

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN060519\  
 Data File : BN006119.D  
 Acq On : 06 Jun 2019 03:14  
 Operator : JU/SJ  
 Sample : K3016-22  
 Misc :  
 ALS Vial : 28 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 A51D9

## Integration Parameters: LSCINT.P

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

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

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

Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN060419MA.M  
 Title : SVOA CALIBRATION

Signal : TIC

| peak<br># | R.T.<br>min | first<br>scan | max<br>scan | last<br>scan | PK<br>TY | peak<br>height | corr.<br>area | corr.<br>% max. | % of<br>total |
|-----------|-------------|---------------|-------------|--------------|----------|----------------|---------------|-----------------|---------------|
| 1         | 3.152       | 24            | 28          | 39           | rVB      | 58987          | 87248         | 1.68%           | 0.102%        |
| 2         | 3.664       | 111           | 115         | 122          | rVB      | 40373          | 55990         | 1.08%           | 0.065%        |
| 3         | 3.781       | 130           | 135         | 142          | rVB      | 97756          | 137593        | 2.65%           | 0.160%        |
| 4         | 4.770       | 298           | 303         | 312          | rVB      | 113096         | 155638        | 3.00%           | 0.182%        |
| 5         | 6.334       | 560           | 569         | 579          | rVB      | 248265         | 372101        | 7.16%           | 0.434%        |
| 6         | 6.640       | 616           | 621         | 625          | rVV5     | 19053          | 33352         | 0.64%           | 0.039%        |
| 7         | 6.817       | 640           | 651         | 664          | rBV      | 778703         | 1422434       | 27.38%          | 1.659%        |
| 8         | 6.981       | 671           | 679         | 691          | rBV      | 637867         | 1045191       | 20.12%          | 1.219%        |
| 9         | 7.087       | 691           | 697         | 703          | rVB      | 44486          | 75130         | 1.45%           | 0.088%        |
| 10        | 7.175       | 706           | 712         | 725          | rBV      | 808306         | 1316499       | 25.34%          | 1.536%        |
| 11        | 7.646       | 785           | 792         | 809          | rBV      | 1326377        | 2180048       | 41.96%          | 2.543%        |
| 12        | 7.864       | 823           | 829         | 836          | rVV      | 105425         | 159291        | 3.07%           | 0.186%        |
| 13        | 8.352       | 905           | 912         | 927          | rBV      | 711076         | 1213080       | 23.35%          | 1.415%        |
| 14        | 8.469       | 927           | 932         | 938          | rVV2     | 20457          | 39466         | 0.76%           | 0.046%        |
| 15        | 8.799       | 980           | 988         | 1001         | rBV      | 762120         | 1291890       | 24.87%          | 1.507%        |
| 16        | 9.222       | 1052          | 1060        | 1065         | rBV      | 26728          | 42087         | 0.81%           | 0.049%        |
| 17        | 9.516       | 1101          | 1110        | 1125         | rBV      | 598032         | 1057647       | 20.36%          | 1.234%        |
| 18        | 9.740       | 1143          | 1148        | 1160         | rVV10    | 27660          | 62063         | 1.19%           | 0.072%        |
| 19        | 10.052      | 1192          | 1201        | 1214         | rBV      | 920780         | 1605900       | 30.91%          | 1.873%        |
| 20        | 10.428      | 1258          | 1265        | 1273         | rBV      | 1986390        | 3317987       | 63.86%          | 3.870%        |
| 21        | 10.569      | 1279          | 1289        | 1302         | rBV      | 317843         | 652458        | 12.56%          | 0.761%        |
| 22        | 13.563      | 1792          | 1798        | 1803         | rBV      | 218764         | 305556        | 5.88%           | 0.356%        |
| 23        | 13.710      | 1817          | 1823        | 1827         | rVV      | 1785282        | 2320475       | 44.66%          | 2.707%        |
| 24        | 13.757      | 1827          | 1831        | 1841         | rVB      | 346952         | 492062        | 9.47%           | 0.574%        |
| 25        | 13.987      | 1860          | 1870        | 1887         | rBV      | 1861177        | 2891117       | 55.65%          | 3.372%        |
| 26        | 14.298      | 1914          | 1923        | 1931         | rBV2     | 2705396        | 4157555       | 80.02%          | 4.849%        |
| 27        | 14.363      | 1931          | 1934        | 1937         | rVV4     | 25454          | 40627         | 0.78%           | 0.047%        |
| 28        | 14.398      | 1937          | 1940        | 1945         | rVB3     | 33428          | 43176         | 0.83%           | 0.050%        |
| 29        | 14.510      | 1953          | 1959        | 1976         | rBV      | 372872         | 768675        | 14.80%          | 0.897%        |
| 30        | 14.663      | 1980          | 1985        | 1990         | rVV3     | 48128          | 108351        | 2.09%           | 0.126%        |
| 31        | 14.728      | 1993          | 1996        | 2001         | rVV5     | 34734          | 63656         | 1.23%           | 0.074%        |
| 32        | 15.298      | 2085          | 2093        | 2100         | rBV      | 2482404        | 3395680       | 65.36%          | 3.961%        |
| 33        | 15.428      | 2109          | 2115        | 2126         | rBV      | 472749         | 695898        | 13.39%          | 0.812%        |
| 34        | 15.540      | 2129          | 2134        | 2140         | rVV4     | 34762          | 54818         | 1.06%           | 0.064%        |

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN060519\  
 Data File : BN006119.D  
 Acq On : 06 Jun 2019 03:14  
 Operator : JU/SJ  
 Sample : K3016-22  
 Misc :  
 ALS Vial : 28 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 A51D9

## Integration Parameters: LSCINT.P

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

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

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

Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN060419MA.M  
 Title : SVOA CALIBRATION

|    |        |      |      |      |      |         |         |        |        |
|----|--------|------|------|------|------|---------|---------|--------|--------|
| 35 | 15.693 | 2155 | 2160 | 2168 | rVB2 | 47961   | 79640   | 1.53%  | 0.093% |
| 36 | 15.875 | 2186 | 2191 | 2196 | rBV5 | 19767   | 34210   | 0.66%  | 0.040% |
| 37 | 16.028 | 2214 | 2217 | 2222 | rBV6 | 24046   | 36132   | 0.70%  | 0.042% |
| 38 | 16.216 | 2236 | 2249 | 2253 | rBV2 | 112891  | 206662  | 3.98%  | 0.241% |
| 39 | 16.328 | 2263 | 2268 | 2272 | rBV2 | 71850   | 105823  | 2.04%  | 0.123% |
| 40 | 16.722 | 2330 | 2335 | 2340 | rBV7 | 26896   | 49105   | 0.95%  | 0.057% |
| 41 | 16.775 | 2340 | 2344 | 2348 | rBV5 | 35359   | 48030   | 0.92%  | 0.056% |
| 42 | 16.934 | 2366 | 2371 | 2377 | rBV2 | 57807   | 83783   | 1.61%  | 0.098% |
| 43 | 17.051 | 2382 | 2391 | 2396 | rBV2 | 3069876 | 4503377 | 86.68% | 5.253% |
| 44 | 17.092 | 2396 | 2398 | 2403 | rVV  | 322894  | 414647  | 7.98%  | 0.484% |
| 45 | 17.151 | 2403 | 2408 | 2421 | rVB  | 2203899 | 3274831 | 63.03% | 3.820% |
| 46 | 17.469 | 2458 | 2462 | 2471 | rVB6 | 24306   | 52423   | 1.01%  | 0.061% |
| 47 | 17.634 | 2488 | 2490 | 2497 | rVB6 | 18778   | 33164   | 0.64%  | 0.039% |
| 48 | 17.834 | 2521 | 2524 | 2531 | rVB7 | 17139   | 40210   | 0.77%  | 0.047% |
| 49 | 17.928 | 2535 | 2540 | 2542 | rBV  | 49342   | 72293   | 1.39%  | 0.084% |
| 50 | 17.957 | 2542 | 2545 | 2554 | rVB3 | 150065  | 322305  | 6.20%  | 0.376% |
| 51 | 18.057 | 2559 | 2562 | 2568 | rVB4 | 49947   | 88599   | 1.71%  | 0.103% |
| 52 | 18.122 | 2568 | 2573 | 2577 | rBV3 | 80330   | 143797  | 2.77%  | 0.168% |
| 53 | 18.163 | 2577 | 2580 | 2586 | rVB3 | 54381   | 84772   | 1.63%  | 0.099% |
| 54 | 18.228 | 2587 | 2591 | 2594 | rBV4 | 54209   | 81285   | 1.56%  | 0.095% |
| 55 | 18.292 | 2597 | 2602 | 2605 | rBV6 | 46339   | 69859   | 1.34%  | 0.081% |
| 56 | 18.404 | 2617 | 2621 | 2624 | rBV3 | 47573   | 73502   | 1.41%  | 0.086% |
| 57 | 18.481 | 2631 | 2634 | 2641 | rVB8 | 54631   | 97593   | 1.88%  | 0.114% |
| 58 | 18.551 | 2642 | 2646 | 2648 | rBV  | 74722   | 94004   | 1.81%  | 0.110% |
| 59 | 18.581 | 2648 | 2651 | 2655 | rVV  | 369235  | 450247  | 8.67%  | 0.525% |
| 60 | 18.681 | 2659 | 2668 | 2672 | rVB3 | 134359  | 256546  | 4.94%  | 0.299% |
| 61 | 18.728 | 2672 | 2676 | 2680 | rBV4 | 87446   | 126211  | 2.43%  | 0.147% |
| 62 | 18.857 | 2693 | 2698 | 2700 | rBV  | 78079   | 129429  | 2.49%  | 0.151% |
| 63 | 18.892 | 2700 | 2704 | 2708 | rVV  | 882518  | 1164582 | 22.42% | 1.358% |
| 64 | 18.928 | 2708 | 2710 | 2718 | rVB2 | 218002  | 265811  | 5.12%  | 0.310% |
| 65 | 19.016 | 2719 | 2725 | 2728 | rBV5 | 62741   | 135188  | 2.60%  | 0.158% |
| 66 | 19.086 | 2734 | 2737 | 2738 | rBV3 | 42103   | 42177   | 0.81%  | 0.049% |
| 67 | 19.122 | 2738 | 2743 | 2746 | rVV  | 620693  | 891943  | 17.17% | 1.040% |
| 68 | 19.151 | 2746 | 2748 | 2753 | rVB  | 533699  | 597721  | 11.50% | 0.697% |
| 69 | 19.216 | 2754 | 2759 | 2762 | rVV  | 391515  | 502191  | 9.67%  | 0.586% |
| 70 | 19.269 | 2762 | 2768 | 2775 | rVB  | 1575239 | 2035848 | 39.18% | 2.375% |
| 71 | 19.398 | 2787 | 2790 | 2794 | rVB6 | 87904   | 128930  | 2.48%  | 0.150% |

Data Path : Z:\SVOASRV\HPCHEM1\BNA\_N\DATA\BN060519\  
Data File : BN006119.D  
Acq On : 06 Jun 2019 03:14  
Operator : JU/SJ  
Sample : K3016-22  
Misc :  
ALS Vial : 28 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
A51D9

## Integration Parameters: LSCINT.P

Integrator: RTE  
Smoothing : OFF  
Sampling : 1  
Start Thrs: 0.2  
Stop Thrs : 0  
Filtering: 5  
Min Area: 0.1 % of largest Peak  
Max Peaks: 100  
Peak Location: TOP

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

Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN060419MA.M  
Title : SVOA CALIBRATION

|     |        |      |      |      |      |         |         |         |        |
|-----|--------|------|------|------|------|---------|---------|---------|--------|
| 72  | 19.457 | 2794 | 2800 | 2809 | rBV2 | 3040664 | 5195505 | 100.00% | 6.060% |
| 73  | 19.557 | 2815 | 2817 | 2823 | rVB4 | 62508   | 94140   | 1.81%   | 0.110% |
| 74  | 19.645 | 2829 | 2832 | 2835 | rBV  | 76606   | 89019   | 1.71%   | 0.104% |
| 75  | 19.716 | 2837 | 2844 | 2849 | rVV2 | 462843  | 866407  | 16.68%  | 1.011% |
| 76  | 19.804 | 2855 | 2859 | 2863 | rBV3 | 183249  | 275108  | 5.30%   | 0.321% |
| 77  | 19.975 | 2880 | 2888 | 2893 | rBV4 | 200435  | 459648  | 8.85%   | 0.536% |
| 78  | 20.022 | 2893 | 2896 | 2900 | rVB3 | 135818  | 162686  | 3.13%   | 0.190% |
| 79  | 20.069 | 2900 | 2904 | 2907 | rBV  | 224183  | 325882  | 6.27%   | 0.380% |
| 80  | 20.133 | 2911 | 2915 | 2918 | rBV5 | 126858  | 208213  | 4.01%   | 0.243% |
| 81  | 20.304 | 2939 | 2944 | 2955 | rVB4 | 1032139 | 1574626 | 30.31%  | 1.837% |
| 82  | 20.451 | 2964 | 2969 | 2974 | rBV  | 2154055 | 2656612 | 51.13%  | 3.099% |
| 83  | 20.675 | 3003 | 3007 | 3014 | rBV5 | 187257  | 408733  | 7.87%   | 0.477% |
| 84  | 20.833 | 3030 | 3034 | 3040 | rBV6 | 149157  | 248020  | 4.77%   | 0.289% |
| 85  | 20.957 | 3052 | 3055 | 3058 | rBV2 | 500357  | 587702  | 11.31%  | 0.685% |
| 86  | 20.992 | 3058 | 3061 | 3069 | rVB2 | 508689  | 753668  | 14.51%  | 0.879% |
| 87  | 21.192 | 3091 | 3095 | 3099 | rBV  | 605501  | 667555  | 12.85%  | 0.779% |
| 88  | 21.269 | 3102 | 3108 | 3113 | rBV  | 3473781 | 5080160 | 97.78%  | 5.925% |
| 89  | 21.427 | 3131 | 3135 | 3142 | rBV2 | 273181  | 380044  | 7.31%   | 0.443% |
| 90  | 21.592 | 3159 | 3163 | 3168 | rBV  | 965619  | 1085624 | 20.90%  | 1.266% |
| 91  | 21.869 | 3206 | 3210 | 3213 | rBV2 | 292971  | 386100  | 7.43%   | 0.450% |
| 92  | 22.327 | 3284 | 3288 | 3295 | rVB  | 332092  | 468891  | 9.02%   | 0.547% |
| 93  | 22.833 | 3369 | 3374 | 3380 | rBV3 | 793782  | 1361819 | 26.21%  | 1.588% |
| 94  | 23.363 | 3458 | 3464 | 3469 | rBV2 | 1615840 | 3138313 | 60.40%  | 3.661% |
| 95  | 23.504 | 3481 | 3488 | 3507 | rVB  | 2414208 | 4758988 | 91.60%  | 5.551% |
| 96  | 24.033 | 3573 | 3578 | 3590 | rVB  | 615401  | 1130345 | 21.76%  | 1.318% |
| 97  | 24.774 | 3699 | 3704 | 3711 | rVB  | 325765  | 646670  | 12.45%  | 0.754% |
| 98  | 25.639 | 3845 | 3851 | 3858 | rBV2 | 208509  | 495008  | 9.53%   | 0.577% |
| 99  | 26.539 | 3997 | 4004 | 4016 | rBV5 | 320462  | 1127072 | 21.69%  | 1.315% |
| 100 | 28.074 | 4257 | 4265 | 4287 | rVB2 | 751580  | 2623963 | 50.50%  | 3.061% |

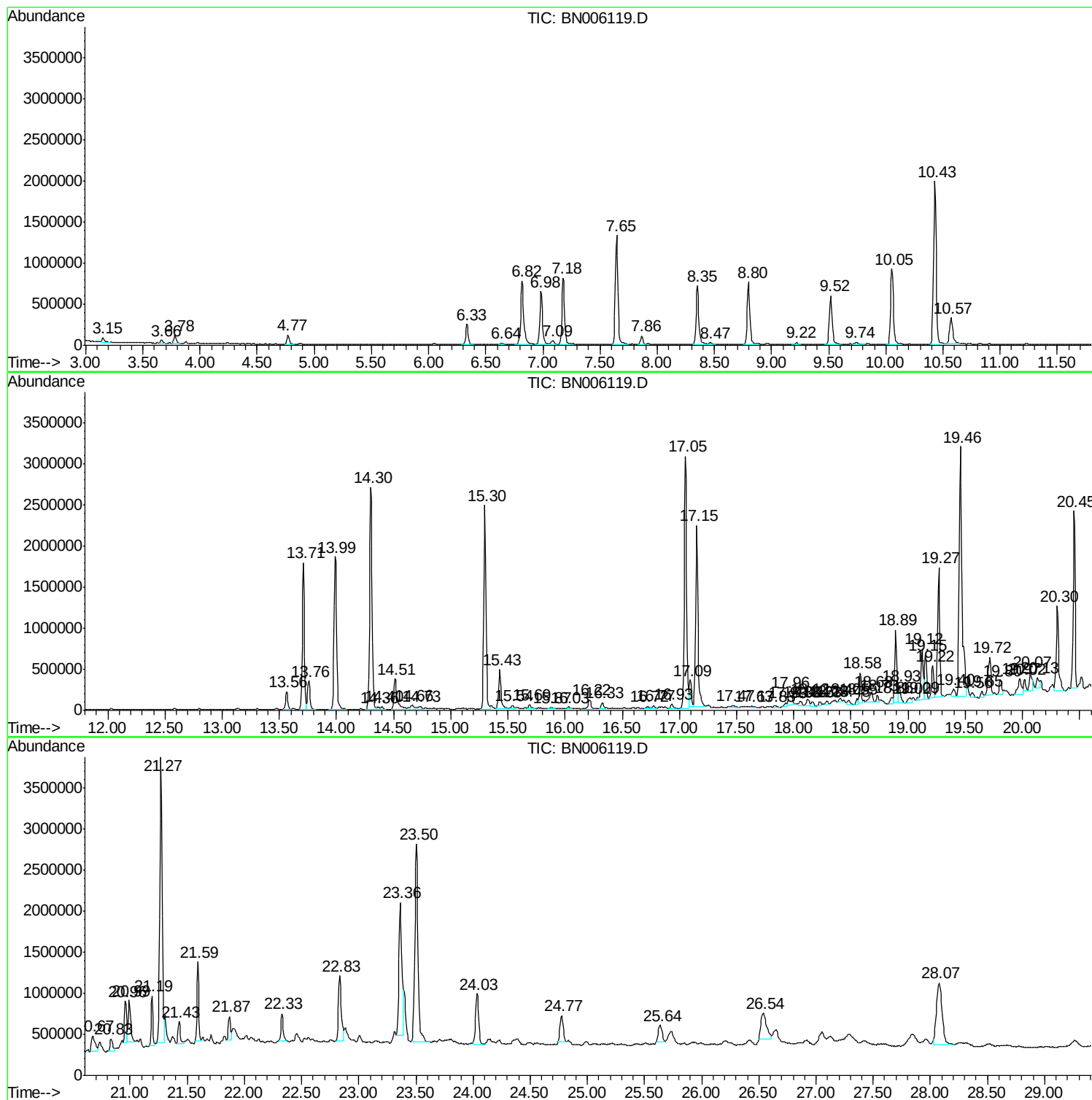
Sum of corrected areas: 85734130

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN060519\  
Data File : BN006119.D  
Acq On : 06 Jun 2019 03:14  
Operator : JU/SJ  
Sample : K3016-22  
Misc :  
ALS Vial : 28 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampled :  
A51D9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN060419MA.M  
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN060519\  
Data File : BN006119.D  
Acq On : 06 Jun 2019 03:14  
Operator : JU/SJ  
Sample : K3016-22  
Misc :  
ALS Vial : 28 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampled :  
A51D9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN060419MA.M  
Quant Title : SVOA CALIBRATION

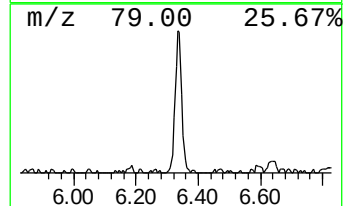
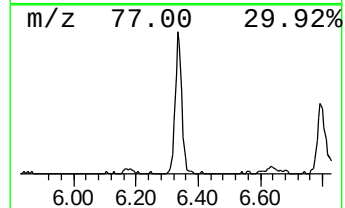
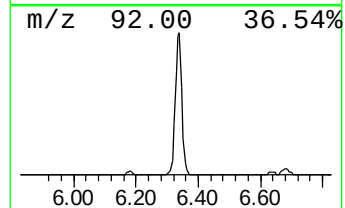
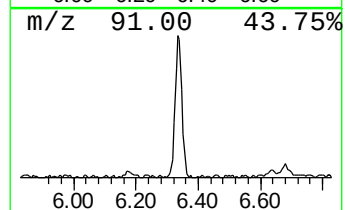
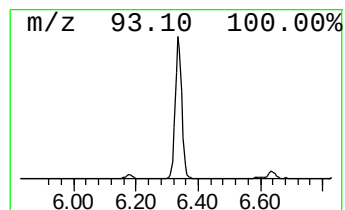
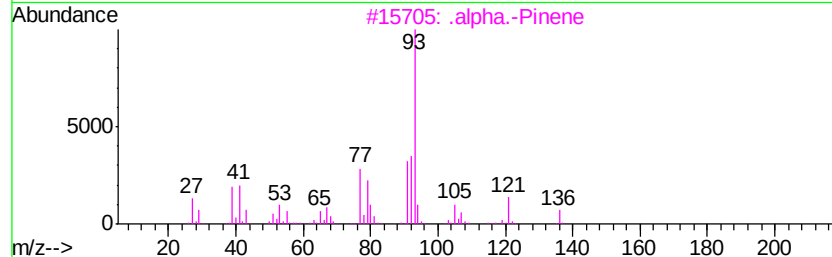
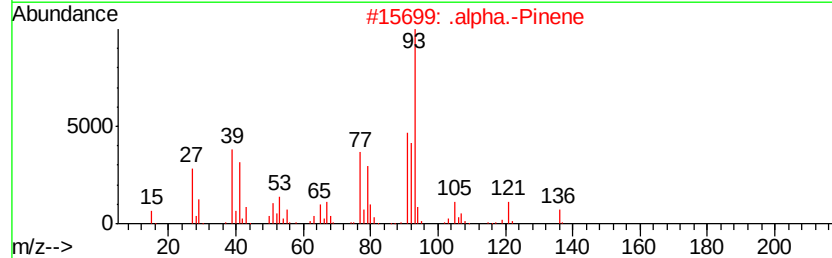
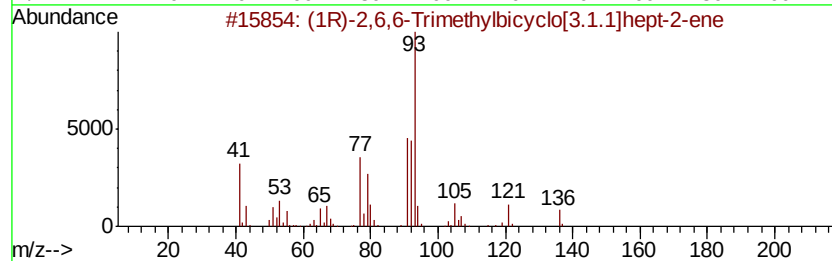
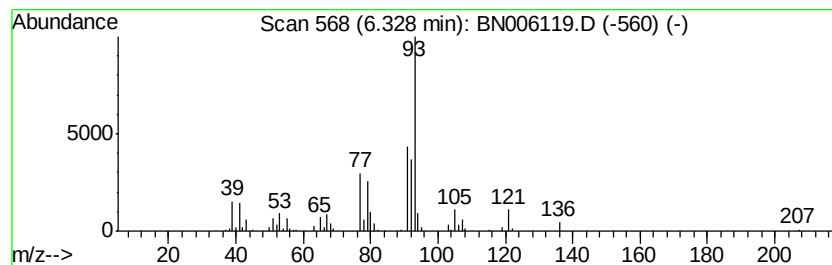
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 1 (1R)-2,6,6-Trimethylbicyclo... Concentration Rank 9

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 6.33 | 3.41 ng/ul | 372101 | 1,4-Dichlorobenzene-d4 | 7.65 |

| Hit# | of | Tentative ID                                 | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------------------------|-----|---------|-------------|------|
| 1    | 5  | (1R)-2,6,6-Trimethylbicyclo[3.1.1]hept-2-ene | 136 | C10H16  | 007785-70-8 | 95   |
| 2    |    | .alpha.-Pinene                               | 136 | C10H16  | 000080-56-8 | 95   |
| 3    |    | .alpha.-Pinene                               | 136 | C10H16  | 000080-56-8 | 95   |
| 4    |    | (1R)-2,6,6-Trimethylbicyclo[3.1.1]hept-2-ene | 136 | C10H16  | 007785-70-8 | 94   |
| 5    |    | Bicyclo[3.1.1]hept-2-ene, 3,6,6-trimethyl-   | 136 | C10H16  | 004889-83-2 | 94   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN060519\  
 Data File : BN006119.D  
 Acq On : 06 Jun 2019 03:14  
 Operator : JU/SJ  
 Sample : K3016-22  
 Misc :  
 ALS Vial : 28 Sample Multiplier: 1

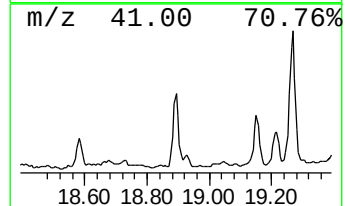
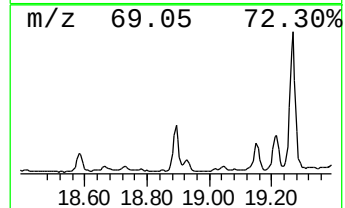
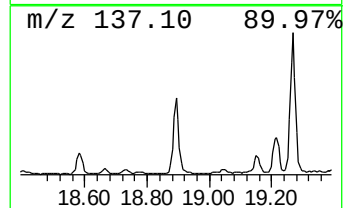
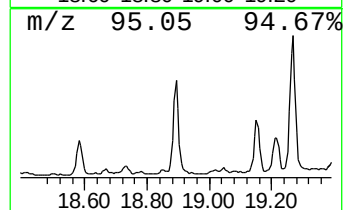
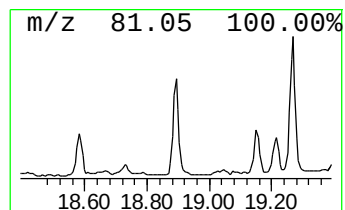
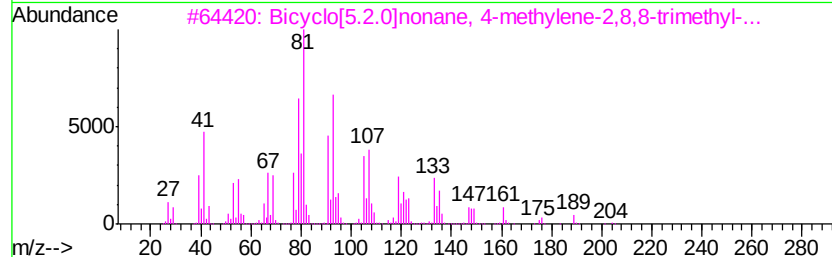
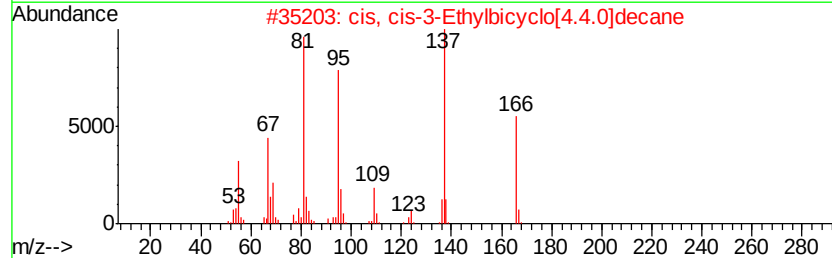
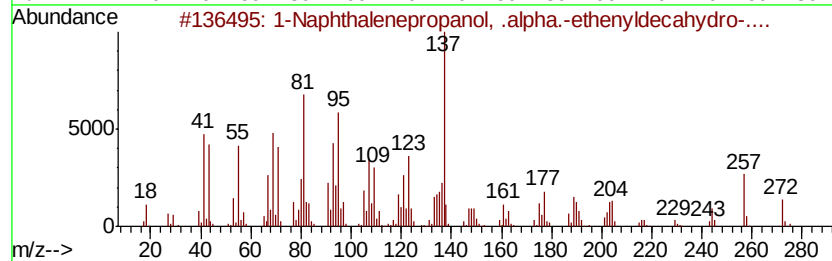
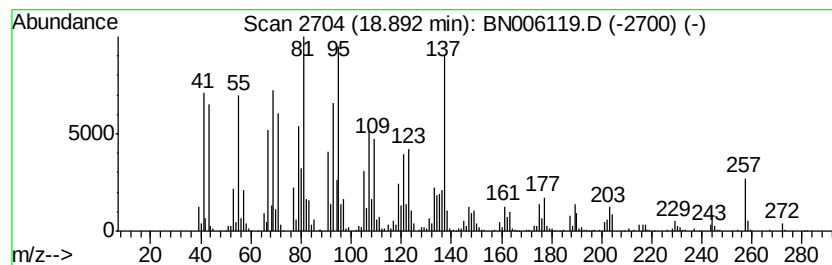
Instrument :  
 BNA\_N  
 ClientSampleId :  
 A51D9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN060419MA.M  
 Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
 Peak Number 2 1-Naphthalenepropanol, .alp... Concentration Rank 5

| R.T.    | EstConc                             | Area         | Relative to ISTD | R.T.      |
|---------|-------------------------------------|--------------|------------------|-----------|
| 18.89   | 5.17 ng/ul                          | 1164580      | Phenanthrene-d10 | 17.05     |
| Hit# of | 5                                   | Tentative ID | MW MolForm       | CAS# Qual |
| 1       | 1-Naphthalenepropanol, .alpha.-e... | 290 C20H34O  | 001438-62-6      | 64        |
| 2       | cis, cis-3-Ethylbicyclo[4.4.0]de... | 166 C12H22   | 066660-42-2      | 38        |
| 3       | Bicyclo[5.2.0]nonane, 4-methylen... | 204 C15H24   | 1000159-38-2     | 30        |
| 4       | 3-Hydroxypyrindine monoacetate      | 137 C7H7NO2  | 017747-43-2      | 25        |
| 5       | p-Menth-8-en-3-ol, acetate          | 196 C12H20O2 | 000089-49-6      | 25        |



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN060519\  
Data File : BN006119.D  
Acq On : 06 Jun 2019 03:14  
Operator : JU/SJ  
Sample : K3016-22  
Misc :  
ALS Vial : 28 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
A51D9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN060419MA.M  
Quant Title : SVOA CALIBRATION

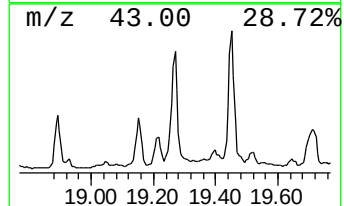
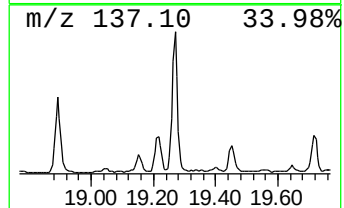
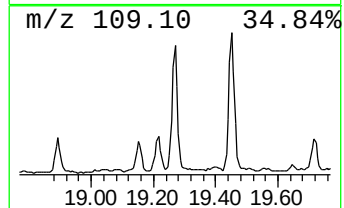
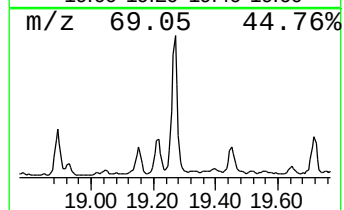
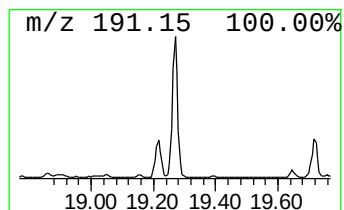
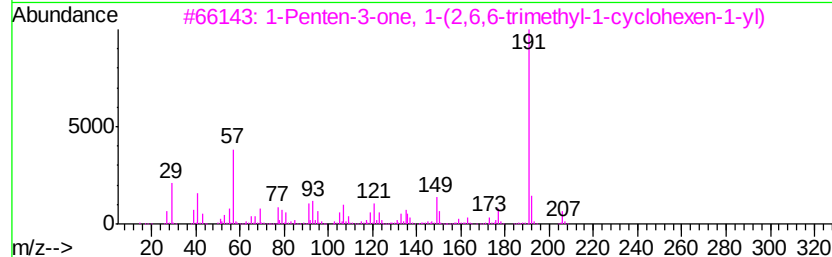
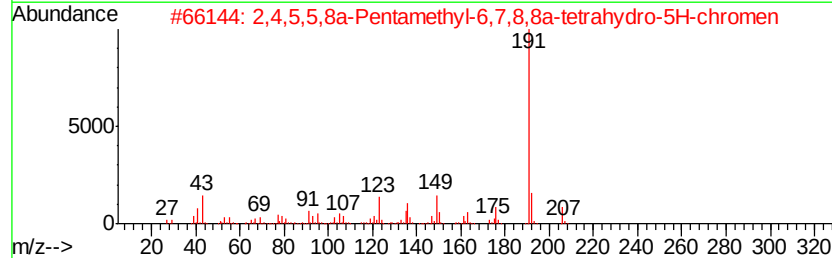
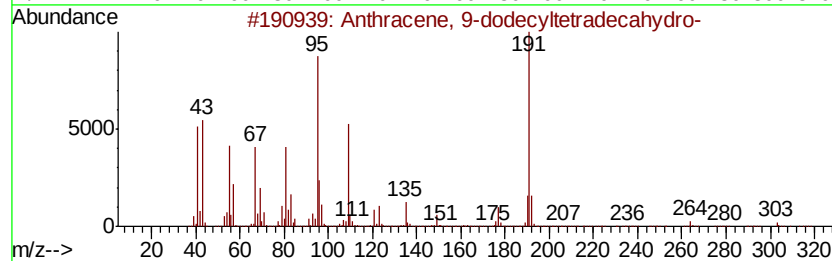
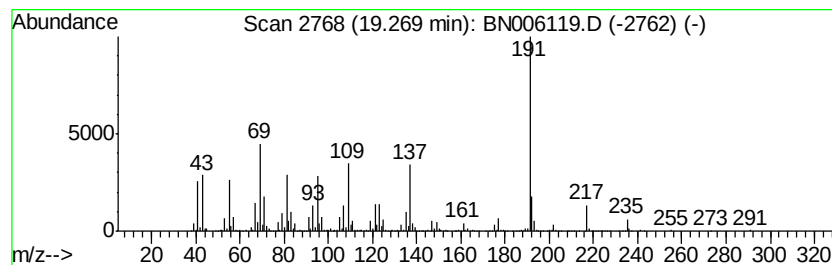
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T.  |
|-------|------------|---------|------------------|-------|
| 19.27 | 8.01 ng/ul | 2035850 | Chrysene-d12     | 21.27 |

| Hit# | of | Tentative ID                         | MW  | MolForm    | CAS#         | Qual |
|------|----|--------------------------------------|-----|------------|--------------|------|
| 1    | 5  | Anthracene, 9-dodecyltetradeca...    | 360 | C26H48     | 055401-75-7  | 45   |
| 2    |    | 2,4,5,5,8a-Pentamethyl-6,7,8,8a-...  | 206 | C14H22O    | 1000195-40-9 | 38   |
| 3    |    | 1-Penten-3-one, 1-(2,6,6-trimeth...  | 206 | C14H22O    | 000127-43-5  | 38   |
| 4    |    | Anthracene, 9-butyl-                 | 234 | C18H18     | 001498-69-7  | 35   |
| 5    |    | 4-Fluoro-2-trifluoromethylbenzoic... | 376 | C20H28F4O2 | 1000338-49-2 | 35   |





Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN060519\  
Data File : BN006119.D  
Acq On : 06 Jun 2019 03:14  
Operator : JU/SJ  
Sample : K3016-22  
Misc :  
ALS Vial : 28 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
A51D9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN060419MA.M  
Quant Title : SVOA CALIBRATION

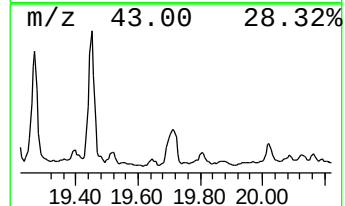
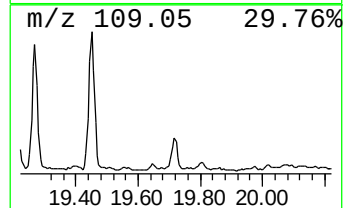
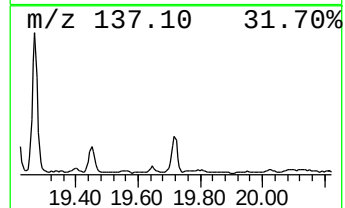
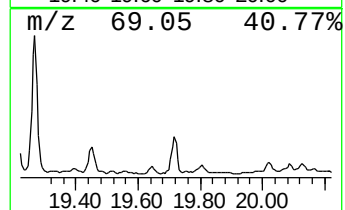
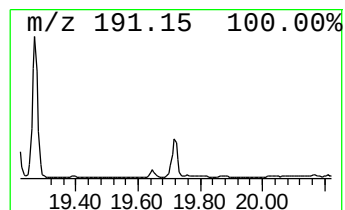
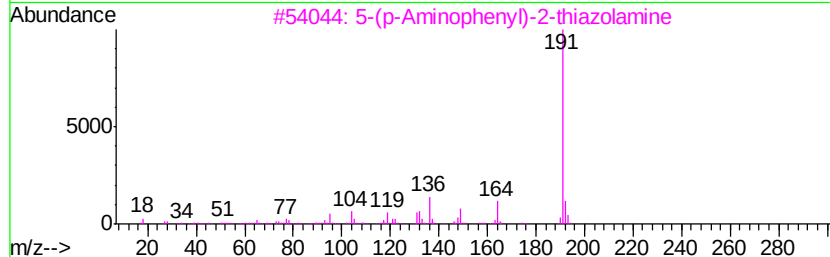
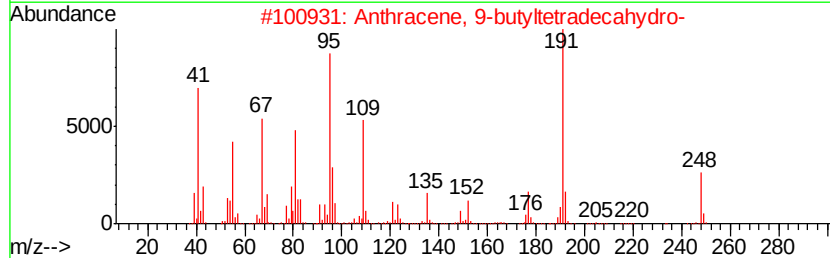
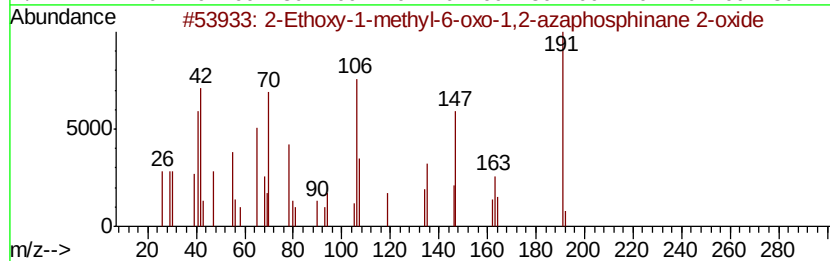
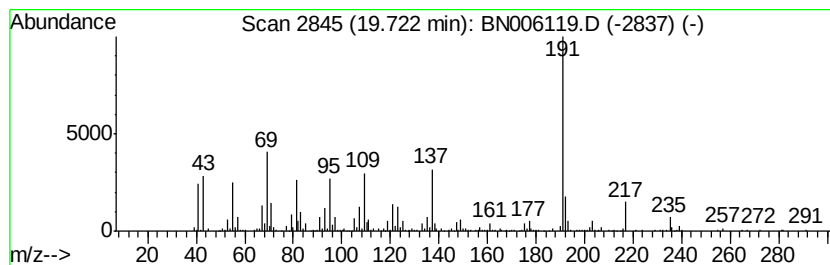
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 4 unknown-02 Concentration Rank 10

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 19.72 | 3.41 ng/ul | 866407 | Chrysene-d12     | 21.27 |

| Hit# | of | Tentative ID                        | MW  | MolForm     | CAS#        | Qual |
|------|----|-------------------------------------|-----|-------------|-------------|------|
| 1    | 5  | 2-Ethoxy-1-methyl-6-oxo-1,2-azap... | 191 | C7H14N03P   | 093989-00-5 | 45   |
| 2    |    | Anthracene, 9-butyltetradecahydro-  | 248 | C18H32      | 055133-89-6 | 45   |
| 3    |    | 5-(p-Aminophenyl)-2-thiazolamine    | 191 | C9H9N3S     | 090349-87-4 | 38   |
| 4    |    | Pyrido[3,2-d]pyrimidine-2,4(1H,3... | 191 | C9H9N3O2    | 024410-25-1 | 38   |
| 5    |    | 4,5-Dihydrothiazol-4-one, 2-(2-m... | 222 | C10H10N2O2S | 022476-61-5 | 38   |





Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN060519\  
Data File : BN006119.D  
Acq On : 06 Jun 2019 03:14  
Operator : JU/SJ  
Sample : K3016-22  
Misc :  
ALS Vial : 28 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleID :  
A51D9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN060419MA.M  
Quant Title : SVOA CALIBRATION

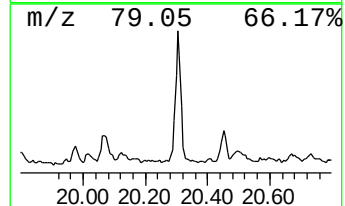
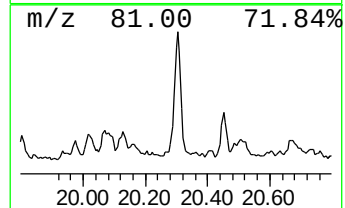
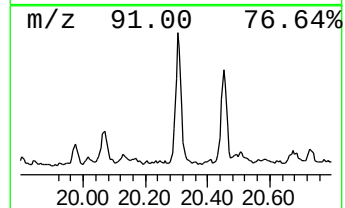
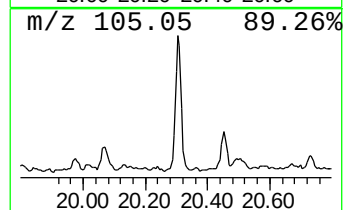
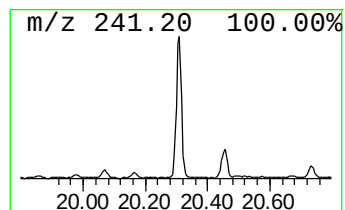
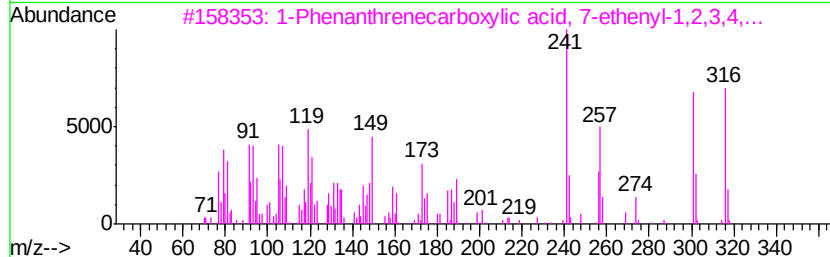
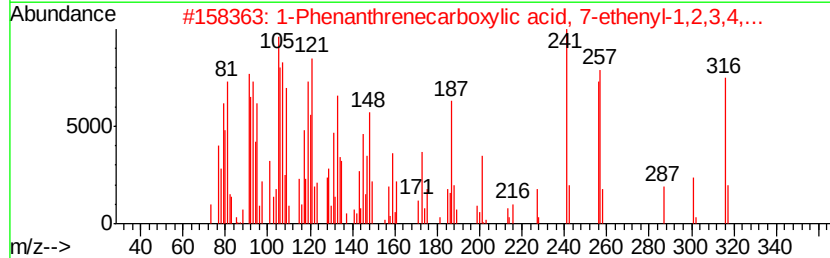
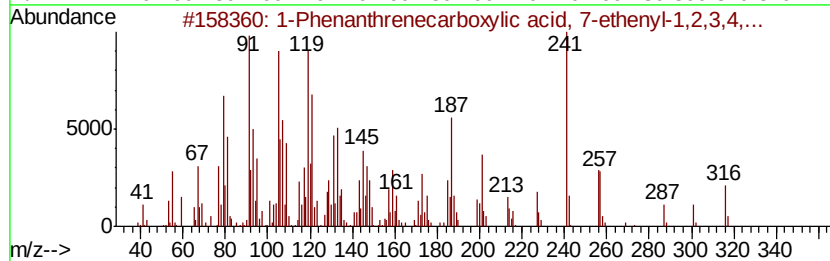
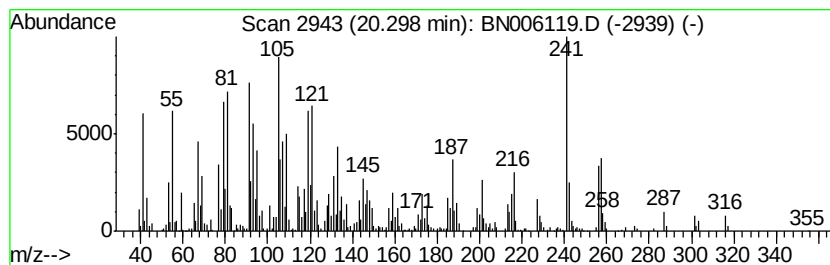
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 5 1-Phenanthrenecarboxylic ac... Concentration Rank 4

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T.  |
|-------|------------|---------|------------------|-------|
| 20.30 | 6.20 ng/ul | 1574630 | Chrysene-d12     | 21.27 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | 1-Phenanthrenecarboxylic acid, 7... | 316 | C21H32O2 | 001686-62-0 | 96   |
| 2    |    | 1-Phenanthrenecarboxylic acid, 7... | 316 | C21H32O2 | 001686-62-0 | 90   |
| 3    |    | 1-Phenanthrenecarboxylic acid, 7... | 316 | C21H32O2 | 003582-26-1 | 62   |
| 4    |    | s-Indacen-1(2H)-one, 3,5,6,7-tet... | 256 | C18H24O  | 038754-94-8 | 42   |
| 5    |    | Bisphenol C                         | 256 | C17H20O2 | 000079-97-0 | 38   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN060519\  
Data File : BN006119.D  
Acq On : 06 Jun 2019 03:14  
Operator : JU/SJ  
Sample : K3016-22  
Misc :  
ALS Vial : 28 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
A51D9

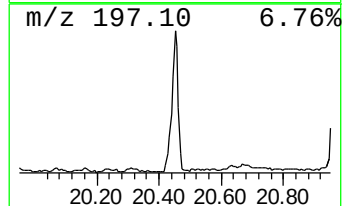
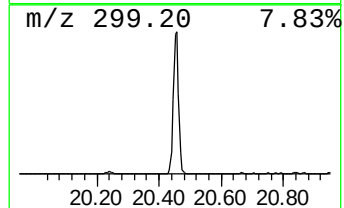
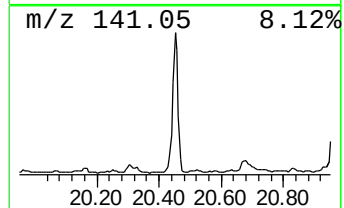
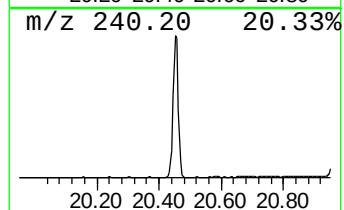
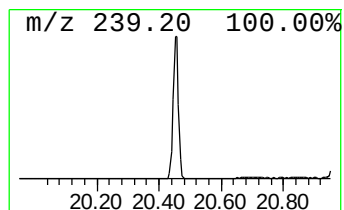
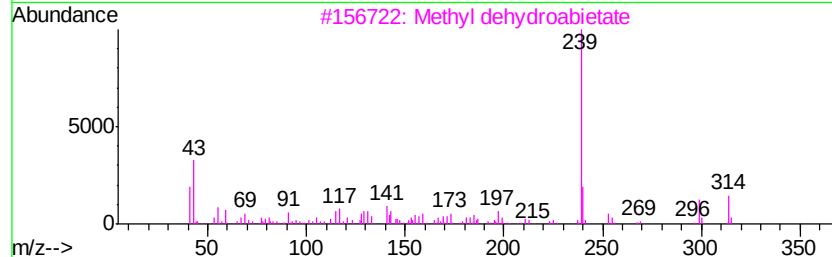
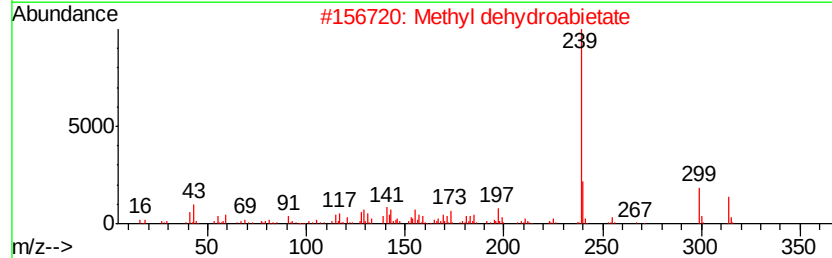
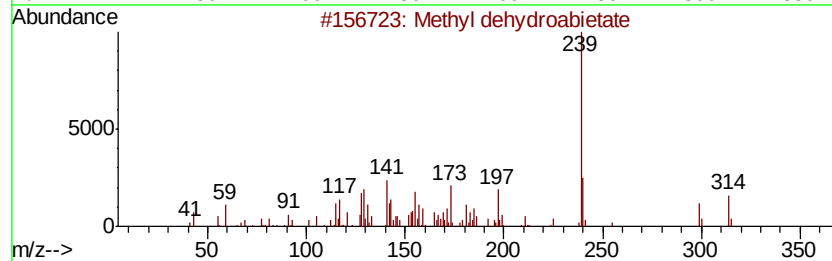
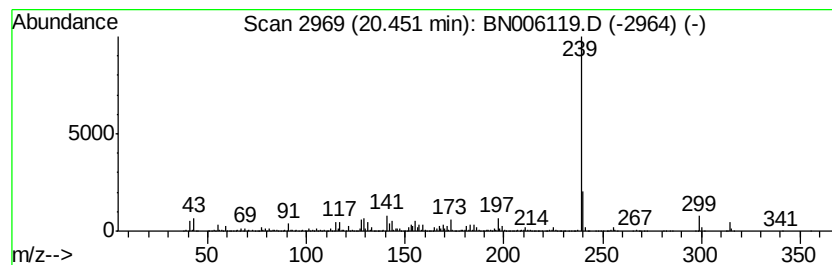
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN060419MA.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 6 Methyl dehydroabietate Concentration Rank 2

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 20.45 | 10.46 ng/ul | 2656610 | Chrysene-d12     | 21.27 |

| Hit# | of | Tentative ID                          | MW  | MolForm    | CAS#         | Qual |
|------|----|---------------------------------------|-----|------------|--------------|------|
| 1    | 5  | Methyl dehydroabietate                | 314 | C21H30O2   | 001235-74-1  | 92   |
| 2    |    | Methyl dehydroabietate                | 314 | C21H30O2   | 001235-74-1  | 91   |
| 3    |    | Methyl dehydroabietate                | 314 | C21H30O2   | 001235-74-1  | 78   |
| 4    |    | 1-[4-(4-Chlorophenyl)thioxomethyl]... | 239 | C12H14ClNS | 003335-29-3  | 59   |
| 5    |    | Phthalic acid, di(3-methylphenyl)...  | 346 | C22H18O4   | 1000315-37-3 | 59   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN060519\  
Data File : BN006119.D  
Acq On : 06 Jun 2019 03:14  
Operator : JU/SJ  
Sample : K3016-22  
Misc :  
ALS Vial : 28 Sample Multiplier: 1

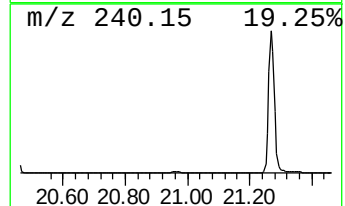
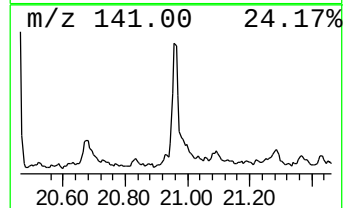
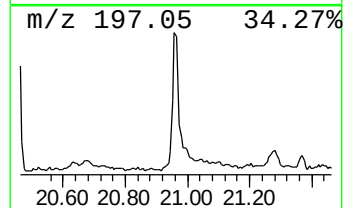
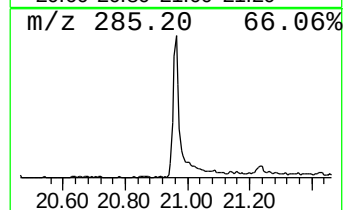
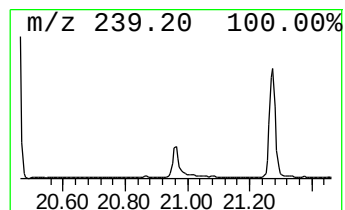
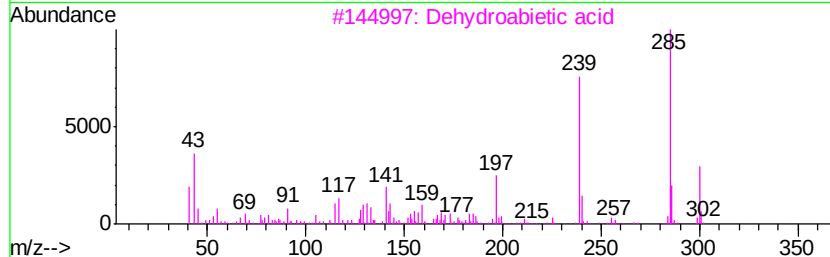
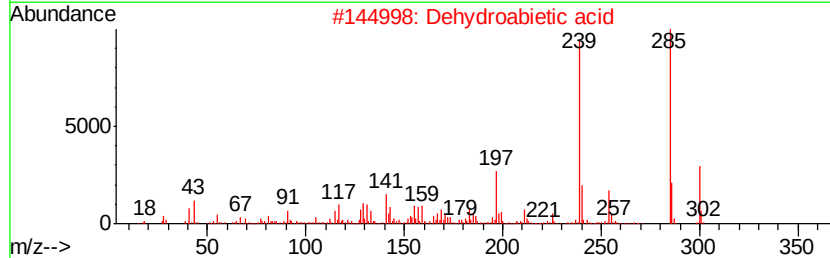
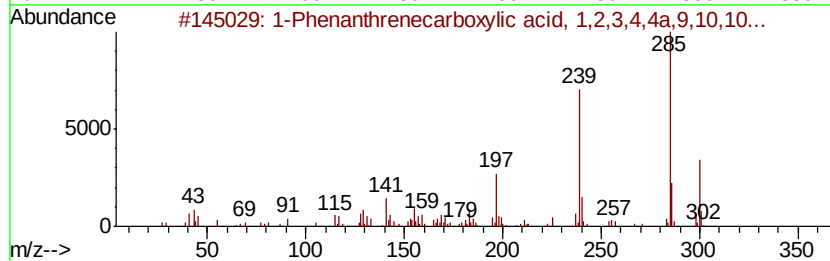
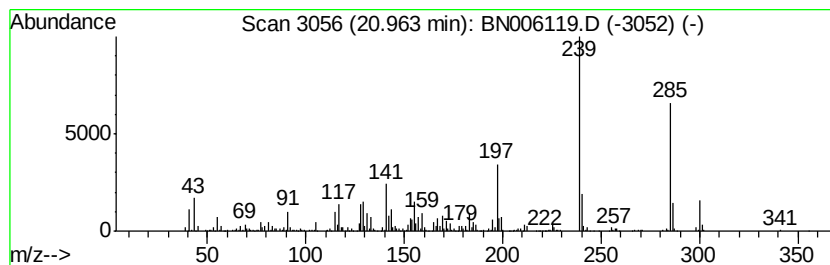
Instrument :  
BNA\_N  
ClientSampleId :  
A51D9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN060419MA.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 7 1-Phenanthrenecarboxylic ac... Concentration Rank 13

| R.T.    | EstConc    | Area                                | Relative to ISTD | R.T.        |             |      |
|---------|------------|-------------------------------------|------------------|-------------|-------------|------|
| 20.96   | 2.31 ng/ul | 587702                              | Chrysene-d12     | 21.27       |             |      |
| Hit# of | 5          | Tentative ID                        | MW               | MolForm     | CAS#        | Qual |
| 1       |            | 1-Phenanthrenecarboxylic acid, 1... | 300              | C20H28O2    | 005155-70-4 | 95   |
| 2       |            | Dehydroabietic acid                 | 300              | C20H28O2    | 001740-19-8 | 94   |
| 3       |            | Dehydroabietic acid                 | 300              | C20H28O2    | 001740-19-8 | 91   |
| 4       |            | Dehydroabietic acid                 | 300              | C20H28O2    | 001740-19-8 | 90   |
| 5       |            | Ethyl 4-hydroxy-7-trifluoromethy... | 285              | C13H10F3NO3 | 000391-02-6 | 53   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN060519\  
Data File : BN006119.D  
Acq On : 06 Jun 2019 03:14  
Operator : JU/SJ  
Sample : K3016-22  
Misc :  
ALS Vial : 28 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
A51D9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN060419MA.M  
Quant Title : SVOA CALIBRATION

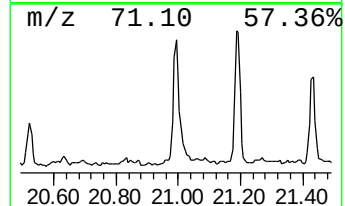
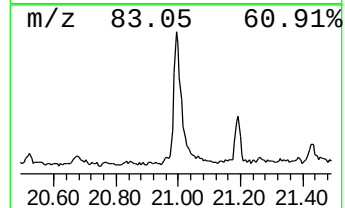
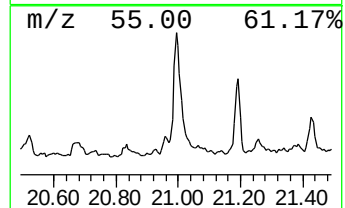
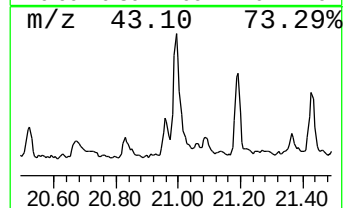
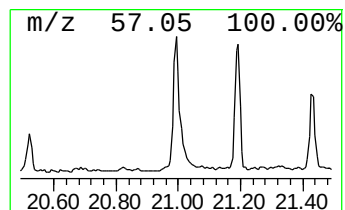
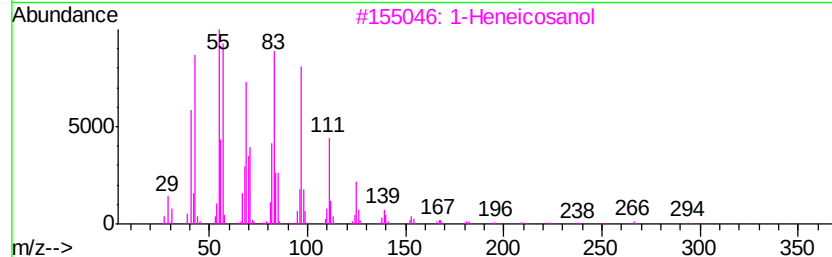
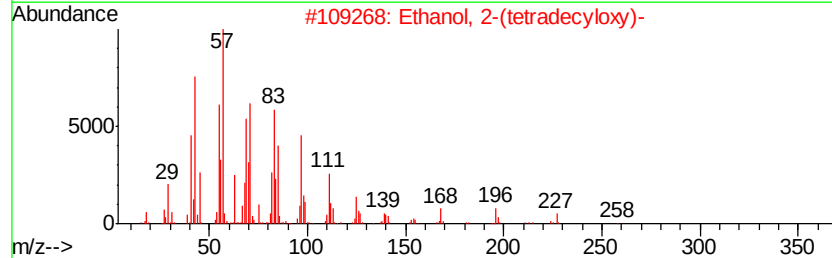
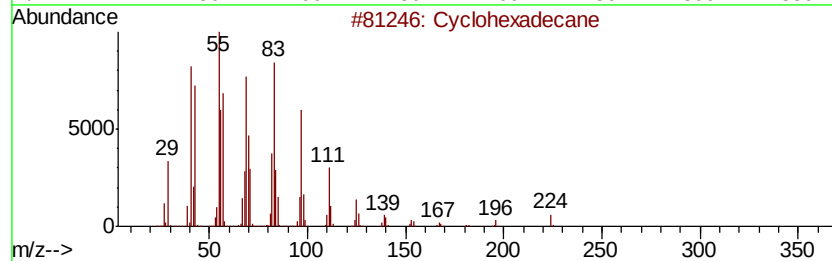
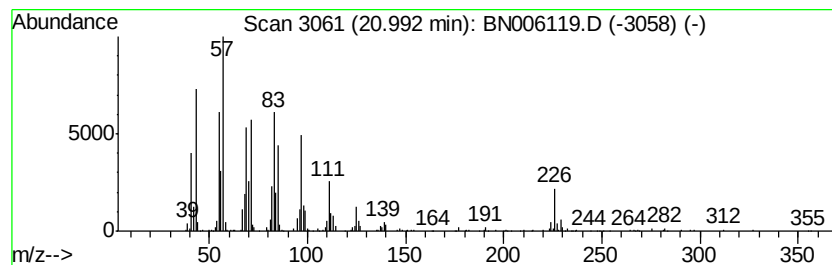
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 8 (DEL) Alkane: Cyclic20.99 Concentration Rank 11

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 20.99 | 2.97 ng/ul | 753668 | Chrysene-d12     | 21.27 |

| Hit# | of | Tentative ID                | MW  | MolForm  | CAS#        | Qual |
|------|----|-----------------------------|-----|----------|-------------|------|
| 1    | 5  | Cyclohexadecane             | 224 | C16H32   | 000295-65-8 | 92   |
| 2    |    | Ethanol, 2-(tetradecyloxy)- | 258 | C16H34O2 | 002136-70-1 | 90   |
| 3    |    | 1-Heneicosanol              | 312 | C21H44O  | 015594-90-8 | 86   |
| 4    |    | Cyclopentadecane            | 210 | C15H30   | 000295-48-7 | 78   |
| 5    |    | Behenic alcohol             | 326 | C22H46O  | 000661-19-8 | 70   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN060519\  
Data File : BN006119.D  
Acq On : 06 Jun 2019 03:14  
Operator : JU/SJ  
Sample : K3016-22  
Misc :  
ALS Vial : 28 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
A51D9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN060419MA.M  
Quant Title : SVOA CALIBRATION

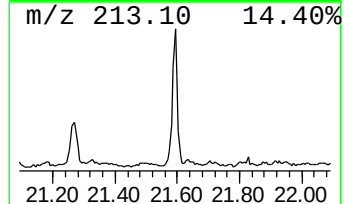
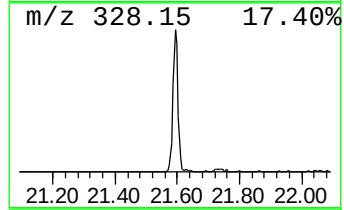
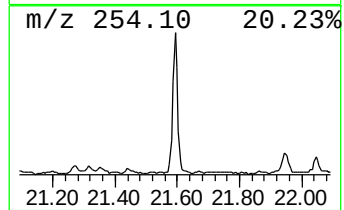
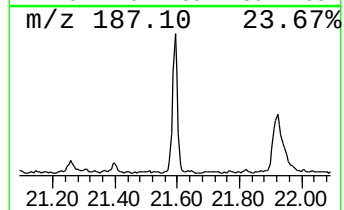
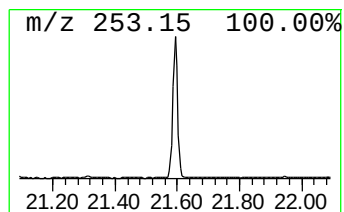
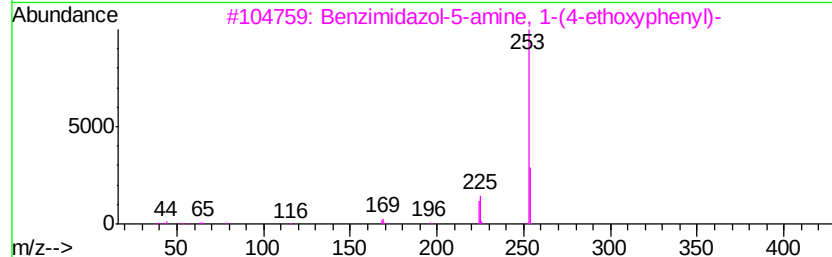
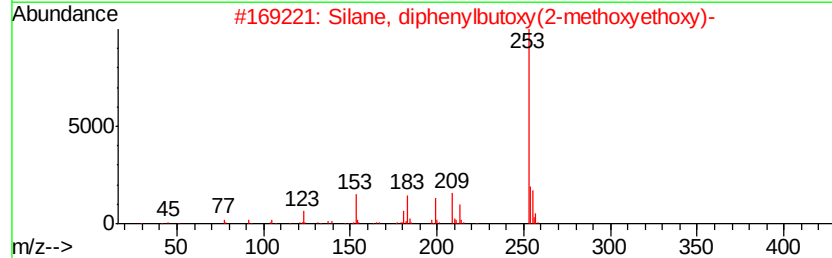
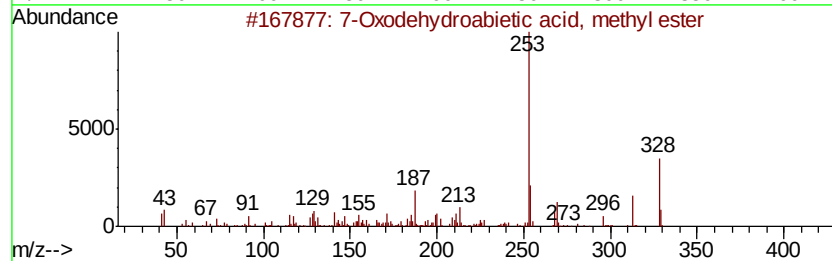
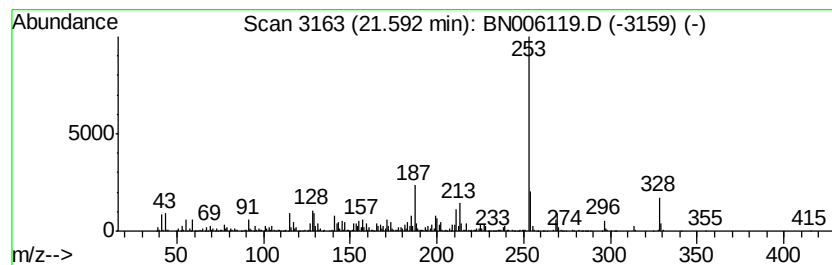
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 9 7-Oxodehydroabiatic acid, m... Concentration Rank 8

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T.  |
|-------|------------|---------|------------------|-------|
| 21.59 | 4.27 ng/ul | 1085620 | Chrysene-d12     | 21.27 |

| Hit# | of | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|-------------------------------------|-----|------------|--------------|------|
| 1    | 5  | 7-Oxodehydroabiatic acid, methyl... | 328 | C21H28O3   | 110936-78-2  | 94   |
| 2    |    | Silane, diphenylbutoxy(2-methoxy... | 330 | C19H26O3Si | 1000367-72-3 | 50   |
| 3    |    | Benzimidazol-5-amine, 1-(4-ethox... | 253 | C15H15N3O  | 007104-62-3  | 46   |
| 4    |    | Imidazole, 2-iodo-5-methyl-4-nitro- | 253 | C4H4IN3O2  | 149510-86-1  | 45   |
| 5    |    | Isophthalic acid, di(2-ethylphen... | 374 | C24H22O4   | 1000356-76-7 | 43   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN060519\  
Data File : BN006119.D  
Acq On : 06 Jun 2019 03:14  
Operator : JU/SJ  
Sample : K3016-22  
Misc :  
ALS Vial : 28 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
A51D9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN060419MA.M  
Quant Title : SVOA CALIBRATION

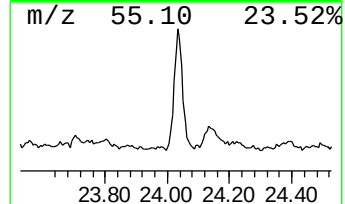
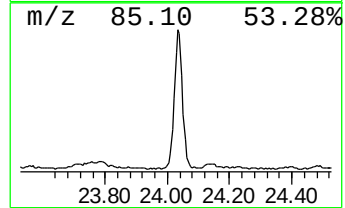
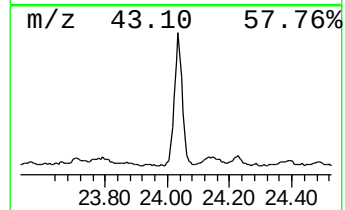
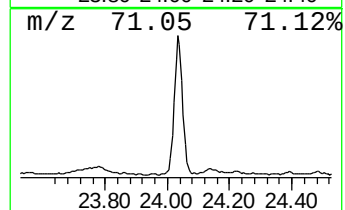
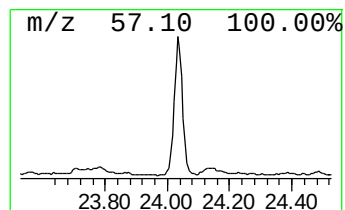
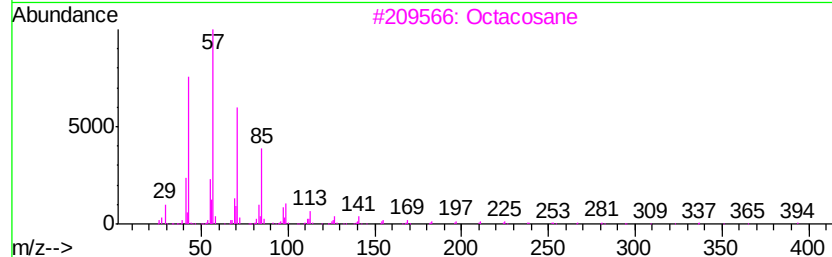
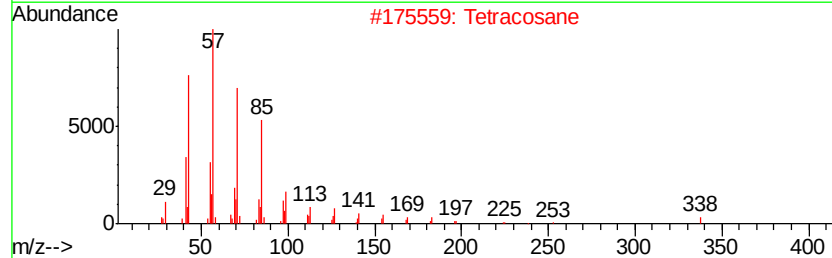
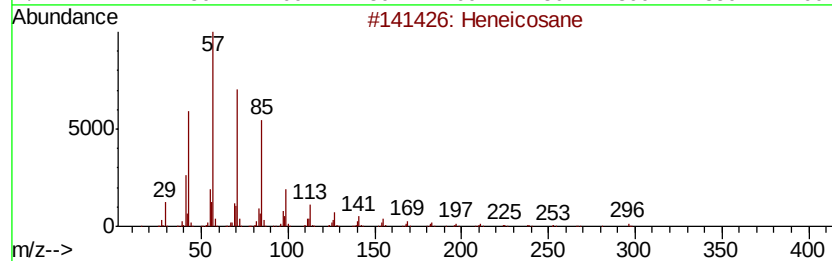
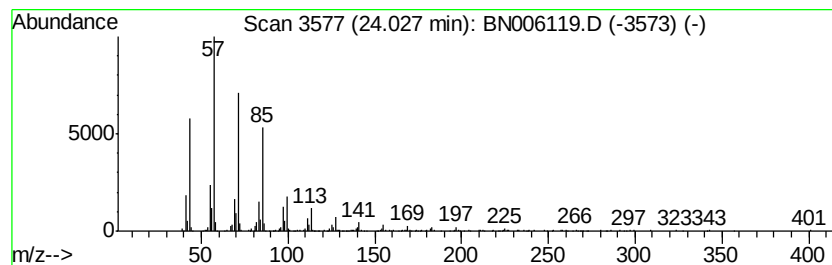
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 6

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T.  |
|-------|------------|---------|------------------|-------|
| 24.03 | 4.75 ng/ul | 1130350 | Perylene-d12     | 23.50 |

| Hit# | of | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------|-----|---------|-------------|------|
| 1    | 5  | Heneicosane           | 296 | C21H44  | 000629-94-7 | 91   |
| 2    |    | Tetracosane           | 338 | C24H50  | 000646-31-1 | 91   |
| 3    |    | Octacosane            | 394 | C28H58  | 000630-02-4 | 91   |
| 4    |    | Heptadecane, 9-octyl- | 352 | C25H52  | 007225-64-1 | 91   |
| 5    |    | Hexacosane            | 366 | C26H54  | 000630-01-3 | 90   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN060519\  
Data File : BN006119.D  
Acq On : 06 Jun 2019 03:14  
Operator : JU/SJ  
Sample : K3016-22  
Misc :  
ALS Vial : 28 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
A51D9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN060419MA.M  
Quant Title : SVOA CALIBRATION

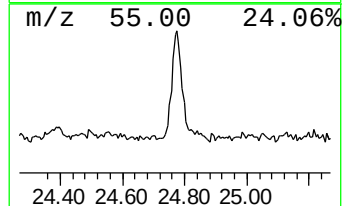
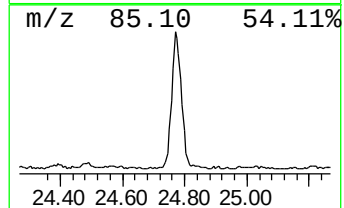
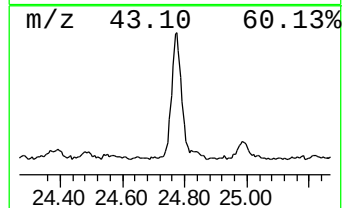
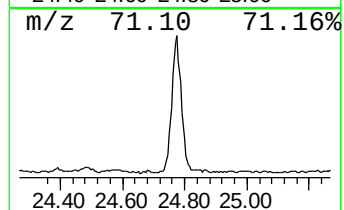
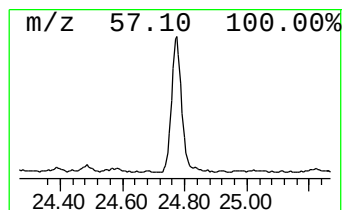
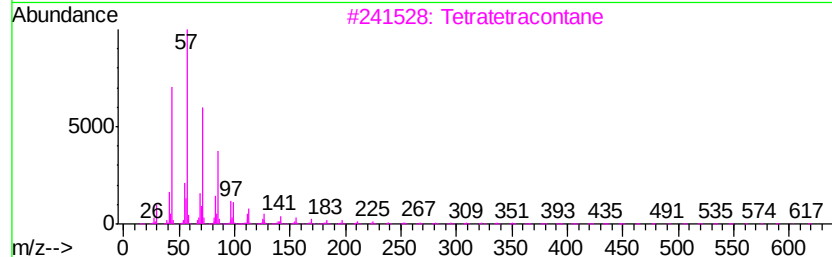
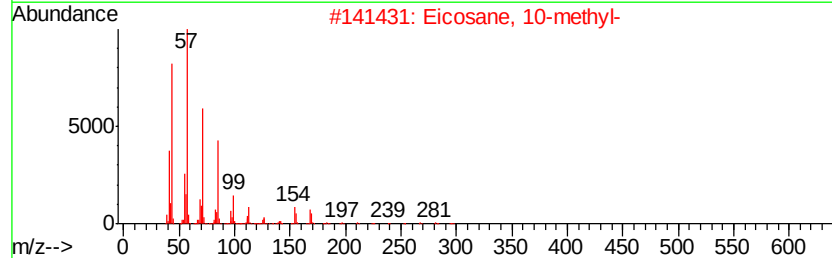
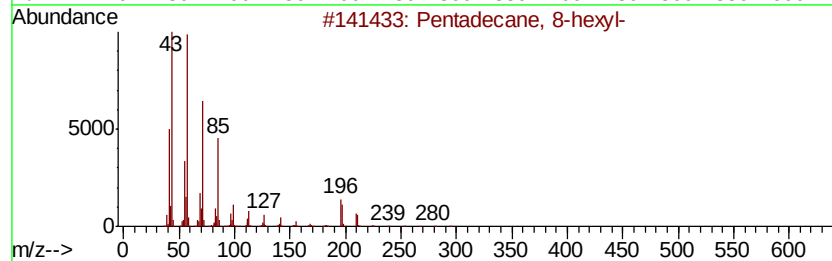
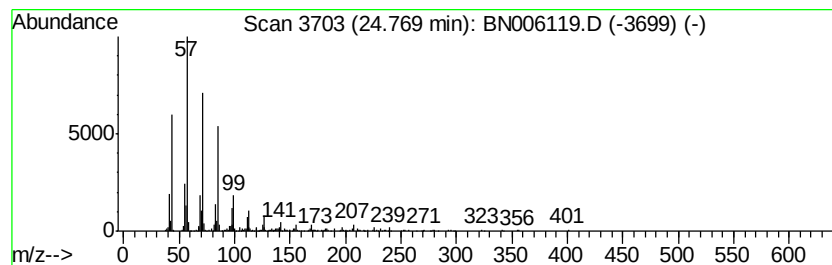
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 12

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 24.77 | 2.72 ng/ul | 646670 | Perylene-d12     | 23.50 |

| Hit# | of | 5 | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------|-----|---------|-------------|------|
| 1    |    |   | Pentadecane, 8-hexyl- | 296 | C21H44  | 013475-75-7 | 94   |
| 2    |    |   | Eicosane, 10-methyl-  | 296 | C21H44  | 054833-23-7 | 94   |
| 3    |    |   | Tetratetracontane     | 619 | C44H90  | 007098-22-8 | 91   |
| 4    |    |   | Octacosane            | 394 | C28H58  | 000630-02-4 | 91   |
| 5    |    |   | Heneicosane           | 296 | C21H44  | 000629-94-7 | 91   |





Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN060519\  
Data File : BN006119.D  
Acq On : 06 Jun 2019 03:14  
Operator : JU/SJ  
Sample : K3016-22  
Misc :  
ALS Vial : 28 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
A51D9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN060419MA.M  
Quant Title : SVOA CALIBRATION

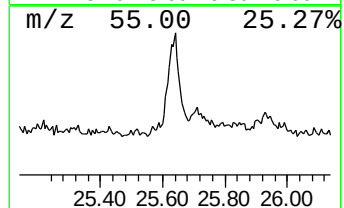
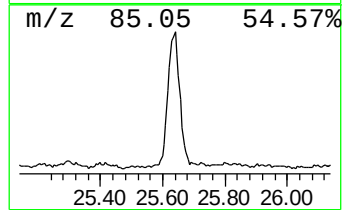
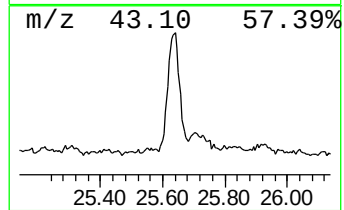
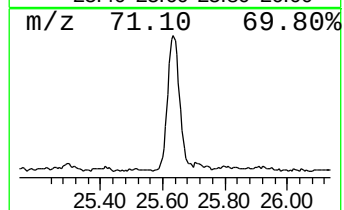
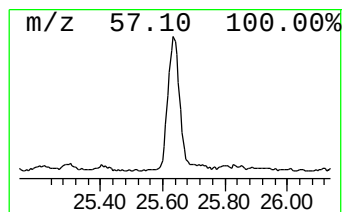
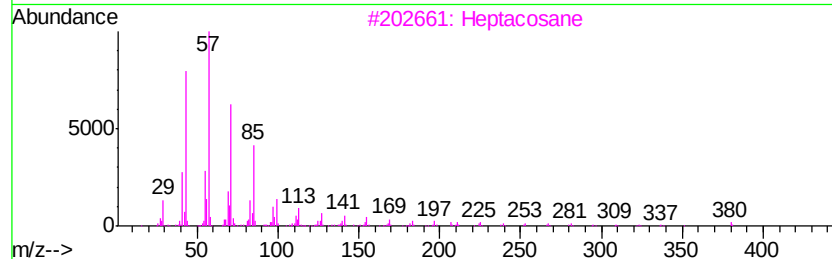
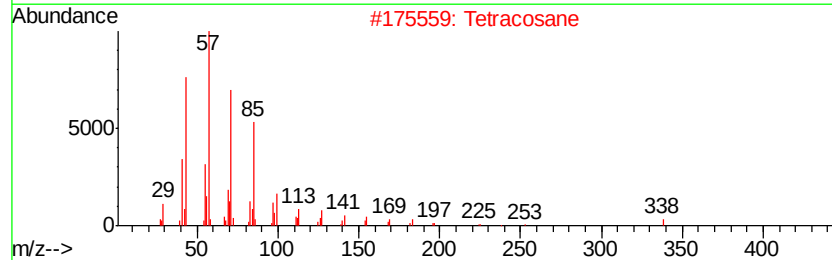
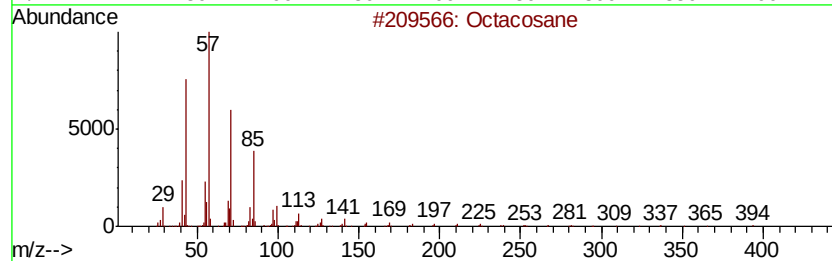
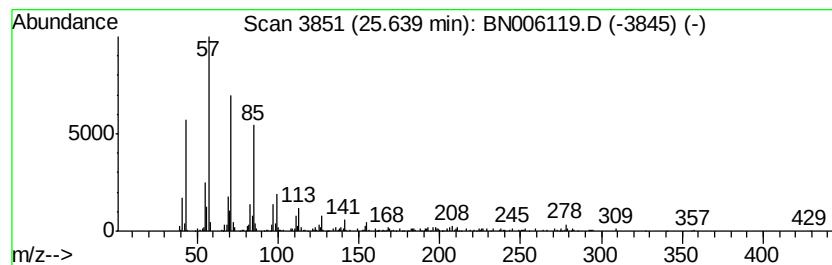
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 14

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 25.64 | 2.08 ng/ul | 495008 | Perylene-d12     | 23.50 |

| Hit# | of | 5 | Tentative ID       | MW  | MolForm | CAS#         | Qual |
|------|----|---|--------------------|-----|---------|--------------|------|
| 1    |    |   | Octacosane         | 394 | C28H58  | 000630-02-4  | 90   |
| 2    |    |   | Tetracosane        | 338 | C24H50  | 000646-31-1  | 87   |
| 3    |    |   | Heptacosane        | 380 | C27H56  | 000593-49-7  | 87   |
| 4    |    |   | 2-methyloctacosane | 408 | C29H60  | 1000376-72-8 | 87   |
| 5    |    |   | Tetradecane        | 198 | C14H30  | 000629-59-4  | 83   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN060519\  
Data File : BN006119.D  
Acq On : 06 Jun 2019 03:14  
Operator : JU/SJ  
Sample : K3016-22  
Misc :  
ALS Vial : 28 Sample Multiplier: 1

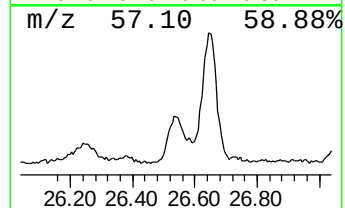
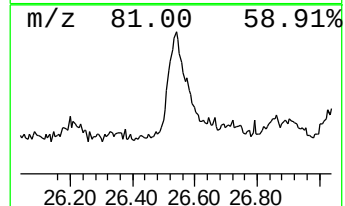
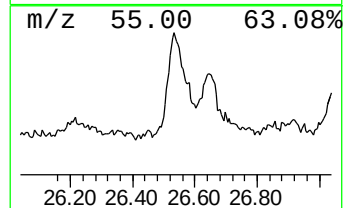
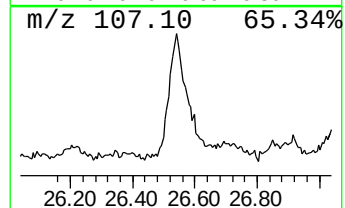
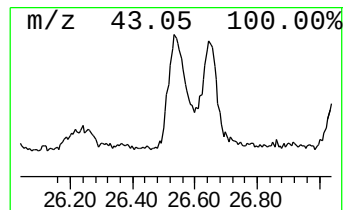
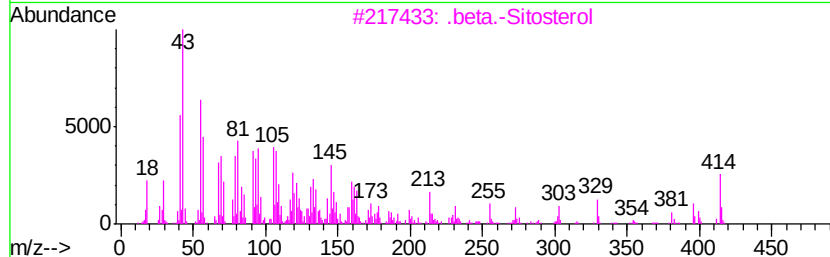
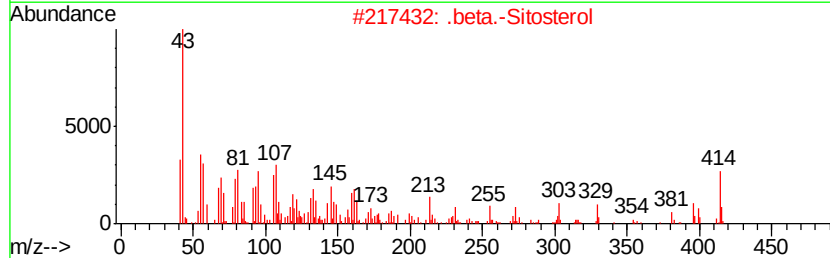
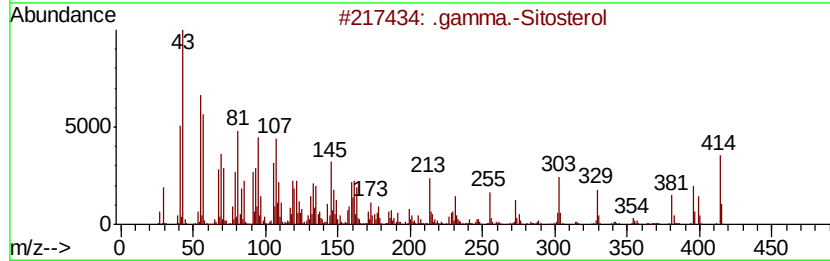
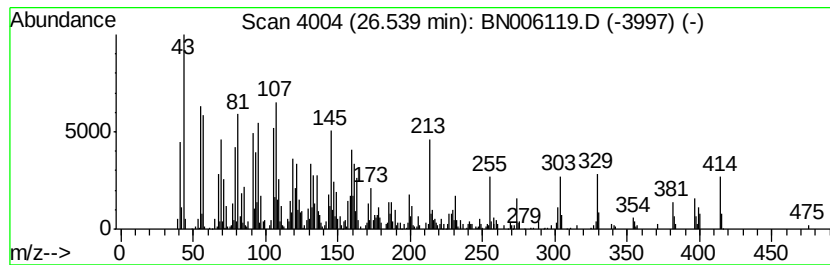
Instrument :  
BNA\_N  
ClientSampleId :  
A51D9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN060419MA.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 13 .gamma.-Sitosterol Concentration Rank 7

| R.T.    | EstConc    | Area                                | Relative to ISTD | R.T.            |
|---------|------------|-------------------------------------|------------------|-----------------|
| 26.54   | 4.74 ng/ul | 1127070                             | Perylene-d12     | 23.50           |
| Hit# of | 5          | Tentative ID                        | MW MolForm       | CAS# Qual       |
| 1       |            | .gamma.-Sitosterol                  | 414 C29H50O      | 000083-47-6 99  |
| 2       |            | .beta.-Sitosterol                   | 414 C29H50O      | 000083-46-5 94  |
| 3       |            | .beta.-Sitosterol                   | 414 C29H50O      | 000083-46-5 94  |
| 4       |            | Cholest-5-ene, 3-ethoxy-, (3.bet... | 414 C29H50O      | 000986-19-6 49  |
| 5       |            | 5-Cholestene-3-ol, 24-methyl-       | 400 C28H48O      | 1000214-17-4 43 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN060519\  
Data File : BN006119.D  
Acq On : 06 Jun 2019 03:14  
Operator : JU/SJ  
Sample : K3016-22  
Misc :  
ALS Vial : 28 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
A51D9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN060419MA.M  
Quant Title : SVOA CALIBRATION

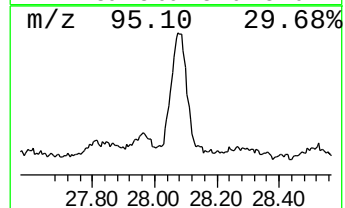
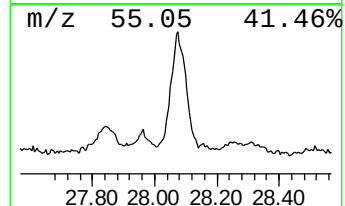
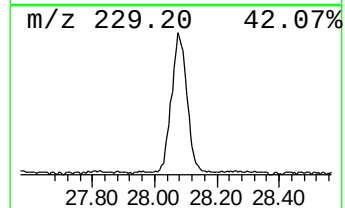
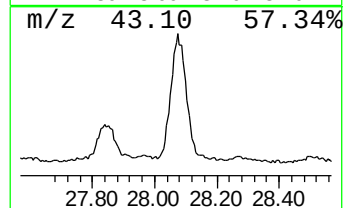
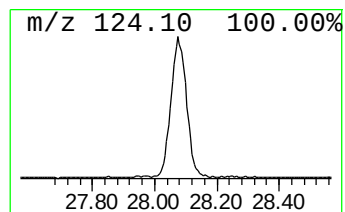
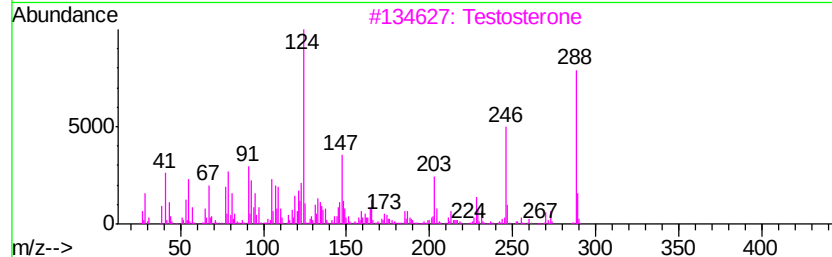
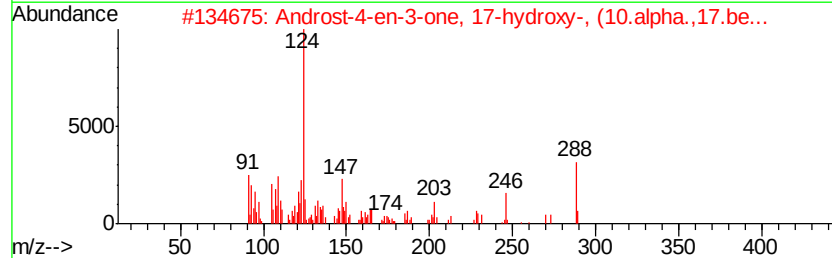
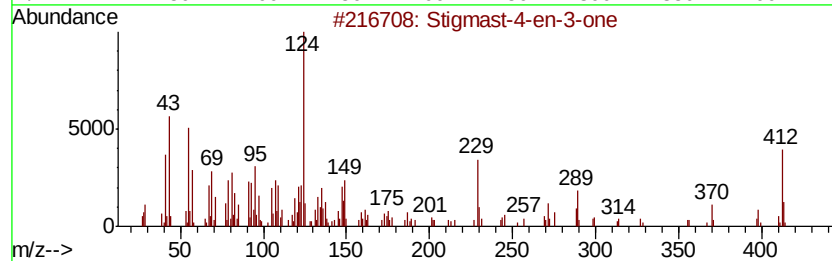
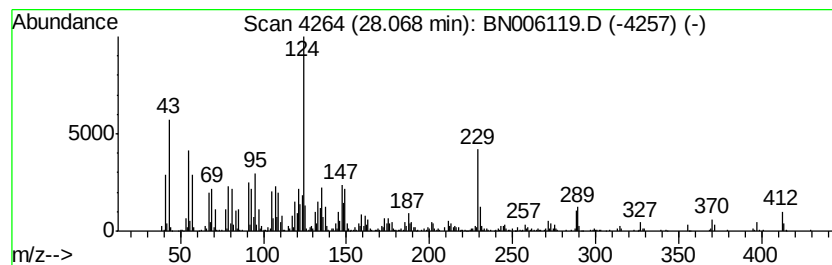
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 14 Stigmast-4-en-3-one Concentration Rank 1

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 28.07 | 11.03 ng/ul | 2623960 | Perylene-d12     | 23.50 |

| Hit# | of | Tentative ID                         | MW  | MolForm     | CAS#        | Qual |
|------|----|--------------------------------------|-----|-------------|-------------|------|
| 1    | 5  | Stigmast-4-en-3-one                  | 412 | C29H48O     | 001058-61-3 | 94   |
| 2    |    | Androst-4-en-3-one, 17-hydroxy-, ... | 288 | C19H28O2    | 000604-39-7 | 50   |
| 3    |    | Testosterone                         | 288 | C19H28O2    | 000058-22-0 | 49   |
| 4    |    | Androst-4-en-3-one, 17-hydroxy-, ... | 288 | C19H28O2    | 000481-30-1 | 35   |
| 5    |    | 3-(4-Fluoroanilino)-1-(3-nitroph...  | 288 | C15H13FN2O3 | 350039-84-8 | 27   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_N\DATA\BN060519\  
Data File : BN006119.D  
Acq On : 06 Jun 2019 03:14  
Operator : JU/SJ  
Sample : K3016-22  
Misc :  
ALS Vial : 28 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
A51D9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN060419MA.M  
Quant Title : SVOA CALIBRATION

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

| TIC Top Hit name     | RT    | EstConc | Units | Response | --Internal Standard-- |       |         |      |
|----------------------|-------|---------|-------|----------|-----------------------|-------|---------|------|
|                      |       |         |       |          | #                     | RT    | Resp    | Conc |
| (1R)-2,6,6-Trimet... | 6.33  | 3.4     | ng/ul | 372101   | 1                     | 7.65  | 2180050 | 20.0 |
| 1-Naphthaleneprop... | 18.89 | 5.2     | ng/ul | 1164580  | 4                     | 17.05 | 4503380 | 20.0 |
| unknown-01           | 19.27 | 8.0     | ng/ul | 2035850  | 5                     | 21.27 | 5080160 | 20.0 |
| unknown-02           | 19.72 | 3.4     | ng/ul | 866407   | 5                     | 21.27 | 5080160 | 20.0 |
| 1-Phenanthrenecar... | 20.30 | 6.2     | ng/ul | 1574630  | 5                     | 21.27 | 5080160 | 20.0 |
| Methyl dehydroabi... | 20.45 | 10.5    | ng/ul | 2656610  | 5                     | 21.27 | 5080160 | 20.0 |
| 1-Phenanthrenecar... | 20.96 | 2.3     | ng/ul | 587702   | 5                     | 21.27 | 5080160 | 20.0 |
| (DEL) Alkane: Cyc... | 20.99 | 3.0     | ng/ul | 753668   | 5                     | 21.27 | 5080160 | 20.0 |
| 7-Oxodehydroabiet... | 21.59 | 4.3     | ng/ul | 1085620  | 5                     | 21.27 | 5080160 | 20.0 |
| (DEL) Alkane: Str... | 24.03 | 4.8     | ng/ul | 1130350  | 6                     | 23.50 | 4758990 | 20.0 |
| (DEL) Alkane: Str... | 24.77 | 2.7     | ng/ul | 646670   | 6                     | 23.50 | 4758990 | 20.0 |
| (DEL) Alkane: Str... | 25.64 | 2.1     | ng/ul | 495008   | 6                     | 23.50 | 4758990 | 20.0 |
| .gamma.-Sitosterol   | 26.54 | 4.7     | ng/ul | 1127070  | 6                     | 23.50 | 4758990 | 20.0 |
| Stigmast-4-en-3-one  | 28.07 | 11.0    | ng/ul | 2623960  | 6                     | 23.50 | 4758990 | 20.0 |