

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP013125\  
 Data File : BP023838.D  
 Acq On : 01 Feb 2025 00:10  
 Operator : RC/JU  
 Sample : Q1224-02  
 Misc :  
 ALS Vial : 22 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 A6302

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-BP012925.MA.M  
 Title : SVOA CALIBRATION

Signal : TIC: BP023838.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         | 3.293       | 32            | 35          | 42           | rVB      | 65919          | 86867         | 0.90%           | 0.083%        |
| 2         | 3.711       | 102           | 106         | 119          | rBV      | 596141         | 899388        | 9.34%           | 0.861%        |
| 3         | 5.170       | 347           | 354         | 368          | rVB3     | 40894          | 97767         | 1.02%           | 0.094%        |
| 4         | 6.222       | 527           | 533         | 545          | rVB      | 39268          | 68473         | 0.71%           | 0.066%        |
| 5         | 6.487       | 574           | 578         | 585          | rVB      | 44863          | 69401         | 0.72%           | 0.066%        |
| 6         | 6.799       | 625           | 631         | 640          | rBV4     | 17212          | 34390         | 0.36%           | 0.033%        |
| 7         | 6.969       | 648           | 660         | 675          | rBV      | 926093         | 1582336       | 16.44%          | 1.515%        |
| 8         | 7.111       | 678           | 684         | 699          | rVB      | 896322         | 1408358       | 14.63%          | 1.348%        |
| 9         | 7.234       | 699           | 705         | 711          | rVB2     | 69932          | 105202        | 1.09%           | 0.101%        |
| 10        | 7.322       | 713           | 720         | 729          | rBV      | 1022632        | 1590985       | 16.53%          | 1.523%        |
| 11        | 7.528       | 742           | 755         | 767          | rBV      | 1328952        | 2609529       | 27.11%          | 2.498%        |
| 12        | 7.775       | 788           | 797         | 804          | rBV      | 1511196        | 2413624       | 25.08%          | 2.310%        |
| 13        | 7.852       | 804           | 810         | 826          | rVB      | 2795381        | 4507309       | 46.83%          | 4.314%        |
| 14        | 8.022       | 832           | 839         | 857          | rVB      | 135065         | 331857        | 3.45%           | 0.318%        |
| 15        | 8.228       | 868           | 874         | 878          | rBV      | 48822          | 74685         | 0.78%           | 0.071%        |
| 16        | 8.281       | 878           | 883         | 892          | rVV2     | 61142          | 109034        | 1.13%           | 0.104%        |
| 17        | 8.487       | 910           | 918         | 925          | rBV      | 853511         | 1310851       | 13.62%          | 1.255%        |
| 18        | 8.563       | 925           | 931         | 935          | rVV4     | 20557          | 42866         | 0.45%           | 0.041%        |
| 19        | 8.922       | 985           | 992         | 1002         | rVV      | 1004613        | 1570634       | 16.32%          | 1.503%        |
| 20        | 9.034       | 1002          | 1011        | 1017         | rVV      | 56432          | 114532        | 1.19%           | 0.110%        |
| 21        | 9.087       | 1017          | 1020        | 1025         | rVB3     | 23006          | 36959         | 0.38%           | 0.035%        |
| 22        | 9.369       | 1063          | 1068        | 1073         | rVB      | 36279          | 58664         | 0.61%           | 0.056%        |
| 23        | 9.487       | 1079          | 1088        | 1093         | rBV      | 152858         | 237995        | 2.47%           | 0.228%        |
| 24        | 9.640       | 1107          | 1114        | 1122         | rBV      | 732369         | 1200073       | 12.47%          | 1.149%        |
| 25        | 9.740       | 1127          | 1131        | 1136         | rBV      | 48193          | 75174         | 0.78%           | 0.072%        |
| 26        | 9.828       | 1136          | 1146        | 1153         | rBV      | 307140         | 630120        | 6.55%           | 0.603%        |
| 27        | 10.187      | 1200          | 1207        | 1217         | rVB      | 1220899        | 1887192       | 19.61%          | 1.806%        |
| 28        | 10.363      | 1230          | 1237        | 1250         | rBV      | 642158         | 1339256       | 13.91%          | 1.282%        |
| 29        | 10.557      | 1263          | 1270        | 1274         | rBV      | 2174368        | 3848155       | 39.98%          | 3.683%        |
| 30        | 10.593      | 1274          | 1276        | 1284         | rVV      | 1120282        | 1597503       | 16.60%          | 1.529%        |
| 31        | 10.693      | 1287          | 1293        | 1300         | rVB      | 164737         | 296372        | 3.08%           | 0.284%        |
| 32        | 11.093      | 1356          | 1361        | 1372         | rVV3     | 76026          | 147692        | 1.53%           | 0.141%        |
| 33        | 11.199      | 1373          | 1379        | 1388         | rVV7     | 23387          | 58358         | 0.61%           | 0.056%        |
| 34        | 11.463      | 1418          | 1424        | 1435         | rBV      | 26678          | 70124         | 0.73%           | 0.067%        |
| 35        | 11.563      | 1435          | 1441        | 1448         | rVV      | 139184         | 255677        | 2.66%           | 0.245%        |
| 36        | 11.646      | 1448          | 1455        | 1463         | rVV      | 597010         | 922716        | 9.59%           | 0.883%        |

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 ALS Vial : 22 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 A6302

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-BP012925.MA.M  
 Title : SVOA CALIBRATION

|    |        |      |      |      |      |         |         |        |        |
|----|--------|------|------|------|------|---------|---------|--------|--------|
| 37 | 12.246 | 1551 | 1557 | 1565 | rVB9 | 13923   | 34237   | 0.36%  | 0.033% |
| 38 | 12.463 | 1590 | 1594 | 1602 | rVB  | 47887   | 80365   | 0.83%  | 0.077% |
| 39 | 12.546 | 1602 | 1608 | 1614 | rBV  | 37321   | 64803   | 0.67%  | 0.062% |
| 40 | 12.740 | 1635 | 1641 | 1649 | rBV2 | 42271   | 100192  | 1.04%  | 0.096% |
| 41 | 12.822 | 1652 | 1655 | 1661 | rVV3 | 18453   | 37379   | 0.39%  | 0.036% |
| 42 | 13.181 | 1708 | 1716 | 1720 | rBV4 | 25195   | 49141   | 0.51%  | 0.047% |
| 43 | 13.228 | 1720 | 1724 | 1731 | rVB  | 33869   | 50741   | 0.53%  | 0.049% |
| 44 | 13.310 | 1731 | 1738 | 1745 | rBV  | 558248  | 810317  | 8.42%  | 0.776% |
| 45 | 13.516 | 1763 | 1773 | 1794 | rBV3 | 849644  | 2670297 | 27.74% | 2.556% |
| 46 | 13.798 | 1813 | 1821 | 1830 | rVV  | 2135359 | 2939527 | 30.54% | 2.814% |
| 47 | 13.928 | 1840 | 1843 | 1849 | rVB  | 21342   | 34669   | 0.36%  | 0.033% |
| 48 | 14.087 | 1861 | 1870 | 1880 | rVB  | 1797754 | 2505645 | 26.03% | 2.398% |
| 49 | 14.257 | 1893 | 1899 | 1905 | rBV5 | 30453   | 59840   | 0.62%  | 0.057% |
| 50 | 14.398 | 1914 | 1923 | 1932 | rBV  | 3798845 | 4900032 | 50.91% | 4.690% |
| 51 | 14.628 | 1954 | 1962 | 1966 | rBV  | 649736  | 929286  | 9.66%  | 0.889% |
| 52 | 14.663 | 1966 | 1968 | 1973 | rVB  | 73434   | 96533   | 1.00%  | 0.092% |
| 53 | 14.745 | 1979 | 1982 | 1988 | rVB3 | 23532   | 38478   | 0.40%  | 0.037% |
| 54 | 14.828 | 1991 | 1996 | 2000 | rBV  | 79975   | 110933  | 1.15%  | 0.106% |
| 55 | 14.934 | 2009 | 2014 | 2017 | rBV  | 100115  | 143918  | 1.50%  | 0.138% |
| 56 | 14.963 | 2017 | 2019 | 2023 | rVV2 | 56185   | 71472   | 0.74%  | 0.068% |
| 57 | 15.222 | 2058 | 2063 | 2070 | rBV3 | 38077   | 72691   | 0.76%  | 0.070% |
| 58 | 15.410 | 2088 | 2095 | 2102 | rBV  | 1982556 | 2647221 | 27.50% | 2.534% |
| 59 | 15.528 | 2108 | 2115 | 2123 | rBV  | 600213  | 900672  | 9.36%  | 0.862% |
| 60 | 15.598 | 2123 | 2127 | 2131 | rVV4 | 21950   | 38381   | 0.40%  | 0.037% |
| 61 | 15.657 | 2132 | 2137 | 2141 | rVB  | 64083   | 88228   | 0.92%  | 0.084% |
| 62 | 15.781 | 2149 | 2158 | 2167 | rBV  | 1852031 | 2582161 | 26.83% | 2.472% |
| 63 | 15.875 | 2169 | 2174 | 2179 | rBV6 | 20071   | 41700   | 0.43%  | 0.040% |
| 64 | 15.969 | 2186 | 2190 | 2201 | rVB6 | 26162   | 69768   | 0.72%  | 0.067% |
| 65 | 16.163 | 2219 | 2223 | 2230 | rVB9 | 21011   | 36206   | 0.38%  | 0.035% |
| 66 | 16.363 | 2247 | 2257 | 2266 | rVB  | 186583  | 343742  | 3.57%  | 0.329% |
| 67 | 16.475 | 2271 | 2276 | 2283 | rVB  | 41626   | 67886   | 0.71%  | 0.065% |
| 68 | 16.581 | 2288 | 2294 | 2313 | rBV5 | 350656  | 1066230 | 11.08% | 1.021% |
| 69 | 16.804 | 2329 | 2332 | 2344 | rBV4 | 36501   | 81360   | 0.85%  | 0.078% |
| 70 | 16.910 | 2345 | 2350 | 2357 | rBV2 | 792431  | 1123231 | 11.67% | 1.075% |
| 71 | 17.010 | 2364 | 2367 | 2375 | rVB4 | 53255   | 87515   | 0.91%  | 0.084% |
| 72 | 17.204 | 2394 | 2400 | 2405 | rBV2 | 3882927 | 5635022 | 58.55% | 5.394% |
| 73 | 17.245 | 2405 | 2407 | 2411 | rVV  | 233545  | 295509  | 3.07%  | 0.283% |
| 74 | 17.304 | 2411 | 2417 | 2426 | rVB  | 1777835 | 2492017 | 25.89% | 2.385% |
| 75 | 17.516 | 2449 | 2453 | 2458 | rBV6 | 46827   | 71827   | 0.75%  | 0.069% |
| 76 | 17.604 | 2463 | 2468 | 2476 | rVV2 | 206466  | 311250  | 3.23%  | 0.298% |

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 ALS Vial : 22 Sample Multiplier: 1

Instrument :  
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 ClientSampleId :  
 A6302

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-BP012925.MA.M  
 Title : SVOA CALIBRATION

|     |        |      |      |      |      |         |         |         |        |
|-----|--------|------|------|------|------|---------|---------|---------|--------|
| 77  | 18.081 | 2543 | 2549 | 2555 | rBV  | 569043  | 982679  | 10.21%  | 0.941% |
| 78  | 18.145 | 2555 | 2560 | 2577 | rVB  | 595668  | 1029364 | 10.70%  | 0.985% |
| 79  | 18.392 | 2598 | 2602 | 2608 | rBV4 | 52118   | 110723  | 1.15%   | 0.106% |
| 80  | 18.457 | 2610 | 2613 | 2620 | rVV6 | 55865   | 109032  | 1.13%   | 0.104% |
| 81  | 18.657 | 2645 | 2647 | 2653 | rVB2 | 56486   | 69327   | 0.72%   | 0.066% |
| 82  | 19.010 | 2704 | 2707 | 2710 | rBV2 | 55452   | 94133   | 0.98%   | 0.090% |
| 83  | 19.110 | 2721 | 2724 | 2730 | rBV  | 56444   | 65667   | 0.68%   | 0.063% |
| 84  | 19.322 | 2754 | 2760 | 2771 | rBV  | 893857  | 1400772 | 14.55%  | 1.341% |
| 85  | 19.469 | 2781 | 2785 | 2792 | rBV2 | 237998  | 364621  | 3.79%   | 0.349% |
| 86  | 19.669 | 2814 | 2819 | 2823 | rBV  | 2081263 | 2818457 | 29.28%  | 2.698% |
| 87  | 19.698 | 2823 | 2824 | 2830 | rVB  | 644535  | 593200  | 6.16%   | 0.568% |
| 88  | 19.828 | 2842 | 2846 | 2850 | rVB  | 74880   | 101215  | 1.05%   | 0.097% |
| 89  | 20.786 | 3007 | 3009 | 3013 | rVB  | 132402  | 119085  | 1.24%   | 0.114% |
| 90  | 21.327 | 3095 | 3101 | 3109 | rBV2 | 1054737 | 1754144 | 18.23%  | 1.679% |
| 91  | 21.416 | 3111 | 3116 | 3124 | rVB  | 937775  | 1342574 | 13.95%  | 1.285% |
| 92  | 21.639 | 3147 | 3154 | 3160 | rBV3 | 4706376 | 9624563 | 100.00% | 9.212% |
| 93  | 22.045 | 3219 | 3223 | 3233 | rVB3 | 476384  | 863389  | 8.97%   | 0.826% |
| 94  | 22.563 | 3306 | 3311 | 3323 | rBV3 | 1031471 | 2245690 | 23.33%  | 2.149% |
| 95  | 23.480 | 3461 | 3467 | 3478 | rVB  | 933217  | 2154168 | 22.38%  | 2.062% |
| 96  | 23.680 | 3497 | 3501 | 3511 | rVB2 | 588357  | 1203605 | 12.51%  | 1.152% |
| 97  | 24.186 | 3582 | 3587 | 3593 | rBV  | 605261  | 1251168 | 13.00%  | 1.198% |
| 98  | 24.780 | 3683 | 3688 | 3696 | rVB  | 603883  | 1394052 | 14.48%  | 1.334% |
| 99  | 25.027 | 3721 | 3730 | 3740 | rVB  | 2216380 | 6259799 | 65.04%  | 5.992% |
| 100 | 25.768 | 3850 | 3856 | 3867 | rVB  | 1221252 | 3055036 | 31.74%  | 2.924% |

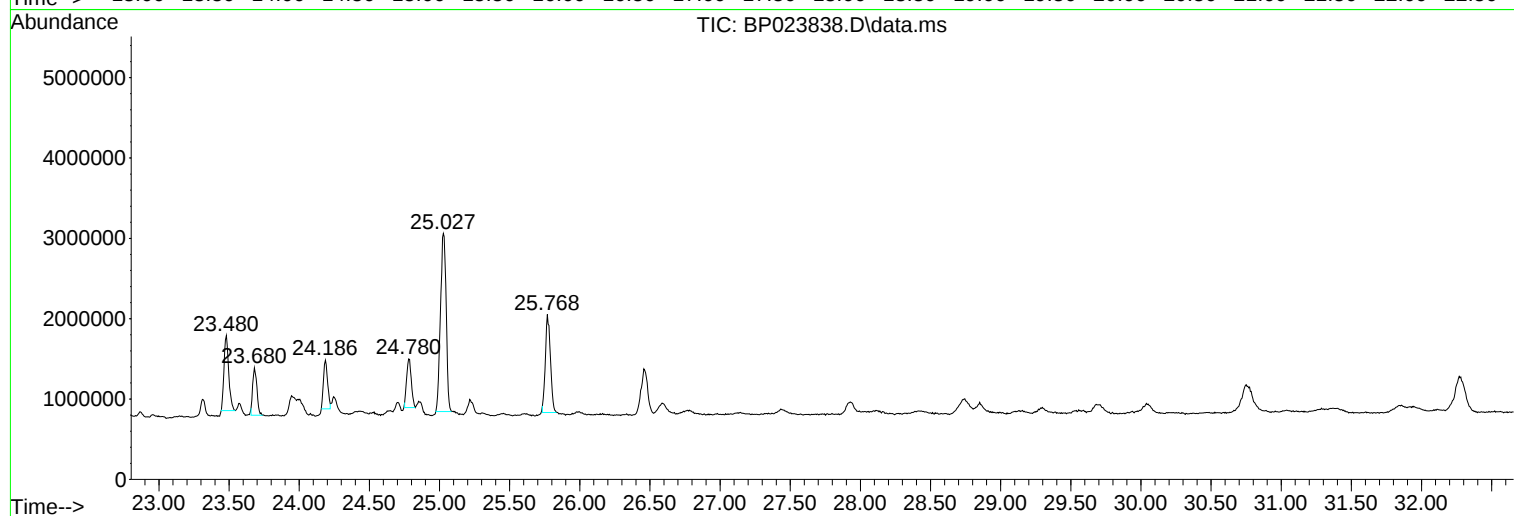
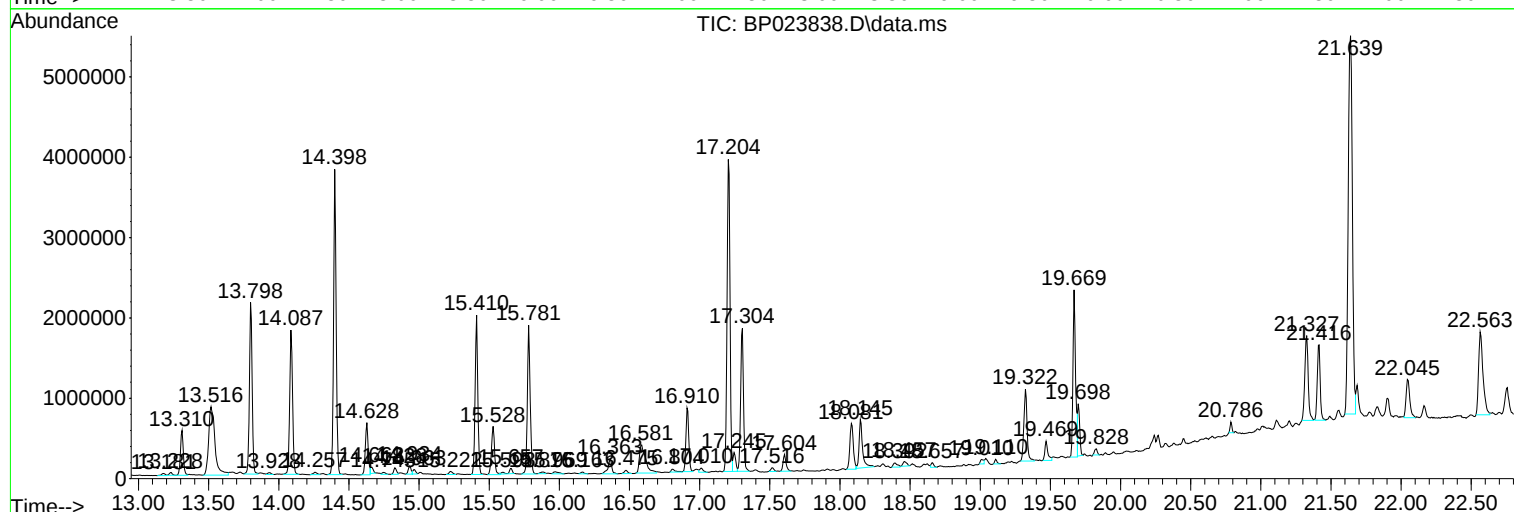
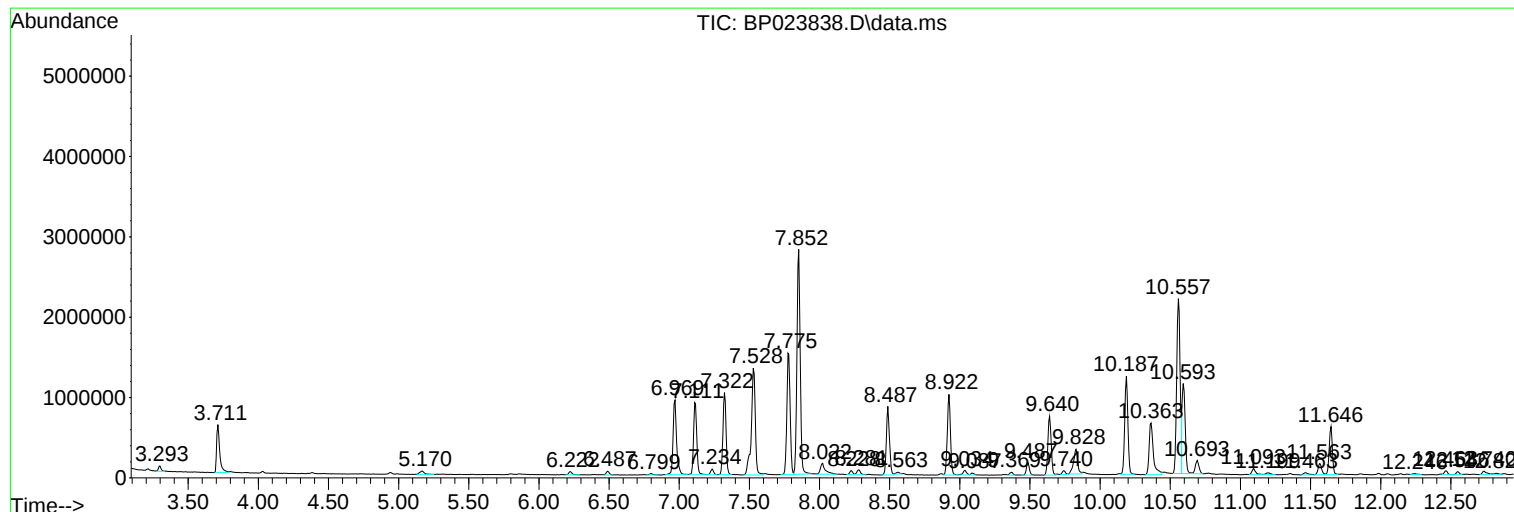
Sum of corrected areas: 104477003

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP013125\  
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Sample : Q1224-02  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
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A6302

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP012925.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\BP013125\  
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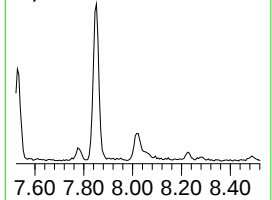
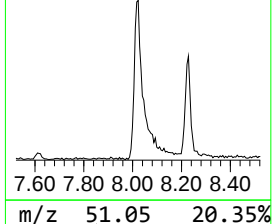
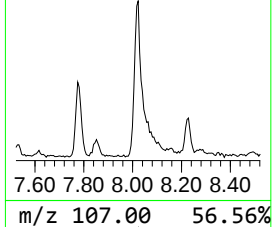
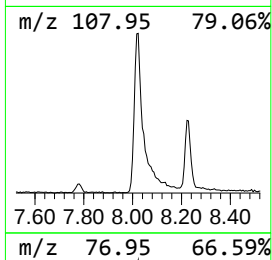
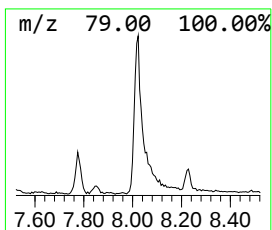
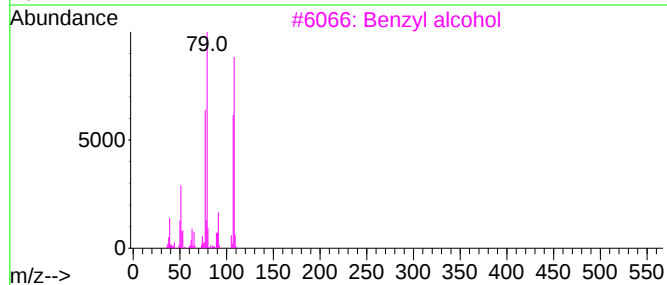
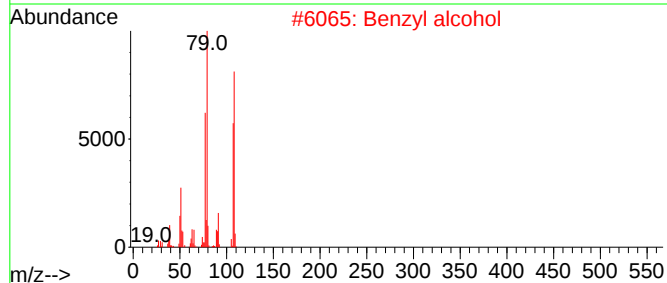
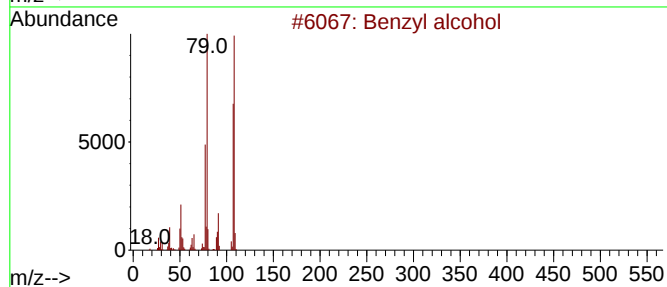
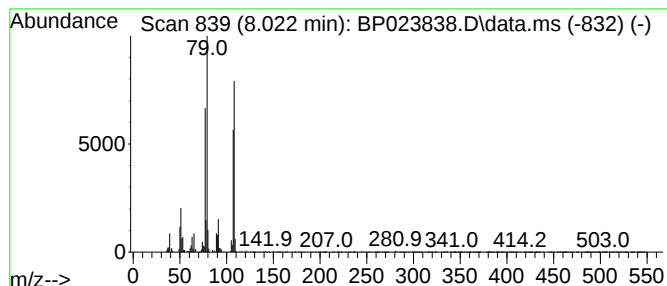
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP012925.MA.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 3 Benzyl alcohol Concentration Rank 20

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 8.022 | 2.75 ng/ul | 331857 | 1,4-Dichlorobenzene-d4 | 7.781 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm    | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|------------|-------------|------|
| 1    |    |   | Benzyl alcohol                      | 108 | C7H8O      | 000100-51-6 | 97   |
| 2    |    |   | Benzyl alcohol                      | 108 | C7H8O      | 000100-51-6 | 96   |
| 3    |    |   | Benzyl alcohol                      | 108 | C7H8O      | 000100-51-6 | 96   |
| 4    |    |   | Benzyl alcohol                      | 108 | C7H8O      | 000100-51-6 | 95   |
| 5    |    |   | N-Cbz-glycylglycine p-nitropheny... | 387 | C18H17N3O7 | 013574-81-7 | 91   |



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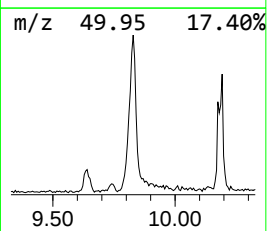
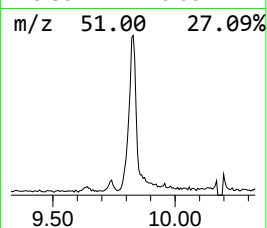
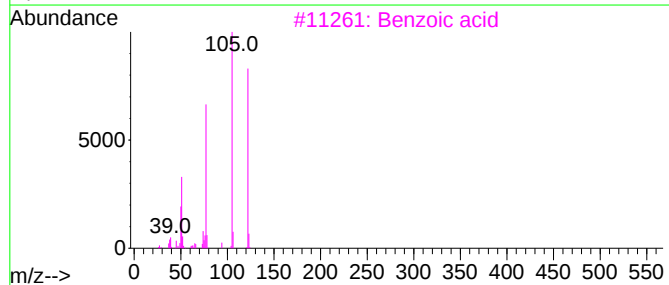
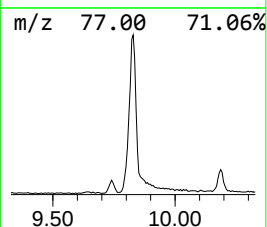
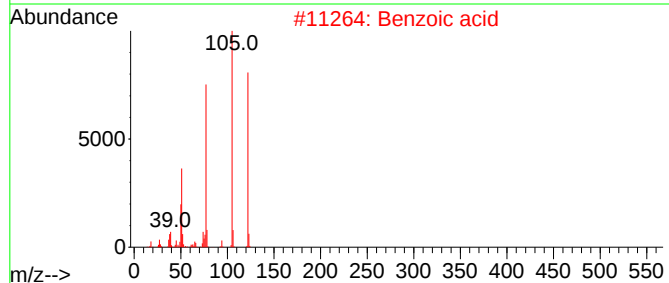
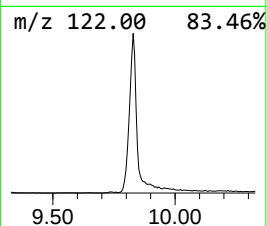
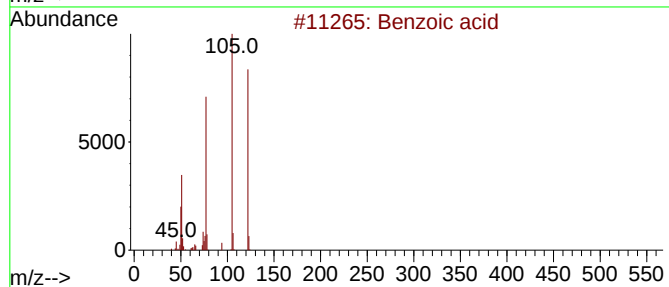
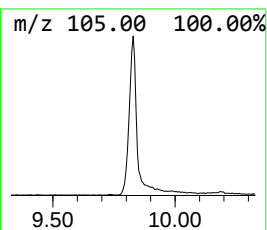
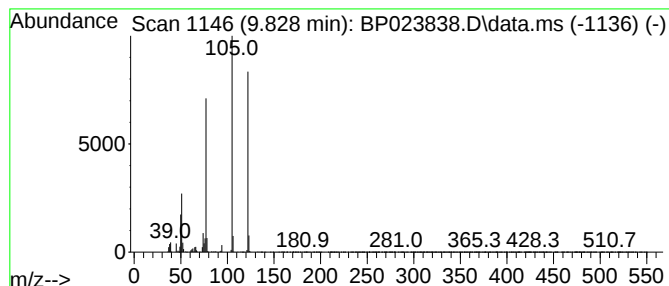
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP012925.MA.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 4 Benzoic acid Concentration Rank 18

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.   |
|-------|------------|--------|------------------|--------|
| 9.828 | 3.27 ng/ul | 630120 | Naphthalene-d8   | 10.557 |

| Hit# | of | 5 | Tentative ID                  | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------|-----|----------|-------------|------|
| 1    |    |   | Benzoic acid                  | 122 | C7H6O2   | 000065-85-0 | 96   |
| 2    |    |   | Benzoic acid                  | 122 | C7H6O2   | 000065-85-0 | 95   |
| 3    |    |   | Benzoic acid                  | 122 | C7H6O2   | 000065-85-0 | 93   |
| 4    |    |   | Benzoic acid                  | 122 | C7H6O2   | 000065-85-0 | 91   |
| 5    |    |   | Benzoic acid, silver(1+) salt | 228 | C7H5AgO2 | 000532-31-0 | 90   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP013125\  
Data File : BP023838.D  
Acq On : 01 Feb 2025 00:10  
Operator : RC/JU  
Sample : Q1224-02  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
A6302

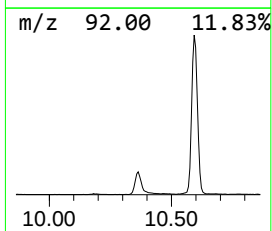
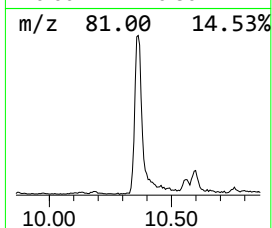
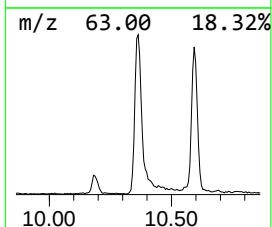
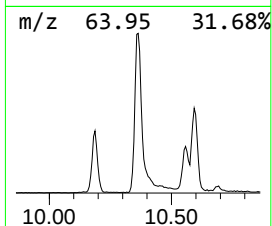
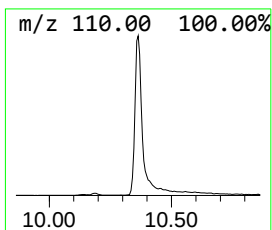
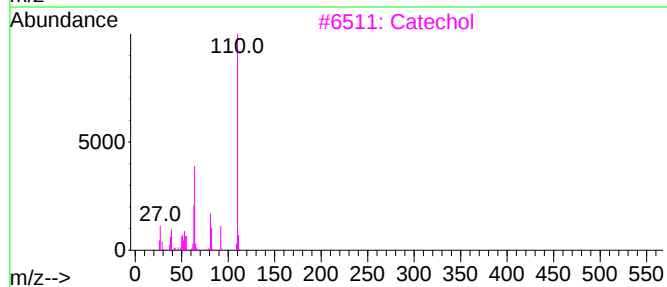
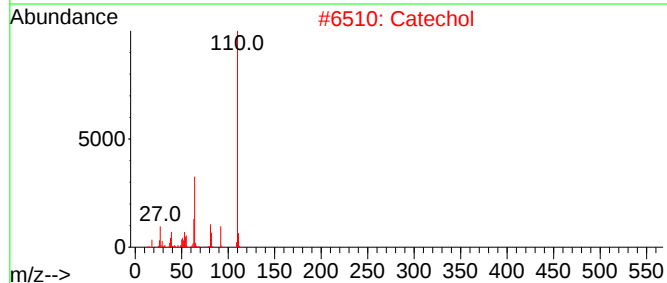
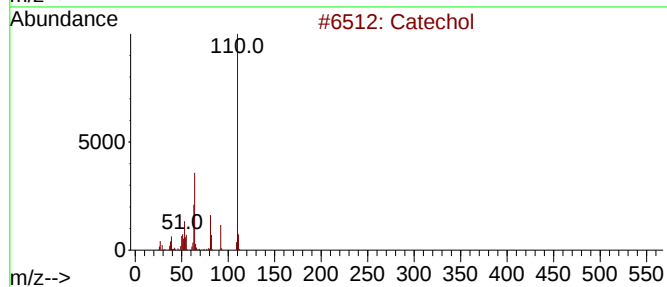
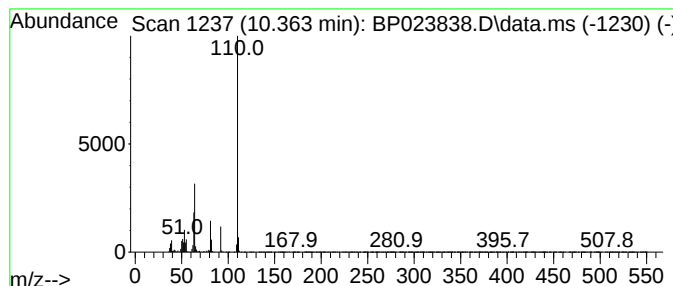
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Quant Title : SVOA CALIBRATION

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

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Peak Number 5 Catechol Concentration Rank 6

| R.T.   | EstConc    | Area    | Relative to ISTD | R.T.   |
|--------|------------|---------|------------------|--------|
| 10.363 | 6.96 ng/ul | 1339260 | Naphthalene-d8   | 10.557 |

| Hit# | of | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------|-----|---------|-------------|------|
| 1    |    |   | Catechol     | 110 | C6H6O2  | 000120-80-9 | 96   |
| 2    |    |   | Catechol     | 110 | C6H6O2  | 000120-80-9 | 91   |
| 3    |    |   | Catechol     | 110 | C6H6O2  | 000120-80-9 | 87   |
| 4    |    |   | Resorcinol   | 110 | C6H6O2  | 000108-46-3 | 59   |
| 5    |    |   | Resorcinol   | 110 | C6H6O2  | 000108-46-3 | 59   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP013125\  
Data File : BP023838.D  
Acq On : 01 Feb 2025 00:10  
Operator : RC/JU  
Sample : Q1224-02  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
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Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP012925.MA.M  
Quant Title : SVOA CALIBRATION

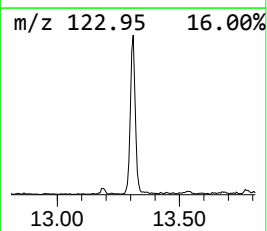
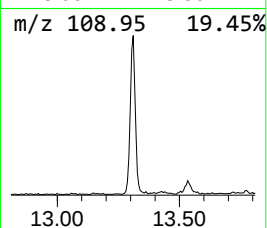
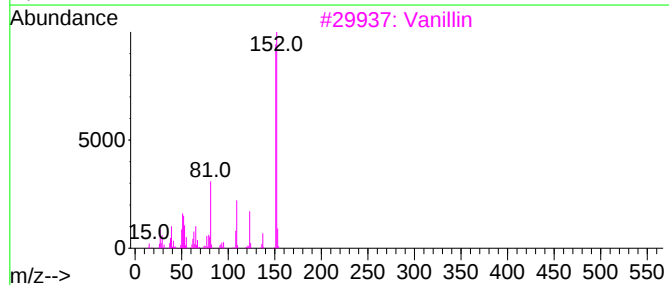
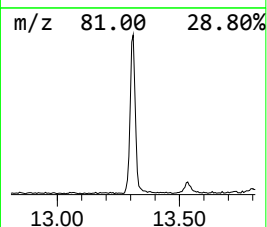
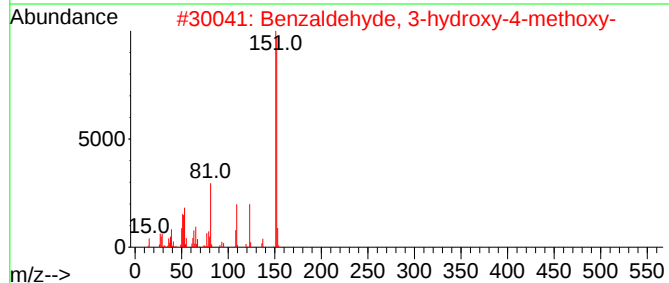
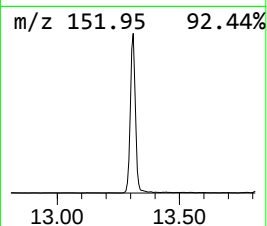
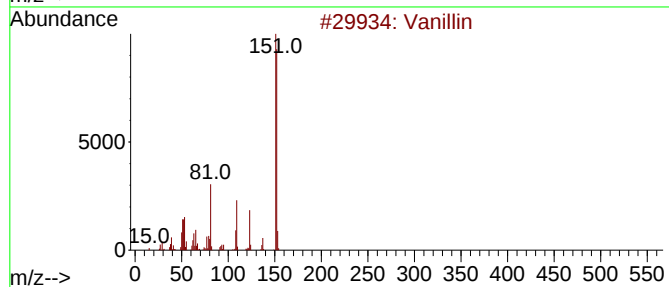
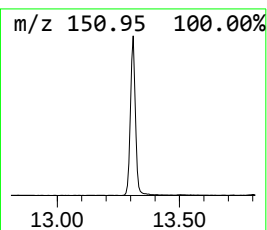
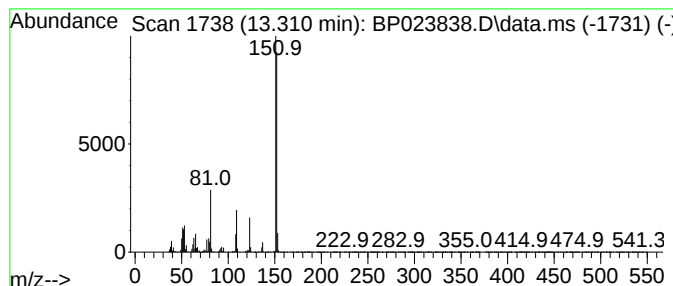
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TIC Integration Parameters: LSCINT.P

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Peak Number 7 Vanillin Concentration Rank 17

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 13.310 | 3.31 ng/ul | 810317 | Acenaphthene-d10 | 14.399 |

| Hit# | of | 5 | Tentative ID                       | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Vanillin                           | 152 | C8H8O3  | 000121-33-5 | 97   |
| 2    |    |   | Benzaldehyde, 3-hydroxy-4-methoxy- | 152 | C8H8O3  | 000621-59-0 | 96   |
| 3    |    |   | Vanillin                           | 152 | C8H8O3  | 000121-33-5 | 96   |
| 4    |    |   | Benzaldehyde, 3-hydroxy-4-methoxy- | 152 | C8H8O3  | 000621-59-0 | 96   |
| 5    |    |   | Benzaldehyde, 3-hydroxy-4-methoxy- | 152 | C8H8O3  | 000621-59-0 | 96   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP013125\  
Data File : BP023838.D  
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Operator : RC/JU  
Sample : Q1224-02  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

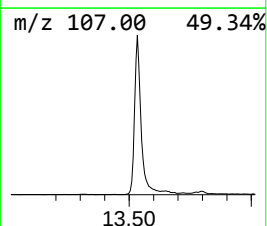
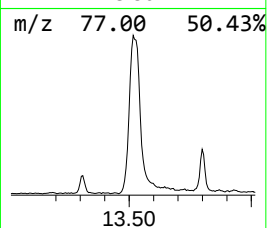
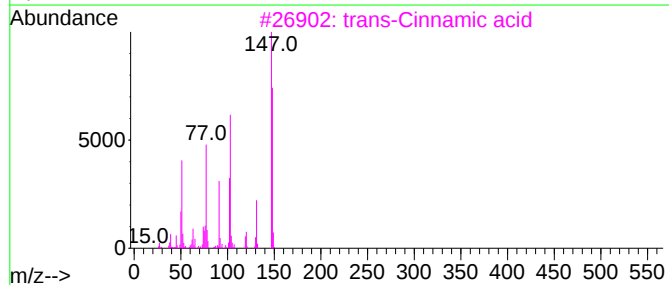
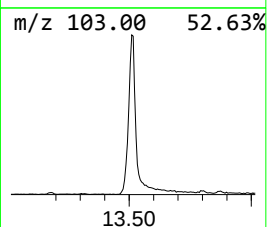
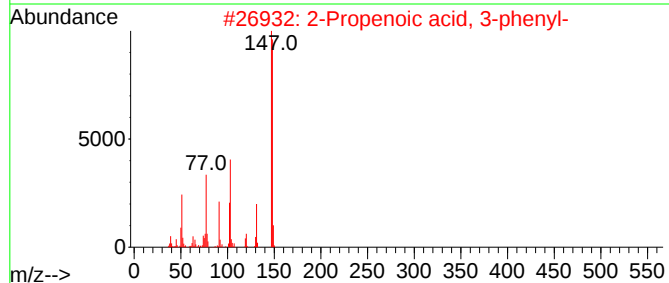
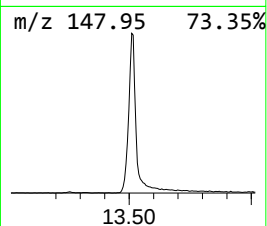
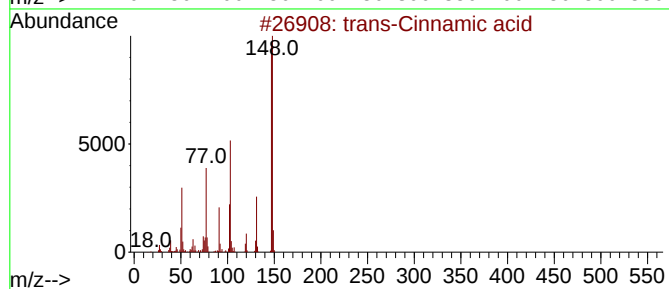
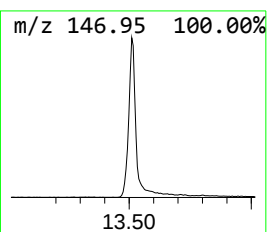
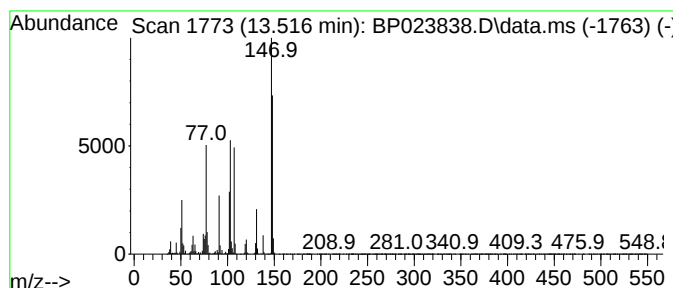
Instrument :  
BNA\_P  
ClientSampleId :  
A6302

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP012925.MA.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 8 trans-Cinnamic acid Concentration Rank 3

| R.T.      | EstConc                     | Area    | Relative to ISTD | R.T.        |      |
|-----------|-----------------------------|---------|------------------|-------------|------|
| 13.516    | 10.90 ng/ul                 | 2670300 | Acenaphthene-d10 | 14.399      |      |
| Hit# of 5 | Tentative ID                | MW      | MolForm          | CAS#        | Qual |
| 1         | trans-Cinnamic acid         | 148     | C9H8O2           | 000140-10-3 | 96   |
| 2         | 2-Propenoic acid, 3-phenyl- | 148     | C9H8O2           | 000621-82-9 | 95   |
| 3         | trans-Cinnamic acid         | 148     | C9H8O2           | 000140-10-3 | 95   |
| 4         | 2-Propenoic acid, 3-phenyl- | 148     | C9H8O2           | 000621-82-9 | 93   |
| 5         | trans-Cinnamic acid         | 148     | C9H8O2           | 000140-10-3 | 93   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP013125\  
Data File : BP023838.D  
Acq On : 01 Feb 2025 00:10  
Operator : RC/JU  
Sample : Q1224-02  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
A6302

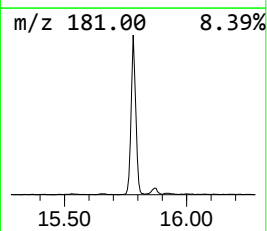
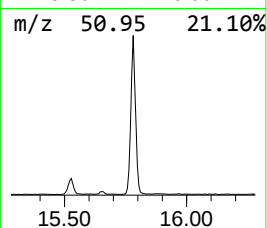
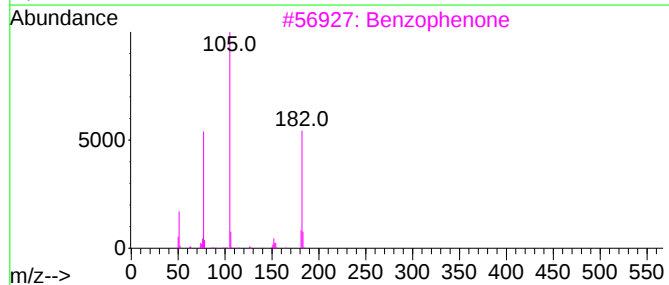
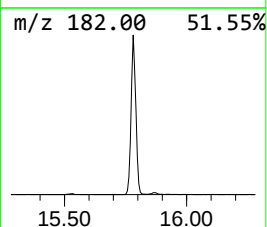
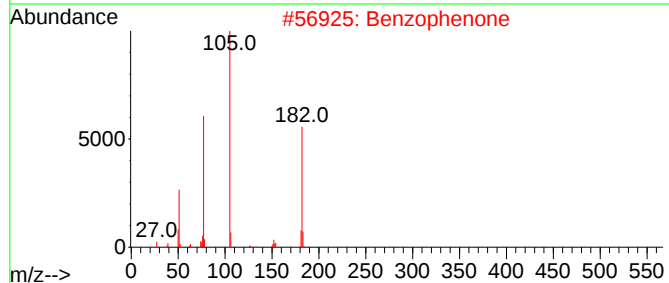
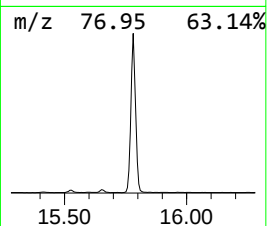
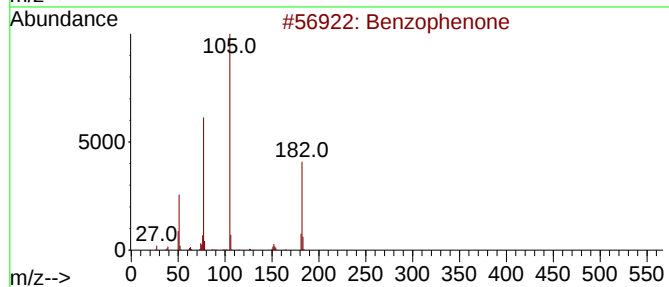
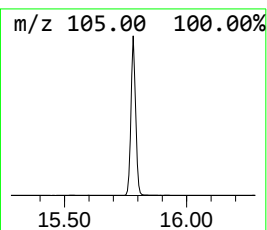
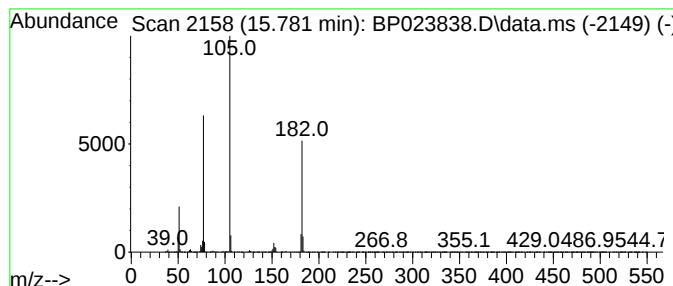
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Quant Title : SVOA CALIBRATION

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

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Peak Number 9 Benzophenone Concentration Rank 4

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 15.781 | 10.54 ng/ul | 2582160 | Acenaphthene-d10 | 14.399 |

| Hit# | of | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------|-----|---------|-------------|------|
| 1    |    |   | Benzophenone | 182 | C13H10O | 000119-61-9 | 97   |
| 2    |    |   | Benzophenone | 182 | C13H10O | 000119-61-9 | 96   |
| 3    |    |   | Benzophenone | 182 | C13H10O | 000119-61-9 | 96   |
| 4    |    |   | Benzophenone | 182 | C13H10O | 000119-61-9 | 95   |
| 5    |    |   | Benzophenone | 182 | C13H10O | 000119-61-9 | 91   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP013125\  
Data File : BP023838.D  
Acq On : 01 Feb 2025 00:10  
Operator : RC/JU  
Sample : Q1224-02  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
A6302

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP012925.MA.M  
Quant Title : SVOA CALIBRATION

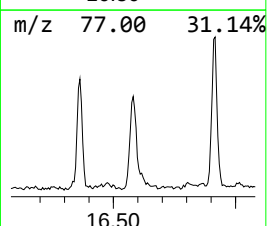
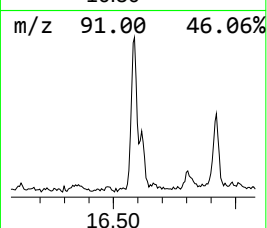
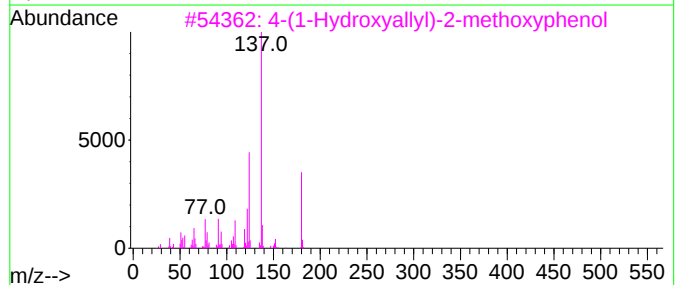
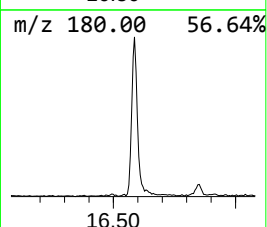
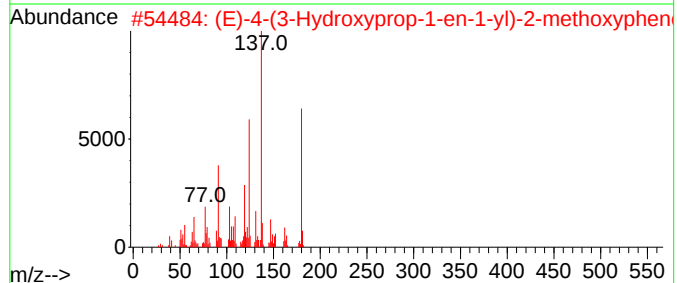
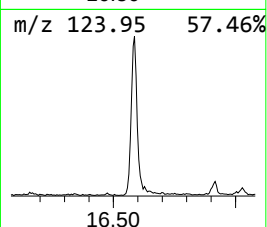
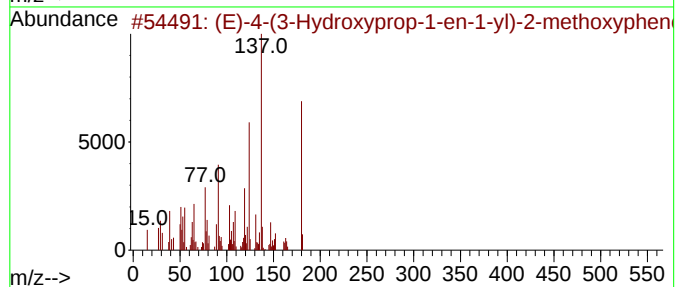
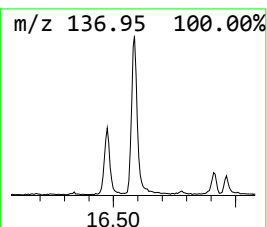
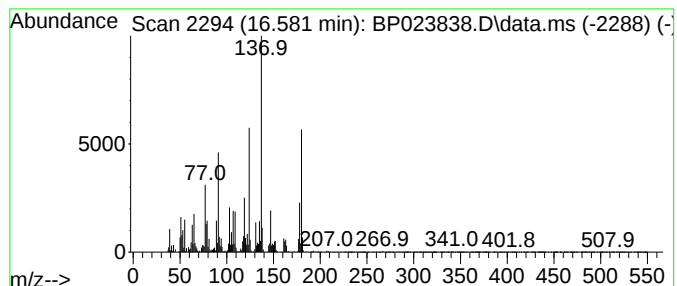
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TIC Integration Parameters: LSCINT.P

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Peak Number 10 (E)-4-(3-Hydroxyprop-1-en-1-yl) Concentration Rank 13

| R.T.   | EstConc    | Area    | Relative to ISTD | R.T.   |
|--------|------------|---------|------------------|--------|
| 16.581 | 3.78 ng/ul | 1066230 | Phenanthrene-d10 | 17.204 |

| Hit# | of                                  | 5   | Tentative ID | MW          | MolForm | CAS# | Qual |
|------|-------------------------------------|-----|--------------|-------------|---------|------|------|
| 1    | (E)-4-(3-Hydroxyprop-1-en-1-yl)-... | 180 | C10H12O3     | 032811-40-8 | 97      |      |      |
| 2    | (E)-4-(3-Hydroxyprop-1-en-1-yl)-... | 180 | C10H12O3     | 032811-40-8 | 90      |      |      |
| 3    | 4-(1-Hydroxyallyl)-2-methoxyphenol  | 180 | C10H12O3     | 112465-50-6 | 64      |      |      |
| 4    | 3,7-Benzofurandiyl, 2,3-dihydro-... | 180 | C10H12O3     | 017781-15-6 | 46      |      |      |
| 5    | 1-Hydroxy-1-(4-methoxyphenyl)pro... | 180 | C10H12O3     | 015482-29-8 | 45      |      |      |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP013125\  
Data File : BP023838.D  
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Operator : RC/JU  
Sample : Q1224-02  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

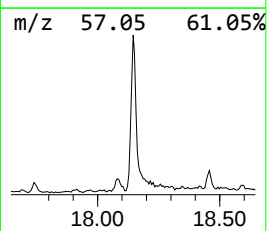
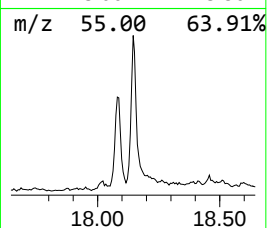
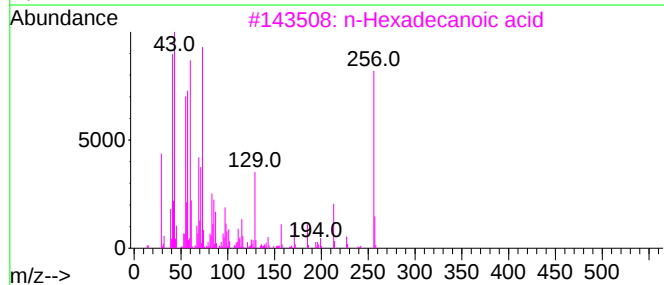
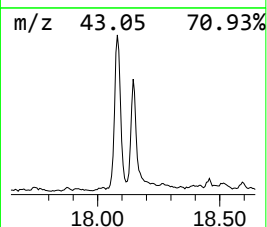
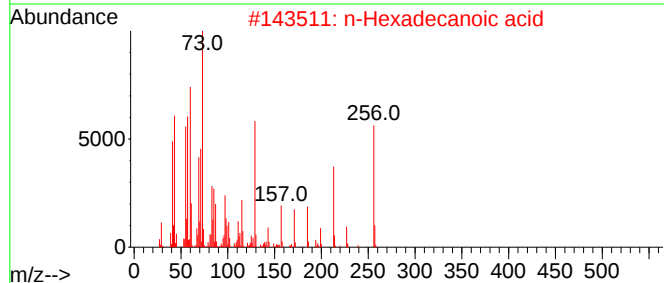
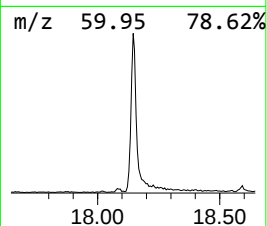
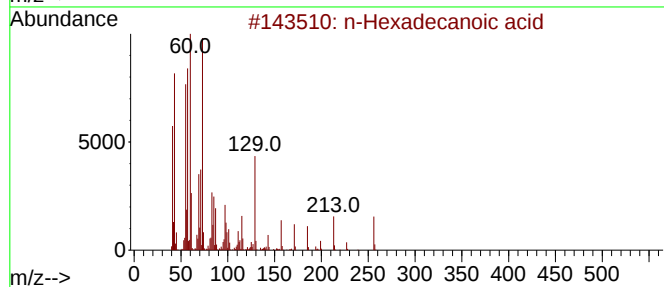
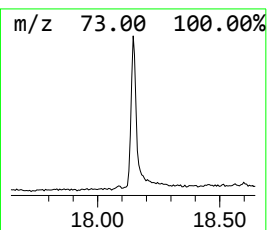
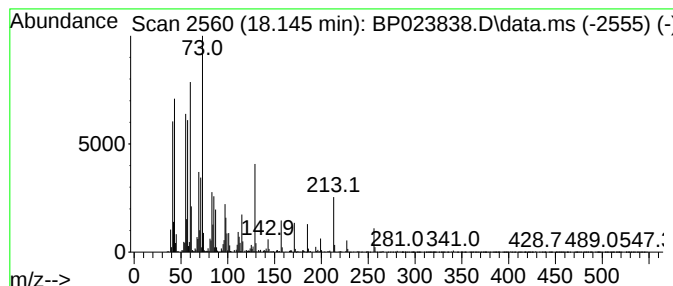
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ClientSampleId :  
A6302

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP012925.MA.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 13 n-Hexadecanoic acid Concentration Rank 14

| R.T.      | EstConc             | Area    | Relative to ISTD | R.T.        |      |
|-----------|---------------------|---------|------------------|-------------|------|
| 18.145    | 3.65 ng/ul          | 1029360 | Phenanthrene-d10 | 17.204      |      |
| Hit# of 5 | Tentative ID        | MW      | MolForm          | CAS#        | Qual |
| 1         | n-Hexadecanoic acid | 256     | C16H32O2         | 000057-10-3 | 99   |
| 2         | n-Hexadecanoic acid | 256     | C16H32O2         | 000057-10-3 | 99   |
| 3         | n-Hexadecanoic acid | 256     | C16H32O2         | 000057-10-3 | 99   |
| 4         | n-Hexadecanoic acid | 256     | C16H32O2         | 000057-10-3 | 99   |
| 5         | Tridecanoic acid    | 214     | C13H26O2         | 000638-53-9 | 94   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP013125\  
Data File : BP023838.D  
Acq On : 01 Feb 2025 00:10  
Operator : RC/JU  
Sample : Q1224-02  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
A6302

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP012925.MA.M  
Quant Title : SVOA CALIBRATION

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

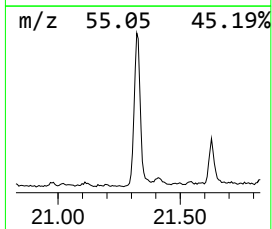
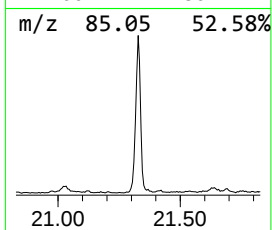
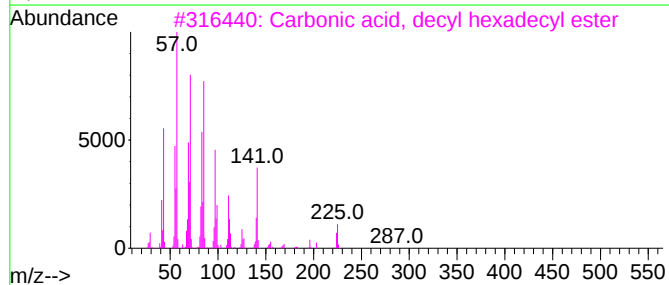
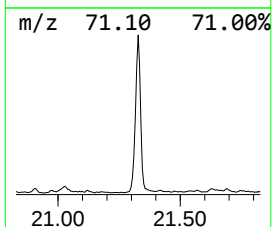
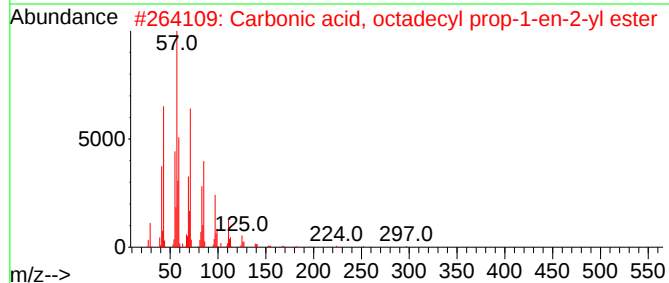
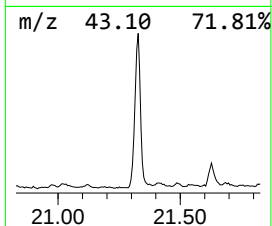
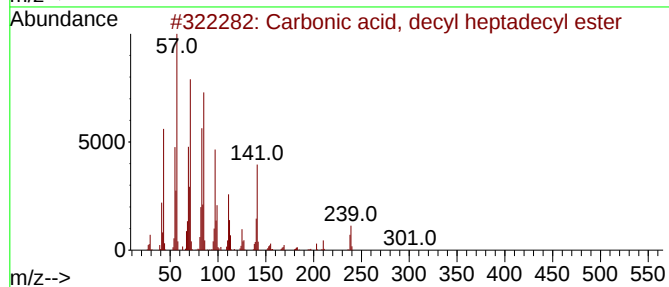
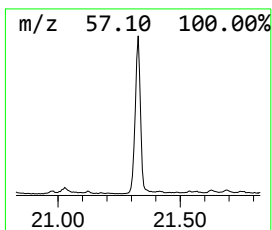
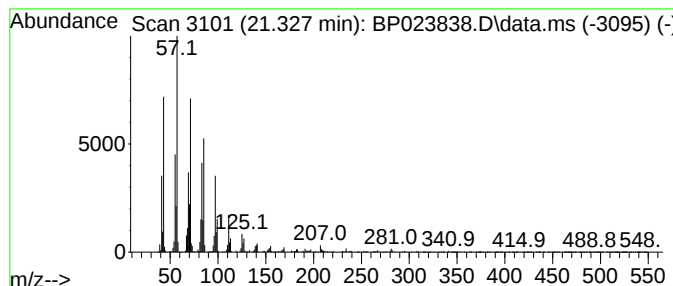
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Peak Number 14 Carbonic acid, decyl heptad... Concentration Rank 15

| R.T.   | EstConc    | Area    | Relative to ISTD | R.T.   |
|--------|------------|---------|------------------|--------|
| 21.328 | 3.65 ng/ul | 1754140 | Chrysene-d12     | 21.645 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Carbonic acid, decyl heptadecyl ... | 440 | C28H56O3 | 1000383-16-6 | 91   |
| 2    |    |   | Carbonic acid, octadecyl prop-1-... | 354 | C22H42O3 | 1000383-11-5 | 91   |
| 3    |    |   | Carbonic acid, decyl hexadecyl e... | 426 | C27H54O3 | 1000383-16-5 | 91   |
| 4    |    |   | Carbonic acid, decyl pentadecyl ... | 412 | C26H52O3 | 1000383-16-4 | 91   |
| 5    |    |   | Carbonic acid, eicosyl vinyl ester  | 368 | C23H44O3 | 1000382-54-3 | 91   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP013125\  
Data File : BP023838.D  
Acq On : 01 Feb 2025 00:10  
Operator : RC/JU  
Sample : Q1224-02  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
A6302

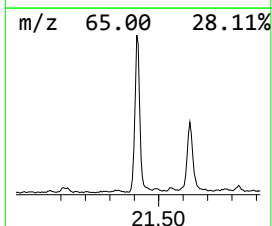
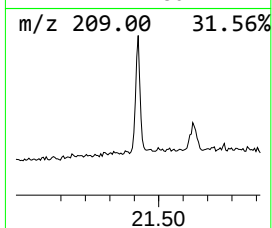
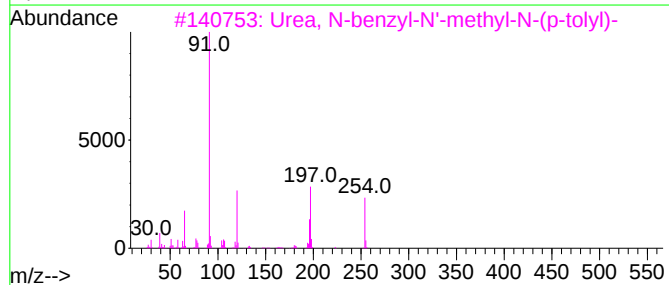
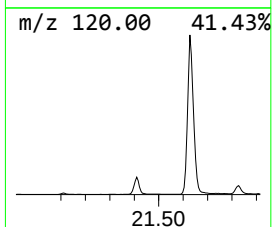
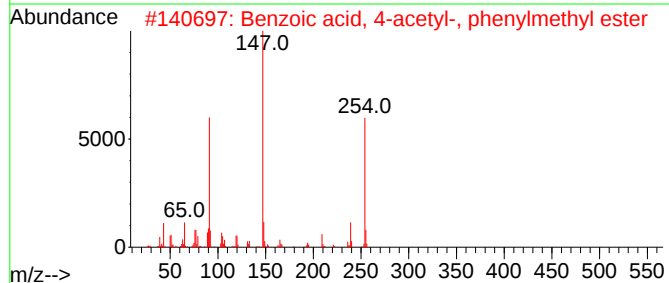
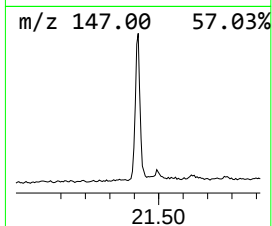
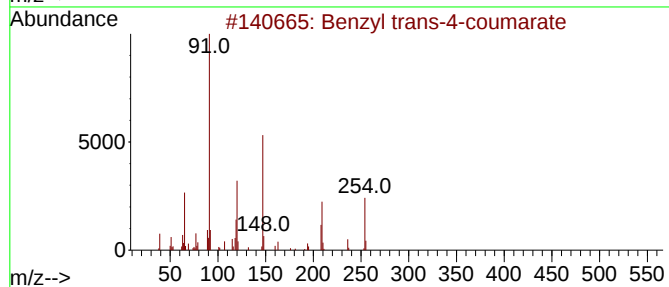
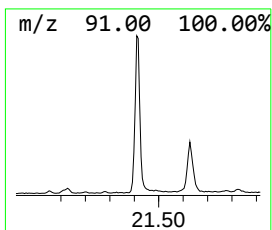
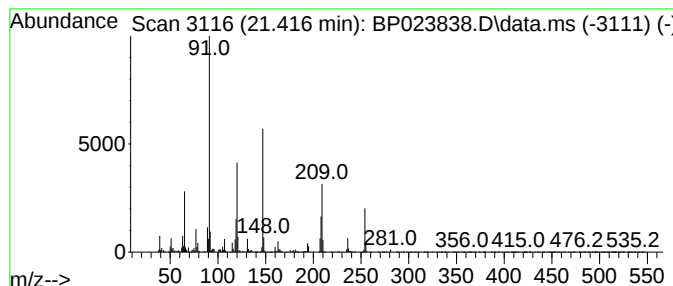
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Quant Title : SVOA CALIBRATION

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

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Peak Number 15 Benzyl trans-4-coumarate Concentration Rank 19

| R.T.   | EstConc    | Area    | Relative to ISTD | R.T.   |
|--------|------------|---------|------------------|--------|
| 21.416 | 2.79 ng/ul | 1342570 | Chrysene-d12     | 21.645 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Benzyl trans-4-coumarate            | 254 | C16H14O3  | 061844-62-0  | 98   |
| 2    |    |   | Benzoic acid, 4-acetyl-, phenylm... | 254 | C16H14O3  | 119838-69-6  | 38   |
| 3    |    |   | Urea, N-benzyl-N'-methyl-N-(p-to... | 254 | C16H18N2O | 1000152-05-0 | 25   |
| 4    |    |   | Pentafluoropropionic acid, benzy... | 254 | C10H7F5O2 | 076008-52-1  | 25   |
| 5    |    |   | 2H-Benzotriazole, 2-ethyl-          | 147 | C8H9N3    | 016584-04-6  | 22   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP013125\  
Data File : BP023838.D  
Acq On : 01 Feb 2025 00:10  
Operator : RC/JU  
Sample : Q1224-02  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
A6302

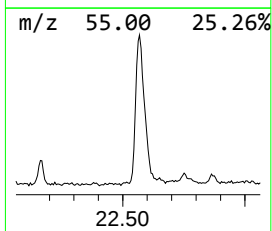
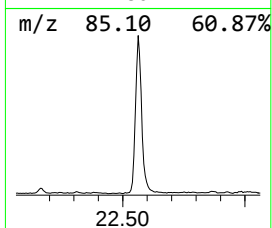
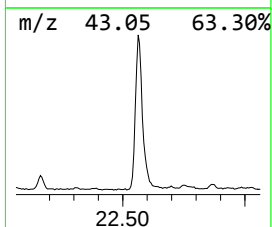
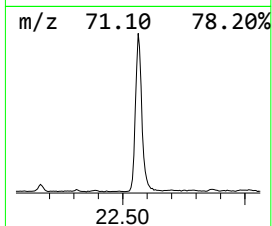
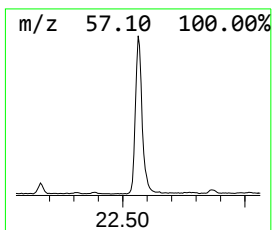
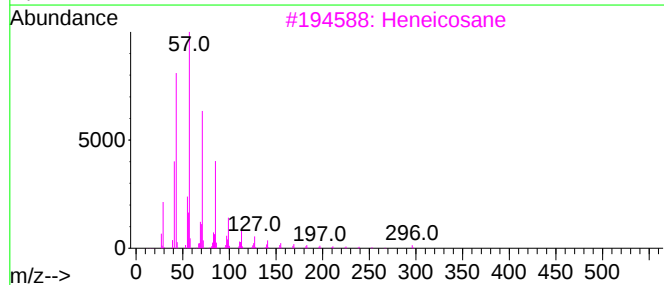
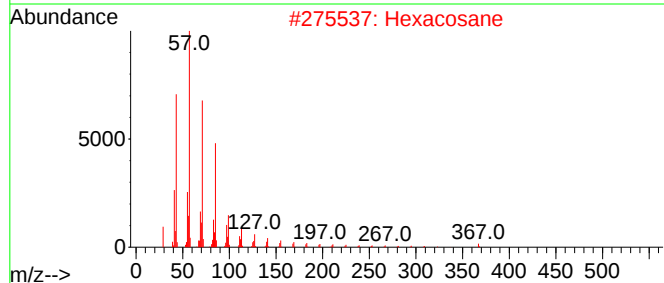
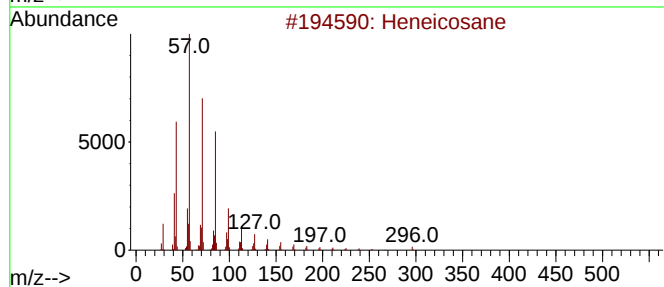
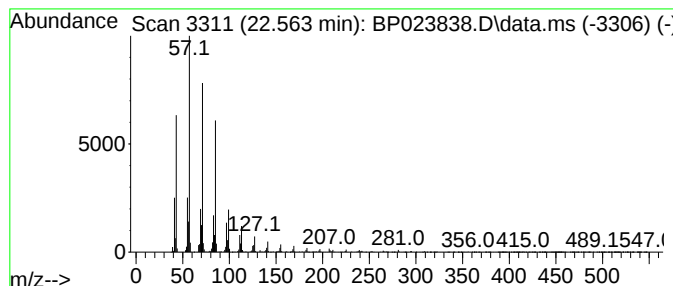
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Quant Title : SVOA CALIBRATION

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

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Peak Number 16 (DEL) Alkane: Straight-Chai... Concentration Rank 9

| R.T.   | EstConc    | Area    | Relative to ISTD | R.T.   |
|--------|------------|---------|------------------|--------|
| 22.563 | 4.67 ng/ul | 2245690 | Chrysene-d12     | 21.645 |

| Hit# | of | 5 | Tentative ID       | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------|-----|---------|-------------|------|
| 1    |    |   | Heneicosane        | 296 | C21H44  | 000629-94-7 | 91   |
| 2    |    |   | Hexacosane         | 366 | C26H54  | 000630-01-3 | 91   |
| 3    |    |   | Heneicosane        | 296 | C21H44  | 000629-94-7 | 91   |
| 4    |    |   | 2-Methylhexacosane | 380 | C27H56  | 001561-02-0 | 90   |
| 5    |    |   | Heptadecane        | 240 | C17H36  | 000629-78-7 | 87   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP013125\  
Data File : BP023838.D  
Acq On : 01 Feb 2025 00:10  
Operator : RC/JU  
Sample : Q1224-02  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
A6302

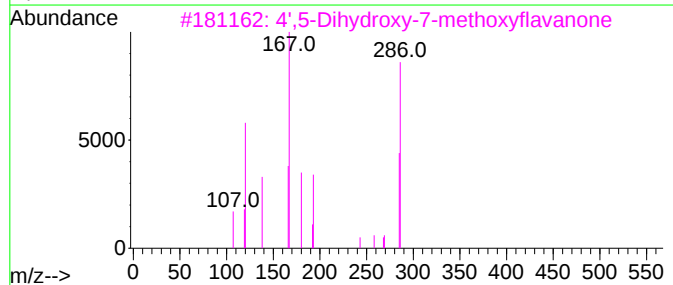
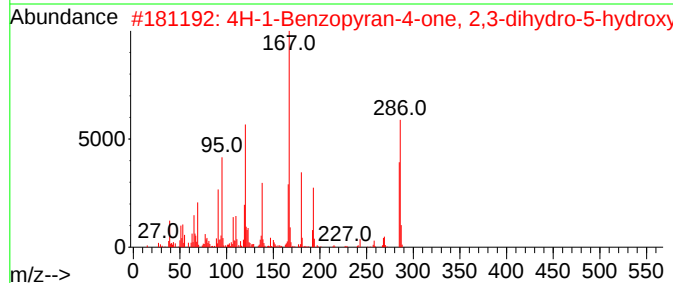
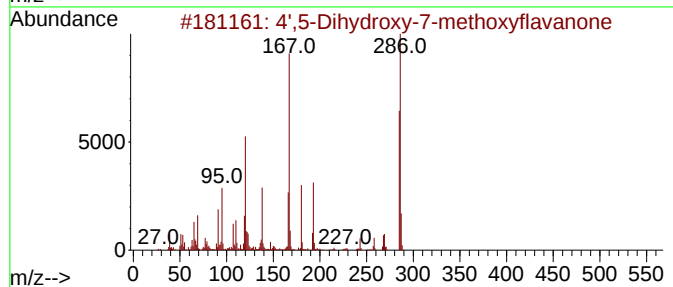
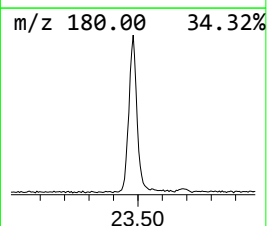
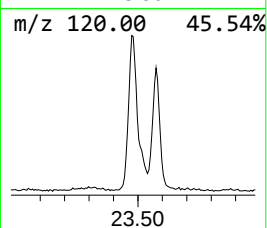
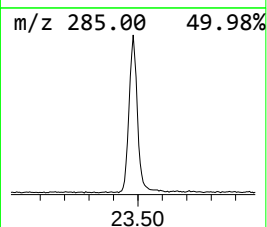
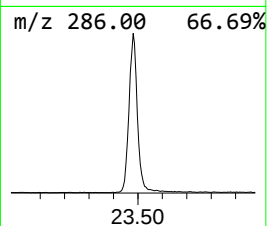
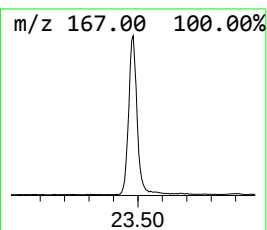
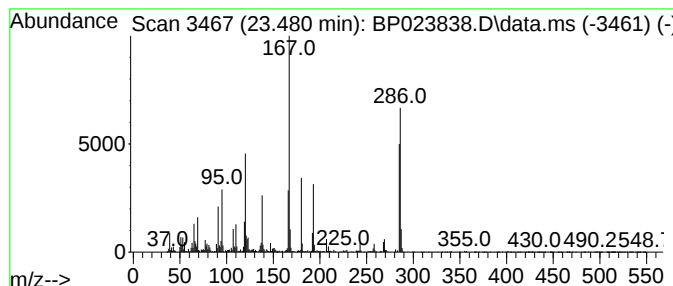
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Quant Title : SVOA CALIBRATION

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

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Peak Number 17 4',5-Dihydroxy-7-methoxyfla... Concentration Rank 7

| R.T.   | EstConc    | Area    | Relative to ISTD | R.T.   |
|--------|------------|---------|------------------|--------|
| 23.480 | 6.88 ng/ul | 2154170 | Perylene-d12     | 25.021 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm   | CAS#        | Qual |
|------|------|-------------------------------------|-----|-----------|-------------|------|
| 1    |      | 4',5-Dihydroxy-7-methoxyflavanone   | 286 | C16H14O5  | 000520-29-6 | 99   |
| 2    |      | 4H-1-Benzopyran-4-one, 2,3-dihyd... | 286 | C16H14O5  | 002957-21-3 | 99   |
| 3    |      | 4',5-Dihydroxy-7-methoxyflavanone   | 286 | C16H14O5  | 000520-29-6 | 96   |
| 4    |      | 2-Propen-1-one, 1-(2,4-dihydroxy... | 286 | C16H14O5  | 069707-17-1 | 86   |
| 5    |      | Sakuranin                           | 448 | C22H24O10 | 000529-39-5 | 72   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP013125\  
Data File : BP023838.D  
Acq On : 01 Feb 2025 00:10  
Operator : RC/JU  
Sample : Q1224-02  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
A6302

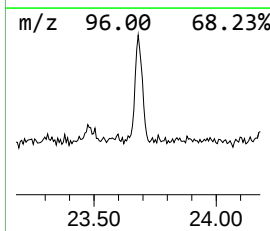
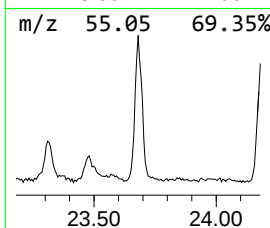
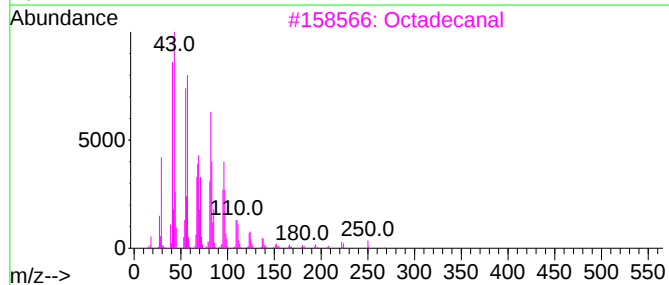
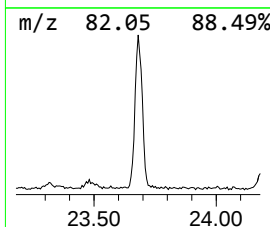
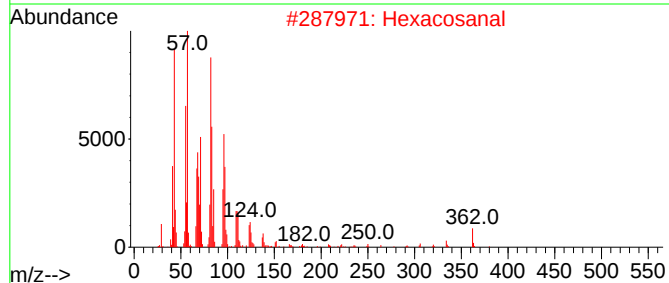
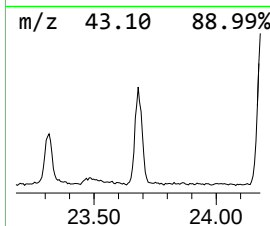
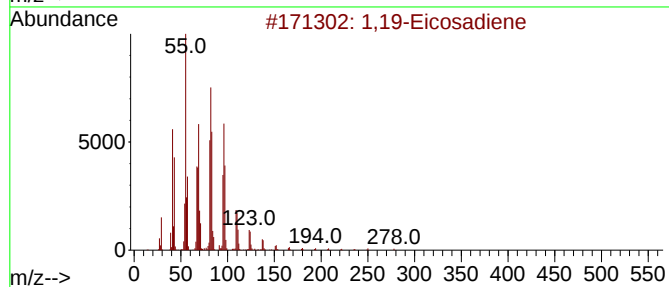
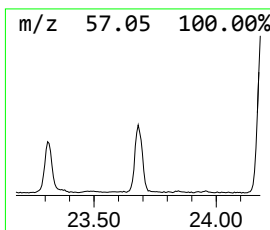
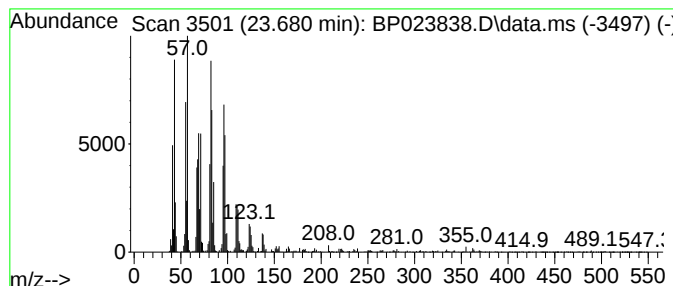
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Quant Title : SVOA CALIBRATION

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

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Peak Number 18 1,19-Eicosadiene Concentration Rank 12

| R.T.   | EstConc    | Area    | Relative to ISTD | R.T.   |
|--------|------------|---------|------------------|--------|
| 23.680 | 3.85 ng/ul | 1203610 | Perylene-d12     | 25.021 |

| Hit# | of 5 | Tentative ID     | MW  | MolForm | CAS#        | Qual |
|------|------|------------------|-----|---------|-------------|------|
| 1    |      | 1,19-Eicosadiene | 278 | C20H38  | 014811-95-1 | 95   |
| 2    |      | Hexacosanal      | 380 | C26H52O | 026627-85-0 | 94   |
| 3    |      | Octadecanal      | 268 | C18H36O | 000638-66-4 | 91   |
| 4    |      | Octadecanal      | 268 | C18H36O | 000638-66-4 | 91   |
| 5    |      | Triacontanal     | 436 | C30H60O | 022725-63-9 | 91   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP013125\  
Data File : BP023838.D  
Acq On : 01 Feb 2025 00:10  
Operator : RC/JU  
Sample : Q1224-02  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
A6302

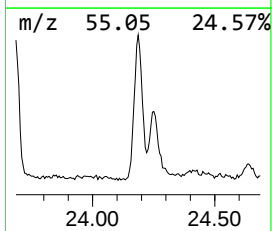
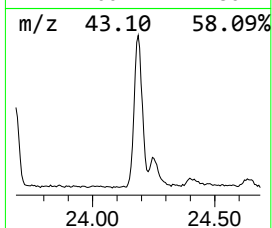
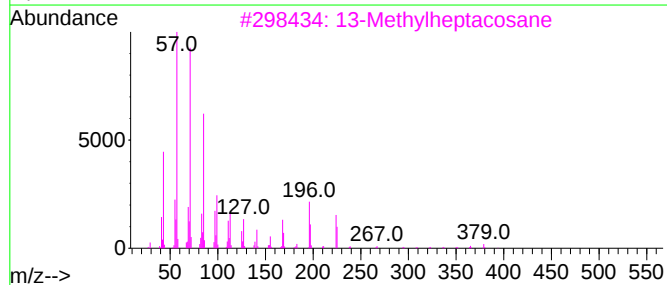
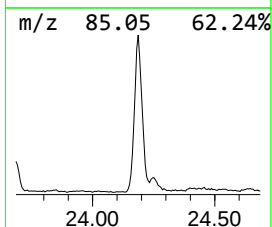
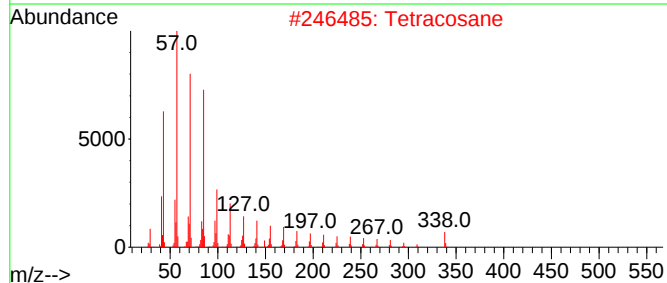
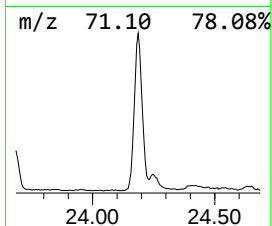
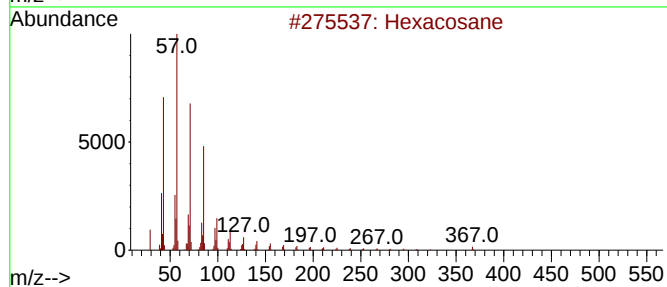
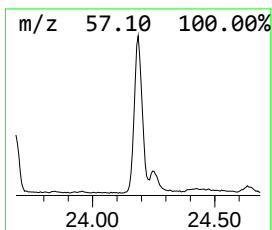
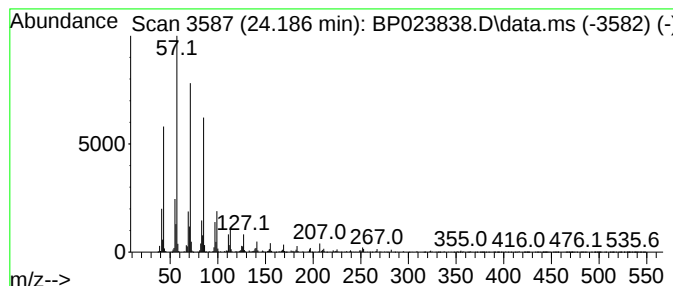
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Quant Title : SVOA CALIBRATION

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

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Peak Number 19 (DEL) Alkane: Straight-Chai... Concentration Rank 10

| R.T.   | EstConc    | Area    | Relative to ISTD | R.T.   |
|--------|------------|---------|------------------|--------|
| 24.186 | 4.00 ng/ul | 1251170 | Perylene-d12     | 25.021 |

| Hit# | of 5 | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|------|-------------------------|-----|---------|-------------|------|
| 1    |      | Hexacosane              | 366 | C26H54  | 000630-01-3 | 94   |
| 2    |      | Tetracosane             | 338 | C24H50  | 000646-31-1 | 94   |
| 3    |      | 13-Methylheptacosane    | 394 | C28H58  | 015689-72-2 | 94   |
| 4    |      | Docosane, 11-butyl-     | 366 | C26H54  | 013475-76-8 | 94   |
| 5    |      | Heneicosane, 11-pentyl- | 366 | C26H54  | 014739-72-1 | 91   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP013125\  
Data File : BP023838.D  
Acq On : 01 Feb 2025 00:10  
Operator : RC/JU  
Sample : Q1224-02  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
A6302

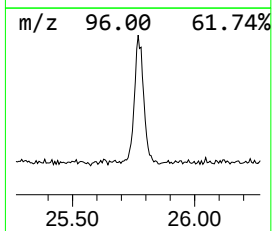
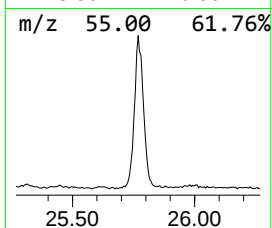
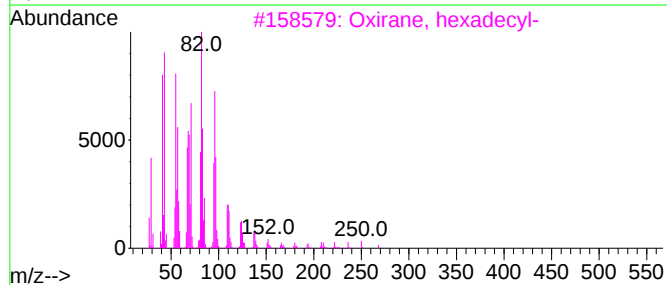
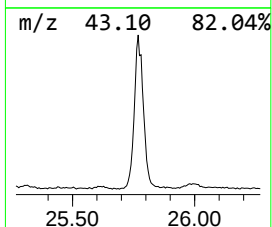
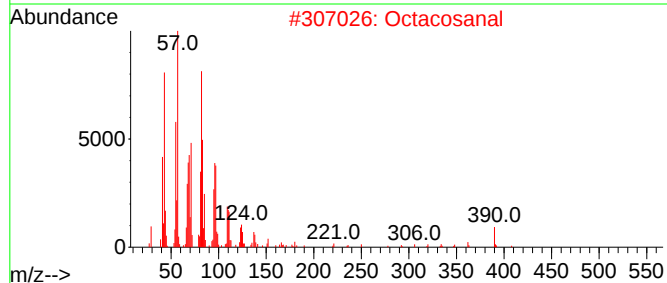
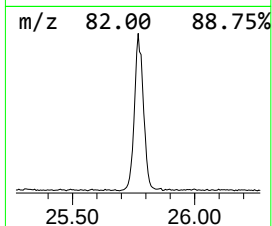
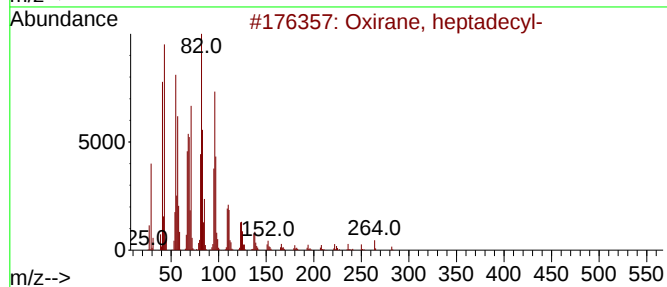
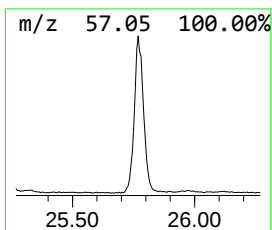
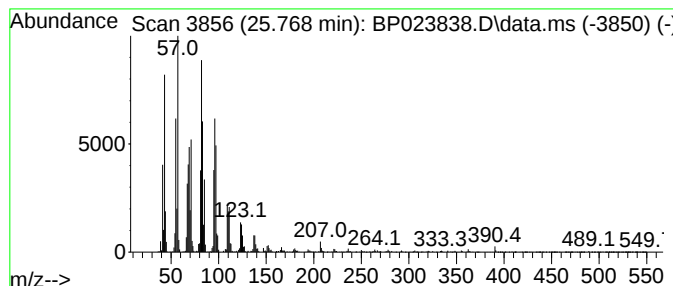
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP012925.MA.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 20 Oxirane, heptadecyl- Concentration Rank 5

| R.T.   | EstConc    | Area    | Relative to ISTD | R.T.   |
|--------|------------|---------|------------------|--------|
| 25.768 | 9.76 ng/ul | 3055040 | Perylene-d12     | 25.021 |

| Hit# | of 5 | Tentative ID         | MW  | MolForm | CAS#        | Qual |
|------|------|----------------------|-----|---------|-------------|------|
| 1    |      | Oxirane, heptadecyl- | 282 | C19H38O | 067860-04-2 | 91   |
| 2    |      | Octacosanal          | 408 | C28H56O | 022725-64-0 | 91   |
| 3    |      | Oxirane, hexadecyl-  | 268 | C18H36O | 007390-81-0 | 91   |
| 4    |      | Oxirane, hexadecyl-  | 268 | C18H36O | 007390-81-0 | 91   |
| 5    |      | Octadecanal          | 268 | C18H36O | 000638-66-4 | 91   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP013125\  
Data File : BP023838.D  
Acq On : 01 Feb 2025 00:10  
Operator : RC/JU  
Sample : Q1224-02  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
A6302

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP012925.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 |
| Benzyl alcohol     | 8.022  | 2.8     | ng/ul | 331857   | 1                     | 7.781  | 2413620 | 20.0 |
| Benzoic acid       | 9.828  | 3.3     | ng/ul | 630120   | 2                     | 10.557 | 3848160 | 20.0 |
| Catechol           | 10.363 | 7.0     | ng/ul | 1339260  | 2                     | 10.557 | 3848160 | 20.0 |
| Vanillin           | 13.310 | 3.3     | ng/ul | 810317   | 3                     | 14.399 | 4900030 | 20.0 |
| trans-Cinnamic ... | 13.516 | 10.9    | ng/ul | 2670300  | 3                     | 14.399 | 4900030 | 20.0 |
| Benzophenone       | 15.781 | 10.5    | ng/ul | 2582160  | 3                     | 14.399 | 4900030 | 20.0 |
| (E)-4-(3-Hydrox... | 16.581 | 3.8     | ng/ul | 1066230  | 4                     | 17.204 | 5635020 | 20.0 |
| n-Hexadecanoic ... | 18.145 | 3.6     | ng/ul | 1029360  | 4                     | 17.204 | 5635020 | 20.0 |
| Carbonic acid, ... | 21.328 | 3.6     | ng/ul | 1754140  | 5                     | 21.645 | 9624560 | 20.0 |
| Benzyl trans-4-... | 21.416 | 2.8     | ng/ul | 1342570  | 5                     | 21.645 | 9624560 | 20.0 |
| (DEL) Alkane: S... | 22.563 | 4.7     | ng/ul | 2245690  | 5                     | 21.645 | 9624560 | 20.0 |
| 4',5-Dihydroxy-... | 23.480 | 6.9     | ng/ul | 2154170  | 6                     | 25.021 | 6259800 | 20.0 |
| 1,19-Eicosadiene   | 23.680 | 3.9     | ng/ul | 1203610  | 6                     | 25.021 | 6259800 | 20.0 |
| (DEL) Alkane: S... | 24.186 | 4.0     | ng/ul | 1251170  | 6                     | 25.021 | 6259800 | 20.0 |
| Oxirane, heptad... | 25.768 | 9.8     | ng/ul | 3055040  | 6                     | 25.021 | 6259800 | 20.0 |