

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
 Data File : BP015549.D
 Acq On : 15 Jun 2023 10:36
 Operator : MA/JU
 Sample : 03088-15
 Misc :
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCGV6

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-BP060723.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP015549.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.404	33	37	39	rBV	30217	39877	0.46%	0.026%
2	3.440	39	43	53	rVB	150590	219491	2.52%	0.143%
3	5.140	327	332	342	rBV	259981	380006	4.37%	0.248%
4	7.222	677	686	691	rBV2	1021701	2566125	29.50%	1.676%
5	7.269	691	694	708	rVV	1000775	1577083	18.13%	1.030%
6	7.410	711	718	728	rVB	502755	781073	8.98%	0.510%
7	7.610	744	752	755	rBV	689555	1130697	13.00%	0.738%
8	7.645	755	758	769	rVB	620765	941460	10.82%	0.615%
9	8.087	826	833	842	rBV	510379	821004	9.44%	0.536%
10	8.804	948	955	965	rBV	866117	1440399	16.56%	0.941%
11	8.904	965	972	984	rVB	730851	1167810	13.42%	0.763%
12	9.263	1026	1033	1037	rBV	674389	1147240	13.19%	0.749%
13	9.310	1037	1041	1052	rVV	553641	917760	10.55%	0.599%
14	9.416	1054	1059	1067	rVB	6443	14127	0.16%	0.009%
15	9.651	1092	1099	1105	rVB	14606	24840	0.29%	0.016%
16	9.992	1149	1157	1160	rBV	626061	1096621	12.61%	0.716%
17	10.022	1160	1162	1180	rVB	554352	858594	9.87%	0.561%
18	10.534	1241	1249	1251	rBV	925524	1464701	16.84%	0.957%
19	10.875	1300	1307	1308	rBV2	30634	48240	0.55%	0.032%
20	10.916	1308	1314	1318	rVV	794249	1327315	15.26%	0.867%
21	10.969	1318	1323	1330	rVV	791574	1309713	15.06%	0.855%
22	11.057	1330	1338	1360	rVV	722077	1430720	16.45%	0.934%
23	11.245	1362	1370	1379	rVB	923474	1489165	17.12%	0.972%
24	12.198	1525	1532	1547	rBV	1596877	2737358	31.47%	1.788%
25	12.433	1564	1572	1578	rBV2	38270	63690	0.73%	0.042%
26	12.498	1579	1583	1590	rVB	17973	29712	0.34%	0.019%
27	12.786	1625	1632	1643	rVV	1425403	2147677	24.69%	1.403%
28	12.933	1647	1657	1668	rVV	1009316	1516817	17.44%	0.991%
29	13.175	1691	1698	1704	rBV	1647285	2547699	29.29%	1.664%
30	13.228	1704	1707	1721	rVV	55239	118568	1.36%	0.077%
31	13.569	1754	1765	1769	rBV	1863305	2839644	32.64%	1.854%
32	13.616	1769	1773	1783	rVB	1029604	1494614	17.18%	0.976%
33	14.139	1856	1862	1866	rBV	1827441	2532369	29.11%	1.654%
34	14.186	1866	1870	1877	rVB	1156328	1582106	18.19%	1.033%
35	14.251	1878	1881	1886	rVB6	14933	20457	0.24%	0.013%
36	14.316	1886	1892	1905	rBV	1146998	1572170	18.07%	1.027%

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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-BP060723.MA.M
 Title : SVOA CALIBRATION

37	14.433	1905	1912	1926	rVB	1899140	2867513	32.96%	1.873%
38	14.604	1935	1941	1947	rVB6	8430	15282	0.18%	0.010%
39	14.733	1954	1963	1969	rBV	1250749	1889619	21.72%	1.234%
40	14.798	1969	1974	1984	rVB	1707283	2421384	27.84%	1.581%
41	14.945	1992	1999	2023	rBV2	1006742	2448798	28.15%	1.599%
42	15.133	2025	2031	2045	rVB	1339801	1914398	22.01%	1.250%
43	15.363	2063	2070	2089	rBV	1233424	1821129	20.94%	1.189%
44	15.563	2099	2104	2109	rBV2	39073	50246	0.58%	0.033%
45	15.727	2126	2132	2136	rBV	2544273	3579182	41.15%	2.337%
46	15.774	2136	2140	2147	rVB	1171903	1644305	18.90%	1.074%
47	15.869	2147	2156	2166	rBV3	1862231	3500658	40.24%	2.286%
48	16.045	2182	2186	2189	rBV2	27997	46439	0.53%	0.030%
49	16.092	2189	2194	2202	rVB	1074156	1438547	16.54%	0.939%
50	16.463	2253	2257	2262	rVB4	15140	24444	0.28%	0.016%
51	16.657	2285	2290	2295	rBV	377205	510756	5.87%	0.334%
52	16.704	2295	2298	2302	rVV	20466	33238	0.38%	0.022%
53	16.792	2306	2313	2319	rVV	1589343	2282396	26.24%	1.491%
54	16.839	2319	2321	2324	rVV4	23571	36728	0.42%	0.024%
55	16.868	2324	2326	2331	rVV3	27741	46166	0.53%	0.030%
56	16.927	2334	2336	2344	rVV4	14308	21374	0.25%	0.014%
57	17.021	2346	2352	2356	rBV6	11530	16886	0.19%	0.011%
58	17.139	2364	2372	2383	rBV	2026011	2935682	33.75%	1.917%
59	17.221	2384	2386	2391	rVB2	16322	21474	0.25%	0.014%
60	17.321	2398	2403	2409	rVB	148163	176290	2.03%	0.115%
61	17.492	2424	2432	2436	rBV	1680580	2478309	28.49%	1.618%
62	17.539	2436	2440	2444	rVV	1805595	2515151	28.91%	1.643%
63	17.598	2444	2450	2453	rVV	2788759	4303751	49.47%	2.811%
64	17.627	2453	2455	2466	rVB	2137918	2626058	30.19%	1.715%
65	17.763	2475	2478	2488	rVB	10552	23749	0.27%	0.016%
66	17.898	2490	2501	2507	rBV	1920131	2668602	30.68%	1.743%
67	17.968	2507	2513	2524	rVB	1985884	2558765	29.41%	1.671%
68	18.098	2531	2535	2538	rBV5	19109	21970	0.25%	0.014%
69	18.139	2538	2542	2547	rVB7	15689	24039	0.28%	0.016%
70	18.233	2554	2558	2563	rBV8	12025	21635	0.25%	0.014%
71	18.357	2573	2579	2586	rBV	973689	1270902	14.61%	0.830%
72	18.427	2586	2591	2599	rVB	1951195	2295135	26.38%	1.499%
73	18.633	2622	2626	2627	rBV2	17899	22663	0.26%	0.015%
74	18.733	2639	2643	2650	rBV7	17061	44497	0.51%	0.029%
75	18.886	2666	2669	2671	rBV4	19177	21951	0.25%	0.014%
76	19.186	2716	2720	2725	rBV7	31523	62938	0.72%	0.041%

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Peak Location: TOP

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77	19.233	2726	2728	2734	rVB2	27988	34437	0.40%	0.022%
78	19.392	2752	2755	2760	rVB4	45424	56518	0.65%	0.037%
79	19.439	2760	2763	2765	rBV2	62001	79936	0.92%	0.052%
80	19.492	2765	2772	2780	rVB	4803229	6579492	75.64%	4.297%
81	19.580	2782	2787	2800	rBV	399666	591331	6.80%	0.386%
82	19.733	2809	2813	2817	rVB	36199	48649	0.56%	0.032%
83	19.821	2820	2828	2830	rVV	3800389	5280852	60.71%	3.449%
84	19.845	2830	2832	2842	rVB	3508723	4394691	50.52%	2.870%
85	20.309	2905	2911	2915	rBV	117000	168883	1.94%	0.110%
86	20.639	2965	2967	2971	rVB	45670	45576	0.52%	0.030%
87	20.698	2973	2977	2982	rVV	2122714	2397029	27.56%	1.565%
88	20.792	2990	2993	2997	rBV	100366	102752	1.18%	0.067%
89	21.251	3067	3071	3086	rBV	860872	1574050	18.09%	1.028%
90	21.456	3101	3106	3111	rBV	3393598	3602881	41.42%	2.353%
91	21.562	3118	3124	3130	rBV2	4827386	8698871	100.00%	5.681%
92	21.621	3130	3134	3139	rVB	3091529	4188583	48.15%	2.735%
93	22.709	3316	3319	3325	rVB2	125461	195110	2.24%	0.127%
94	23.303	3416	3420	3421	rBV	155454	217417	2.50%	0.142%
95	23.345	3421	3427	3438	rVB	2326553	4382688	50.38%	2.862%
96	23.968	3525	3533	3537	rBV2	2427058	5109816	58.74%	3.337%
97	24.015	3537	3541	3550	rVV	1854077	3726827	42.84%	2.434%
98	24.133	3554	3561	3571	rVB	1151927	2519786	28.97%	1.646%
99	24.756	3662	3667	3675	rVB	242188	514525	5.91%	0.336%
100	26.839	4011	4021	4034	rVB2	1918017	6518521	74.94%	4.257%

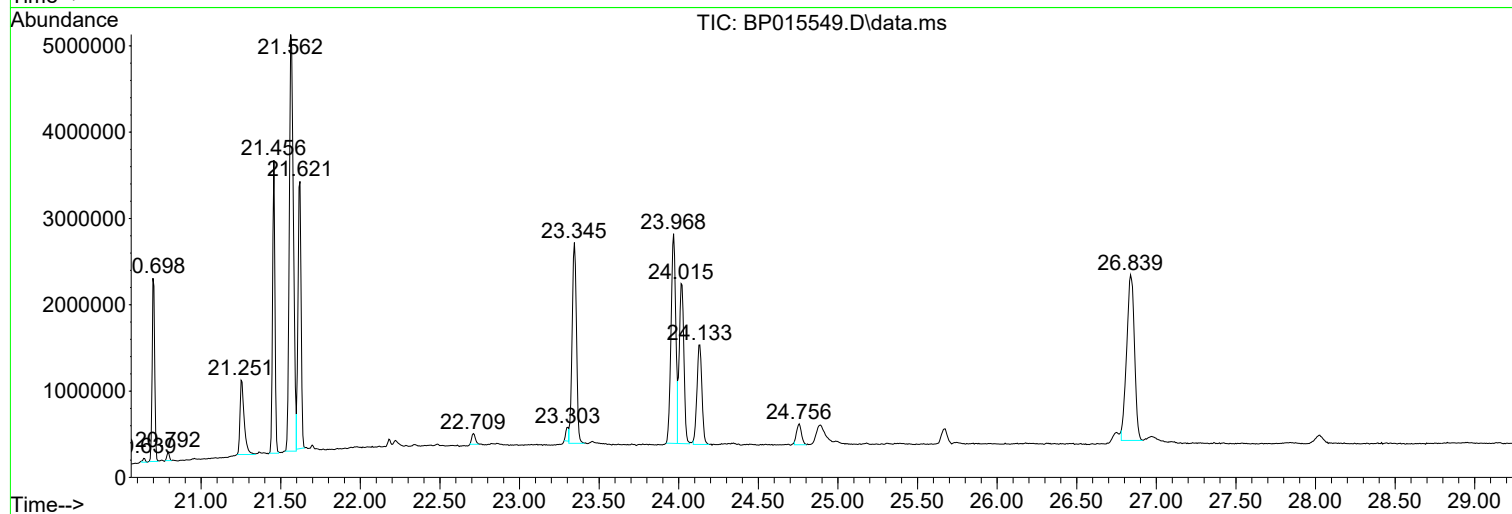
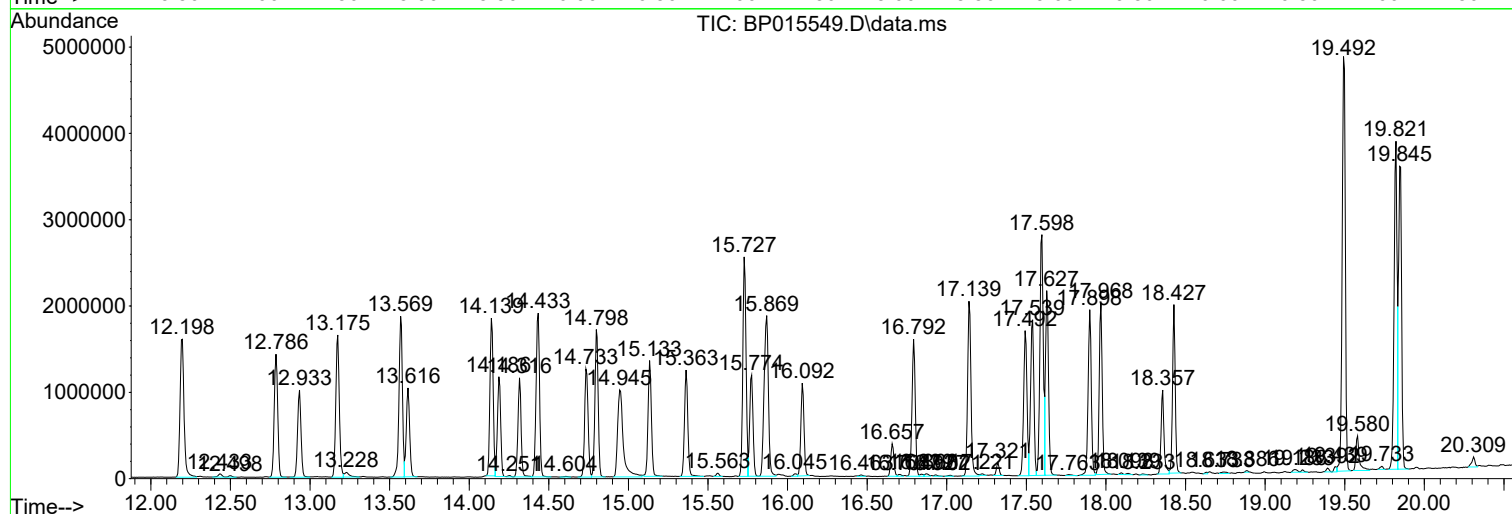
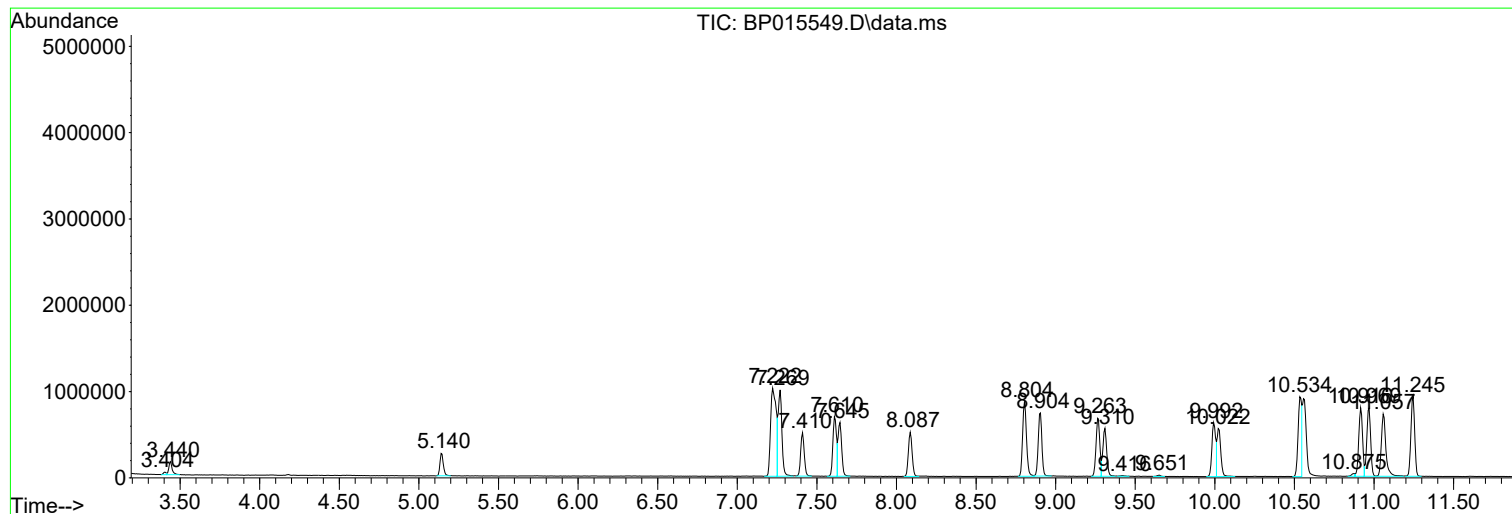
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

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



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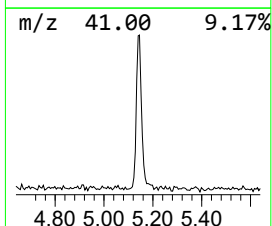
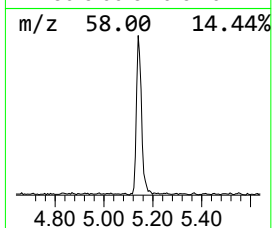
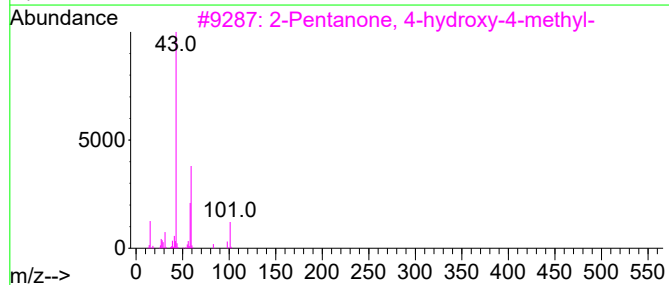
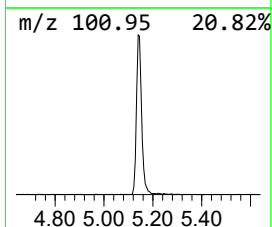
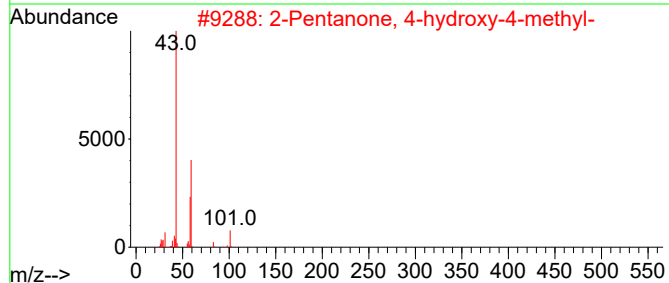
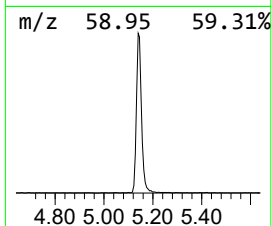
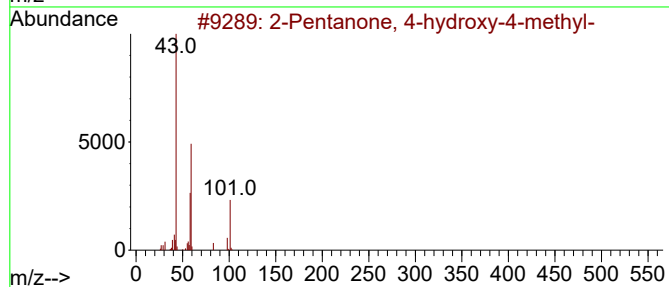
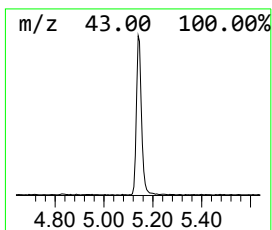
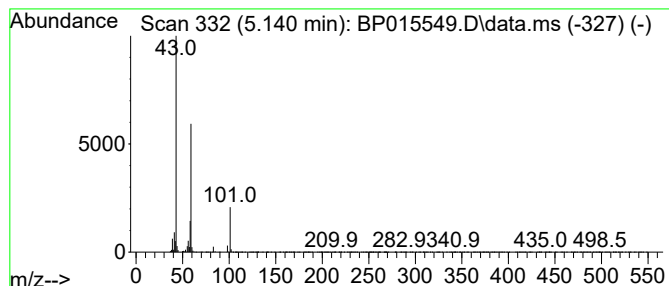
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.140	9.26 ng/ul	380006	1,4-Dichlorobenzene-d4	8.087

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56		
2	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	40		
3	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	40		
4	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	35		
5	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	32		



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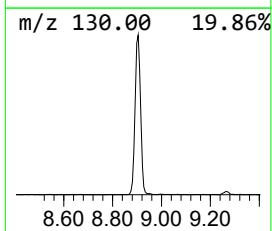
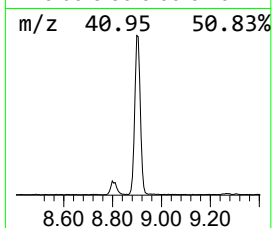
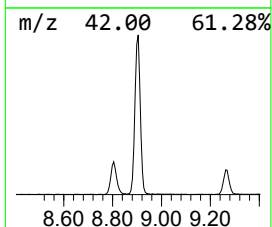
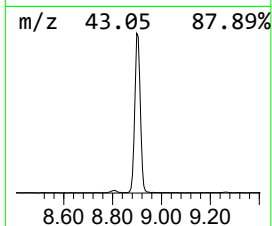
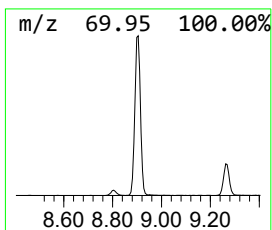
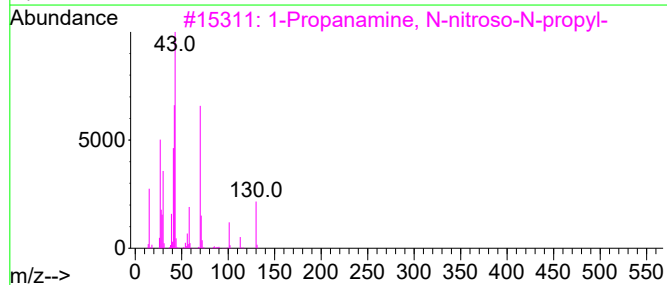
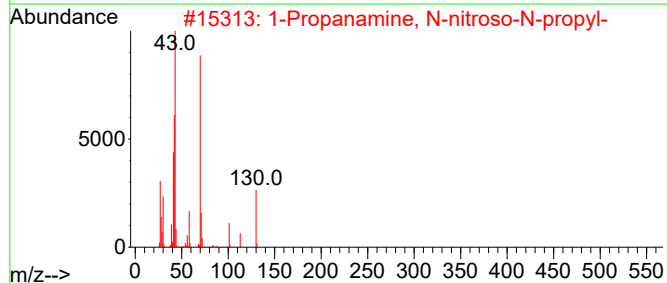
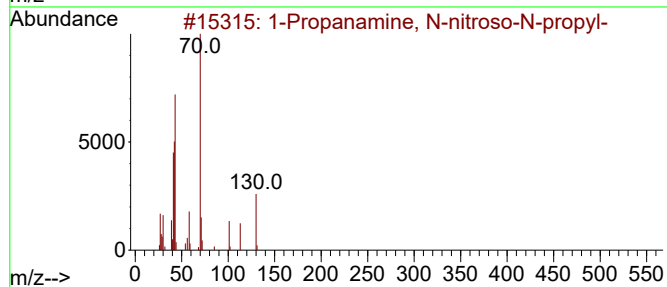
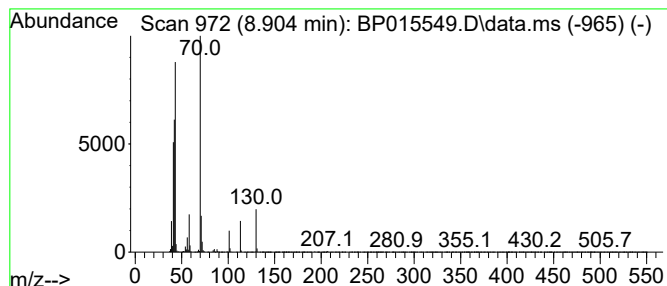
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 1-Propanamine, N-nitroso-N-... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.904	28.45 ng/ul	1167810	1,4-Dichlorobenzene-d4	8.087

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Propanamine, N-nitroso-N-propyl-	130	C6H14N2O	000621-64-7	81
2			1-Propanamine, N-nitroso-N-propyl-	130	C6H14N2O	000621-64-7	76
3			1-Propanamine, N-nitroso-N-propyl-	130	C6H14N2O	000621-64-7	76
4			Oxazolidine, 2,2-dimethyl-	101	C5H11NO	020515-62-2	47
5			Pentane, 2-chloro-	106	C5H11Cl	000625-29-6	25



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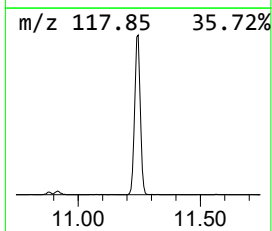
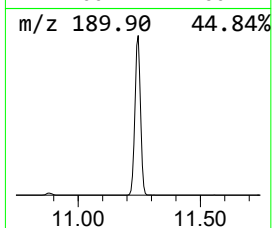
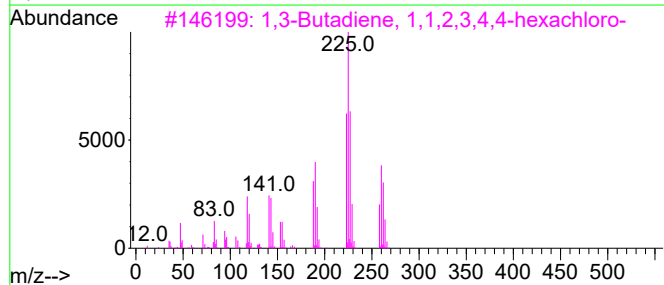
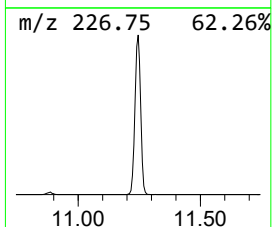
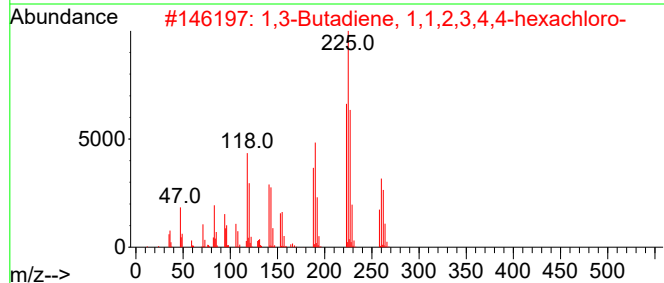
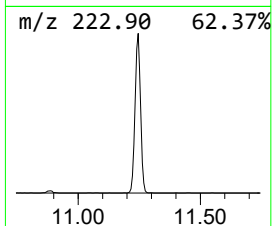
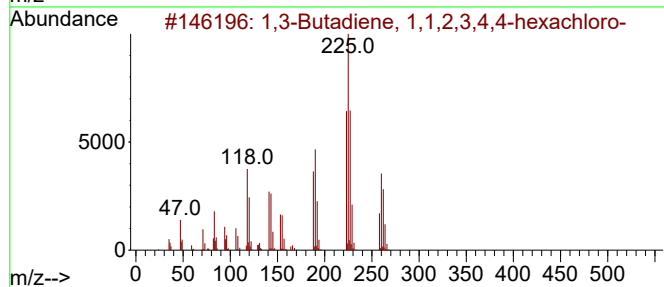
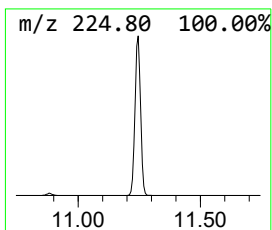
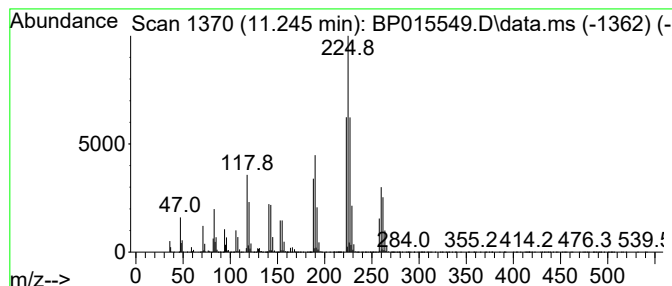
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.245	22.44 ng/ul	1489170	Naphthalene-d8	10.916

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Butadiene, 1,1,2,3,4,4-hexac...	258	C4Cl6	000087-68-3	99		
2	1,3-Butadiene, 1,1,2,3,4,4-hexac...	258	C4Cl6	000087-68-3	99		
3	1,3-Butadiene, 1,1,2,3,4,4-hexac...	258	C4Cl6	000087-68-3	99		
4	1,3-Butadiene, 1,1,2,3,4,4-hexac...	258	C4Cl6	000087-68-3	95		
5	Tin(IV) chloride	260	Cl4Sn	007646-78-8	17		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
Data File : BP015549.D
Acq On : 15 Jun 2023 10:36
Operator : MA/JU
Sample : 03088-15
Misc :
ALS Vial : 44 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCGV6

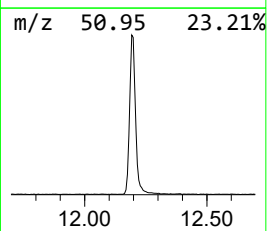
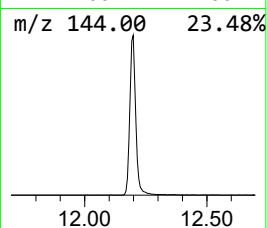
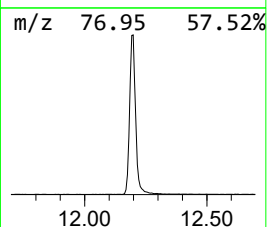
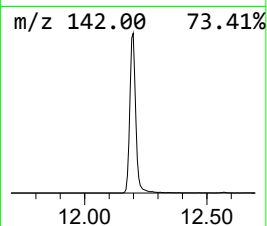
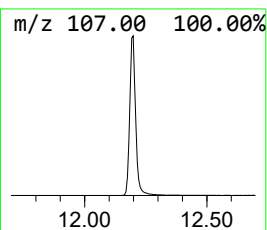
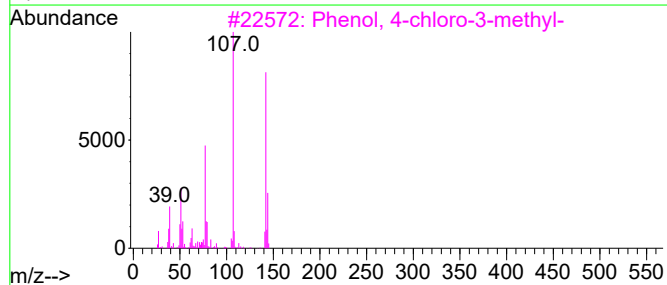
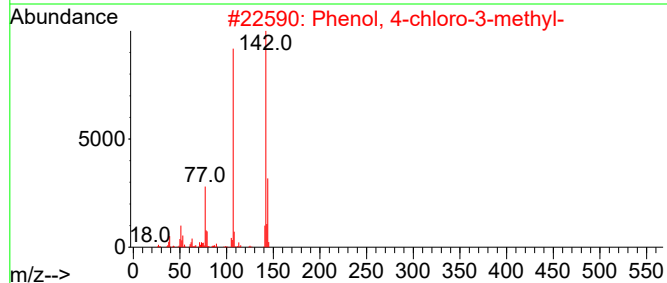
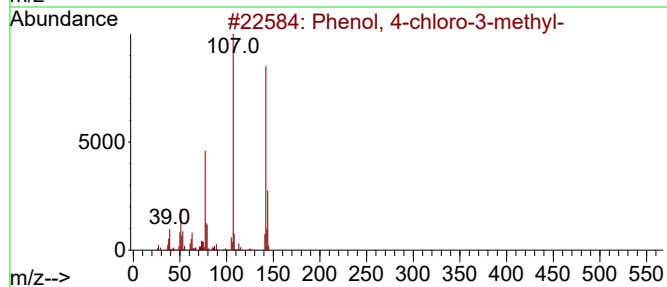
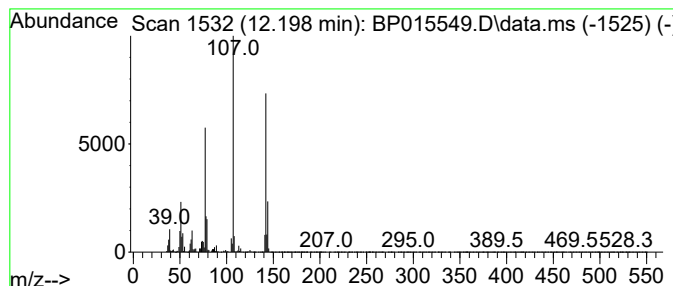
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Phenol, 4-chloro-3-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.198	41.25 ng/ul	2737360	Naphthalene-d8	10.916

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 4-chloro-3-methyl-	142	C7H7ClO	000059-50-7	97
2			Phenol, 4-chloro-3-methyl-	142	C7H7ClO	000059-50-7	97
3			Phenol, 4-chloro-3-methyl-	142	C7H7ClO	000059-50-7	96
4			Phenol, 4-chloro-3-methyl-	142	C7H7ClO	000059-50-7	94
5			Phenol, 4-chloro-3-methyl-	142	C7H7ClO	000059-50-7	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
Data File : BP015549.D
Acq On : 15 Jun 2023 10:36
Operator : MA/JU
Sample : 03088-15
Misc :
ALS Vial : 44 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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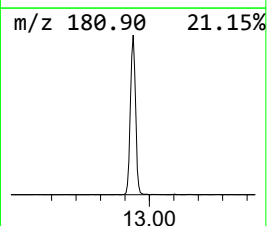
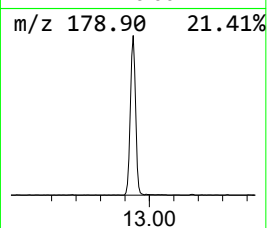
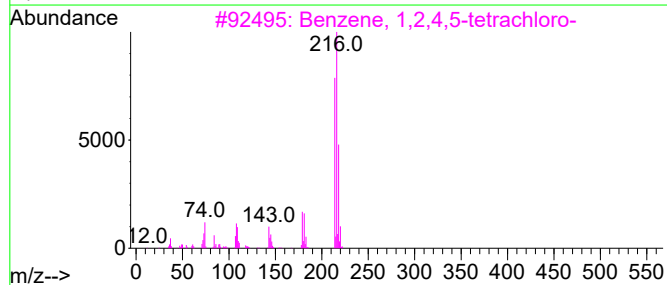
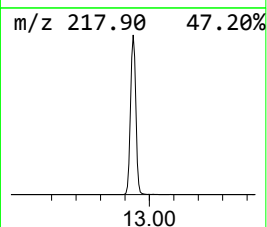
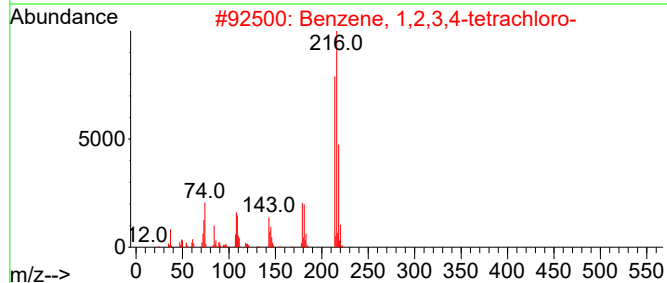
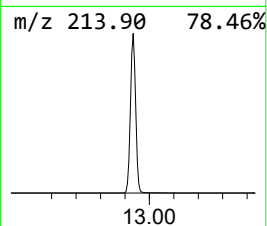
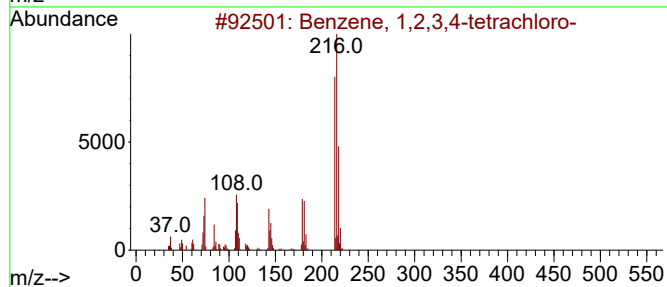
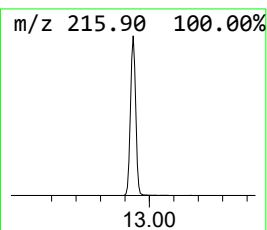
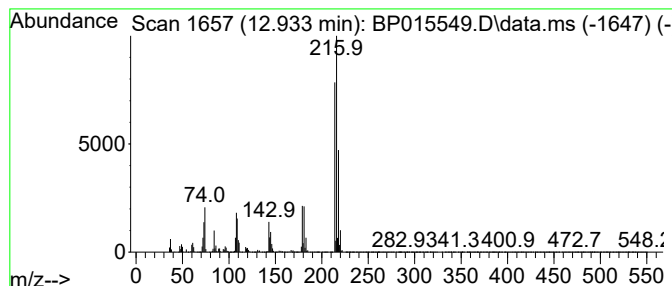
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Benzene, 1,2,3,4-tetrachloro- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.933	16.05 ng/ul	1516820	Acenaphthene-d10	14.733

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3,4-tetrachloro-	214	C6H2Cl4	000634-66-2	99
2			Benzene, 1,2,3,4-tetrachloro-	214	C6H2Cl4	000634-66-2	99
3			Benzene, 1,2,4,5-tetrachloro-	214	C6H2Cl4	000095-94-3	99
4			Benzene, 1,2,4,5-tetrachloro-	214	C6H2Cl4	000095-94-3	99
5			Benzene, 1,2,4,5-tetrachloro-	214	C6H2Cl4	000095-94-3	99



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
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Sample : 03088-15
Misc :
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ClientSampleId :
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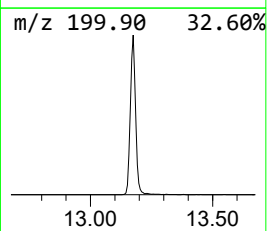
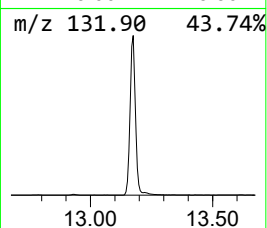
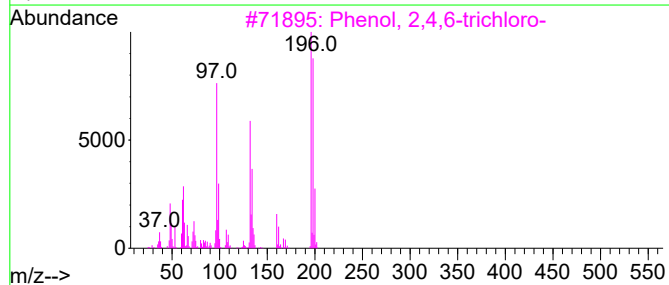
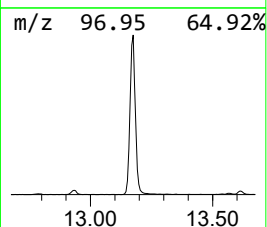
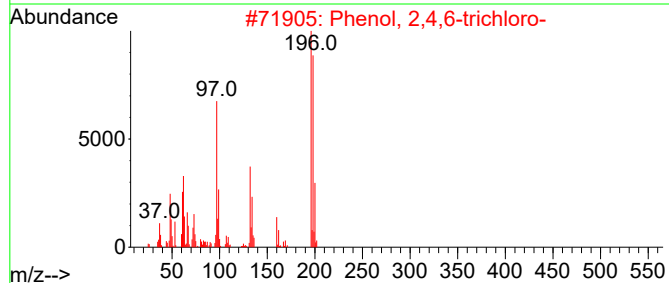
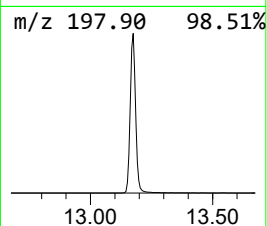
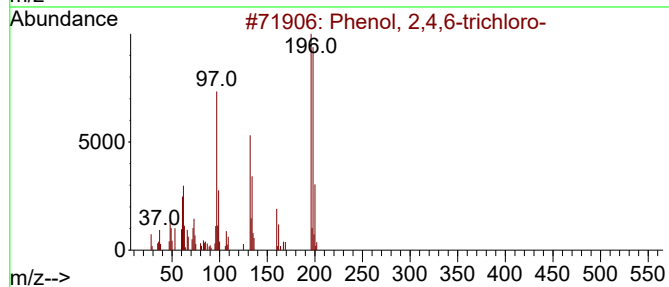
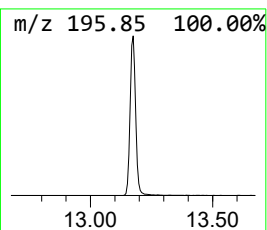
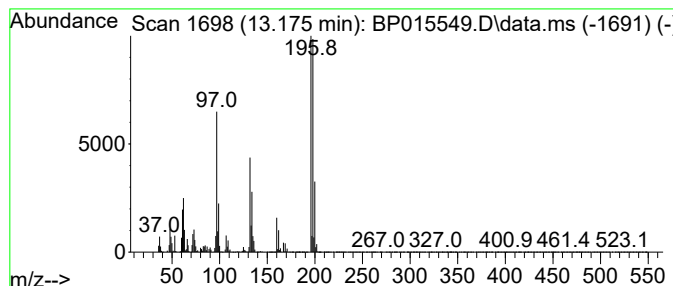
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Phenol, 2,4,6-trichloro- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.175	26.97 ng/ul	2547700	Acenaphthene-d10	14.733

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 2,4,6-trichloro-	196	C6H3Cl3O	000088-06-2	99
2			Phenol, 2,4,6-trichloro-	196	C6H3Cl3O	000088-06-2	95
3			Phenol, 2,4,6-trichloro-	196	C6H3Cl3O	000088-06-2	94
4			Phenol, 2,4,5-trichloro-	196	C6H3Cl3O	000095-95-4	94
5			Phenol, 2,4,6-trichloro-	196	C6H3Cl3O	000088-06-2	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
Data File : BP015549.D
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Sample : 03088-15
Misc :
ALS Vial : 44 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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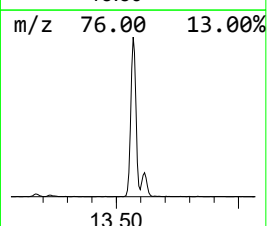
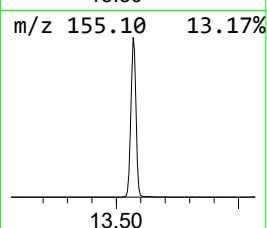
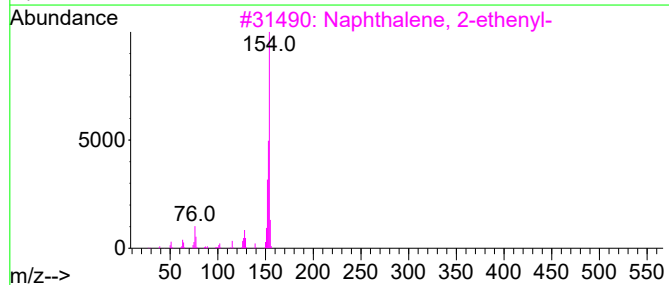
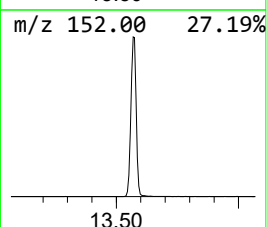
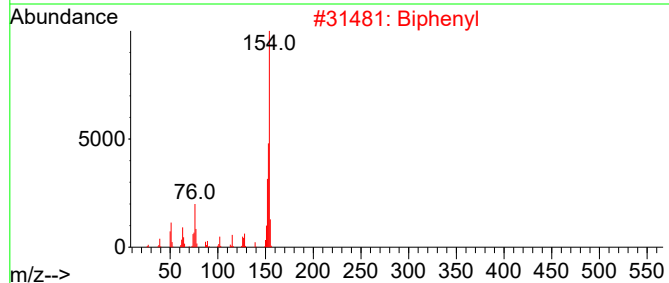
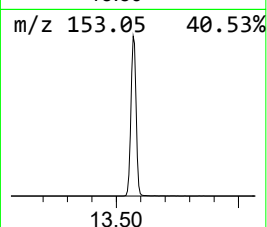
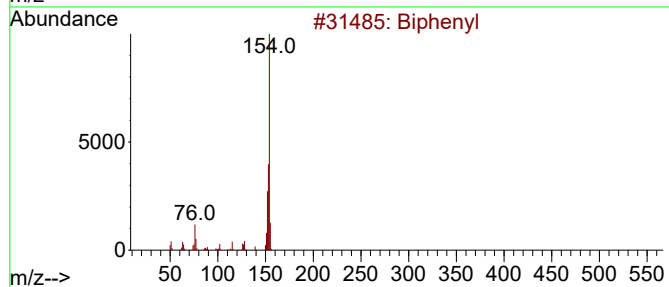
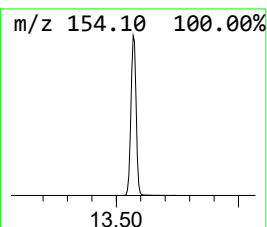
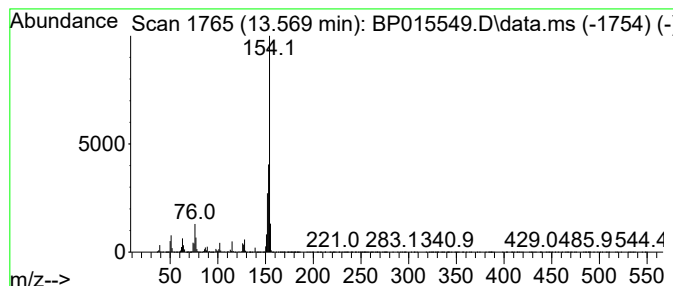
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Biphenyl Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.569	30.06 ng/ul	2839640	Acenaphthene-d10	14.733

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Biphenyl	154	C12H10	000092-52-4	95
2			Biphenyl	154	C12H10	000092-52-4	93
3			Naphthalene, 2-ethenyl-	154	C12H10	000827-54-3	93
4			Biphenyl	154	C12H10	000092-52-4	93
5			Biphenyl	154	C12H10	000092-52-4	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
Data File : BP015549.D
Acq On : 15 Jun 2023 10:36
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Sample : 03088-15
Misc :
ALS Vial : 44 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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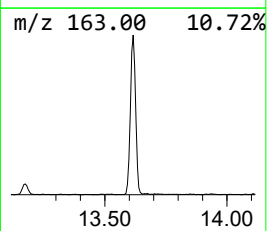
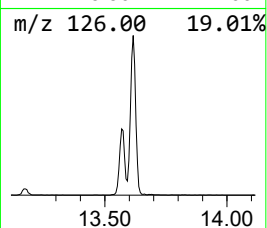
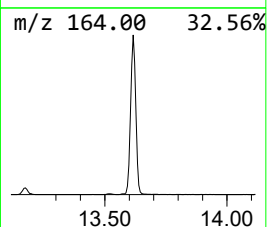
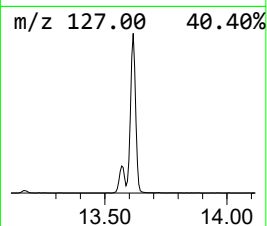
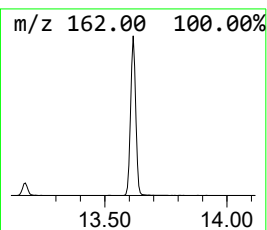
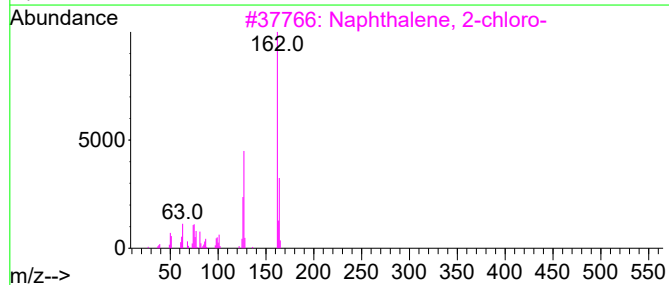
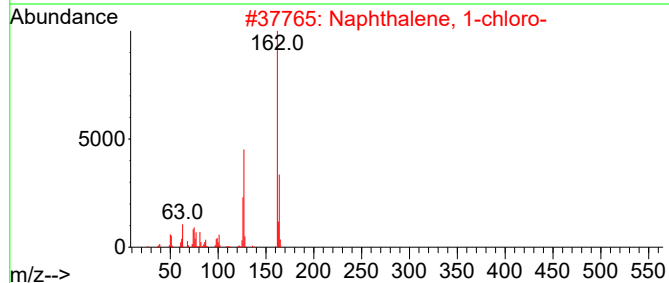
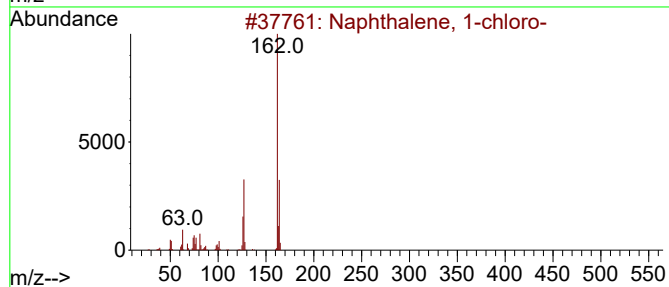
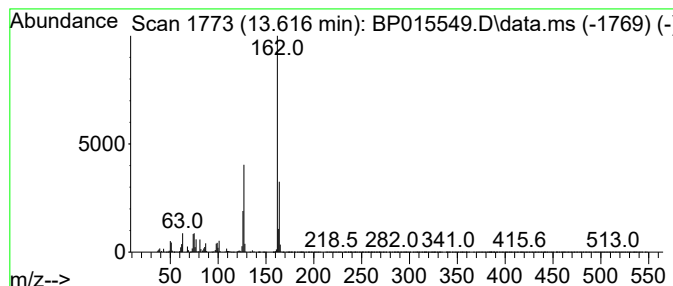
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Naphthalene, 1-chloro- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.616	15.82 ng/ul	1494610	Acenaphthene-d10	14.733

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1-chloro-	162	C10H7Cl	000090-13-1	97
2			Naphthalene, 1-chloro-	162	C10H7Cl	000090-13-1	95
3			Naphthalene, 2-chloro-	162	C10H7Cl	000091-58-7	94
4			Naphthalene, 1-chloro-	162	C10H7Cl	000090-13-1	94
5			Naphthalene, 1-chloro-	162	C10H7Cl	000090-13-1	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
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Misc :
ALS Vial : 44 Sample Multiplier: 1

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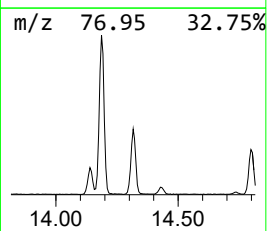
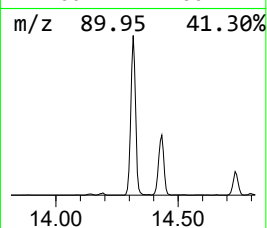
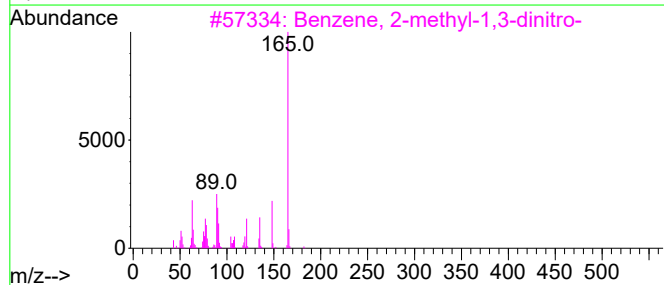
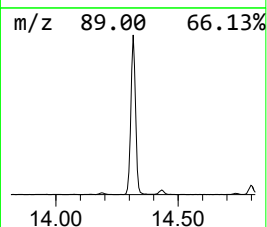
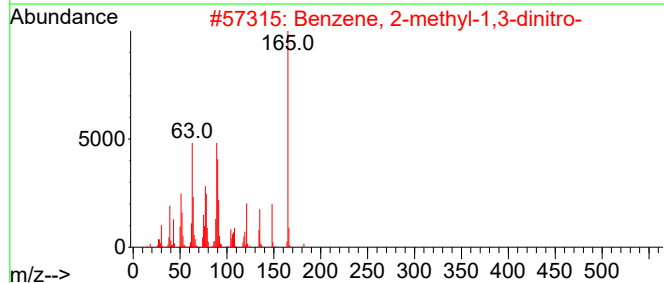
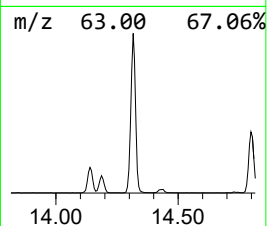
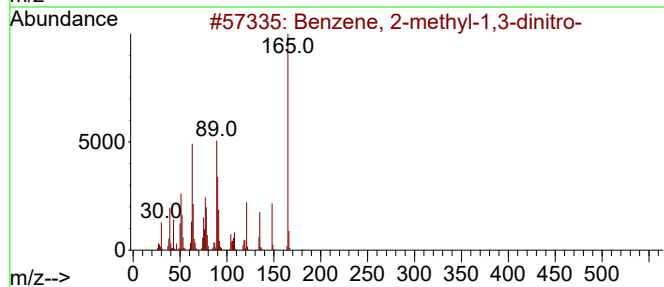
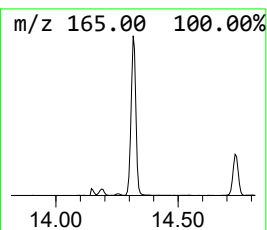
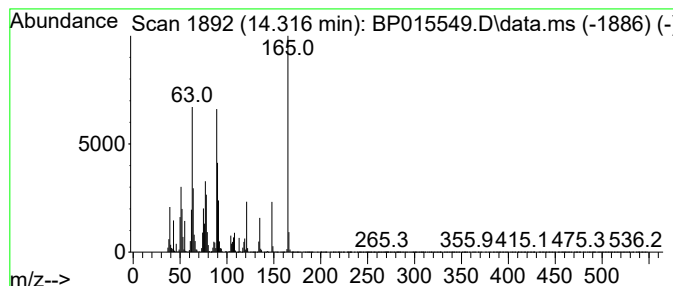
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Benzene, 2-methyl-1,3-dinitro- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.316	16.64 ng/ul	1572170	Acenaphthene-d10	14.733

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-methyl-1,3-dinitro-	182	C7H6N2O4	000606-20-2	64
2			Benzene, 2-methyl-1,3-dinitro-	182	C7H6N2O4	000606-20-2	64
3			Benzene, 2-methyl-1,3-dinitro-	182	C7H6N2O4	000606-20-2	52
4			Benzene, 1-(chloromethyl)-2-nitro-	171	C7H6ClNO2	000612-23-7	37
5			1H-Benzotriazole, 4-nitro-	164	C6H4N4O2	006299-39-4	33



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
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Acq On : 15 Jun 2023 10:36
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Misc :
ALS Vial : 44 Sample Multiplier: 1

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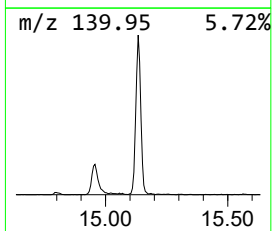
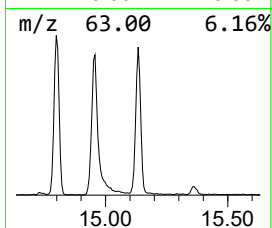
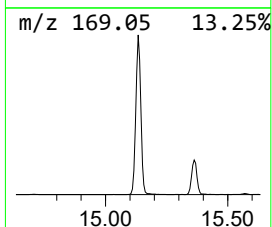
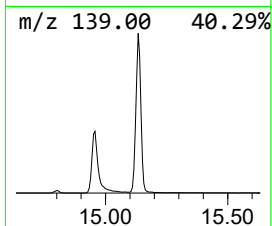
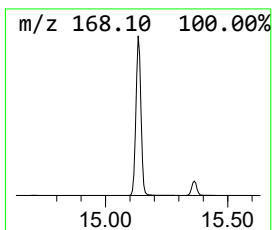
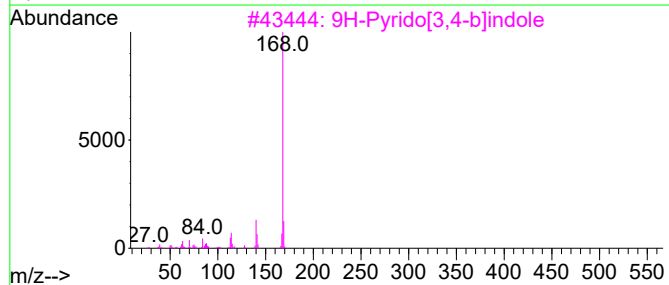
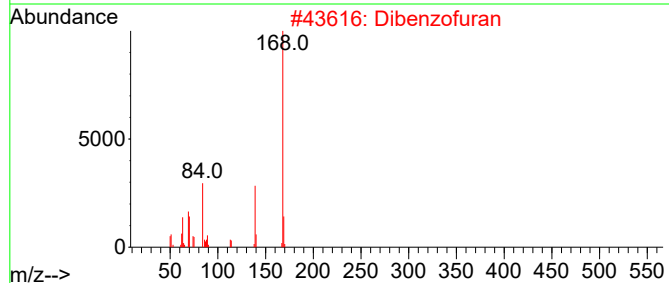
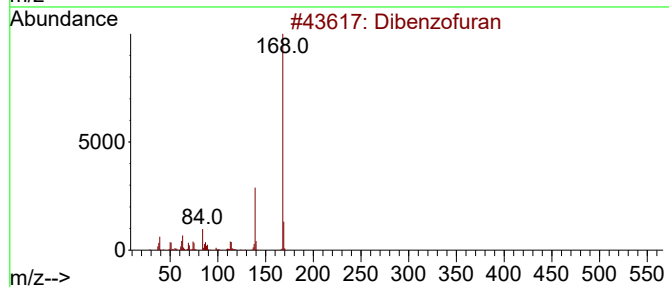
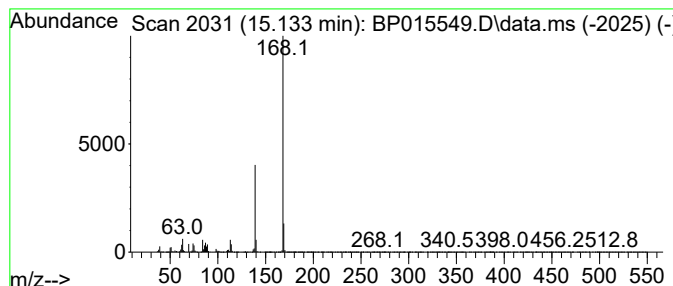
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Dibenzo furan Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.133	20.26 ng/ul	1914400	Acenaphthene-d10	14.733

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dibenzofuran	168	C12H8O	000132-64-9	91
2			Dibenzofuran	168	C12H8O	000132-64-9	78
3			9H-Pyrido[3,4-b]indole	168	C11H8N2	000244-63-3	53
4			1-Naphthalenecarbonitrile, 8-amino-	168	C11H8N2	038515-13-8	53
5			9H-Pyrido[3,4-b]indole	168	C11H8N2	000244-63-3	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
Data File : BP015549.D
Acq On : 15 Jun 2023 10:36
Operator : MA/JU
Sample : 03088-15
Misc :
ALS Vial : 44 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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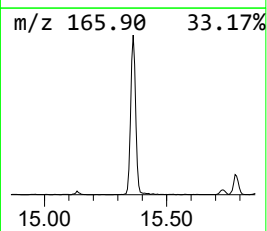
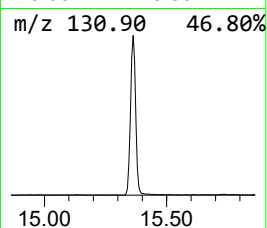
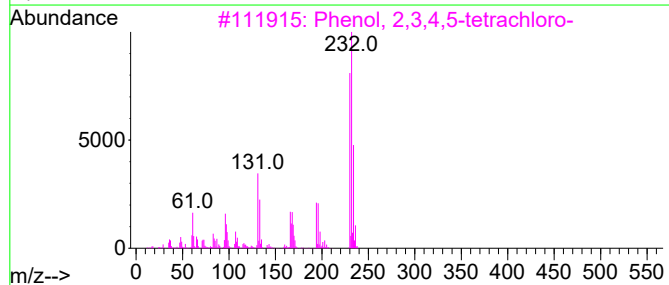
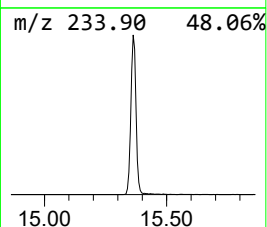
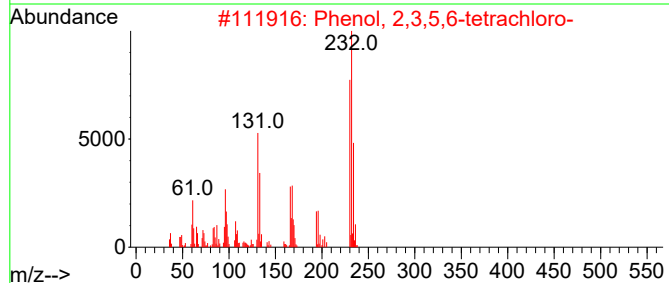
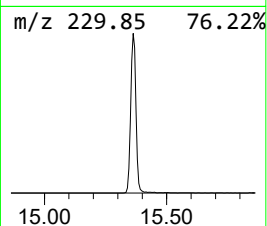
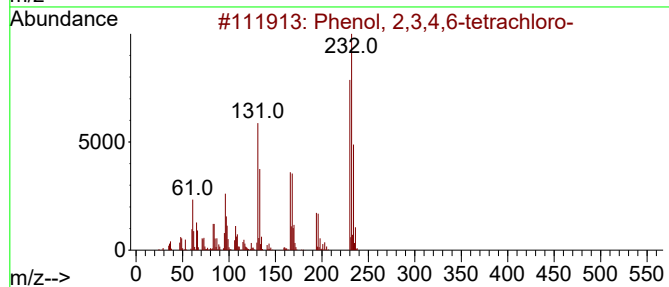
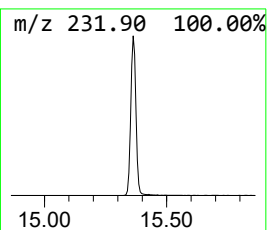
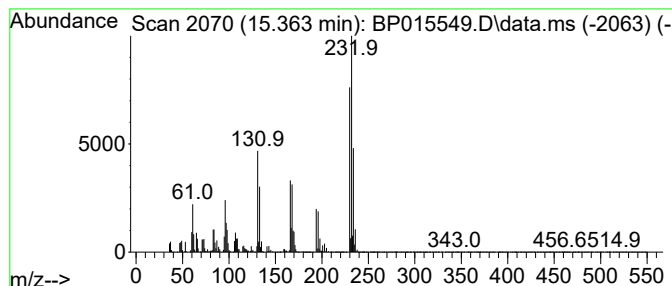
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 Phenol, 2,3,4,6-tetrachloro- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.363	19.28 ng/ul	1821130	Acenaphthene-d10	14.733

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 2,3,4,6-tetrachloro-	230	C6H2Cl4O	000058-90-2	99
2			Phenol, 2,3,5,6-tetrachloro-	230	C6H2Cl4O	000935-95-5	99
3			Phenol, 2,3,4,5-tetrachloro-	230	C6H2Cl4O	004901-51-3	98
4			Phenol, 2,3,4,6-tetrachloro-	230	C6H2Cl4O	000058-90-2	98
5			Phenol, 2,3,4,6-tetrachloro-	230	C6H2Cl4O	000058-90-2	98



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
Data File : BP015549.D
Acq On : 15 Jun 2023 10:36
Operator : MA/JU
Sample : 03088-15
Misc :
ALS Vial : 44 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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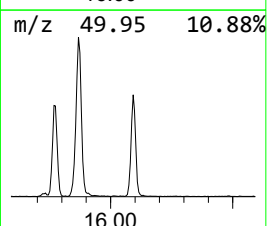
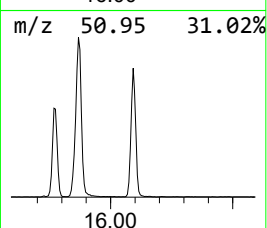
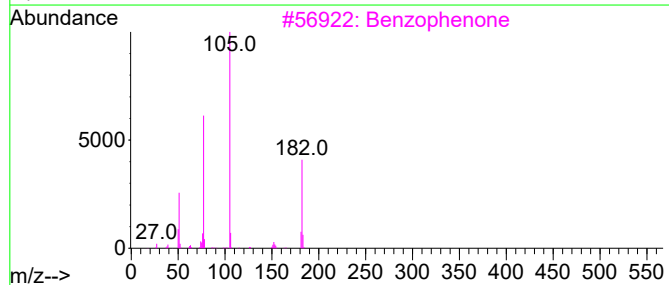
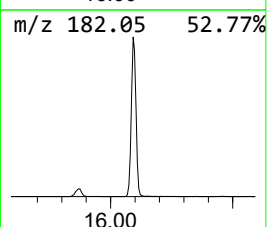
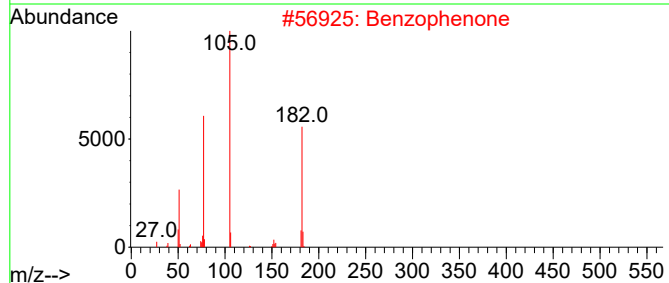
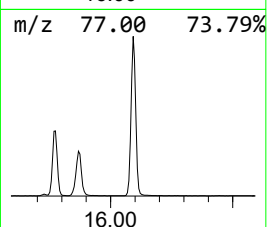
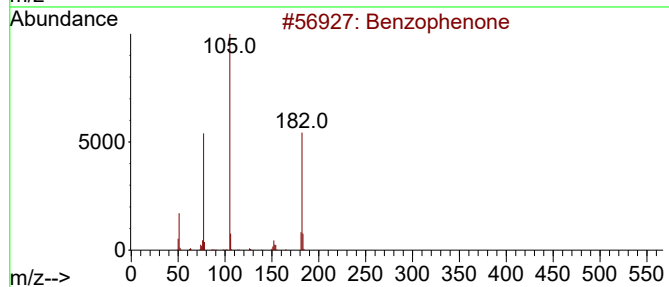
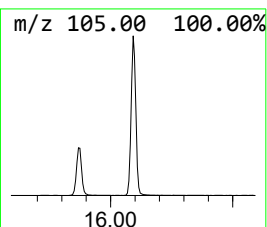
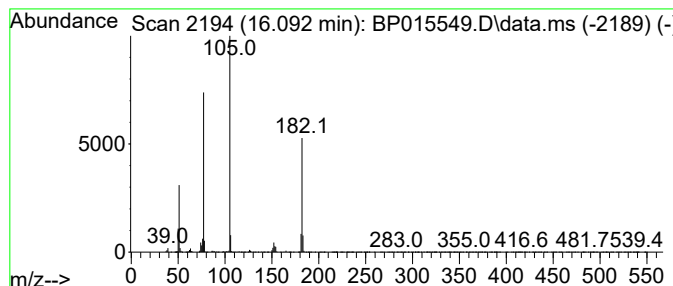
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 Benzophenone Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.092	15.23 ng/ul	1438550	Acenaphthene-d10	14.733

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	96
4			Benzophenone	182	C13H10O	000119-61-9	91
5			Benzophenone	182	C13H10O	000119-61-9	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
Data File : BP015549.D
Acq On : 15 Jun 2023 10:36
Operator : MA/JU
Sample : 03088-15
Misc :
ALS Vial : 44 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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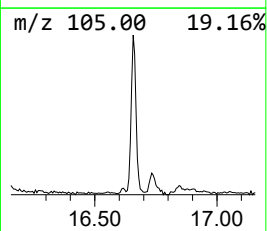
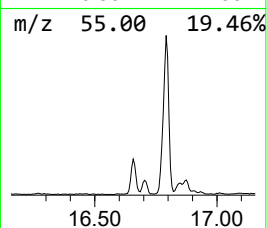
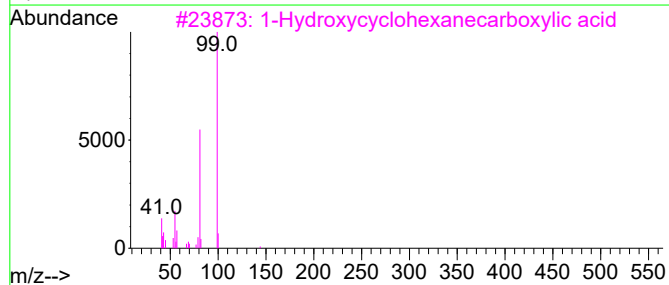
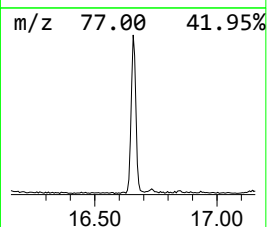
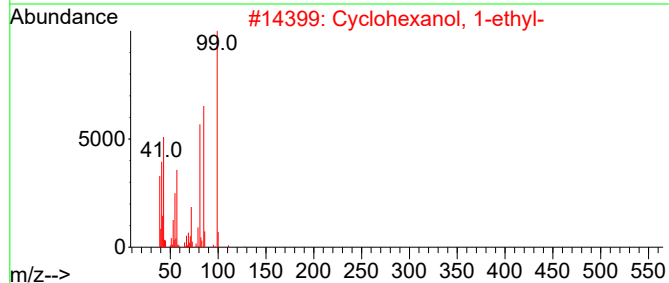
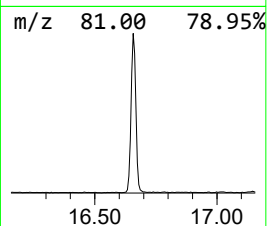
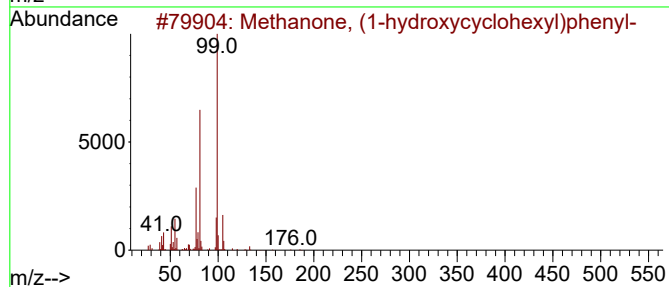
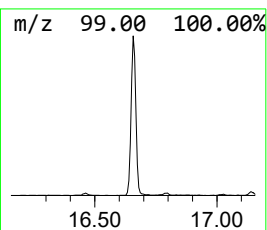
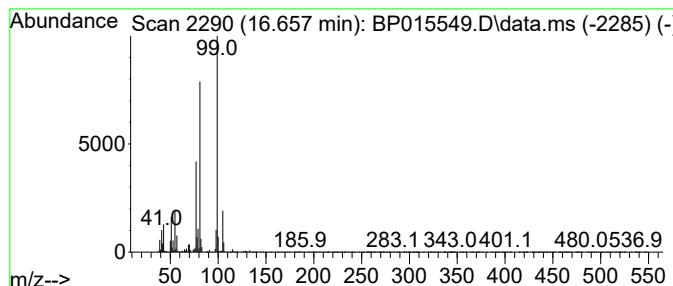
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 Methanone, (1-hydroxycyclohexyl) Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.657	4.12 ng/ul	510756	Phenanthrene-d10	17.492

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methanone, (1-hydroxycyclohexyl)...	204	C13H16O2	000947-19-3	90
2			Cyclohexanol, 1-ethyl-	128	C8H16O	001940-18-7	50
3			1-Hydroxycyclohexanecarboxylic acid	144	C7H12O3	001123-28-0	50
4			Methanone, (1-hydroxycyclohexyl)...	204	C13H16O2	000947-19-3	49
5			1,1'-Bicyclohexyl-1,1'-diol	198	C12H22O2	002888-11-1	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
Data File : BP015549.D
Acq On : 15 Jun 2023 10:36
Operator : MA/JU
Sample : 03088-15
Misc :
ALS Vial : 44 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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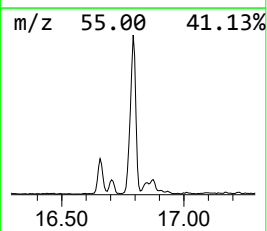
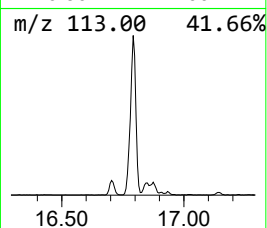
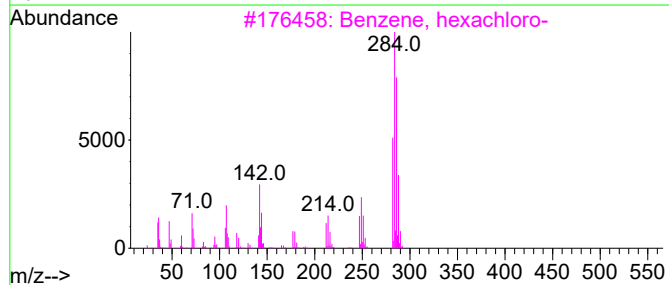
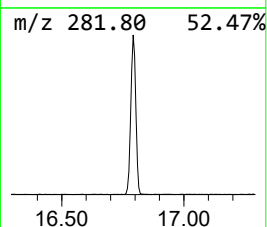
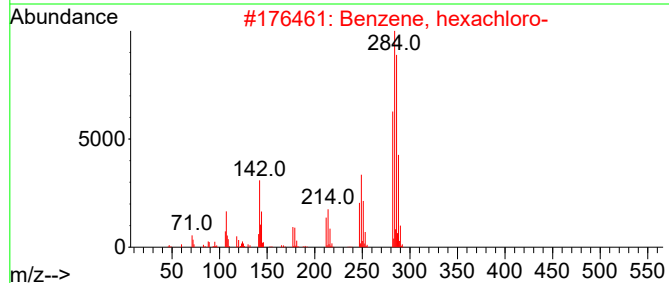
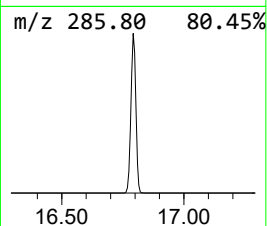
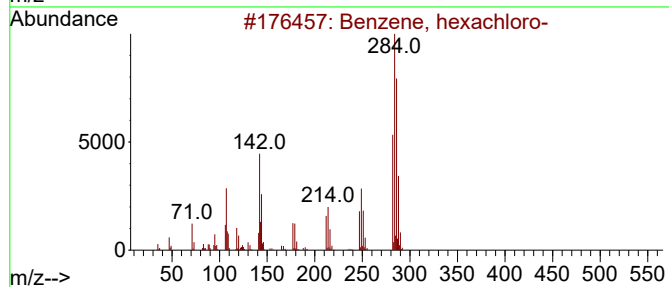
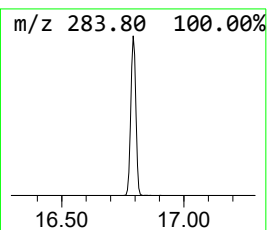
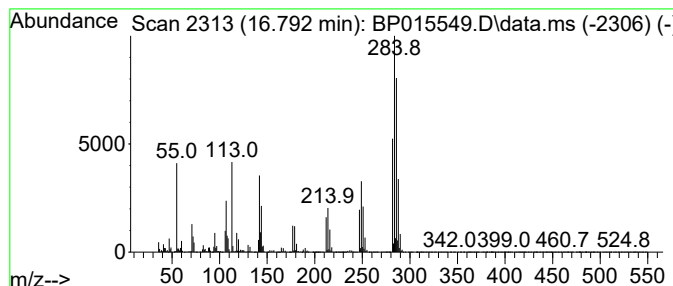
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 Benzene, hexachloro- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.792	18.42 ng/ul	2282400	Phenanthrene-d10	17.492

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, hexachloro-	282	C6Cl6	000118-74-1	99
2			Benzene, hexachloro-	282	C6Cl6	000118-74-1	99
3			Benzene, hexachloro-	282	C6Cl6	000118-74-1	99
4			1,3-Cyclopentadiene, 1,2,3,4-tet...	282	C6Cl6	006317-25-5	98
5			Benzene, hexachloro-	282	C6Cl6	000118-74-1	98



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
Data File : BP015549.D
Acq On : 15 Jun 2023 10:36
Operator : MA/JU
Sample : 03088-15
Misc :
ALS Vial : 44 Sample Multiplier: 1

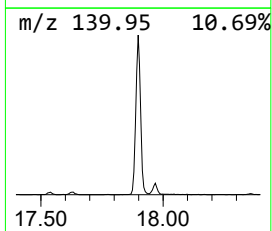
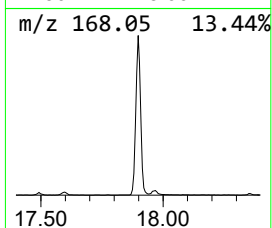
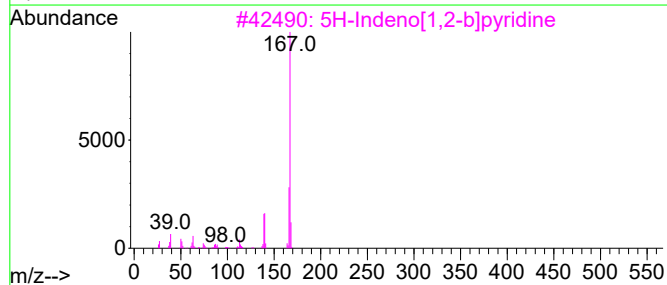
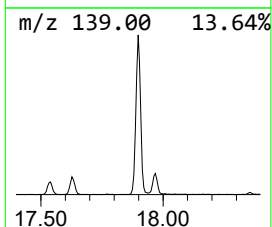
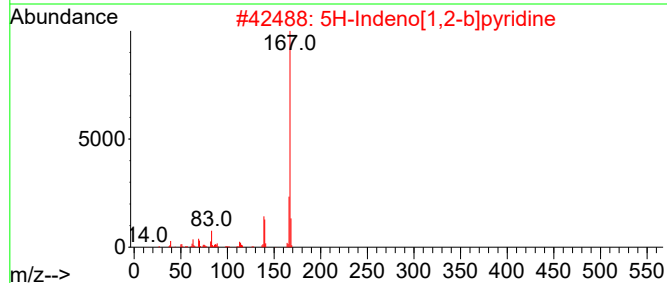
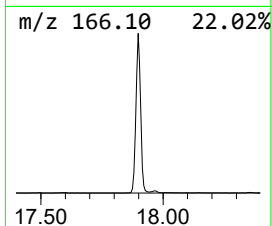
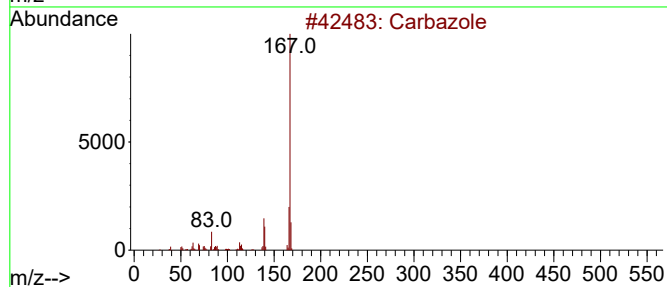
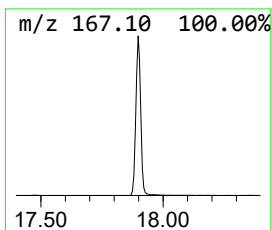
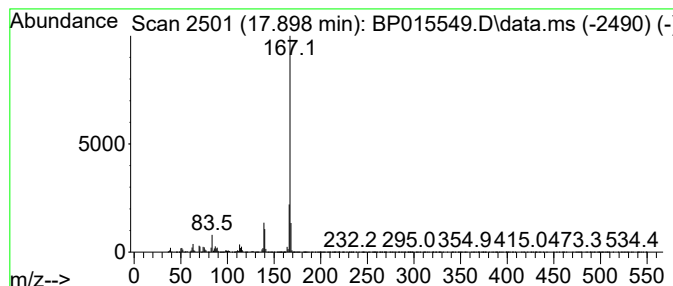
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 Carba zole Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD		R.T.	
17.898	21.54 ng/ul	2668600	Phenanthrene-d10		17.492	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Carbazole		167	C12H9N	000086-74-8	96
2	5H-Indeno[1,2-b]pyridine		167	C12H9N	000244-99-5	95
3	5H-Indeno[1,2-b]pyridine		167	C12H9N	000244-99-5	94
4	Carbazole		167	C12H9N	000086-74-8	91
5	9H-Carbazole-9-methanol		197	C13H11NO	002409-36-1	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
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Acq On : 15 Jun 2023 10:36
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Sample : 03088-15
Misc :
ALS Vial : 44 Sample Multiplier: 1

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ClientSampleId :
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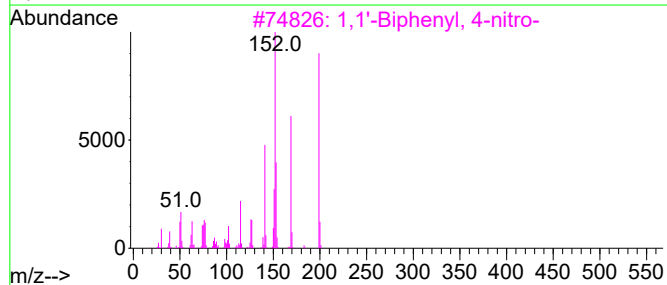
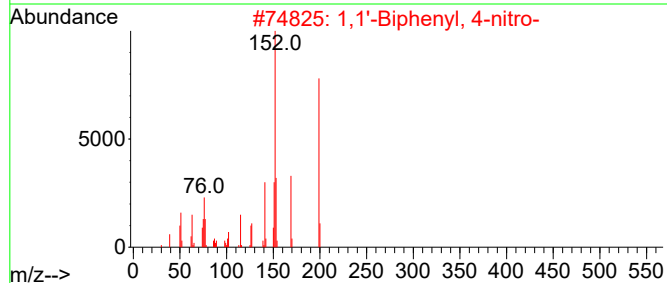
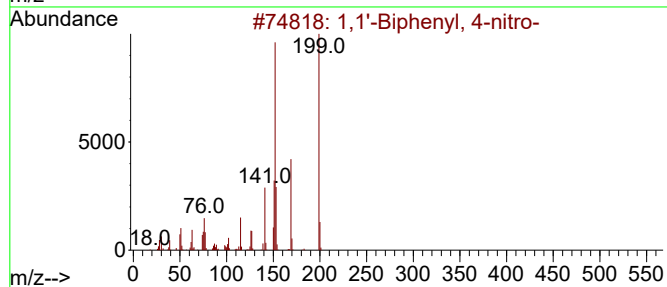
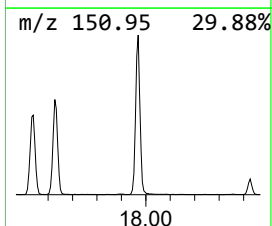
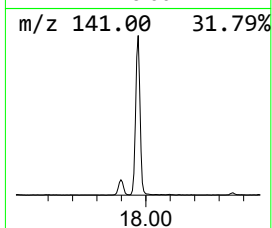
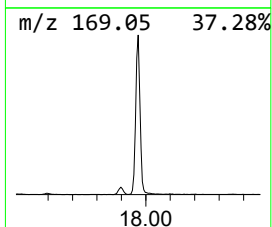
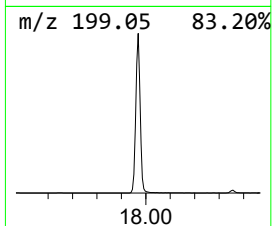
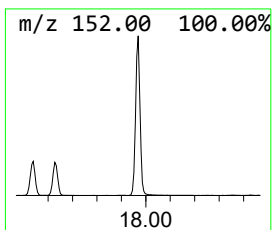
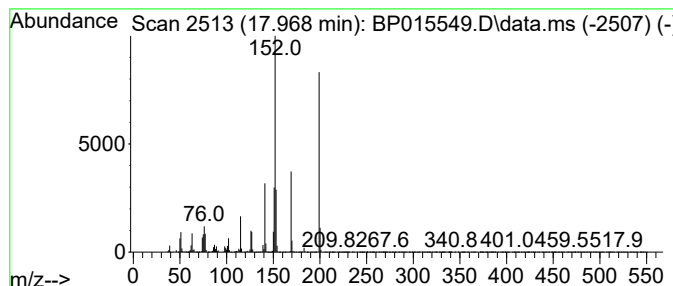
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 1,1'-Biphenyl, 4-nitro- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.968	20.65 ng/ul	2558770	Phenanthrene-d10	17.492

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 4-nitro-			199	C12H9NO2	000092-93-3	99
2	1,1'-Biphenyl, 4-nitro-			199	C12H9NO2	000092-93-3	98
3	1,1'-Biphenyl, 4-nitro-			199	C12H9NO2	000092-93-3	97
4	1,1'-Biphenyl, 4-nitro-			199	C12H9NO2	000092-93-3	94
5	1,1'-Biphenyl, 3-nitro-			199	C12H9NO2	002113-58-8	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
Data File : BP015549.D
Acq On : 15 Jun 2023 10:36
Operator : MA/JU
Sample : 03088-15
Misc :
ALS Vial : 44 Sample Multiplier: 1

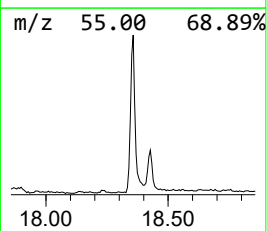
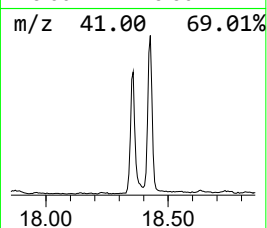
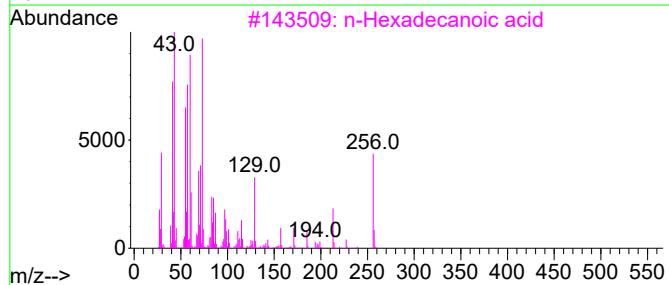
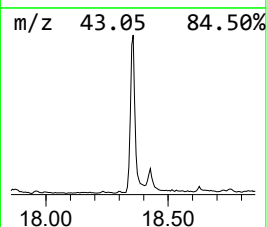
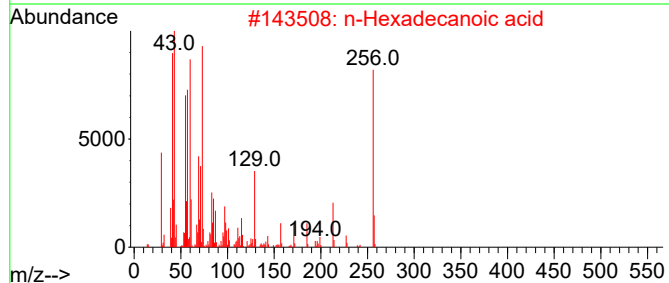
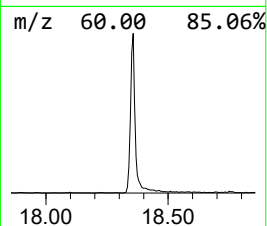
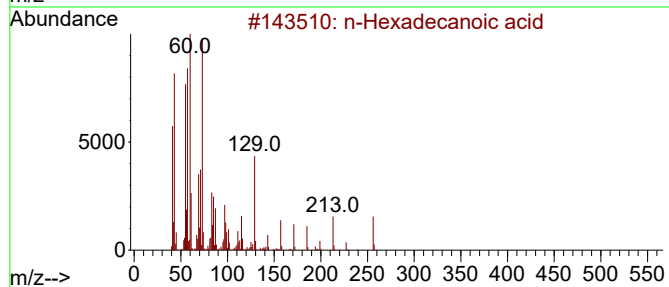
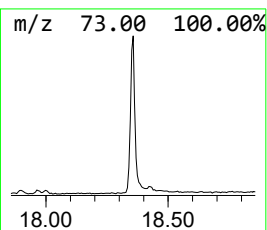
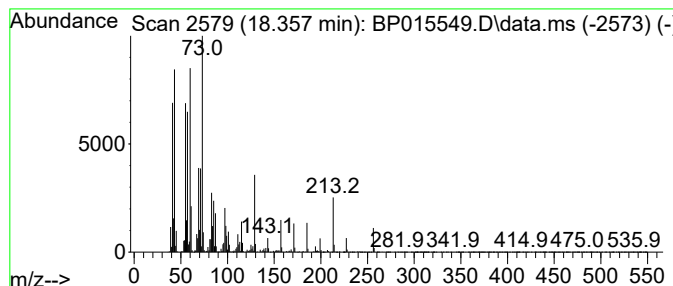
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Quant Title : SVOA CALIBRATION

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

Peak Number 17 n-Hexadecanoic acid Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.357	10.26 ng/ul	1270900	Phenanthrene-d10	17.492	
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	98
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
5	Pentadecanoic acid	242	C15H30O2	001002-84-2	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
Data File : BP015549.D
Acq On : 15 Jun 2023 10:36
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Sample : 03088-15
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Instrument :
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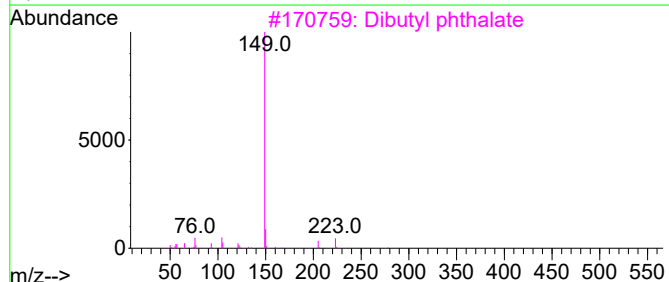
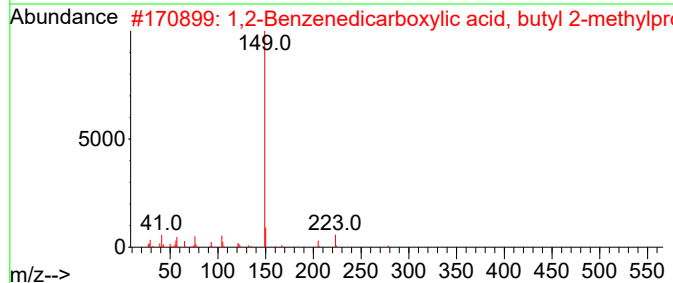
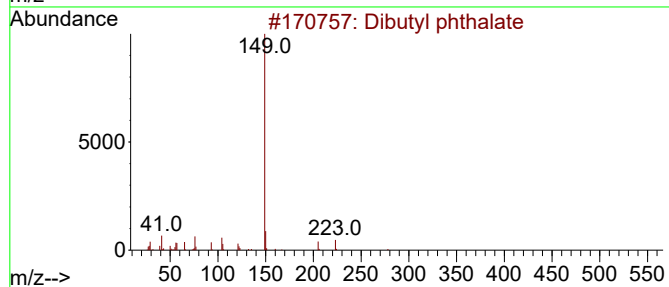
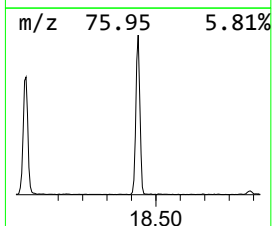
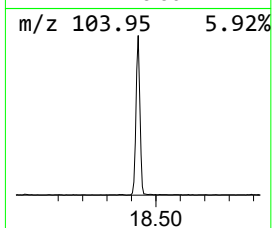
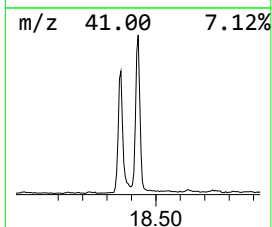
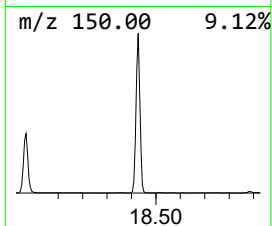
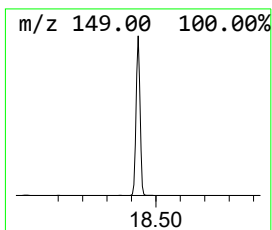
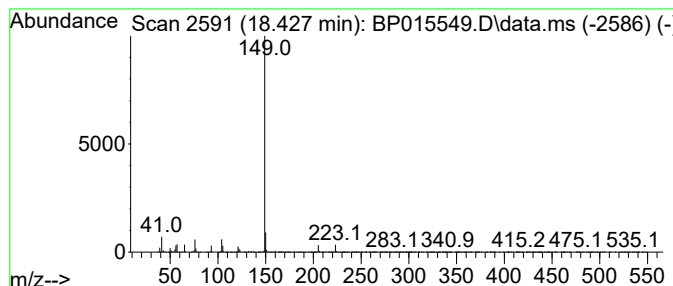
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Quant Title : SVOA CALIBRATION

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

Peak Number 18 Dibutyl phthalate Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.427	18.52 ng/ul	2295140	Phenanthrene-d10	17.492

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dibutyl phthalate	278	C16H22O4	000084-74-2	95
2			1,2-Benzenedicarboxylic acid, bu...	278	C16H22O4	017851-53-5	94
3			Dibutyl phthalate	278	C16H22O4	000084-74-2	91
4			1,4-Dibutyl benzene-1,4-dicarbox...	278	C16H22O4	001962-75-0	90
5			1,2-Benzenedicarboxylic acid, bi...	278	C16H22O4	000084-69-5	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
Data File : BP015549.D
Acq On : 15 Jun 2023 10:36
Operator : MA/JU
Sample : 03088-15
Misc :
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Instrument :
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ClientSampleId :
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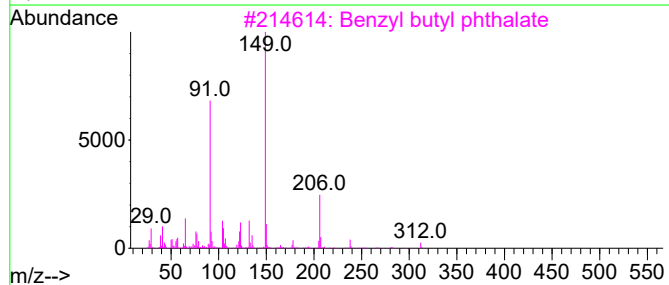
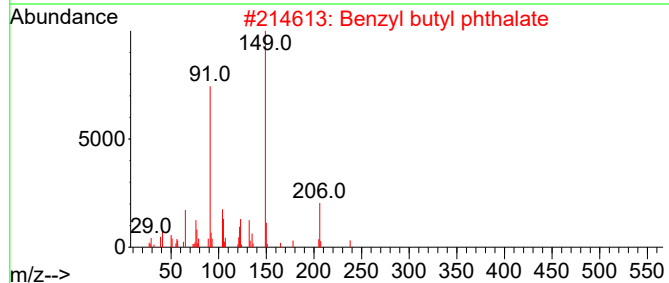
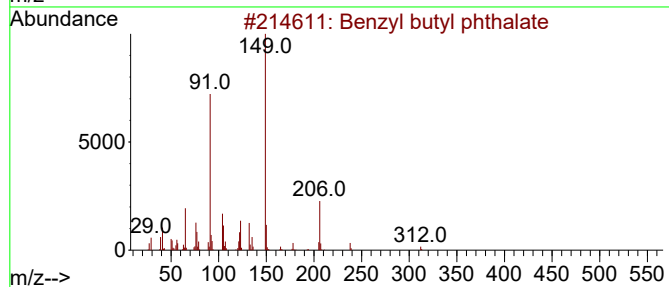
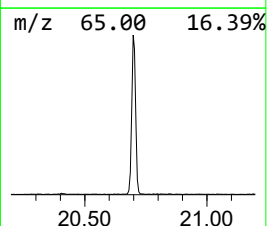
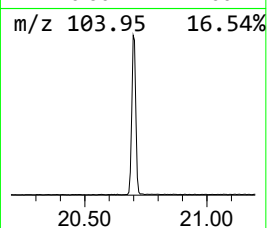
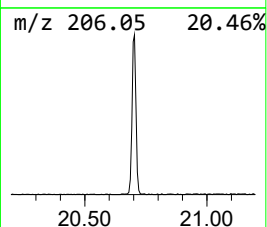
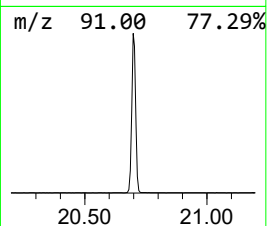
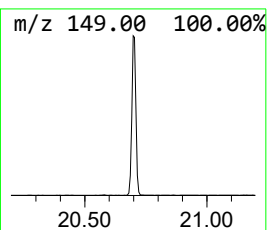
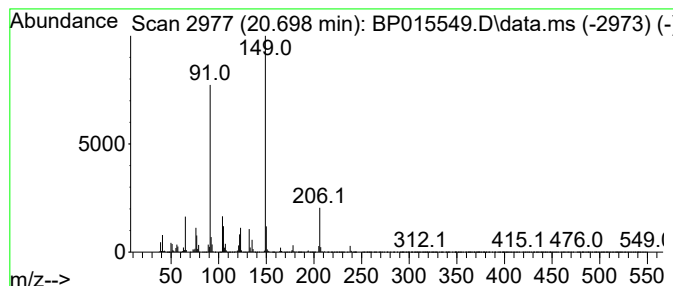
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Quant Title : SVOA CALIBRATION

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

Peak Number 19 Benzyl butyl phthalate Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.698	5.51 ng/ul	2397030	Chrysene-d12	21.580

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzyl butyl phthalate	312	C19H20O4	000085-68-7	94
2			Benzyl butyl phthalate	312	C19H20O4	000085-68-7	91
3			Benzyl butyl phthalate	312	C19H20O4	000085-68-7	91
4			Benzyl butyl phthalate	312	C19H20O4	000085-68-7	72
5			Phthalic acid, 3-methoxy-4-nitro...	457	C25H31NO7	1000382-53-7	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
Data File : BP015549.D
Acq On : 15 Jun 2023 10:36
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Sample : 03088-15
Misc :
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Instrument :
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ClientSampleId :
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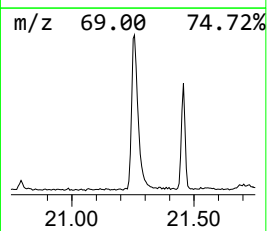
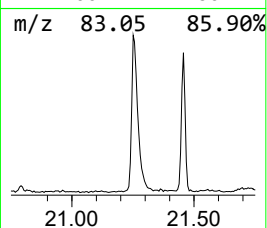
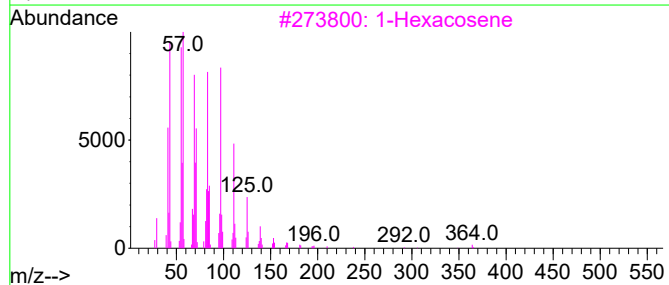
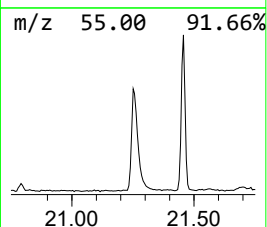
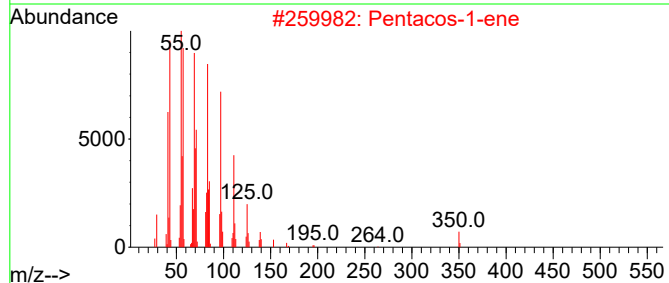
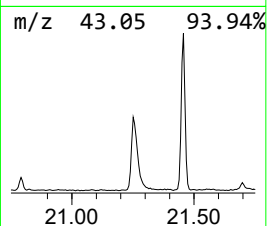
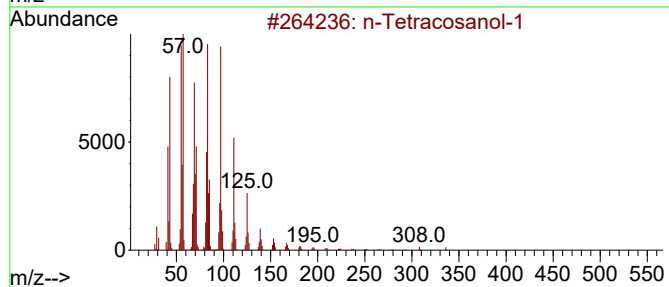
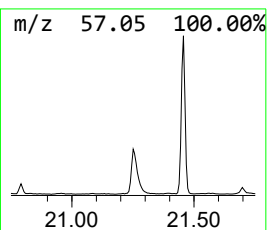
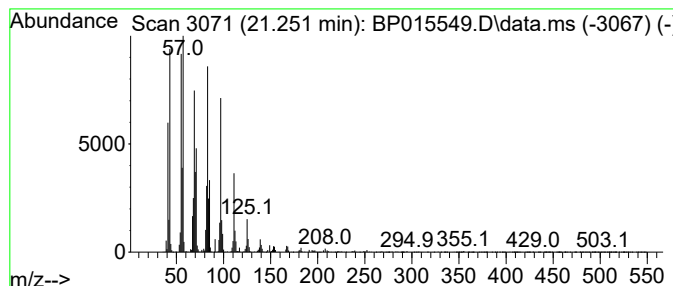
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Quant Title : SVOA CALIBRATION

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

Peak Number 20 n-Tetracosanol-1 Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.251	3.62 ng/ul	1574050	Chrysene-d12	21.580

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Tetracosanol-1	354	C24H50O	000506-51-4	94
2			Pentacos-1-ene	350	C25H50	016980-85-1	94
3			1-Hexacosene	364	C26H52	018835-33-1	94
4			Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	94
5			Cyclopentadecane	210	C15H30	000295-48-7	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
Data File : BP015549.D
Acq On : 15 Jun 2023 10:36
Operator : MA/JU
Sample : 03088-15
Misc :
ALS Vial : 44 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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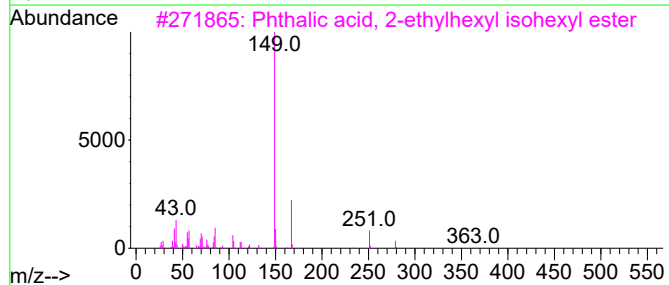
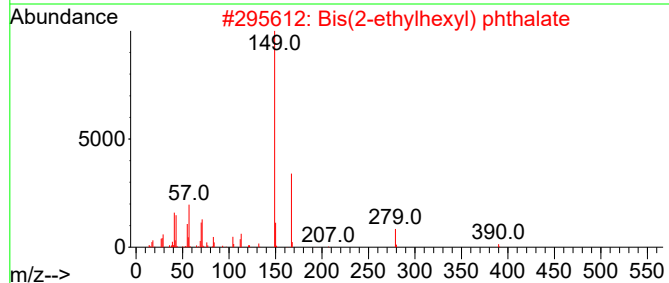
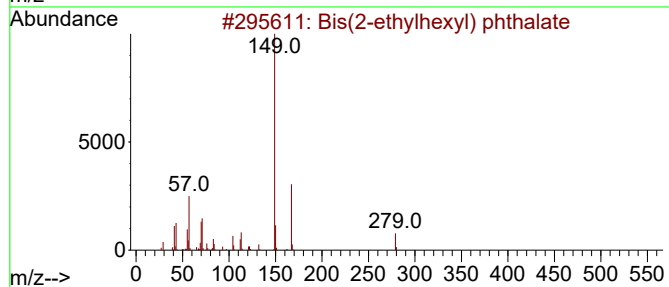
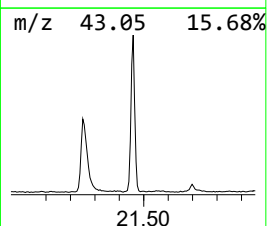
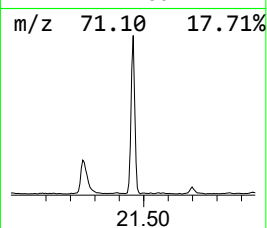
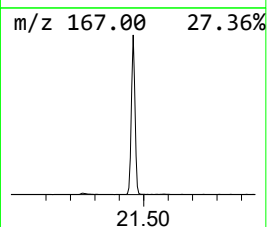
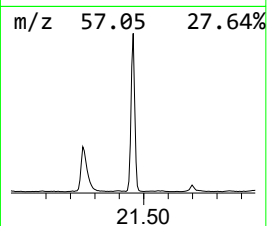
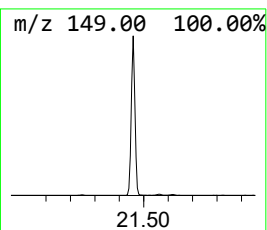
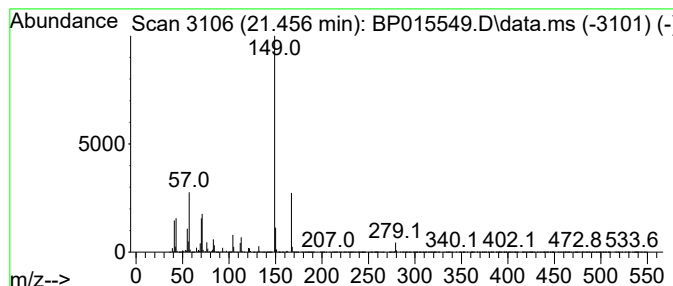
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Quant Title : SVOA CALIBRATION

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

Peak Number 21 Bis(2-ethylhexyl) phthalate Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.456	8.28 ng/ul	3602880	Chrysene-d12	21.580

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bis(2-ethylhexyl) phthalate	390	C24H38O4	000117-81-7	91
2			Bis(2-ethylhexyl) phthalate	390	C24H38O4	000117-81-7	91
3			Phthalic acid, 2-ethylhexyl isoh...	362	C22H34O4	1000308-98-5	86
4			Phthalic acid, di(oct-3-yl) ester	390	C24H38O4	1000377-72-3	80
5			1,2-Benzenedicarboxylic acid, bu...	334	C20H30O4	000085-69-8	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
Data File : BP015549.D
Acq On : 15 Jun 2023 10:36
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Sample : 03088-15
Misc :
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Instrument :
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ClientSampleId :
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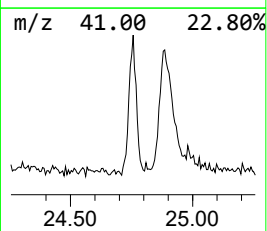
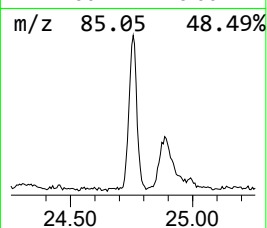
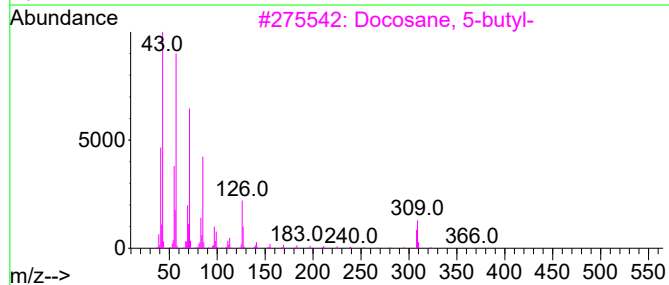
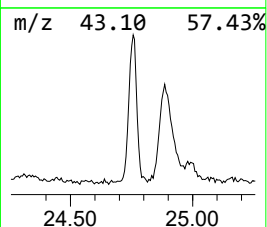
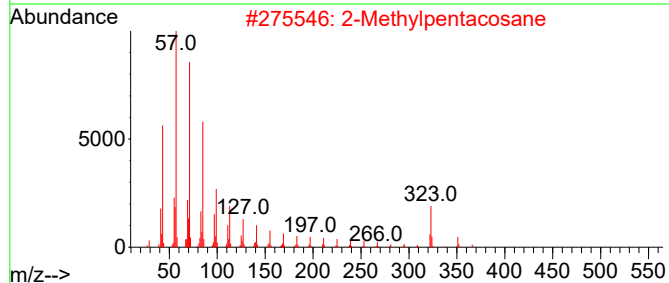
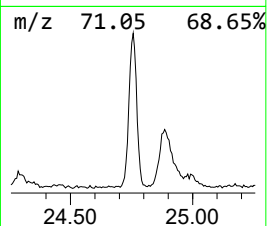
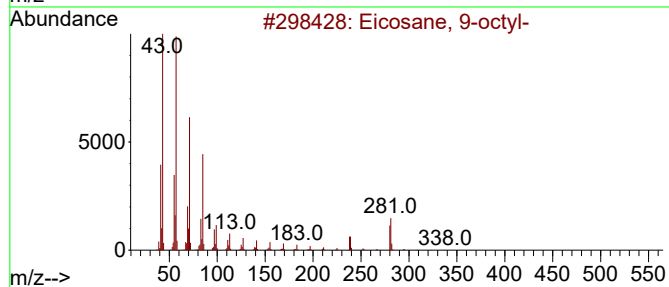
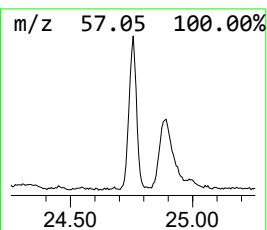
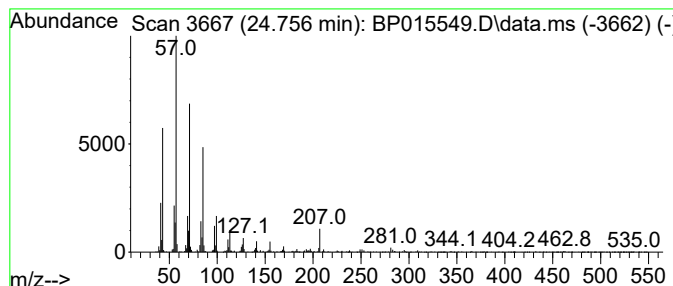
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Quant Title : SVOA CALIBRATION

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

Peak Number 22 (DEL) Alkane: Straight-Chai... Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.756	4.08 ng/ul	514525	Perylene-d12	24.127

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane, 9-octyl-	394	C28H58	013475-77-9	93
2			2-Methylpentacosane	366	C26H54	000629-87-8	91
3			Docosane, 5-butyl-	366	C26H54	055282-16-1	91
4			Heptadecane, 9-hexyl-	324	C23H48	055124-79-3	91
5			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
 Data File : BP015549.D
 Acq On : 15 Jun 2023 10:36
 Operator : MA/JU
 Sample : 03088-15
 Misc :
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCGV6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.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
2-Pentanone, 4-...	5.140	9.3	ng/ul	380006	1	8.087	821004	20.0
1-Propanamine, ...	8.904	28.4	ng/ul	1167810	1	8.087	821004	20.0
1,3-Butadiene, ...	11.245	22.4	ng/ul	1489170	2	10.916	1327320	20.0
Phenol, 4-chlor...	12.198	41.3	ng/ul	2737360	2	10.916	1327320	20.0
Benzene, 1,2,3,...	12.933	16.1	ng/ul	1516820	3	14.733	1889620	20.0
Phenol, 2,4,6-t...	13.175	27.0	ng/ul	2547700	3	14.733	1889620	20.0
Biphenyl	13.569	30.1	ng/ul	2839640	3	14.733	1889620	20.0
Naphthalene, 1-...	13.616	15.8	ng/ul	1494610	3	14.733	1889620	20.0
Benzene, 2-meth...	14.316	16.6	ng/ul	1572170	3	14.733	1889620	20.0
Dibenzo furan	15.133	20.3	ng/ul	1914400	3	14.733	1889620	20.0
Phenol, 2,3,4,6...	15.363	19.3	ng/ul	1821130	3	14.733	1889620	20.0
Benzophenone	16.092	15.2	ng/ul	1438550	3	14.733	1889620	20.0
Methanone, (1-h...	16.657	4.1	ng/ul	510756	4	17.492	2478310	20.0
Benzene, hexach...	16.792	18.4	ng/ul	2282400	4	17.492	2478310	20.0
Carba zole	17.898	21.5	ng/ul	2668600	4	17.492	2478310	20.0
1,1'-Biphenyl, ...	17.968	20.6	ng/ul	2558770	4	17.492	2478310	20.0
n-Hexadecanoic ...	18.357	10.3	ng/ul	1270900	4	17.492	2478310	20.0
Dibutyl phthalate	18.427	18.5	ng/ul	2295140	4	17.492	2478310	20.0
Benzyl butyl ph...	20.698	5.5	ng/ul	2397030	5	21.580	8698870	20.0
n-Tetracosanol-1	21.251	3.6	ng/ul	1574050	5	21.580	8698870	20.0
Bis(2-ethylhexy...	21.456	8.3	ng/ul	3602880	5	21.580	8698870	20.0
(DEL) Alkane: S...	24.756	4.1	ng/ul	514525	6	24.127	2519790	20.0