

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102424\
 Data File : BP022588.D
 Acq On : 24 Oct 2024 15:12
 Operator : RC/JU
 Sample : P4436-02
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A4E97

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: BP022588.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.499	422	427	431	rBV2	14633	22511	0.12%	0.009%
2	6.034	510	518	522	rBV3	39720	76547	0.42%	0.032%
3	6.075	523	525	530	rVB2	14861	19613	0.11%	0.008%
4	6.481	588	594	598	rBV2	19745	29483	0.16%	0.012%
5	6.528	598	602	611	rVB	27944	45932	0.25%	0.019%
6	6.869	645	660	675	rVB2	435336	1132210	6.24%	0.474%
7	6.993	675	681	688	rBV	204309	307075	1.69%	0.129%
8	7.087	690	697	709	rVB	1821288	2770724	15.27%	1.161%
9	7.228	709	721	731	rBV2	1062800	2126100	11.72%	0.891%
10	7.687	784	799	812	rBV	3011341	5313623	29.28%	2.226%
11	7.999	843	852	862	rBV	3104761	4885491	26.92%	2.047%
12	8.364	907	914	921	rBV	335408	553907	3.05%	0.232%
13	8.446	921	928	941	rVV	2240311	3576054	19.71%	1.498%
14	8.711	965	973	981	rVB	1825038	2966527	16.35%	1.243%
15	8.793	981	987	990	rVV	282182	440698	2.43%	0.185%
16	8.834	990	994	1001	rVB	492604	769882	4.24%	0.323%
17	9.193	1048	1055	1061	rVB4	17822	30735	0.17%	0.013%
18	9.358	1073	1083	1099	rBV	2690294	4595409	25.33%	1.925%
19	9.505	1099	1108	1109	rVV	242901	356429	1.96%	0.149%
20	9.534	1109	1113	1123	rVV	496864	853659	4.70%	0.358%
21	9.658	1125	1134	1142	rVB	101820	170872	0.94%	0.072%
22	9.828	1155	1163	1172	rVB	2375729	3567181	19.66%	1.495%
23	10.069	1193	1204	1217	rBV2	1661791	3423779	18.87%	1.434%
24	10.410	1251	1262	1269	rBV	471063	814358	4.49%	0.341%
25	10.552	1279	1286	1300	rVB	188651	362115	2.00%	0.152%
26	10.740	1307	1318	1325	rBV	3047146	4866511	26.82%	2.039%
27	11.705	1472	1482	1498	rBV	795332	1305426	7.19%	0.547%
28	11.963	1520	1526	1537	rVB3	24292	53145	0.29%	0.022%
29	12.069	1537	1544	1558	rBV	1558401	2414028	13.30%	1.011%
30	12.487	1610	1615	1621	rVV	38953	53561	0.30%	0.022%
31	12.687	1631	1649	1656	rBV	3899428	6125312	33.76%	2.566%
32	12.763	1656	1662	1676	rVB	853266	1410783	7.78%	0.591%
33	13.187	1729	1734	1741	rVV2	27570	40826	0.22%	0.017%
34	13.675	1809	1817	1828	rVB	686380	1146266	6.32%	0.480%
35	13.857	1840	1848	1855	rBV	5062671	6462452	35.62%	2.708%
36	13.957	1859	1865	1868	rVV2	824934	1296137	7.14%	0.543%

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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
 Title : SVOA CALIBRATION

37	13.987	1868	1870	1877	rVB	638704	820696	4.52%	0.344%
38	14.187	1900	1904	1909	rVB3	16167	25341	0.14%	0.011%
39	14.275	1912	1919	1923	rBV	733444	1132635	6.24%	0.475%
40	14.334	1923	1929	1935	rVV	4554034	6296350	34.70%	2.638%
41	14.410	1936	1942	1949	rVB2	42187	68524	0.38%	0.029%
42	14.522	1952	1961	1969	rBV2	1602944	2556794	14.09%	1.071%
43	14.598	1969	1974	1977	rVV	92591	143064	0.79%	0.060%
44	14.657	1977	1984	1990	rVV	3158277	4671694	25.75%	1.957%
45	14.710	1990	1993	2004	rVB4	22411	37898	0.21%	0.016%
46	14.910	2023	2027	2038	rVB	68181	104714	0.58%	0.044%
47	15.163	2063	2070	2076	rVB7	15311	34461	0.19%	0.014%
48	15.281	2082	2090	2094	rBV	1160265	1781709	9.82%	0.746%
49	15.340	2094	2100	2108	rVV2	7768147	13299390	73.30%	5.572%
50	15.428	2108	2115	2126	rVV	257357	507027	2.79%	0.212%
51	15.534	2126	2133	2138	rVV10	7292	19542	0.11%	0.008%
52	15.593	2138	2143	2147	rVV	107672	151684	0.84%	0.064%
53	15.640	2147	2151	2158	rVB	320164	443504	2.44%	0.186%
54	15.934	2198	2201	2206	rVB3	18434	24426	0.13%	0.010%
55	16.010	2206	2214	2231	rVB5	77094	195404	1.08%	0.082%
56	16.245	2244	2254	2262	rBV	1641546	2153352	11.87%	0.902%
57	16.363	2267	2274	2281	rBV	8610437	13179377	72.63%	5.522%
58	16.428	2281	2285	2290	rVB3	53659	89016	0.49%	0.037%
59	16.475	2290	2293	2298	rBV6	12379	20009	0.11%	0.008%
60	16.722	2325	2335	2347	rBV	6220883	9657470	53.22%	4.046%
61	16.881	2357	2362	2370	rVB5	48671	106179	0.59%	0.044%
62	17.069	2387	2394	2398	rBV	821046	1471583	8.11%	0.617%
63	17.122	2398	2403	2409	rVV	2082992	3289745	18.13%	1.378%
64	17.187	2409	2414	2416	rVV	1368354	1902802	10.49%	0.797%
65	17.222	2416	2420	2425	rVB	2233414	3109243	17.14%	1.303%
66	17.504	2462	2468	2475	rBV4	24190	46001	0.25%	0.019%
67	17.704	2497	2502	2510	rBV4	54864	94755	0.52%	0.040%
68	17.781	2510	2515	2522	rVB	98530	141151	0.78%	0.059%
69	17.963	2542	2546	2549	rBV5	17614	25558	0.14%	0.011%
70	18.004	2549	2553	2559	rVV	431480	603164	3.32%	0.253%
71	18.081	2559	2566	2573	rVB	5577074	9081361	50.05%	3.805%
72	18.175	2579	2582	2588	rVB5	21958	34992	0.19%	0.015%
73	18.316	2600	2606	2614	rBV4	53740	103195	0.57%	0.043%
74	18.392	2614	2619	2627	rVV5	49825	128084	0.71%	0.054%
75	18.481	2629	2634	2640	rVV4	26135	62148	0.34%	0.026%
76	18.557	2641	2647	2652	rVB	245990	399404	2.20%	0.167%

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Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

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Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

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Stop Thrs : 0

Peak Location: TOP

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Peak separation: 5

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 Title : SVOA CALIBRATION

77	18.739	2676	2678	2683	rVB6	25221	31905	0.18%	0.013%
78	18.881	2697	2702	2706	rBV	49952	91473	0.50%	0.038%
79	18.998	2720	2722	2725	rVB3	26036	23050	0.13%	0.010%
80	19.098	2736	2739	2742	rBV3	29246	35418	0.20%	0.015%
81	19.145	2743	2747	2750	rVV4	49711	76402	0.42%	0.032%
82	19.210	2750	2758	2767	rVB	8573965	13336747	73.50%	5.588%
83	19.339	2776	2780	2785	rBV2	89459	151265	0.83%	0.063%
84	19.457	2796	2800	2805	rBV	88564	125071	0.69%	0.052%
85	19.551	2810	2816	2818	rVV	1644267	2583992	14.24%	1.083%
86	19.586	2818	2822	2832	rVB	3952964	5971970	32.91%	2.502%
87	20.286	2938	2941	2946	rBV	129181	148976	0.82%	0.062%
88	20.533	2979	2983	2988	rBV	2090189	2369386	13.06%	0.993%
89	21.180	3089	3093	3105	rVB3	348698	629186	3.47%	0.264%
90	21.339	3117	3120	3127	rVB2	103543	143497	0.79%	0.060%
91	21.422	3128	3134	3140	rVV	7735105	10873220	59.92%	4.555%
92	21.498	3140	3147	3153	rVV2	6888012	13673160	75.36%	5.729%
93	21.569	3153	3159	3165	rVB	11714626	18144904	100.00%	7.602%
94	23.757	3520	3531	3541	rVB	7417283	17487998	96.38%	7.327%
95	24.551	3657	3666	3671	rBV	862660	2415357	13.31%	1.012%
96	24.627	3671	3679	3692	rVB	1307733	3539825	19.51%	1.483%
97	24.774	3696	3704	3715	rVB	611520	1685839	9.29%	0.706%
98	28.474	4325	4333	4334	rBV	406619	690989	3.81%	0.289%
99	28.533	4337	4343	4344	rBV	1103047	1316892	7.26%	0.552%
100	29.609	4518	4526	4527	rBV	403242	739555	4.08%	0.310%

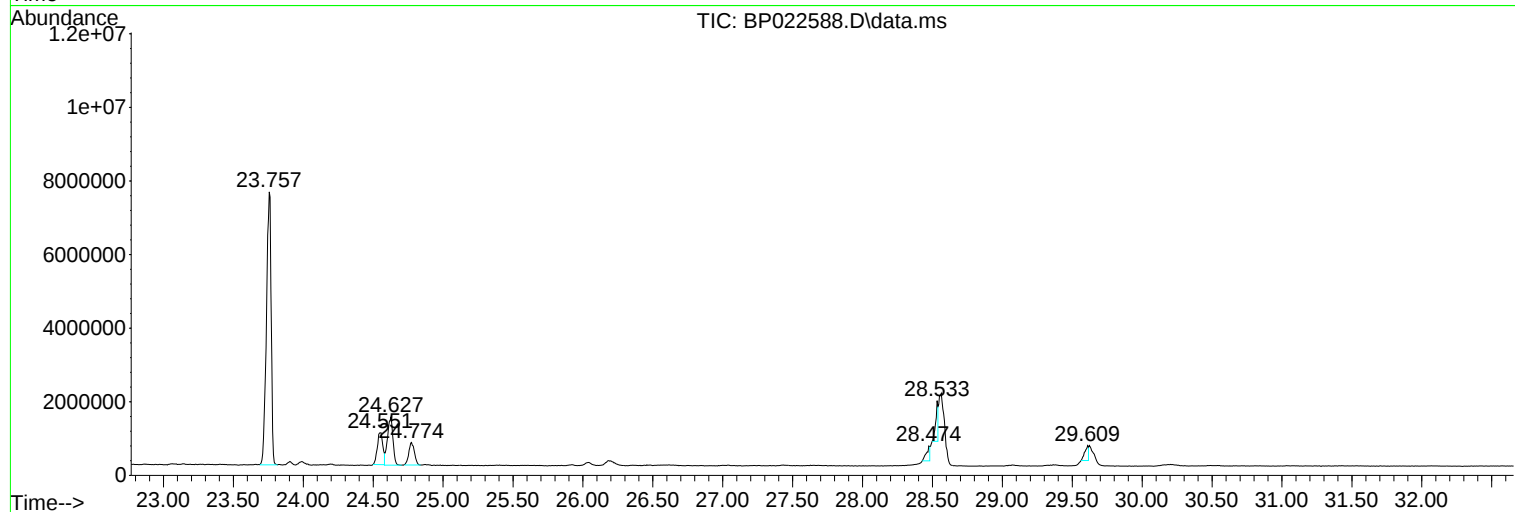
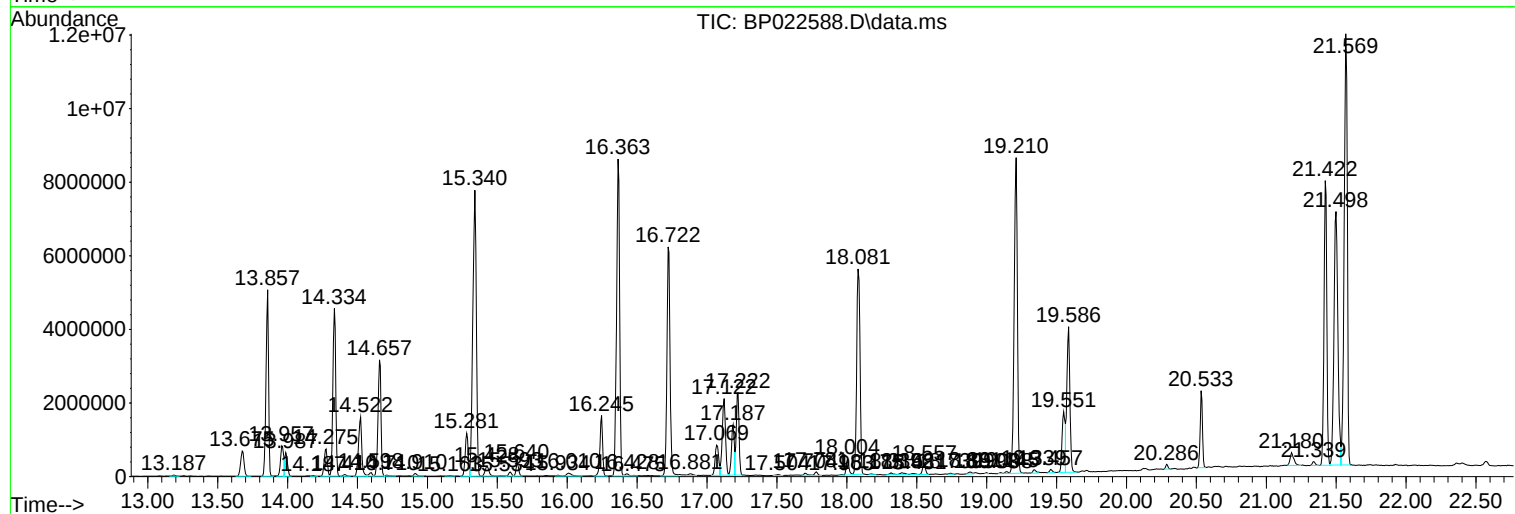
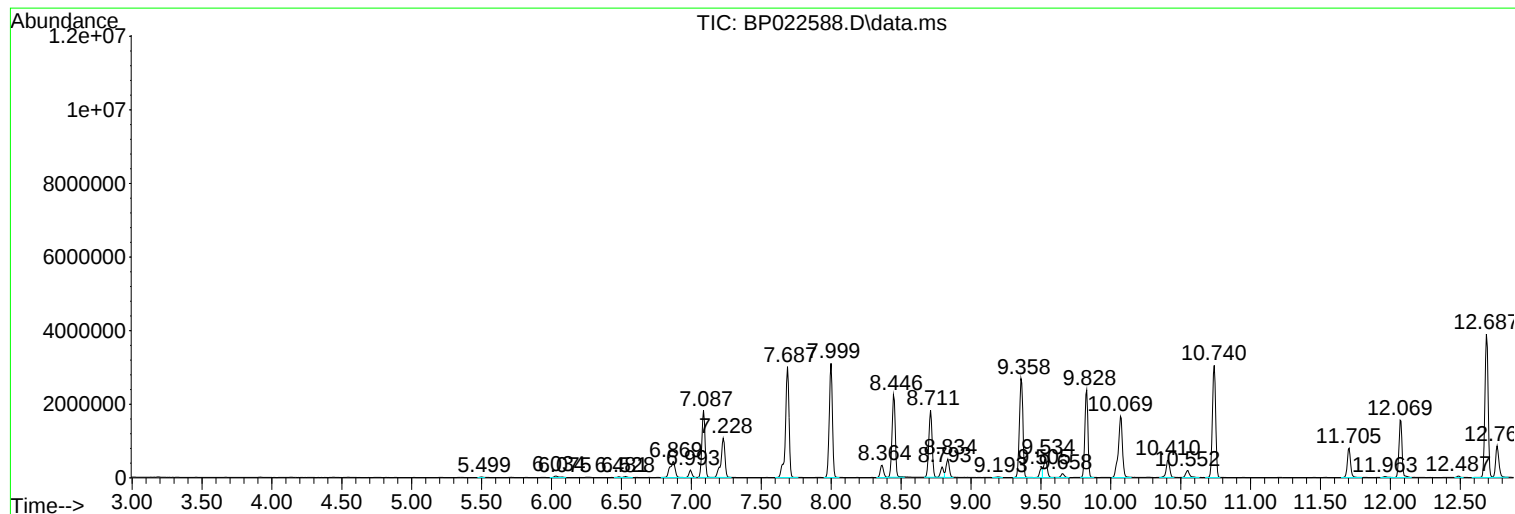
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Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION

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



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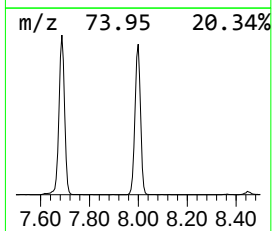
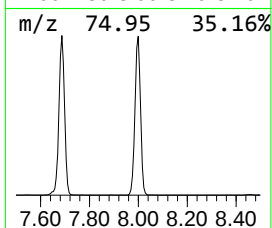
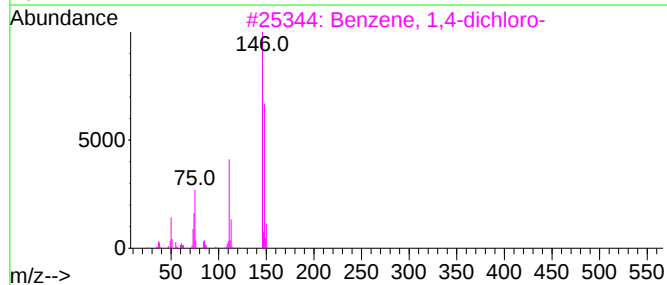
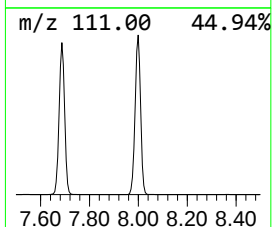
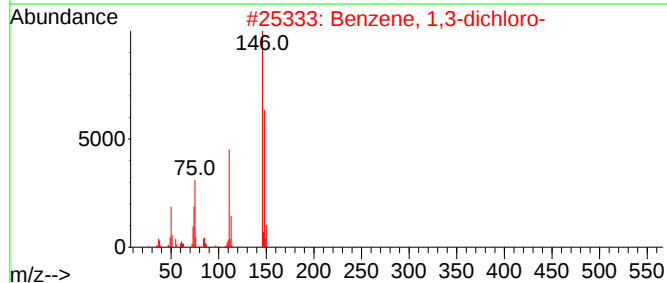
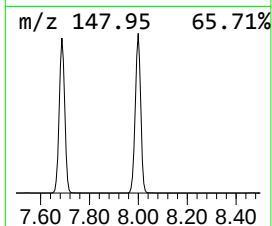
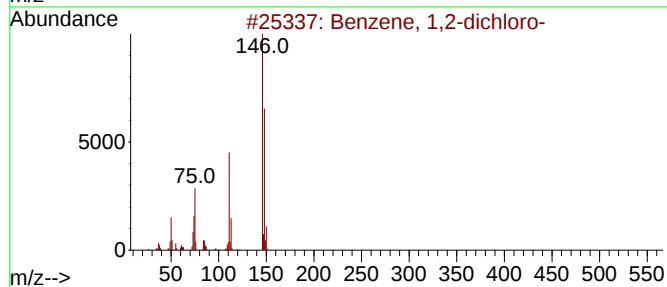
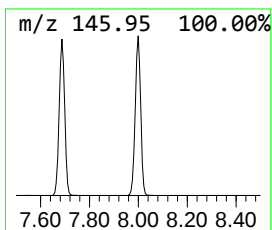
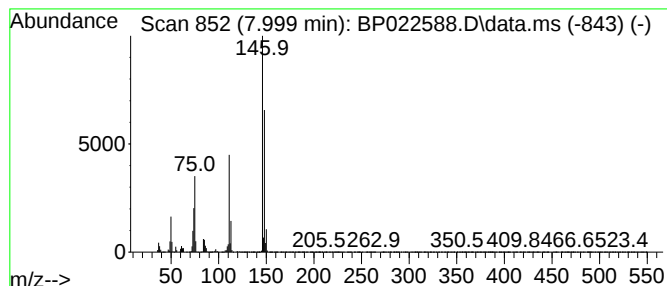
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.999	18.39 ng/ul	4885490	1,4-Dichlorobenzene-d4	7.652

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, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	97
5			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	97



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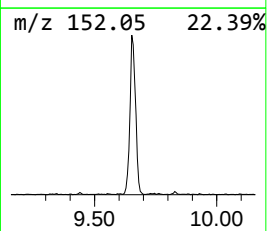
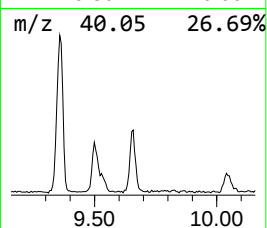
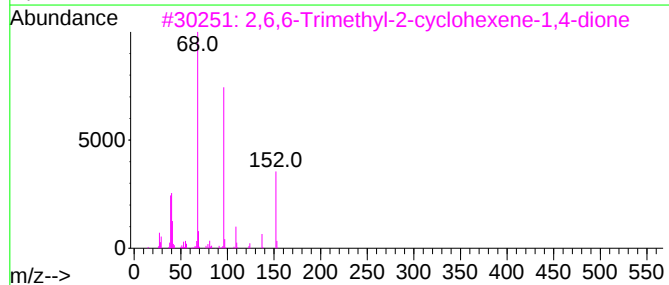
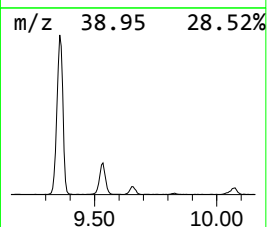
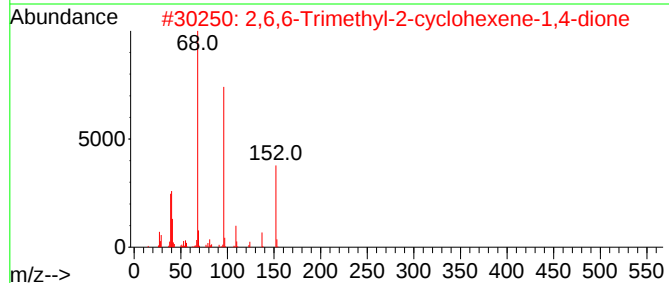
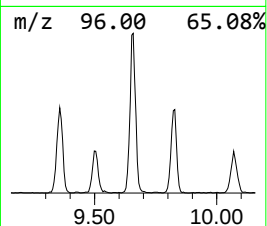
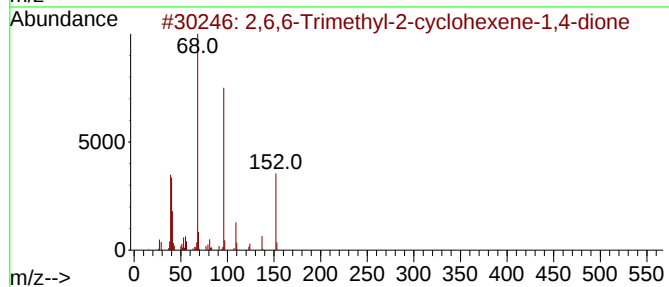
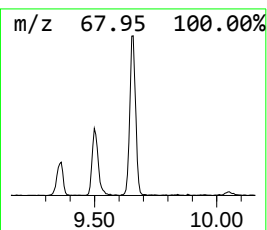
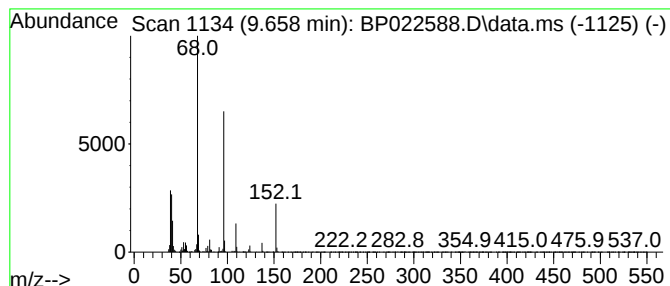
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.658	4.20 ng/ul	170872	Naphthalene-d8	10.411

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,6,6-Trimethyl-2-cyclohexene-1,...	152	C9H12O2	001125-21-9	95		
2	2,6,6-Trimethyl-2-cyclohexene-1,...	152	C9H12O2	001125-21-9	91		
3	2,6,6-Trimethyl-2-cyclohexene-1,...	152	C9H12O2	001125-21-9	91		
4	2,6,6-Trimethyl-2-cyclohexene-1,...	152	C9H12O2	001125-21-9	87		
5	2-Cyclohexen-1-one	96	C6H8O	000930-68-7	53		



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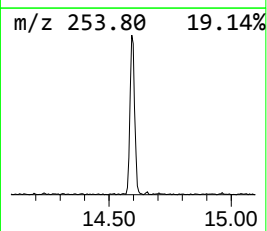
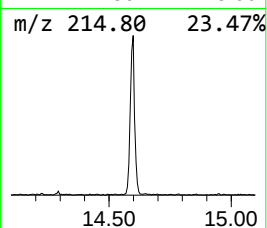
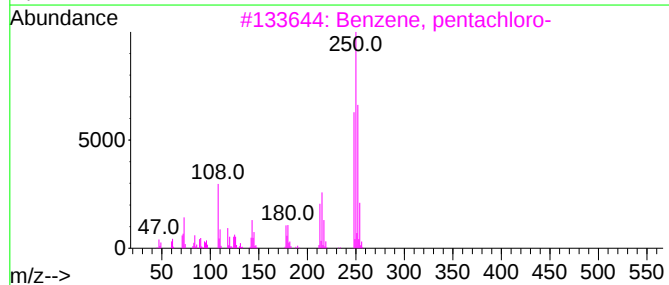
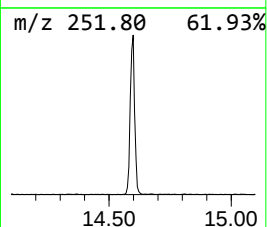
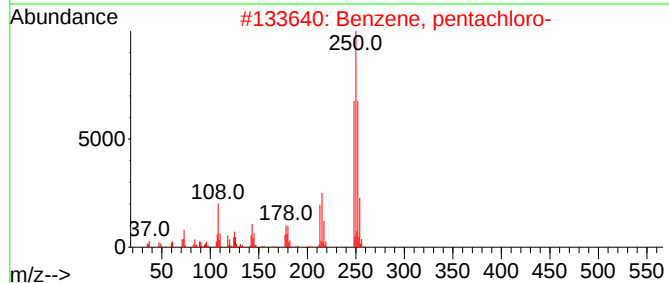
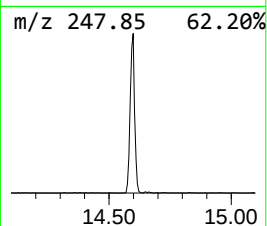
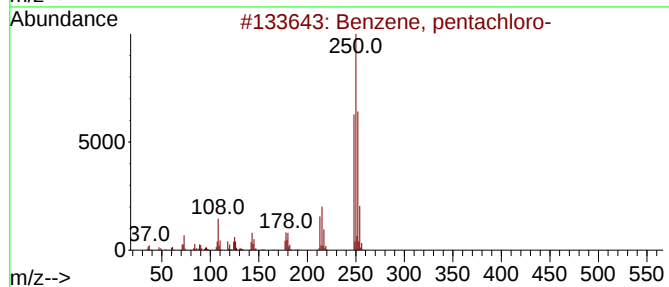
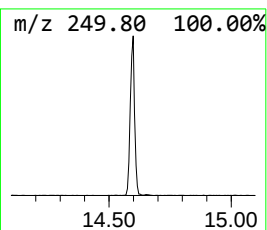
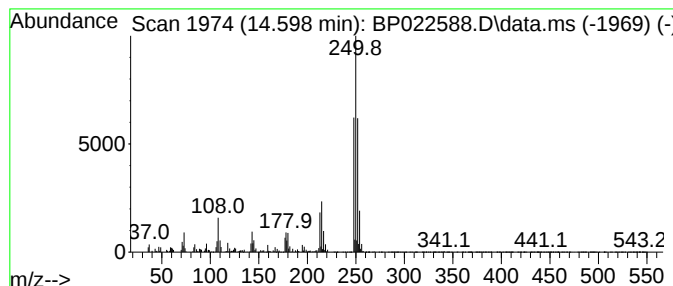
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Benzene, pentachloro- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.599	2.53 ng/ul	143064	Acenaphthene-d10	14.275

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	99
4			Benzene, pentachloro-	248	C6HCl5	000608-93-5	96
5			Benzene, pentachloro-	248	C6HCl5	000608-93-5	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102424\
Data File : BP022588.D
Acq On : 24 Oct 2024 15:12
Operator : RC/JU
Sample : P4436-02
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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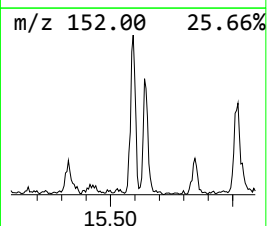
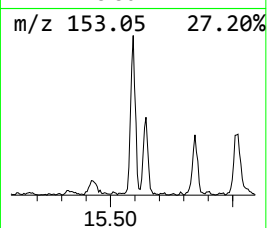
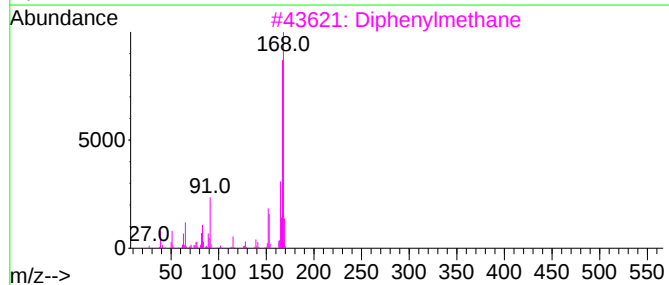
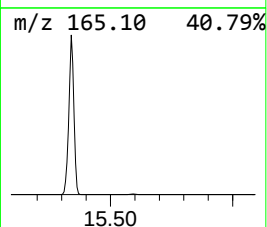
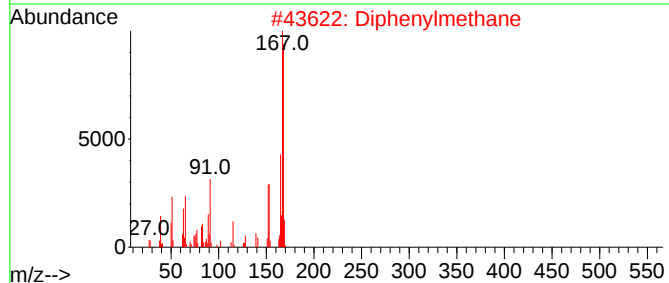
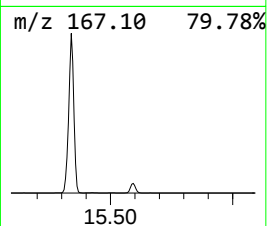
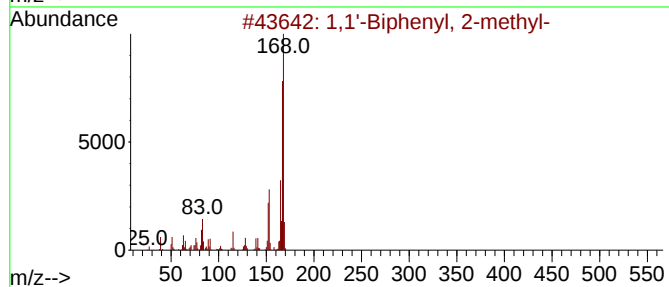
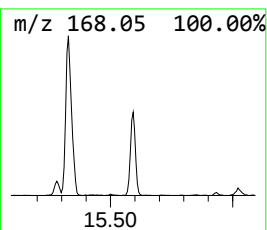
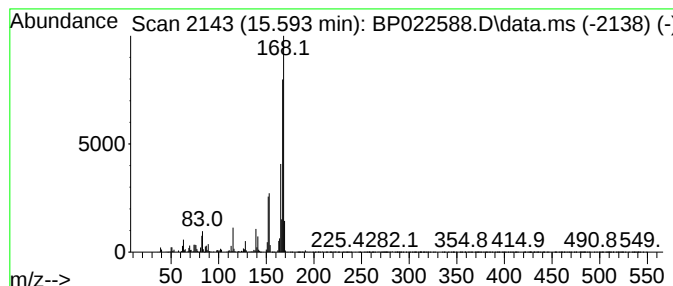
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 1,1'-Biphenyl, 2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.592	2.68 ng/ul	151684	Acenaphthene-d10	14.275

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,1'-Biphenyl, 2-methyl-	168	C13H12	000643-58-3	94
2			Diphenylmethane	168	C13H12	000101-81-5	91
3			Diphenylmethane	168	C13H12	000101-81-5	91
4			Diphenylmethane	168	C13H12	000101-81-5	90
5			1,1'-Biphenyl, 2-methyl-	168	C13H12	000643-58-3	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102424\
Data File : BP022588.D
Acq On : 24 Oct 2024 15:12
Operator : RC/JU
Sample : P4436-02
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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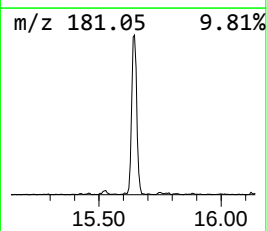
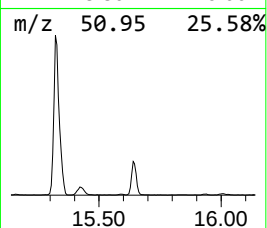
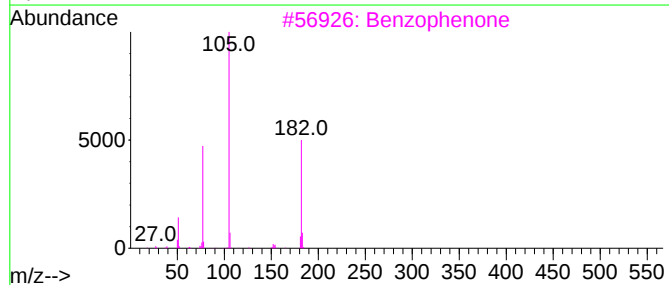
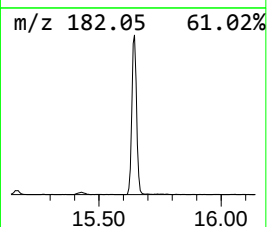
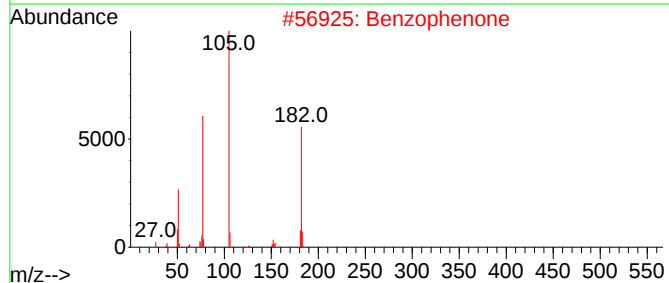
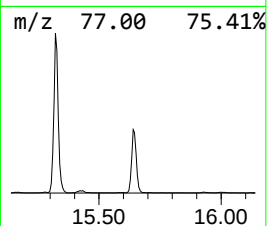
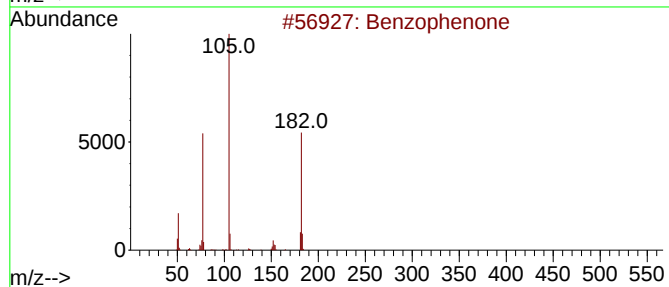
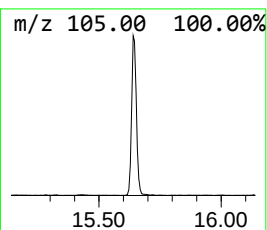
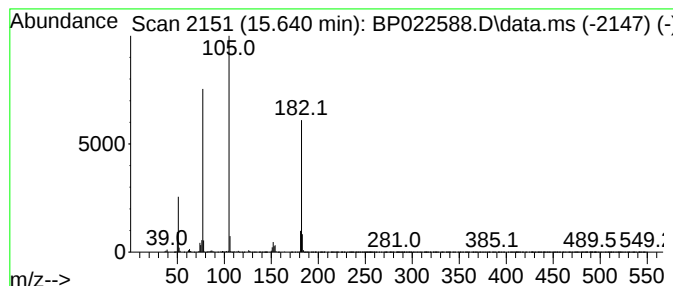
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Benzophenone Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.640	7.83 ng/ul	443504	Acenaphthene-d10	14.275

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	91
4			Benzophenone	182	C13H10O	000119-61-9	81
5			1,2,4-Trioxolane, 3,3,5-triphenyl-	304	C20H16O3	023246-12-0	78



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102424\
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Acq On : 24 Oct 2024 15:12
Operator : RC/JU
Sample : P4436-02
Misc :
ALS Vial : 8 Sample Multiplier: 1

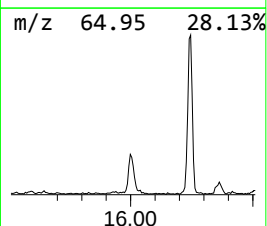
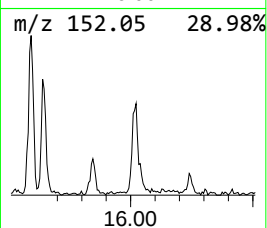
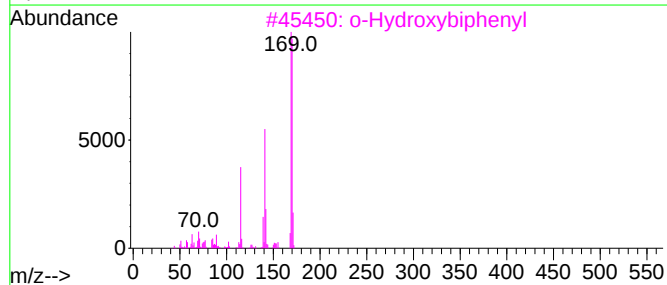
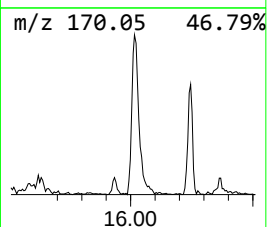
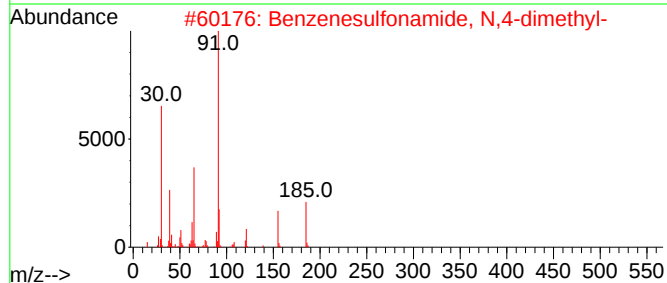
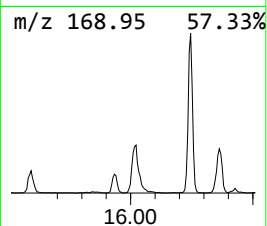
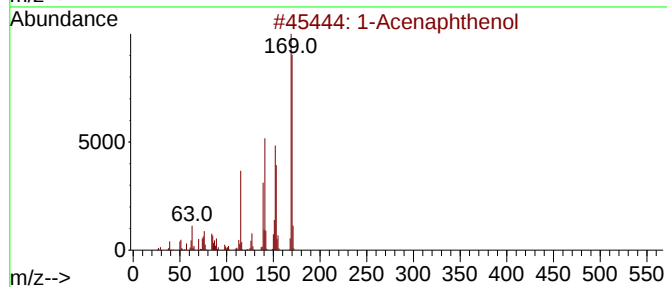
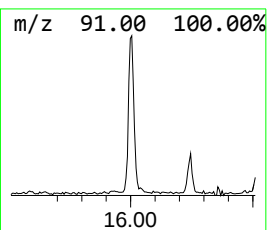
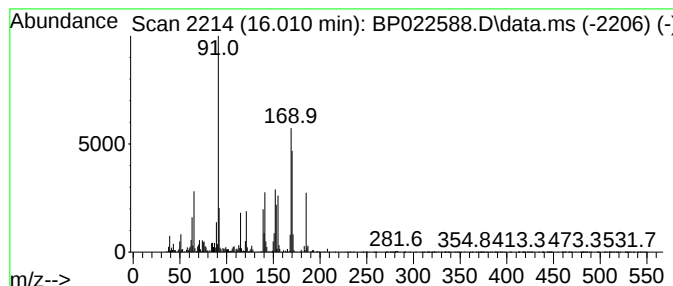
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 1-Acenaphthenol Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD		R.T.	
16.010	2.66 ng/ul	195404	Phenanthrene-d10		17.069	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	1-Acenaphthenol		170	C12H10O	006306-07-6	86
2	Benzenesulfonamide, N,4-dimethyl-		185	C8H11NO2S	000640-61-9	44
3	o-Hydroxybiphenyl		170	C12H10O	000090-43-7	38
4	Benzenesulfonamide, N,4-dimethyl-		185	C8H11NO2S	000640-61-9	35
5	1,8-Naphthalenedimethanol		188	C12H12O2	002026-08-6	32



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102424\
Data File : BP022588.D
Acq On : 24 Oct 2024 15:12
Operator : RC/JU
Sample : P4436-02
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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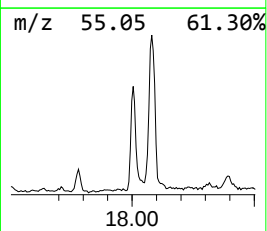
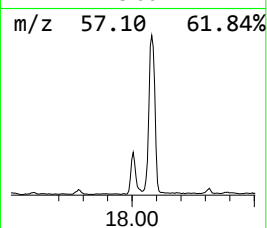
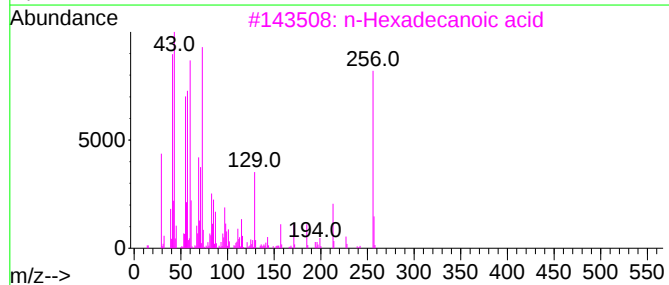
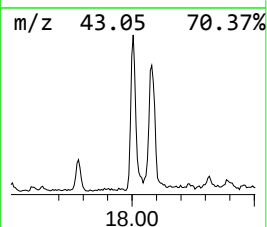
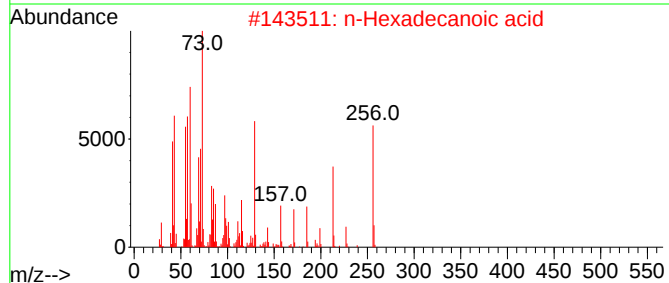
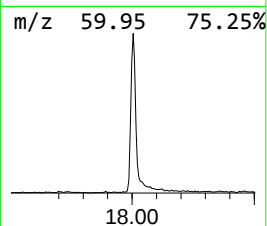
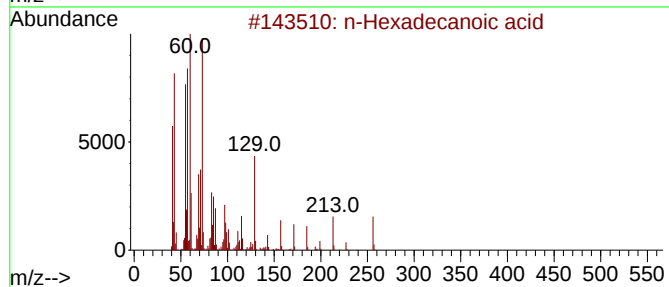
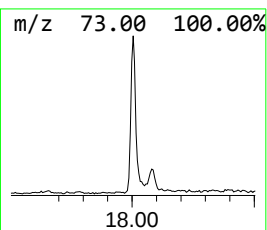
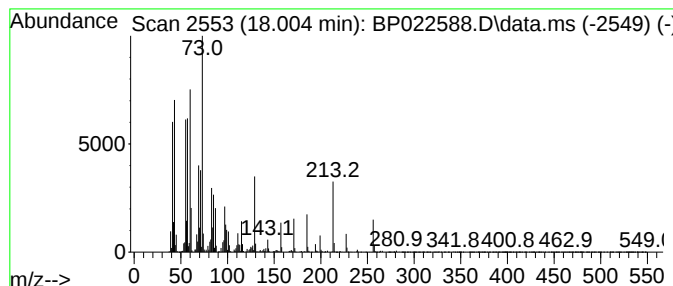
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.004	8.20 ng/ul	603164	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	99
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
5			Tridecanoic acid	214	C13H26O2	000638-53-9	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102424\
Data File : BP022588.D
Acq On : 24 Oct 2024 15:12
Operator : RC/JU
Sample : P4436-02
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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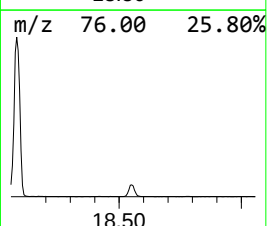
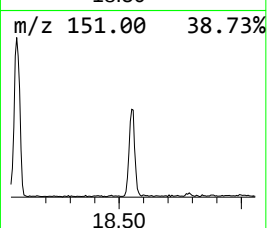
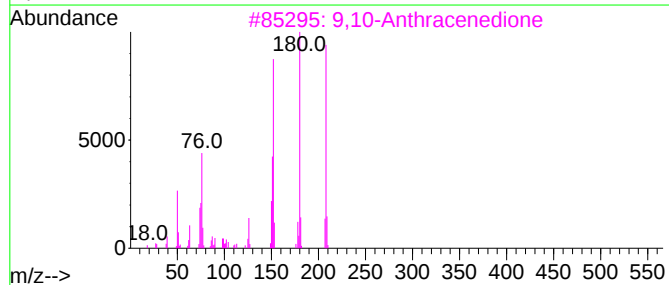
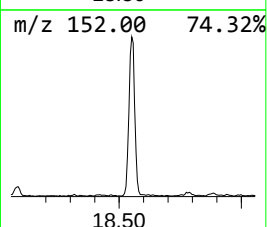
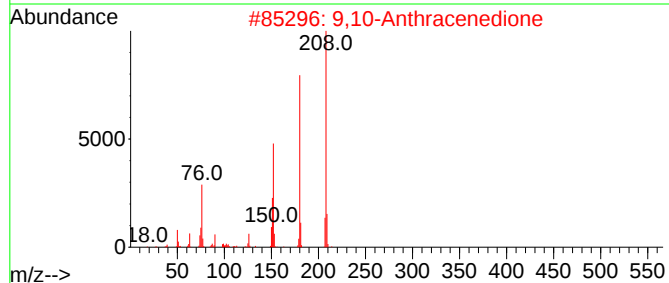
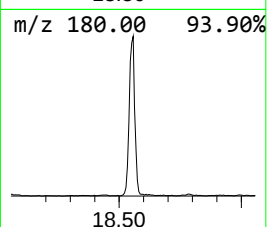
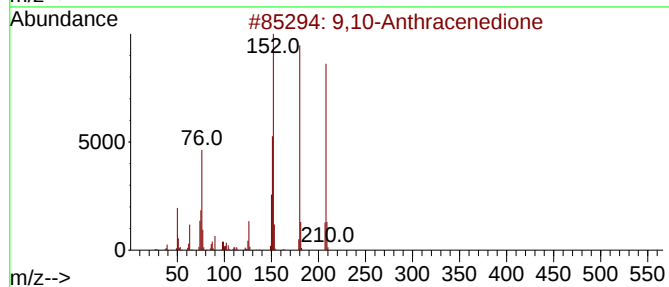
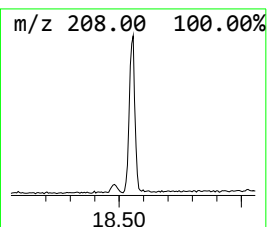
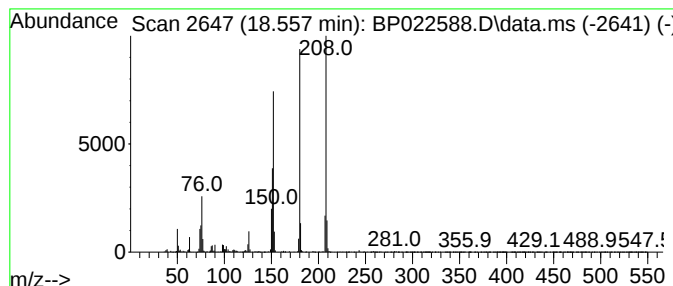
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 9,10-Anthracenedione Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.557	5.43 ng/ul	399404	Phenanthrene-d10	17.069

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,10-Anthracenedione			208	C14H8O2	000084-65-1	98
2	9,10-Anthracenedione			208	C14H8O2	000084-65-1	97
3	9,10-Anthracenedione			208	C14H8O2	000084-65-1	97
4	9,10-Anthracenedione			208	C14H8O2	000084-65-1	95
5	Benzo[c]cinoline			180	C12H8N2	000230-17-1	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102424\
Data File : BP022588.D
Acq On : 24 Oct 2024 15:12
Operator : RC/JU
Sample : P4436-02
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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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\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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2,6,6-Trimethyl...	9.658	4.2	ng/ul	170872	2	10.411	814358	20.0
Benzene, pentac...	14.599	2.5	ng/ul	143064	3	14.275	1132640	20.0
1,1'-Biphenyl, ...	15.592	2.7	ng/ul	151684	3	14.275	1132640	20.0
Benzophenone	15.640	7.8	ng/ul	443504	3	14.275	1132640	20.0
1-Acenaphthenol	16.010	2.7	ng/ul	195404	4	17.069	1471580	20.0
n-Hexadecanoic ...	18.004	8.2	ng/ul	603164	4	17.069	1471580	20.0
9,10-Anthracene...	18.557	5.4	ng/ul	399404	4	17.069	1471580	20.0