

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102124\
 Data File : BP022537.D
 Acq On : 22 Oct 2024 13:08
 Operator : RC/JU
 Sample : P4414-02
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A4E88

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: BP022537.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.440	242	247	255	rBV	34747	55302	0.27%	0.021%
2	6.058	516	522	530	rBV5	10186	26452	0.13%	0.010%
3	6.875	645	661	674	rVB2	480502	1237558	6.08%	0.459%
4	7.005	674	683	689	rBV	217942	345243	1.70%	0.128%
5	7.093	691	698	709	rVB	2070605	3048290	14.98%	1.131%
6	7.234	717	722	732	rVB	1204296	2006302	9.86%	0.744%
7	7.658	785	794	795	rBV	358368	492140	2.42%	0.183%
8	7.693	795	800	809	rVB	3418944	5406767	26.58%	2.006%
9	8.005	844	853	862	rBV	3437507	5470042	26.89%	2.029%
10	8.369	907	915	922	rBV	371480	595192	2.93%	0.221%
11	8.458	922	930	961	rVV	2178801	3661581	18.00%	1.358%
12	8.716	964	974	982	rVV	1962395	3199991	15.73%	1.187%
13	8.799	982	988	992	rVV	275333	508048	2.50%	0.188%
14	8.846	992	996	1005	rVB	516226	823302	4.05%	0.305%
15	9.205	1050	1057	1062	rBV3	19519	33158	0.16%	0.012%
16	9.363	1071	1084	1096	rBV	2597835	4701781	23.11%	1.744%
17	9.516	1102	1110	1111	rBV	266387	399691	1.96%	0.148%
18	9.540	1111	1114	1124	rVV	556162	871874	4.29%	0.323%
19	9.663	1128	1135	1142	rVB	104606	173182	0.85%	0.064%
20	9.834	1156	1164	1176	rVB	2378924	3713671	18.25%	1.378%
21	10.075	1194	1205	1220	rBV2	1732210	3696733	18.17%	1.371%
22	10.416	1252	1263	1270	rBV	445207	786879	3.87%	0.292%
23	10.552	1278	1286	1303	rBV	226991	433420	2.13%	0.161%
24	10.746	1307	1319	1330	rVB	3263289	5354203	26.32%	1.986%
25	11.710	1475	1483	1496	rBV	865550	1375656	6.76%	0.510%
26	11.969	1521	1527	1538	rVB3	26469	53883	0.26%	0.020%
27	12.081	1538	1546	1560	rBV	1614381	2584347	12.70%	0.959%
28	12.493	1607	1616	1623	rBV	40030	66149	0.33%	0.025%
29	12.699	1639	1651	1658	rBV	3918330	6577886	32.33%	2.440%
30	12.775	1658	1664	1685	rVB	1063613	1530414	7.52%	0.568%
31	13.099	1713	1719	1723	rBV2	20456	32724	0.16%	0.012%
32	13.199	1730	1736	1743	rVV	20171	35753	0.18%	0.013%
33	13.263	1743	1747	1751	rVV	16501	21592	0.11%	0.008%
34	13.534	1787	1793	1798	rBV4	15497	22893	0.11%	0.008%
35	13.681	1811	1818	1830	rBV	693809	1196939	5.88%	0.444%
36	13.869	1842	1850	1854	rBV	4778192	6695738	32.91%	2.484%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
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37	13.963	1857	1866	1869	rVV	873446	1452993	7.14%	0.539%
38	13.993	1869	1871	1880	rVB	695017	914641	4.50%	0.339%
39	14.281	1911	1920	1924	rBV	771133	1125523	5.53%	0.418%
40	14.340	1924	1930	1937	rVV	5528798	6968768	34.25%	2.585%
41	14.410	1937	1942	1950	rVB	42807	73011	0.36%	0.027%
42	14.528	1950	1962	1970	rBV2	1739943	2697850	13.26%	1.001%
43	14.604	1970	1975	1978	rVV	99058	152306	0.75%	0.057%
44	14.663	1978	1985	1991	rVV	3159268	4842037	23.80%	1.796%
45	14.916	2024	2028	2038	rVB	72806	109908	0.54%	0.041%
46	15.163	2063	2070	2079	rVB8	16913	43977	0.22%	0.016%
47	15.287	2084	2091	2094	rVV	1123506	1856650	9.13%	0.689%
48	15.345	2094	2101	2107	rVV2	8497163	14516007	71.35%	5.385%
49	15.422	2107	2114	2126	rVB	180369	350076	1.72%	0.130%
50	15.598	2140	2144	2148	rBV	121518	156688	0.77%	0.058%
51	15.651	2148	2153	2160	rVB	449047	573366	2.82%	0.213%
52	15.845	2182	2186	2193	rBV4	9489	22558	0.11%	0.008%
53	15.945	2199	2203	2207	rBV3	21477	26582	0.13%	0.010%
54	16.022	2211	2216	2231	rVB4	52202	112931	0.56%	0.042%
55	16.251	2245	2255	2262	rBV2	1730895	2377645	11.69%	0.882%
56	16.369	2267	2275	2289	rVB	10046009	14422873	70.89%	5.351%
57	16.728	2325	2336	2346	rBV	7152700	10419346	51.21%	3.865%
58	16.887	2358	2363	2373	rVB7	39822	96999	0.48%	0.036%
59	17.069	2388	2394	2398	rBV	935464	1504886	7.40%	0.558%
60	17.122	2398	2403	2408	rVV	2116244	3536784	17.38%	1.312%
61	17.181	2408	2413	2416	rVV2	1375731	2170398	10.67%	0.805%
62	17.216	2416	2419	2426	rVB	2199834	3312374	16.28%	1.229%
63	17.504	2462	2468	2474	rVB5	26970	56640	0.28%	0.021%
64	17.710	2498	2503	2510	rVV4	77319	134072	0.66%	0.050%
65	17.775	2510	2514	2521	rVB	144764	180360	0.89%	0.067%
66	17.969	2543	2547	2550	rBV5	20373	31933	0.16%	0.012%
67	18.010	2550	2554	2560	rBV	328818	432019	2.12%	0.160%
68	18.086	2560	2567	2571	rVV	7969072	9757244	47.96%	3.620%
69	18.169	2579	2581	2590	rVB6	25828	47294	0.23%	0.018%
70	18.322	2599	2607	2611	rBV4	55727	123923	0.61%	0.046%
71	18.398	2615	2620	2629	rVV5	77022	196627	0.97%	0.073%
72	18.481	2629	2634	2640	rVV4	45835	106221	0.52%	0.039%
73	18.539	2640	2644	2650	rVB	255822	367062	1.80%	0.136%
74	18.751	2677	2680	2684	rVB3	41105	55745	0.27%	0.021%
75	18.792	2685	2687	2691	rVB3	19784	21210	0.10%	0.008%
76	18.875	2696	2701	2706	rBV	86216	172090	0.85%	0.064%

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77	18.922	2706	2709	2714	rVV3	32377	55034	0.27%	0.020%
78	18.969	2715	2717	2720	rVV4	19511	22730	0.11%	0.008%
79	18.998	2720	2722	2726	rVB3	61197	60785	0.30%	0.023%
80	19.104	2737	2740	2743	rBV3	26977	39880	0.20%	0.015%
81	19.139	2743	2746	2750	rVV3	56854	79435	0.39%	0.029%
82	19.216	2752	2759	2766	rVB	10865814	14186752	69.73%	5.263%
83	19.339	2776	2780	2789	rBV	123467	190387	0.94%	0.071%
84	19.463	2798	2801	2807	rVB	72739	97658	0.48%	0.036%
85	19.557	2811	2817	2819	rBV	2014586	2840283	13.96%	1.054%
86	19.586	2819	2822	2832	rVB	4533618	6307553	31.00%	2.340%
87	20.292	2937	2942	2946	rBV2	222970	325291	1.60%	0.121%
88	20.533	2979	2983	2988	rVB	2366877	2575674	12.66%	0.956%
89	21.192	3090	3095	3104	rVB	332352	578540	2.84%	0.215%
90	21.339	3116	3120	3126	rBV2	215098	293332	1.44%	0.109%
91	21.427	3129	3135	3139	rVV	8601675	11982814	58.90%	4.446%
92	21.498	3139	3147	3153	rVV	8148718	14846163	72.97%	5.508%
93	21.575	3153	3160	3169	rVB	12412614	20344703	100.00%	7.548%
94	22.574	3326	3330	3338	rVB4	195459	375300	1.84%	0.139%
95	23.763	3520	3532	3542	rVB	7691083	19042472	93.60%	7.065%
96	24.545	3657	3665	3671	rBV	908516	2628083	12.92%	0.975%
97	24.621	3672	3678	3690	rVB	1590397	3908858	19.21%	1.450%
98	24.768	3695	3703	3714	rVB	663495	1786149	8.78%	0.663%
99	28.568	4324	4349	4364	rVB2	2231156	11342927	55.75%	4.208%
100	29.615	4517	4527	4550	rVB	624377	2882847	14.17%	1.070%

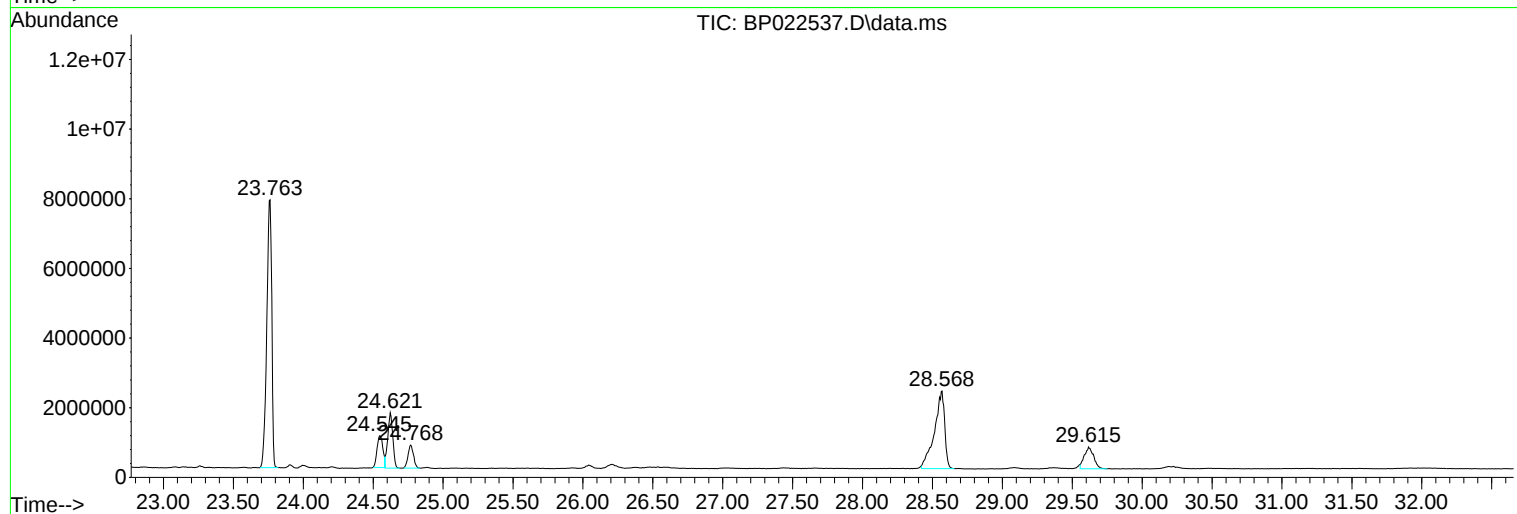
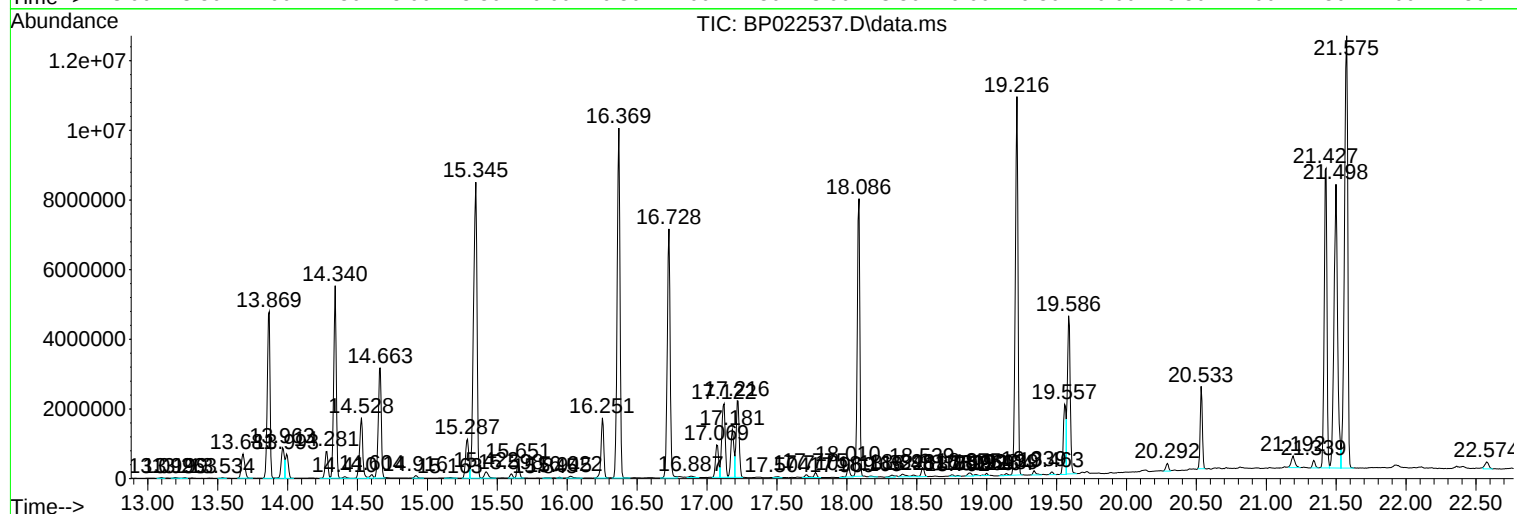
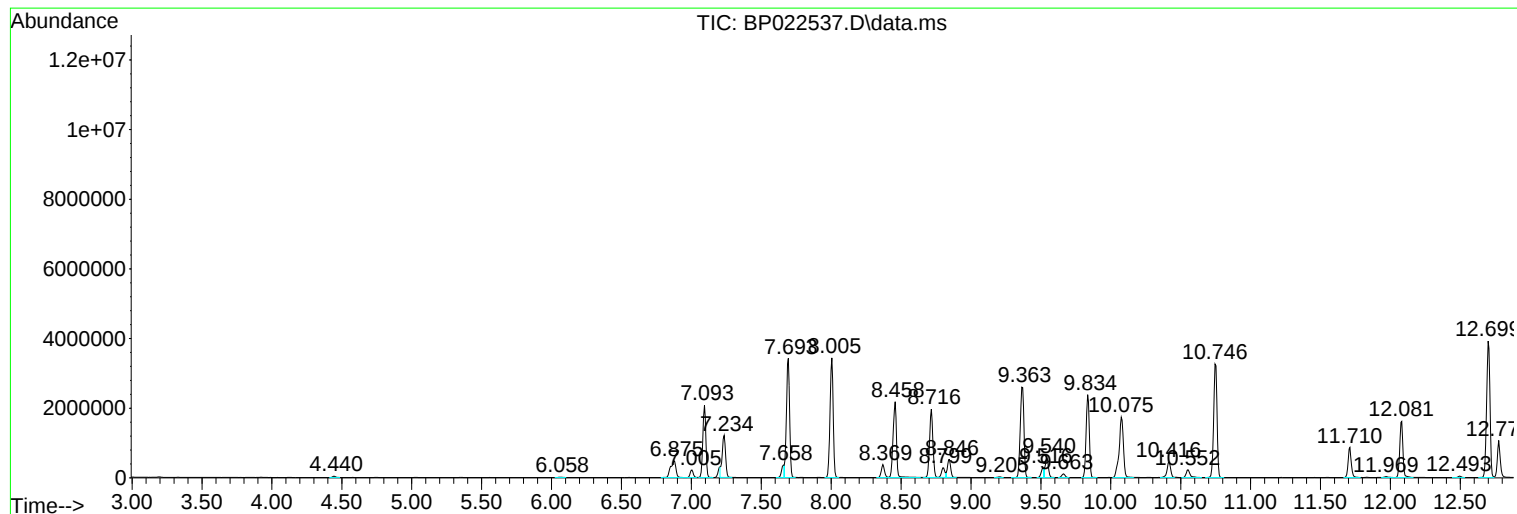
Sum of corrected areas: 269548043

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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



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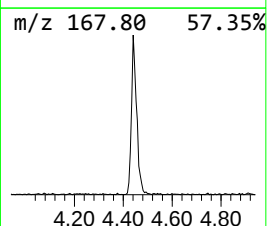
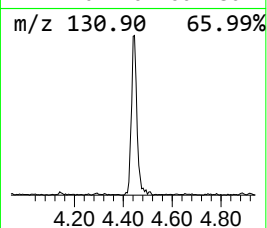
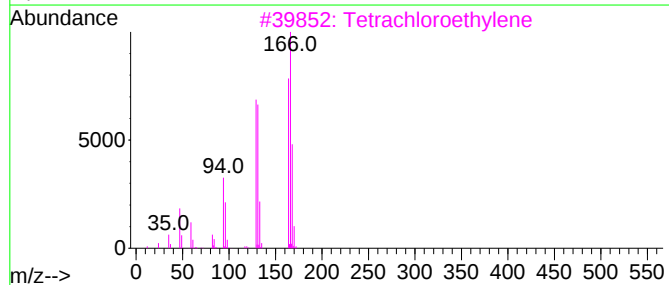
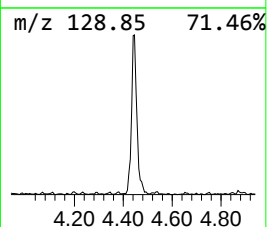
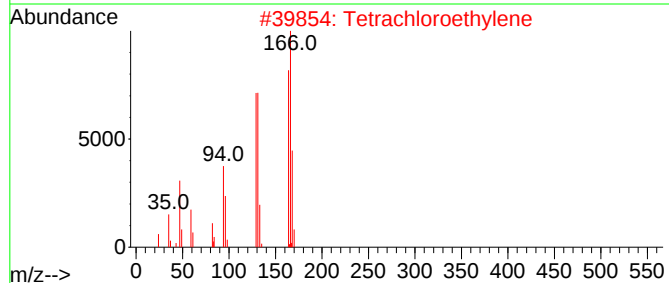
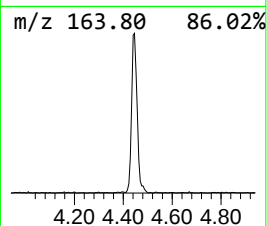
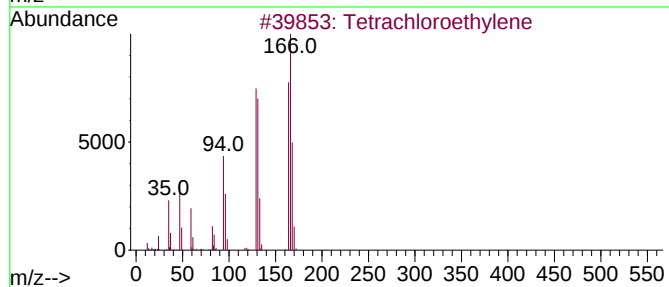
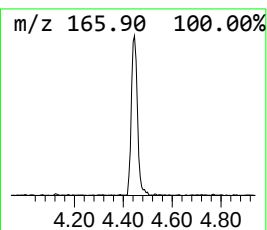
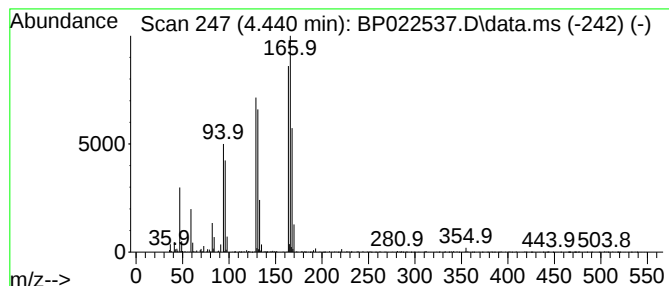
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Tetrachloroethylene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.440	2.25 ng/ul	55302	1,4-Dichlorobenzene-d4	7.658

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetrachloroethylene	164	C2Cl4	000127-18-4	95
2			Tetrachloroethylene	164	C2Cl4	000127-18-4	95
3			Tetrachloroethylene	164	C2Cl4	000127-18-4	94
4			Tetrachloroethylene	164	C2Cl4	000127-18-4	93
5			Pyrimidine, 5-fluoro-2,4-dichloro-	166	C4HC12FN2	002927-71-1	35



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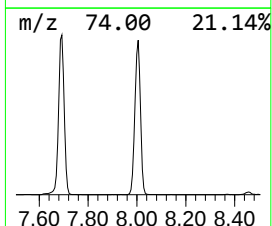
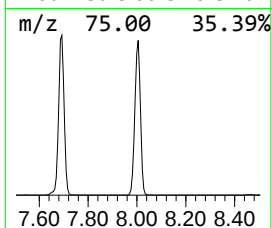
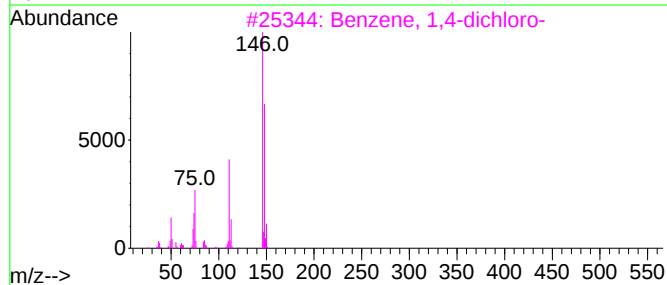
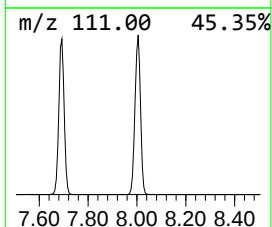
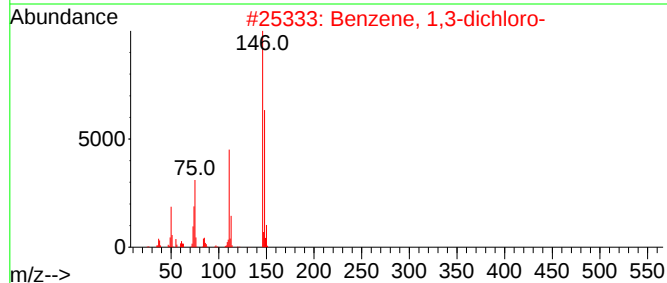
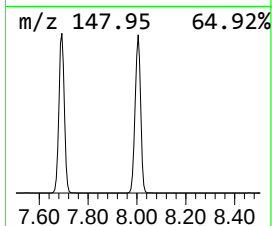
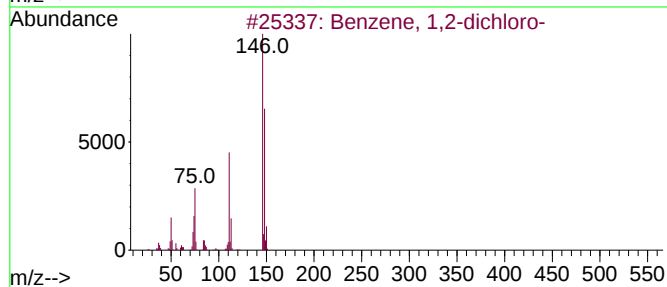
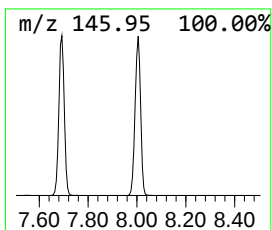
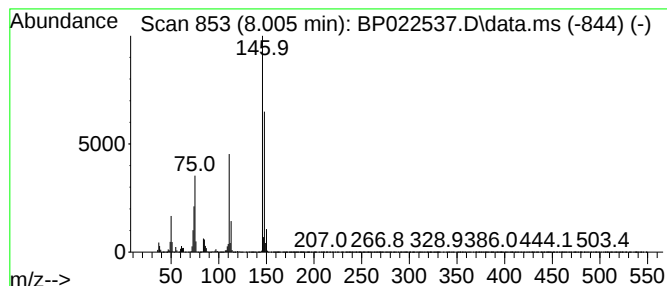
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzene, 1,2-dichloro- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.005	222.30 ng/ul	5470040	1,4-Dichlorobenzene-d4	7.658

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	98
2			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	98
3			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	98
4			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	97
5			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	96



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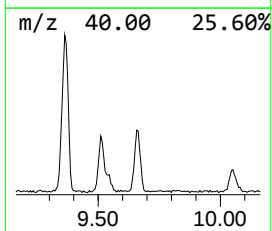
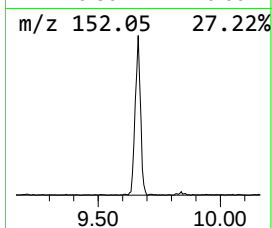
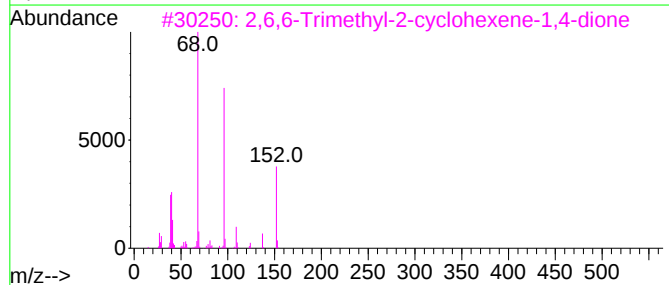
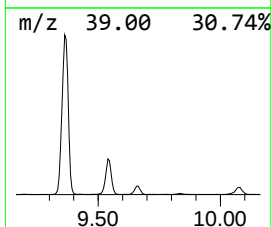
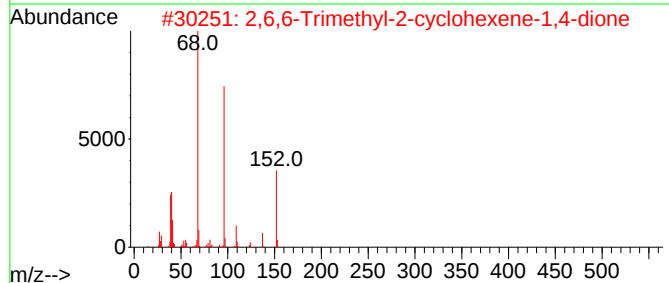
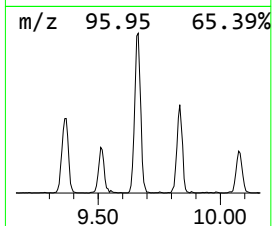
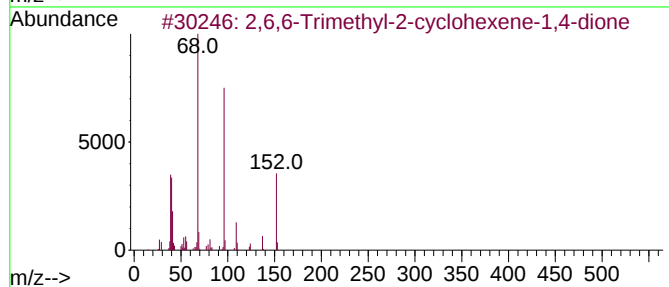
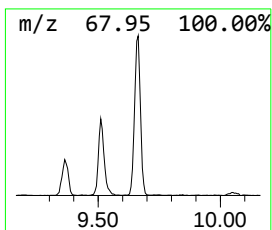
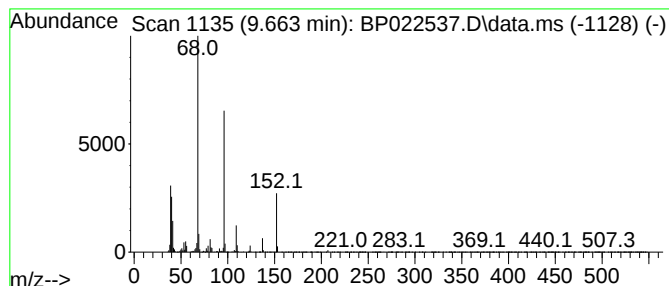
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 2,6,6-Trimethyl-2-cyclohexe... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.663	4.40 ng/ul	173182	Naphthalene-d8	10.416

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,6,6-Trimethyl-2-cyclohexene-1,...	152	C9H12O2	001125-21-9	97		
2	2,6,6-Trimethyl-2-cyclohexene-1,...	152	C9H12O2	001125-21-9	95		
3	2,6,6-Trimethyl-2-cyclohexene-1,...	152	C9H12O2	001125-21-9	94		
4	2,6,6-Trimethyl-2-cyclohexene-1,...	152	C9H12O2	001125-21-9	91		
5	2-Cyclohexen-1-one	96	C6H8O	000930-68-7	53		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102124\
Data File : BP022537.D
Acq On : 22 Oct 2024 13:08
Operator : RC/JU
Sample : P4414-02
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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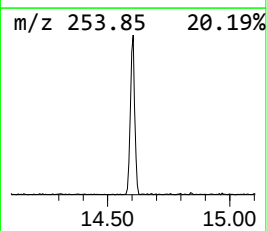
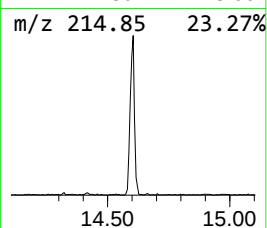
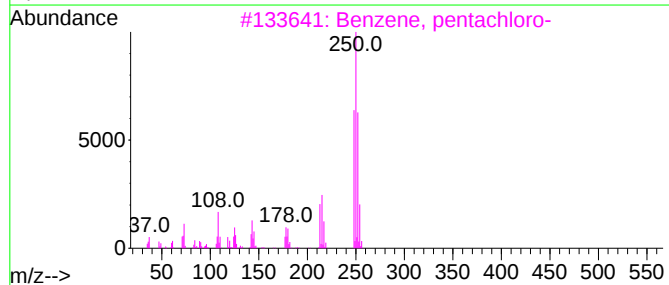
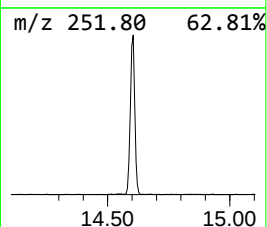
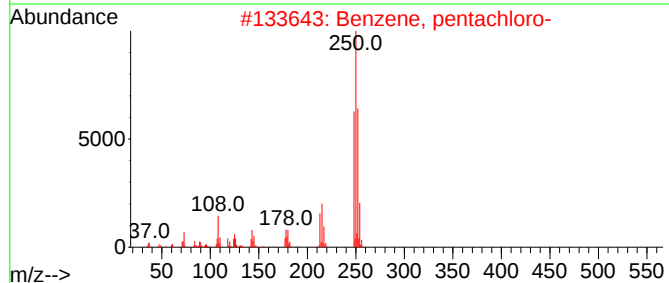
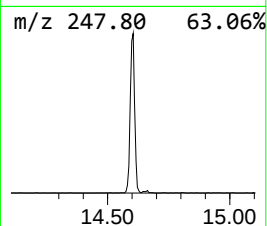
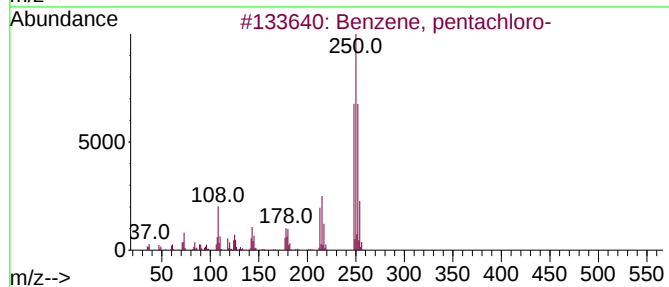
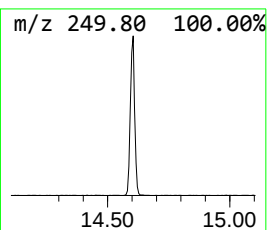
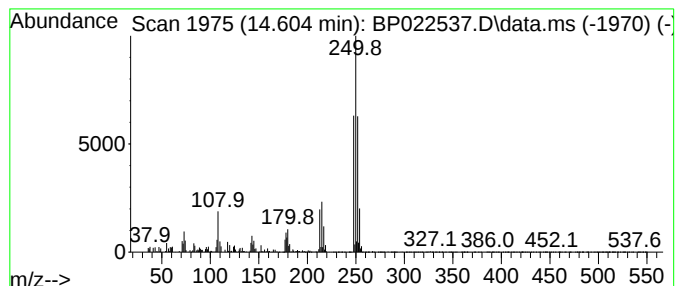
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Benzene, pentachloro- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.604	2.71 ng/ul	152306	Acenaphthene-d10	14.281

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, pentachloro-	248	C6HCl5	000608-93-5	99
2			Benzene, pentachloro-	248	C6HCl5	000608-93-5	99
3			Benzene, pentachloro-	248	C6HCl5	000608-93-5	96
4			Benzene, pentachloro-	248	C6HCl5	000608-93-5	91
5			Benzene, pentachloro-	248	C6HCl5	000608-93-5	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102124\
Data File : BP022537.D
Acq On : 22 Oct 2024 13:08
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Sample : P4414-02
Misc :
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Instrument :
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ClientSampleId :
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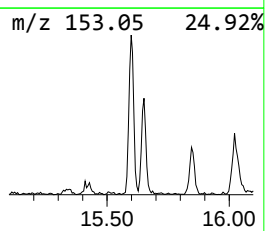
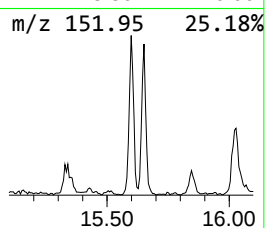
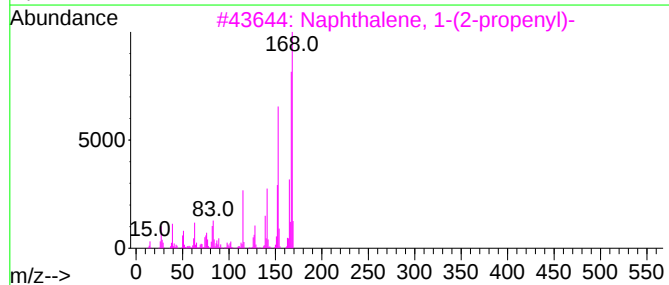
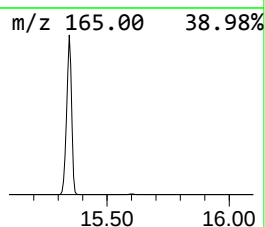
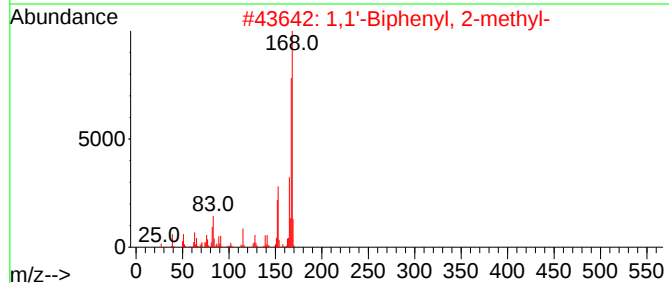
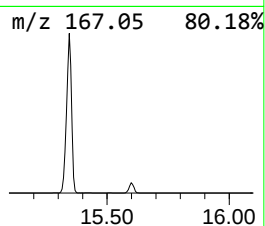
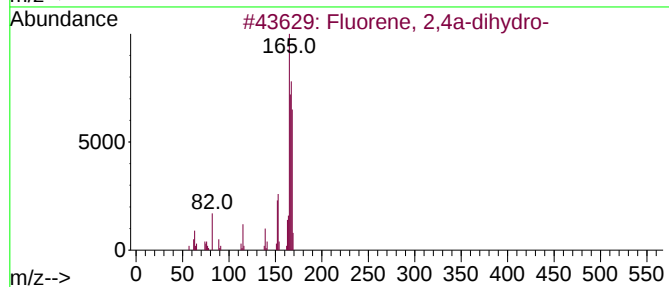
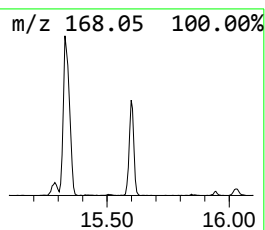
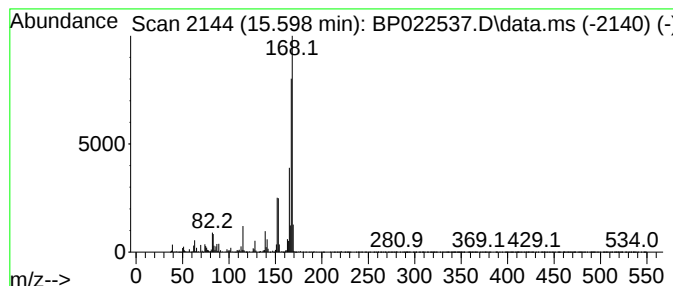
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Fluorene, 2,4a-dihydro- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.598	2.78 ng/ul	156688	Acenaphthene-d10	14.281

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Fluorene, 2,4a-dihydro-	168	C13H12	059247-36-8	94
2			1,1'-Biphenyl, 2-methyl-	168	C13H12	000643-58-3	94
3			Naphthalene, 1-(2-propenyl)-	168	C13H12	002489-86-3	93
4			1,1'-Biphenyl, 4-methyl-	168	C13H12	000644-08-6	91
5			Diphenylmethane	168	C13H12	000101-81-5	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102124\
Data File : BP022537.D
Acq On : 22 Oct 2024 13:08
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Sample : P4414-02
Misc :
ALS Vial : 38 Sample Multiplier: 1

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ClientSampleId :
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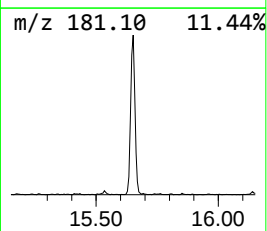
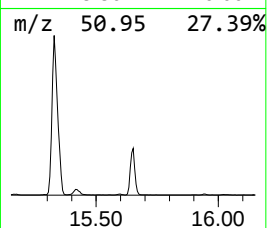
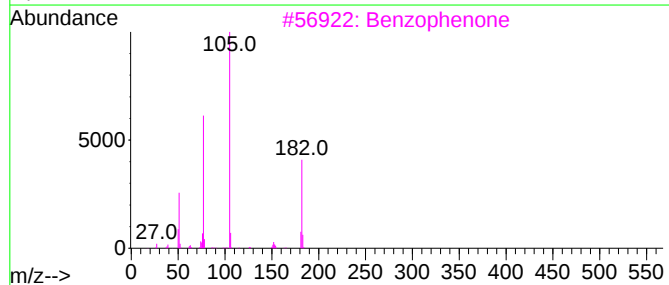
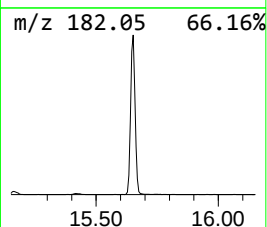
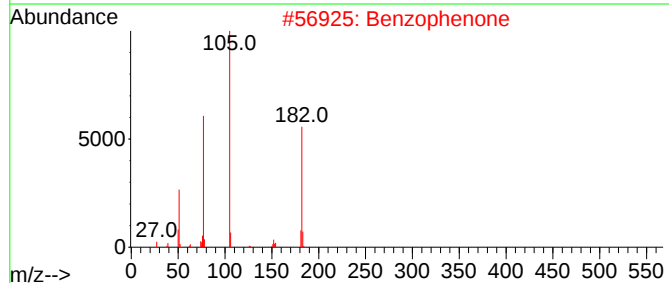
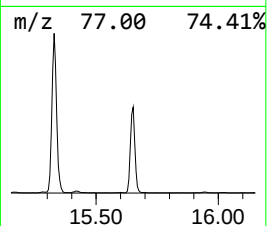
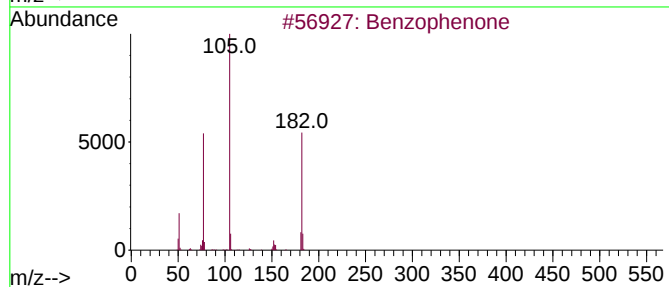
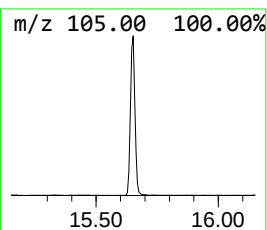
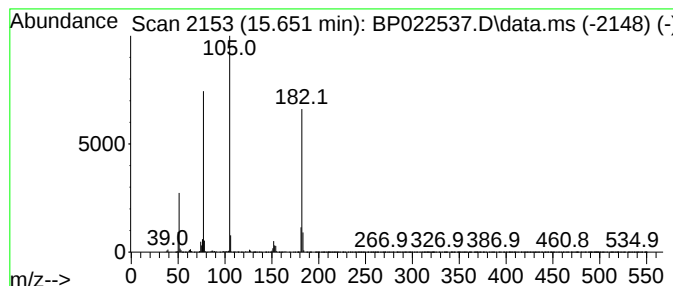
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Benzophenone Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.651	10.19 ng/ul	573366	Acenaphthene-d10	14.281

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	97
2			Benzophenone	182	C13H10O	000119-61-9	96
3			Benzophenone	182	C13H10O	000119-61-9	94
4			Benzophenone	182	C13H10O	000119-61-9	91
5			Benzophenone	182	C13H10O	000119-61-9	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102124\
Data File : BP022537.D
Acq On : 22 Oct 2024 13:08
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Sample : P4414-02
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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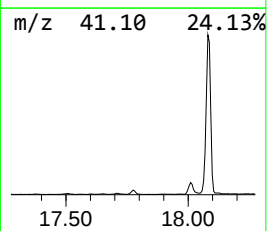
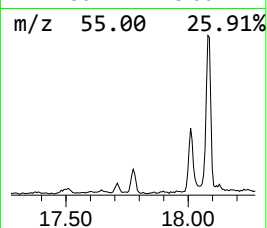
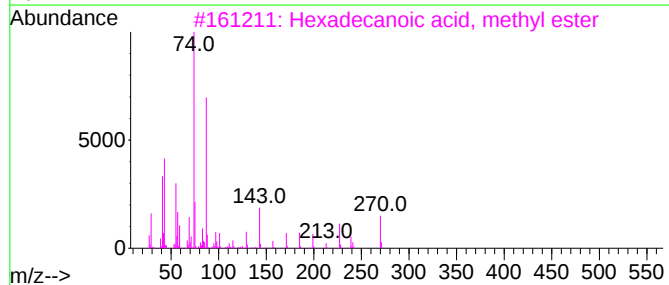
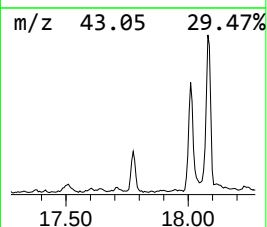
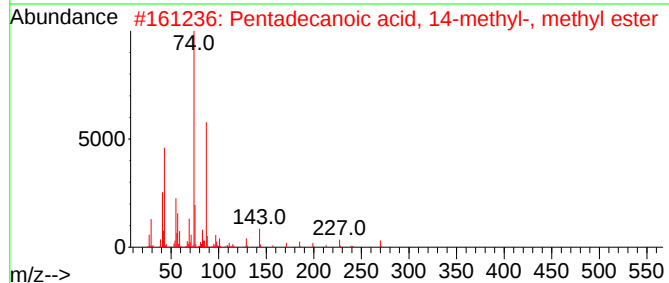
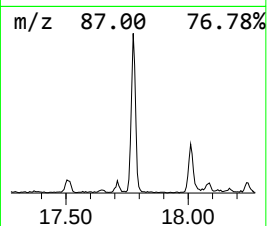
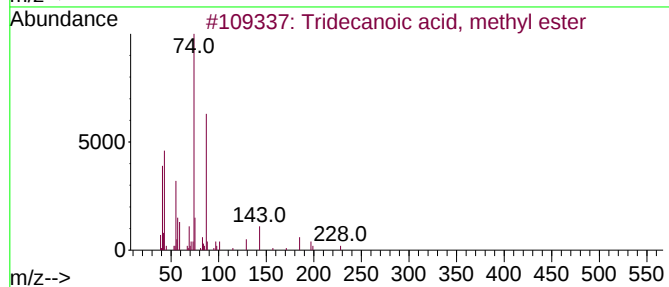
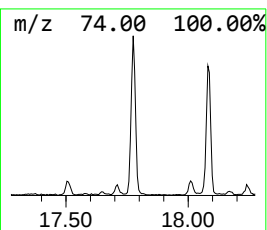
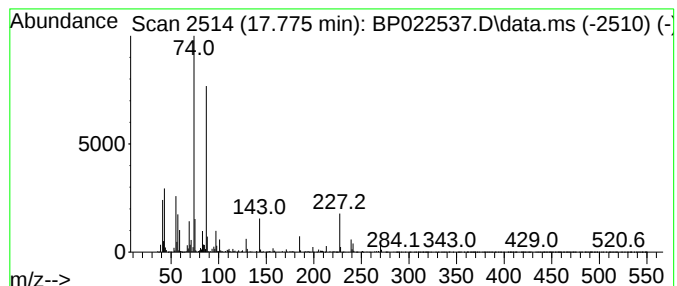
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Tridecanoic acid, methyl ester Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.775	2.40 ng/ul	180360	Phenanthrene-d10	17.069

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	93
2			Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	93
3			Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	93
4			Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	91
5			Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102124\
Data File : BP022537.D
Acq On : 22 Oct 2024 13:08
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Sample : P4414-02
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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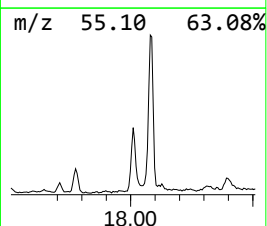
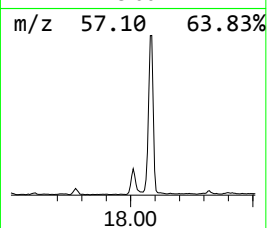
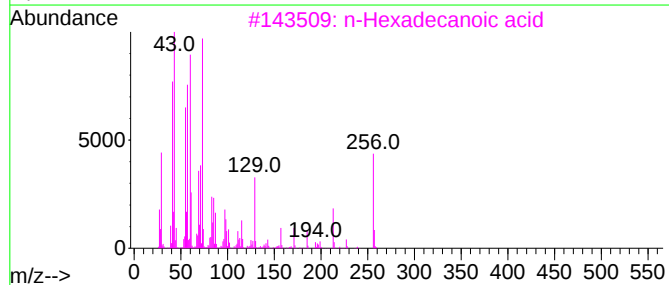
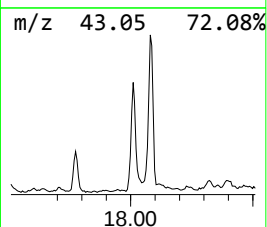
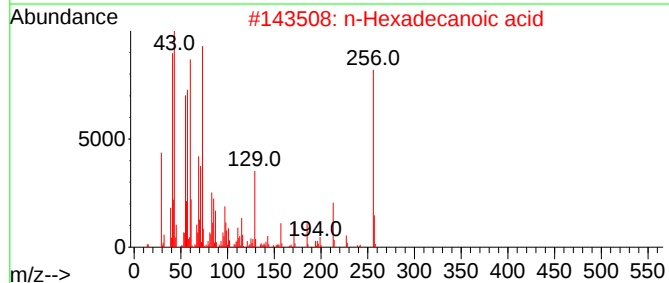
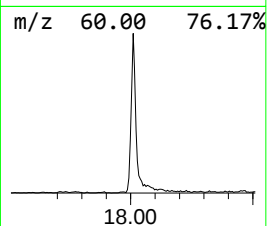
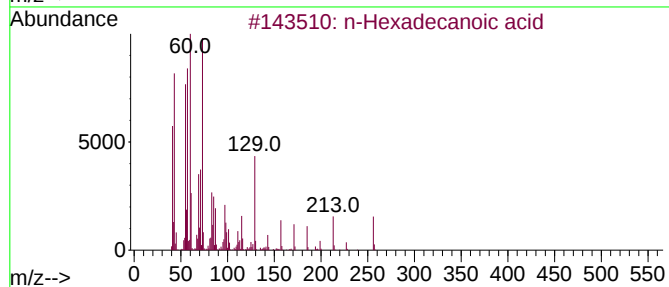
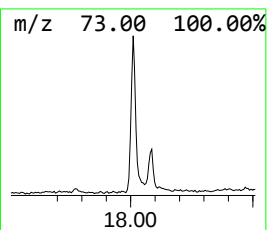
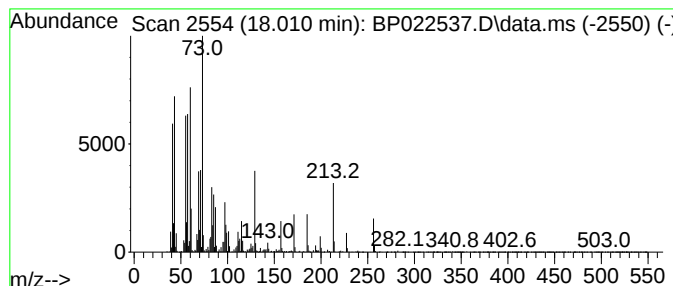
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.010	5.74 ng/ul	432019	Phenanthrene-d10	17.069

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
4			Tridecanoic acid	214	C13H26O2	000638-53-9	90
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102124\
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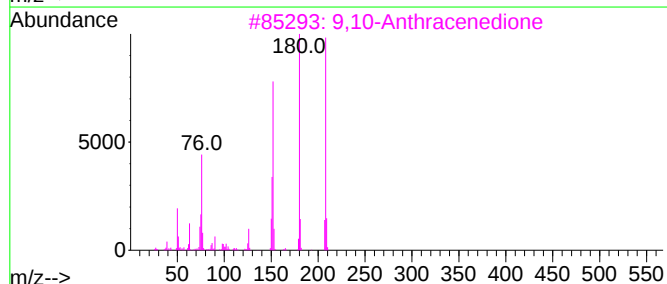
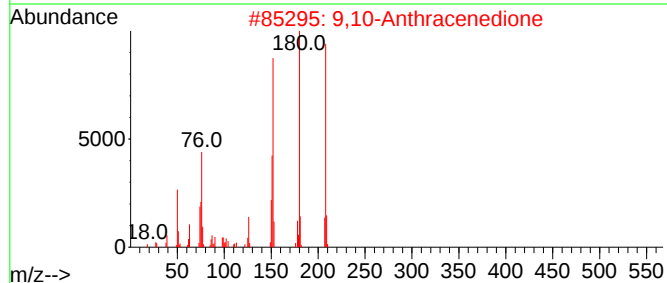
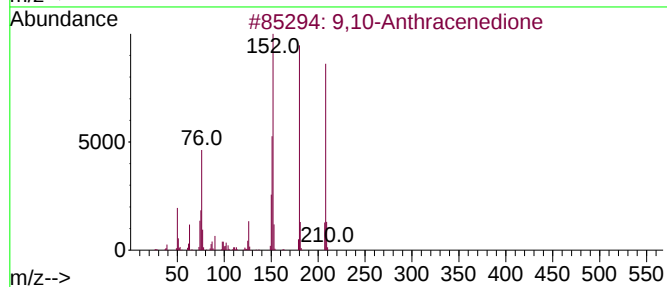
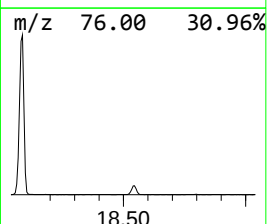
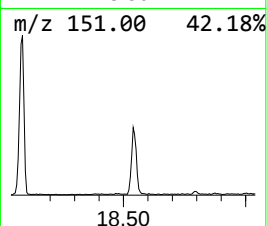
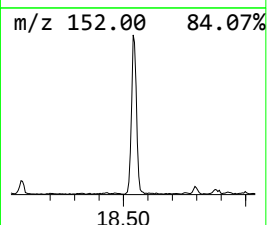
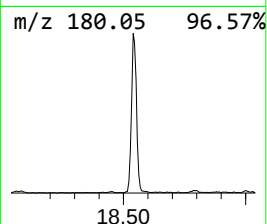
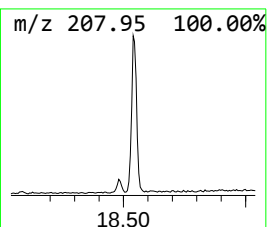
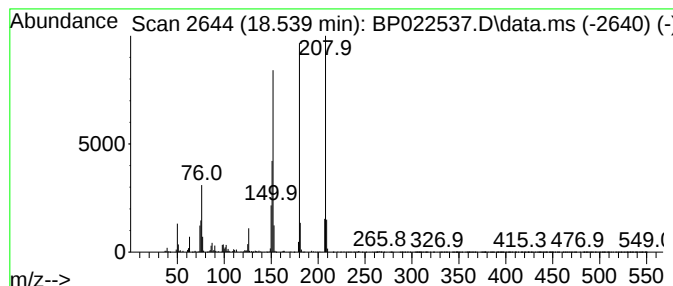
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 9,10-Anthracenedione Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.539	4.88 ng/ul	367062	Phenanthrene-d10	17.069

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,10-Anthracenedione			208	C14H8O2	000084-65-1	99
2	9,10-Anthracenedione			208	C14H8O2	000084-65-1	98
3	9,10-Anthracenedione			208	C14H8O2	000084-65-1	94
4	9,10-Anthracenedione			208	C14H8O2	000084-65-1	90
5	9H-Fluoren-9-one			180	C13H8O	000486-25-9	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102124\
Data File : BP022537.D
Acq On : 22 Oct 2024 13:08
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Sample : P4414-02
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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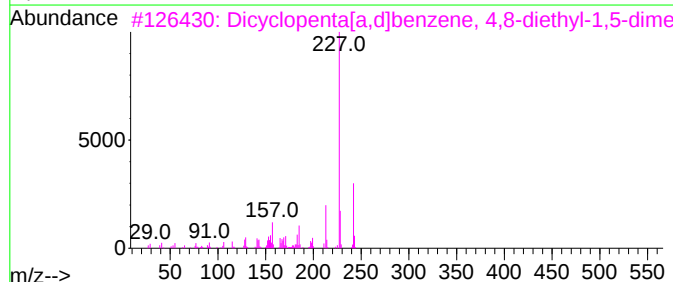
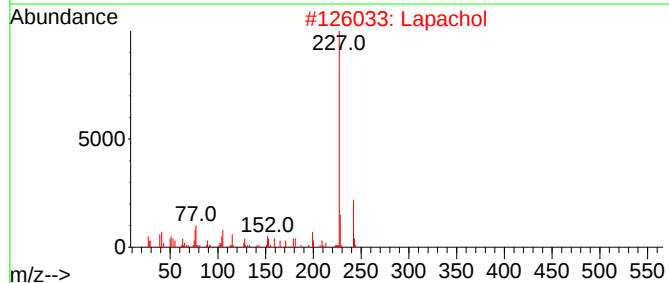
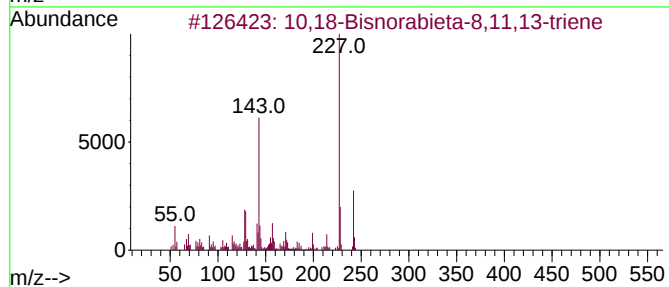
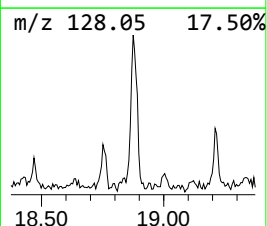
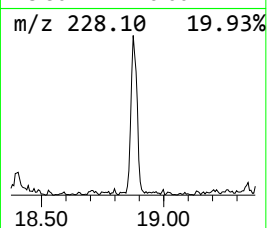
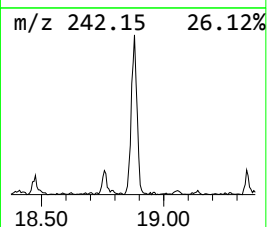
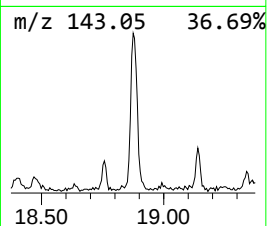
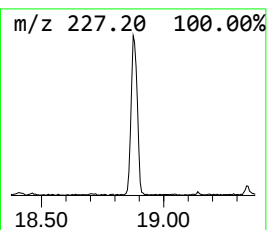
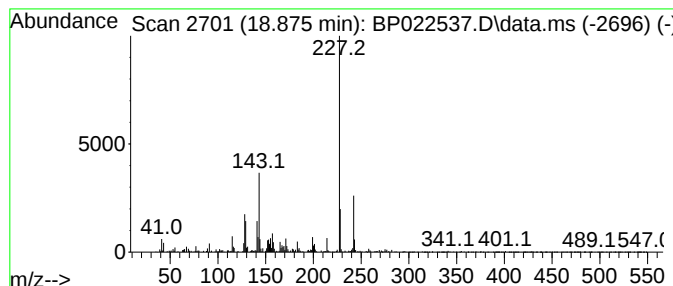
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 10,18-Bisnorabieta-8,11,13-... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.875	2.29 ng/ul	172090	Phenanthrene-d10	17.069

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	10,18-Bisnorabieta-8,11,13-triene	242	C18H26	032624-67-2	94		
2	Lapachol	242	C15H14O3	000084-79-7	70		
3	Dicyclopenta[a,d]benzene, 4,8-di...	242	C18H26	1000156-41-6	64		
4	s-Indacene-1,7-dione, 2,3,5,6-te...	242	C16H18O2	055591-17-8	58		
5	5-Methoxy-2-methyl-4-oxo-1,2,3,4...	242	C14H14N2O2	1000227-06-0	52		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102124\
Data File : BP022537.D
Acq On : 22 Oct 2024 13:08
Operator : RC/JU
Sample : P4414-02
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4E88

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
Tetrachloroethy...	4.440	2.3	ng/ul	55302	1	7.658	492140	20.0
Benzene, 1,2-di...	8.005	222.3	ng/ul	5470040	1	7.658	492140	20.0
2,6,6-Trimethyl...	9.663	4.4	ng/ul	173182	2	10.416	786879	20.0
Benzene, pentac...	14.604	2.7	ng/ul	152306	3	14.281	1125520	20.0
Fluorene, 2,4a-...	15.598	2.8	ng/ul	156688	3	14.281	1125520	20.0
Benzophenone	15.651	10.2	ng/ul	573366	3	14.281	1125520	20.0
Tridecanoic aci...	17.775	2.4	ng/ul	180360	4	17.069	1504890	20.0
n-Hexadecanoic ...	18.010	5.7	ng/ul	432019	4	17.069	1504890	20.0
9,10-Anthracene...	18.539	4.9	ng/ul	367062	4	17.069	1504890	20.0
10,18-Bisnorabi...	18.875	2.3	ng/ul	172090	4	17.069	1504890	20.0