

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101624\
Data File : BP022412.D
Acq On : 16 Oct 2024 12:00
Operator : RC/JU
Sample : P4329-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4E79

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 0.1 % of largest Peak

Max Peaks: 100

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: BP022412.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	6.887	647	663	677	rVB2	523824	1323551	6.08%	0.467%
2	7.011	677	684	693	rBV	226556	359809	1.65%	0.127%
3	7.105	693	700	710	rVB	2039749	3153744	14.48%	1.114%
4	7.246	712	724	732	rBV2	1348483	2534157	11.64%	0.895%
5	7.705	786	802	812	rBV	3475654	6121573	28.11%	2.162%
6	8.016	846	855	865	rBV	3462729	5764462	26.47%	2.036%
7	8.381	910	917	924	rBV	429756	671207	3.08%	0.237%
8	8.475	924	933	958	rVV	2521557	4241442	19.48%	1.498%
9	8.734	969	977	984	rBV	2193290	3577381	16.43%	1.263%
10	8.810	984	990	993	rVV	310088	522691	2.40%	0.185%
11	8.852	993	997	1006	rVB	556541	909813	4.18%	0.321%
12	9.222	1051	1060	1065	rBV2	21563	43165	0.20%	0.015%
13	9.381	1075	1087	1102	rBV	3393242	5565504	25.56%	1.966%
14	9.522	1102	1111	1112	rVV	286930	388080	1.78%	0.137%
15	9.552	1112	1116	1128	rVV	606873	1091796	5.01%	0.386%
16	9.675	1130	1137	1145	rVB	129542	211557	0.97%	0.075%
17	9.852	1158	1167	1176	rBV	2590406	4185537	19.22%	1.478%
18	10.087	1196	1207	1221	rBV2	2303397	4263053	19.57%	1.506%
19	10.434	1256	1266	1272	rBV	397182	693889	3.19%	0.245%
20	10.569	1281	1289	1304	rVB	269368	545950	2.51%	0.193%
21	10.757	1310	1321	1335	rBV	4249987	6186224	28.41%	2.185%
22	11.722	1477	1485	1496	rBV	837346	1459165	6.70%	0.515%
23	11.981	1524	1529	1540	rVB2	28676	61955	0.28%	0.022%
24	12.093	1540	1548	1562	rBV	1867697	2864670	13.15%	1.012%
25	12.510	1612	1619	1624	rBV	47267	67981	0.31%	0.024%
26	12.710	1645	1653	1659	rBV	5013146	7199417	33.06%	2.543%
27	12.775	1659	1664	1679	rVB	1008894	1552688	7.13%	0.548%
28	13.104	1715	1720	1724	rBV	25443	32751	0.15%	0.012%
29	13.198	1730	1736	1744	rBV	25592	40924	0.19%	0.014%
30	13.269	1744	1748	1753	rVV2	13335	22136	0.10%	0.008%
31	13.322	1753	1757	1766	rVB3	12950	23609	0.11%	0.008%
32	13.698	1813	1821	1830	rBV	760202	1264830	5.81%	0.447%
33	13.869	1843	1850	1855	rBV	5179572	6883866	31.61%	2.431%
34	13.975	1862	1868	1871	rBV	996314	1597582	7.34%	0.564%
35	14.004	1871	1873	1881	rVB	687221	809234	3.72%	0.286%
36	14.287	1912	1921	1926	rBV	692633	996243	4.57%	0.352%

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

37	14.351	1926	1932	1940	rVV	5025547	7153595	32.85%	2.527%
38	14.422	1940	1944	1951	rVB3	61523	86547	0.40%	0.031%
39	14.528	1954	1962	1971	rBV2	1817474	2848800	13.08%	1.006%
40	14.610	1971	1976	1979	rVV	90239	143186	0.66%	0.051%

41	14.669	1979	1986	1992	rVV	3250194	4827862	22.17%	1.705%
42	14.722	1992	1995	2005	rVB4	19399	36189	0.17%	0.013%
43	14.928	2025	2030	2041	rVB	72394	113190	0.52%	0.040%
44	15.110	2054	2061	2065	rBV2	16072	28663	0.13%	0.010%
45	15.169	2066	2071	2076	rVB5	18560	33264	0.15%	0.012%

46	15.292	2086	2092	2096	rBV	1369830	1927674	8.85%	0.681%
47	15.351	2096	2102	2109	rVV2	8168285	14523533	66.69%	5.130%
48	15.434	2109	2116	2127	rVV	326956	591539	2.72%	0.209%
49	15.610	2140	2146	2149	rVV	135266	166489	0.76%	0.059%
50	15.657	2149	2154	2164	rVB	545598	771411	3.54%	0.272%

51	15.863	2182	2189	2194	rBV7	15353	28307	0.13%	0.010%
52	15.945	2200	2203	2208	rVB3	22258	26911	0.12%	0.010%
53	16.034	2212	2218	2228	rVB6	42375	96494	0.44%	0.034%
54	16.257	2246	2256	2263	rBV2	1768287	2467643	11.33%	0.872%
55	16.381	2269	2277	2287	rVB	9410829	14578977	66.94%	5.149%

56	16.739	2327	2338	2349	rBV	6945395	10644860	48.88%	3.760%
57	16.892	2359	2364	2372	rVB3	66122	128453	0.59%	0.045%
58	17.092	2391	2398	2401	rVV	763727	1343450	6.17%	0.474%
59	17.139	2401	2406	2411	rVV	2106845	3546566	16.28%	1.253%
60	17.198	2411	2416	2419	rVV	1529075	2266572	10.41%	0.801%

61	17.234	2419	2422	2437	rVV	2752154	3321284	15.25%	1.173%
62	17.369	2440	2445	2447	rVV4	14507	25483	0.12%	0.009%
63	17.522	2465	2471	2476	rBV5	26317	50859	0.23%	0.018%
64	17.716	2499	2504	2508	rBV3	67799	103466	0.48%	0.037%
65	17.798	2513	2518	2524	rVB	99431	132585	0.61%	0.047%

66	17.981	2545	2549	2551	rBV5	18514	25028	0.11%	0.009%
67	18.022	2551	2556	2562	rVV	538420	799131	3.67%	0.282%
68	18.104	2562	2570	2579	rVB	7162641	10236072	47.00%	3.615%
69	18.333	2603	2609	2616	rBV2	81392	150066	0.69%	0.053%
70	18.404	2616	2621	2629	rVV3	65609	138834	0.64%	0.049%

71	18.486	2630	2635	2642	rVV3	45128	86528	0.40%	0.031%
72	18.563	2642	2648	2654	rVB	224156	353330	1.62%	0.125%
73	18.657	2661	2664	2667	rVB5	21150	23247	0.11%	0.008%
74	18.751	2677	2680	2685	rVB4	37347	57641	0.26%	0.020%
75	18.804	2686	2689	2695	rVB3	26689	38040	0.17%	0.013%

76	18.904	2699	2706	2717	rBV4	81638	228298	1.05%	0.081%
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Integrator: RTE

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77	19.010	2722	2724	2727	rVB4	29146	33493	0.15%	0.012%
78	19.116	2738	2742	2745	rBV2	28996	33244	0.15%	0.012%
79	19.157	2745	2749	2752	rBV2	53829	83513	0.38%	0.029%
80	19.228	2752	2761	2770	rVB	8777016	14756239	67.76%	5.212%
81	19.357	2778	2783	2791	rBV	177053	294257	1.35%	0.104%
82	19.475	2799	2803	2807	rBV	93748	113993	0.52%	0.040%
83	19.563	2812	2818	2820	rVV	1808828	2695903	12.38%	0.952%
84	19.598	2820	2824	2835	rVB	4088800	6695870	30.75%	2.365%
85	19.692	2838	2840	2843	rVV4	35985	44578	0.20%	0.016%
86	19.728	2843	2846	2851	rVB2	52014	63692	0.29%	0.022%
87	20.298	2939	2943	2948	rBV	202180	271467	1.25%	0.096%
88	20.545	2980	2985	2991	rBV	2347239	2931211	13.46%	1.035%
89	21.198	3091	3096	3106	rVB	436328	698037	3.21%	0.247%
90	21.351	3119	3122	3128	rVB	140952	181401	0.83%	0.064%
91	21.433	3130	3136	3142	rVV	9940191	13156629	60.41%	4.647%
92	21.510	3142	3149	3155	rVV	8839372	16032701	73.62%	5.663%
93	21.580	3155	3161	3167	rVB	13994313	21778234	100.00%	7.692%
94	23.774	3521	3534	3547	rVB	8388207	21129888	97.02%	7.463%
95	24.563	3658	3668	3673	rBV	1059222	3087580	14.18%	1.090%
96	24.627	3673	3679	3693	rVB	1759344	4421885	20.30%	1.562%
97	24.786	3697	3706	3716	rVB2	664960	1840550	8.45%	0.650%
98	28.551	4326	4346	4347	rBV2	1860335	5233402	24.03%	1.848%
99	28.574	4348	4350	4368	rVB	2592235	6175069	28.35%	2.181%

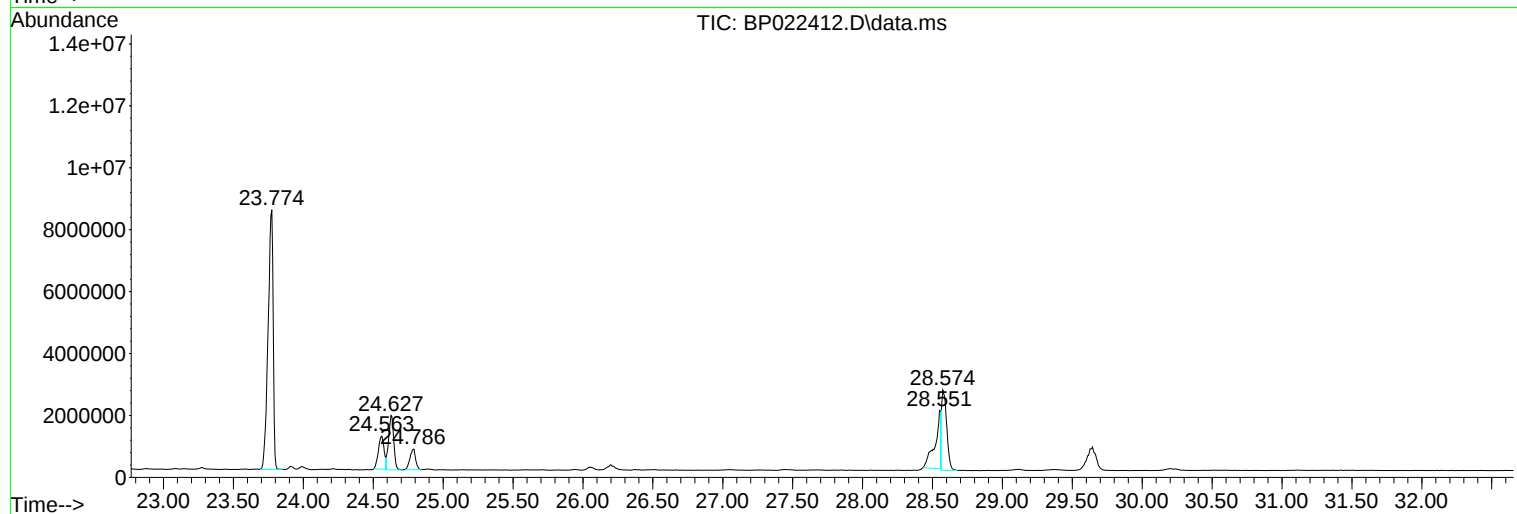
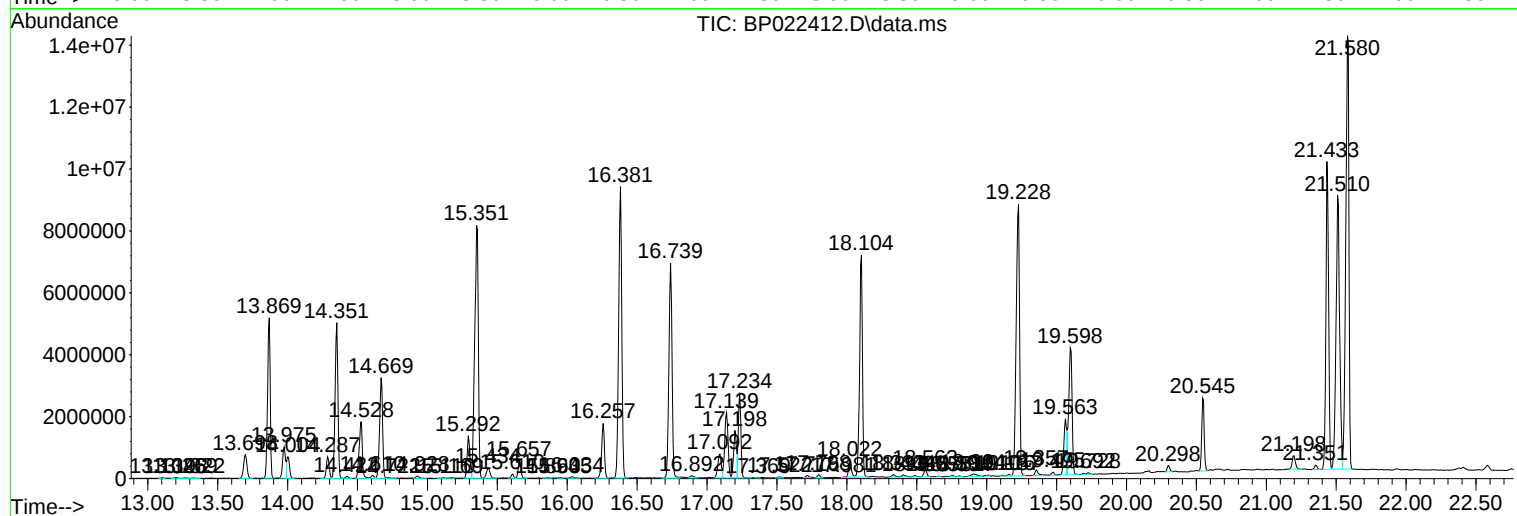
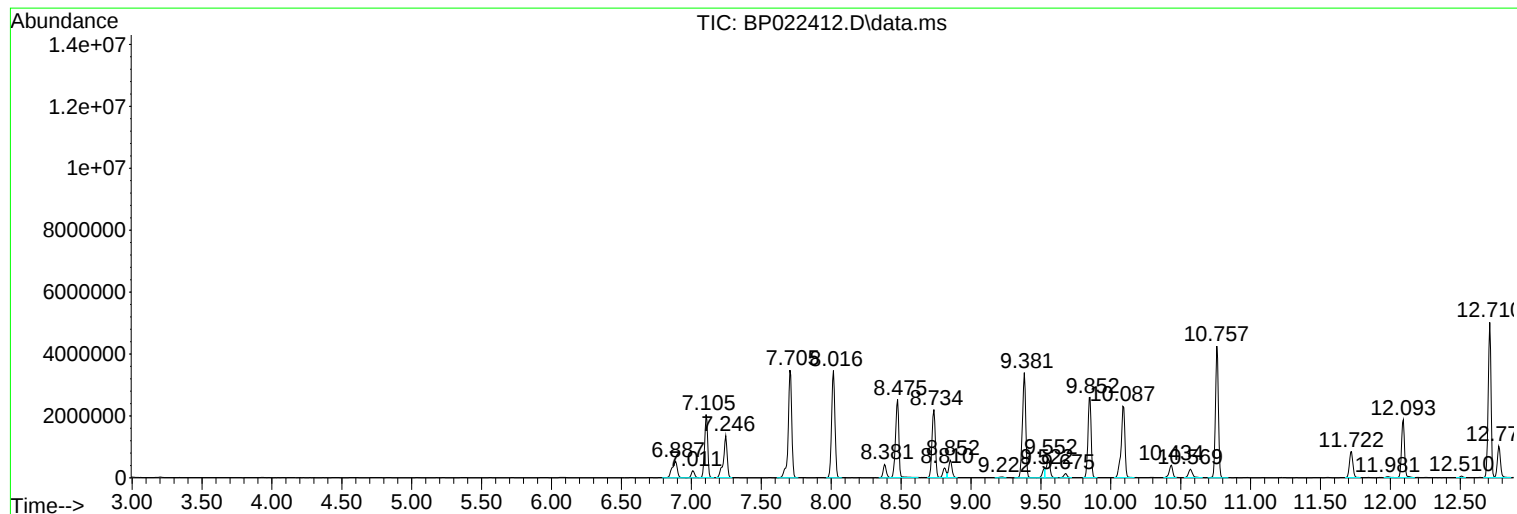
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ALS Vial : 4 Sample Multiplier: 1

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\NIST20.L
TIC Integration Parameters: LSCINT.P



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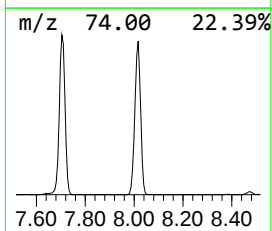
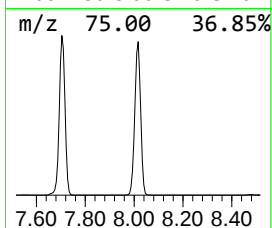
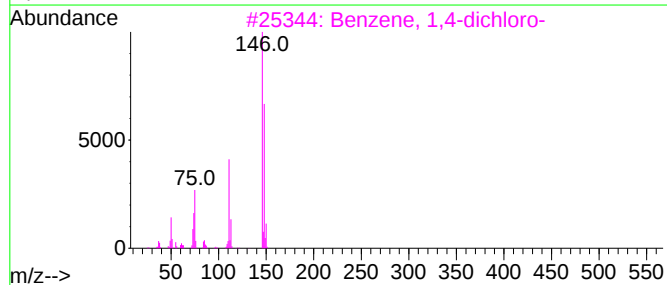
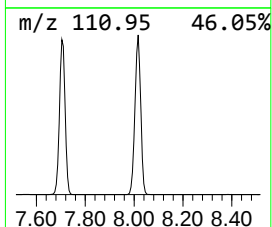
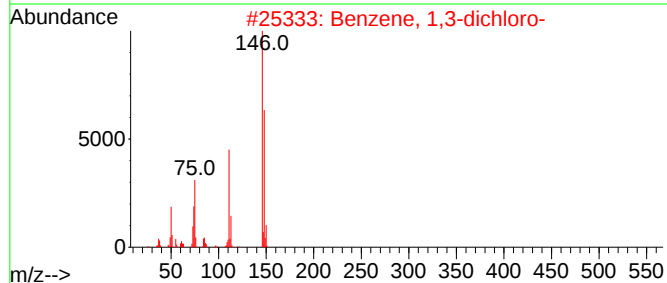
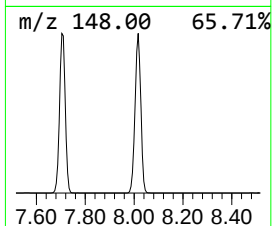
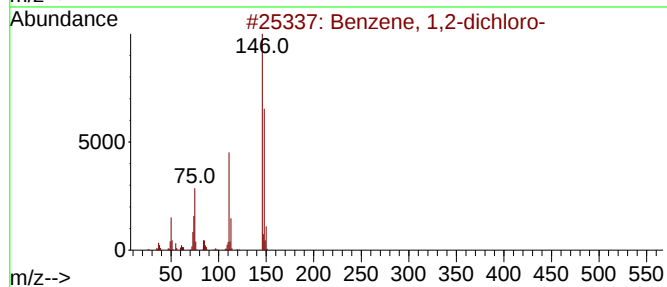
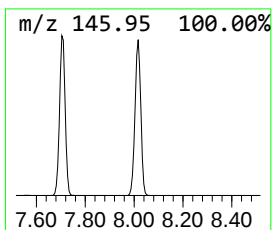
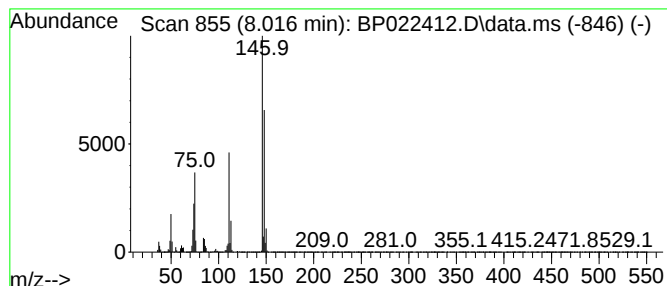
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION

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.
8.016	18.83 ng/ul	5764460	1,4-Dichlorobenzene-d4	7.669

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	97
4			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	97
5			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	96



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ALS Vial : 4 Sample Multiplier: 1

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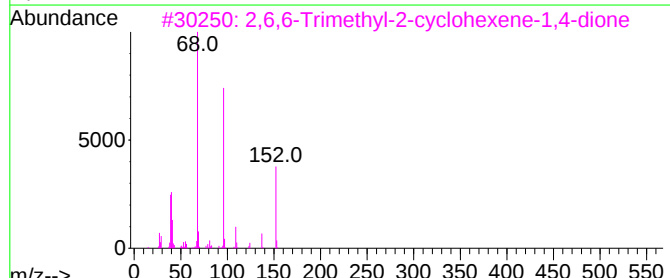
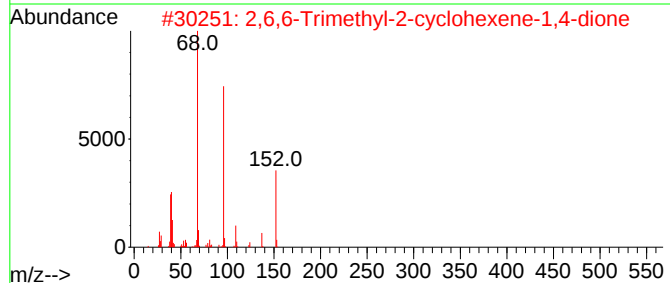
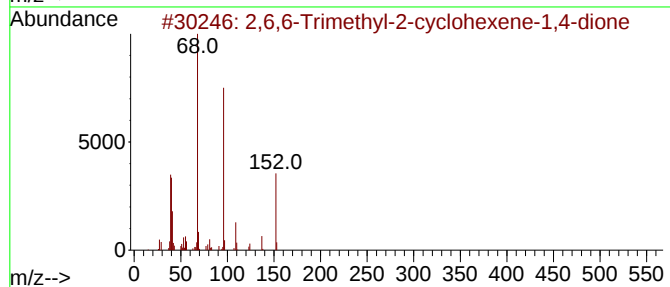
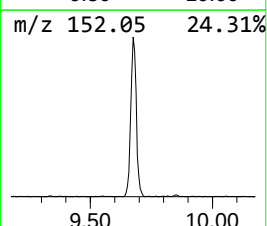
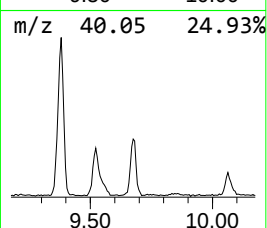
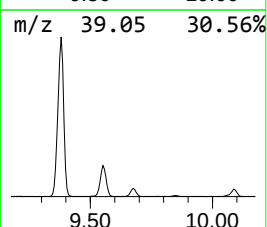
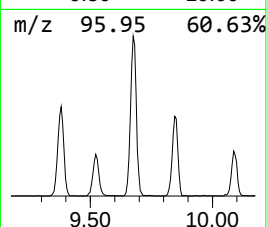
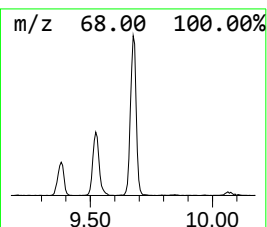
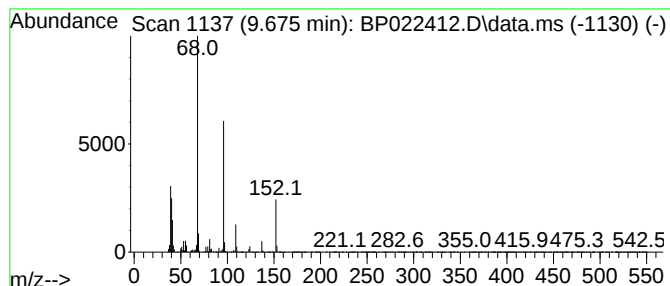
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 2,6,6-Trimethyl-2-cyclohexe... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.675	6.10 ng/ul	211557	Naphthalene-d8	10.434

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



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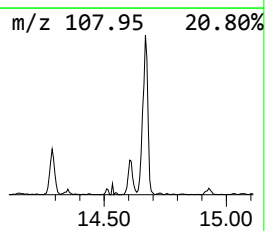
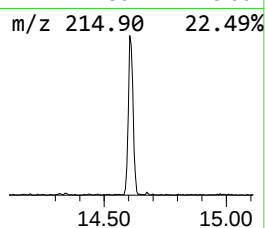
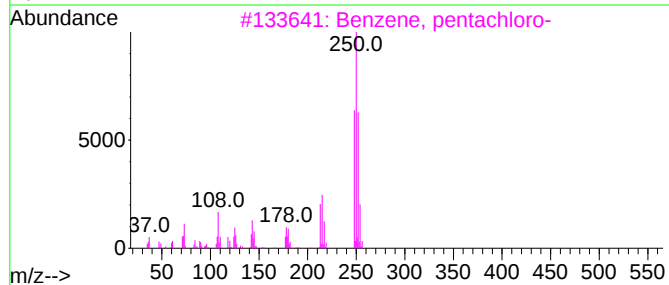
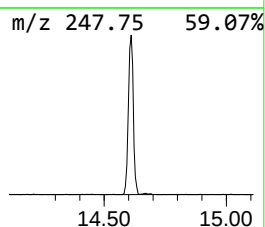
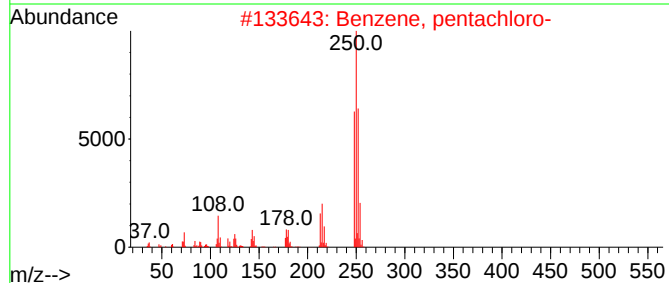
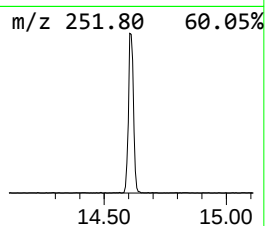
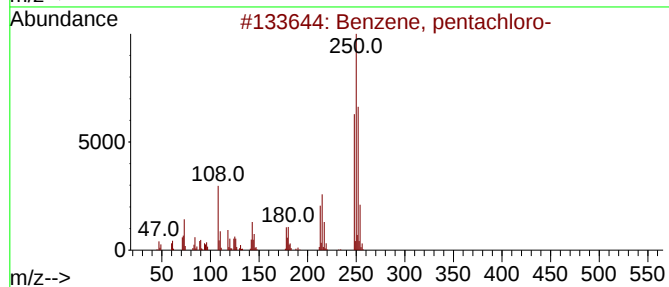
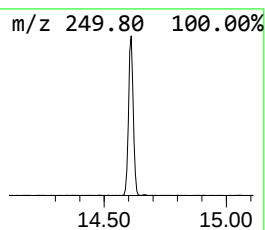
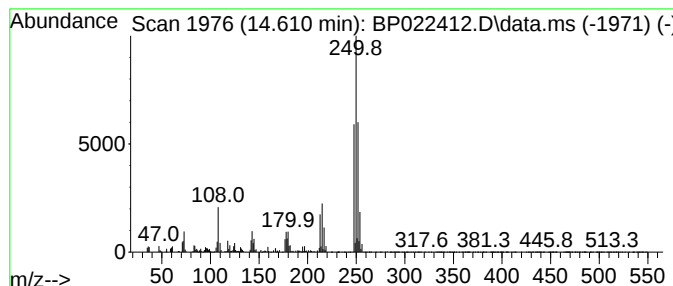
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Quant Title : SVOA CALIBRATION

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.
14.610	2.87 ng/ul	143186	Acenaphthene-d10	14.287

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	98
3			Benzene, pentachloro-	248	C6HCl5	000608-93-5	96
4			Benzene, pentachloro-	248	C6HCl5	000608-93-5	94
5			Benzene, pentachloro-	248	C6HCl5	000608-93-5	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101624\
Data File : BP022412.D
Acq On : 16 Oct 2024 12:00
Operator : RC/JU
Sample : P4329-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4E79

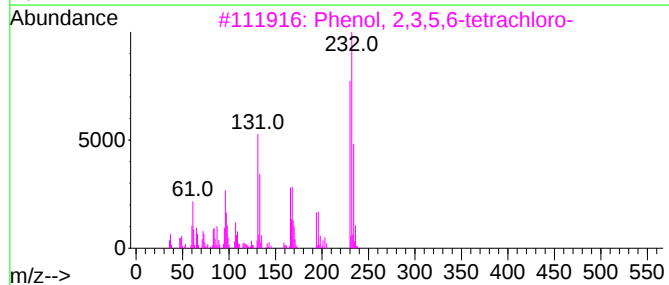
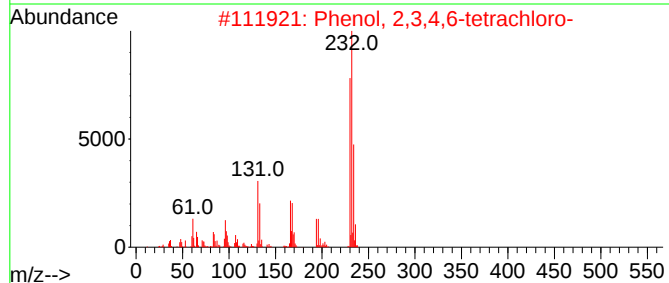
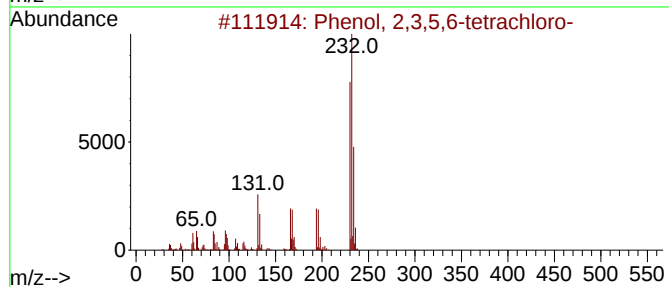
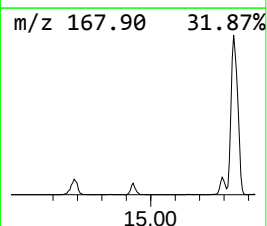
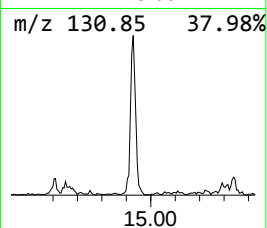
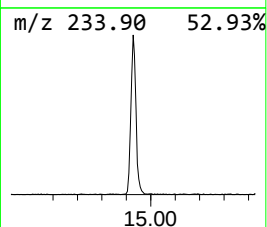
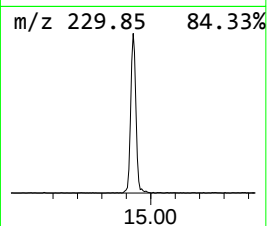
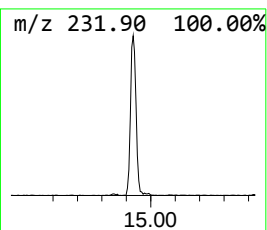
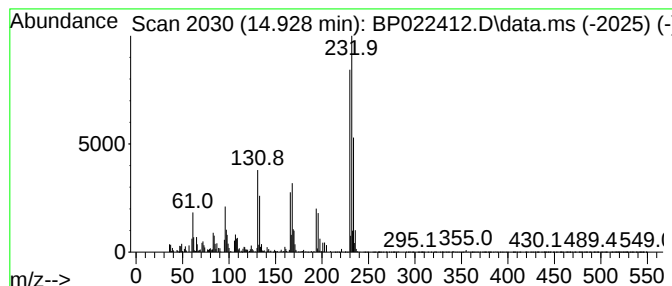
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Phenol, 2,3,5,6-tetrachloro- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.928	2.27 ng/ul	113190	Acenaphthene-d10	14.287

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 2,3,5,6-tetrachloro-	230	C6H2Cl4O	000935-95-5	97
2			Phenol, 2,3,4,6-tetrachloro-	230	C6H2Cl4O	000058-90-2	96
3			Phenol, 2,3,5,6-tetrachloro-	230	C6H2Cl4O	000935-95-5	96
4			Benzene, 1,2,3,5-tetrachloro-4-e...	258	C8H6Cl4O	056818-02-1	95
5			Phenol, 2,3,5,6-tetrachloro-	230	C6H2Cl4O	000935-95-5	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101624\
Data File : BP022412.D
Acq On : 16 Oct 2024 12:00
Operator : RC/JU
Sample : P4329-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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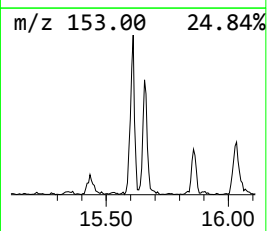
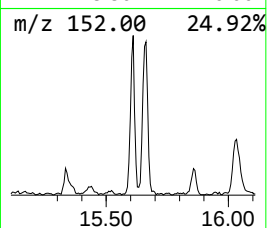
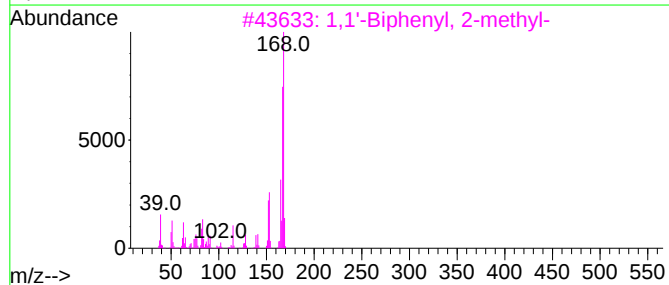
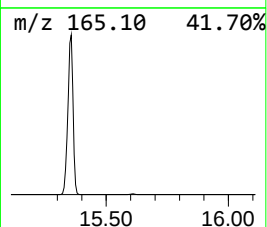
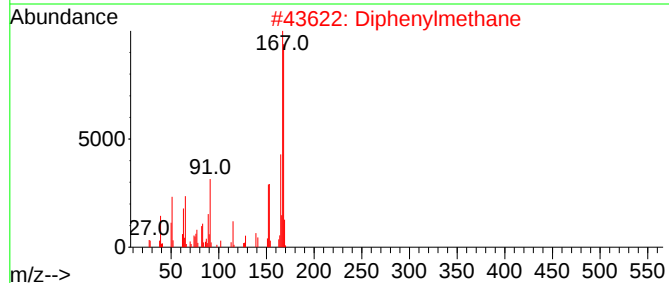
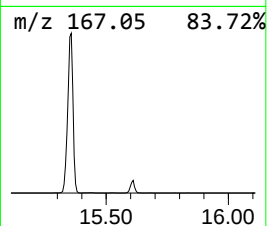
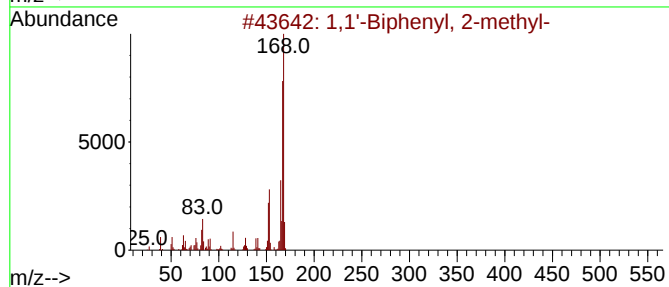
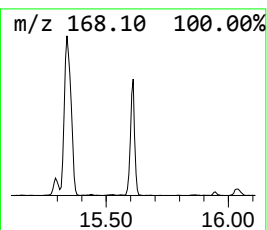
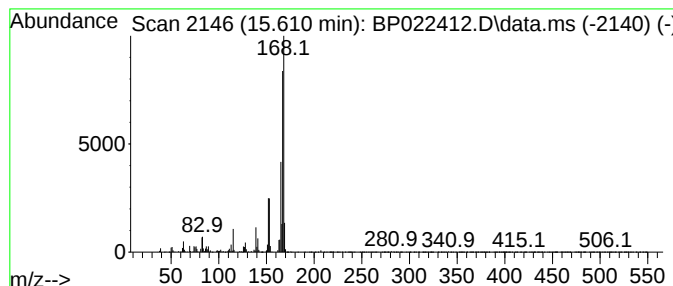
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 1,1'-Biphenyl, 2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.610	3.34 ng/ul	166489	Acenaphthene-d10	14.287

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			1,1'-Biphenyl, 2-methyl-	168	C13H12	000643-58-3	90
4			1,1'-Biphenyl, 4-methyl-	168	C13H12	000644-08-6	87
5			Diphenylmethane	168	C13H12	000101-81-5	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101624\
Data File : BP022412.D
Acq On : 16 Oct 2024 12:00
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Sample : P4329-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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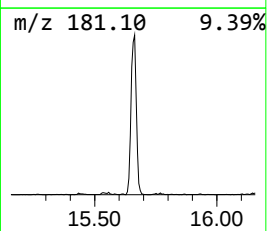
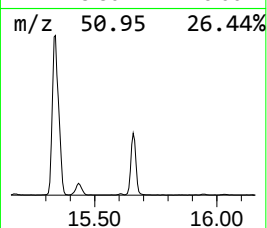
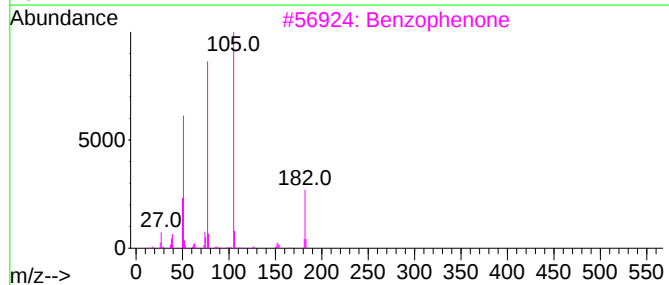
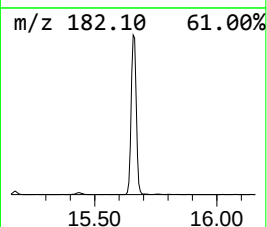
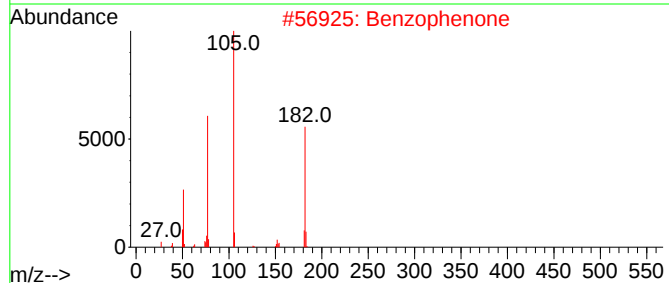
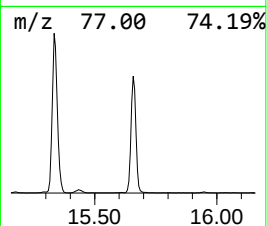
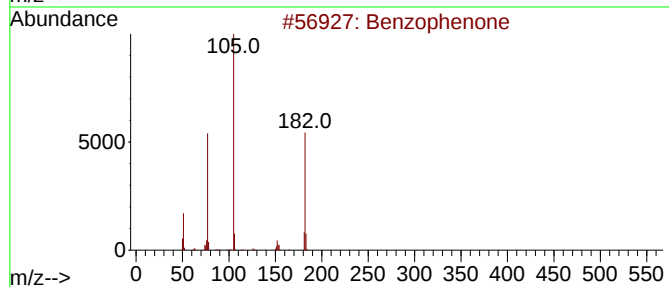
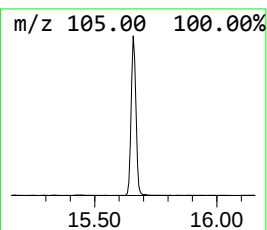
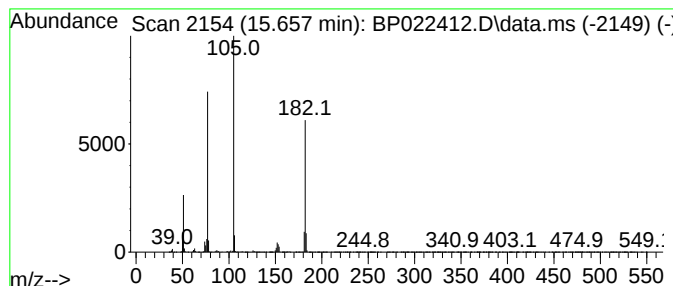
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Benzophenone Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.657	15.49 ng/ul	771411	Acenaphthene-d10	14.287

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	95
4			Benzophenone	182	C13H10O	000119-61-9	91
5			Benzophenone	182	C13H10O	000119-61-9	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101624\
Data File : BP022412.D
Acq On : 16 Oct 2024 12:00
Operator : RC/JU
Sample : P4329-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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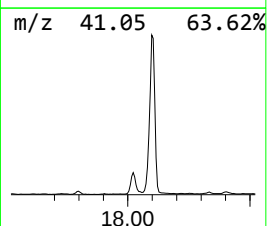
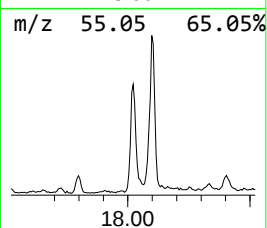
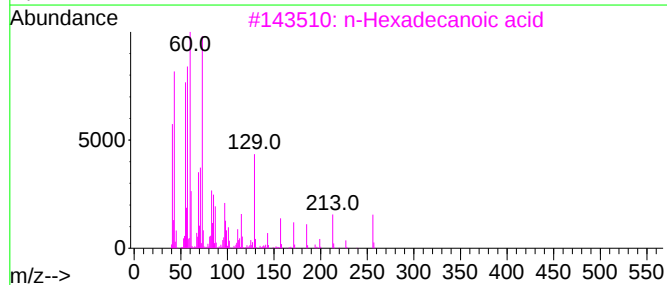
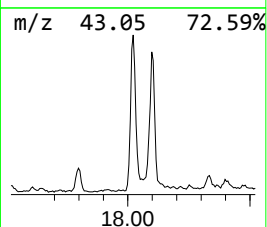
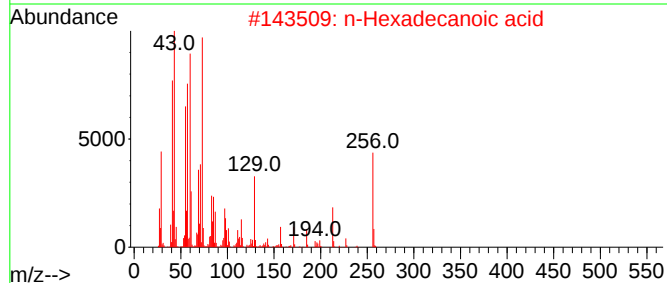
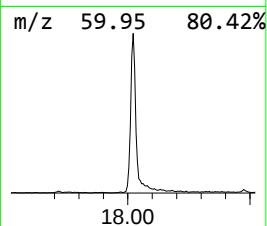
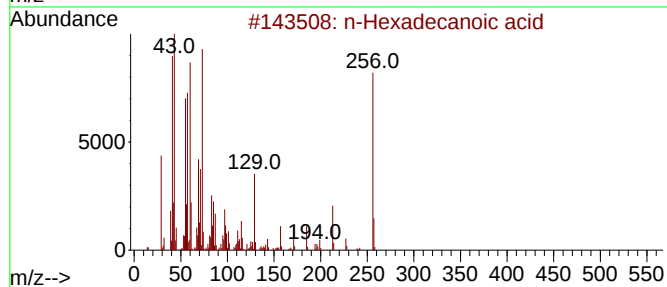
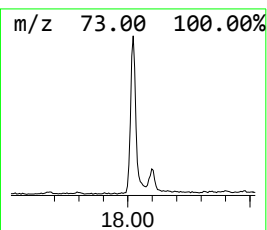
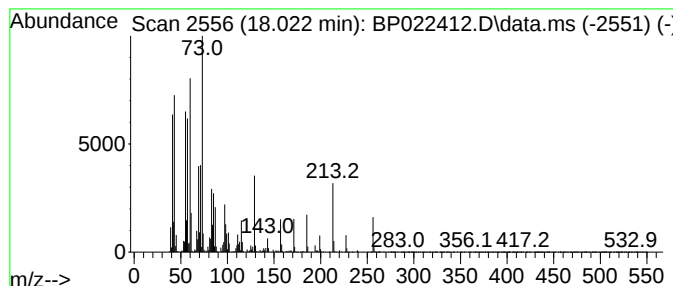
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.022	11.90 ng/ul	799131	Phenanthrene-d10	17.092

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93
4			Tridecanoic acid	214	C13H26O2	000638-53-9	89
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101624\
Data File : BP022412.D
Acq On : 16 Oct 2024 12:00
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Sample : P4329-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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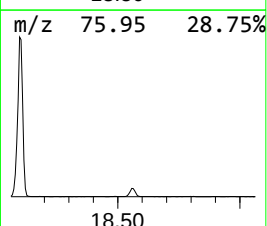
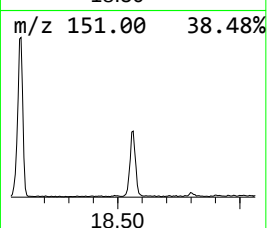
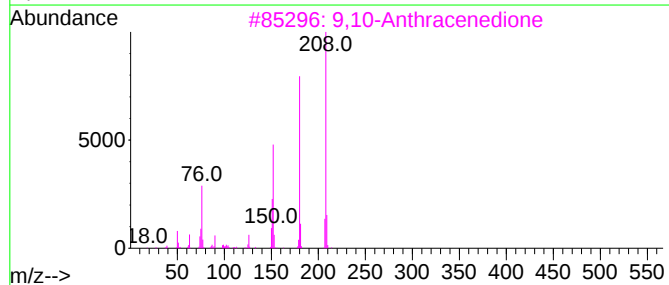
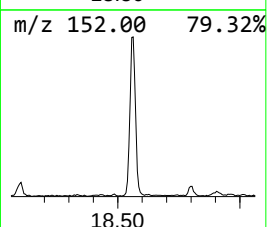
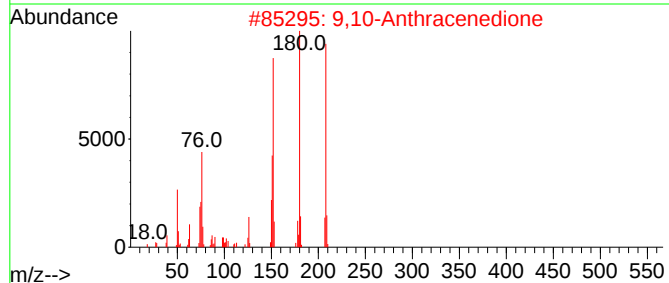
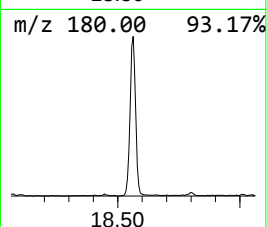
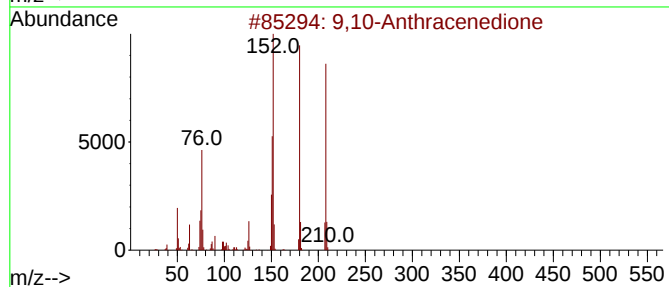
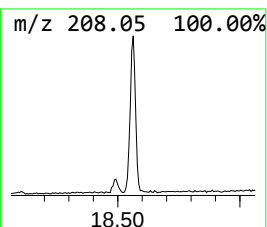
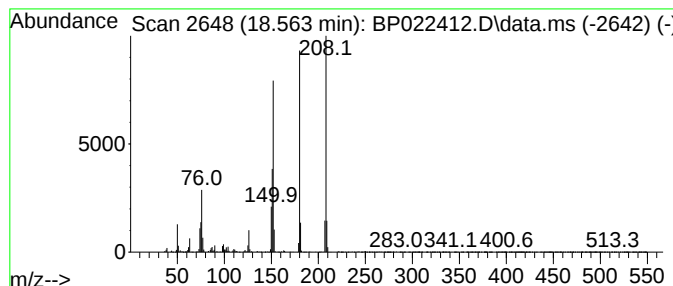
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 9,10-Anthracenedione Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.563	5.26 ng/ul	353330	Phenanthrene-d10	17.092

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	98
3		9,10-Anthracenedione	208	C14H8O2	000084-65-1	94
4		9,10-Anthracenedione	208	C14H8O2	000084-65-1	90
5		Benzo[c]cinnoline	180	C12H8N2	000230-17-1	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101624\
Data File : BP022412.D
Acq On : 16 Oct 2024 12:00
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Sample : P4329-02
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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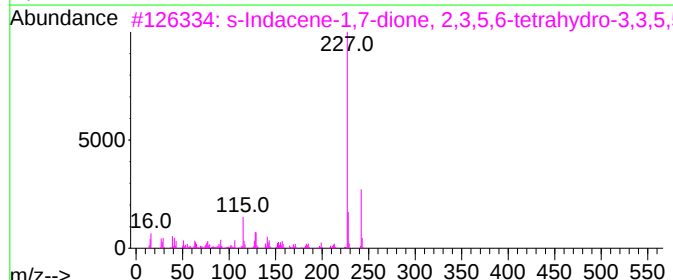
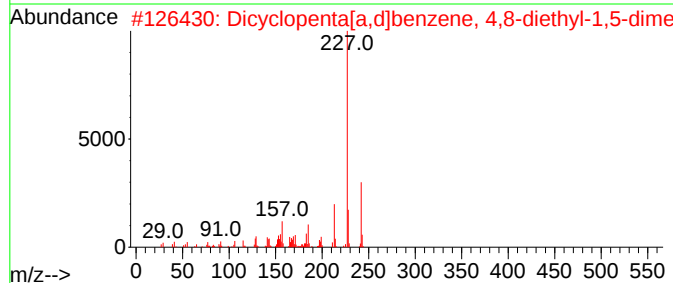
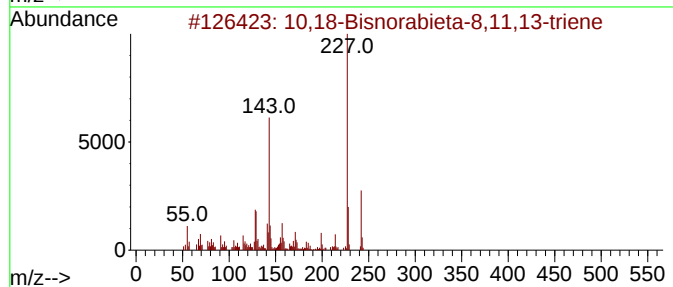
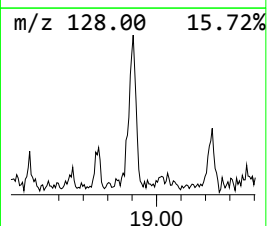
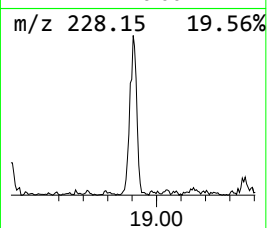
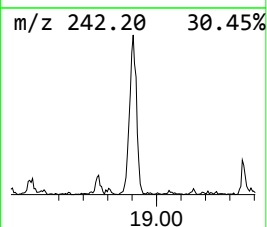
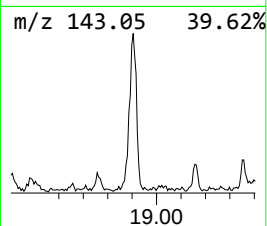
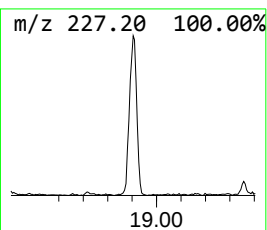
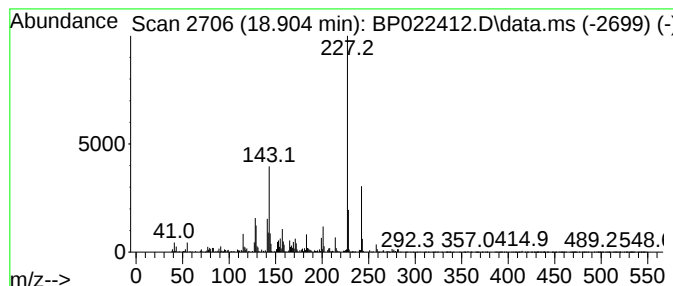
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 10,18-Bisnorabieta-8,11,13-... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.904	3.40 ng/ul	228298	Phenanthrene-d10	17.092

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		10,18-Bisnorabieta-8,11,13-triene	242	C18H26	032624-67-2	95
2		Dicyclopenta[a,d]benzene, 4,8-di...	242	C18H26	1000156-41-6	83
3		s-Indacene-1,7-dione, 2,3,5,6-te...	242	C16H18O2	055591-17-8	60
4		3-Cyano-7-(diethylamino)coumarin	242	C14H14N2O2	051473-74-6	50
5		7-Diethylamino-2-oxo-2H-chromene...	242	C14H14N2O2	332411-59-3	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101624\
Data File : BP022412.D
Acq On : 16 Oct 2024 12:00
Operator : RC/JU
Sample : P4329-02
Misc :
ALS Vial : 4 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
Benzene, 1,2-di...	8.016	18.8	ng/ul	5764460	1	7.669	6121570	20.0
2,6,6-Trimethyl...	9.675	6.1	ng/ul	211557	2	10.434	693889	20.0
Benzene, pentac...	14.610	2.9	ng/ul	143186	3	14.287	996243	20.0
Phenol, 2,3,5,6...	14.928	2.3	ng/ul	113190	3	14.287	996243	20.0
1,1'-Biphenyl, ...	15.610	3.3	ng/ul	166489	3	14.287	996243	20.0
Benzophenone	15.657	15.5	ng/ul	771411	3	14.287	996243	20.0
n-Hexadecanoic ...	18.022	11.9	ng/ul	799131	4	17.092	1343450	20.0
9,10-Anthracene...	18.563	5.3	ng/ul	353330	4	17.092	1343450	20.0
10,18-Bisnorabi...	18.904	3.4	ng/ul	228298	4	17.092	1343450	20.0