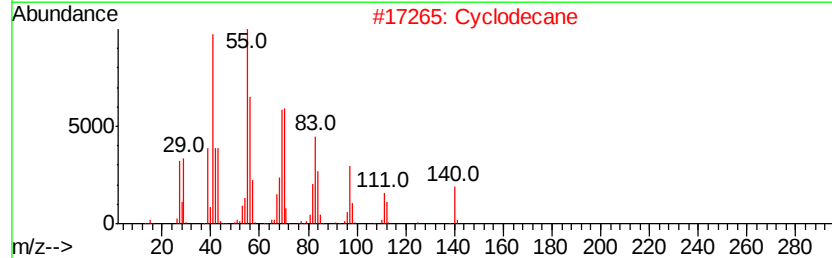
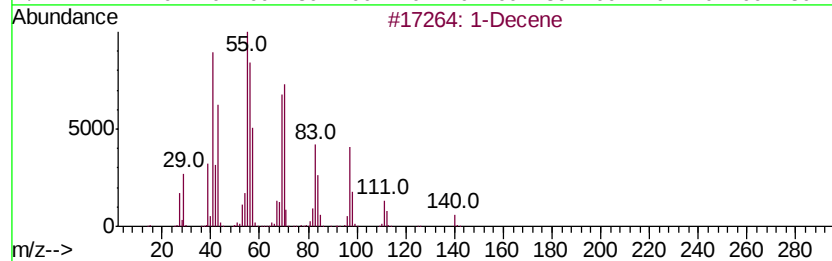
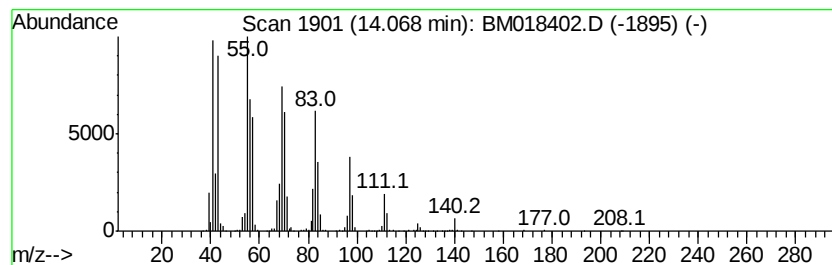
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ClientSampleId :  
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM122718.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 7 1-Decene Concentration Rank 4

| R.T.    | EstConc                             | Area         | Relative to ISTD | R.T.           |
|---------|-------------------------------------|--------------|------------------|----------------|
| 14.07   | 9.82 ng                             | 2112310      | Acenaphthene-d10 | 14.43          |
| Hit# of | 5                                   | Tentative ID | MW MolForm       | CAS# Qual      |
| 1       | 1-Decene                            |              | 140 C10H20       | 000872-05-9 93 |
| 2       | Cyclodecane                         |              | 140 C10H20       | 000293-96-9 93 |
| 3       | 2-Butenedioic acid (Z)-, monodod... |              | 284 C16H28O4     | 002424-61-5 91 |
| 4       | Cyclododecane                       |              | 168 C12H24       | 000294-62-2 91 |
| 5       | 1-Tridecanol                        |              | 200 C13H28O      | 000112-70-9 91 |



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Operator : JU/SJ  
Sample : J6569-03  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
381128115-4

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

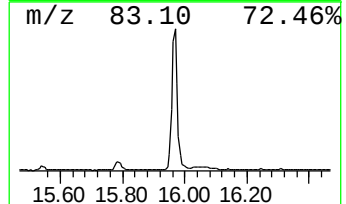
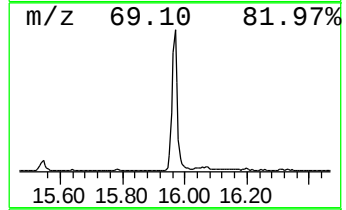
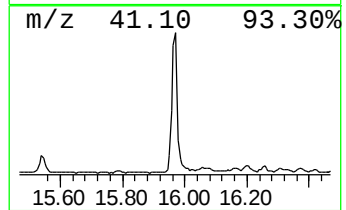
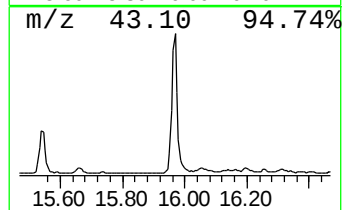
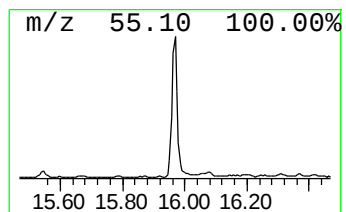
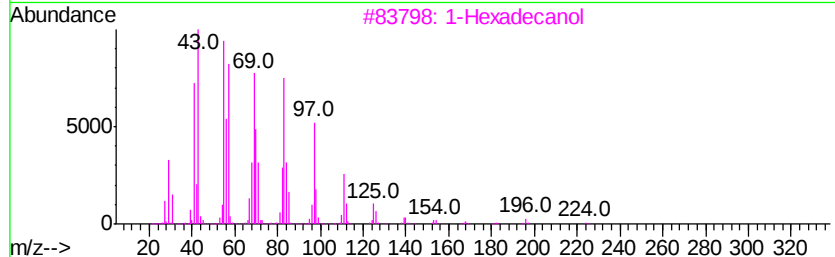
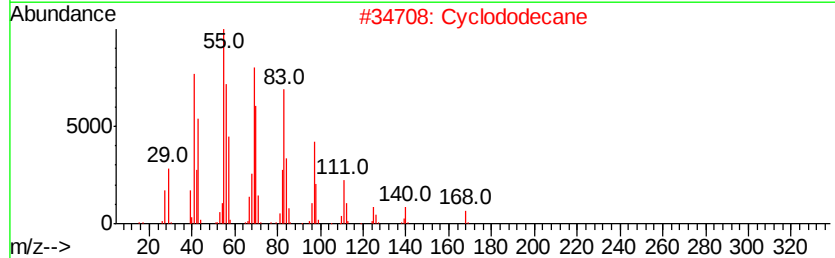
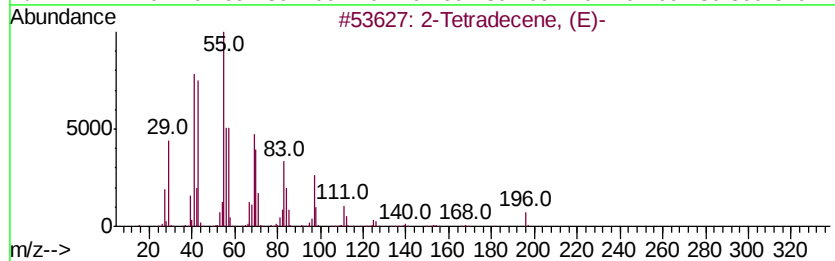
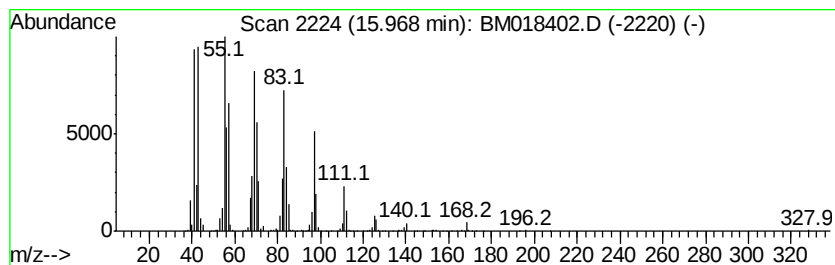
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TIC Integration Parameters: LSCINT.P

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Peak Number 8 2-Tetradecene, (E)- Concentration Rank 5

| R.T.  | EstConc | Area    | Relative to ISTD | R.T.  |
|-------|---------|---------|------------------|-------|
| 15.97 | 8.08 ng | 2042740 | Phenanthrene-d10 | 17.17 |

| Hit# | of | Tentative ID        | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------|-----|---------|-------------|------|
| 1    | 5  | 2-Tetradecene, (E)- | 196 | C14H28  | 035953-53-8 | 97   |
| 2    |    | Cyclododecane       | 168 | C12H24  | 000294-62-2 | 97   |
| 3    |    | 1-Hexadecanol       | 242 | C16H34O | 036653-82-4 | 93   |
| 4    |    | 1-Pentadecanol      | 228 | C15H32O | 000629-76-5 | 91   |
| 5    |    | 1-Tridecanol        | 200 | C13H28O | 000112-70-9 | 91   |



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Acq On : 27 Dec 2018 18:41  
Operator : JU/SJ  
Sample : J6569-03  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

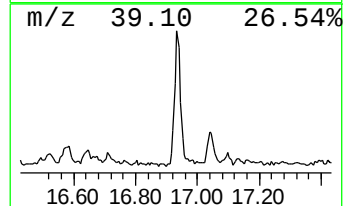
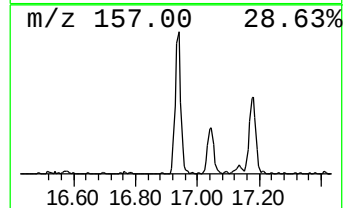
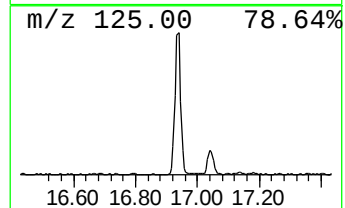
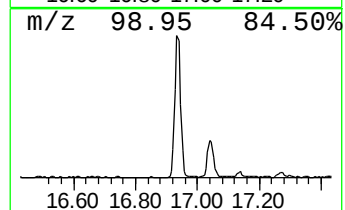
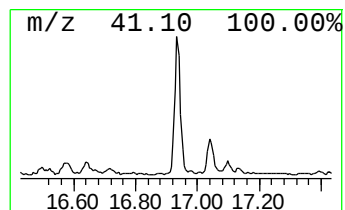
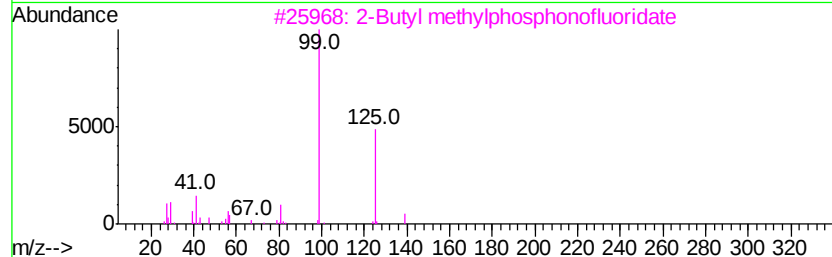
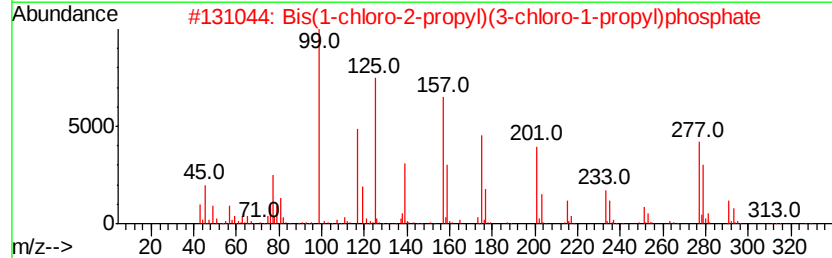
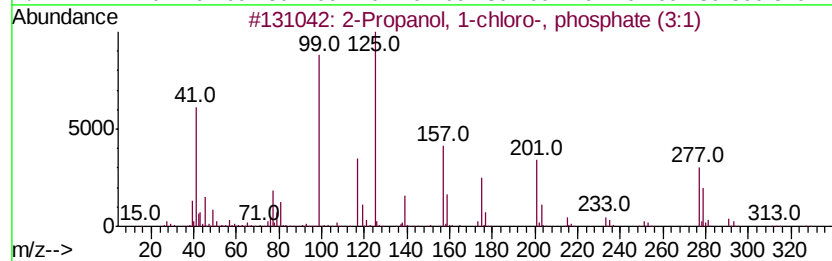
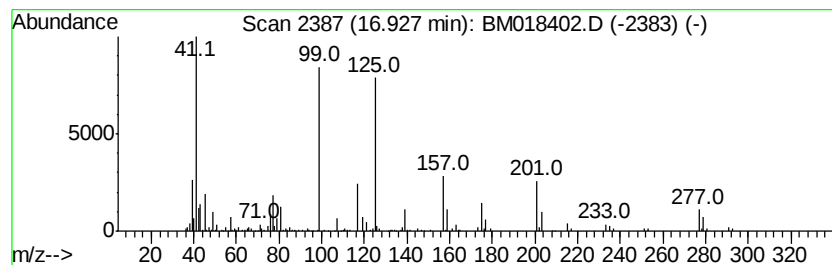
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ClientSampleId :  
381128115-4

Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM122718.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 9 2-Propanol, 1-chloro-, phos... Concentration Rank 12

| R.T.    | EstConc                             | Area         | Relative to ISTD | R.T.         |      |      |
|---------|-------------------------------------|--------------|------------------|--------------|------|------|
| 16.93   | 3.48 ng                             | 879420       | Phenanthrene-d10 | 17.17        |      |      |
| Hit# of | 5                                   | Tentative ID | MW               | MolForm      | CAS# | Qual |
| 1       | 2-Propanol, 1-chloro-, phosphate... | 326          | C9H18Cl3O4P      | 013674-84-5  | 94   |      |
| 2       | Bis(1-chloro-2-propyl)(3-chloro-... | 326          | C9H18Cl3O4P      | 137909-40-1  | 46   |      |
| 3       | 2-Butyl methylphosphonofluoridate   | 154          | C5H12F02P        | 000352-52-3  | 17   |      |
| 4       | Oxamide, N-methyl-N'-(2-(4-piper... | 214          | C9H18N4O2        | 294877-47-7  | 14   |      |
| 5       | Cholestan-22(26)-isoepoxy           | 386          | C27H46O          | 1000252-59-7 | 10   |      |



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Data File : BM018402.D  
Acq On : 27 Dec 2018 18:41  
Operator : JU/SJ  
Sample : J6569-03  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

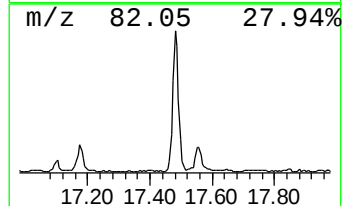
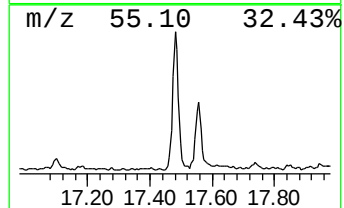
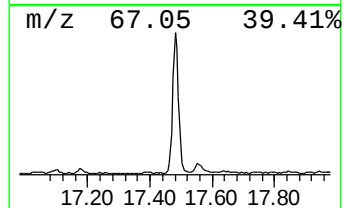
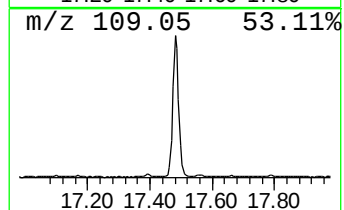
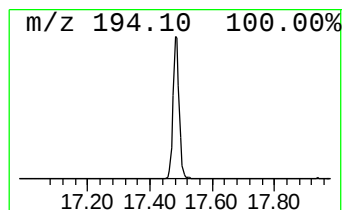
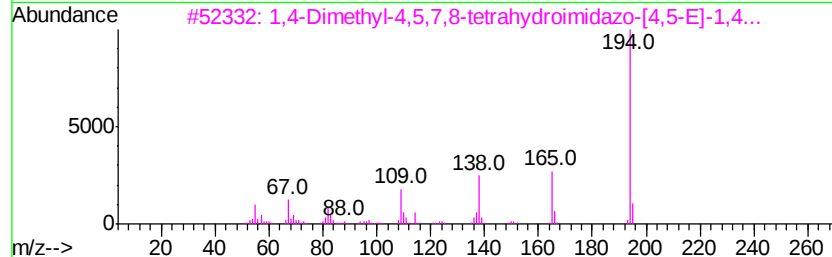
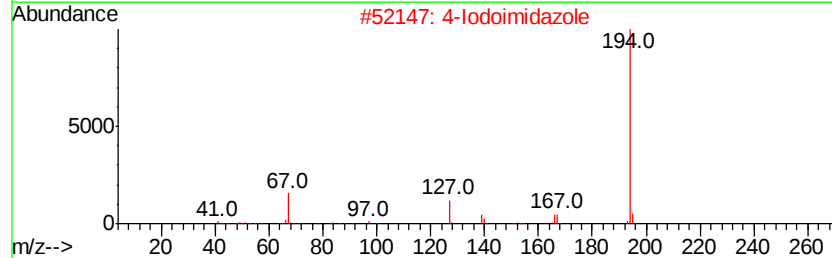
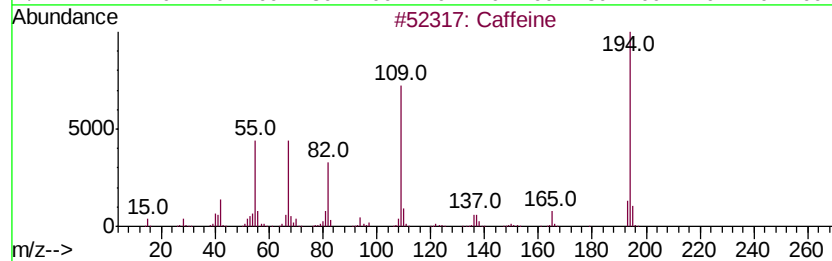
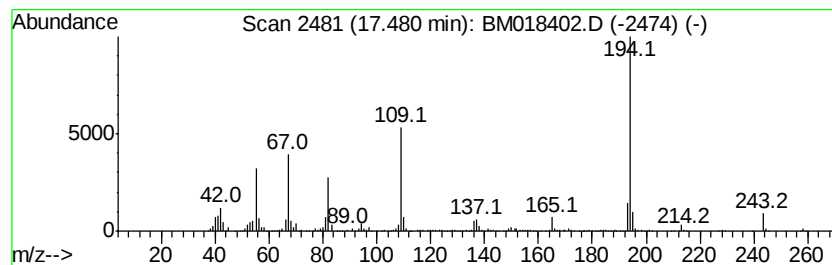
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ClientSampleId :  
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 10 Caffeine Concentration Rank 7

| R.T.      | EstConc                             | Area    | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------------|---------|------------------|-------------|------|
| 17.48     | 5.95 ng                             | 1502860 | Phenanthrene-d10 | 17.17       |      |
| Hit# of 5 | Tentative ID                        | MW      | MolForm          | CAS#        | Qual |
| 1         | Caffeine                            | 194     | C8H10N4O2        | 000058-08-2 | 98   |
| 2         | 4-Iodoimidazole                     | 194     | C3H3IN2          | 071759-89-2 | 49   |
| 3         | 1,4-Dimethyl-4,5,7,8-tetrahydroi... | 194     | C8H10N4O2        | 130063-15-9 | 43   |
| 4         | 3(2H)-Benzofuranone, 4,6-dimethoxy- | 194     | C10H10O4         | 004225-35-8 | 38   |
| 5         | 3,3'-Dimethyl-5-methoxy-4,5'-bii... | 194     | C9H10N2O3        | 019668-80-5 | 32   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM122718\  
Data File : BM018402.D  
Acq On : 27 Dec 2018 18:41  
Operator : JU/SJ  
Sample : J6569-03  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
381128115-4

Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM122718.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

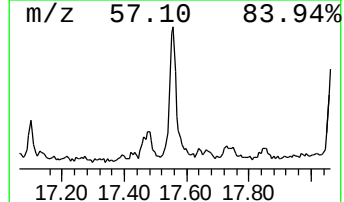
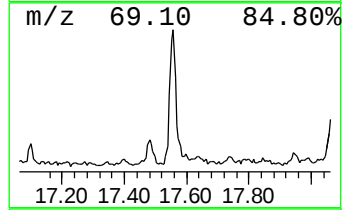
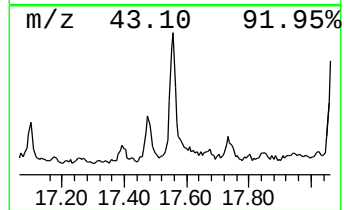
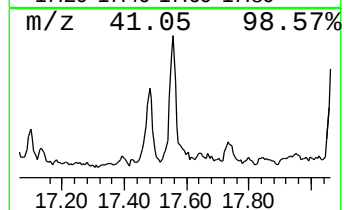
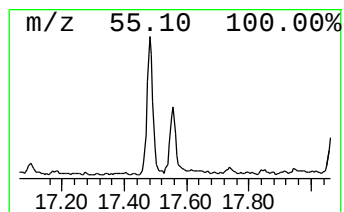
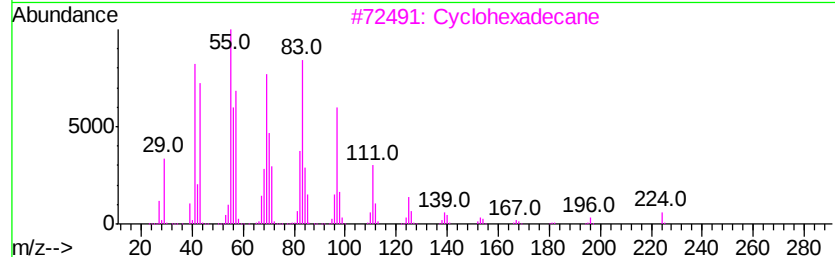
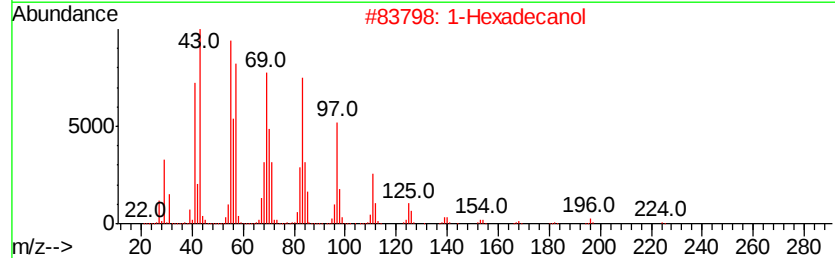
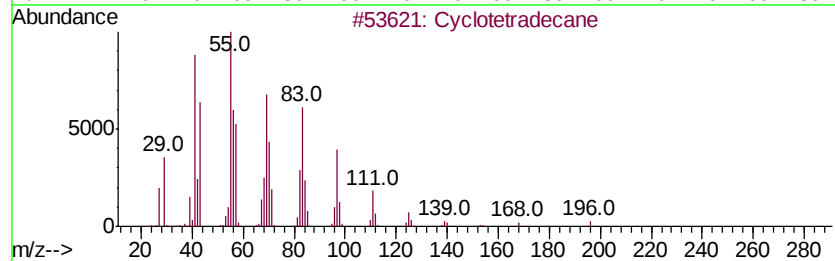
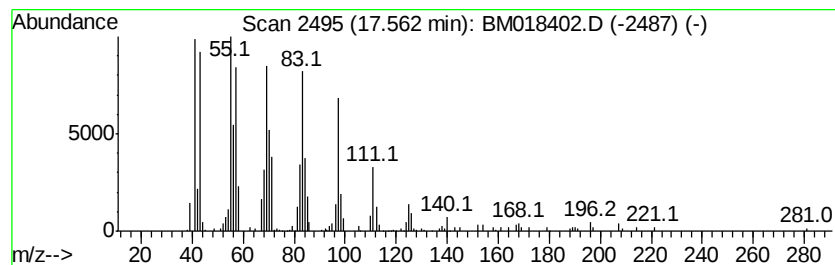
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 11 Cyclotetradecane Concentration Rank 17

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 17.56 | 2.13 ng | 538820 | Phenanthrene-d10 | 17.17 |

| Hit# | of | 5 | Tentative ID     | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------|-----|---------|-------------|------|
| 1    |    |   | Cyclotetradecane | 196 | C14H28  | 000295-17-0 | 98   |
| 2    |    |   | 1-Hexadecanol    | 242 | C16H34O | 036653-82-4 | 95   |
| 3    |    |   | Cyclohexadecane  | 224 | C16H32  | 000295-65-8 | 94   |
| 4    |    |   | 1-Pentadecanol   | 228 | C15H32O | 000629-76-5 | 91   |
| 5    |    |   | 1-Pentadecene    | 210 | C15H30  | 013360-61-7 | 91   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM122718\  
Data File : BM018402.D  
Acq On : 27 Dec 2018 18:41  
Operator : JU/SJ  
Sample : J6569-03  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
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ClientSampleId :  
381128115-4

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

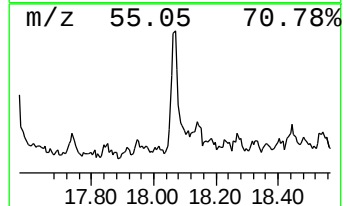
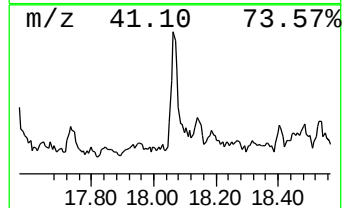
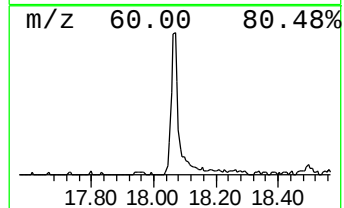
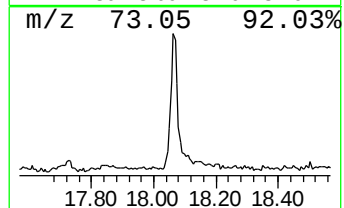
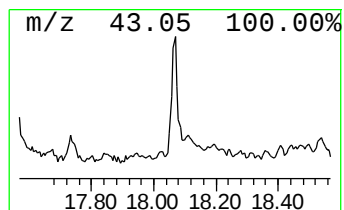
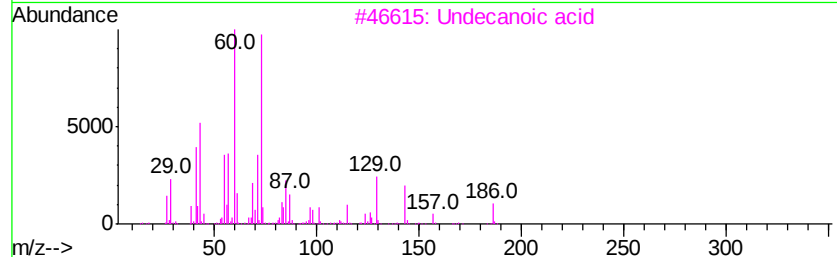
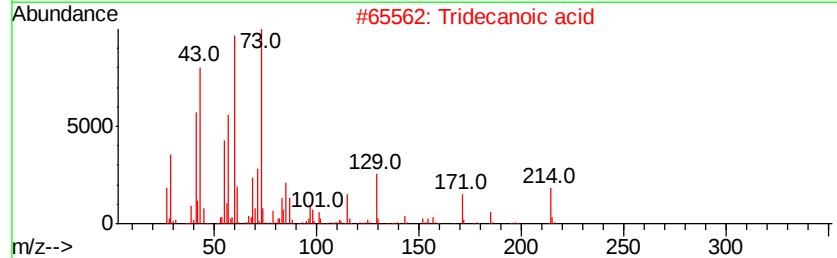
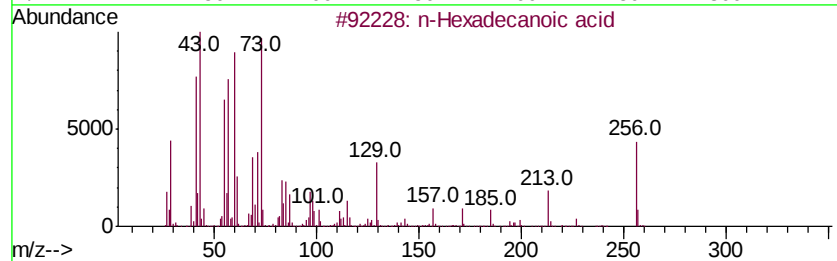
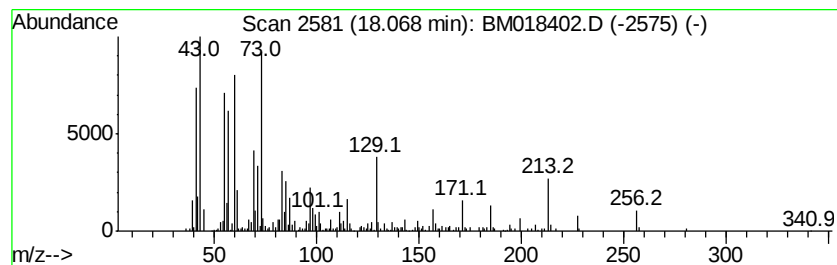
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 12 n-Hexadecanoic acid Concentration Rank 16

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 18.07 | 2.15 ng | 542855 | Phenanthrene-d10 | 17.17 |

| Hit# | of | Tentative ID        | MW  | MolForm  | CAS#        | Qual |
|------|----|---------------------|-----|----------|-------------|------|
| 1    | 5  | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 96   |
| 2    |    | Tridecanoic acid    | 214 | C13H26O2 | 000638-53-9 | 64   |
| 3    |    | Undecanoic acid     | 186 | C11H22O2 | 000112-37-8 | 62   |
| 4    |    | Tetradecanoic acid  | 228 | C14H28O2 | 000544-63-8 | 50   |
| 5    |    | Nonanoic acid       | 158 | C9H18O2  | 000112-05-0 | 50   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM122718\  
Data File : BM018402.D  
Acq On : 27 Dec 2018 18:41  
Operator : JU/SJ  
Sample : J6569-03  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

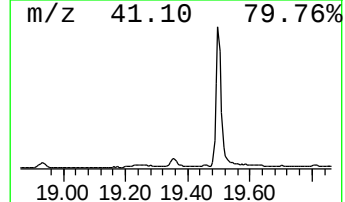
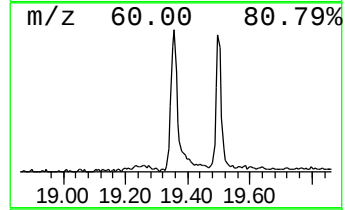
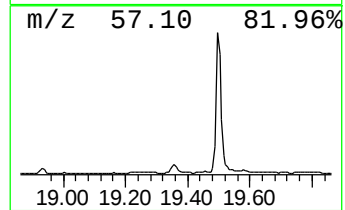
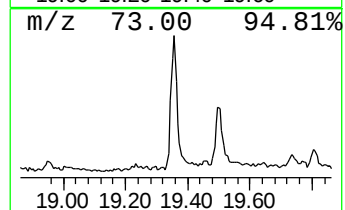
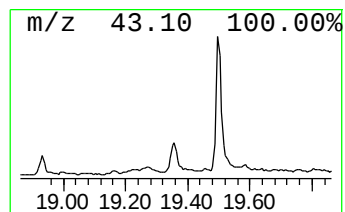
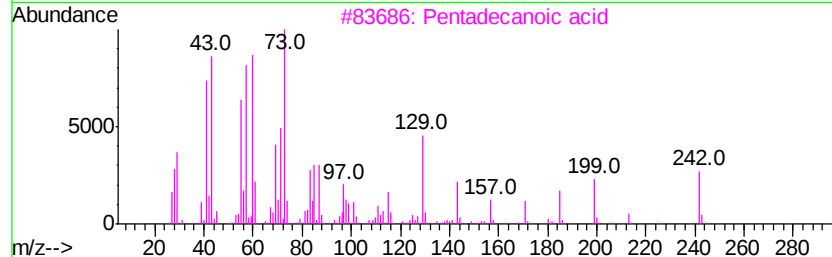
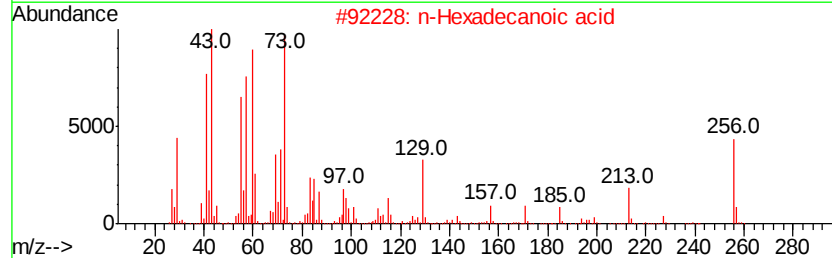
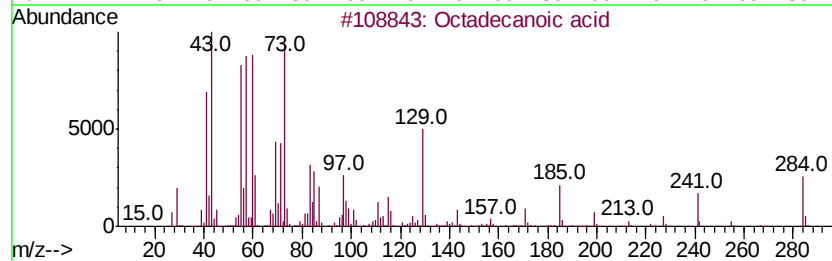
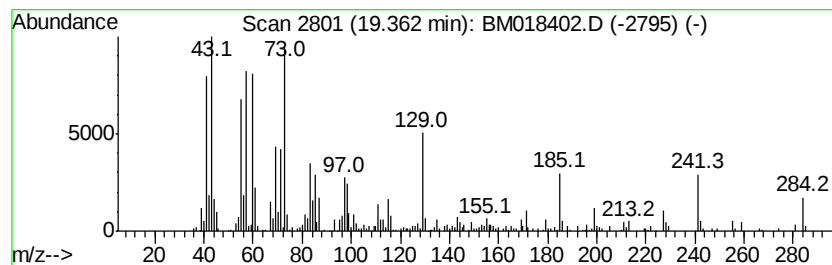
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TIC Integration Parameters: LSCINT.P

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Peak Number 13 Octadecanoic acid Concentration Rank 19

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 19.36 | 2.04 ng | 673503 | Chrysene-d12     | 21.37 |

| Hit# | of | 5 | Tentative ID                       | MW  | MolForm  | CAS#        | Qual |
|------|----|---|------------------------------------|-----|----------|-------------|------|
| 1    |    |   | Octadecanoic acid                  | 284 | C18H36O2 | 000057-11-4 | 99   |
| 2    |    |   | n-Hexadecanoic acid                | 256 | C16H32O2 | 000057-10-3 | 83   |
| 3    |    |   | Pentadecanoic acid                 | 242 | C15H30O2 | 001002-84-2 | 76   |
| 4    |    |   | Tetradecanoic acid                 | 228 | C14H28O2 | 000544-63-8 | 74   |
| 5    |    |   | Octadecanoic acid, 2-(2-hydroxy... | 372 | C22H44O4 | 000106-11-6 | 59   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM122718\  
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Operator : JU/SJ  
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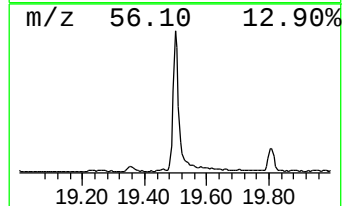
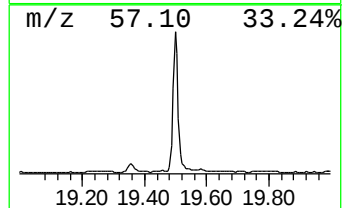
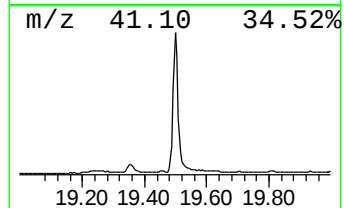
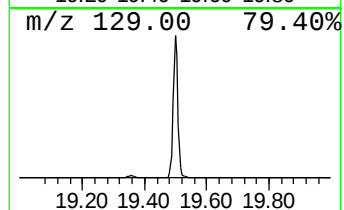
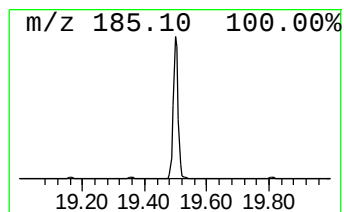
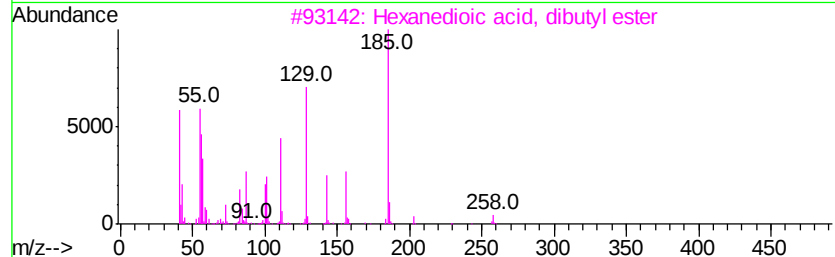
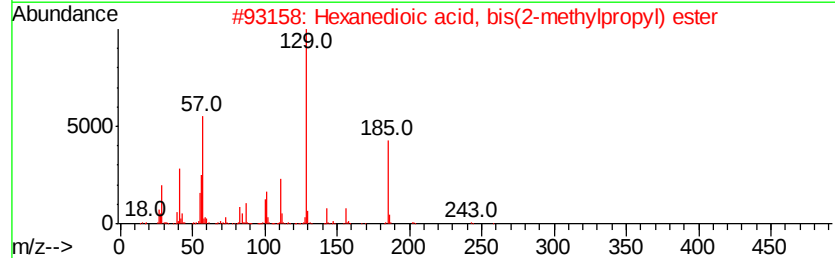
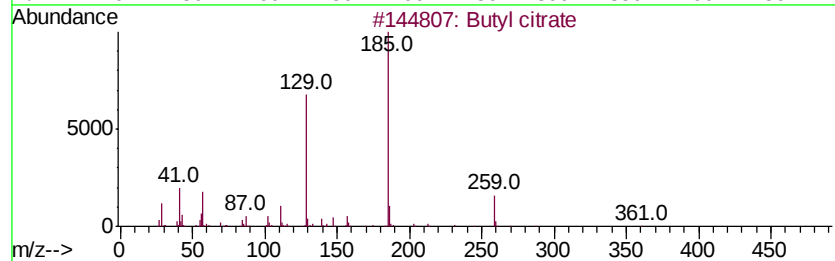
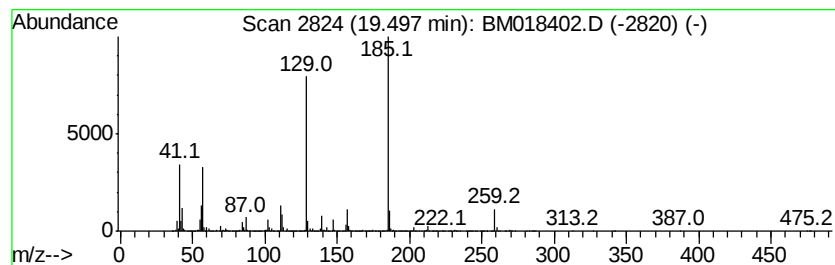
Instrument :  
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ClientSampleId :  
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM122718.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 14 Butyl citrate Concentration Rank 2

| R.T.    | EstConc                             | Area         | Relative to ISTD | R.T.      |
|---------|-------------------------------------|--------------|------------------|-----------|
| 19.50   | 20.52 ng                            | 6764670      | Chrysene-d12     | 21.37     |
| Hit# of | 5                                   | Tentative ID | MW MolForm       | CAS# Qual |
| 1       | Butyl citrate                       | 360 C18H32O7 | 000077-94-1      | 87        |
| 2       | Hexanedioic acid, bis(2-methylpr... | 258 C14H26O4 | 000141-04-8      | 50        |
| 3       | Hexanedioic acid, dibutyl ester     | 258 C14H26O4 | 000105-99-7      | 39        |
| 4       | Hexanedioic acid, bis(1-methylpr... | 258 C14H26O4 | 038447-22-2      | 38        |
| 5       | 1,3,5-Triazine, 2-amino-4-ethyla... | 185 C6H11N5S | 004147-58-4      | 38        |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM122718\  
Data File : BM018402.D  
Acq On : 27 Dec 2018 18:41  
Operator : JU/SJ  
Sample : J6569-03  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
381128115-4

Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM122718.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

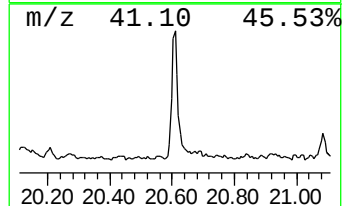
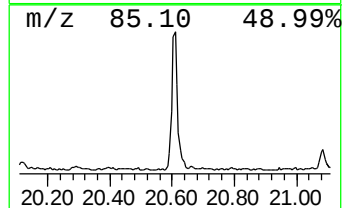
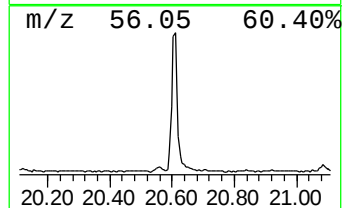
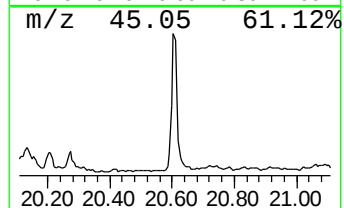
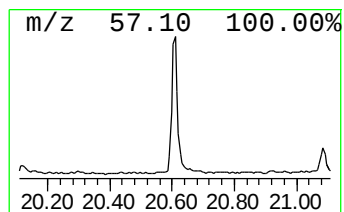
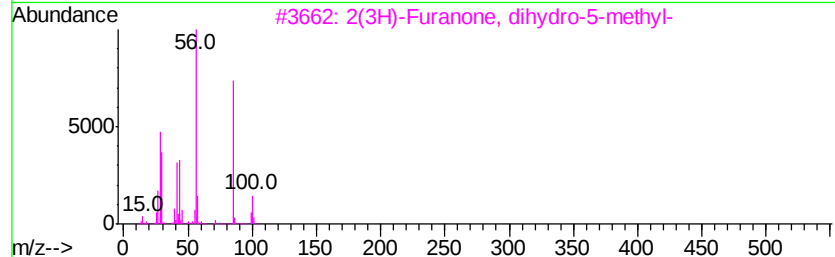
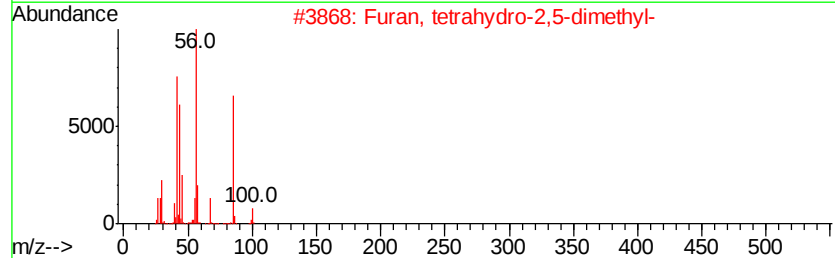
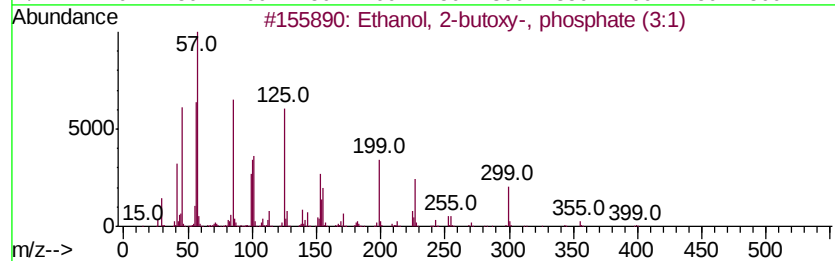
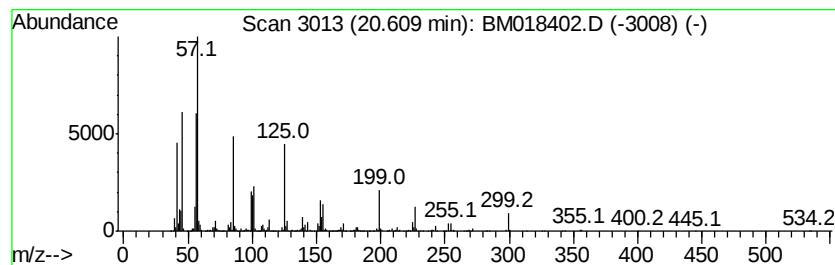
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 15 Ethanol, 2-butoxy-, phospho... Concentration Rank 9

| R.T.  | EstConc | Area    | Relative to ISTD | R.T.  |
|-------|---------|---------|------------------|-------|
| 20.61 | 4.22 ng | 1389820 | Chrysene-d12     | 21.37 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#        | Qual |
|------|----|-------------------------------------|-----|-----------|-------------|------|
| 1    | 5  | Ethanol, 2-butoxy-, phosphate (3:1) | 398 | C18H39O7P | 000078-51-3 | 81   |
| 2    |    | Furan, tetrahydro-2,5-dimethyl-     | 100 | C6H12O    | 001003-38-9 | 22   |
| 3    |    | 2(3H)-Furanone, dihydro-5-methyl-   | 100 | C5H8O2    | 000108-29-2 | 22   |
| 4    |    | Pentane, 2,2-dimethyl-              | 100 | C7H16     | 000590-35-2 | 14   |
| 5    |    | 2-Propanone, ethylhydrazone         | 100 | C5H12N2   | 007422-99-3 | 12   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM122718\  
Data File : BM018402.D  
Acq On : 27 Dec 2018 18:41  
Operator : JU/SJ  
Sample : J6569-03  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
381128115-4

Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM122718.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

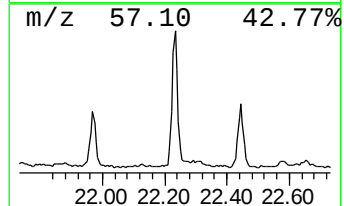
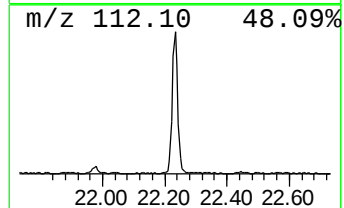
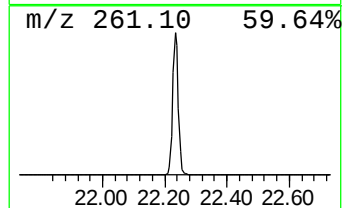
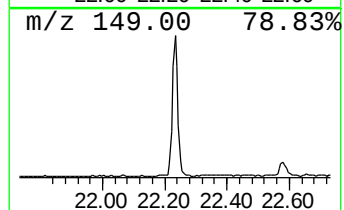
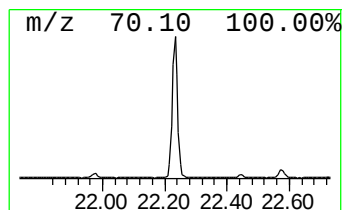
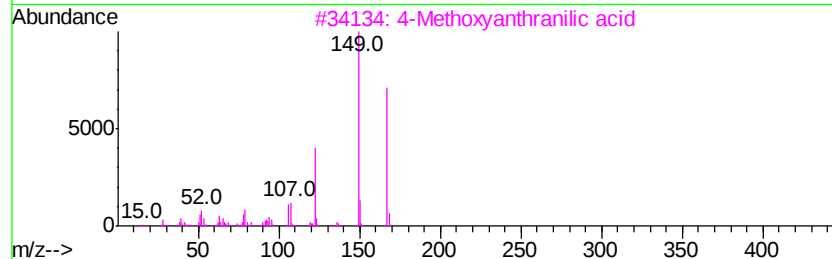
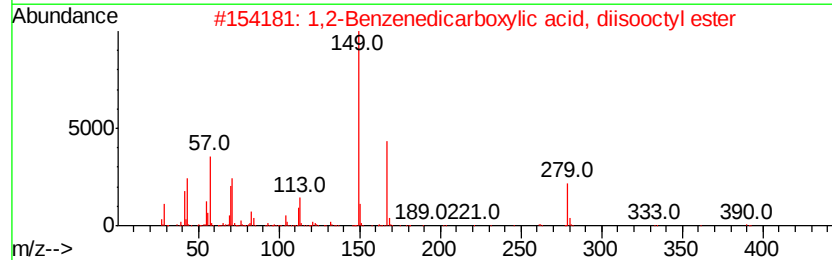
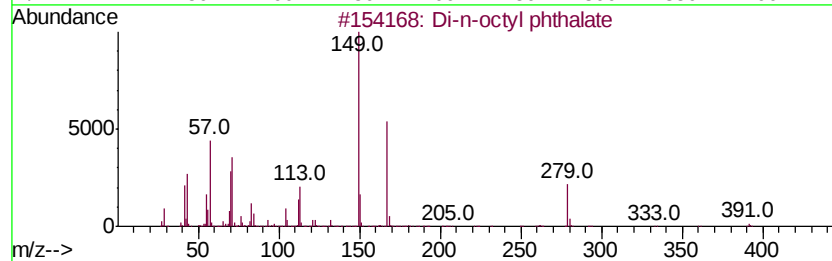
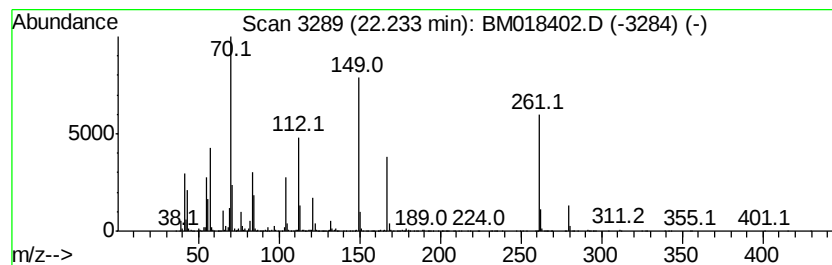
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 16 unknown22.23 Concentration Rank 6

| R.T.  | EstConc | Area    | Relative to ISTD | R.T.  |
|-------|---------|---------|------------------|-------|
| 22.23 | 7.66 ng | 2525030 | Chrysene-d12     | 21.37 |

| Hit# | of | Tentative ID                        | MW  | MolForm       | CAS#         | Qual |
|------|----|-------------------------------------|-----|---------------|--------------|------|
| 1    | 5  | Di-n-octyl phthalate                | 390 | C24H38O4      | 000117-84-0  | 30   |
| 2    |    | 1,2-Benzenedicarboxylic acid, di... | 390 | C24H38O4      | 027554-26-3  | 27   |
| 3    |    | 4-Methoxyanthranilic acid           | 167 | C8H9NO3       | 004294-95-5  | 22   |
| 4    |    | Didodecyl phthalate                 | 502 | C32H54O4      | 002432-90-8  | 22   |
| 5    |    | 9-(2',2'-Dimethylpropanoilhydraz... | 576 | C30H42Cl2N4O3 | 1000111-04-6 | 16   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM122718\  
Data File : BM018402.D  
Acq On : 27 Dec 2018 18:41  
Operator : JU/SJ  
Sample : J6569-03  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

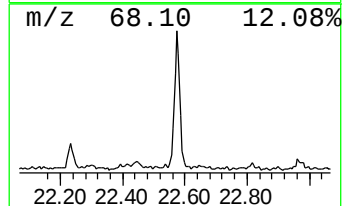
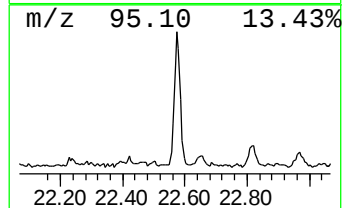
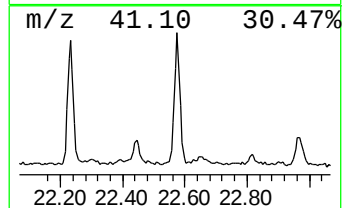
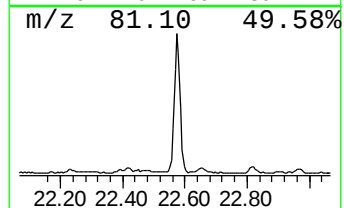
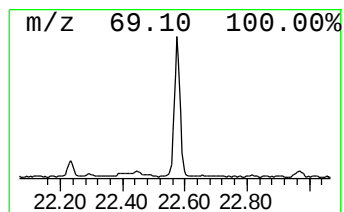
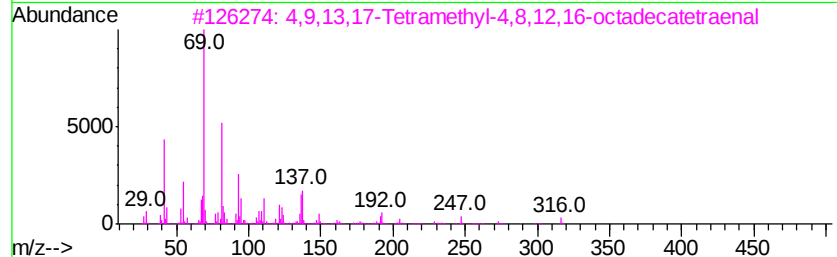
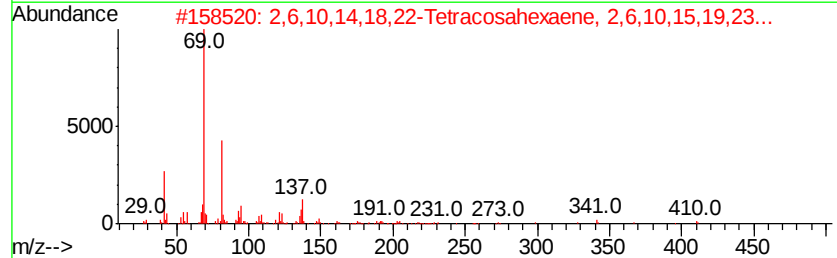
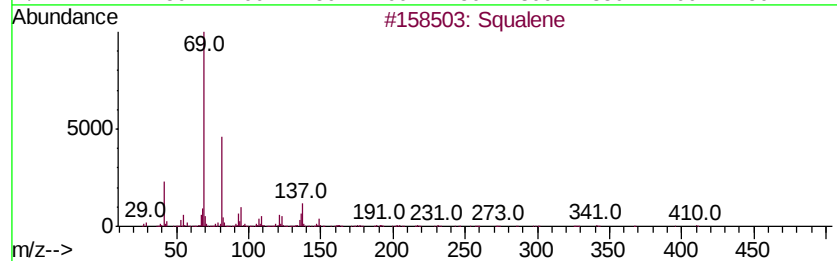
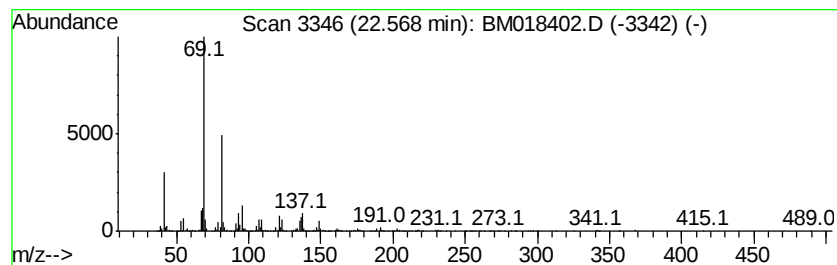
Instrument :  
BNA\_M  
ClientSampleId :  
381128115-4

Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM122718.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 17 Squalene Concentration Rank 10

| R.T.    | EstConc                             | Area         | Relative to ISTD | R.T.           |
|---------|-------------------------------------|--------------|------------------|----------------|
| 22.57   | 4.02 ng                             | 1402020      | Perylene-d12     | 23.66          |
| Hit# of | 5                                   | Tentative ID | MW MolForm       | CAS# Qual      |
| 1       | Squalene                            |              | 410 C30H50       | 007683-64-9 91 |
| 2       | 2,6,10,14,18,22-Tetracosahexaene... |              | 410 C30H50       | 000111-02-4 91 |
| 3       | 4,9,13,17-Tetramethyl-4,8,12,16-... |              | 316 C22H36O      | 056882-09-8 83 |
| 4       | 3,7,11-Tridecatrienitrile, 4,8...   |              | 231 C16H25N      | 006006-01-5 78 |
| 5       | 7,11-Dimethyldodeca-2,6,10-trien... |              | 208 C14H24O      | 125529-10-4 72 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM122718\  
Data File : BM018402.D  
Acq On : 27 Dec 2018 18:41  
Operator : JU/SJ  
Sample : J6569-03  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
381128115-4

Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM122718.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

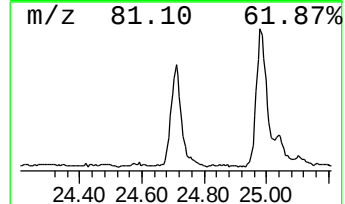
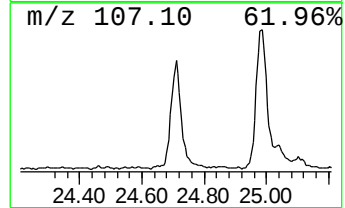
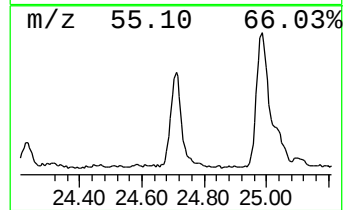
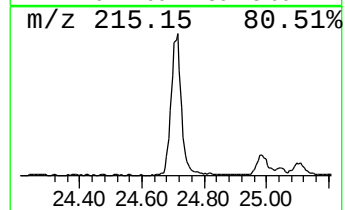
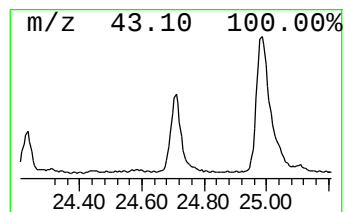
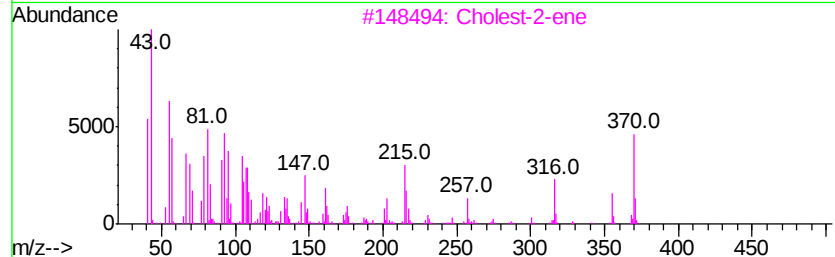
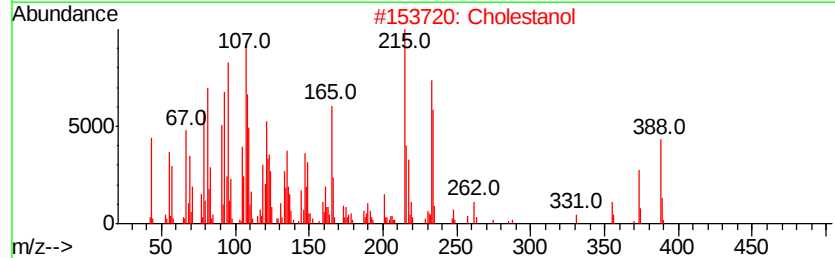
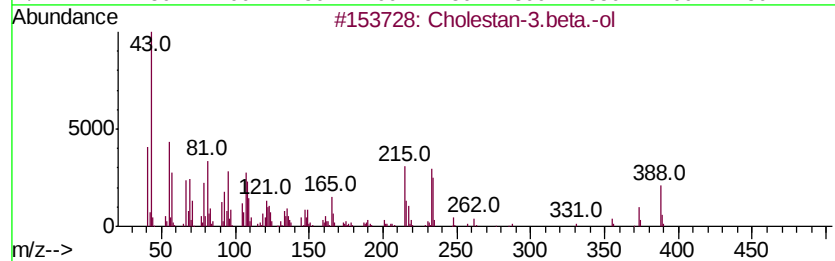
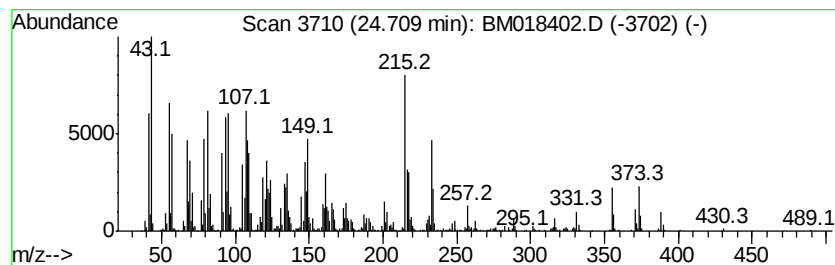
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 18 Cholestan-3.beta.-ol Concentration Rank 3

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.  |
|-------|----------|---------|------------------|-------|
| 24.71 | 11.05 ng | 3854030 | Perylene-d12     | 23.66 |

| Hit# | of | Tentative ID                       | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Cholestan-3.beta.-ol               | 388 | C27H48O | 017608-41-2 | 90   |
| 2    |    | Cholestanol                        | 388 | C27H48O | 000080-97-7 | 72   |
| 3    |    | Cholest-2-ene                      | 370 | C27H46  | 015910-23-3 | 55   |
| 4    |    | Cholestan-3-ol                     | 388 | C27H48O | 027409-41-2 | 55   |
| 5    |    | Cholestan-3-ol, (3.beta.,5.beta.)- | 388 | C27H48O | 000360-68-9 | 47   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM122718\  
Data File : BM018402.D  
Acq On : 27 Dec 2018 18:41  
Operator : JU/SJ  
Sample : J6569-03  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
381128115-4

Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM122718.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

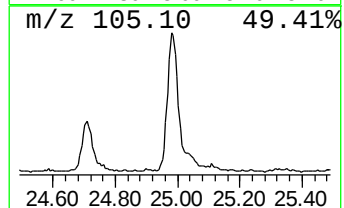
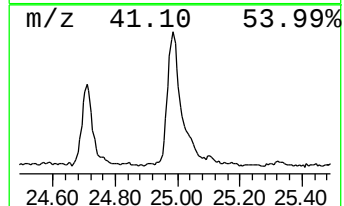
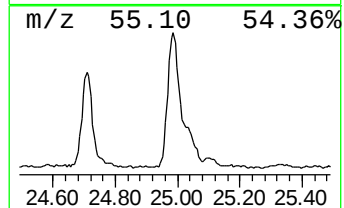
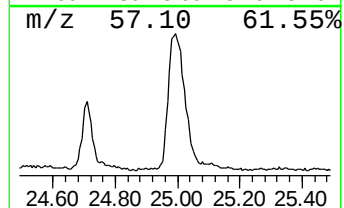
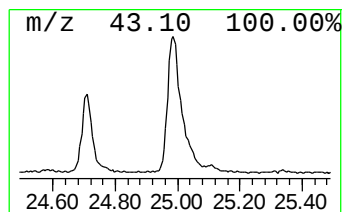
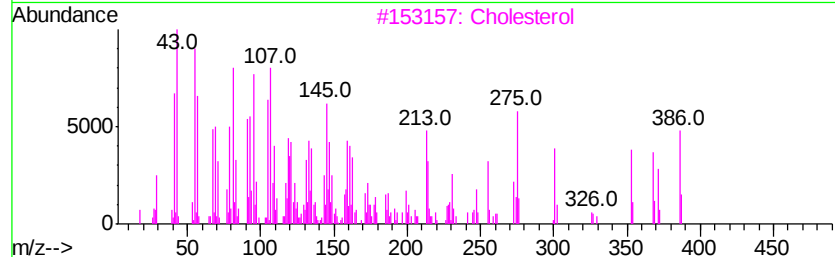
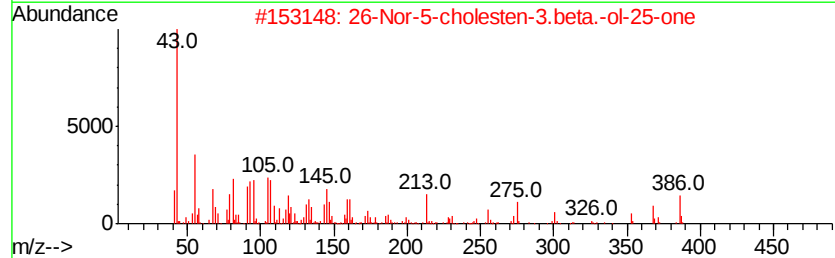
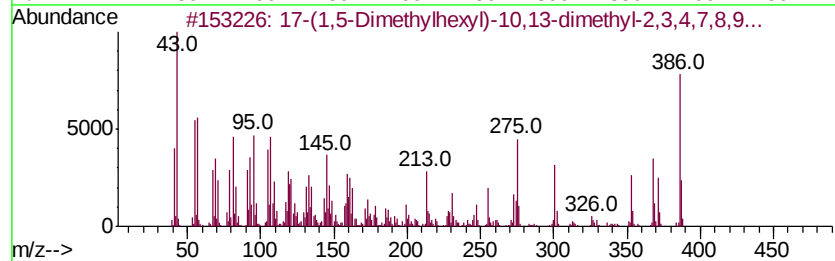
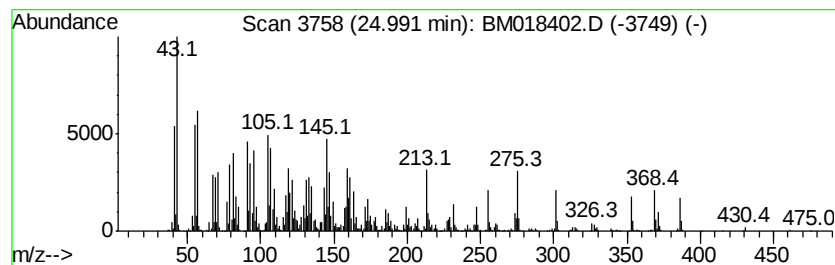
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 19 17-(1,5-Dimethylhexyl)-10,1... Concentration Rank 1

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.  |
|-------|----------|---------|------------------|-------|
| 24.99 | 21.61 ng | 7536690 | Perylene-d12     | 23.66 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|-------------------------------------|-----|----------|--------------|------|
| 1    | 5  | 17-(1,5-Dimethylhexyl)-10,13-dim... | 386 | C27H46O  | 1000210-38-4 | 99   |
| 2    |    | 26-Nor-5-cholesten-3.beta.-ol-25... | 386 | C26H42O2 | 007494-34-0  | 99   |
| 3    |    | Cholesterol                         | 386 | C27H46O  | 000057-88-5  | 96   |
| 4    |    | Cholestane-3,5-diol, 5-acetate, ... | 446 | C29H50O3 | 041721-93-1  | 64   |
| 5    |    | Cholesterol isocaproate             | 484 | C33H56O2 | 1000252-03-0 | 50   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_M\DATA\BM122718\  
 Data File : BM018402.D  
 Acq On : 27 Dec 2018 18:41  
 Operator : JU/SJ  
 Sample : J6569-03  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 381128115-4

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\8270-BM122718.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

| TIC Top Hit name     | RT    | EstConc | Units | Response | --Internal Standard-- |       |         |      |
|----------------------|-------|---------|-------|----------|-----------------------|-------|---------|------|
|                      |       |         |       |          | #                     | RT    | Resp    | Conc |
| 3-Ethyl-2-methyl-... | 6.13  | 2.7     | ng    | 251129   | 1                     | 7.78  | 1846900 | 20.0 |
| Cyclohexane, 1,2-... | 6.62  | 3.1     | ng    | 283103   | 1                     | 7.78  | 1846900 | 20.0 |
| Benzenemethanamin... | 8.17  | 2.1     | ng    | 191329   | 1                     | 7.78  | 1846900 | 20.0 |
| Benzenemethanol, ... | 8.88  | 3.8     | ng    | 353695   | 1                     | 7.78  | 1846900 | 20.0 |
| unknown11.25         | 11.25 | 2.6     | ng    | 374901   | 2                     | 10.57 | 2903310 | 20.0 |
| unknown13.32         | 13.32 | 5.0     | ng    | 1065700  | 3                     | 14.43 | 4303620 | 20.0 |
| 1-Decene             | 14.07 | 9.8     | ng    | 2112310  | 3                     | 14.43 | 4303620 | 20.0 |
| 2-Tetradecene, (E)-  | 15.97 | 8.1     | ng    | 2042740  | 4                     | 17.17 | 5054770 | 20.0 |
| 2-Propanol, 1-chl... | 16.93 | 3.5     | ng    | 879420   | 4                     | 17.17 | 5054770 | 20.0 |
| Caffeine             | 17.48 | 6.0     | ng    | 1502860  | 4                     | 17.17 | 5054770 | 20.0 |
| Cyclotetradecane     | 17.56 | 2.1     | ng    | 538820   | 4                     | 17.17 | 5054770 | 20.0 |
| n-Hexadecanoic acid  | 18.07 | 2.1     | ng    | 542855   | 4                     | 17.17 | 5054770 | 20.0 |
| Octadecanoic acid    | 19.36 | 2.0     | ng    | 673503   | 5                     | 21.37 | 6593600 | 20.0 |
| Butyl citrate        | 19.50 | 20.5    | ng    | 6764670  | 5                     | 21.37 | 6593600 | 20.0 |
| Ethanol, 2-butoxy... | 20.61 | 4.2     | ng    | 1389820  | 5                     | 21.37 | 6593600 | 20.0 |
| unknown22.23         | 22.23 | 7.7     | ng    | 2525030  | 5                     | 21.37 | 6593600 | 20.0 |
| Squalene             | 22.57 | 4.0     | ng    | 1402020  | 6                     | 23.66 | 6975740 | 20.0 |
| Cholestan-3.beta.-ol | 24.71 | 11.1    | ng    | 3854030  | 6                     | 23.66 | 6975740 | 20.0 |
| 17-(1,5-Dimethylh... | 24.99 | 21.6    | ng    | 7536690  | 6                     | 23.66 | 6975740 | 20.0 |