

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP060723\  
 Data File : BP015404.D  
 Acq On : 07 Jun 2023 18:08  
 Operator : MA/JU  
 Sample : 02950-01  
 Misc :  
 ALS Vial : 10 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 EW978

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % 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\_P\Methods\SFAM-EPA-BP060723.MA.M  
 Title : SVOA CALIBRATION

Signal : TIC: BP015404.D\data.ms

| 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         | 8.099       | 830           | 835         | 843          | rVB      | 896644         | 1420131       | 7.29%           | 0.342%        |
| 2         | 10.928      | 1308          | 1316        | 1322         | rVV      | 1297717        | 2459295       | 12.63%          | 0.592%        |
| 3         | 12.799      | 1625          | 1634        | 1643         | rBV      | 3655290        | 5795425       | 29.76%          | 1.395%        |
| 4         | 13.798      | 1799          | 1804        | 1813         | rVB      | 988185         | 1722626       | 8.84%           | 0.415%        |
| 5         | 13.922      | 1817          | 1825        | 1832         | rVB      | 1900133        | 3460334       | 17.77%          | 0.833%        |
| 6         | 14.069      | 1843          | 1850        | 1856         | rVV      | 5748743        | 10087554      | 51.80%          | 2.429%        |
| 7         | 14.145      | 1856          | 1863        | 1868         | rVV      | 1258173        | 2152118       | 11.05%          | 0.518%        |
| 8         | 14.204      | 1868          | 1873        | 1878         | rVB2     | 972626         | 1499338       | 7.70%           | 0.361%        |
| 9         | 14.304      | 1884          | 1890        | 1893         | rVV      | 3410065        | 5213319       | 26.77%          | 1.255%        |
| 10        | 14.334      | 1893          | 1895        | 1900         | rVB      | 1085214        | 1259898       | 6.47%           | 0.303%        |
| 11        | 14.440      | 1908          | 1913        | 1914         | rBV      | 1208367        | 1445164       | 7.42%           | 0.348%        |
| 12        | 14.469      | 1914          | 1918        | 1928         | rVB      | 2646159        | 4188221       | 21.50%          | 1.008%        |
| 13        | 14.710      | 1953          | 1959        | 1962         | rBV      | 1397484        | 2410826       | 12.38%          | 0.580%        |
| 14        | 14.745      | 1962          | 1965        | 1970         | rVB      | 2013951        | 2772467       | 14.24%          | 0.667%        |
| 15        | 14.810      | 1970          | 1976        | 1982         | rBV      | 6550456        | 10514149      | 53.99%          | 2.531%        |
| 16        | 14.951      | 1993          | 2000        | 2008         | rVB      | 2774086        | 6484286       | 33.29%          | 1.561%        |
| 17        | 15.034      | 2008          | 2014        | 2022         | rBV2     | 1394641        | 2568998       | 13.19%          | 0.618%        |
| 18        | 15.140      | 2022          | 2032        | 2039         | rVB3     | 3317085        | 6806956       | 34.95%          | 1.639%        |
| 19        | 15.216      | 2039          | 2045        | 2050         | rBV      | 2876980        | 4140103       | 21.26%          | 0.997%        |
| 20        | 15.357      | 2061          | 2069        | 2074         | rBV      | 2613672        | 5055374       | 25.96%          | 1.217%        |
| 21        | 15.410      | 2074          | 2078        | 2082         | rVV      | 2247070        | 2931705       | 15.05%          | 0.706%        |
| 22        | 15.528      | 2090          | 2098        | 2101         | rBV2     | 1974498        | 4284832       | 22.00%          | 1.032%        |
| 23        | 15.563      | 2101          | 2104        | 2109         | rVB2     | 1326092        | 1895381       | 9.73%           | 0.456%        |
| 24        | 15.616      | 2109          | 2113        | 2118         | rBV      | 1030399        | 1476362       | 7.58%           | 0.355%        |
| 25        | 15.675      | 2118          | 2123        | 2124         | rBV2     | 744571         | 1074911       | 5.52%           | 0.259%        |
| 26        | 15.740      | 2130          | 2134        | 2138         | rVB      | 2809450        | 3860806       | 19.82%          | 0.929%        |
| 27        | 15.792      | 2138          | 2143        | 2153         | rVB2     | 3993927        | 7409431       | 38.04%          | 1.784%        |
| 28        | 15.887      | 2153          | 2159        | 2162         | rBV2     | 2436543        | 3828408       | 19.66%          | 0.922%        |
| 29        | 15.957      | 2167          | 2171        | 2175         | rVV4     | 1073908        | 1930816       | 9.91%           | 0.465%        |
| 30        | 16.004      | 2175          | 2179        | 2183         | rVB2     | 2228669        | 3152917       | 16.19%          | 0.759%        |
| 31        | 16.045      | 2183          | 2186        | 2189         | rBV2     | 813816         | 1118770       | 5.74%           | 0.269%        |
| 32        | 16.204      | 2208          | 2213        | 2216         | rBV3     | 1484036        | 2437856       | 12.52%          | 0.587%        |
| 33        | 16.234      | 2216          | 2218        | 2225         | rVB      | 1317339        | 1679274       | 8.62%           | 0.404%        |
| 34        | 16.457      | 2248          | 2256        | 2262         | rBV4     | 3286977        | 6995633       | 35.92%          | 1.684%        |
| 35        | 16.598      | 2276          | 2280        | 2284         | rBV      | 849764         | 1380440       | 7.09%           | 0.332%        |
| 36        | 16.798      | 2306          | 2314        | 2319         | rVV2     | 2446593        | 6263144       | 32.16%          | 1.508%        |

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 ALS Vial : 10 Sample Multiplier: 1

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 BNA\_P  
 ClientSampleId :  
 EW978

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % 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\_P\Methods\SFAM-EPA-BP060723.MA.M  
 Title : SVOA CALIBRATION

|    |        |      |      |      |      |          |          |         |        |
|----|--------|------|------|------|------|----------|----------|---------|--------|
| 37 | 16.851 | 2319 | 2323 | 2328 | rVV  | 2826450  | 4410724  | 22.65%  | 1.062% |
| 38 | 16.904 | 2329 | 2332 | 2335 | rVV3 | 701311   | 1139494  | 5.85%   | 0.274% |
| 39 | 16.951 | 2336 | 2340 | 2345 | rVB  | 1943239  | 2813370  | 14.45%  | 0.677% |
| 40 | 17.157 | 2372 | 2375 | 2380 | rVB2 | 1505454  | 2122778  | 10.90%  | 0.511% |
| 41 | 17.234 | 2384 | 2388 | 2392 | rVV3 | 546428   | 1060543  | 5.45%   | 0.255% |
| 42 | 17.287 | 2393 | 2397 | 2399 | rVV  | 1183045  | 1800617  | 9.25%   | 0.434% |
| 43 | 17.322 | 2399 | 2403 | 2410 | rVB2 | 2389121  | 3800530  | 19.51%  | 0.915% |
| 44 | 17.504 | 2429 | 2434 | 2437 | rBV  | 2122420  | 3136491  | 16.10%  | 0.755% |
| 45 | 17.557 | 2437 | 2443 | 2447 | rVV  | 12446255 | 19475792 | 100.00% | 4.689% |
| 46 | 17.604 | 2447 | 2451 | 2454 | rVV2 | 1655684  | 2483392  | 12.75%  | 0.598% |
| 47 | 17.639 | 2454 | 2457 | 2461 | rVV  | 5206754  | 6635727  | 34.07%  | 1.598% |
| 48 | 17.692 | 2461 | 2466 | 2471 | rVV2 | 1268297  | 2647543  | 13.59%  | 0.637% |
| 49 | 17.739 | 2471 | 2474 | 2476 | rVV2 | 892319   | 1006887  | 5.17%   | 0.242% |
| 50 | 17.769 | 2476 | 2479 | 2488 | rVB3 | 635434   | 1562837  | 8.02%   | 0.376% |
| 51 | 17.904 | 2498 | 2502 | 2511 | rBV3 | 699072   | 1872196  | 9.61%   | 0.451% |
| 52 | 18.075 | 2527 | 2531 | 2536 | rVB  | 2391054  | 3320343  | 17.05%  | 0.799% |
| 53 | 18.216 | 2550 | 2555 | 2562 | rBV2 | 1511905  | 2749198  | 14.12%  | 0.662% |
| 54 | 18.363 | 2575 | 2580 | 2584 | rBV  | 6971726  | 10316457 | 52.97%  | 2.484% |
| 55 | 18.416 | 2584 | 2589 | 2593 | rVV  | 7521861  | 10499559 | 53.91%  | 2.528% |
| 56 | 18.492 | 2597 | 2602 | 2606 | rVV  | 3680937  | 5542971  | 28.46%  | 1.334% |
| 57 | 18.551 | 2606 | 2612 | 2616 | rVV2 | 8963382  | 17328335 | 88.97%  | 4.172% |
| 58 | 18.592 | 2616 | 2619 | 2624 | rVB  | 6489779  | 8030723  | 41.23%  | 1.933% |
| 59 | 18.745 | 2641 | 2645 | 2649 | rVB  | 1018576  | 1452712  | 7.46%   | 0.350% |
| 60 | 18.839 | 2657 | 2661 | 2665 | rBV2 | 2677331  | 3477484  | 17.86%  | 0.837% |
| 61 | 18.963 | 2678 | 2682 | 2685 | rBV2 | 631018   | 965917   | 4.96%   | 0.233% |
| 62 | 19.051 | 2693 | 2697 | 2698 | rBV2 | 1086749  | 1322633  | 6.79%   | 0.318% |
| 63 | 19.075 | 2698 | 2701 | 2705 | rVV  | 2490288  | 3573487  | 18.35%  | 0.860% |
| 64 | 19.128 | 2706 | 2710 | 2713 | rVV  | 2113871  | 3078657  | 15.81%  | 0.741% |
| 65 | 19.163 | 2713 | 2716 | 2720 | rVB2 | 1419653  | 1764602  | 9.06%   | 0.425% |
| 66 | 19.245 | 2725 | 2730 | 2734 | rVV2 | 6372764  | 10072570 | 51.72%  | 2.425% |
| 67 | 19.304 | 2734 | 2740 | 2743 | rVV3 | 3431311  | 7699851  | 39.54%  | 1.854% |
| 68 | 19.333 | 2743 | 2745 | 2748 | rVV  | 2530386  | 2725215  | 13.99%  | 0.656% |
| 69 | 19.392 | 2748 | 2755 | 2767 | rVB3 | 2367252  | 9069131  | 46.57%  | 2.183% |
| 70 | 19.504 | 2768 | 2774 | 2779 | rBV  | 8376259  | 12624620 | 64.82%  | 3.039% |
| 71 | 19.557 | 2779 | 2783 | 2786 | rVV2 | 957184   | 1379559  | 7.08%   | 0.332% |
| 72 | 19.592 | 2786 | 2789 | 2793 | rVB  | 1437255  | 1806535  | 9.28%   | 0.435% |
| 73 | 19.645 | 2794 | 2798 | 2808 | rVB  | 3373283  | 5416031  | 27.81%  | 1.304% |
| 74 | 19.733 | 2808 | 2813 | 2815 | rBV3 | 899423   | 1501881  | 7.71%   | 0.362% |
| 75 | 19.763 | 2816 | 2818 | 2824 | rVB  | 1169560  | 1331345  | 6.84%   | 0.321% |
| 76 | 19.828 | 2825 | 2829 | 2831 | rBV3 | 2950001  | 4258608  | 21.87%  | 1.025% |

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Instrument :  
 BNA\_P  
 ClientSampleId :  
 EW978

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % 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\_P\Methods\SFAM-EPA-BP060723.MA.M  
 Title : SVOA CALIBRATION

|     |        |      |      |      |      |          |          |        |        |
|-----|--------|------|------|------|------|----------|----------|--------|--------|
| 77  | 19.863 | 2831 | 2835 | 2841 | rVB  | 11539612 | 16301651 | 83.70% | 3.925% |
| 78  | 19.922 | 2841 | 2845 | 2850 | rVB  | 1722858  | 2507475  | 12.87% | 0.604% |
| 79  | 19.969 | 2850 | 2853 | 2855 | rBV2 | 740423   | 971431   | 4.99%  | 0.234% |
| 80  | 20.075 | 2868 | 2871 | 2877 | rVB2 | 948025   | 1516494  | 7.79%  | 0.365% |
| 81  | 20.169 | 2881 | 2887 | 2895 | rVB4 | 1963096  | 4717642  | 24.22% | 1.136% |
| 82  | 20.333 | 2905 | 2915 | 2922 | rVB  | 4001229  | 9395433  | 48.24% | 2.262% |
| 83  | 20.433 | 2928 | 2932 | 2937 | rVV  | 2906873  | 3913031  | 20.09% | 0.942% |
| 84  | 20.486 | 2938 | 2941 | 2945 | rVB  | 2542584  | 3159635  | 16.22% | 0.761% |
| 85  | 20.580 | 2954 | 2957 | 2961 | rVB  | 902128   | 1194981  | 6.14%  | 0.288% |
| 86  | 20.628 | 2961 | 2965 | 2969 | rVV  | 2748451  | 3855332  | 19.80% | 0.928% |
| 87  | 20.669 | 2969 | 2972 | 2983 | rVB  | 3405299  | 5432854  | 27.90% | 1.308% |
| 88  | 21.227 | 3064 | 3067 | 3070 | rBV  | 1156248  | 1465277  | 7.52%  | 0.353% |
| 89  | 21.263 | 3070 | 3073 | 3077 | rVB  | 1756073  | 2227574  | 11.44% | 0.536% |
| 90  | 21.575 | 3121 | 3126 | 3132 | rBV2 | 5213519  | 11962769 | 61.42% | 2.880% |
| 91  | 21.627 | 3132 | 3135 | 3144 | rVB  | 4730630  | 6577159  | 33.77% | 1.583% |
| 92  | 21.751 | 3153 | 3156 | 3161 | rBV2 | 658336   | 1095536  | 5.63%  | 0.264% |
| 93  | 21.810 | 3163 | 3166 | 3175 | rVB3 | 859608   | 1601822  | 8.22%  | 0.386% |
| 94  | 22.192 | 3228 | 3231 | 3237 | rVB  | 1608874  | 2168424  | 11.13% | 0.522% |
| 95  | 22.245 | 3237 | 3240 | 3245 | rVB2 | 1138656  | 1529744  | 7.85%  | 0.368% |
| 96  | 23.357 | 3424 | 3429 | 3442 | rVB  | 1090506  | 3847186  | 19.75% | 0.926% |
| 97  | 23.916 | 3519 | 3524 | 3529 | rBV  | 847984   | 1590335  | 8.17%  | 0.383% |
| 98  | 23.974 | 3529 | 3534 | 3538 | rVV2 | 1036091  | 2108067  | 10.82% | 0.508% |
| 99  | 24.027 | 3539 | 3543 | 3551 | rVB  | 1714802  | 3260632  | 16.74% | 0.785% |
| 100 | 24.139 | 3557 | 3562 | 3568 | rVB  | 1271876  | 2510066  | 12.89% | 0.604% |

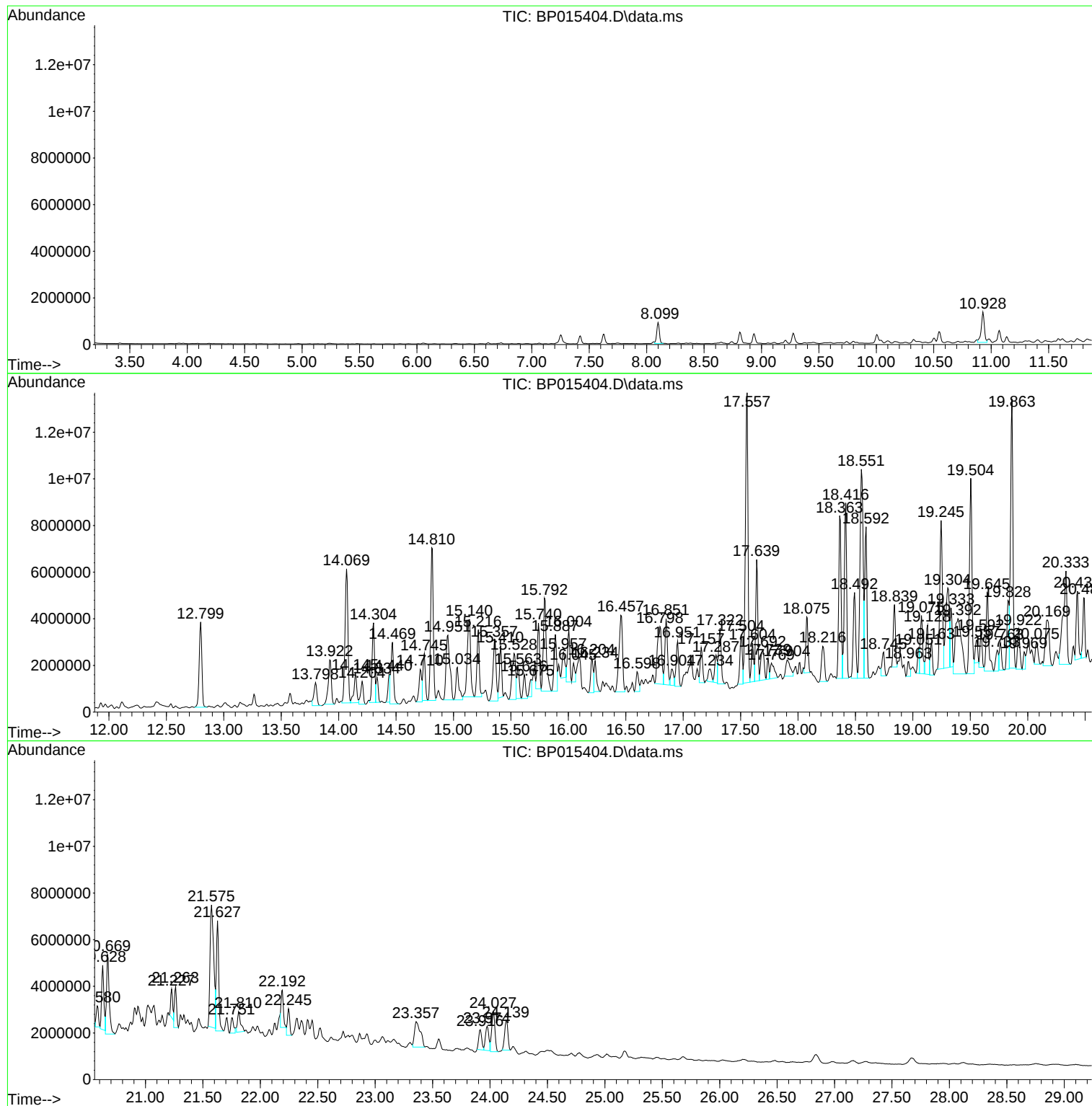
Sum of corrected areas: 415365163

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Instrument :  
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Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP060723.MA.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP060723\  
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Instrument :  
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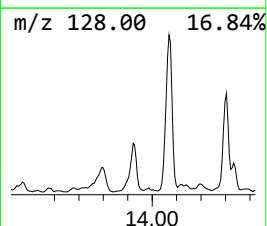
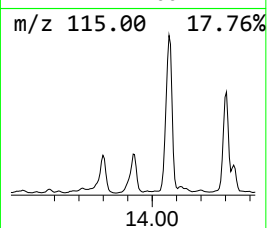
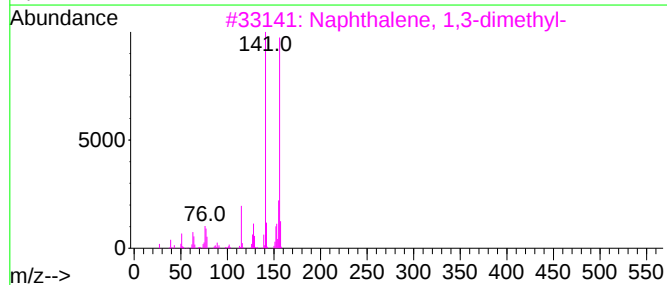
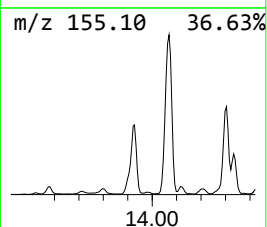
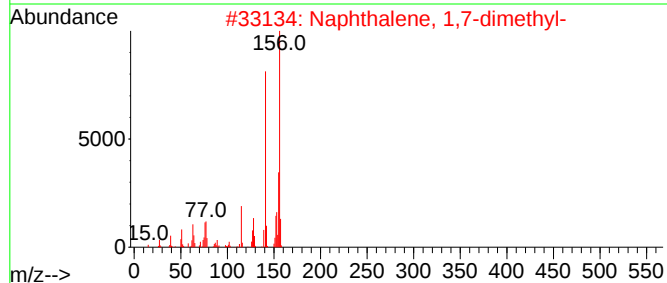
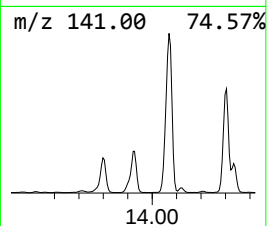
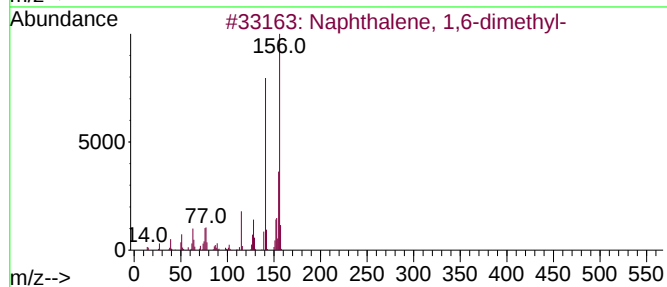
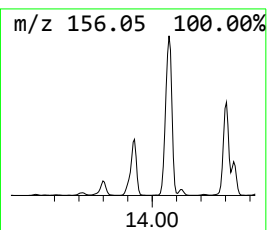
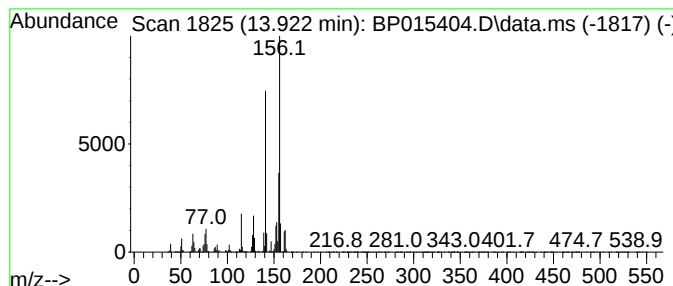
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP060723.MA.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 2 Naphthalene, 1,6-dimethyl- Concentration Rank 18

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 13.922 | 24.96 ng/ul | 3460330 | Acenaphthene-d10 | 14.746 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, 1,6-dimethyl- | 156 | C12H12  | 000575-43-9 | 98   |
| 2    |    |   | Naphthalene, 1,7-dimethyl- | 156 | C12H12  | 000575-37-1 | 96   |
| 3    |    |   | Naphthalene, 1,3-dimethyl- | 156 | C12H12  | 000575-41-7 | 96   |
| 4    |    |   | Naphthalene, 1,7-dimethyl- | 156 | C12H12  | 000575-37-1 | 96   |
| 5    |    |   | Naphthalene, 2,7-dimethyl- | 156 | C12H12  | 000582-16-1 | 96   |



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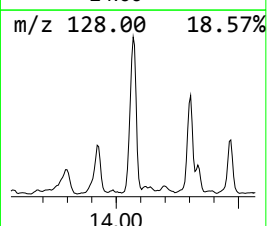
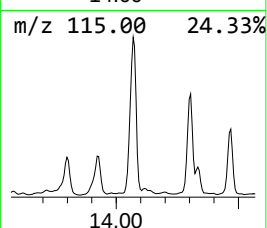
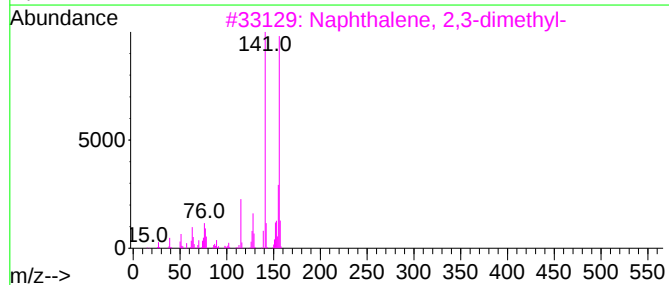
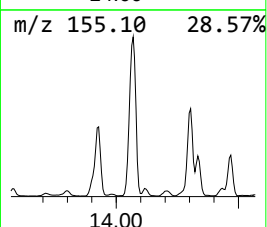
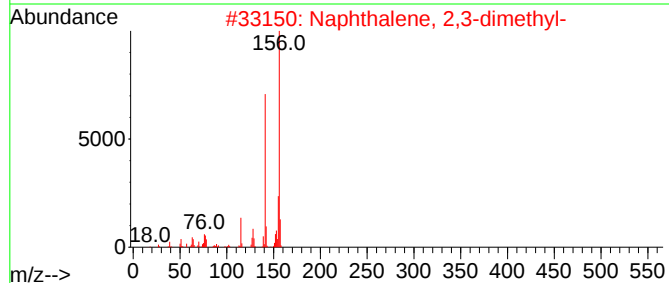
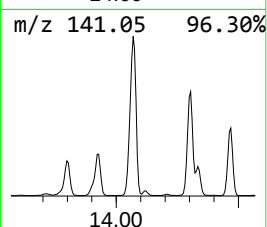
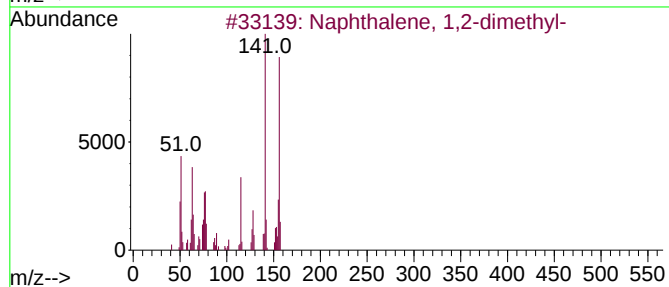
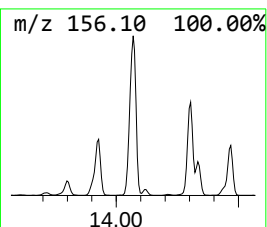
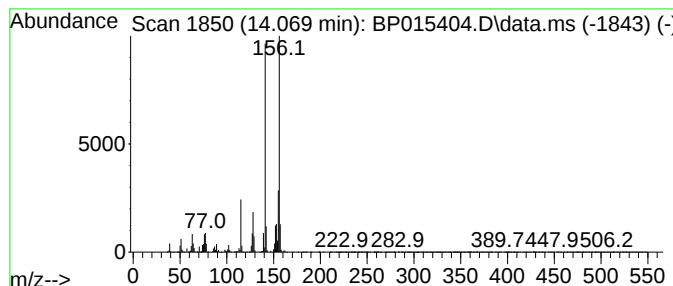
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP060723.MA.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 3 Naphthalene, 1,2-dimethyl- Concentration Rank 2

| R.T.   | EstConc     | Area     | Relative to ISTD | R.T.   |
|--------|-------------|----------|------------------|--------|
| 14.069 | 72.77 ng/ul | 10087600 | Acenaphthene-d10 | 14.746 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, 1,2-dimethyl- | 156 | C12H12  | 000573-98-8 | 97   |
| 2    |    |   | Naphthalene, 2,3-dimethyl- | 156 | C12H12  | 000581-40-8 | 97   |
| 3    |    |   | Naphthalene, 2,3-dimethyl- | 156 | C12H12  | 000581-40-8 | 96   |
| 4    |    |   | Naphthalene, 1,5-dimethyl- | 156 | C12H12  | 000571-61-9 | 96   |
| 5    |    |   | Naphthalene, 1,6-dimethyl- | 156 | C12H12  | 000575-43-9 | 96   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP060723\  
Data File : BP015404.D  
Acq On : 07 Jun 2023 18:08  
Operator : MA/JU  
Sample : 02950-01  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
EW978

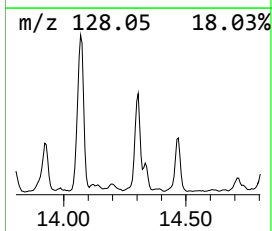
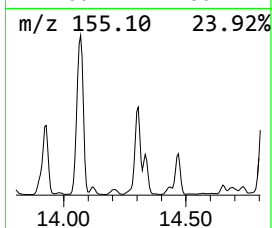
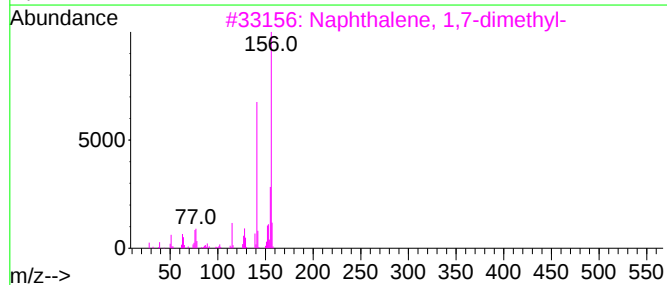
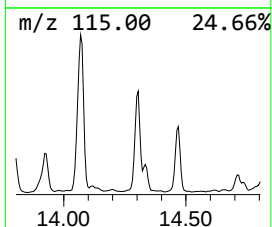
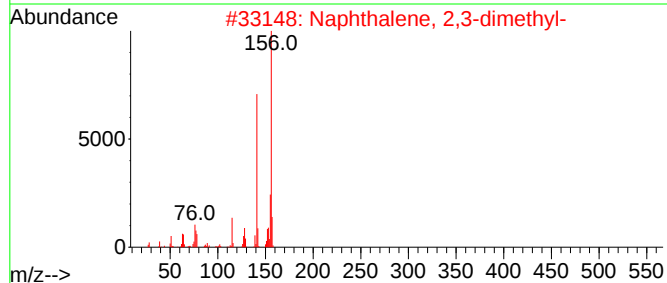
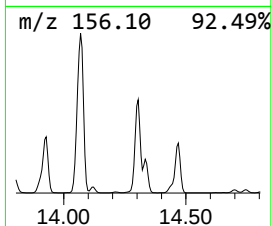
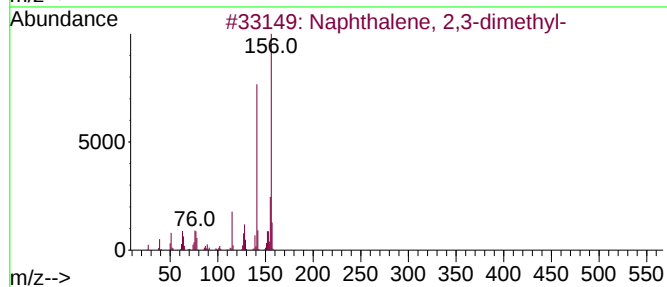
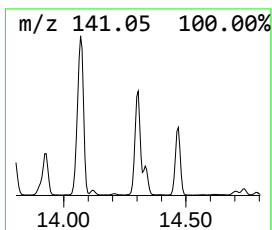
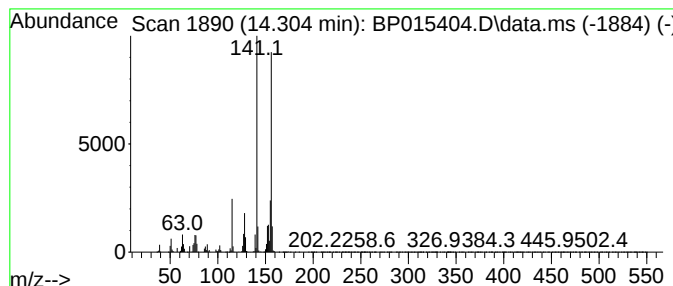
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 5 Naphthalene, 2,3-dimethyl- Concentration Rank 12

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 14.304 | 37.61 ng/ul | 5213320 | Acenaphthene-d10 | 14.746 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, 2,3-dimethyl- | 156 | C12H12  | 000581-40-8 | 97   |
| 2    |    |   | Naphthalene, 2,3-dimethyl- | 156 | C12H12  | 000581-40-8 | 97   |
| 3    |    |   | Naphthalene, 1,7-dimethyl- | 156 | C12H12  | 000575-37-1 | 97   |
| 4    |    |   | Naphthalene, 2,3-dimethyl- | 156 | C12H12  | 000581-40-8 | 97   |
| 5    |    |   | Naphthalene, 1,5-dimethyl- | 156 | C12H12  | 000571-61-9 | 97   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP060723\  
Data File : BP015404.D  
Acq On : 07 Jun 2023 18:08  
Operator : MA/JU  
Sample : 02950-01  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
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Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP060723.MA.M  
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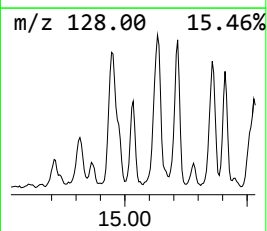
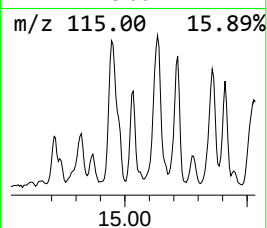
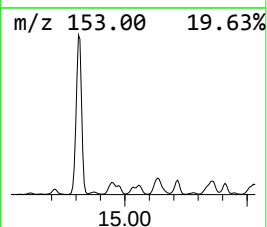
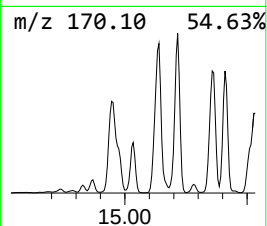
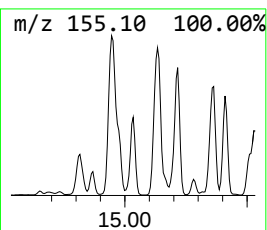
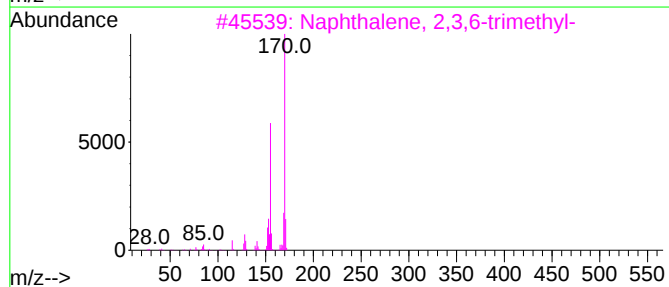
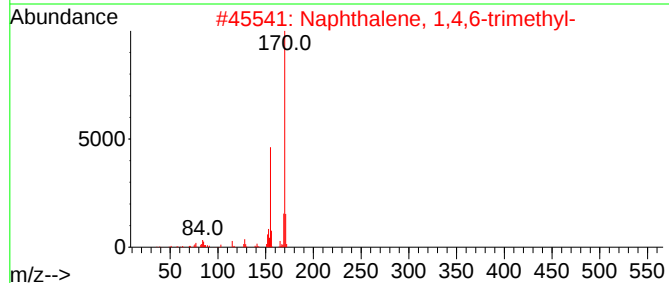
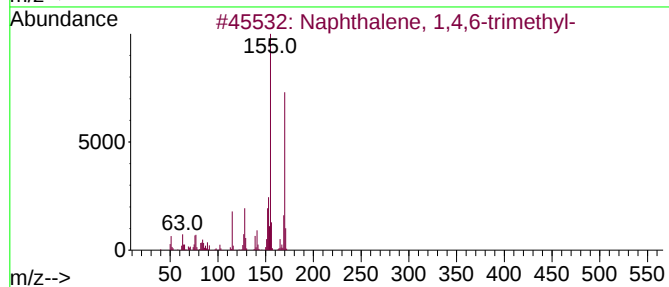
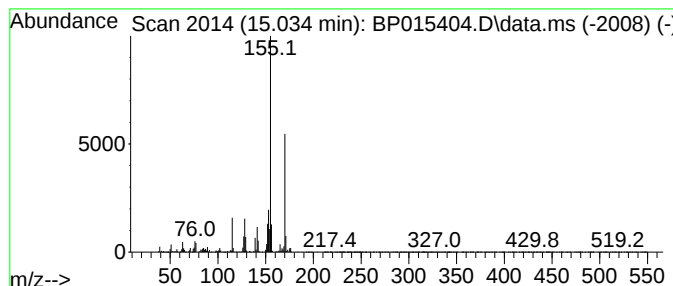
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TIC Integration Parameters: LSCINT.P

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Peak Number 7 Naphthalene, 1,4,6-trimethyl- Concentration Rank 26

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 15.034 | 18.53 ng/ul | 2569000 | Acenaphthene-d10 | 14.746 |

| Hit# | of | 5 | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, 1,4,6-trimethyl- | 170 | C13H14  | 002131-42-2 | 95   |
| 2    |    |   | Naphthalene, 1,4,6-trimethyl- | 170 | C13H14  | 002131-42-2 | 94   |
| 3    |    |   | Naphthalene, 2,3,6-trimethyl- | 170 | C13H14  | 000829-26-5 | 94   |
| 4    |    |   | Naphthalene, 1,6,7-trimethyl- | 170 | C13H14  | 002245-38-7 | 94   |
| 5    |    |   | Naphthalene, 2,3,6-trimethyl- | 170 | C13H14  | 000829-26-5 | 94   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP060723\  
Data File : BP015404.D  
Acq On : 07 Jun 2023 18:08  
Operator : MA/JU  
Sample : 02950-01  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
EW978

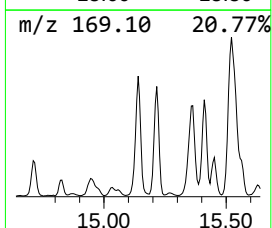
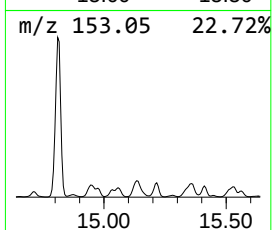
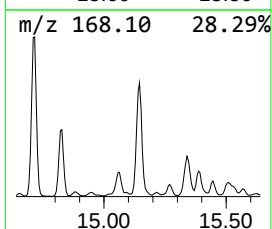
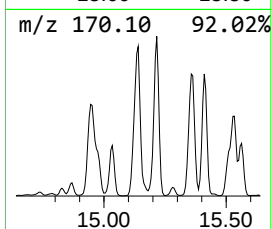
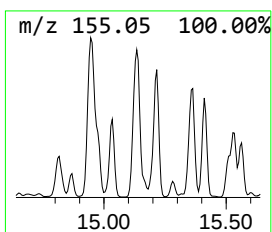
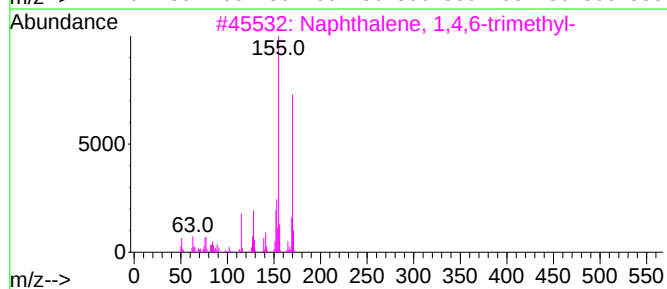
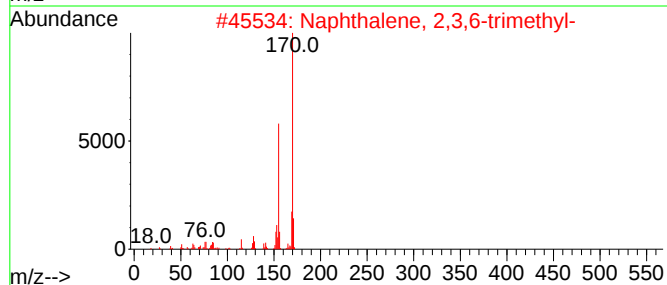
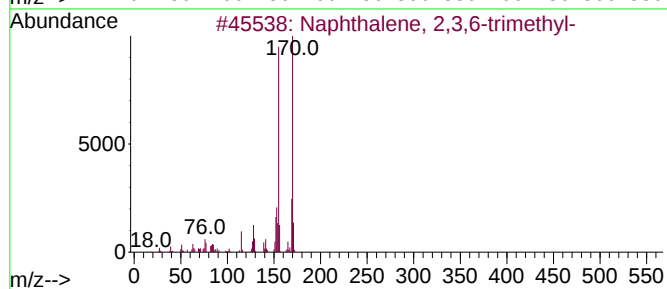
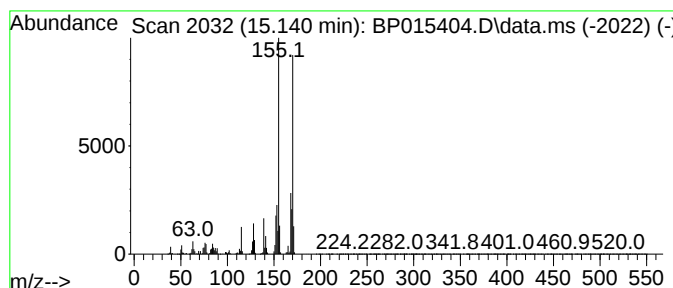
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 8 Naphthalene, 2,3,6-trimethyl- Concentration Rank 8

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 15.140 | 49.10 ng/ul | 6806960 | Acenaphthene-d10 | 14.746 |

| Hit# | of | 5 | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, 2,3,6-trimethyl- | 170 | C13H14  | 000829-26-5 | 98   |
| 2    |    |   | Naphthalene, 2,3,6-trimethyl- | 170 | C13H14  | 000829-26-5 | 96   |
| 3    |    |   | Naphthalene, 1,4,6-trimethyl- | 170 | C13H14  | 002131-42-2 | 95   |
| 4    |    |   | Naphthalene, 2,3,6-trimethyl- | 170 | C13H14  | 000829-26-5 | 95   |
| 5    |    |   | Naphthalene, 1,6,7-trimethyl- | 170 | C13H14  | 002245-38-7 | 95   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP060723\  
Data File : BP015404.D  
Acq On : 07 Jun 2023 18:08  
Operator : MA/JU  
Sample : 02950-01  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
EW978

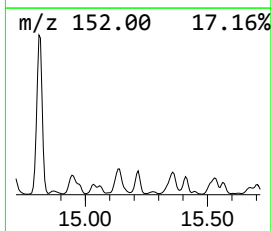
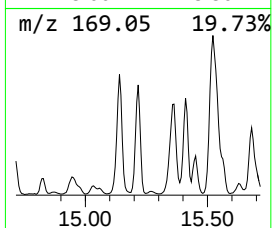
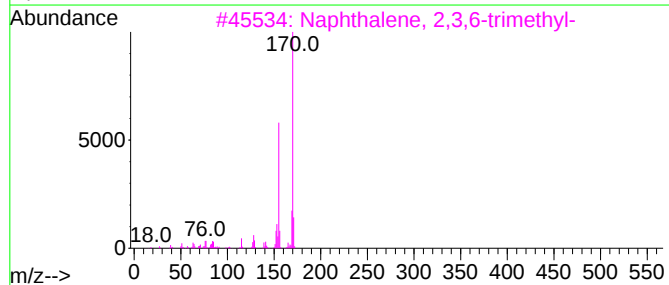
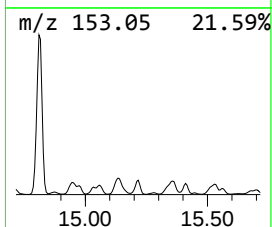
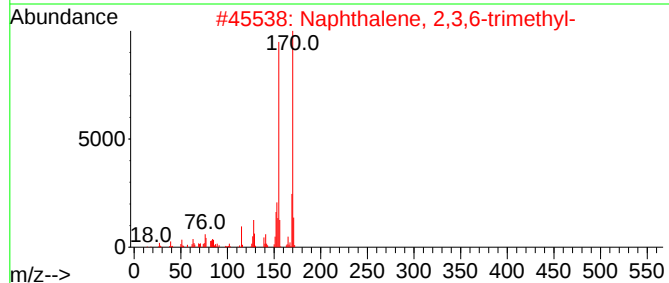
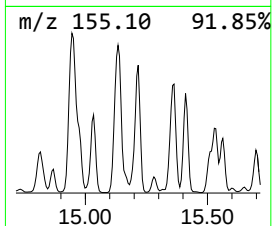
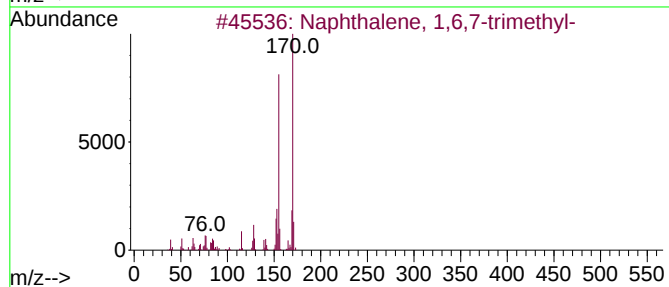
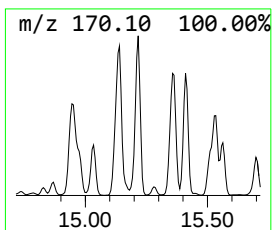
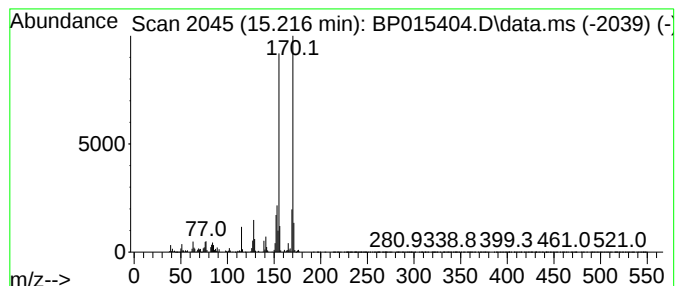
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 9 Naphthalene, 1,6,7-trimethyl- Concentration Rank 16

| R.T.      | EstConc                       | Area    | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------|---------|------------------|-------------|------|
| 15.216    | 29.87 ng/ul                   | 4140100 | Acenaphthene-d10 | 14.746      |      |
| Hit# of 5 | Tentative ID                  | MW      | MolForm          | CAS#        | Qual |
| 1         | Naphthalene, 1,6,7-trimethyl- | 170     | C13H14           | 002245-38-7 | 97   |
| 2         | Naphthalene, 2,3,6-trimethyl- | 170     | C13H14           | 000829-26-5 | 97   |
| 3         | Naphthalene, 2,3,6-trimethyl- | 170     | C13H14           | 000829-26-5 | 97   |
| 4         | Naphthalene, 1,4,6-trimethyl- | 170     | C13H14           | 002131-42-2 | 96   |
| 5         | Naphthalene, 2,3,6-trimethyl- | 170     | C13H14           | 000829-26-5 | 94   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP060723\  
Data File : BP015404.D  
Acq On : 07 Jun 2023 18:08  
Operator : MA/JU  
Sample : 02950-01  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
EW978

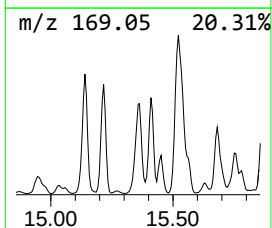
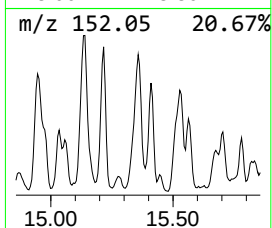
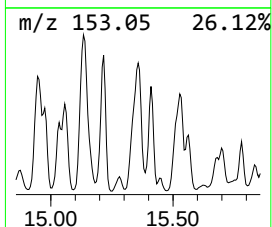
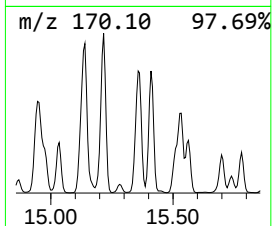
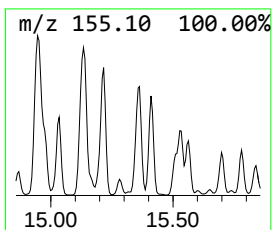
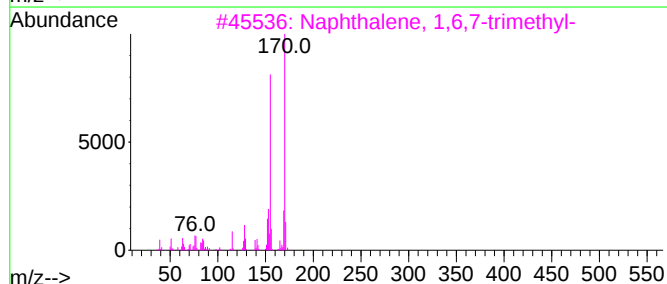
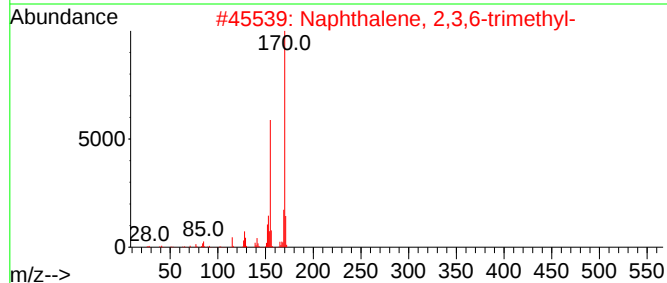
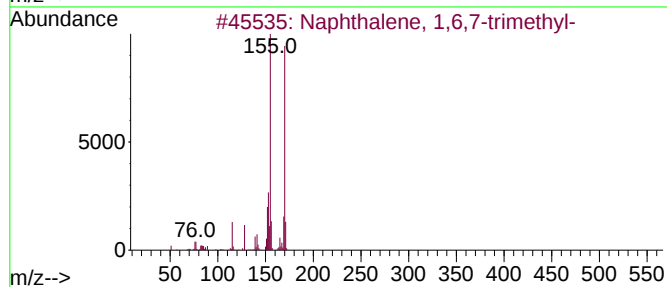
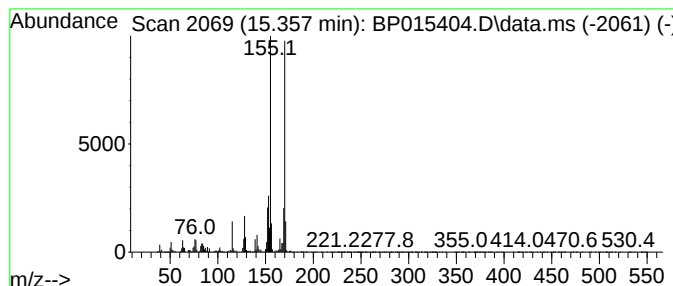
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 10 unknown-01 Concentration Rank 13

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 15.357 | 36.47 ng/ul | 5055370 | Acenaphthene-d10 | 14.746 |

| Hit# | of 5 | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|------|-------------------------------|-----|---------|-------------|------|
| 1    |      | Naphthalene, 1,6,7-trimethyl- | 170 | C13H14  | 002245-38-7 | 98   |
| 2    |      | Naphthalene, 2,3,6-trimethyl- | 170 | C13H14  | 000829-26-5 | 97   |
| 3    |      | Naphthalene, 1,6,7-trimethyl- | 170 | C13H14  | 002245-38-7 | 97   |
| 4    |      | Naphthalene, 2,3,6-trimethyl- | 170 | C13H14  | 000829-26-5 | 96   |
| 5    |      | Naphthalene, 2,3,6-trimethyl- | 170 | C13H14  | 000829-26-5 | 96   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP060723\  
Data File : BP015404.D  
Acq On : 07 Jun 2023 18:08  
Operator : MA/JU  
Sample : 02950-01  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
EW978

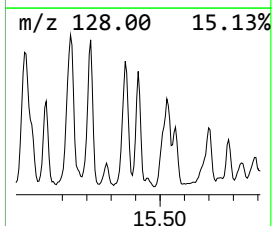
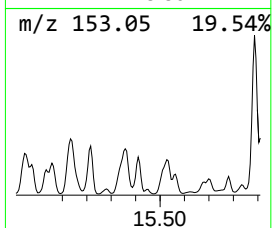
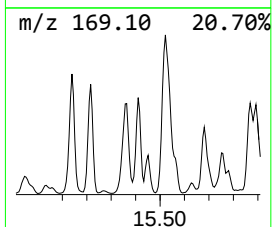
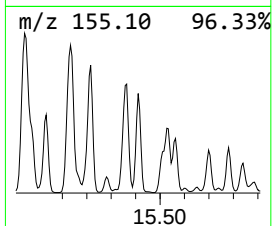
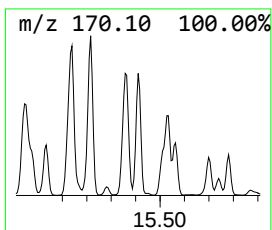
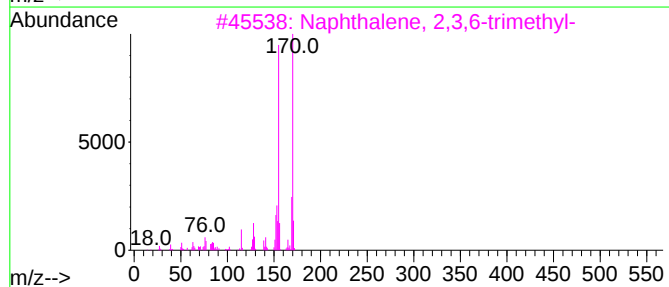
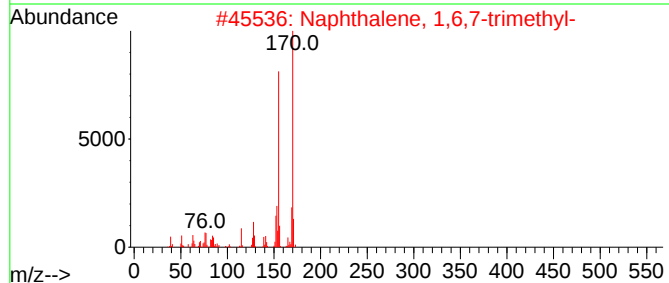
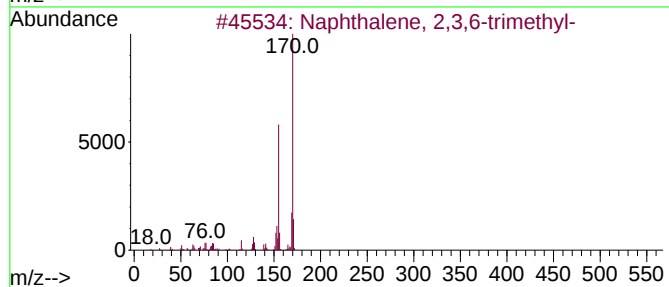
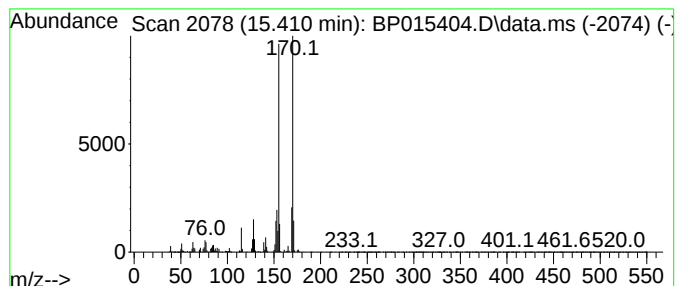
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 11 unknown-02 Concentration Rank 24

| R.T.      | EstConc                       | Area    | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------|---------|------------------|-------------|------|
| 15.410    | 21.15 ng/ul                   | 2931710 | Acenaphthene-d10 | 14.746      |      |
| Hit# of 5 | Tentative ID                  | MW      | MolForm          | CAS#        | Qual |
| 1         | Naphthalene, 2,3,6-trimethyl- | 170     | C13H14           | 000829-26-5 | 97   |
| 2         | Naphthalene, 1,6,7-trimethyl- | 170     | C13H14           | 002245-38-7 | 97   |
| 3         | Naphthalene, 2,3,6-trimethyl- | 170     | C13H14           | 000829-26-5 | 97   |
| 4         | Naphthalene, 2,3,6-trimethyl- | 170     | C13H14           | 000829-26-5 | 97   |
| 5         | Naphthalene, 1,4,6-trimethyl- | 170     | C13H14           | 002131-42-2 | 96   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP060723\  
Data File : BP015404.D  
Acq On : 07 Jun 2023 18:08  
Operator : MA/JU  
Sample : 02950-01  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

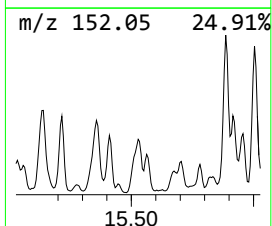
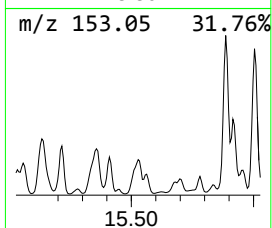
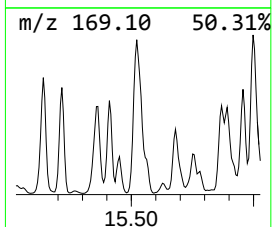
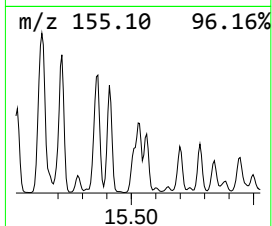
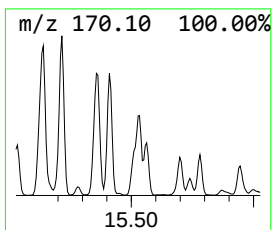
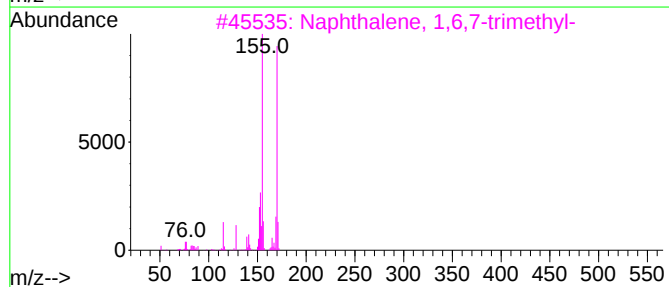
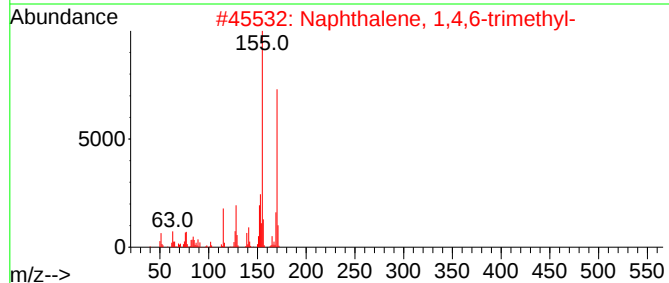
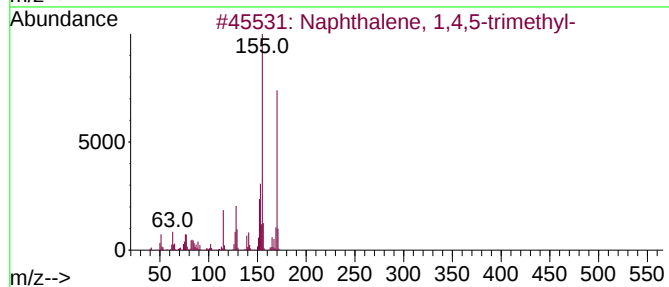
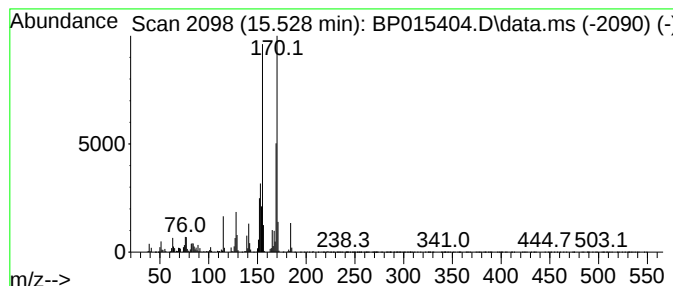
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BNA\_P  
ClientSampleId :  
EW978

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP060723.MA.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 12 Naphthalene, 1,4,5-trimethyl- Concentration Rank 15

| R.T.      | EstConc                       | Area    | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------|---------|------------------|-------------|------|
| 15.528    | 30.91 ng/ul                   | 4284830 | Acenaphthene-d10 | 14.746      |      |
| Hit# of 5 | Tentative ID                  | MW      | MolForm          | CAS#        | Qual |
| 1         | Naphthalene, 1,4,5-trimethyl- | 170     | C13H14           | 002131-41-1 | 95   |
| 2         | Naphthalene, 1,4,6-trimethyl- | 170     | C13H14           | 002131-42-2 | 94   |
| 3         | Naphthalene, 1,6,7-trimethyl- | 170     | C13H14           | 002245-38-7 | 94   |
| 4         | 4,6,8-Trimethylazulene        | 170     | C13H14           | 000941-81-1 | 90   |
| 5         | Naphthalene, 1,6,7-trimethyl- | 170     | C13H14           | 002245-38-7 | 89   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP060723\  
Data File : BP015404.D  
Acq On : 07 Jun 2023 18:08  
Operator : MA/JU  
Sample : 02950-01  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

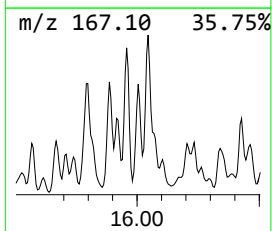
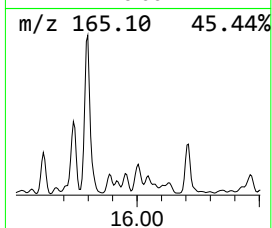
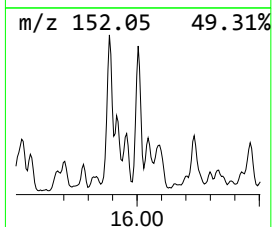
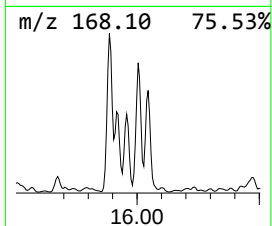
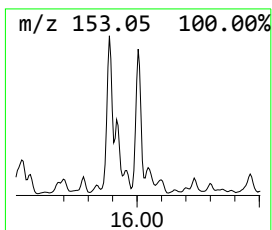
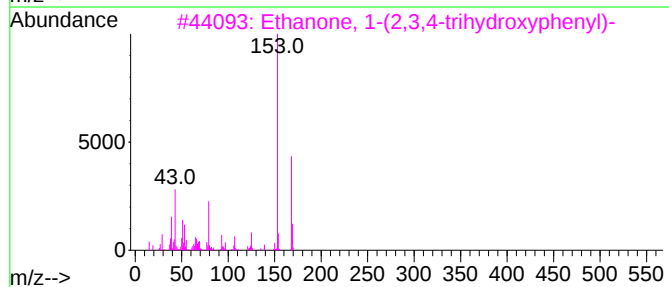
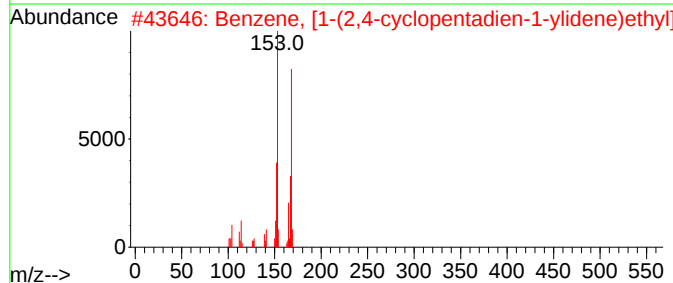
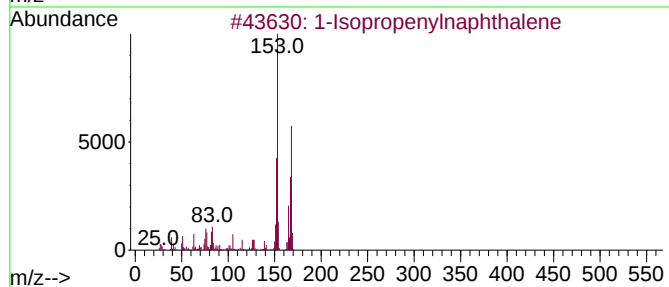
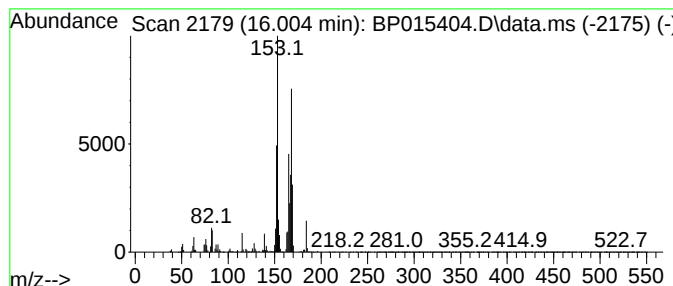
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ClientSampleId :  
EW978

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP060723.MA.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 17 1-Isopropenylnaphthalene Concentration Rank 21

| R.T.      | EstConc                             | Area    | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------------|---------|------------------|-------------|------|
| 16.004    | 22.74 ng/ul                         | 3152920 | Acenaphthene-d10 | 14.746      |      |
| Hit# of 5 | Tentative ID                        | MW      | MolForm          | CAS#        | Qual |
| 1         | 1-Isopropenylnaphthalene            | 168     | C13H12           | 001855-47-6 | 62   |
| 2         | Benzene, [1-(2,4-cyclopentadien-... | 168     | C13H12           | 002320-32-3 | 49   |
| 3         | Ethanone, 1-(2,3,4-trihydroxyphe... | 168     | C8H8O4           | 000528-21-2 | 43   |
| 4         | Nordiphenamid                       | 225     | C15H15NO         | 000954-21-2 | 38   |
| 5         | 1,1'-Biphenyl, 3-methyl-            | 168     | C13H12           | 000643-93-6 | 38   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP060723\  
Data File : BP015404.D  
Acq On : 07 Jun 2023 18:08  
Operator : MA/JU  
Sample : 02950-01  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
EW978

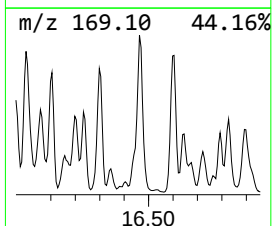
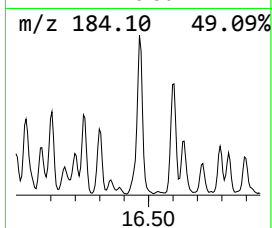
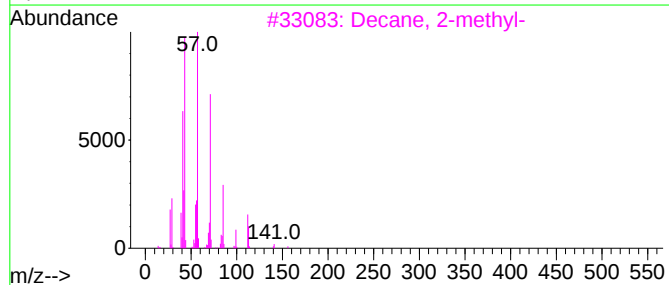
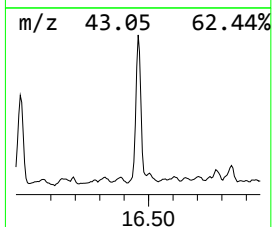
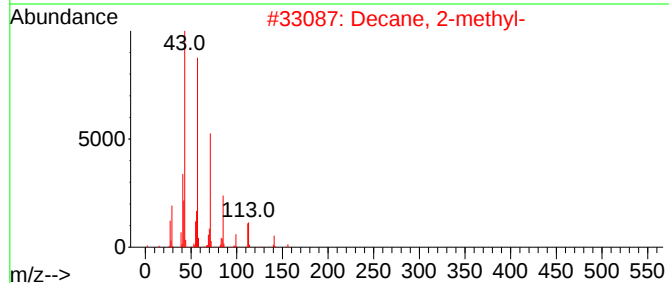
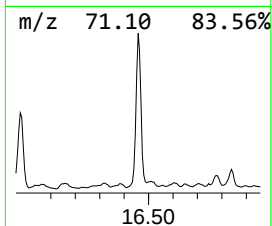
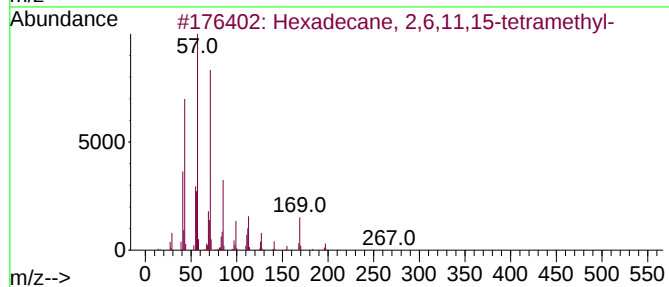
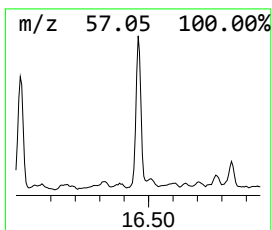
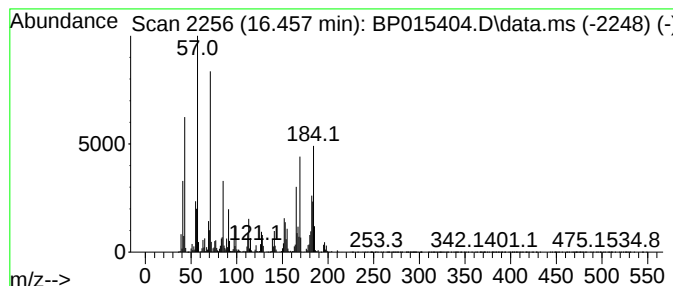
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 21 (DEL) Alkane: Straight-Chai... Concentration Rank 10

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 16.457 | 44.61 ng/ul | 6995630 | Phenanthrene-d10 | 17.504 |

| Hit# | of | 5 | Tentative ID                       | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Hexadecane, 2,6,11,15-tetramethyl- | 282 | C20H42  | 000504-44-9 | 42   |
| 2    |    |   | Decane, 2-methyl-                  | 156 | C11H24  | 006975-98-0 | 38   |
| 3    |    |   | Decane, 2-methyl-                  | 156 | C11H24  | 006975-98-0 | 38   |
| 4    |    |   | Hexadecane, 7,9-dimethyl-          | 254 | C18H38  | 021164-95-4 | 30   |
| 5    |    |   | Hexadecane, 2,6,10,14-tetramethyl- | 282 | C20H42  | 000638-36-8 | 30   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP060723\  
Data File : BP015404.D  
Acq On : 07 Jun 2023 18:08  
Operator : MA/JU  
Sample : 02950-01  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
EW978

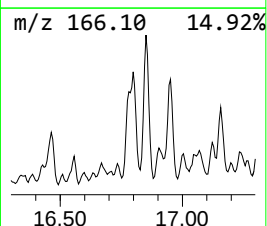
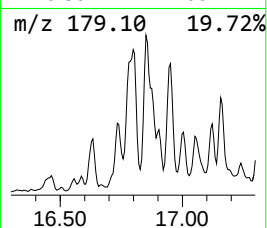
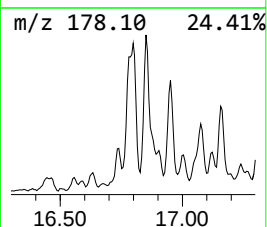
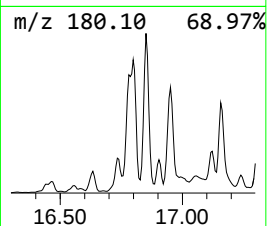
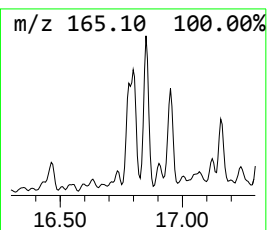
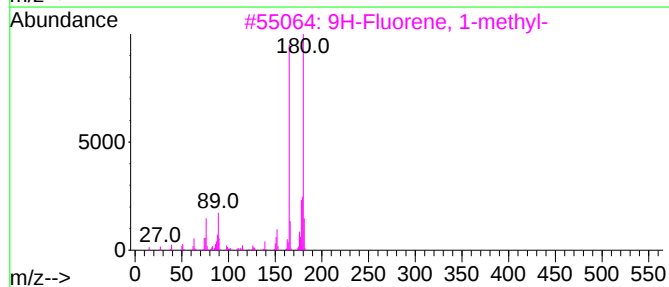
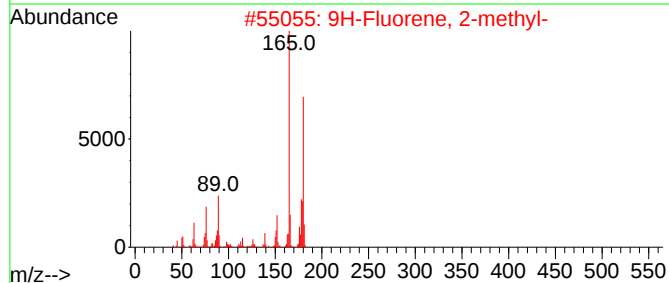
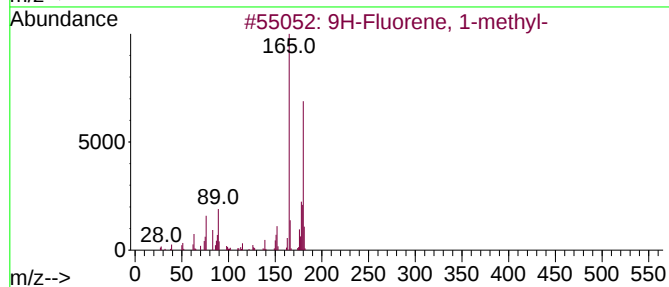
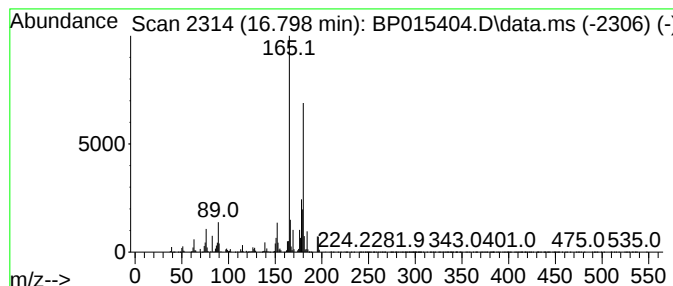
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 23 9H-Fluorene, 1-methyl- Concentration Rank 11

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 16.798 | 39.94 ng/ul | 6263140 | Phenanthrene-d10 | 17.504 |

| Hit# | of 5                   | Tentative ID | MW     | MolForm     | CAS# | Qual |
|------|------------------------|--------------|--------|-------------|------|------|
| 1    | 9H-Fluorene, 1-methyl- | 180          | C14H12 | 001730-37-6 | 95   |      |
| 2    | 9H-Fluorene, 2-methyl- | 180          | C14H12 | 001430-97-3 | 95   |      |
| 3    | 9H-Fluorene, 1-methyl- | 180          | C14H12 | 001730-37-6 | 94   |      |
| 4    | 9H-Fluorene, 1-methyl- | 180          | C14H12 | 001730-37-6 | 93   |      |
| 5    | 9H-Fluorene, 2-methyl- | 180          | C14H12 | 001430-97-3 | 93   |      |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP060723\  
Data File : BP015404.D  
Acq On : 07 Jun 2023 18:08  
Operator : MA/JU  
Sample : 02950-01  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
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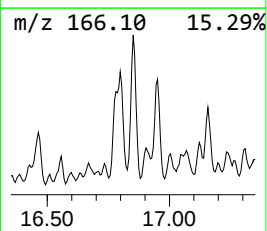
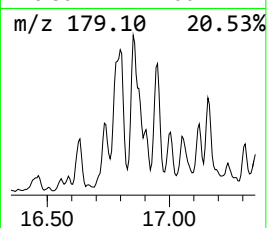
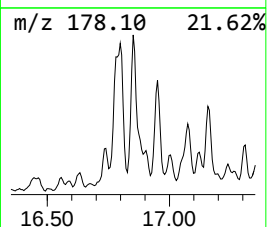
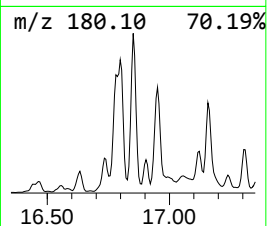
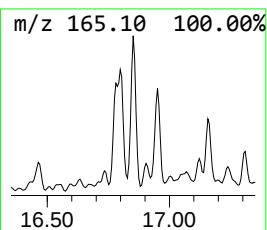
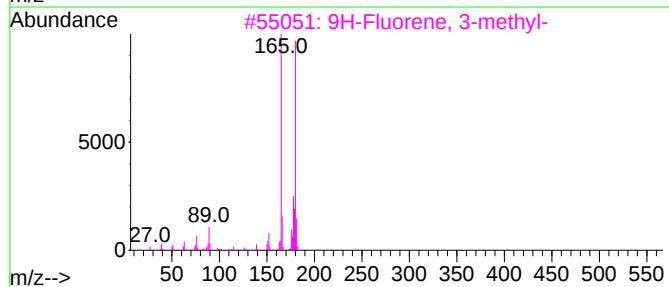
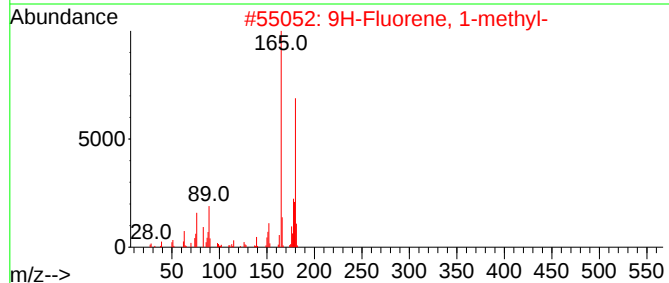
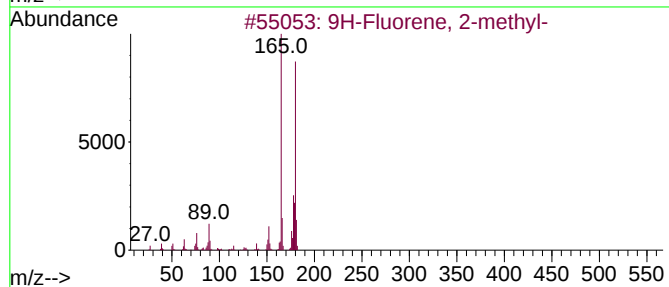
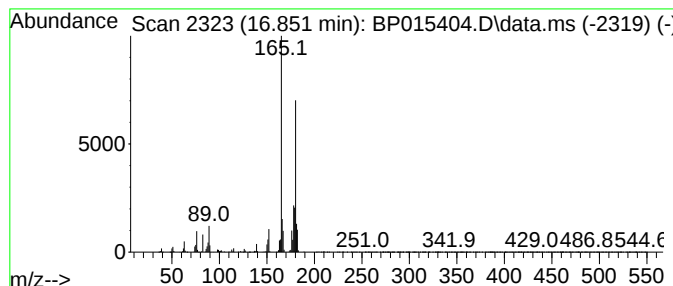
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 24 9H-Fluorene, 2-methyl- Concentration Rank 17

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 16.851 | 28.13 ng/ul | 4410720 | Phenanthrene-d10 | 17.504 |

| Hit# | of | 5 | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------------|-----|---------|-------------|------|
| 1    |    |   | 9H-Fluorene, 2-methyl- | 180 | C14H12  | 001430-97-3 | 97   |
| 2    |    |   | 9H-Fluorene, 1-methyl- | 180 | C14H12  | 001730-37-6 | 96   |
| 3    |    |   | 9H-Fluorene, 3-methyl- | 180 | C14H12  | 002523-39-9 | 95   |
| 4    |    |   | 9H-Fluorene, 1-methyl- | 180 | C14H12  | 001730-37-6 | 95   |
| 5    |    |   | 9H-Fluorene, 1-methyl- | 180 | C14H12  | 001730-37-6 | 92   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP060723\  
Data File : BP015404.D  
Acq On : 07 Jun 2023 18:08  
Operator : MA/JU  
Sample : 02950-01  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
EW978

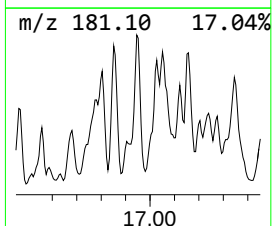
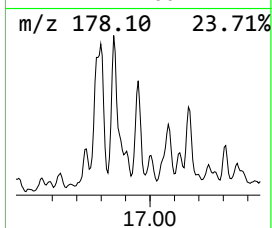
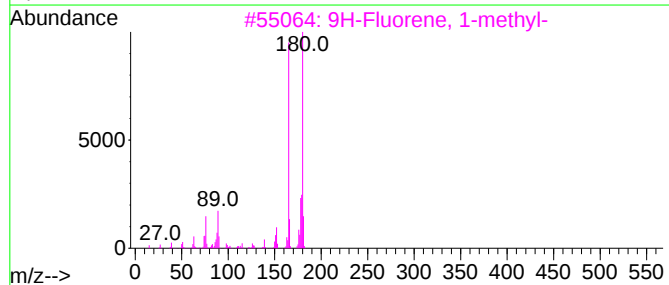
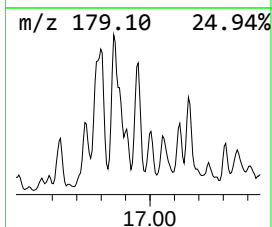
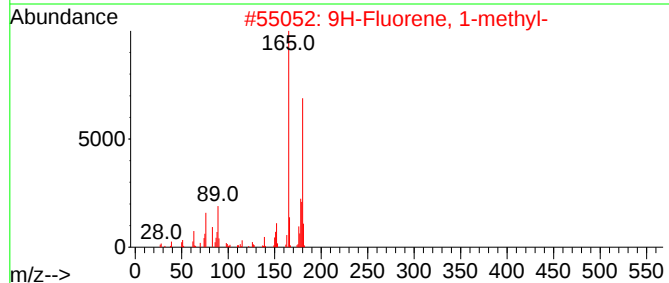
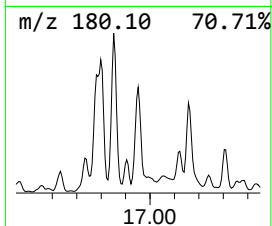
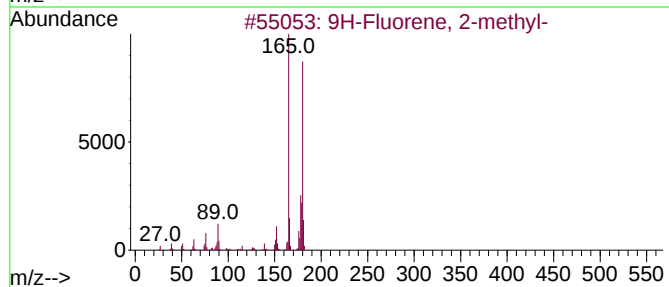
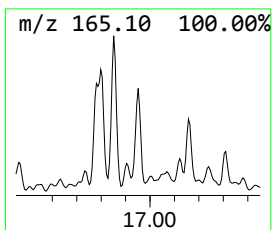
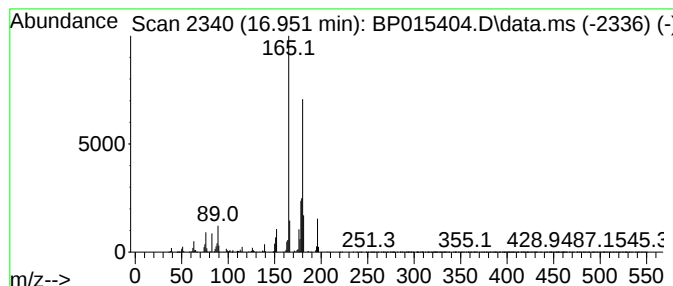
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 26 9H-Fluorene, 4-methyl- Concentration Rank 27

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 16.951 | 17.94 ng/ul | 2813370 | Phenanthrene-d10 | 17.504 |

| Hit# | of 5                   | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|------------------------|--------------|-----|---------|-------------|------|
| 1    | 9H-Fluorene, 2-methyl- |              | 180 | C14H12  | 001430-97-3 | 95   |
| 2    | 9H-Fluorene, 1-methyl- |              | 180 | C14H12  | 001730-37-6 | 95   |
| 3    | 9H-Fluorene, 1-methyl- |              | 180 | C14H12  | 001730-37-6 | 91   |
| 4    | 9H-Fluorene, 4-methyl- |              | 180 | C14H12  | 001556-99-6 | 90   |
| 5    | 9H-Fluorene, 3-methyl- |              | 180 | C14H12  | 002523-39-9 | 76   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP060723\  
Data File : BP015404.D  
Acq On : 07 Jun 2023 18:08  
Operator : MA/JU  
Sample : 02950-01  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
EW978

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP060723.MA.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

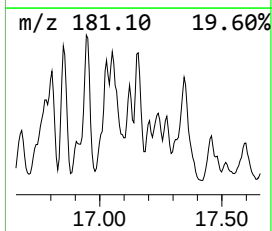
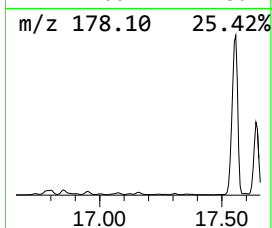
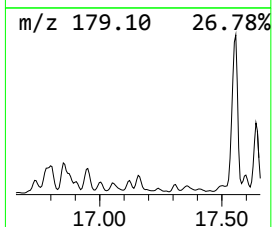
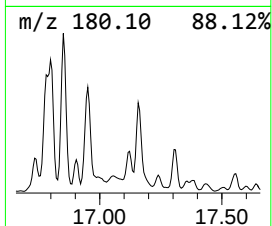
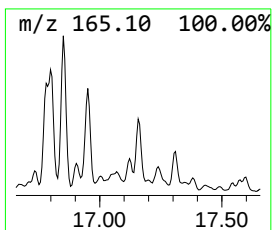
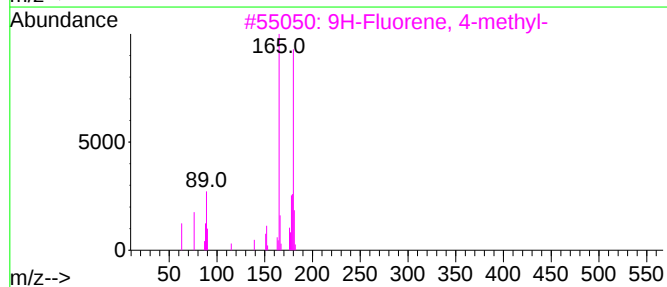
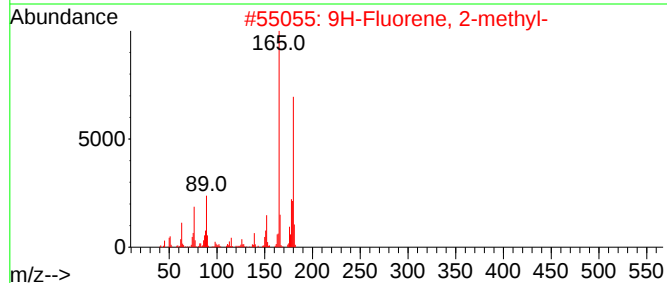
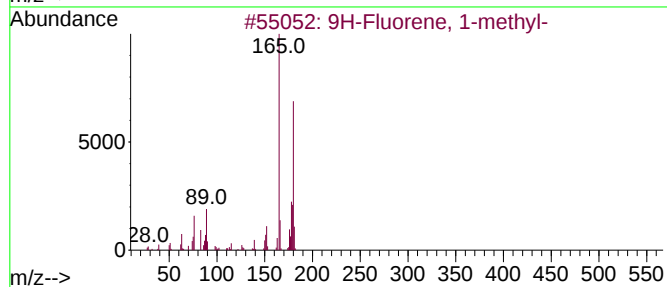
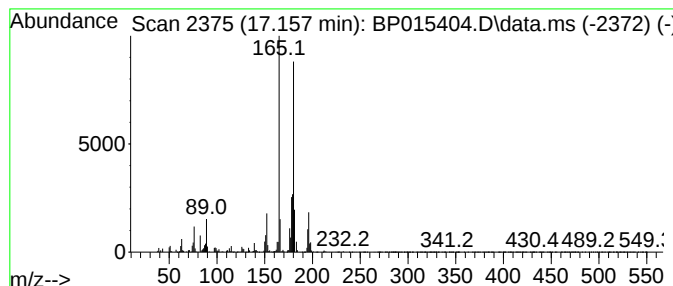
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Peak Number 27 unknown-03

Concentration Rank 35

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 17.157 | 13.54 ng/ul | 2122780 | Phenanthrene-d10 | 17.504 |

| Hit# | of                     | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|------------------------|---|--------------|-----|---------|-------------|------|
| 1    | 9H-Fluorene, 1-methyl- |   |              | 180 | C14H12  | 001730-37-6 | 94   |
| 2    | 9H-Fluorene, 2-methyl- |   |              | 180 | C14H12  | 001430-97-3 | 94   |
| 3    | 9H-Fluorene, 4-methyl- |   |              | 180 | C14H12  | 001556-99-6 | 90   |
| 4    | 9H-Fluorene, 2-methyl- |   |              | 180 | C14H12  | 001430-97-3 | 90   |
| 5    | 9H-Fluorene, 1-methyl- |   |              | 180 | C14H12  | 001730-37-6 | 90   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP060723\  
Data File : BP015404.D  
Acq On : 07 Jun 2023 18:08  
Operator : MA/JU  
Sample : 02950-01  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
EW978

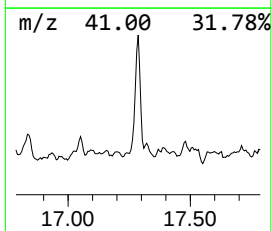
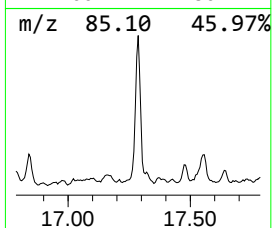
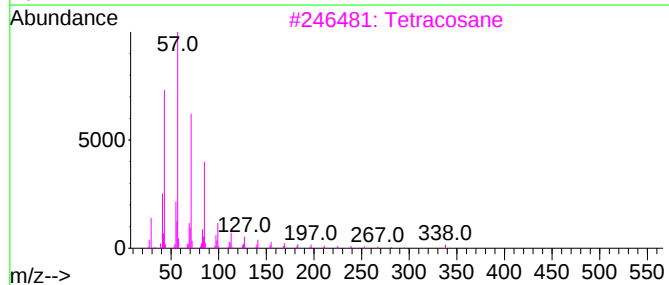
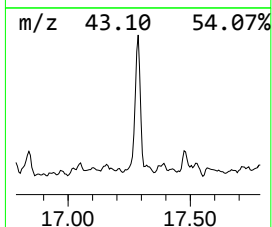
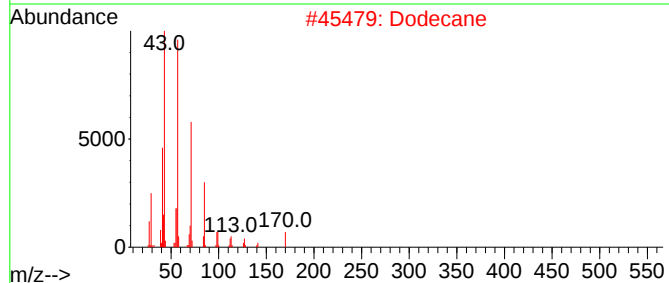
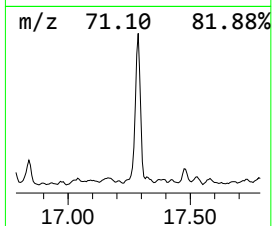
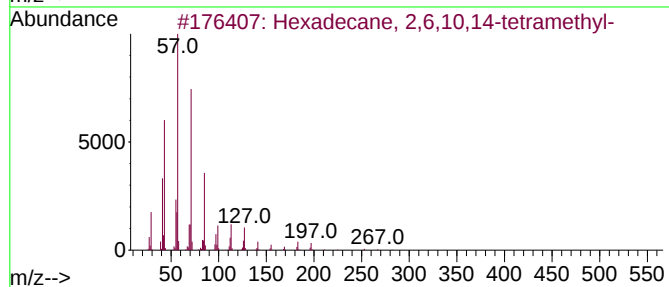
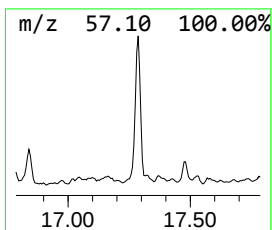
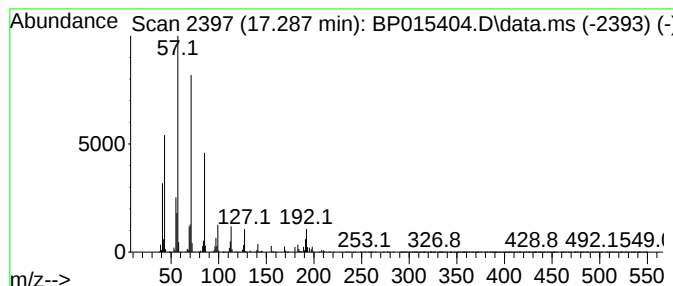
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 29 (DEL) Alkane: Straight-Chai... Concentration Rank 39

| R.T.      | EstConc                            | Area    | Relative to ISTD | R.T.        |      |
|-----------|------------------------------------|---------|------------------|-------------|------|
| 17.287    | 11.48 ng/ul                        | 1800620 | Phenanthrene-d10 | 17.504      |      |
| Hit# of 5 | Tentative ID                       | MW      | MolForm          | CAS#        | Qual |
| 1         | Hexadecane, 2,6,10,14-tetramethyl- | 282     | C20H42           | 000638-36-8 | 96   |
| 2         | Dodecane                           | 170     | C12H26           | 000112-40-3 | 90   |
| 3         | Tetracosane                        | 338     | C24H50           | 000646-31-1 | 86   |
| 4         | Octadecane                         | 254     | C18H38           | 000593-45-3 | 86   |
| 5         | Hexadecane, 1-iodo-                | 352     | C16H33I          | 000544-77-4 | 86   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP060723\  
Data File : BP015404.D  
Acq On : 07 Jun 2023 18:08  
Operator : MA/JU  
Sample : 02950-01  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

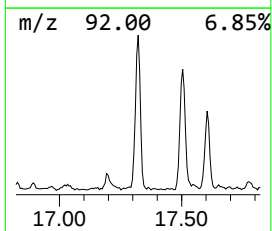
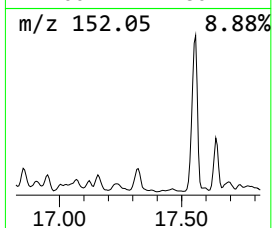
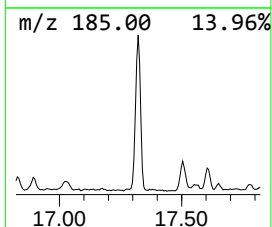
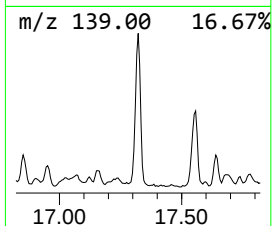
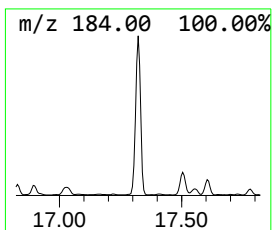
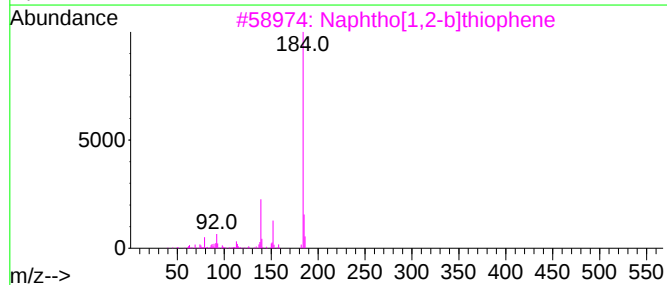
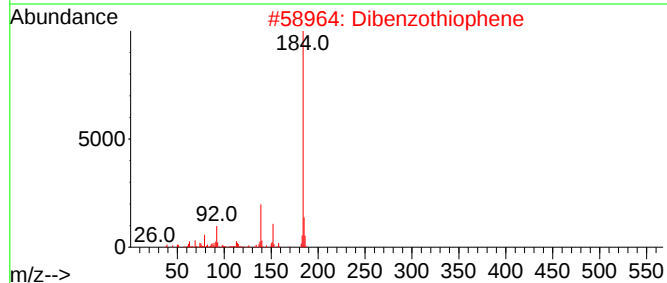
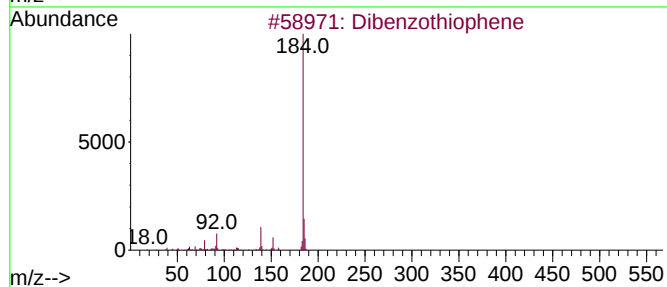
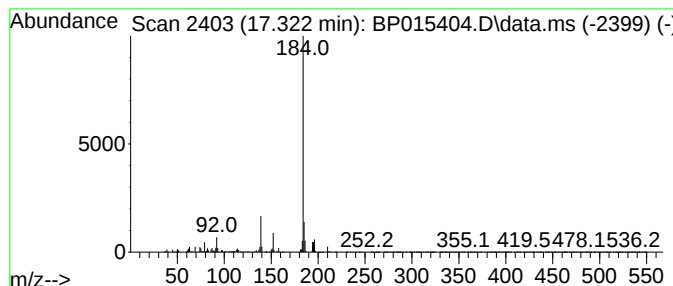
Instrument :  
BNA\_P  
ClientSampleId :  
EW978

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP060723.MA.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 30 Dibenzothiophene Concentration Rank 19

| R.T.      | EstConc                 | Area    | Relative to ISTD |         | R.T.        |      |
|-----------|-------------------------|---------|------------------|---------|-------------|------|
| 17.322    | 24.23 ng/ul             | 3800530 | Phenanthrene-d10 |         | 17.504      |      |
| Hit# of 5 | Tentative ID            |         | MW               | MolForm | CAS#        | Qual |
| 1         | Dibenzothiophene        |         | 184              | C12H8S  | 000132-65-0 | 96   |
| 2         | Dibenzothiophene        |         | 184              | C12H8S  | 000132-65-0 | 95   |
| 3         | Naphtho[1,2-b]thiophene |         | 184              | C12H8S  | 000234-41-3 | 93   |
| 4         | Dibenzothiophene        |         | 184              | C12H8S  | 000132-65-0 | 91   |
| 5         | Dibenzothiophene        |         | 184              | C12H8S  | 000132-65-0 | 91   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP060723\  
Data File : BP015404.D  
Acq On : 07 Jun 2023 18:08  
Operator : MA/JU  
Sample : 02950-01  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
EW978

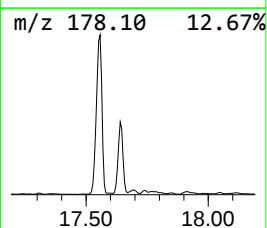
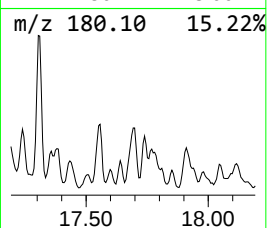
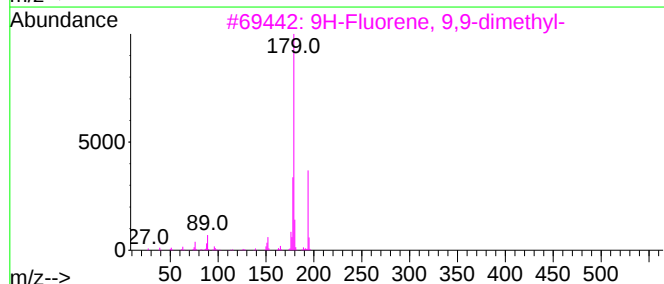
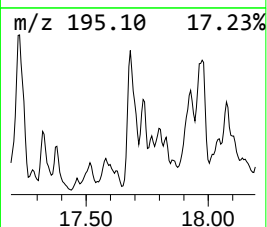
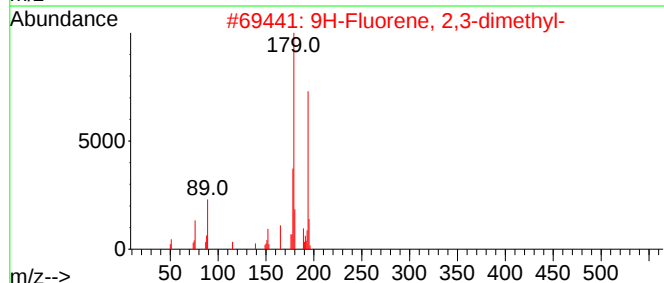
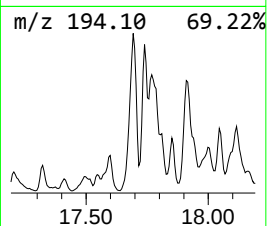
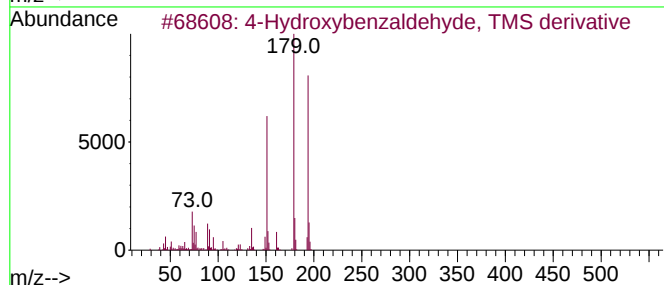
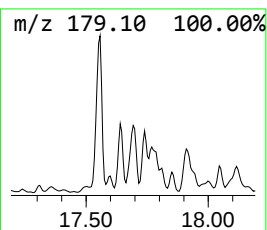
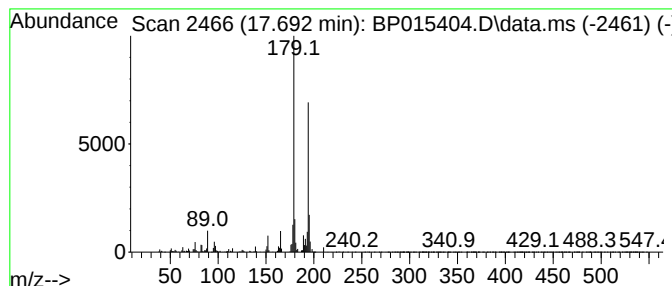
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 31 4-Hydroxybenzaldehyde, TMS ... Concentration Rank 30

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 17.692 | 16.88 ng/ul | 2647540 | Phenanthrene-d10 | 17.504 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm    | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|------------|-------------|------|
| 1    |    |   | 4-Hydroxybenzaldehyde, TMS deriv... | 194 | C10H14O2Si | 001012-12-0 | 80   |
| 2    |    |   | 9H-Fluorene, 2,3-dimethyl-          | 194 | C15H14     | 004612-63-9 | 78   |
| 3    |    |   | 9H-Fluorene, 9,9-dimethyl-          | 194 | C15H14     | 004569-45-3 | 64   |
| 4    |    |   | 4-Acetyl-7-hydroxybenzo-2,1,3-th... | 194 | C8H6N2O2S  | 120893-38-1 | 64   |
| 5    |    |   | Benzene, 1-ethyl-4-[(trimethylsi... | 194 | C11H18OSi  | 017993-90-7 | 59   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP060723\  
Data File : BP015404.D  
Acq On : 07 Jun 2023 18:08  
Operator : MA/JU  
Sample : 02950-01  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

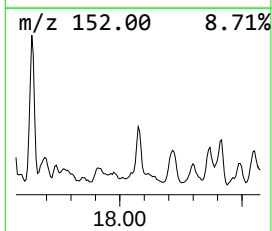
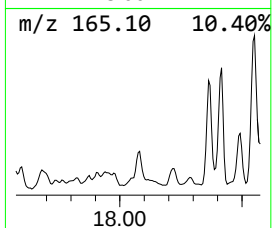
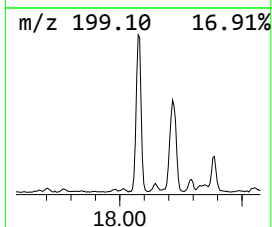
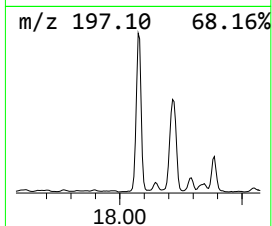
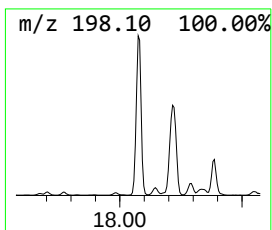
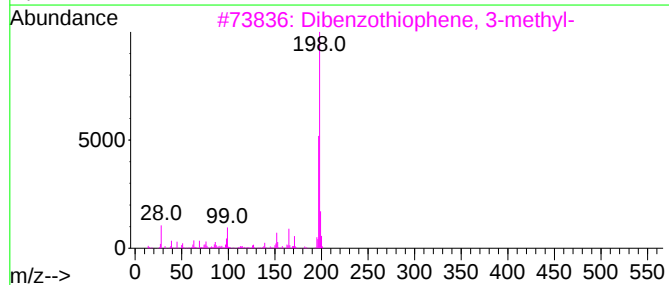
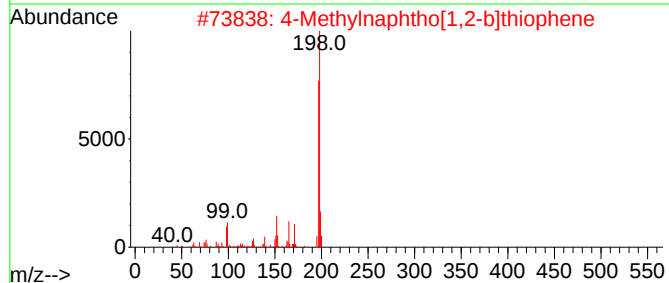
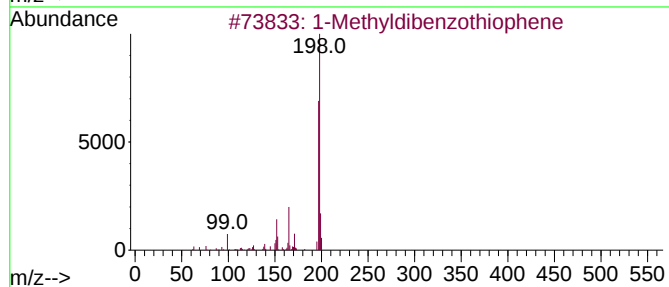
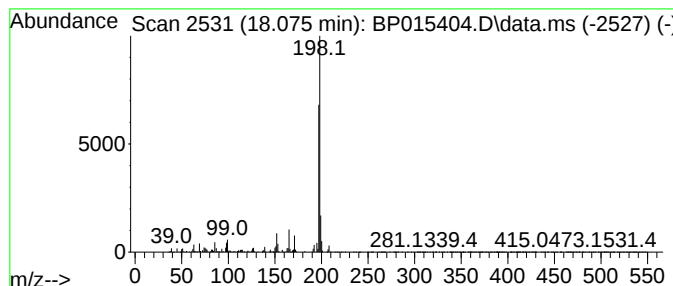
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BNA\_P  
ClientSampleId :  
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Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP060723.MA.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 35 1-Methyldibenzothiophene Concentration Rank 23

| R.T.      | EstConc                         | Area    | Relative to ISTD | R.T.         |      |
|-----------|---------------------------------|---------|------------------|--------------|------|
| 18.075    | 21.17 ng/ul                     | 3320340 | Phenanthrene-d10 | 17.504       |      |
| Hit# of 5 | Tentative ID                    | MW      | MolForm          | CAS#         | Qual |
| 1         | 1-Methyldibenzothiophene        | 198     | C13H10S          | 031317-07-4  | 97   |
| 2         | 4-Methylnaphtho[1,2-b]thiophene | 198     | C13H10S          | 067388-11-8  | 96   |
| 3         | Dibenzothiophene, 3-methyl-     | 198     | C13H10S          | 016587-52-3  | 93   |
| 4         | Dibenzothiophene, 4-methyl-     | 198     | C13H10S          | 007372-88-5  | 93   |
| 5         | 2-Methyldibenzothiophene        | 198     | C13H10S          | 1000383-55-1 | 91   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP060723\  
Data File : BP015404.D  
Acq On : 07 Jun 2023 18:08  
Operator : MA/JU  
Sample : 02950-01  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
EW978

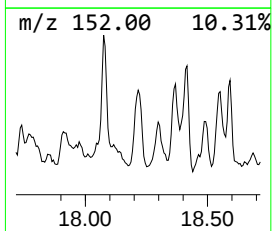
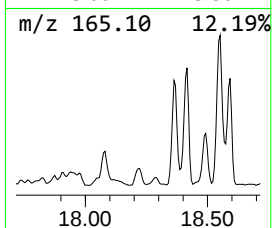
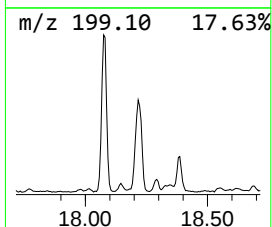
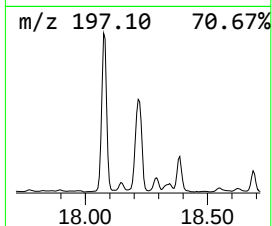
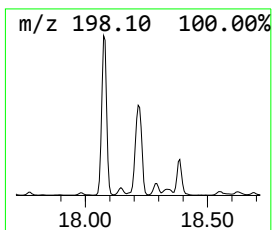
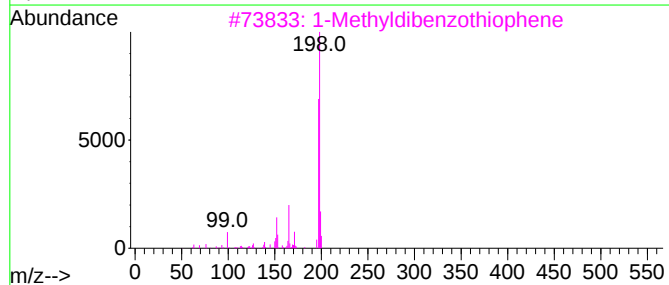
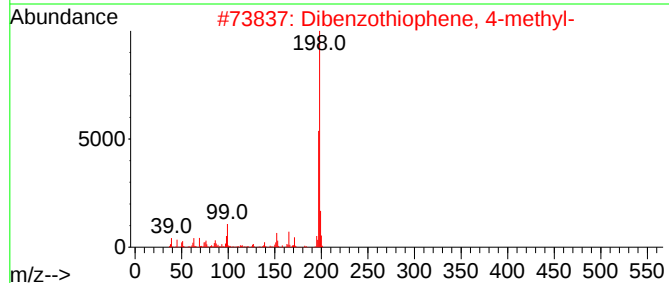
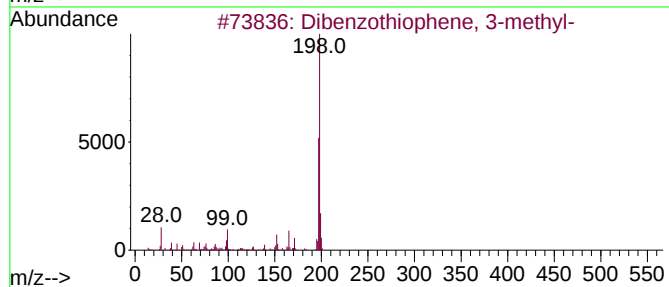
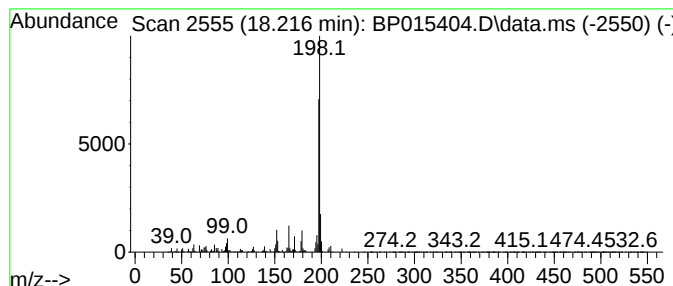
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 36 Dibenzothiophene, 3-methyl- Concentration Rank 28

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 18.216 | 17.53 ng/ul | 2749200 | Phenanthrene-d10 | 17.504 |

| Hit# | of | 5 | Tentative ID                    | MW  | MolForm | CAS#         | Qual |
|------|----|---|---------------------------------|-----|---------|--------------|------|
| 1    |    |   | Dibenzothiophene, 3-methyl-     | 198 | C13H10S | 016587-52-3  | 95   |
| 2    |    |   | Dibenzothiophene, 4-methyl-     | 198 | C13H10S | 007372-88-5  | 95   |
| 3    |    |   | 1-Methyldibenzothiophene        | 198 | C13H10S | 031317-07-4  | 94   |
| 4    |    |   | 4-Methylnaphtho[1,2-b]thiophene | 198 | C13H10S | 067388-11-8  | 93   |
| 5    |    |   | 2-Methyldibenzothiophene        | 198 | C13H10S | 1000383-55-1 | 90   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP060723\  
Data File : BP015404.D  
Acq On : 07 Jun 2023 18:08  
Operator : MA/JU  
Sample : 02950-01  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
EW978

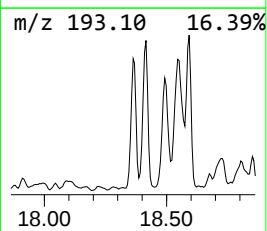
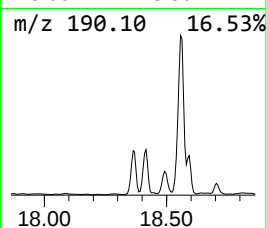
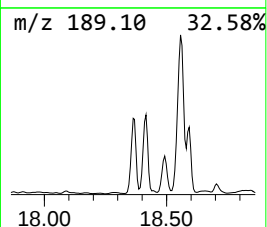
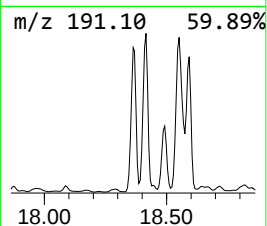
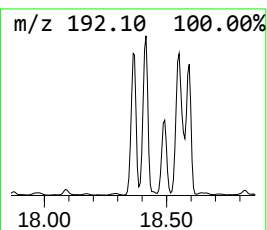
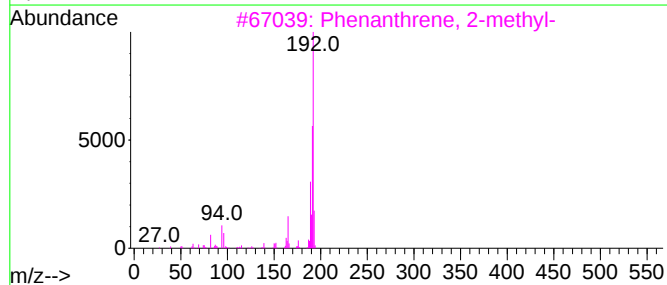
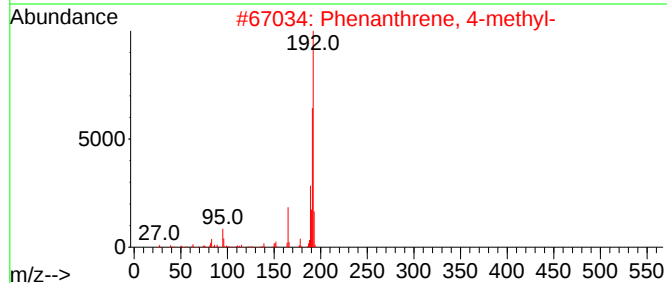
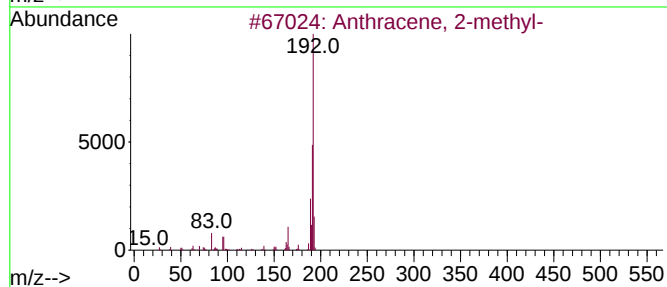
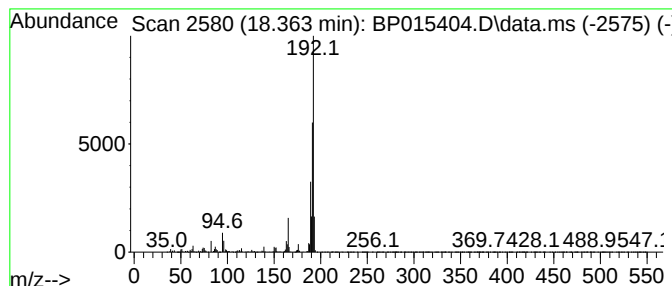
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 37 Anthracene, 2-methyl- Concentration Rank 4

| R.T.   | EstConc     | Area     | Relative to ISTD | R.T.   |
|--------|-------------|----------|------------------|--------|
| 18.363 | 65.78 ng/ul | 10316500 | Phenanthrene-d10 | 17.504 |

| Hit# | of | 5 | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------|-----|---------|-------------|------|
| 1    |    |   | Anthracene, 2-methyl-   | 192 | C15H12  | 000613-12-7 | 95   |
| 2    |    |   | Phenanthrene, 4-methyl- | 192 | C15H12  | 000832-64-4 | 95   |
| 3    |    |   | Phenanthrene, 2-methyl- | 192 | C15H12  | 002531-84-2 | 94   |
| 4    |    |   | Phenanthrene, 2-methyl- | 192 | C15H12  | 002531-84-2 | 93   |
| 5    |    |   | Phenanthrene, 1-methyl- | 192 | C15H12  | 000832-69-9 | 93   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP060723\  
Data File : BP015404.D  
Acq On : 07 Jun 2023 18:08  
Operator : MA/JU  
Sample : 02950-01  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
EW978

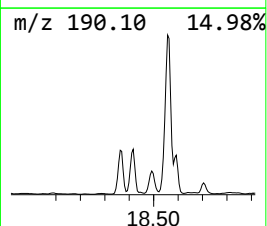
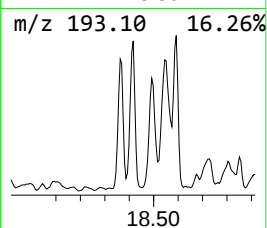
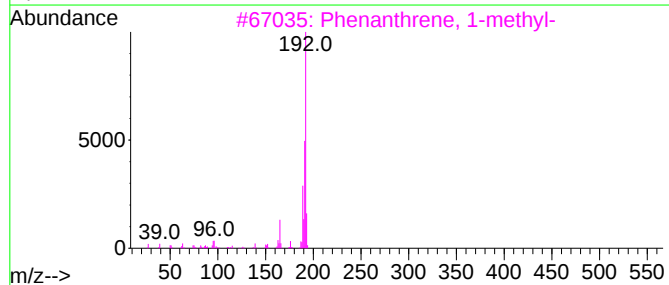
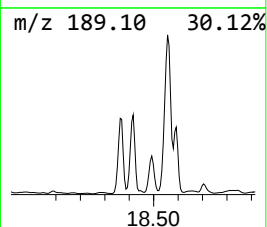
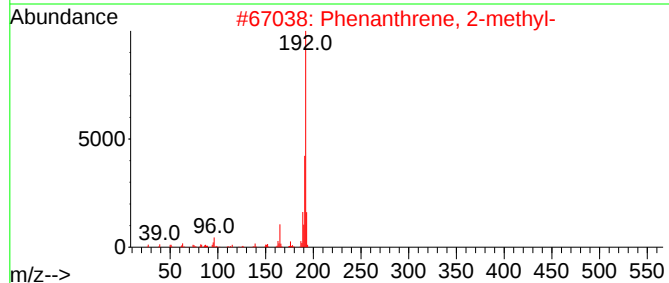
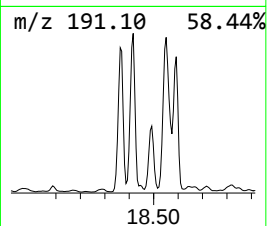
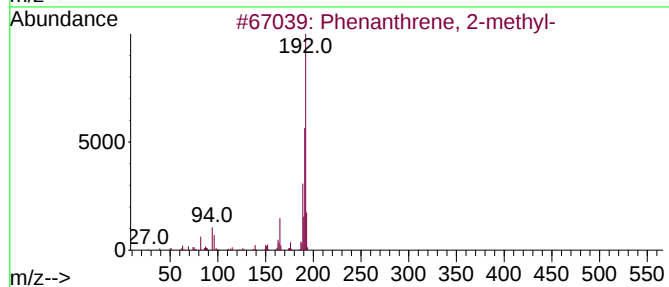
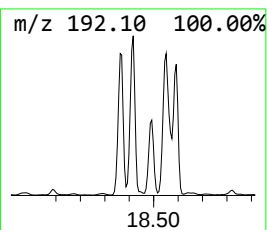
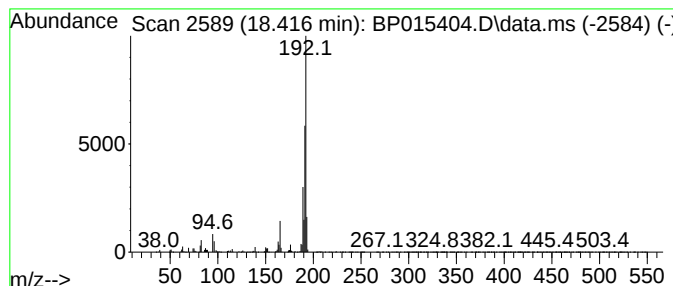
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 38 Phenanthrene, 2-methyl- Concentration Rank 3

| R.T.   | EstConc     | Area     | Relative to ISTD | R.T.   |
|--------|-------------|----------|------------------|--------|
| 18.416 | 66.95 ng/ul | 10499600 | Phenanthrene-d10 | 17.504 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Phenanthrene, 2-methyl-             | 192 | C15H12  | 002531-84-2 | 98   |
| 2    |    |   | Phenanthrene, 2-methyl-             | 192 | C15H12  | 002531-84-2 | 97   |
| 3    |    |   | Phenanthrene, 1-methyl-             | 192 | C15H12  | 000832-69-9 | 97   |
| 4    |    |   | 1H-Cyclopropa[1]phenanthrene,1a,... | 192 | C15H12  | 000949-41-7 | 96   |
| 5    |    |   | Phenanthrene, 1-methyl-             | 192 | C15H12  | 000832-69-9 | 95   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP060723\  
Data File : BP015404.D  
Acq On : 07 Jun 2023 18:08  
Operator : MA/JU  
Sample : 02950-01  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
EW978

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP060723.MA.M  
Quant Title : SVOA CALIBRATION

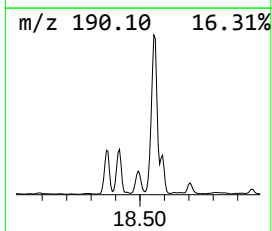
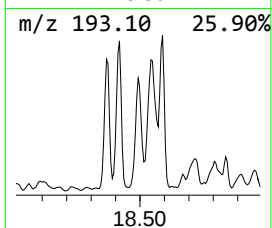
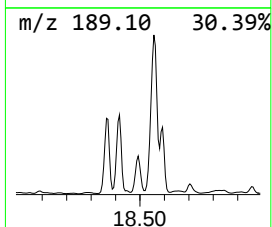
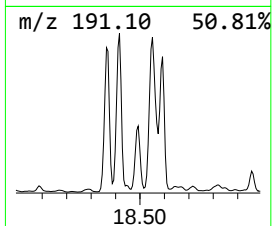
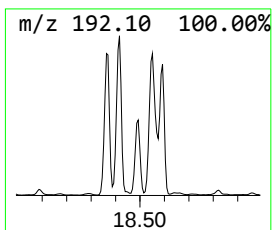
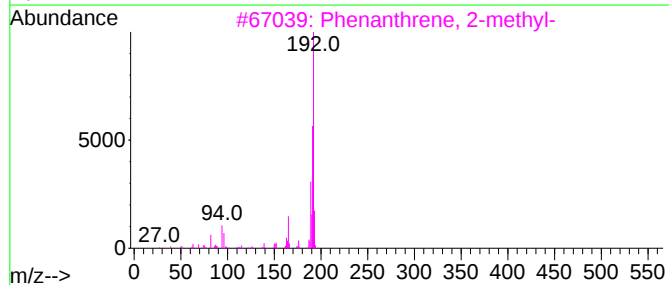
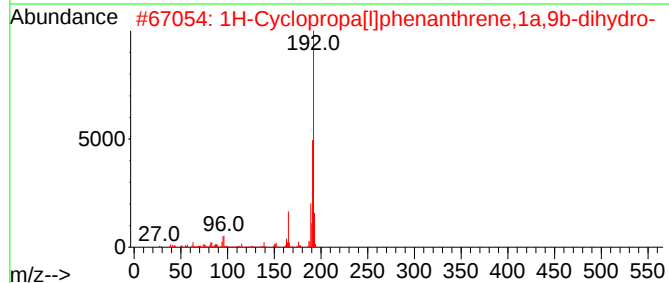
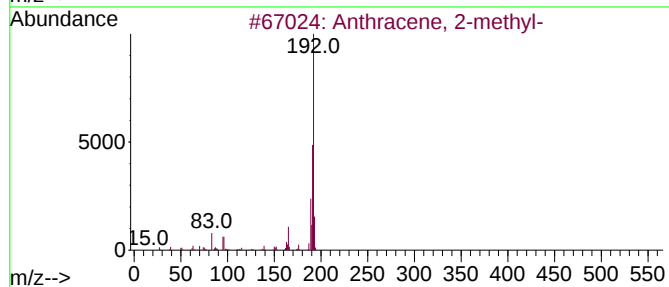
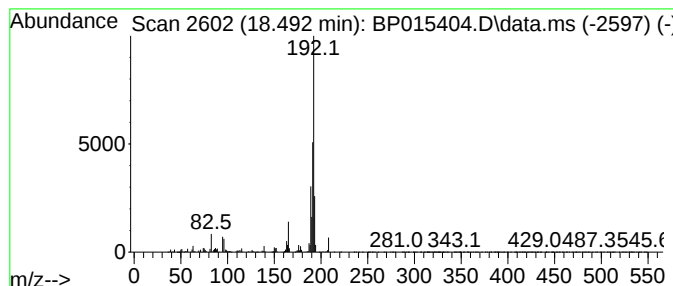
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TIC Integration Parameters: LSCINT.P

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Peak Number 39 1H-Cyclopropa[l]phenanthren... Concentration Rank 14

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 18.492 | 35.34 ng/ul | 5542970 | Phenanthrene-d10 | 17.504 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|------|-------------------------------------|-----|---------|-------------|------|
| 1    |      | Anthracene, 2-methyl-               | 192 | C15H12  | 000613-12-7 | 96   |
| 2    |      | 1H-Cyclopropa[l]phenanthrene,1a,... | 192 | C15H12  | 000949-41-7 | 96   |
| 3    |      | Phenanthrene, 2-methyl-             | 192 | C15H12  | 002531-84-2 | 95   |
| 4    |      | Phenanthrene, 1-methyl-             | 192 | C15H12  | 000832-69-9 | 95   |
| 5    |      | Anthracene, 1-methyl-               | 192 | C15H12  | 000610-48-0 | 95   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP060723\  
Data File : BP015404.D  
Acq On : 07 Jun 2023 18:08  
Operator : MA/JU  
Sample : 02950-01  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
EW978

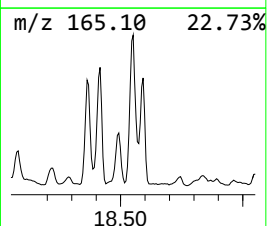
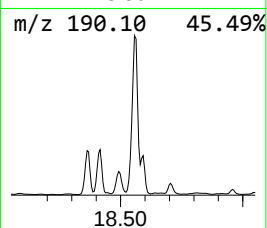
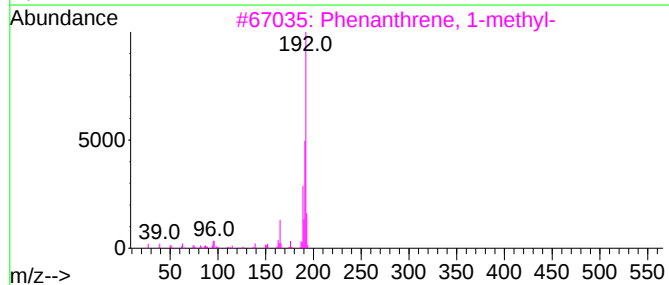
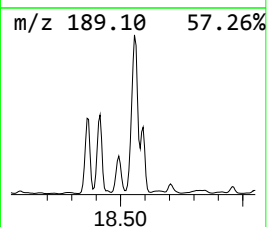
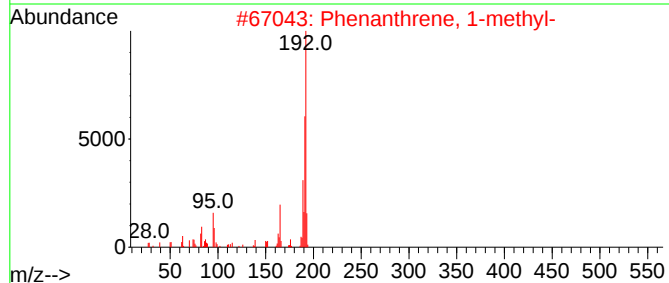
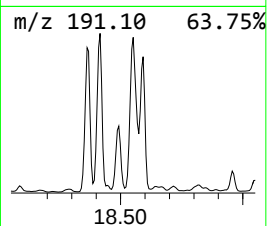
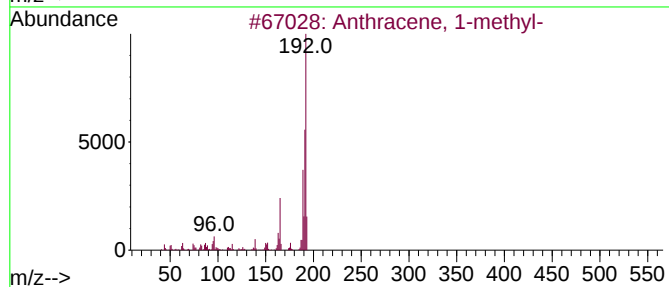
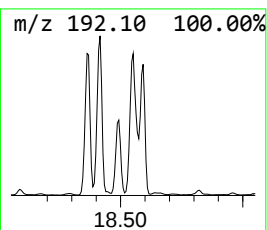
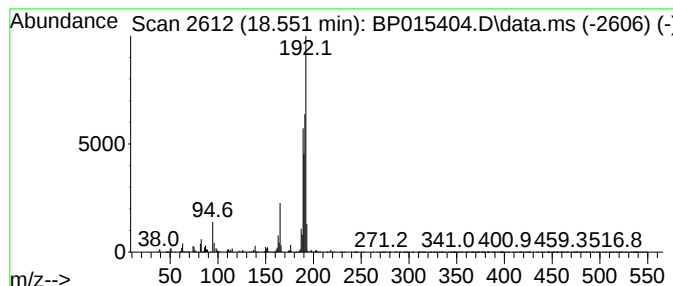
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 40 Anthracene, 1-methyl- Concentration Rank 1

| R.T.   | EstConc      | Area     | Relative to ISTD | R.T.   |
|--------|--------------|----------|------------------|--------|
| 18.551 | 110.50 ng/ul | 17328300 | Phenanthrene-d10 | 17.504 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Anthracene, 1-methyl-               | 192 | C15H12  | 000610-48-0 | 60   |
| 2    |    |   | Phenanthrene, 1-methyl-             | 192 | C15H12  | 000832-69-9 | 60   |
| 3    |    |   | Phenanthrene, 1-methyl-             | 192 | C15H12  | 000832-69-9 | 55   |
| 4    |    |   | Anthracene, 1-methyl-               | 192 | C15H12  | 000610-48-0 | 55   |
| 5    |    |   | 1H-Cyclopropa[1]phenanthrene,1a,... | 192 | C15H12  | 000949-41-7 | 53   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP060723\  
Data File : BP015404.D  
Acq On : 07 Jun 2023 18:08  
Operator : MA/JU  
Sample : 02950-01  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

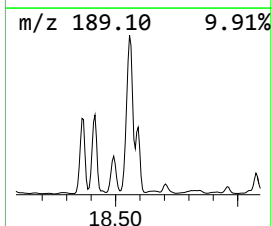
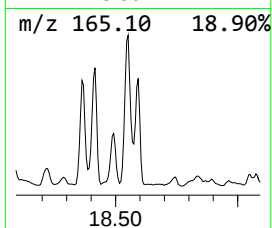
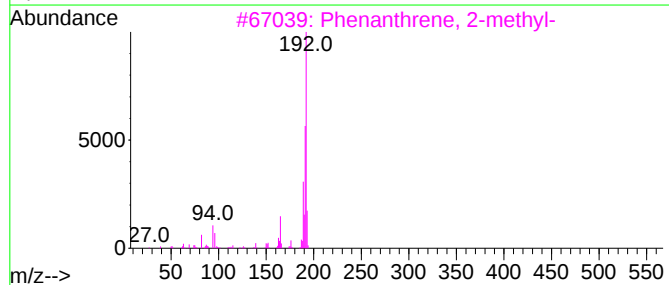
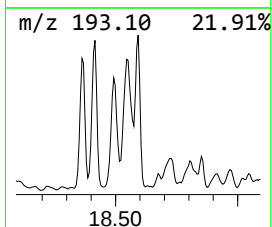
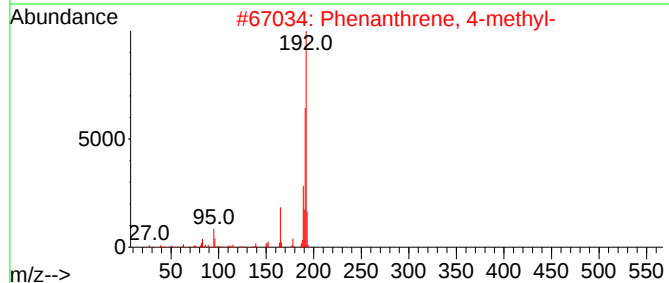
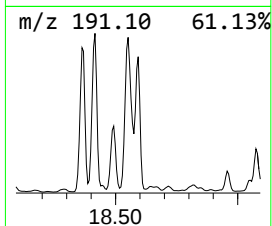
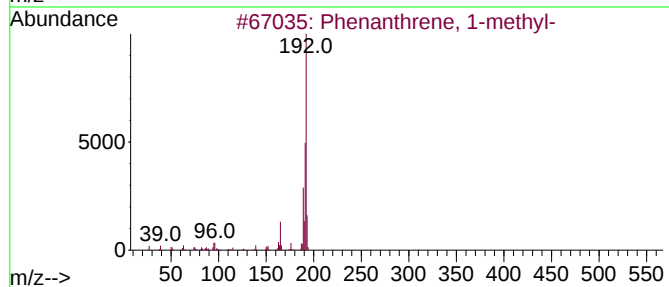
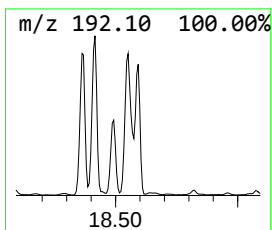
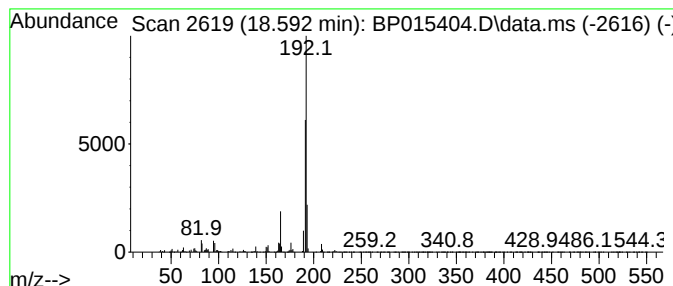
Instrument :  
BNA\_P  
ClientSampleId :  
EW978

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP060723.MA.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 41 Phenanthrene, 1-methyl- Concentration Rank 7

| R.T.      | EstConc                 | Area    | Relative to ISTD |         |             | R.T.   |
|-----------|-------------------------|---------|------------------|---------|-------------|--------|
| 18.592    | 51.21 ng/ul             | 8030720 | Phenanthrene-d10 |         |             | 17.504 |
| Hit# of 5 | Tentative ID            |         | MW               | MolForm | CAS#        | Qual   |
| 1         | Phenanthrene, 1-methyl- |         | 192              | C15H12  | 000832-69-9 | 93     |
| 2         | Phenanthrene, 4-methyl- |         | 192              | C15H12  | 000832-64-4 | 91     |
| 3         | Phenanthrene, 2-methyl- |         | 192              | C15H12  | 002531-84-2 | 90     |
| 4         | Anthracene, 1-methyl-   |         | 192              | C15H12  | 000610-48-0 | 90     |
| 5         | Anthracene, 9-methyl-   |         | 192              | C15H12  | 000779-02-2 | 90     |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP060723\  
Data File : BP015404.D  
Acq On : 07 Jun 2023 18:08  
Operator : MA/JU  
Sample : 02950-01  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
EW978

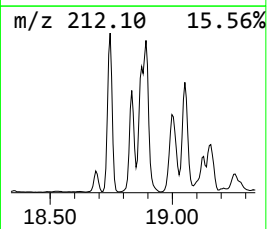
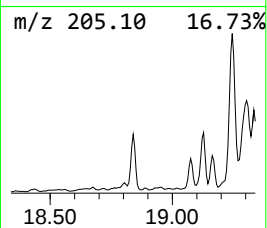
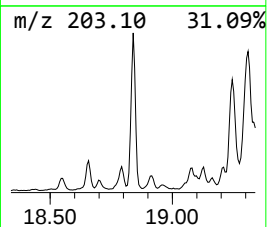
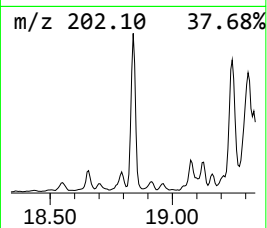
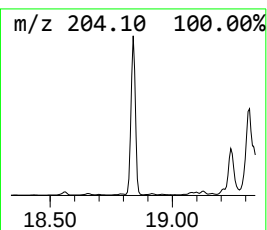
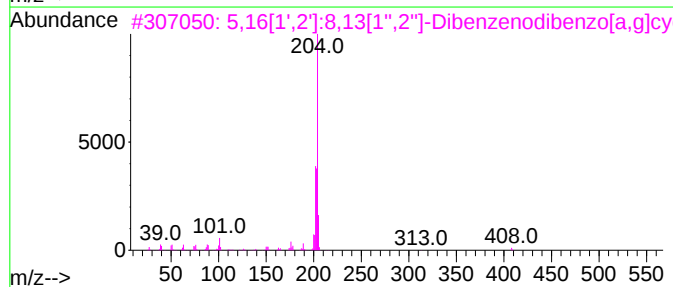
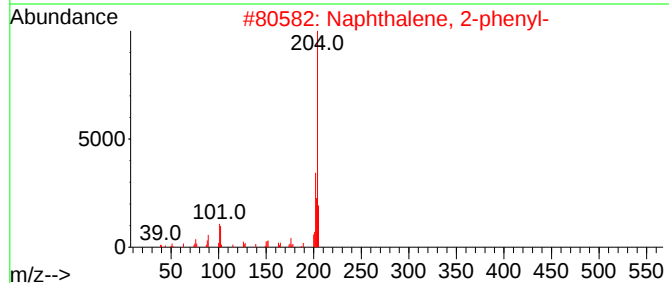
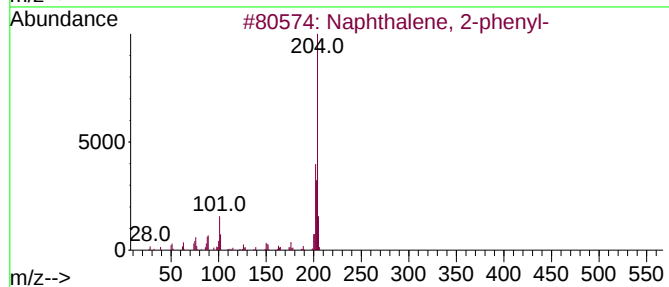
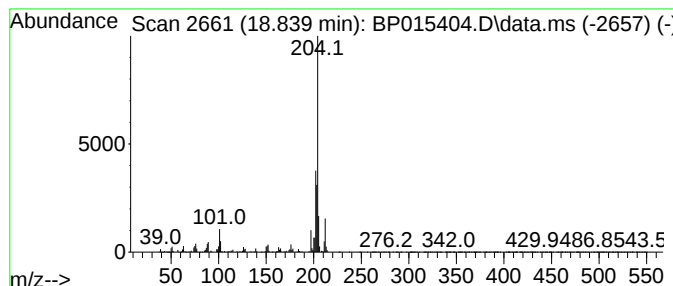
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 43 Naphthalene, 2-phenyl- Concentration Rank 22

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 18.839 | 22.17 ng/ul | 3477480 | Phenanthrene-d10 | 17.504 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, 2-phenyl-              | 204 | C16H12  | 000612-94-2 | 96   |
| 2    |    |   | Naphthalene, 2-phenyl-              | 204 | C16H12  | 000612-94-2 | 83   |
| 3    |    |   | 5,16[1',2']:8,13[1'',2'']-Dibenz... | 408 | C32H24  | 005672-97-9 | 72   |
| 4    |    |   | Naphthalene, 2-phenyl-              | 204 | C16H12  | 000612-94-2 | 70   |
| 5    |    |   | Naphthalene, 2-phenyl-              | 204 | C16H12  | 000612-94-2 | 70   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP060723\  
Data File : BP015404.D  
Acq On : 07 Jun 2023 18:08  
Operator : MA/JU  
Sample : 02950-01  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

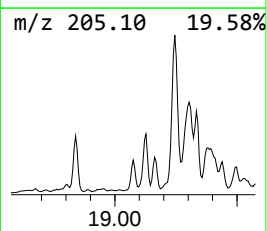
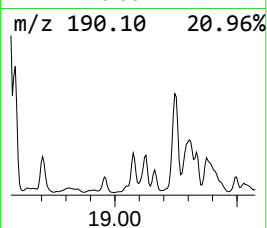
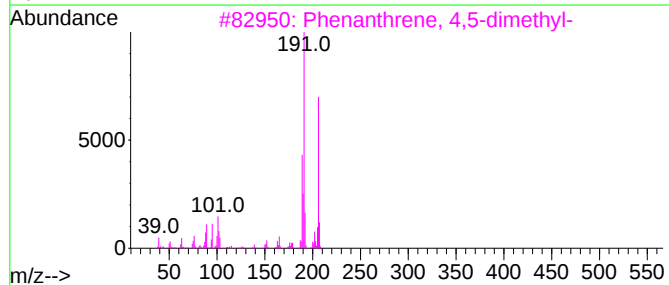
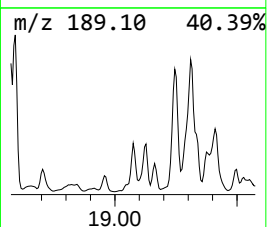
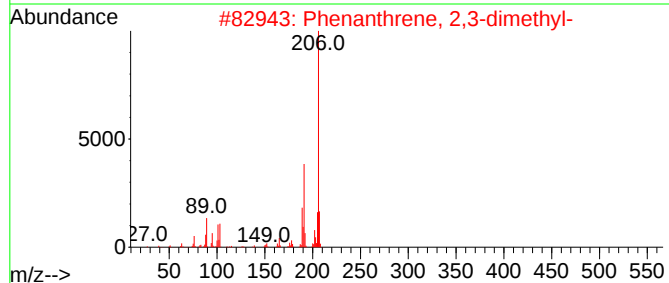
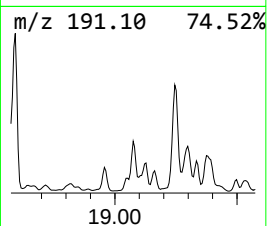
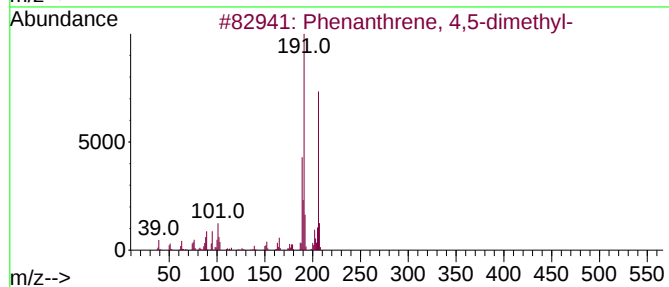
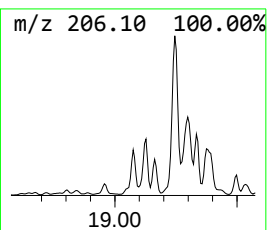
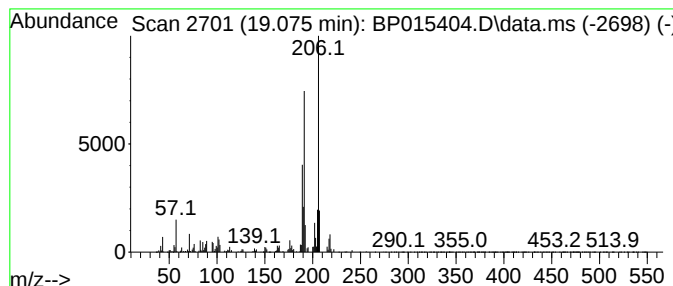
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BNA\_P  
ClientSampleId :  
EW978

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP060723.MA.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 46 Phenanthrene, 4,5-dimethyl- Concentration Rank 20

| R.T.      | EstConc                     | Area    | Relative to ISTD |         | R.T.        |      |
|-----------|-----------------------------|---------|------------------|---------|-------------|------|
| 19.075    | 22.79 ng/ul                 | 3573490 | Phenanthrene-d10 |         | 17.504      |      |
| Hit# of 5 | Tentative ID                |         | MW               | MolForm | CAS#        | Qual |
| 1         | Phenanthrene, 4,5-dimethyl- |         | 206              | C16H14  | 003674-69-9 | 93   |
| 2         | Phenanthrene, 2,3-dimethyl- |         | 206              | C16H14  | 003674-65-5 | 93   |
| 3         | Phenanthrene, 4,5-dimethyl- |         | 206              | C16H14  | 003674-69-9 | 87   |
| 4         | Phenanthrene, 4,5-dimethyl- |         | 206              | C16H14  | 003674-69-9 | 81   |
| 5         | 9,10-Dimethylanthracene     |         | 206              | C16H14  | 000781-43-1 | 76   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP060723\  
Data File : BP015404.D  
Acq On : 07 Jun 2023 18:08  
Operator : MA/JU  
Sample : 02950-01  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

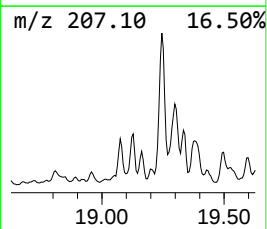
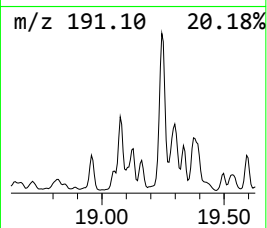
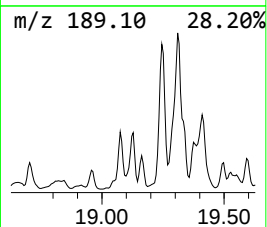
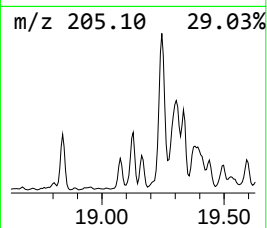
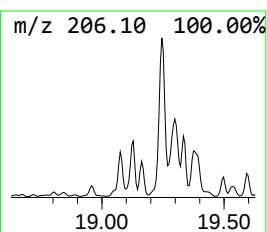
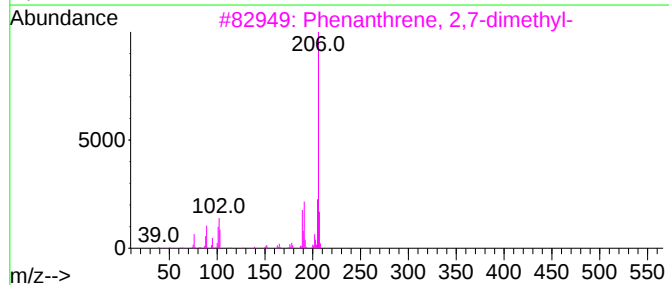
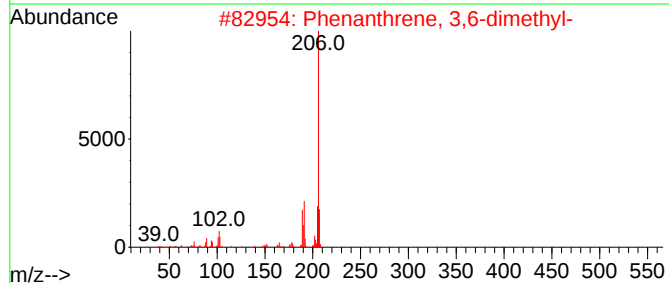
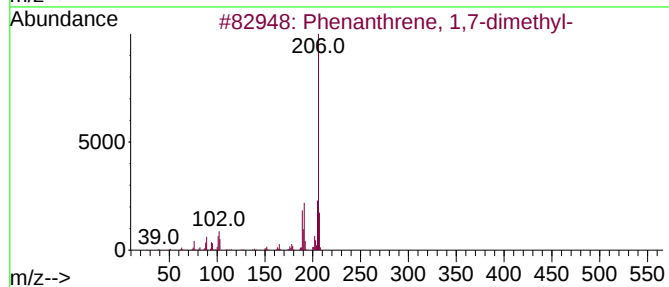
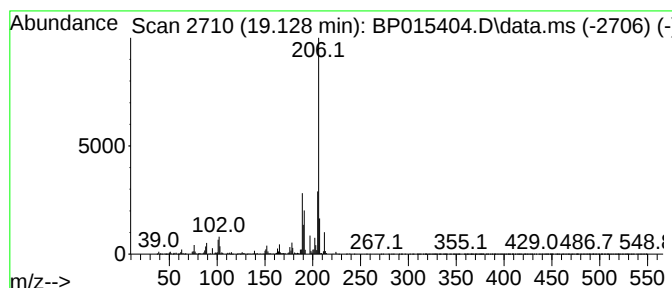
Instrument :  
BNA\_P  
ClientSampleId :  
EW978

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP060723.MA.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 47 Phenanthrene, 1,7-dimethyl- Concentration Rank 25

| R.T.      | EstConc                     | Area    | Relative to ISTD | R.T.        |      |
|-----------|-----------------------------|---------|------------------|-------------|------|
| 19.128    | 19.63 ng/ul                 | 3078660 | Phenanthrene-d10 | 17.504      |      |
| Hit# of 5 | Tentative ID                | MW      | MolForm          | CAS#        | Qual |
| 1         | Phenanthrene, 1,7-dimethyl- | 206     | C16H14           | 000483-87-4 | 90   |
| 2         | Phenanthrene, 3,6-dimethyl- | 206     | C16H14           | 001576-67-6 | 90   |
| 3         | Phenanthrene, 2,7-dimethyl- | 206     | C16H14           | 001576-69-8 | 87   |
| 4         | Phenanthrene, 3,6-dimethyl- | 206     | C16H14           | 001576-67-6 | 81   |
| 5         | di-p-Tolylacetylene         | 206     | C16H14           | 002789-88-0 | 74   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP060723\  
Data File : BP015404.D  
Acq On : 07 Jun 2023 18:08  
Operator : MA/JU  
Sample : 02950-01  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
EW978

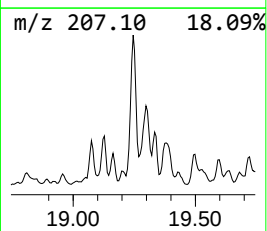
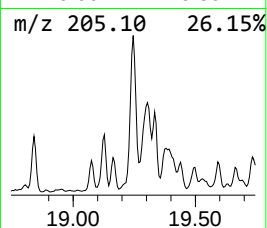
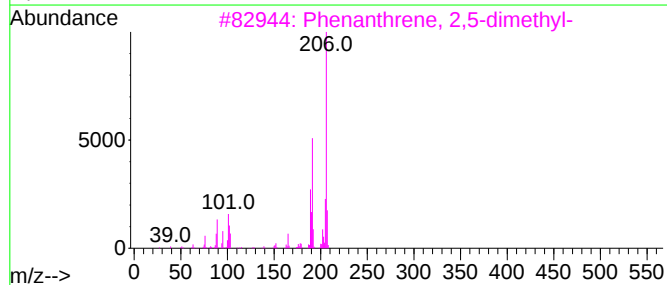
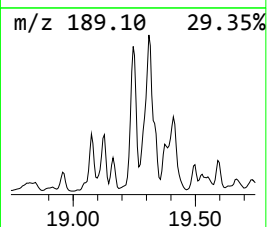
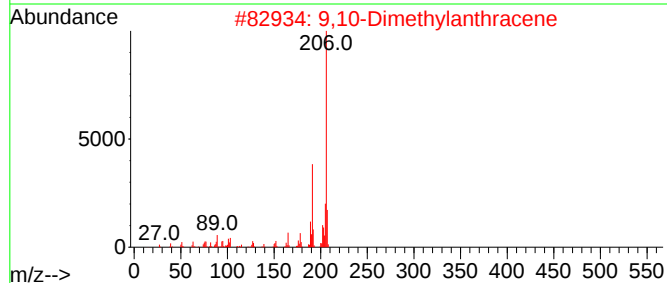
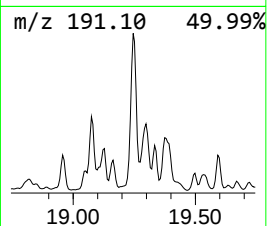
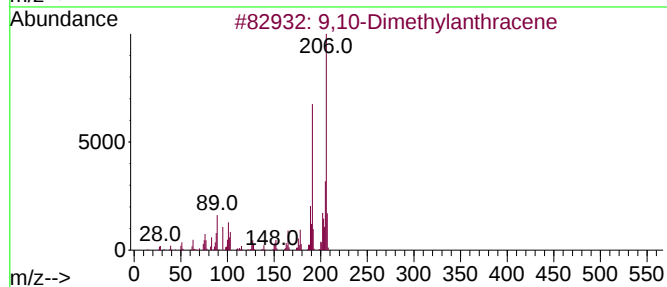
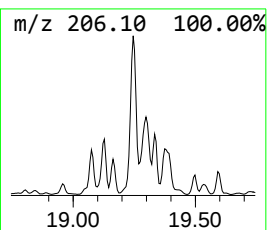
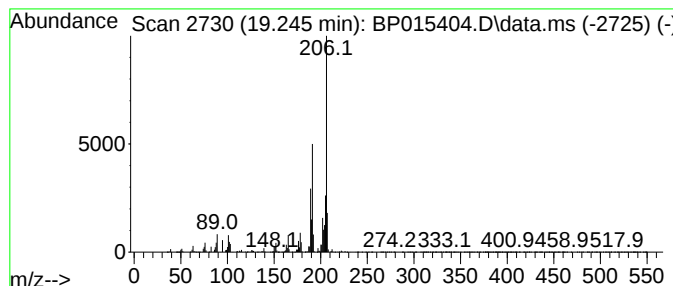
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 49 9,10-Dimethylantracene Concentration Rank 5

| R.T.   | EstConc     | Area     | Relative to ISTD | R.T.   |
|--------|-------------|----------|------------------|--------|
| 19.245 | 64.23 ng/ul | 10072600 | Phenanthrene-d10 | 17.504 |

| Hit# | of | 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------------|-----|---------|-------------|------|
| 1    |    |   | 9,10-Dimethylantracene      | 206 | C16H14  | 000781-43-1 | 97   |
| 2    |    |   | 9,10-Dimethylantracene      | 206 | C16H14  | 000781-43-1 | 95   |
| 3    |    |   | Phenanthrene, 2,5-dimethyl- | 206 | C16H14  | 003674-66-6 | 95   |
| 4    |    |   | Phenanthrene, 3,6-dimethyl- | 206 | C16H14  | 001576-67-6 | 91   |
| 5    |    |   | 9,10-Dimethylantracene      | 206 | C16H14  | 000781-43-1 | 91   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP060723\  
Data File : BP015404.D  
Acq On : 07 Jun 2023 18:08  
Operator : MA/JU  
Sample : 02950-01  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
EW978

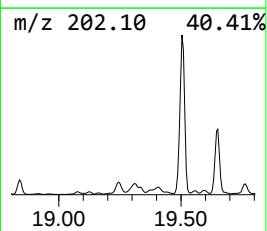
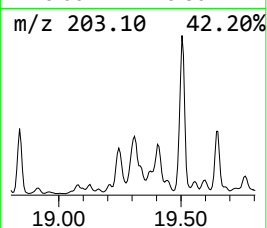
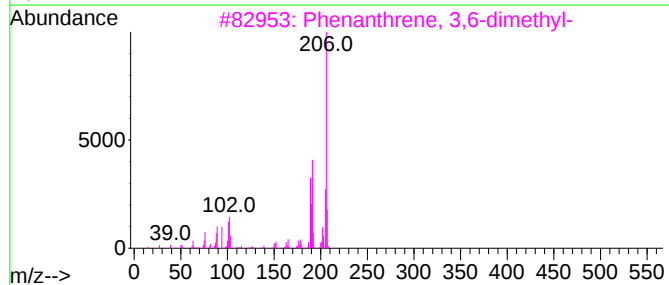
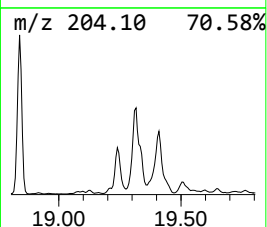
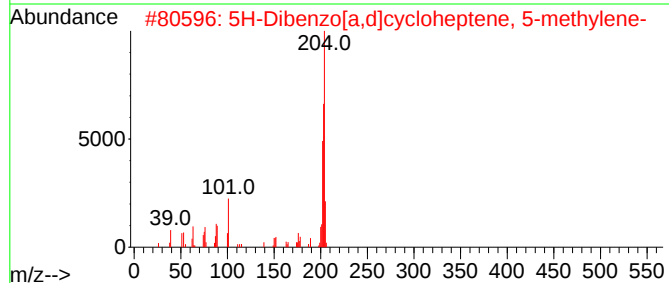
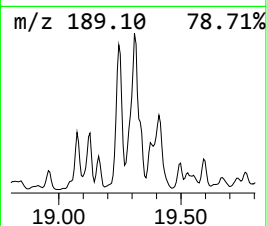
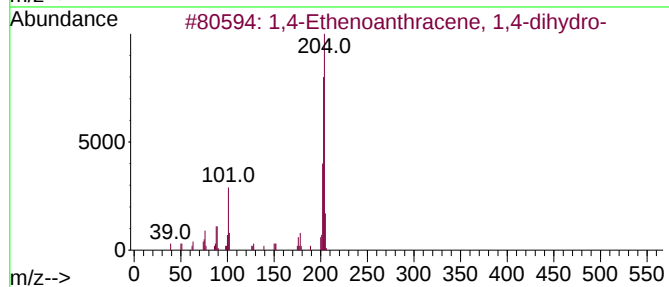
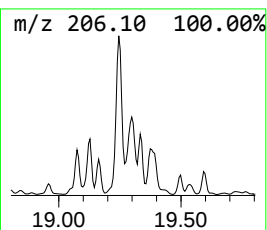
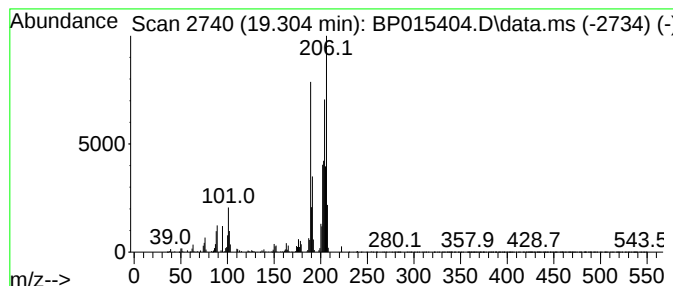
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 50 1,4-Ethenoanthracene, 1,4-d... Concentration Rank 9

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 19.304 | 49.10 ng/ul | 7699850 | Phenanthrene-d10 | 17.504 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | 1,4-Ethenoanthracene, 1,4-dihydro-  | 204 | C16H12  | 027765-96-4 | 55   |
| 2    |    |   | 5H-Dibenzo[a,d]cycloheptene, 5-m... | 204 | C16H12  | 002975-79-3 | 53   |
| 3    |    |   | Phenanthrene, 3,6-dimethyl-         | 206 | C16H14  | 001576-67-6 | 46   |
| 4    |    |   | Pyrene, 4,5,9,10-tetrahydro-        | 206 | C16H14  | 000781-17-9 | 46   |
| 5    |    |   | 1,3-Butadiene, 1,4-diphenyl-, (E... | 206 | C16H14  | 000538-81-8 | 41   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP060723\  
Data File : BP015404.D  
Acq On : 07 Jun 2023 18:08  
Operator : MA/JU  
Sample : 02950-01  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
EW978

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP060723.MA.M  
Quant Title : SVOA CALIBRATION

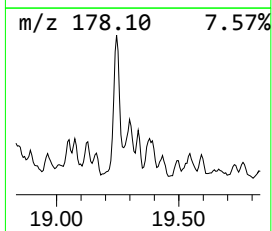
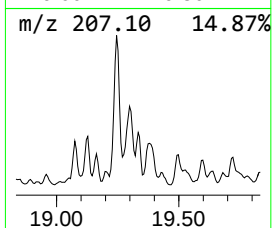
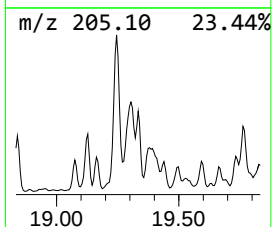
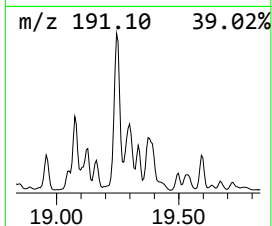
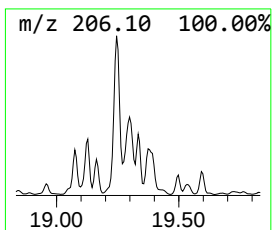
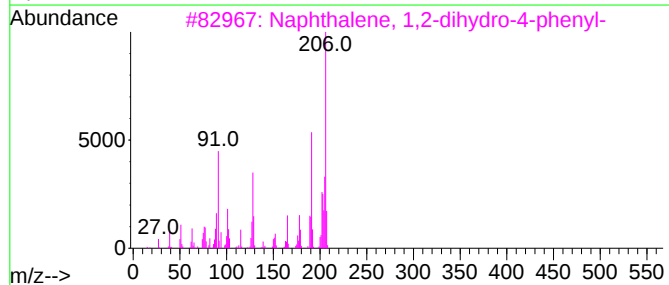
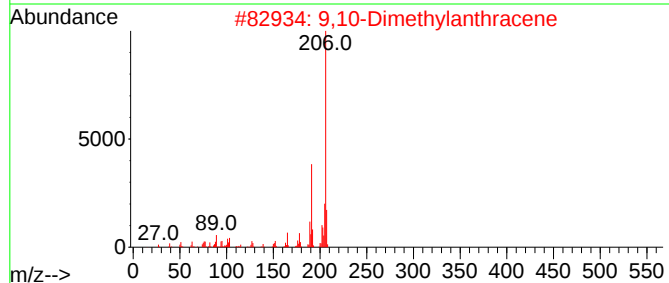
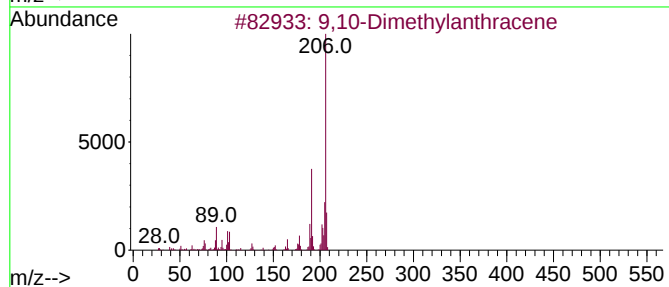
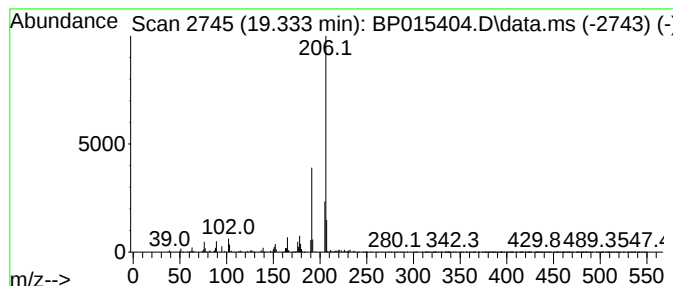
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 51 Naphthalene, 1,2-dihydro-4-... Concentration Rank 29

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 19.334 | 17.38 ng/ul | 2725220 | Phenanthrene-d10 | 17.504 |

| Hit# | of                                 | 5   | Tentative ID | MW          | MolForm | CAS# | Qual |
|------|------------------------------------|-----|--------------|-------------|---------|------|------|
| 1    | 9,10-Dimethylantracene             | 206 | C16H14       | 000781-43-1 | 91      |      |      |
| 2    | 9,10-Dimethylantracene             | 206 | C16H14       | 000781-43-1 | 87      |      |      |
| 3    | Naphthalene, 1,2-dihydro-4-phenyl- | 206 | C16H14       | 007469-40-1 | 83      |      |      |
| 4    | Anthracene, 1,4-dimethyl-          | 206 | C16H14       | 000781-92-0 | 80      |      |      |
| 5    | Phenanthrene, 3,6-dimethyl-        | 206 | C16H14       | 001576-67-6 | 78      |      |      |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP060723\  
Data File : BP015404.D  
Acq On : 07 Jun 2023 18:08  
Operator : MA/JU  
Sample : 02950-01  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
EW978

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP060723.MA.M  
Quant Title : SVOA CALIBRATION

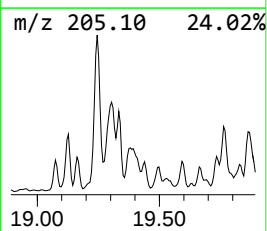
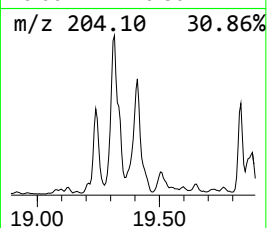
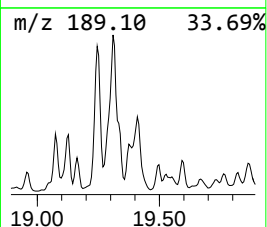
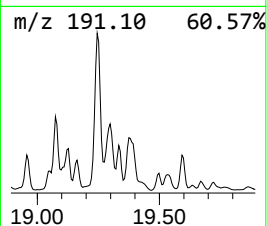
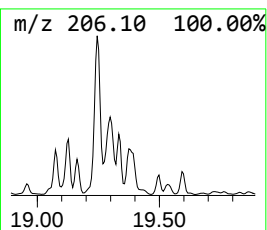
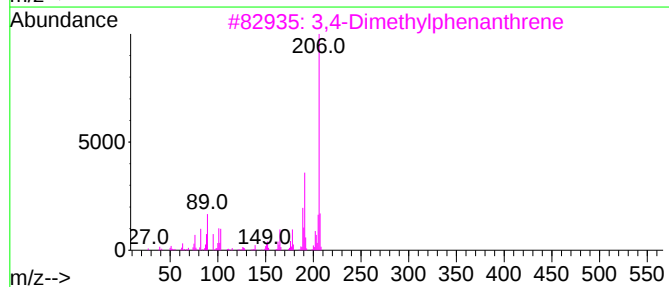
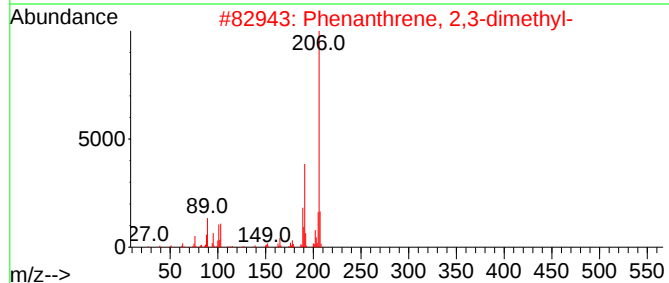
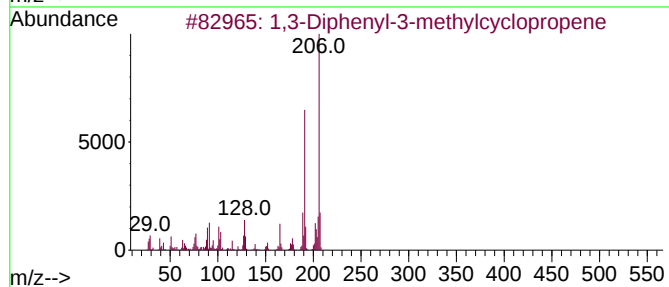
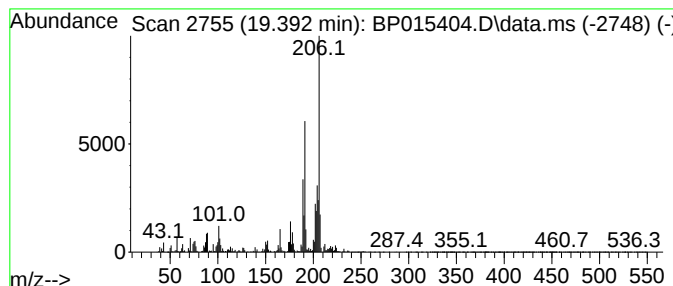
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 52 1,3-Diphenyl-3-methylcyclop... Concentration Rank 6

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 19.392 | 57.83 ng/ul | 9069130 | Phenanthrene-d10 | 17.504 |

| Hit# | of 5 | Tentative ID                      | MW  | MolForm | CAS#         | Qual |
|------|------|-----------------------------------|-----|---------|--------------|------|
| 1    |      | 1,3-Diphenyl-3-methylcyclopropene | 206 | C16H14  | 086544-79-8  | 93   |
| 2    |      | Phenanthrene, 2,3-dimethyl-       | 206 | C16H14  | 003674-65-5  | 90   |
| 3    |      | 3,4-Dimethylphenanthrene          | 206 | C16H14  | 1000452-24-5 | 89   |
| 4    |      | Phenanthrene, 2,5-dimethyl-       | 206 | C16H14  | 003674-66-6  | 89   |
| 5    |      | 9,10-Dimethylanthracene           | 206 | C16H14  | 000781-43-1  | 89   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP060723\  
 Data File : BP015404.D  
 Acq On : 07 Jun 2023 18:08  
 Operator : MA/JU  
 Sample : 02950-01  
 Misc :  
 ALS Vial : 10 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 EW978

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP060723.MA.M  
 Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST0.L  
 TIC Integration Parameters: LSCINT.P

| TIC Top Hit name   | RT     | EstConc | Units | Response | --Internal Standard-- |        |         |      |
|--------------------|--------|---------|-------|----------|-----------------------|--------|---------|------|
|                    |        |         |       |          | #                     | RT     | Resp    | Conc |
| Naphthalene, 1,... | 13.922 | 25.0    | ng/ul | 3460330  | 3                     | 14.746 | 2772470 | 20.0 |
| Naphthalene, 1,... | 14.069 | 72.8    | ng/ul | 10087600 | 3                     | 14.746 | 2772470 | 20.0 |
| Naphthalene, 2,... | 14.304 | 37.6    | ng/ul | 5213320  | 3                     | 14.746 | 2772470 | 20.0 |
| Naphthalene, 1,... | 15.034 | 18.5    | ng/ul | 2569000  | 3                     | 14.746 | 2772470 | 20.0 |
| Naphthalene, 2,... | 15.140 | 49.1    | ng/ul | 6806960  | 3                     | 14.746 | 2772470 | 20.0 |
| Naphthalene, 1,... | 15.216 | 29.9    | ng/ul | 4140100  | 3                     | 14.746 | 2772470 | 20.0 |
| unknown-01         | 15.357 | 36.5    | ng/ul | 5055370  | 3                     | 14.746 | 2772470 | 20.0 |
| unknown-02         | 15.410 | 21.1    | ng/ul | 2931710  | 3                     | 14.746 | 2772470 | 20.0 |
| Naphthalene, 1,... | 15.528 | 30.9    | ng/ul | 4284830  | 3                     | 14.746 | 2772470 | 20.0 |
| 1-Isopropenylna... | 16.004 | 22.7    | ng/ul | 3152920  | 3                     | 14.746 | 2772470 | 20.0 |
| (DEL) Alkane: S... | 16.457 | 44.6    | ng/ul | 6995630  | 4                     | 17.504 | 3136490 | 20.0 |
| 9H-Fluorene, 1-... | 16.798 | 39.9    | ng/ul | 6263140  | 4                     | 17.504 | 3136490 | 20.0 |
| 9H-Fluorene, 2-... | 16.851 | 28.1    | ng/ul | 4410720  | 4                     | 17.504 | 3136490 | 20.0 |
| 9H-Fluorene, 4-... | 16.951 | 17.9    | ng/ul | 2813370  | 4                     | 17.504 | 3136490 | 20.0 |
| unknown-03         | 17.157 | 13.5    | ng/ul | 2122780  | 4                     | 17.504 | 3136490 | 20.0 |
| (DEL) Alkane: S... | 17.287 | 11.5    | ng/ul | 1800620  | 4                     | 17.504 | 3136490 | 20.0 |
| Dibenzothiophene   | 17.322 | 24.2    | ng/ul | 3800530  | 4                     | 17.504 | 3136490 | 20.0 |
| 4-Hydroxybenzal... | 17.692 | 16.9    | ng/ul | 2647540  | 4                     | 17.504 | 3136490 | 20.0 |
| 1-Methyldibenzo... | 18.075 | 21.2    | ng/ul | 3320340  | 4                     | 17.504 | 3136490 | 20.0 |
| Dibenzothiophen... | 18.216 | 17.5    | ng/ul | 2749200  | 4                     | 17.504 | 3136490 | 20.0 |
| Anthracene, 2-m... | 18.363 | 65.8    | ng/ul | 10316500 | 4                     | 17.504 | 3136490 | 20.0 |
| Phenanthrene, 2... | 18.416 | 67.0    | ng/ul | 10499600 | 4                     | 17.504 | 3136490 | 20.0 |
| 1H-Cyclopropa[1... | 18.492 | 35.3    | ng/ul | 5542970  | 4                     | 17.504 | 3136490 | 20.0 |
| Anthracene, 1-m... | 18.551 | 110.5   | ng/ul | 17328300 | 4                     | 17.504 | 3136490 | 20.0 |
| Phenanthrene, 1... | 18.592 | 51.2    | ng/ul | 8030720  | 4                     | 17.504 | 3136490 | 20.0 |
| Naphthalene, 2-... | 18.839 | 22.2    | ng/ul | 3477480  | 4                     | 17.504 | 3136490 | 20.0 |
| Phenanthrene, 4... | 19.075 | 22.8    | ng/ul | 3573490  | 4                     | 17.504 | 3136490 | 20.0 |
| Phenanthrene, 1... | 19.128 | 19.6    | ng/ul | 3078660  | 4                     | 17.504 | 3136490 | 20.0 |
| 9,10-Dimethylan... | 19.245 | 64.2    | ng/ul | 10072600 | 4                     | 17.504 | 3136490 | 20.0 |
| 1,4-Ethenoanthr... | 19.304 | 49.1    | ng/ul | 7699850  | 4                     | 17.504 | 3136490 | 20.0 |
| Naphthalene, 1...  | 19.334 | 17.4    | ng/ul | 2725220  | 4                     | 17.504 | 3136490 | 20.0 |
| 1,3-Diphenyl-3-... | 19.392 | 57.8    | ng/ul | 9069130  | 4                     | 17.504 | 3136490 | 20.0 |