

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011722\  
 Data File : BG052058.D  
 Acq On : 18 Jan 2022 18:42  
 Operator : CG/JU  
 Sample : N1100-15ME  
 Misc :  
 ALS Vial : 44 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 BGLS8ME

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\_G\Methods\SFAM-EPA-BG010622.M  
 Title : SVOA CALIBRATION

Signal : TIC: BG052058.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         | 5.709       | 373           | 378         | 386          | rVV      | 575961         | 948505        | 6.57%           | 1.203%        |
| 2         | 5.850       | 395           | 402         | 420          | rVB      | 1338244        | 3056382       | 21.18%          | 3.878%        |
| 3         | 6.226       | 460           | 466         | 475          | rVB      | 753946         | 1229877       | 8.52%           | 1.560%        |
| 4         | 6.708       | 541           | 548         | 555          | rBV      | 516679         | 902880        | 6.26%           | 1.146%        |
| 5         | 6.866       | 570           | 575         | 588          | rVB3     | 93087          | 236275        | 1.64%           | 0.300%        |
| 6         | 7.213       | 624           | 634         | 640          | rBV2     | 252836         | 538379        | 3.73%           | 0.683%        |
| 7         | 7.319       | 647           | 652         | 657          | rVB      | 442851         | 695676        | 4.82%           | 0.883%        |
| 8         | 7.383       | 657           | 663         | 669          | rVB2     | 443398         | 746790        | 5.18%           | 0.948%        |
| 9         | 7.454       | 669           | 675         | 683          | rVB      | 314905         | 519679        | 3.60%           | 0.659%        |
| 10        | 7.624       | 699           | 704         | 709          | rBV      | 190578         | 327250        | 2.27%           | 0.415%        |
| 11        | 7.859       | 738           | 744         | 747          | rBV      | 919593         | 1560110       | 10.81%          | 1.979%        |
| 12        | 7.889       | 747           | 749         | 759          | rVB      | 894397         | 1199315       | 8.31%           | 1.522%        |
| 13        | 8.153       | 784           | 794         | 800          | rBV7     | 71650          | 203052        | 1.41%           | 0.258%        |
| 14        | 8.218       | 800           | 805         | 813          | rVB2     | 243250         | 450887        | 3.12%           | 0.572%        |
| 15        | 8.364       | 825           | 830         | 837          | rVB2     | 254876         | 464455        | 3.22%           | 0.589%        |
| 16        | 8.629       | 866           | 875         | 880          | rVB3     | 142364         | 315746        | 2.19%           | 0.401%        |
| 17        | 8.793       | 896           | 903         | 907          | rVV3     | 236215         | 571388        | 3.96%           | 0.725%        |
| 18        | 8.893       | 914           | 920         | 926          | rBV3     | 334287         | 773188        | 5.36%           | 0.981%        |
| 19        | 8.940       | 926           | 928         | 934          | rVB2     | 134481         | 205582        | 1.42%           | 0.261%        |
| 20        | 9.011       | 934           | 940         | 945          | rBV      | 161407         | 270381        | 1.87%           | 0.343%        |
| 21        | 9.263       | 978           | 983         | 990          | rVB      | 119765         | 223292        | 1.55%           | 0.283%        |
| 22        | 9.363       | 990           | 1000        | 1005         | rBV3     | 257495         | 541019        | 3.75%           | 0.686%        |
| 23        | 9.481       | 1012          | 1020        | 1026         | rVV      | 1037001        | 1780365       | 12.34%          | 2.259%        |
| 24        | 9.616       | 1032          | 1043        | 1052         | rBV9     | 116502         | 369957        | 2.56%           | 0.469%        |
| 25        | 9.745       | 1052          | 1065        | 1072         | rVV8     | 119407         | 422823        | 2.93%           | 0.536%        |
| 26        | 9.898       | 1086          | 1091        | 1096         | rVV3     | 155626         | 296069        | 2.05%           | 0.376%        |
| 27        | 9.956       | 1096          | 1101        | 1105         | rVV      | 176241         | 314753        | 2.18%           | 0.399%        |
| 28        | 10.150      | 1125          | 1134        | 1138         | rBV5     | 129483         | 282026        | 1.95%           | 0.358%        |
| 29        | 10.321      | 1158          | 1163        | 1172         | rVB2     | 125501         | 329590        | 2.28%           | 0.418%        |
| 30        | 10.479      | 1184          | 1190        | 1198         | rBV6     | 265436         | 674279        | 4.67%           | 0.856%        |
| 31        | 10.591      | 1203          | 1209        | 1217         | rVV3     | 120082         | 410352        | 2.84%           | 0.521%        |
| 32        | 10.697      | 1223          | 1227        | 1236         | rVB3     | 123895         | 266291        | 1.85%           | 0.338%        |
| 33        | 11.043      | 1280          | 1286        | 1291         | rVV      | 604349         | 1022771       | 7.09%           | 1.298%        |
| 34        | 11.155      | 1295          | 1305        | 1311         | rVV      | 3988980        | 8339947       | 57.80%          | 10.582%       |
| 35        | 11.231      | 1311          | 1318        | 1326         | rVV3     | 165396         | 553537        | 3.84%           | 0.702%        |
| 36        | 11.989      | 1440          | 1447        | 1454         | rBV3     | 117606         | 240441        | 1.67%           | 0.305%        |

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 ClientSampleId :  
 BGLS8ME

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\_G\Methods\SFAM-EPA-BG010622.M  
 Title : SVOA CALIBRATION

|    |        |      |      |      |      |         |         |        |        |
|----|--------|------|------|------|------|---------|---------|--------|--------|
| 37 | 12.089 | 1458 | 1464 | 1468 | rVV  | 183705  | 290596  | 2.01%  | 0.369% |
| 38 | 12.476 | 1522 | 1530 | 1536 | rBV  | 590533  | 940918  | 6.52%  | 1.194% |
| 39 | 12.735 | 1562 | 1574 | 1582 | rVB  | 1094604 | 1939143 | 13.44% | 2.460% |
| 40 | 12.946 | 1602 | 1610 | 1619 | rVB  | 481323  | 822160  | 5.70%  | 1.043% |
| 41 | 13.416 | 1686 | 1690 | 1698 | rVB  | 165787  | 245892  | 1.70%  | 0.312% |
| 42 | 13.704 | 1731 | 1739 | 1747 | rBV2 | 856674  | 1478712 | 10.25% | 1.876% |
| 43 | 13.922 | 1771 | 1776 | 1784 | rVB  | 109494  | 253588  | 1.76%  | 0.322% |
| 44 | 14.051 | 1792 | 1798 | 1800 | rBV2 | 197448  | 331998  | 2.30%  | 0.421% |
| 45 | 14.209 | 1819 | 1825 | 1830 | rBV  | 293890  | 524386  | 3.63%  | 0.665% |
| 46 | 14.268 | 1830 | 1835 | 1840 | rVB2 | 335992  | 570174  | 3.95%  | 0.723% |
| 47 | 14.356 | 1846 | 1850 | 1853 | rVB2 | 181029  | 240058  | 1.66%  | 0.305% |
| 48 | 14.585 | 1883 | 1889 | 1896 | rBV  | 212536  | 427029  | 2.96%  | 0.542% |
| 49 | 14.767 | 1915 | 1920 | 1925 | rVB  | 704315  | 882500  | 6.12%  | 1.120% |
| 50 | 14.891 | 1937 | 1941 | 1945 | rVB  | 187228  | 286645  | 1.99%  | 0.364% |
| 51 | 14.955 | 1946 | 1952 | 1960 | rBV4 | 115322  | 305526  | 2.12%  | 0.388% |
| 52 | 15.096 | 1967 | 1976 | 1984 | rVB3 | 93876   | 318635  | 2.21%  | 0.404% |
| 53 | 15.279 | 2001 | 2007 | 2012 | rVV4 | 291262  | 552203  | 3.83%  | 0.701% |
| 54 | 15.361 | 2016 | 2021 | 2031 | rVV5 | 112634  | 349793  | 2.42%  | 0.444% |
| 55 | 15.678 | 2067 | 2075 | 2077 | rBV7 | 123807  | 251756  | 1.74%  | 0.319% |
| 56 | 15.713 | 2077 | 2081 | 2086 | rVB  | 580573  | 754052  | 5.23%  | 0.957% |
| 57 | 15.819 | 2094 | 2099 | 2105 | rBV3 | 104177  | 208732  | 1.45%  | 0.265% |
| 58 | 15.878 | 2105 | 2109 | 2113 | rBV  | 262068  | 373763  | 2.59%  | 0.474% |
| 59 | 16.124 | 2144 | 2151 | 2155 | rBV2 | 168647  | 365965  | 2.54%  | 0.464% |
| 60 | 16.577 | 2223 | 2228 | 2230 | rBV  | 558908  | 721151  | 5.00%  | 0.915% |
| 61 | 16.600 | 2230 | 2232 | 2237 | rVB2 | 309168  | 381915  | 2.65%  | 0.485% |
| 62 | 17.017 | 2299 | 2303 | 2311 | rVB2 | 296950  | 503099  | 3.49%  | 0.638% |
| 63 | 17.370 | 2359 | 2363 | 2367 | rVB  | 388439  | 468386  | 3.25%  | 0.594% |
| 64 | 17.517 | 2383 | 2388 | 2392 | rBV2 | 158797  | 268237  | 1.86%  | 0.340% |
| 65 | 17.640 | 2403 | 2409 | 2412 | rBV  | 224067  | 379441  | 2.63%  | 0.481% |
| 66 | 17.734 | 2421 | 2425 | 2430 | rVB2 | 275494  | 413026  | 2.86%  | 0.524% |
| 67 | 17.846 | 2440 | 2444 | 2451 | rVB2 | 168858  | 320024  | 2.22%  | 0.406% |
| 68 | 18.110 | 2485 | 2489 | 2496 | rVB  | 373589  | 496134  | 3.44%  | 0.629% |
| 69 | 18.533 | 2553 | 2561 | 2570 | rBV  | 1723874 | 3366368 | 23.33% | 4.271% |
| 70 | 18.797 | 2602 | 2606 | 2610 | rVB  | 255483  | 299883  | 2.08%  | 0.380% |
| 71 | 19.068 | 2648 | 2652 | 2661 | rVB5 | 186019  | 286094  | 1.98%  | 0.363% |
| 72 | 19.420 | 2706 | 2712 | 2716 | rVB2 | 316676  | 492743  | 3.41%  | 0.625% |
| 73 | 19.608 | 2740 | 2744 | 2752 | rVB2 | 217665  | 309121  | 2.14%  | 0.392% |
| 74 | 19.772 | 2768 | 2772 | 2781 | rVV2 | 272953  | 350539  | 2.43%  | 0.445% |
| 75 | 19.890 | 2788 | 2792 | 2800 | rVB  | 147681  | 222852  | 1.54%  | 0.283% |
| 76 | 20.019 | 2806 | 2814 | 2827 | rVB2 | 491804  | 954216  | 6.61%  | 1.211% |

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 ALS Vial : 44 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 BGLS8ME

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\_G\Methods\SFAM-EPA-BG010622.M  
 Title : SVOA CALIBRATION

|     |        |      |      |      |      |         |          |         |         |
|-----|--------|------|------|------|------|---------|----------|---------|---------|
| 77  | 20.524 | 2896 | 2900 | 2904 | rVB  | 244688  | 296098   | 2.05%   | 0.376%  |
| 78  | 20.918 | 2958 | 2967 | 2972 | rVB  | 574357  | 926745   | 6.42%   | 1.176%  |
| 79  | 21.030 | 2982 | 2986 | 2989 | rVB  | 203162  | 249438   | 1.73%   | 0.316%  |
| 80  | 21.282 | 3025 | 3029 | 3036 | rVB  | 144917  | 208945   | 1.45%   | 0.265%  |
| 81  | 21.564 | 3072 | 3077 | 3083 | rVB3 | 389408  | 579770   | 4.02%   | 0.736%  |
| 82  | 21.840 | 3115 | 3124 | 3134 | rVB  | 8400690 | 14429881 | 100.00% | 18.309% |
| 83  | 21.964 | 3137 | 3145 | 3156 | rVB3 | 269919  | 648833   | 4.50%   | 0.823%  |
| 84  | 22.146 | 3168 | 3176 | 3187 | rBV2 | 303885  | 617706   | 4.28%   | 0.784%  |
| 85  | 22.598 | 3250 | 3253 | 3257 | rVB2 | 168027  | 223461   | 1.55%   | 0.284%  |
| 86  | 22.804 | 3282 | 3288 | 3295 | rBV2 | 270090  | 504357   | 3.50%   | 0.640%  |
| 87  | 23.027 | 3320 | 3326 | 3329 | rBV  | 167425  | 344335   | 2.39%   | 0.437%  |
| 88  | 23.062 | 3330 | 3332 | 3337 | rVV2 | 153455  | 243293   | 1.69%   | 0.309%  |
| 89  | 23.133 | 3338 | 3344 | 3355 | rVB  | 213611  | 549487   | 3.81%   | 0.697%  |
| 90  | 23.321 | 3370 | 3376 | 3383 | rBV4 | 140235  | 305379   | 2.12%   | 0.387%  |
| 91  | 23.556 | 3411 | 3416 | 3423 | rVB  | 202200  | 395339   | 2.74%   | 0.502%  |
| 92  | 24.443 | 3561 | 3567 | 3576 | rBV2 | 171620  | 365247   | 2.53%   | 0.463%  |
| 93  | 24.836 | 3629 | 3634 | 3649 | rVB  | 181931  | 580794   | 4.02%   | 0.737%  |
| 94  | 25.148 | 3680 | 3687 | 3692 | rBV3 | 152182  | 390838   | 2.71%   | 0.496%  |
| 95  | 25.471 | 3732 | 3742 | 3757 | rVB4 | 175484  | 766845   | 5.31%   | 0.973%  |
| 96  | 26.734 | 3951 | 3957 | 3970 | rVB3 | 93982   | 312162   | 2.16%   | 0.396%  |
| 97  | 27.221 | 4032 | 4040 | 4050 | rBV2 | 75031   | 231985   | 1.61%   | 0.294%  |
| 98  | 27.750 | 4118 | 4130 | 4142 | rVB7 | 165403  | 719512   | 4.99%   | 0.913%  |
| 99  | 28.237 | 4201 | 4213 | 4225 | rBV4 | 49862   | 210015   | 1.46%   | 0.266%  |
| 100 | 31.480 | 4757 | 4765 | 4784 | rVB9 | 98957   | 483491   | 3.35%   | 0.613%  |

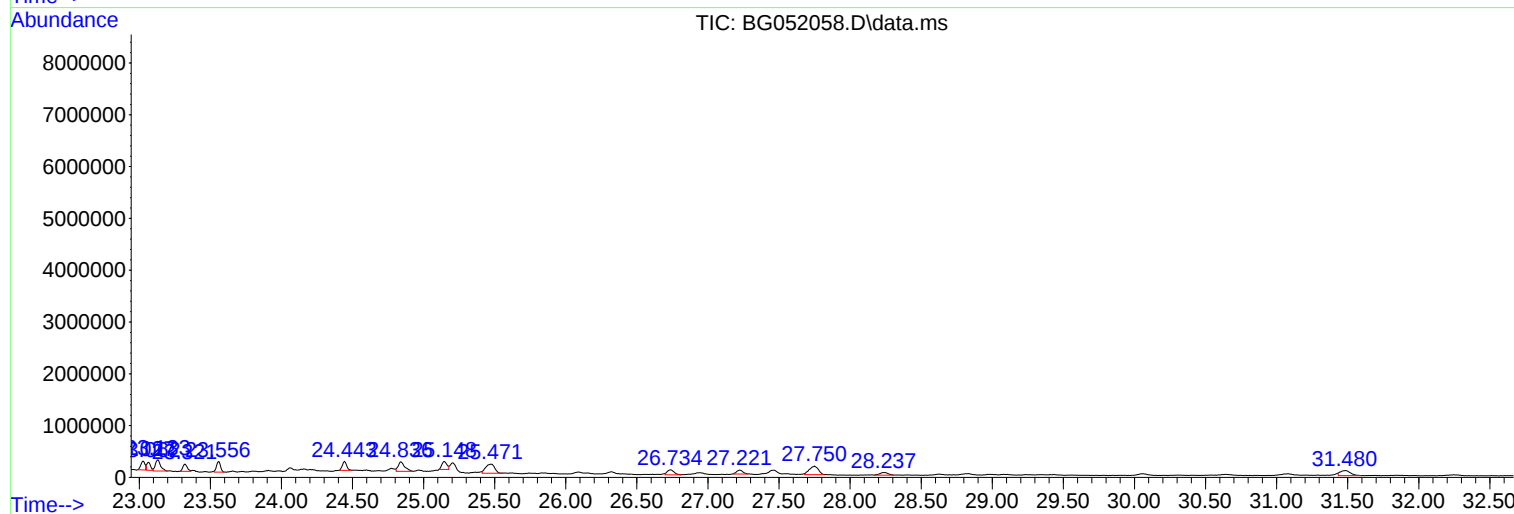
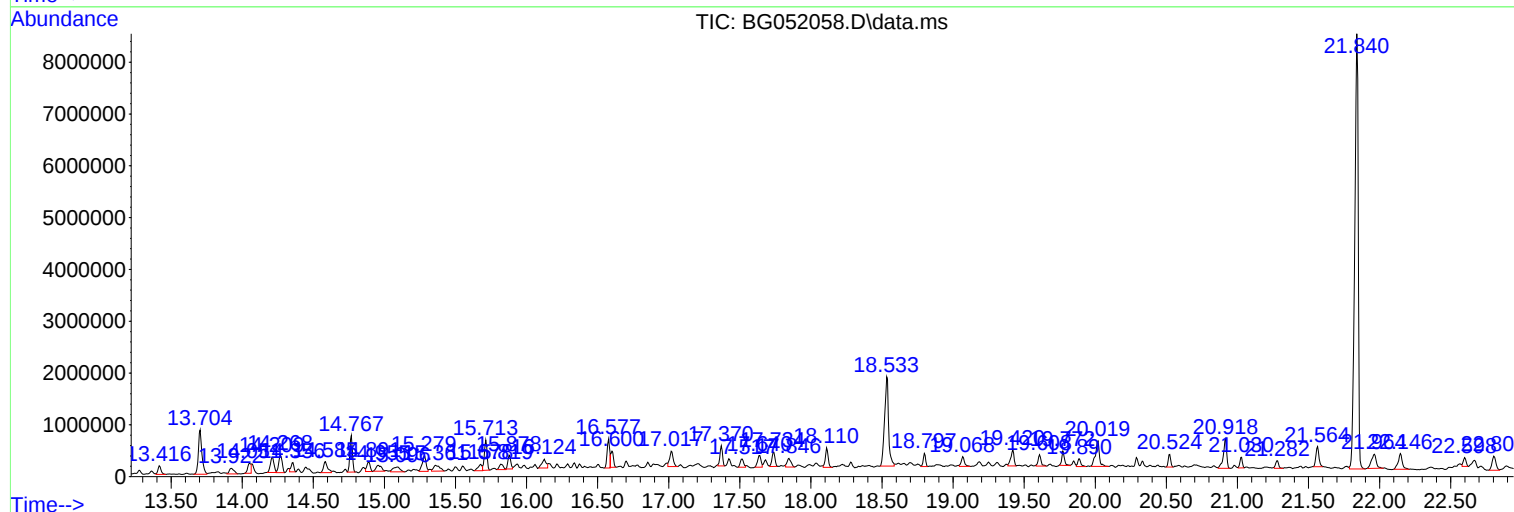
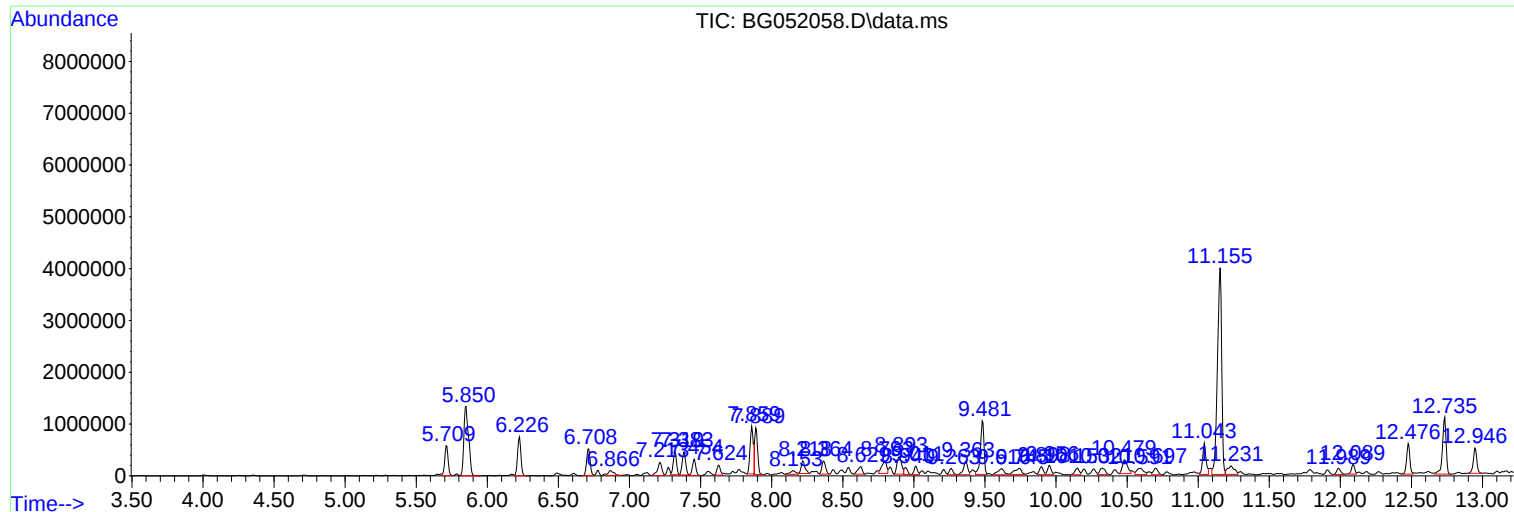
Sum of corrected areas: 78814718

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Instrument :  
BNA\_G  
ClientSampleId :  
BGLS8ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010622.M  
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011722\  
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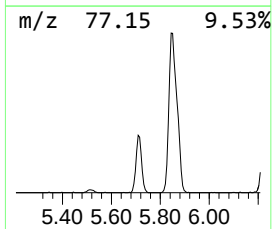
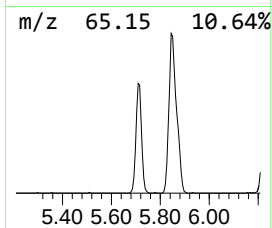
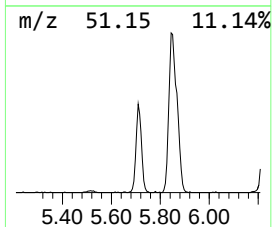
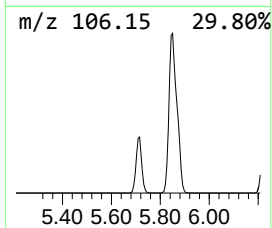
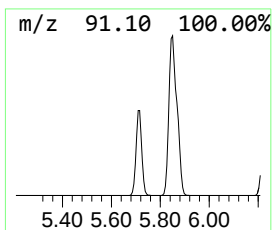
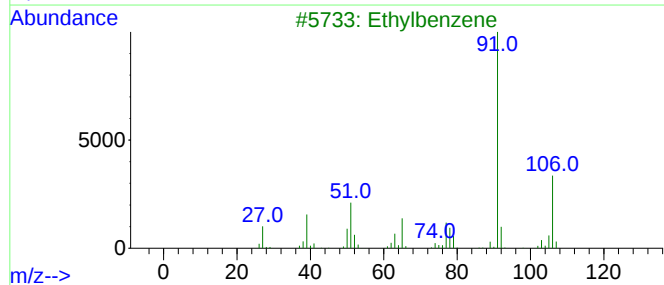
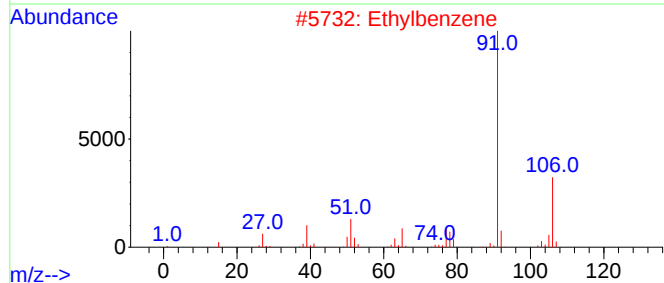
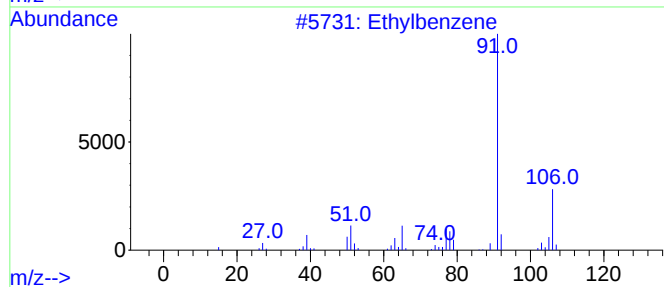
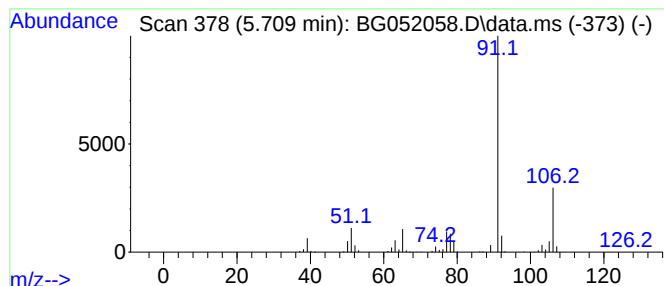
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010622.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 1 Ethylbenzene Concentration Rank 9

| R.T.  | EstConc     | Area   | Relative to ISTD       | R.T.  |
|-------|-------------|--------|------------------------|-------|
| 5.709 | 42.07 ng/ul | 948505 | 1,4-Dichlorobenzene-d4 | 8.247 |

| Hit# | of | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------|-----|---------|-------------|------|
| 1    |    |   | Ethylbenzene | 106 | C8H10   | 000100-41-4 | 95   |
| 2    |    |   | Ethylbenzene | 106 | C8H10   | 000100-41-4 | 94   |
| 3    |    |   | Ethylbenzene | 106 | C8H10   | 000100-41-4 | 91   |
| 4    |    |   | Ethylbenzene | 106 | C8H10   | 000100-41-4 | 91   |
| 5    |    |   | Ethylbenzene | 106 | C8H10   | 000100-41-4 | 91   |



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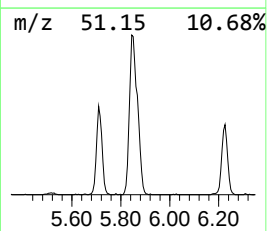
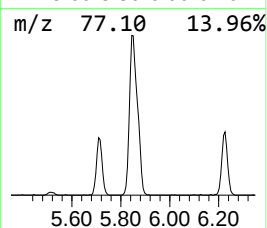
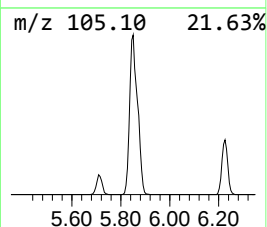
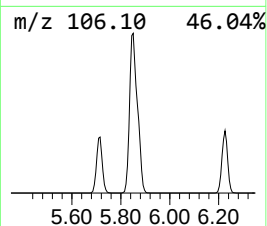
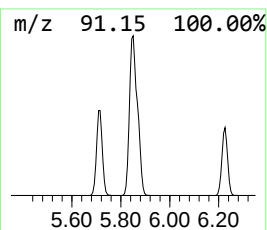
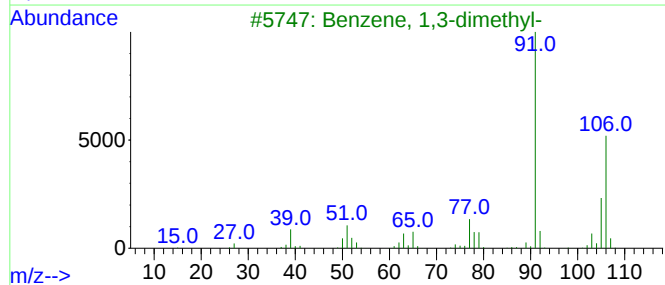
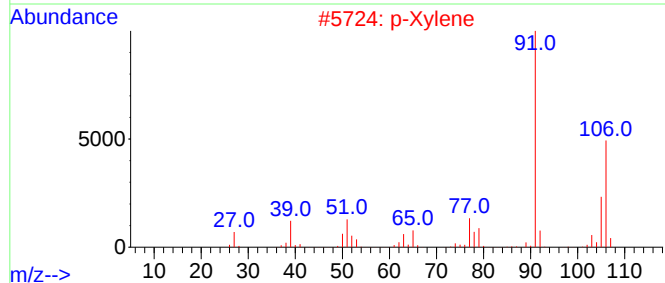
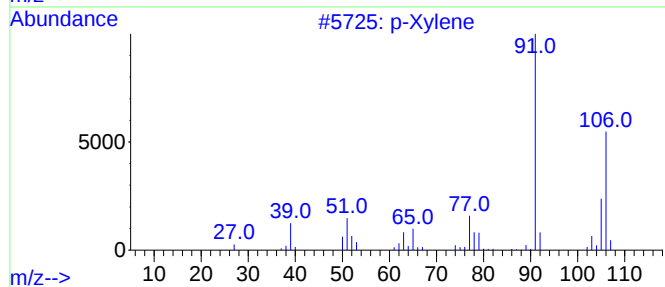
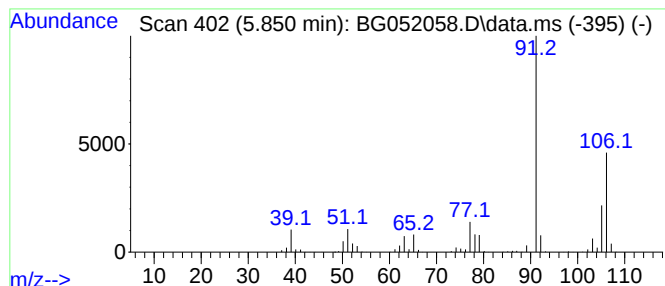
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010622.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 2 p-Xylene Concentration Rank 2

| R.T.  | EstConc      | Area    | Relative to ISTD       | R.T.  |
|-------|--------------|---------|------------------------|-------|
| 5.850 | 135.57 ng/ul | 3056380 | 1,4-Dichlorobenzene-d4 | 8.247 |

| Hit# | of                     | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|------------------------|---|--------------|-----|---------|-------------|------|
| 1    | p-Xylene               |   |              | 106 | C8H10   | 000106-42-3 | 97   |
| 2    | p-Xylene               |   |              | 106 | C8H10   | 000106-42-3 | 97   |
| 3    | Benzene, 1,3-dimethyl- |   |              | 106 | C8H10   | 000108-38-3 | 97   |
| 4    | Benzene, 1,3-dimethyl- |   |              | 106 | C8H10   | 000108-38-3 | 97   |
| 5    | o-Xylene               |   |              | 106 | C8H10   | 000095-47-6 | 97   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011722\  
Data File : BG052058.D  
Acq On : 18 Jan 2022 18:42  
Operator : CG/JU  
Sample : N1100-15ME  
Misc :  
ALS Vial : 44 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLS8ME

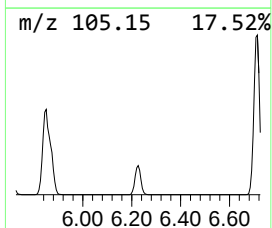
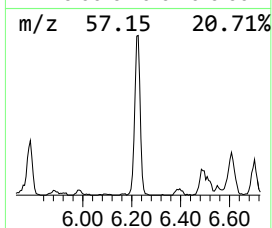
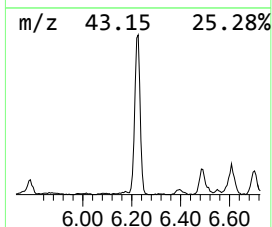
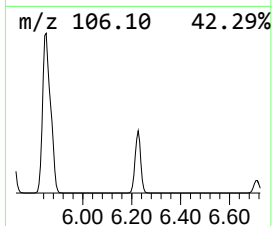
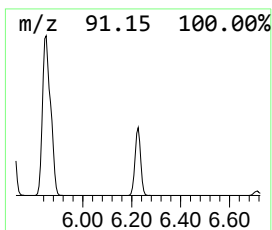
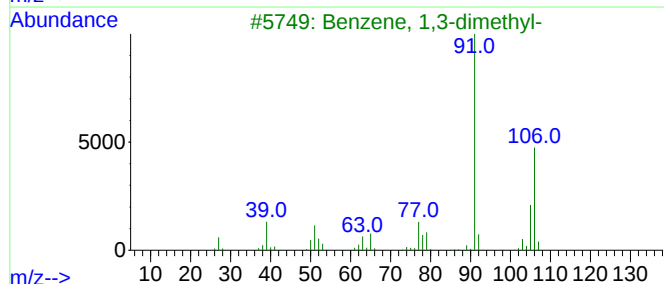
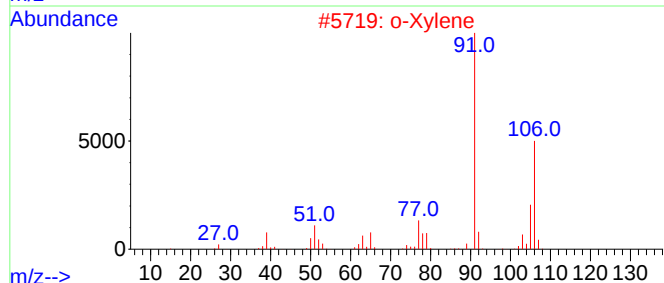
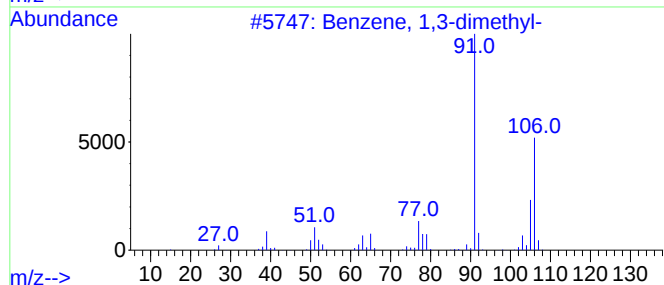
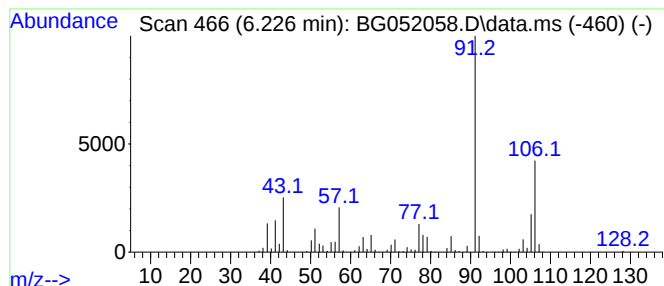
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Quant Title : SVOA CALIBRATION

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

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Peak Number 3 Benzene, 1,3-dimethyl- Concentration Rank 6

| R.T.  | EstConc     | Area    | Relative to ISTD       | R.T.  |
|-------|-------------|---------|------------------------|-------|
| 6.226 | 54.55 ng/ul | 1229880 | 1,4-Dichlorobenzene-d4 | 8.247 |

| Hit# | of | 5 | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1,3-dimethyl- | 106 | C8H10   | 000108-38-3 | 97   |
| 2    |    |   | o-Xylene               | 106 | C8H10   | 000095-47-6 | 97   |
| 3    |    |   | Benzene, 1,3-dimethyl- | 106 | C8H10   | 000108-38-3 | 97   |
| 4    |    |   | p-Xylene               | 106 | C8H10   | 000106-42-3 | 97   |
| 5    |    |   | o-Xylene               | 106 | C8H10   | 000095-47-6 | 97   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011722\  
Data File : BG052058.D  
Acq On : 18 Jan 2022 18:42  
Operator : CG/JU  
Sample : N1100-15ME  
Misc :  
ALS Vial : 44 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLS8ME

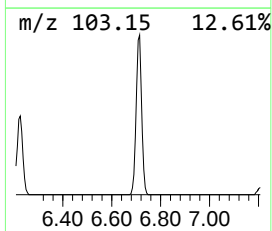
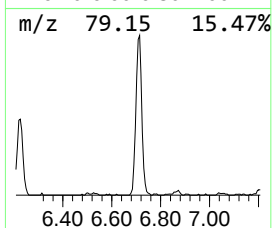
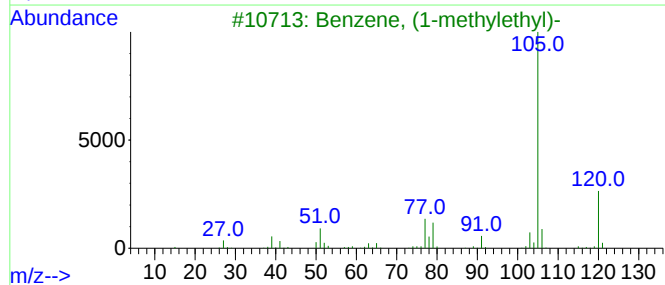
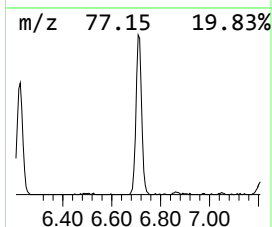
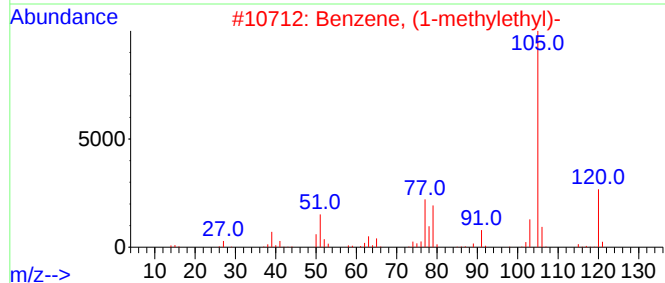
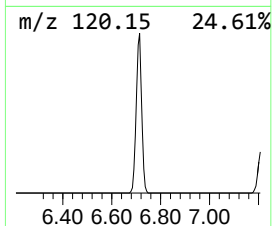
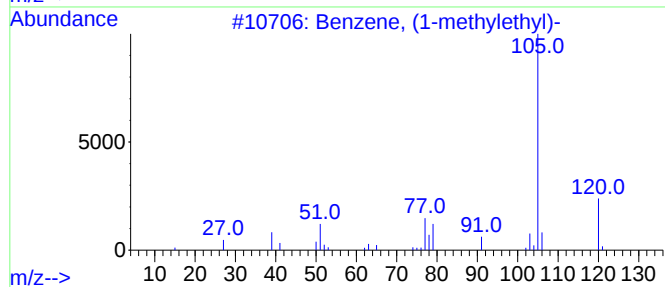
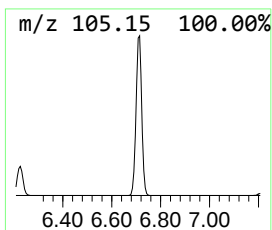
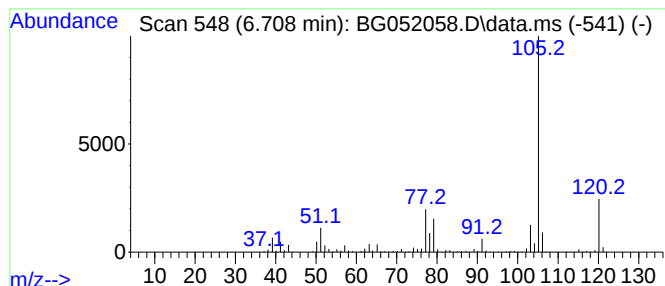
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Quant Title : SVOA CALIBRATION

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

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Peak Number 4 Benzene, (1-methylethyl)- Concentration Rank 10

| R.T.  | EstConc     | Area   | Relative to ISTD       | R.T.  |
|-------|-------------|--------|------------------------|-------|
| 6.708 | 40.05 ng/ul | 902880 | 1,4-Dichlorobenzene-d4 | 8.247 |

| Hit# | of | 5 | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|------|----|---|---------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, (1-methylethyl)- | 120 | C9H12   | 000098-82-8 | 95   |
| 2    |    |   | Benzene, (1-methylethyl)- | 120 | C9H12   | 000098-82-8 | 94   |
| 3    |    |   | Benzene, (1-methylethyl)- | 120 | C9H12   | 000098-82-8 | 91   |
| 4    |    |   | Benzene, 1,2,3-trimethyl- | 120 | C9H12   | 000526-73-8 | 91   |
| 5    |    |   | Benzene, (1-methylethyl)- | 120 | C9H12   | 000098-82-8 | 90   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011722\  
Data File : BG052058.D  
Acq On : 18 Jan 2022 18:42  
Operator : CG/JU  
Sample : N1100-15ME  
Misc :  
ALS Vial : 44 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLS8ME

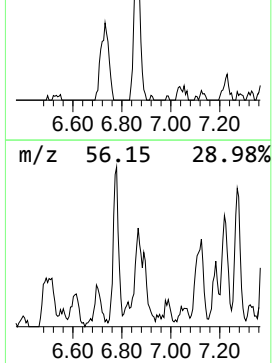
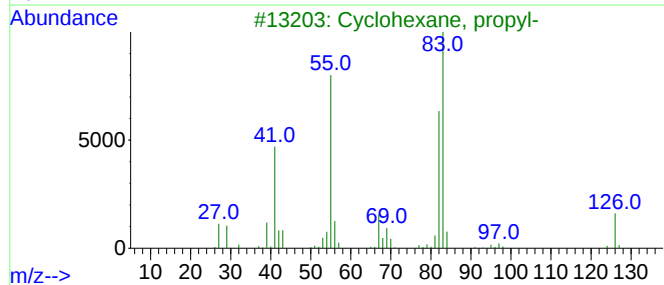
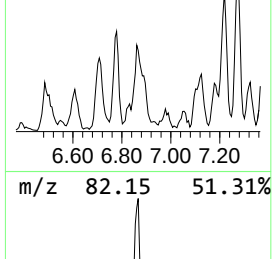
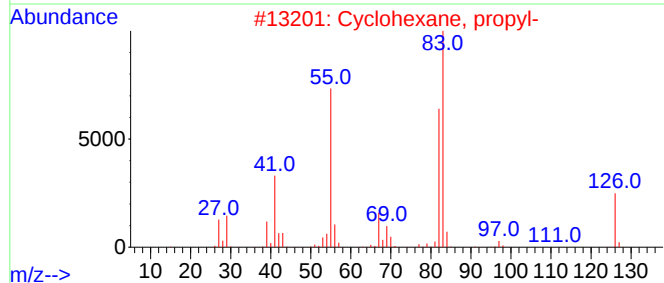
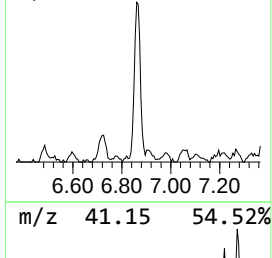
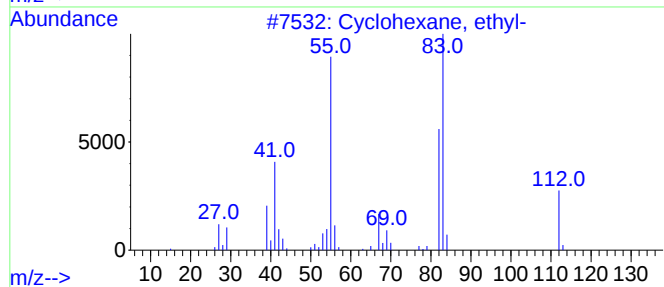
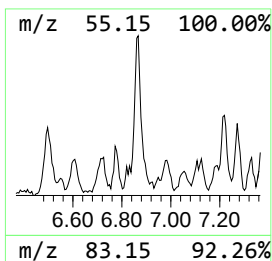
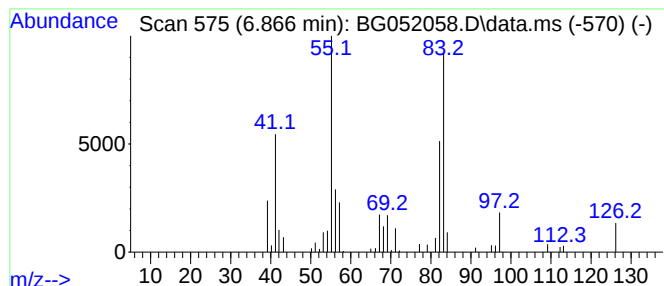
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Quant Title : SVOA CALIBRATION

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

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Peak Number 5 (DEL) Alkane: Cyclic6.866 Concentration Rank 53

| R.T.  | EstConc     | Area   | Relative to ISTD       | R.T.  |
|-------|-------------|--------|------------------------|-------|
| 6.866 | 10.48 ng/ul | 236275 | 1,4-Dichlorobenzene-d4 | 8.247 |

| Hit# | of | 5 | Tentative ID         | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------|-----|---------|-------------|------|
| 1    |    |   | Cyclohexane, ethyl-  | 112 | C8H16   | 001678-91-7 | 81   |
| 2    |    |   | Cyclohexane, propyl- | 126 | C9H18   | 001678-92-8 | 80   |
| 3    |    |   | Cyclohexane, propyl- | 126 | C9H18   | 001678-92-8 | 80   |
| 4    |    |   | Cyclohexane, propyl- | 126 | C9H18   | 001678-92-8 | 72   |
| 5    |    |   | Cyclohexane, propyl- | 126 | C9H18   | 001678-92-8 | 72   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011722\  
Data File : BG052058.D  
Acq On : 18 Jan 2022 18:42  
Operator : CG/JU  
Sample : N1100-15ME  
Misc :  
ALS Vial : 44 Sample Multiplier: 1

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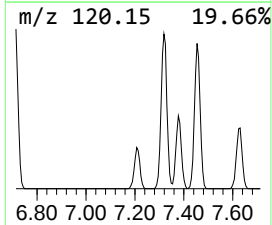
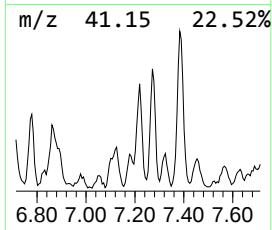
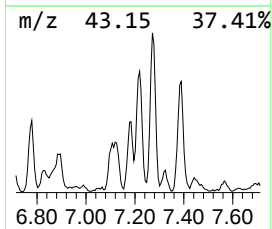
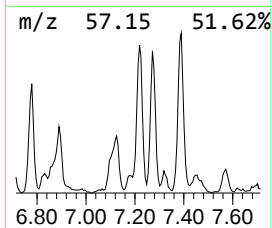
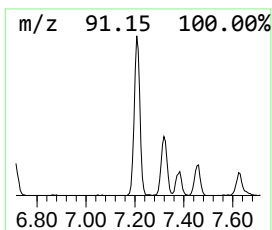
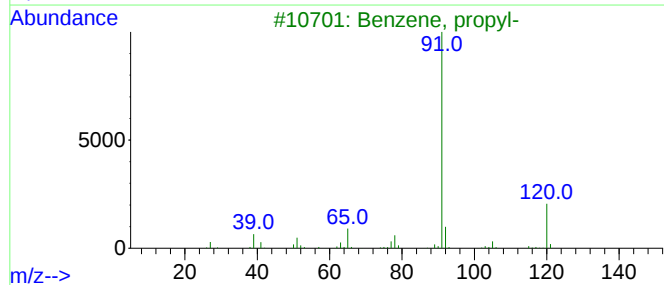
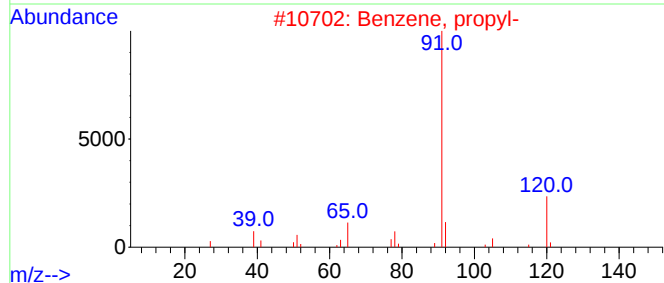
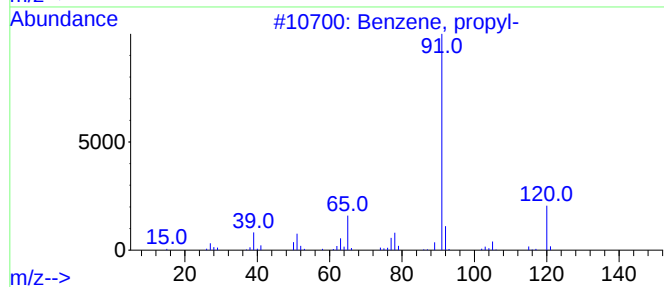
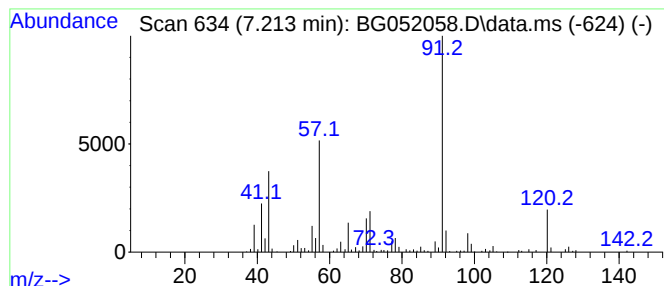
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Quant Title : SVOA CALIBRATION

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

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Peak Number 6 Benzene, propyl- Concentration Rank 23

| R.T.  | EstConc     | Area   | Relative to ISTD       | R.T.  |
|-------|-------------|--------|------------------------|-------|
| 7.213 | 23.88 ng/ul | 538379 | 1,4-Dichlorobenzene-d4 | 8.247 |

| Hit# | of | 5 | Tentative ID        | MW  | MolForm | CAS#        | Qual |
|------|----|---|---------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, propyl-    | 120 | C9H12   | 000103-65-1 | 91   |
| 2    |    |   | Benzene, propyl-    | 120 | C9H12   | 000103-65-1 | 60   |
| 3    |    |   | Benzene, propyl-    | 120 | C9H12   | 000103-65-1 | 60   |
| 4    |    |   | Benzene, propyl-    | 120 | C9H12   | 000103-65-1 | 53   |
| 5    |    |   | Benzeneacetaldehyde | 120 | C8H8O   | 000122-78-1 | 45   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011722\  
Data File : BG052058.D  
Acq On : 18 Jan 2022 18:42  
Operator : CG/JU  
Sample : N1100-15ME  
Misc :  
ALS Vial : 44 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLS8ME

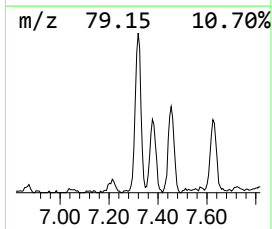
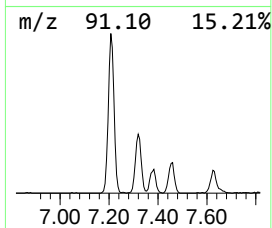
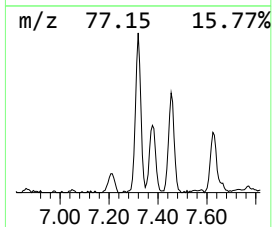
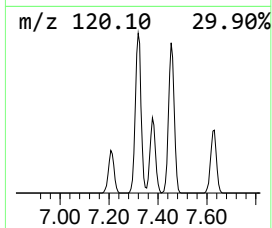
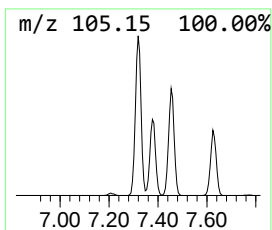
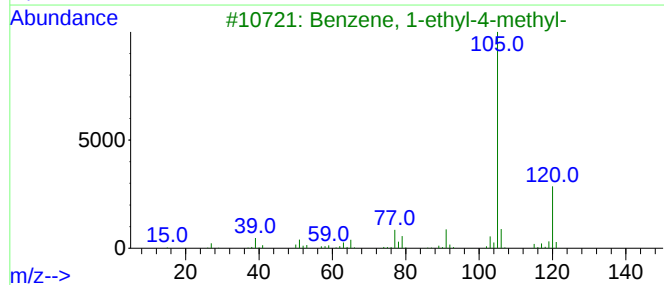
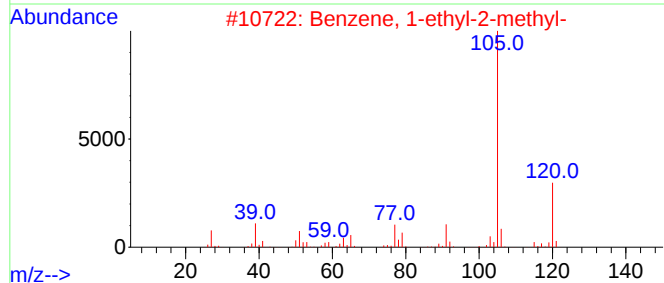
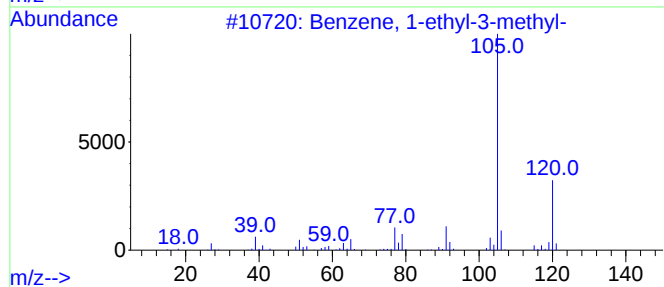
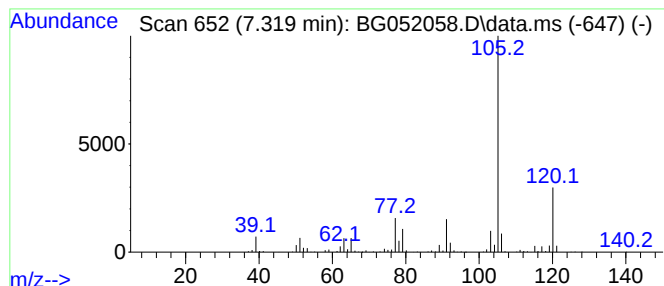
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Quant Title : SVOA CALIBRATION

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

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Peak Number 7 Benzene, 1-ethyl-3-methyl- Concentration Rank 14

| R.T.  | EstConc     | Area   | Relative to ISTD       | R.T.  |
|-------|-------------|--------|------------------------|-------|
| 7.319 | 30.86 ng/ul | 695676 | 1,4-Dichlorobenzene-d4 | 8.247 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 94   |
| 2    |    |   | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 94   |
| 3    |    |   | Benzene, 1-ethyl-4-methyl- | 120 | C9H12   | 000622-96-8 | 91   |
| 4    |    |   | Benzene, 1,2,3-trimethyl-  | 120 | C9H12   | 000526-73-8 | 91   |
| 5    |    |   | Benzene, 1,2,3-trimethyl-  | 120 | C9H12   | 000526-73-8 | 91   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011722\  
Data File : BG052058.D  
Acq On : 18 Jan 2022 18:42  
Operator : CG/JU  
Sample : N1100-15ME  
Misc :  
ALS Vial : 44 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLS8ME

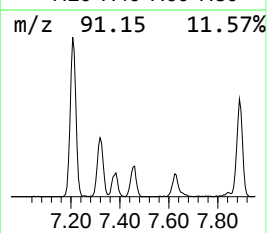
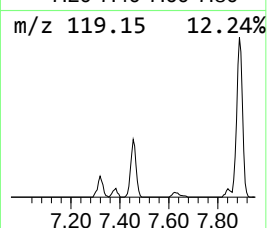
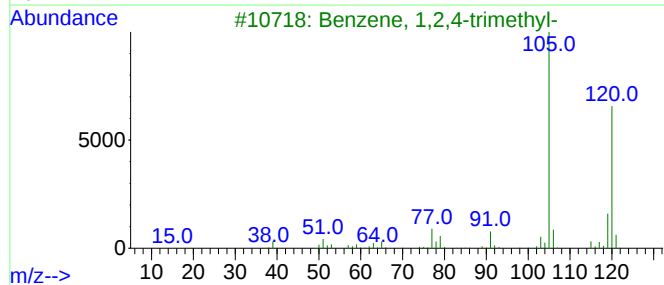
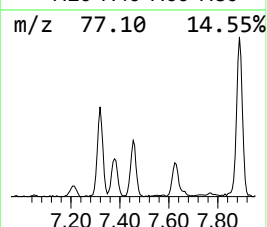
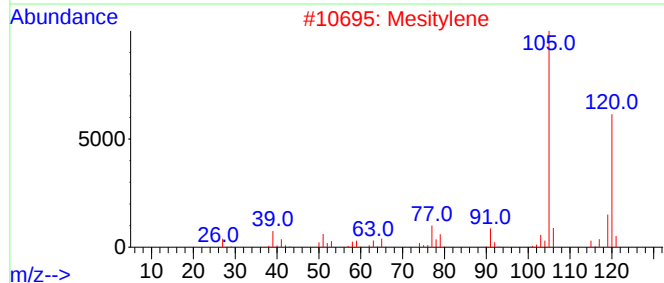
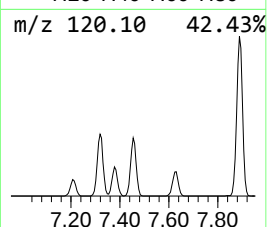
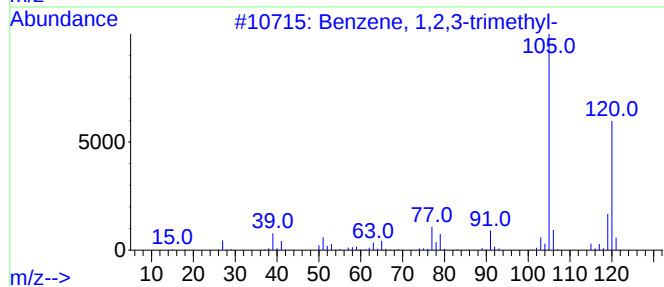
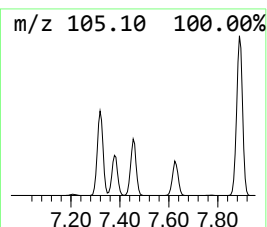
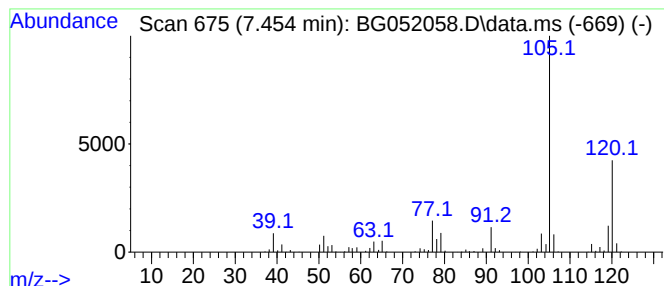
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Quant Title : SVOA CALIBRATION

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

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Peak Number 8 Benzene, 1,2,3-trimethyl- Concentration Rank 25

| R.T.  | EstConc     | Area   | Relative to ISTD       | R.T.  |
|-------|-------------|--------|------------------------|-------|
| 7.454 | 23.05 ng/ul | 519679 | 1,4-Dichlorobenzene-d4 | 8.247 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1,2,3-trimethyl-  | 120 | C9H12   | 000526-73-8 | 95   |
| 2    |    |   | Mesitylene                 | 120 | C9H12   | 000108-67-8 | 94   |
| 3    |    |   | Benzene, 1,2,4-trimethyl-  | 120 | C9H12   | 000095-63-6 | 94   |
| 4    |    |   | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 94   |
| 5    |    |   | Mesitylene                 | 120 | C9H12   | 000108-67-8 | 94   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011722\  
Data File : BG052058.D  
Acq On : 18 Jan 2022 18:42  
Operator : CG/JU  
Sample : N1100-15ME  
Misc :  
ALS Vial : 44 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLS8ME

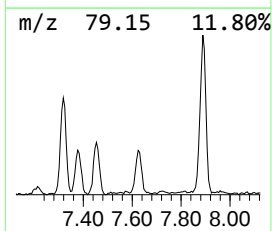
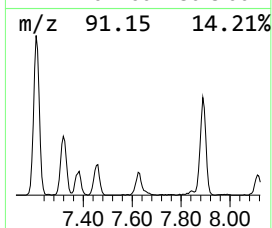
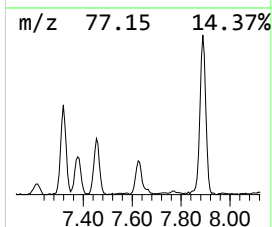
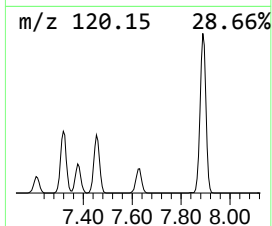
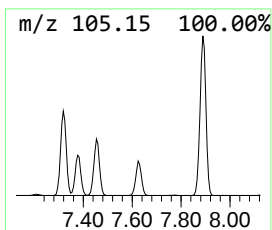
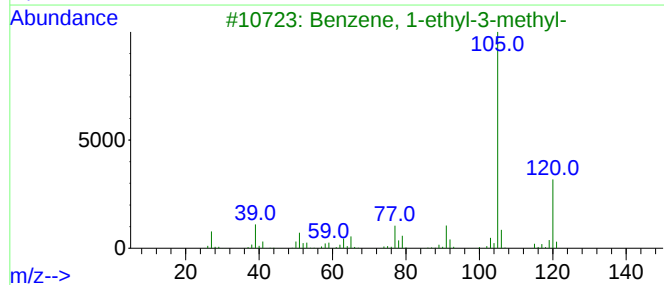
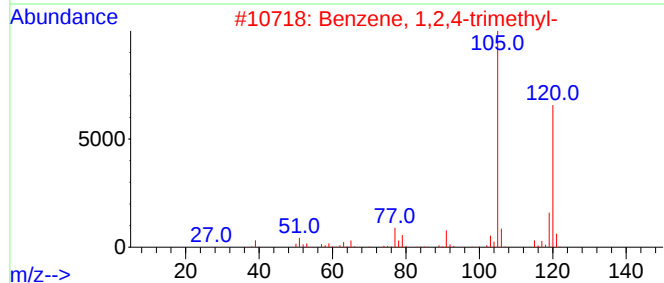
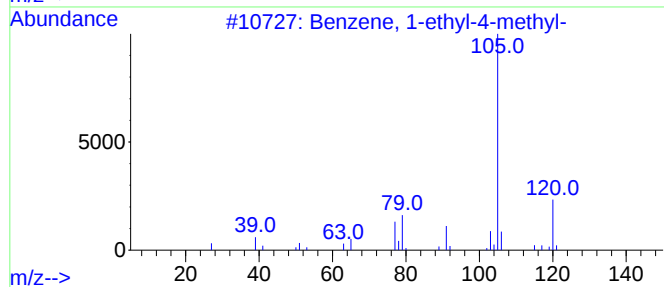
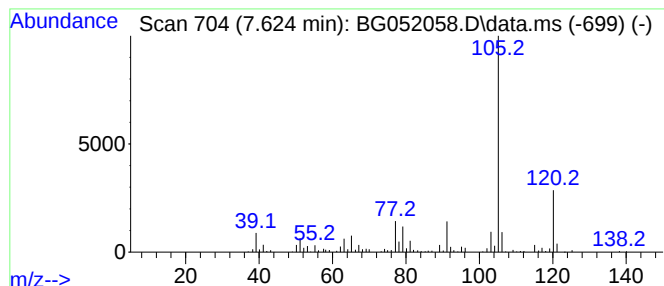
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Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 9 Benzene, 1-ethyl-4-methyl- Concentration Rank 45

| R.T.  | EstConc     | Area   | Relative to ISTD       | R.T.  |
|-------|-------------|--------|------------------------|-------|
| 7.624 | 14.52 ng/ul | 327250 | 1,4-Dichlorobenzene-d4 | 8.247 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1-ethyl-4-methyl- | 120 | C9H12   | 000622-96-8 | 93   |
| 2    |    |   | Benzene, 1,2,4-trimethyl-  | 120 | C9H12   | 000095-63-6 | 91   |
| 3    |    |   | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 91   |
| 4    |    |   | Benzene, 1,2,3-trimethyl-  | 120 | C9H12   | 000526-73-8 | 91   |
| 5    |    |   | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 91   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011722\  
Data File : BG052058.D  
Acq On : 18 Jan 2022 18:42  
Operator : CG/JU  
Sample : N1100-15ME  
Misc :  
ALS Vial : 44 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLS8ME

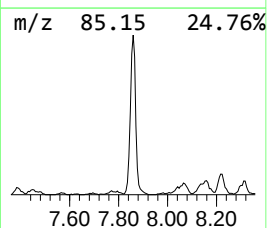
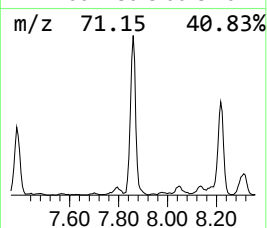
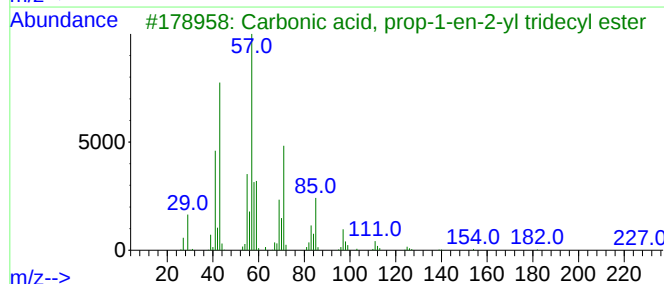
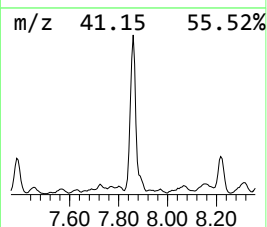
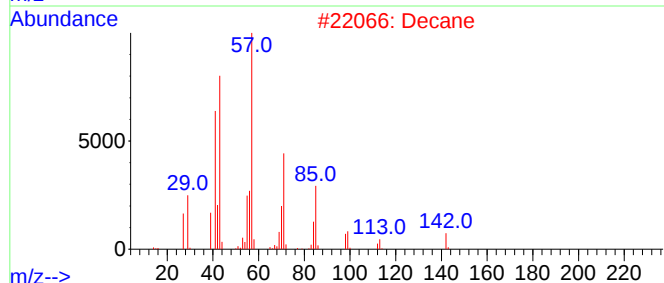
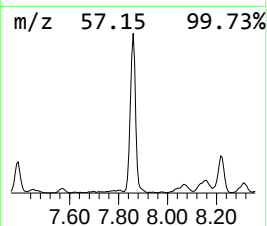
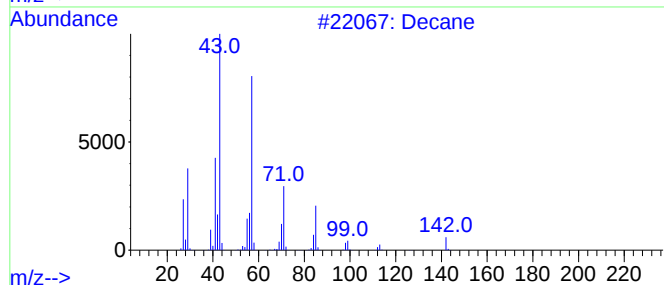
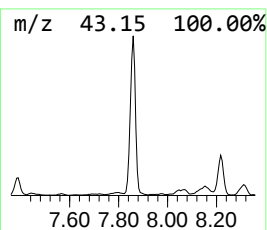
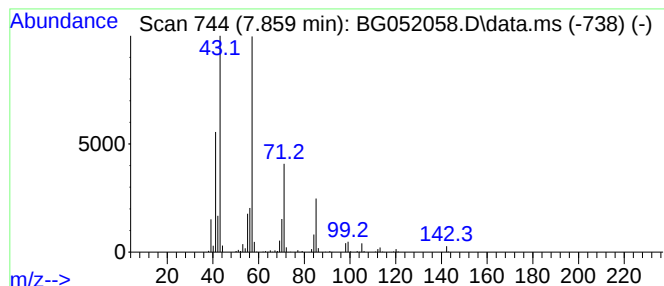
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Quant Title : SVOA CALIBRATION

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

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Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 4

| R.T.  | EstConc     | Area    | Relative to ISTD       | R.T.  |
|-------|-------------|---------|------------------------|-------|
| 7.859 | 69.20 ng/ul | 1560110 | 1,4-Dichlorobenzene-d4 | 8.247 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Decane                              | 142 | C10H22   | 000124-18-5  | 90   |
| 2    |    |   | Decane                              | 142 | C10H22   | 000124-18-5  | 86   |
| 3    |    |   | Carbonic acid, prop-1-en-2-yl tr... | 284 | C17H32O3 | 1000382-90-8 | 83   |
| 4    |    |   | 1-Iodo-2-methylundecane             | 296 | C12H25I  | 073105-67-6  | 83   |
| 5    |    |   | Nonane                              | 128 | C9H20    | 000111-84-2  | 80   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011722\  
Data File : BG052058.D  
Acq On : 18 Jan 2022 18:42  
Operator : CG/JU  
Sample : N1100-15ME  
Misc :  
ALS Vial : 44 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLS8ME

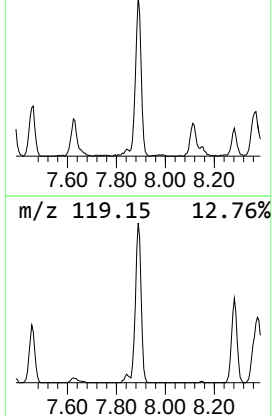
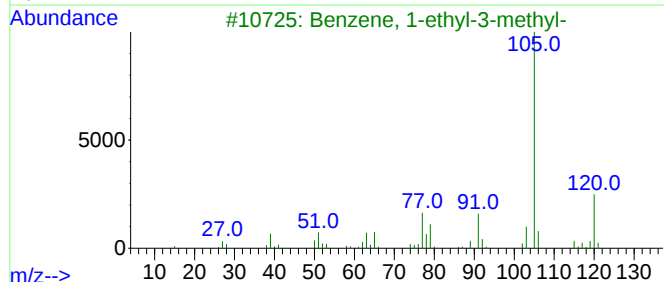
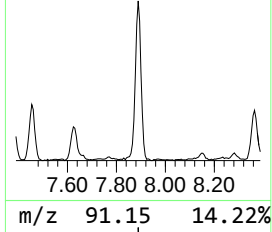
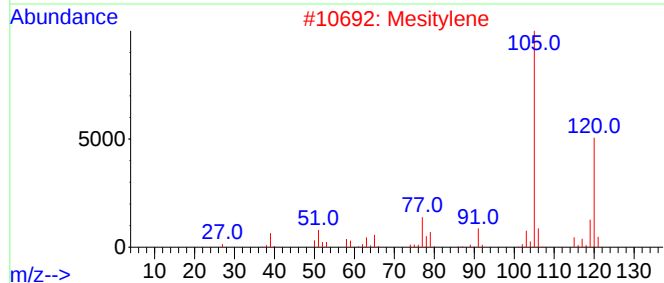
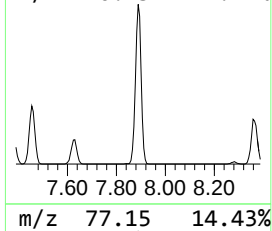
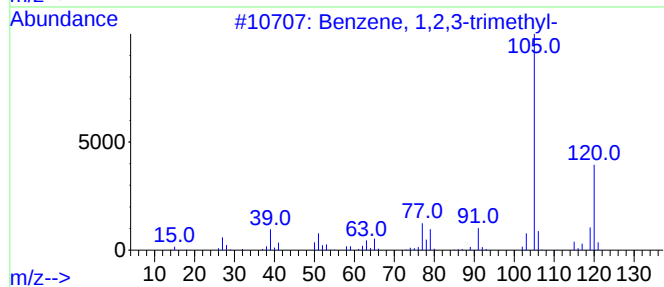
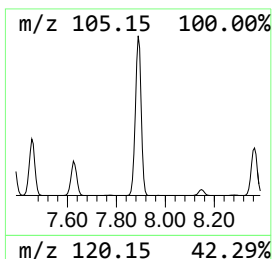
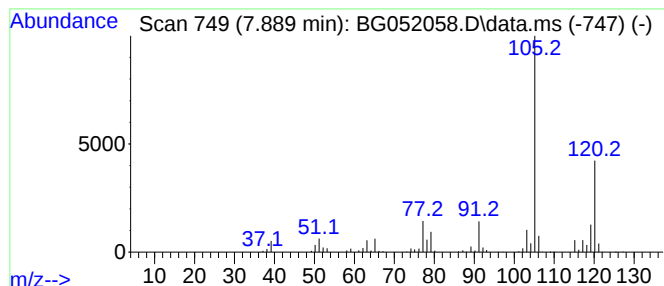
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Quant Title : SVOA CALIBRATION

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

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Peak Number 11 Mesitylene Concentration Rank 7

| R.T.  | EstConc     | Area    | Relative to ISTD       | R.T.  |
|-------|-------------|---------|------------------------|-------|
| 7.889 | 53.20 ng/ul | 1199320 | 1,4-Dichlorobenzene-d4 | 8.247 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1,2,3-trimethyl-  | 120 | C9H12   | 000526-73-8 | 97   |
| 2    |    |   | Mesitylene                 | 120 | C9H12   | 000108-67-8 | 94   |
| 3    |    |   | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 94   |
| 4    |    |   | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 93   |
| 5    |    |   | Benzene, 1,2,3-trimethyl-  | 120 | C9H12   | 000526-73-8 | 91   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011722\  
Data File : BG052058.D  
Acq On : 18 Jan 2022 18:42  
Operator : CG/JU  
Sample : N1100-15ME  
Misc :  
ALS Vial : 44 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLS8ME

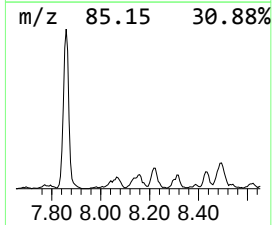
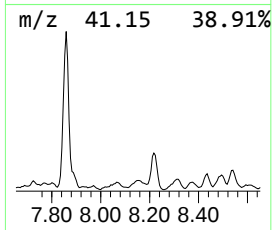
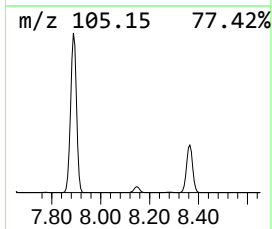
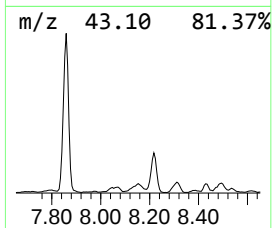
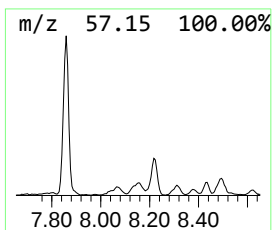
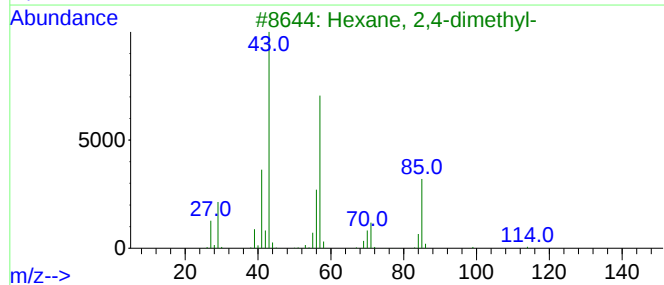
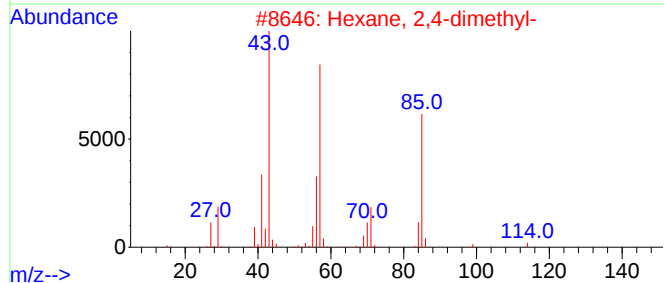
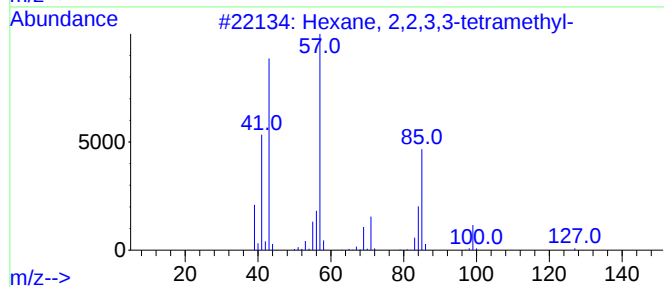
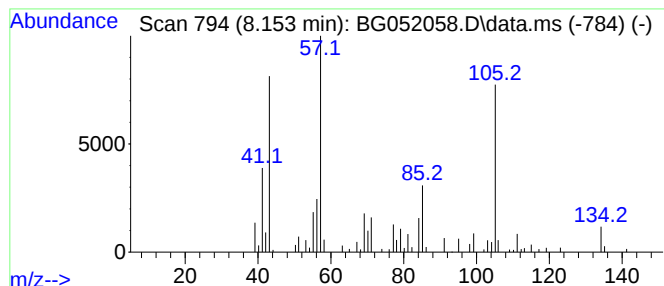
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Quant Title : SVOA CALIBRATION

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

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Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 59

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 8.153 | 9.01 ng/ul | 203052 | 1,4-Dichlorobenzene-d4 | 8.247 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Hexane, 2,2,3,3-tetramethyl-        | 142 | C10H22   | 013475-81-5  | 38   |
| 2    |    |   | Hexane, 2,4-dimethyl-               | 114 | C8H18    | 000589-43-5  | 35   |
| 3    |    |   | Hexane, 2,4-dimethyl-               | 114 | C8H18    | 000589-43-5  | 35   |
| 4    |    |   | Oxalic acid, hexyl isohexyl ester   | 258 | C14H26O4 | 1010309-33-0 | 35   |
| 5    |    |   | Benzeneacetaldehyde, .alpha.-met... | 134 | C9H10O   | 000093-53-8  | 30   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011722\  
Data File : BG052058.D  
Acq On : 18 Jan 2022 18:42  
Operator : CG/JU  
Sample : N1100-15ME  
Misc :  
ALS Vial : 44 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLS8ME

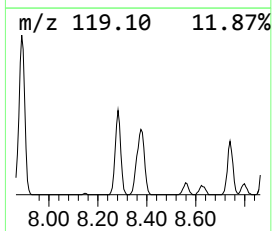
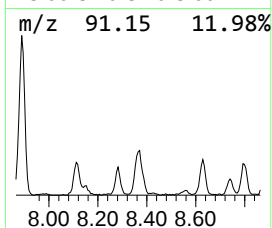
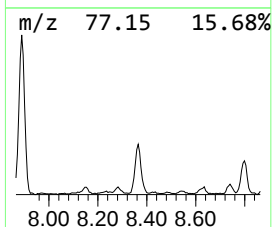
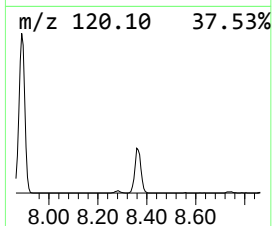
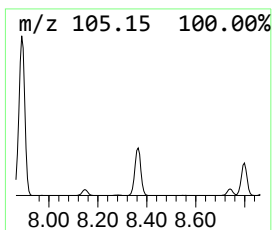
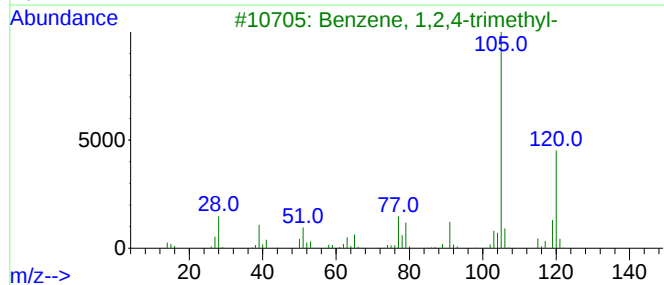
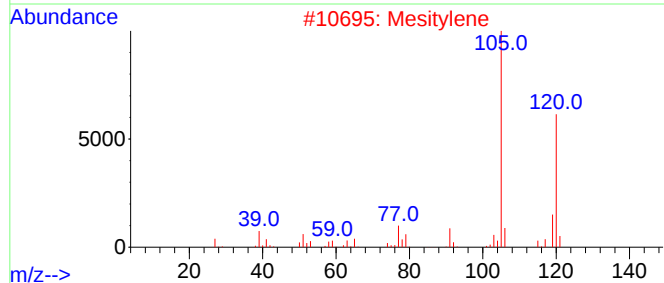
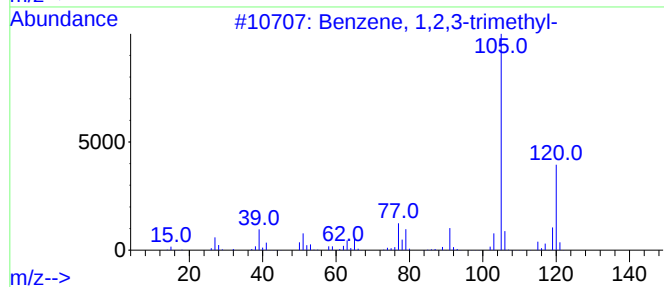
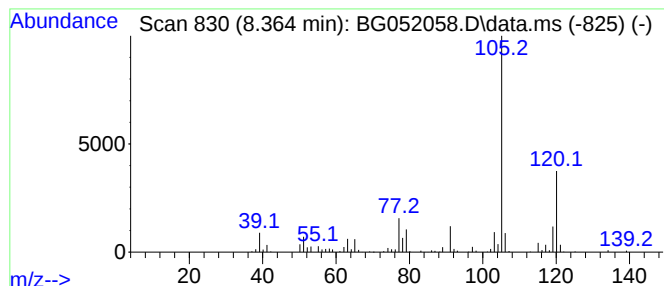
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Quant Title : SVOA CALIBRATION

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

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Peak Number 13 Benzene, 1,2,4-trimethyl- Concentration Rank 26

| R.T.  | EstConc     | Area   | Relative to ISTD       | R.T.  |
|-------|-------------|--------|------------------------|-------|
| 8.364 | 20.60 ng/ul | 464455 | 1,4-Dichlorobenzene-d4 | 8.247 |

| Hit# | of | 5 | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|------|----|---|---------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1,2,3-trimethyl- | 120 | C9H12   | 000526-73-8 | 97   |
| 2    |    |   | Mesitylene                | 120 | C9H12   | 000108-67-8 | 94   |
| 3    |    |   | Benzene, 1,2,4-trimethyl- | 120 | C9H12   | 000095-63-6 | 94   |
| 4    |    |   | Benzene, 1,2,3-trimethyl- | 120 | C9H12   | 000526-73-8 | 94   |
| 5    |    |   | Mesitylene                | 120 | C9H12   | 000108-67-8 | 94   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011722\  
Data File : BG052058.D  
Acq On : 18 Jan 2022 18:42  
Operator : CG/JU  
Sample : N1100-15ME  
Misc :  
ALS Vial : 44 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLS8ME

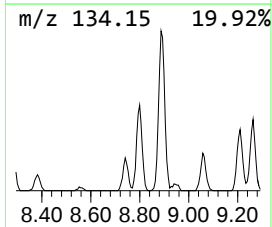
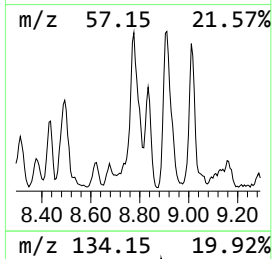
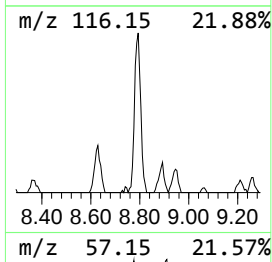
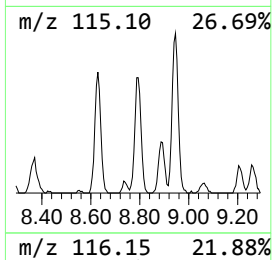
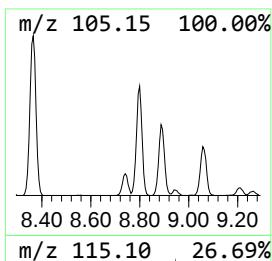
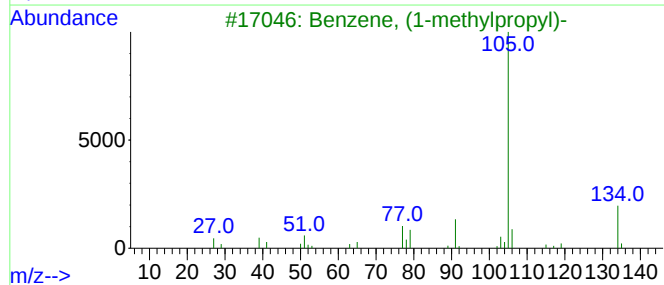
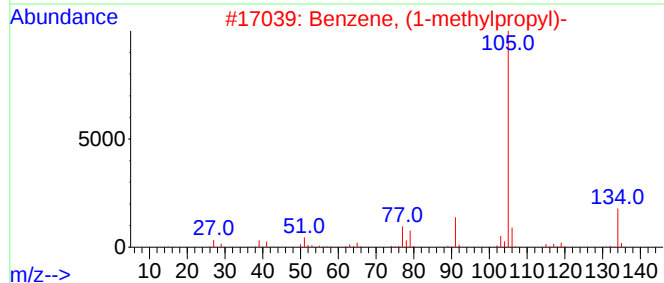
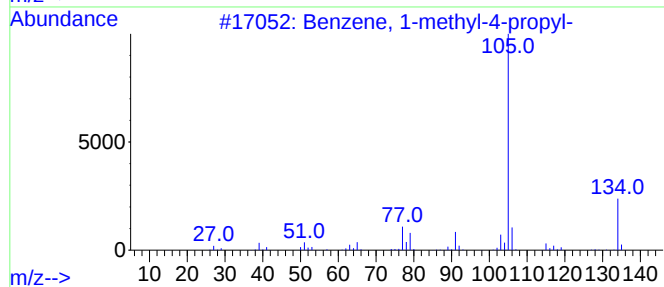
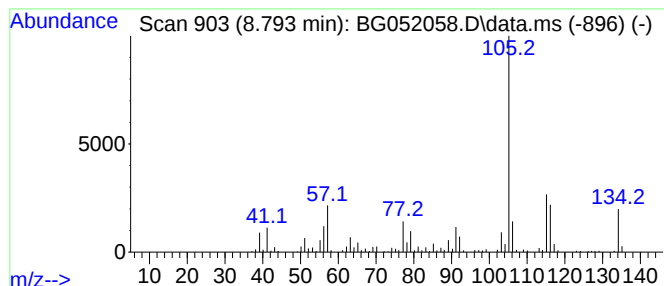
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Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 15 Benzene, 1-methyl-4-propyl- Concentration Rank 19

| R.T.  | EstConc     | Area   | Relative to ISTD       | R.T.  |
|-------|-------------|--------|------------------------|-------|
| 8.793 | 25.35 ng/ul | 571388 | 1,4-Dichlorobenzene-d4 | 8.247 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1-methyl-4-propyl-         | 134 | C10H14  | 001074-55-1 | 64   |
| 2    |    |   | Benzene, (1-methylpropyl)-          | 134 | C10H14  | 000135-98-8 | 60   |
| 3    |    |   | Benzene, (1-methylpropyl)-          | 134 | C10H14  | 000135-98-8 | 60   |
| 4    |    |   | Benzene, 1-methyl-3-propyl-         | 134 | C10H14  | 001074-43-7 | 58   |
| 5    |    |   | Benzeneacetaldehyde, .alpha.-met... | 134 | C9H10O  | 000093-53-8 | 55   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011722\  
Data File : BG052058.D  
Acq On : 18 Jan 2022 18:42  
Operator : CG/JU  
Sample : N1100-15ME  
Misc :  
ALS Vial : 44 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLS8ME

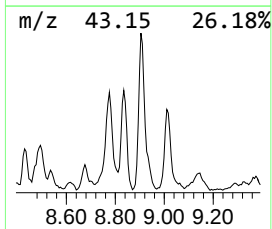
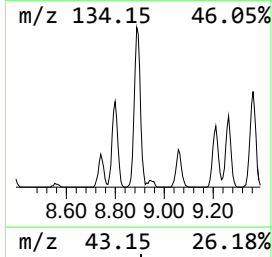
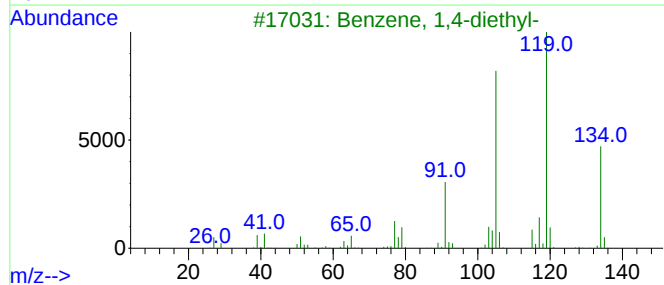
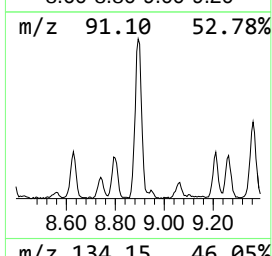
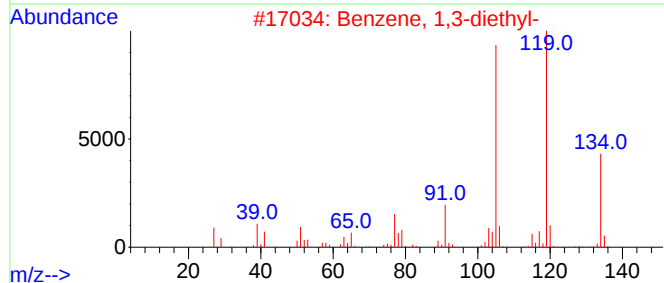
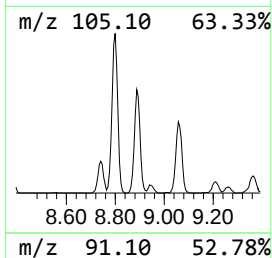
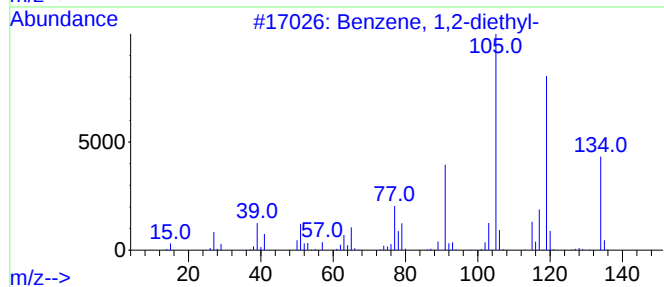
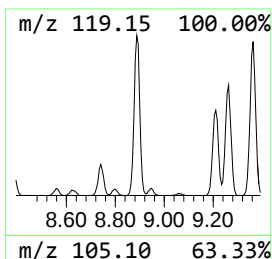
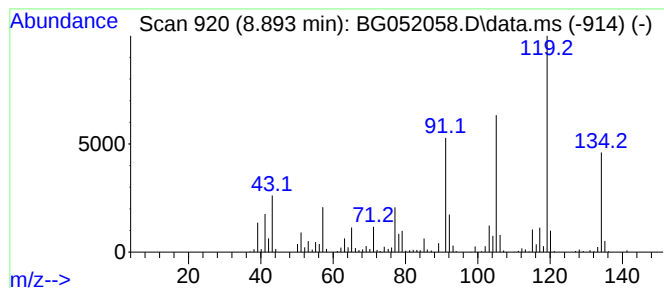
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Quant Title : SVOA CALIBRATION

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

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Peak Number 16 Benzene, 1,2-diethyl- Concentration Rank 13

| R.T.  | EstConc     | Area   | Relative to ISTD       | R.T.  |
|-------|-------------|--------|------------------------|-------|
| 8.893 | 34.30 ng/ul | 773188 | 1,4-Dichlorobenzene-d4 | 8.247 |

| Hit# | of | 5 | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1,2-diethyl-          | 134 | C10H14  | 000135-01-3 | 94   |
| 2    |    |   | Benzene, 1,3-diethyl-          | 134 | C10H14  | 000141-93-5 | 94   |
| 3    |    |   | Benzene, 1,4-diethyl-          | 134 | C10H14  | 000105-05-5 | 93   |
| 4    |    |   | Benzene, 1-ethyl-2,3-dimethyl- | 134 | C10H14  | 000933-98-2 | 91   |
| 5    |    |   | Benzene, 1-ethyl-3,5-dimethyl- | 134 | C10H14  | 000934-74-7 | 91   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011722\  
Data File : BG052058.D  
Acq On : 18 Jan 2022 18:42  
Operator : CG/JU  
Sample : N1100-15ME  
Misc :  
ALS Vial : 44 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLS8ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010622.M  
Quant Title : SVOA CALIBRATION

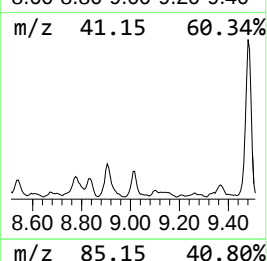
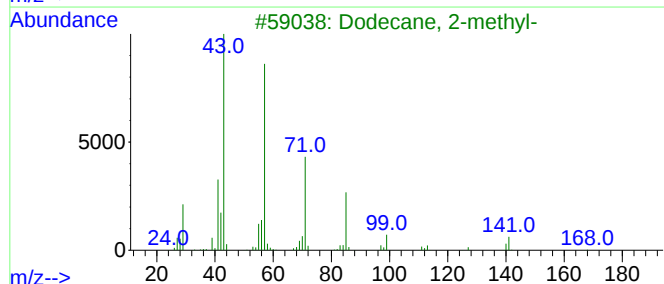
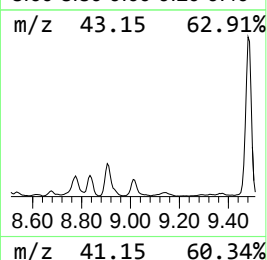
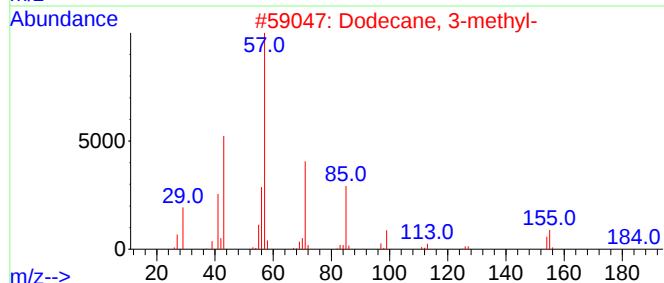
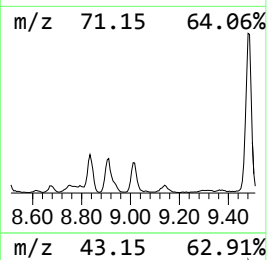
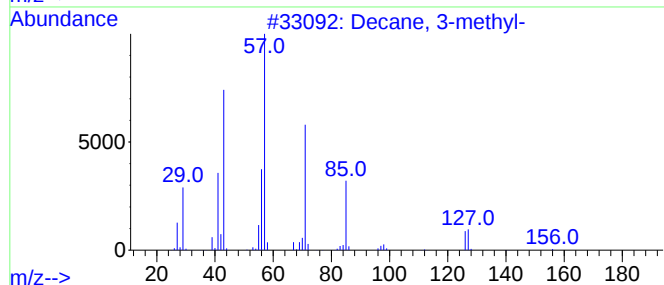
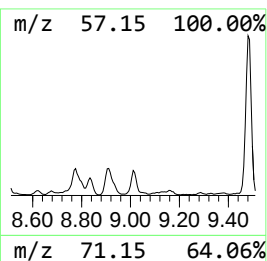
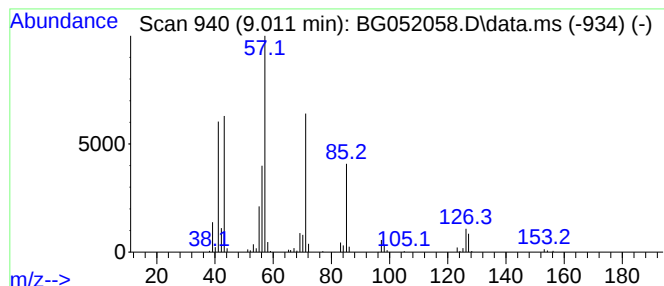
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 17 (DEL) Alkane: Straight-Chai... Concentration Rank 51

| R.T.  | EstConc     | Area   | Relative to ISTD       | R.T.  |
|-------|-------------|--------|------------------------|-------|
| 9.011 | 11.99 ng/ul | 270381 | 1,4-Dichlorobenzene-d4 | 8.247 |

| Hit# | of | 5 | Tentative ID                | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-----------------------------|-----|----------|--------------|------|
| 1    |    |   | Decane, 3-methyl-           | 156 | C11H24   | 013151-34-3  | 93   |
| 2    |    |   | Dodecane, 3-methyl-         | 184 | C13H28   | 017312-57-1  | 72   |
| 3    |    |   | Dodecane, 2-methyl-         | 184 | C13H28   | 001560-97-0  | 53   |
| 4    |    |   | 2-Nonanol, trimethylacetate | 228 | C14H28O2 | 1000463-21-6 | 50   |
| 5    |    |   | Nonane, 1-iodo-             | 254 | C9H19I   | 004282-42-2  | 50   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011722\  
Data File : BG052058.D  
Acq On : 18 Jan 2022 18:42  
Operator : CG/JU  
Sample : N1100-15ME  
Misc :  
ALS Vial : 44 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLS8ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010622.M  
Quant Title : SVOA CALIBRATION

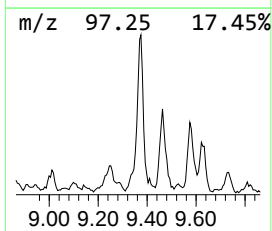
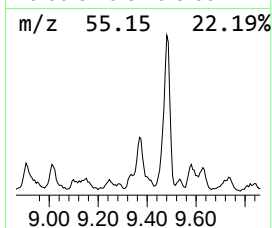
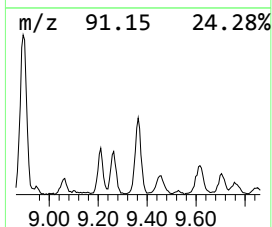
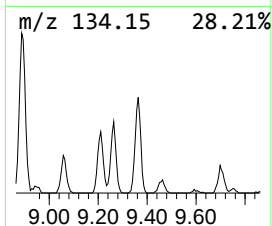
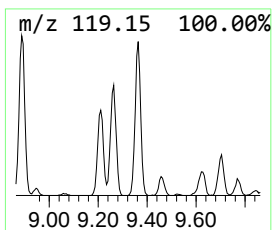
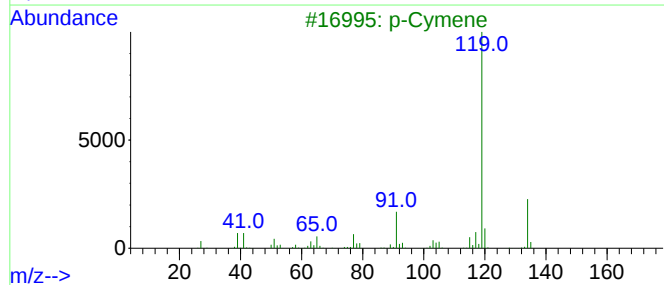
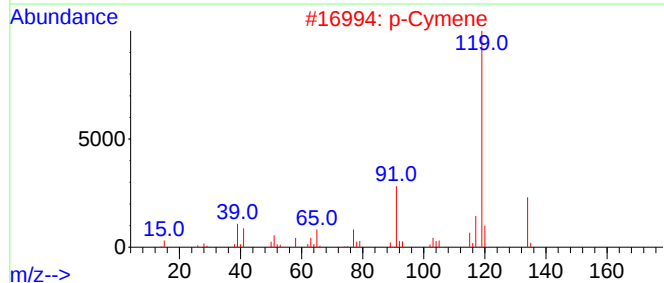
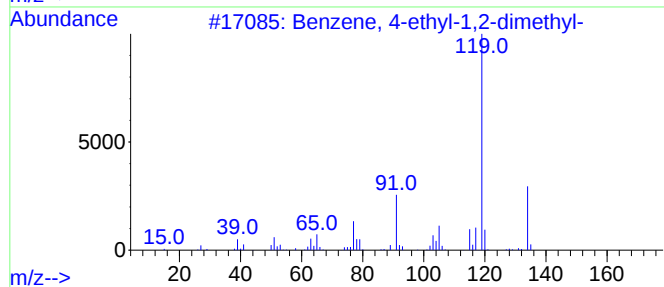
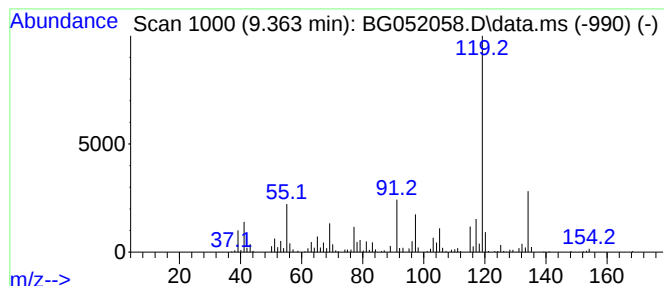
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 19 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 22

| R.T.  | EstConc     | Area   | Relative to ISTD       | R.T.  |
|-------|-------------|--------|------------------------|-------|
| 9.363 | 24.00 ng/ul | 541019 | 1,4-Dichlorobenzene-d4 | 8.247 |

| Hit# | of | 5 | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 4-ethyl-1,2-dimethyl- | 134 | C10H14  | 000934-80-5 | 96   |
| 2    |    |   | p-Cymene                       | 134 | C10H14  | 000099-87-6 | 94   |
| 3    |    |   | p-Cymene                       | 134 | C10H14  | 000099-87-6 | 93   |
| 4    |    |   | Benzene, 1-ethyl-2,4-dimethyl- | 134 | C10H14  | 000874-41-9 | 91   |
| 5    |    |   | Benzene, 2-ethyl-1,3-dimethyl- | 134 | C10H14  | 002870-04-4 | 91   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011722\  
Data File : BG052058.D  
Acq On : 18 Jan 2022 18:42  
Operator : CG/JU  
Sample : N1100-15ME  
Misc :  
ALS Vial : 44 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLS8ME

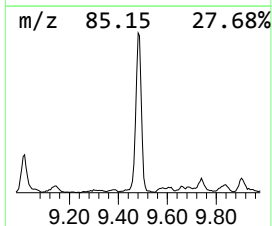
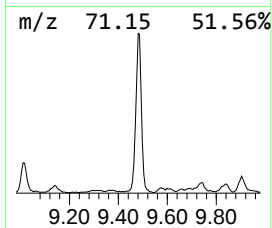
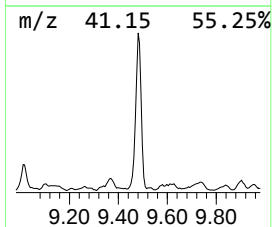
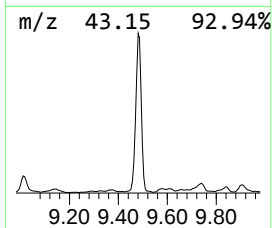
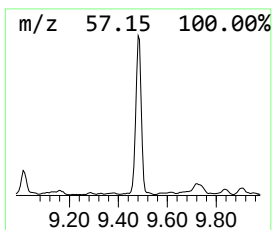
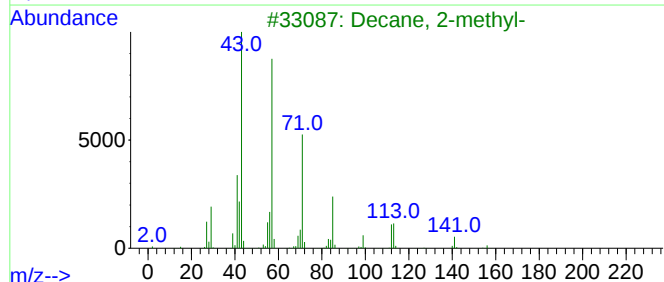
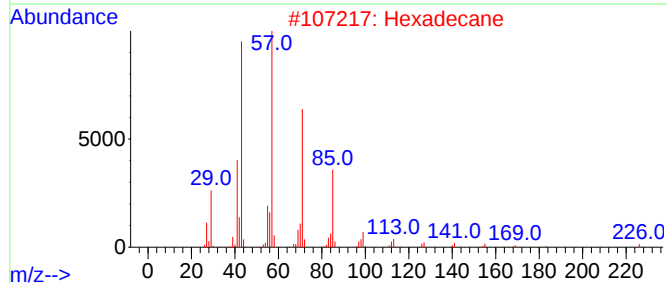
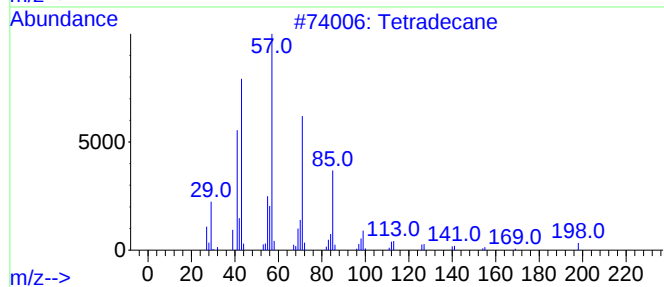
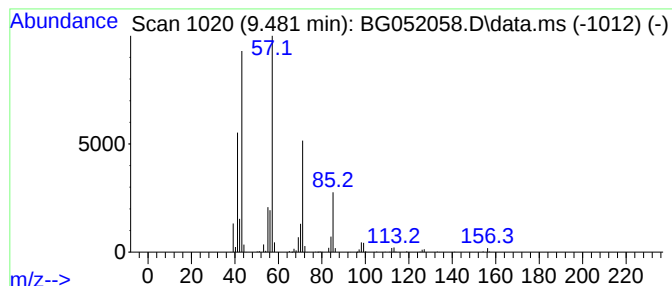
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Quant Title : SVOA CALIBRATION

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

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Peak Number 20 (DEL) Alkane: Straight-Chai... Concentration Rank 3

| R.T.  | EstConc     | Area    | Relative to ISTD       | R.T.  |
|-------|-------------|---------|------------------------|-------|
| 9.481 | 78.97 ng/ul | 1780370 | 1,4-Dichlorobenzene-d4 | 8.247 |

| Hit# | of 5 | Tentative ID      | MW  | MolForm | CAS#        | Qual |
|------|------|-------------------|-----|---------|-------------|------|
| 1    |      | Tetradecane       | 198 | C14H30  | 000629-59-4 | 90   |
| 2    |      | Hexadecane        | 226 | C16H34  | 000544-76-3 | 80   |
| 3    |      | Decane, 2-methyl- | 156 | C11H24  | 006975-98-0 | 80   |
| 4    |      | Undecane          | 156 | C11H24  | 001120-21-4 | 72   |
| 5    |      | Heptadecane       | 240 | C17H36  | 000629-78-7 | 64   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011722\  
Data File : BG052058.D  
Acq On : 18 Jan 2022 18:42  
Operator : CG/JU  
Sample : N1100-15ME  
Misc :  
ALS Vial : 44 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLS8ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010622.M  
Quant Title : SVOA CALIBRATION

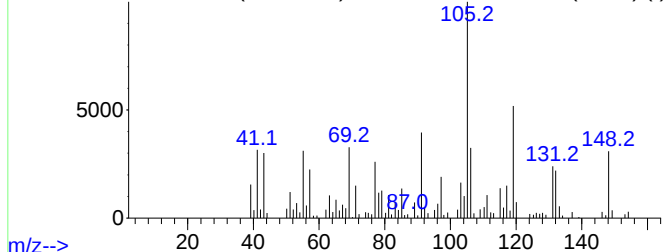
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TIC Integration Parameters: LSCINT.P

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Peak Number 21 unknown-01 Concentration Rank 38

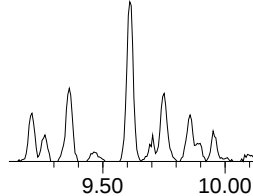
| R.T.  | EstConc     | Area   | Relative to ISTD       | R.T.  |
|-------|-------------|--------|------------------------|-------|
| 9.616 | 16.41 ng/ul | 369957 | 1,4-Dichlorobenzene-d4 | 8.247 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | 3,4-Dihydro-2-[1H]quinoxalinone     | 148 | C8H8N2O | 059564-59-9 | 35   |
| 2    |    |   | Benzenepropanal, .beta.-methyl-     | 148 | C10H12O | 016251-77-7 | 30   |
| 3    |    |   | Benzene, 1-methyl-4-(2-methylpro... | 148 | C11H16  | 005161-04-6 | 30   |
| 4    |    |   | Benzene, 2-ethyl-1,4-dimethyl-      | 134 | C10H14  | 001758-88-9 | 25   |
| 5    |    |   | Benzene, (1,2-dimethylpropyl)-      | 148 | C11H16  | 004481-30-5 | 18   |

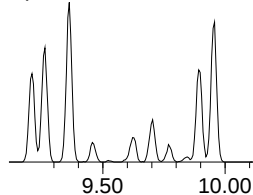
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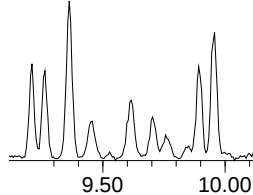
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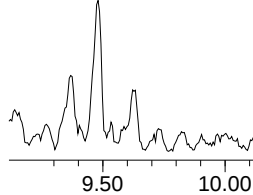
m/z 119.15 51.85%



m/z 91.15 39.55%

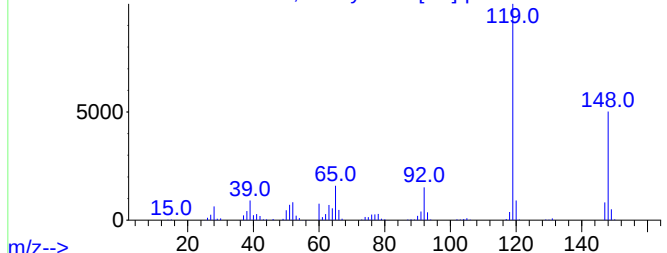


m/z 69.15 32.79%

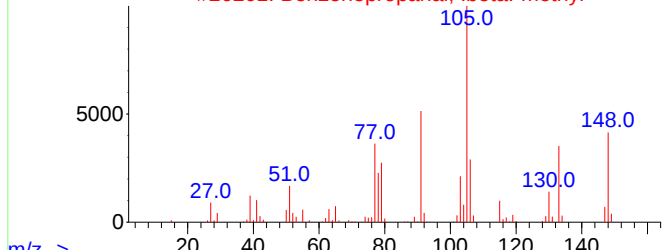


m/z 106.15 32.46%

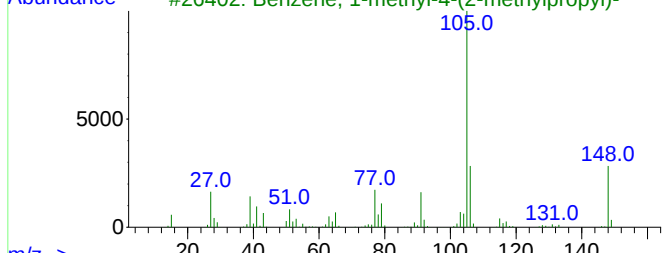
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m/z--> Abundance #26261: Benzenepropanal, .beta.-methyl-



m/z--> Abundance #26402: Benzene, 1-methyl-4-(2-methylpropyl)-



m/z--> Abundance

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011722\  
Data File : BG052058.D  
Acq On : 18 Jan 2022 18:42  
Operator : CG/JU  
Sample : N1100-15ME  
Misc :  
ALS Vial : 44 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLS8ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010622.M  
Quant Title : SVOA CALIBRATION

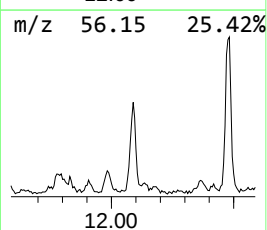
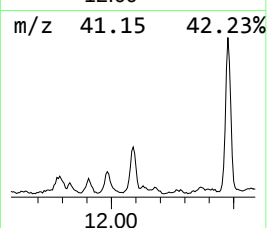
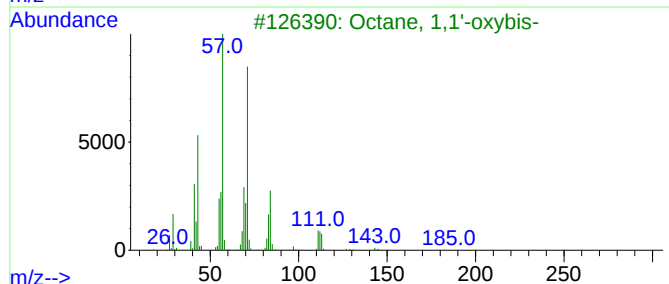
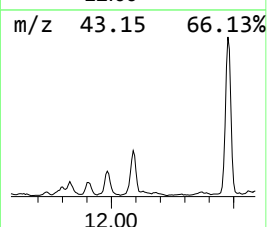
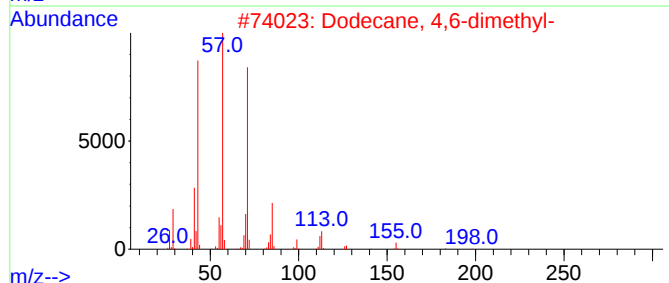
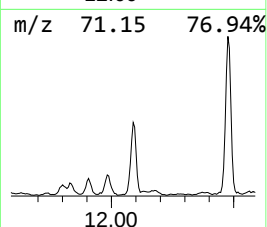
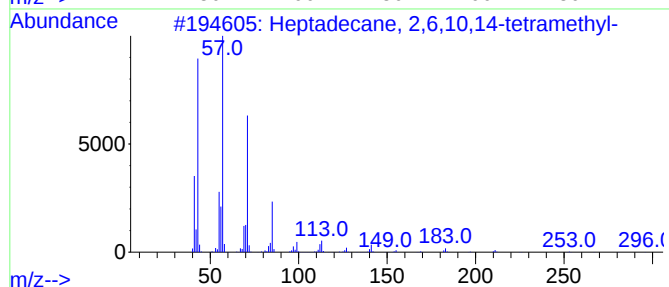
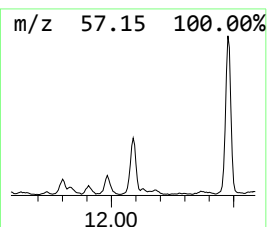
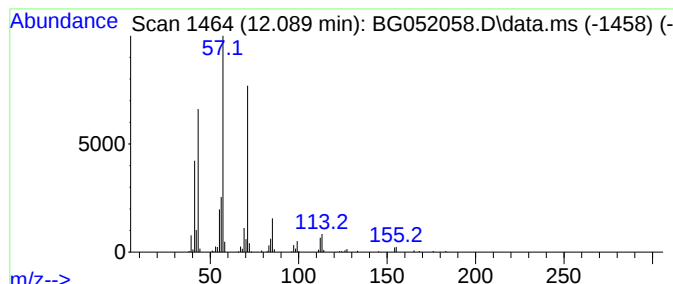
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 28 (DEL) Alkane: Straight-Chai... Concentration Rank 71

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 12.089 | 5.68 ng/ul | 290596 | Naphthalene-d8   | 11.090 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Heptadecane, 2,6,10,14-tetramethyl- | 296 | C21H44   | 018344-37-1  | 72   |
| 2    |    |   | Dodecane, 4,6-dimethyl-             | 198 | C14H30   | 061141-72-8  | 72   |
| 3    |    |   | Octane, 1,1'-oxybis-                | 242 | C16H34O  | 000629-82-3  | 72   |
| 4    |    |   | Oxalic acid, bis(2-ethylhexyl) e... | 314 | C18H34O4 | 1000309-39-0 | 64   |
| 5    |    |   | Octane, 2,6-dimethyl-               | 142 | C10H22   | 002051-30-1  | 64   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011722\  
Data File : BG052058.D  
Acq On : 18 Jan 2022 18:42  
Operator : CG/JU  
Sample : N1100-15ME  
Misc :  
ALS Vial : 44 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLS8ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010622.M  
Quant Title : SVOA CALIBRATION

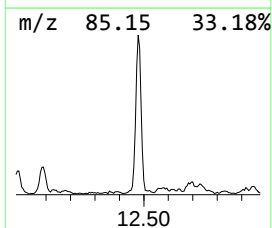
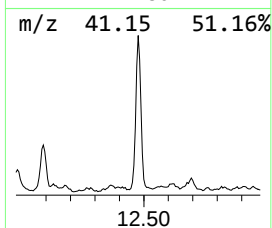
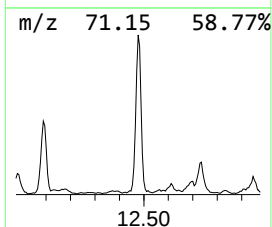
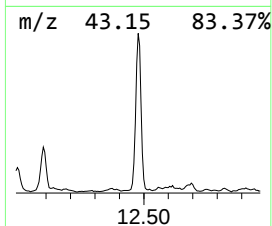
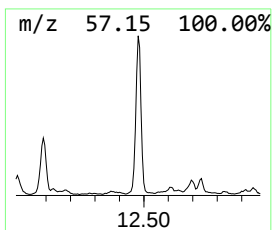
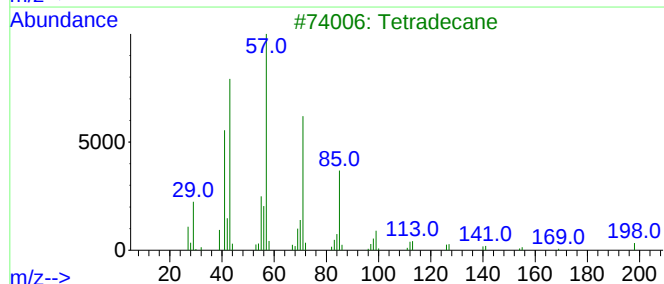
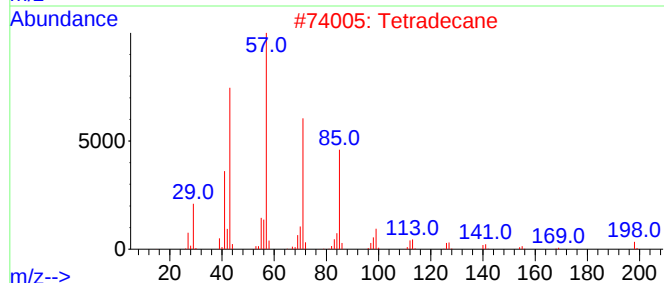
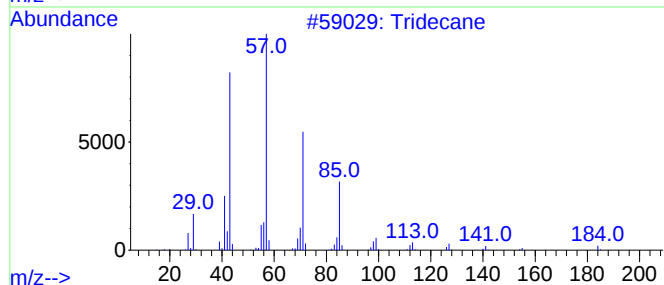
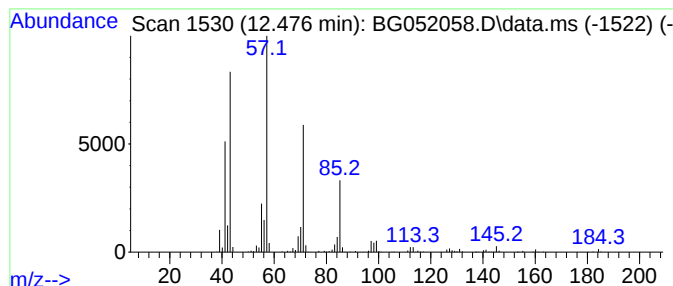
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 29 (DEL) Alkane: Straight-Chai... Concentration Rank 31

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 12.477 | 18.40 ng/ul | 940918 | Naphthalene-d8   | 11.090 |

| Hit# | of 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|------|--------------|-----|---------|-------------|------|
| 1    |      | Tridecane    | 184 | C13H28  | 000629-50-5 | 91   |
| 2    |      | Tetradecane  | 198 | C14H30  | 000629-59-4 | 86   |
| 3    |      | Tetradecane  | 198 | C14H30  | 000629-59-4 | 86   |
| 4    |      | Hexadecane   | 226 | C16H34  | 000544-76-3 | 86   |
| 5    |      | Tridecane    | 184 | C13H28  | 000629-50-5 | 81   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011722\  
Data File : BG052058.D  
Acq On : 18 Jan 2022 18:42  
Operator : CG/JU  
Sample : N1100-15ME  
Misc :  
ALS Vial : 44 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLS8ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010622.M  
Quant Title : SVOA CALIBRATION

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

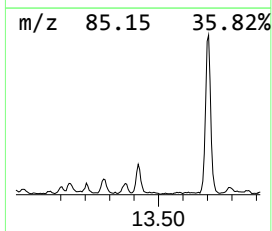
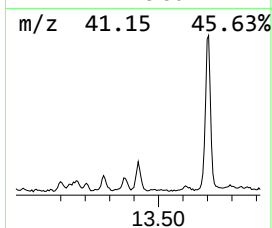
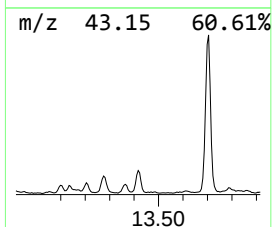
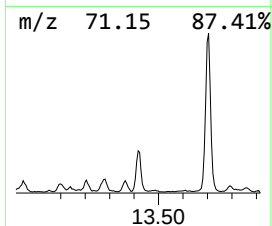
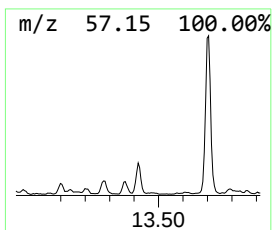
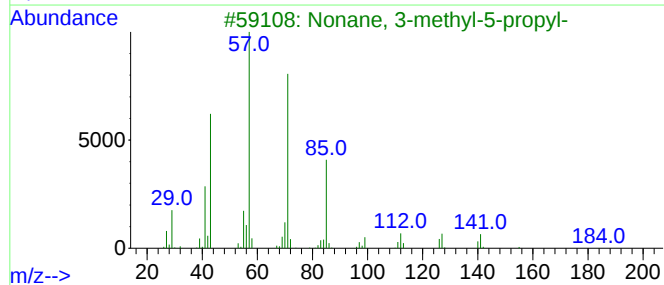
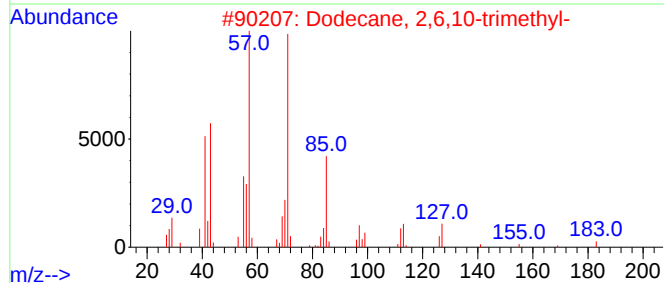
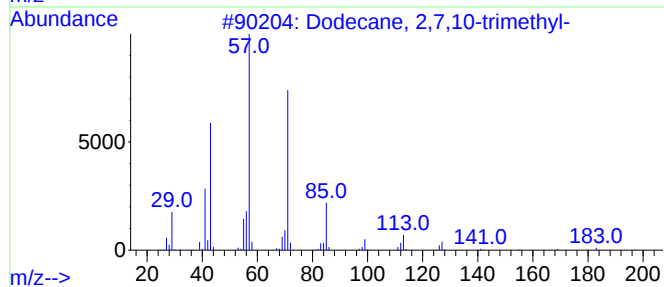
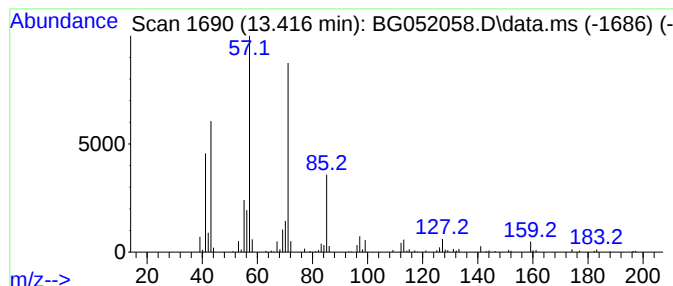
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Peak Number 30 (DEL) Alkane: Straight-Chai... Concentration Rank 35

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 13.416 | 17.16 ng/ul | 245892 | Acenaphthene-d10 | 14.891 |

| Hit# | of 5 | Tentative ID                       | MW  | MolForm | CAS#        | Qual |
|------|------|------------------------------------|-----|---------|-------------|------|
| 1    |      | Dodecane, 2,7,10-trimethyl-        | 212 | C15H32  | 074645-98-0 | 80   |
| 2    |      | Dodecane, 2,6,10-trimethyl-        | 212 | C15H32  | 003891-98-3 | 78   |
| 3    |      | Nonane, 3-methyl-5-propyl-         | 184 | C13H28  | 031081-18-2 | 78   |
| 4    |      | Butane, 2,2-dimethyl-              | 86  | C6H14   | 000075-83-2 | 53   |
| 5    |      | Hexadecane, 2,6,10,14-tetramethyl- | 282 | C20H42  | 000638-36-8 | 53   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011722\  
Data File : BG052058.D  
Acq On : 18 Jan 2022 18:42  
Operator : CG/JU  
Sample : N1100-15ME  
Misc :  
ALS Vial : 44 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLS8ME

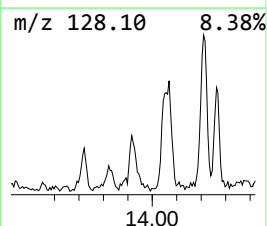
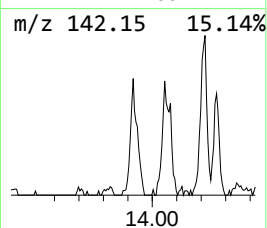
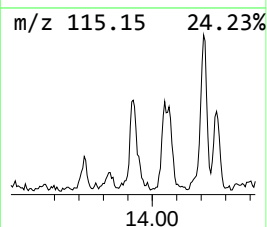
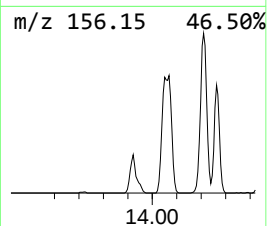
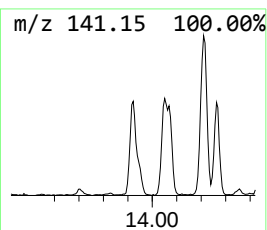
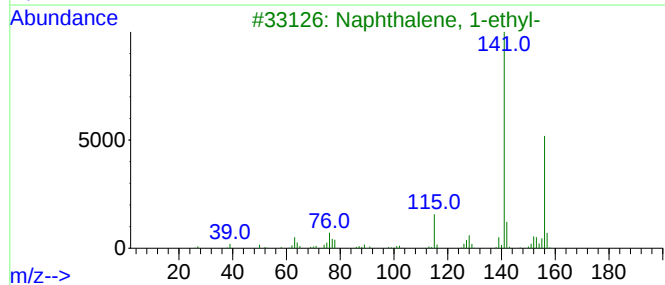
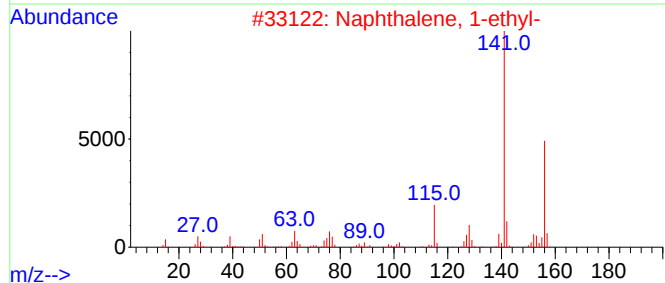
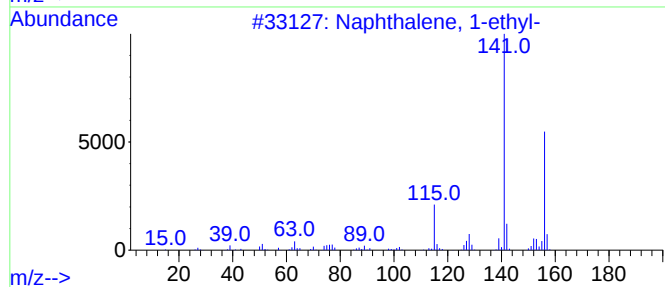
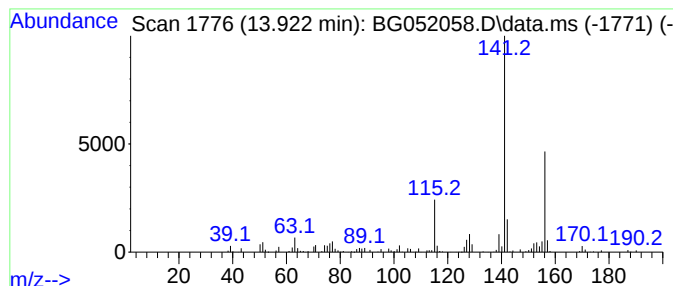
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Quant Title : SVOA CALIBRATION

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

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Peak Number 31 Naphthalene, 1-ethyl- Concentration Rank 33

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 13.922 | 17.69 ng/ul | 253588 | Acenaphthene-d10 | 14.891 |

| Hit# | of | 5 | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, 1-ethyl- | 156 | C12H12  | 001127-76-0 | 95   |
| 2    |    |   | Naphthalene, 1-ethyl- | 156 | C12H12  | 001127-76-0 | 94   |
| 3    |    |   | Naphthalene, 1-ethyl- | 156 | C12H12  | 001127-76-0 | 94   |
| 4    |    |   | Naphthalene, 2-ethyl- | 156 | C12H12  | 000939-27-5 | 94   |
| 5    |    |   | Naphthalene, 2-ethyl- | 156 | C12H12  | 000939-27-5 | 93   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011722\  
Data File : BG052058.D  
Acq On : 18 Jan 2022 18:42  
Operator : CG/JU  
Sample : N1100-15ME  
Misc :  
ALS Vial : 44 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLS8ME

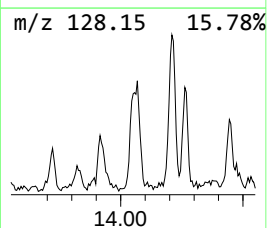
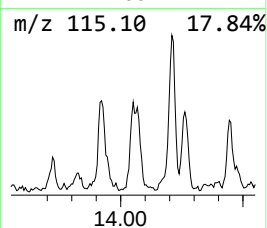
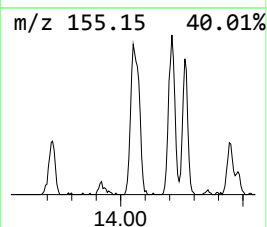
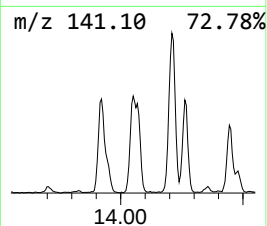
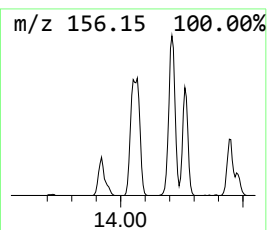
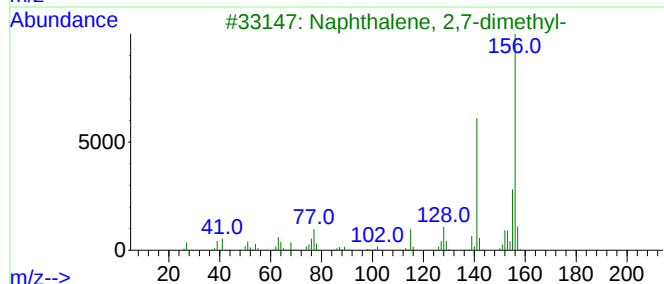
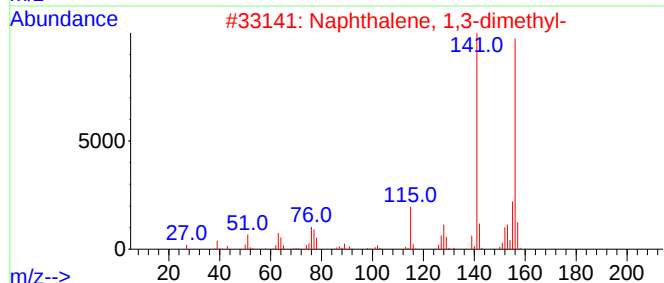
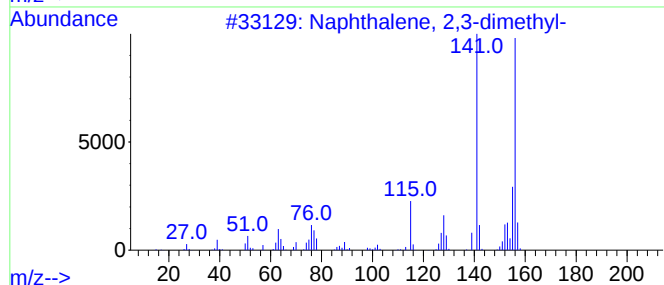
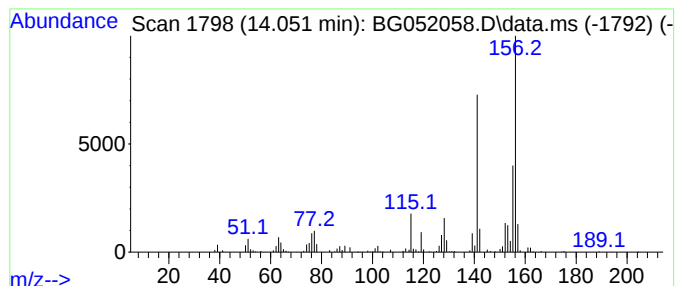
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Quant Title : SVOA CALIBRATION

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

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Peak Number 32 Naphthalene, 2,3-dimethyl- Concentration Rank 24

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 14.051 | 23.16 ng/ul | 331998 | Acenaphthene-d10 | 14.891 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, 2,3-dimethyl- | 156 | C12H12  | 000581-40-8 | 96   |
| 2    |    |   | Naphthalene, 1,3-dimethyl- | 156 | C12H12  | 000575-41-7 | 96   |
| 3    |    |   | Naphthalene, 2,7-dimethyl- | 156 | C12H12  | 000582-16-1 | 96   |
| 4    |    |   | Naphthalene, 2,6-dimethyl- | 156 | C12H12  | 000581-42-0 | 96   |
| 5    |    |   | Naphthalene, 1,4-dimethyl- | 156 | C12H12  | 000571-58-4 | 96   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011722\  
Data File : BG052058.D  
Acq On : 18 Jan 2022 18:42  
Operator : CG/JU  
Sample : N1100-15ME  
Misc :  
ALS Vial : 44 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLS8ME

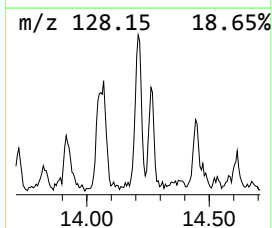
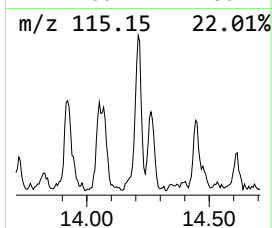
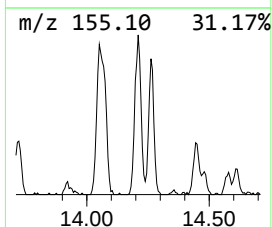
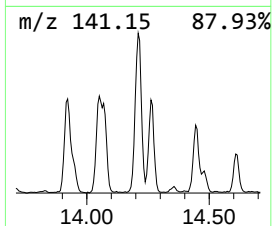
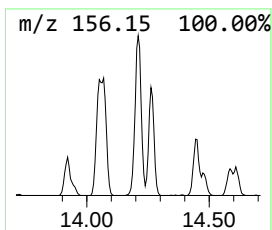
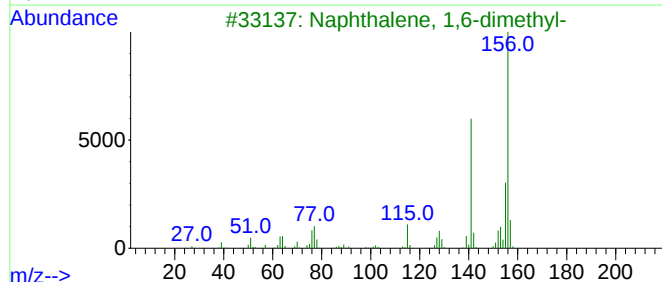
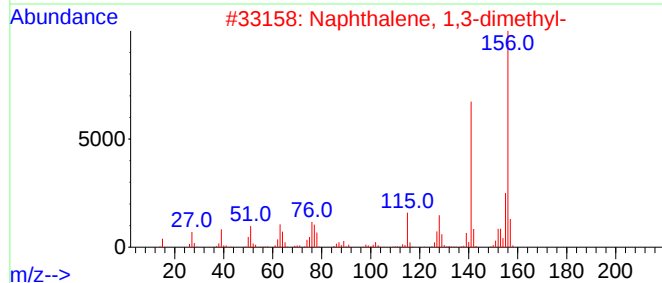
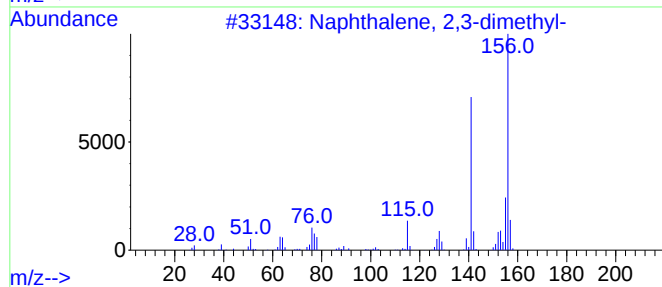
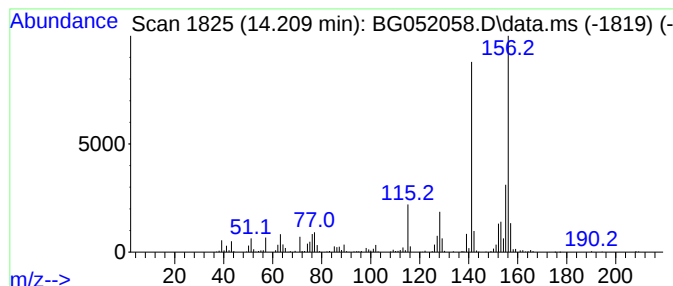
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Quant Title : SVOA CALIBRATION

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

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Peak Number 33 Naphthalene, 1,3-dimethyl- Concentration Rank 12

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 14.209 | 36.59 ng/ul | 524386 | Acenaphthene-d10 | 14.891 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, 2,3-dimethyl- | 156 | C12H12  | 000581-40-8 | 97   |
| 2    |    |   | Naphthalene, 1,3-dimethyl- | 156 | C12H12  | 000575-41-7 | 97   |
| 3    |    |   | Naphthalene, 1,6-dimethyl- | 156 | C12H12  | 000575-43-9 | 96   |
| 4    |    |   | Naphthalene, 2,6-dimethyl- | 156 | C12H12  | 000581-42-0 | 96   |
| 5    |    |   | Naphthalene, 2,7-dimethyl- | 156 | C12H12  | 000582-16-1 | 96   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011722\  
Data File : BG052058.D  
Acq On : 18 Jan 2022 18:42  
Operator : CG/JU  
Sample : N1100-15ME  
Misc :  
ALS Vial : 44 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLS8ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010622.M  
Quant Title : SVOA CALIBRATION

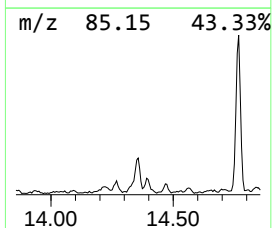
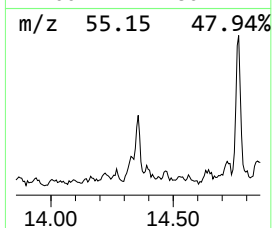
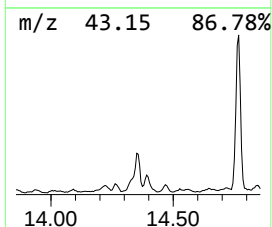
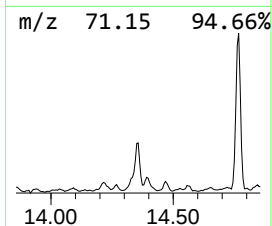
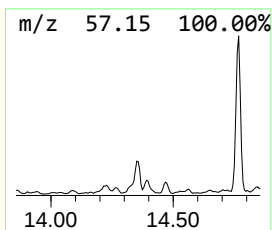
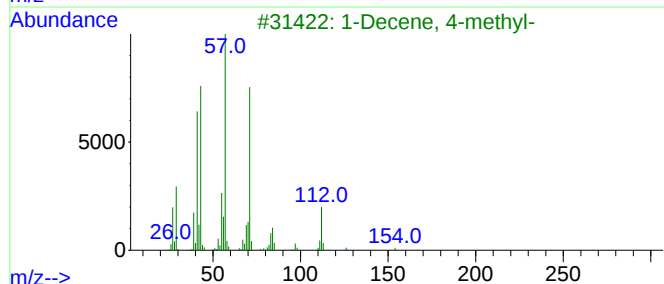
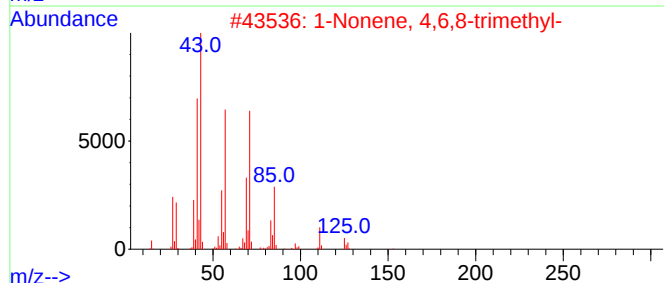
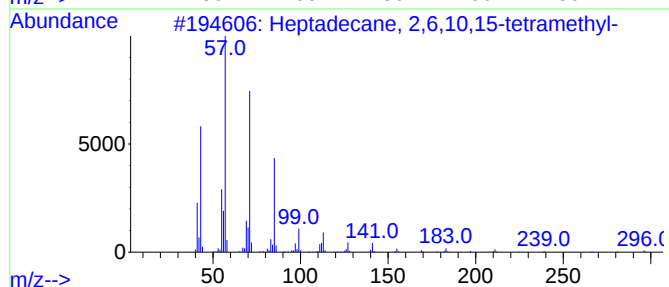
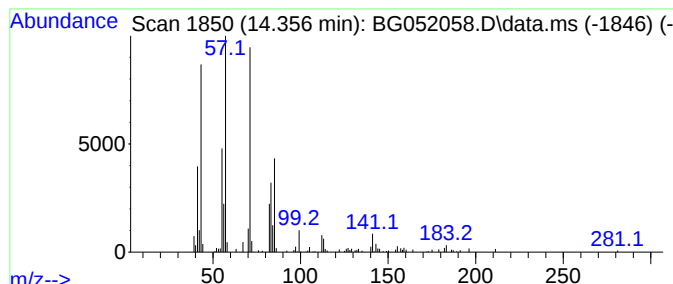
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 34 (DEL) Alkane: Straight-Chai... Concentration Rank 37

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 14.356 | 16.75 ng/ul | 240058 | Acenaphthene-d10 | 14.891 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|------|-------------------------------------|-----|---------|-------------|------|
| 1    |      | Heptadecane, 2,6,10,15-tetramethyl- | 296 | C21H44  | 054833-48-6 | 64   |
| 2    |      | 1-Nonene, 4,6,8-trimethyl-          | 168 | C12H24  | 054410-98-9 | 53   |
| 3    |      | 1-Decene, 4-methyl-                 | 154 | C11H22  | 013151-29-6 | 49   |
| 4    |      | Octane, 2,3,3-trimethyl-            | 156 | C11H24  | 062016-30-2 | 43   |
| 5    |      | Decane, 2,3,5-trimethyl-            | 184 | C13H28  | 062238-11-3 | 38   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011722\  
Data File : BG052058.D  
Acq On : 18 Jan 2022 18:42  
Operator : CG/JU  
Sample : N1100-15ME  
Misc :  
ALS Vial : 44 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLS8ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010622.M  
Quant Title : SVOA CALIBRATION

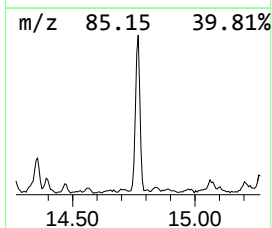
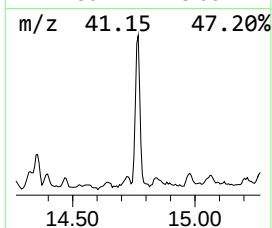
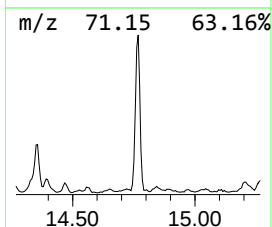
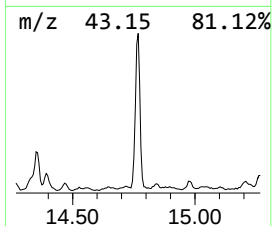
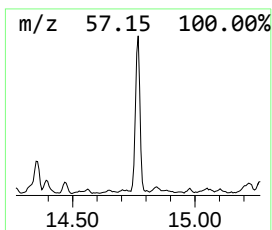
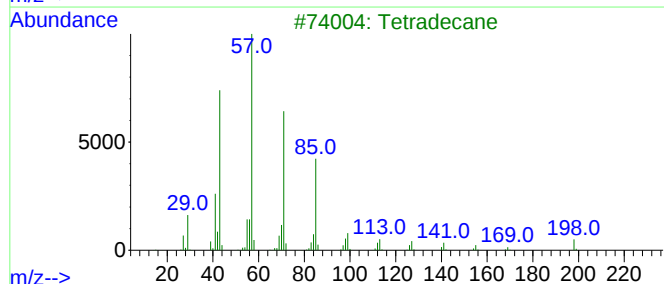
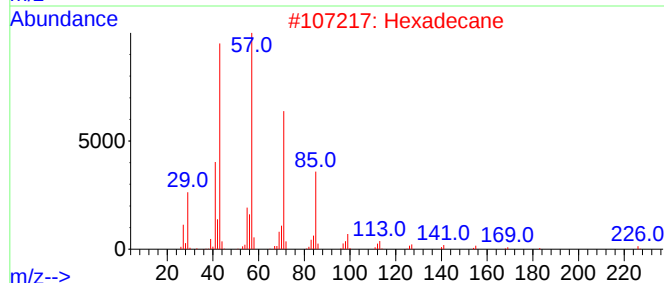
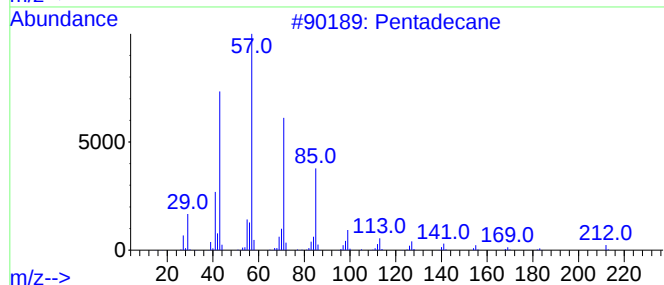
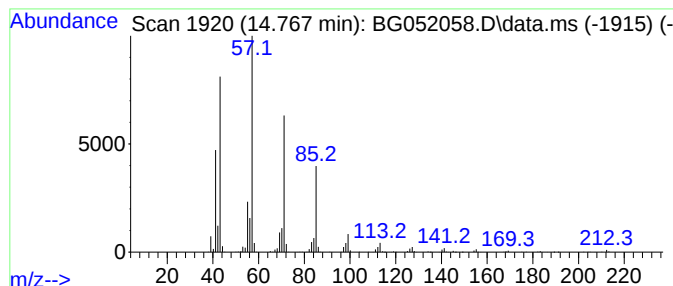
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 35 (DEL) Alkane: Straight-Chai... Concentration Rank 5

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 14.768 | 61.57 ng/ul | 882500 | Acenaphthene-d10 | 14.891 |

| Hit# | of 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|------|--------------|-----|---------|-------------|------|
| 1    |      | Pentadecane  | 212 | C15H32  | 000629-62-9 | 94   |
| 2    |      | Hexadecane   | 226 | C16H34  | 000544-76-3 | 91   |
| 3    |      | Tetradecane  | 198 | C14H30  | 000629-59-4 | 90   |
| 4    |      | Tridecane    | 184 | C13H28  | 000629-50-5 | 90   |
| 5    |      | Tetradecane  | 198 | C14H30  | 000629-59-4 | 90   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011722\  
Data File : BG052058.D  
Acq On : 18 Jan 2022 18:42  
Operator : CG/JU  
Sample : N1100-15ME  
Misc :  
ALS Vial : 44 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLS8ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010622.M  
Quant Title : SVOA CALIBRATION

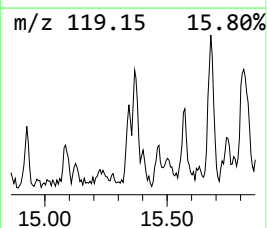
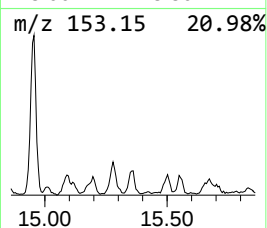
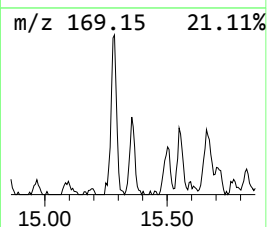
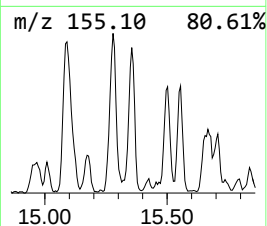
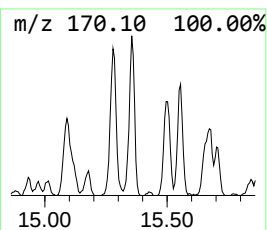
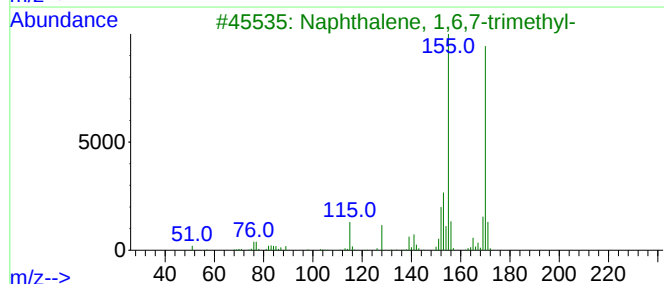
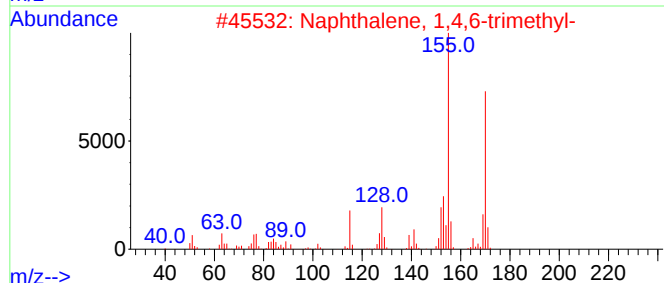
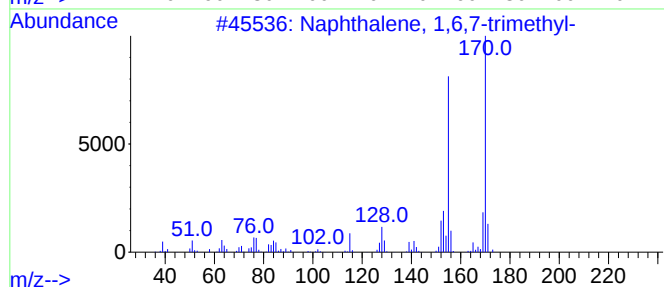
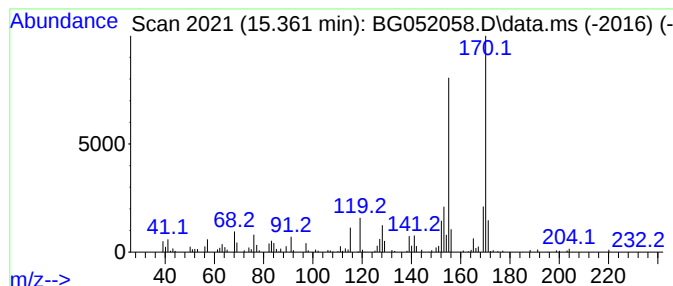
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 36 Naphthalene, 1,6,7-trimethyl- Concentration Rank 21

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 15.361 | 24.41 ng/ul | 349793 | Acenaphthene-d10 | 14.891 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|---------|--------------|------|
| 1    |    |   | Naphthalene, 1,6,7-trimethyl-       | 170 | C13H14  | 002245-38-7  | 96   |
| 2    |    |   | Naphthalene, 1,4,6-trimethyl-       | 170 | C13H14  | 002131-42-2  | 96   |
| 3    |    |   | Naphthalene, 1,6,7-trimethyl-       | 170 | C13H14  | 002245-38-7  | 95   |
| 4    |    |   | 1-(2,2-Dimethylcyclopropyl)-2-ph... | 170 | C13H14  | 1000294-91-1 | 94   |
| 5    |    |   | Naphthalene, 2,3,6-trimethyl-       | 170 | C13H14  | 000829-26-5  | 94   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011722\  
Data File : BG052058.D  
Acq On : 18 Jan 2022 18:42  
Operator : CG/JU  
Sample : N1100-15ME  
Misc :  
ALS Vial : 44 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLS8ME

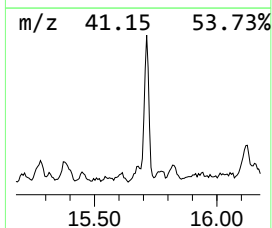
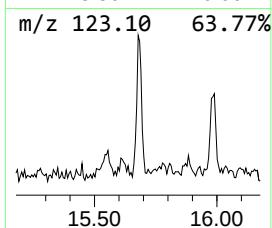
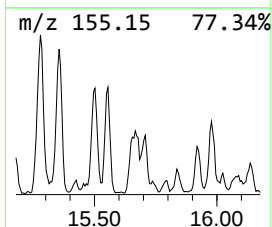
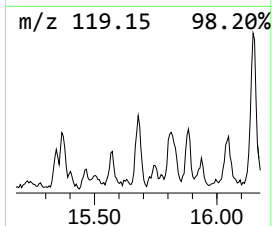
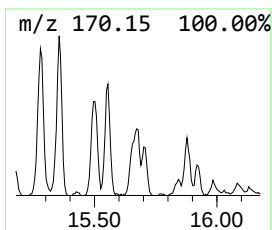
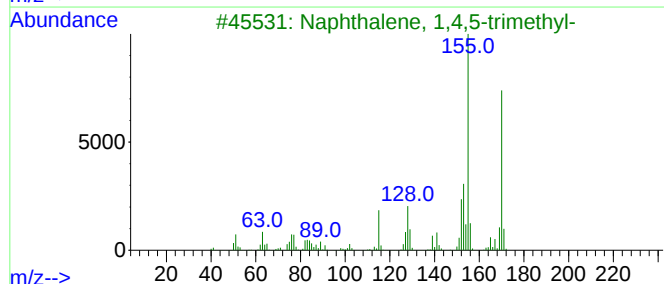
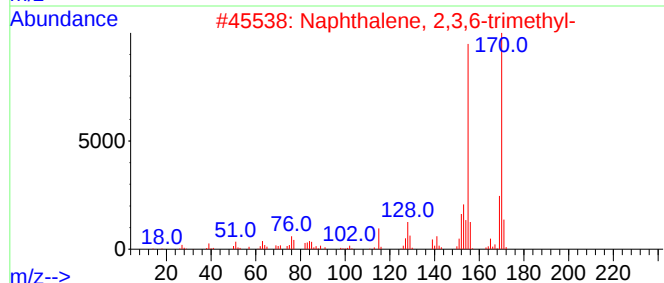
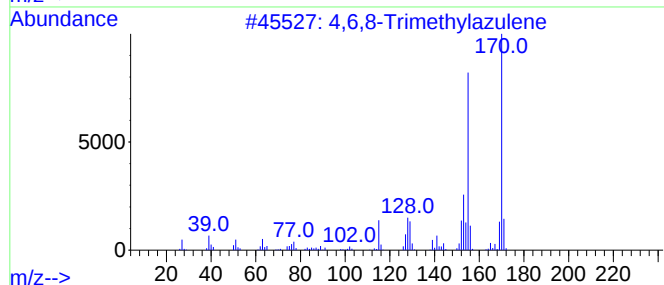
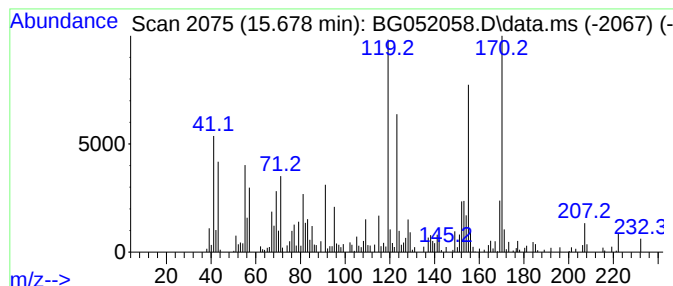
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010622.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 37 4,6,8-Trimethylazulene Concentration Rank 34

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 15.678 | 17.57 ng/ul | 251756 | Acenaphthene-d10 | 14.891 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|---------|--------------|------|
| 1    |    |   | 4,6,8-Trimethylazulene              | 170 | C13H14  | 000941-81-1  | 90   |
| 2    |    |   | Naphthalene, 2,3,6-trimethyl-       | 170 | C13H14  | 000829-26-5  | 90   |
| 3    |    |   | Naphthalene, 1,4,5-trimethyl-       | 170 | C13H14  | 002131-41-1  | 89   |
| 4    |    |   | Naphthalene, 1,6,7-trimethyl-       | 170 | C13H14  | 002245-38-7  | 87   |
| 5    |    |   | 1-(2,2-Dimethylcyclopropyl)-2-ph... | 170 | C13H14  | 1000294-91-1 | 86   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011722\  
Data File : BG052058.D  
Acq On : 18 Jan 2022 18:42  
Operator : CG/JU  
Sample : N1100-15ME  
Misc :  
ALS Vial : 44 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLS8ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010622.M  
Quant Title : SVOA CALIBRATION

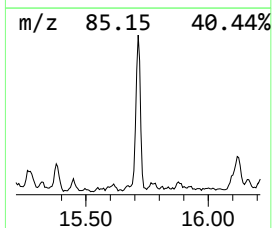
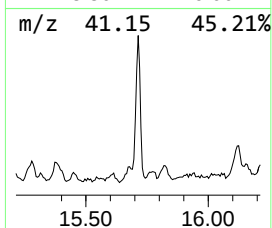
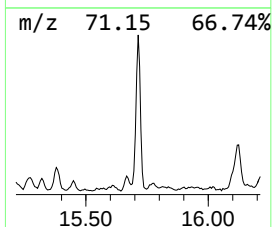
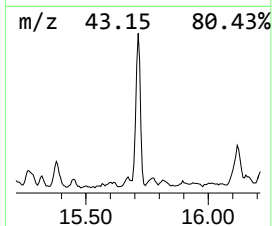
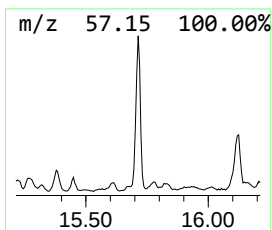
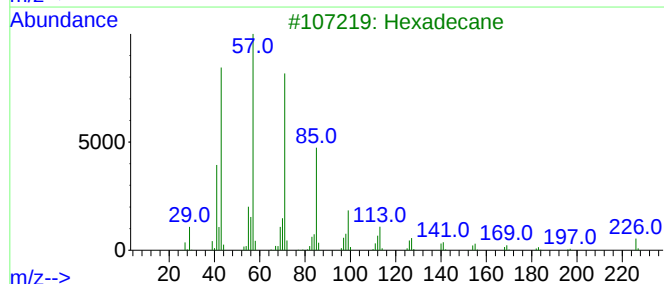
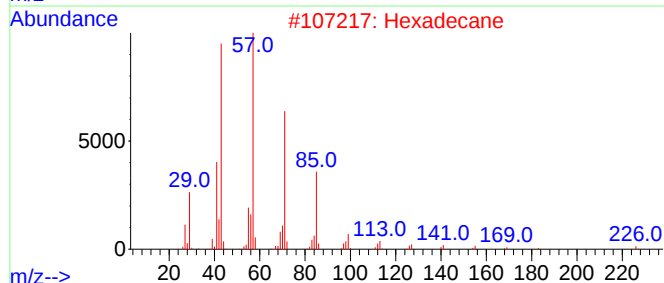
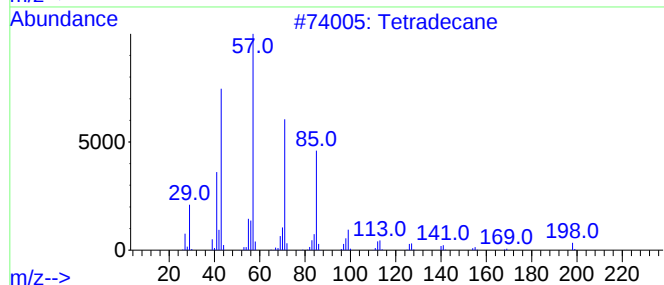
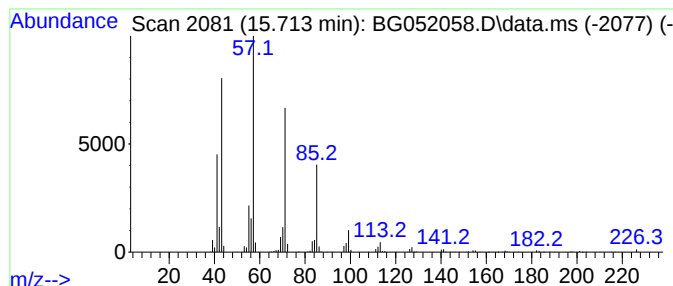
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 38 (DEL) Alkane: Straight-Chai... Concentration Rank 8

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 15.713 | 52.61 ng/ul | 754052 | Acenaphthene-d10 | 14.891 |

| Hit# | of 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|------|--------------|-----|---------|-------------|------|
| 1    |      | Tetradecane  | 198 | C14H30  | 000629-59-4 | 97   |
| 2    |      | Hexadecane   | 226 | C16H34  | 000544-76-3 | 95   |
| 3    |      | Hexadecane   | 226 | C16H34  | 000544-76-3 | 94   |
| 4    |      | Pentadecane  | 212 | C15H32  | 000629-62-9 | 90   |
| 5    |      | Tridecane    | 184 | C13H28  | 000629-50-5 | 90   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011722\  
Data File : BG052058.D  
Acq On : 18 Jan 2022 18:42  
Operator : CG/JU  
Sample : N1100-15ME  
Misc :  
ALS Vial : 44 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLS8ME

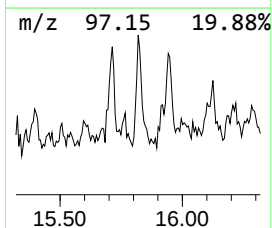
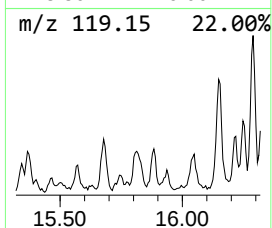
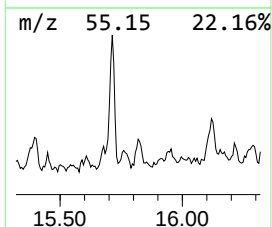
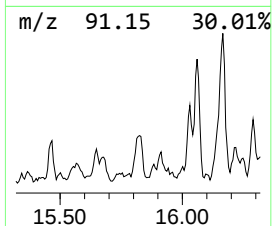
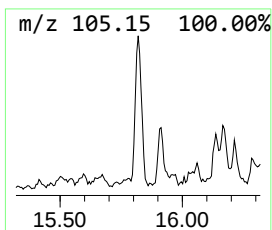
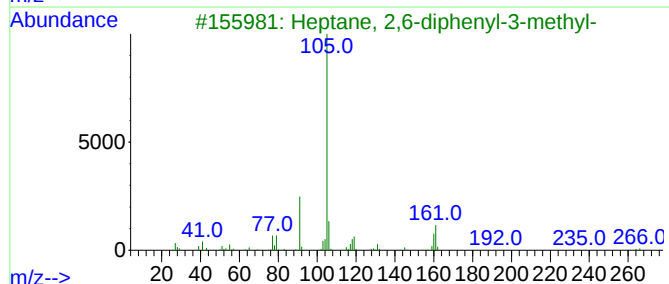
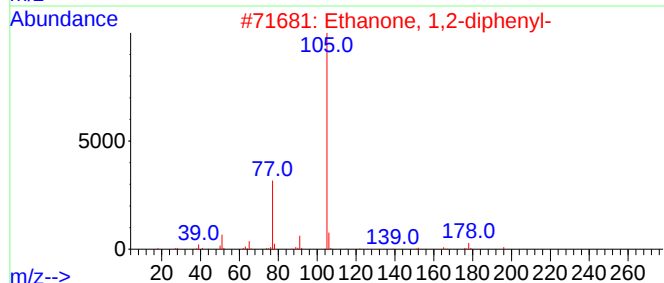
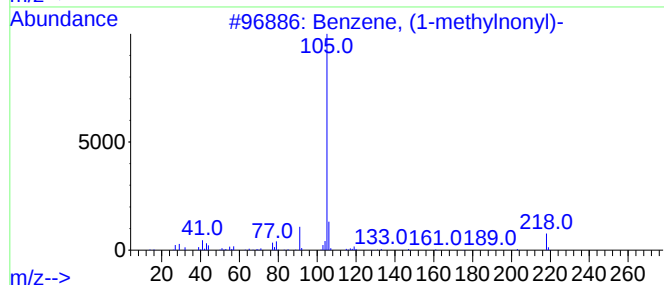
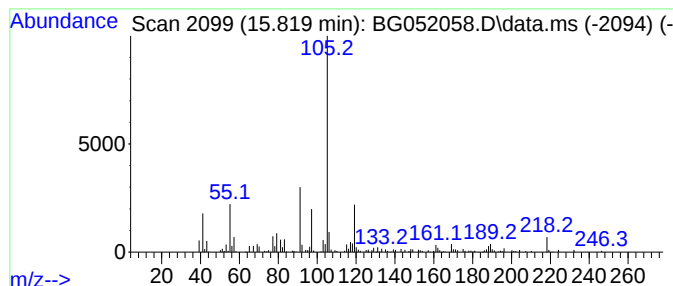
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Quant Title : SVOA CALIBRATION

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

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Peak Number 39 unknown-02 Concentration Rank 44

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 15.819 | 14.56 ng/ul | 208732 | Acenaphthene-d10 | 14.891 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|------------|--------------|------|
| 1    |    |   | Benzene, (1-methylnonyl)-           | 218 | C16H26     | 004537-13-7  | 27   |
| 2    |    |   | Ethanone, 1,2-diphenyl-             | 196 | C14H12O    | 000451-40-1  | 22   |
| 3    |    |   | Heptane, 2,6-diphenyl-3-methyl-     | 266 | C20H26     | 1000161-22-5 | 22   |
| 4    |    |   | Ethyl 2-cyano-3-methyl-2-(0-meth... | 259 | C16H21NO2  | 029107-35-5  | 22   |
| 5    |    |   | N-(4-Methylthiazol-2-yl)benzamide   | 218 | C11H10N2OS | 037529-67-2  | 22   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011722\  
Data File : BG052058.D  
Acq On : 18 Jan 2022 18:42  
Operator : CG/JU  
Sample : N1100-15ME  
Misc :  
ALS Vial : 44 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLS8ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010622.M  
Quant Title : SVOA CALIBRATION

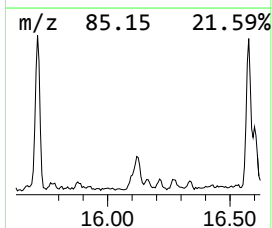
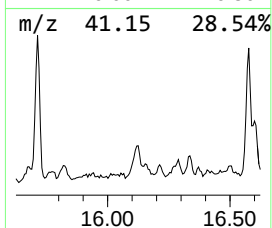
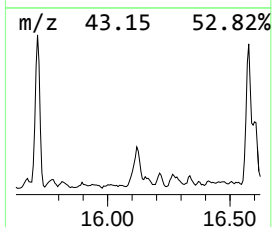
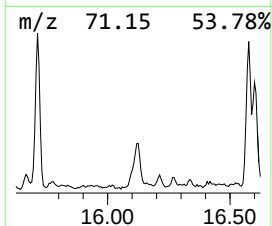
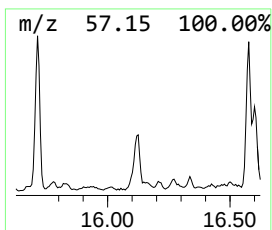
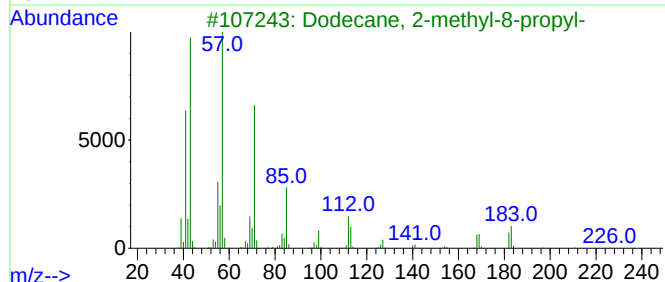
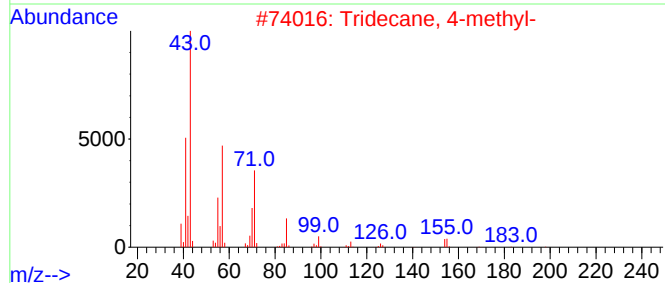
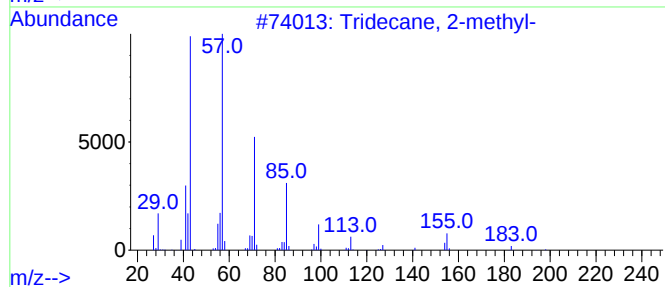
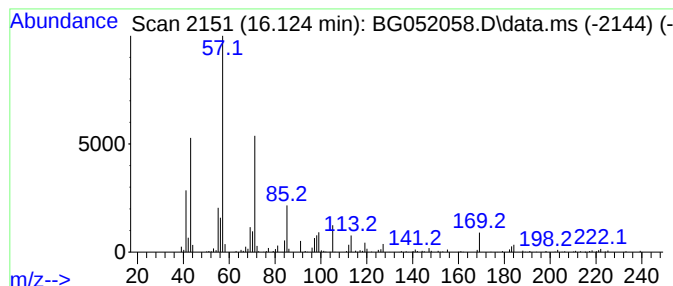
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 40 (DEL) Alkane: Straight-Chai... Concentration Rank 18

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 16.125 | 25.53 ng/ul | 365965 | Acenaphthene-d10 | 14.891 |

| Hit# | of 5 | Tentative ID                 | MW  | MolForm | CAS#        | Qual |
|------|------|------------------------------|-----|---------|-------------|------|
| 1    |      | Tridecane, 2-methyl-         | 198 | C14H30  | 001560-96-9 | 74   |
| 2    |      | Tridecane, 4-methyl-         | 198 | C14H30  | 026730-12-1 | 64   |
| 3    |      | Dodecane, 2-methyl-8-propyl- | 226 | C16H34  | 055045-07-3 | 59   |
| 4    |      | Tridecane, 3-methyl-         | 198 | C14H30  | 006418-41-3 | 58   |
| 5    |      | Heptadecane, 7-methyl-       | 254 | C18H38  | 020959-33-5 | 58   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011722\  
Data File : BG052058.D  
Acq On : 18 Jan 2022 18:42  
Operator : CG/JU  
Sample : N1100-15ME  
Misc :  
ALS Vial : 44 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLS8ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010622.M  
Quant Title : SVOA CALIBRATION

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

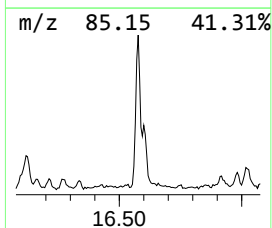
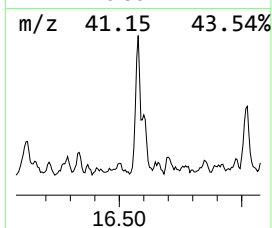
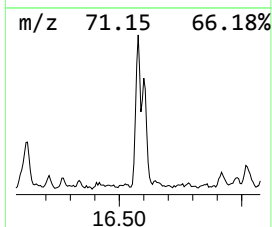
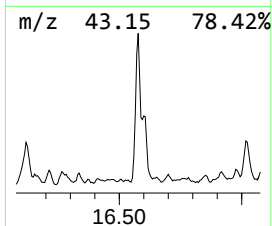
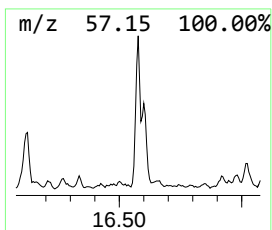
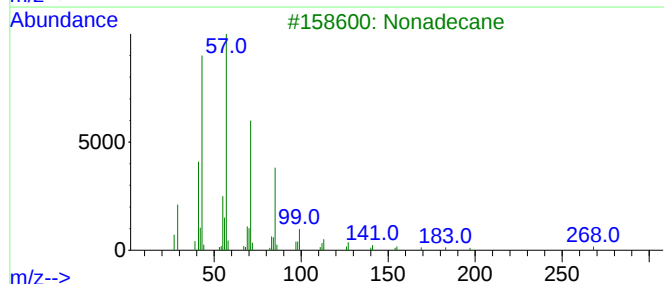
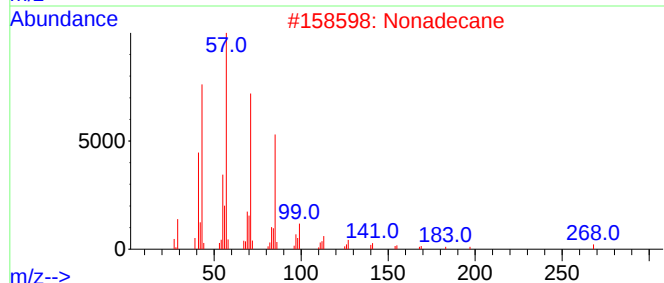
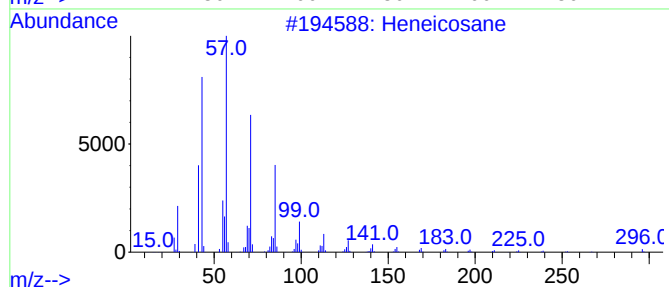
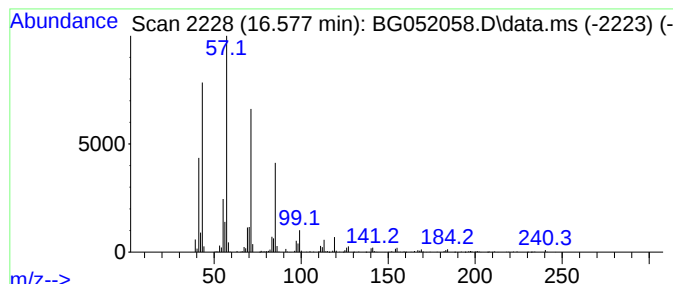
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Peak Number 41 (DEL) Alkane: Straight-Chai... Concentration Rank 11

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 16.577 | 38.01 ng/ul | 721151 | Phenanthrene-d10 | 17.640 |

| Hit# | of 5 | Tentative ID         | MW  | MolForm | CAS#        | Qual |
|------|------|----------------------|-----|---------|-------------|------|
| 1    |      | Heneicosane          | 296 | C21H44  | 000629-94-7 | 91   |
| 2    |      | Nonadecane           | 268 | C19H40  | 000629-92-5 | 91   |
| 3    |      | Nonadecane           | 268 | C19H40  | 000629-92-5 | 91   |
| 4    |      | Dodecane, 2-methyl-  | 184 | C13H28  | 001560-97-0 | 90   |
| 5    |      | Eicosane, 10-methyl- | 296 | C21H44  | 054833-23-7 | 90   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011722\  
Data File : BG052058.D  
Acq On : 18 Jan 2022 18:42  
Operator : CG/JU  
Sample : N1100-15ME  
Misc :  
ALS Vial : 44 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLS8ME

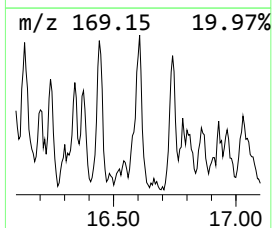
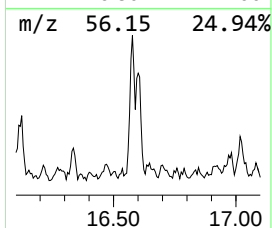
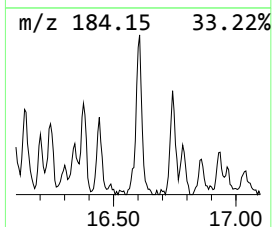
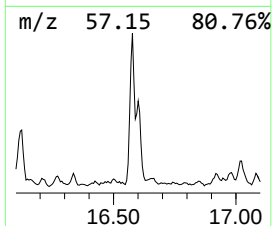
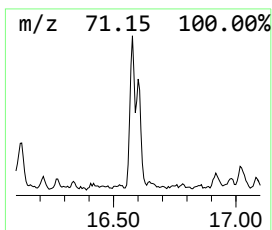
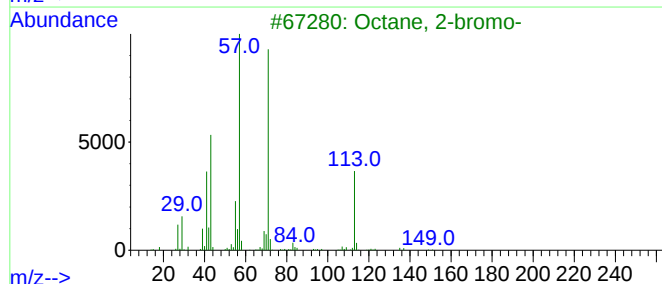
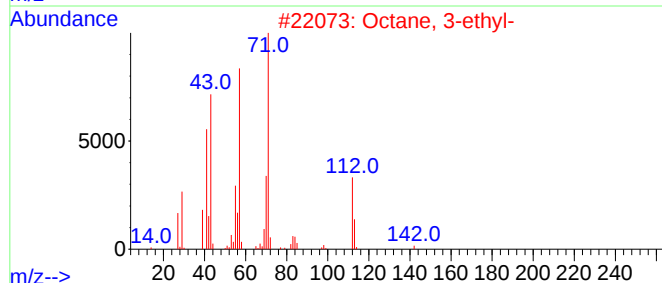
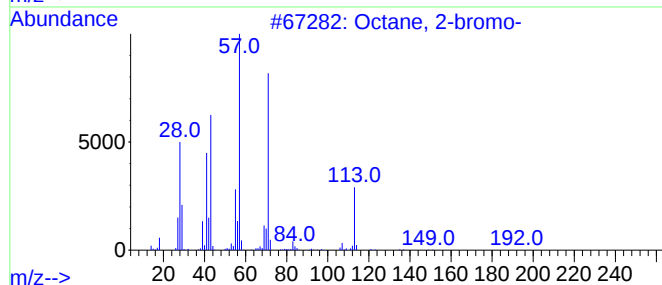
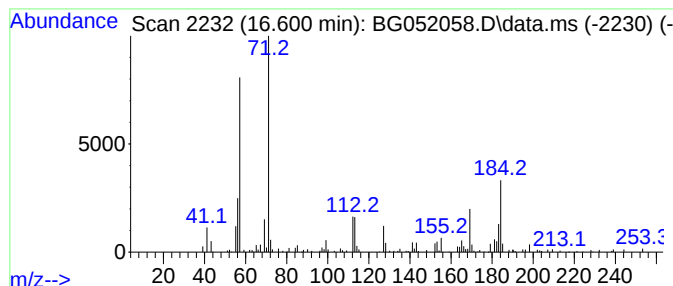
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Quant Title : SVOA CALIBRATION

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

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Peak Number 42 unknown-03 Concentration Rank 27

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 16.600 | 20.13 ng/ul | 381915 | Phenanthrene-d10 | 17.640 |

| Hit# | of | 5 | Tentative ID         | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------|-----|---------|-------------|------|
| 1    |    |   | Octane, 2-bromo-     | 192 | C8H17Br | 000557-35-7 | 47   |
| 2    |    |   | Octane, 3-ethyl-     | 142 | C10H22  | 005881-17-4 | 46   |
| 3    |    |   | Octane, 2-bromo-     | 192 | C8H17Br | 000557-35-7 | 43   |
| 4    |    |   | Nonane, 3-methyl-    | 142 | C10H22  | 005911-04-6 | 38   |
| 5    |    |   | Octane, 1,1'-oxybis- | 242 | C16H34O | 000629-82-3 | 38   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011722\  
Data File : BG052058.D  
Acq On : 18 Jan 2022 18:42  
Operator : CG/JU  
Sample : N1100-15ME  
Misc :  
ALS Vial : 44 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLS8ME

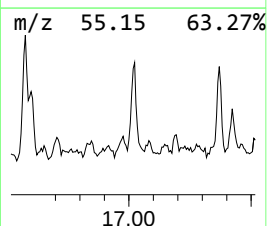
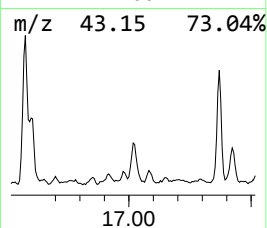
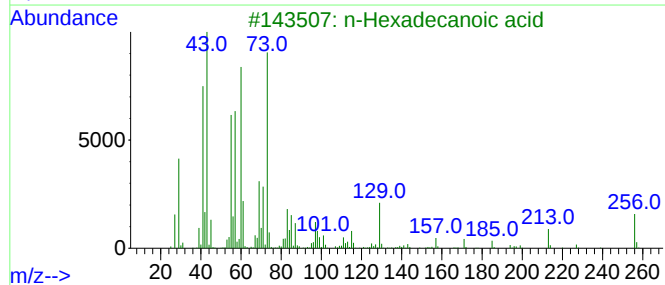
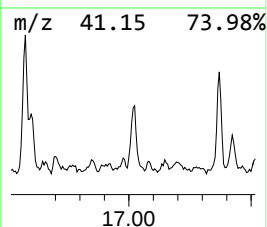
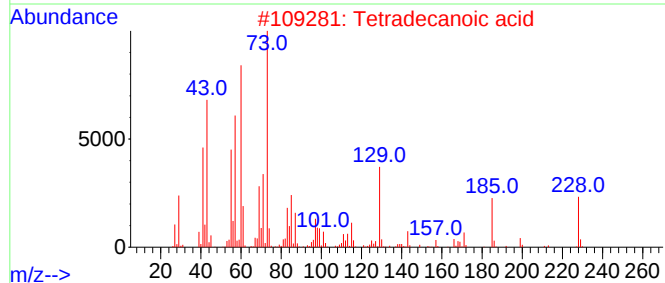
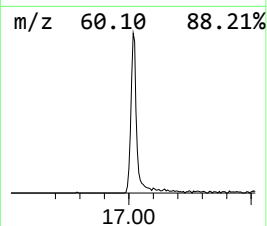
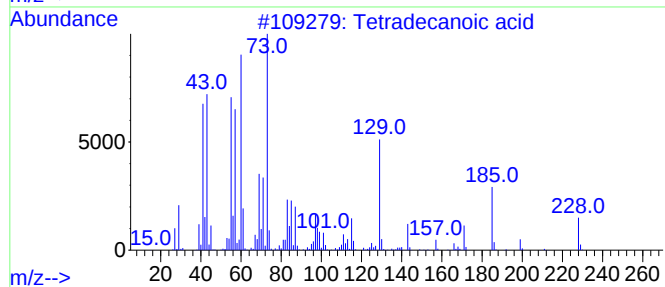
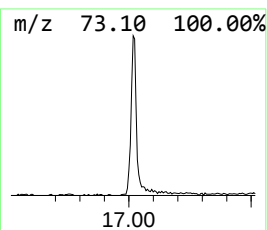
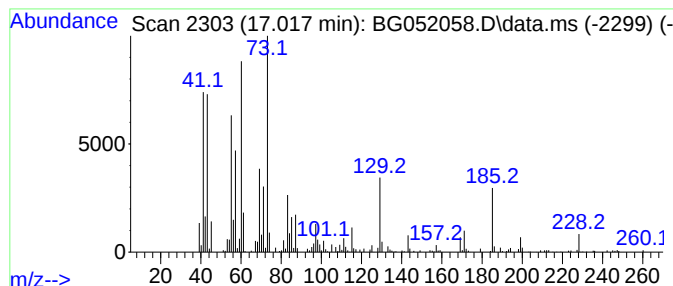
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Quant Title : SVOA CALIBRATION

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

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Peak Number 43 Tetradecanoic acid Concentration Rank 15

| R.T.      | EstConc             | Area   | Relative to ISTD | R.T.        |      |
|-----------|---------------------|--------|------------------|-------------|------|
| 17.017    | 26.52 ng/ul         | 503099 | Phenanthrene-d10 | 17.640      |      |
| Hit# of 5 | Tentative ID        | MW     | MolForm          | CAS#        | Qual |
| 1         | Tetradecanoic acid  | 228    | C14H28O2         | 000544-63-8 | 90   |
| 2         | Tetradecanoic acid  | 228    | C14H28O2         | 000544-63-8 | 87   |
| 3         | n-Hexadecanoic acid | 256    | C16H32O2         | 000057-10-3 | 64   |
| 4         | Dodecanoic acid     | 200    | C12H24O2         | 000143-07-7 | 64   |
| 5         | Dodecanoic acid     | 200    | C12H24O2         | 000143-07-7 | 62   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011722\  
Data File : BG052058.D  
Acq On : 18 Jan 2022 18:42  
Operator : CG/JU  
Sample : N1100-15ME  
Misc :  
ALS Vial : 44 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLS8ME

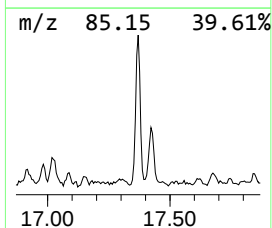
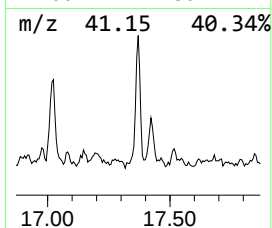
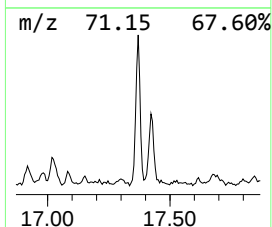
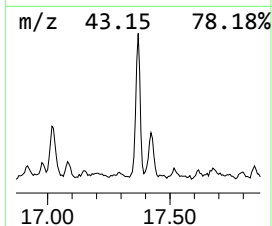
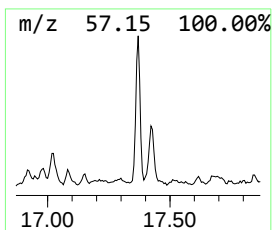
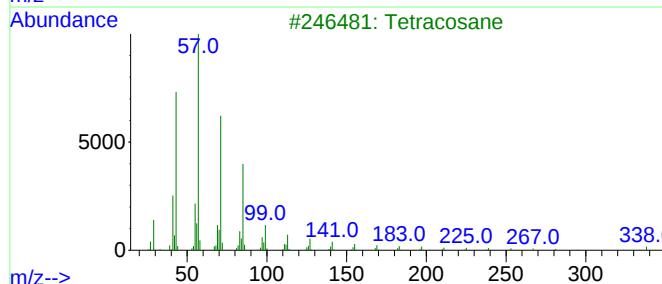
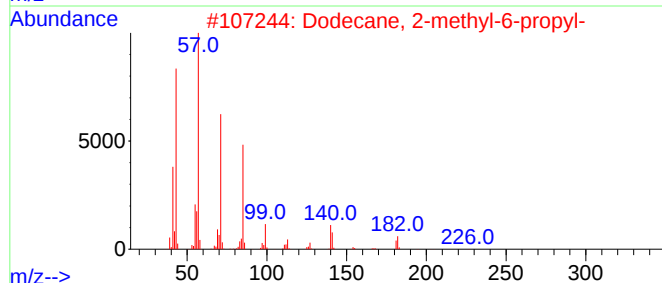
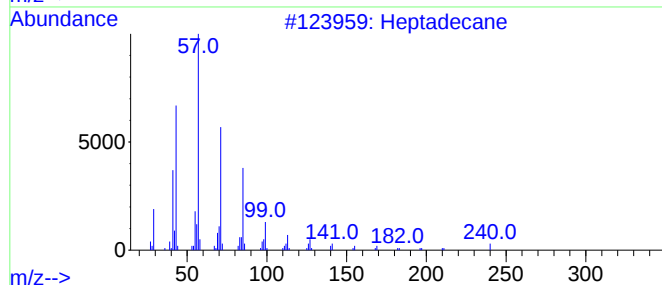
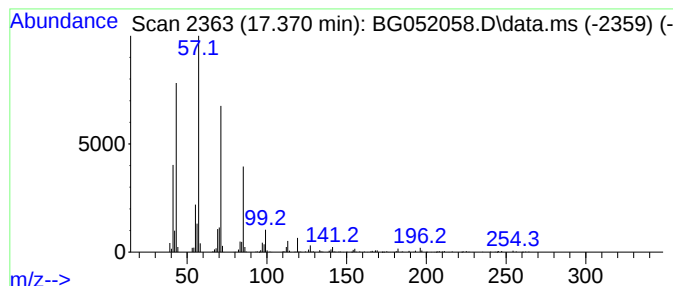
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Quant Title : SVOA CALIBRATION

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

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Peak Number 44 (DEL) Alkane: Straight-Chai... Concentration Rank 20

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 17.370 | 24.69 ng/ul | 468386 | Phenanthrene-d10 | 17.640 |

| Hit# | of 5 | Tentative ID                 | MW  | MolForm | CAS#        | Qual |
|------|------|------------------------------|-----|---------|-------------|------|
| 1    |      | Heptadecane                  | 240 | C17H36  | 000629-78-7 | 94   |
| 2    |      | Dodecane, 2-methyl-6-propyl- | 226 | C16H34  | 055045-08-4 | 93   |
| 3    |      | Tetracosane                  | 338 | C24H50  | 000646-31-1 | 90   |
| 4    |      | Heptadecane                  | 240 | C17H36  | 000629-78-7 | 87   |
| 5    |      | Pentadecane                  | 212 | C15H32  | 000629-62-9 | 86   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011722\  
Data File : BG052058.D  
Acq On : 18 Jan 2022 18:42  
Operator : CG/JU  
Sample : N1100-15ME  
Misc :  
ALS Vial : 44 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLS8ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010622.M  
Quant Title : SVOA CALIBRATION

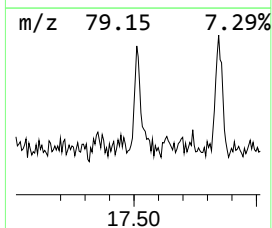
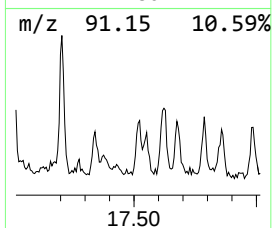
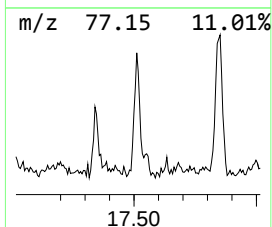
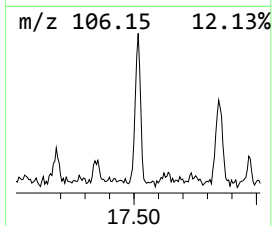
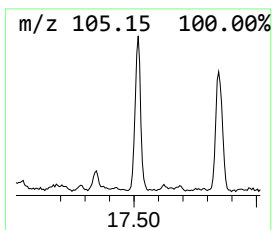
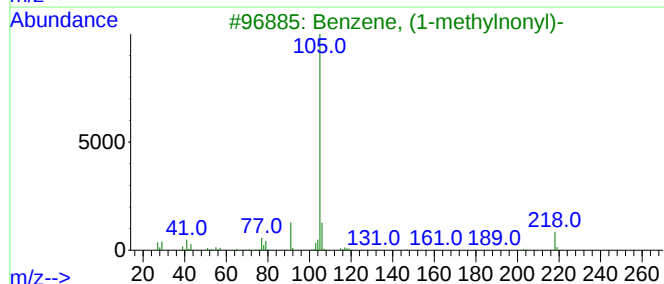
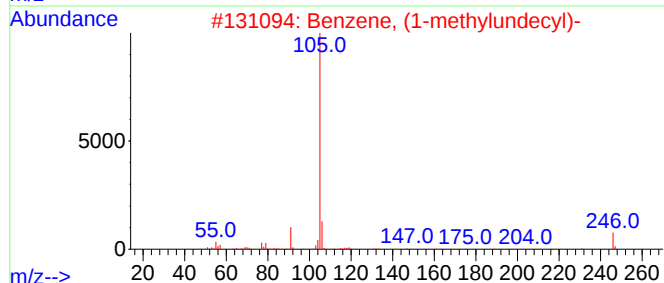
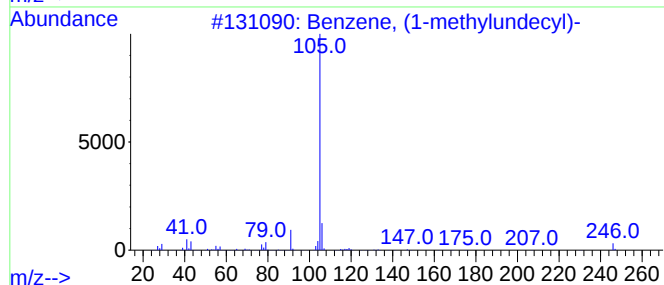
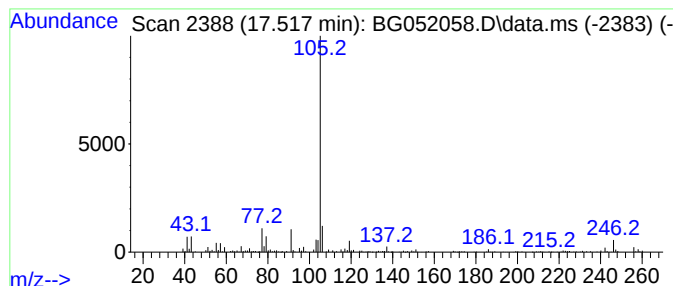
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 45 Benzene, (1-methylundecyl)- Concentration Rank 46

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 17.517 | 14.14 ng/ul | 268237 | Phenanthrene-d10 | 17.640 |

| Hit# | of | 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, (1-methylundecyl)- | 246 | C18H30  | 002719-61-1 | 81   |
| 2    |    |   | Benzene, (1-methylundecyl)- | 246 | C18H30  | 002719-61-1 | 72   |
| 3    |    |   | Benzene, (1-methylnonyl)-   | 218 | C16H26  | 004537-13-7 | 59   |
| 4    |    |   | Benzene, (1-methylundecyl)- | 246 | C18H30  | 002719-61-1 | 58   |
| 5    |    |   | Benzene, 1-methyl-4-propyl- | 134 | C10H14  | 001074-55-1 | 53   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011722\  
Data File : BG052058.D  
Acq On : 18 Jan 2022 18:42  
Operator : CG/JU  
Sample : N1100-15ME  
Misc :  
ALS Vial : 44 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLS8ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010622.M  
Quant Title : SVOA CALIBRATION

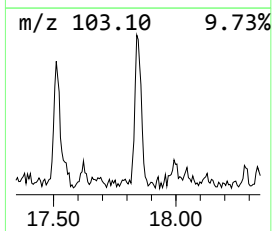
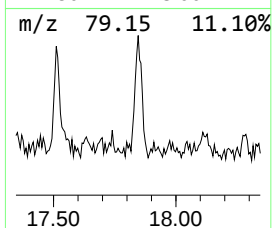
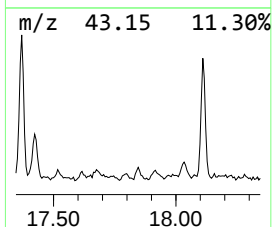
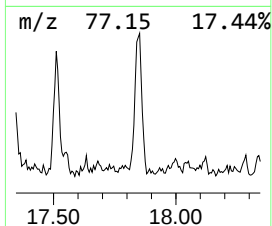
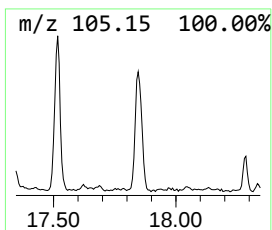
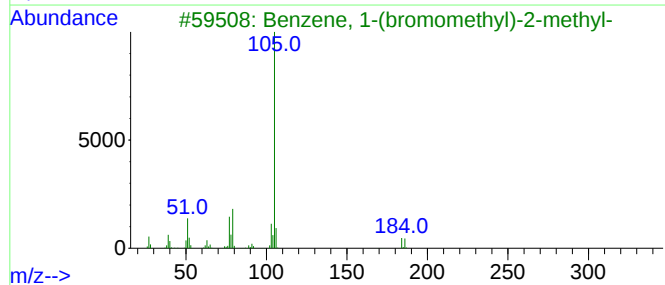
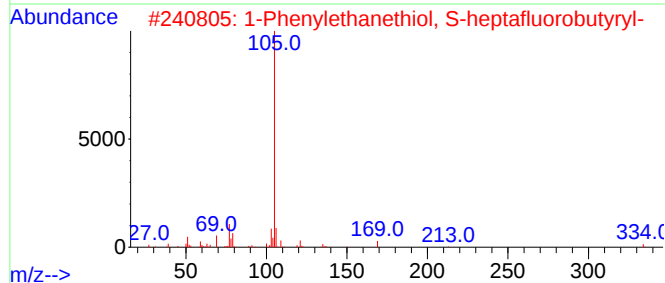
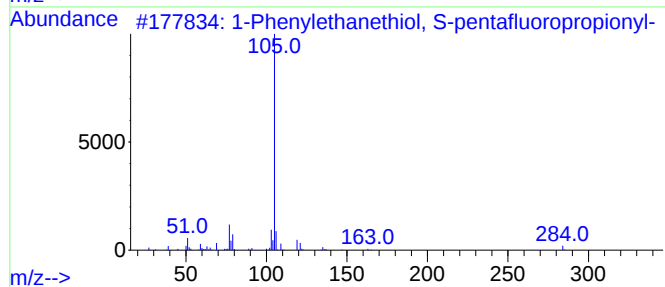
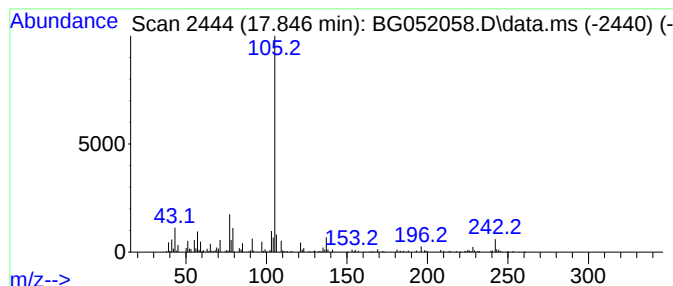
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 46 1-Phenylethanethiol, S-pent... Concentration Rank 36

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 17.846 | 16.87 ng/ul | 320024 | Phenanthrene-d10 | 17.640 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|------------|--------------|------|
| 1    |    |   | 1-Phenylethanethiol, S-pentafluo... | 284 | C11H9F5OS  | 1000469-38-3 | 64   |
| 2    |    |   | 1-Phenylethanethiol, S-heptaflu...  | 334 | C12H9F7OS  | 1000469-38-2 | 59   |
| 3    |    |   | Benzene, 1-(bromomethyl)-2-methyl-  | 184 | C8H9Br     | 000089-92-9  | 53   |
| 4    |    |   | Ethane, 1-(9-borabicyclo[3.3.1]n... | 242 | C16H23BO   | 1000160-80-6 | 50   |
| 5    |    |   | Etomidate                           | 244 | C14H16N2O2 | 033125-97-2  | 49   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011722\  
Data File : BG052058.D  
Acq On : 18 Jan 2022 18:42  
Operator : CG/JU  
Sample : N1100-15ME  
Misc :  
ALS Vial : 44 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLS8ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010622.M  
Quant Title : SVOA CALIBRATION

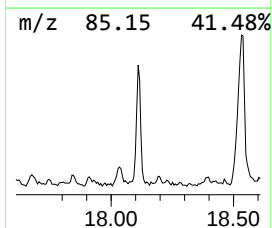
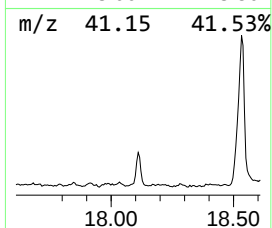
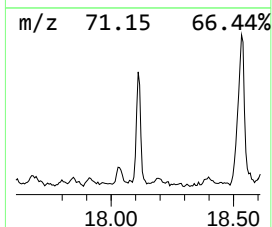
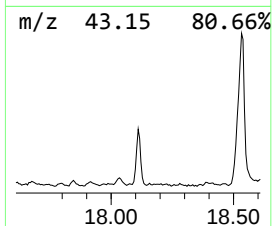
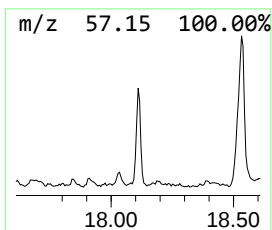
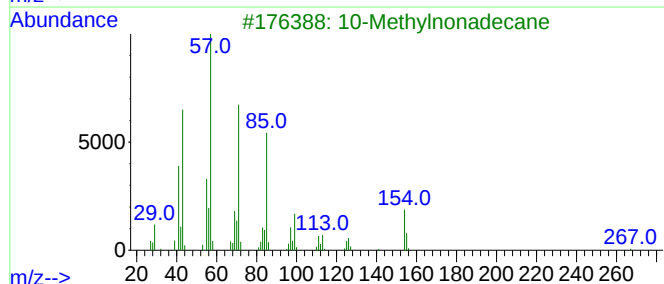
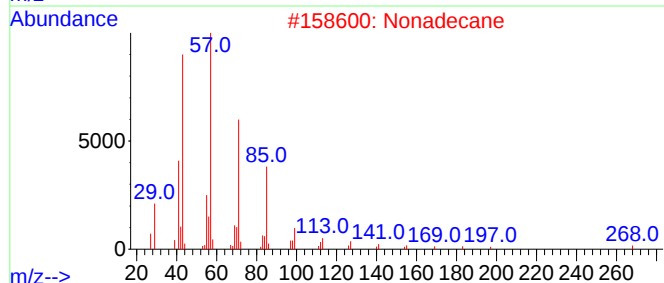
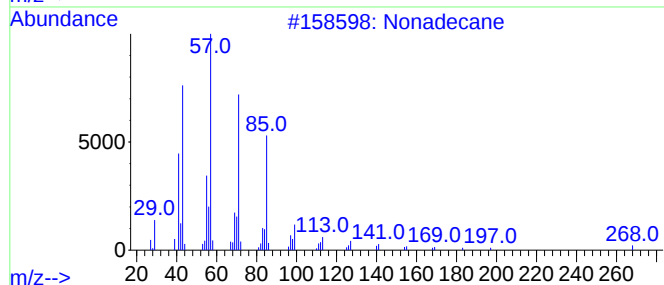
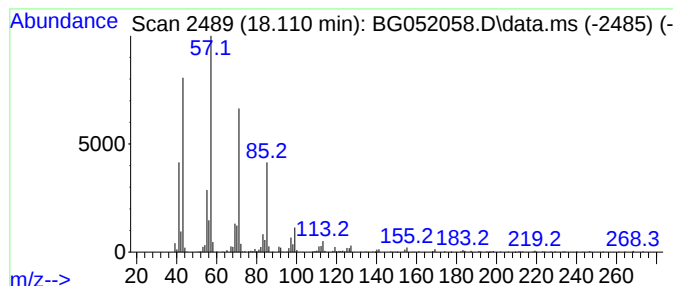
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 47 (DEL) Alkane: Straight-Chai... Concentration Rank 16

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 18.110 | 26.15 ng/ul | 496134 | Phenanthrene-d10 | 17.640 |

| Hit# | of | 5 | Tentative ID        | MW  | MolForm | CAS#        | Qual |
|------|----|---|---------------------|-----|---------|-------------|------|
| 1    |    |   | Nonadecane          | 268 | C19H40  | 000629-92-5 | 97   |
| 2    |    |   | Nonadecane          | 268 | C19H40  | 000629-92-5 | 91   |
| 3    |    |   | 10-Methylnonadecane | 282 | C20H42  | 056862-62-5 | 91   |
| 4    |    |   | Nonane, 5-butyl-    | 184 | C13H28  | 017312-63-9 | 90   |
| 5    |    |   | Hexadecane          | 226 | C16H34  | 000544-76-3 | 90   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011722\  
Data File : BG052058.D  
Acq On : 18 Jan 2022 18:42  
Operator : CG/JU  
Sample : N1100-15ME  
Misc :  
ALS Vial : 44 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLS8ME

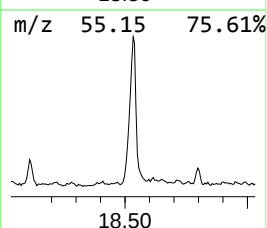
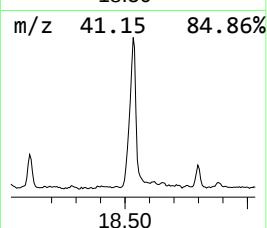
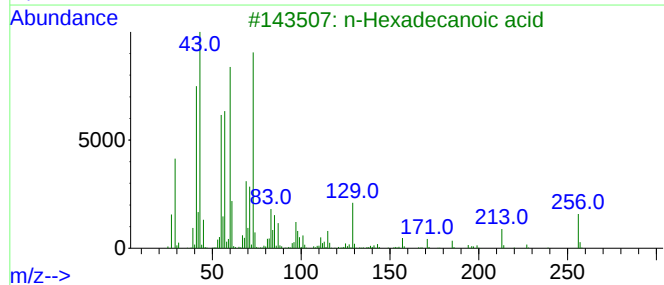
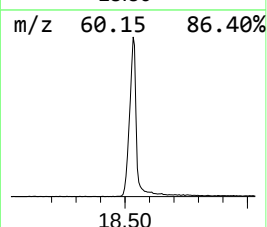
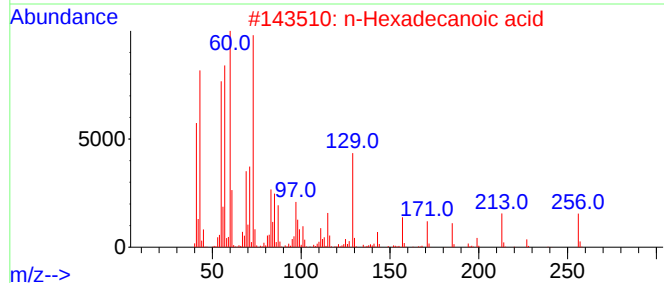
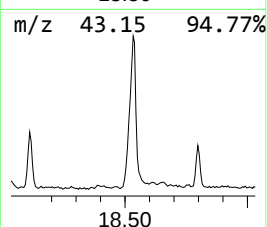
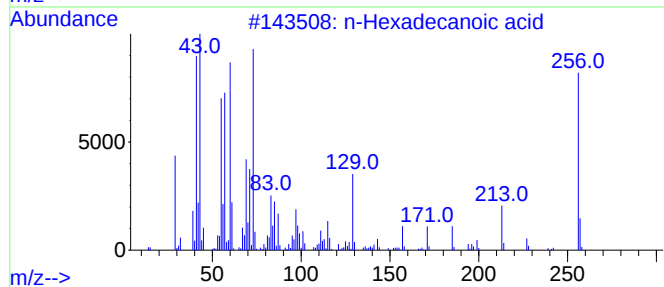
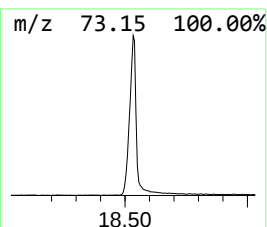
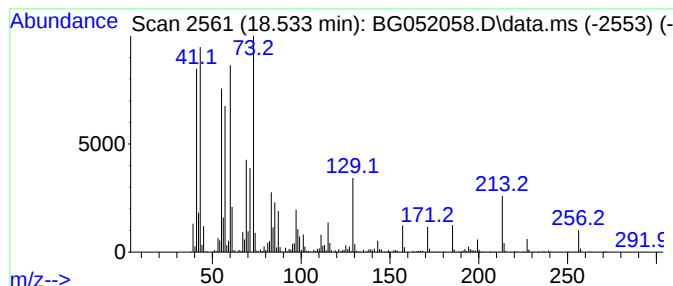
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Quant Title : SVOA CALIBRATION

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

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Peak Number 48 n-Hexadecanoic acid Concentration Rank 1

| R.T.   | EstConc      | Area    | Relative to ISTD | R.T.   |
|--------|--------------|---------|------------------|--------|
| 18.533 | 177.44 ng/ul | 3366370 | Phenanthrene-d10 | 17.640 |

| Hit# | of | 5 | Tentative ID        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|---------------------|-----|----------|-------------|------|
| 1    |    |   | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 98   |
| 2    |    |   | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 97   |
| 3    |    |   | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 95   |
| 4    |    |   | Tridecanoic acid    | 214 | C13H26O2 | 000638-53-9 | 95   |
| 5    |    |   | Tetradecanoic acid  | 228 | C14H28O2 | 000544-63-8 | 89   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011722\  
Data File : BG052058.D  
Acq On : 18 Jan 2022 18:42  
Operator : CG/JU  
Sample : N1100-15ME  
Misc :  
ALS Vial : 44 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLS8ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010622.M  
Quant Title : SVOA CALIBRATION

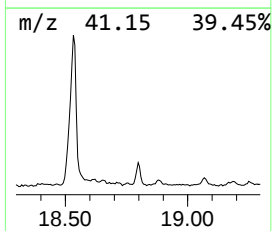
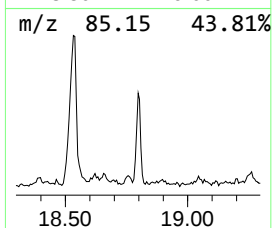
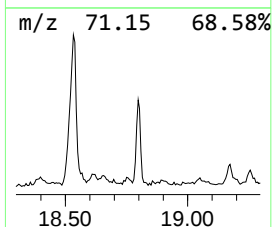
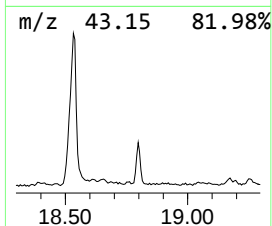
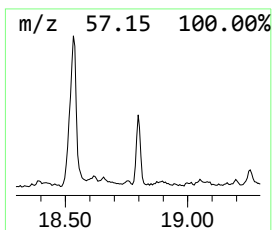
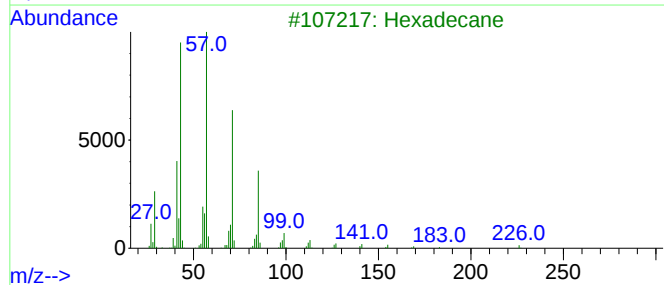
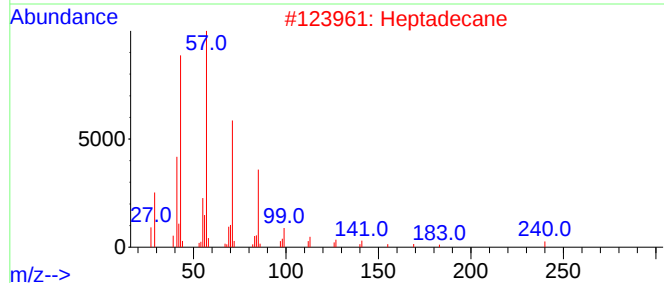
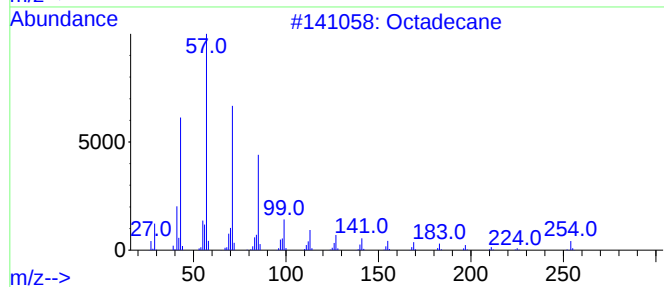
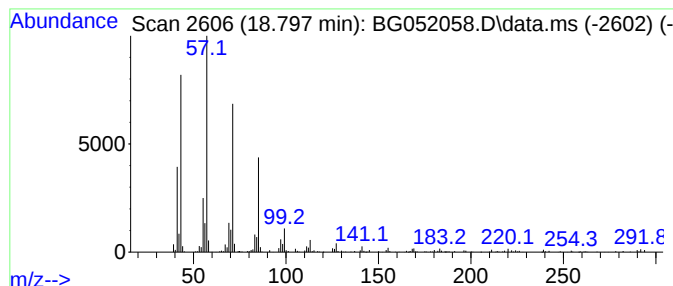
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 49 (DEL) Alkane: Straight-Chai... Concentration Rank 40

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 18.797 | 15.81 ng/ul | 299883 | Phenanthrene-d10 | 17.640 |

| Hit# | of 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|------|--------------|-----|---------|-------------|------|
| 1    |      | Octadecane   | 254 | C18H38  | 000593-45-3 | 97   |
| 2    |      | Heptadecane  | 240 | C17H36  | 000629-78-7 | 96   |
| 3    |      | Hexadecane   | 226 | C16H34  | 000544-76-3 | 95   |
| 4    |      | Hexadecane   | 226 | C16H34  | 000544-76-3 | 93   |
| 5    |      | Nonadecane   | 268 | C19H40  | 000629-92-5 | 91   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011722\  
Data File : BG052058.D  
Acq On : 18 Jan 2022 18:42  
Operator : CG/JU  
Sample : N1100-15ME  
Misc :  
ALS Vial : 44 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLS8ME

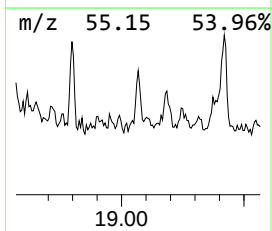
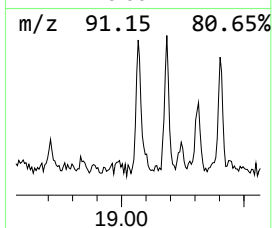
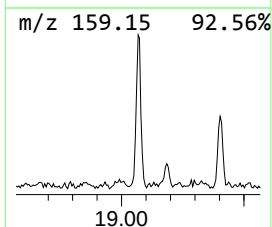
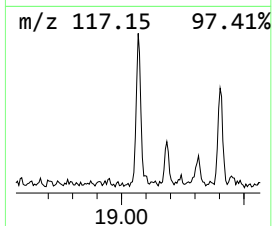
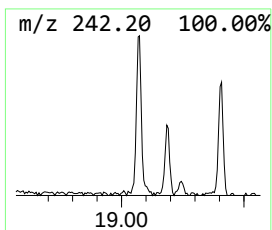
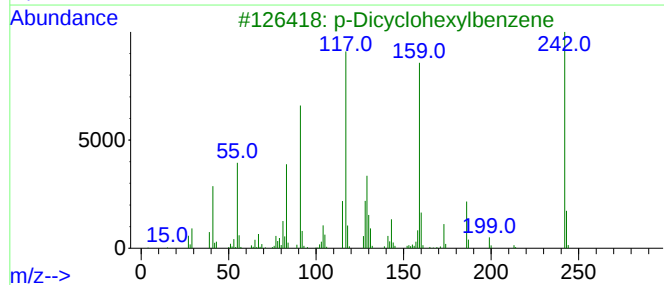
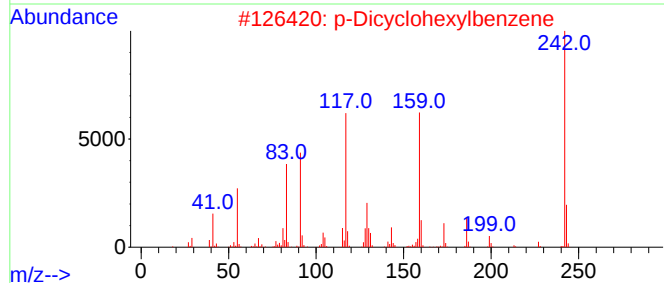
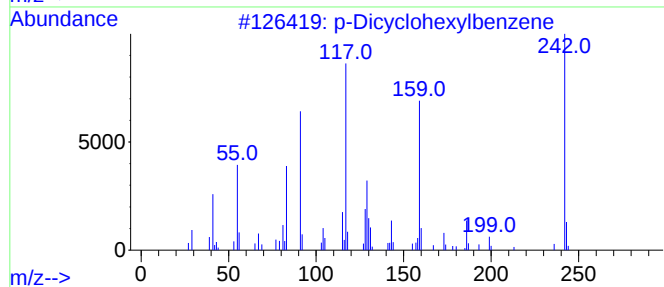
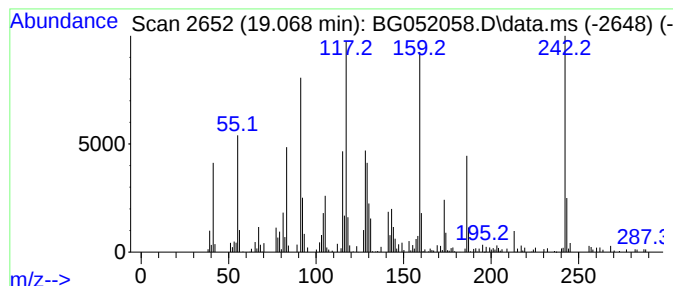
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Quant Title : SVOA CALIBRATION

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

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Peak Number 50 p-Dicyclohexylbenzene Concentration Rank 43

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 19.067 | 15.08 ng/ul | 286094 | Phenanthrene-d10 | 17.640 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | p-Dicyclohexylbenzene               | 242 | C18H26   | 001087-02-1  | 90   |
| 2    |    |   | p-Dicyclohexylbenzene               | 242 | C18H26   | 001087-02-1  | 89   |
| 3    |    |   | p-Dicyclohexylbenzene               | 242 | C18H26   | 001087-02-1  | 81   |
| 4    |    |   | Bicyclohexyl, 4-phenyl-             | 242 | C18H26   | 020273-27-2  | 68   |
| 5    |    |   | 2-Phenylpropionic acid, 4'-(2-me... | 218 | C14H18O2 | 1010454-94-3 | 43   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011722\  
Data File : BG052058.D  
Acq On : 18 Jan 2022 18:42  
Operator : CG/JU  
Sample : N1100-15ME  
Misc :  
ALS Vial : 44 Sample Multiplier: 1

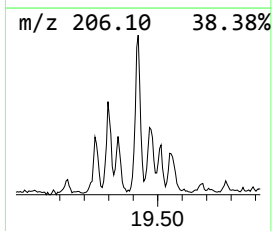
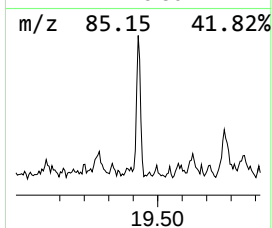
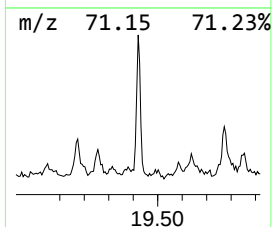
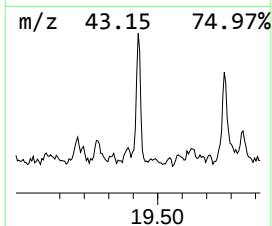
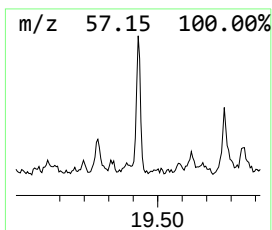
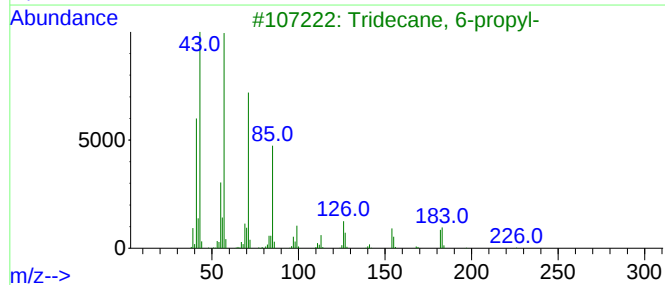
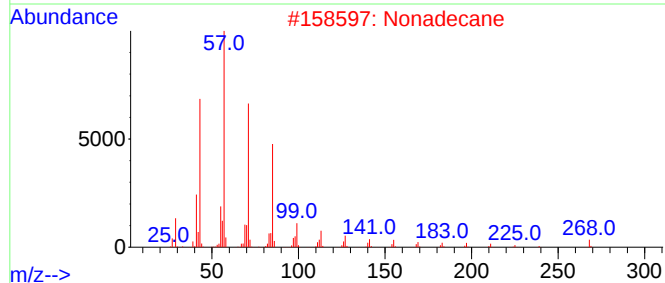
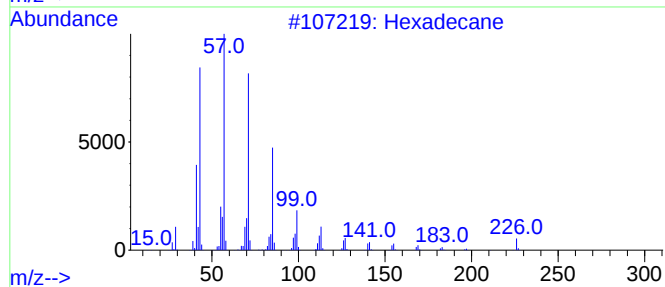
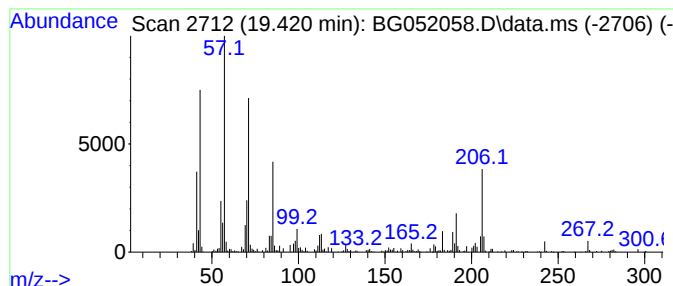
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BNA\_G  
ClientSampleId :  
BGLS8ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010622.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 51 (DEL) Alkane: Straight-Chai... Concentration Rank 17

| R.T.      | EstConc              | Area   | Relative to ISTD |         | R.T.        |      |
|-----------|----------------------|--------|------------------|---------|-------------|------|
| 19.420    | 25.97 ng/ul          | 492743 | Phenanthrene-d10 |         | 17.640      |      |
| Hit# of 5 | Tentative ID         |        | MW               | MolForm | CAS#        | Qual |
| 1         | Hexadecane           |        | 226              | C16H34  | 000544-76-3 | 90   |
| 2         | Nonadecane           |        | 268              | C19H40  | 000629-92-5 | 78   |
| 3         | Tridecane, 6-propyl- |        | 226              | C16H34  | 055045-10-8 | 78   |
| 4         | Pentadecane          |        | 212              | C15H32  | 000629-62-9 | 70   |
| 5         | Heneicosane          |        | 296              | C21H44  | 000629-94-7 | 64   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011722\  
Data File : BG052058.D  
Acq On : 18 Jan 2022 18:42  
Operator : CG/JU  
Sample : N1100-15ME  
Misc :  
ALS Vial : 44 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLS8ME

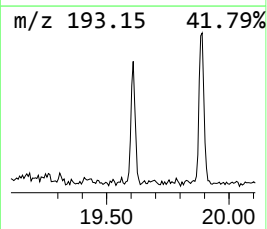
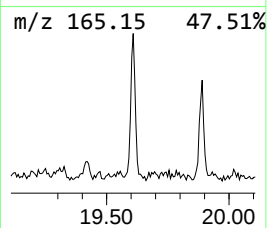
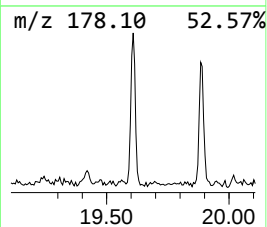
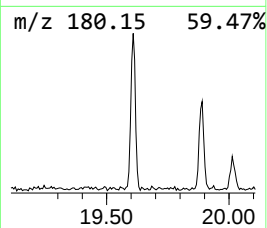
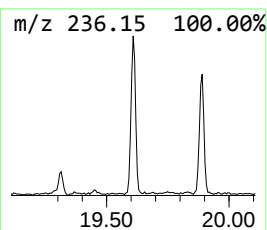
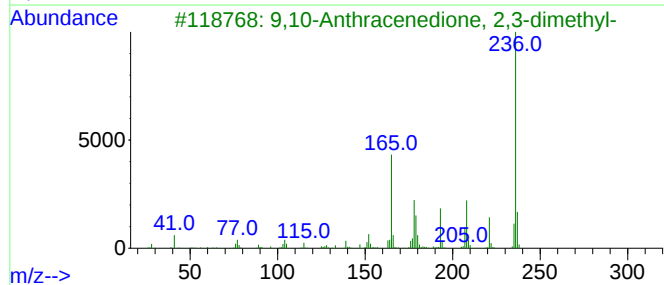
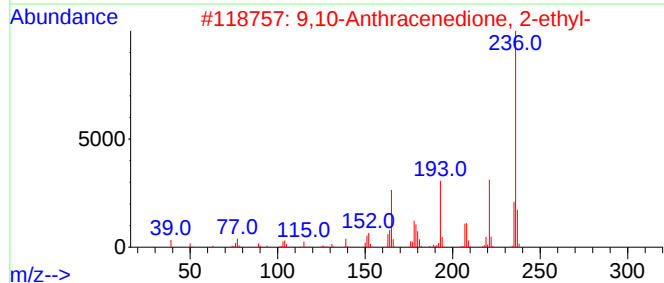
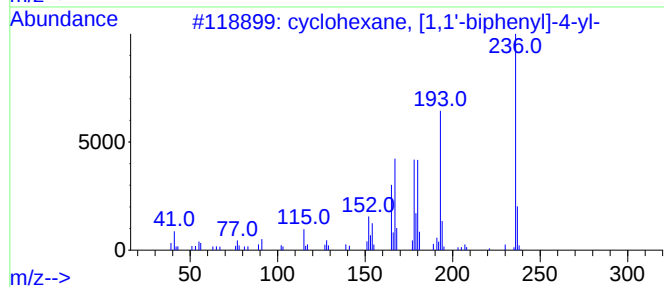
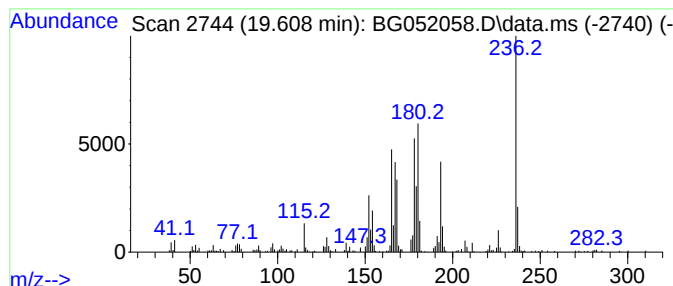
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Quant Title : SVOA CALIBRATION

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

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Peak Number 52 cyclohexane, [1,1'-biphenyl]... Concentration Rank 39

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 19.608 | 16.29 ng/ul | 309121 | Phenanthrene-d10 | 17.640 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | cyclohexane, [1,1'-biphenyl]-4-yl-  | 236 | C18H20    | 1000401-12-4 | 89   |
| 2    |    |   | 9,10-Anthracenedione, 2-ethyl-      | 236 | C16H12O2  | 000084-51-5  | 46   |
| 3    |    |   | 9,10-Anthracenedione, 2,3-dimethyl- | 236 | C16H12O2  | 006531-35-7  | 38   |
| 4    |    |   | 7-Methyl-4-phenylcoumarin           | 236 | C16H12O2  | 007758-71-6  | 27   |
| 5    |    |   | 6-Ethynyl-1,3-dimethylperimidin-... | 236 | C15H12N2O | 1000443-36-0 | 22   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011722\  
Data File : BG052058.D  
Acq On : 18 Jan 2022 18:42  
Operator : CG/JU  
Sample : N1100-15ME  
Misc :  
ALS Vial : 44 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLS8ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010622.M  
Quant Title : SVOA CALIBRATION

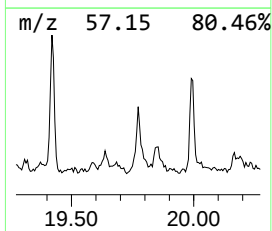
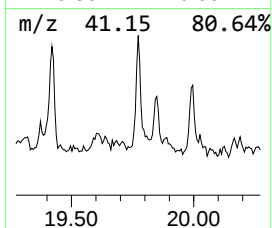
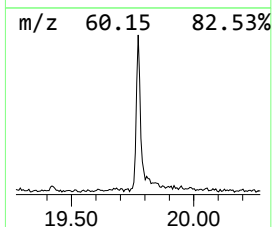
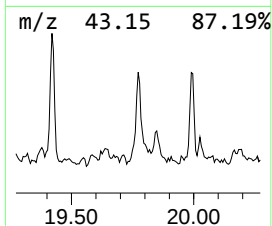
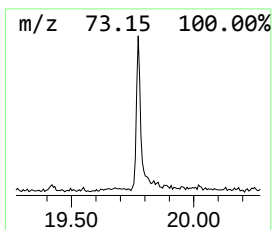
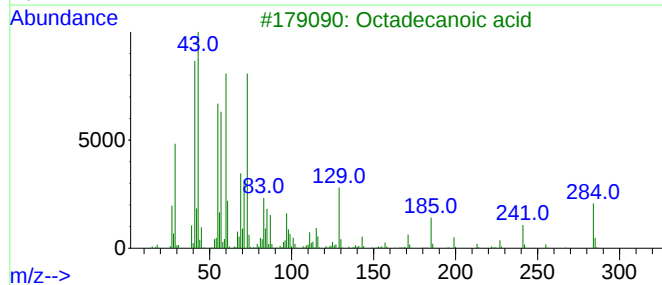
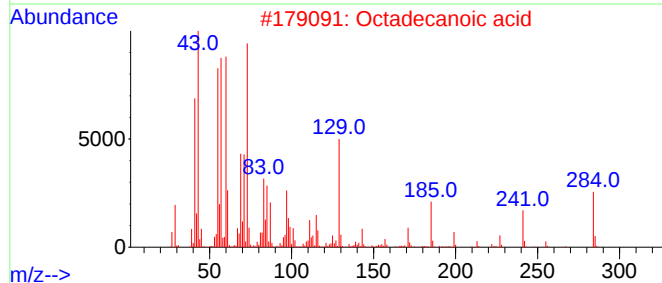
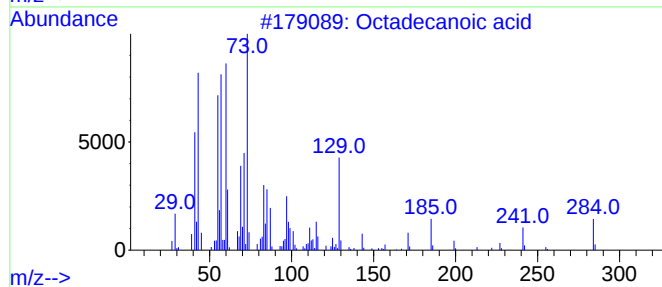
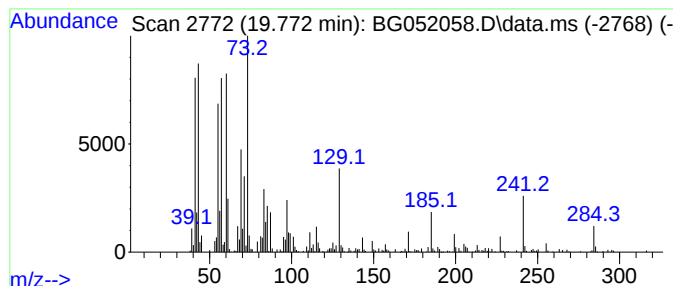
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 53 Octadecanoic acid Concentration Rank 30

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 19.773 | 18.48 ng/ul | 350539 | Phenanthrene-d10 | 17.640 |

| Hit# | of | 5 | Tentative ID       | MW  | MolForm  | CAS#        | Qual |
|------|----|---|--------------------|-----|----------|-------------|------|
| 1    |    |   | Octadecanoic acid  | 284 | C18H36O2 | 000057-11-4 | 99   |
| 2    |    |   | Octadecanoic acid  | 284 | C18H36O2 | 000057-11-4 | 99   |
| 3    |    |   | Octadecanoic acid  | 284 | C18H36O2 | 000057-11-4 | 94   |
| 4    |    |   | Octadecanoic acid  | 284 | C18H36O2 | 000057-11-4 | 93   |
| 5    |    |   | Pentadecanoic acid | 242 | C15H30O2 | 001002-84-2 | 93   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011722\  
Data File : BG052058.D  
Acq On : 18 Jan 2022 18:42  
Operator : CG/JU  
Sample : N1100-15ME  
Misc :  
ALS Vial : 44 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLS8ME

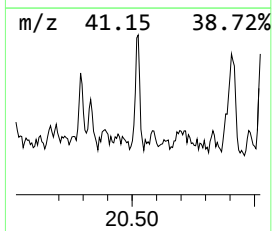
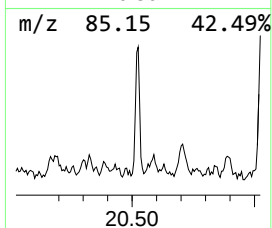
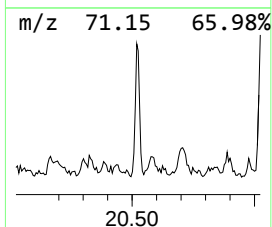
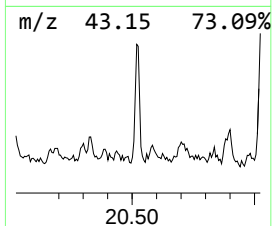
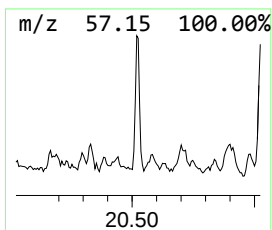
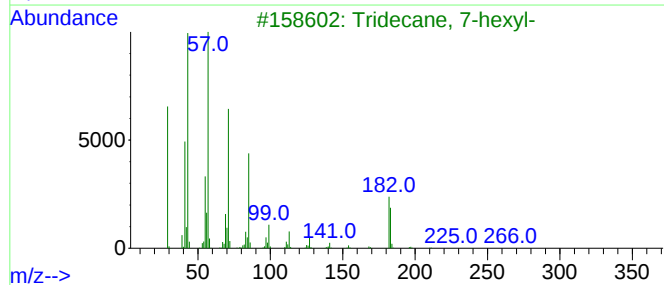
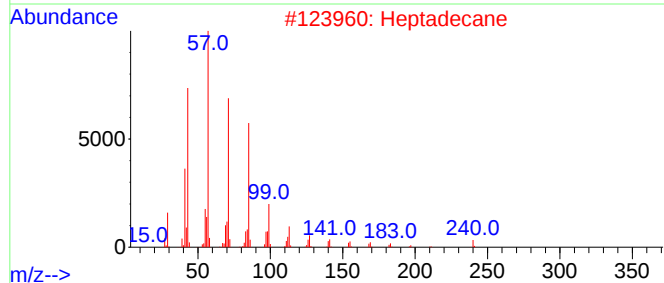
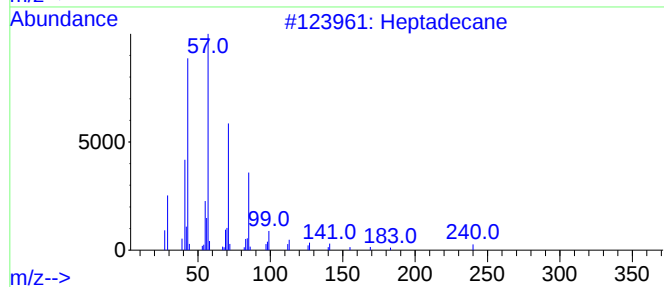
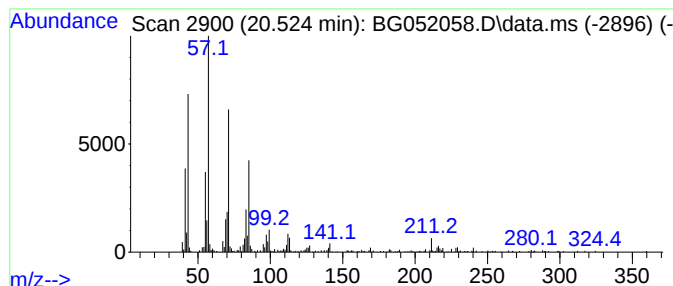
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010622.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 55 (DEL) Alkane: Straight-Chai... Concentration Rank 58

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 20.524 | 9.13 ng/ul | 296098 | Chrysene-d12     | 21.958 |

| Hit# | of | 5 | Tentative ID        | MW  | MolForm | CAS#        | Qual |
|------|----|---|---------------------|-----|---------|-------------|------|
| 1    |    |   | Heptadecane         | 240 | C17H36  | 000629-78-7 | 93   |
| 2    |    |   | Heptadecane         | 240 | C17H36  | 000629-78-7 | 92   |
| 3    |    |   | Tridecane, 7-hexyl- | 268 | C19H40  | 007225-66-3 | 89   |
| 4    |    |   | Octacosane          | 394 | C28H58  | 000630-02-4 | 80   |
| 5    |    |   | Tricosane           | 324 | C23H48  | 000638-67-5 | 76   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011722\  
Data File : BG052058.D  
Acq On : 18 Jan 2022 18:42  
Operator : CG/JU  
Sample : N1100-15ME  
Misc :  
ALS Vial : 44 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLS8ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010622.M  
Quant Title : SVOA CALIBRATION

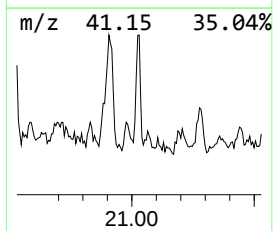
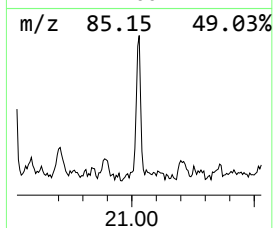
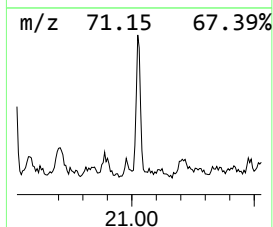
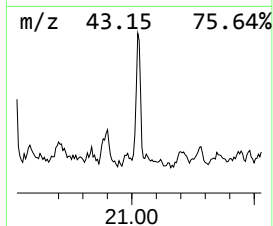
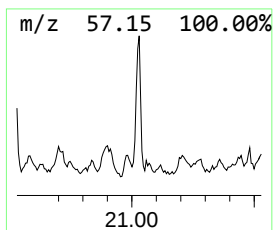
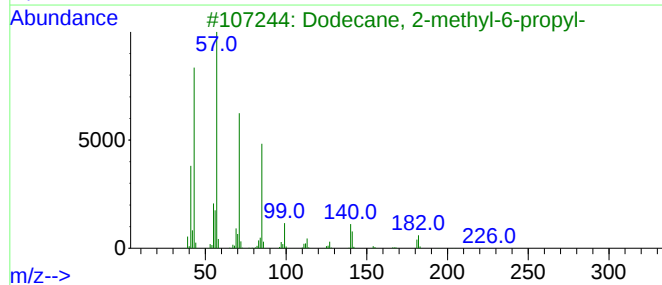
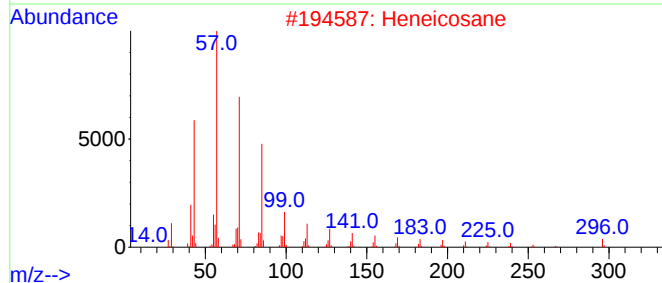
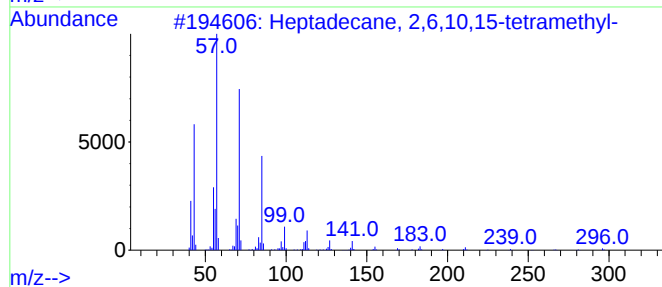
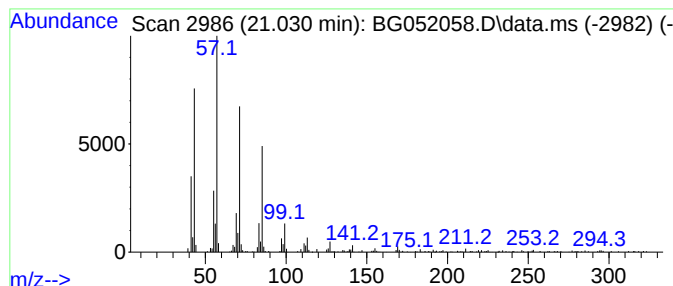
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 56 (DEL) Alkane: Straight-Chai... Concentration Rank 63

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 21.030 | 7.69 ng/ul | 249438 | Chrysene-d12     | 21.958 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Heptadecane, 2,6,10,15-tetramethyl- | 296 | C21H44  | 054833-48-6 | 97   |
| 2    |    |   | Heneicosane                         | 296 | C21H44  | 000629-94-7 | 94   |
| 3    |    |   | Dodecane, 2-methyl-6-propyl-        | 226 | C16H34  | 055045-08-4 | 94   |
| 4    |    |   | Heptadecane                         | 240 | C17H36  | 000629-78-7 | 93   |
| 5    |    |   | Heneicosane                         | 296 | C21H44  | 000629-94-7 | 91   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011722\  
Data File : BG052058.D  
Acq On : 18 Jan 2022 18:42  
Operator : CG/JU  
Sample : N1100-15ME  
Misc :  
ALS Vial : 44 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLS8ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010622.M  
Quant Title : SVOA CALIBRATION

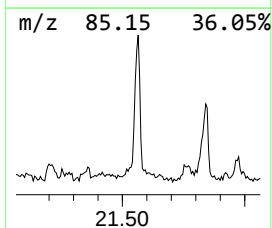
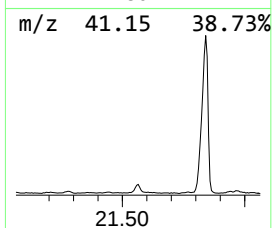
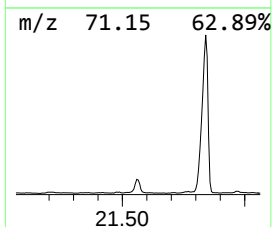
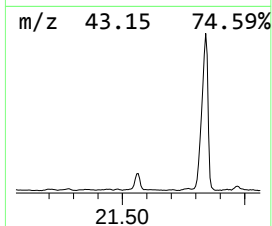
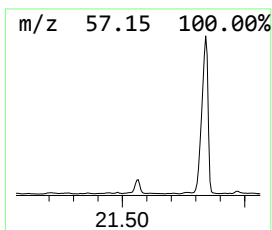
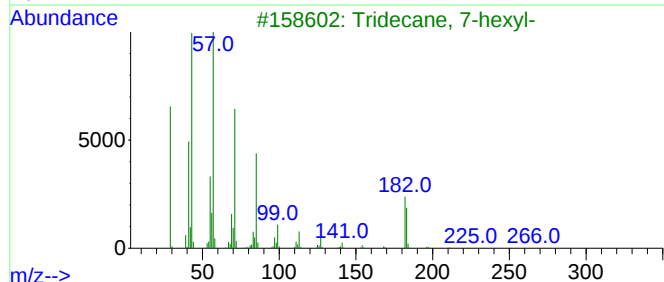
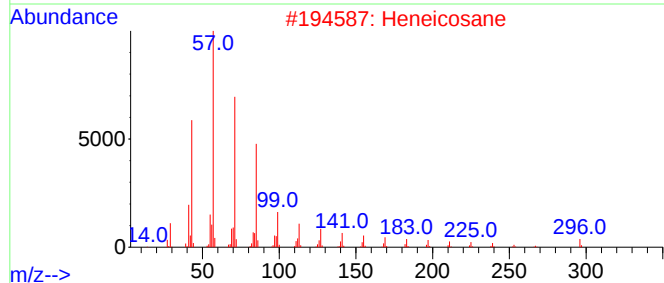
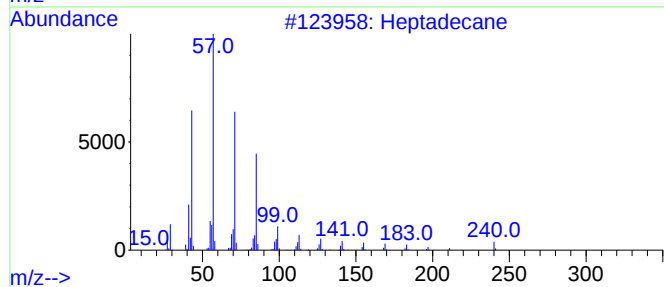
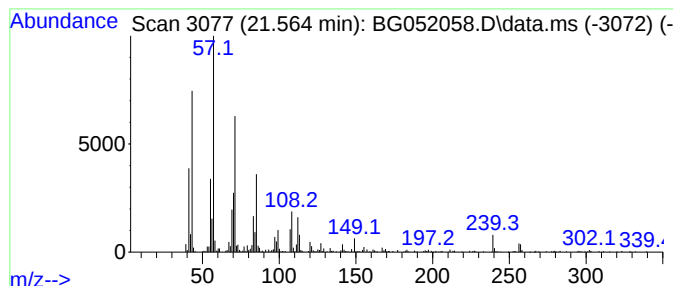
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 58 (DEL) Alkane: Straight-Chai... Concentration Rank 32

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 21.564 | 17.87 ng/ul | 579770 | Chrysene-d12     | 21.958 |

| Hit# | of | 5 | Tentative ID        | MW  | MolForm | CAS#        | Qual |
|------|----|---|---------------------|-----|---------|-------------|------|
| 1    |    |   | Heptadecane         | 240 | C17H36  | 000629-78-7 | 86   |
| 2    |    |   | Heneicosane         | 296 | C21H44  | 000629-94-7 | 86   |
| 3    |    |   | Tridecane, 7-hexyl- | 268 | C19H40  | 007225-66-3 | 83   |
| 4    |    |   | Octadecane          | 254 | C18H38  | 000593-45-3 | 78   |
| 5    |    |   | Octadecane          | 254 | C18H38  | 000593-45-3 | 70   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011722\  
Data File : BG052058.D  
Acq On : 18 Jan 2022 18:42  
Operator : CG/JU  
Sample : N1100-15ME  
Misc :  
ALS Vial : 44 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLS8ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010622.M  
Quant Title : SVOA CALIBRATION

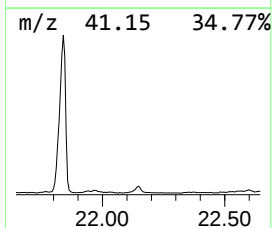
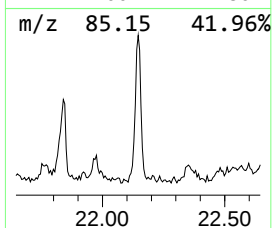
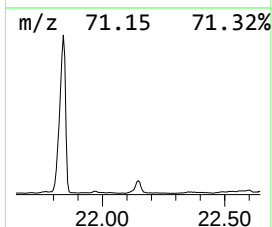
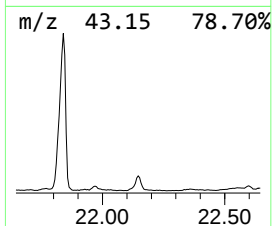
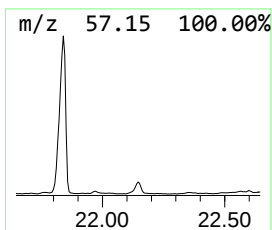
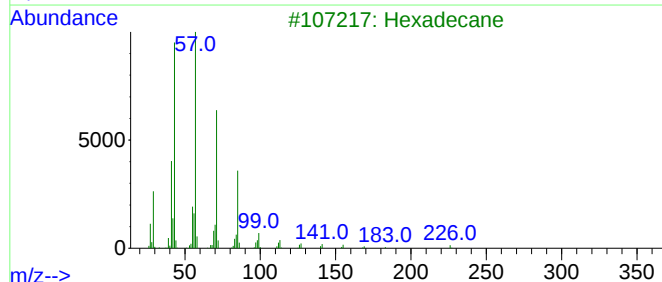
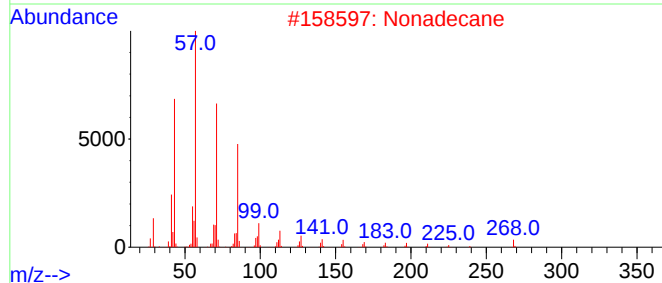
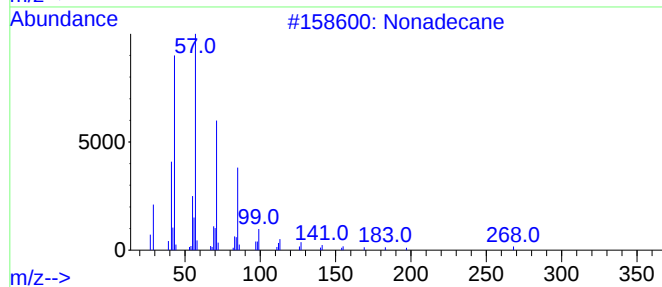
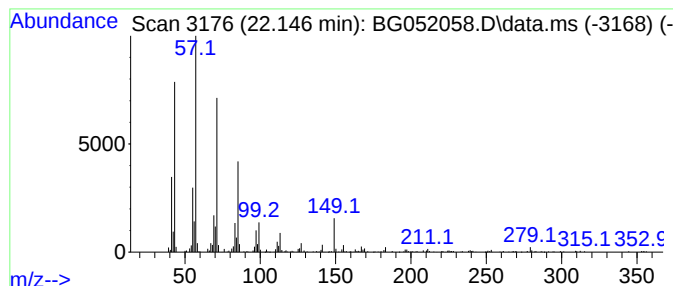
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 59 (DEL) Alkane: Straight-Chai... Concentration Rank 28

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 22.146 | 19.04 ng/ul | 617706 | Chrysene-d12     | 21.958 |

| Hit# | of | 5 | Tentative ID       | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------|-----|---------|-------------|------|
| 1    |    |   | Nonadecane         | 268 | C19H40  | 000629-92-5 | 95   |
| 2    |    |   | Nonadecane         | 268 | C19H40  | 000629-92-5 | 94   |
| 3    |    |   | Hexadecane         | 226 | C16H34  | 000544-76-3 | 93   |
| 4    |    |   | Heptadecane        | 240 | C17H36  | 000629-78-7 | 93   |
| 5    |    |   | Tridecane, 1-iodo- | 310 | C13H27I | 035599-77-0 | 93   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011722\  
Data File : BG052058.D  
Acq On : 18 Jan 2022 18:42  
Operator : CG/JU  
Sample : N1100-15ME  
Misc :  
ALS Vial : 44 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLS8ME

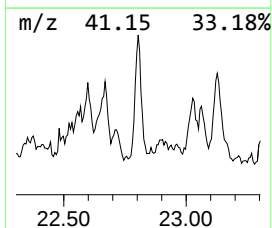
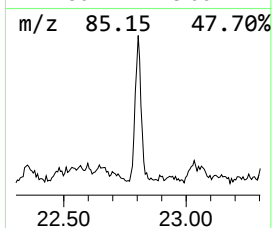
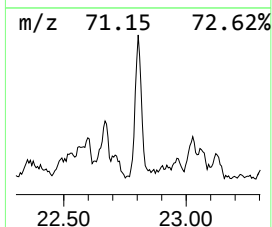
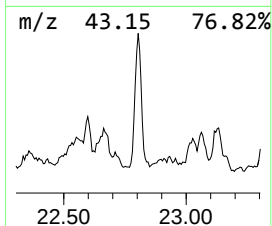
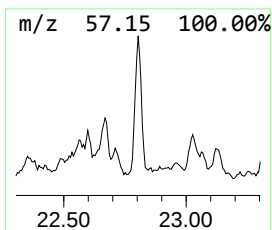
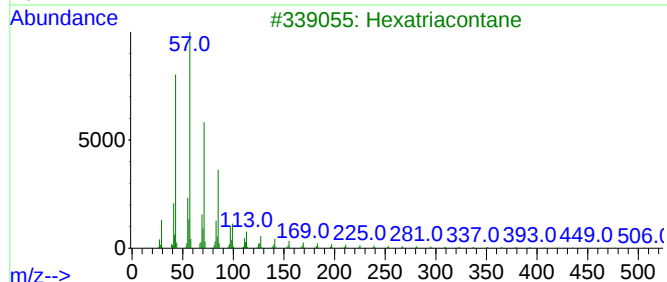
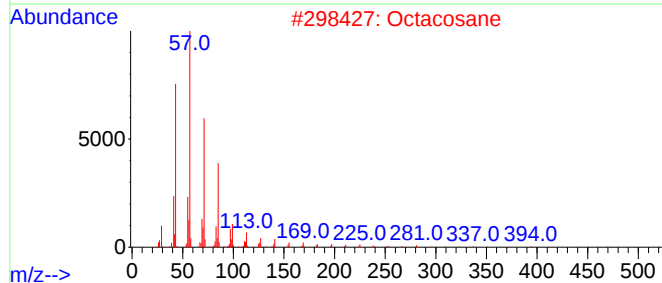
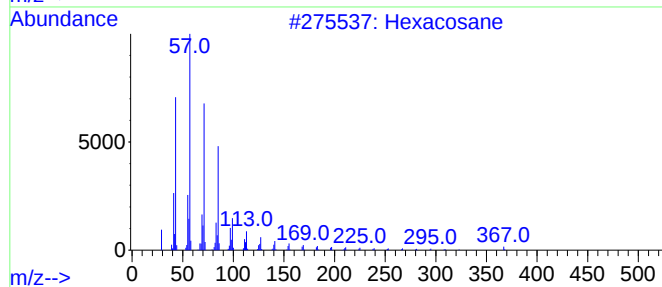
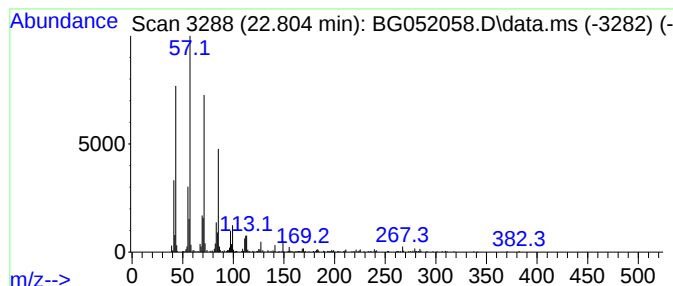
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Quant Title : SVOA CALIBRATION

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

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Peak Number 61 (DEL) Alkane: Straight-Chai... Concentration Rank 41

| R.T.      | EstConc         | Area   | Relative to ISTD | R.T.        |      |
|-----------|-----------------|--------|------------------|-------------|------|
| 22.804    | 15.55 ng/ul     | 504357 | Chrysene-d12     | 21.958      |      |
| Hit# of 5 | Tentative ID    | MW     | MolForm          | CAS#        | Qual |
| 1         | Hexacosane      | 366    | C26H54           | 000630-01-3 | 90   |
| 2         | Octacosane      | 394    | C28H58           | 000630-02-4 | 90   |
| 3         | Hexatriacontane | 507    | C36H74           | 000630-06-8 | 87   |
| 4         | Heptadecane     | 240    | C17H36           | 000629-78-7 | 87   |
| 5         | Tetracosane     | 338    | C24H50           | 000646-31-1 | 87   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011722\  
Data File : BG052058.D  
Acq On : 18 Jan 2022 18:42  
Operator : CG/JU  
Sample : N1100-15ME  
Misc :  
ALS Vial : 44 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLS8ME

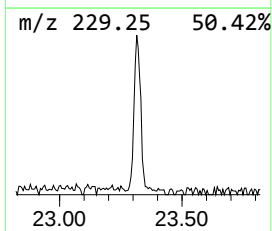
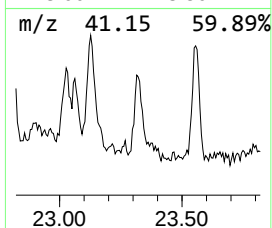
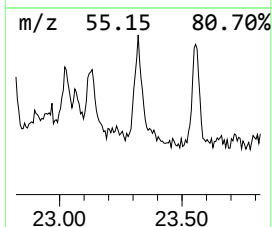
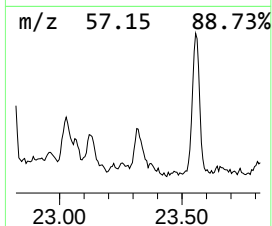
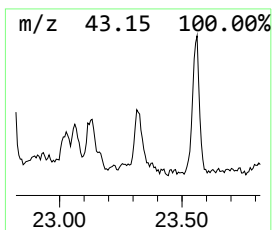
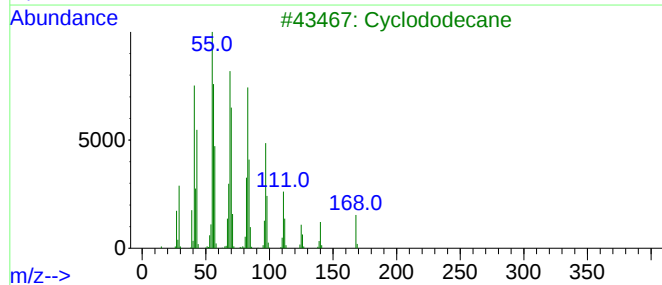
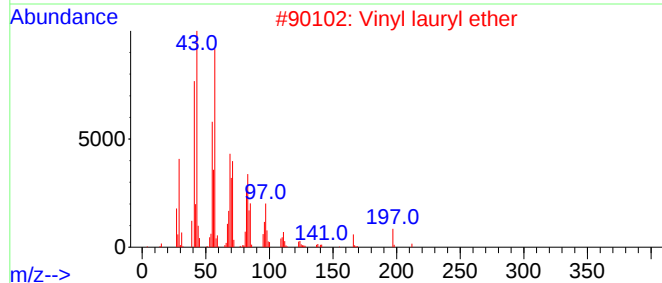
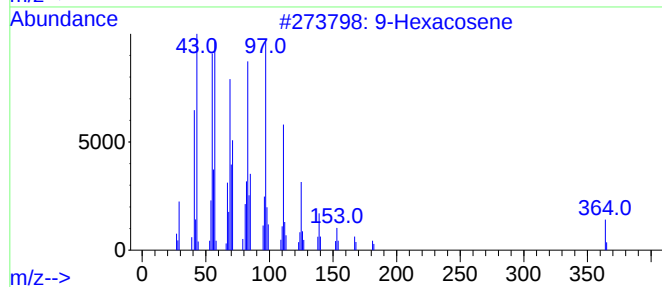
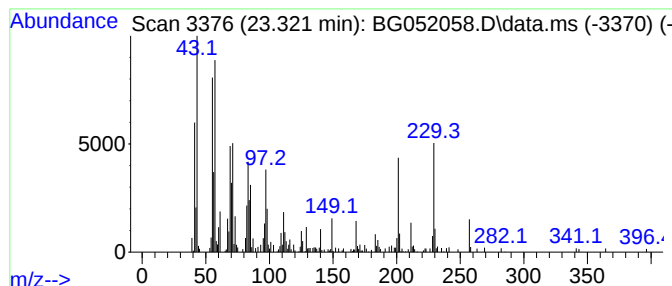
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Quant Title : SVOA CALIBRATION

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

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Peak Number 64 (DEL) Alkane: Straight-Chai... Concentration Rank 57

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 23.321 | 9.41 ng/ul | 305379 | Chrysene-d12     | 21.958 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm     | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-------------|--------------|------|
| 1    |    |   | 9-Hexacosene                        | 364 | C26H52      | 071502-22-2  | 64   |
| 2    |    |   | Vinyl lauryl ether                  | 212 | C14H28O     | 000765-14-0  | 64   |
| 3    |    |   | Cyclododecane                       | 168 | C12H24      | 000294-62-2  | 60   |
| 4    |    |   | Butyl dodecyl ether                 | 242 | C16H34O     | 1000406-40-3 | 53   |
| 5    |    |   | Pentadecafluorooctanoic acid, oc... | 666 | C26H37F15O2 | 1000406-04-8 | 49   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011722\  
Data File : BG052058.D  
Acq On : 18 Jan 2022 18:42  
Operator : CG/JU  
Sample : N1100-15ME  
Misc :  
ALS Vial : 44 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLS8ME

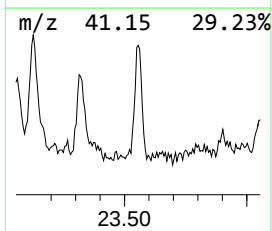
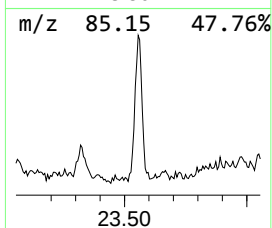
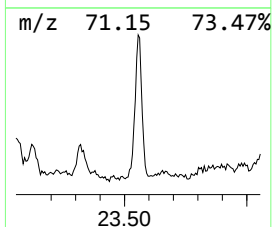
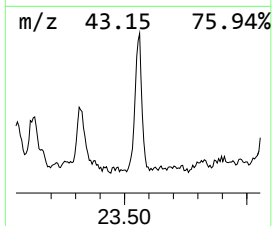
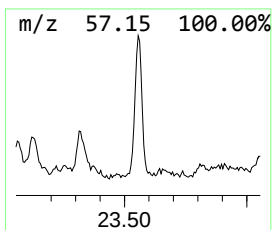
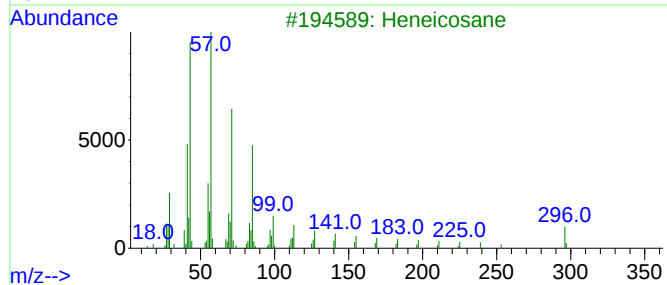
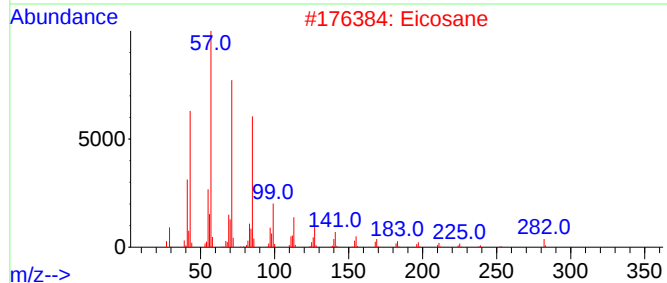
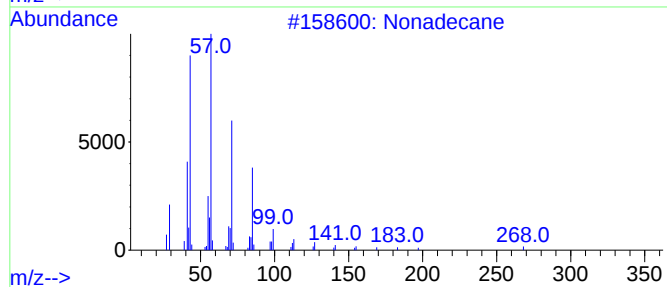
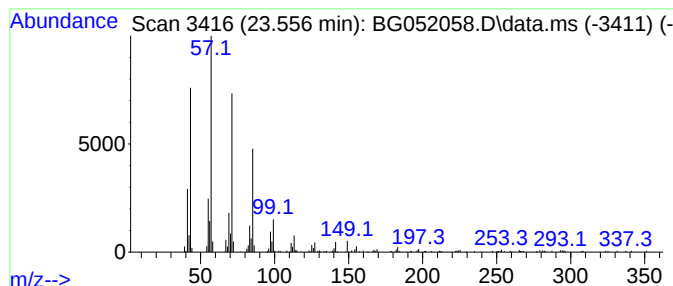
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Quant Title : SVOA CALIBRATION

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

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Peak Number 65 (DEL) Alkane: Straight-Chai... Concentration Rank 50

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 23.556 | 12.19 ng/ul | 395339 | Chrysene-d12     | 21.958 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Nonadecane                          | 268 | C19H40  | 000629-92-5 | 95   |
| 2    |    |   | Eicosane                            | 282 | C20H42  | 000112-95-8 | 95   |
| 3    |    |   | Heneicosane                         | 296 | C21H44  | 000629-94-7 | 93   |
| 4    |    |   | Heptadecane, 2,6,10,15-tetramethyl- | 296 | C21H44  | 054833-48-6 | 93   |
| 5    |    |   | Docosane, 11-butyl-                 | 366 | C26H54  | 013475-76-8 | 91   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011722\  
Data File : BG052058.D  
Acq On : 18 Jan 2022 18:42  
Operator : CG/JU  
Sample : N1100-15ME  
Misc :  
ALS Vial : 44 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLS8ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010622.M  
Quant Title : SVOA CALIBRATION

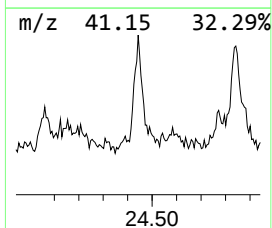
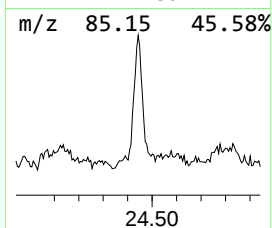
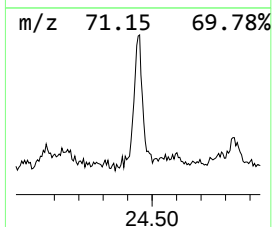
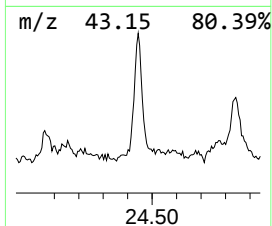
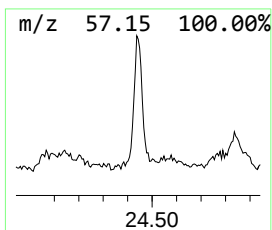
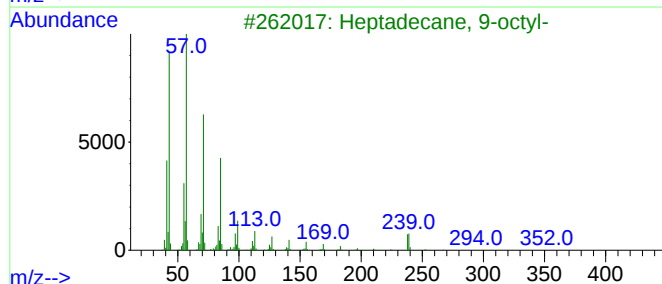
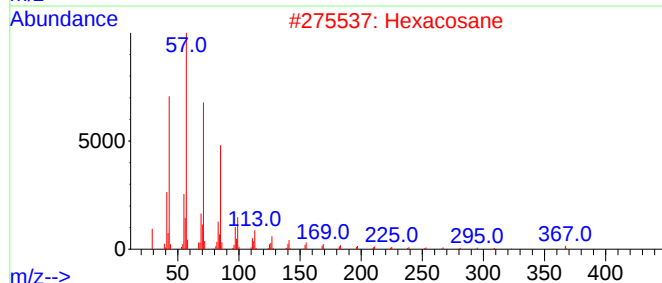
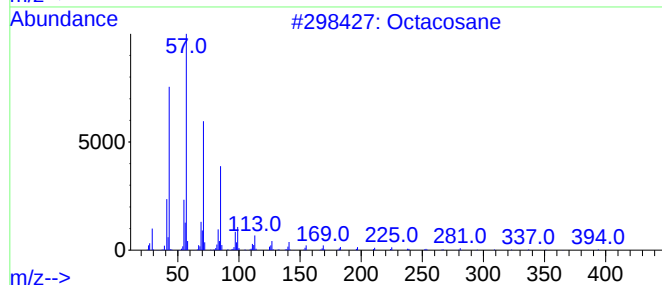
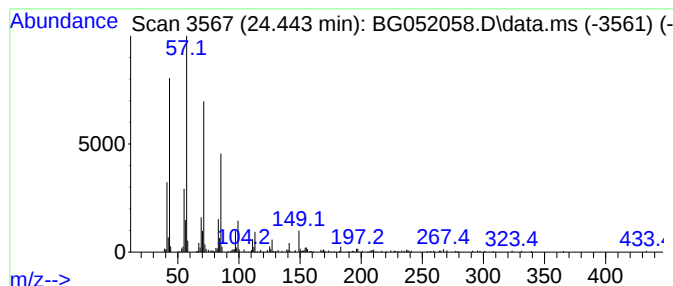
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 66 (DEL) Alkane: Straight-Chai... Concentration Rank 56

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 24.443 | 9.53 ng/ul | 365247 | Perylene-d12     | 25.447 |

| Hit# | of | 5 | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------------|-----|---------|-------------|------|
| 1    |    |   | Octacosane             | 394 | C28H58  | 000630-02-4 | 87   |
| 2    |    |   | Hexacosane             | 366 | C26H54  | 000630-01-3 | 87   |
| 3    |    |   | Heptadecane, 9-octyl-  | 352 | C25H52  | 007225-64-1 | 83   |
| 4    |    |   | Tetracosane            | 338 | C24H50  | 000646-31-1 | 83   |
| 5    |    |   | Heneicosane, 11-decyl- | 437 | C31H64  | 055320-06-4 | 83   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011722\  
Data File : BG052058.D  
Acq On : 18 Jan 2022 18:42  
Operator : CG/JU  
Sample : N1100-15ME  
Misc :  
ALS Vial : 44 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLS8ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010622.M  
Quant Title : SVOA CALIBRATION

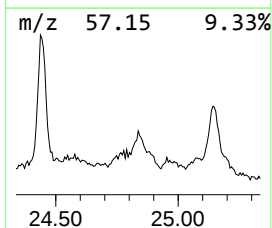
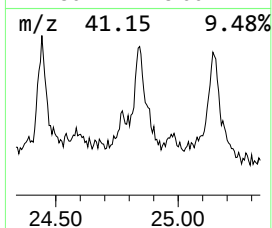
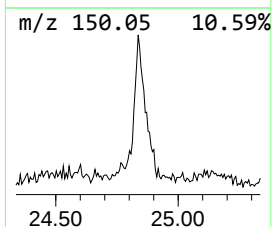
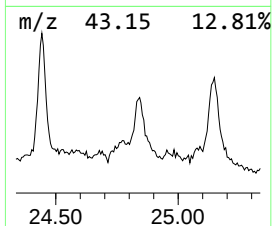
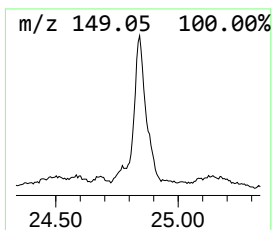
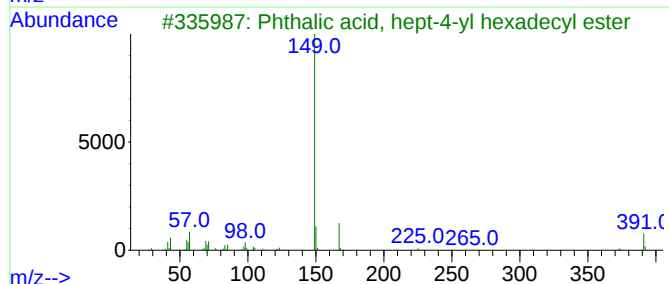
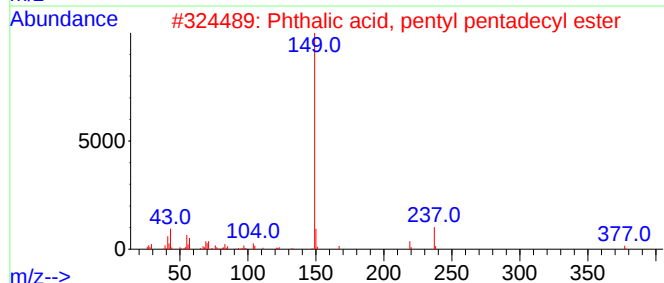
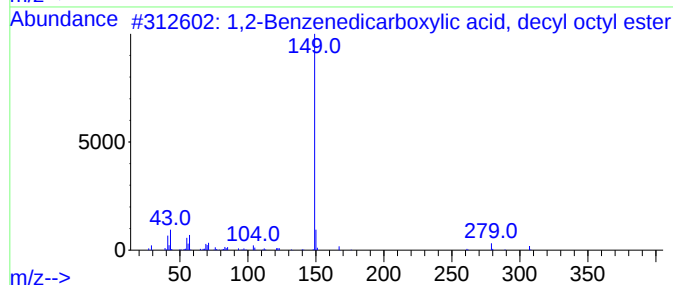
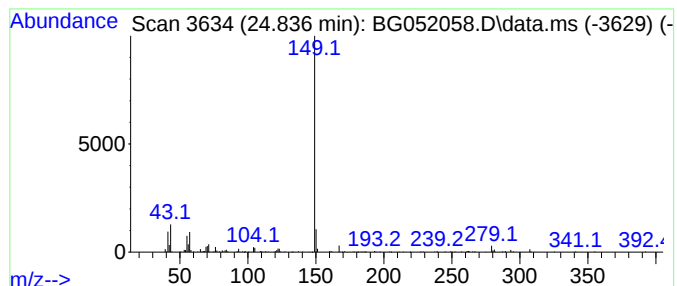
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 67 1,2-Benzenedicarboxylic aci... Concentration Rank 42

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 24.836 | 15.15 ng/ul | 580794 | Perylene-d12     | 25.447 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | 1,2-Benzenedicarboxylic acid, de... | 418 | C26H42O4 | 000119-07-3  | 83   |
| 2    |    |   | Phthalic acid, pentyl pentadecyl... | 446 | C28H46O4 | 1000308-92-9 | 72   |
| 3    |    |   | Phthalic acid, hept-4-yl hexadec... | 488 | C31H52O4 | 1000356-79-7 | 72   |
| 4    |    |   | Didodecyl phthalate                 | 502 | C32H54O4 | 002432-90-8  | 72   |
| 5    |    |   | Di-isononyl phthlate                | 418 | C26H42O4 | 020548-62-3  | 72   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011722\  
Data File : BG052058.D  
Acq On : 18 Jan 2022 18:42  
Operator : CG/JU  
Sample : N1100-15ME  
Misc :  
ALS Vial : 44 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLS8ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010622.M  
Quant Title : SVOA CALIBRATION

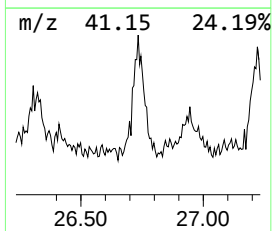
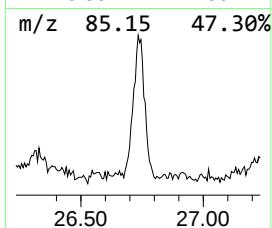
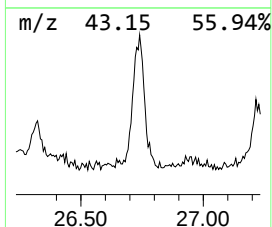
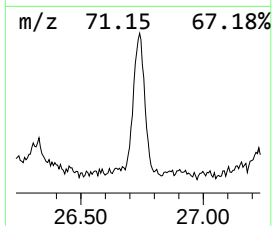
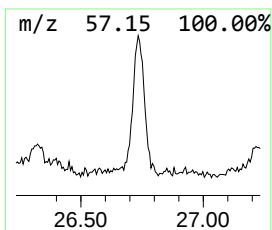
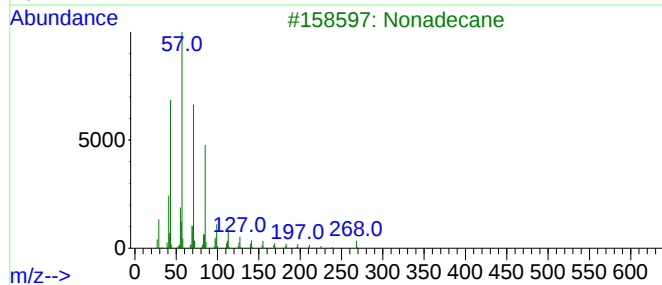
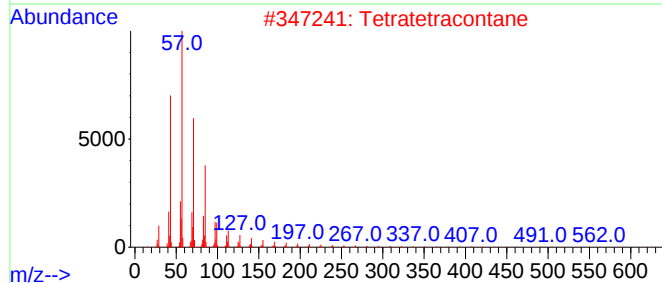
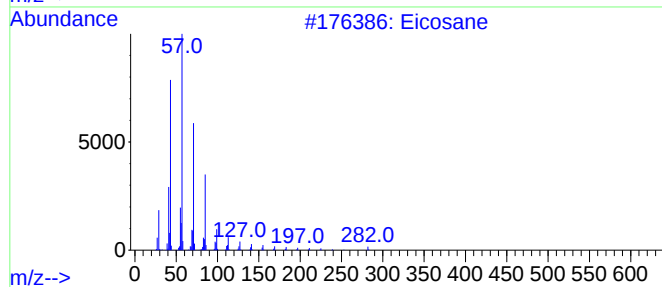
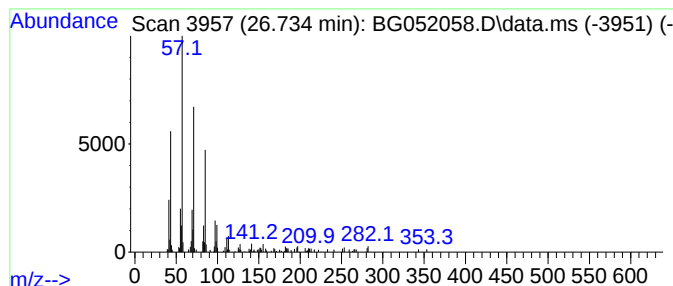
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 69 (DEL) Alkane: Straight-Chai... Concentration Rank 61

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 26.734 | 8.14 ng/ul | 312162 | Perylene-d12     | 25.447 |

| Hit# | of | 5 | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------|-----|---------|-------------|------|
| 1    |    |   | Eicosane              | 282 | C20H42  | 000112-95-8 | 91   |
| 2    |    |   | Tetratetracontane     | 619 | C44H90  | 007098-22-8 | 90   |
| 3    |    |   | Nonadecane            | 268 | C19H40  | 000629-92-5 | 87   |
| 4    |    |   | Octacosane            | 394 | C28H58  | 000630-02-4 | 87   |
| 5    |    |   | Octacosane, 2-methyl- | 408 | C29H60  | 001560-98-1 | 87   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011722\  
Data File : BG052058.D  
Acq On : 18 Jan 2022 18:42  
Operator : CG/JU  
Sample : N1100-15ME  
Misc :  
ALS Vial : 44 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLS8ME

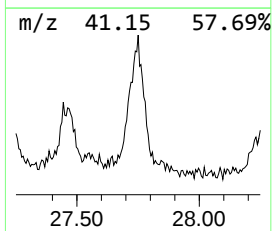
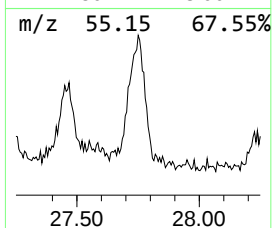
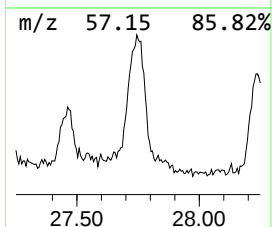
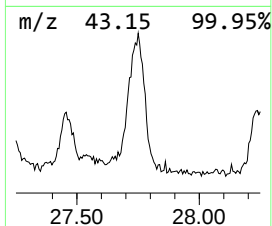
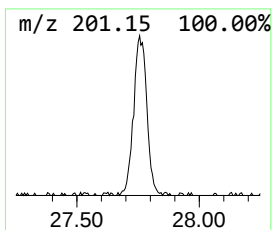
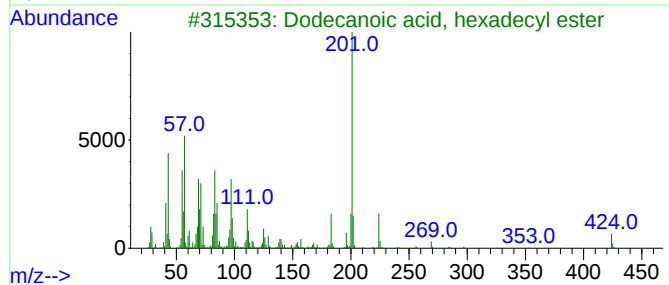
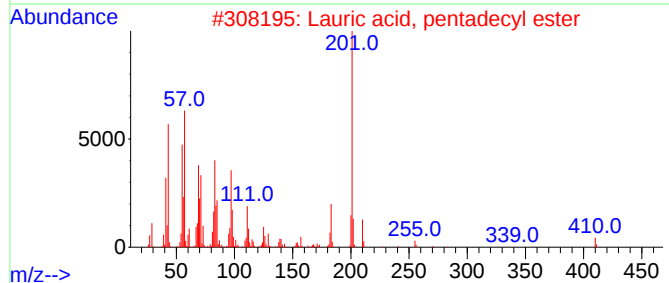
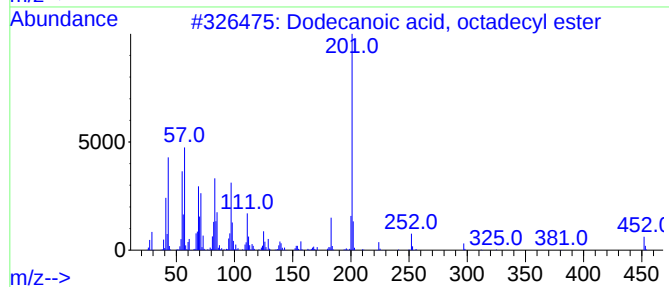
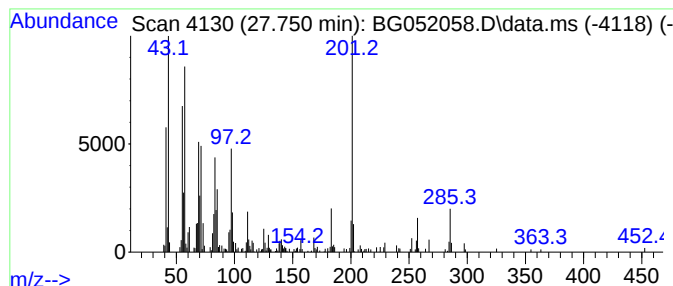
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010622.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 71 Dodecanoic acid, octadecyl ... Concentration Rank 29

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 27.750 | 18.77 ng/ul | 719512 | Perylene-d12     | 25.447 |

| Hit# | of | 5 | Tentative ID                      | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-----------------------------------|-----|----------|--------------|------|
| 1    |    |   | Dodecanoic acid, octadecyl ester  | 452 | C30H60O2 | 003234-84-2  | 97   |
| 2    |    |   | Lauric acid, pentadecyl ester     | 410 | C27H54O2 | 1000438-93-8 | 90   |
| 3    |    |   | Dodecanoic acid, hexadecyl ester  | 424 | C28H56O2 | 020834-06-4  | 90   |
| 4    |    |   | Dodecanoic acid, hexadecyl ester  | 424 | C28H56O2 | 020834-06-4  | 81   |
| 5    |    |   | Dodecanoic acid, tetradecyl ester | 396 | C26H52O2 | 022412-97-1  | 81   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011722\  
Data File : BG052058.D  
Acq On : 18 Jan 2022 18:42  
Operator : CG/JU  
Sample : N1100-15ME  
Misc :  
ALS Vial : 44 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLS8ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010622.M  
Quant Title : SVOA CALIBRATION

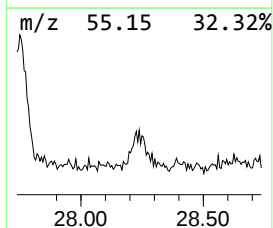
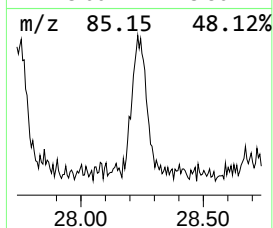
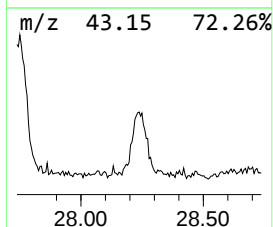
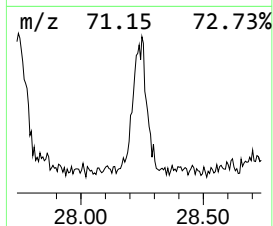
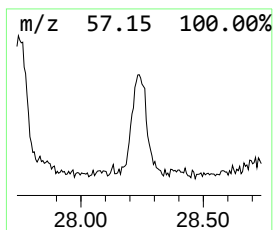
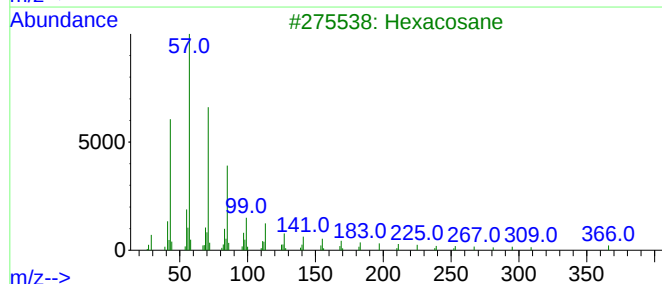
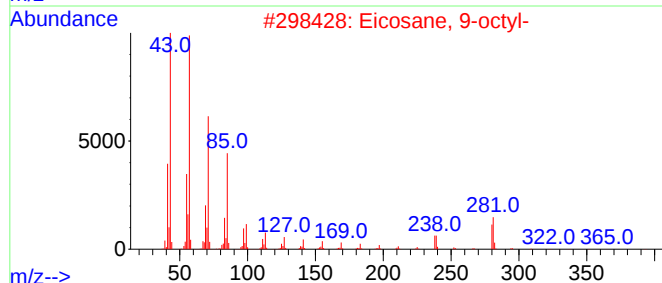
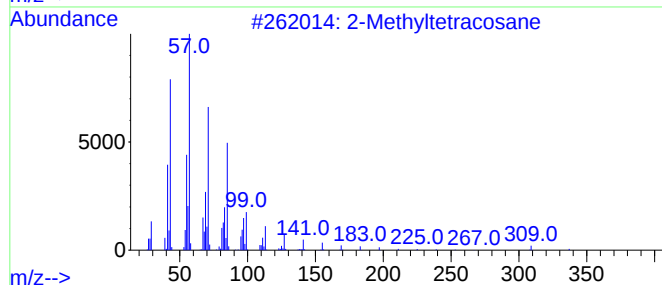
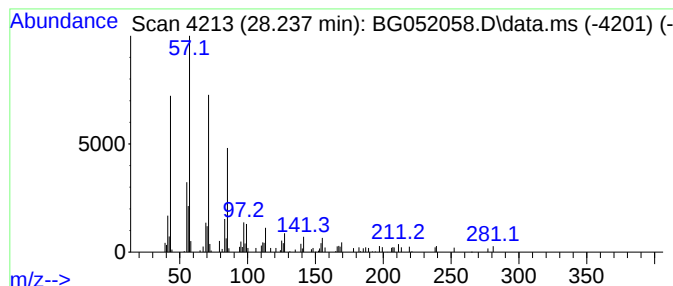
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 72 (DEL) Alkane: Straight-Chai... Concentration Rank 72

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 28.238 | 5.48 ng/ul | 210015 | Perylene-d12     | 25.447 |

| Hit# | of | 5 | Tentative ID        | MW  | MolForm | CAS#        | Qual |
|------|----|---|---------------------|-----|---------|-------------|------|
| 1    |    |   | 2-Methyltetracosane | 352 | C25H52  | 001560-78-7 | 91   |
| 2    |    |   | Eicosane, 9-octyl-  | 394 | C28H58  | 013475-77-9 | 91   |
| 3    |    |   | Hexacosane          | 366 | C26H54  | 000630-01-3 | 87   |
| 4    |    |   | Heptacosane         | 380 | C27H56  | 000593-49-7 | 87   |
| 5    |    |   | Hexacosane          | 366 | C26H54  | 000630-01-3 | 87   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG011722\  
 Data File : BG052058.D  
 Acq On : 18 Jan 2022 18:42  
 Operator : CG/JU  
 Sample : N1100-15ME  
 Misc :  
 ALS Vial : 44 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 BGLS8ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG010622.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 |
| Ethylbenzene       | 5.709  | 42.1    | ng/ul | 948505   | 1                     | 8.247  | 450887  | 20.0 |
| p-Xylene           | 5.850  | 135.6   | ng/ul | 3056380  | 1                     | 8.247  | 450887  | 20.0 |
| Benzene, 1,3-di... | 6.226  | 54.5    | ng/ul | 1229880  | 1                     | 8.247  | 450887  | 20.0 |
| Benzene, (1-met... | 6.708  | 40.0    | ng/ul | 902880   | 1                     | 8.247  | 450887  | 20.0 |
| (DEL) Alkane: C... | 6.866  | 10.5    | ng/ul | 236275   | 1                     | 8.247  | 450887  | 20.0 |
| Benzene, propyl-   | 7.213  | 23.9    | ng/ul | 538379   | 1                     | 8.247  | 450887  | 20.0 |
| Benzene, 1-ethy... | 7.319  | 30.9    | ng/ul | 695676   | 1                     | 8.247  | 450887  | 20.0 |
| Benzene, 1,2,3-... | 7.454  | 23.1    | ng/ul | 519679   | 1                     | 8.247  | 450887  | 20.0 |
| Benzene, 1-ethy... | 7.624  | 14.5    | ng/ul | 327250   | 1                     | 8.247  | 450887  | 20.0 |
| (DEL) Alkane: S... | 7.859  | 69.2    | ng/ul | 1560110  | 1                     | 8.247  | 450887  | 20.0 |
| Mesitylene         | 7.889  | 53.2    | ng/ul | 1199320  | 1                     | 8.247  | 450887  | 20.0 |
| (DEL) Alkane: S... | 8.153  | 9.0     | ng/ul | 203052   | 1                     | 8.247  | 450887  | 20.0 |
| Benzene, 1,2,4-... | 8.364  | 20.6    | ng/ul | 464455   | 1                     | 8.247  | 450887  | 20.0 |
| Benzene, 1-meth... | 8.793  | 25.4    | ng/ul | 571388   | 1                     | 8.247  | 450887  | 20.0 |
| Benzene, 1,2-di... | 8.893  | 34.3    | ng/ul | 773188   | 1                     | 8.247  | 450887  | 20.0 |
| (DEL) Alkane: S... | 9.011  | 12.0    | ng/ul | 270381   | 1                     | 8.247  | 450887  | 20.0 |
| Benzene, 4-ethy... | 9.363  | 24.0    | ng/ul | 541019   | 1                     | 8.247  | 450887  | 20.0 |
| (DEL) Alkane: S... | 9.481  | 79.0    | ng/ul | 1780370  | 1                     | 8.247  | 450887  | 20.0 |
| unknown-01         | 9.616  | 16.4    | ng/ul | 369957   | 1                     | 8.247  | 450887  | 20.0 |
| (DEL) Alkane: S... | 12.089 | 5.7     | ng/ul | 290596   | 2                     | 11.090 | 1022770 | 20.0 |
| (DEL) Alkane: S... | 12.477 | 18.4    | ng/ul | 940918   | 2                     | 11.090 | 1022770 | 20.0 |
| (DEL) Alkane: S... | 13.416 | 17.2    | ng/ul | 245892   | 3                     | 14.891 | 286645  | 20.0 |
| Naphthalene, 1-... | 13.922 | 17.7    | ng/ul | 253588   | 3                     | 14.891 | 286645  | 20.0 |
| Naphthalene, 2,... | 14.051 | 23.2    | ng/ul | 331998   | 3                     | 14.891 | 286645  | 20.0 |
| Naphthalene, 1,... | 14.209 | 36.6    | ng/ul | 524386   | 3                     | 14.891 | 286645  | 20.0 |
| (DEL) Alkane: S... | 14.356 | 16.8    | ng/ul | 240058   | 3                     | 14.891 | 286645  | 20.0 |
| (DEL) Alkane: S... | 14.768 | 61.6    | ng/ul | 882500   | 3                     | 14.891 | 286645  | 20.0 |
| Naphthalene, 1,... | 15.361 | 24.4    | ng/ul | 349793   | 3                     | 14.891 | 286645  | 20.0 |
| 4,6,8-Trimethyl... | 15.678 | 17.6    | ng/ul | 251756   | 3                     | 14.891 | 286645  | 20.0 |
| (DEL) Alkane: S... | 15.713 | 52.6    | ng/ul | 754052   | 3                     | 14.891 | 286645  | 20.0 |
| unknown-02         | 15.819 | 14.6    | ng/ul | 208732   | 3                     | 14.891 | 286645  | 20.0 |
| (DEL) Alkane: S... | 16.125 | 25.5    | ng/ul | 365965   | 3                     | 14.891 | 286645  | 20.0 |
| (DEL) Alkane: S... | 16.577 | 38.0    | ng/ul | 721151   | 4                     | 17.640 | 379441  | 20.0 |
| unknown-03         | 16.600 | 20.1    | ng/ul | 381915   | 4                     | 17.640 | 379441  | 20.0 |
| Tetradecanoic acid | 17.017 | 26.5    | ng/ul | 503099   | 4                     | 17.640 | 379441  | 20.0 |
| (DEL) Alkane: S... | 17.370 | 24.7    | ng/ul | 468386   | 4                     | 17.640 | 379441  | 20.0 |
| Benzene, (1-met... | 17.517 | 14.1    | ng/ul | 268237   | 4                     | 17.640 | 379441  | 20.0 |
| 1-Phenylethanet... | 17.846 | 16.9    | ng/ul | 320024   | 4                     | 17.640 | 379441  | 20.0 |
| (DEL) Alkane: S... | 18.110 | 26.1    | ng/ul | 496134   | 4                     | 17.640 | 379441  | 20.0 |
| n-Hexadecanoic ... | 18.533 | 177.4   | ng/ul | 3366370  | 4                     | 17.640 | 379441  | 20.0 |
| (DEL) Alkane: S... | 18.797 | 15.8    | ng/ul | 299883   | 4                     | 17.640 | 379441  | 20.0 |
| p-Dicyclohexylb... | 19.067 | 15.1    | ng/ul | 286094   | 4                     | 17.640 | 379441  | 20.0 |
| (DEL) Alkane: S... | 19.420 | 26.0    | ng/ul | 492743   | 4                     | 17.640 | 379441  | 20.0 |
| cyclohexane, [1... | 19.608 | 16.3    | ng/ul | 309121   | 4                     | 17.640 | 379441  | 20.0 |
| Octadecanoic acid  | 19.773 | 18.5    | ng/ul | 350539   | 4                     | 17.640 | 379441  | 20.0 |
| (DEL) Alkane: S... | 20.524 | 9.1     | ng/ul | 296098   | 5                     | 21.958 | 648833  | 20.0 |
| (DEL) Alkane: S... | 21.030 | 7.7     | ng/ul | 249438   | 5                     | 21.958 | 648833  | 20.0 |
| (DEL) Alkane: S... | 21.564 | 17.9    | ng/ul | 579770   | 5                     | 21.958 | 648833  | 20.0 |
| (DEL) Alkane: S... | 22.146 | 19.0    | ng/ul | 617706   | 5                     | 21.958 | 648833  | 20.0 |
| (DEL) Alkane: S... | 22.804 | 15.6    | ng/ul | 504357   | 5                     | 21.958 | 648833  | 20.0 |
| (DEL) Alkane: S... | 23.321 | 9.4     | ng/ul | 305379   | 5                     | 21.958 | 648833  | 20.0 |
| (DEL) Alkane: S... | 23.556 | 12.2    | ng/ul | 395339   | 5                     | 21.958 | 648833  | 20.0 |

|                    |        |      |       |        |   |        |        |      |
|--------------------|--------|------|-------|--------|---|--------|--------|------|
| (DEL) Alkane: S... | 24.443 | 9.5  | ng/ul | 365247 | 6 | 25.447 | 766845 | 20.0 |
| 1,2-Benzenedica... | 24.836 | 15.2 | ng/ul | 580794 | 6 | 25.447 | 766845 | 20.0 |
| (DEL) Alkane: S... | 26.734 | 8.1  | ng/ul | 312162 | 6 | 25.447 | 766845 | 20.0 |
| Dodecanoic acid... | 27.750 | 18.8 | ng/ul | 719512 | 6 | 25.447 | 766845 | 20.0 |
| (DEL) Alkane: S... | 28.238 | 5.5  | ng/ul | 210015 | 6 | 25.447 | 766845 | 20.0 |

Instrument :

BNA\_G

ClientSampleId :

BGLS8ME