

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012324\
 Data File : BP019420.D
 Acq On : 23 Jan 2024 15:02
 Operator : MA/JU
 Sample : P1157-05
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AB2

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-BP010824.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP019420.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.129	3	7	12	rVB	949216	955548	49.14%	5.480%
2	3.464	61	64	68	rVB	21588	25763	1.32%	0.148%
3	3.946	140	146	151	rBV	43471	59313	3.05%	0.340%
4	4.058	162	165	168	rBV5	3942	4591	0.24%	0.026%
5	4.387	218	221	225	rVB2	8544	12628	0.65%	0.072%
6	4.487	235	238	244	rVB8	2739	5200	0.27%	0.030%
7	4.587	251	255	259	rVB6	4204	6438	0.33%	0.037%
8	4.699	269	274	277	rBV4	7388	10003	0.51%	0.057%
9	4.787	285	289	295	rVB	16174	22755	1.17%	0.131%
10	4.852	295	300	304	rVB8	3872	7599	0.39%	0.044%
11	5.205	356	360	362	rBV4	3835	4820	0.25%	0.028%
12	5.270	367	371	378	rBV	27540	39942	2.05%	0.229%
13	5.405	389	394	403	rVB2	13477	25003	1.29%	0.143%
14	5.775	452	457	465	rVB	15878	23407	1.20%	0.134%
15	6.028	494	500	503	rBV7	3547	6045	0.31%	0.035%
16	6.252	531	538	546	rBV2	21031	36191	1.86%	0.208%
17	6.311	546	548	554	rVB7	2877	4121	0.21%	0.024%
18	6.722	613	618	619	rBV4	2431	4167	0.21%	0.024%
19	6.752	619	623	628	rVB7	5951	10749	0.55%	0.062%
20	6.805	628	632	634	rBV5	4729	6273	0.32%	0.036%
21	6.852	636	640	644	rBV3	7478	10630	0.55%	0.061%
22	6.987	658	663	670	rVB2	9718	16398	0.84%	0.094%
23	7.140	684	689	691	rBV6	6350	9004	0.46%	0.052%
24	7.234	699	705	706	rBV5	3996	6855	0.35%	0.039%
25	7.252	706	708	712	rVB5	4127	3996	0.21%	0.023%
26	7.311	714	718	724	rVV9	2816	4949	0.25%	0.028%
27	7.381	726	730	732	rVV2	6962	9949	0.51%	0.057%
28	7.411	732	735	742	rVB	14313	20616	1.06%	0.118%
29	7.658	769	777	784	rBV	5363	15749	0.81%	0.090%
30	7.769	787	796	810	rBV	528102	847617	43.59%	4.861%
31	7.887	810	816	817	rBV4	5443	9031	0.46%	0.052%
32	8.140	853	859	865	rVB6	5896	9869	0.51%	0.057%
33	8.316	883	889	891	rVV7	2720	4537	0.23%	0.026%
34	8.405	900	904	910	rBV8	5052	8921	0.46%	0.051%
35	8.875	977	984	992	rVB7	5114	11702	0.60%	0.067%
36	8.987	1000	1003	1008	rVB3	5606	8199	0.42%	0.047%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010824.MA.M
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37	9.399	1067	1073	1077	rBV8	3932	7893	0.41%	0.045%
38	9.452	1079	1082	1088	rVB6	4004	7008	0.36%	0.040%
39	9.963	1163	1169	1176	rVB10	4172	8499	0.44%	0.049%
40	10.187	1202	1207	1212	rVB9	2921	6193	0.32%	0.036%
41	10.234	1212	1215	1218	rBV5	2531	3750	0.19%	0.022%
42	10.334	1228	1232	1236	rBV6	2223	3574	0.18%	0.020%
43	10.428	1244	1248	1255	rBV10	3555	9935	0.51%	0.057%
44	10.563	1262	1271	1276	rBV	670401	1127883	58.00%	6.469%
45	10.610	1276	1279	1286	rVB	84095	145142	7.46%	0.832%
46	11.228	1377	1384	1388	rBV10	2769	6279	0.32%	0.036%
47	11.587	1441	1445	1448	rBV5	2402	3976	0.20%	0.023%
48	11.987	1509	1513	1517	rBV6	4080	6701	0.34%	0.038%
49	12.234	1549	1555	1564	rBV2	24974	51653	2.66%	0.296%
50	12.452	1584	1592	1601	rBV	18861	37718	1.94%	0.216%
51	12.610	1617	1619	1621	rBV3	3075	3499	0.18%	0.020%
52	12.752	1640	1643	1645	rBV4	2703	3576	0.18%	0.021%
53	12.834	1653	1657	1663	rVB9	5109	9622	0.49%	0.055%
54	12.887	1663	1666	1671	rBV6	4077	7706	0.40%	0.044%
55	12.934	1673	1674	1680	rVV6	4543	6806	0.35%	0.039%
56	13.004	1680	1686	1690	rVV9	3907	8168	0.42%	0.047%
57	13.181	1710	1716	1719	rBV8	7973	13890	0.71%	0.080%
58	13.252	1723	1728	1736	rVV4	9492	24104	1.24%	0.138%
59	13.587	1780	1785	1786	rBV4	6914	12061	0.62%	0.069%
60	13.734	1805	1810	1815	rBV5	13700	26729	1.37%	0.153%
61	13.787	1817	1819	1821	rVV2	7759	8823	0.45%	0.051%
62	13.816	1821	1824	1830	rVB4	22744	34684	1.78%	0.199%
63	13.893	1833	1837	1840	rBV4	9528	9356	0.48%	0.054%
64	13.934	1840	1844	1848	rBV6	6641	12617	0.65%	0.072%
65	14.004	1854	1856	1860	rVB5	5506	6573	0.34%	0.038%
66	14.140	1874	1879	1885	rBV8	19480	40123	2.06%	0.230%
67	14.228	1890	1894	1900	rVB3	34197	53441	2.75%	0.307%
68	14.310	1903	1908	1912	rBV5	12876	20242	1.04%	0.116%
69	14.363	1913	1917	1919	rBV4	15156	22092	1.14%	0.127%
70	14.416	1919	1926	1932	rVV	1085647	1607748	82.68%	9.221%
71	14.475	1932	1936	1945	rVV3	25565	47596	2.45%	0.273%
72	14.628	1958	1962	1969	rBV10	8836	20299	1.04%	0.116%
73	14.816	1990	1994	2003	rVB	43315	72252	3.72%	0.414%
74	15.040	2028	2032	2033	rBV4	11112	15430	0.79%	0.088%
75	15.193	2053	2058	2064	rBV4	37171	65149	3.35%	0.374%
76	15.469	2100	2105	2109	rBV	63044	93470	4.81%	0.536%

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77	15.675	2136	2140	2145	rBV6	30958	48283	2.48%	0.277%
78	15.916	2178	2181	2188	rBV6	16455	38446	1.98%	0.221%
79	16.151	2214	2221	2228	rBV7	36897	88684	4.56%	0.509%
80	16.463	2271	2274	2279	rVB4	29546	38136	1.96%	0.219%
81	17.169	2389	2394	2398	rBV	1351346	1944576	100.00%	11.153%
82	17.210	2398	2401	2408	rVB	389709	561849	28.89%	3.222%
83	18.228	2570	2574	2579	rBV2	236508	317587	16.33%	1.821%
84	18.287	2581	2584	2588	rVB	113979	145297	7.47%	0.833%
85	18.504	2618	2621	2624	rBV2	82098	115187	5.92%	0.661%
86	19.028	2706	2710	2711	rBV	134679	156846	8.07%	0.900%
87	19.051	2711	2714	2720	rVB2	198921	323809	16.65%	1.857%
88	19.192	2733	2738	2743	rBV	483359	633715	32.59%	3.635%
89	19.328	2757	2761	2767	rVB	295712	380284	19.56%	2.181%
90	19.386	2768	2771	2775	rVV	119645	133030	6.84%	0.763%
91	19.516	2790	2793	2794	rBV	74140	77021	3.96%	0.442%
92	19.545	2795	2798	2801	rVB	161827	201913	10.38%	1.158%
93	19.763	2831	2835	2841	rVB	286615	358378	18.43%	2.055%
94	20.016	2874	2878	2883	rBV4	177068	292911	15.06%	1.680%
95	20.069	2884	2887	2893	rVB	201134	229240	11.79%	1.315%
96	20.286	2921	2924	2928	rVB	161075	182683	9.39%	1.048%
97	20.598	2975	2977	2989	rVB6	157788	219617	11.29%	1.260%
98	21.275	3087	3092	3096	rBV	1376427	1852227	95.25%	10.623%
99	22.751	3338	3343	3350	rVB	904185	1453522	74.75%	8.337%
100	23.622	3485	3491	3502	rVB	845733	1701522	87.50%	9.759%

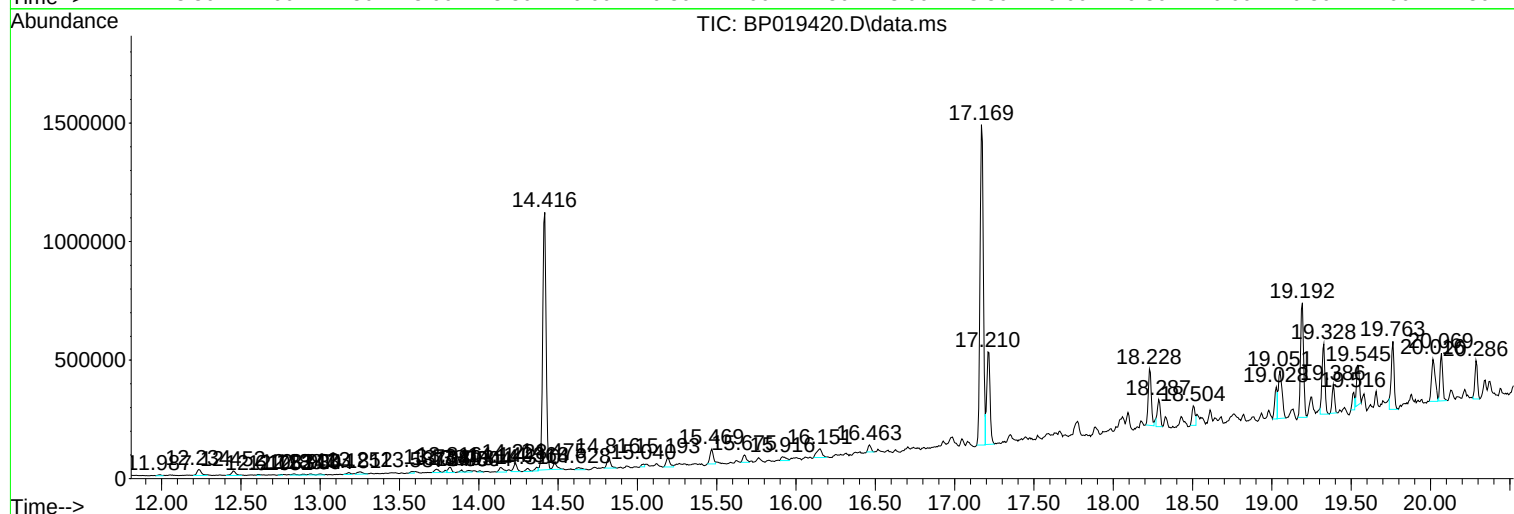
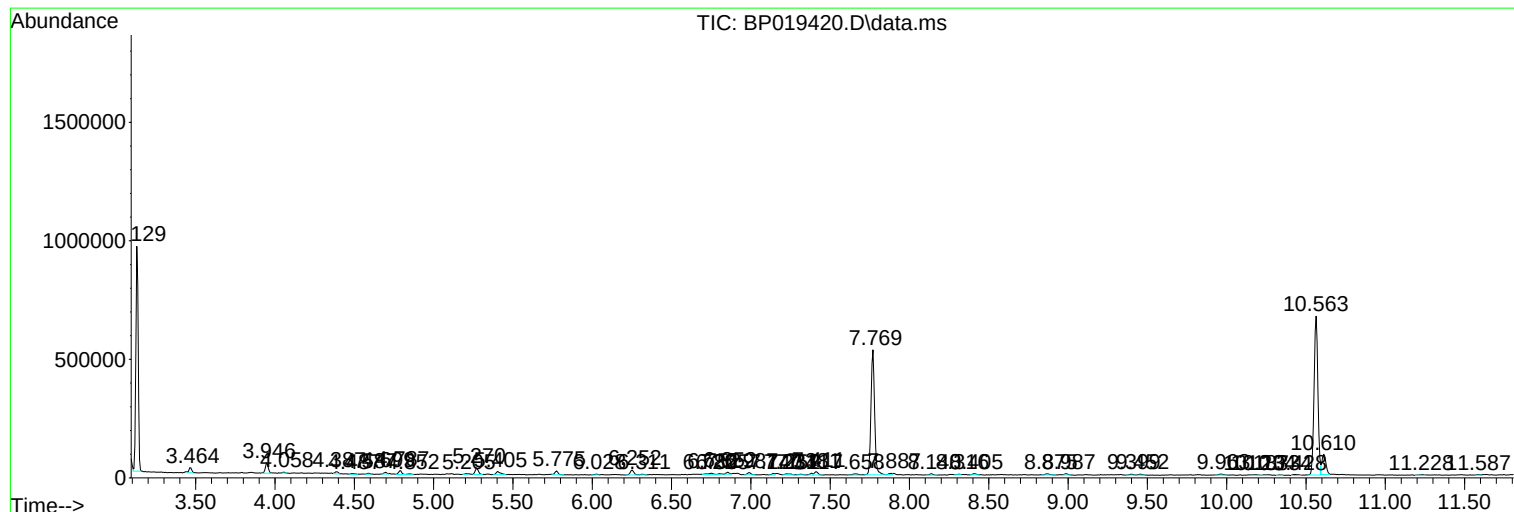
Sum of corrected areas: 17435601

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010824.MA.M
Quant Title : SVOA CALIBRATION

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



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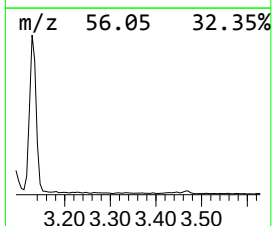
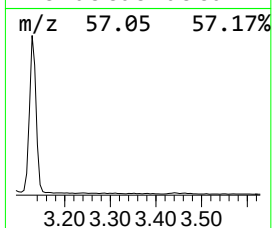
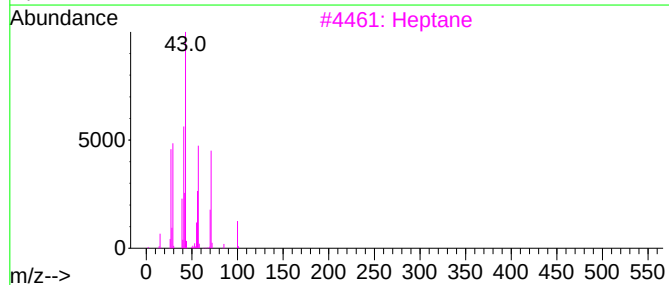
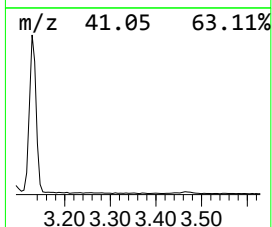
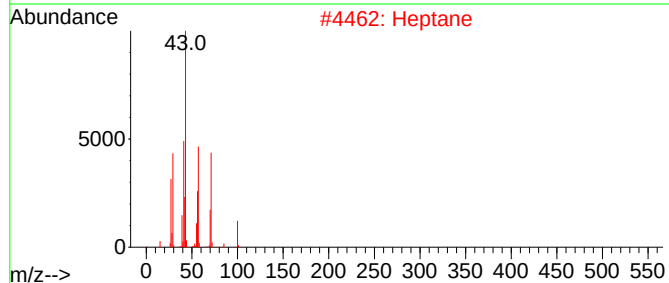
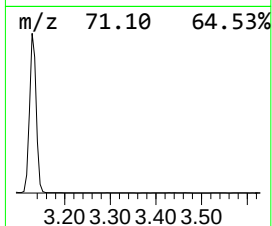
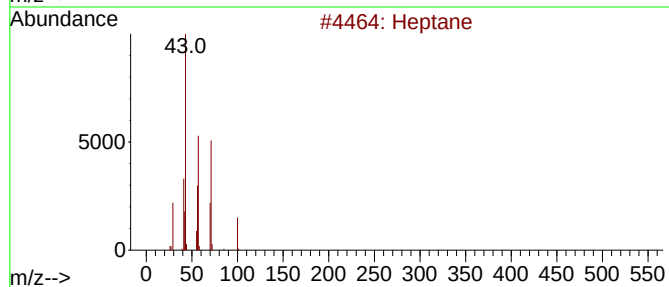
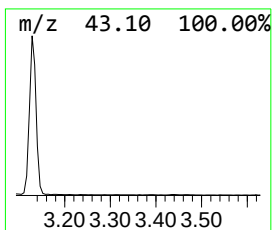
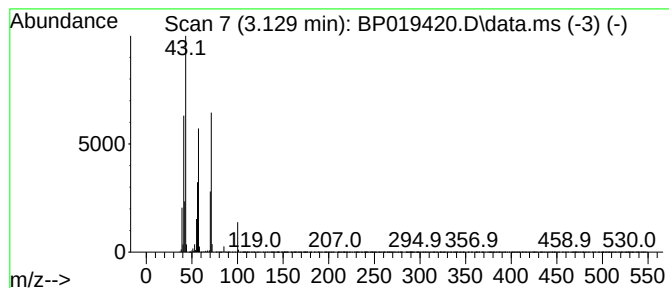
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010824.MA.M
Quant Title : SVOA CALIBRATION

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.
3.129	22.55 ng/ul	955548	1,4-Dichlorobenzene-d4	7.769

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane			100	C7H16	000142-82-5	91
2	Heptane			100	C7H16	000142-82-5	90
3	Heptane			100	C7H16	000142-82-5	90
4	Heptane			100	C7H16	000142-82-5	87
5	Hexane, 3-methyl-			100	C7H16	000589-34-4	43



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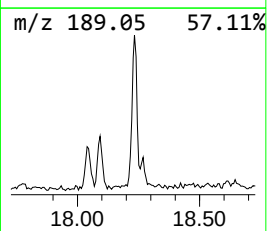
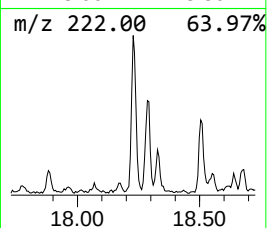
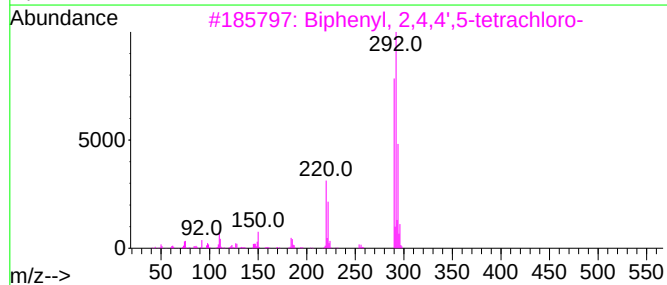
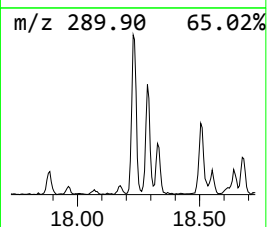
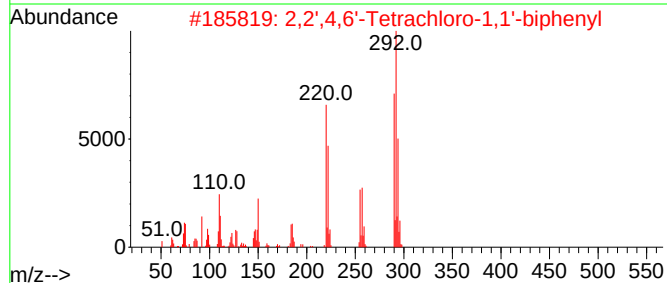
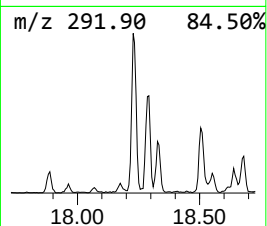
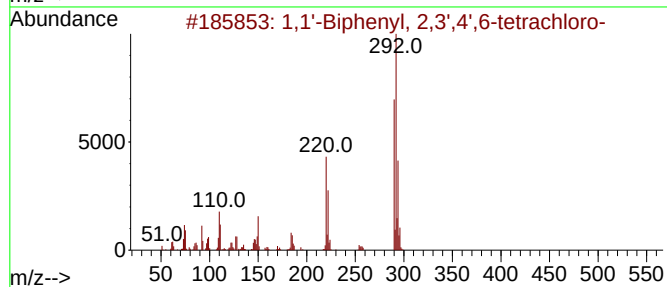
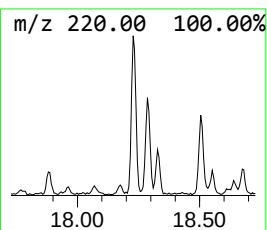
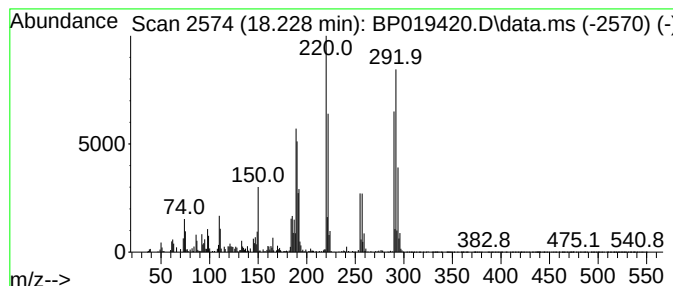
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010824.MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.228	3.27 ng/ul	317587	Phenanthrene-d10	17.169

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,3',4',6-tetrach...	290	C12H6Cl4	041464-46-4	99		
2	2,2',4,6'-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	068194-04-7	99		
3	Biphenyl, 2,4,4',5-tetrachloro-	290	C12H6Cl4	032690-93-0	99		
4	1,1'-Biphenyl, 2,2',4,4'-tetrach...	290	C12H6Cl4	002437-79-8	99		
5	1,1'-Biphenyl, 2,2',4,4'-tetrach...	290	C12H6Cl4	002437-79-8	99		



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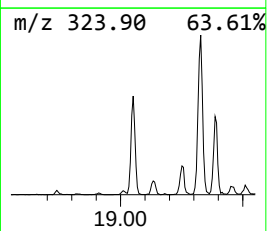
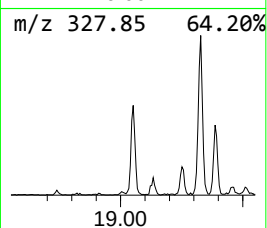
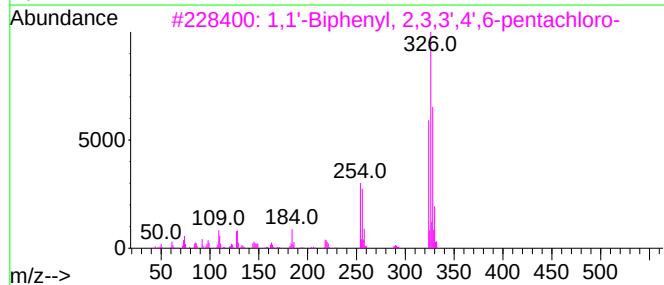
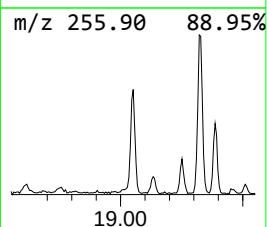
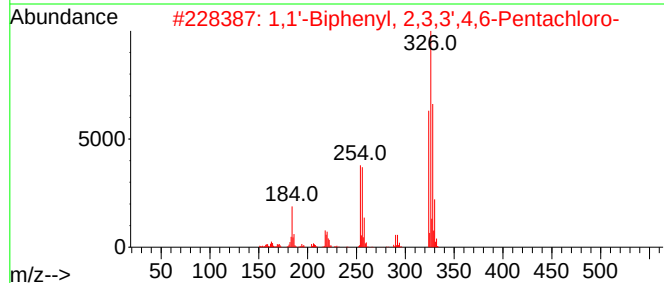
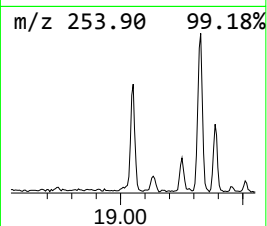
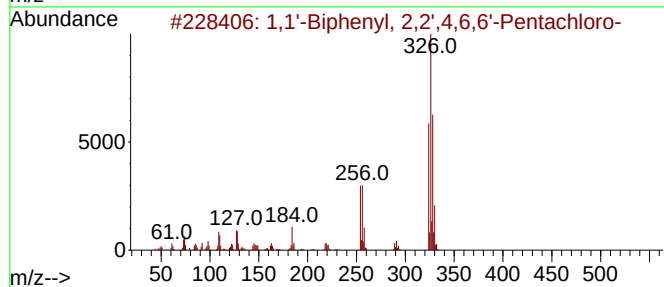
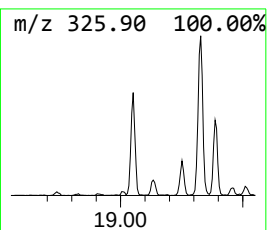
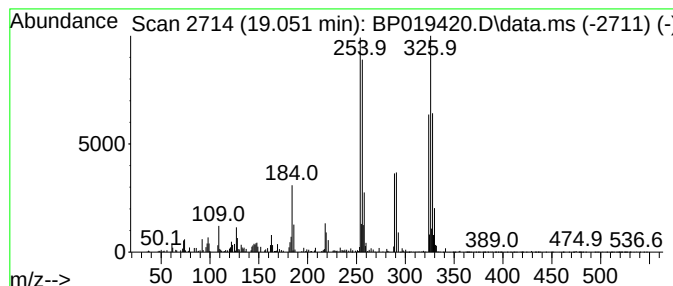
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010824.MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.051	3.33 ng/ul	323809	Phenanthrene-d10	17.169

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',4,6,6'-Penta...	324	C12H5Cl5	056558-16-8	99		
2	1,1'-Biphenyl, 2,3,3',4,6-Pentac...	324	C12H5Cl5	074472-35-8	99		
3	1,1'-Biphenyl, 2,3,3',4',6-penta...	324	C12H5Cl5	038380-03-9	99		
4	2,3,3',5,6-Pentachloro-1,1'-biph...	324	C12H5Cl5	074472-36-9	99		
5	2,3',4,4',5'-Pentachloro-1,1'-bi...	324	C12H5Cl5	065510-44-3	99		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012324\
Data File : BP019420.D
Acq On : 23 Jan 2024 15:02
Operator : MA/JU
Sample : P1157-05
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AB2

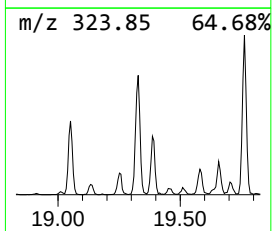
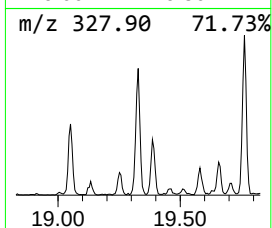
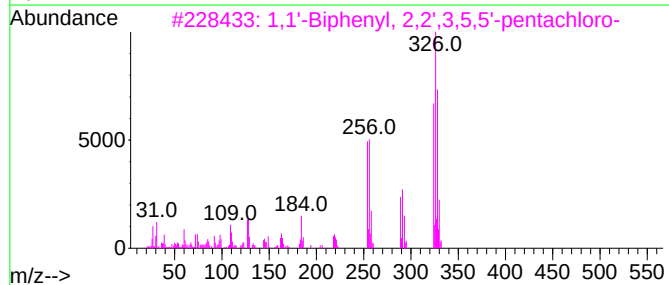
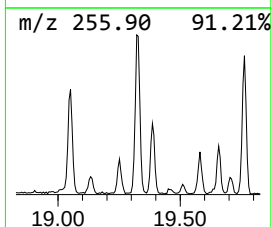
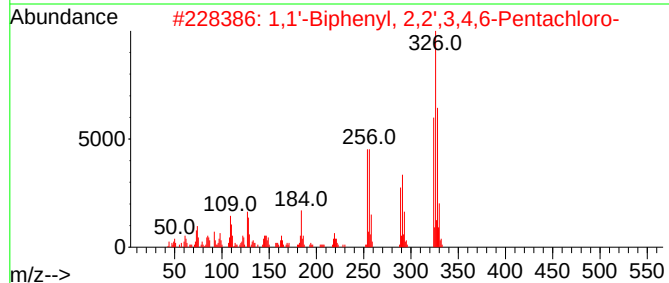
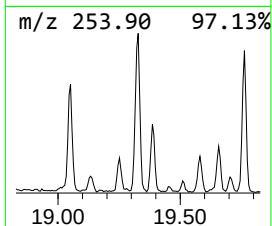
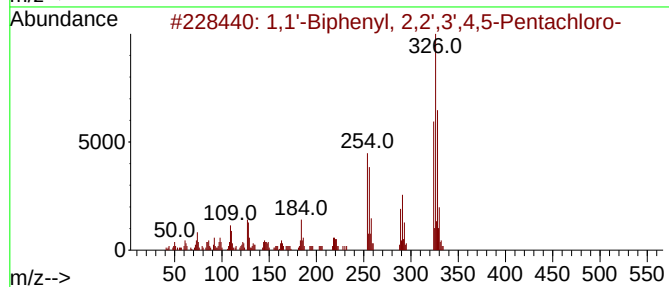
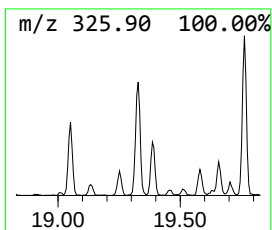
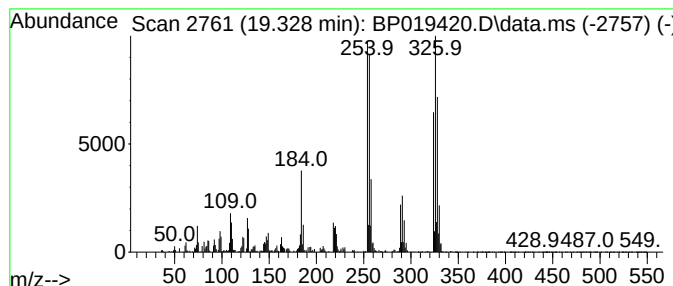
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.328	4.11 ng/ul	380284	Chrysene-d12	21.275

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',3',4,5-Penta...	324	C12H5Cl5	041464-51-1	99		
2	1,1'-Biphenyl, 2,2',3,4,6-Pentac...	324	C12H5Cl5	055215-17-3	99		
3	1,1'-Biphenyl, 2,2',3,5,5'-penta...	324	C12H5Cl5	052663-61-3	99		
4	2,3,3',4',5'- Pentachloro-1,1'-b...	324	C12H5Cl5	076842-07-4	99		
5	1,1'-Biphenyl, 2,2',4,6,6'-Penta...	324	C12H5Cl5	056558-16-8	99		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012324\
Data File : BP019420.D
Acq On : 23 Jan 2024 15:02
Operator : MA/JU
Sample : P1157-05
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AB2

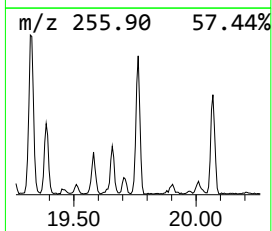
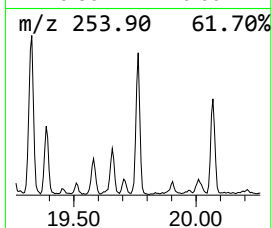
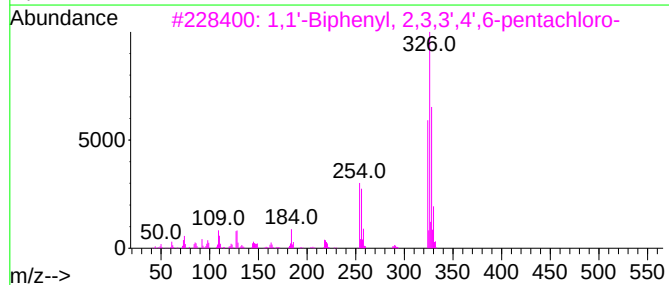
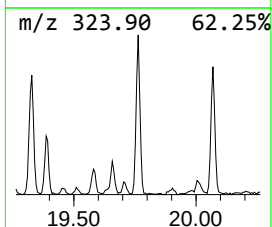
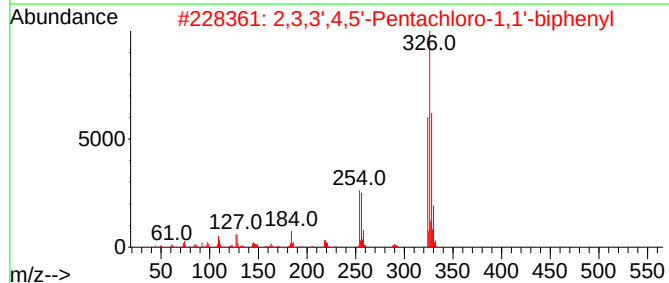
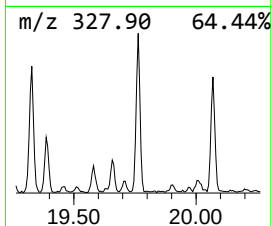
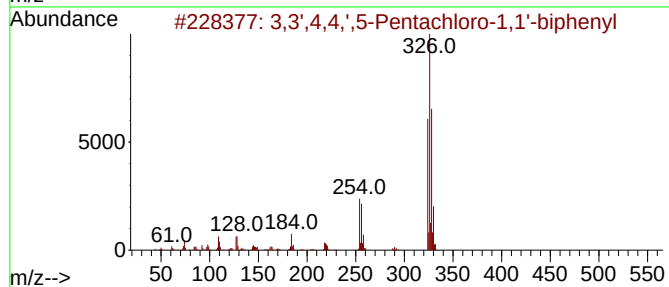
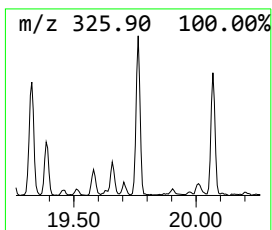
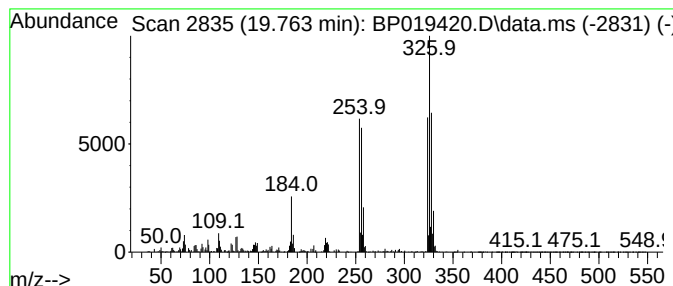
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 3,3',4,4',5-Pentachloro-1,... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.763	3.87 ng/ul	358378	Chrysene-d12	21.275

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,3',4,4',5-Pentachloro-1,1'-bi...	324	C12H5Cl5	057465-28-8	98		
2	2,3,3',4,5'-Pentachloro-1,1'-bip...	324	C12H5Cl5	070362-41-3	98		
3	1,1'-Biphenyl, 2,3,3',4',6-penta...	324	C12H5Cl5	038380-03-9	97		
4	1,1'-Biphenyl, 2,3,3',4,4'-penta...	324	C12H5Cl5	032598-14-4	97		
5	3,3',4,5,5'-Pentachloro-1,1'-bip...	324	C12H5Cl5	039635-33-1	96		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012324\
Data File : BP019420.D
Acq On : 23 Jan 2024 15:02
Operator : MA/JU
Sample : P1157-05
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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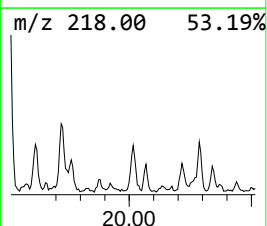
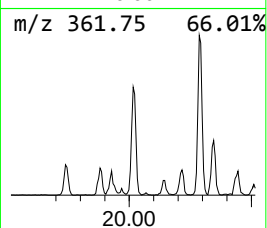
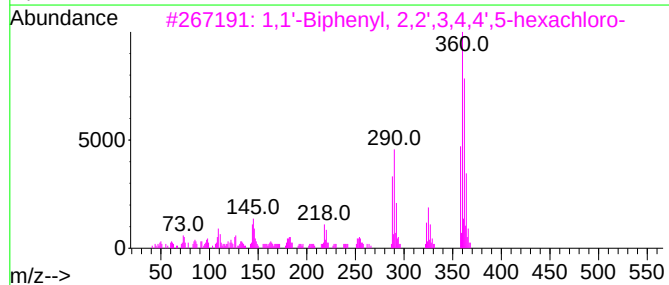
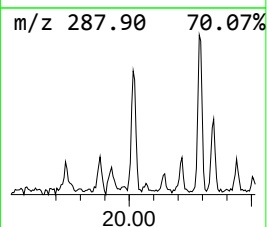
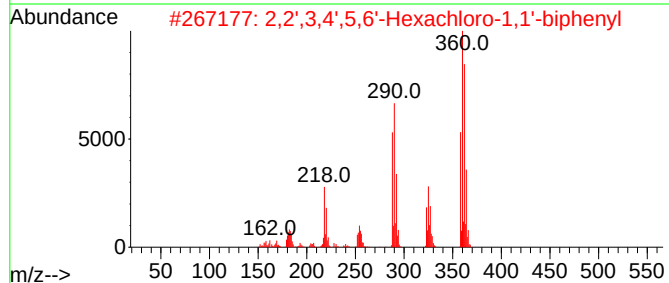
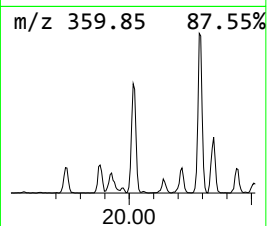
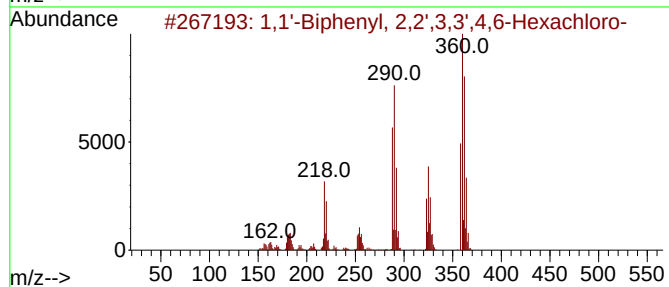
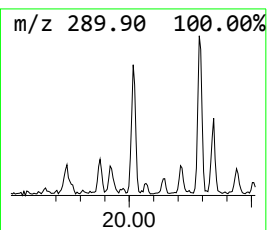
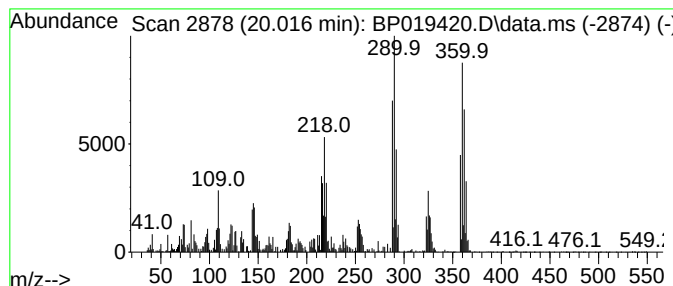
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 1,1'-Biphenyl, 2,2',3,3',4,... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.016	3.16 ng/ul	292911	Chrysene-d12	21.275

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',3,3',4,6-Hex...	358	C12H4Cl6	061798-70-7	99		
2	2,2',3,4',5,6'-Hexachloro-1,1'-b...	358	C12H4Cl6	074472-41-6	99		
3	1,1'-Biphenyl, 2,2',3,4,4',5-hex...	358	C12H4Cl6	035694-06-5	99		
4	2,2',3,4,5,6-Hexachloro-1,1'-bip...	358	C12H4Cl6	041411-61-4	99		
5	1,1'-Biphenyl, 2,2',3,3',4,5'-he...	358	C12H4Cl6	052663-66-8	99		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012324\
Data File : BP019420.D
Acq On : 23 Jan 2024 15:02
Operator : MA/JU
Sample : P1157-05
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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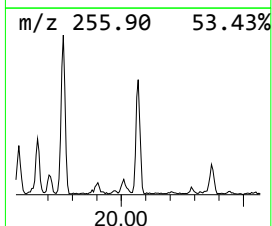
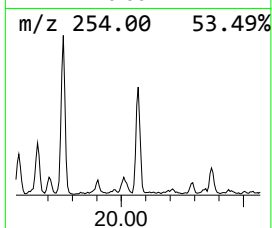
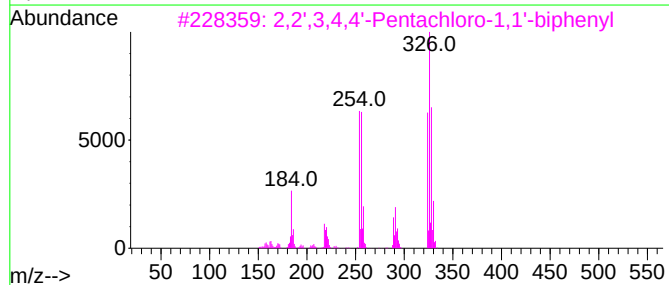
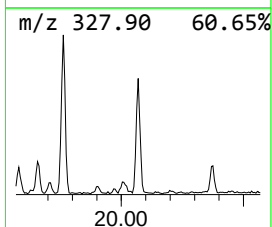
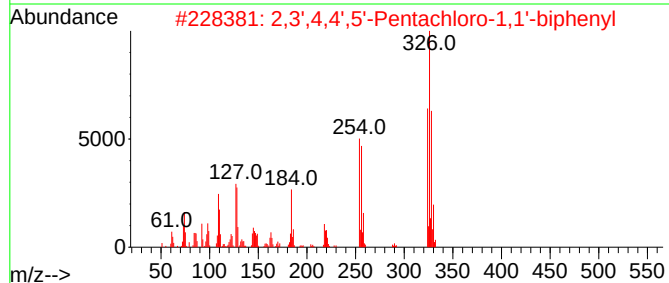
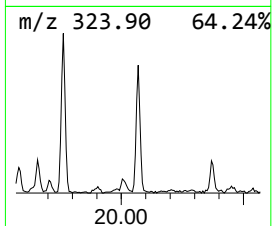
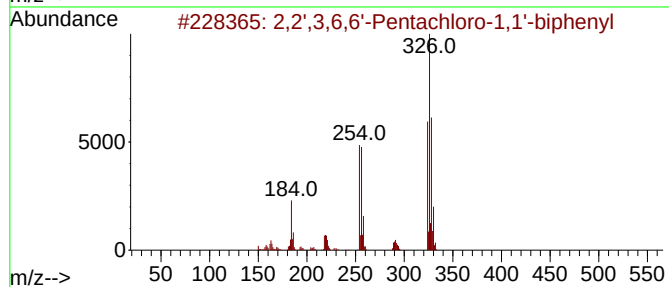
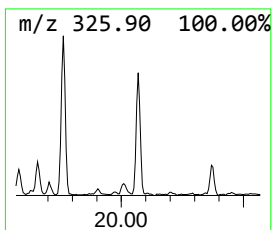
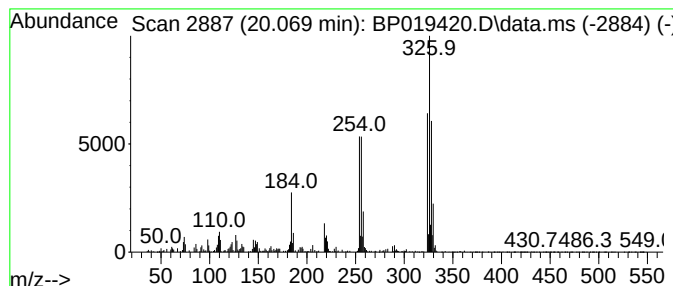
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 2,2',3,6,6'-Pentachloro-1,1... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.069	2.48 ng/ul	229240	Chrysene-d12	21.275

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2',3,6,6'-Pentachloro-1,1'-bip...	324	C12H5Cl5	073575-54-9	97		
2	2,3',4,4',5'-Pentachloro-1,1'-bi...	324	C12H5Cl5	065510-44-3	96		
3	2,2',3,4,4'-Pentachloro-1,1'-bip...	324	C12H5Cl5	065510-45-4	96		
4	2,3,3',4,5'-Pentachloro-1,1'-bip...	324	C12H5Cl5	070362-41-3	96		
5	1,1'-Biphenyl, 2,3,3',4,6-Pentac...	324	C12H5Cl5	074472-35-8	96		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012324\
Data File : BP019420.D
Acq On : 23 Jan 2024 15:02
Operator : MA/JU
Sample : P1157-05
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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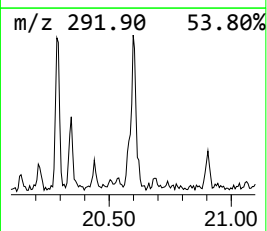
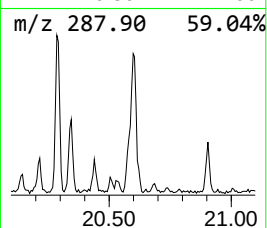
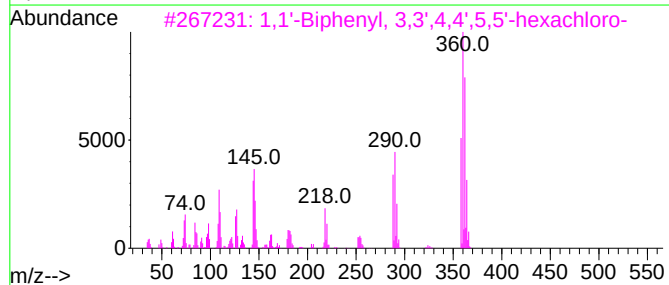
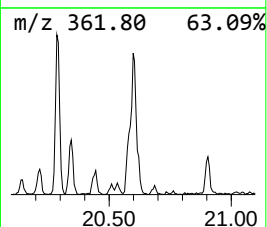
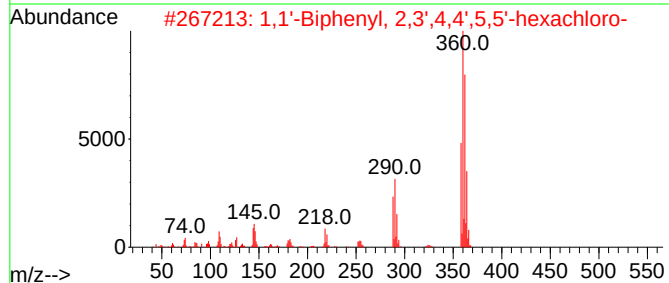
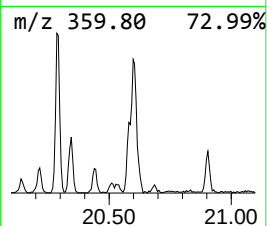
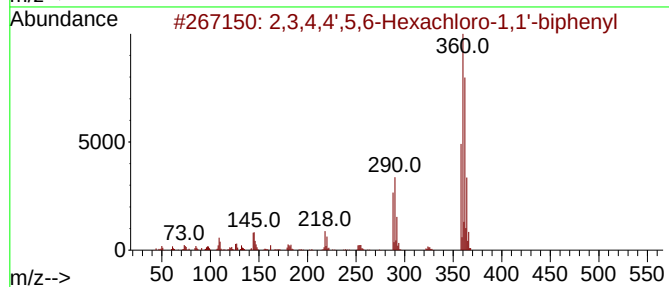
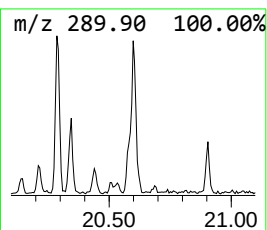
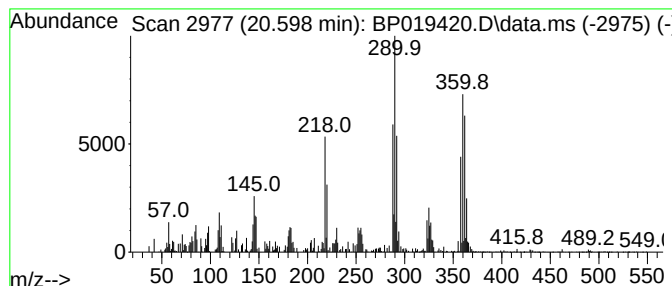
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 2,3,4,4',5,6-Hexachloro-1,1... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.598	2.37 ng/ul	219617	Chrysene-d12	21.275

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	95		
2	1,1'-Biphenyl, 2,3',4,4',5,5'-he...	358	C12H4Cl6	052663-72-6	95		
3	1,1'-Biphenyl, 3,3',4,4',5,5'-he...	358	C12H4Cl6	032774-16-6	94		
4	1,1'-Biphenyl, 2,3,3',4,4',5-hex...	358	C12H4Cl6	038380-08-4	93		
5	2,3,3',4,5',6-Hexachloro-1,1'-bi...	358	C12H4Cl6	074472-43-8	91		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012324\
Data File : BP019420.D
Acq On : 23 Jan 2024 15:02
Operator : MA/JU
Sample : P1157-05
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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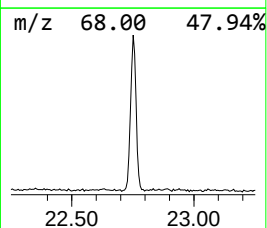
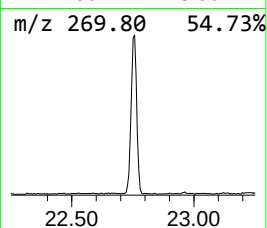
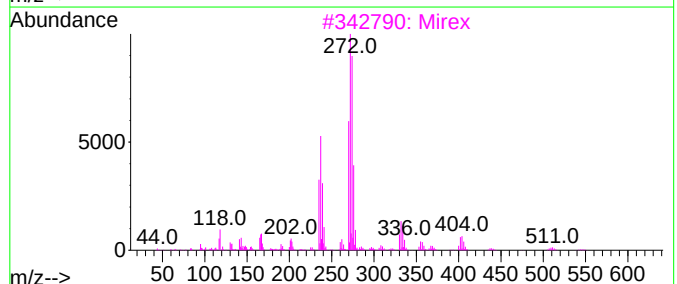
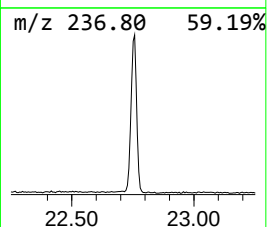
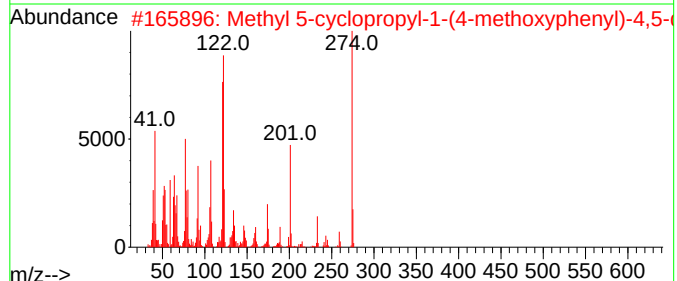
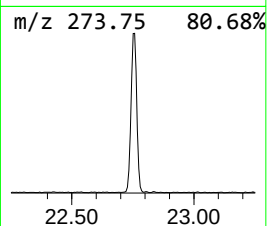
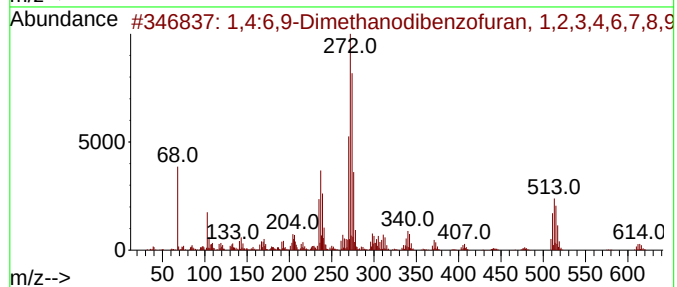
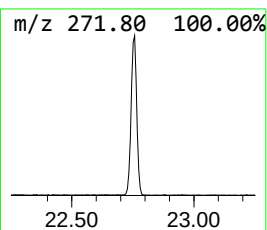
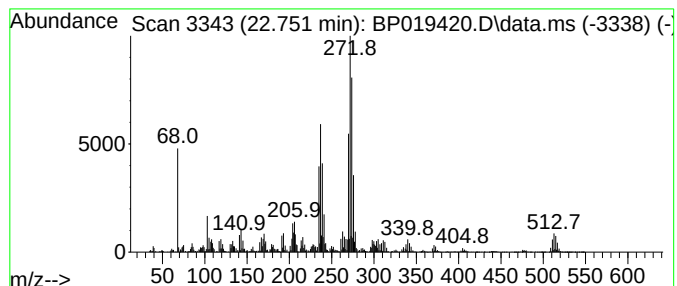
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 1,4:6,9-Dimethanodibenzofur... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.751	17.09 ng/ul	1453520	Perylene-d12	23.622

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,4:6,9-Dimethanodibenzofuran, 1...	608	C14H4Cl120	031107-44-5	93		
2	Methyl 5-cyclopropyl-1-(4-methox...	274	C15H18N2O3	1265897-79-7	90		
3	Mirex	540	C10Cl12	002385-85-5	87		
4	Mirex	540	C10Cl12	002385-85-5	87		
5	Mirex	540	C10Cl12	002385-85-5	78		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012324\
Data File : BP019420.D
Acq On : 23 Jan 2024 15:02
Operator : MA/JU
Sample : P1157-05
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AB2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010824.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
(DEL) Alkane: S...	3.129	22.6	ng/ul	955548	1	7.769	847617	20.0
1,1'-Biphenyl, ...	18.228	3.3	ng/ul	317587	4	17.169	1944580	20.0
1,1'-Biphenyl, ...	19.051	3.3	ng/ul	323809	4	17.169	1944580	20.0
1,1'-Biphenyl, ...	19.328	4.1	ng/ul	380284	5	21.275	1852230	20.0
3,3',4,4',5-Pe...	19.763	3.9	ng/ul	358378	5	21.275	1852230	20.0
1,1'-Biphenyl, ...	20.016	3.2	ng/ul	292911	5	21.275	1852230	20.0
2,2',3,6,6'-Pen...	20.069	2.5	ng/ul	229240	5	21.275	1852230	20.0
2,3,4,4',5,6-He...	20.598	2.4	ng/ul	219617	5	21.275	1852230	20.0
1,4:6,9-Dimetha...	22.751	17.1	ng/ul	1453520	6	23.622	1701520	20.0