

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121824\
 Data File : BP023555.D
 Acq On : 20 Dec 2024 15:42
 Operator : RC/JU
 Sample : AR1660 50PPM
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 AR1660 50PPM

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-BP112624.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP023555.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.334	56	59	66	rVB	23012	27172	0.72%	0.101%
2	3.805	136	139	145	rVB	13193	18780	0.50%	0.070%
3	7.516	763	770	787	rBV	775742	1466172	38.84%	5.457%
4	10.263	1229	1237	1259	rBV	988451	1891752	50.11%	7.042%
5	14.140	1886	1896	1915	rBV	1445529	2583616	68.44%	9.617%
6	14.275	1915	1919	1927	rVB2	23941	40493	1.07%	0.151%
7	15.116	2056	2062	2068	rBV5	5824	12018	0.32%	0.045%
8	15.404	2104	2111	2122	rBV	161744	265461	7.03%	0.988%
9	15.875	2183	2191	2201	rBV	41139	71574	1.90%	0.266%
10	16.040	2213	2219	2225	rBV2	74722	117534	3.11%	0.437%
11	16.157	2232	2239	2252	rBV	375141	583424	15.45%	2.172%
12	16.457	2285	2290	2297	rBV	54197	71727	1.90%	0.267%
13	16.734	2333	2337	2345	rBV3	8758	18890	0.50%	0.070%
14	16.810	2345	2350	2354	rBV	522706	716385	18.98%	2.667%
15	16.851	2354	2357	2366	rVB	185781	276361	7.32%	1.029%
16	16.951	2366	2374	2390	rBV2	1987858	3299952	87.41%	12.283%
17	17.122	2397	2403	2410	rVB2	268908	431431	11.43%	1.606%
18	17.334	2434	2439	2443	rBV7	6162	11088	0.29%	0.041%
19	17.422	2448	2454	2457	rBV	80558	126243	3.34%	0.470%
20	17.451	2457	2459	2466	rVB	40109	46828	1.24%	0.174%
21	17.563	2471	2478	2491	rVB	558661	1216767	32.23%	4.529%
22	17.704	2494	2502	2509	rVB	347282	536654	14.22%	1.998%
23	17.769	2509	2513	2517	rBV5	19430	24530	0.65%	0.091%
24	17.828	2518	2523	2528	rBV	166574	244455	6.48%	0.910%
25	17.869	2528	2530	2541	rVB3	56189	118892	3.15%	0.443%
26	17.992	2547	2551	2557	rVB4	23828	33412	0.89%	0.124%
27	18.057	2557	2562	2568	rBV	237935	354312	9.39%	1.319%
28	18.116	2568	2572	2576	rVV	160817	225926	5.98%	0.841%
29	18.157	2576	2579	2586	rVB	151757	190573	5.05%	0.709%
30	18.351	2607	2612	2618	rBV	200089	328352	8.70%	1.222%
31	18.404	2618	2621	2627	rVB3	69706	106008	2.81%	0.395%
32	18.481	2629	2634	2636	rBV	91213	144485	3.83%	0.538%
33	18.498	2636	2637	2640	rVV	84875	77518	2.05%	0.289%
34	18.539	2640	2644	2649	rVB	155938	214151	5.67%	0.797%
35	18.634	2656	2660	2666	rVB5	39054	62016	1.64%	0.231%
36	18.863	2696	2699	2704	rBV	48096	71084	1.88%	0.265%

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Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

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Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

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Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
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37	18.934	2704	2711	2723	rVB4	168673	399494	10.58%	1.487%
38	19.239	2758	2763	2770	rBV	172043	261582	6.93%	0.974%
39	19.681	2835	2838	2841	rBV3	60431	79710	2.11%	0.297%
40	19.710	2841	2843	2848	rVB	83161	86768	2.30%	0.323%
41	19.833	2860	2864	2868	rBV2	144186	196012	5.19%	0.730%
42	19.886	2869	2873	2879	rVB4	55017	103410	2.74%	0.385%
43	19.986	2885	2890	2894	rBV	394447	495258	13.12%	1.843%
44	20.281	2935	2940	2944	rBV	421960	519244	13.75%	1.933%
45	20.333	2945	2949	2954	rBV4	98230	149754	3.97%	0.557%
46	20.445	2964	2968	2973	rBV3	183633	271136	7.18%	1.009%
47	20.604	2989	2995	3002	rVB	221246	410197	10.87%	1.527%
48	20.763	3018	3022	3027	rBV2	280968	344517	9.13%	1.282%
49	20.828	3030	3033	3037	rVV3	136936	151639	4.02%	0.564%
50	21.063	3069	3073	3078	rBV2	213732	275250	7.29%	1.025%
51	21.133	3081	3085	3090	rVB3	106587	173003	4.58%	0.644%
52	21.380	3122	3127	3140	rBV2	1971698	3775154	100.00%	14.052%
53	24.557	3660	3667	3685	rVB	1251977	3147541	83.38%	11.716%

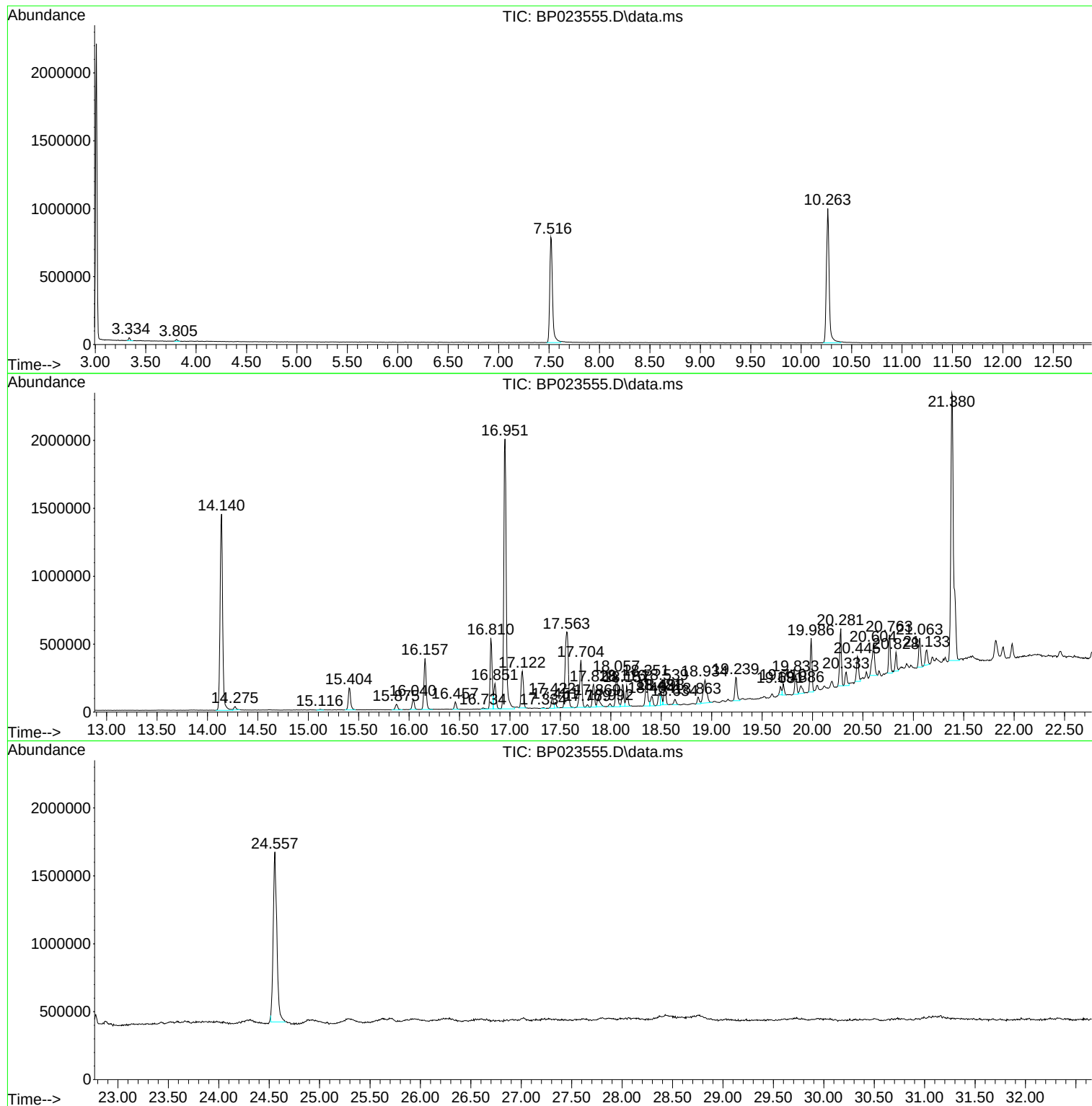
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Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
AR1660 50PPM

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

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



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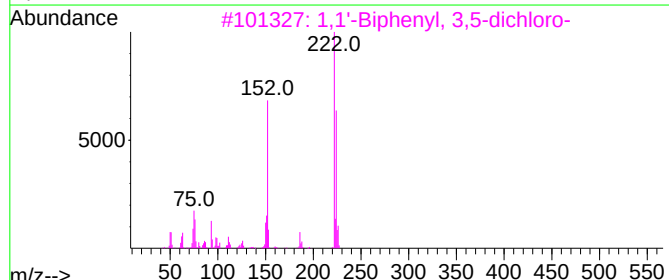
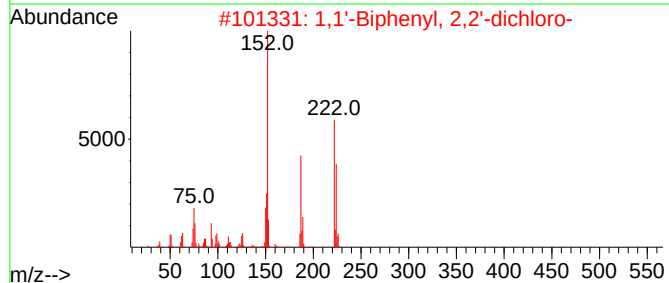
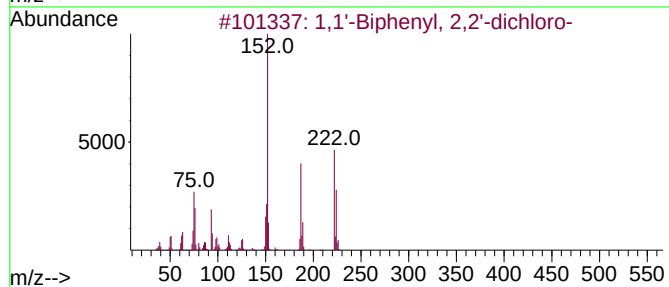
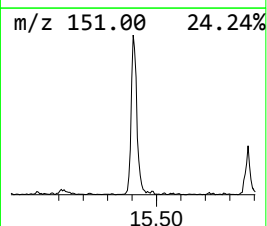
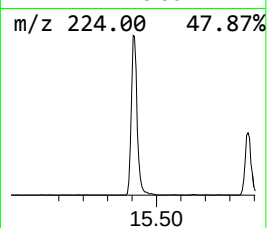
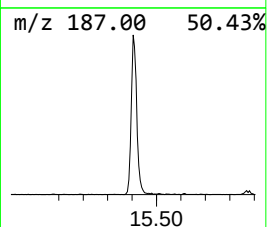
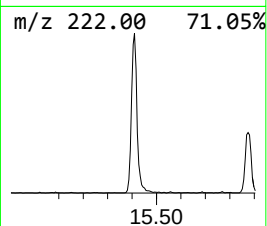
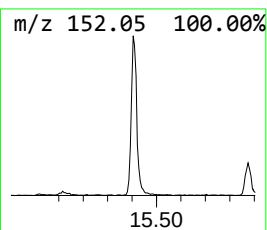
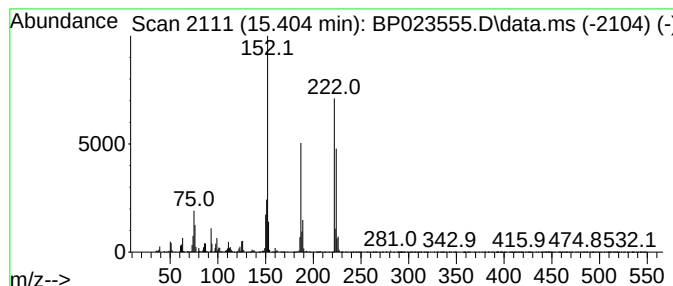
Instrument :
BNA_P
ClientSampleId :
AR1660 50PPM

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 1,1'-Biphenyl, 2,2'-dichloro- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.404	2.05 ng/ul	265461	Acenaphthene-d10	14.140	
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	98
3	1,1'-Biphenyl, 3,5-dichloro-	222	C12H8Cl2	034883-41-5	95
4	1,1'-Biphenyl, 2,4'-dichloro-	222	C12H8Cl2	034883-43-7	95
5	1,1'-Biphenyl, 2,2'-dichloro-	222	C12H8Cl2	013029-08-8	95



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ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
AR1660 50PPM

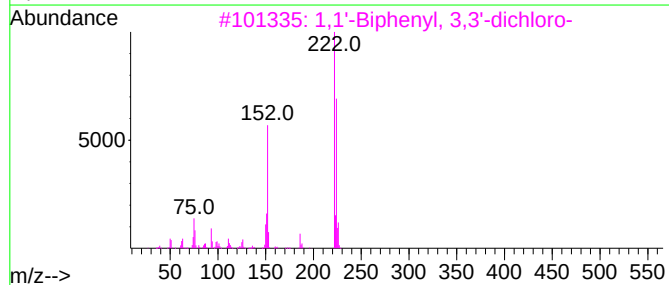
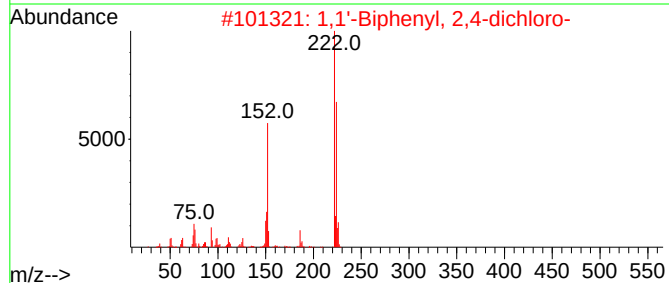
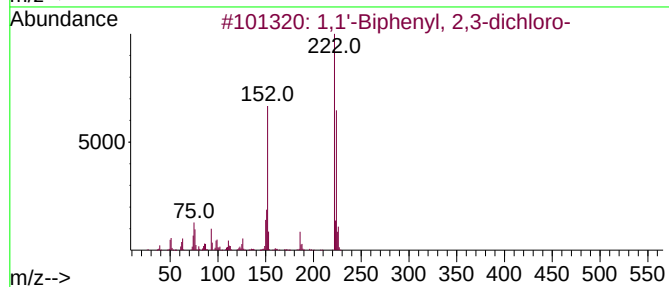
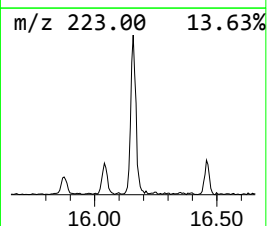
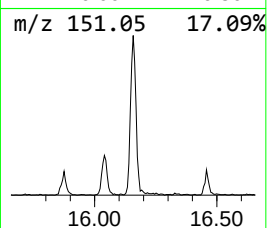
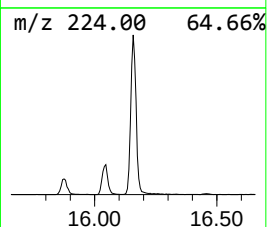
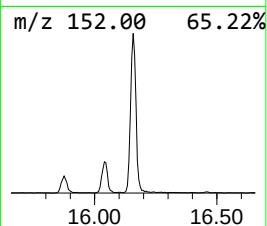
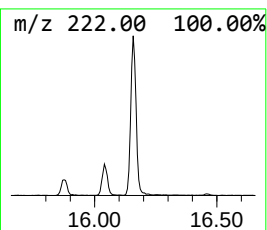
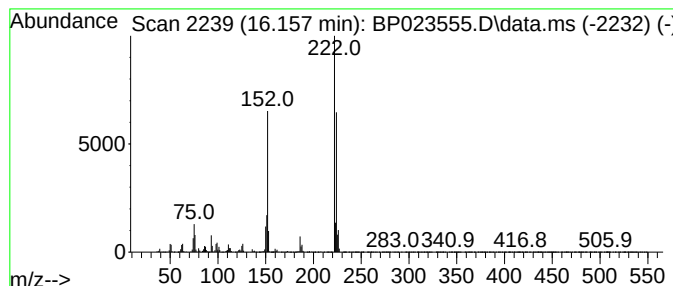
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.157	3.54 ng/ul	583424	Phenanthrene-d10	16.951

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, 3,3'-dichloro-	222	C12H8Cl2	002050-67-1	98		
4	1,1'-Biphenyl, 2,5-dichloro-	222	C12H8Cl2	034883-39-1	98		
5	1,1'-Biphenyl, 2,4-dichloro-	222	C12H8Cl2	033284-50-3	98		



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Misc :
ALS Vial : 6 Sample Multiplier: 1

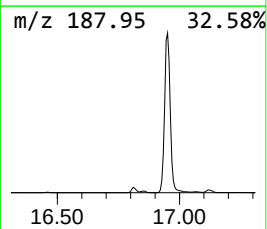
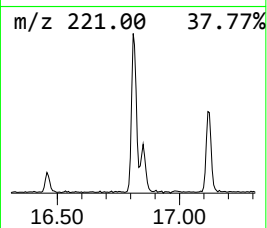
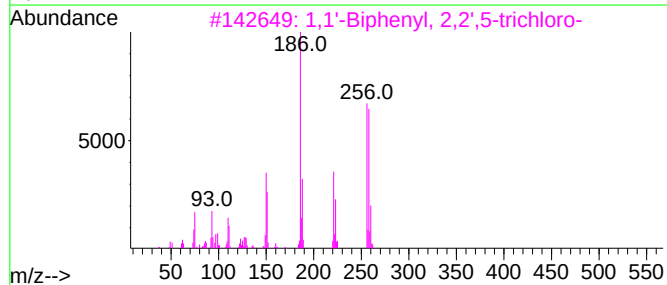
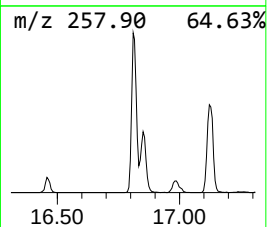
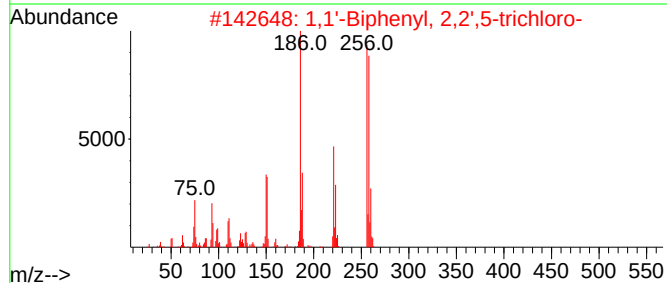
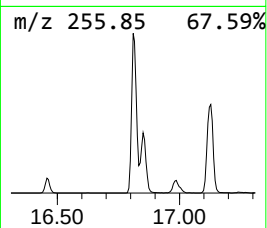
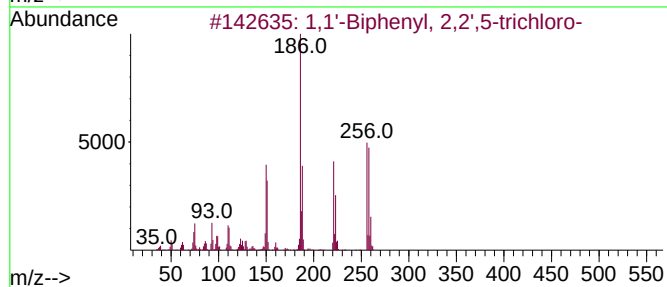
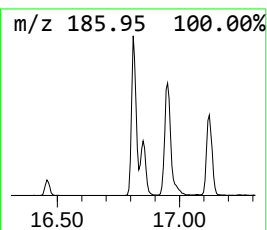
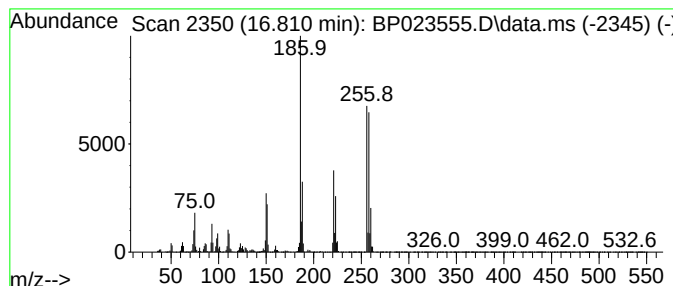
Instrument :
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ClientSampleId :
AR1660 50PPM

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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
16.810	4.34 ng/ul	716385	Phenanthrene-d10	16.951	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',5-trichloro-	256	C12H7Cl3	037680-65-2	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, 2,4,4'-trichloro-	256	C12H7Cl3	007012-37-5	99
5	2,2',3-Trichloro-1,1'-biphenyl	256	C12H7Cl3	038444-78-9	99



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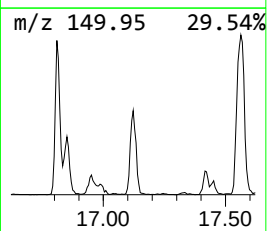
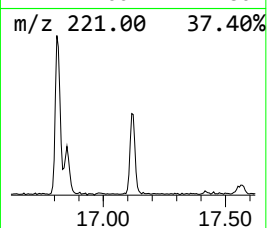
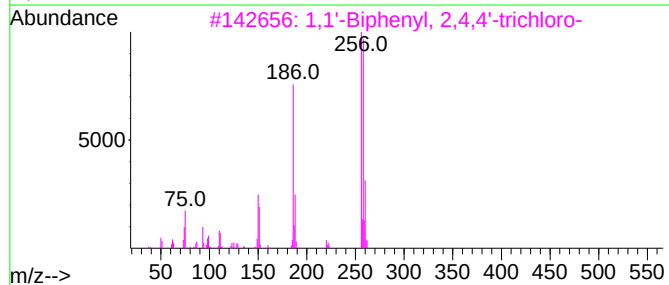
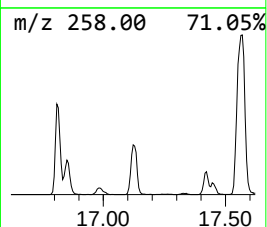
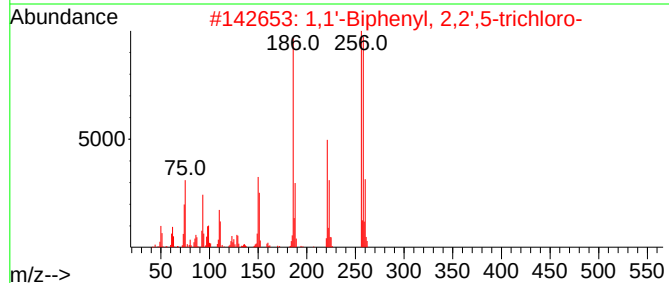
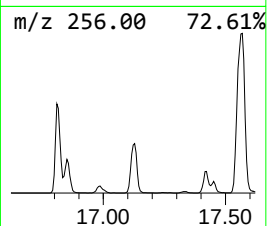
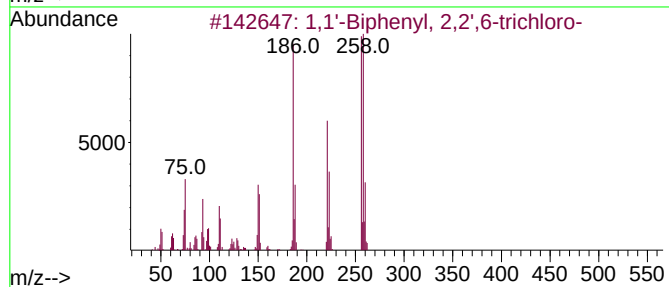
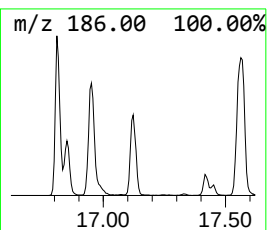
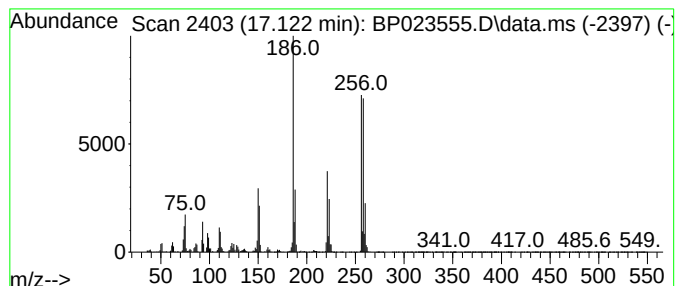
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 1,1'-Biphenyl, 2,2',6-trich... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.122	2.61 ng/ul	431431	Phenanthrene-d10	16.951

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	2,2',3-Trichloro-1,1'-biphenyl	256	C12H7Cl3	038444-78-9	99		
5	1,1'-Biphenyl, 2,2',5-trichloro-	256	C12H7Cl3	037680-65-2	99		



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AR1660 50PPM

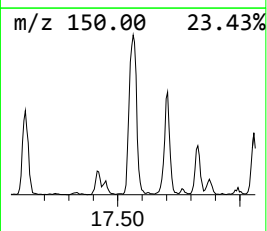
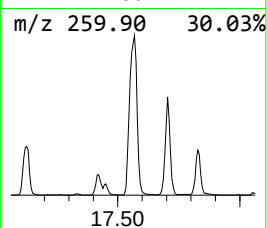
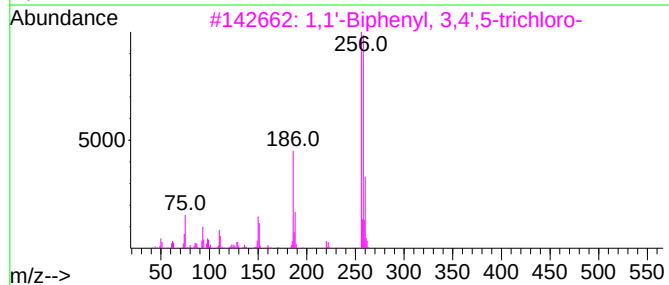
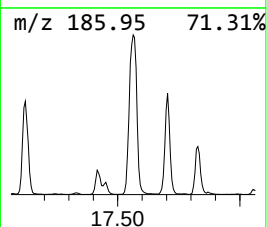
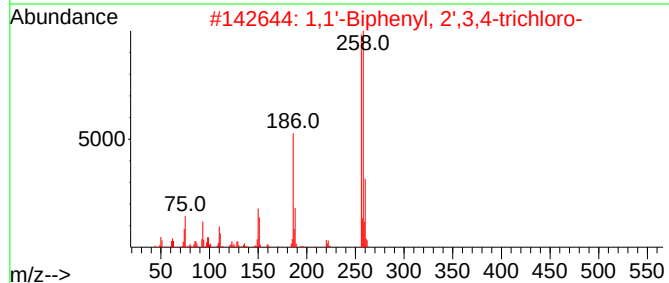
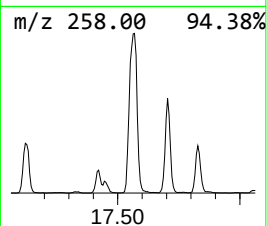
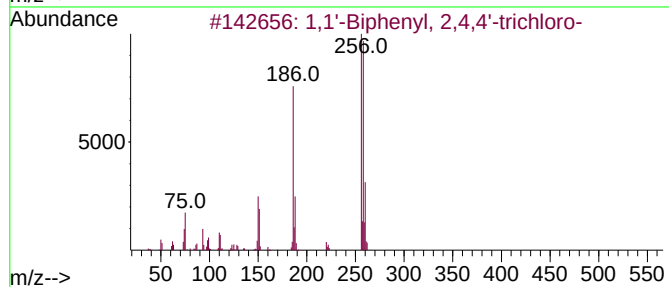
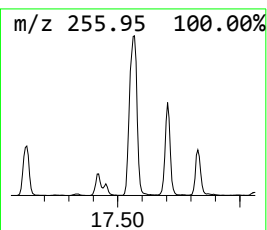
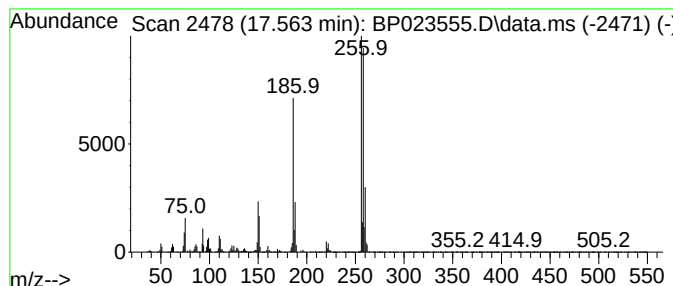
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.563	7.37 ng/ul	1216770	Phenanthrene-d10	16.951

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',3,4-trichloro-	256	C12H7Cl3	038444-86-9	99		
3	1,1'-Biphenyl, 3,4',5-trichloro-	256	C12H7Cl3	038444-88-1	99		
4	1,1'-Biphenyl, 2,3',5-trichloro-	256	C12H7Cl3	038444-81-4	99		
5	1,1'-Biphenyl, 3,3',5-trichloro-	256	C12H7Cl3	038444-87-0	99		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121824\
Data File : BP023555.D
Acq On : 20 Dec 2024 15:42
Operator : RC/JU
Sample : AR1660 50PPM
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
AR1660 50PPM

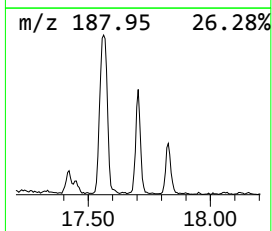
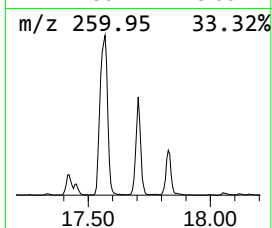
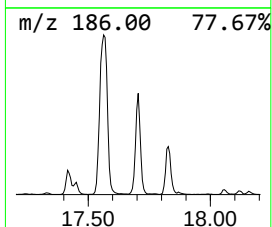
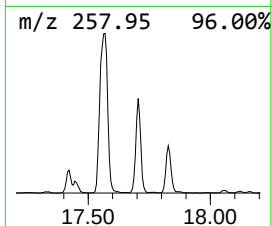
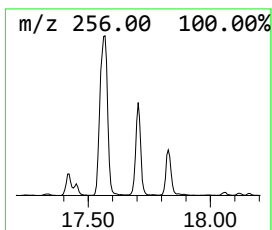
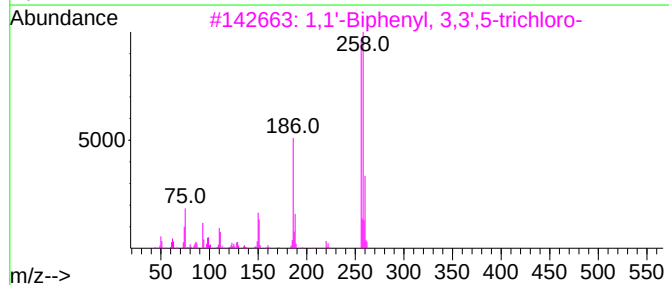
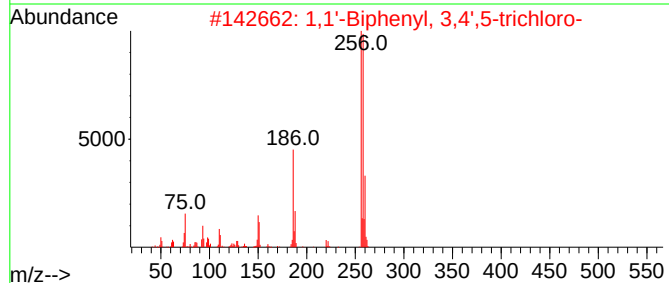
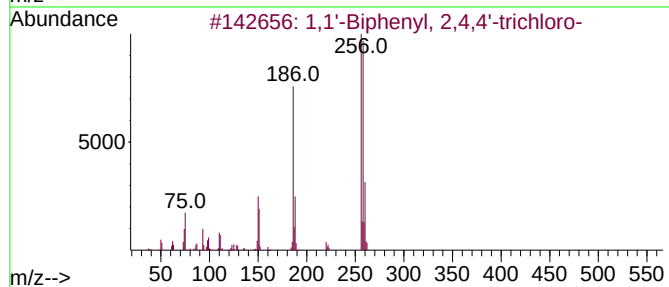
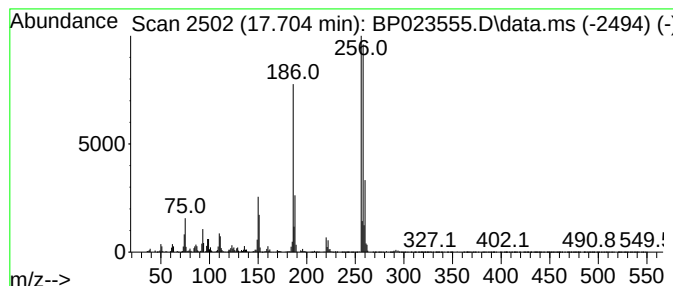
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 1,1'-Biphenyl, 3,4',5-trich... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.704	3.25 ng/ul	536654	Phenanthrene-d10	16.951

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, 3,4',5-trichloro-	256	C12H7Cl3	038444-88-1	99	
3	1,1'-Biphenyl, 3,3',5-trichloro-	256	C12H7Cl3	038444-87-0	99	
4	1,1'-Biphenyl, 2,4',5-trichloro-	256	C12H7Cl3	016606-02-3	99	
5	1,1'-Biphenyl, 2,3',5-trichloro-	256	C12H7Cl3	038444-81-4	99	



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121824\
Data File : BP023555.D
Acq On : 20 Dec 2024 15:42
Operator : RC/JU
Sample : AR1660 50PPM
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
AR1660 50PPM

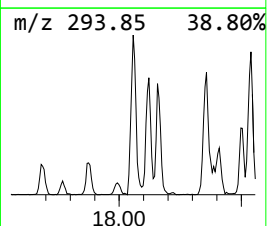
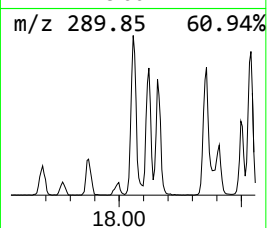
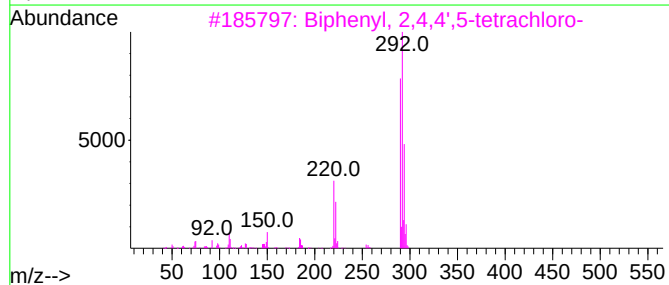
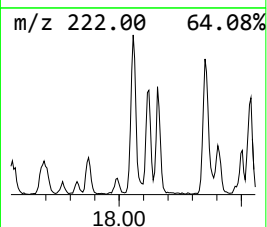
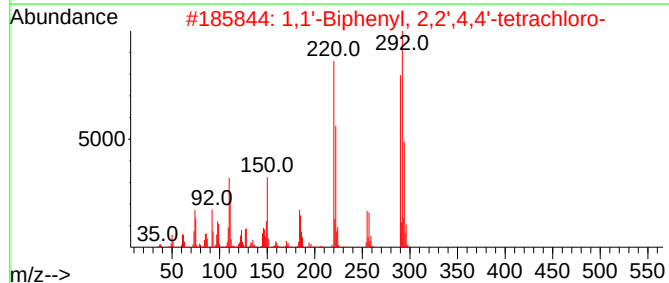
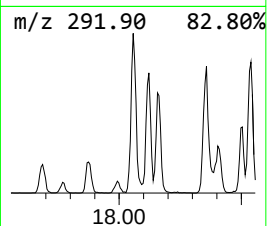
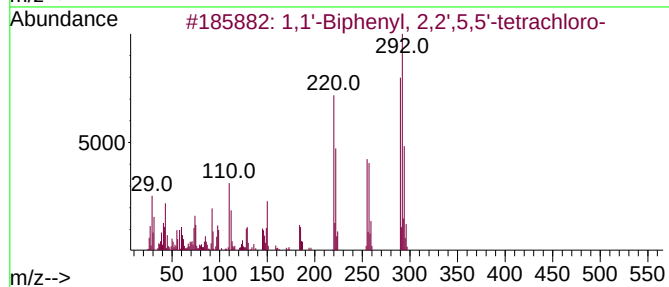
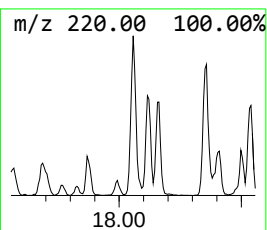
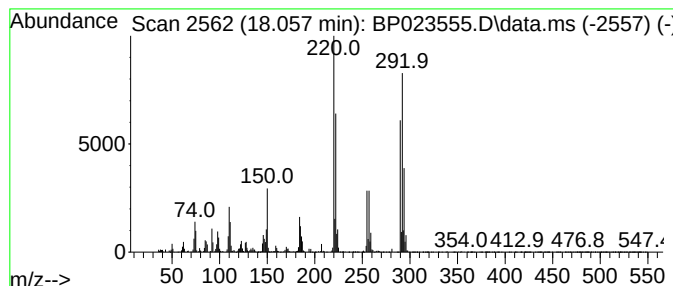
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 1,1'-Biphenyl, 2,2',5,5'-te... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.057	2.15 ng/ul	354312	Phenanthrene-d10	16.951

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',5,5'-tetrach...	290	C12H6Cl4	035693-99-3	99		
2	1,1'-Biphenyl, 2,2',4,4'-tetrach...	290	C12H6Cl4	002437-79-8	99		
3	Biphenyl, 2,4,4',5-tetrachloro-	290	C12H6Cl4	032690-93-0	99		
4	2,2',3,6-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-45-7	99		
5	1,1'-Biphenyl, 2,3',4,6-Tetrachl...	290	C12H6Cl4	060233-24-1	99		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121824\
Data File : BP023555.D
Acq On : 20 Dec 2024 15:42
Operator : RC/JU
Sample : AR1660 50PPM
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
AR1660 50PPM

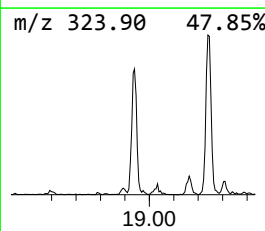
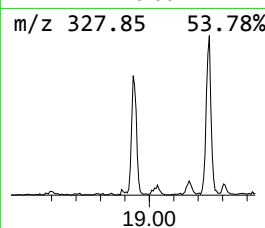
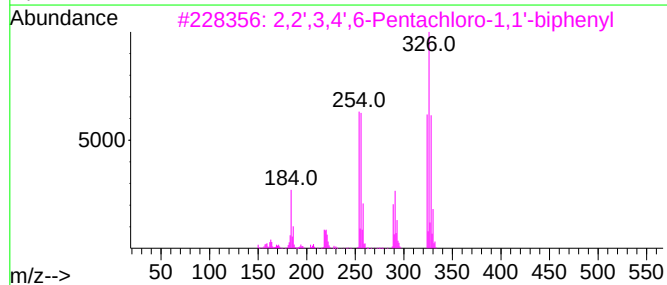
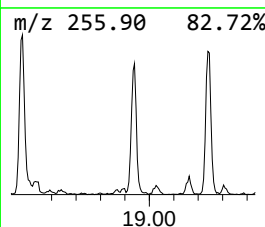
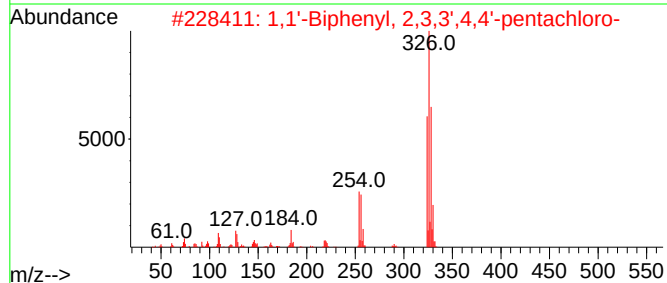
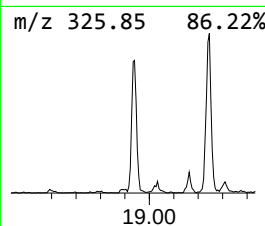
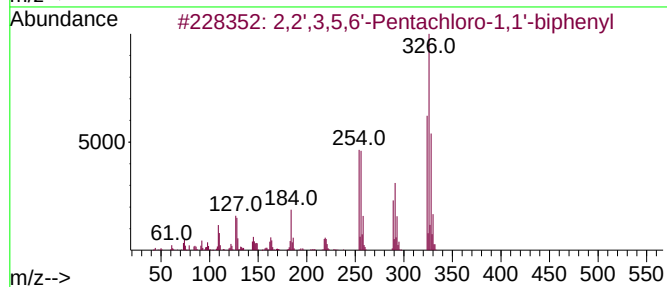
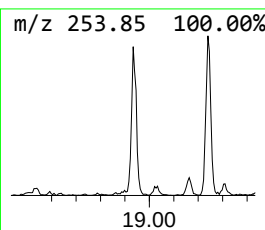
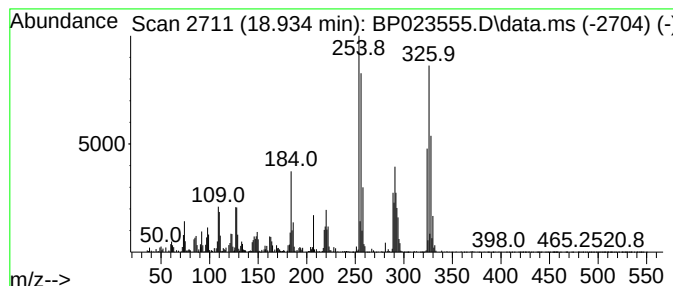
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 2,2',3,5,6'-Pentachloro-1,1... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.933	2.42 ng/ul	399494	Phenanthrene-d10	16.951

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, 2,3,3',4,4'-penta...	324	C12H5Cl5	032598-14-4	99		
3	2,2',3,4',6-Pentachloro-1,1'-bip...	324	C12H5Cl5	068194-05-8	99		
4	1,1'-Biphenyl, 2,2',3',4,6-Penta...	324	C12H5Cl5	060233-25-2	99		
5	1,1'-Biphenyl, 2,2',3,4,6-Pentac...	324	C12H5Cl5	055215-17-3	99		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121824\
Data File : BP023555.D
Acq On : 20 Dec 2024 15:42
Operator : RC/JU
Sample : AR1660 50PPM
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
AR1660 50PPM

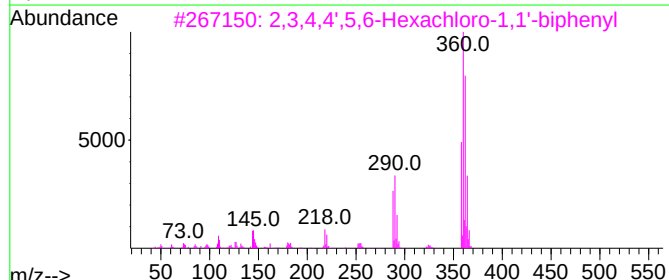
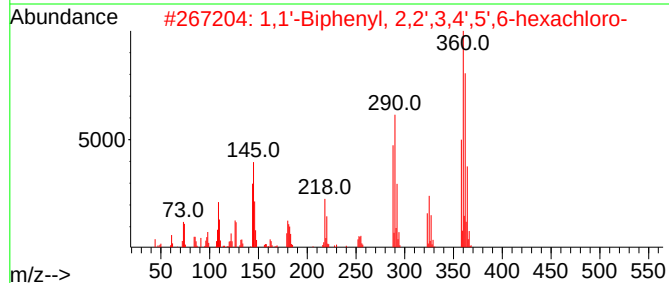
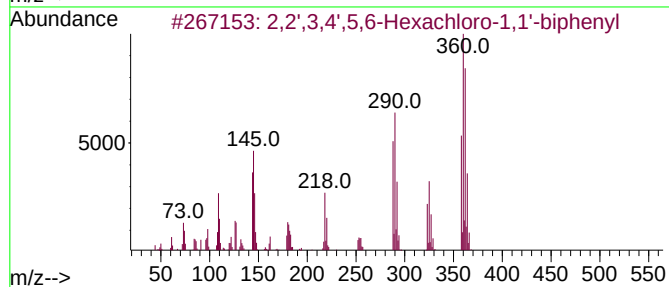
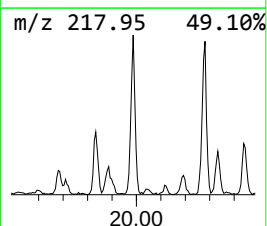
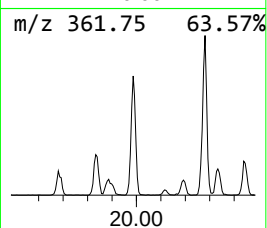
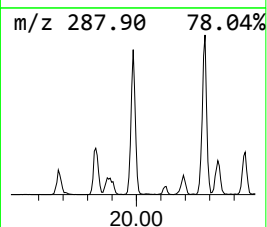
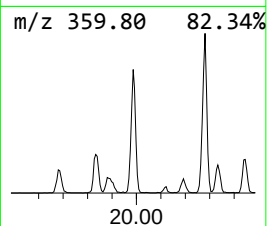
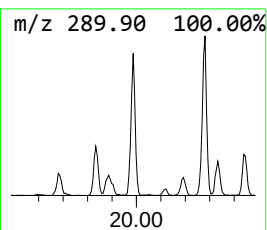
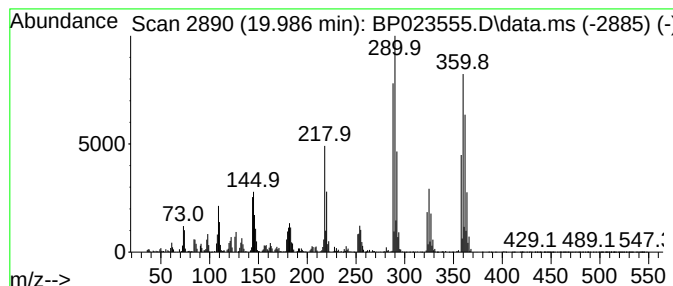
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 2,2',3,4',5,6-Hexachloro-1,... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.986	2.62 ng/ul	495258	Chrysene-d12	21.381

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2',3,4',5,6-Hexachloro-1,1'-bi...	358	C12H4Cl6	068194-13-8	99		
2	1,1'-Biphenyl, 2,2',3,4',5',6-he...	358	C12H4Cl6	038380-04-0	99		
3	2,3,4,4',5,6-Hexachloro-1,1'-bip...	358	C12H4Cl6	041411-63-6	99		
4	1,1'-Biphenyl, 2,2',4,4',5',6-He...	358	C12H4Cl6	060145-22-4	99		
5	2,2',3,5,6,6'-Hexachloro-1,1'-bi...	358	C12H4Cl6	068194-09-2	99		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121824\
Data File : BP023555.D
Acq On : 20 Dec 2024 15:42
Operator : RC/JU
Sample : AR1660 50PPM
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
AR1660 50PPM

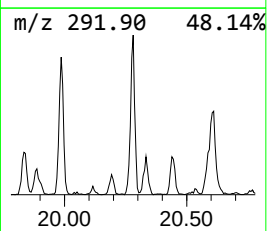
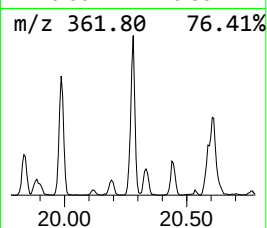
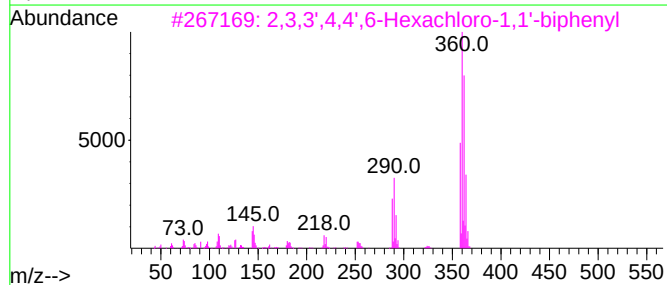
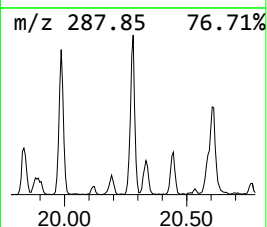
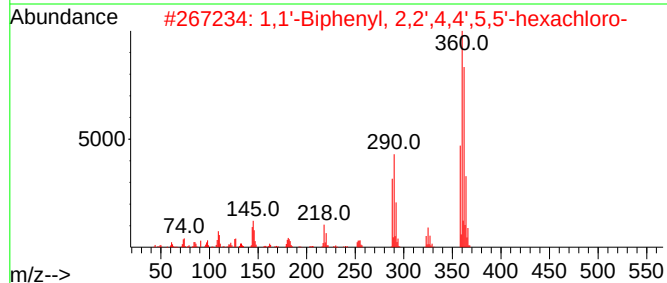
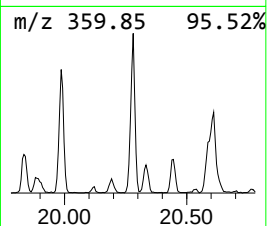
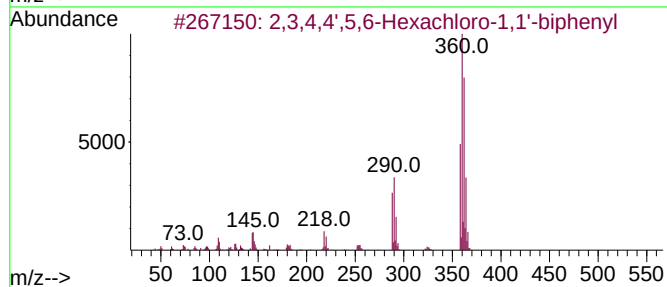
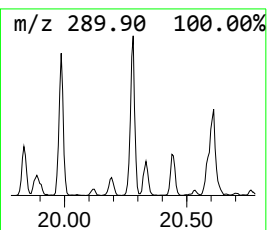
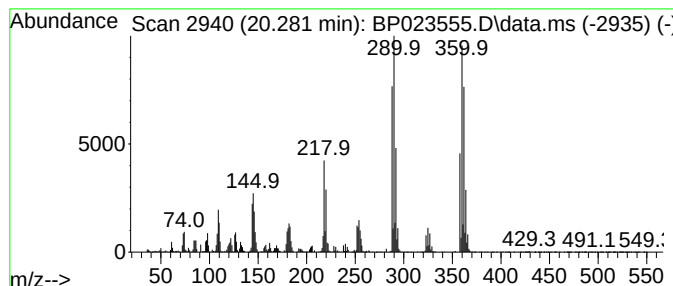
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 2,3,4,4',5,6-Hexachloro-1,1... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.281	2.75 ng/ul	519244	Chrysene-d12	21.381

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,3,4,4',5,6-Hexachloro-1,1'-bip...	358	C12H4Cl6	041411-63-6	99		
2	1,1'-Biphenyl, 2,2',4,4',5,5'-he...	358	C12H4Cl6	035065-27-1	99		
3	2,3,3',4,4',6-Hexachloro-1,1'-bi...	358	C12H4Cl6	074472-42-7	99		
4	2,3,3',4',5,6-Hexachloro-1,1'-bi...	358	C12H4Cl6	074472-44-9	99		
5	2,3,3',4,5,5'-Hexachloro-1,1'-bi...	358	C12H4Cl6	039635-35-3	99		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121824\
Data File : BP023555.D
Acq On : 20 Dec 2024 15:42
Operator : RC/JU
Sample : AR1660 50PPM
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
AR1660 50PPM

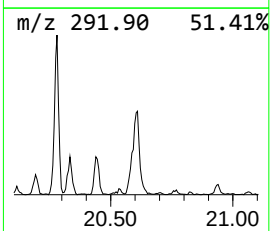
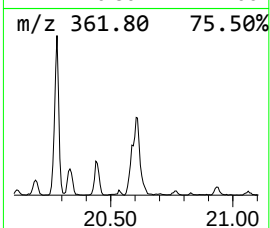
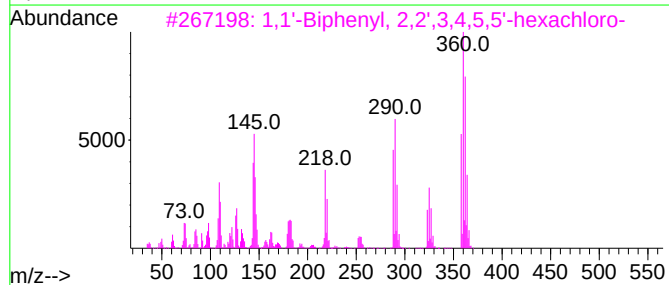
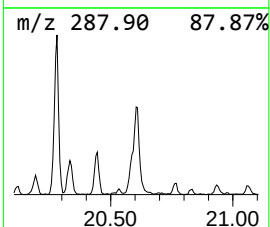
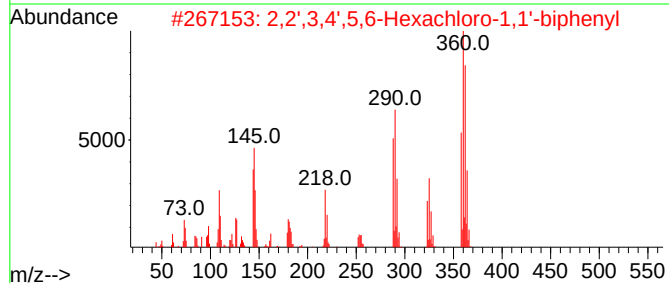
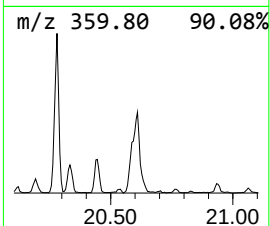
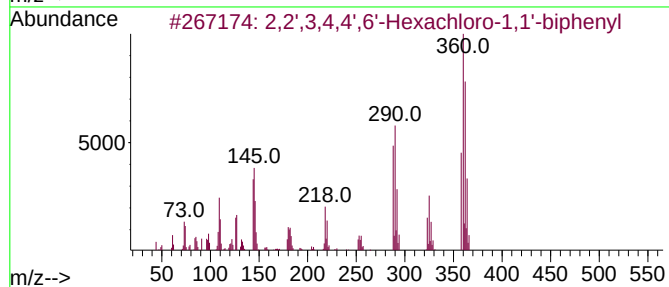
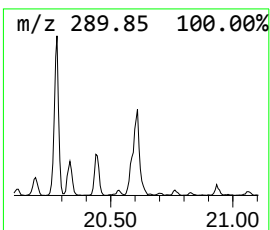
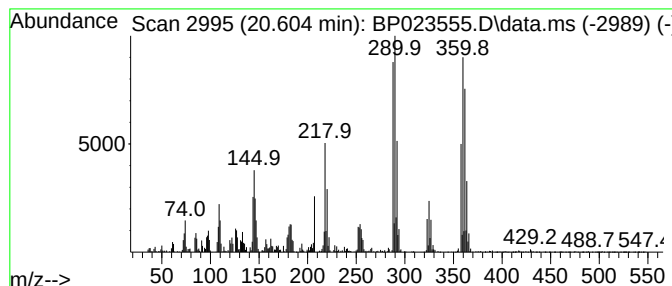
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 11 2,2',3,4,4',6'-Hexachloro-1... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.604	2.17 ng/ul	410197	Chrysene-d12	21.381

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2',3,4,4',6'-Hexachloro-1,1'-b...	358	C12H4Cl6	059291-64-4	99		
2	2,2',3,4',5,6-Hexachloro-1,1'-bi...	358	C12H4Cl6	068194-13-8	99		
3	1,1'-Biphenyl, 2,2',3,4,5,5'-hex...	358	C12H4Cl6	052712-04-6	99		
4	2,3,4,4',5,6-Hexachloro-1,1'-bip...	358	C12H4Cl6	041411-63-6	99		
5	2,2',3,5,6,6'-Hexachloro-1,1'-bi...	358	C12H4Cl6	068194-09-2	99		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121824\
Data File : BP023555.D
Acq On : 20 Dec 2024 15:42
Operator : RC/JU
Sample : AR1660 50PPM
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
AR1660 50PPM

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
1,1'-Biphenyl, ...	15.404	2.0	ng/ul	265461	3	14.140	2583620	20.0
1,1'-Biphenyl, ...	16.157	3.5	ng/ul	583424	4	16.951	3299950	20.0
1,1'-Biphenyl, ...	16.810	4.3	ng/ul	716385	4	16.951	3299950	20.0
1,1'-Biphenyl, ...	17.122	2.6	ng/ul	431431	4	16.951	3299950	20.0
1,1'-Biphenyl, ...	17.563	7.4	ng/ul	1216770	4	16.951	3299950	20.0
1,1'-Biphenyl, ...	17.704	3.3	ng/ul	536654	4	16.951	3299950	20.0
1,1'-Biphenyl, ...	18.057	2.1	ng/ul	354312	4	16.951	3299950	20.0
2,2',3,5,6'-Pen...	18.933	2.4	ng/ul	399494	4	16.951	3299950	20.0
2,2',3,4',5,6-H...	19.986	2.6	ng/ul	495258	5	21.381	3775150	20.0
2,3,4,4',5,6-He...	20.281	2.8	ng/ul	519244	5	21.381	3775150	20.0
2,2',3,4,4',6'-...	20.604	2.2	ng/ul	410197	5	21.381	3775150	20.0