

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP122422\  
 Data File : BP013333.D  
 Acq On : 25 Dec 2022 10:58  
 Operator : CG/JU  
 Sample : N6187-08  
 Misc : GCMS CONFIRMATION  
 ALS Vial : 46 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 A43X5

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120722.M  
 Title : SVOA CALIBRATION

Signal : TIC: BP013333.D\data.ms

| peak<br># | R.T.<br>min | first<br>scan | max<br>scan | last<br>scan | PK<br>TY | peak<br>height | corr.<br>area | corr.<br>% max. | % of<br>total |
|-----------|-------------|---------------|-------------|--------------|----------|----------------|---------------|-----------------|---------------|
| 1         | 3.211       | 3             | 4           | 7            | rVB      | 282946         | 176713        | 0.31%           | 0.044%        |
| 2         | 3.252       | 7             | 11          | 15           | rVB      | 11824960       | 12677748      | 22.59%          | 3.131%        |
| 3         | 3.599       | 66            | 70          | 75           | rVB      | 403886         | 491286        | 0.88%           | 0.121%        |
| 4         | 4.081       | 148           | 152         | 157          | rVB      | 52145          | 63509         | 0.11%           | 0.016%        |
| 5         | 7.934       | 800           | 807         | 818          | rBV      | 2095721        | 3333926       | 5.94%           | 0.823%        |
| 6         | 10.752      | 1271          | 1286        | 1296         | rBV      | 2767205        | 4618018       | 8.23%           | 1.140%        |
| 7         | 14.581      | 1929          | 1937        | 1950         | rBV      | 3745354        | 5282018       | 9.41%           | 1.304%        |
| 8         | 14.704      | 1950          | 1958        | 1965         | rVV      | 1339968        | 1951954       | 3.48%           | 0.482%        |
| 9         | 15.428      | 2075          | 2081        | 2088         | rBV      | 72651          | 107056        | 0.19%           | 0.026%        |
| 10        | 15.516      | 2090          | 2096        | 2103         | rBV      | 348901         | 468758        | 0.84%           | 0.116%        |
| 11        | 15.839      | 2144          | 2151        | 2157         | rBV      | 11147547       | 15122247      | 26.94%          | 3.735%        |
| 12        | 16.281      | 2219          | 2226        | 2233         | rBV      | 2736336        | 3885541       | 6.92%           | 0.960%        |
| 13        | 16.445      | 2248          | 2254        | 2262         | rBV      | 5039932        | 6951573       | 12.38%          | 1.717%        |
| 14        | 16.569      | 2268          | 2275        | 2289         | rBV      | 20473460       | 28828564      | 51.36%          | 7.119%        |
| 15        | 16.863      | 2319          | 2325        | 2332         | rBV      | 3203326        | 4422868       | 7.88%           | 1.092%        |
| 16        | 17.110      | 2363          | 2367        | 2373         | rVB2     | 61065          | 88942         | 0.16%           | 0.022%        |
| 17        | 17.222      | 2378          | 2386        | 2389         | rBV      | 24404364       | 37157879      | 66.20%          | 9.176%        |
| 18        | 17.257      | 2389          | 2392        | 2398         | rVV      | 10888484       | 13442929      | 23.95%          | 3.320%        |
| 19        | 17.328      | 2398          | 2404        | 2419         | rVB3     | 7177327        | 16252423      | 28.95%          | 4.014%        |
| 20        | 17.516      | 2429          | 2436        | 2450         | rVB      | 15485925       | 23023458      | 41.02%          | 5.686%        |
| 21        | 17.628      | 2450          | 2455        | 2463         | rBV      | 103726         | 190422        | 0.34%           | 0.047%        |
| 22        | 17.704      | 2463          | 2468        | 2476         | rBV2     | 327635         | 506309        | 0.90%           | 0.125%        |
| 23        | 17.792      | 2477          | 2483        | 2486         | rBV      | 4795481        | 6585523       | 11.73%          | 1.626%        |
| 24        | 17.822      | 2486          | 2488        | 2494         | rVB      | 2251479        | 2383555       | 4.25%           | 0.589%        |
| 25        | 17.881      | 2494          | 2498        | 2500         | rBV      | 51554          | 60631         | 0.11%           | 0.015%        |
| 26        | 17.939      | 2500          | 2508        | 2515         | rVV      | 26440617       | 56131742      | 100.00%         | 13.862%       |
| 27        | 18.063      | 2521          | 2529        | 2535         | rBV2     | 17419160       | 26083389      | 46.47%          | 6.441%        |
| 28        | 18.128      | 2535          | 2540        | 2544         | rVV      | 981720         | 1212413       | 2.16%           | 0.299%        |
| 29        | 18.181      | 2544          | 2549        | 2554         | rVV      | 9965207        | 12652894      | 22.54%          | 3.125%        |
| 30        | 18.233      | 2554          | 2558        | 2564         | rVB      | 3568392        | 4389214       | 7.82%           | 1.084%        |
| 31        | 18.339      | 2570          | 2576        | 2580         | rBV      | 1375624        | 1844787       | 3.29%           | 0.456%        |
| 32        | 18.392      | 2580          | 2585        | 2591         | rVV      | 12849749       | 16807299      | 29.94%          | 4.151%        |
| 33        | 18.451      | 2591          | 2595        | 2598         | rVV      | 9435852        | 11627292      | 20.71%          | 2.871%        |
| 34        | 18.492      | 2598          | 2602        | 2607         | rVB      | 8073919        | 9819304       | 17.49%          | 2.425%        |
| 35        | 18.669      | 2625          | 2632        | 2636         | rBV      | 11821972       | 16581170      | 29.54%          | 4.095%        |
| 36        | 18.710      | 2636          | 2639        | 2644         | rVV      | 4169379        | 5533055       | 9.86%           | 1.366%        |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP122422\  
 Data File : BP013333.D  
 Acq On : 25 Dec 2022 10:58  
 Operator : CG/JU  
 Sample : N6187-08  
 Misc : GCMS CONFIRMATION  
 ALS Vial : 46 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 A43X5

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120722.M  
 Title : SVOA CALIBRATION

|    |        |      |      |      |      |         |          |        |        |
|----|--------|------|------|------|------|---------|----------|--------|--------|
| 37 | 18.763 | 2644 | 2648 | 2651 | rVV  | 4951368 | 6172048  | 11.00% | 1.524% |
| 38 | 18.798 | 2651 | 2654 | 2657 | rVV  | 3762627 | 4772179  | 8.50%  | 1.179% |
| 39 | 18.833 | 2657 | 2660 | 2667 | rVV  | 8498272 | 10711415 | 19.08% | 2.645% |
| 40 | 18.898 | 2667 | 2671 | 2673 | rVV  | 126328  | 170982   | 0.30%  | 0.042% |
| 41 | 18.933 | 2673 | 2677 | 2682 | rVV  | 2583093 | 3102884  | 5.53%  | 0.766% |
| 42 | 19.004 | 2684 | 2689 | 2696 | rVB  | 332352  | 432895   | 0.77%  | 0.107% |
| 43 | 19.075 | 2696 | 2701 | 2705 | rBV  | 309981  | 396749   | 0.71%  | 0.098% |
| 44 | 19.133 | 2706 | 2711 | 2715 | rVV  | 2348330 | 2852133  | 5.08%  | 0.704% |
| 45 | 19.180 | 2715 | 2719 | 2722 | rVV  | 3848648 | 5309629  | 9.46%  | 1.311% |
| 46 | 19.216 | 2722 | 2725 | 2731 | rVV3 | 3279797 | 4810614  | 8.57%  | 1.188% |
| 47 | 19.292 | 2734 | 2738 | 2742 | rVV  | 346852  | 429498   | 0.77%  | 0.106% |
| 48 | 19.328 | 2742 | 2744 | 2753 | rVB4 | 52328   | 64569    | 0.12%  | 0.016% |
| 49 | 19.422 | 2754 | 2760 | 2767 | rBV  | 789503  | 1717308  | 3.06%  | 0.424% |
| 50 | 19.480 | 2767 | 2770 | 2777 | rVV  | 631122  | 819293   | 1.46%  | 0.202% |
| 51 | 19.545 | 2777 | 2781 | 2786 | rVB  | 175760  | 221418   | 0.39%  | 0.055% |
| 52 | 19.739 | 2810 | 2814 | 2819 | rBV2 | 115900  | 165038   | 0.29%  | 0.041% |
| 53 | 19.810 | 2823 | 2826 | 2830 | rVV  | 165051  | 190063   | 0.34%  | 0.047% |
| 54 | 19.857 | 2832 | 2834 | 2839 | rVB2 | 92514   | 105926   | 0.19%  | 0.026% |
| 55 | 19.916 | 2840 | 2844 | 2848 | rBV  | 278645  | 327219   | 0.58%  | 0.081% |
| 56 | 19.963 | 2849 | 2852 | 2857 | rVV2 | 104199  | 135179   | 0.24%  | 0.033% |
| 57 | 20.057 | 2866 | 2868 | 2873 | rVB4 | 72892   | 70430    | 0.13%  | 0.017% |
| 58 | 20.222 | 2892 | 2896 | 2901 | rBV2 | 234241  | 297997   | 0.53%  | 0.074% |
| 59 | 20.522 | 2944 | 2947 | 2952 | rBV2 | 151860  | 175212   | 0.31%  | 0.043% |
| 60 | 21.433 | 3097 | 3102 | 3112 | rBV  | 3762668 | 4932312  | 8.79%  | 1.218% |
| 61 | 23.916 | 3517 | 3524 | 3535 | rVB2 | 2792510 | 5792180  | 10.32% | 1.430% |

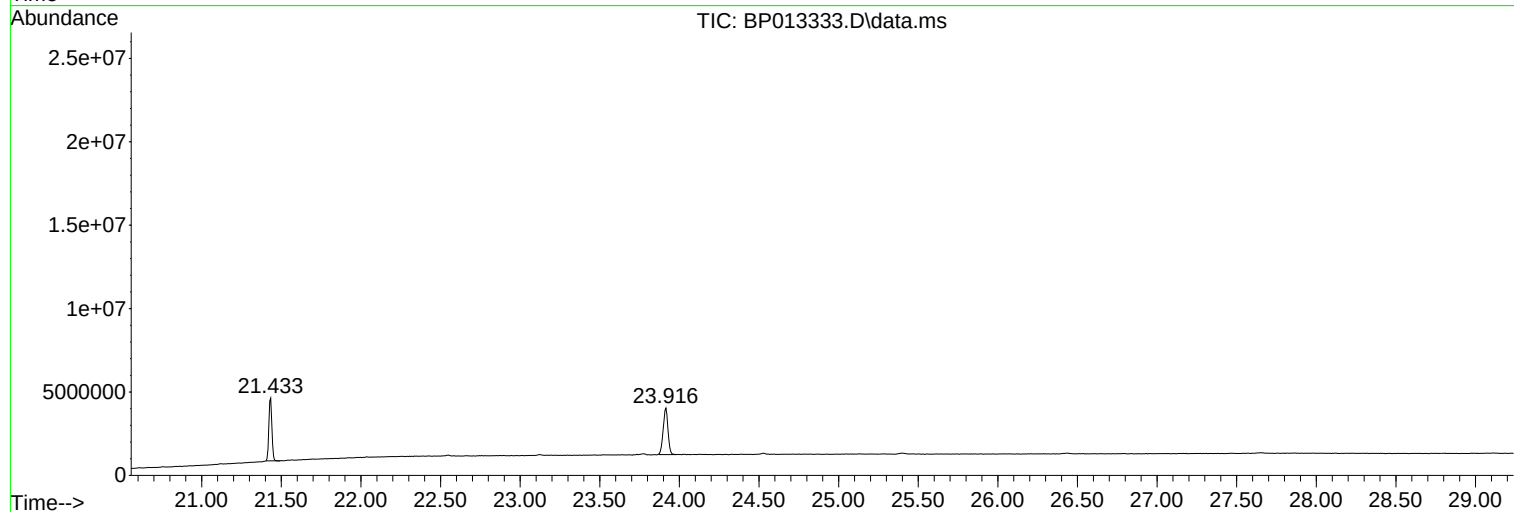
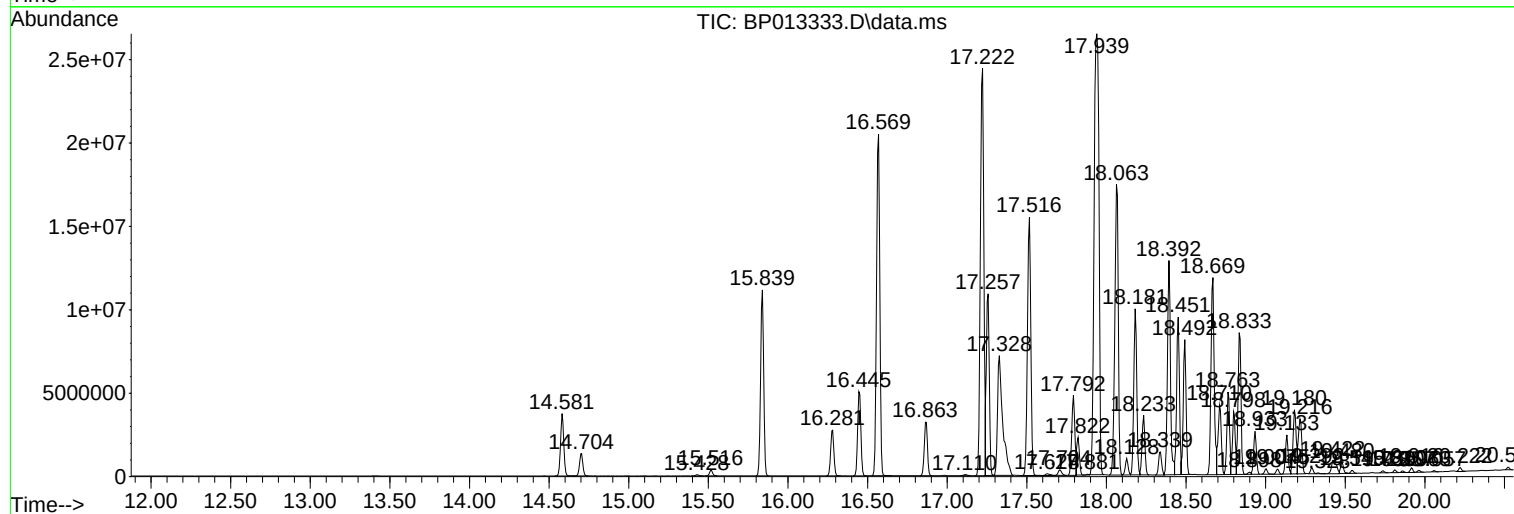
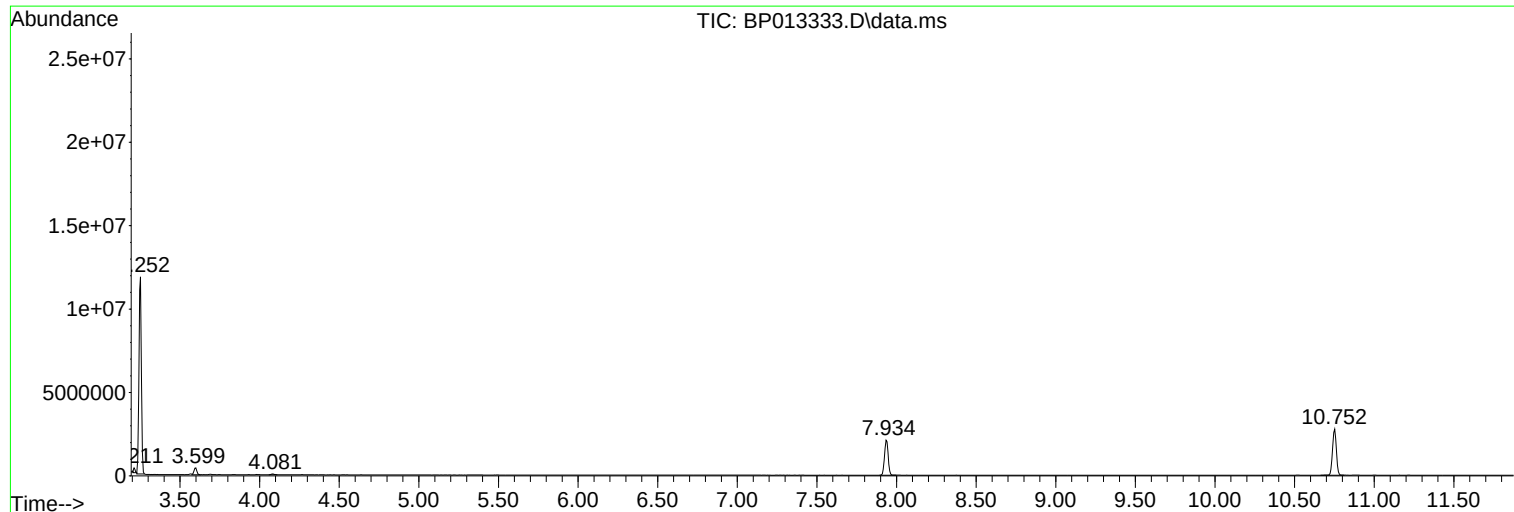
Sum of corrected areas: 404931579

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP122422\  
Data File : BP013333.D  
Acq On : 25 Dec 2022 10:58  
Operator : CG/JU  
Sample : N6187-08  
Misc : GCMS CONFIRMATION  
ALS Vial : 46 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
A43X5

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120722.M  
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP122422\  
Data File : BP013333.D  
Acq On : 25 Dec 2022 10:58  
Operator : CG/JU  
Sample : N6187-08  
Misc : GCMS CONFIRMATION  
ALS Vial : 46 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
A43X5

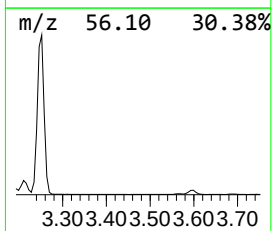
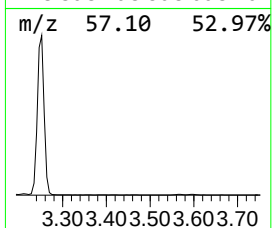
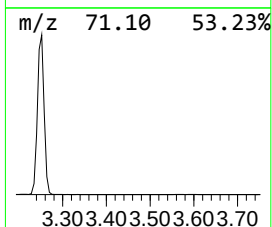
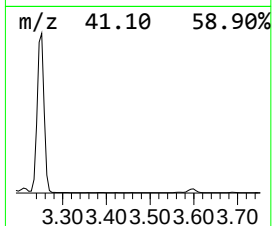
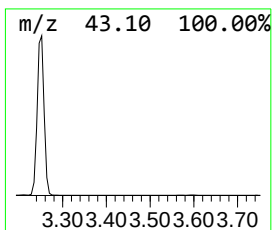
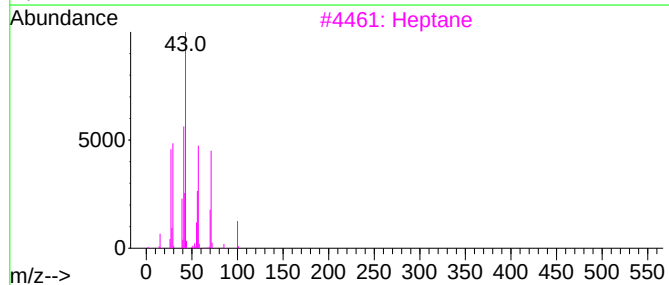
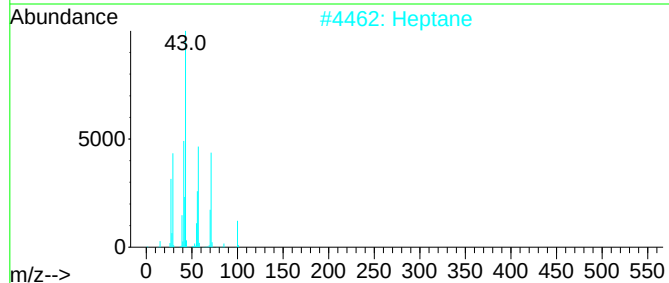
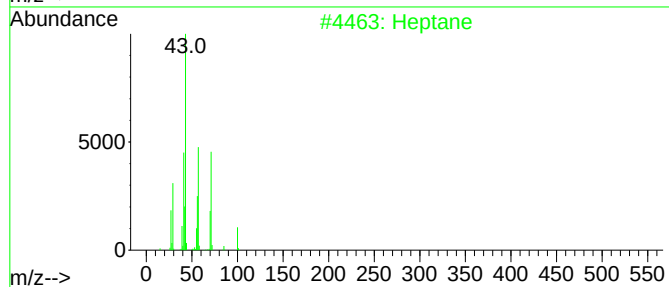
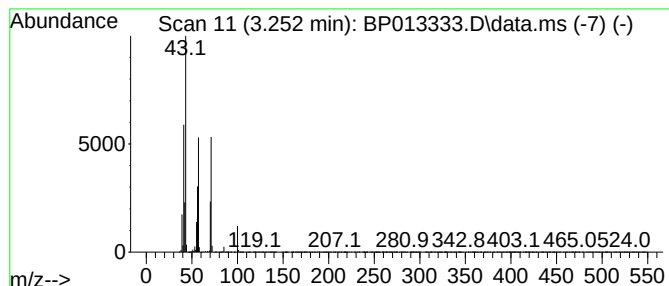
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120722.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 1

| R.T.  | EstConc     | Area     | Relative to ISTD       | R.T.  |
|-------|-------------|----------|------------------------|-------|
| 3.252 | 76.05 ng/ul | 12677700 | 1,4-Dichlorobenzene-d4 | 7.934 |

| Hit# | of                                 | 5 | Tentative ID | MW  | MolForm  | CAS#         | Qual |
|------|------------------------------------|---|--------------|-----|----------|--------------|------|
| 1    | Heptane                            |   |              | 100 | C7H16    | 000142-82-5  | 95   |
| 2    | Heptane                            |   |              | 100 | C7H16    | 000142-82-5  | 94   |
| 3    | Heptane                            |   |              | 100 | C7H16    | 000142-82-5  | 91   |
| 4    | Heptane                            |   |              | 100 | C7H16    | 000142-82-5  | 83   |
| 5    | Oxalic acid, isobutyl pentyl ester |   |              | 216 | C11H20O4 | 1000309-37-0 | 64   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP122422\  
Data File : BP013333.D  
Acq On : 25 Dec 2022 10:58  
Operator : CG/JU  
Sample : N6187-08  
Misc : GCMS CONFIRMATION  
ALS Vial : 46 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
A43X5

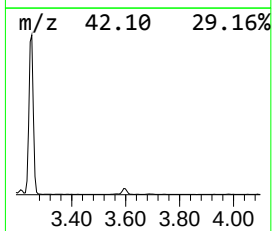
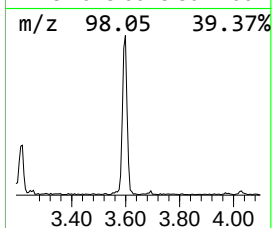
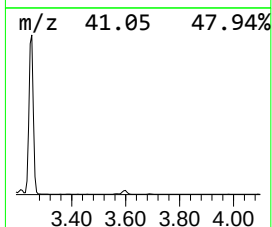
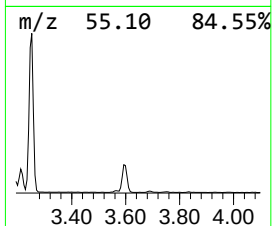
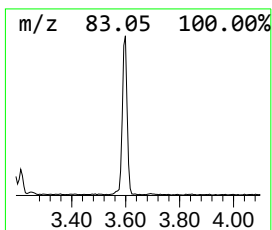
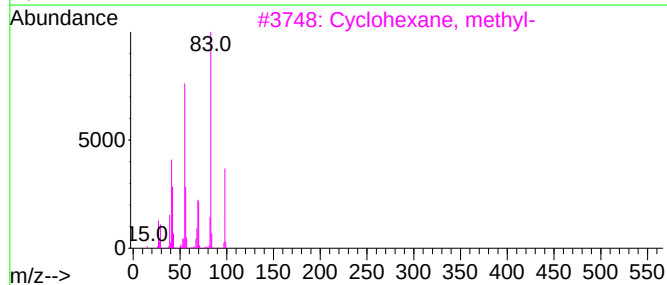
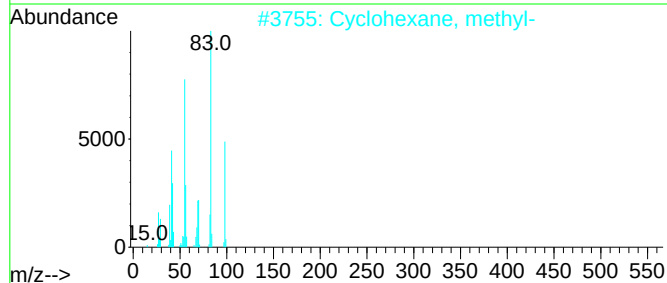
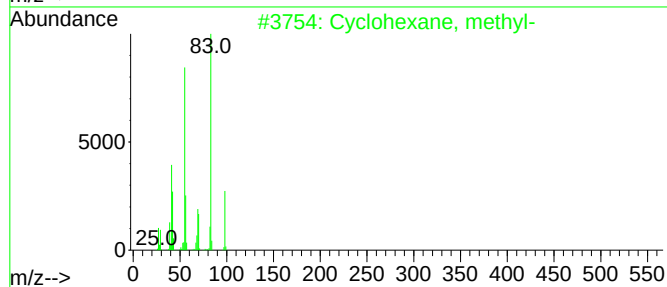
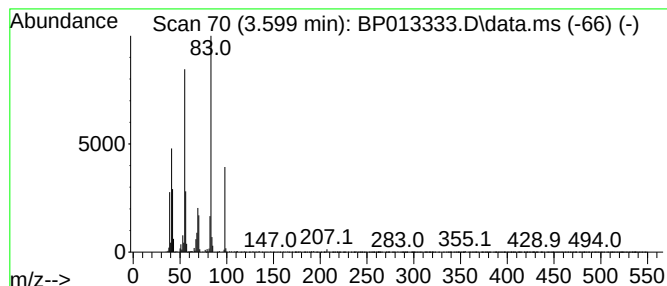
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120722.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 2 (DEL) Alkane: Cyclic3.599 Concentration Rank 30

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 3.599 | 2.95 ng/ul | 491286 | 1,4-Dichlorobenzene-d4 | 7.934 |

| Hit# | of | 5 | Tentative ID         | MW | MolForm | CAS#        | Qual |
|------|----|---|----------------------|----|---------|-------------|------|
| 1    |    |   | Cyclohexane, methyl- | 98 | C7H14   | 000108-87-2 | 94   |
| 2    |    |   | Cyclohexane, methyl- | 98 | C7H14   | 000108-87-2 | 94   |
| 3    |    |   | Cyclohexane, methyl- | 98 | C7H14   | 000108-87-2 | 94   |
| 4    |    |   | Cyclohexane, methyl- | 98 | C7H14   | 000108-87-2 | 94   |
| 5    |    |   | Cyclohexane, methyl- | 98 | C7H14   | 000108-87-2 | 94   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP122422\  
Data File : BP013333.D  
Acq On : 25 Dec 2022 10:58  
Operator : CG/JU  
Sample : N6187-08  
Misc : GCMS CONFIRMATION  
ALS Vial : 46 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
A43X5

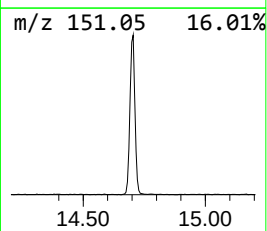
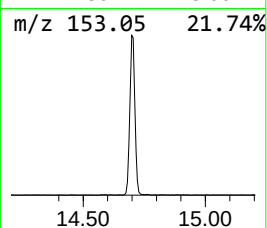
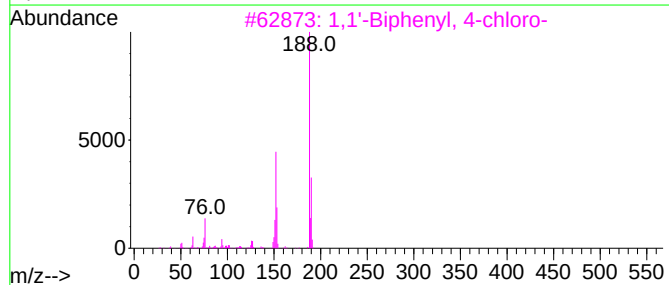
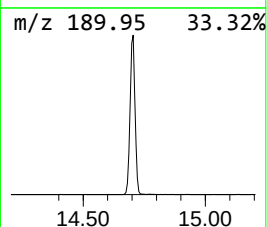
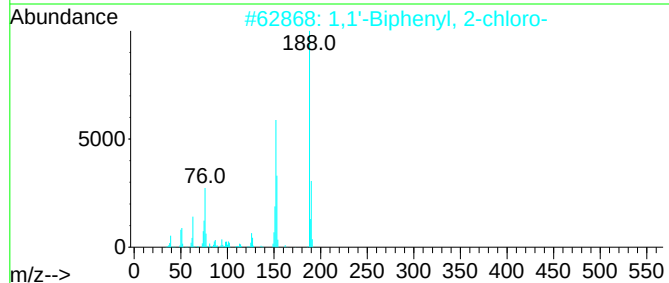
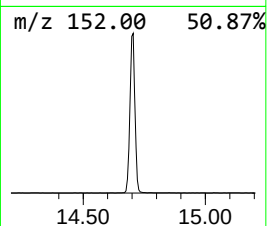
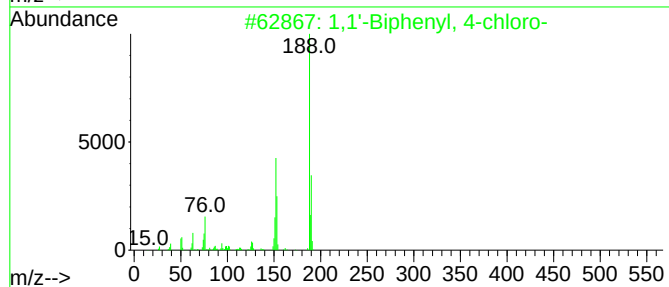
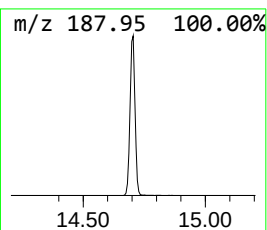
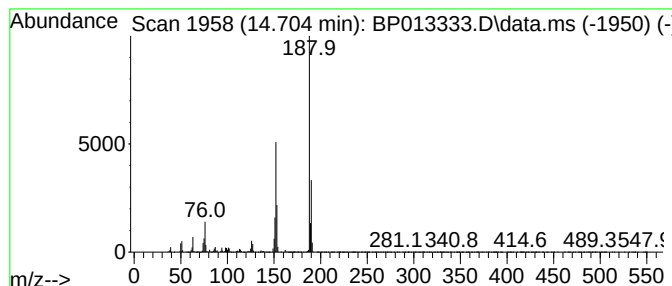
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120722.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 3 1,1'-Biphenyl, 4-chloro- Concentration Rank 18

| R.T.   | EstConc    | Area    | Relative to ISTD | R.T.   |
|--------|------------|---------|------------------|--------|
| 14.704 | 7.39 ng/ul | 1951950 | Acenaphthene-d10 | 14.581 |

| Hit# | of                       | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|--------------------------|---|--------------|-----|---------|-------------|------|
| 1    | 1,1'-Biphenyl, 4-chloro- |   |              | 188 | C12H9Cl | 002051-62-9 | 98   |
| 2    | 1,1'-Biphenyl, 2-chloro- |   |              | 188 | C12H9Cl | 002051-60-7 | 97   |
| 3    | 1,1'-Biphenyl, 4-chloro- |   |              | 188 | C12H9Cl | 002051-62-9 | 95   |
| 4    | 1,1'-Biphenyl, 4-chloro- |   |              | 188 | C12H9Cl | 002051-62-9 | 95   |
| 5    | 1,1'-Biphenyl, 2-chloro- |   |              | 188 | C12H9Cl | 002051-60-7 | 93   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP122422\  
Data File : BP013333.D  
Acq On : 25 Dec 2022 10:58  
Operator : CG/JU  
Sample : N6187-08  
Misc : GCMS CONFIRMATION  
ALS Vial : 46 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
A43X5

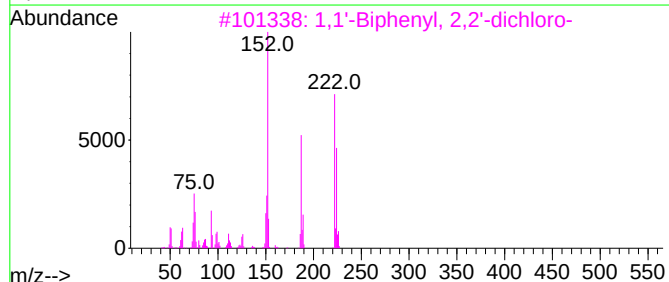
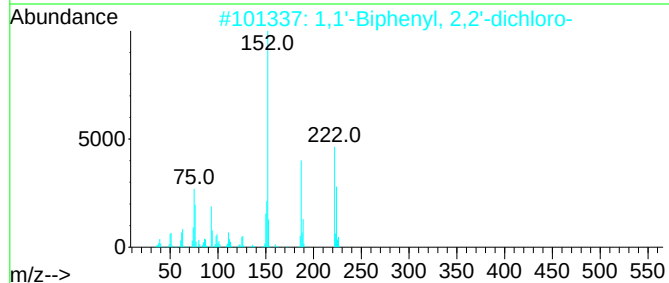
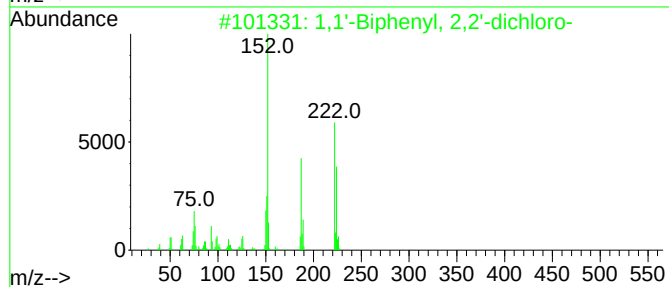
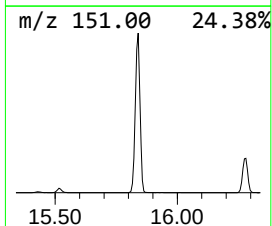
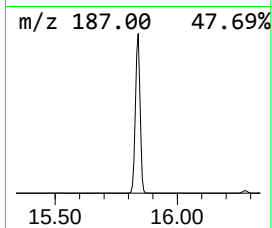
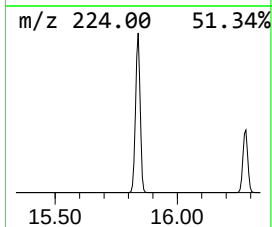
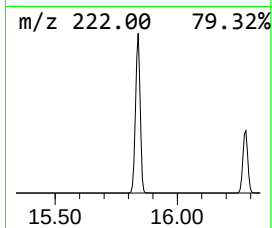
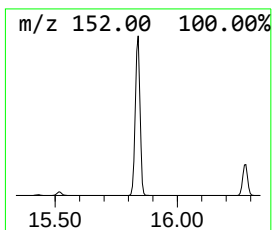
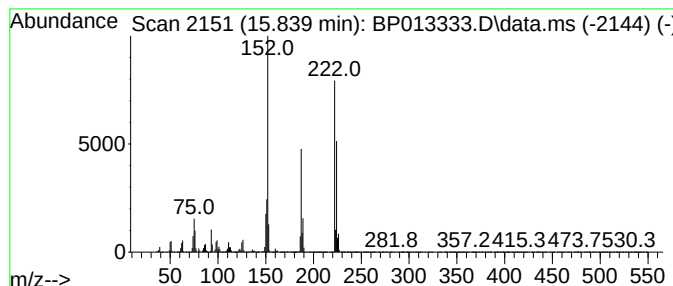
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120722.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 4 1,1'-Biphenyl, 2,2'-dichloro- Concentration Rank 3

| R.T.   | EstConc     | Area     | Relative to ISTD | R.T.   |
|--------|-------------|----------|------------------|--------|
| 15.839 | 57.26 ng/ul | 15122200 | Acenaphthene-d10 | 14.581 |

| Hit# | of                            | 5   | Tentative ID | MW          | MolForm | CAS# | Qual |
|------|-------------------------------|-----|--------------|-------------|---------|------|------|
| 1    | 1,1'-Biphenyl, 2,2'-dichloro- | 222 | C12H8Cl2     | 013029-08-8 | 99      |      |      |
| 2    | 1,1'-Biphenyl, 2,2'-dichloro- | 222 | C12H8Cl2     | 013029-08-8 | 99      |      |      |
| 3    | 1,1'-Biphenyl, 2,2'-dichloro- | 222 | C12H8Cl2     | 013029-08-8 | 98      |      |      |
| 4    | 1,1'-Biphenyl, 2,4'-dichloro- | 222 | C12H8Cl2     | 034883-43-7 | 98      |      |      |
| 5    | 1,1'-Biphenyl, 3,5-dichloro-  | 222 | C12H8Cl2     | 034883-41-5 | 98      |      |      |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP122422\  
Data File : BP013333.D  
Acq On : 25 Dec 2022 10:58  
Operator : CG/JU  
Sample : N6187-08  
Misc : GCMS CONFIRMATION  
ALS Vial : 46 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
A43X5

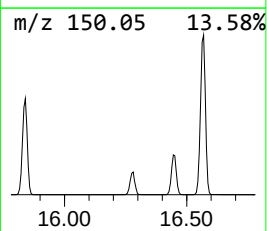
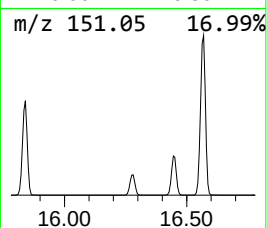
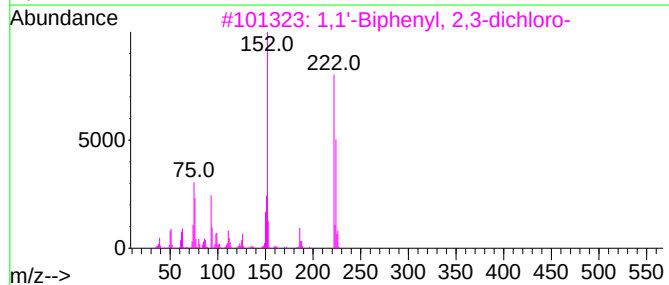
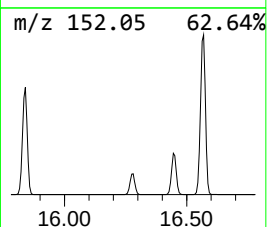
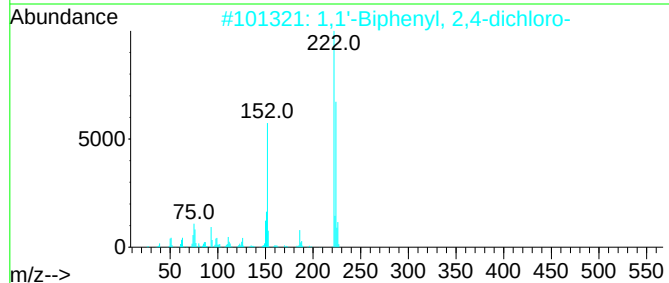
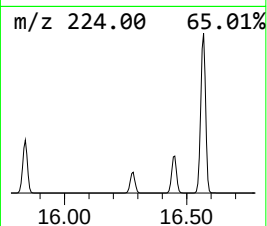
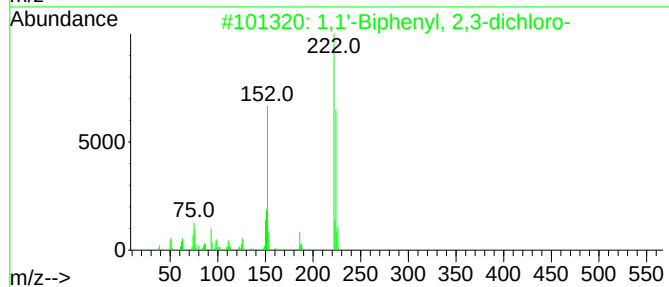
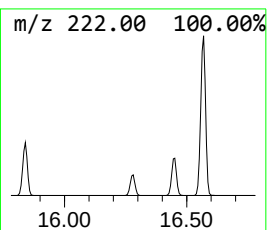
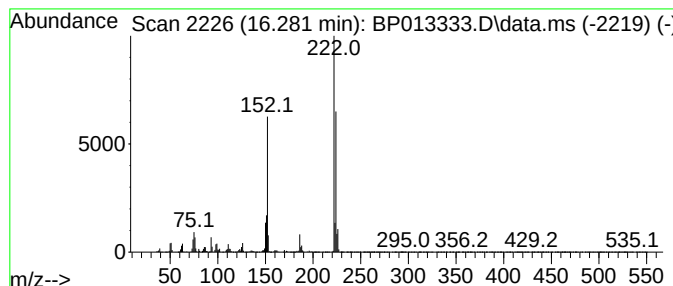
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120722.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 5 1,1'-Biphenyl, 2,3-dichloro- Concentration Rank 26

| R.T.   | EstConc    | Area    | Relative to ISTD | R.T.   |
|--------|------------|---------|------------------|--------|
| 16.281 | 4.78 ng/ul | 3885540 | Phenanthrene-d10 | 17.345 |

| Hit# | of                            | 5   | Tentative ID | MW          | MolForm | CAS# | Qual |
|------|-------------------------------|-----|--------------|-------------|---------|------|------|
| 1    | 1,1'-Biphenyl, 2,3-dichloro-  | 222 | C12H8Cl2     | 016605-91-7 | 99      |      |      |
| 2    | 1,1'-Biphenyl, 2,4-dichloro-  | 222 | C12H8Cl2     | 033284-50-3 | 99      |      |      |
| 3    | 1,1'-Biphenyl, 2,3-dichloro-  | 222 | C12H8Cl2     | 016605-91-7 | 97      |      |      |
| 4    | 1,1'-Biphenyl, 2,4-dichloro-  | 222 | C12H8Cl2     | 033284-50-3 | 97      |      |      |
| 5    | 1,1'-Biphenyl, 4,4'-dichloro- | 222 | C12H8Cl2     | 002050-68-2 | 97      |      |      |





Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP122422\  
Data File : BP013333.D  
Acq On : 25 Dec 2022 10:58  
Operator : CG/JU  
Sample : N6187-08  
Misc : GCMS CONFIRMATION  
ALS Vial : 46 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
A43X5

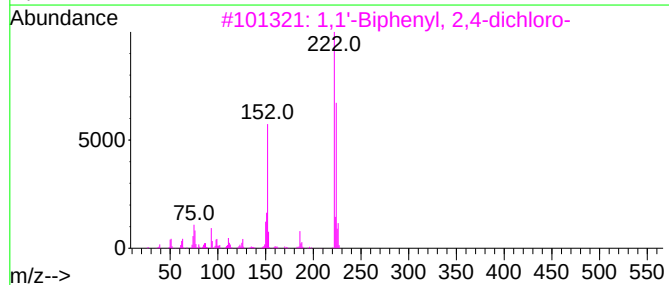
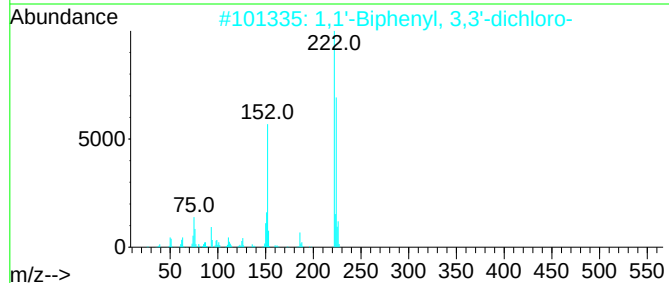
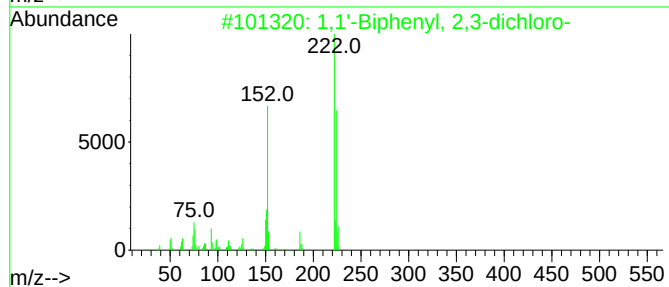
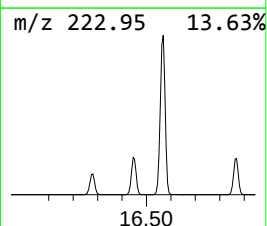
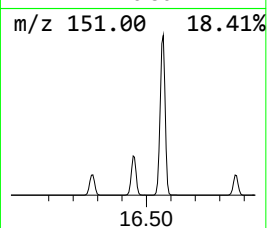
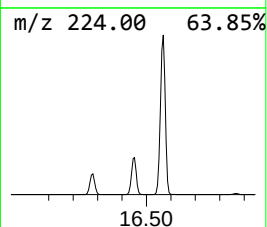
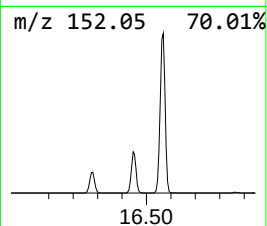
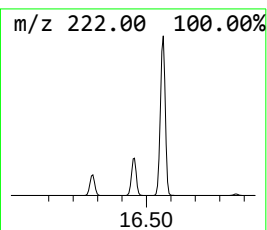
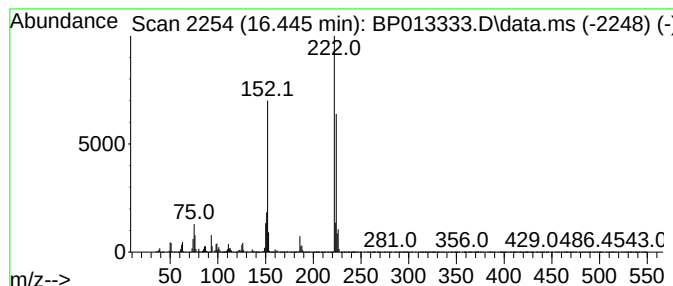
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120722.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*

Peak Number 6 1,1'-Biphenyl, 3,3'-dichloro- Concentration Rank 15

| R.T.      | EstConc                       | Area    | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------|---------|------------------|-------------|------|
| 16.445    | 8.55 ng/ul                    | 6951570 | Phenanthrene-d10 | 17.345      |      |
| Hit# of 5 | Tentative ID                  | MW      | MolForm          | CAS#        | Qual |
| 1         | 1,1'-Biphenyl, 2,3-dichloro-  | 222     | C12H8Cl2         | 016605-91-7 | 99   |
| 2         | 1,1'-Biphenyl, 3,3'-dichloro- | 222     | C12H8Cl2         | 002050-67-1 | 99   |
| 3         | 1,1'-Biphenyl, 2,4-dichloro-  | 222     | C12H8Cl2         | 033284-50-3 | 99   |
| 4         | 1,1'-Biphenyl, 2,5-dichloro-  | 222     | C12H8Cl2         | 034883-39-1 | 98   |
| 5         | 1,1'-Biphenyl, 2,4'-dichloro- | 222     | C12H8Cl2         | 034883-43-7 | 98   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP122422\  
Data File : BP013333.D  
Acq On : 25 Dec 2022 10:58  
Operator : CG/JU  
Sample : N6187-08  
Misc : GCMS CONFIRMATION  
ALS Vial : 46 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
A43X5

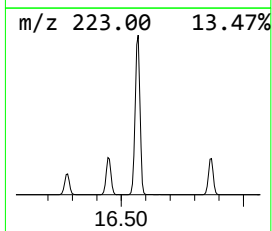
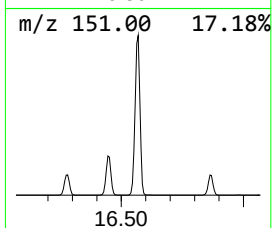
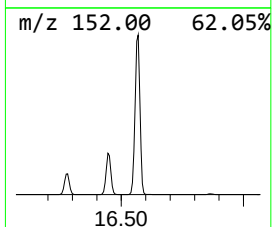
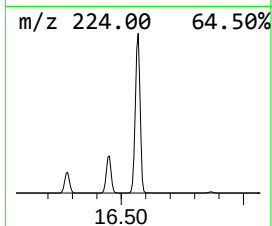
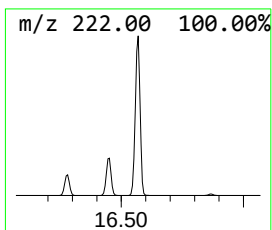
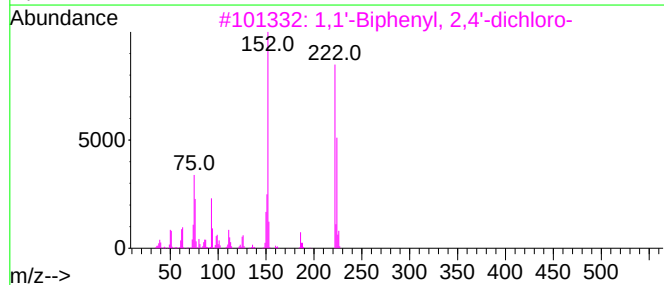
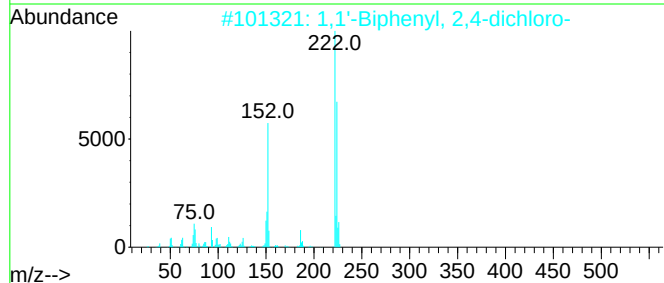
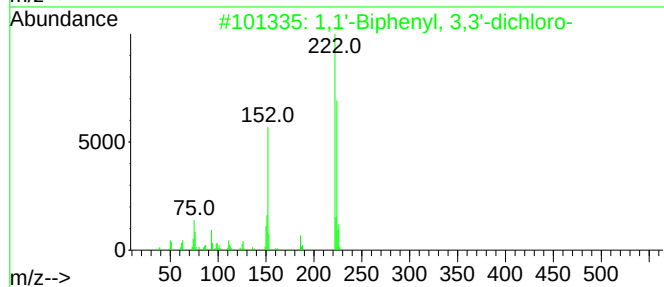
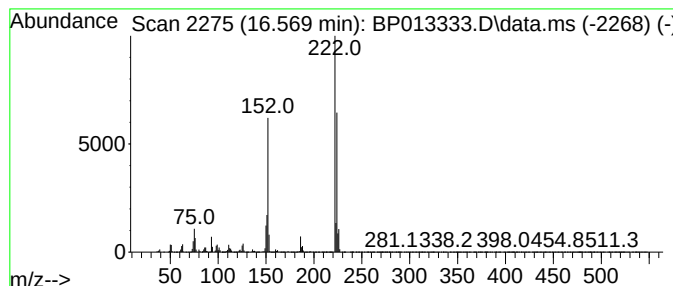
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120722.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 7 1,1'-Biphenyl, 2,4-dichloro- Concentration Rank 5

| R.T.   | EstConc     | Area     | Relative to ISTD | R.T.   |
|--------|-------------|----------|------------------|--------|
| 16.569 | 35.48 ng/ul | 28828600 | Phenanthrene-d10 | 17.345 |

| Hit# | of                            | 5   | Tentative ID | MW          | MolForm | CAS# | Qual |
|------|-------------------------------|-----|--------------|-------------|---------|------|------|
| 1    | 1,1'-Biphenyl, 3,3'-dichloro- | 222 | C12H8Cl2     | 002050-67-1 | 99      |      |      |
| 2    | 1,1'-Biphenyl, 2,4-dichloro-  | 222 | C12H8Cl2     | 033284-50-3 | 99      |      |      |
| 3    | 1,1'-Biphenyl, 2,4'-dichloro- | 222 | C12H8Cl2     | 034883-43-7 | 98      |      |      |
| 4    | 1,1'-Biphenyl, 4,4'-dichloro- | 222 | C12H8Cl2     | 002050-68-2 | 98      |      |      |
| 5    | 1,1'-Biphenyl, 3,3'-dichloro- | 222 | C12H8Cl2     | 002050-67-1 | 97      |      |      |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP122422\  
Data File : BP013333.D  
Acq On : 25 Dec 2022 10:58  
Operator : CG/JU  
Sample : N6187-08  
Misc : GCMS CONFIRMATION  
ALS Vial : 46 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
A43X5

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120722.M  
Quant Title : SVOA CALIBRATION

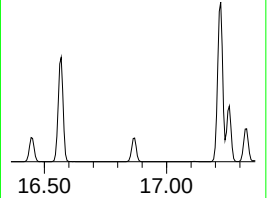
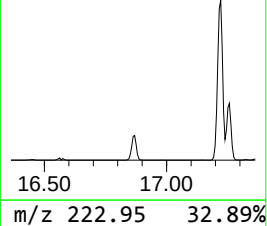
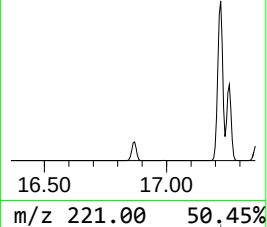
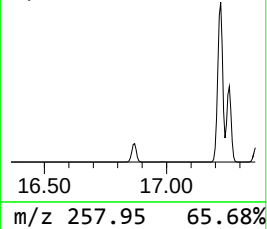
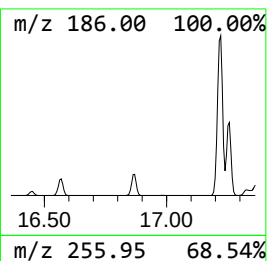
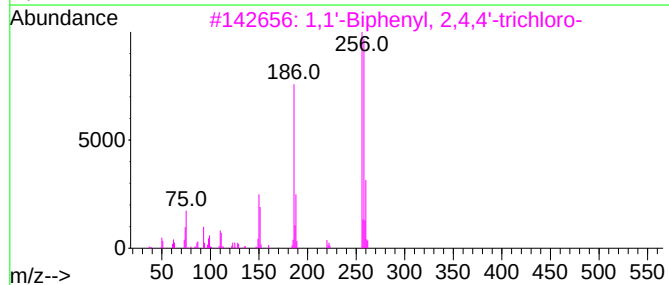
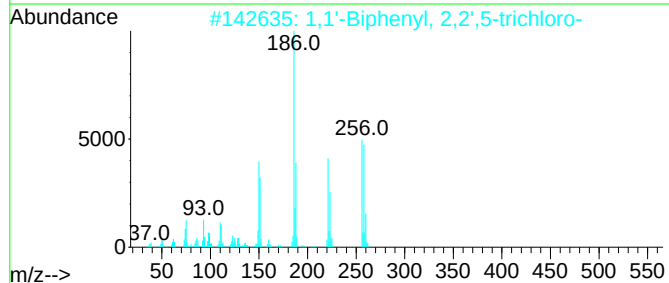
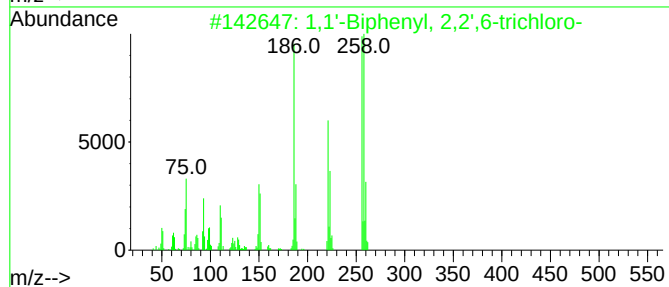
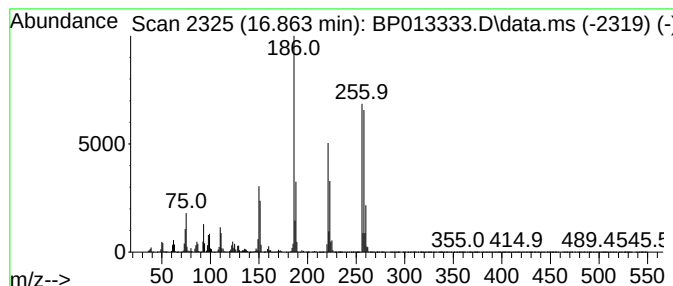
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*

Peak Number 8 1,1'-Biphenyl, 2,2',6-trich... Concentration Rank 24

| R.T.   | EstConc    | Area    | Relative to ISTD | R.T.   |
|--------|------------|---------|------------------|--------|
| 16.863 | 5.44 ng/ul | 4422870 | Phenanthrene-d10 | 17.345 |

| Hit# | of                               | 5   | Tentative ID | MW          | MolForm | CAS# | Qual |
|------|----------------------------------|-----|--------------|-------------|---------|------|------|
| 1    | 1,1'-Biphenyl, 2,2',6-trichloro- | 256 | C12H7Cl3     | 038444-73-4 | 99      |      |      |
| 2    | 1,1'-Biphenyl, 2,2',5-trichloro- | 256 | C12H7Cl3     | 037680-65-2 | 99      |      |      |
| 3    | 1,1'-Biphenyl, 2,4,4'-trichloro- | 256 | C12H7Cl3     | 007012-37-5 | 99      |      |      |
| 4    | 1,1'-Biphenyl, 2,2',5-trichloro- | 256 | C12H7Cl3     | 037680-65-2 | 99      |      |      |
| 5    | 1,1'-Biphenyl, 2,2',5-trichloro- | 256 | C12H7Cl3     | 037680-65-2 | 98      |      |      |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP122422\  
Data File : BP013333.D  
Acq On : 25 Dec 2022 10:58  
Operator : CG/JU  
Sample : N6187-08  
Misc : GCMS CONFIRMATION  
ALS Vial : 46 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
A43X5

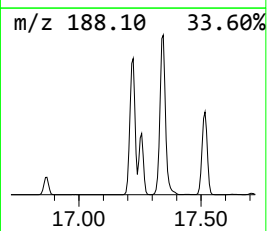
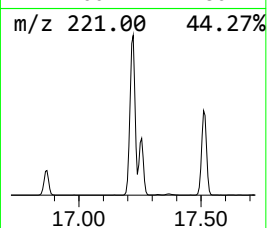
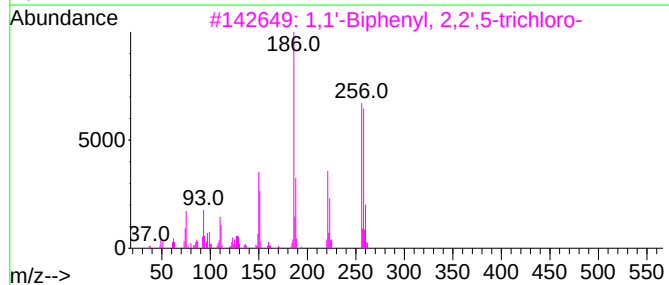
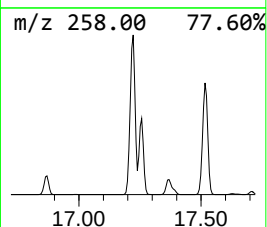
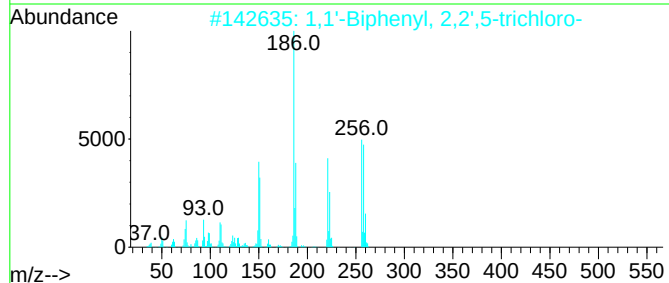
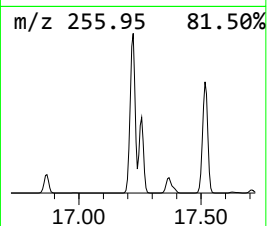
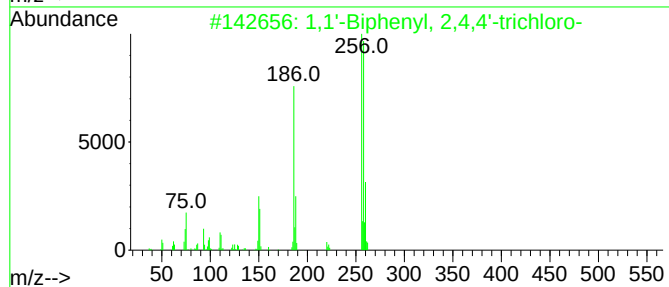
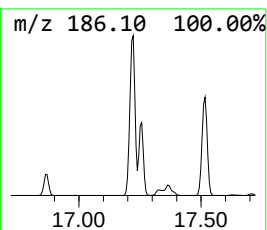
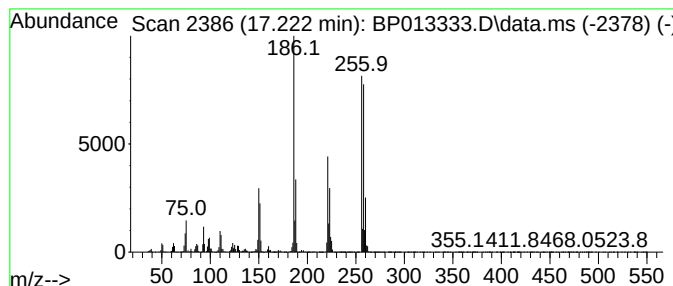
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120722.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 9 1,1'-Biphenyl, 2,4,4'-trich... Concentration Rank 4

| R.T.   | EstConc     | Area     | Relative to ISTD | R.T.   |
|--------|-------------|----------|------------------|--------|
| 17.222 | 45.73 ng/ul | 37157900 | Phenanthrene-d10 | 17.345 |

| Hit# | of                               | 5   | Tentative ID | MW          | MolForm | CAS# | Qual |
|------|----------------------------------|-----|--------------|-------------|---------|------|------|
| 1    | 1,1'-Biphenyl, 2,4,4'-trichloro- | 256 | C12H7Cl3     | 007012-37-5 | 99      |      |      |
| 2    | 1,1'-Biphenyl, 2,2',5-trichloro- | 256 | C12H7Cl3     | 037680-65-2 | 99      |      |      |
| 3    | 1,1'-Biphenyl, 2,2',5-trichloro- | 256 | C12H7Cl3     | 037680-65-2 | 99      |      |      |
| 4    | 1,1'-Biphenyl, 2,2',6-trichloro- | 256 | C12H7Cl3     | 038444-73-4 | 99      |      |      |
| 5    | 2,2',3-Trichloro-1,1'-biphenyl   | 256 | C12H7Cl3     | 038444-78-9 | 99      |      |      |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP122422\  
Data File : BP013333.D  
Acq On : 25 Dec 2022 10:58  
Operator : CG/JU  
Sample : N6187-08  
Misc : GCMS CONFIRMATION  
ALS Vial : 46 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
A43X5

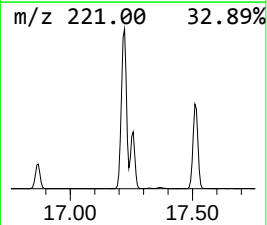
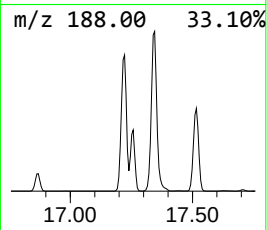
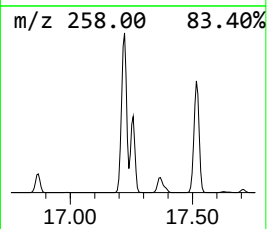
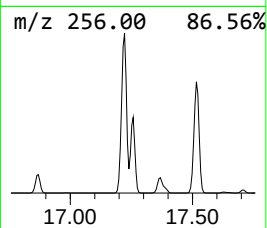
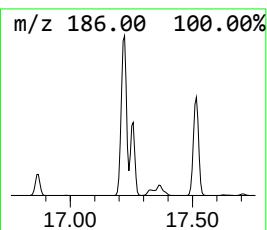
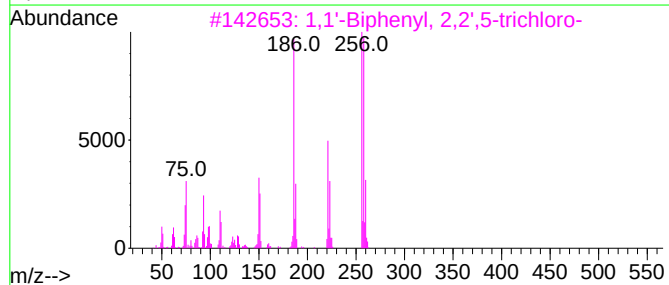
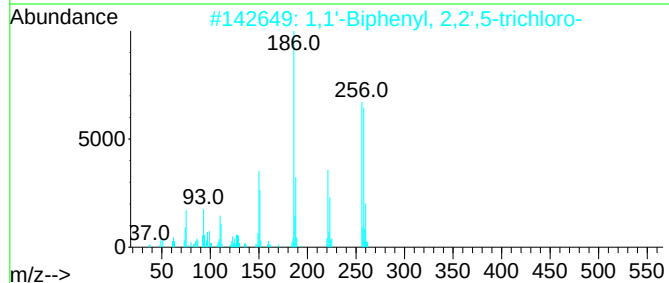
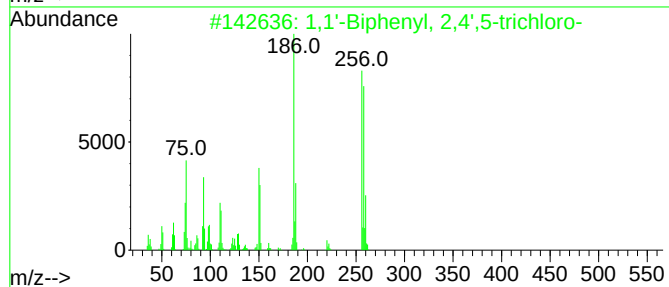
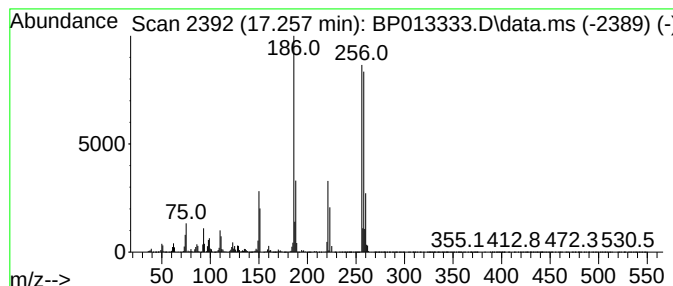
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120722.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 10 1,1'-Biphenyl, 2,4',5-trich... Concentration Rank 10

| R.T.   | EstConc     | Area     | Relative to ISTD | R.T.   |
|--------|-------------|----------|------------------|--------|
| 17.257 | 16.54 ng/ul | 13442900 | Phenanthrene-d10 | 17.345 |

| Hit# | of                               | 5   | Tentative ID | MW          | MolForm | CAS# | Qual |
|------|----------------------------------|-----|--------------|-------------|---------|------|------|
| 1    | 1,1'-Biphenyl, 2,4',5-trichloro- | 256 | C12H7Cl3     | 016606-02-3 | 97      |      |      |
| 2    | 1,1'-Biphenyl, 2,2',5-trichloro- | 256 | C12H7Cl3     | 037680-65-2 | 96      |      |      |
| 3    | 1,1'-Biphenyl, 2,2',5-trichloro- | 256 | C12H7Cl3     | 037680-65-2 | 96      |      |      |
| 4    | 1,1'-Biphenyl, 2,4',5-trichloro- | 256 | C12H7Cl3     | 016606-02-3 | 95      |      |      |
| 5    | 1,1'-Biphenyl, 2,4,4'-trichloro- | 256 | C12H7Cl3     | 007012-37-5 | 95      |      |      |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP122422\  
Data File : BP013333.D  
Acq On : 25 Dec 2022 10:58  
Operator : CG/JU  
Sample : N6187-08  
Misc : GCMS CONFIRMATION  
ALS Vial : 46 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
A43X5

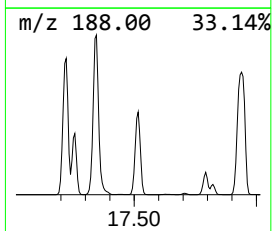
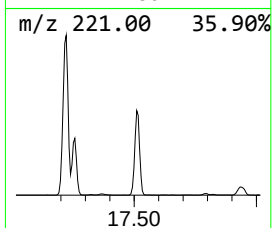
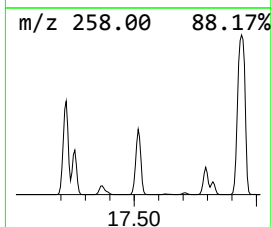
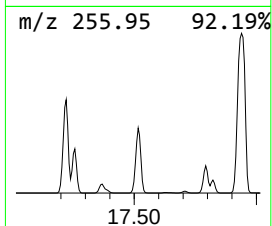
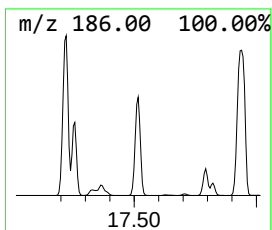
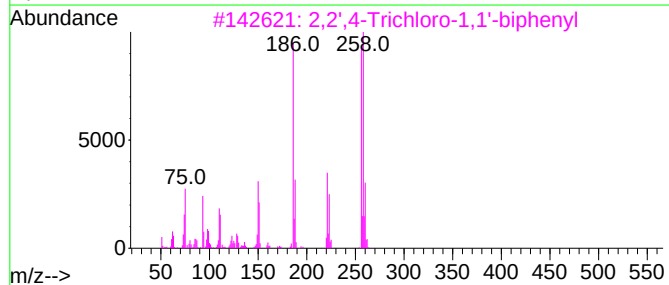
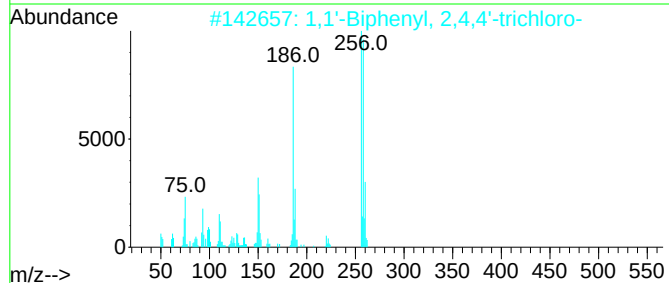
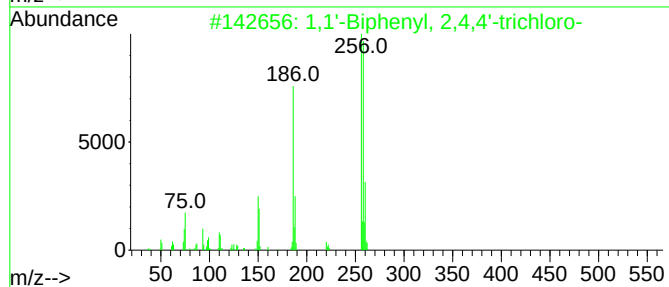
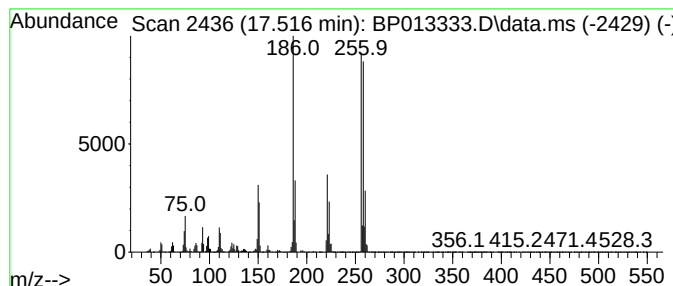
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120722.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 11 2,2',4-Trichloro-1,1'-biphenyl Concentration Rank 7

| R.T.   | EstConc     | Area     | Relative to ISTD | R.T.   |
|--------|-------------|----------|------------------|--------|
| 17.516 | 28.33 ng/ul | 23023500 | Phenanthrene-d10 | 17.345 |

| Hit# | of 5                             | Tentative ID | MW       | MolForm     | CAS# | Qual |
|------|----------------------------------|--------------|----------|-------------|------|------|
| 1    | 1,1'-Biphenyl, 2,4,4'-trichloro- | 256          | C12H7Cl3 | 007012-37-5 | 99   |      |
| 2    | 1,1'-Biphenyl, 2,4,4'-trichloro- | 256          | C12H7Cl3 | 007012-37-5 | 99   |      |
| 3    | 2,2',4-Trichloro-1,1'-biphenyl   | 256          | C12H7Cl3 | 037680-66-3 | 99   |      |
| 4    | 1,1'-Biphenyl, 2,2',5-trichloro- | 256          | C12H7Cl3 | 037680-65-2 | 99   |      |
| 5    | 1,1'-Biphenyl, 2,2',5-trichloro- | 256          | C12H7Cl3 | 037680-65-2 | 99   |      |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP122422\  
Data File : BP013333.D  
Acq On : 25 Dec 2022 10:58  
Operator : CG/JU  
Sample : N6187-08  
Misc : GCMS CONFIRMATION  
ALS Vial : 46 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
A43X5

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120722.M  
Quant Title : SVOA CALIBRATION

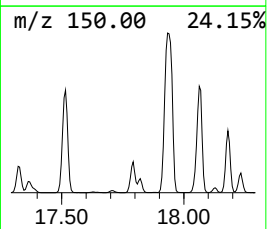
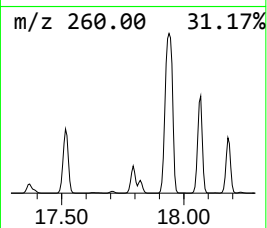
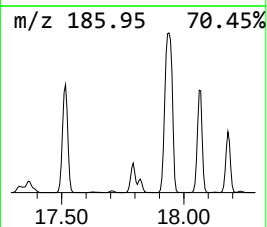
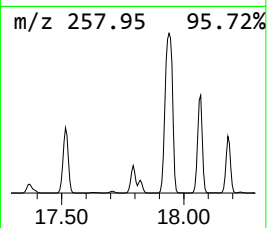
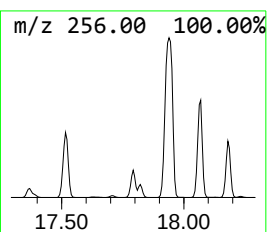
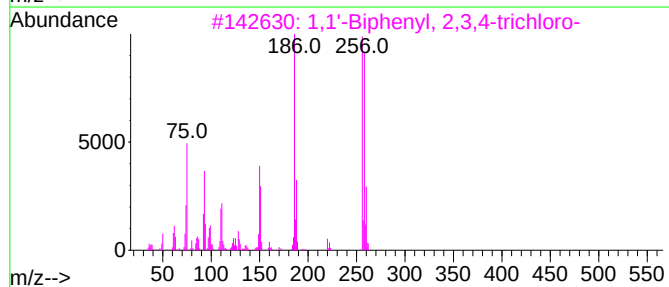
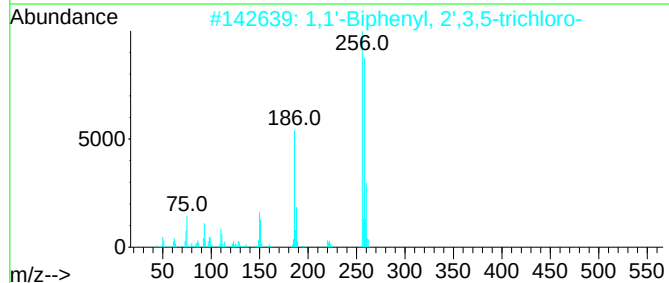
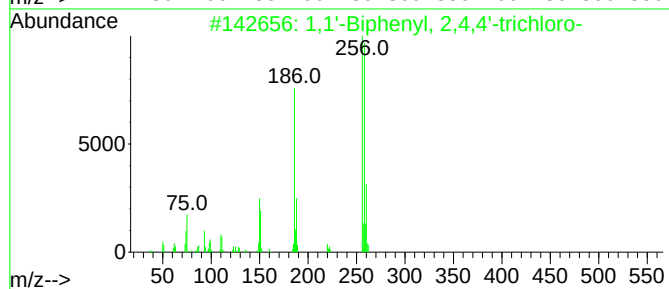
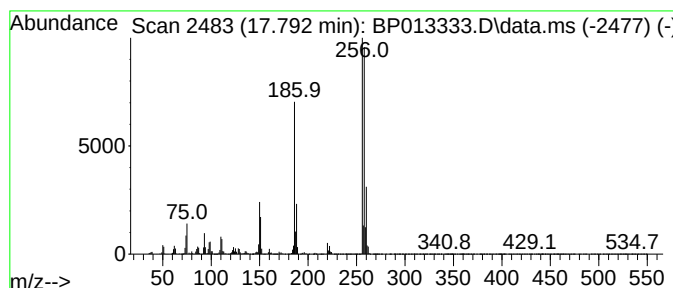
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*

Peak Number 12 1,1'-Biphenyl, 2',3,5-trich... Concentration Rank 16

| R.T.   | EstConc    | Area    | Relative to ISTD | R.T.   |
|--------|------------|---------|------------------|--------|
| 17.792 | 8.10 ng/ul | 6585520 | Phenanthrene-d10 | 17.345 |

| Hit# | of 5                             | Tentative ID | MW       | MolForm     | CAS# | Qual |
|------|----------------------------------|--------------|----------|-------------|------|------|
| 1    | 1,1'-Biphenyl, 2,4,4'-trichloro- | 256          | C12H7Cl3 | 007012-37-5 | 99   |      |
| 2    | 1,1'-Biphenyl, 2',3,5-trichloro- | 256          | C12H7Cl3 | 037680-68-5 | 99   |      |
| 3    | 1,1'-Biphenyl, 2,3,4-trichloro-  | 256          | C12H7Cl3 | 055702-46-0 | 99   |      |
| 4    | 1,1'-Biphenyl, 2,3',5-trichloro- | 256          | C12H7Cl3 | 038444-81-4 | 99   |      |
| 5    | 1,1'-Biphenyl, 2,4',5-trichloro- | 256          | C12H7Cl3 | 016606-02-3 | 99   |      |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP122422\  
Data File : BP013333.D  
Acq On : 25 Dec 2022 10:58  
Operator : CG/JU  
Sample : N6187-08  
Misc : GCMS CONFIRMATION  
ALS Vial : 46 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
A43X5

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120722.M  
Quant Title : SVOA CALIBRATION

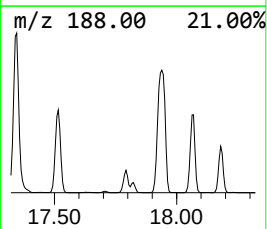
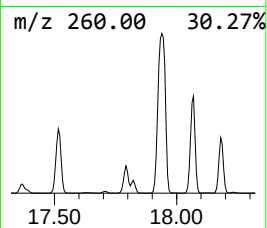
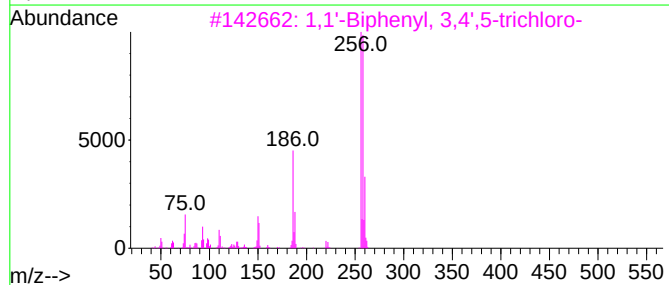
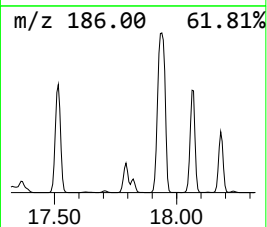
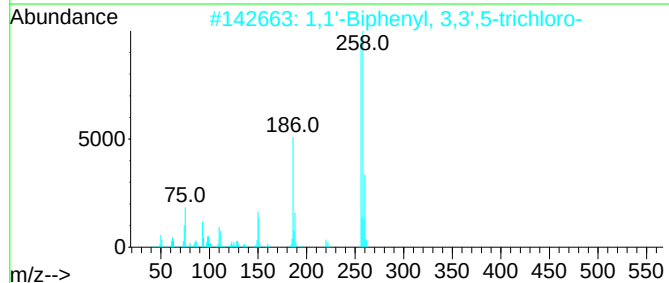
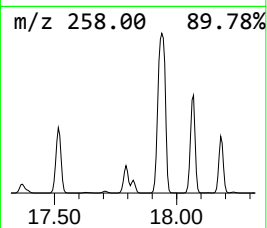
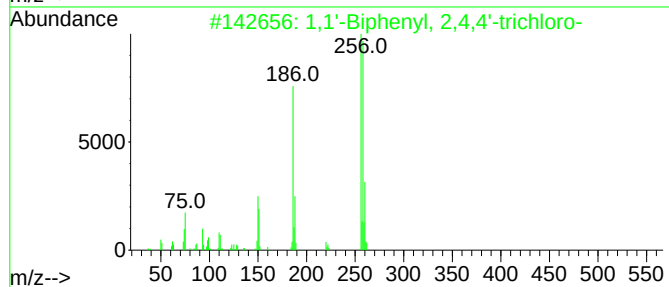
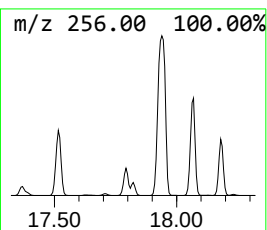
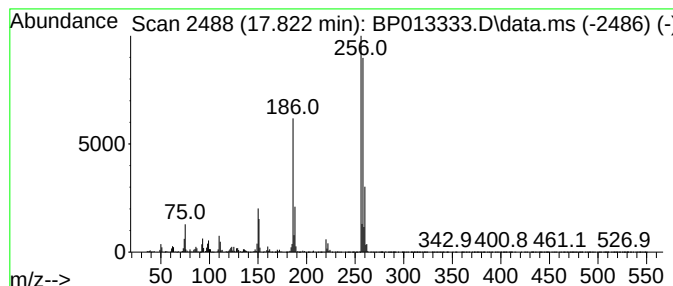
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*

Peak Number 13 1,1'-Biphenyl, 3,3',5-trich... Concentration Rank 31

| R.T.   | EstConc    | Area    | Relative to ISTD | R.T.   |
|--------|------------|---------|------------------|--------|
| 17.822 | 2.93 ng/ul | 2383560 | Phenanthrene-d10 | 17.345 |

| Hit# | of 5                             | Tentative ID | MW       | MolForm     | CAS# | Qual |
|------|----------------------------------|--------------|----------|-------------|------|------|
| 1    | 1,1'-Biphenyl, 2,4,4'-trichloro- | 256          | C12H7Cl3 | 007012-37-5 | 99   |      |
| 2    | 1,1'-Biphenyl, 3,3',5-trichloro- | 256          | C12H7Cl3 | 038444-87-0 | 99   |      |
| 3    | 1,1'-Biphenyl, 3,4',5-trichloro- | 256          | C12H7Cl3 | 038444-88-1 | 99   |      |
| 4    | 1,1'-Biphenyl, 2,3,3'-trichloro- | 256          | C12H7Cl3 | 038444-84-7 | 99   |      |
| 5    | 1,1'-Biphenyl, 2,3',5-trichloro- | 256          | C12H7Cl3 | 038444-81-4 | 99   |      |





Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP122422\  
Data File : BP013333.D  
Acq On : 25 Dec 2022 10:58  
Operator : CG/JU  
Sample : N6187-08  
Misc : GCMS CONFIRMATION  
ALS Vial : 46 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
A43X5

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120722.M  
Quant Title : SVOA CALIBRATION

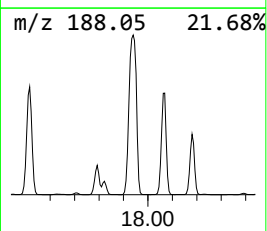
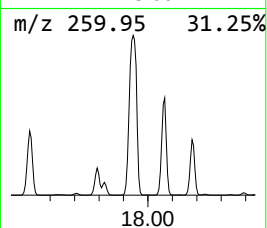
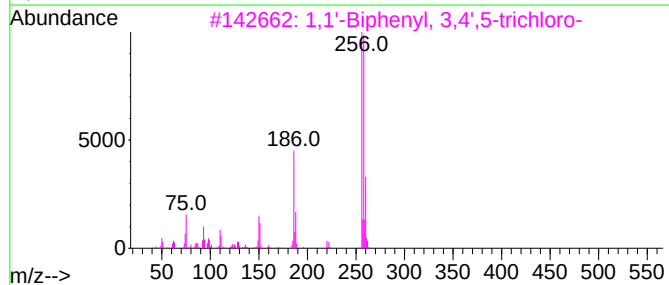
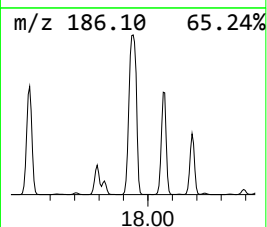
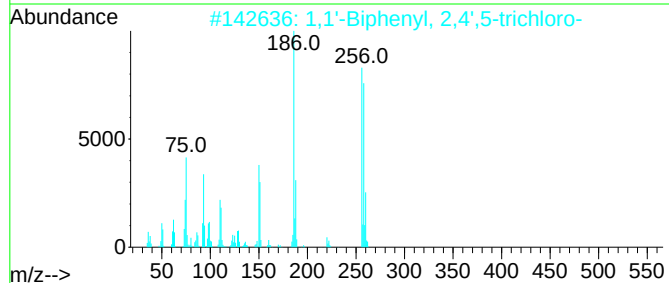
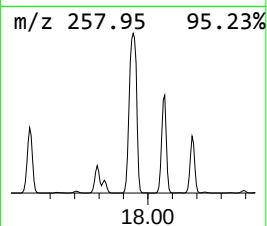
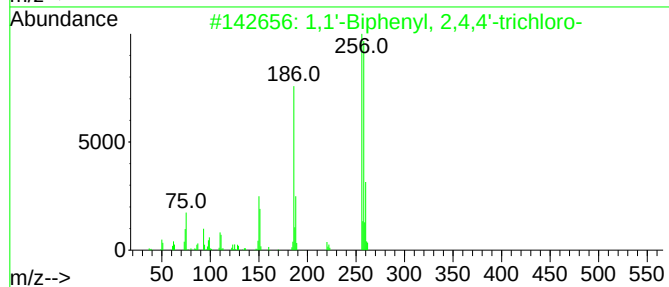
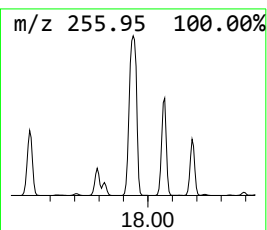
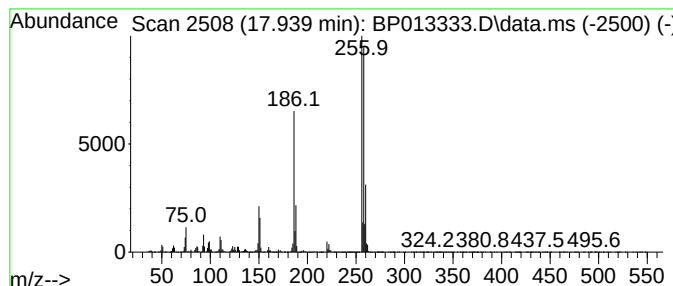
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*

Peak Number 14 1,1'-Biphenyl, 3,4',5-trich... Concentration Rank 2

| R.T.   | EstConc     | Area     | Relative to ISTD | R.T.   |
|--------|-------------|----------|------------------|--------|
| 17.939 | 69.07 ng/ul | 56131700 | Phenanthrene-d10 | 17.345 |

| Hit# | of 5                             | Tentative ID | MW       | MolForm     | CAS# | Qual |
|------|----------------------------------|--------------|----------|-------------|------|------|
| 1    | 1,1'-Biphenyl, 2,4,4'-trichloro- | 256          | C12H7Cl3 | 007012-37-5 | 99   |      |
| 2    | 1,1'-Biphenyl, 2,4',5-trichloro- | 256          | C12H7Cl3 | 016606-02-3 | 99   |      |
| 3    | 1,1'-Biphenyl, 3,4',5-trichloro- | 256          | C12H7Cl3 | 038444-88-1 | 99   |      |
| 4    | 1,1'-Biphenyl, 2,3,3'-trichloro- | 256          | C12H7Cl3 | 038444-84-7 | 99   |      |
| 5    | 1,1'-Biphenyl, 3,3',5-trichloro- | 256          | C12H7Cl3 | 038444-87-0 | 99   |      |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP122422\  
Data File : BP013333.D  
Acq On : 25 Dec 2022 10:58  
Operator : CG/JU  
Sample : N6187-08  
Misc : GCMS CONFIRMATION  
ALS Vial : 46 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
A43X5

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120722.M  
Quant Title : SVOA CALIBRATION

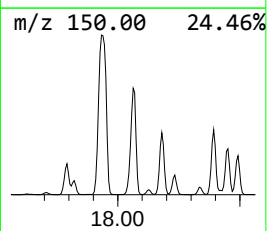
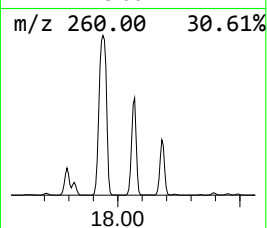
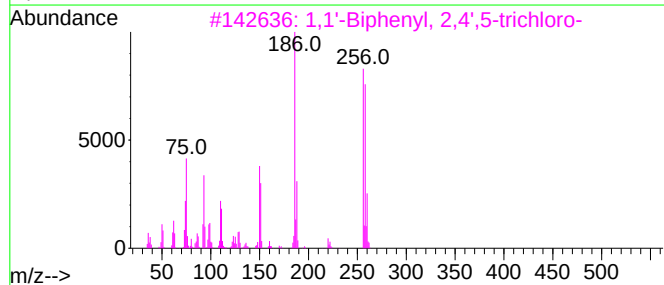
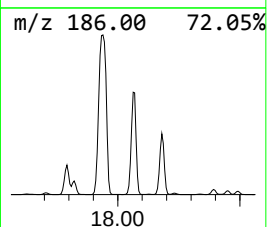
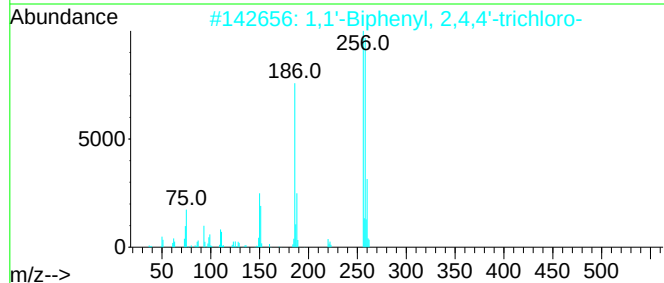
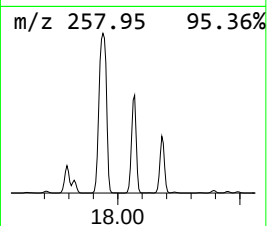
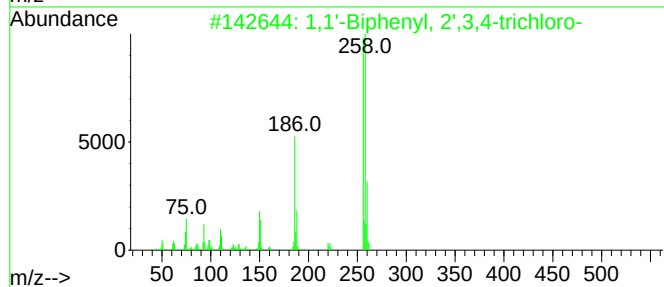
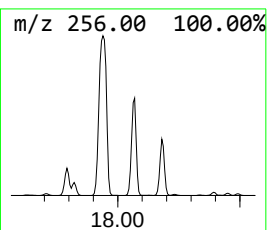
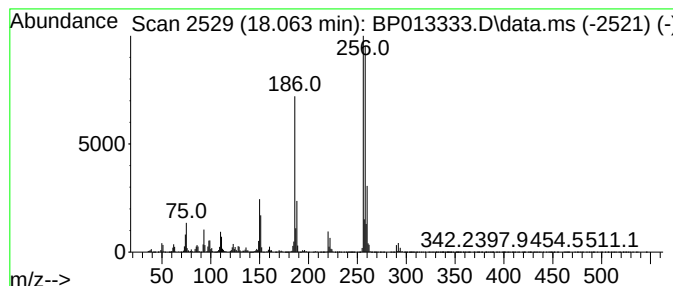
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*

Peak Number 15 1,1'-Biphenyl, 2',3,4-trich... Concentration Rank 6

| R.T.   | EstConc     | Area     | Relative to ISTD | R.T.   |
|--------|-------------|----------|------------------|--------|
| 18.063 | 32.10 ng/ul | 26083400 | Phenanthrene-d10 | 17.345 |

| Hit# | of                               | 5   | Tentative ID | MW          | MolForm | CAS# | Qual |
|------|----------------------------------|-----|--------------|-------------|---------|------|------|
| 1    | 1,1'-Biphenyl, 2',3,4-trichloro- | 256 | C12H7Cl3     | 038444-86-9 | 99      |      |      |
| 2    | 1,1'-Biphenyl, 2,4,4'-trichloro- | 256 | C12H7Cl3     | 007012-37-5 | 99      |      |      |
| 3    | 1,1'-Biphenyl, 2,4',5-trichloro- | 256 | C12H7Cl3     | 016606-02-3 | 99      |      |      |
| 4    | 1,1'-Biphenyl, 3,4',5-trichloro- | 256 | C12H7Cl3     | 038444-88-1 | 99      |      |      |
| 5    | 1,1'-Biphenyl, 2,3,3'-trichloro- | 256 | C12H7Cl3     | 038444-84-7 | 99      |      |      |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP122422\  
Data File : BP013333.D  
Acq On : 25 Dec 2022 10:58  
Operator : CG/JU  
Sample : N6187-08  
Misc : GCMS CONFIRMATION  
ALS Vial : 46 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
A43X5

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120722.M  
Quant Title : SVOA CALIBRATION

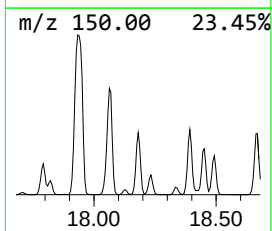
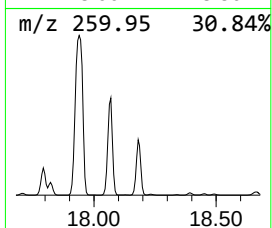
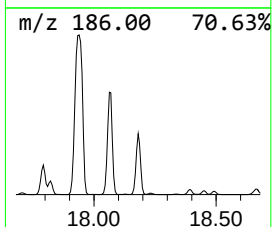
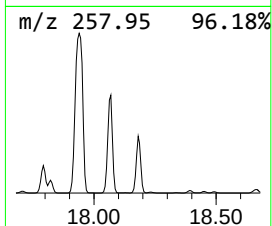
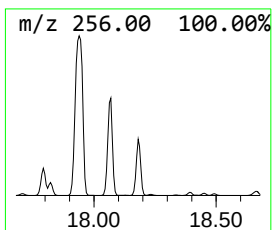
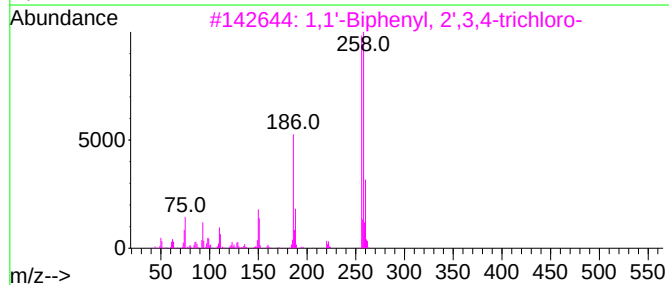
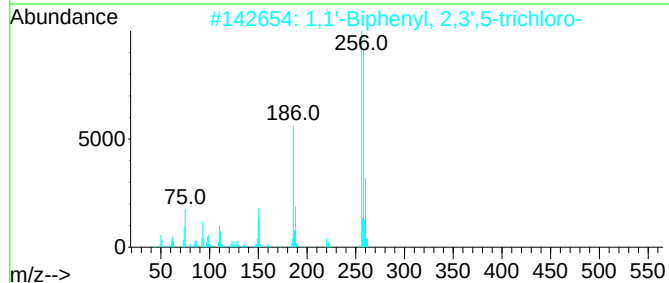
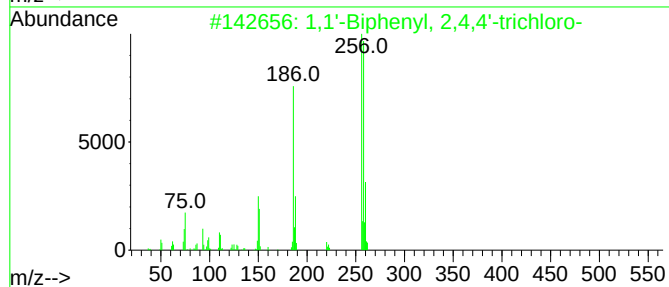
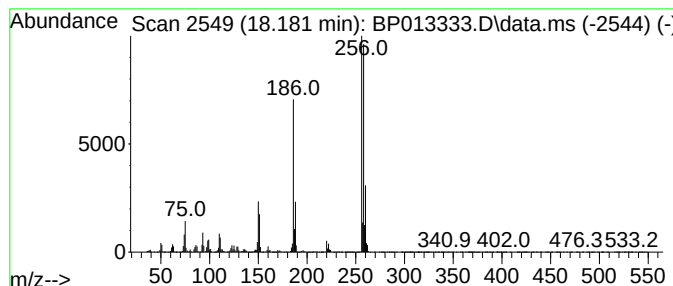
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*

Peak Number 16 1,1'-Biphenyl, 2,3,3'-trich... Concentration Rank 11

| R.T.   | EstConc     | Area     | Relative to ISTD | R.T.   |
|--------|-------------|----------|------------------|--------|
| 18.180 | 15.57 ng/ul | 12652900 | Phenanthrene-d10 | 17.345 |

| Hit# | of                               | 5   | Tentative ID | MW          | MolForm | CAS# | Qual |
|------|----------------------------------|-----|--------------|-------------|---------|------|------|
| 1    | 1,1'-Biphenyl, 2,4,4'-trichloro- | 256 | C12H7Cl3     | 007012-37-5 | 99      |      |      |
| 2    | 1,1'-Biphenyl, 2,3',5-trichloro- | 256 | C12H7Cl3     | 038444-81-4 | 99      |      |      |
| 3    | 1,1'-Biphenyl, 2',3,4-trichloro- | 256 | C12H7Cl3     | 038444-86-9 | 99      |      |      |
| 4    | 1,1'-Biphenyl, 3,4',5-trichloro- | 256 | C12H7Cl3     | 038444-88-1 | 99      |      |      |
| 5    | 1,1'-Biphenyl, 2,3,3'-trichloro- | 256 | C12H7Cl3     | 038444-84-7 | 99      |      |      |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP122422\  
Data File : BP013333.D  
Acq On : 25 Dec 2022 10:58  
Operator : CG/JU  
Sample : N6187-08  
Misc : GCMS CONFIRMATION  
ALS Vial : 46 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
A43X5

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120722.M  
Quant Title : SVOA CALIBRATION

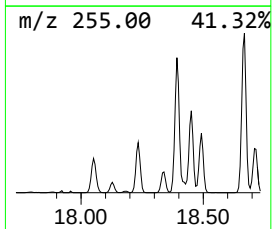
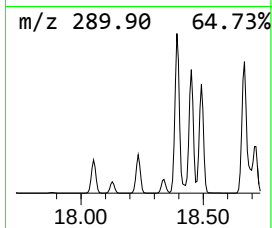
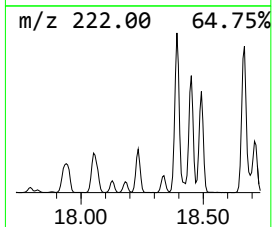
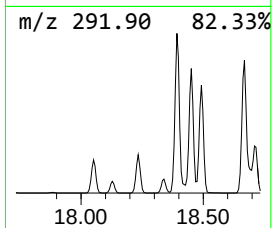
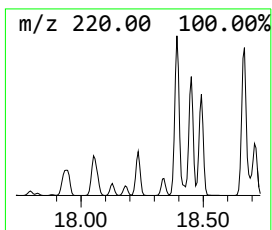
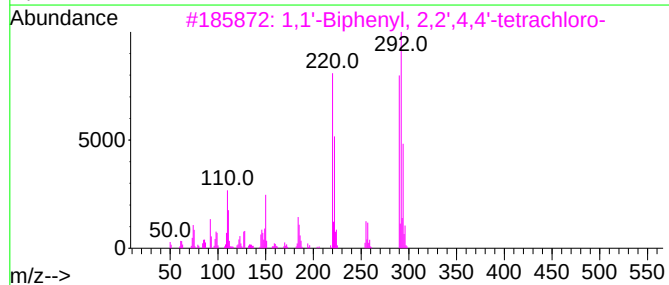
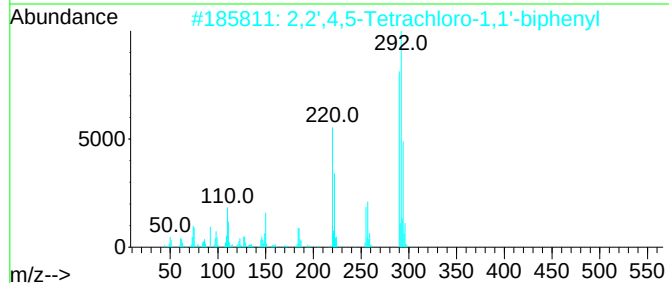
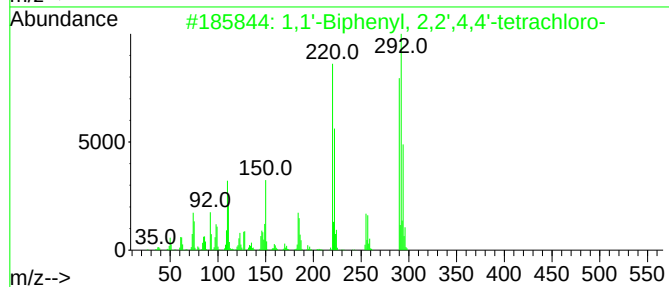
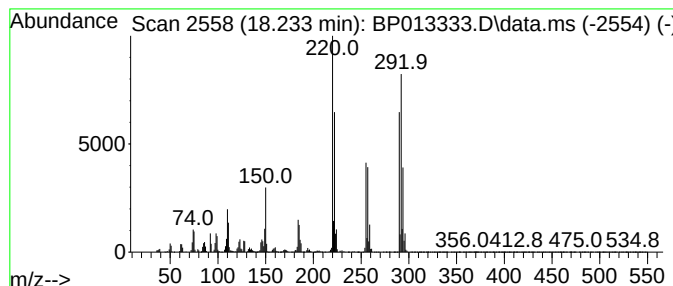
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*

Peak Number 17 1,1'-Biphenyl, 2,2',4,4'-te... Concentration Rank 25

| R.T.   | EstConc    | Area    | Relative to ISTD | R.T.   |
|--------|------------|---------|------------------|--------|
| 18.233 | 5.40 ng/ul | 4389210 | Phenanthrene-d10 | 17.345 |

| Hit# | of                                  | 5   | Tentative ID | MW          | MolForm | CAS# | Qual |
|------|-------------------------------------|-----|--------------|-------------|---------|------|------|
| 1    | 1,1'-Biphenyl, 2,2',4,4'-tetrach... | 290 | C12H6Cl4     | 002437-79-8 | 99      |      |      |
| 2    | 2,2',4,5-Tetrachloro-1,1'-biphenyl  | 290 | C12H6Cl4     | 070362-47-9 | 99      |      |      |
| 3    | 1,1'-Biphenyl, 2,2',4,4'-tetrach... | 290 | C12H6Cl4     | 002437-79-8 | 99      |      |      |
| 4    | 2,2',4,6'-Tetrachloro-1,1'-biphenyl | 290 | C12H6Cl4     | 068194-04-7 | 99      |      |      |
| 5    | 2,2',3,6-Tetrachloro-1,1'-biphenyl  | 290 | C12H6Cl4     | 070362-45-7 | 99      |      |      |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP122422\  
Data File : BP013333.D  
Acq On : 25 Dec 2022 10:58  
Operator : CG/JU  
Sample : N6187-08  
Misc : GCMS CONFIRMATION  
ALS Vial : 46 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
A43X5

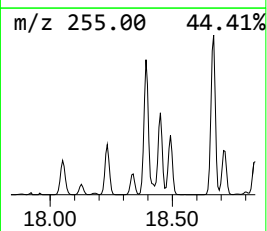
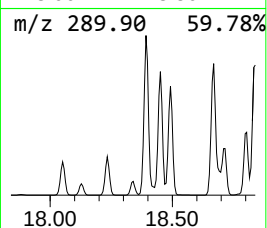
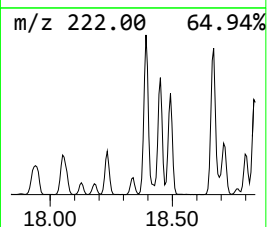
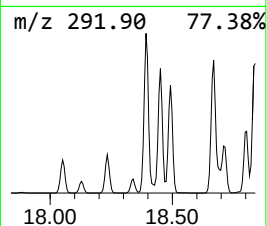
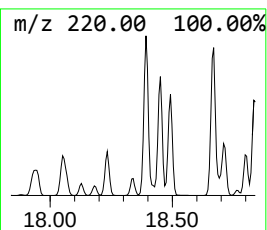
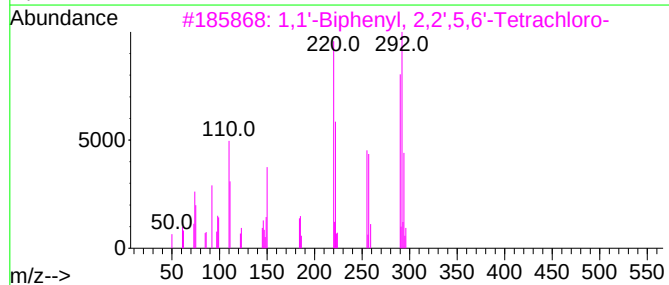
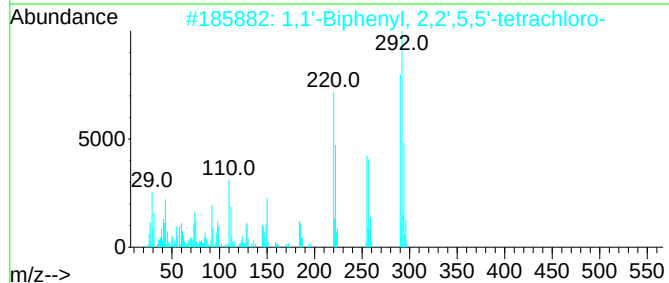
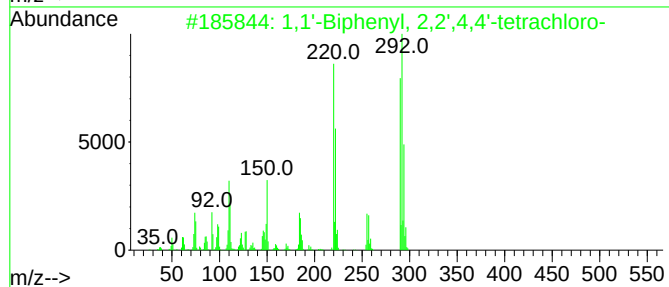
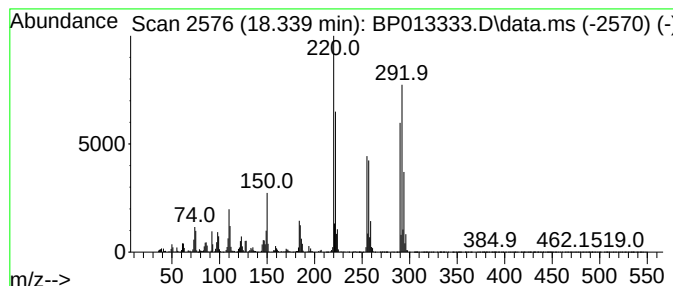
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120722.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*

Peak Number 18 1,1'-Biphenyl, 2,2',5,5'-te... Concentration Rank 32

| R.T.   | EstConc                             | Area         | Relative to ISTD | R.T.        |      |      |
|--------|-------------------------------------|--------------|------------------|-------------|------|------|
| 18.339 | 2.27 ng/ul                          | 1844790      | Phenanthrene-d10 | 17.345      |      |      |
| Hit#   | of 5                                | Tentative ID | MW               | MolForm     | CAS# | Qual |
| 1      | 1,1'-Biphenyl, 2,2',4,4'-tetrach... | 290          | C12H6Cl4         | 002437-79-8 | 99   |      |
| 2      | 1,1'-Biphenyl, 2,2',5,5'-tetrach... | 290          | C12H6Cl4         | 035693-99-3 | 99   |      |
| 3      | 1,1'-Biphenyl, 2,2',5,6'-Tetrach... | 290          | C12H6Cl4         | 041464-41-9 | 99   |      |
| 4      | 2,2',4,6'-Tetrachloro-1,1'-biphenyl | 290          | C12H6Cl4         | 068194-04-7 | 99   |      |
| 5      | 1,1'-Biphenyl, 2,2',4,6-Tetrachl... | 290          | C12H6Cl4         | 062796-65-0 | 99   |      |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP122422\  
Data File : BP013333.D  
Acq On : 25 Dec 2022 10:58  
Operator : CG/JU  
Sample : N6187-08  
Misc : GCMS CONFIRMATION  
ALS Vial : 46 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
A43X5

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120722.M  
Quant Title : SVOA CALIBRATION

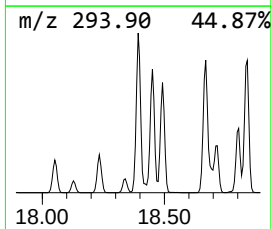
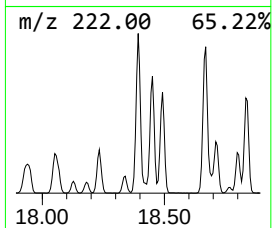
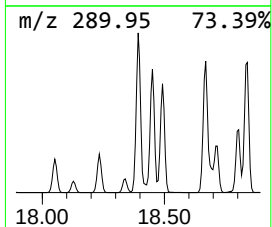
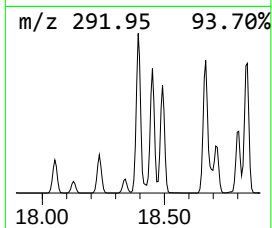
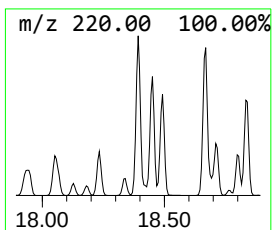
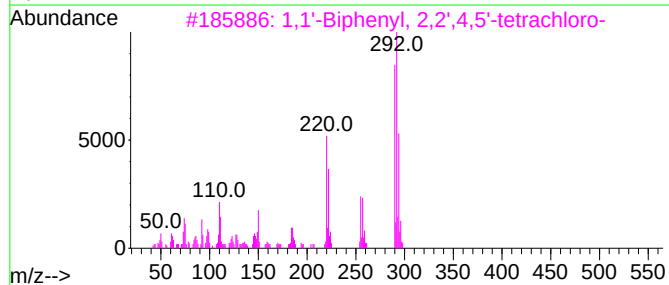
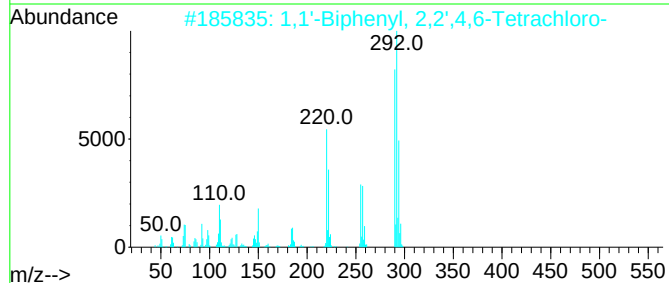
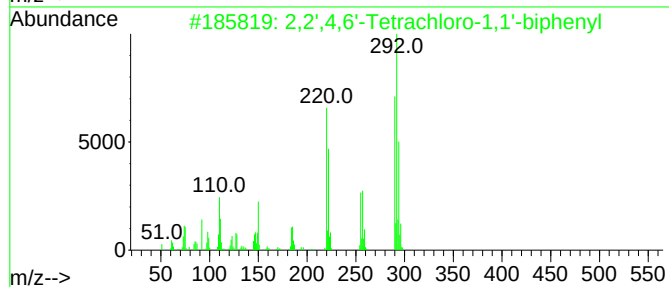
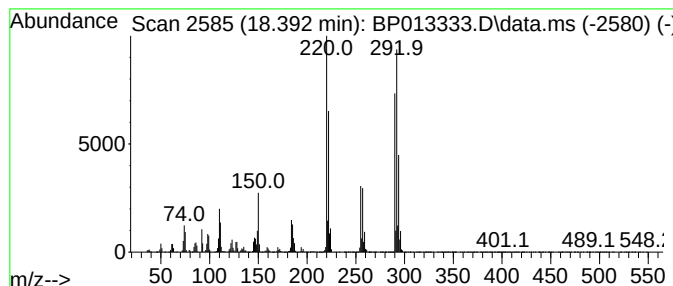
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*

Peak Number 19 2,2',4,6'-Tetrachloro-1,1'-... Concentration Rank 8

| R.T.   | EstConc     | Area     | Relative to ISTD | R.T.   |
|--------|-------------|----------|------------------|--------|
| 18.392 | 20.68 ng/ul | 16807300 | Phenanthrene-d10 | 17.345 |

| Hit# | of                                  | 5   | Tentative ID | MW          | MolForm | CAS# | Qual |
|------|-------------------------------------|-----|--------------|-------------|---------|------|------|
| 1    | 2,2',4,6'-Tetrachloro-1,1'-biphenyl | 290 | C12H6Cl4     | 068194-04-7 | 99      |      |      |
| 2    | 1,1'-Biphenyl, 2,2',4,6-Tetrachl... | 290 | C12H6Cl4     | 062796-65-0 | 99      |      |      |
| 3    | 1,1'-Biphenyl, 2,2',4,5'-tetrach... | 290 | C12H6Cl4     | 041464-40-8 | 99      |      |      |
| 4    | 1,1'-Biphenyl, 2,2',3,5'-tetrach... | 290 | C12H6Cl4     | 041464-39-5 | 99      |      |      |
| 5    | 2,2',4,5-Tetrachloro-1,1'-biphenyl  | 290 | C12H6Cl4     | 070362-47-9 | 99      |      |      |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP122422\  
Data File : BP013333.D  
Acq On : 25 Dec 2022 10:58  
Operator : CG/JU  
Sample : N6187-08  
Misc : GCMS CONFIRMATION  
ALS Vial : 46 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
A43X5

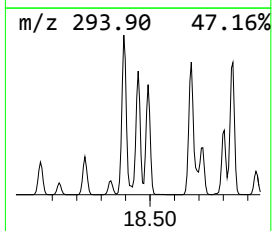
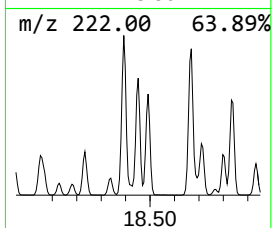
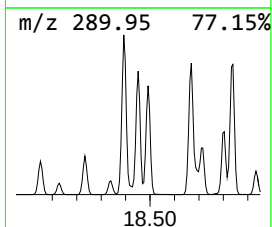
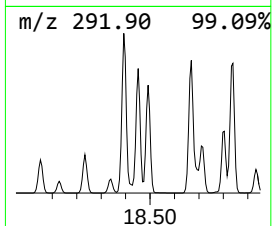
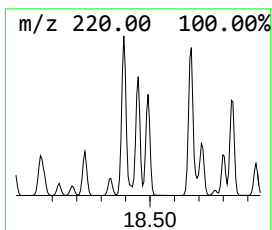
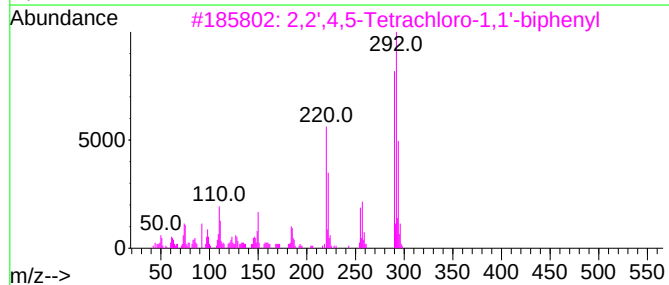
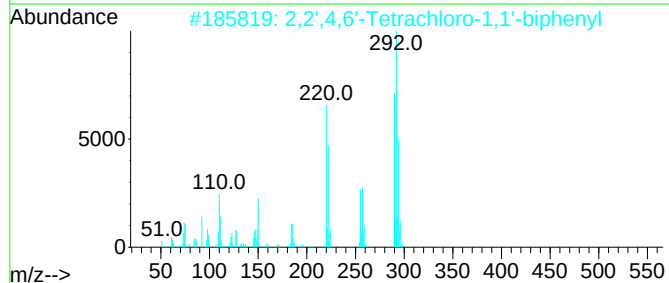
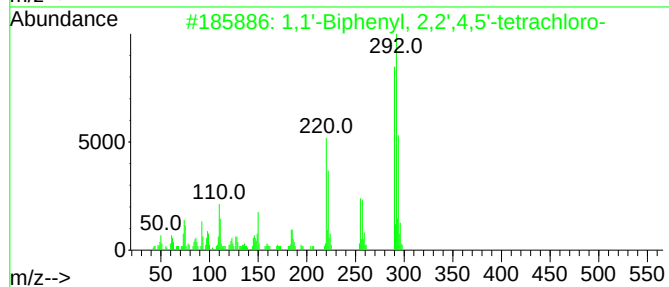
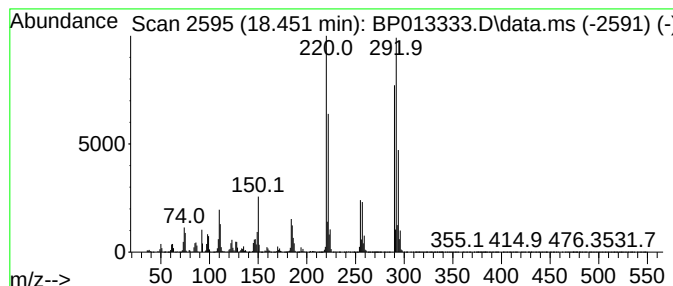
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120722.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*

Peak Number 20 1,1'-Biphenyl, 2,2',4,5'-te... Concentration Rank 12

| R.T.   | EstConc                  | Area                            | Relative to ISTD | R.T.     |             |      |
|--------|--------------------------|---------------------------------|------------------|----------|-------------|------|
| 18.451 | 14.31 ng/ul              | 11627300                        | Phenanthrene-d10 | 17.345   |             |      |
| Hit#   | of 5                     | Tentative ID                    | MW               | MolForm  | CAS#        | Qual |
| 1      | 1,1'                     | -Biphenyl, 2,2',4,5'-tetrach... | 290              | C12H6Cl4 | 041464-40-8 | 99   |
| 2      | 2,2',4,6'                | -Tetrachloro-1,1'-biphenyl      | 290              | C12H6Cl4 | 068194-04-7 | 99   |
| 3      | 2,2',4,5-                | Tetrachloro-1,1'-biphenyl       | 290              | C12H6Cl4 | 070362-47-9 | 99   |
| 4      | 1,1'-Biphenyl, 2,2',4,4' | -tetrach...                     | 290              | C12H6Cl4 | 002437-79-8 | 99   |
| 5      | 2,2',4,5-                | Tetrachloro-1,1'-biphenyl       | 290              | C12H6Cl4 | 070362-47-9 | 99   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP122422\  
Data File : BP013333.D  
Acq On : 25 Dec 2022 10:58  
Operator : CG/JU  
Sample : N6187-08  
Misc : GCMS CONFIRMATION  
ALS Vial : 46 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
A43X5

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120722.M  
Quant Title : SVOA CALIBRATION

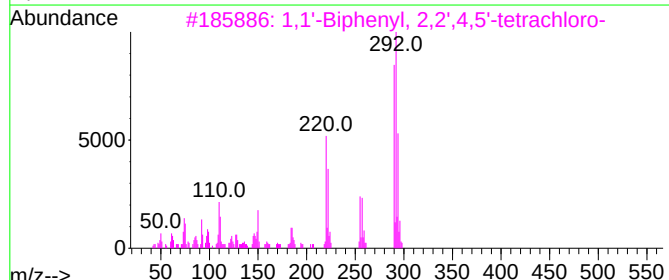
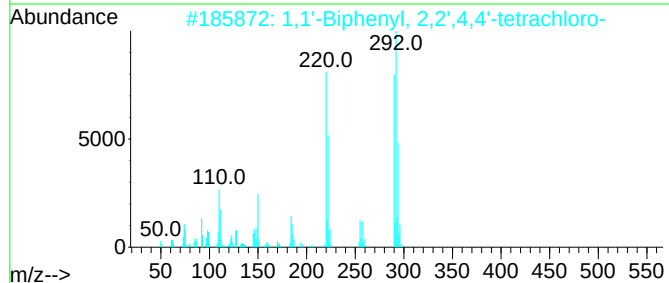
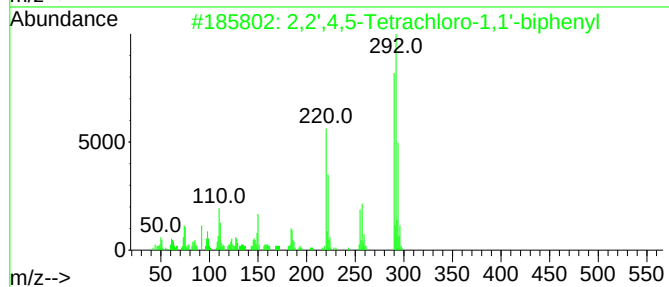
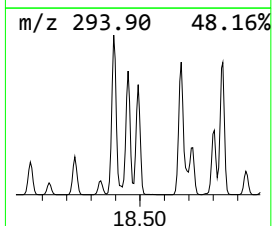
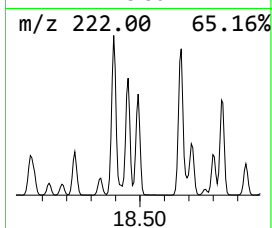
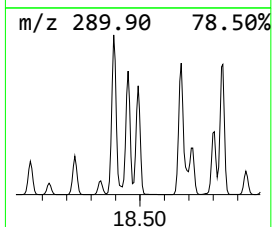
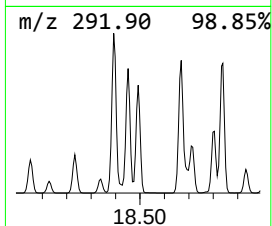
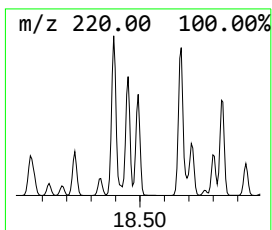
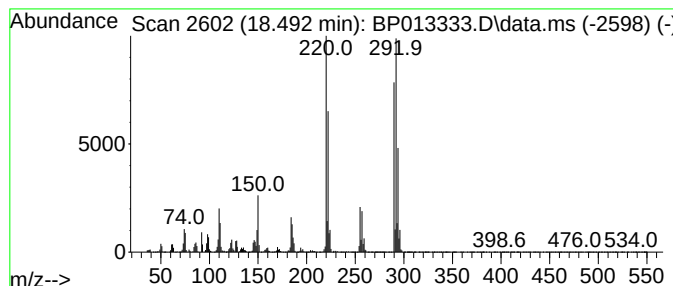
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*

Peak Number 21 2,2',4,5-Tetrachloro-1,1'-b... Concentration Rank 14

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 18.492 | 12.08 ng/ul | 9819300 | Phenanthrene-d10 | 17.345 |

| Hit# | of                                  | 5   | Tentative ID | MW          | MolForm | CAS# | Qual |
|------|-------------------------------------|-----|--------------|-------------|---------|------|------|
| 1    | 2,2',4,5-Tetrachloro-1,1'-biphenyl  | 290 | C12H6Cl4     | 070362-47-9 | 99      |      |      |
| 2    | 1,1'-Biphenyl, 2,2',4,4'-tetrach... | 290 | C12H6Cl4     | 002437-79-8 | 99      |      |      |
| 3    | 1,1'-Biphenyl, 2,2',4,5'-tetrach... | 290 | C12H6Cl4     | 041464-40-8 | 99      |      |      |
| 4    | 1,1'-Biphenyl, 2,2',4,4'-tetrach... | 290 | C12H6Cl4     | 002437-79-8 | 99      |      |      |
| 5    | 1,1'-Biphenyl, 2,2',6,6'-tetrach... | 290 | C12H6Cl4     | 015968-05-5 | 99      |      |      |





Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP122422\  
Data File : BP013333.D  
Acq On : 25 Dec 2022 10:58  
Operator : CG/JU  
Sample : N6187-08  
Misc : GCMS CONFIRMATION  
ALS Vial : 46 Sample Multiplier: 1

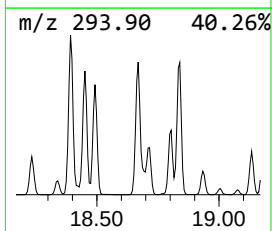
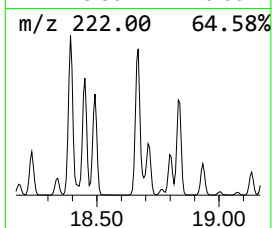
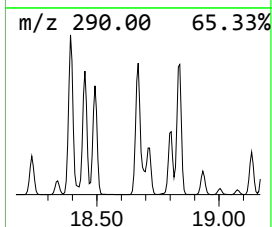
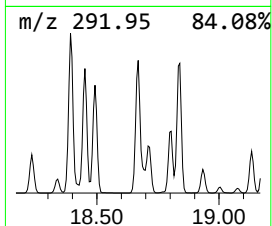
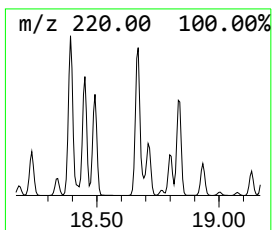
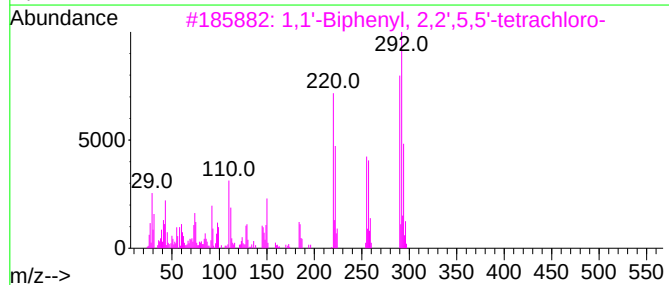
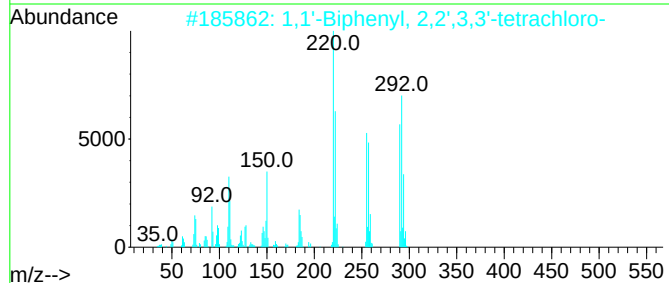
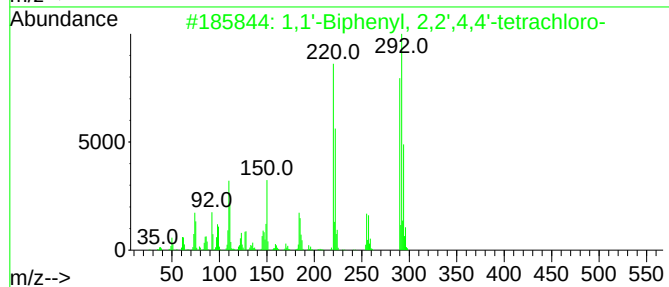
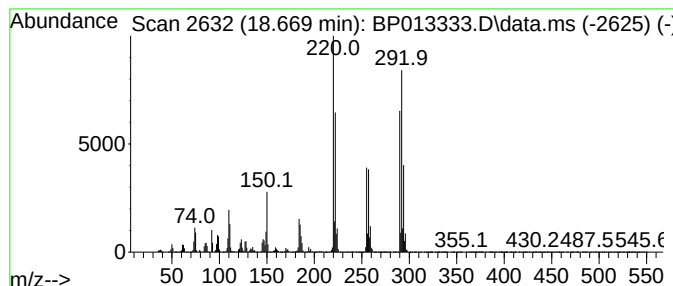
Instrument :  
BNA\_P  
ClientSampleId :  
A43X5

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120722.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 22 1,1'-Biphenyl, 2,2',3,3'-te... Concentration Rank 9

| R.T.   | EstConc                             | Area         | Relative to ISTD | R.T.        |      |      |
|--------|-------------------------------------|--------------|------------------|-------------|------|------|
| 18.669 | 20.40 ng/ul                         | 16581200     | Phenanthrene-d10 | 17.345      |      |      |
| Hit#   | of 5                                | Tentative ID | MW               | MolForm     | CAS# | Qual |
| 1      | 1,1'-Biphenyl, 2,2',4,4'-tetrach... | 290          | C12H6Cl4         | 002437-79-8 | 99   |      |
| 2      | 1,1'-Biphenyl, 2,2',3,3'-tetrach... | 290          | C12H6Cl4         | 038444-93-8 | 99   |      |
| 3      | 1,1'-Biphenyl, 2,2',5,5'-tetrach... | 290          | C12H6Cl4         | 035693-99-3 | 99   |      |
| 4      | 2,2',4,5-Tetrachloro-1,1'-biphenyl  | 290          | C12H6Cl4         | 070362-47-9 | 99   |      |
| 5      | 1,1'-Biphenyl, 2,2',3,4'-tetrach... | 290          | C12H6Cl4         | 036559-22-5 | 99   |      |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP122422\  
Data File : BP013333.D  
Acq On : 25 Dec 2022 10:58  
Operator : CG/JU  
Sample : N6187-08  
Misc : GCMS CONFIRMATION  
ALS Vial : 46 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
A43X5

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120722.M  
Quant Title : SVOA CALIBRATION

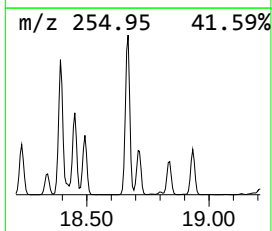
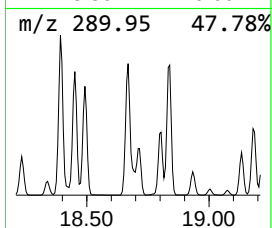
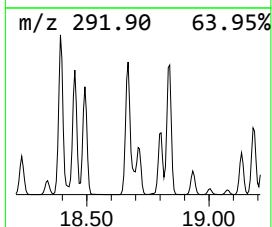
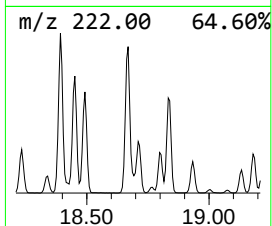
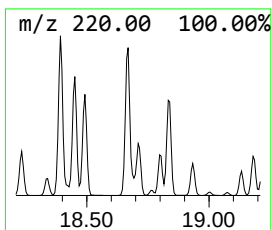
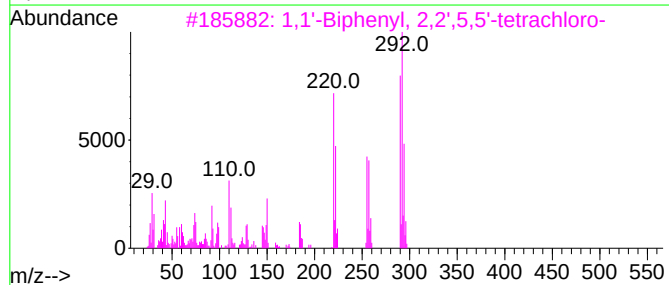
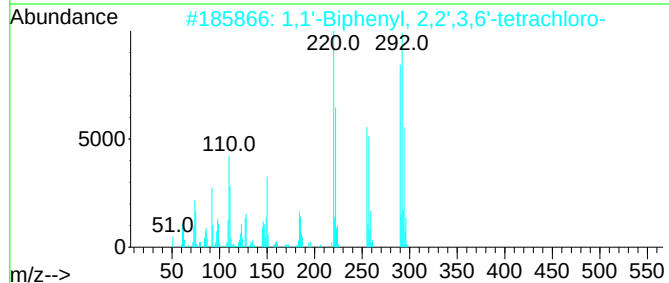
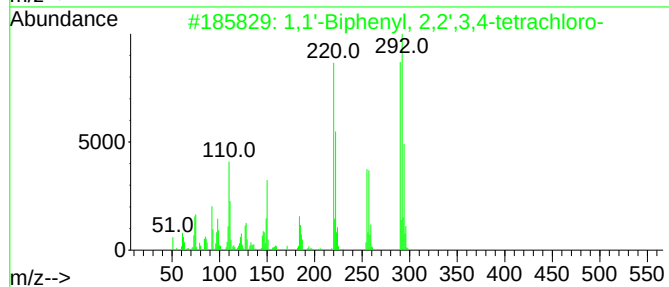
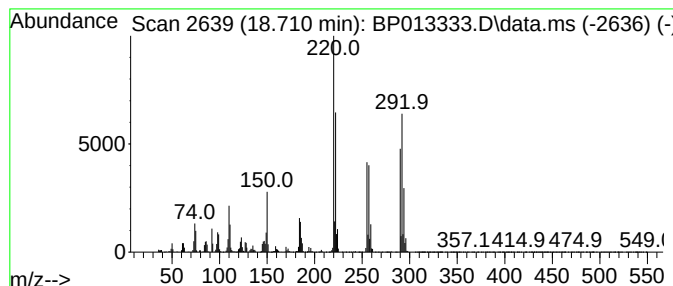
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*

Peak Number 23 1,1'-Biphenyl, 2,2',3,4-tet... Concentration Rank 20

| R.T.   | EstConc    | Area    | Relative to ISTD | R.T.   |
|--------|------------|---------|------------------|--------|
| 18.710 | 6.81 ng/ul | 5533060 | Phenanthrene-d10 | 17.345 |

| Hit# | of                                  | 5   | Tentative ID | MW          | MolForm | CAS# | Qual |
|------|-------------------------------------|-----|--------------|-------------|---------|------|------|
| 1    | 1,1'-Biphenyl, 2,2',3,4-tetrachl... | 290 | C12H6Cl4     | 052663-59-9 | 99      |      |      |
| 2    | 1,1'-Biphenyl, 2,2',3,6'-tetrach... | 290 | C12H6Cl4     | 041464-47-5 | 99      |      |      |
| 3    | 1,1'-Biphenyl, 2,2',5,5'-tetrach... | 290 | C12H6Cl4     | 035693-99-3 | 99      |      |      |
| 4    | 1,1'-Biphenyl, 2,2',3,3'-tetrach... | 290 | C12H6Cl4     | 038444-93-8 | 99      |      |      |
| 5    | 1,1'-Biphenyl, 2,2',4,6-Tetrachl... | 290 | C12H6Cl4     | 062796-65-0 | 99      |      |      |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP122422\  
Data File : BP013333.D  
Acq On : 25 Dec 2022 10:58  
Operator : CG/JU  
Sample : N6187-08  
Misc : GCMS CONFIRMATION  
ALS Vial : 46 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
A43X5

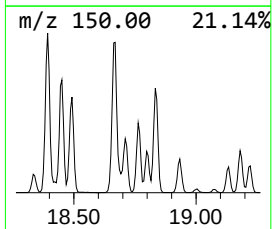
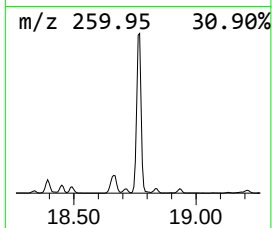
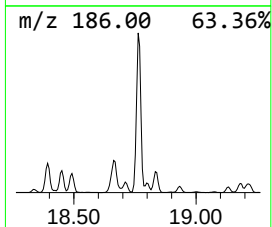
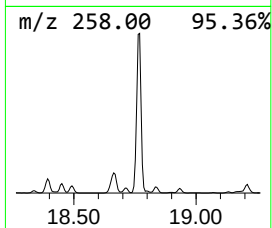
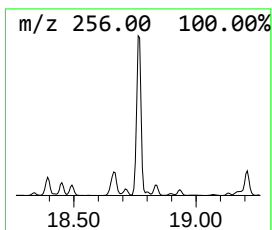
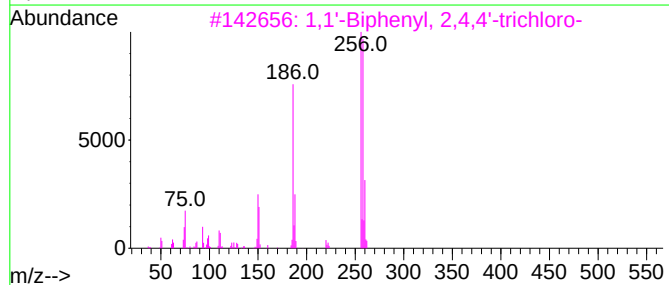
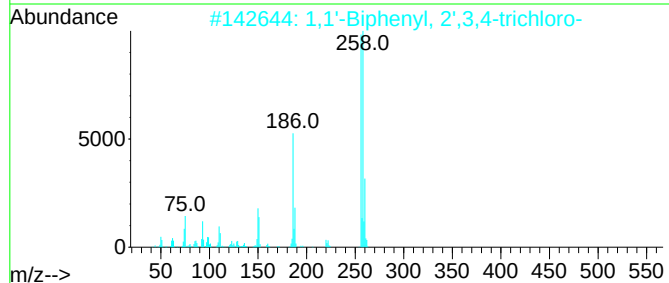
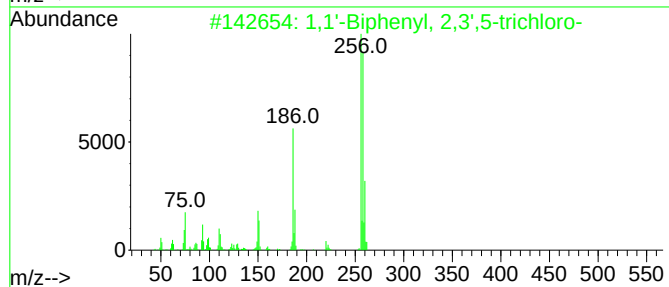
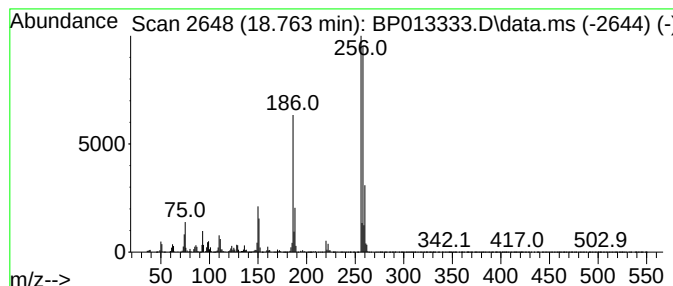
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120722.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 24 unknown-01 Concentration Rank 17

| R.T.   | EstConc    | Area    | Relative to ISTD | R.T.   |
|--------|------------|---------|------------------|--------|
| 18.763 | 7.60 ng/ul | 6172050 | Phenanthrene-d10 | 17.345 |

| Hit# | of                               | 5   | Tentative ID | MW          | MolForm | CAS# | Qual |
|------|----------------------------------|-----|--------------|-------------|---------|------|------|
| 1    | 1,1'-Biphenyl, 2,3',5-trichloro- | 256 | C12H7Cl3     | 038444-81-4 | 99      |      |      |
| 2    | 1,1'-Biphenyl, 2',3,4-trichloro- | 256 | C12H7Cl3     | 038444-86-9 | 99      |      |      |
| 3    | 1,1'-Biphenyl, 2,4,4'-trichloro- | 256 | C12H7Cl3     | 007012-37-5 | 99      |      |      |
| 4    | 1,1'-Biphenyl, 2,4,4'-trichloro- | 256 | C12H7Cl3     | 007012-37-5 | 99      |      |      |
| 5    | 1,1'-Biphenyl, 2,4',5-trichloro- | 256 | C12H7Cl3     | 016606-02-3 | 99      |      |      |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP122422\  
Data File : BP013333.D  
Acq On : 25 Dec 2022 10:58  
Operator : CG/JU  
Sample : N6187-08  
Misc : GCMS CONFIRMATION  
ALS Vial : 46 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
A43X5

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120722.M  
Quant Title : SVOA CALIBRATION

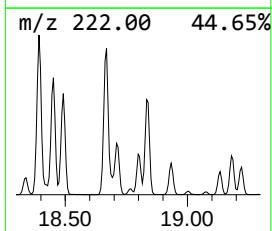
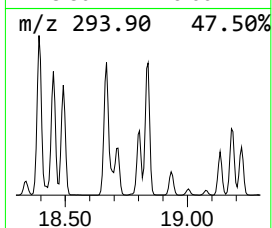
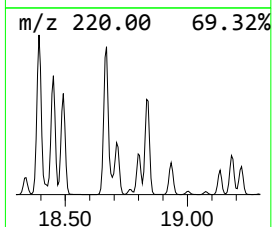
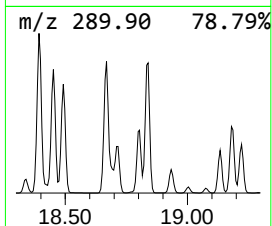
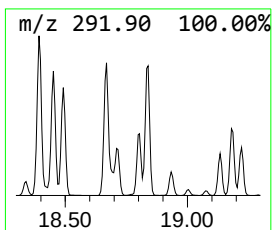
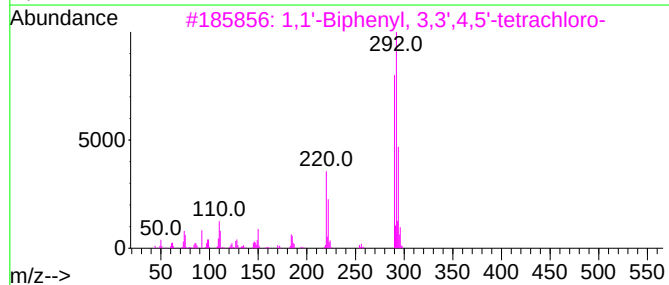
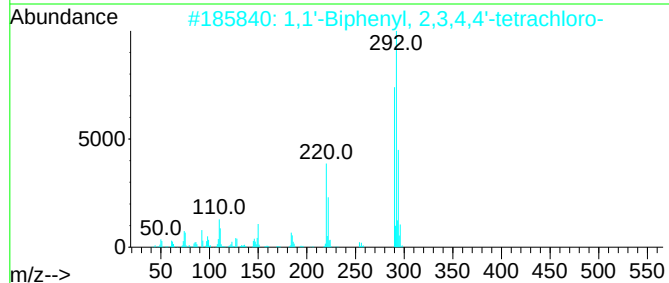
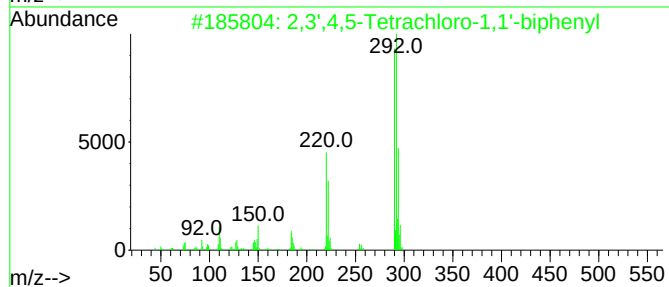
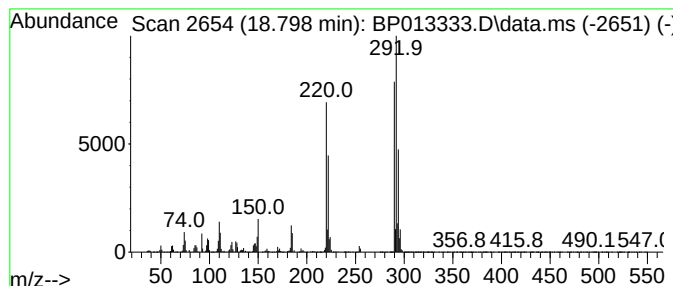
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*

Peak Number 25 2,3',4,5-Tetrachloro-1,1'-b... Concentration Rank 23

| R.T.   | EstConc    | Area    | Relative to ISTD | R.T.   |
|--------|------------|---------|------------------|--------|
| 18.798 | 5.87 ng/ul | 4772180 | Phenanthrene-d10 | 17.345 |

| Hit# | of                                  | 5   | Tentative ID | MW          | MolForm | CAS# | Qual |
|------|-------------------------------------|-----|--------------|-------------|---------|------|------|
| 1    | 2,3',4,5-Tetrachloro-1,1'-biphenyl  | 290 | C12H6Cl4     | 073575-53-8 | 99      |      |      |
| 2    | 1,1'-Biphenyl, 2,3,4,4'-tetrachl... | 290 | C12H6Cl4     | 033025-41-1 | 99      |      |      |
| 3    | 1,1'-Biphenyl, 3,3',4,5'-tetrach... | 290 | C12H6Cl4     | 041464-48-6 | 99      |      |      |
| 4    | 1,1'-Biphenyl, 2,2',3,4-tetrachl... | 290 | C12H6Cl4     | 052663-59-9 | 99      |      |      |
| 5    | 2,3,3',4-Tetrachloro-1,1'-biphenyl  | 290 | C12H6Cl4     | 074338-24-2 | 99      |      |      |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP122422\  
Data File : BP013333.D  
Acq On : 25 Dec 2022 10:58  
Operator : CG/JU  
Sample : N6187-08  
Misc : GCMS CONFIRMATION  
ALS Vial : 46 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
A43X5

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120722.M  
Quant Title : SVOA CALIBRATION

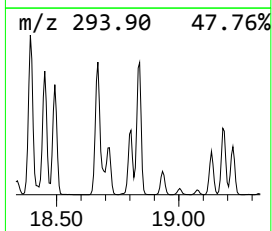
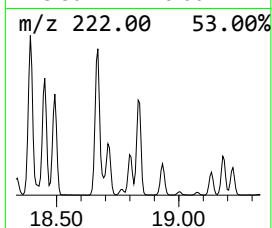
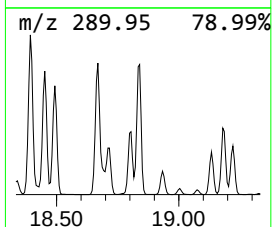
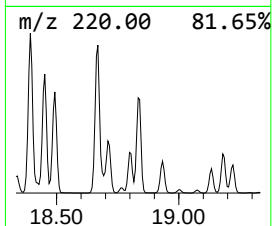
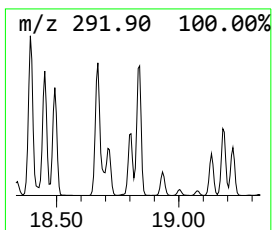
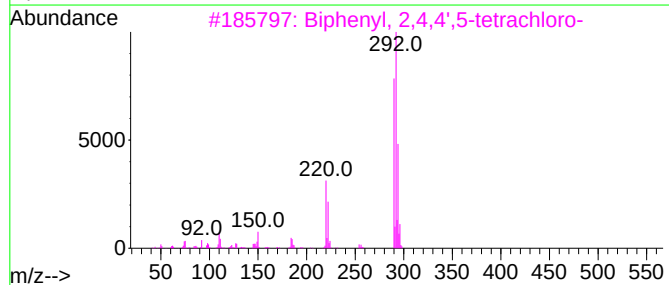
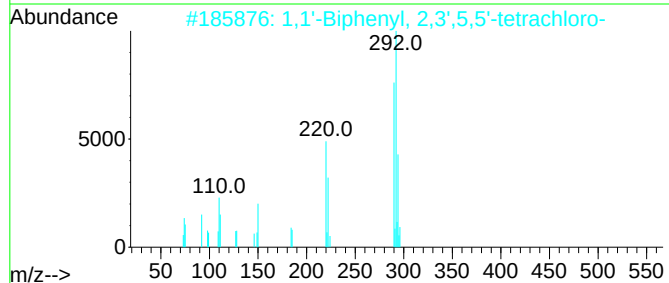
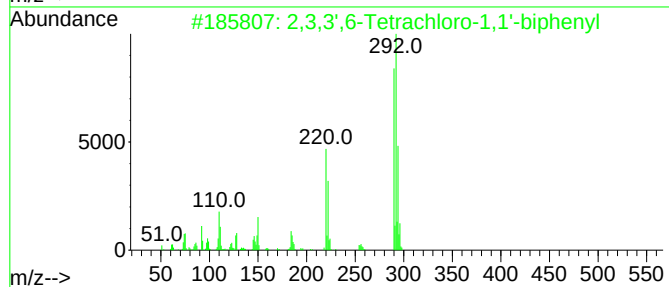
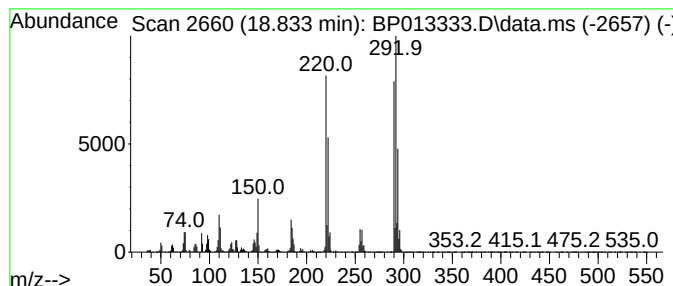
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*

Peak Number 26 2,3,3',6-Tetrachloro-1,1'-b... Concentration Rank 13

| R.T.   | EstConc     | Area     | Relative to ISTD | R.T.   |
|--------|-------------|----------|------------------|--------|
| 18.833 | 13.18 ng/ul | 10711400 | Phenanthrene-d10 | 17.345 |

| Hit# | of                                  | 5   | Tentative ID | MW          | MolForm | CAS# | Qual |
|------|-------------------------------------|-----|--------------|-------------|---------|------|------|
| 1    | 2,3,3',6-Tetrachloro-1,1'-biphenyl  | 290 | C12H6Cl4     | 074472-33-6 | 99      |      |      |
| 2    | 1,1'-Biphenyl, 2,3',5,5'-tetrach... | 290 | C12H6Cl4     | 041464-42-0 | 99      |      |      |
| 3    | Biphenyl, 2,4,4',5-tetrachloro-     | 290 | C12H6Cl4     | 032690-93-0 | 99      |      |      |
| 4    | 2,3',5',6-Tetrachloro-1,1'-biphenyl | 290 | C12H6Cl4     | 074338-23-1 | 99      |      |      |
| 5    | 1,1'-Biphenyl, 2,3,3',5'-tetrach... | 290 | C12H6Cl4     | 041464-49-7 | 99      |      |      |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP122422\  
Data File : BP013333.D  
Acq On : 25 Dec 2022 10:58  
Operator : CG/JU  
Sample : N6187-08  
Misc : GCMS CONFIRMATION  
ALS Vial : 46 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
A43X5

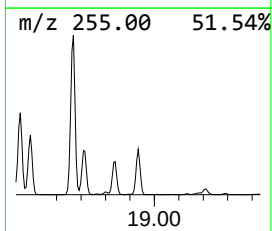
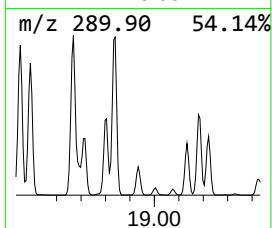
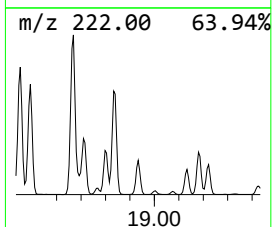
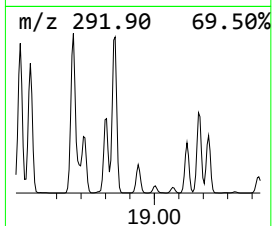
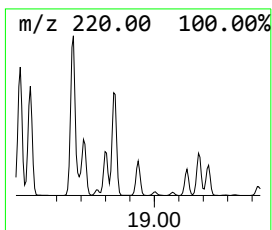
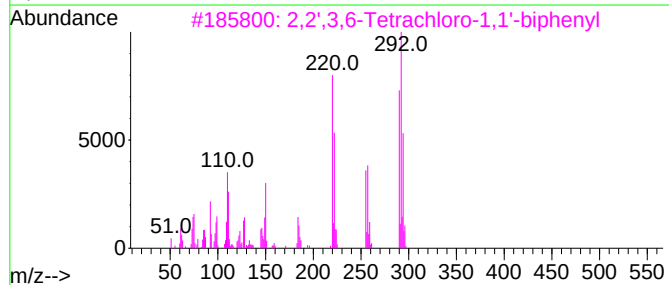
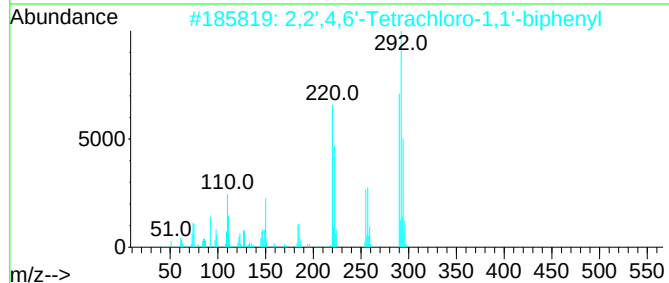
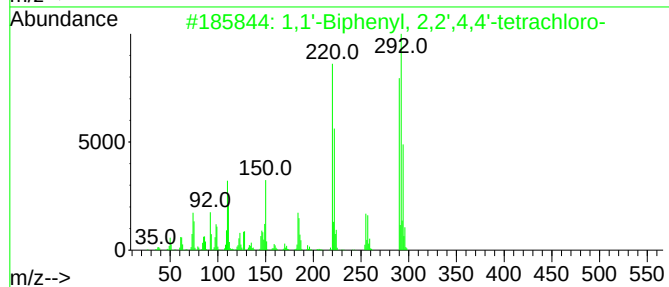
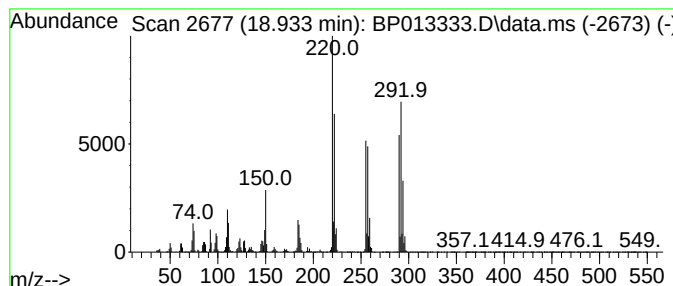
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120722.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*

Peak Number 27 2,2',3,6-Tetrachloro-1,1'-b... Concentration Rank 27

| R.T.      | EstConc                             | Area    | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------------|---------|------------------|-------------|------|
| 18.933    | 3.82 ng/ul                          | 3102880 | Phenanthrene-d10 | 17.345      |      |
| Hit# of 5 | Tentative ID                        | MW      | MolForm          | CAS#        | Qual |
| 1         | 1,1'-Biphenyl, 2,2',4,4'-tetrach... | 290     | C12H6Cl4         | 002437-79-8 | 99   |
| 2         | 2,2',4,6'-Tetrachloro-1,1'-biphenyl | 290     | C12H6Cl4         | 068194-04-7 | 99   |
| 3         | 2,2',3,6-Tetrachloro-1,1'-biphenyl  | 290     | C12H6Cl4         | 070362-45-7 | 99   |
| 4         | 2,2',4,5-Tetrachloro-1,1'-biphenyl  | 290     | C12H6Cl4         | 070362-47-9 | 99   |
| 5         | 2,2',4,5-Tetrachloro-1,1'-biphenyl  | 290     | C12H6Cl4         | 070362-47-9 | 99   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP122422\  
Data File : BP013333.D  
Acq On : 25 Dec 2022 10:58  
Operator : CG/JU  
Sample : N6187-08  
Misc : GCMS CONFIRMATION  
ALS Vial : 46 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
A43X5

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120722.M  
Quant Title : SVOA CALIBRATION

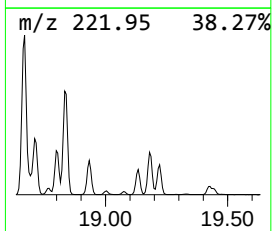
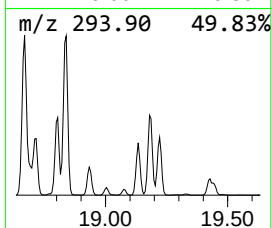
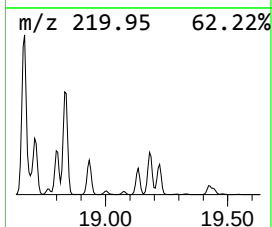
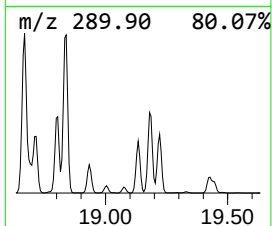
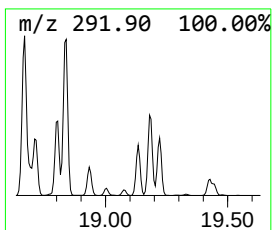
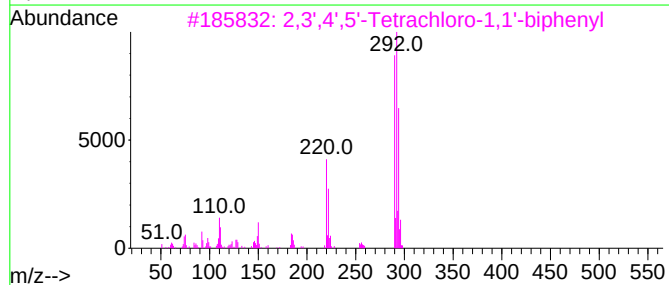
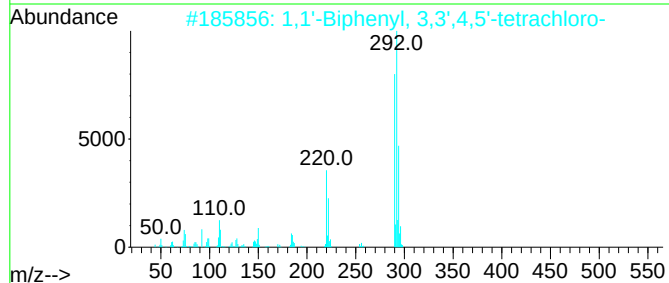
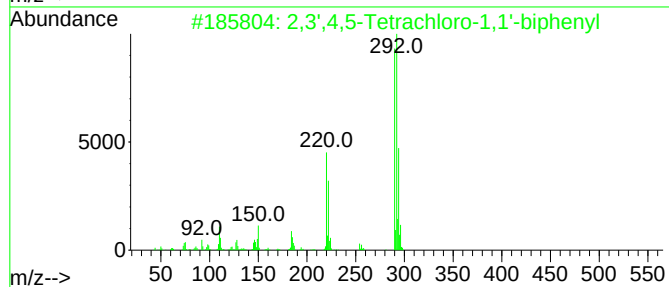
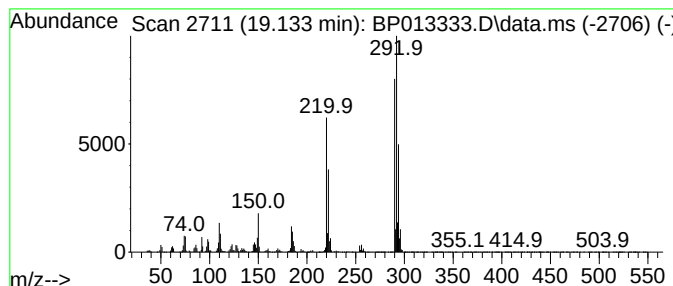
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*

Peak Number 28 1,1'-Biphenyl, 3,3',4,5'-te... Concentration Rank 28

| R.T.   | EstConc    | Area    | Relative to ISTD | R.T.   |
|--------|------------|---------|------------------|--------|
| 19.133 | 3.51 ng/ul | 2852130 | Phenanthrene-d10 | 17.345 |

| Hit# | of                                  | 5   | Tentative ID | MW          | MolForm | CAS# | Qual |
|------|-------------------------------------|-----|--------------|-------------|---------|------|------|
| 1    | 2,3',4,5-Tetrachloro-1,1'-biphenyl  | 290 | C12H6Cl4     | 073575-53-8 | 99      |      |      |
| 2    | 1,1'-Biphenyl, 3,3',4,5'-tetrach... | 290 | C12H6Cl4     | 041464-48-6 | 99      |      |      |
| 3    | 2,3',4',5'-Tetrachloro-1,1'-biph... | 290 | C12H6Cl4     | 070362-48-0 | 99      |      |      |
| 4    | 2,3,3',4-Tetrachloro-1,1'-biphenyl  | 290 | C12H6Cl4     | 074338-24-2 | 99      |      |      |
| 5    | 1,1'-Biphenyl, 2,3,3',4'-tetrach... | 290 | C12H6Cl4     | 041464-43-1 | 99      |      |      |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP122422\  
Data File : BP013333.D  
Acq On : 25 Dec 2022 10:58  
Operator : CG/JU  
Sample : N6187-08  
Misc : GCMS CONFIRMATION  
ALS Vial : 46 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
A43X5

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120722.M  
Quant Title : SVOA CALIBRATION

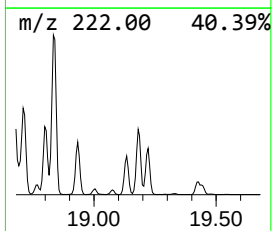
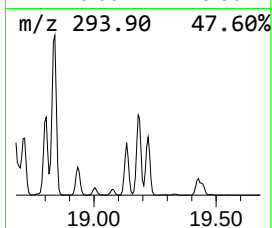
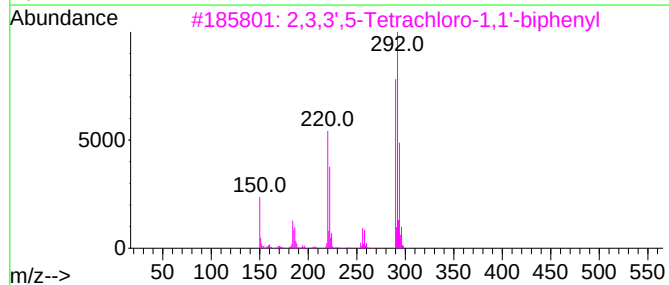
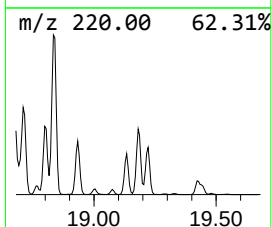
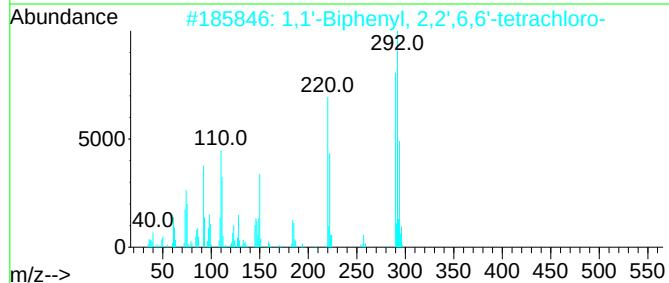
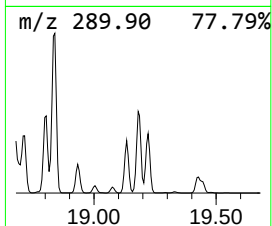
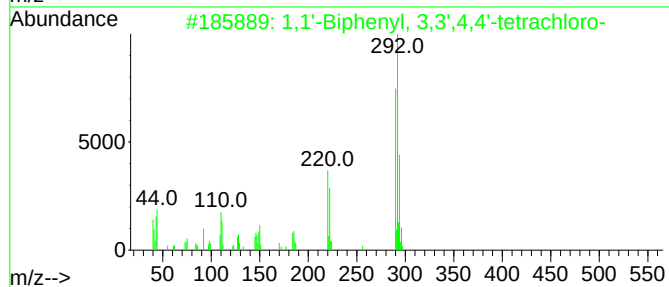
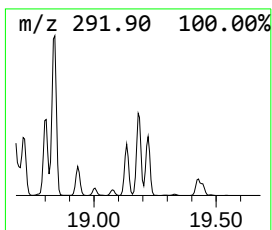
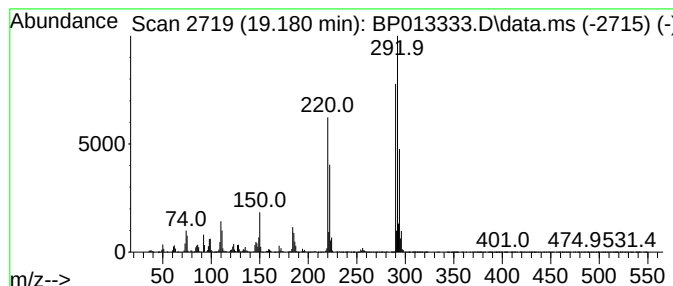
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*

Peak Number 29 1,1'-Biphenyl, 3,3',4,4'-te... Concentration Rank 21

| R.T.   | EstConc    | Area    | Relative to ISTD | R.T.   |
|--------|------------|---------|------------------|--------|
| 19.180 | 6.53 ng/ul | 5309630 | Phenanthrene-d10 | 17.345 |

| Hit# | of 5                                | Tentative ID | MW       | MolForm     | CAS# | Qual |
|------|-------------------------------------|--------------|----------|-------------|------|------|
| 1    | 1,1'-Biphenyl, 3,3',4,4'-tetrach... | 290          | C12H6Cl4 | 032598-13-3 | 99   |      |
| 2    | 1,1'-Biphenyl, 2,2',6,6'-tetrach... | 290          | C12H6Cl4 | 015968-05-5 | 99   |      |
| 3    | 2,3,3',5-Tetrachloro-1,1'-biphenyl  | 290          | C12H6Cl4 | 070424-67-8 | 96   |      |
| 4    | 2,3,4,6-Tetrachloro-1,1'-biphenyl   | 290          | C12H6Cl4 | 054230-22-7 | 96   |      |
| 5    | 2,3',4',5'-Tetrachloro-1,1'-biph... | 290          | C12H6Cl4 | 070362-48-0 | 96   |      |





Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP122422\  
Data File : BP013333.D  
Acq On : 25 Dec 2022 10:58  
Operator : CG/JU  
Sample : N6187-08  
Misc : GCMS CONFIRMATION  
ALS Vial : 46 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
A43X5

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120722.M  
Quant Title : SVOA CALIBRATION

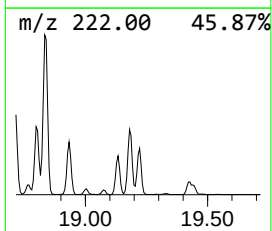
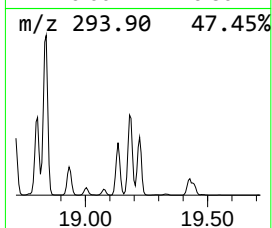
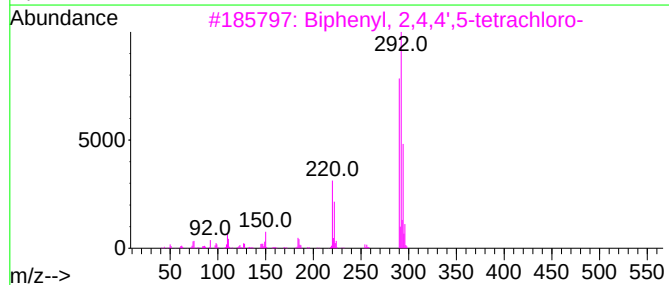
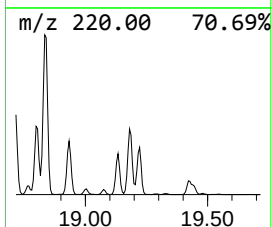
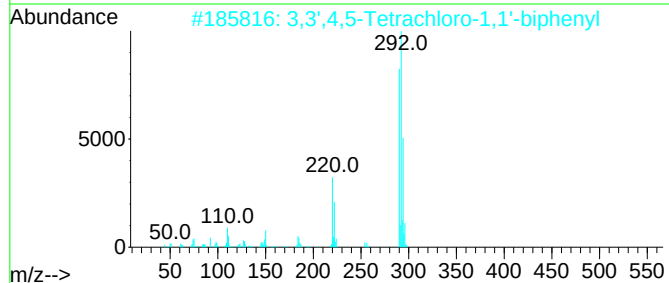
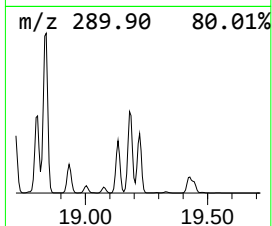
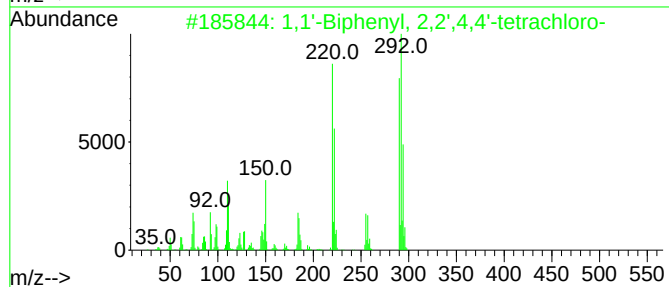
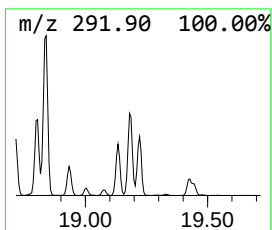
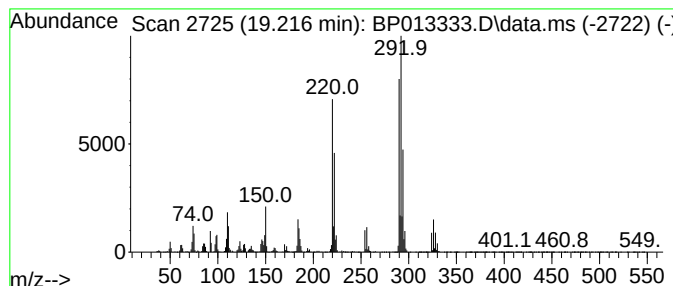
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*

Peak Number 30 3,3',4,5-Tetrachloro-1,1'-b... Concentration Rank 22

| R.T.   | EstConc    | Area    | Relative to ISTD | R.T.   |
|--------|------------|---------|------------------|--------|
| 19.216 | 5.92 ng/ul | 4810610 | Phenanthrene-d10 | 17.345 |

| Hit# | of                                  | 5   | Tentative ID | MW          | MolForm | CAS# | Qual |
|------|-------------------------------------|-----|--------------|-------------|---------|------|------|
| 1    | 1,1'-Biphenyl, 2,2',4,4'-tetrach... | 290 | C12H6Cl4     | 002437-79-8 | 99      |      |      |
| 2    | 3,3',4,5-Tetrachloro-1,1'-biphenyl  | 290 | C12H6Cl4     | 070362-49-1 | 99      |      |      |
| 3    | Biphenyl, 2,4,4',5-tetrachloro-     | 290 | C12H6Cl4     | 032690-93-0 | 99      |      |      |
| 4    | 1,1'-Biphenyl, 2,3',4',6-tetrach... | 290 | C12H6Cl4     | 041464-46-4 | 99      |      |      |
| 5    | 2,3',4,5-Tetrachloro-1,1'-biphenyl  | 290 | C12H6Cl4     | 073575-53-8 | 99      |      |      |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP122422\  
Data File : BP013333.D  
Acq On : 25 Dec 2022 10:58  
Operator : CG/JU  
Sample : N6187-08  
Misc : GCMS CONFIRMATION  
ALS Vial : 46 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
A43X5

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120722.M  
Quant Title : SVOA CALIBRATION

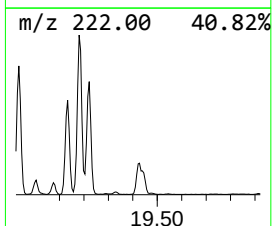
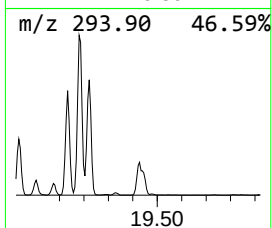
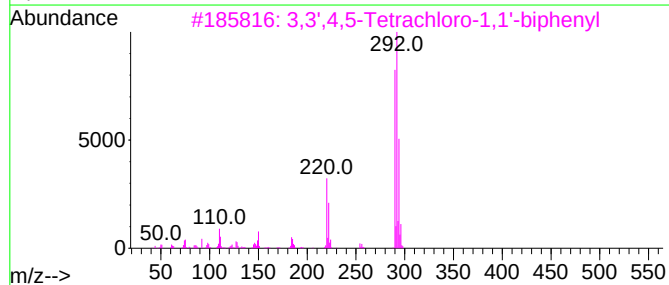
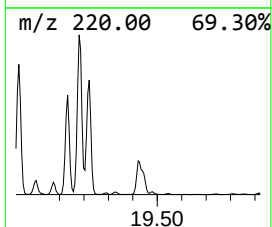
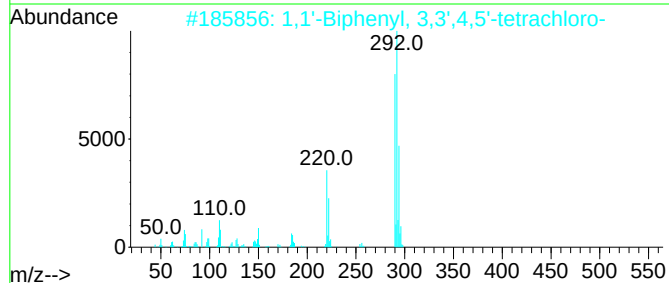
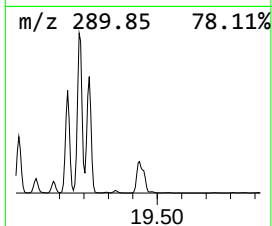
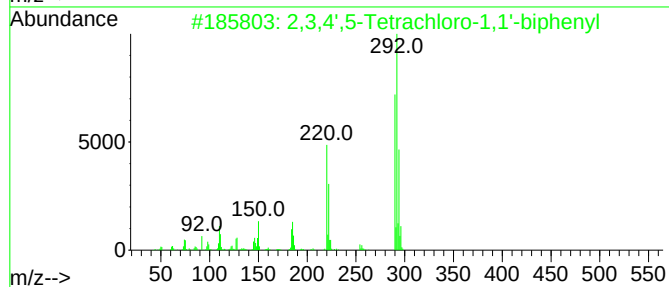
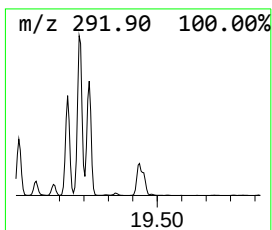
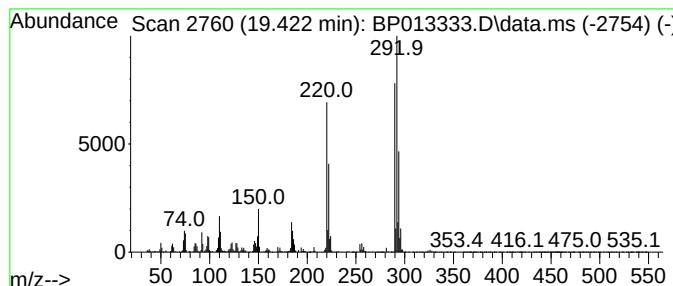
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*

Peak Number 31 Biphenyl, 2,4,4',5-tetrachl... Concentration Rank 19

| R.T.   | EstConc    | Area    | Relative to ISTD | R.T.   |
|--------|------------|---------|------------------|--------|
| 19.422 | 6.96 ng/ul | 1717310 | Chrysene-d12     | 21.433 |

| Hit# | of                                  | 5   | Tentative ID | MW          | MolForm | CAS# | Qual |
|------|-------------------------------------|-----|--------------|-------------|---------|------|------|
| 1    | 2,3,4',5-Tetrachloro-1,1'-biphenyl  | 290 | C12H6Cl4     | 074472-34-7 | 99      |      |      |
| 2    | 1,1'-Biphenyl, 3,3',4,5'-tetrach... | 290 | C12H6Cl4     | 041464-48-6 | 99      |      |      |
| 3    | 3,3',4,5-Tetrachloro-1,1'-biphenyl  | 290 | C12H6Cl4     | 070362-49-1 | 99      |      |      |
| 4    | Biphenyl, 2,4,4',5-tetrachloro-     | 290 | C12H6Cl4     | 032690-93-0 | 99      |      |      |
| 5    | 1,1'-Biphenyl, 2,2',4,4'-tetrach... | 290 | C12H6Cl4     | 002437-79-8 | 99      |      |      |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP122422\  
Data File : BP013333.D  
Acq On : 25 Dec 2022 10:58  
Operator : CG/JU  
Sample : N6187-08  
Misc : GCMS CONFIRMATION  
ALS Vial : 46 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
A43X5

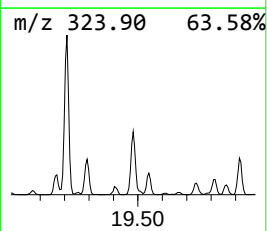
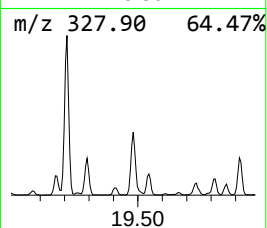
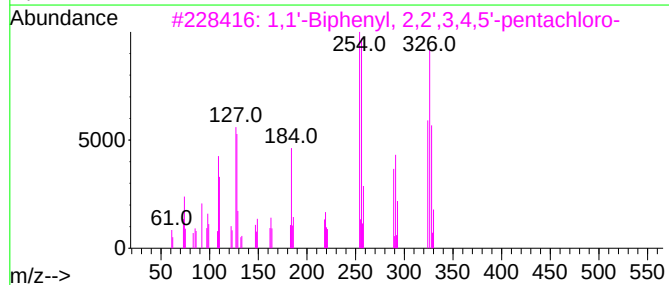
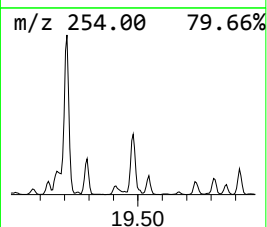
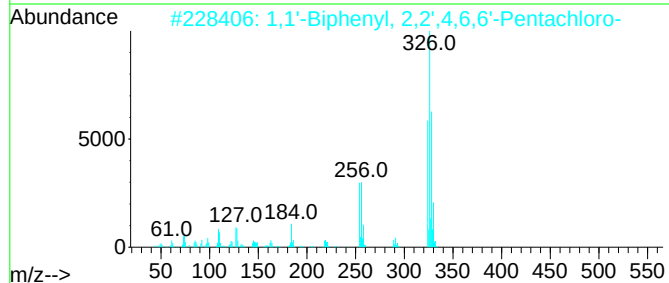
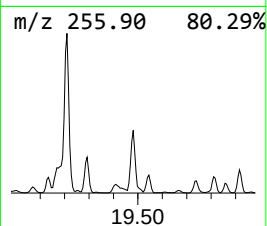
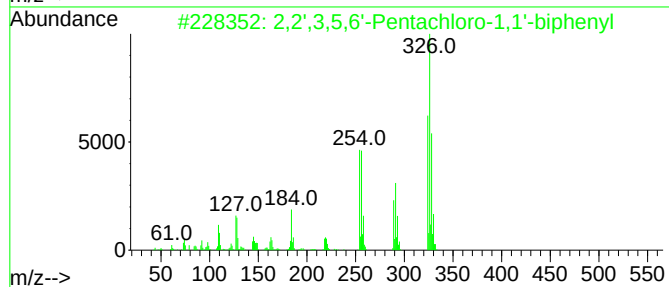
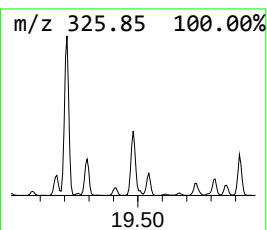
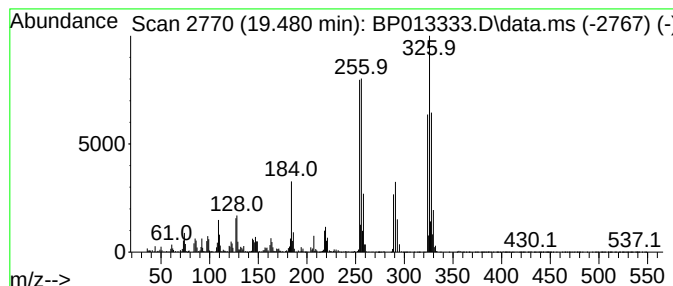
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120722.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*

Peak Number 32 2,2',3,5,6'-Pentachloro-1,1... Concentration Rank 29

| R.T.      | EstConc                              | Area   | Relative to ISTD | R.T.        |      |
|-----------|--------------------------------------|--------|------------------|-------------|------|
| 19.480    | 3.32 ng/ul                           | 819293 | Chrysene-d12     | 21.433      |      |
| Hit# of 5 | Tentative ID                         | MW     | MolForm          | CAS#        | Qual |
| 1         | 2,2',3,5,6'-Pentachloro-1,1'-bip...  | 324    | C12H5Cl5         | 073575-55-0 | 99   |
| 2         | 1,1'-Biphenyl, 2,2',4,6,6'-Penta...  | 324    | C12H5Cl5         | 056558-16-8 | 99   |
| 3         | 1,1'-Biphenyl, 2,2',3,4,5'-penta...  | 324    | C12H5Cl5         | 038380-02-8 | 99   |
| 4         | 1,1'-Biphenyl, 2,2',3,5',6'-penta... | 324    | C12H5Cl5         | 038379-99-6 | 99   |
| 5         | 1,1'-Biphenyl, 2,2',3',4,5'-Penta... | 324    | C12H5Cl5         | 041464-51-1 | 99   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP122422\  
 Data File : BP013333.D  
 Acq On : 25 Dec 2022 10:58  
 Operator : CG/JU  
 Sample : N6187-08  
 Misc : GCMS CONFIRMATION  
 ALS Vial : 46 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 A43X5

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\SFAM-EPA-BP120722.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 |
| (DEL) Alkane: S... | 3.252  | 76.0    | ng/ul | 12677700 | 1                     | 7.934  | 3333930  | 20.0 |
| (DEL) Alkane: C... | 3.599  | 3.0     | ng/ul | 491286   | 1                     | 7.934  | 3333930  | 20.0 |
| 1,1'-Biphenyl, ... | 14.704 | 7.4     | ng/ul | 1951950  | 3                     | 14.581 | 5282020  | 20.0 |
| 1,1'-Biphenyl, ... | 15.839 | 57.3    | ng/ul | 15122200 | 3                     | 14.581 | 5282020  | 20.0 |
| 1,1'-Biphenyl, ... | 16.281 | 4.8     | ng/ul | 3885540  | 4                     | 17.345 | 16252400 | 20.0 |
| 1,1'-Biphenyl, ... | 16.445 | 8.6     | ng/ul | 6951570  | 4                     | 17.345 | 16252400 | 20.0 |
| 1,1'-Biphenyl, ... | 16.569 | 35.5    | ng/ul | 28828600 | 4                     | 17.345 | 16252400 | 20.0 |
| 1,1'-Biphenyl, ... | 16.863 | 5.4     | ng/ul | 4422870  | 4                     | 17.345 | 16252400 | 20.0 |
| 1,1'-Biphenyl, ... | 17.222 | 45.7    | ng/ul | 37157900 | 4                     | 17.345 | 16252400 | 20.0 |
| 1,1'-Biphenyl, ... | 17.257 | 16.5    | ng/ul | 13442900 | 4                     | 17.345 | 16252400 | 20.0 |
| 2,2',4-Trichlor... | 17.516 | 28.3    | ng/ul | 23023500 | 4                     | 17.345 | 16252400 | 20.0 |
| 1,1'-Biphenyl, ... | 17.792 | 8.1     | ng/ul | 6585520  | 4                     | 17.345 | 16252400 | 20.0 |
| 1,1'-Biphenyl, ... | 17.822 | 2.9     | ng/ul | 2383560  | 4                     | 17.345 | 16252400 | 20.0 |
| 1,1'-Biphenyl, ... | 17.939 | 69.1    | ng/ul | 56131700 | 4                     | 17.345 | 16252400 | 20.0 |
| 1,1'-Biphenyl, ... | 18.063 | 32.1    | ng/ul | 26083400 | 4                     | 17.345 | 16252400 | 20.0 |
| 1,1'-Biphenyl, ... | 18.180 | 15.6    | ng/ul | 12652900 | 4                     | 17.345 | 16252400 | 20.0 |
| 1,1'-Biphenyl, ... | 18.233 | 5.4     | ng/ul | 4389210  | 4                     | 17.345 | 16252400 | 20.0 |
| 1,1'-Biphenyl, ... | 18.339 | 2.3     | ng/ul | 1844790  | 4                     | 17.345 | 16252400 | 20.0 |
| 2,2',4,6'-Tetra... | 18.392 | 20.7    | ng/ul | 16807300 | 4                     | 17.345 | 16252400 | 20.0 |
| 1,1'-Biphenyl, ... | 18.451 | 14.3    | ng/ul | 11627300 | 4                     | 17.345 | 16252400 | 20.0 |
| 2,2',4,5-Tetrac... | 18.492 | 12.1    | ng/ul | 9819300  | 4                     | 17.345 | 16252400 | 20.0 |
| 1,1'-Biphenyl, ... | 18.669 | 20.4    | ng/ul | 16581200 | 4                     | 17.345 | 16252400 | 20.0 |
| 1,1'-Biphenyl, ... | 18.710 | 6.8     | ng/ul | 5533060  | 4                     | 17.345 | 16252400 | 20.0 |
| unknown-01         | 18.763 | 7.6     | ng/ul | 6172050  | 4                     | 17.345 | 16252400 | 20.0 |
| 2,3',4,5-Tetrac... | 18.798 | 5.9     | ng/ul | 4772180  | 4                     | 17.345 | 16252400 | 20.0 |
| 2,3,3',6-Tetrac... | 18.833 | 13.2    | ng/ul | 10711400 | 4                     | 17.345 | 16252400 | 20.0 |
| 2,2',3,6-Tetrac... | 18.933 | 3.8     | ng/ul | 3102880  | 4                     | 17.345 | 16252400 | 20.0 |
| 1,1'-Biphenyl, ... | 19.133 | 3.5     | ng/ul | 2852130  | 4                     | 17.345 | 16252400 | 20.0 |
| 1,1'-Biphenyl, ... | 19.180 | 6.5     | ng/ul | 5309630  | 4                     | 17.345 | 16252400 | 20.0 |
| 3,3',4,5-Tetrac... | 19.216 | 5.9     | ng/ul | 4810610  | 4                     | 17.345 | 16252400 | 20.0 |
| Biphenyl, 2,4,4... | 19.422 | 7.0     | ng/ul | 1717310  | 5                     | 21.433 | 4932310  | 20.0 |
| 2,2',3,5,6'-Pen... | 19.480 | 3.3     | ng/ul | 819293   | 5                     | 21.433 | 4932310  | 20.0 |