

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122721\  
 Data File : BG051828.D  
 Acq On : 28 Dec 2021 14:05  
 Operator : CG/JU  
 Sample : M5215-01DL 2X  
 Misc :  
 ALS Vial : 35 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 BGLD3DL

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-BG121421.M  
 Title : SVOA CALIBRATION

Signal : TIC: BG051828.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         | 5.766       | 378           | 388         | 395          | rVB      | 7215890        | 14744470      | 14.16%          | 2.577%        |
| 2         | 5.919       | 403           | 414         | 422          | rVB      | 13936095       | 40123117      | 38.54%          | 7.014%        |
| 3         | 6.271       | 468           | 474         | 484          | rVB2     | 6452499        | 11341974      | 10.89%          | 1.983%        |
| 4         | 6.530       | 511           | 518         | 526          | rBV4     | 598322         | 1503929       | 1.44%           | 0.263%        |
| 5         | 6.653       | 532           | 539         | 546          | rVB2     | 461724         | 924323        | 0.89%           | 0.162%        |
| 6         | 6.747       | 546           | 555         | 563          | rBV6     | 658806         | 1579641       | 1.52%           | 0.276%        |
| 7         | 6.818       | 563           | 567         | 572          | rVB      | 831057         | 1202131       | 1.15%           | 0.210%        |
| 8         | 6.906       | 578           | 582         | 597          | rVB3     | 880911         | 2269793       | 2.18%           | 0.397%        |
| 9         | 7.164       | 619           | 626         | 632          | rVB2     | 445132         | 1068786       | 1.03%           | 0.187%        |
| 10        | 7.258       | 632           | 642         | 648          | rBV3     | 1299180        | 3118264       | 2.99%           | 0.545%        |
| 11        | 7.323       | 648           | 653         | 656          | rBV      | 995933         | 1501109       | 1.44%           | 0.262%        |
| 12        | 7.364       | 656           | 660         | 665          | rVB      | 2015884        | 3155410       | 3.03%           | 0.552%        |
| 13        | 7.429       | 665           | 671         | 677          | rVB2     | 2074605        | 3659725       | 3.52%           | 0.640%        |
| 14        | 7.499       | 677           | 683         | 691          | rVB      | 1535426        | 2619848       | 2.52%           | 0.458%        |
| 15        | 7.611       | 691           | 702         | 707          | rBV4     | 302693         | 846437        | 0.81%           | 0.148%        |
| 16        | 7.670       | 707           | 712         | 717          | rBV      | 1047558        | 1663142       | 1.60%           | 0.291%        |
| 17        | 7.910       | 746           | 753         | 755          | rBV      | 3976585        | 6558646       | 6.30%           | 1.146%        |
| 18        | 7.940       | 755           | 758         | 766          | rVB      | 3271798        | 4742023       | 4.55%           | 0.829%        |
| 19        | 8.192       | 791           | 801         | 808          | rBV7     | 356014         | 1015658       | 0.98%           | 0.178%        |
| 20        | 8.263       | 808           | 813         | 820          | rVB2     | 1183447        | 2012179       | 1.93%           | 0.352%        |
| 21        | 8.410       | 833           | 838         | 845          | rVB2     | 1086226        | 1949193       | 1.87%           | 0.341%        |
| 22        | 8.586       | 863           | 868         | 874          | rVB2     | 608244         | 1025354       | 0.98%           | 0.179%        |
| 23        | 8.839       | 904           | 911         | 915          | rVV3     | 842364         | 2116835       | 2.03%           | 0.370%        |
| 24        | 8.938       | 922           | 928         | 941          | rBV4     | 1122249        | 2827020       | 2.72%           | 0.494%        |
| 25        | 9.056       | 941           | 948         | 952          | rBV2     | 675782         | 1118647       | 1.07%           | 0.196%        |
| 26        | 9.408       | 998           | 1008        | 1014         | rBV2     | 736051         | 1667161       | 1.60%           | 0.291%        |
| 27        | 9.532       | 1019          | 1029        | 1034         | rVB      | 2837216        | 5094679       | 4.89%           | 0.891%        |
| 28        | 9.661       | 1040          | 1051        | 1060         | rBV8     | 367958         | 1322659       | 1.27%           | 0.231%        |
| 29        | 10.524      | 1191          | 1198        | 1204         | rVV4     | 534048         | 1150453       | 1.10%           | 0.201%        |
| 30        | 11.083      | 1287          | 1293        | 1303         | rVV2     | 1650655        | 3018162       | 2.90%           | 0.528%        |
| 31        | 11.182      | 1303          | 1310        | 1316         | rVV      | 2017936        | 3798278       | 3.65%           | 0.664%        |
| 32        | 11.271      | 1316          | 1325        | 1331         | rVB      | 487937         | 974163        | 0.94%           | 0.170%        |
| 33        | 12.128      | 1465          | 1471        | 1475         | rBV      | 714047         | 1179991       | 1.13%           | 0.206%        |
| 34        | 12.516      | 1529          | 1537        | 1547         | rBV      | 2302943        | 3701257       | 3.55%           | 0.647%        |
| 35        | 12.769      | 1575          | 1580        | 1587         | rVB      | 1459626        | 2428856       | 2.33%           | 0.425%        |
| 36        | 12.980      | 1611          | 1616        | 1625         | rVB      | 777830         | 1340152       | 1.29%           | 0.234%        |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122721\  
 Data File : BG051828.D  
 Acq On : 28 Dec 2021 14:05  
 Operator : CG/JU  
 Sample : M5215-01DL 2X  
 Misc :  
 ALS Vial : 35 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 BGLD3DL

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-BG121421.M  
 Title : SVOA CALIBRATION

|    |        |      |      |      |      |         |         |       |        |
|----|--------|------|------|------|------|---------|---------|-------|--------|
| 37 | 13.203 | 1646 | 1654 | 1657 | rVV3 | 369355  | 865951  | 0.83% | 0.151% |
| 38 | 13.456 | 1692 | 1697 | 1703 | rVB  | 1004243 | 1467637 | 1.41% | 0.257% |
| 39 | 13.750 | 1738 | 1747 | 1754 | rBV  | 5472373 | 8839669 | 8.49% | 1.545% |
| 40 | 13.955 | 1778 | 1782 | 1792 | rVB2 | 363810  | 911624  | 0.88% | 0.159% |

|    |        |      |      |      |      |         |         |       |        |
|----|--------|------|------|------|------|---------|---------|-------|--------|
| 41 | 14.084 | 1799 | 1804 | 1815 | rVB2 | 683717  | 2047178 | 1.97% | 0.358% |
| 42 | 14.249 | 1826 | 1832 | 1837 | rBV4 | 1213582 | 2152043 | 2.07% | 0.376% |
| 43 | 14.302 | 1837 | 1841 | 1846 | rVB2 | 1170460 | 1746165 | 1.68% | 0.305% |
| 44 | 14.360 | 1846 | 1851 | 1853 | rBV4 | 787361  | 1266393 | 1.22% | 0.221% |
| 45 | 14.390 | 1853 | 1856 | 1860 | rVB  | 1679189 | 2138264 | 2.05% | 0.374% |

|    |        |      |      |      |      |         |         |       |        |
|----|--------|------|------|------|------|---------|---------|-------|--------|
| 46 | 14.813 | 1922 | 1928 | 1935 | rVB  | 6077335 | 8863279 | 8.51% | 1.549% |
| 47 | 14.883 | 1935 | 1940 | 1944 | rBV4 | 771257  | 1484274 | 1.43% | 0.259% |
| 48 | 15.136 | 1977 | 1983 | 1989 | rVB2 | 561035  | 1300802 | 1.25% | 0.227% |
| 49 | 15.312 | 2007 | 2013 | 2017 | rBV3 | 1166336 | 2591161 | 2.49% | 0.453% |
| 50 | 15.418 | 2029 | 2031 | 2039 | rVB2 | 1017962 | 1742217 | 1.67% | 0.305% |

|    |        |      |      |      |      |         |         |       |        |
|----|--------|------|------|------|------|---------|---------|-------|--------|
| 51 | 15.535 | 2047 | 2051 | 2056 | rBV2 | 537698  | 946615  | 0.91% | 0.165% |
| 52 | 15.588 | 2056 | 2060 | 2064 | rBV2 | 627919  | 877488  | 0.84% | 0.153% |
| 53 | 15.764 | 2084 | 2090 | 2094 | rVB  | 5815420 | 8699767 | 8.36% | 1.521% |
| 54 | 15.853 | 2102 | 2105 | 2112 | rVB2 | 843279  | 1528803 | 1.47% | 0.267% |
| 55 | 16.164 | 2150 | 2158 | 2162 | rBV  | 2582897 | 4937300 | 4.74% | 0.863% |

|    |        |      |      |      |      |         |          |       |        |
|----|--------|------|------|------|------|---------|----------|-------|--------|
| 56 | 16.252 | 2170 | 2173 | 2180 | rVB6 | 730534  | 1072515  | 1.03% | 0.187% |
| 57 | 16.323 | 2180 | 2185 | 2190 | rBV3 | 565438  | 1293049  | 1.24% | 0.226% |
| 58 | 16.375 | 2190 | 2194 | 2197 | rVV3 | 1077915 | 1304015  | 1.25% | 0.228% |
| 59 | 16.411 | 2197 | 2200 | 2204 | rVB2 | 796432  | 907653   | 0.87% | 0.159% |
| 60 | 16.628 | 2231 | 2237 | 2239 | rBV  | 6602505 | 10107789 | 9.71% | 1.767% |

|    |        |      |      |      |      |         |         |       |        |
|----|--------|------|------|------|------|---------|---------|-------|--------|
| 61 | 16.651 | 2239 | 2241 | 2246 | rVB  | 4744054 | 5705845 | 5.48% | 0.997% |
| 62 | 16.740 | 2252 | 2256 | 2260 | rBV  | 1691558 | 2726456 | 2.62% | 0.477% |
| 63 | 16.892 | 2279 | 2282 | 2286 | rVB2 | 976382  | 1239982 | 1.19% | 0.217% |
| 64 | 17.022 | 2301 | 2304 | 2307 | rBV3 | 779287  | 1153557 | 1.11% | 0.202% |
| 65 | 17.421 | 2367 | 2372 | 2376 | rBV  | 5288563 | 7044033 | 6.77% | 1.231% |

|    |        |      |      |      |      |         |         |       |        |
|----|--------|------|------|------|------|---------|---------|-------|--------|
| 66 | 17.474 | 2376 | 2381 | 2385 | rBV  | 2325759 | 3221660 | 3.09% | 0.563% |
| 67 | 17.562 | 2392 | 2396 | 2409 | rVB  | 1460243 | 2746660 | 2.64% | 0.480% |
| 68 | 17.662 | 2410 | 2413 | 2419 | rBV4 | 692970  | 1565169 | 1.50% | 0.274% |
| 69 | 17.721 | 2420 | 2423 | 2431 | rVB  | 1038931 | 1649038 | 1.58% | 0.288% |
| 70 | 18.161 | 2493 | 2498 | 2503 | rVB  | 5625025 | 7327513 | 7.04% | 1.281% |

|    |        |      |      |      |      |          |          |        |         |
|----|--------|------|------|------|------|----------|----------|--------|---------|
| 71 | 18.320 | 2522 | 2525 | 2532 | rVB3 | 1123580  | 1706839  | 1.64%  | 0.298%  |
| 72 | 18.672 | 2562 | 2585 | 2589 | rBV3 | 13217577 | 57894875 | 55.61% | 10.120% |
| 73 | 18.849 | 2610 | 2615 | 2619 | rVB  | 4078047  | 4920405  | 4.73%  | 0.860%  |
| 74 | 19.113 | 2657 | 2660 | 2669 | rVB4 | 1594391  | 2307580  | 2.22%  | 0.403%  |
| 75 | 19.230 | 2677 | 2680 | 2686 | rVB4 | 866435   | 1102217  | 1.06%  | 0.193%  |

|    |        |      |      |      |      |         |         |       |        |
|----|--------|------|------|------|------|---------|---------|-------|--------|
| 76 | 19.465 | 2714 | 2720 | 2723 | rVB2 | 3734230 | 5283490 | 5.07% | 0.924% |
|----|--------|------|------|------|------|---------|---------|-------|--------|

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122721\  
 Data File : BG051828.D  
 Acq On : 28 Dec 2021 14:05  
 Operator : CG/JU  
 Sample : M5215-01DL 2X  
 Misc :  
 ALS Vial : 35 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 BGLD3DL

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-BG121421.M  
 Title : SVOA CALIBRATION

|     |        |      |      |      |      |          |           |         |         |
|-----|--------|------|------|------|------|----------|-----------|---------|---------|
| 77  | 19.647 | 2747 | 2751 | 2754 | rBV  | 2031101  | 2353315   | 2.26%   | 0.411%  |
| 78  | 19.712 | 2758 | 2762 | 2765 | rBV  | 4265744  | 6557461   | 6.30%   | 1.146%  |
| 79  | 19.871 | 2782 | 2789 | 2795 | rVB  | 5722057  | 10920746  | 10.49%  | 1.909%  |
| 80  | 19.924 | 2796 | 2798 | 2807 | rVB  | 1466729  | 2140571   | 2.06%   | 0.374%  |
| 81  | 20.035 | 2813 | 2817 | 2823 | rVB2 | 2697276  | 3343153   | 3.21%   | 0.584%  |
| 82  | 20.329 | 2864 | 2867 | 2870 | rBV  | 1377395  | 1613477   | 1.55%   | 0.282%  |
| 83  | 20.364 | 2871 | 2873 | 2880 | rVB  | 893336   | 1101207   | 1.06%   | 0.192%  |
| 84  | 20.558 | 2902 | 2906 | 2910 | rBV  | 3659488  | 4253414   | 4.09%   | 0.744%  |
| 85  | 21.016 | 2968 | 2984 | 2987 | rBV  | 23255566 | 67915895  | 65.23%  | 11.872% |
| 86  | 21.069 | 2987 | 2993 | 2997 | rVB  | 3092778  | 3734109   | 3.59%   | 0.653%  |
| 87  | 21.328 | 3033 | 3037 | 3042 | rVB  | 2059043  | 2811362   | 2.70%   | 0.491%  |
| 88  | 21.486 | 3062 | 3064 | 3069 | rVB  | 979962   | 1203974   | 1.16%   | 0.210%  |
| 89  | 21.609 | 3080 | 3085 | 3098 | rVB  | 4053260  | 6565575   | 6.31%   | 1.148%  |
| 90  | 21.944 | 3124 | 3142 | 3148 | rBV2 | 25610857 | 104115806 | 100.00% | 18.200% |
| 91  | 22.203 | 3178 | 3186 | 3191 | rBV2 | 3352224  | 6924847   | 6.65%   | 1.210%  |
| 92  | 22.661 | 3261 | 3264 | 3268 | rBV  | 895793   | 1269525   | 1.22%   | 0.222%  |
| 93  | 22.702 | 3268 | 3271 | 3272 | rBV  | 710379   | 773115    | 0.74%   | 0.135%  |
| 94  | 22.867 | 3294 | 3299 | 3305 | rVB  | 2637677  | 4639029   | 4.46%   | 0.811%  |
| 95  | 23.096 | 3332 | 3338 | 3341 | rBV  | 1573408  | 3253145   | 3.12%   | 0.569%  |
| 96  | 23.201 | 3350 | 3356 | 3368 | rVB  | 2541382  | 6298644   | 6.05%   | 1.101%  |
| 97  | 23.630 | 3423 | 3429 | 3437 | rVB2 | 1805308  | 3721042   | 3.57%   | 0.650%  |
| 98  | 24.141 | 3509 | 3516 | 3527 | rVB  | 1161699  | 3148519   | 3.02%   | 0.550%  |
| 99  | 24.535 | 3577 | 3583 | 3590 | rBV  | 1285866  | 2741144   | 2.63%   | 0.479%  |
| 100 | 27.349 | 4053 | 4062 | 4078 | rVB2 | 897437   | 3620133   | 3.48%   | 0.633%  |

Sum of corrected areas: 572069643



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122721\  
Data File : BG051828.D  
Acq On : 28 Dec 2021 14:05  
Operator : CG/JU  
Sample : M5215-01DL 2X  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD3DL

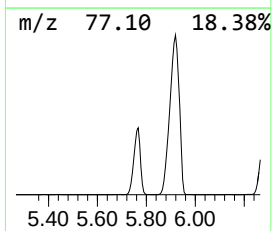
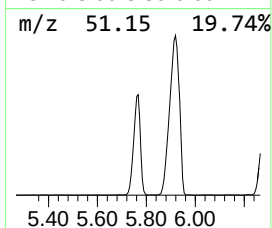
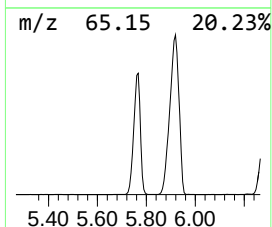
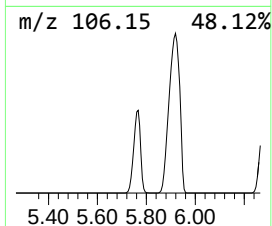
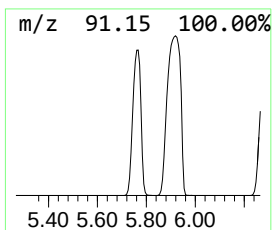
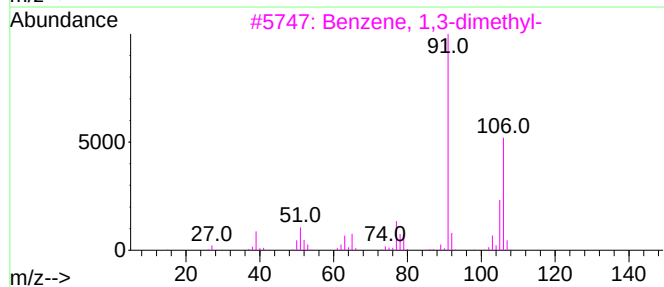
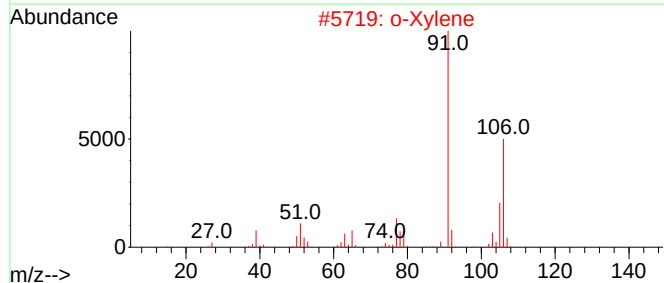
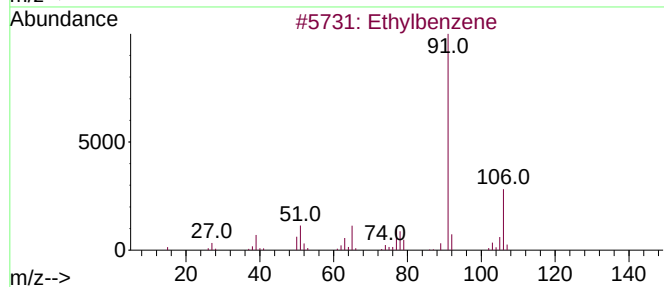
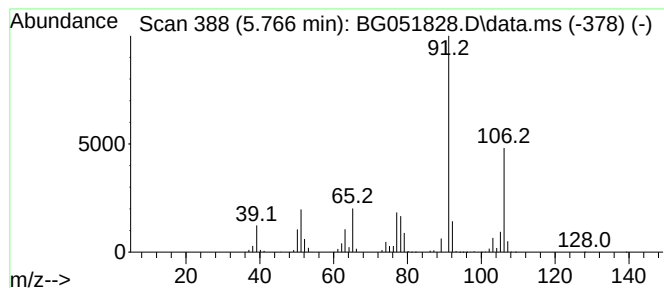
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 1 Ethylbenzene Concentration Rank 3

| R.T.  | EstConc      | Area     | Relative to ISTD       | R.T.  |
|-------|--------------|----------|------------------------|-------|
| 5.766 | 146.55 ng/ul | 14744500 | 1,4-Dichlorobenzene-d4 | 8.286 |

| Hit# | of | 5 | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------------|-----|---------|-------------|------|
| 1    |    |   | Ethylbenzene           | 106 | C8H10   | 000100-41-4 | 91   |
| 2    |    |   | o-Xylene               | 106 | C8H10   | 000095-47-6 | 87   |
| 3    |    |   | Benzene, 1,3-dimethyl- | 106 | C8H10   | 000108-38-3 | 87   |
| 4    |    |   | p-Xylene               | 106 | C8H10   | 000106-42-3 | 87   |
| 5    |    |   | Ethylbenzene           | 106 | C8H10   | 000100-41-4 | 83   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122721\  
Data File : BG051828.D  
Acq On : 28 Dec 2021 14:05  
Operator : CG/JU  
Sample : M5215-01DL 2X  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD3DL

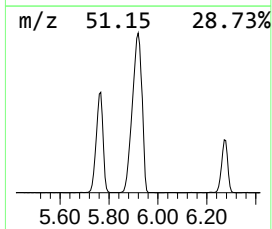
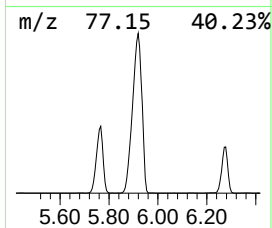
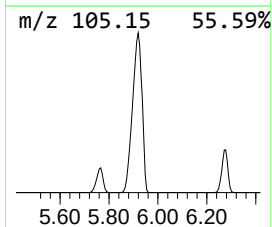
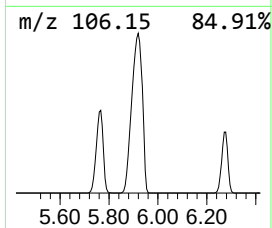
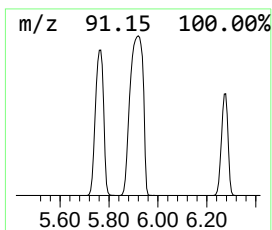
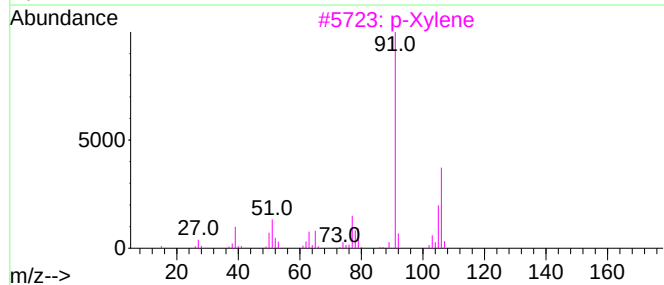
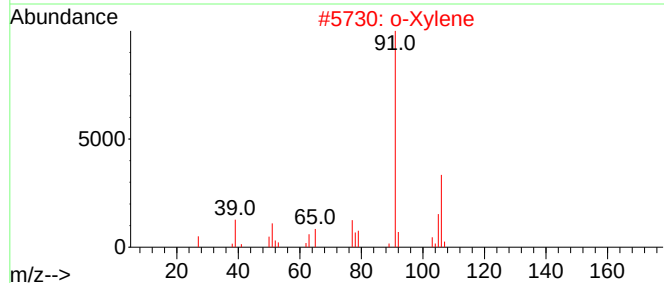
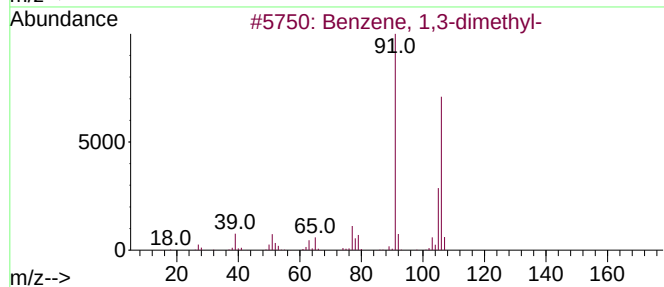
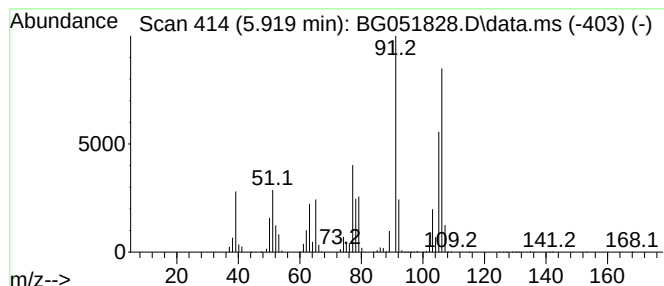
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

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

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

| R.T.  | EstConc      | Area     | Relative to ISTD       | R.T.  |
|-------|--------------|----------|------------------------|-------|
| 5.919 | 398.80 ng/ul | 40123100 | 1,4-Dichlorobenzene-d4 | 8.286 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1,3-dimethyl-              | 106 | C8H10   | 000108-38-3 | 90   |
| 2    |    |   | o-Xylene                            | 106 | C8H10   | 000095-47-6 | 87   |
| 3    |    |   | p-Xylene                            | 106 | C8H10   | 000106-42-3 | 64   |
| 4    |    |   | 3,5-Octadiyne                       | 106 | C8H10   | 016387-70-5 | 58   |
| 5    |    |   | 1,3-Cyclopentadiene, 5-(1-methyl... | 106 | C8H10   | 002175-91-9 | 53   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122721\  
Data File : BG051828.D  
Acq On : 28 Dec 2021 14:05  
Operator : CG/JU  
Sample : M5215-01DL 2X  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD3DL

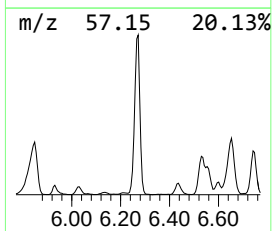
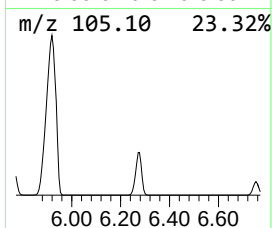
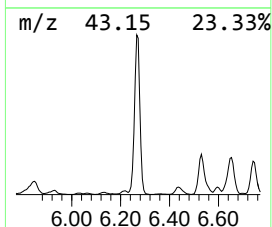
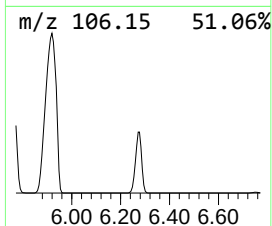
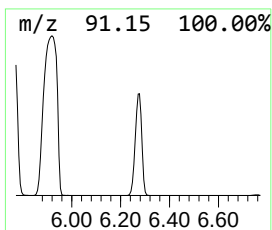
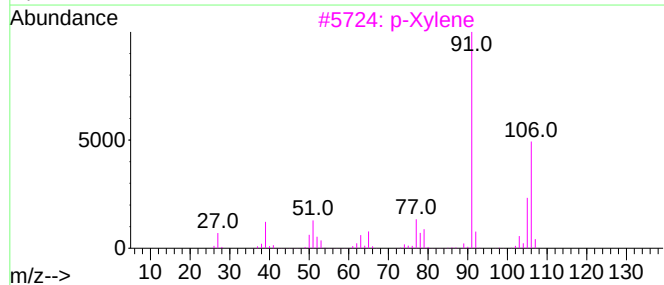
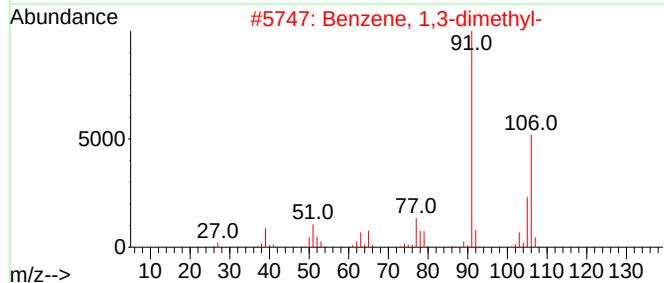
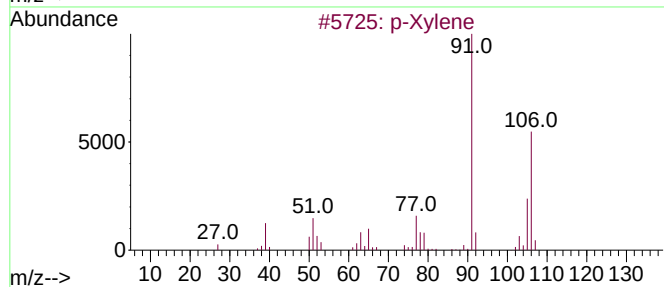
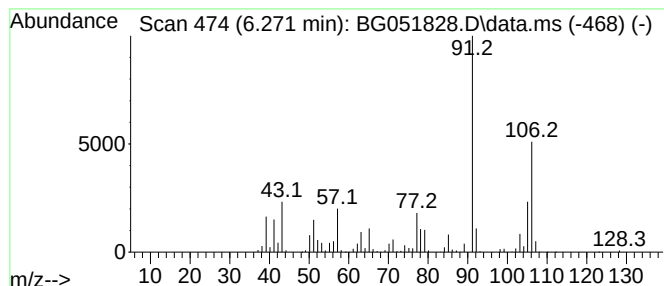
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 3 p-Xylene Concentration Rank 6

| R.T.  | EstConc      | Area     | Relative to ISTD       | R.T.  |
|-------|--------------|----------|------------------------|-------|
| 6.271 | 112.73 ng/ul | 11342000 | 1,4-Dichlorobenzene-d4 | 8.286 |

| Hit# | of                     | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|------------------------|---|--------------|-----|---------|-------------|------|
| 1    | p-Xylene               |   |              | 106 | C8H10   | 000106-42-3 | 97   |
| 2    | Benzene, 1,3-dimethyl- |   |              | 106 | C8H10   | 000108-38-3 | 97   |
| 3    | p-Xylene               |   |              | 106 | C8H10   | 000106-42-3 | 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\BG122721\  
Data File : BG051828.D  
Acq On : 28 Dec 2021 14:05  
Operator : CG/JU  
Sample : M5215-01DL 2X  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD3DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

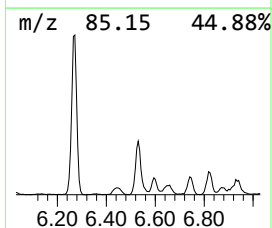
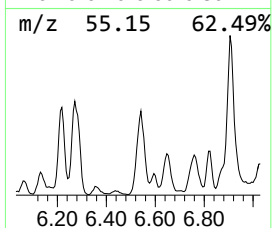
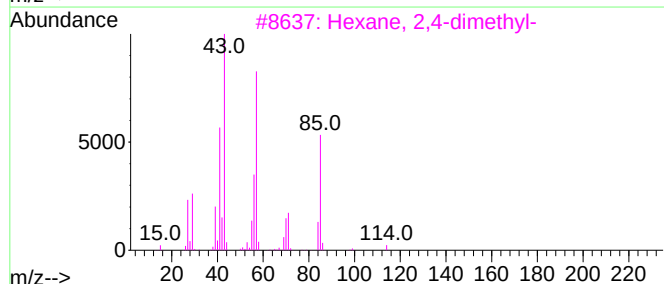
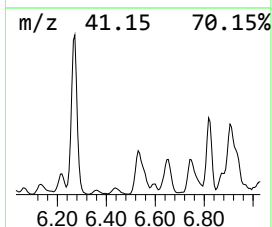
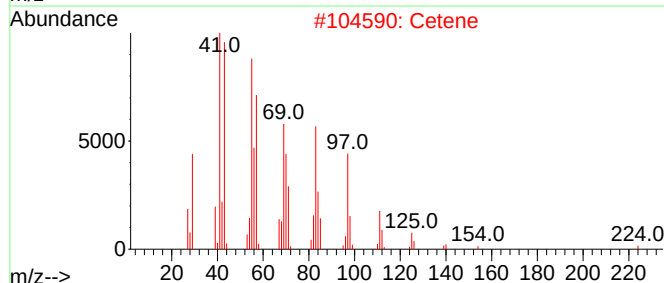
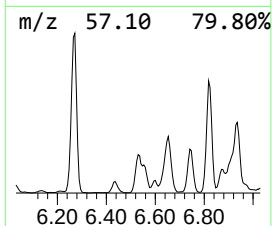
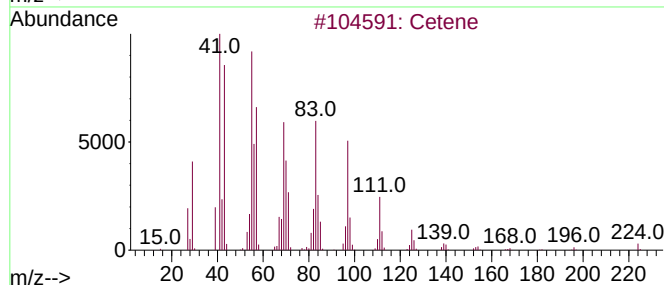
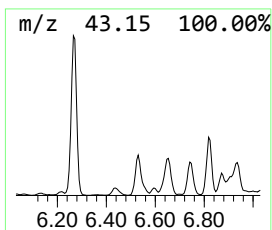
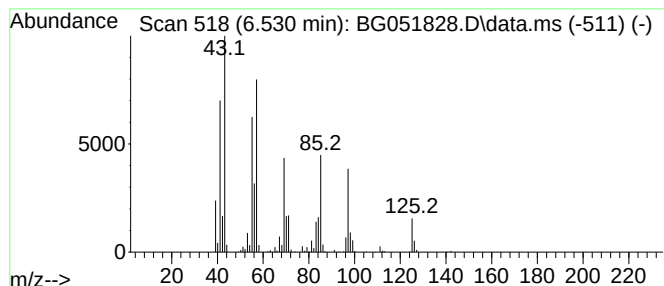
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 43

| R.T.  | EstConc     | Area    | Relative to ISTD       | R.T.  |
|-------|-------------|---------|------------------------|-------|
| 6.530 | 14.95 ng/ul | 1503930 | 1,4-Dichlorobenzene-d4 | 8.286 |

| Hit# | of | 5 | Tentative ID             | MW  | MolForm  | CAS#        | Qual |
|------|----|---|--------------------------|-----|----------|-------------|------|
| 1    |    |   | Cetene                   | 224 | C16H32   | 000629-73-2 | 47   |
| 2    |    |   | Cetene                   | 224 | C16H32   | 000629-73-2 | 47   |
| 3    |    |   | Hexane, 2,4-dimethyl-    | 114 | C8H18    | 000589-43-5 | 46   |
| 4    |    |   | Hexyl isobutyl carbonate | 202 | C11H22O3 | 959067-98-2 | 43   |
| 5    |    |   | Hexane, 2,4-dimethyl-    | 114 | C8H18    | 000589-43-5 | 43   |





Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122721\  
Data File : BG051828.D  
Acq On : 28 Dec 2021 14:05  
Operator : CG/JU  
Sample : M5215-01DL 2X  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD3DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

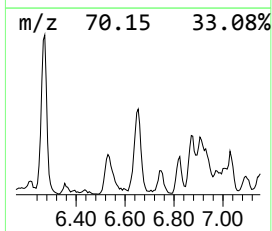
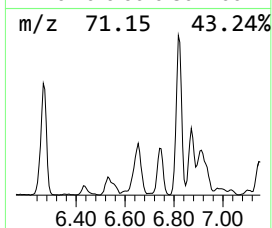
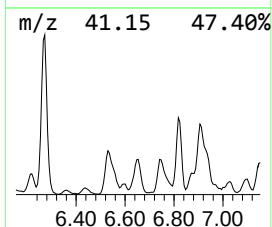
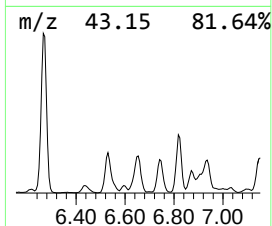
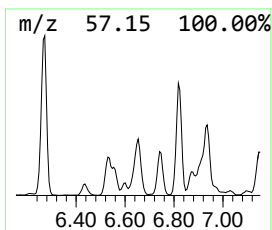
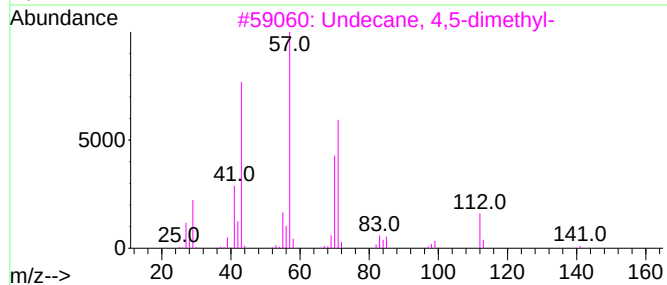
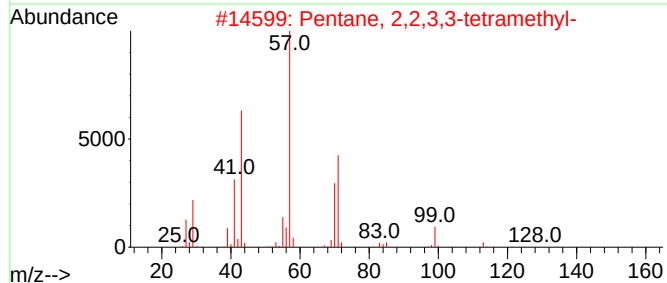
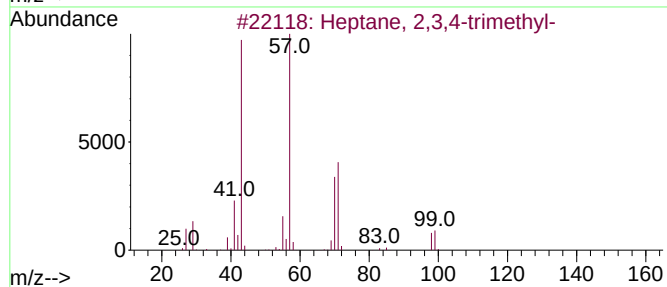
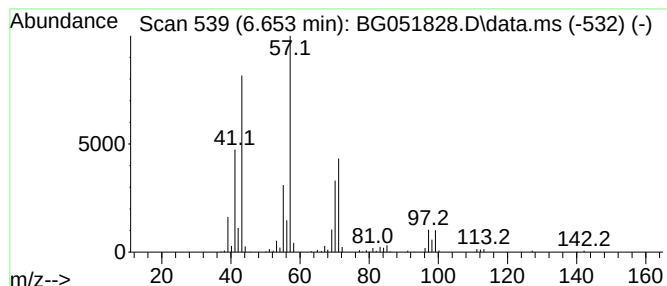
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 58

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 6.653 | 9.19 ng/ul | 924323 | 1,4-Dichlorobenzene-d4 | 8.286 |

| Hit# | of | 5 | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------|-----|---------|-------------|------|
| 1    |    |   | Heptane, 2,3,4-trimethyl-      | 142 | C10H22  | 052896-95-4 | 64   |
| 2    |    |   | Pentane, 2,2,3,3-tetramethyl-  | 128 | C9H20   | 007154-79-2 | 64   |
| 3    |    |   | Undecane, 4,5-dimethyl-        | 184 | C13H28  | 017312-79-7 | 59   |
| 4    |    |   | Nonane, 4-methyl-              | 142 | C10H22  | 017301-94-9 | 53   |
| 5    |    |   | Pentane, 3-ethyl-2,2-dimethyl- | 128 | C9H20   | 016747-32-3 | 50   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122721\  
Data File : BG051828.D  
Acq On : 28 Dec 2021 14:05  
Operator : CG/JU  
Sample : M5215-01DL 2X  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD3DL

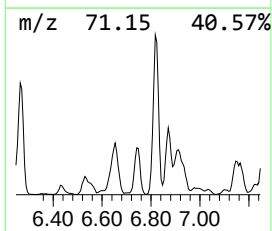
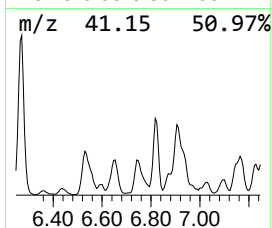
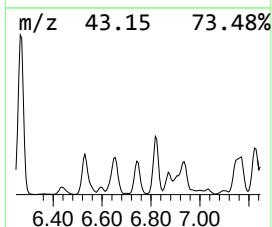
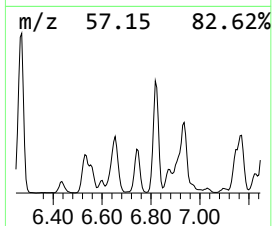
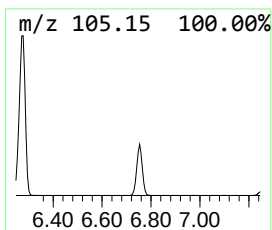
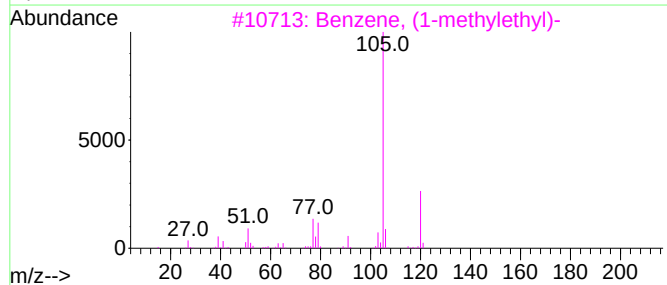
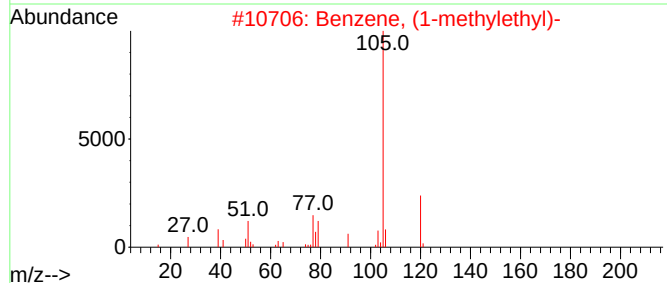
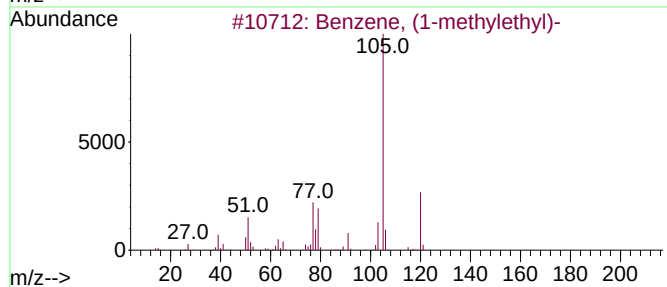
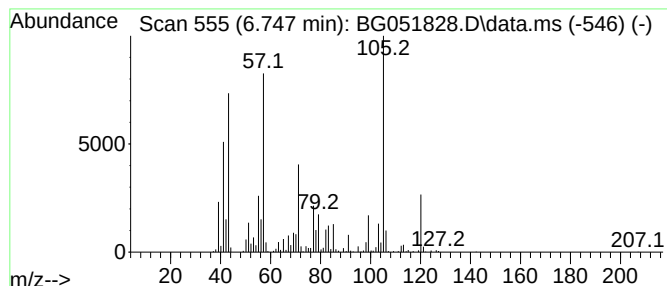
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 6 Benzene, (1-methylethyl)- Concentration Rank 42

| R.T.  | EstConc     | Area    | Relative to ISTD       | R.T.  |
|-------|-------------|---------|------------------------|-------|
| 6.747 | 15.70 ng/ul | 1579640 | 1,4-Dichlorobenzene-d4 | 8.286 |

| Hit# | of | 5 | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|------|----|---|---------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, (1-methylethyl)- | 120 | C9H12   | 000098-82-8 | 94   |
| 2    |    |   | Benzene, (1-methylethyl)- | 120 | C9H12   | 000098-82-8 | 89   |
| 3    |    |   | Benzene, (1-methylethyl)- | 120 | C9H12   | 000098-82-8 | 50   |
| 4    |    |   | Benzene, (1-methylethyl)- | 120 | C9H12   | 000098-82-8 | 50   |
| 5    |    |   | Benzene, 1,2,3-trimethyl- | 120 | C9H12   | 000526-73-8 | 45   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122721\  
Data File : BG051828.D  
Acq On : 28 Dec 2021 14:05  
Operator : CG/JU  
Sample : M5215-01DL 2X  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD3DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

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

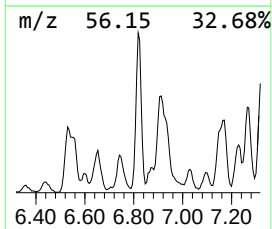
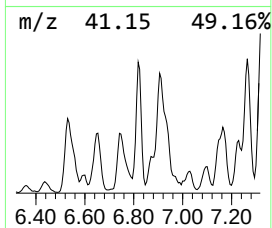
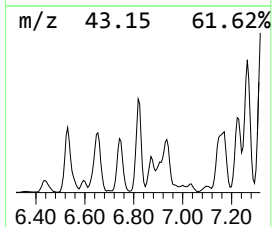
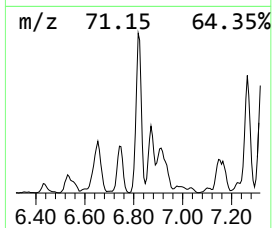
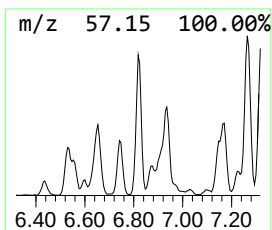
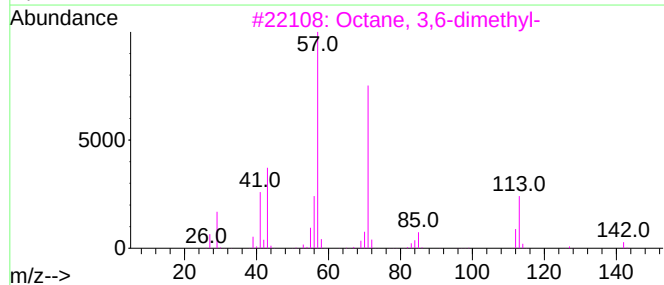
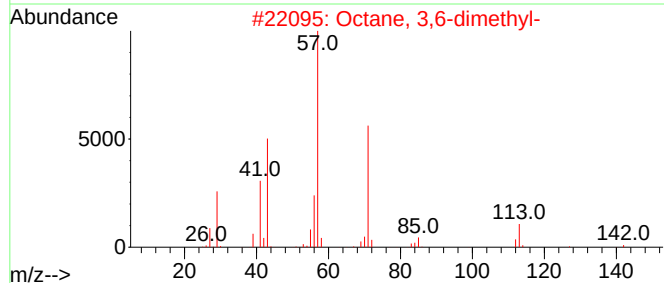
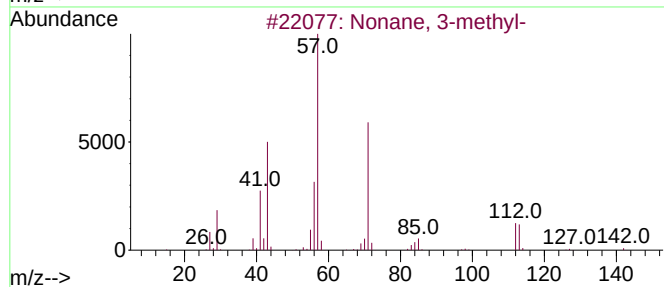
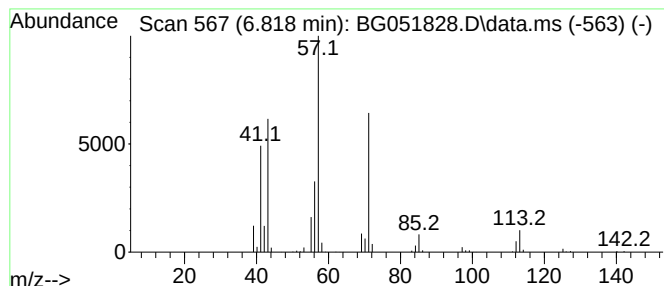
\*\*\*\*\*

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 51

| R.T.  | EstConc     | Area    | Relative to ISTD       | R.T.  |
|-------|-------------|---------|------------------------|-------|
| 6.818 | 11.95 ng/ul | 1202130 | 1,4-Dichlorobenzene-d4 | 8.286 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|------|-------------------------------------|-----|-----------|--------------|------|
| 1    |      | Nonane, 3-methyl-                   | 142 | C10H22    | 005911-04-6  | 90   |
| 2    |      | Octane, 3,6-dimethyl-               | 142 | C10H22    | 015869-94-0  | 86   |
| 3    |      | Octane, 3,6-dimethyl-               | 142 | C10H22    | 015869-94-0  | 78   |
| 4    |      | Nonane, 3-methyl-                   | 142 | C10H22    | 005911-04-6  | 72   |
| 5    |      | Sulfurous acid, butyl 2-ethylhex... | 250 | C12H26O3S | 1000309-18-8 | 64   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122721\  
Data File : BG051828.D  
Acq On : 28 Dec 2021 14:05  
Operator : CG/JU  
Sample : M5215-01DL 2X  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD3DL

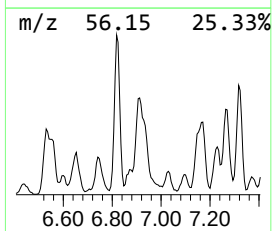
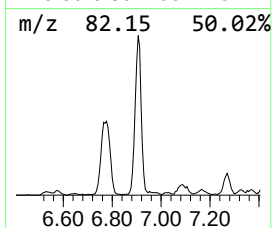
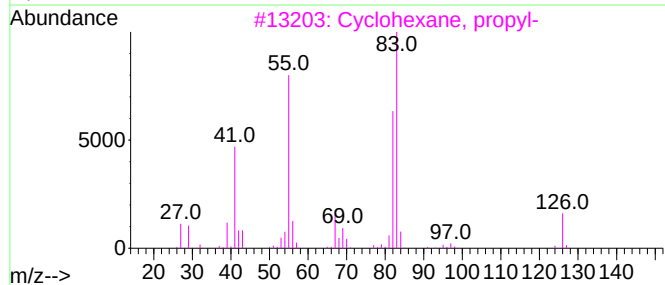
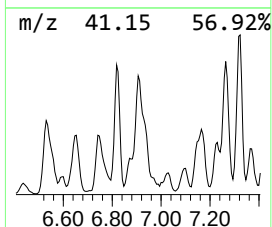
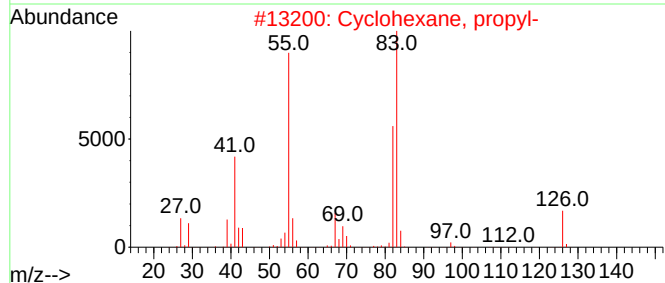
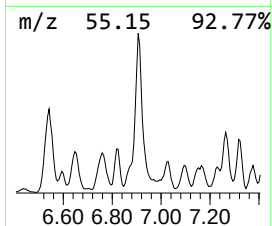
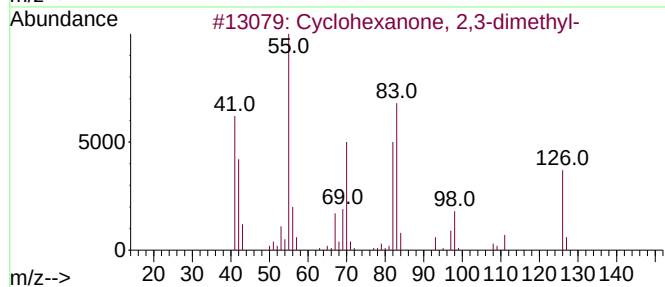
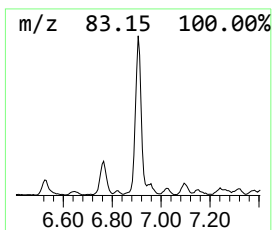
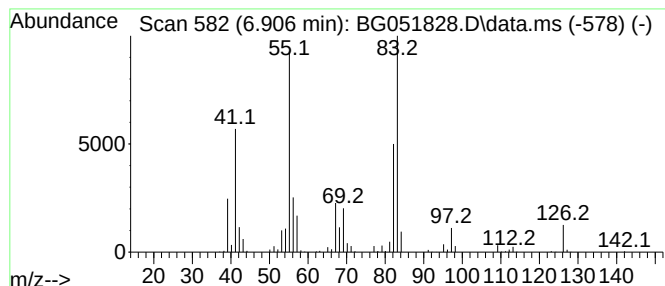
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 8 Cyclohexanone, 2,3-dimethyl- Concentration Rank 30

| R.T.  | EstConc     | Area    | Relative to ISTD       | R.T.  |
|-------|-------------|---------|------------------------|-------|
| 6.906 | 22.56 ng/ul | 2269790 | 1,4-Dichlorobenzene-d4 | 8.286 |

| Hit# | of | 5 | Tentative ID                 | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------------------|-----|---------|-------------|------|
| 1    |    |   | Cyclohexanone, 2,3-dimethyl- | 126 | C8H14O  | 013395-76-1 | 83   |
| 2    |    |   | Cyclohexane, propyl-         | 126 | C9H18   | 001678-92-8 | 72   |
| 3    |    |   | Cyclohexane, propyl-         | 126 | C9H18   | 001678-92-8 | 72   |
| 4    |    |   | Cyclohexane, propyl-         | 126 | C9H18   | 001678-92-8 | 72   |
| 5    |    |   | Cyclohexane, propyl-         | 126 | C9H18   | 001678-92-8 | 64   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122721\  
Data File : BG051828.D  
Acq On : 28 Dec 2021 14:05  
Operator : CG/JU  
Sample : M5215-01DL 2X  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD3DL

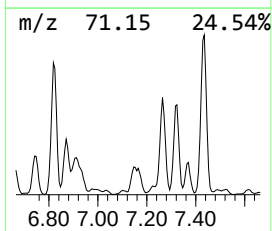
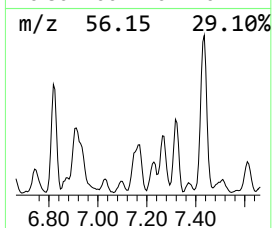
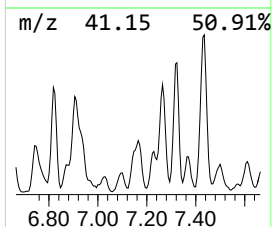
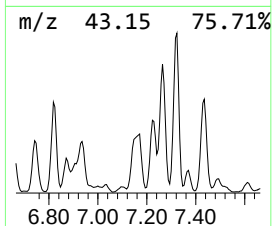
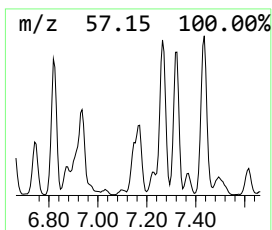
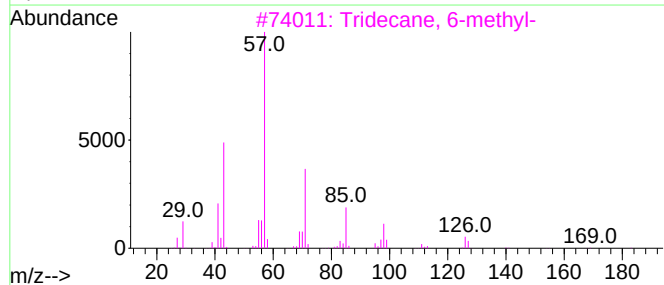
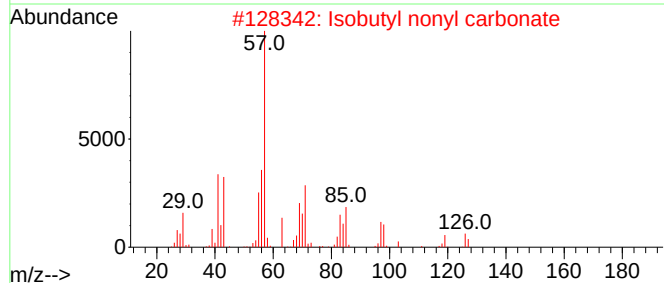
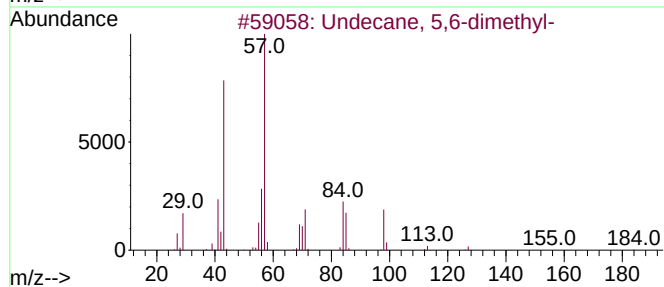
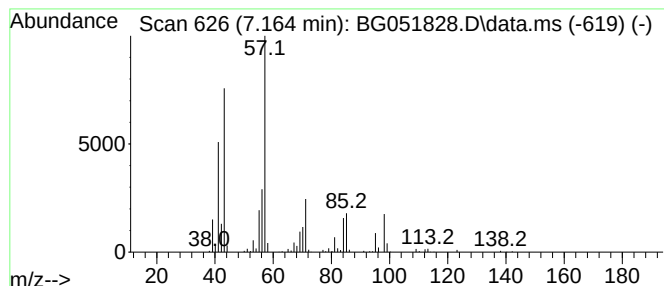
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 55

| R.T.      | EstConc                  | Area    | Relative to ISTD       | R.T.        |      |
|-----------|--------------------------|---------|------------------------|-------------|------|
| 7.164     | 10.62 ng/ul              | 1068790 | 1,4-Dichlorobenzene-d4 | 8.286       |      |
| Hit# of 5 | Tentative ID             | MW      | MolForm                | CAS#        | Qual |
| 1         | Undecane, 5,6-dimethyl-  | 184     | C13H28                 | 017615-91-7 | 80   |
| 2         | Isobutyl nonyl carbonate | 244     | C14H28O3               | 959311-27-4 | 64   |
| 3         | Tridecane, 6-methyl-     | 198     | C14H30                 | 013287-21-3 | 59   |
| 4         | Hexane, 2,4-dimethyl-    | 114     | C8H18                  | 000589-43-5 | 58   |
| 5         | Octane, 2,5,6-trimethyl- | 156     | C11H24                 | 062016-14-2 | 53   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122721\  
Data File : BG051828.D  
Acq On : 28 Dec 2021 14:05  
Operator : CG/JU  
Sample : M5215-01DL 2X  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD3DL

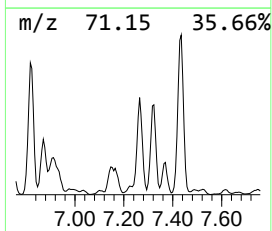
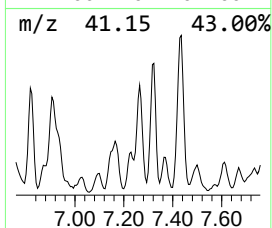
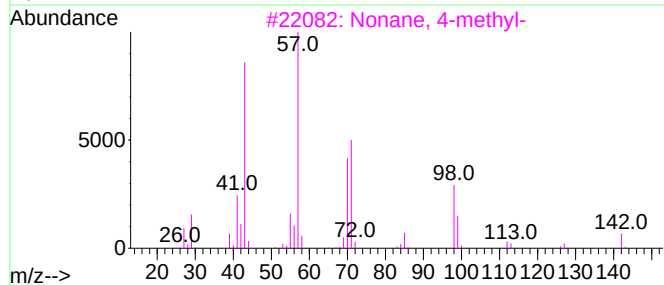
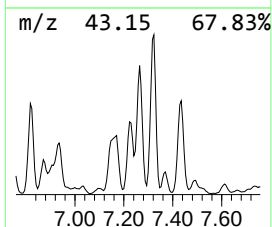
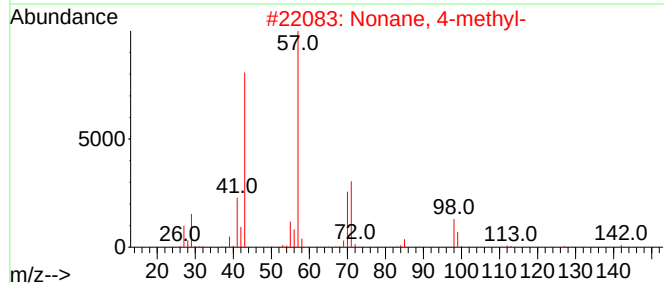
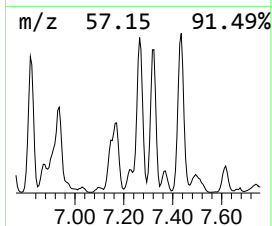
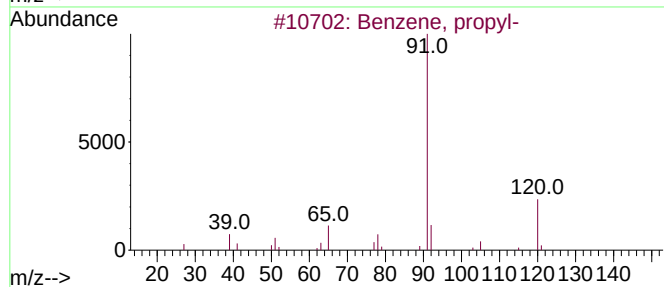
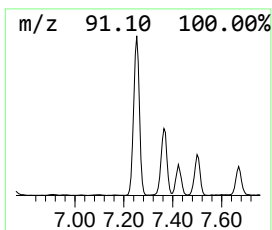
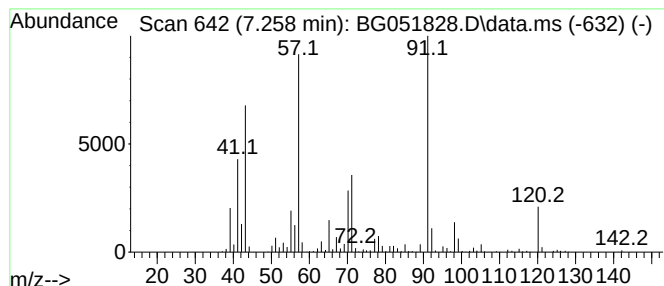
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 10 Benzene, propyl- Concentration Rank 19

| R.T.  | EstConc     | Area    | Relative to ISTD       | R.T.  |
|-------|-------------|---------|------------------------|-------|
| 7.258 | 30.99 ng/ul | 3118260 | 1,4-Dichlorobenzene-d4 | 8.286 |

| Hit# | of | 5 | Tentative ID      | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, propyl-  | 120 | C9H12   | 000103-65-1 | 80   |
| 2    |    |   | Nonane, 4-methyl- | 142 | C10H22  | 017301-94-9 | 60   |
| 3    |    |   | Nonane, 4-methyl- | 142 | C10H22  | 017301-94-9 | 47   |
| 4    |    |   | Benzene, propyl-  | 120 | C9H12   | 000103-65-1 | 42   |
| 5    |    |   | Benzene, propyl-  | 120 | C9H12   | 000103-65-1 | 38   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122721\  
Data File : BG051828.D  
Acq On : 28 Dec 2021 14:05  
Operator : CG/JU  
Sample : M5215-01DL 2X  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD3DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

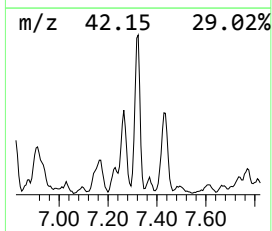
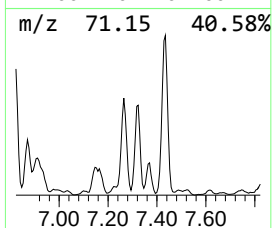
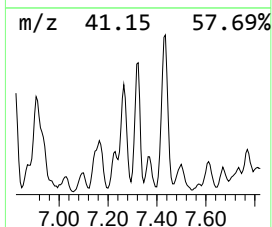
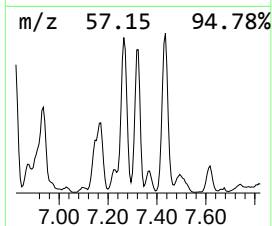
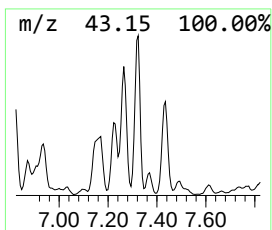
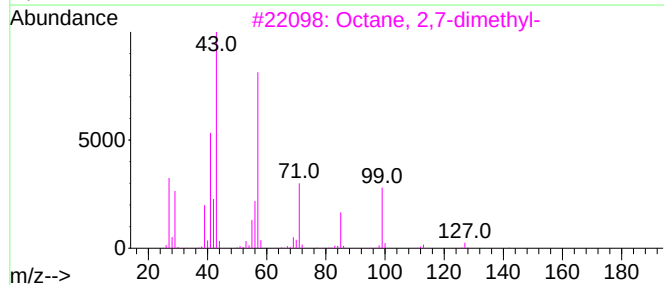
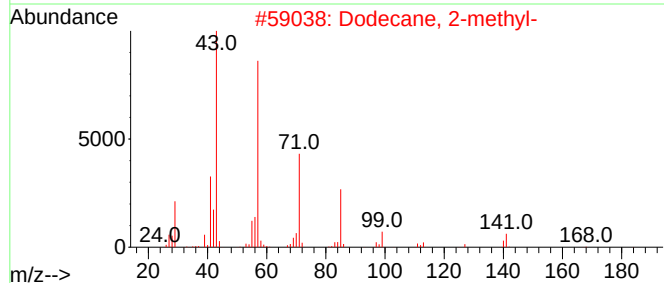
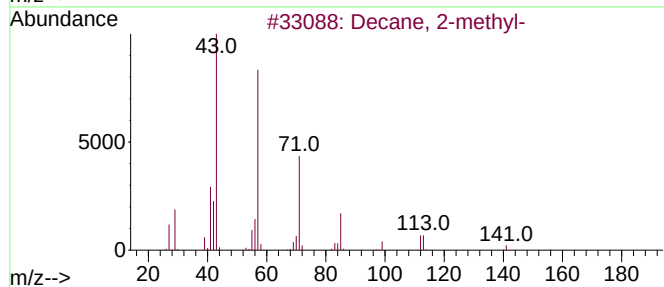
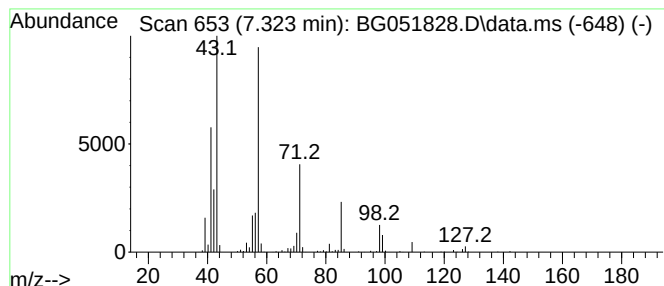
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 44

| R.T.  | EstConc     | Area    | Relative to ISTD       | R.T.  |
|-------|-------------|---------|------------------------|-------|
| 7.323 | 14.92 ng/ul | 1501110 | 1,4-Dichlorobenzene-d4 | 8.286 |

| Hit# | of | 5 | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------|-----|---------|-------------|------|
| 1    |    |   | Decane, 2-methyl-     | 156 | C11H24  | 006975-98-0 | 72   |
| 2    |    |   | Dodecane, 2-methyl-   | 184 | C13H28  | 001560-97-0 | 72   |
| 3    |    |   | Octane, 2,7-dimethyl- | 142 | C10H22  | 001072-16-8 | 72   |
| 4    |    |   | Octane, 2-methyl-     | 128 | C9H20   | 003221-61-2 | 59   |
| 5    |    |   | Decane                | 142 | C10H22  | 000124-18-5 | 59   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122721\  
Data File : BG051828.D  
Acq On : 28 Dec 2021 14:05  
Operator : CG/JU  
Sample : M5215-01DL 2X  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD3DL

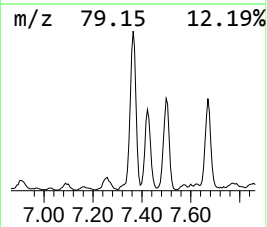
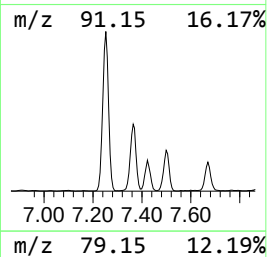
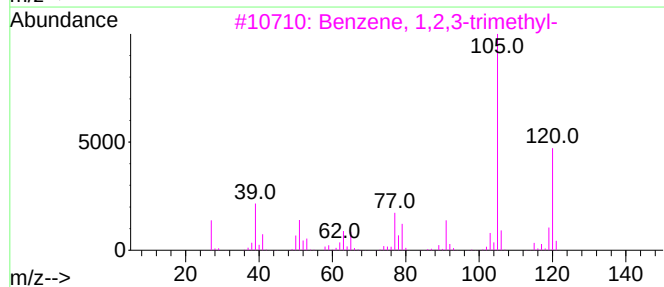
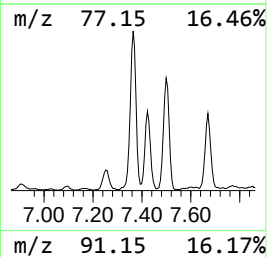
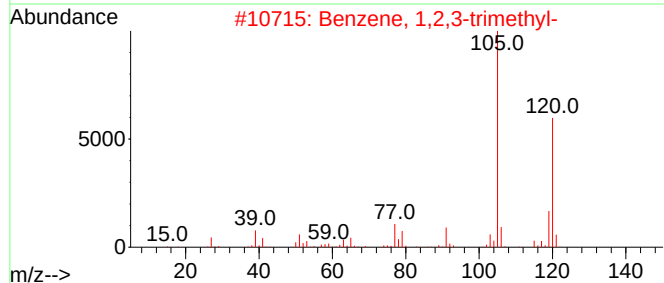
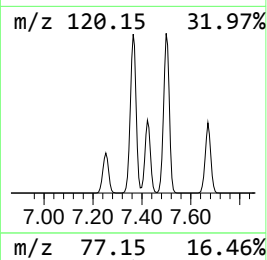
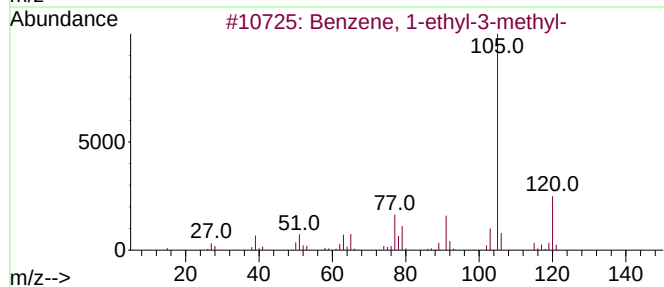
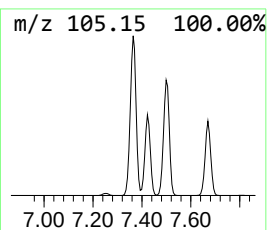
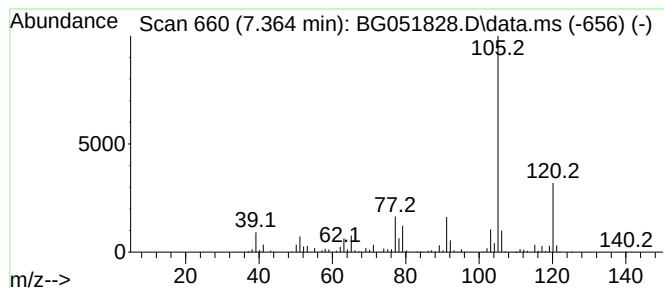
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 12 Benzene, 1-ethyl-3-methyl- Concentration Rank 18

| R.T.  | EstConc     | Area    | Relative to ISTD       | R.T.  |
|-------|-------------|---------|------------------------|-------|
| 7.364 | 31.36 ng/ul | 3155410 | 1,4-Dichlorobenzene-d4 | 8.286 |

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





Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122721\  
Data File : BG051828.D  
Acq On : 28 Dec 2021 14:05  
Operator : CG/JU  
Sample : M5215-01DL 2X  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD3DL

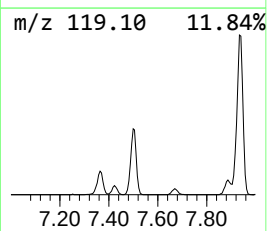
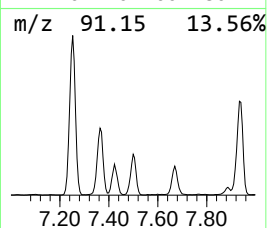
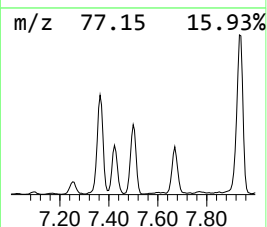
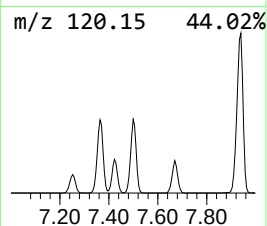
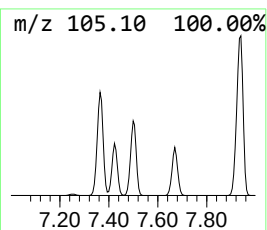
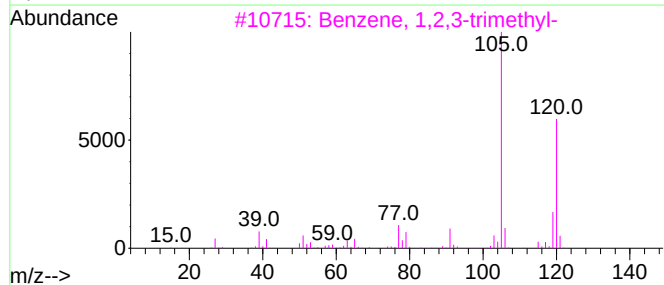
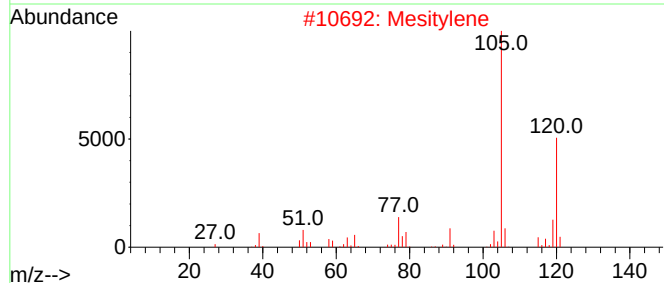
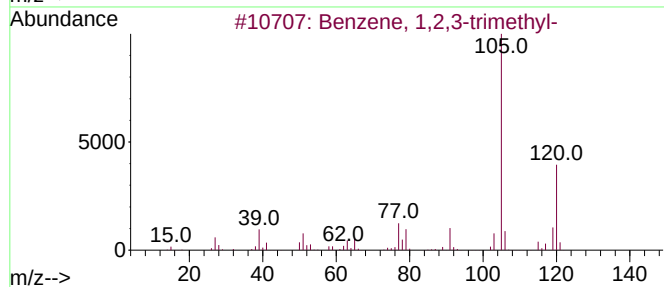
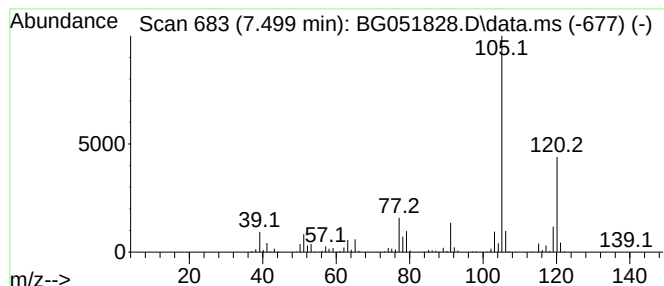
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 13 Benzene, 1,2,3-trimethyl- Concentration Rank 26

| R.T.  | EstConc     | Area    | Relative to ISTD       | R.T.  |
|-------|-------------|---------|------------------------|-------|
| 7.499 | 26.04 ng/ul | 2619850 | 1,4-Dichlorobenzene-d4 | 8.286 |

| 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 | 95   |
| 3    |    |   | Benzene, 1,2,3-trimethyl- | 120 | C9H12   | 000526-73-8 | 95   |
| 4    |    |   | Mesitylene                | 120 | C9H12   | 000108-67-8 | 94   |
| 5    |    |   | Benzene, 1,2,3-trimethyl- | 120 | C9H12   | 000526-73-8 | 94   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122721\  
Data File : BG051828.D  
Acq On : 28 Dec 2021 14:05  
Operator : CG/JU  
Sample : M5215-01DL 2X  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD3DL

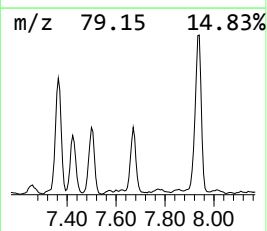
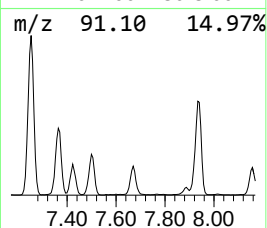
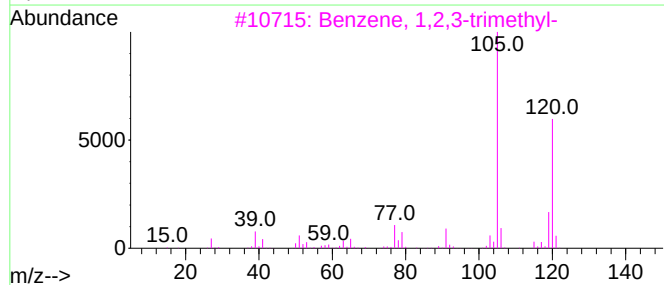
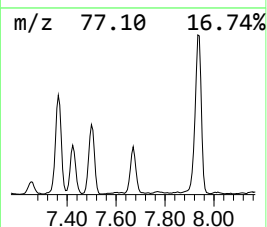
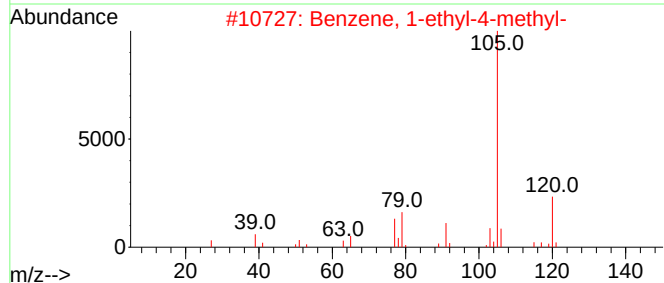
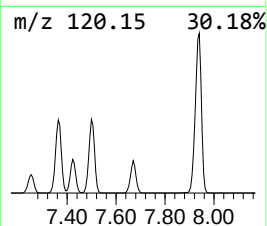
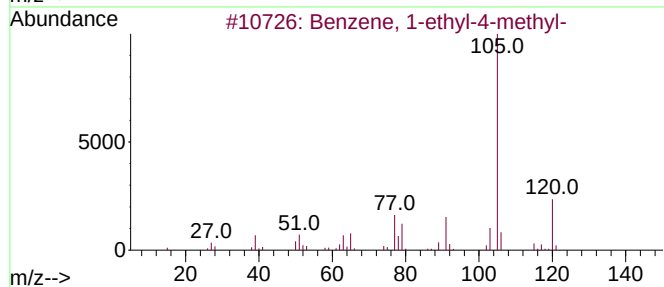
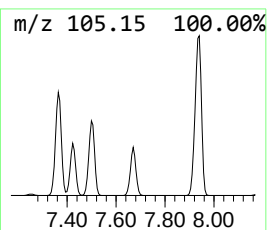
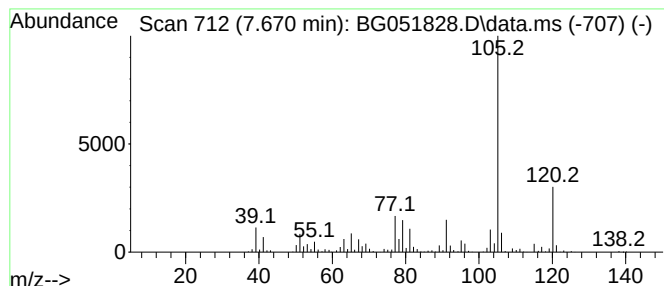
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 14 Benzene, 1-ethyl-4-methyl- Concentration Rank 39

| R.T.  | EstConc     | Area    | Relative to ISTD       | R.T.  |
|-------|-------------|---------|------------------------|-------|
| 7.670 | 16.53 ng/ul | 1663140 | 1,4-Dichlorobenzene-d4 | 8.286 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1-ethyl-4-methyl- | 120 | C9H12   | 000622-96-8 | 95   |
| 2    |    |   | Benzene, 1-ethyl-4-methyl- | 120 | C9H12   | 000622-96-8 | 94   |
| 3    |    |   | Benzene, 1,2,3-trimethyl-  | 120 | C9H12   | 000526-73-8 | 90   |
| 4    |    |   | Mesitylene                 | 120 | C9H12   | 000108-67-8 | 81   |
| 5    |    |   | Benzene, 1-ethyl-4-methyl- | 120 | C9H12   | 000622-96-8 | 81   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122721\  
Data File : BG051828.D  
Acq On : 28 Dec 2021 14:05  
Operator : CG/JU  
Sample : M5215-01DL 2X  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD3DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

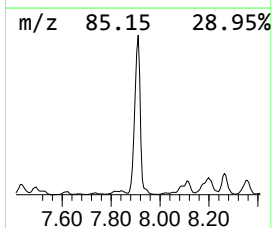
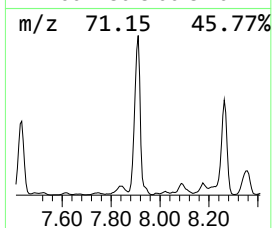
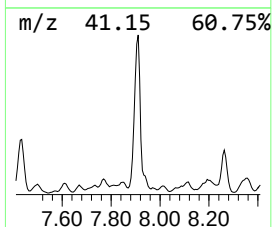
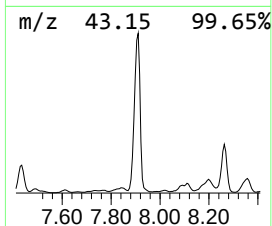
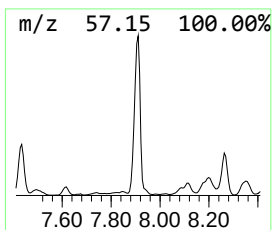
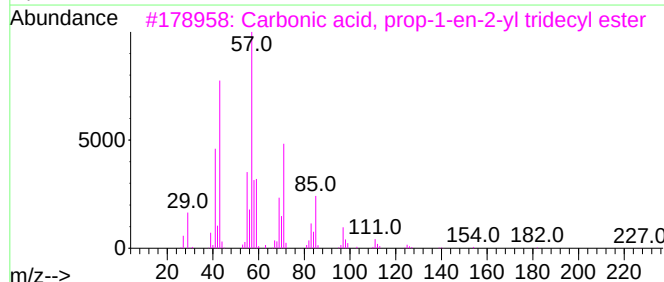
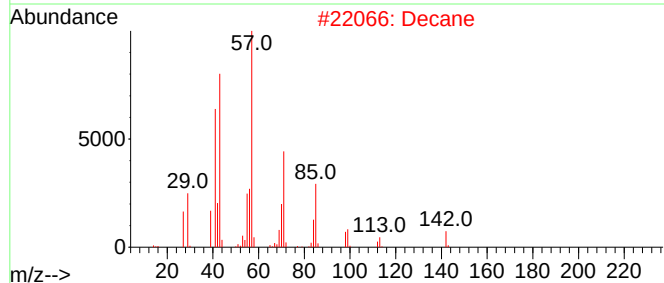
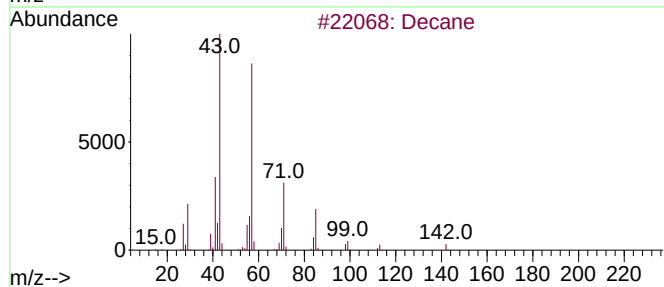
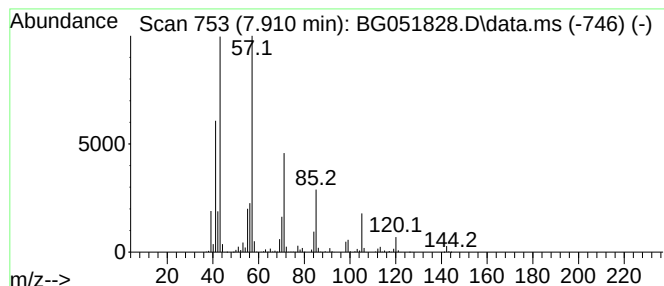
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*

Peak Number 15 (DEL) Alkane: Straight-Chai... Concentration Rank 11

| R.T.  | EstConc     | Area    | Relative to ISTD       | R.T.  |
|-------|-------------|---------|------------------------|-------|
| 7.910 | 65.19 ng/ul | 6558650 | 1,4-Dichlorobenzene-d4 | 8.286 |

| 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 | 86   |
| 4    |    |   | Decane                              | 142 | C10H22   | 000124-18-5  | 78   |
| 5    |    |   | Octane, 4-ethyl-                    | 142 | C10H22   | 015869-86-0  | 72   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122721\  
Data File : BG051828.D  
Acq On : 28 Dec 2021 14:05  
Operator : CG/JU  
Sample : M5215-01DL 2X  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD3DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

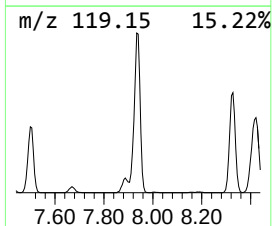
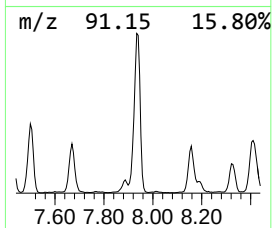
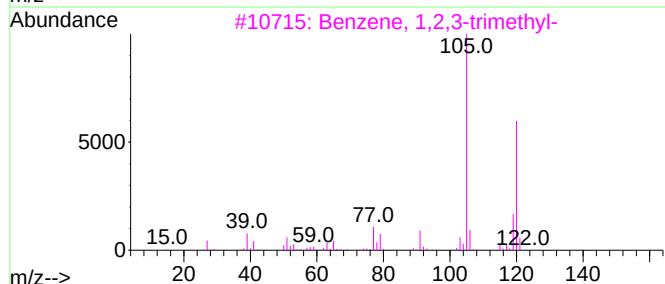
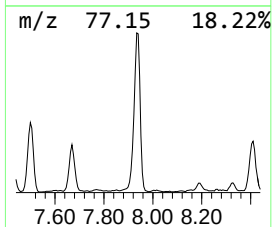
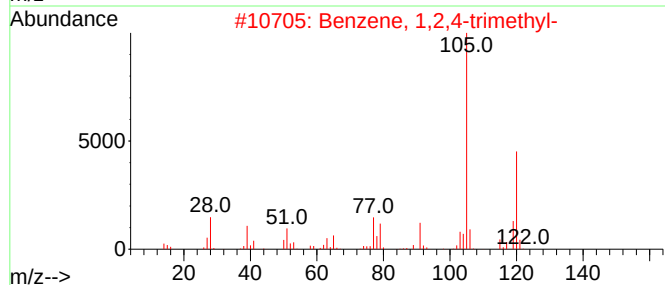
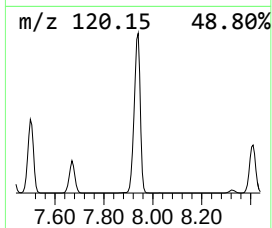
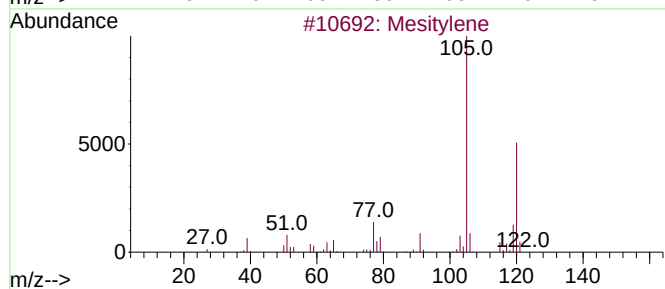
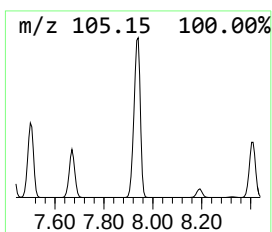
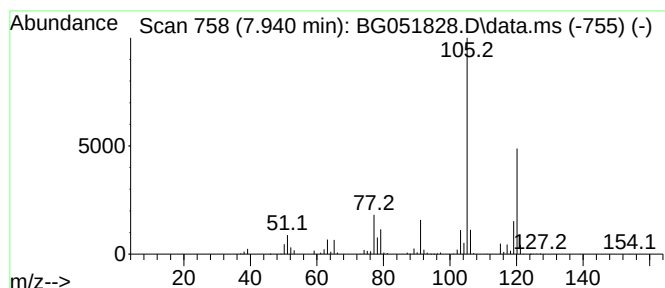
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 16 Mesitylene Concentration Rank 14

| R.T.  | EstConc     | Area    | Relative to ISTD       | R.T.  |
|-------|-------------|---------|------------------------|-------|
| 7.940 | 47.13 ng/ul | 4742020 | 1,4-Dichlorobenzene-d4 | 8.286 |

| Hit# | of | 5 | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|------|----|---|---------------------------|-----|---------|-------------|------|
| 1    |    |   | Mesitylene                | 120 | C9H12   | 000108-67-8 | 95   |
| 2    |    |   | Benzene, 1,2,4-trimethyl- | 120 | C9H12   | 000095-63-6 | 94   |
| 3    |    |   | Benzene, 1,2,3-trimethyl- | 120 | C9H12   | 000526-73-8 | 94   |
| 4    |    |   | Benzene, 1,2,4-trimethyl- | 120 | C9H12   | 000095-63-6 | 93   |
| 5    |    |   | Benzene, 1,2,4-trimethyl- | 120 | C9H12   | 000095-63-6 | 91   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122721\  
Data File : BG051828.D  
Acq On : 28 Dec 2021 14:05  
Operator : CG/JU  
Sample : M5215-01DL 2X  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD3DL

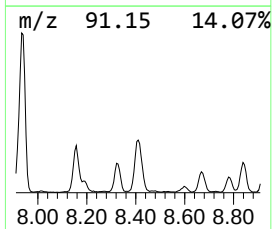
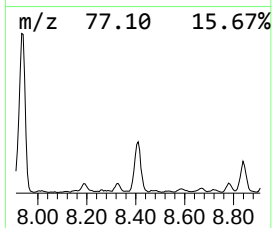
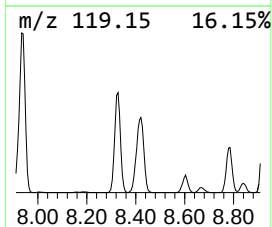
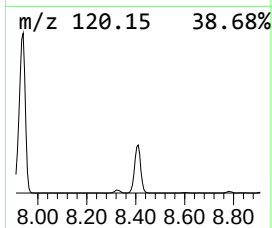
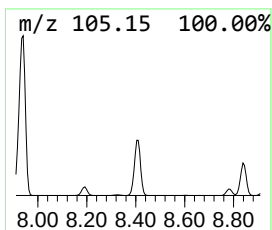
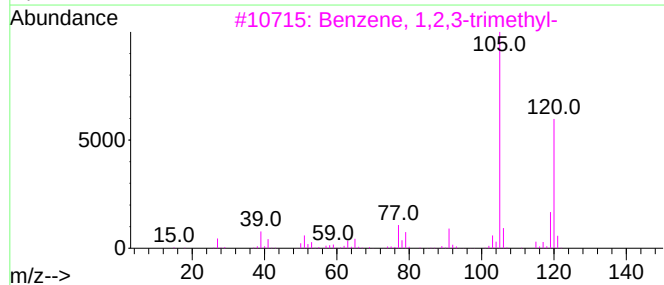
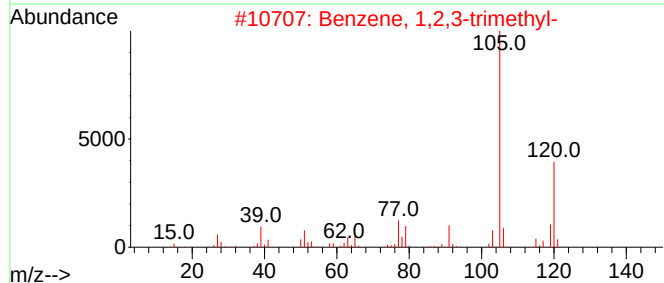
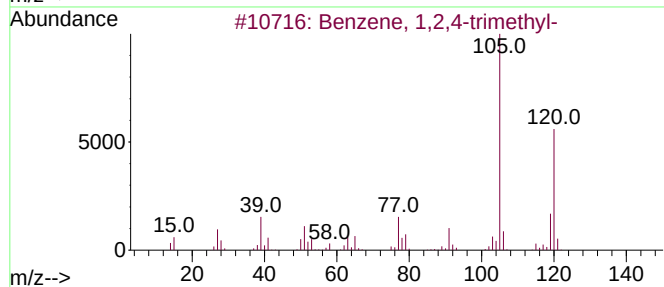
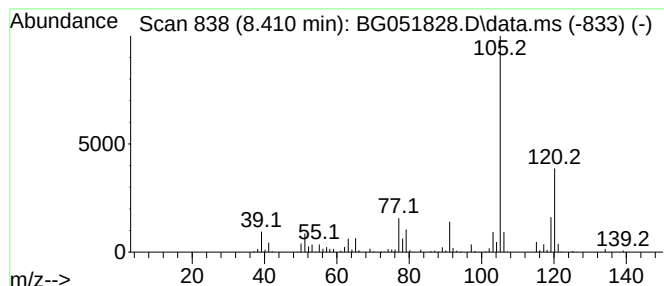
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 18 Benzene, 1,2,4-trimethyl- Concentration Rank 36

| R.T.  | EstConc     | Area    | Relative to ISTD       | R.T.  |
|-------|-------------|---------|------------------------|-------|
| 8.410 | 19.37 ng/ul | 1949190 | 1,4-Dichlorobenzene-d4 | 8.286 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1,2,4-trimethyl-  | 120 | C9H12   | 000095-63-6 | 95   |
| 2    |    |   | Benzene, 1,2,3-trimethyl-  | 120 | C9H12   | 000526-73-8 | 95   |
| 3    |    |   | Benzene, 1,2,3-trimethyl-  | 120 | C9H12   | 000526-73-8 | 94   |
| 4    |    |   | Benzene, 1,2,3-trimethyl-  | 120 | C9H12   | 000526-73-8 | 94   |
| 5    |    |   | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 93   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122721\  
Data File : BG051828.D  
Acq On : 28 Dec 2021 14:05  
Operator : CG/JU  
Sample : M5215-01DL 2X  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD3DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

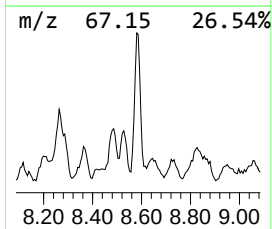
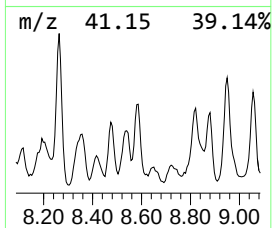
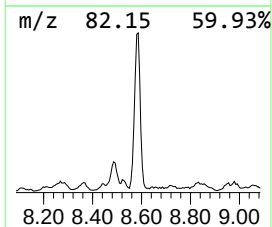
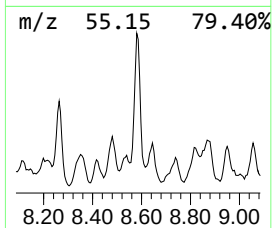
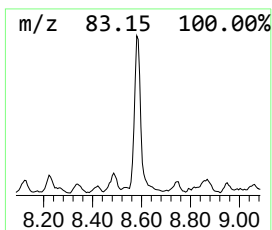
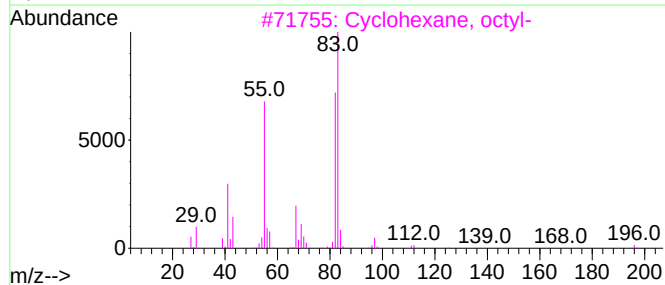
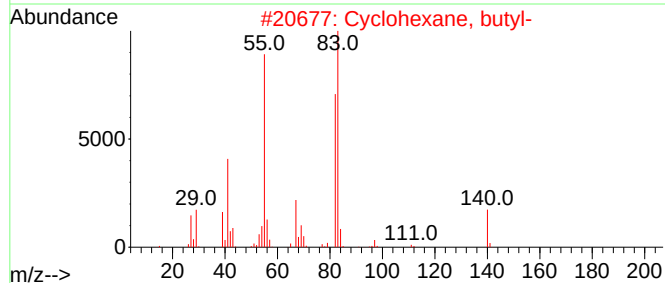
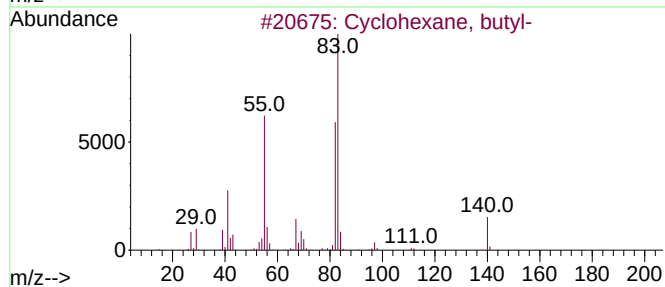
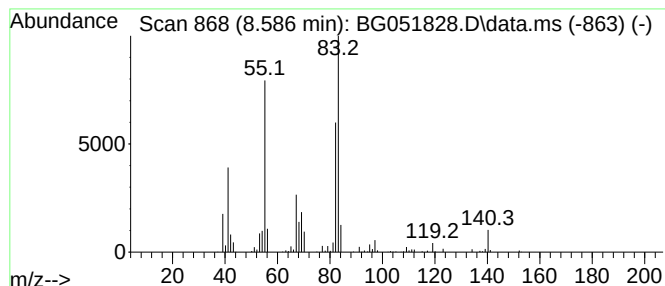
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 19 (DEL) Alkane: Cyclic8.586 Concentration Rank 56

| R.T.  | EstConc     | Area    | Relative to ISTD       | R.T.  |
|-------|-------------|---------|------------------------|-------|
| 8.586 | 10.19 ng/ul | 1025350 | 1,4-Dichlorobenzene-d4 | 8.286 |

| Hit# | of 5 | Tentative ID         | MW  | MolForm | CAS#        | Qual |
|------|------|----------------------|-----|---------|-------------|------|
| 1    |      | Cyclohexane, butyl-  | 140 | C10H20  | 001678-93-9 | 87   |
| 2    |      | Cyclohexane, butyl-  | 140 | C10H20  | 001678-93-9 | 87   |
| 3    |      | Cyclohexane, octyl-  | 196 | C14H28  | 001795-15-9 | 86   |
| 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\BG122721\  
Data File : BG051828.D  
Acq On : 28 Dec 2021 14:05  
Operator : CG/JU  
Sample : M5215-01DL 2X  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD3DL

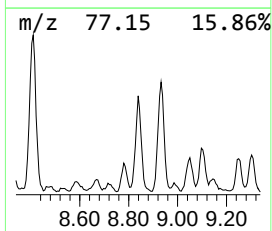
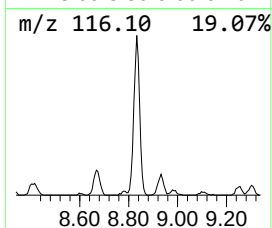
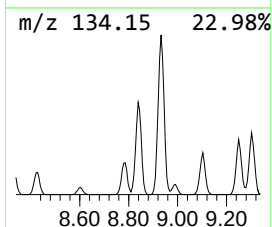
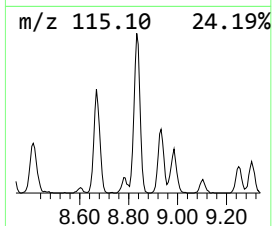
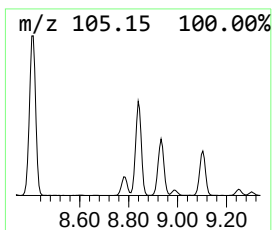
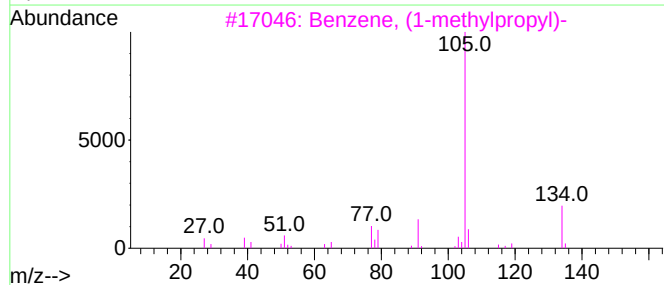
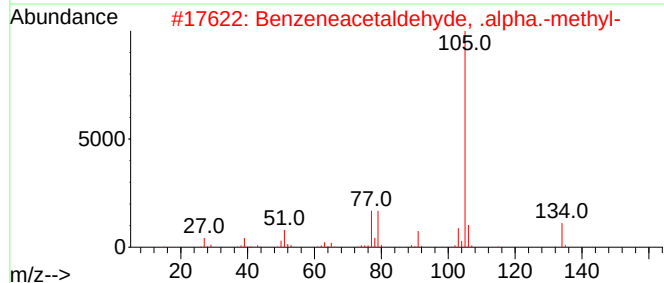
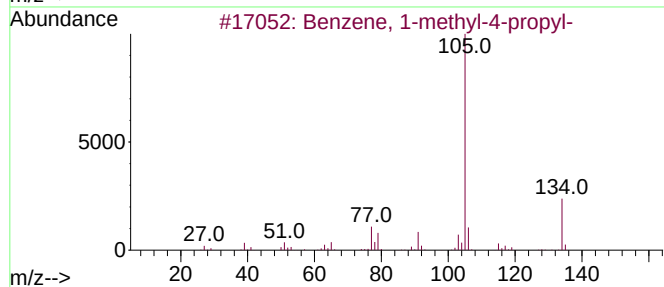
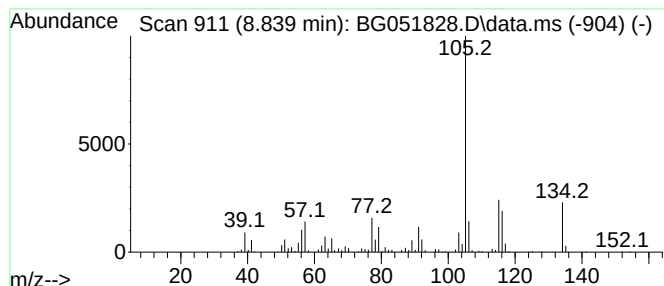
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 20 Benzene, 1-methyl-4-propyl- Concentration Rank 32

| R.T.  | EstConc     | Area    | Relative to ISTD       | R.T.  |
|-------|-------------|---------|------------------------|-------|
| 8.839 | 21.04 ng/ul | 2116840 | 1,4-Dichlorobenzene-d4 | 8.286 |

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



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122721\  
Data File : BG051828.D  
Acq On : 28 Dec 2021 14:05  
Operator : CG/JU  
Sample : M5215-01DL 2X  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD3DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

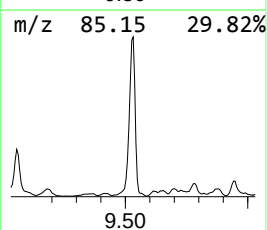
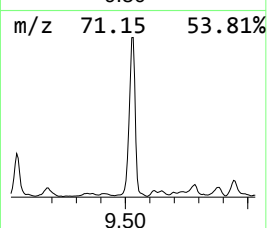
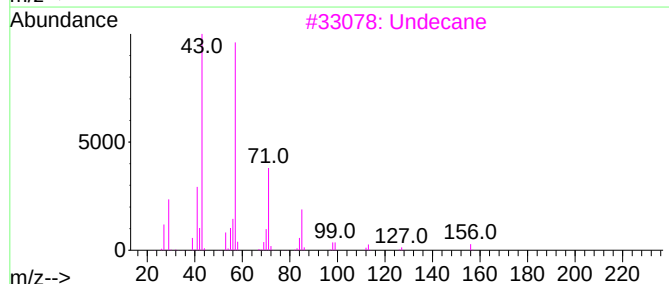
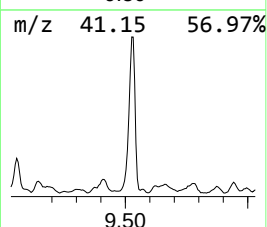
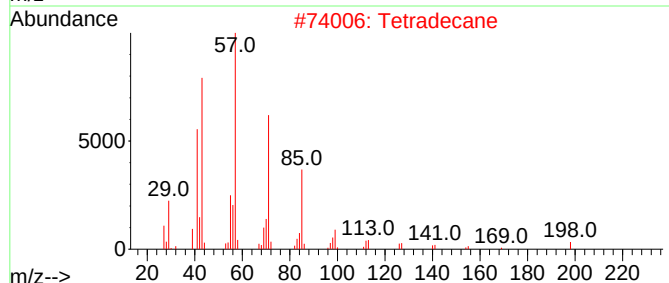
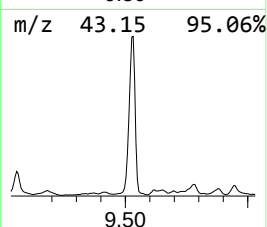
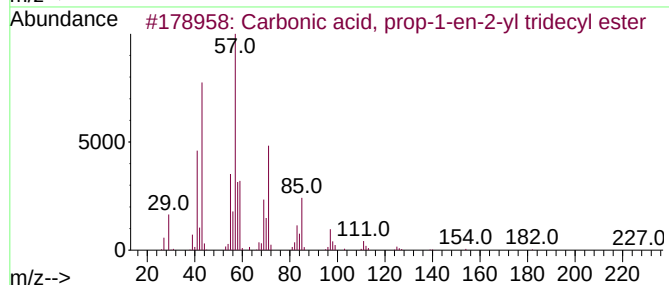
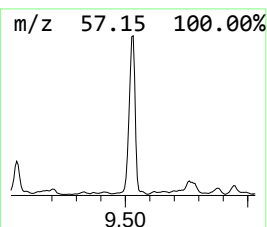
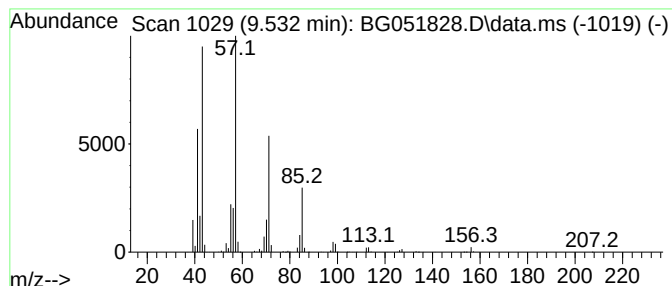
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*

Peak Number 21 Carbonic acid, prop-1-en-2-... Concentration Rank 13

| R.T.  | EstConc     | Area    | Relative to ISTD       | R.T.  |
|-------|-------------|---------|------------------------|-------|
| 9.532 | 50.64 ng/ul | 5094680 | 1,4-Dichlorobenzene-d4 | 8.286 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Carbonic acid, prop-1-en-2-yl tr... | 284 | C17H32O3 | 1000382-90-8 | 90   |
| 2    |    |   | Tetradecane                         | 198 | C14H30   | 000629-59-4  | 72   |
| 3    |    |   | Undecane                            | 156 | C11H24   | 001120-21-4  | 70   |
| 4    |    |   | Decane, 3,8-dimethyl-               | 170 | C12H26   | 017312-55-9  | 59   |
| 5    |    |   | 3-Ethyl-3-methylheptane             | 142 | C10H22   | 017302-01-1  | 59   |





Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122721\  
Data File : BG051828.D  
Acq On : 28 Dec 2021 14:05  
Operator : CG/JU  
Sample : M5215-01DL 2X  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

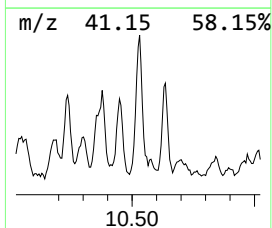
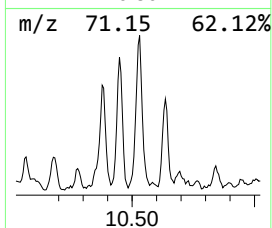
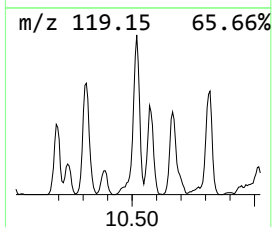
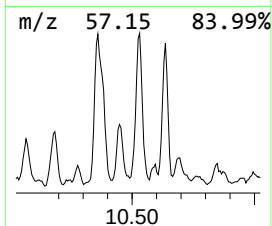
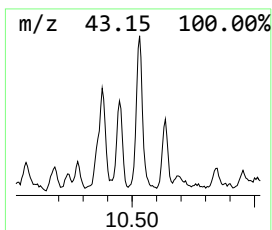
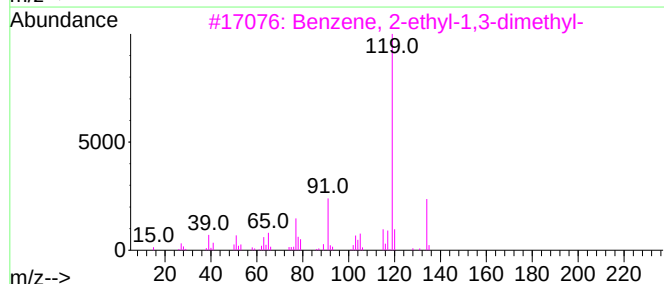
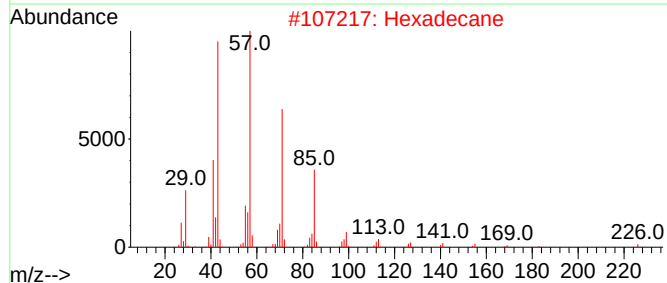
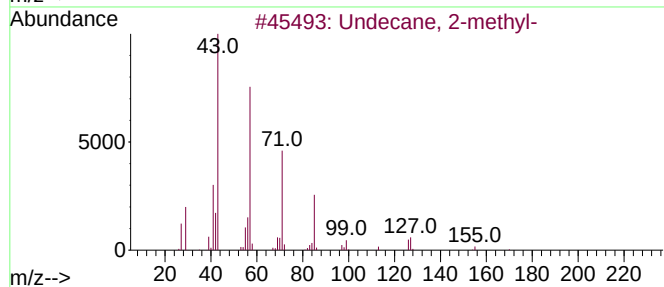
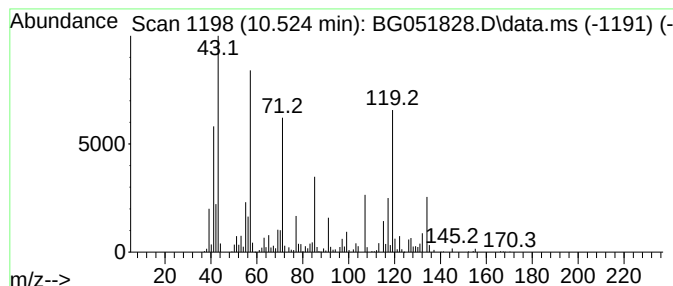
Instrument :  
BNA\_G  
ClientSampleId :  
BGLD3DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 23 (DEL) Alkane: Straight-Chai... Concentration Rank 60

| R.T.      | EstConc                             | Area    | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------------|---------|------------------|-------------|------|
| 10.525    | 7.62 ng/ul                          | 1150450 | Naphthalene-d8   | 11.124      |      |
| Hit# of 5 | Tentative ID                        | MW      | MolForm          | CAS#        | Qual |
| 1         | Undecane, 2-methyl-                 | 170     | C12H26           | 007045-71-8 | 50   |
| 2         | Hexadecane                          | 226     | C16H34           | 000544-76-3 | 38   |
| 3         | Benzene, 2-ethyl-1,3-dimethyl-      | 134     | C10H14           | 002870-04-4 | 35   |
| 4         | p-Cymene                            | 134     | C10H14           | 000099-87-6 | 35   |
| 5         | Benzene, 1-methyl-3-(1-methyleth... | 134     | C10H14           | 000535-77-3 | 35   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122721\  
Data File : BG051828.D  
Acq On : 28 Dec 2021 14:05  
Operator : CG/JU  
Sample : M5215-01DL 2X  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD3DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

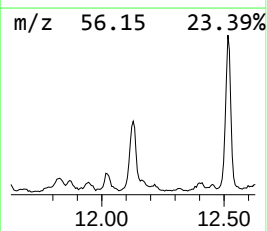
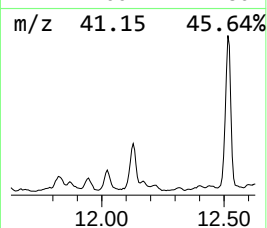
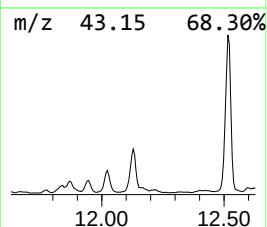
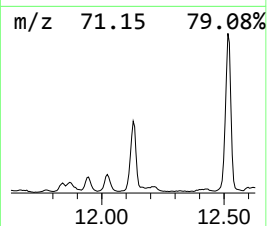
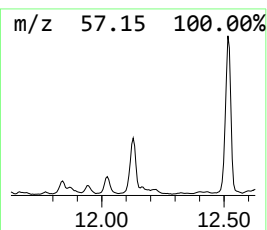
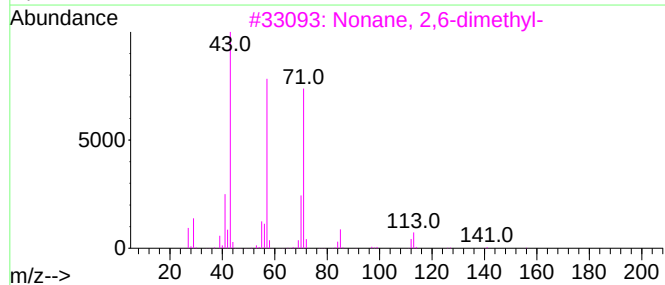
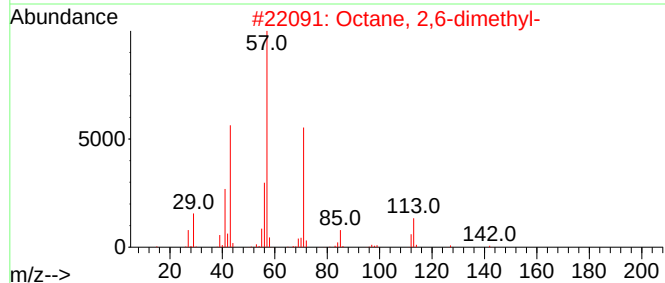
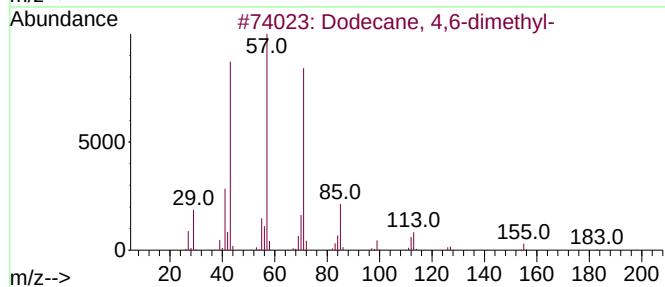
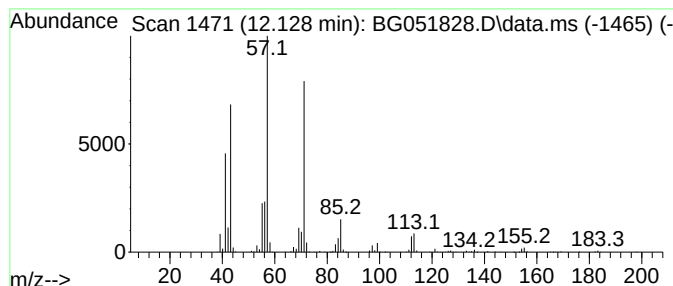
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 24 (DEL) Alkane: Straight-Chai... Concentration Rank 59

| R.T.   | EstConc    | Area    | Relative to ISTD | R.T.   |
|--------|------------|---------|------------------|--------|
| 12.128 | 7.82 ng/ul | 1179990 | Naphthalene-d8   | 11.124 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|------|-------------------------------------|-----|-----------|--------------|------|
| 1    |      | Dodecane, 4,6-dimethyl-             | 198 | C14H30    | 061141-72-8  | 91   |
| 2    |      | Octane, 2,6-dimethyl-               | 142 | C10H22    | 002051-30-1  | 90   |
| 3    |      | Nonane, 2,6-dimethyl-               | 156 | C11H24    | 017302-28-2  | 80   |
| 4    |      | Sulfurous acid, 2-ethylhexyl non... | 320 | C17H36O3S | 1010309-19-2 | 78   |
| 5    |      | Octane, 2,3,7-trimethyl-            | 156 | C11H24    | 062016-34-6  | 72   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122721\  
Data File : BG051828.D  
Acq On : 28 Dec 2021 14:05  
Operator : CG/JU  
Sample : M5215-01DL 2X  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

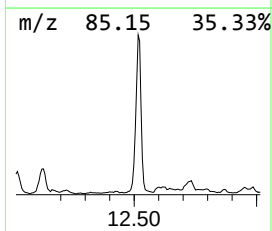
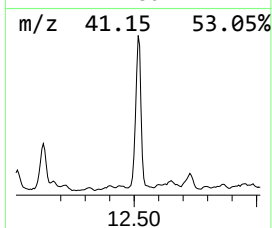
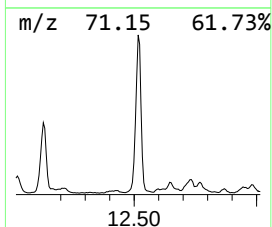
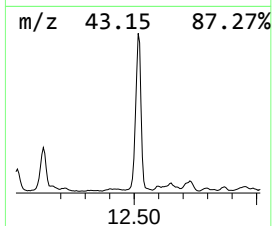
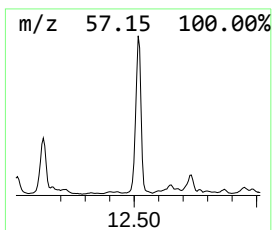
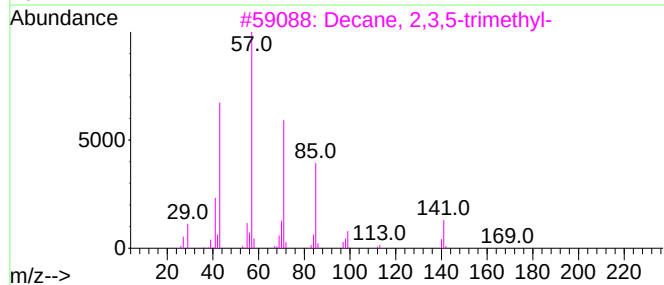
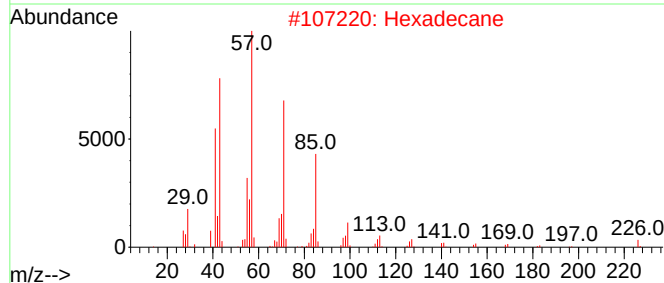
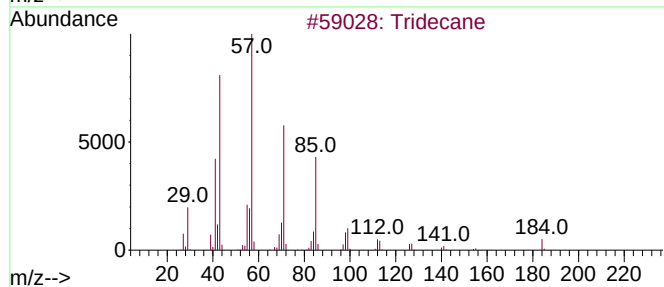
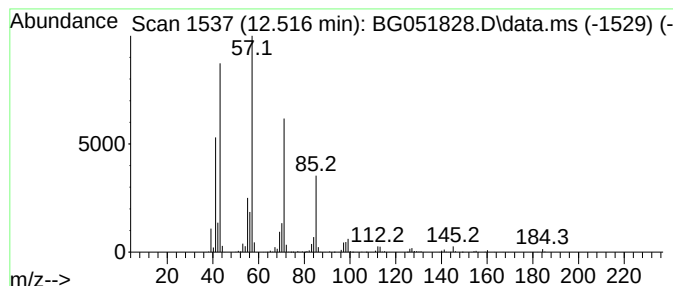
Instrument :  
BNA\_G  
ClientSampleId :  
BGLD3DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 25 (DEL) Alkane: Straight-Chai... Concentration Rank 27

| R.T.      | EstConc                  | Area    | Relative to ISTD | R.T.        |      |
|-----------|--------------------------|---------|------------------|-------------|------|
| 12.516    | 24.53 ng/ul              | 3701260 | Naphthalene-d8   | 11.124      |      |
| Hit# of 5 | Tentative ID             | MW      | MolForm          | CAS#        | Qual |
| 1         | Tridecane                | 184     | C13H28           | 000629-50-5 | 93   |
| 2         | Hexadecane               | 226     | C16H34           | 000544-76-3 | 91   |
| 3         | Decane, 2,3,5-trimethyl- | 184     | C13H28           | 062238-11-3 | 90   |
| 4         | Hexadecane               | 226     | C16H34           | 000544-76-3 | 86   |
| 5         | Tetradecane              | 198     | C14H30           | 000629-59-4 | 86   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122721\  
Data File : BG051828.D  
Acq On : 28 Dec 2021 14:05  
Operator : CG/JU  
Sample : M5215-01DL 2X  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD3DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

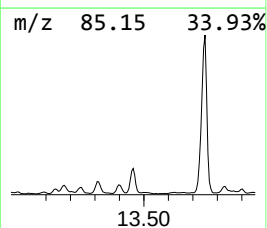
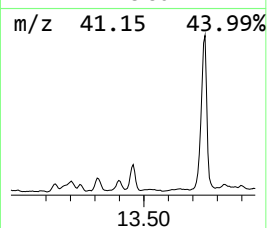
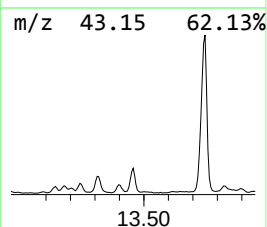
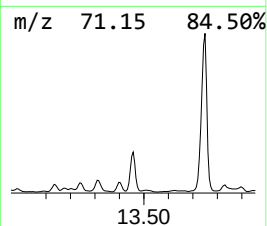
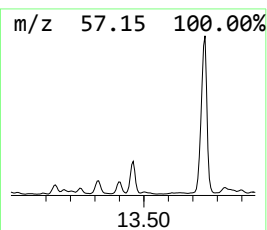
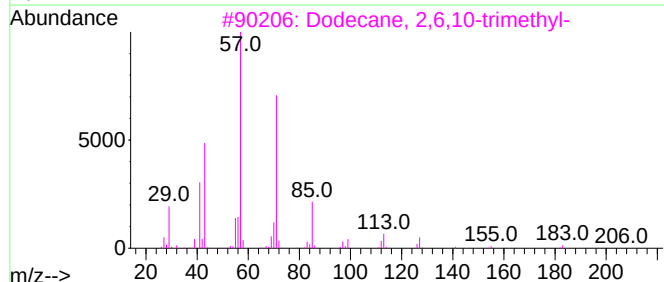
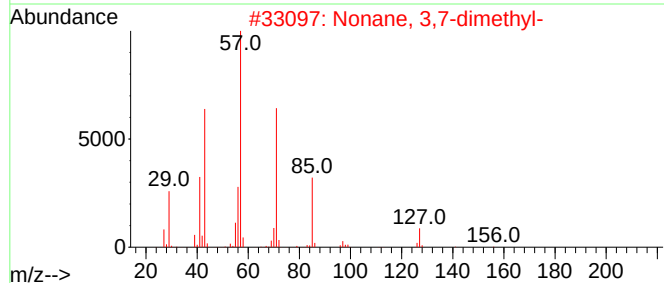
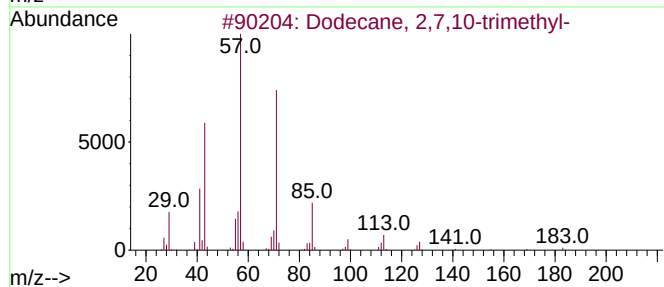
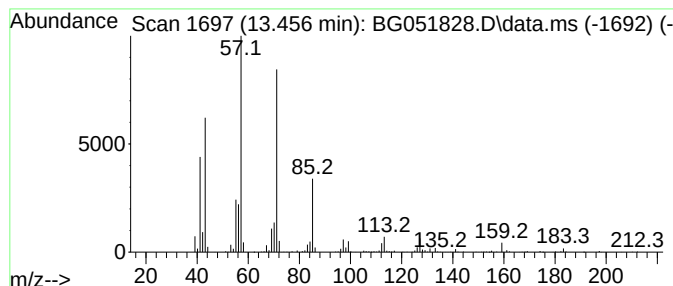
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*

Peak Number 27 (DEL) Alkane: Straight-Chai... Concentration Rank 35

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 13.456 | 19.78 ng/ul | 1467640 | Acenaphthene-d10 | 14.919 |

| Hit# | of | 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------------|-----|---------|-------------|------|
| 1    |    |   | Dodecane, 2,7,10-trimethyl- | 212 | C15H32  | 074645-98-0 | 90   |
| 2    |    |   | Nonane, 3,7-dimethyl-       | 156 | C11H24  | 017302-32-8 | 87   |
| 3    |    |   | Dodecane, 2,6,10-trimethyl- | 212 | C15H32  | 003891-98-3 | 86   |
| 4    |    |   | Dodecane, 2,6,11-trimethyl- | 212 | C15H32  | 031295-56-4 | 83   |
| 5    |    |   | Dodecane, 2,6,11-trimethyl- | 212 | C15H32  | 031295-56-4 | 80   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122721\  
Data File : BG051828.D  
Acq On : 28 Dec 2021 14:05  
Operator : CG/JU  
Sample : M5215-01DL 2X  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD3DL

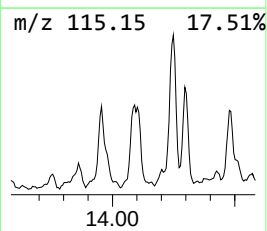
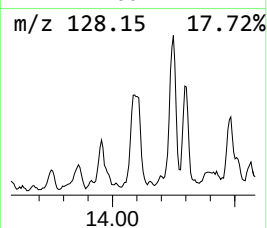
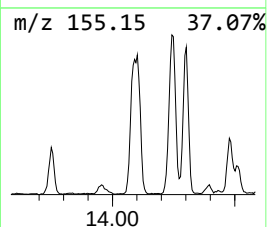
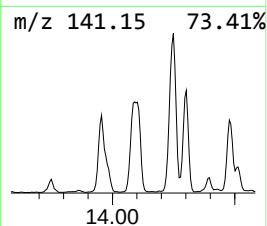
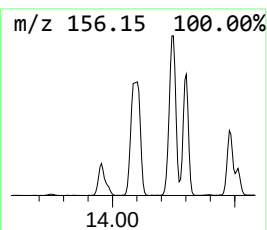
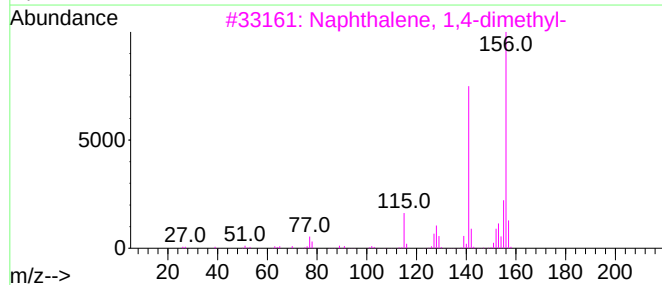
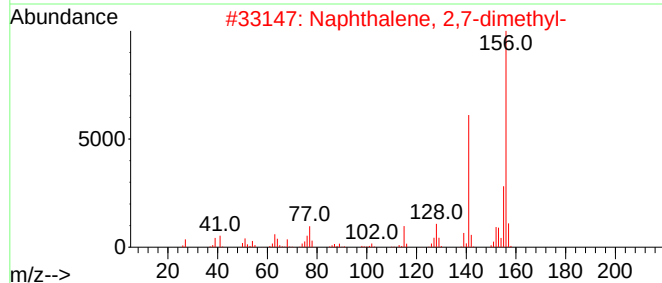
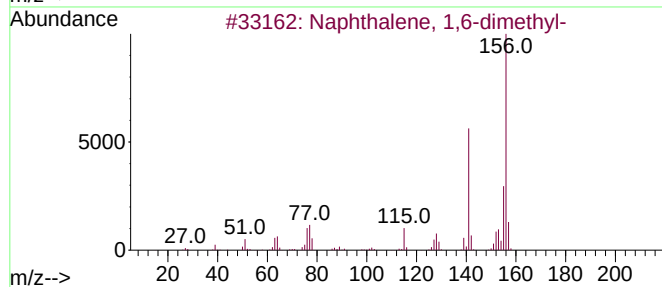
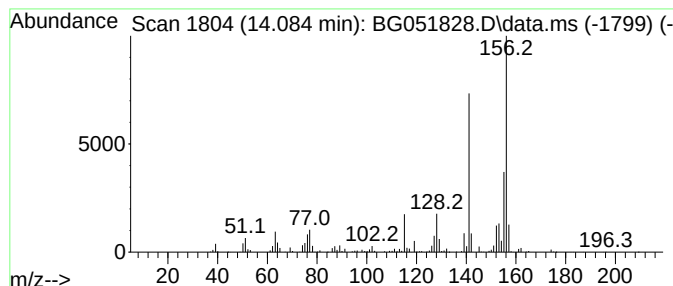
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 29 Naphthalene, 1,6-dimethyl- Concentration Rank 24

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 14.084 | 27.58 ng/ul | 2047180 | Acenaphthene-d10 | 14.919 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, 1,6-dimethyl- | 156 | C12H12  | 000575-43-9 | 97   |
| 2    |    |   | Naphthalene, 2,7-dimethyl- | 156 | C12H12  | 000582-16-1 | 97   |
| 3    |    |   | Naphthalene, 1,4-dimethyl- | 156 | C12H12  | 000571-58-4 | 97   |
| 4    |    |   | Naphthalene, 1,4-dimethyl- | 156 | C12H12  | 000571-58-4 | 97   |
| 5    |    |   | Naphthalene, 1,6-dimethyl- | 156 | C12H12  | 000575-43-9 | 96   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122721\  
Data File : BG051828.D  
Acq On : 28 Dec 2021 14:05  
Operator : CG/JU  
Sample : M5215-01DL 2X  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

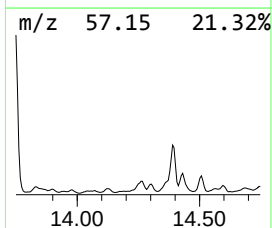
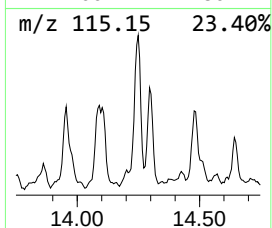
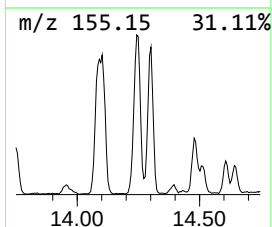
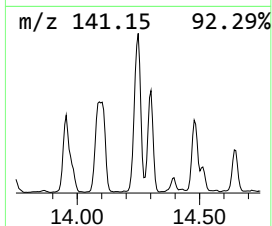
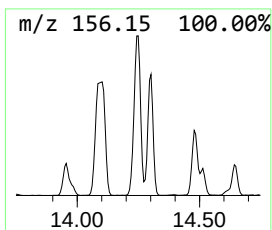
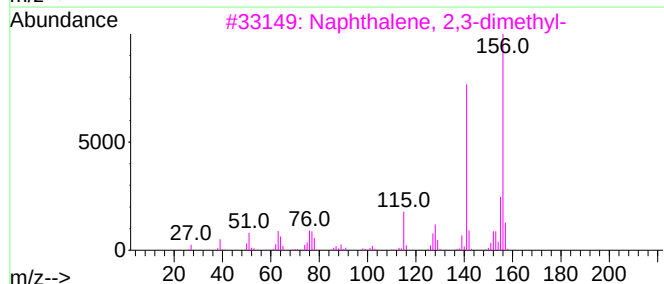
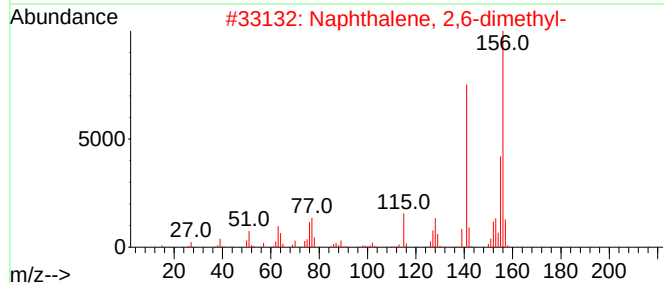
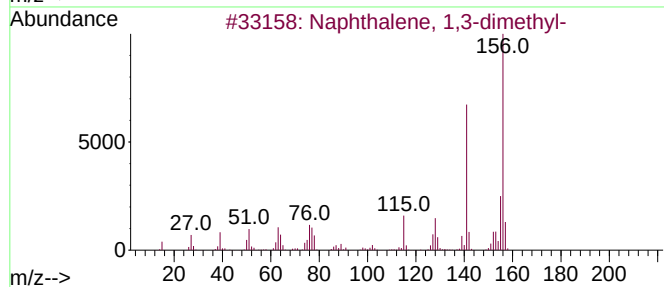
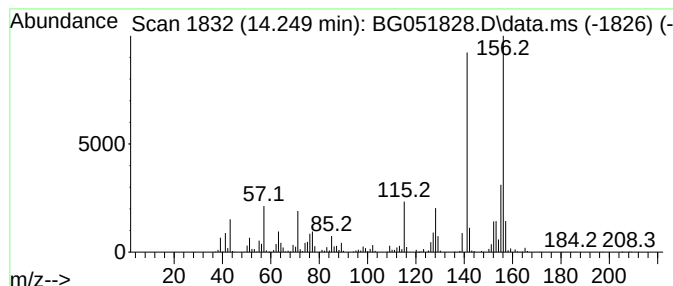
Instrument :  
BNA\_G  
ClientSampleId :  
BGLD3DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 30 Naphthalene, 1,3-dimethyl- Concentration Rank 22

| R.T.      | EstConc                    | Area    | Relative to ISTD | R.T.        |      |
|-----------|----------------------------|---------|------------------|-------------|------|
| 14.249    | 29.00 ng/ul                | 2152040 | Acenaphthene-d10 | 14.919      |      |
| Hit# of 5 | Tentative ID               | MW      | MolForm          | CAS#        | Qual |
| 1         | Naphthalene, 1,3-dimethyl- | 156     | C12H12           | 000575-41-7 | 97   |
| 2         | Naphthalene, 2,6-dimethyl- | 156     | C12H12           | 000581-42-0 | 96   |
| 3         | Naphthalene, 2,3-dimethyl- | 156     | C12H12           | 000581-40-8 | 96   |
| 4         | Naphthalene, 1,4-dimethyl- | 156     | C12H12           | 000571-58-4 | 96   |
| 5         | Naphthalene, 1,6-dimethyl- | 156     | C12H12           | 000575-43-9 | 96   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122721\  
Data File : BG051828.D  
Acq On : 28 Dec 2021 14:05  
Operator : CG/JU  
Sample : M5215-01DL 2X  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD3DL

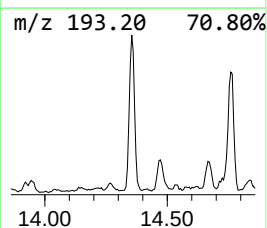
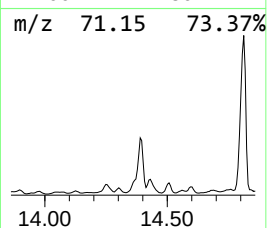
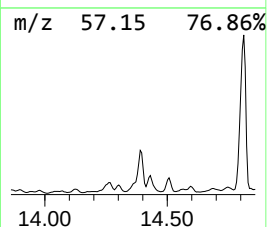
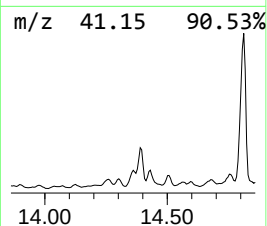
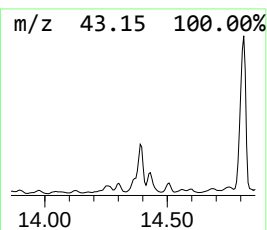
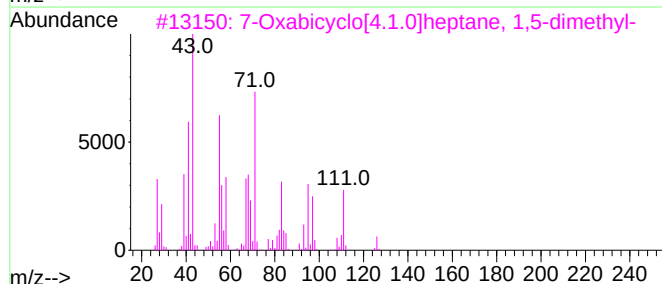
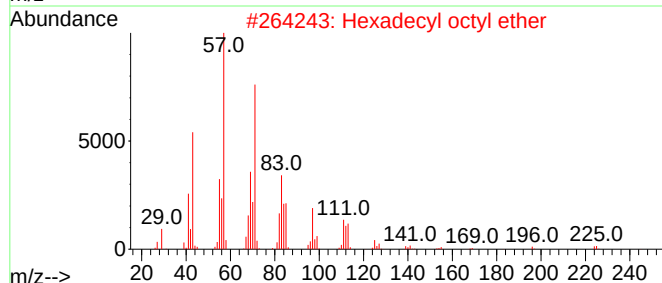
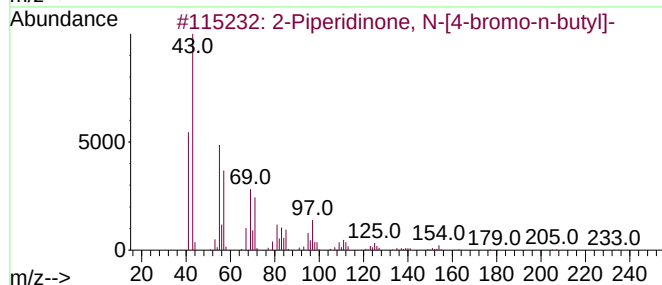
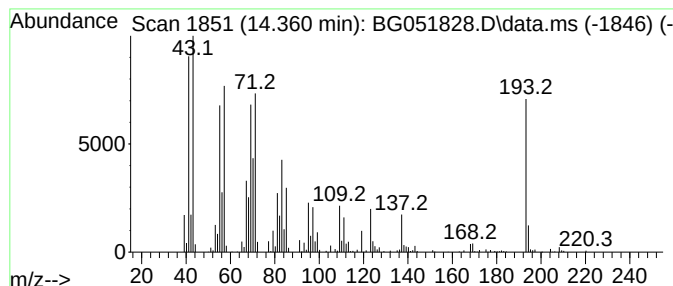
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 31 unknown-01 Concentration Rank 37

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 14.361 | 17.06 ng/ul | 1266390 | Acenaphthene-d10 | 14.919 |

| Hit# | of                                  | 5   | Tentative ID | MW           | MolForm | CAS# | Qual |
|------|-------------------------------------|-----|--------------|--------------|---------|------|------|
| 1    | 2-Piperidinone, N-[4-bromo-n-but... | 233 | C9H16BrNO    | 195194-80-0  | 49      |      |      |
| 2    | Hexadecyl octyl ether               | 354 | C24H50O      | 1000406-38-6 | 46      |      |      |
| 3    | 7-Oxabicyclo[4.1.0]heptane, 1,5-... | 126 | C8H14O       | 162239-52-3  | 46      |      |      |
| 4    | Nonane, 1-bromo-                    | 206 | C9H19Br      | 000693-58-3  | 30      |      |      |
| 5    | 2,6,10-Trimethylundeca-1,3-diene    | 194 | C14H26       | 1000222-09-1 | 25      |      |      |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122721\  
Data File : BG051828.D  
Acq On : 28 Dec 2021 14:05  
Operator : CG/JU  
Sample : M5215-01DL 2X  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

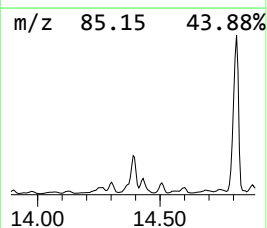
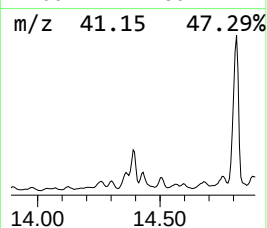
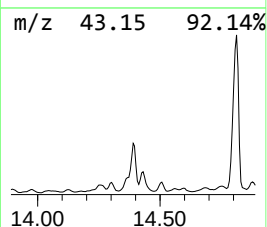
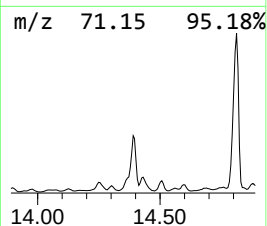
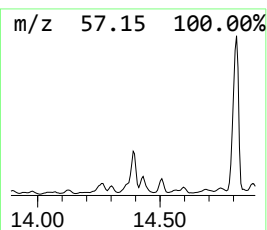
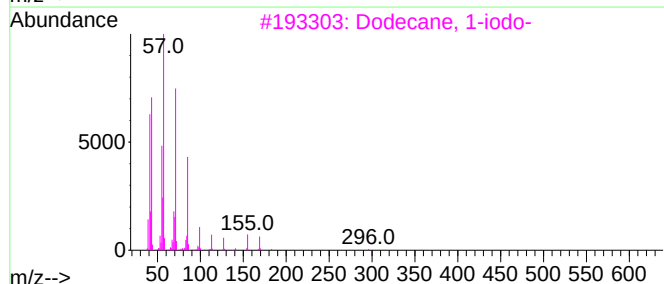
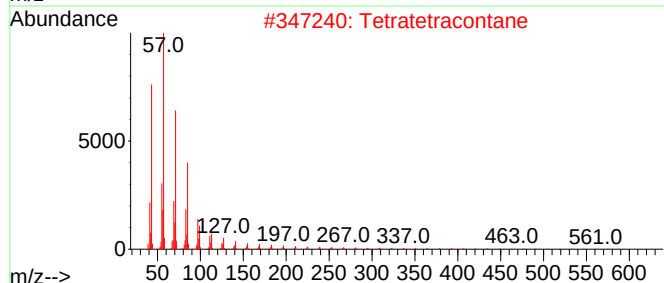
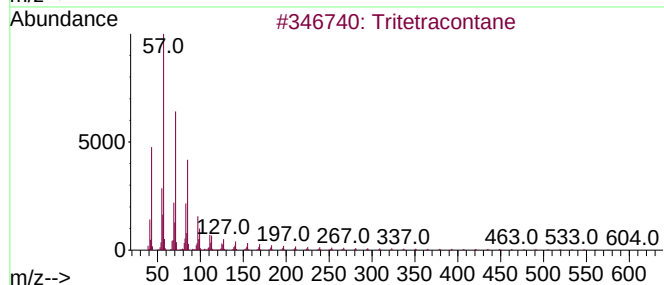
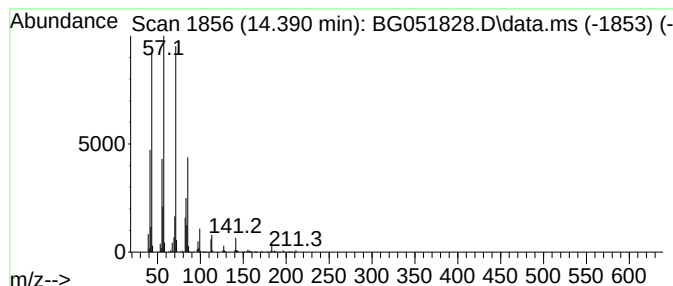
Instrument :  
BNA\_G  
ClientSampleId :  
BGLD3DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 32 (DEL) Alkane: Straight-Chai... Concentration Rank 23

| R.T.      | EstConc             | Area    | Relative to ISTD |         | R.T.        |      |
|-----------|---------------------|---------|------------------|---------|-------------|------|
| 14.390    | 28.81 ng/ul         | 2138260 | Acenaphthene-d10 |         | 14.919      |      |
| Hit# of 5 | Tentative ID        |         | MW               | MolForm | CAS#        | Qual |
| 1         | Tritetracontane     |         | 605              | C43H88  | 007098-21-7 | 90   |
| 2         | Tetratetracontane   |         | 619              | C44H90  | 007098-22-8 | 90   |
| 3         | Dodecane, 1-iodo-   |         | 296              | C12H25I | 004292-19-7 | 80   |
| 4         | Decane, 5-propyl-   |         | 184              | C13H28  | 017312-62-8 | 64   |
| 5         | Dodecane, 4-methyl- |         | 184              | C13H28  | 006117-97-1 | 62   |





Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122721\  
Data File : BG051828.D  
Acq On : 28 Dec 2021 14:05  
Operator : CG/JU  
Sample : M5215-01DL 2X  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD3DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

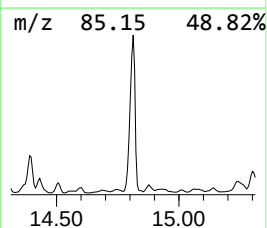
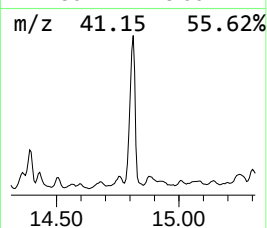
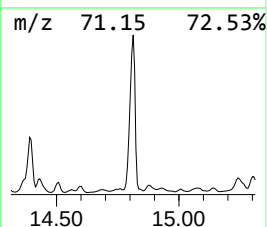
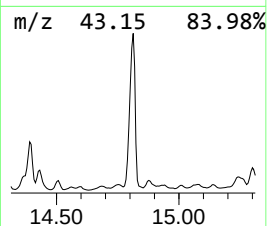
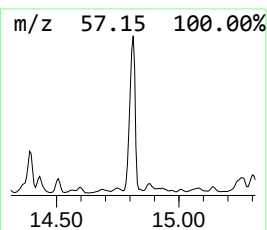
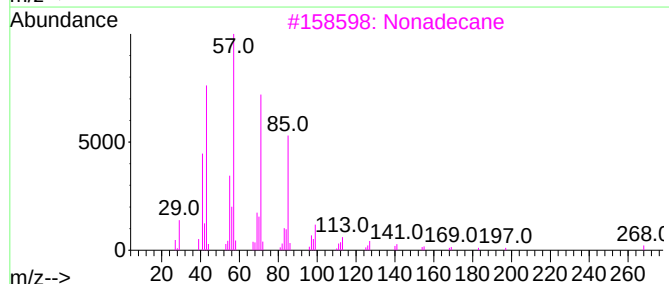
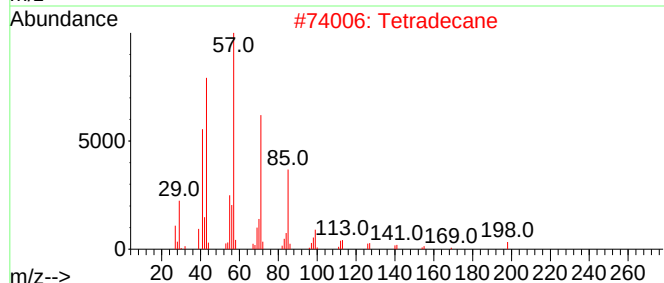
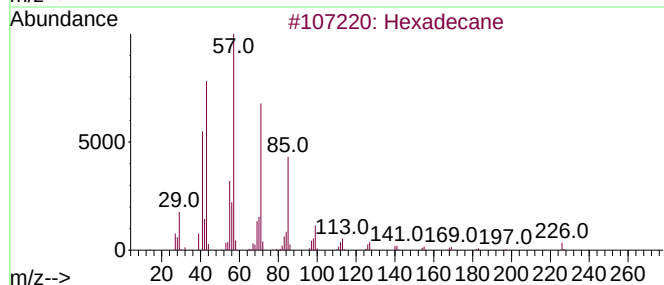
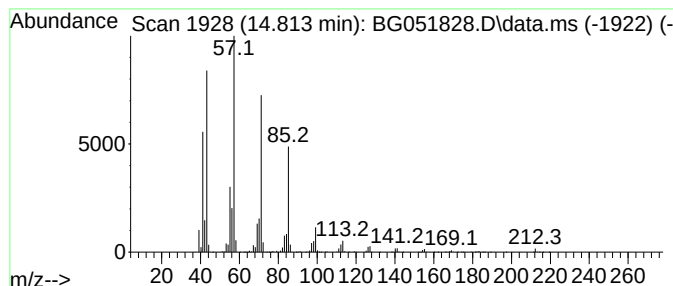
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 33 (DEL) Alkane: Straight-Chai... Concentration Rank 5

| R.T.   | EstConc      | Area    | Relative to ISTD | R.T.   |
|--------|--------------|---------|------------------|--------|
| 14.813 | 119.43 ng/ul | 8863280 | Acenaphthene-d10 | 14.919 |

| Hit# | of 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|------|--------------|-----|---------|-------------|------|
| 1    |      | Hexadecane   | 226 | C16H34  | 000544-76-3 | 91   |
| 2    |      | Tetradecane  | 198 | C14H30  | 000629-59-4 | 91   |
| 3    |      | Nonadecane   | 268 | C19H40  | 000629-92-5 | 90   |
| 4    |      | Nonadecane   | 268 | C19H40  | 000629-92-5 | 90   |
| 5    |      | Tridecane    | 184 | C13H28  | 000629-50-5 | 90   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122721\  
Data File : BG051828.D  
Acq On : 28 Dec 2021 14:05  
Operator : CG/JU  
Sample : M5215-01DL 2X  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD3DL

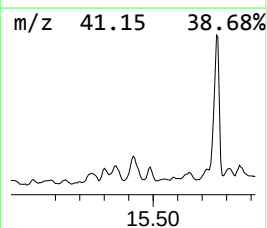
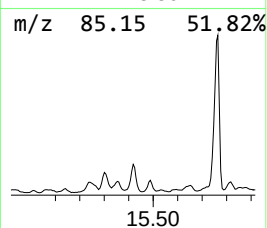
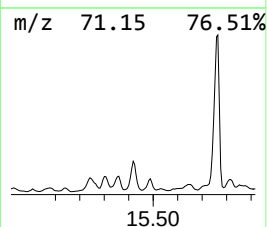
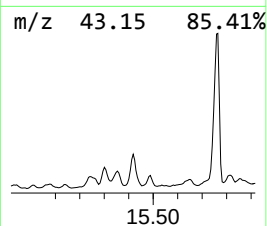
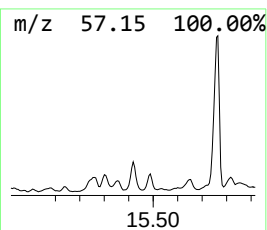
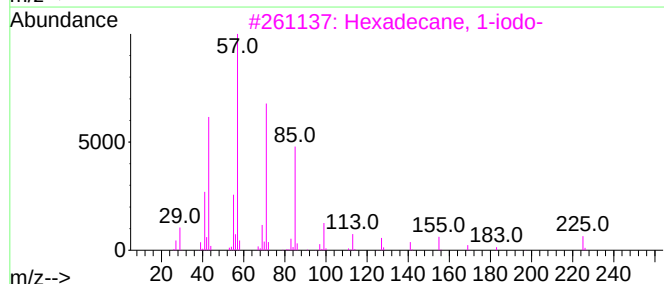
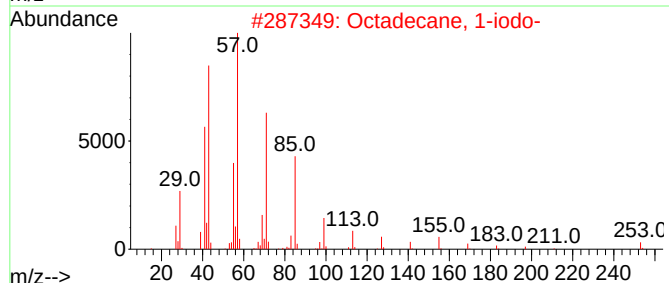
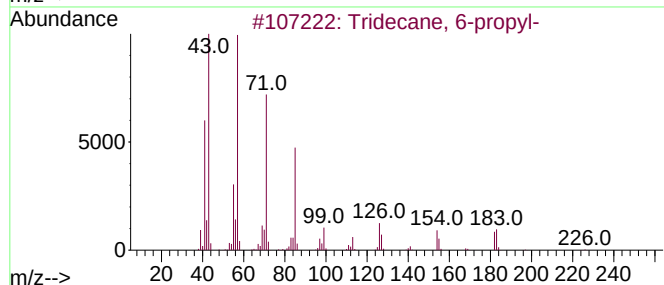
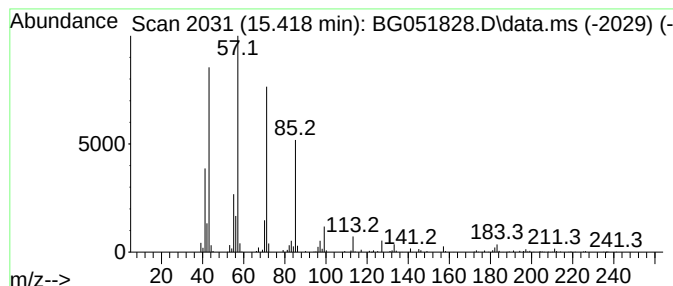
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*

Peak Number 34 (DEL) Alkane: Straight-Chai... Concentration Rank 28

| R.T.      | EstConc              | Area    | Relative to ISTD | R.T.        |      |
|-----------|----------------------|---------|------------------|-------------|------|
| 15.418    | 23.48 ng/ul          | 1742220 | Acenaphthene-d10 | 14.919      |      |
| Hit# of 5 | Tentative ID         | MW      | MolForm          | CAS#        | Qual |
| 1         | Tridecane, 6-propyl- | 226     | C16H34           | 055045-10-8 | 91   |
| 2         | Octadecane, 1-iodo-  | 380     | C18H37I          | 000629-93-6 | 90   |
| 3         | Hexadecane, 1-iodo-  | 352     | C16H33I          | 000544-77-4 | 90   |
| 4         | Nonadecane           | 268     | C19H40           | 000629-92-5 | 86   |
| 5         | Docosane             | 310     | C22H46           | 000629-97-0 | 83   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122721\  
Data File : BG051828.D  
Acq On : 28 Dec 2021 14:05  
Operator : CG/JU  
Sample : M5215-01DL 2X  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD3DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

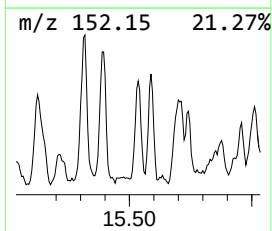
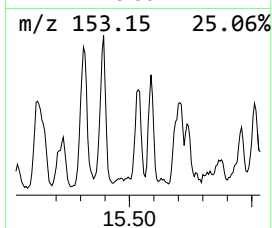
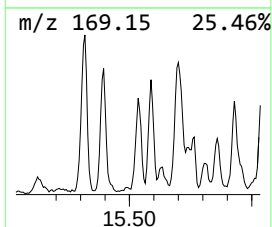
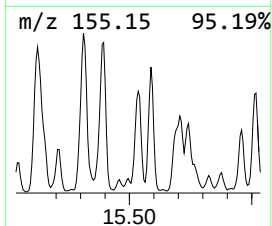
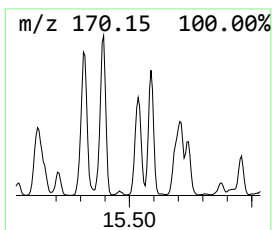
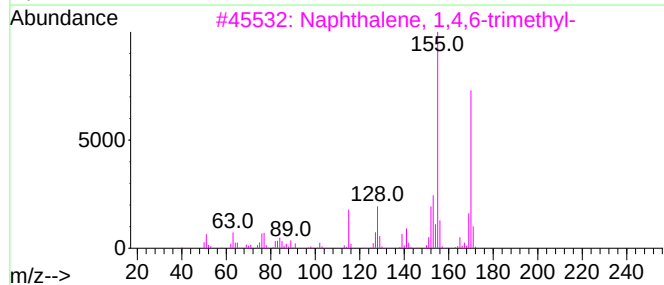
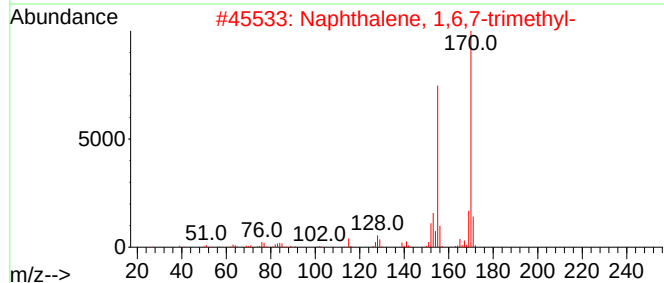
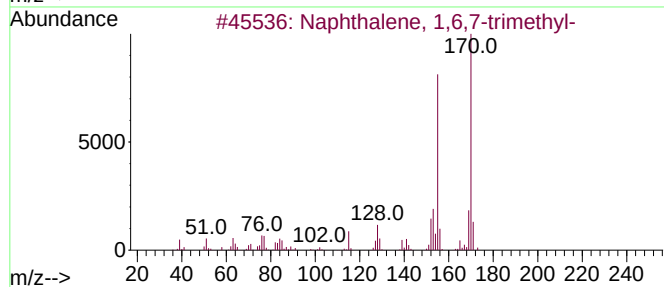
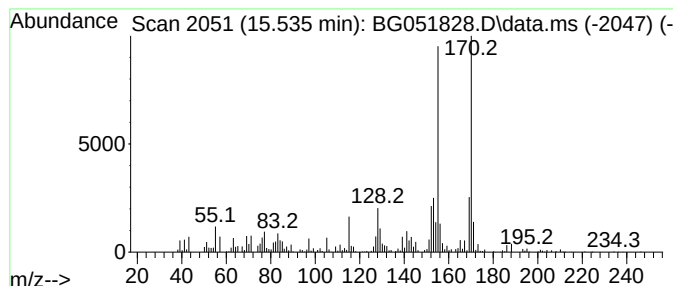
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 35 Naphthalene, 1,6,7-trimethyl- Concentration Rank 49

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 15.535 | 12.76 ng/ul | 946615 | Acenaphthene-d10 | 14.919 |

| Hit# | of | 5 | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, 1,6,7-trimethyl- | 170 | C13H14  | 002245-38-7 | 97   |
| 2    |    |   | Naphthalene, 1,6,7-trimethyl- | 170 | C13H14  | 002245-38-7 | 97   |
| 3    |    |   | Naphthalene, 1,4,6-trimethyl- | 170 | C13H14  | 002131-42-2 | 97   |
| 4    |    |   | Naphthalene, 2,3,6-trimethyl- | 170 | C13H14  | 000829-26-5 | 97   |
| 5    |    |   | Naphthalene, 2,3,6-trimethyl- | 170 | C13H14  | 000829-26-5 | 96   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122721\  
Data File : BG051828.D  
Acq On : 28 Dec 2021 14:05  
Operator : CG/JU  
Sample : M5215-01DL 2X  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

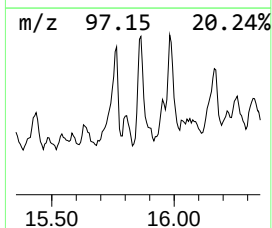
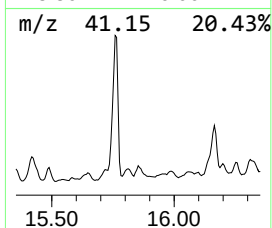
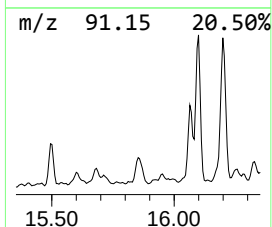
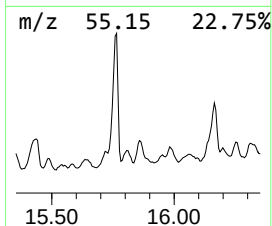
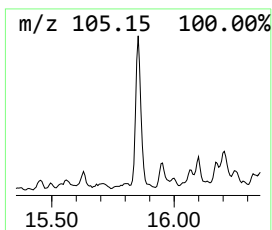
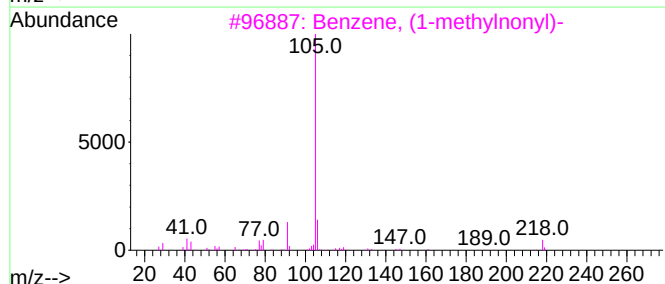
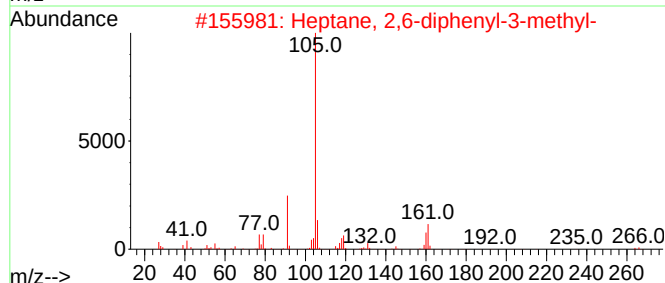
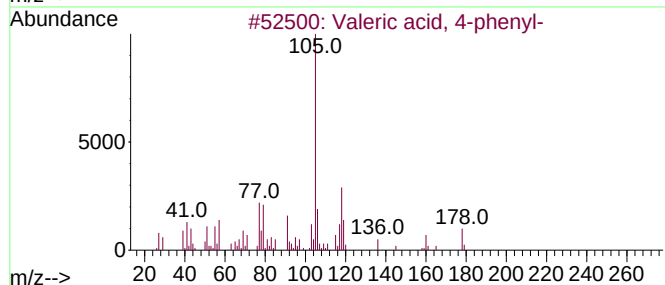
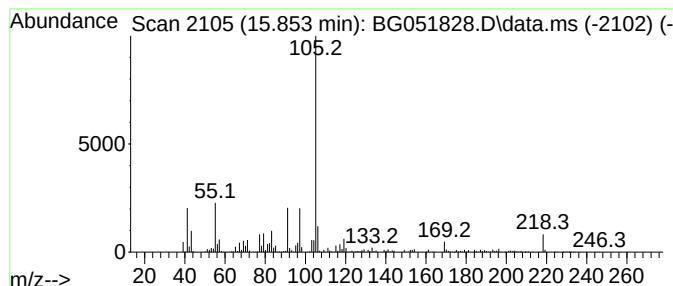
Instrument :  
BNA\_G  
ClientSampleId :  
BGLD3DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 37 unknown-02 Concentration Rank 33

| R.T.      | EstConc                         | Area    | Relative to ISTD |          |              | R.T.   |
|-----------|---------------------------------|---------|------------------|----------|--------------|--------|
| 15.853    | 20.60 ng/ul                     | 1528800 | Acenaphthene-d10 |          |              | 14.919 |
| Hit# of 5 | Tentative ID                    |         | MW               | MolForm  | CAS#         | Qual   |
| 1         | Valeric acid, 4-phenyl-         |         | 178              | C11H14O2 | 016433-43-5  | 43     |
| 2         | Heptane, 2,6-diphenyl-3-methyl- |         | 266              | C20H26   | 1000161-22-5 | 38     |
| 3         | Benzene, (1-methylnonyl)-       |         | 218              | C16H26   | 004537-13-7  | 38     |
| 4         | Benzene, (1-methylnonyl)-       |         | 218              | C16H26   | 004537-13-7  | 38     |
| 5         | Benzene, (1-methylnonyl)-       |         | 218              | C16H26   | 004537-13-7  | 35     |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122721\  
Data File : BG051828.D  
Acq On : 28 Dec 2021 14:05  
Operator : CG/JU  
Sample : M5215-01DL 2X  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD3DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

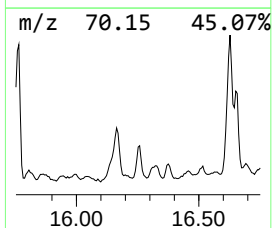
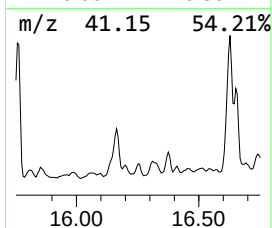
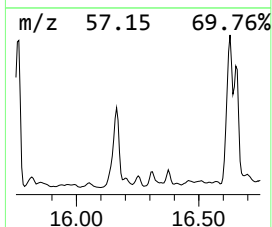
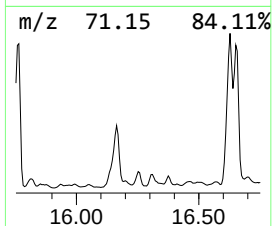
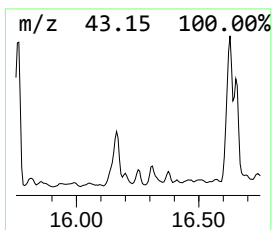
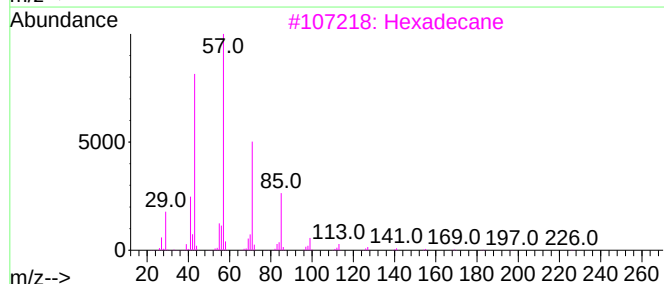
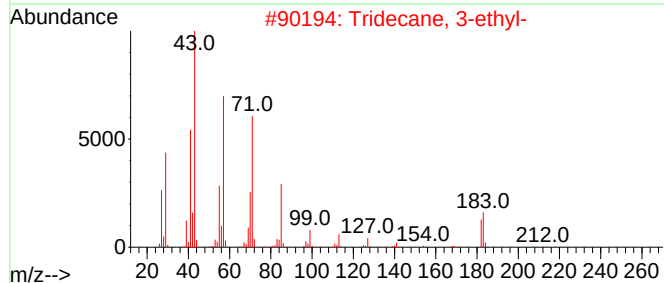
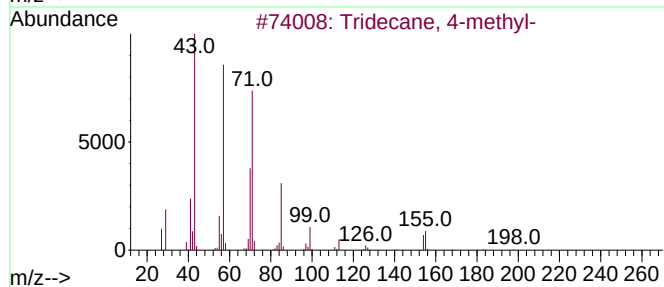
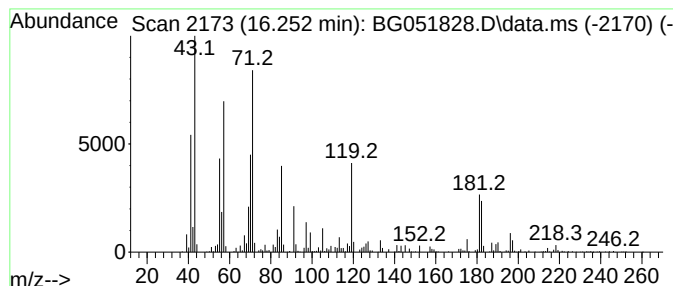
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*

Peak Number 38 (DEL) Alkane: Straight-Chai... Concentration Rank 46

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 16.252 | 14.45 ng/ul | 1072520 | Acenaphthene-d10 | 14.919 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Tridecane, 4-methyl-                | 198 | C14H30    | 026730-12-1  | 60   |
| 2    |    |   | Tridecane, 3-ethyl-                 | 212 | C15H32    | 013286-73-2  | 60   |
| 3    |    |   | Hexadecane                          | 226 | C16H34    | 000544-76-3  | 50   |
| 4    |    |   | Dodecane, 4-methyl-                 | 184 | C13H28    | 006117-97-1  | 47   |
| 5    |    |   | Sulfurous acid, pentyl undecyl e... | 306 | C16H34O3S | 1000309-14-4 | 46   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122721\  
Data File : BG051828.D  
Acq On : 28 Dec 2021 14:05  
Operator : CG/JU  
Sample : M5215-01DL 2X  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD3DL

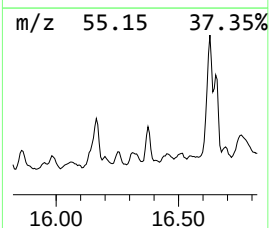
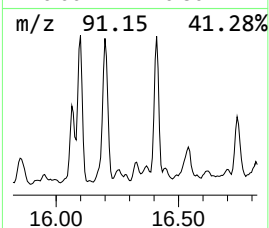
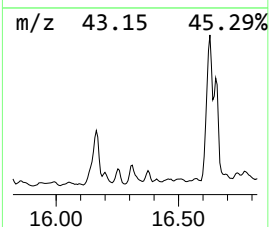
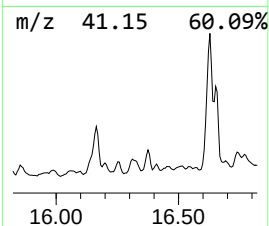
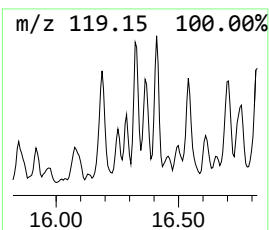
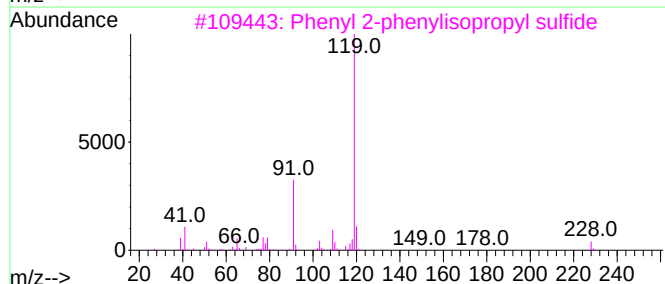
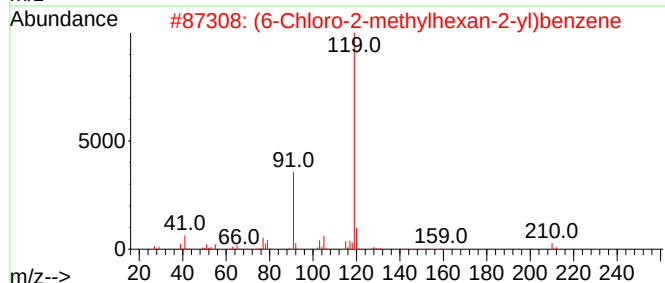
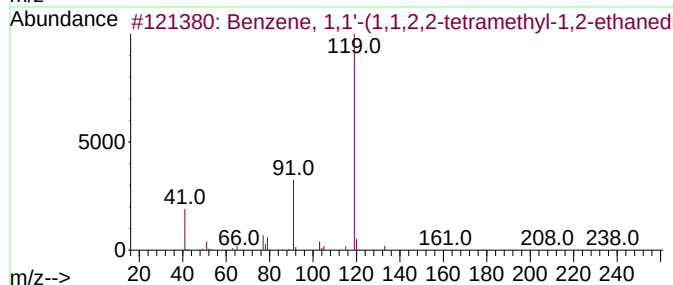
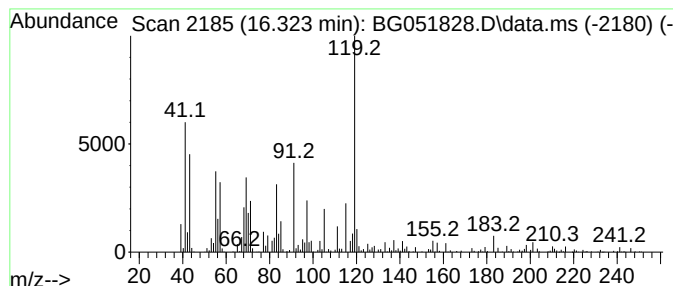
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 39 unknown-03 Concentration Rank 40

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 16.323 | 16.52 ng/ul | 1293050 | Phenanthrene-d10 | 17.680 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|------------|--------------|------|
| 1    |    |   | Benzene, 1,1'-(1,1,2,2-tetrameth... | 238 | C18H22     | 001889-67-4  | 42   |
| 2    |    |   | (6-Chloro-2-methylhexan-2-yl)ben... | 210 | C13H19Cl   | 1417621-45-4 | 42   |
| 3    |    |   | Phenyl 2-phenylisopropyl sulfide    | 228 | C15H16S    | 004148-93-0  | 30   |
| 4    |    |   | 1H-Pyrrolo[1,2-c]imidazole-1,3(2... | 216 | C12H12N2O2 | 002221-09-2  | 30   |
| 5    |    |   | 2,4,4,6-Tetramethyl-6-phenyl-1-h... | 230 | C17H26     | 1000293-27-9 | 27   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122721\  
Data File : BG051828.D  
Acq On : 28 Dec 2021 14:05  
Operator : CG/JU  
Sample : M5215-01DL 2X  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD3DL

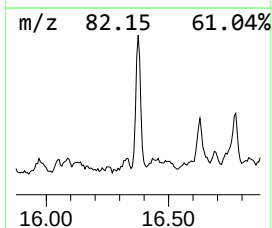
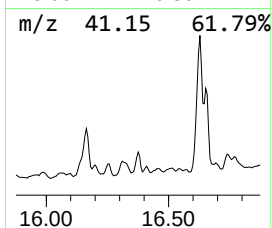
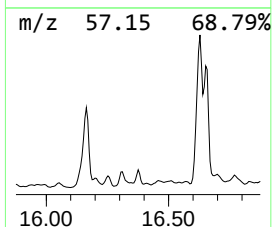
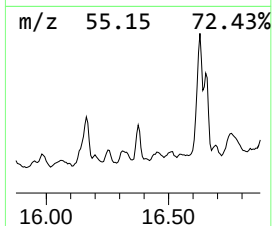
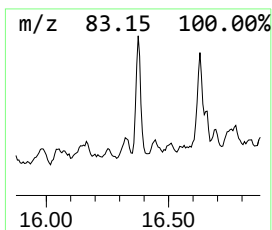
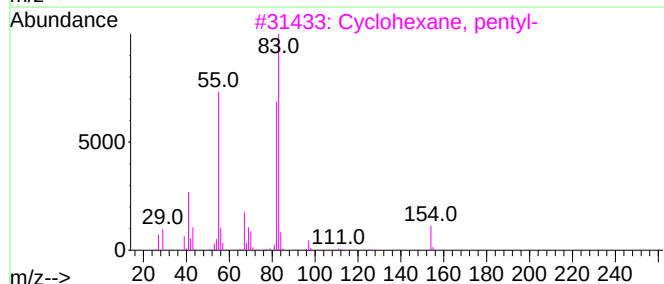
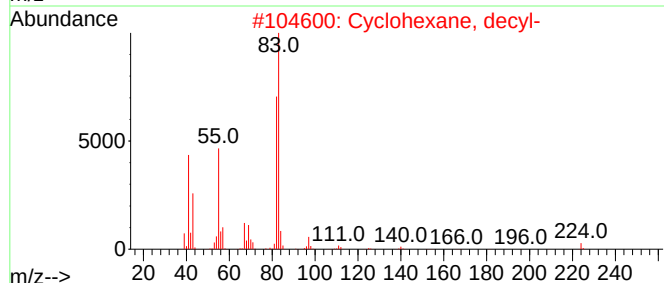
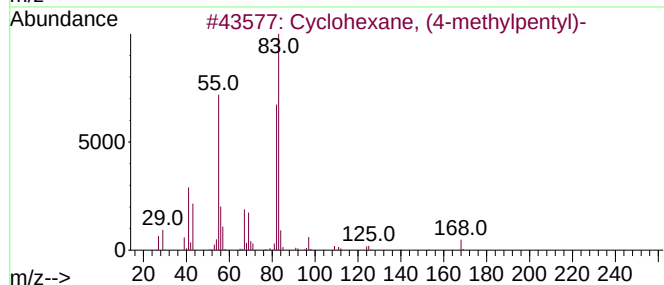
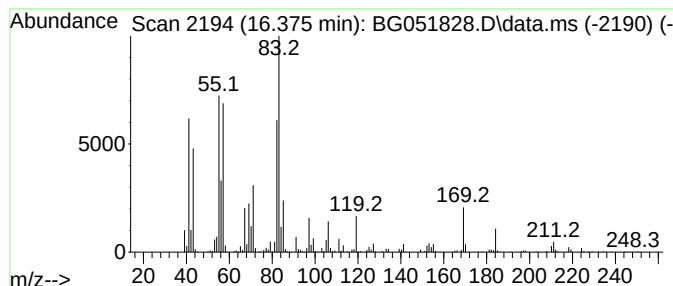
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 40 (DEL) Alkane: Cyclic16.375 Concentration Rank 38

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 16.375 | 16.66 ng/ul | 1304020 | Phenanthrene-d10 | 17.680 |

| Hit# | of | 5 | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------|-----|---------|-------------|------|
| 1    |    |   | Cyclohexane, (4-methylpentyl)- | 168 | C12H24  | 061142-20-9 | 49   |
| 2    |    |   | Cyclohexane, decyl-            | 224 | C16H32  | 001795-16-0 | 49   |
| 3    |    |   | Cyclohexane, pentyl-           | 154 | C11H22  | 004292-92-6 | 47   |
| 4    |    |   | Cyclohexane, pentyl-           | 154 | C11H22  | 004292-92-6 | 47   |
| 5    |    |   | Cyclohexane, butyl-            | 140 | C10H20  | 001678-93-9 | 47   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122721\  
Data File : BG051828.D  
Acq On : 28 Dec 2021 14:05  
Operator : CG/JU  
Sample : M5215-01DL 2X  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD3DL

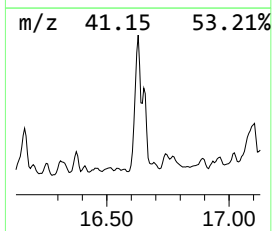
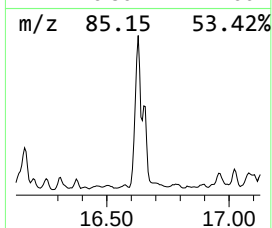
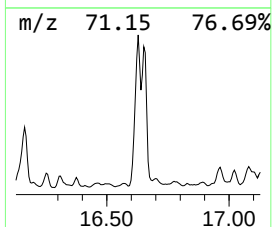
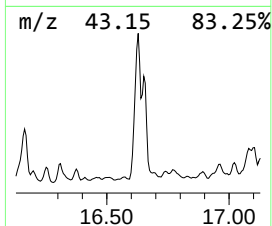
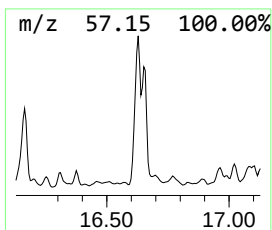
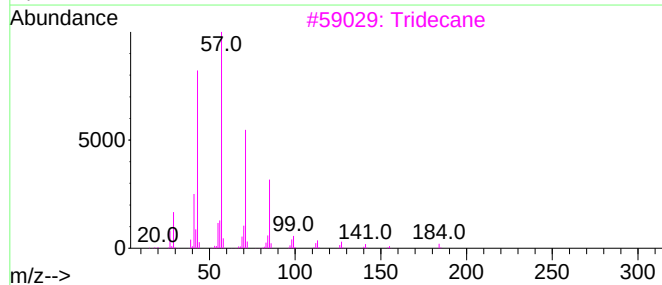
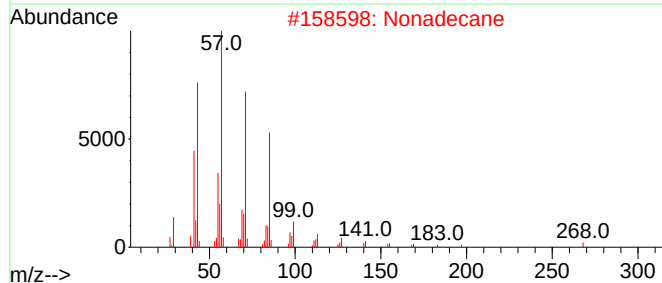
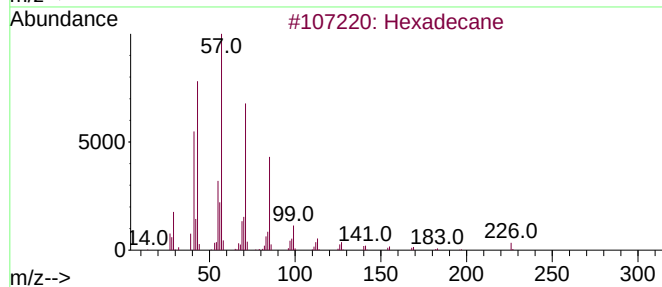
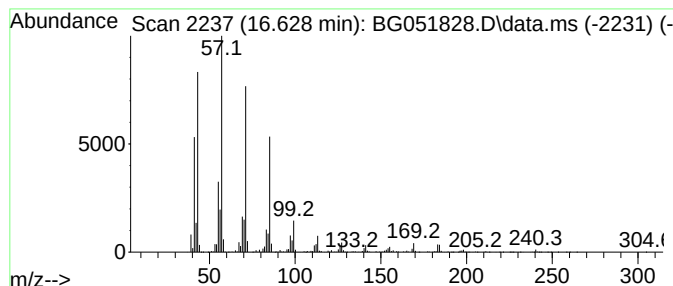
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 42 (DEL) Alkane: Straight-Chai... Concentration Rank 4

| R.T.   | EstConc      | Area     | Relative to ISTD | R.T.   |
|--------|--------------|----------|------------------|--------|
| 16.628 | 129.16 ng/ul | 10107800 | Phenanthrene-d10 | 17.680 |

| Hit# | of 5 | Tentative ID                       | MW  | MolForm  | CAS#         | Qual |
|------|------|------------------------------------|-----|----------|--------------|------|
| 1    |      | Hexadecane                         | 226 | C16H34   | 000544-76-3  | 90   |
| 2    |      | Nonadecane                         | 268 | C19H40   | 000629-92-5  | 90   |
| 3    |      | Tridecane                          | 184 | C13H28   | 000629-50-5  | 90   |
| 4    |      | Carbonic acid, eicosyl vinyl ester | 368 | C23H44O3 | 1000382-54-3 | 87   |
| 5    |      | Pentadecane                        | 212 | C15H32   | 000629-62-9  | 87   |





Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122721\  
Data File : BG051828.D  
Acq On : 28 Dec 2021 14:05  
Operator : CG/JU  
Sample : M5215-01DL 2X  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD3DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

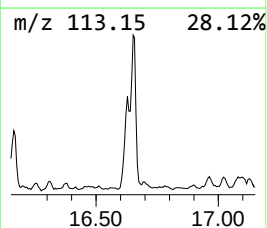
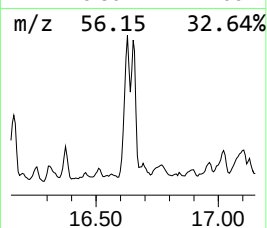
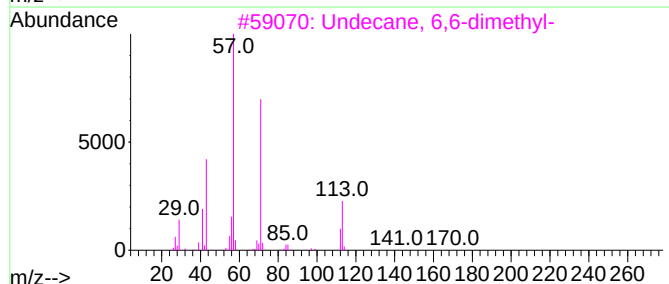
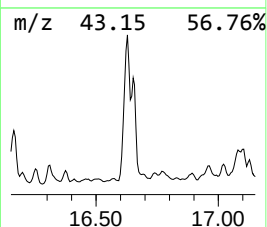
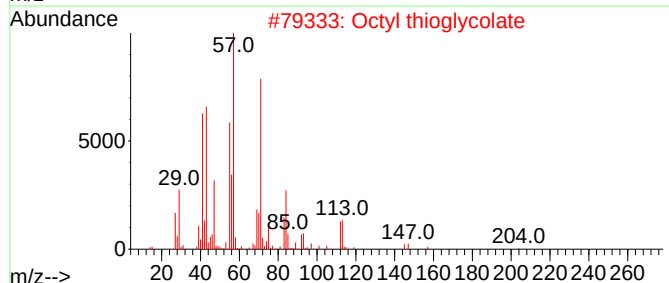
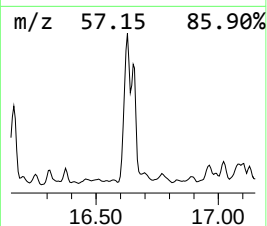
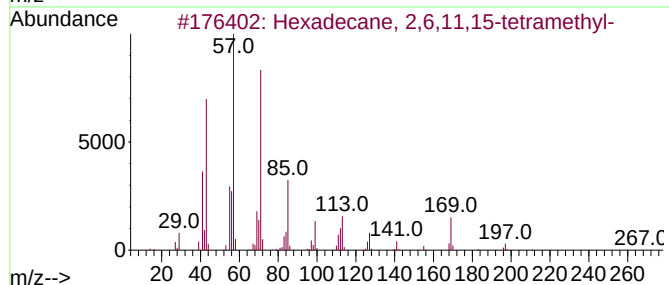
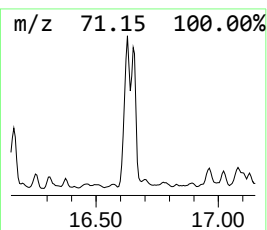
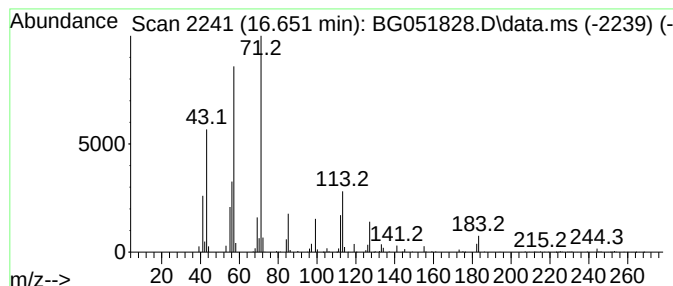
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 43 (DEL) Alkane: Straight-Chai... Concentration Rank 9

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 16.651 | 72.91 ng/ul | 5705850 | Phenanthrene-d10 | 17.680 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|------|-------------------------------------|-----|-----------|--------------|------|
| 1    |      | Hexadecane, 2,6,11,15-tetramethyl-  | 282 | C20H42    | 000504-44-9  | 72   |
| 2    |      | Octyl thioglycolate                 | 204 | C10H20O2S | 007664-80-4  | 64   |
| 3    |      | Undecane, 6,6-dimethyl-             | 184 | C13H28    | 017312-76-4  | 59   |
| 4    |      | Sulfurous acid, 2-ethylhexyl pen... | 404 | C23H48O3S | 1000309-19-8 | 59   |
| 5    |      | Sulfurous acid, dodecyl 2-ethylh... | 362 | C20H42O3S | 1000309-19-5 | 59   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122721\  
Data File : BG051828.D  
Acq On : 28 Dec 2021 14:05  
Operator : CG/JU  
Sample : M5215-01DL 2X  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD3DL

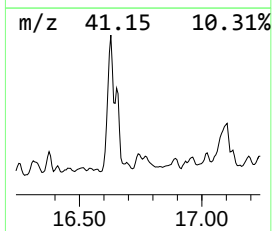
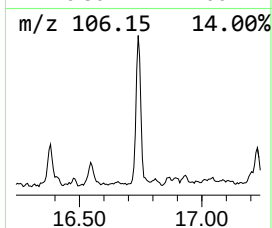
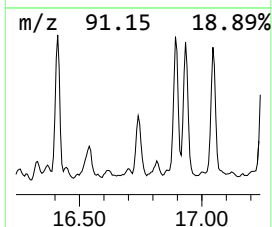
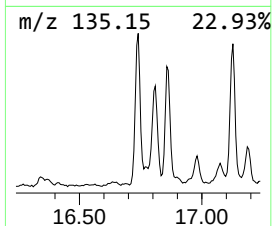
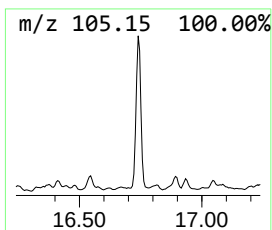
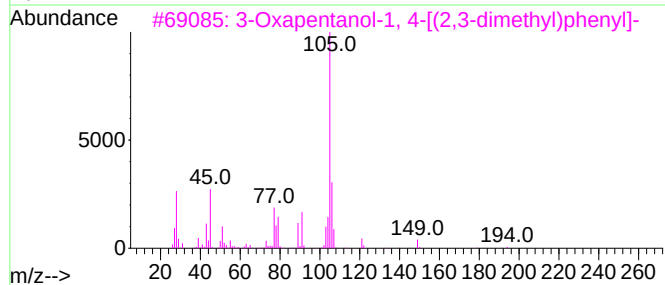
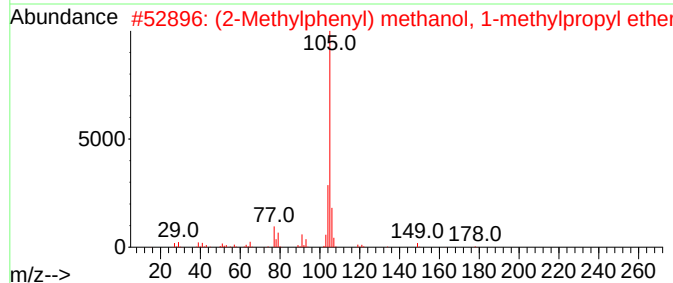
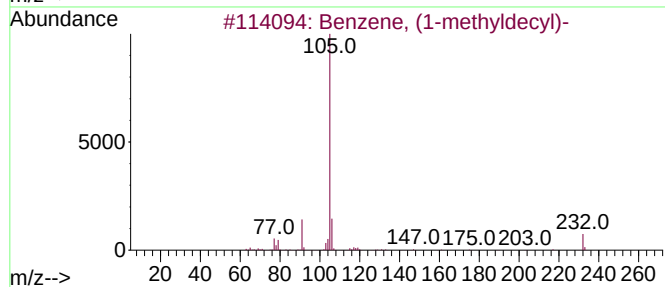
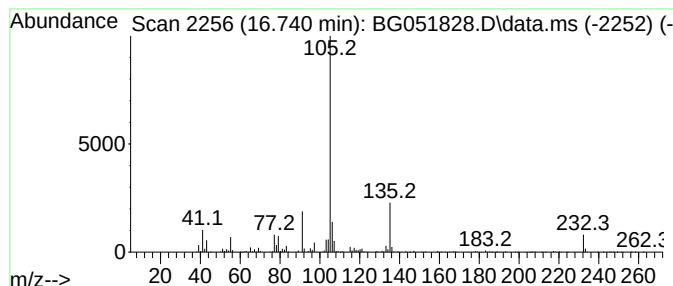
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 44 Benzene, (1-methyldecyl)- Concentration Rank 17

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 16.740 | 34.84 ng/ul | 2726460 | Phenanthrene-d10 | 17.680 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Benzene, (1-methyldecyl)-           | 232 | C17H28   | 004536-88-3  | 92   |
| 2    |    |   | (2-Methylphenyl) methanol, 1-met... | 178 | C12H18O  | 1000374-44-3 | 45   |
| 3    |    |   | 3-Oxapentanol-1, 4-[(2,3-dimethy... | 194 | C12H18O2 | 1000063-22-7 | 43   |
| 4    |    |   | 2-Indanol                           | 134 | C9H10O   | 004254-29-9  | 43   |
| 5    |    |   | Disulfide, bis(p-methylbenzyl)      | 274 | C16H18S2 | 020193-94-6  | 43   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122721\  
Data File : BG051828.D  
Acq On : 28 Dec 2021 14:05  
Operator : CG/JU  
Sample : M5215-01DL 2X  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD3DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

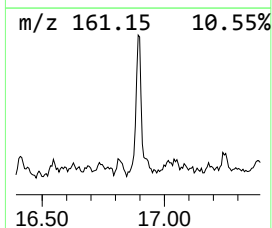
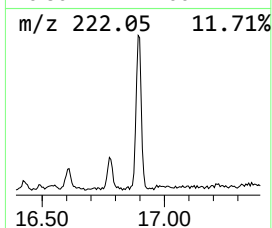
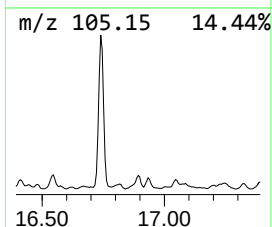
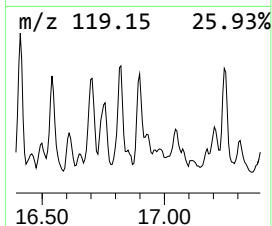
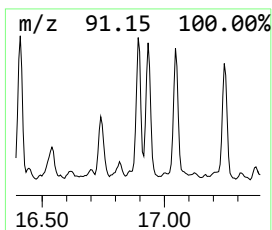
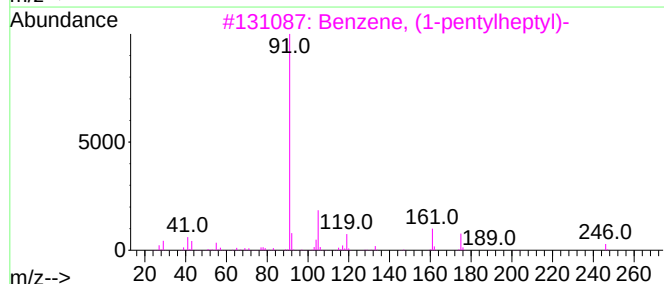
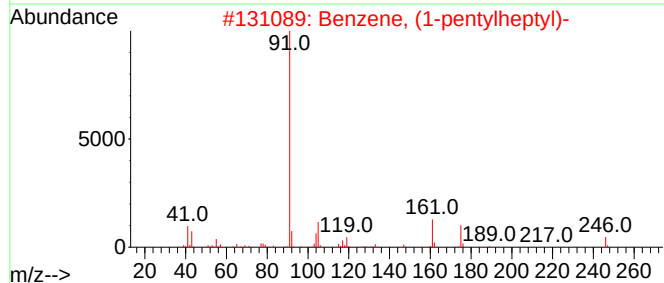
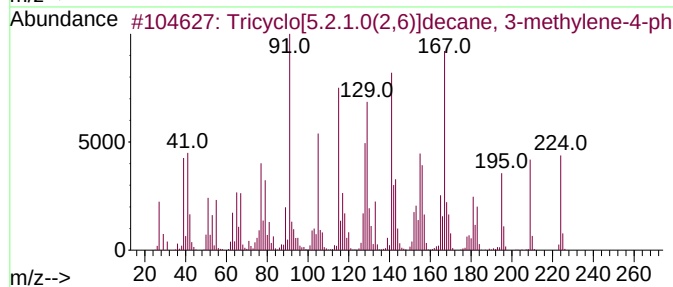
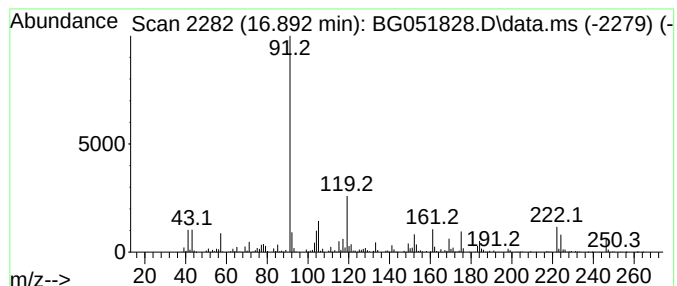
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*

Peak Number 45 Tricyclo[5.2.1.0(2,6)]decan... Concentration Rank 41

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 16.892 | 15.84 ng/ul | 1239980 | Phenanthrene-d10 | 17.680 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|------------|--------------|------|
| 1    |    |   | Tricyclo[5.2.1.0(2,6)]decane, 3-... | 224 | C17H20     | 1000150-37-1 | 90   |
| 2    |    |   | Benzene, (1-pentylheptyl)-          | 246 | C18H30     | 002719-62-2  | 52   |
| 3    |    |   | Benzene, (1-pentylheptyl)-          | 246 | C18H30     | 002719-62-2  | 50   |
| 4    |    |   | 2-phenylbutyric acid, 2,2,2-trif... | 246 | C12H13F3O2 | 1000376-55-3 | 46   |
| 5    |    |   | Aziridine, 1-phenyl-                | 119 | C8H9N      | 000696-18-4  | 43   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122721\  
Data File : BG051828.D  
Acq On : 28 Dec 2021 14:05  
Operator : CG/JU  
Sample : M5215-01DL 2X  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

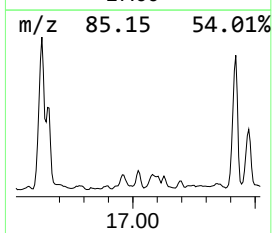
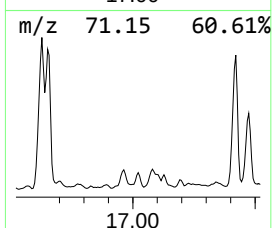
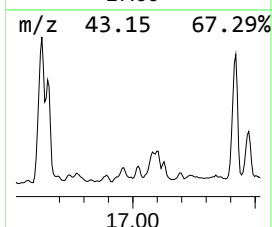
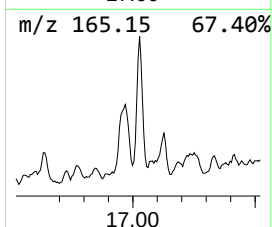
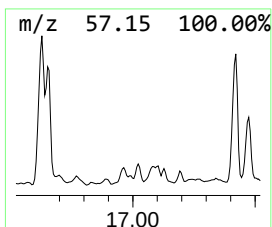
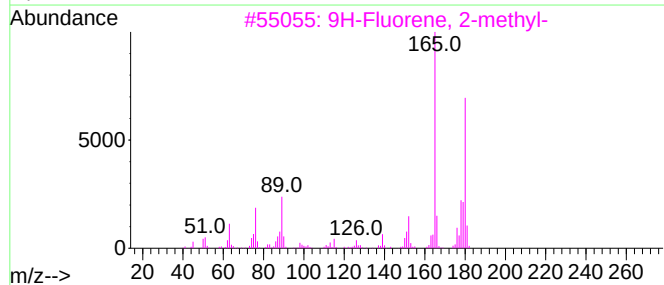
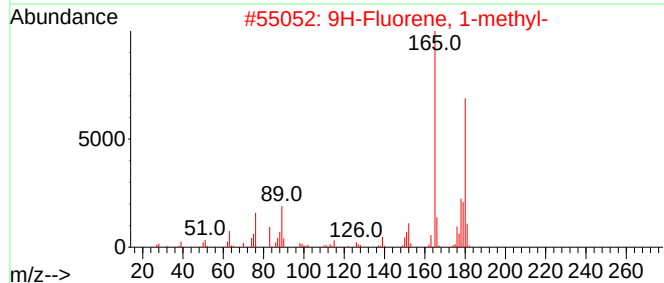
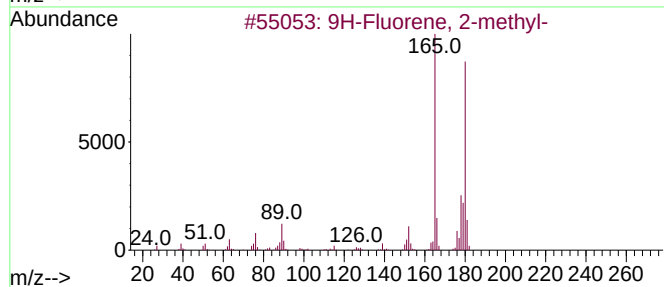
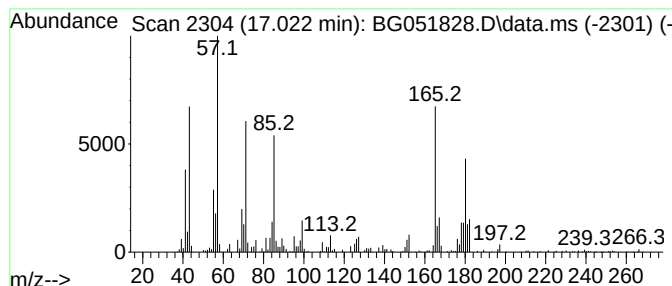
Instrument :  
BNA\_G  
ClientSampleId :  
BGLD3DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 46 9H-Fluorene, 2-methyl- Concentration Rank 45

| R.T.      | EstConc                | Area    | Relative to ISTD |         | R.T.        |      |
|-----------|------------------------|---------|------------------|---------|-------------|------|
| 17.022    | 14.74 ng/ul            | 1153560 | Phenanthrene-d10 |         | 17.680      |      |
| Hit# of 5 | Tentative ID           |         | MW               | MolForm | CAS#        | Qual |
| 1         | 9H-Fluorene, 2-methyl- |         | 180              | C14H12  | 001430-97-3 | 86   |
| 2         | 9H-Fluorene, 1-methyl- |         | 180              | C14H12  | 001730-37-6 | 52   |
| 3         | 9H-Fluorene, 2-methyl- |         | 180              | C14H12  | 001430-97-3 | 52   |
| 4         | 9H-Fluorene, 9-methyl- |         | 180              | C14H12  | 002523-37-7 | 46   |
| 5         | 9H-Fluorene, 3-methyl- |         | 180              | C14H12  | 002523-39-9 | 46   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122721\  
Data File : BG051828.D  
Acq On : 28 Dec 2021 14:05  
Operator : CG/JU  
Sample : M5215-01DL 2X  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD3DL

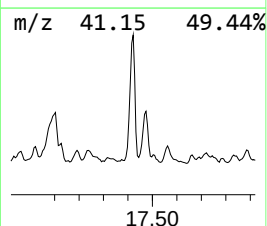
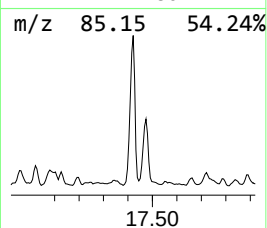
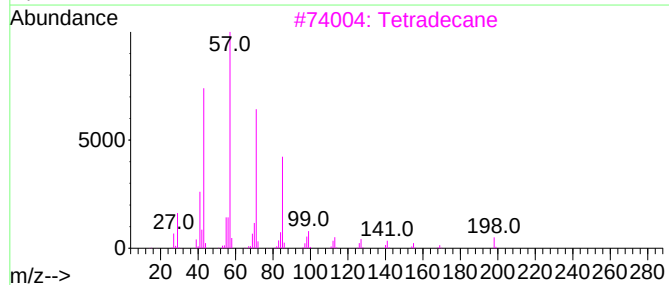
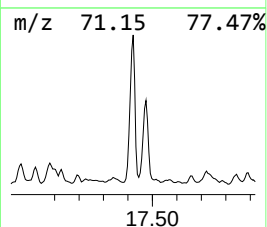
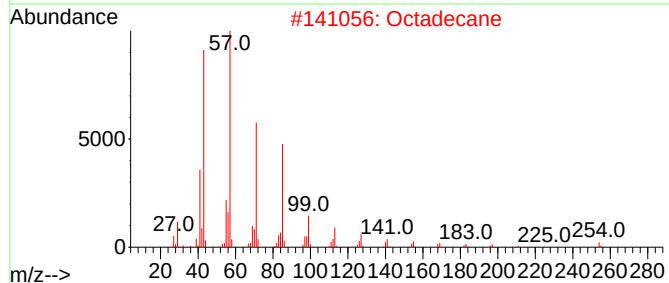
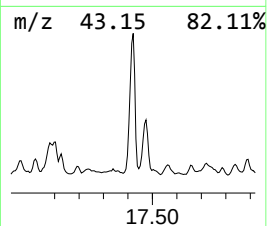
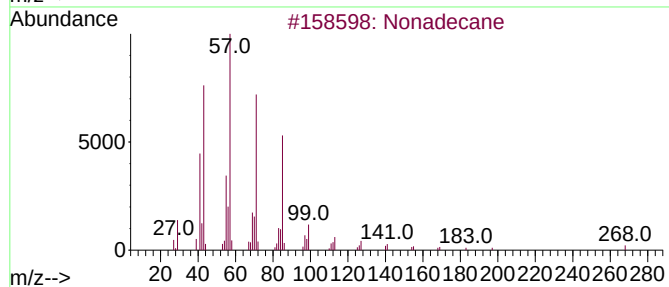
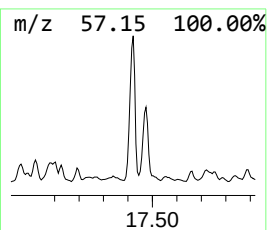
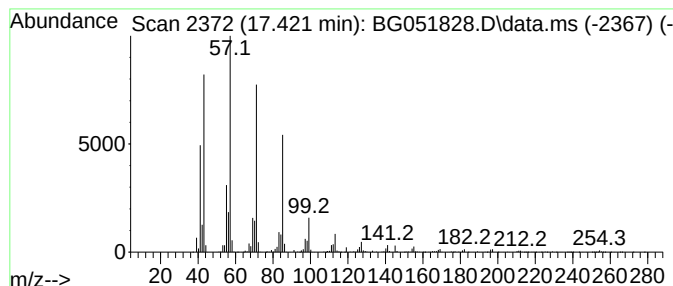
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 47 (DEL) Alkane: Straight-Chai... Concentration Rank 8

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 17.421 | 90.01 ng/ul | 7044030 | Phenanthrene-d10 | 17.680 |

| Hit# | of | 5 | Tentative ID                 | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------------------|-----|---------|-------------|------|
| 1    |    |   | Nonadecane                   | 268 | C19H40  | 000629-92-5 | 91   |
| 2    |    |   | Octadecane                   | 254 | C18H38  | 000593-45-3 | 91   |
| 3    |    |   | Tetradecane                  | 198 | C14H30  | 000629-59-4 | 91   |
| 4    |    |   | Dodecane, 2-methyl-6-propyl- | 226 | C16H34  | 055045-08-4 | 91   |
| 5    |    |   | Tetradecane, 4-ethyl-        | 226 | C16H34  | 055045-14-2 | 91   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122721\  
Data File : BG051828.D  
Acq On : 28 Dec 2021 14:05  
Operator : CG/JU  
Sample : M5215-01DL 2X  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD3DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

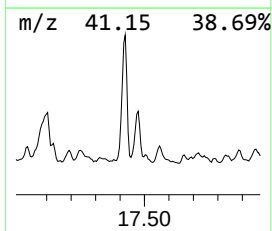
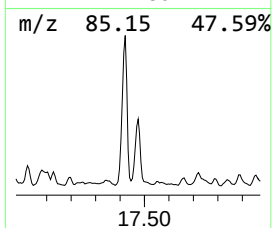
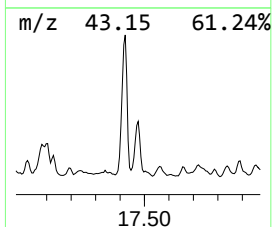
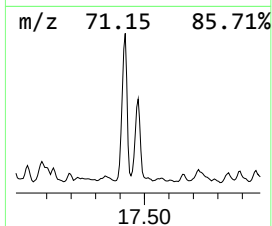
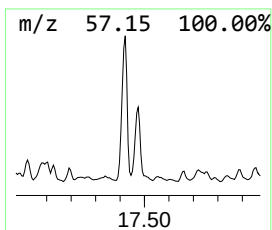
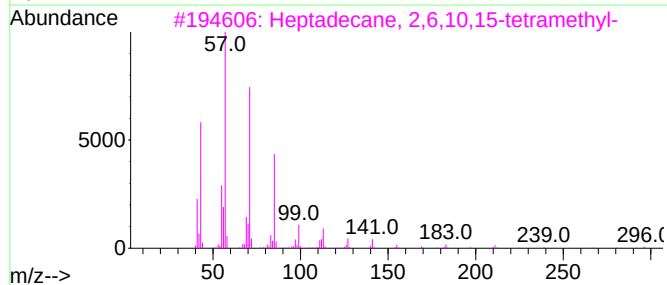
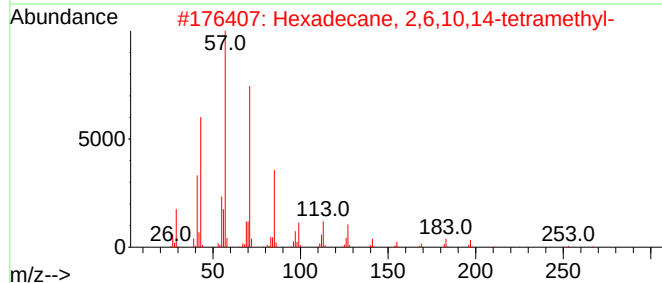
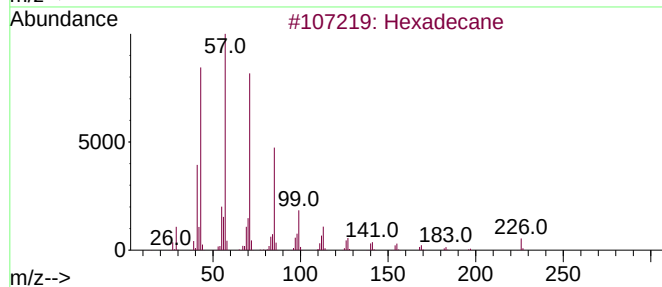
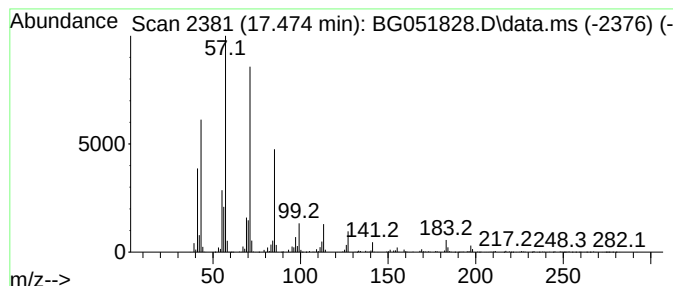
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*

Peak Number 48 (DEL) Alkane: Straight-Chai... Concentration Rank 15

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 17.474 | 41.17 ng/ul | 3221660 | Phenanthrene-d10 | 17.680 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|---------|--------------|------|
| 1    |    |   | Hexadecane                          | 226 | C16H34  | 000544-76-3  | 91   |
| 2    |    |   | Hexadecane, 2,6,10,14-tetramethyl-  | 282 | C20H42  | 000638-36-8  | 91   |
| 3    |    |   | Heptadecane, 2,6,10,15-tetramethyl- | 296 | C21H44  | 054833-48-6  | 91   |
| 4    |    |   | Tridecane                           | 184 | C13H28  | 000629-50-5  | 90   |
| 5    |    |   | 3-Ethyl-2,6,10-trimethylundecane    | 226 | C16H34  | 1000432-25-9 | 86   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122721\  
Data File : BG051828.D  
Acq On : 28 Dec 2021 14:05  
Operator : CG/JU  
Sample : M5215-01DL 2X  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD3DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

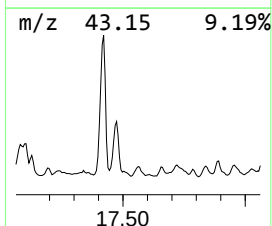
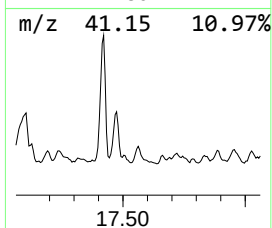
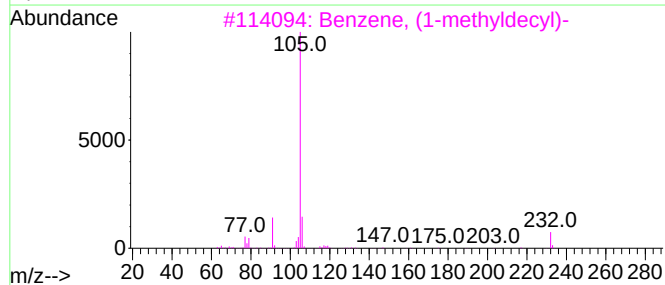
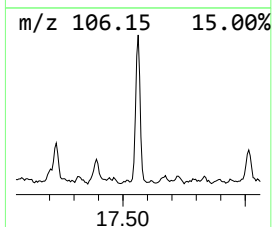
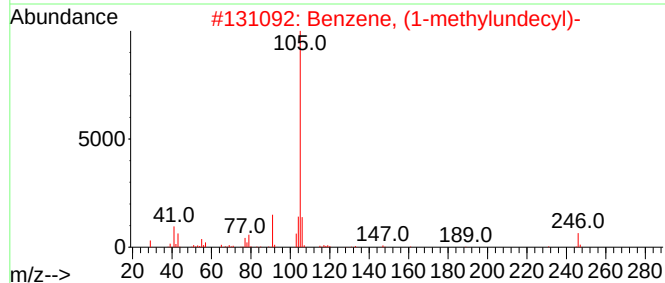
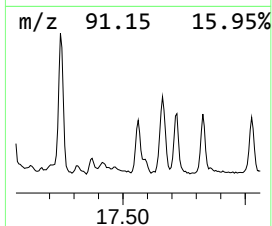
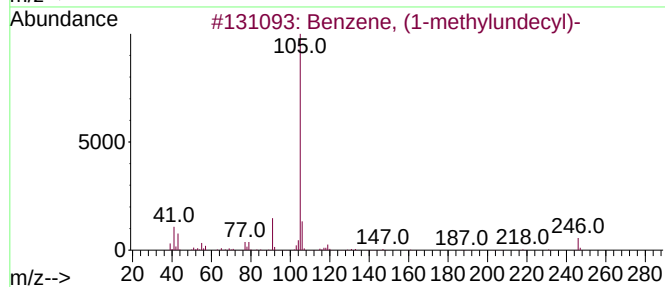
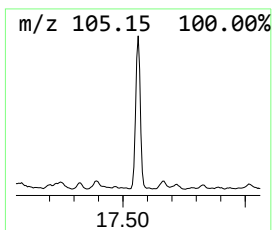
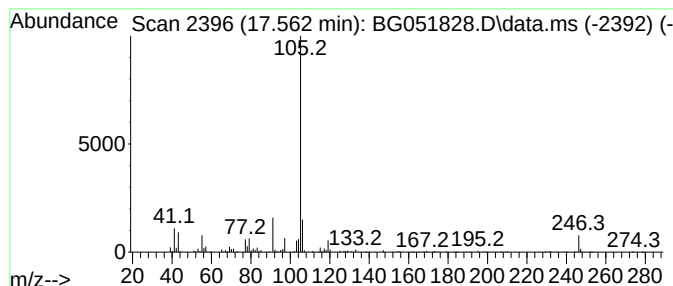
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 49 Benzene, (1-methylundecyl)- Concentration Rank 16

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 17.562 | 35.10 ng/ul | 2746660 | Phenanthrene-d10 | 17.680 |

| Hit# | of 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|------|-----------------------------|-----|---------|-------------|------|
| 1    |      | Benzene, (1-methylundecyl)- | 246 | C18H30  | 002719-61-1 | 91   |
| 2    |      | Benzene, (1-methylundecyl)- | 246 | C18H30  | 002719-61-1 | 89   |
| 3    |      | Benzene, (1-methyldecyl)-   | 232 | C17H28  | 004536-88-3 | 76   |
| 4    |      | Benzene, (1-methylundecyl)- | 246 | C18H30  | 002719-61-1 | 64   |
| 5    |      | Benzene, (1-methylheptyl)-  | 190 | C14H22  | 000777-22-0 | 59   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122721\  
Data File : BG051828.D  
Acq On : 28 Dec 2021 14:05  
Operator : CG/JU  
Sample : M5215-01DL 2X  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

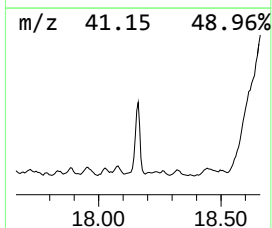
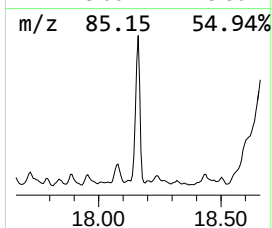
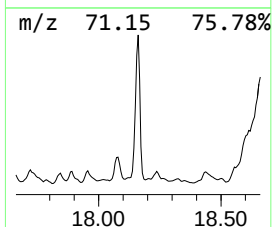
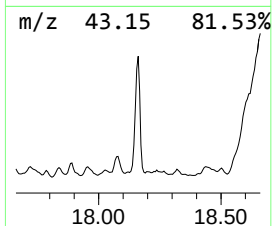
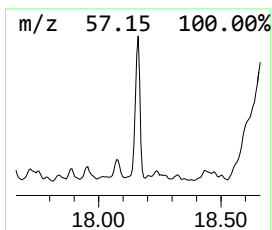
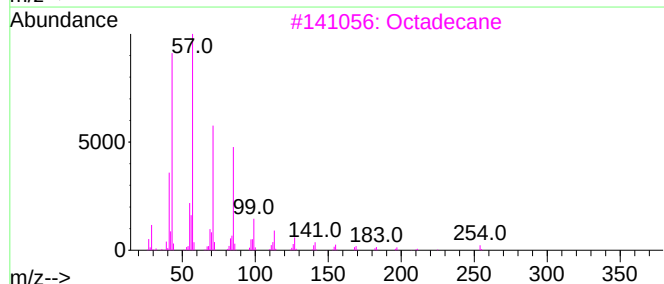
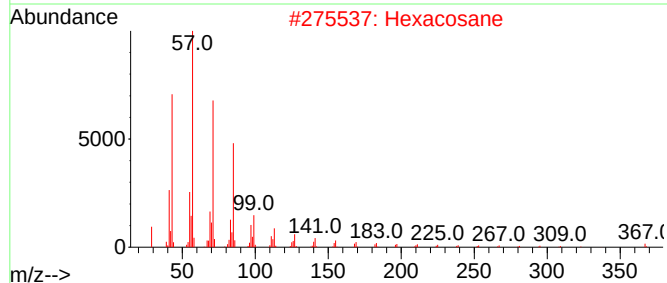
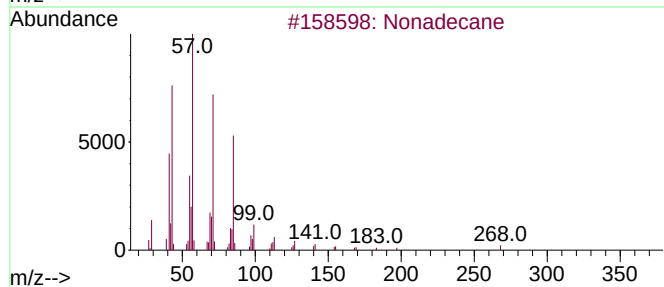
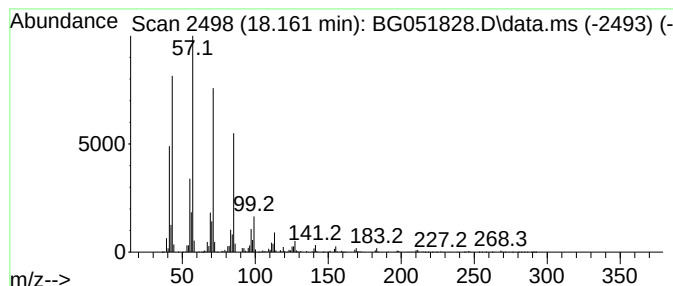
Instrument :  
BNA\_G  
ClientSampleId :  
BGLD3DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 50 (DEL) Alkane: Straight-Chai... Concentration Rank 7

| R.T.      | EstConc      | Area    | Relative to ISTD |         | R.T.        |      |
|-----------|--------------|---------|------------------|---------|-------------|------|
| 18.161    | 93.63 ng/ul  | 7327510 | Phenanthrene-d10 |         | 17.680      |      |
| Hit# of 5 | Tentative ID |         | MW               | MolForm | CAS#        | Qual |
| 1         | Nonadecane   |         | 268              | C19H40  | 000629-92-5 | 91   |
| 2         | Hexacosane   |         | 366              | C26H54  | 000630-01-3 | 91   |
| 3         | Octadecane   |         | 254              | C18H38  | 000593-45-3 | 91   |
| 4         | Heneicosane  |         | 296              | C21H44  | 000629-94-7 | 91   |
| 5         | Heptadecane  |         | 240              | C17H36  | 000629-78-7 | 91   |





Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122721\  
Data File : BG051828.D  
Acq On : 28 Dec 2021 14:05  
Operator : CG/JU  
Sample : M5215-01DL 2X  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD3DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

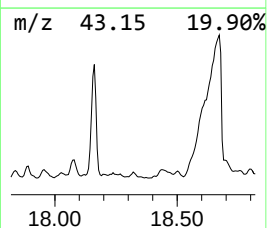
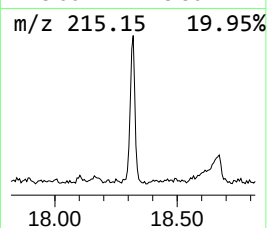
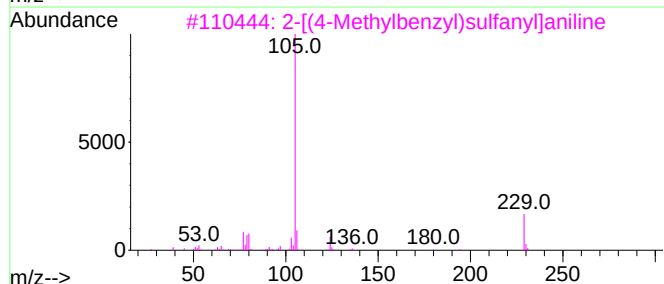
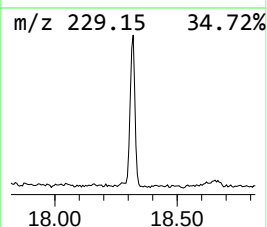
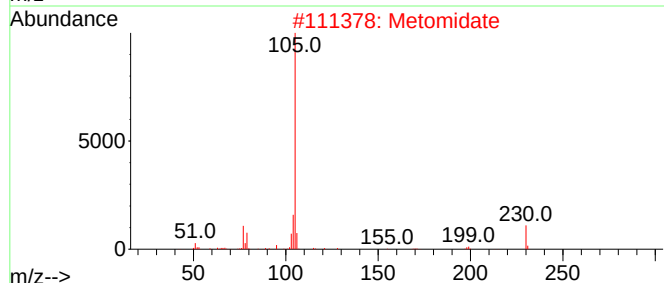
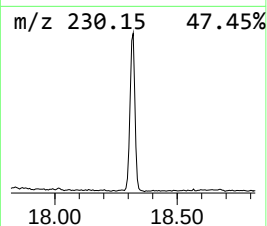
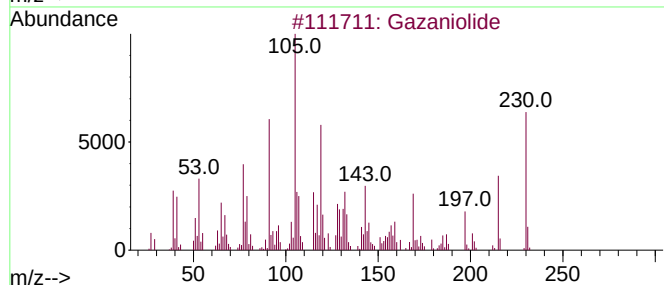
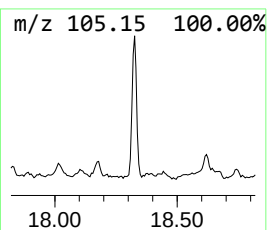
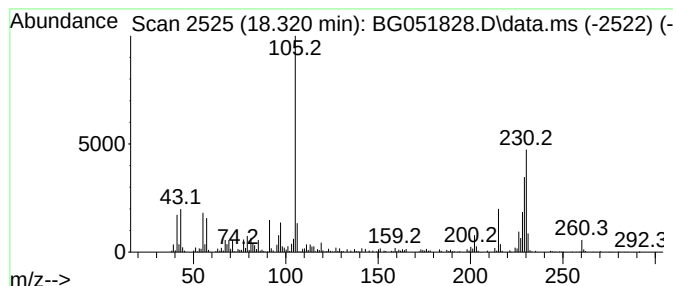
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 51 unknown-04 Concentration Rank 31

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 18.320 | 21.81 ng/ul | 1706840 | Phenanthrene-d10 | 17.680 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|------------|--------------|------|
| 1    |    |   | Gazaniolide                         | 230 | C15H18O2   | 071609-02-4  | 37   |
| 2    |    |   | Metomidate                          | 230 | C13H14N2O2 | 005377-20-8  | 25   |
| 3    |    |   | 2-[(4-Methylbenzyl)sulfanyl]aniline | 229 | C14H15NS   | 136620-24-1  | 22   |
| 4    |    |   | 2H-Pyran-3-aldimien, 5,6-dihydro... | 215 | C14H17NO   | 1000153-18-8 | 14   |
| 5    |    |   | N-Benzoyl derivative of p-fluoro... | 215 | C13H10FNO  | 000366-75-6  | 12   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122721\  
Data File : BG051828.D  
Acq On : 28 Dec 2021 14:05  
Operator : CG/JU  
Sample : M5215-01DL 2X  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD3DL

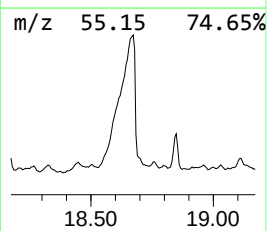
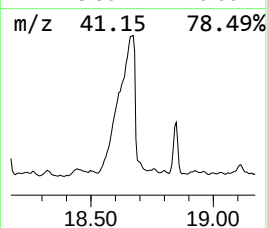
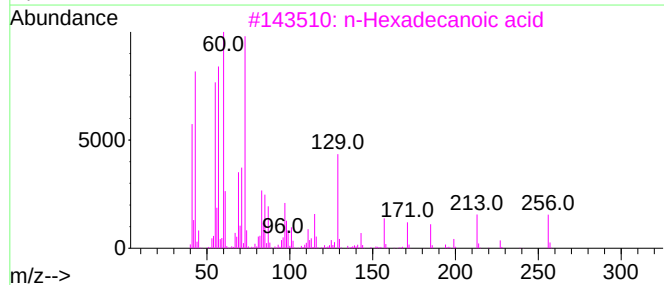
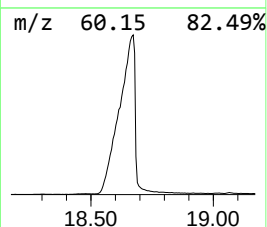
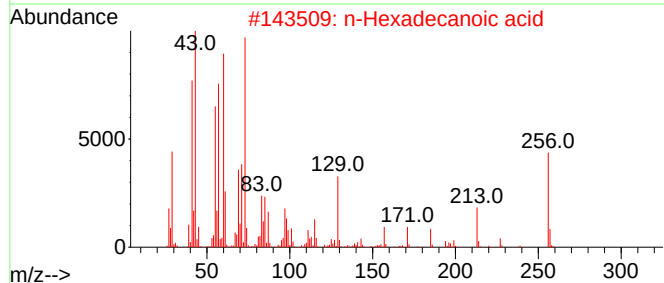
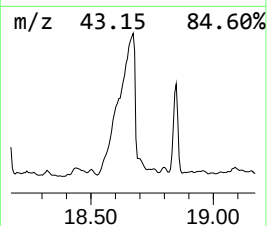
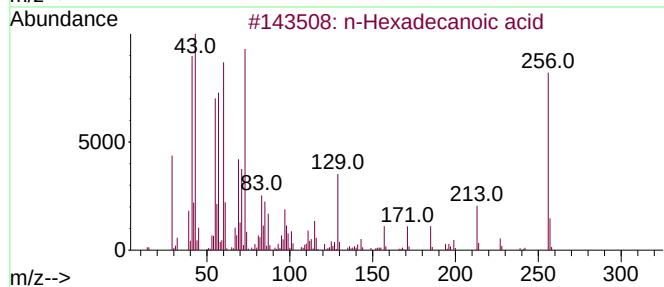
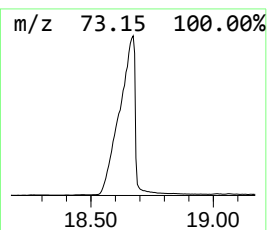
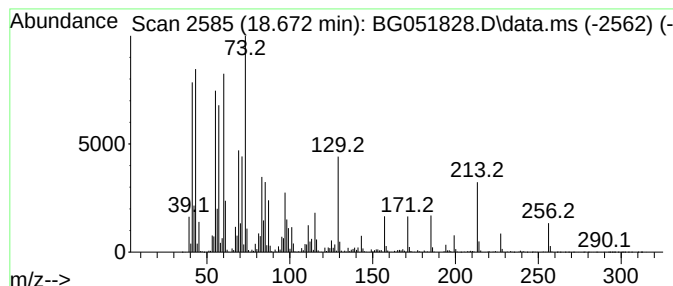
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 52 n-Hexadecanoic acid Concentration Rank 1

| R.T.   | EstConc      | Area     | Relative to ISTD | R.T.   |
|--------|--------------|----------|------------------|--------|
| 18.672 | 739.79 ng/ul | 57894900 | Phenanthrene-d10 | 17.680 |

| Hit# | of | 5 | Tentative ID        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|---------------------|-----|----------|-------------|------|
| 1    |    |   | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 99   |
| 2    |    |   | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 97   |
| 3    |    |   | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 96   |
| 4    |    |   | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 93   |
| 5    |    |   | Pentadecanoic acid  | 242 | C15H30O2 | 001002-84-2 | 90   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122721\  
Data File : BG051828.D  
Acq On : 28 Dec 2021 14:05  
Operator : CG/JU  
Sample : M5215-01DL 2X  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD3DL

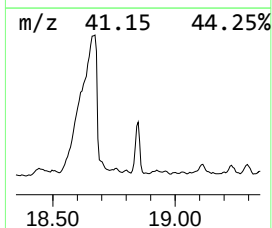
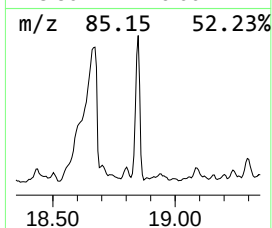
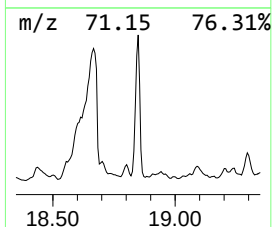
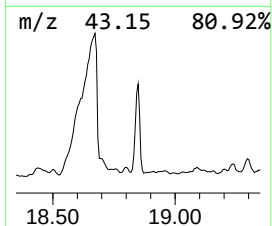
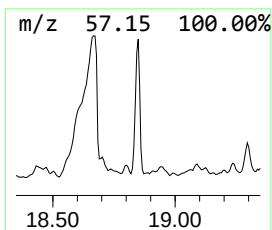
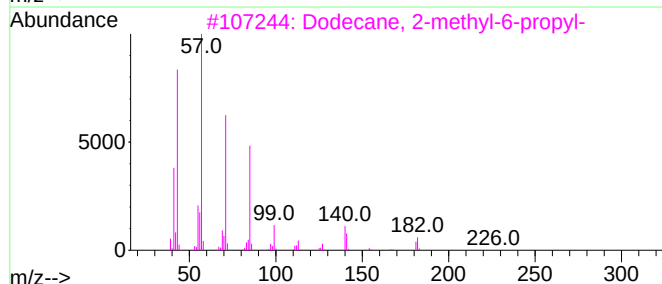
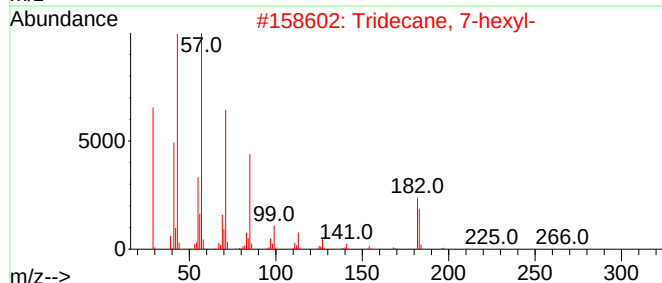
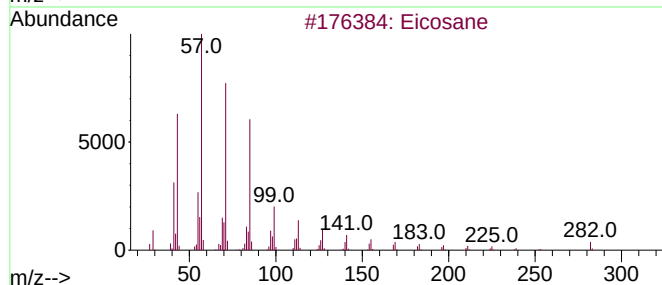
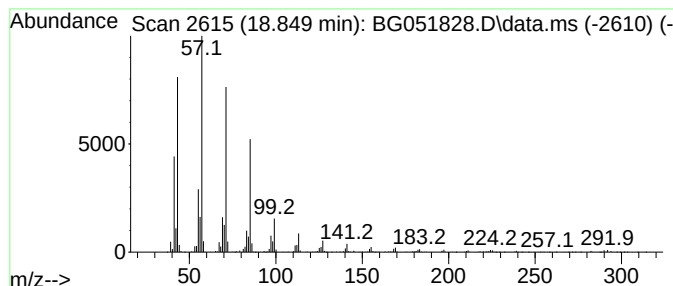
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*

Peak Number 53 (DEL) Alkane: Straight-Chai... Concentration Rank 12

| R.T.      | EstConc                      | Area    | Relative to ISTD |         | R.T.        |      |
|-----------|------------------------------|---------|------------------|---------|-------------|------|
| 18.849    | 62.87 ng/ul                  | 4920410 | Phenanthrene-d10 |         | 17.680      |      |
| Hit# of 5 | Tentative ID                 |         | MW               | MolForm | CAS#        | Qual |
| 1         | Eicosane                     |         | 282              | C20H42  | 000112-95-8 | 96   |
| 2         | Tridecane, 7-hexyl-          |         | 268              | C19H40  | 007225-66-3 | 95   |
| 3         | Dodecane, 2-methyl-6-propyl- |         | 226              | C16H34  | 055045-08-4 | 93   |
| 4         | Nonadecane                   |         | 268              | C19H40  | 000629-92-5 | 91   |
| 5         | Eicosane                     |         | 282              | C20H42  | 000112-95-8 | 91   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122721\  
Data File : BG051828.D  
Acq On : 28 Dec 2021 14:05  
Operator : CG/JU  
Sample : M5215-01DL 2X  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD3DL

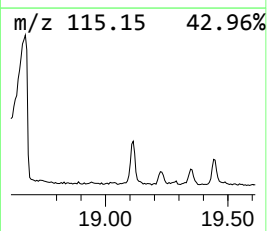
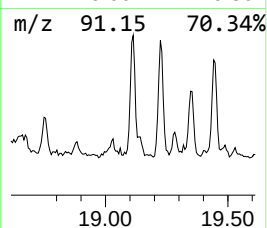
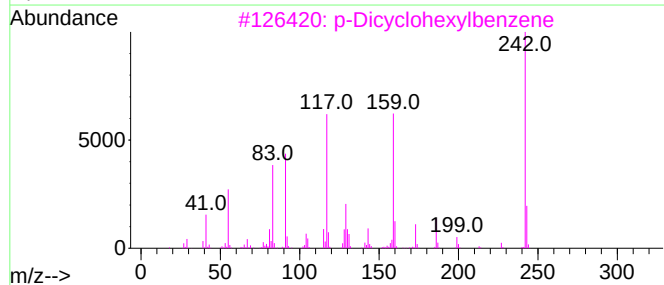
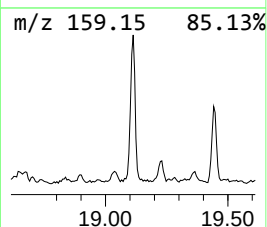
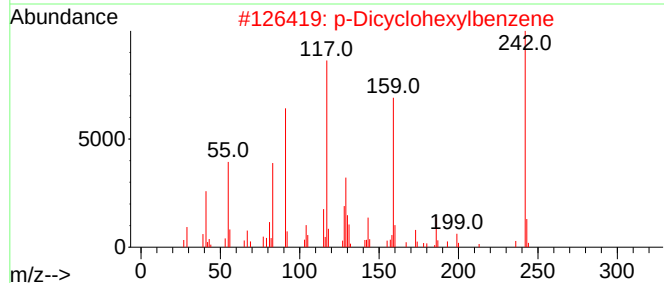
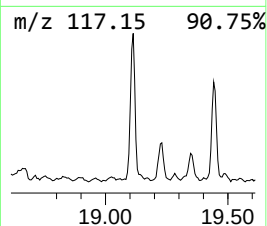
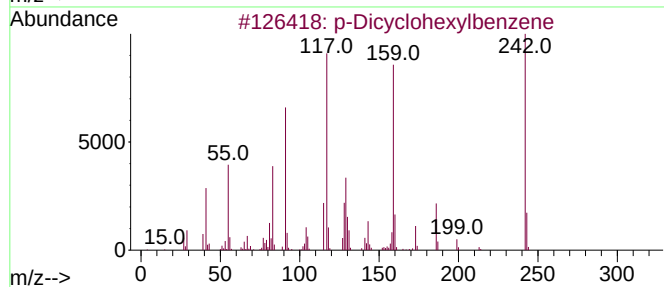
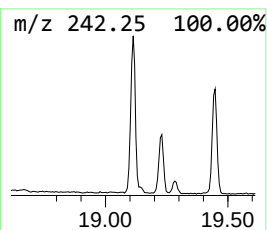
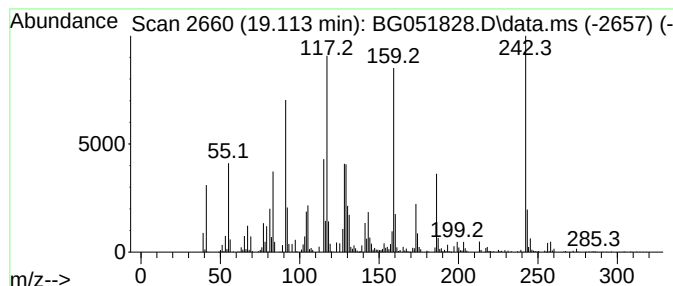
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 54 p-Dicyclohexylbenzene Concentration Rank 21

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 19.113 | 29.49 ng/ul | 2307580 | Phenanthrene-d10 | 17.680 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | p-Dicyclohexylbenzene               | 242 | C18H26  | 001087-02-1 | 98   |
| 2    |    |   | p-Dicyclohexylbenzene               | 242 | C18H26  | 001087-02-1 | 90   |
| 3    |    |   | p-Dicyclohexylbenzene               | 242 | C18H26  | 001087-02-1 | 89   |
| 4    |    |   | Bicyclohexyl, 4-phenyl-             | 242 | C18H26  | 020273-27-2 | 58   |
| 5    |    |   | Naphthalene, 1,2,3,4-tetrahydro-... | 174 | C13H18  | 030316-36-0 | 30   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122721\  
Data File : BG051828.D  
Acq On : 28 Dec 2021 14:05  
Operator : CG/JU  
Sample : M5215-01DL 2X  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD3DL

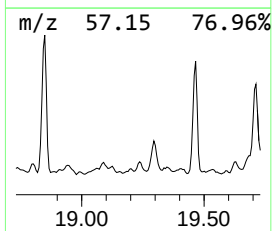
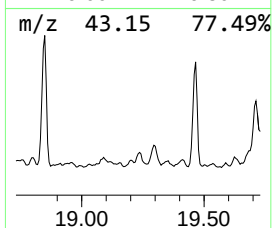
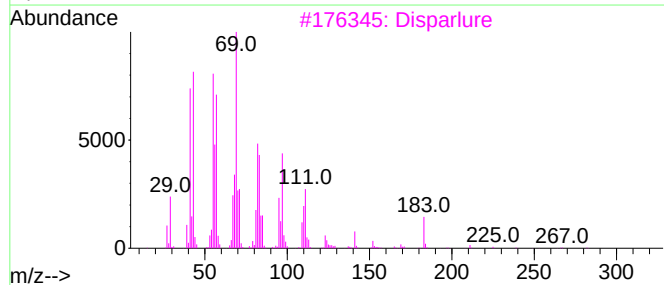
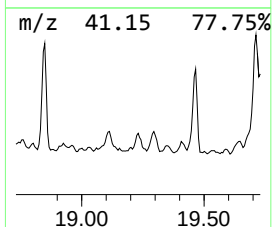
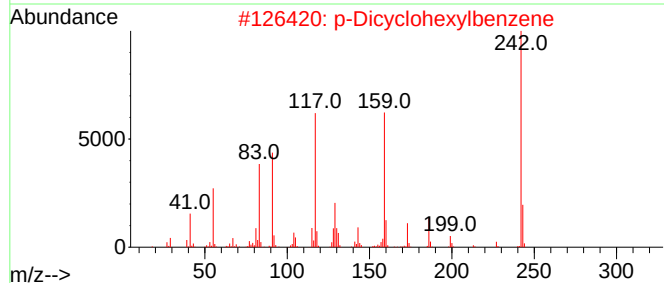
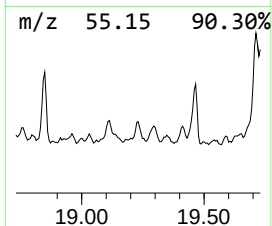
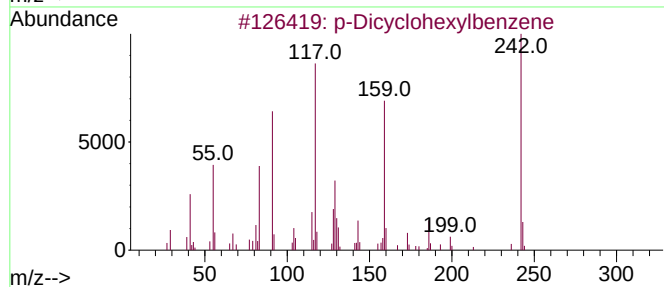
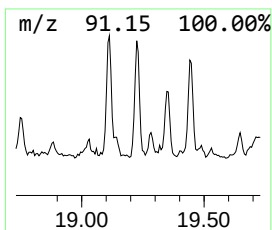
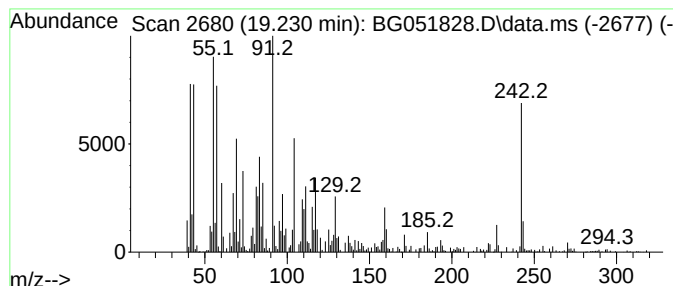
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 55 unknown-05 Concentration Rank 47

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 19.230 | 14.08 ng/ul | 1102220 | Phenanthrene-d10 | 17.680 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm       | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|---------------|--------------|------|
| 1    |    |   | p-Dicyclohexylbenzene               | 242 | C18H26        | 001087-02-1  | 38   |
| 2    |    |   | p-Dicyclohexylbenzene               | 242 | C18H26        | 001087-02-1  | 18   |
| 3    |    |   | Disparlure                          | 282 | C19H38O       | 029804-22-6  | 15   |
| 4    |    |   | 10-Chloro-1-decanol, chlorodiflu... | 304 | C12H20Cl2F2O2 | 1000447-45-2 | 14   |
| 5    |    |   | Benzene, cyclohexyl-                | 160 | C12H16        | 000827-52-1  | 11   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122721\  
Data File : BG051828.D  
Acq On : 28 Dec 2021 14:05  
Operator : CG/JU  
Sample : M5215-01DL 2X  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD3DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

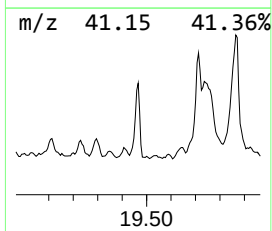
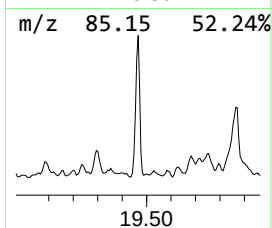
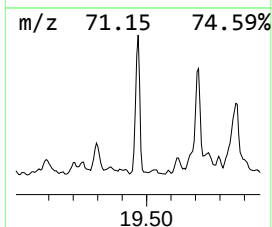
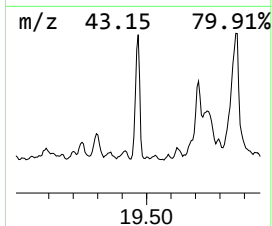
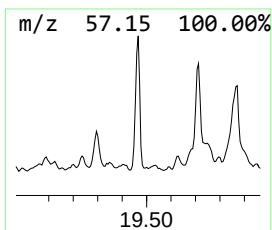
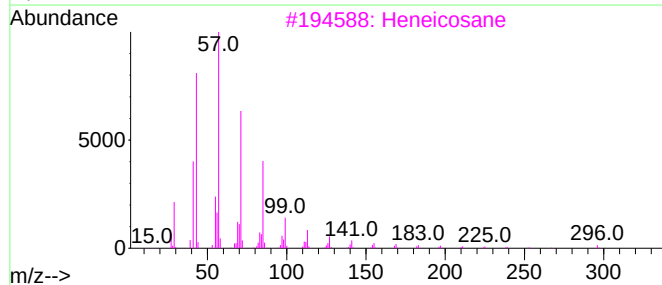
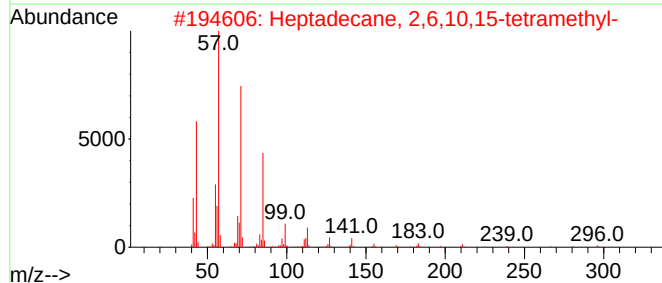
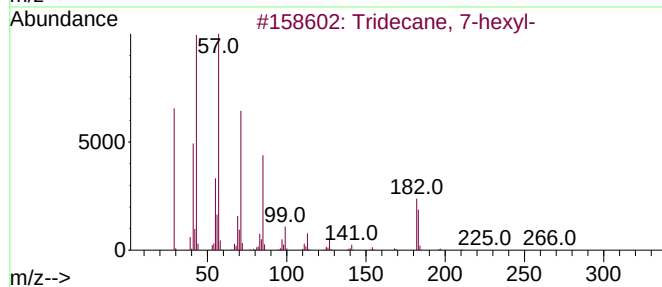
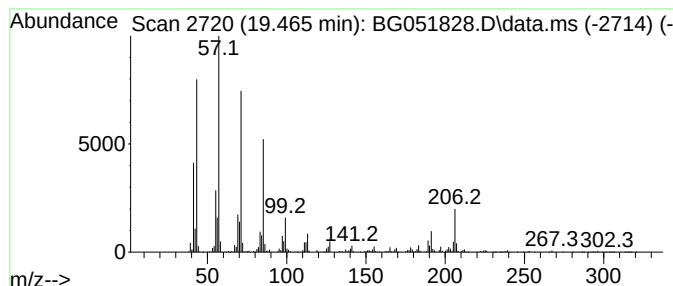
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 56 (DEL) Alkane: Straight-Chai... Concentration Rank 10

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 19.465 | 67.51 ng/ul | 5283490 | Phenanthrene-d10 | 17.680 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|------|-------------------------------------|-----|---------|-------------|------|
| 1    |      | Tridecane, 7-hexyl-                 | 268 | C19H40  | 007225-66-3 | 93   |
| 2    |      | Heptadecane, 2,6,10,15-tetramethyl- | 296 | C21H44  | 054833-48-6 | 93   |
| 3    |      | Heneicosane                         | 296 | C21H44  | 000629-94-7 | 89   |
| 4    |      | Nonadecane                          | 268 | C19H40  | 000629-92-5 | 76   |
| 5    |      | Tetracosane                         | 338 | C24H50  | 000646-31-1 | 68   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122721\  
Data File : BG051828.D  
Acq On : 28 Dec 2021 14:05  
Operator : CG/JU  
Sample : M5215-01DL 2X  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD3DL

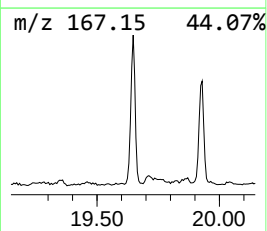
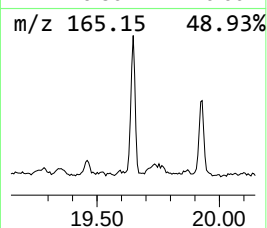
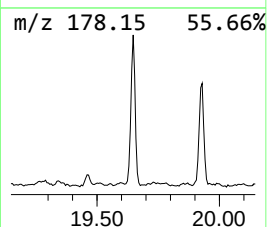
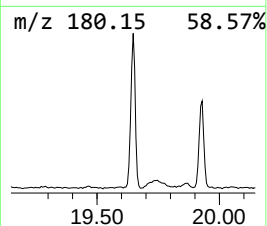
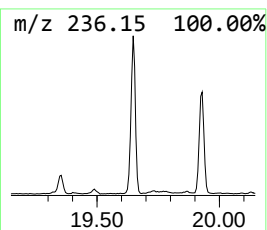
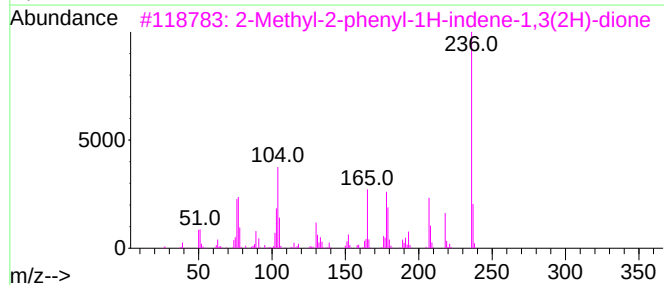
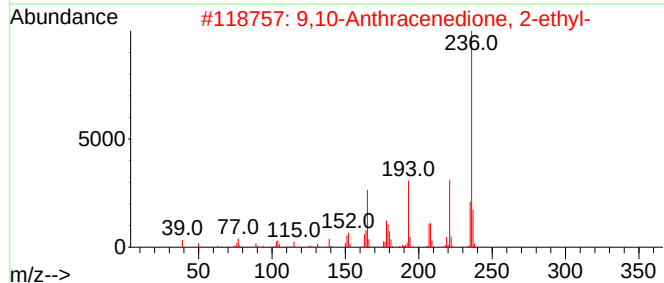
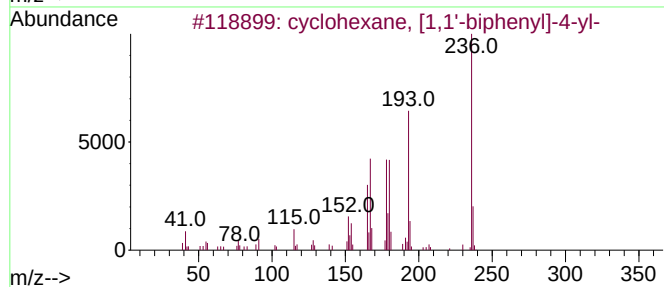
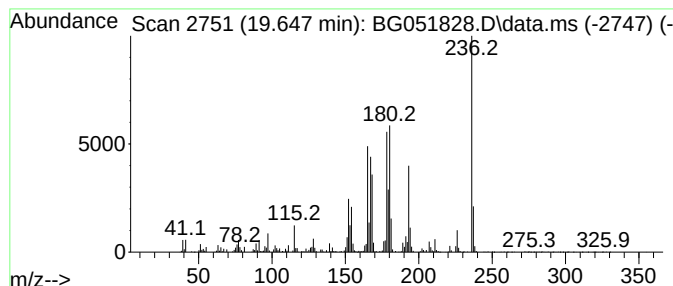
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 57 cyclohexane, [1,1'-biphenyl]... Concentration Rank 20

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 19.647 | 30.07 ng/ul | 2353320 | Phenanthrene-d10 | 17.680 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm      | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|--------------|--------------|------|
| 1    |    |   | cyclohexane, [1,1'-biphenyl]-4-yl-  | 236 | C18H20       | 1000401-12-4 | 93   |
| 2    |    |   | 9,10-Anthracenedione, 2-ethyl-      | 236 | C16H12O2     | 000084-51-5  | 43   |
| 3    |    |   | 2-Methyl-2-phenyl-1H-indene-1,3(... | 236 | C16H12O2     | 1010332-18-2 | 35   |
| 4    |    |   | Ritodrine 0,0',0"-tris-trimethyl... | 503 | C26H45NO3Si3 | 335627-90-2  | 35   |
| 5    |    |   | 7-Methyl-4-phenylcoumarin           | 236 | C16H12O2     | 007758-71-6  | 27   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122721\  
Data File : BG051828.D  
Acq On : 28 Dec 2021 14:05  
Operator : CG/JU  
Sample : M5215-01DL 2X  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD3DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

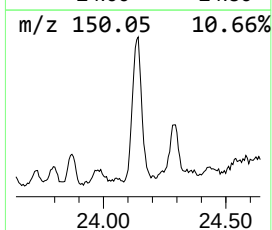
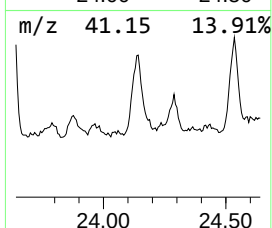
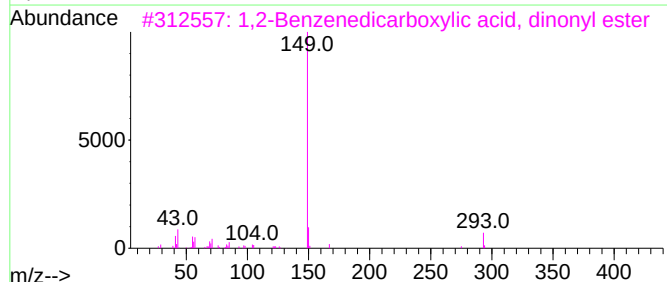
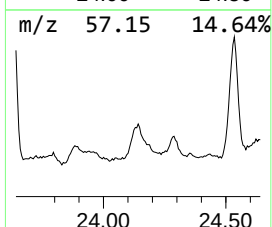
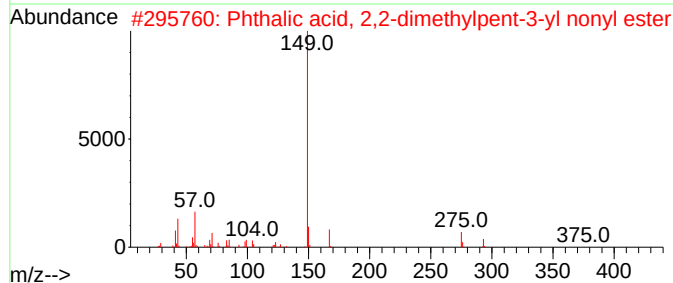
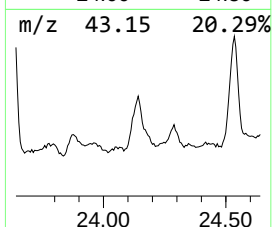
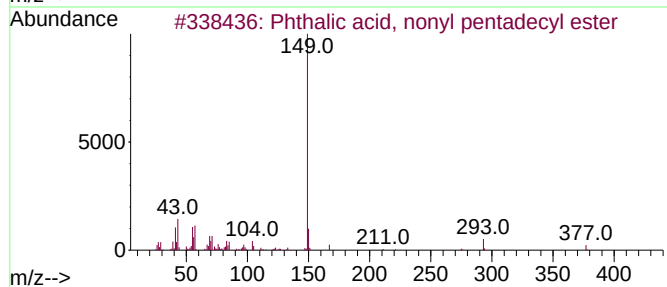
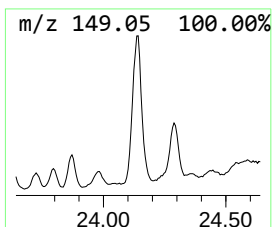
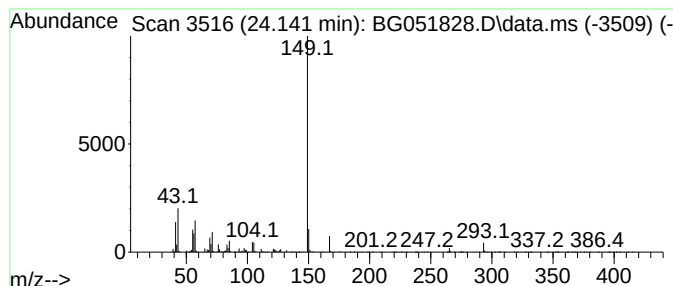
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*

Peak Number 59 Phthalic acid, nonyl pentad... Concentration Rank 29

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 24.141 | 22.97 ng/ul | 3148520 | Perylene-d12     | 25.545 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Phthalic acid, nonyl pentadecyl ... | 502 | C32H54O4 | 1000308-92-7 | 91   |
| 2    |    |   | Phthalic acid, 2,2-dimethylpent-... | 390 | C24H38O4 | 1000415-53-2 | 90   |
| 3    |    |   | 1,2-Benzenedicarboxylic acid, di... | 418 | C26H42O4 | 000084-76-4  | 90   |
| 4    |    |   | Phthalic acid, decyl 2-ethylhexy... | 418 | C26H42O4 | 1000309-00-0 | 86   |
| 5    |    |   | Phthalic acid, dodecyl octyl ester  | 446 | C28H46O4 | 1010308-91-7 | 86   |





Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122721\  
Data File : BG051828.D  
Acq On : 28 Dec 2021 14:05  
Operator : CG/JU  
Sample : M5215-01DL 2X  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD3DL

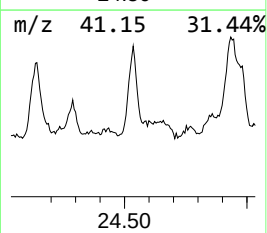
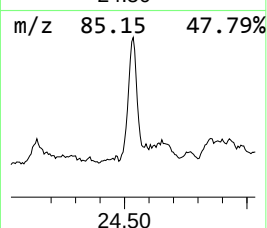
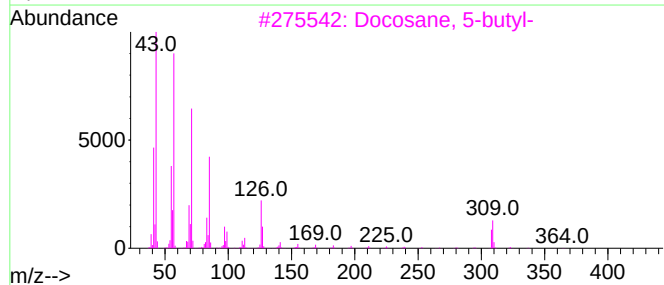
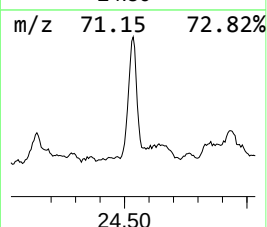
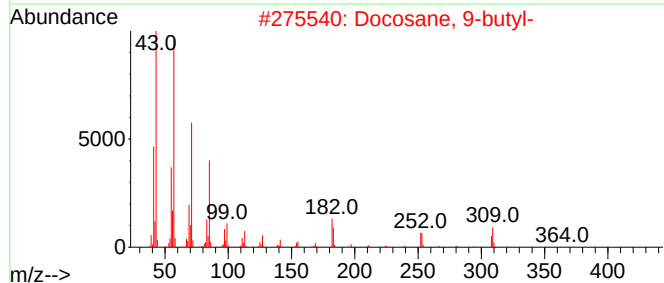
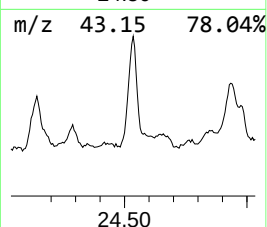
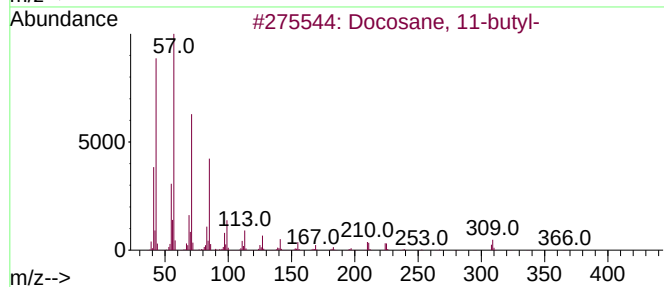
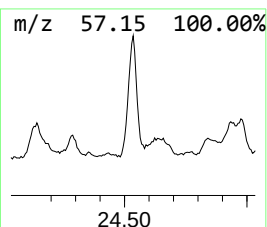
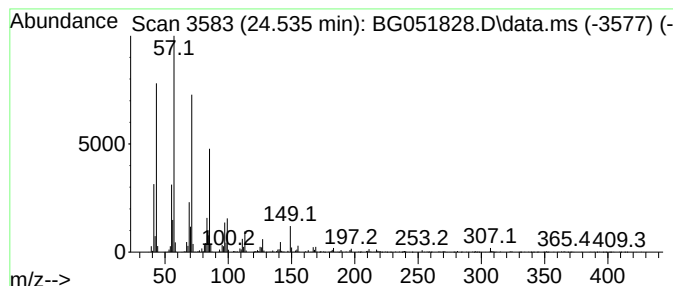
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*

Peak Number 60 (DEL) Alkane: Straight-Chai... Concentration Rank 34

| R.T.      | EstConc               | Area    | Relative to ISTD | R.T.        |      |
|-----------|-----------------------|---------|------------------|-------------|------|
| 24.535    | 20.00 ng/ul           | 2741140 | Perylene-d12     | 25.545      |      |
| Hit# of 5 | Tentative ID          | MW      | MolForm          | CAS#        | Qual |
| 1         | Docosane, 11-butyl-   | 366     | C26H54           | 013475-76-8 | 93   |
| 2         | Docosane, 9-butyl-    | 366     | C26H54           | 055282-14-9 | 93   |
| 3         | Docosane, 5-butyl-    | 366     | C26H54           | 055282-16-1 | 91   |
| 4         | Hexacosane            | 366     | C26H54           | 000630-01-3 | 90   |
| 5         | Octacosane, 2-methyl- | 408     | C29H60           | 001560-98-1 | 90   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122721\  
Data File : BG051828.D  
Acq On : 28 Dec 2021 14:05  
Operator : CG/JU  
Sample : M5215-01DL 2X  
Misc :  
ALS Vial : 35 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD3DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

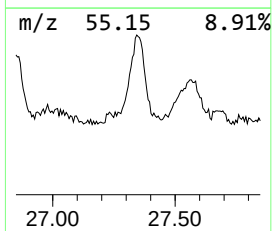
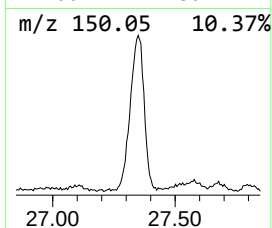
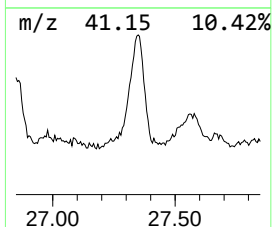
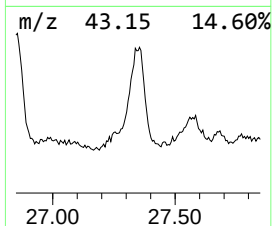
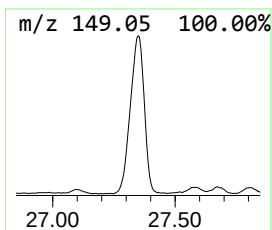
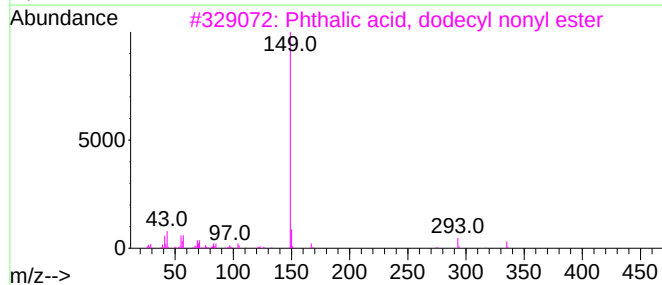
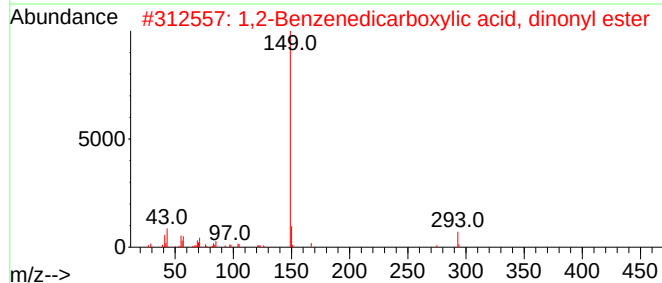
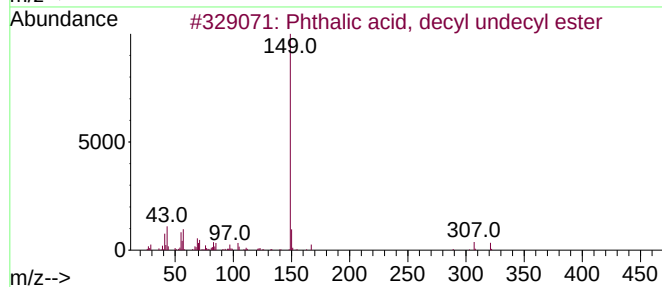
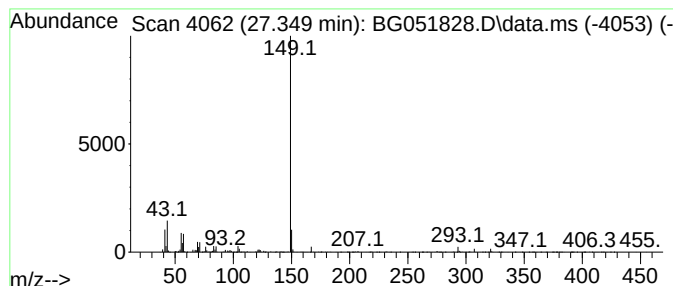
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*

Peak Number 61 Phthalic acid, decyl undecy... Concentration Rank 25

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 27.349 | 26.41 ng/ul | 3620130 | Perylene-d12     | 25.545 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Phthalic acid, decyl undecyl ester  | 460 | C29H48O4 | 1000308-91-9 | 86   |
| 2    |    |   | 1,2-Benzenedicarboxylic acid, di... | 418 | C26H42O4 | 000084-76-4  | 80   |
| 3    |    |   | Phthalic acid, dodecyl nonyl ester  | 460 | C29H48O4 | 1000308-92-6 | 80   |
| 4    |    |   | Phthalic acid, hexyl octyl ester    | 362 | C22H34O4 | 1000424-04-5 | 80   |
| 5    |    |   | 1,2-Benzenedicarboxylic acid, di... | 474 | C30H50O4 | 003648-20-2  | 80   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122721\  
 Data File : BG051828.D  
 Acq On : 28 Dec 2021 14:05  
 Operator : CG/JU  
 Sample : M5215-01DL 2X  
 Misc :  
 ALS Vial : 35 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 BGLD3DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.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       | 5.766  | 146.6   | ng/ul | 14744500 | 1                     | 8.286  | 2012180 | 20.0 |
| Benzene, 1,3-di... | 5.919  | 398.8   | ng/ul | 40123100 | 1                     | 8.286  | 2012180 | 20.0 |
| p-Xylene           | 6.271  | 112.7   | ng/ul | 11342000 | 1                     | 8.286  | 2012180 | 20.0 |
| (DEL) Alkane: S... | 6.530  | 14.9    | ng/ul | 1503930  | 1                     | 8.286  | 2012180 | 20.0 |
| (DEL) Alkane: S... | 6.653  | 9.2     | ng/ul | 924323   | 1                     | 8.286  | 2012180 | 20.0 |
| Benzene, (1-met... | 6.747  | 15.7    | ng/ul | 1579640  | 1                     | 8.286  | 2012180 | 20.0 |
| (DEL) Alkane: S... | 6.818  | 11.9    | ng/ul | 1202130  | 1                     | 8.286  | 2012180 | 20.0 |
| Cyclohexanone, ... | 6.906  | 22.6    | ng/ul | 2269790  | 1                     | 8.286  | 2012180 | 20.0 |
| (DEL) Alkane: S... | 7.164  | 10.6    | ng/ul | 1068790  | 1                     | 8.286  | 2012180 | 20.0 |
| Benzene, propyl-   | 7.258  | 31.0    | ng/ul | 3118260  | 1                     | 8.286  | 2012180 | 20.0 |
| (DEL) Alkane: S... | 7.323  | 14.9    | ng/ul | 1501110  | 1                     | 8.286  | 2012180 | 20.0 |
| Benzene, 1-ethy... | 7.364  | 31.4    | ng/ul | 3155410  | 1                     | 8.286  | 2012180 | 20.0 |
| Benzene, 1,2,3-... | 7.499  | 26.0    | ng/ul | 2619850  | 1                     | 8.286  | 2012180 | 20.0 |
| Benzene, 1-ethy... | 7.670  | 16.5    | ng/ul | 1663140  | 1                     | 8.286  | 2012180 | 20.0 |
| (DEL) Alkane: S... | 7.910  | 65.2    | ng/ul | 6558650  | 1                     | 8.286  | 2012180 | 20.0 |
| Mesitylene         | 7.940  | 47.1    | ng/ul | 4742020  | 1                     | 8.286  | 2012180 | 20.0 |
| Benzene, 1,2,4-... | 8.410  | 19.4    | ng/ul | 1949190  | 1                     | 8.286  | 2012180 | 20.0 |
| (DEL) Alkane: C... | 8.586  | 10.2    | ng/ul | 1025350  | 1                     | 8.286  | 2012180 | 20.0 |
| Benzene, 1-meth... | 8.839  | 21.0    | ng/ul | 2116840  | 1                     | 8.286  | 2012180 | 20.0 |
| Carbonic acid, ... | 9.532  | 50.6    | ng/ul | 5094680  | 1                     | 8.286  | 2012180 | 20.0 |
| (DEL) Alkane: S... | 10.525 | 7.6     | ng/ul | 1150450  | 2                     | 11.124 | 3018160 | 20.0 |
| (DEL) Alkane: S... | 12.128 | 7.8     | ng/ul | 1179990  | 2                     | 11.124 | 3018160 | 20.0 |
| (DEL) Alkane: S... | 12.516 | 24.5    | ng/ul | 3701260  | 2                     | 11.124 | 3018160 | 20.0 |
| (DEL) Alkane: S... | 13.456 | 19.8    | ng/ul | 1467640  | 3                     | 14.919 | 1484270 | 20.0 |
| Naphthalene, 1,... | 14.084 | 27.6    | ng/ul | 2047180  | 3                     | 14.919 | 1484270 | 20.0 |
| Naphthalene, 1,... | 14.249 | 29.0    | ng/ul | 2152040  | 3                     | 14.919 | 1484270 | 20.0 |
| unknown-01         | 14.361 | 17.1    | ng/ul | 1266390  | 3                     | 14.919 | 1484270 | 20.0 |
| (DEL) Alkane: S... | 14.390 | 28.8    | ng/ul | 2138260  | 3                     | 14.919 | 1484270 | 20.0 |
| (DEL) Alkane: S... | 14.813 | 119.4   | ng/ul | 8863280  | 3                     | 14.919 | 1484270 | 20.0 |
| (DEL) Alkane: S... | 15.418 | 23.5    | ng/ul | 1742220  | 3                     | 14.919 | 1484270 | 20.0 |
| Naphthalene, 1,... | 15.535 | 12.8    | ng/ul | 946615   | 3                     | 14.919 | 1484270 | 20.0 |
| unknown-02         | 15.853 | 20.6    | ng/ul | 1528800  | 3                     | 14.919 | 1484270 | 20.0 |
| (DEL) Alkane: S... | 16.252 | 14.4    | ng/ul | 1072520  | 3                     | 14.919 | 1484270 | 20.0 |
| unknown-03         | 16.323 | 16.5    | ng/ul | 1293050  | 4                     | 17.680 | 1565170 | 20.0 |
| (DEL) Alkane: C... | 16.375 | 16.7    | ng/ul | 1304020  | 4                     | 17.680 | 1565170 | 20.0 |
| (DEL) Alkane: S... | 16.628 | 129.2   | ng/ul | 10107800 | 4                     | 17.680 | 1565170 | 20.0 |
| (DEL) Alkane: S... | 16.651 | 72.9    | ng/ul | 5705850  | 4                     | 17.680 | 1565170 | 20.0 |
| Benzene, (1-met... | 16.740 | 34.8    | ng/ul | 2726460  | 4                     | 17.680 | 1565170 | 20.0 |
| Tricyclo[5.2.1...  | 16.892 | 15.8    | ng/ul | 1239980  | 4                     | 17.680 | 1565170 | 20.0 |
| 9H-Fluorene, 2-... | 17.022 | 14.7    | ng/ul | 1153560  | 4                     | 17.680 | 1565170 | 20.0 |
| (DEL) Alkane: S... | 17.421 | 90.0    | ng/ul | 7044030  | 4                     | 17.680 | 1565170 | 20.0 |
| (DEL) Alkane: S... | 17.474 | 41.2    | ng/ul | 3221660  | 4                     | 17.680 | 1565170 | 20.0 |
| Benzene, (1-met... | 17.562 | 35.1    | ng/ul | 2746660  | 4                     | 17.680 | 1565170 | 20.0 |
| (DEL) Alkane: S... | 18.161 | 93.6    | ng/ul | 7327510  | 4                     | 17.680 | 1565170 | 20.0 |
| unknown-04         | 18.320 | 21.8    | ng/ul | 1706840  | 4                     | 17.680 | 1565170 | 20.0 |
| n-Hexadecanoic ... | 18.672 | 739.8   | ng/ul | 57894900 | 4                     | 17.680 | 1565170 | 20.0 |
| (DEL) Alkane: S... | 18.849 | 62.9    | ng/ul | 4920410  | 4                     | 17.680 | 1565170 | 20.0 |
| p-Dicyclohexylb... | 19.113 | 29.5    | ng/ul | 2307580  | 4                     | 17.680 | 1565170 | 20.0 |
| unknown-05         | 19.230 | 14.1    | ng/ul | 1102220  | 4                     | 17.680 | 1565170 | 20.0 |
| (DEL) Alkane: S... | 19.465 | 67.5    | ng/ul | 5283490  | 4                     | 17.680 | 1565170 | 20.0 |
| cyclohexane, [1... | 19.647 | 30.1    | ng/ul | 2353320  | 4                     | 17.680 | 1565170 | 20.0 |
| Phthalic acid, ... | 24.141 | 23.0    | ng/ul | 3148520  | 6                     | 25.545 | 2741140 | 20.0 |

|                    |        |      |       |         |   |        |         |      |
|--------------------|--------|------|-------|---------|---|--------|---------|------|
| (DEL) Alkane: S... | 24.535 | 20.0 | ng/ul | 2741140 | 6 | 25.545 | 2741140 | 20.0 |
| Phthalic acid, ... | 27.349 | 26.4 | ng/ul | 3620130 | 6 | 25.545 | 2741140 | 20.0 |

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD3DL