

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100924\
 Data File : BP022295.D
 Acq On : 09 Oct 2024 18:24
 Operator : RC/JU
 Sample : P4162-16
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A4EJ1

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-BP100724.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP022295.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.228	37	41	45	rVB	20824	25834	0.65%	0.058%
2	3.640	107	111	121	rBV	240214	361653	9.07%	0.814%
3	3.964	162	166	170	rVB	13338	17682	0.44%	0.040%
4	4.869	315	320	325	rBV3	6630	12596	0.32%	0.028%
5	5.828	481	483	488	rVB5	2923	4523	0.11%	0.010%
6	5.999	508	512	516	rVB6	4208	6467	0.16%	0.015%
7	6.752	632	640	644	rBV10	4681	8897	0.22%	0.020%
8	6.893	653	664	678	rBV	411602	720919	18.08%	1.623%
9	7.046	683	690	702	rBV	293165	484420	12.15%	1.091%
10	7.252	718	725	736	rBV	426013	690435	17.32%	1.554%
11	7.705	796	802	810	rBV	488312	807861	20.26%	1.819%
12	7.775	810	814	819	rVB7	5079	7766	0.19%	0.017%
13	8.410	914	922	929	rBV	549092	866046	21.72%	1.950%
14	8.846	988	996	1005	rBV	402570	678909	17.03%	1.528%
15	8.999	1020	1022	1028	rVB7	2433	4367	0.11%	0.010%
16	9.263	1060	1067	1072	rBV2	12403	21054	0.53%	0.047%
17	9.405	1084	1091	1098	rBV2	28824	51672	1.30%	0.116%
18	9.563	1109	1118	1131	rBV	360014	626708	15.72%	1.411%
19	9.810	1151	1160	1164	rBV	3412	7401	0.19%	0.017%
20	9.969	1181	1187	1190	rBV7	2236	4132	0.10%	0.009%
21	10.099	1200	1209	1223	rBV	592176	1055731	26.48%	2.377%
22	10.469	1262	1272	1286	rBV	649222	1186758	29.77%	2.672%
23	10.604	1286	1295	1309	rBV	471816	794851	19.94%	1.789%
24	10.999	1354	1362	1370	rBV8	12368	23593	0.59%	0.053%
25	11.269	1401	1408	1412	rBV8	4892	10746	0.27%	0.024%
26	11.487	1439	1445	1450	rVB7	3760	6131	0.15%	0.014%
27	11.740	1482	1488	1496	rBV6	7030	13948	0.35%	0.031%
28	11.987	1521	1530	1534	rBV7	8264	21101	0.53%	0.048%
29	12.051	1534	1541	1549	rVB3	11209	23023	0.58%	0.052%
30	12.122	1549	1553	1556	rBV6	4014	5782	0.15%	0.013%
31	12.346	1586	1591	1597	rVB9	4071	5944	0.15%	0.013%
32	12.598	1630	1634	1638	rVB7	2591	4288	0.11%	0.010%
33	12.840	1671	1675	1683	rVB5	6415	10463	0.26%	0.024%
34	13.128	1721	1724	1727	rBV5	3826	5603	0.14%	0.013%
35	13.298	1746	1753	1761	rBV3	17122	36868	0.92%	0.083%
36	13.457	1773	1780	1792	rBV	23905	72203	1.81%	0.163%

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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-BP100724.MA.M
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37	13.634	1805	1810	1815	rBV7	7255	13461	0.34%	0.030%
38	13.722	1820	1825	1832	rVV	1262782	1659531	41.62%	3.736%
39	13.798	1836	1838	1842	rVB4	8812	10322	0.26%	0.023%
40	14.010	1863	1874	1887	rBV	1270905	2009397	50.40%	4.523%
41	14.157	1895	1899	1905	rVB8	5287	10170	0.26%	0.023%
42	14.216	1905	1909	1915	rVB9	4884	8375	0.21%	0.019%
43	14.322	1918	1927	1934	rBV	1030242	1796140	45.05%	4.043%
44	14.392	1934	1939	1944	rBV2	15873	26498	0.66%	0.060%
45	14.445	1946	1948	1953	rVB3	11146	14281	0.36%	0.032%
46	14.540	1957	1964	1978	rBV	363568	777950	19.51%	1.751%
47	14.734	1993	1997	1998	rBV2	9575	11900	0.30%	0.027%
48	14.987	2038	2040	2045	rVB5	7166	9571	0.24%	0.022%
49	15.128	2060	2064	2070	rBV9	9832	19435	0.49%	0.044%
50	15.316	2090	2096	2103	rBV	1830007	2635918	66.11%	5.934%
51	15.381	2103	2107	2113	rVB2	16786	26145	0.66%	0.059%
52	15.457	2115	2120	2128	rBV	416598	539180	13.52%	1.214%
53	15.581	2137	2141	2153	rVB2	73969	129116	3.24%	0.291%
54	15.704	2155	2162	2169	rVB	284494	459622	11.53%	1.035%
55	16.210	2244	2248	2253	rBV	28421	43891	1.10%	0.099%
56	16.322	2263	2267	2273	rVB	70009	100650	2.52%	0.227%
57	16.634	2315	2320	2327	rVB	196286	279227	7.00%	0.629%
58	16.986	2376	2380	2383	rBV2	64366	88641	2.22%	0.200%
59	17.022	2383	2386	2395	rVB	129544	193165	4.84%	0.435%
60	17.116	2395	2402	2406	rBV	1701205	2445791	61.35%	5.506%
61	17.157	2406	2409	2413	rVV	340334	469154	11.77%	1.056%
62	17.216	2413	2419	2428	rVV2	1883221	3198219	80.22%	7.200%
63	17.292	2428	2432	2446	rVB	340396	523249	13.12%	1.178%
64	17.586	2478	2482	2484	rBV	102769	129948	3.26%	0.293%
65	17.610	2484	2486	2495	rVB2	71916	119862	3.01%	0.270%
66	17.739	2500	2508	2517	rVB	224828	474097	11.89%	1.067%
67	17.863	2524	2529	2534	rBV4	88041	147007	3.69%	0.331%
68	17.939	2537	2542	2545	rBV2	61642	86778	2.18%	0.195%
69	17.992	2547	2551	2553	rBV	50105	80220	2.01%	0.181%
70	18.057	2557	2562	2572	rVB3	542826	873385	21.91%	1.966%
71	18.228	2586	2591	2595	rBV2	244963	373410	9.37%	0.

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Integrator: RTE
Smoothing : OFF Filtering: 5
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Title : SVOA CALIBRATION

77	18.692	2667	2670	2674	rVB	129570	173904	4.36%	0.391%
78	19.028	2724	2727	2728	rBV2	58926	60002	1.50%	0.135%
79	19.251	2760	2765	2772	rVB	417866	640690	16.07%	1.442%
80	19.386	2784	2788	2794	rVB5	167813	280535	7.04%	0.632%
81	19.592	2817	2823	2828	rBV	2595196	3986910	100.00%	8.975%
82	19.798	2856	2858	2861	rBV3	75627	86296	2.16%	0.194%
83	20.686	3003	3009	3010	rBV5	196020	310804	7.80%	0.700%
84	21.210	3095	3098	3105	rVB2	449184	680113	17.06%	1.531%
85	21.551	3150	3156	3161	rBV2	1519696	2625788	65.86%	5.911%
86	24.621	3671	3678	3687	rBV2	1347491	3261271	81.80%	7.342%
87	24.857	3710	3718	3728	rVB	998991	2922861	73.31%	6.580%

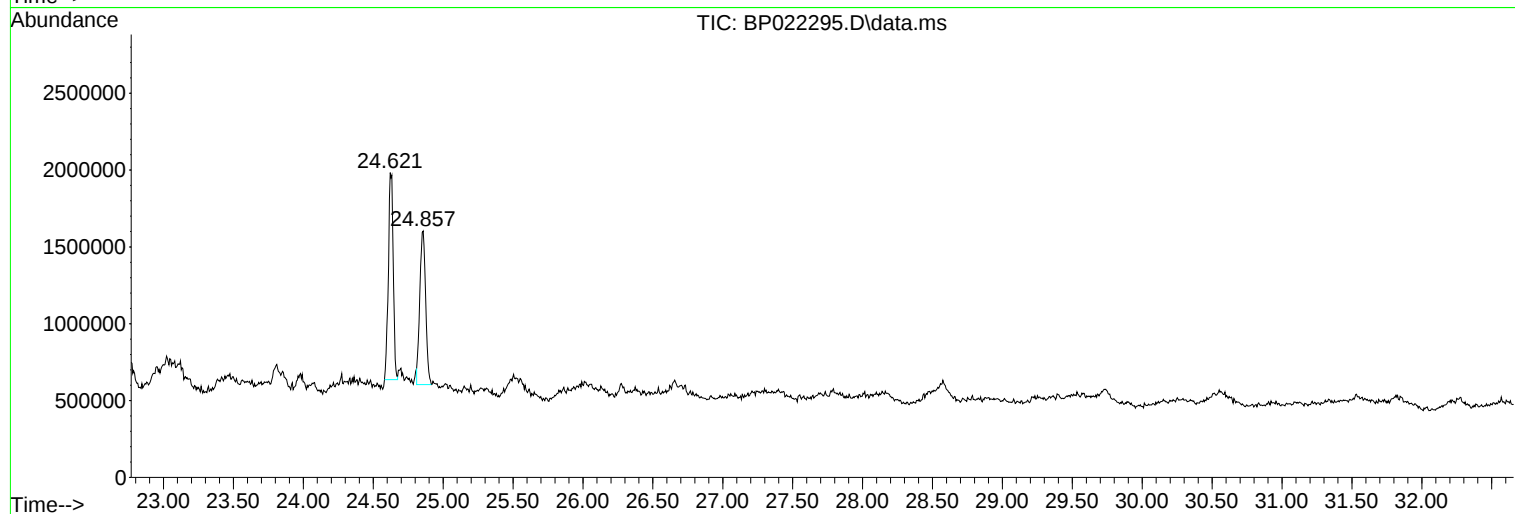
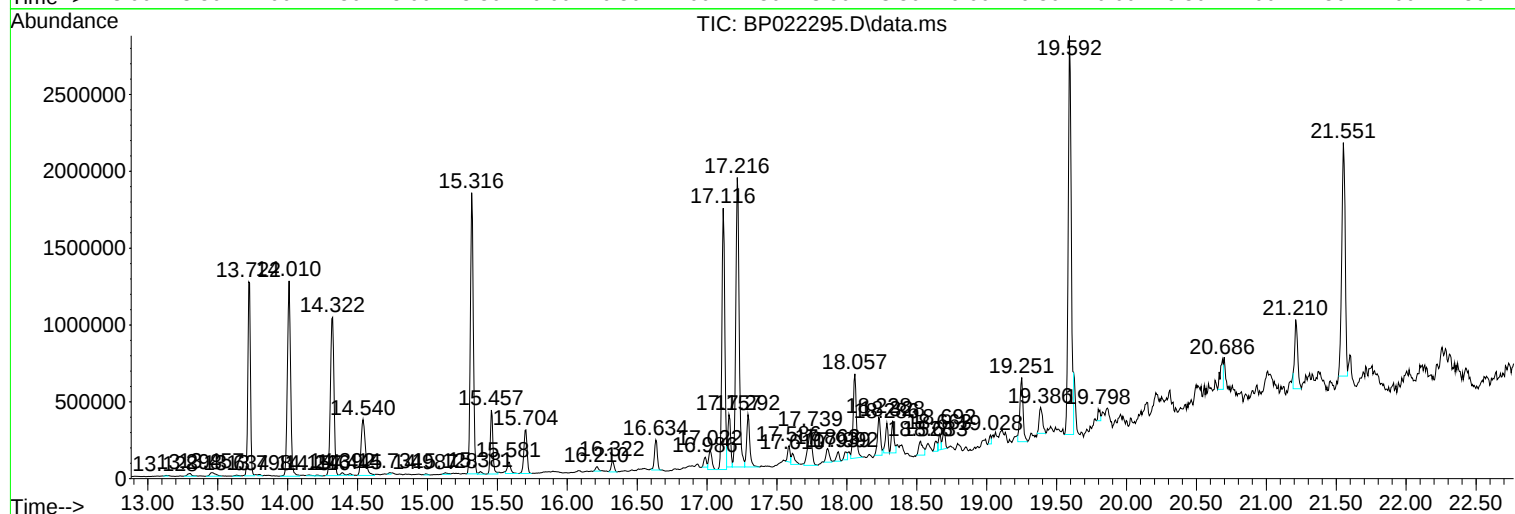
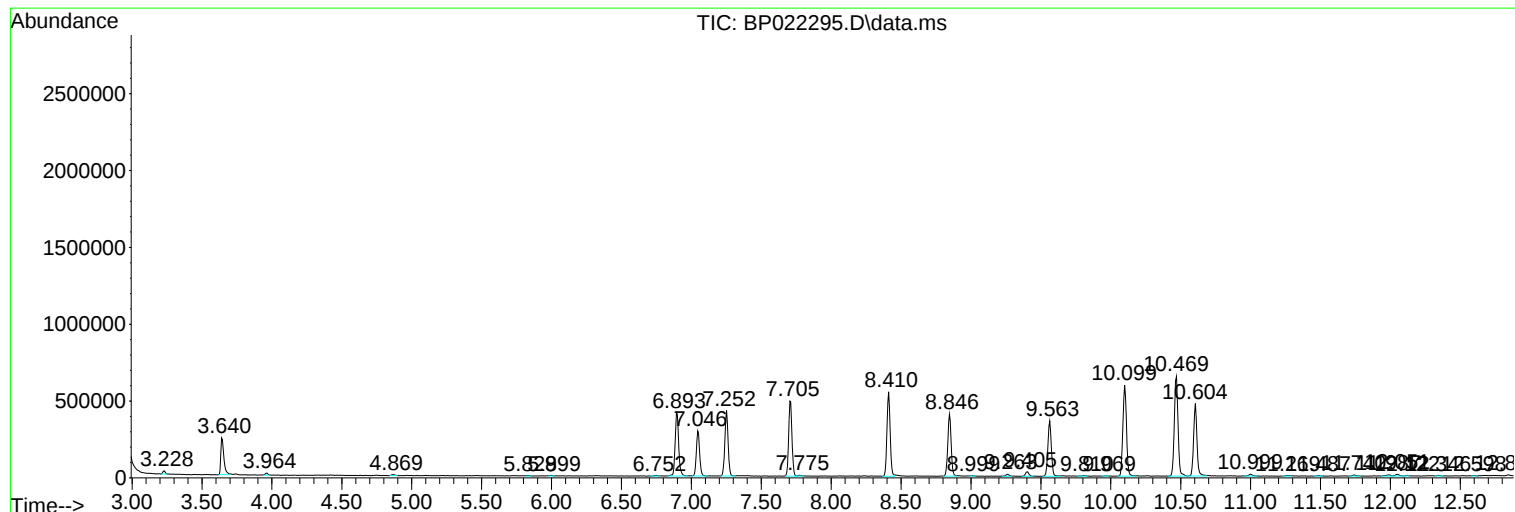
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100924\
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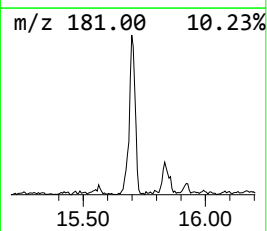
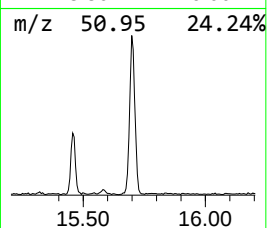
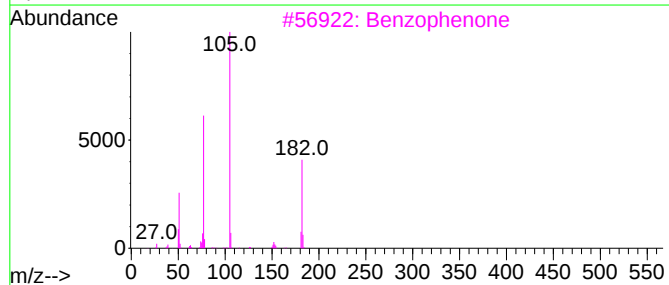
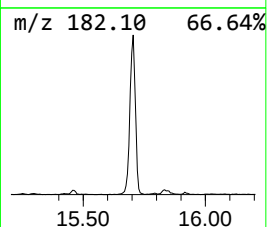
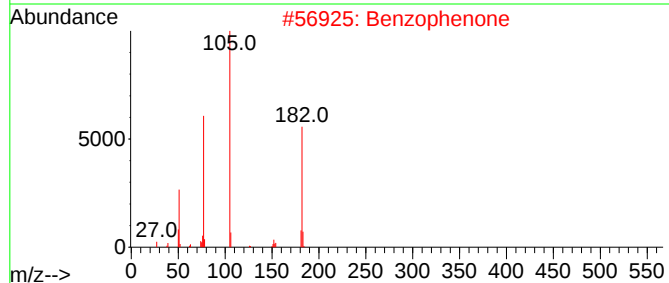
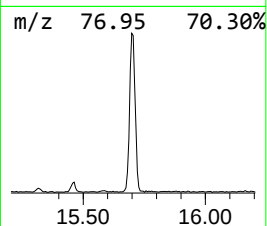
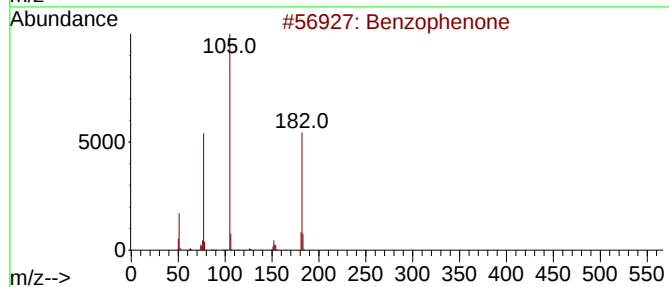
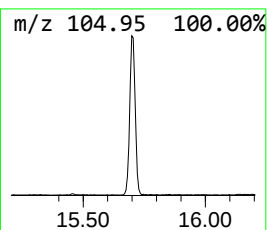
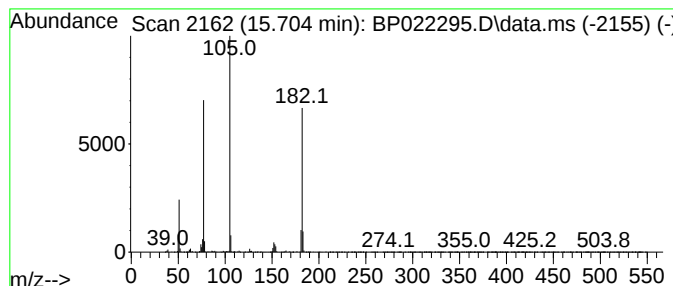
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzophenone Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.704	5.12 ng/ul	459622	Acenaphthene-d10	14.316

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			Benzophenone	182	C13H10O	000119-61-9	96
3			Benzophenone	182	C13H10O	000119-61-9	95
4			Benzophenone	182	C13H10O	000119-61-9	91
5			Benzophenone	182	C13H10O	000119-61-9	81



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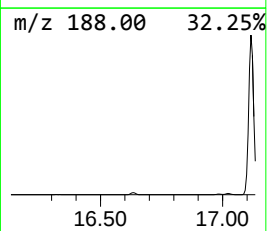
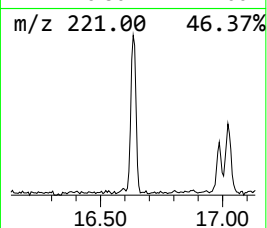
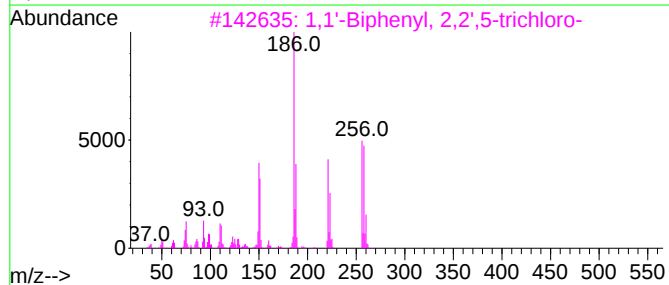
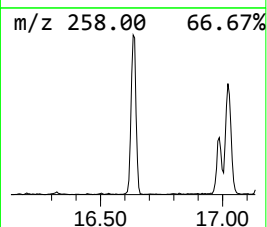
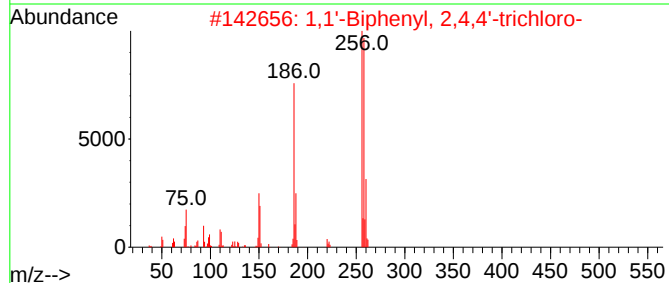
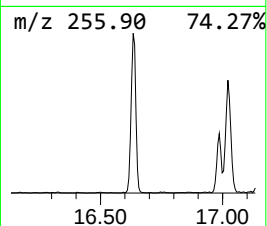
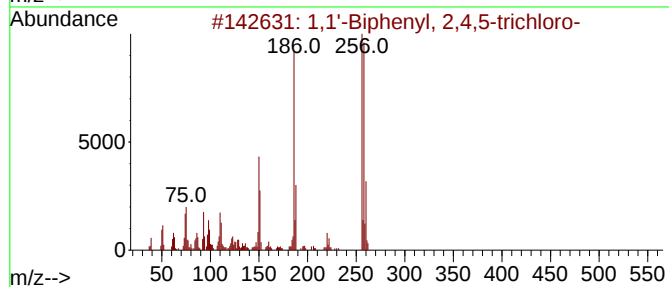
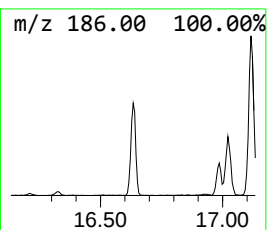
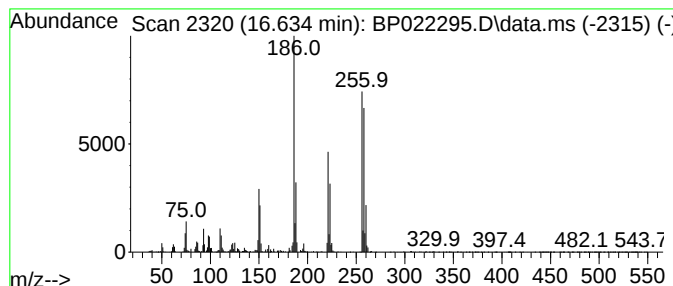
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 1,1'-Biphenyl, 2,4,5-trichl... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.634	2.28 ng/ul	279227	Phenanthrene-d10	17.116

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,4,5-trichloro-	256	C12H7Cl3	015862-07-4	99		
2	1,1'-Biphenyl, 2,4,4'-trichloro-	256	C12H7Cl3	007012-37-5	99		
3	1,1'-Biphenyl, 2,2',5-trichloro-	256	C12H7Cl3	037680-65-2	99		
4	2,2',3-Trichloro-1,1'-biphenyl	256	C12H7Cl3	038444-78-9	99		
5	1,1'-Biphenyl, 2,3,3'-trichloro-	256	C12H7Cl3	038444-84-7	99		



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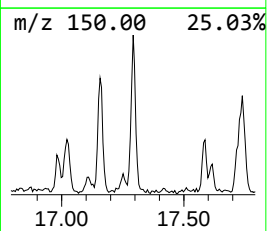
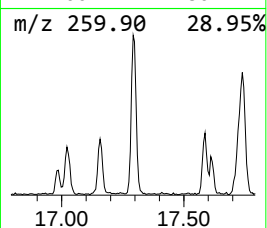
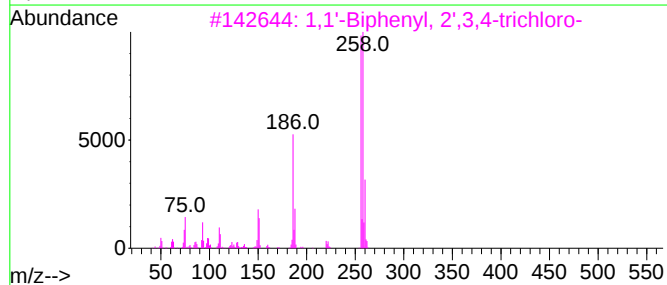
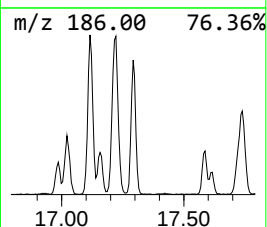
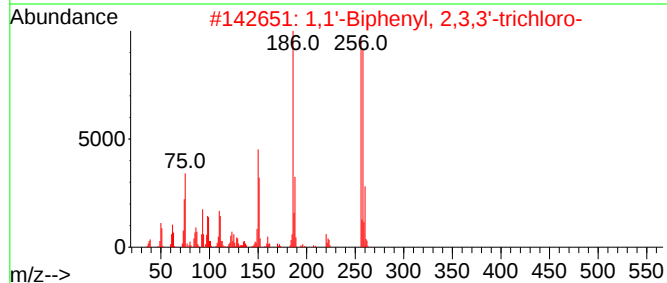
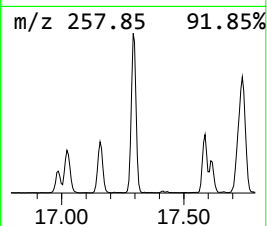
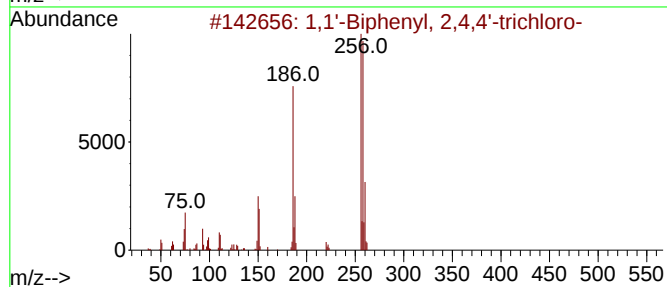
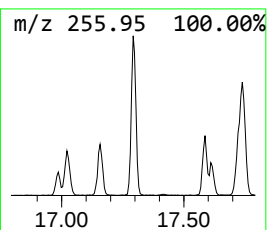
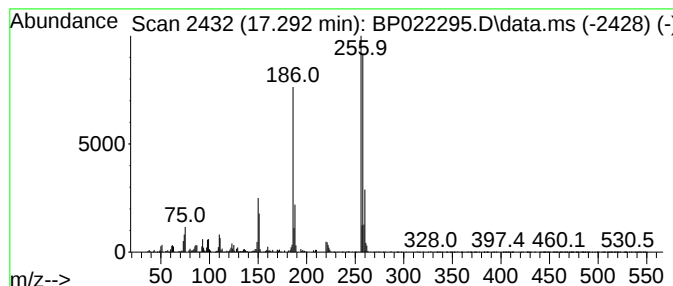
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.292	4.28 ng/ul	523249	Phenanthrene-d10	17.116

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, 2,3,3'-trichloro-	256	C12H7Cl3	038444-84-7	99		
3	1,1'-Biphenyl, 2',3,4-trichloro-	256	C12H7Cl3	038444-86-9	99		
4	1,1'-Biphenyl, 2,3',5-trichloro-	256	C12H7Cl3	038444-81-4	99		
5	1,1'-Biphenyl, 2,4,4'-trichloro-	256	C12H7Cl3	007012-37-5	99		



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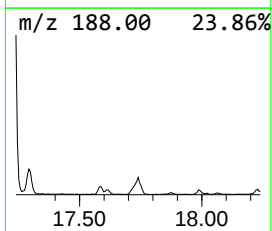
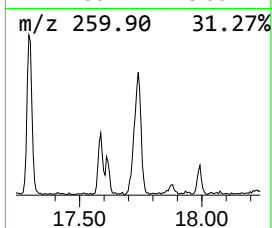
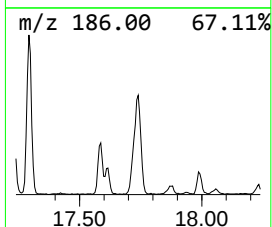
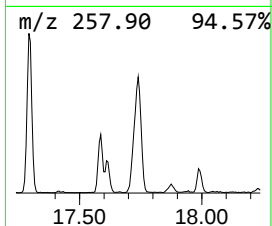
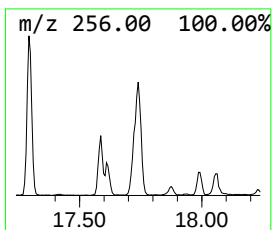
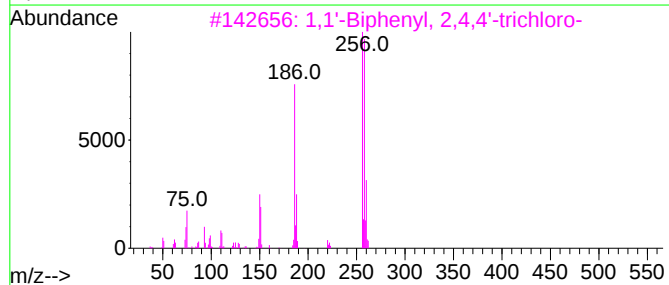
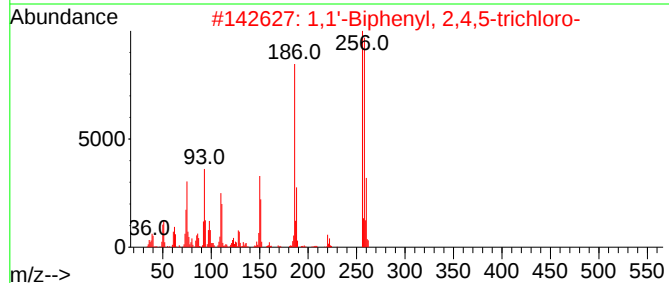
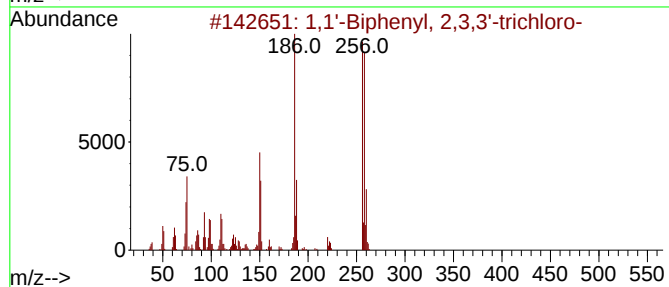
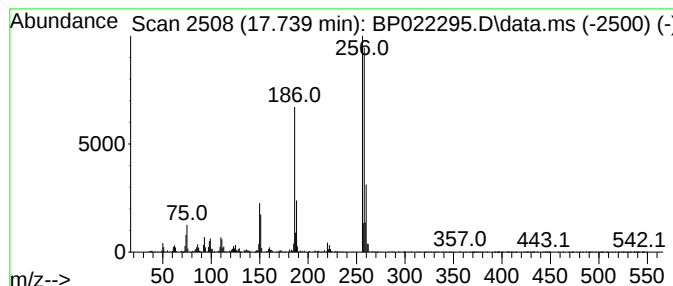
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BNA_P
ClientSampleId :
A4EJ1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.739	3.88 ng/ul	474097	Phenanthrene-d10	17.116	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,3,3'-trichloro-	256	C12H7Cl3	038444-84-7	99
2	1,1'-Biphenyl, 2,4,5-trichloro-	256	C12H7Cl3	015862-07-4	98
3	1,1'-Biphenyl, 2,4,4'-trichloro-	256	C12H7Cl3	007012-37-5	97
4	1,1'-Biphenyl, 2',3,4-trichloro-	256	C12H7Cl3	038444-86-9	96
5	1,1'-Biphenyl, 2,3',5-trichloro-	256	C12H7Cl3	038444-81-4	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100924\
Data File : BP022295.D
Acq On : 09 Oct 2024 18:24
Operator : RC/JU
Sample : P4162-16
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4EJ1

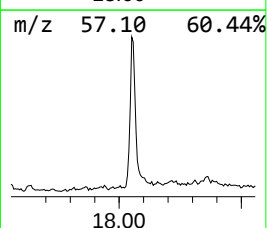
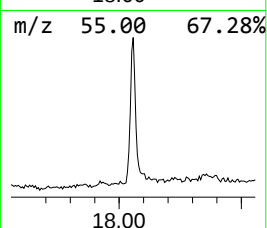
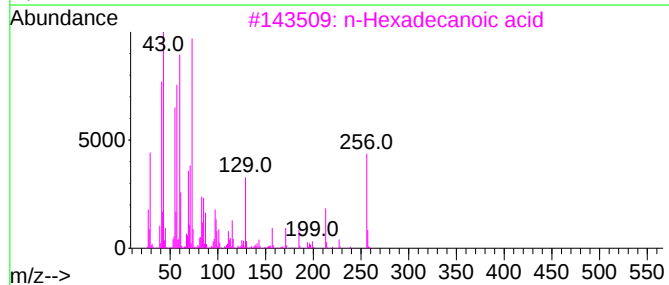
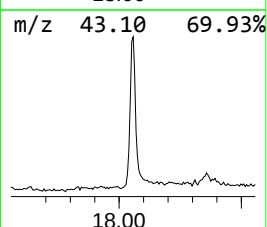
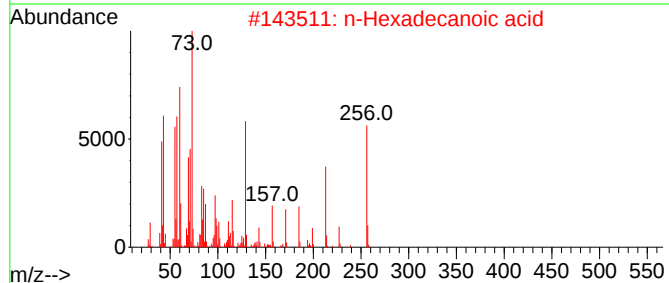
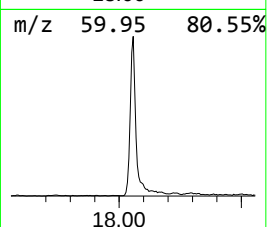
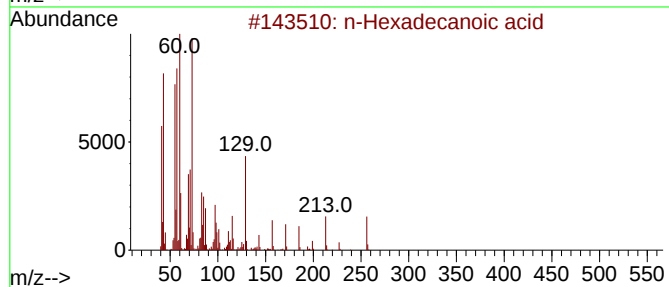
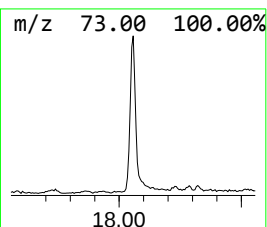
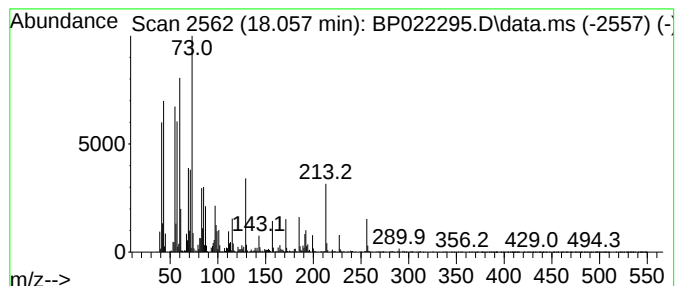
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 n-Hexadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.057	7.14 ng/ul	873385	Phenanthrene-d10	17.116

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
5			Tridecanoic acid	214	C13H26O2	000638-53-9	92



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100924\
Data File : BP022295.D
Acq On : 09 Oct 2024 18:24
Operator : RC/JU
Sample : P4162-16
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4EJ1

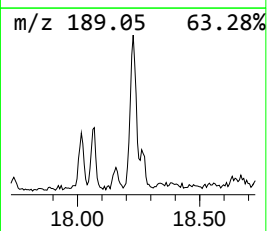
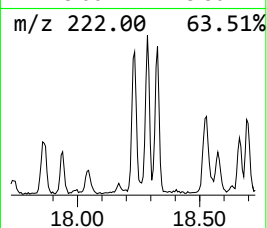
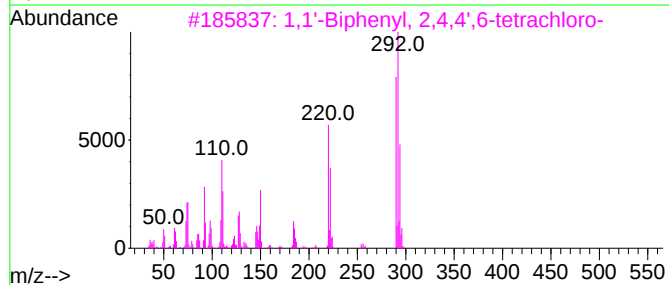
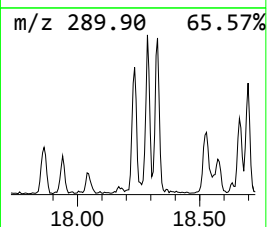
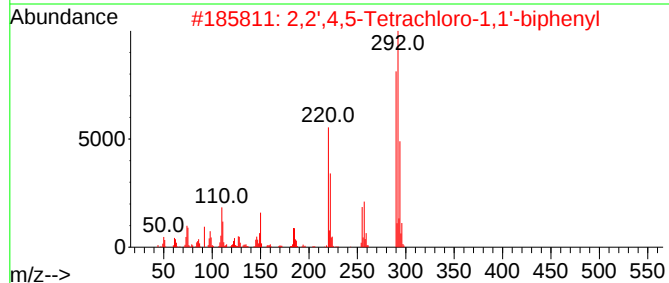
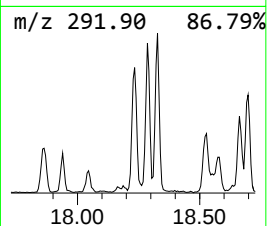
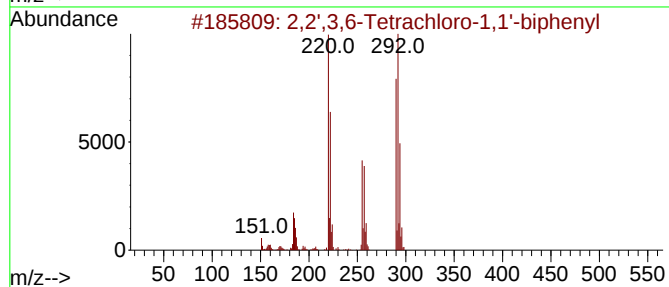
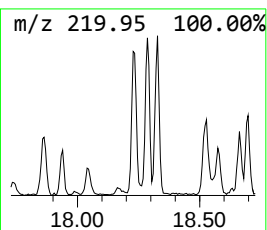
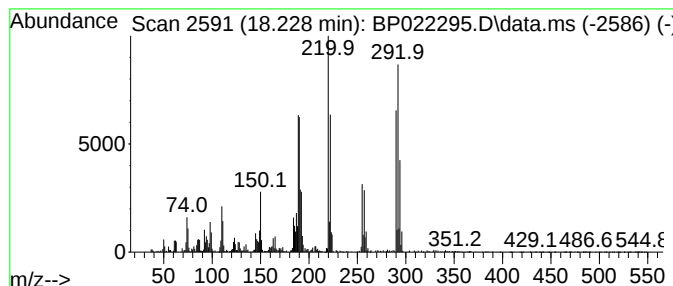
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.228	3.05 ng/ul	373410	Phenanthrene-d10	17.116

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2',3,6-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-45-7	95		
2	2,2',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-47-9	95		
3	1,1'-Biphenyl, 2,4,4',6-tetrachl...	290	C12H6Cl4	032598-12-2	95		
4	2,3',4,5'-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	073575-52-7	95		
5	1,1'-Biphenyl, 2,2',3,5-Tetrachl...	290	C12H6Cl4	070362-46-8	95		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100924\
Data File : BP022295.D
Acq On : 09 Oct 2024 18:24
Operator : RC/JU
Sample : P4162-16
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4EJ1

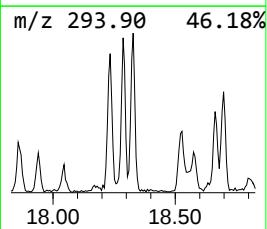
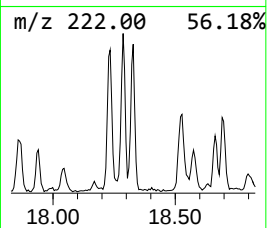
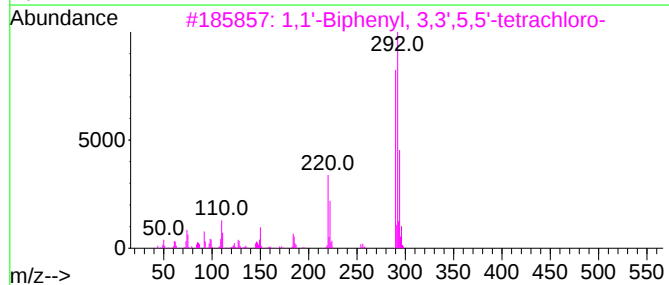
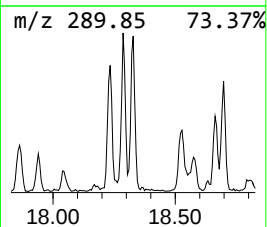
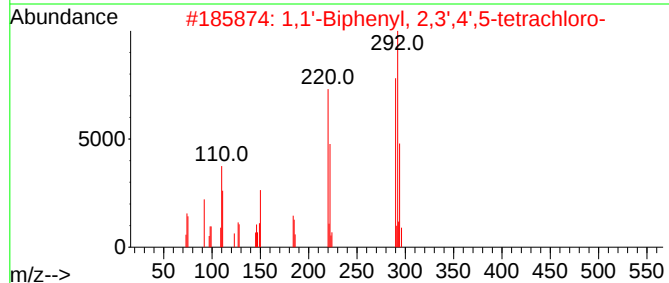
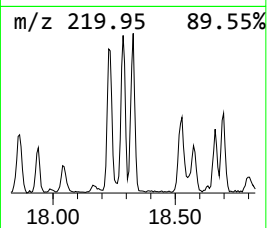
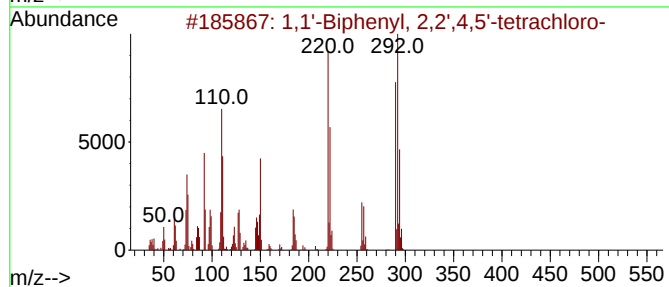
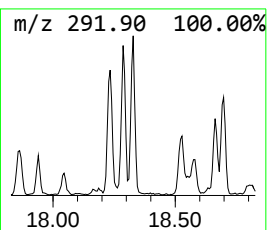
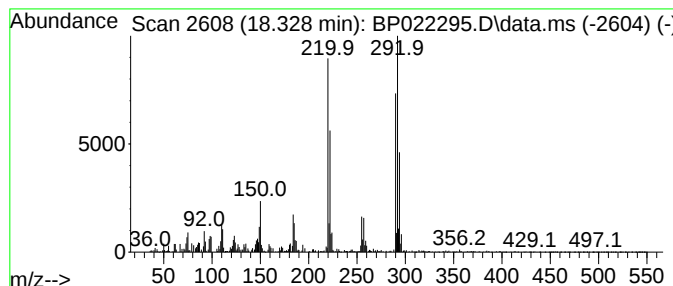
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 1,1'-Biphenyl, 2,2',4,5'-te... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.328	2.22 ng/ul	271865	Phenanthrene-d10	17.116

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',4,5'-tetrach...	290	C12H6Cl4	041464-40-8	97		
2	1,1'-Biphenyl, 2,3',4',5-tetrach...	290	C12H6Cl4	032598-11-1	96		
3	1,1'-Biphenyl, 3,3',5,5'-tetrach...	290	C12H6Cl4	033284-52-5	96		
4	1,1'-Biphenyl, 2,3,3',5'-tetrach...	290	C12H6Cl4	041464-49-7	96		
5	1,1'-Biphenyl, 2,3,4,4'-tetrachl...	290	C12H6Cl4	033025-41-1	96		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100924\
Data File : BP022295.D
Acq On : 09 Oct 2024 18:24
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Sample : P4162-16
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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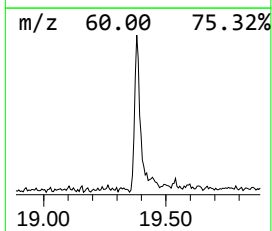
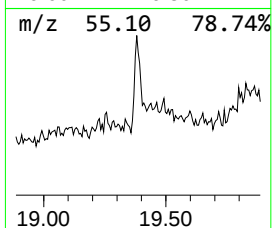
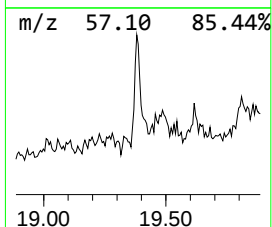
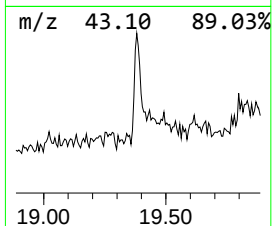
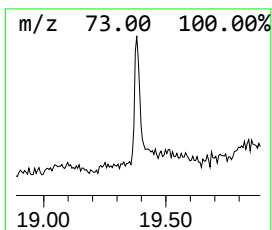
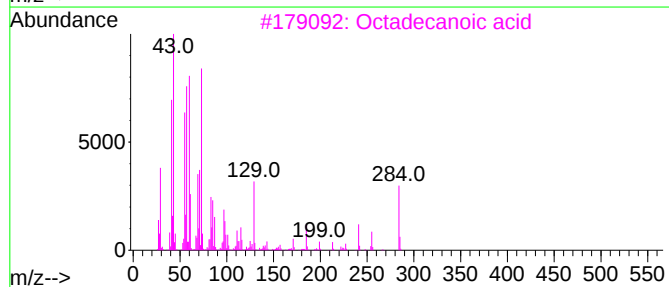
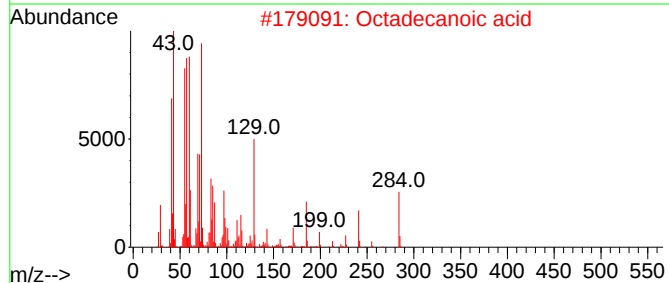
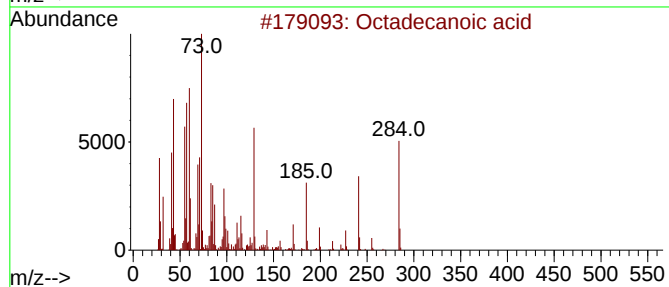
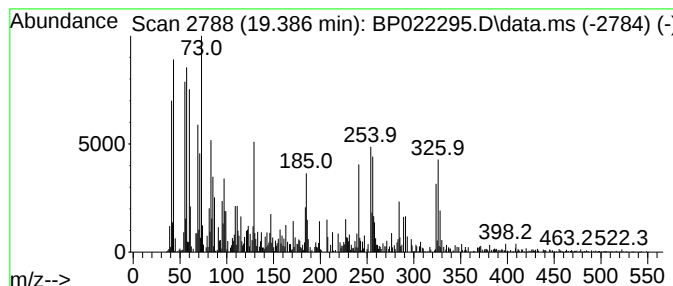
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Octadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.386	2.14 ng/ul	280535	Chrysene-d12	21.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	95
2			Octadecanoic acid	284	C18H36O2	000057-11-4	94
3			Octadecanoic acid	284	C18H36O2	000057-11-4	86
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	66
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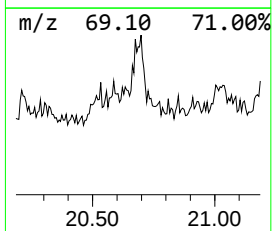
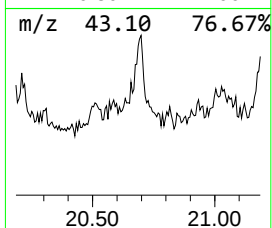
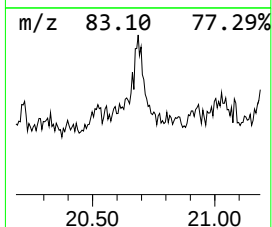
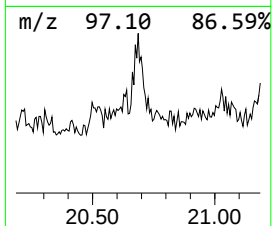
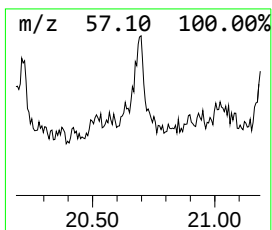
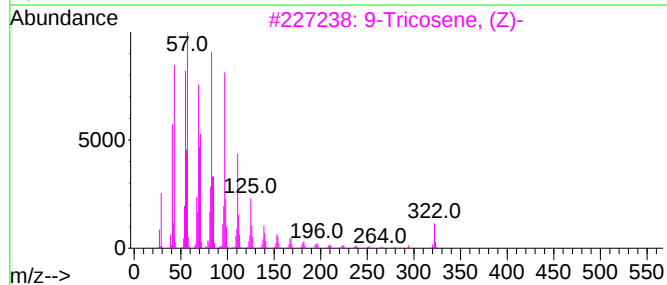
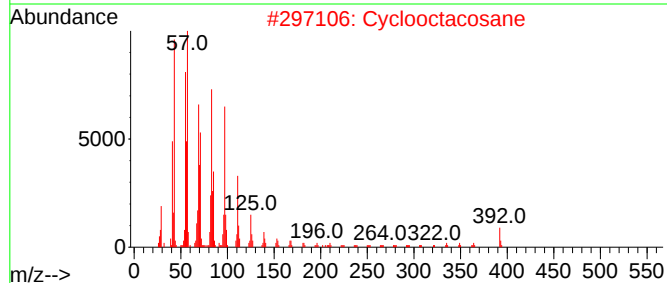
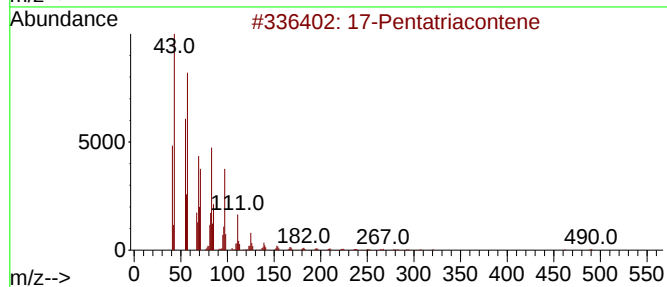
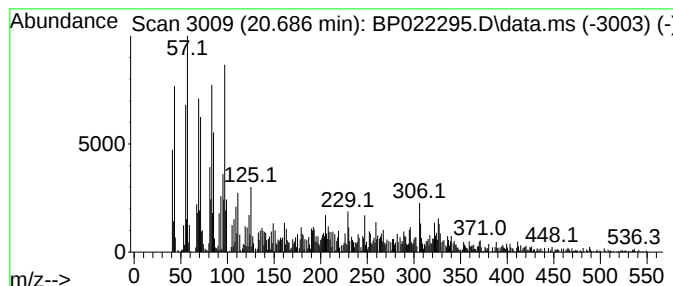
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Data File : BP022295.D
Acq On : 09 Oct 2024 18:24
Operator : RC/JU
Sample : P4162-16
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4EJ1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 7



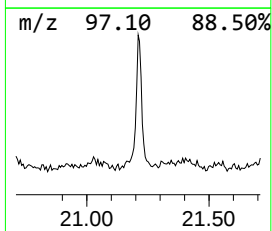
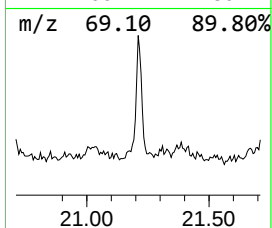
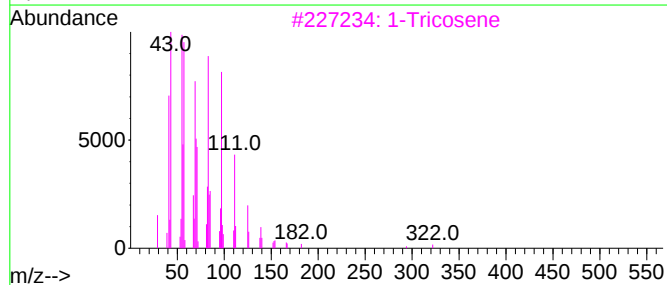
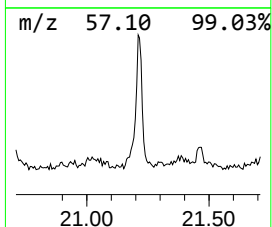
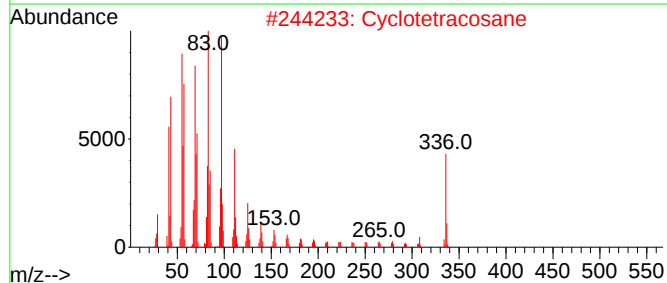
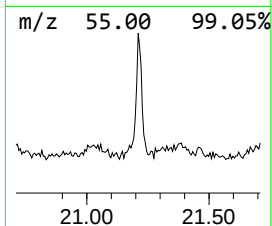
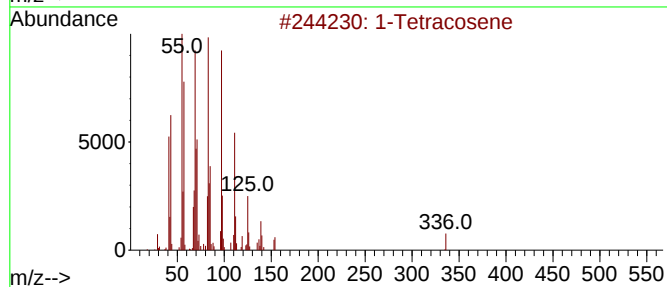
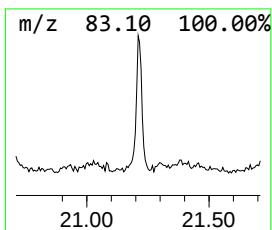
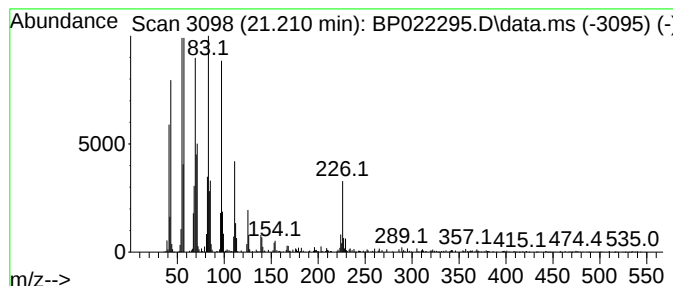
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Acq On : 09 Oct 2024 18:24
Operator : RC/JU
Sample : P4162-16
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4EJ1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 2



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100924\
Data File : BP022295.D
Acq On : 09 Oct 2024 18:24
Operator : RC/JU
Sample : P4162-16
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4EJ1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.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
Benzophenone	15.704	5.1	ng/ul	459622	3	14.316	1796140	20.0
1,1'-Biphenyl, ...	16.634	2.3	ng/ul	279227	4	17.116	2445790	20.0
1,1'-Biphenyl, ...	17.292	4.3	ng/ul	523249	4	17.116	2445790	20.0
1,1'-Biphenyl, ...	17.739	3.9	ng/ul	474097	4	17.116	2445790	20.0
n-Hexadecanoic ...	18.057	7.1	ng/ul	873385	4	17.116	2445790	20.0
2,2',3,6-Tetrac...	18.228	3.0	ng/ul	373410	4	17.116	2445790	20.0
1,1'-Biphenyl, ...	18.328	2.2	ng/ul	271865	4	17.116	2445790	20.0
Octadecanoic acid	19.386	2.1	ng/ul	280535	5	21.551	2625790	20.0
(DEL) Alkane: S...	20.686	2.4	ng/ul	310804	5	21.551	2625790	20.0
(DEL) Alkane: S...	21.210	5.2	ng/ul	680113	5	21.551	2625790	20.0