

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101124\
 Data File : BM048018.D
 Acq On : 11 Oct 2024 19:55
 Operator : RC/JU
 Sample : P4237-15
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4EN7

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_M\Methods\SFAM-EPA-BM092724.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BM048018.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.240	40	43	50	rVB	26187	33649	0.60%	0.058%
2	3.669	113	116	121	rBV4	19451	26892	0.48%	0.046%
3	3.757	127	131	136	rVB2	31109	39362	0.70%	0.067%
4	3.981	166	169	174	rVB	23596	32962	0.59%	0.057%
5	4.781	300	305	309	rBV3	16402	26575	0.48%	0.046%
6	4.892	320	324	331	rVB	43313	63654	1.14%	0.109%
7	6.045	516	520	530	rVB4	30942	46676	0.84%	0.080%
8	6.957	665	675	689	rBV	329164	638907	11.44%	1.096%
9	7.116	695	702	712	rBV	272286	447594	8.01%	0.768%
10	7.316	730	736	743	rBV	356159	586761	10.51%	1.006%
11	7.781	808	815	823	rBV	762544	1233410	22.08%	2.115%
12	7.851	823	827	837	rVB4	19415	38587	0.69%	0.066%
13	8.492	929	936	943	rBV	352145	581991	10.42%	0.998%
14	8.604	950	955	963	rVB	73574	123874	2.22%	0.212%
15	8.933	1005	1011	1023	rBV	301238	520922	9.33%	0.893%
16	9.504	1103	1108	1116	rVB2	30978	52338	0.94%	0.090%
17	9.657	1126	1134	1141	rBV	250678	432928	7.75%	0.742%
18	9.822	1156	1162	1167	rBV7	13647	26257	0.47%	0.045%
19	10.198	1220	1226	1238	rBV	364725	662198	11.86%	1.136%
20	10.574	1282	1290	1296	rBV	1058090	1738469	31.13%	2.981%
21	10.622	1296	1298	1306	rVB	36070	54664	0.98%	0.094%
22	10.716	1306	1314	1328	rBV2	65955	169677	3.04%	0.291%
23	11.121	1377	1383	1392	rBV2	14567	36726	0.66%	0.063%
24	11.369	1416	1425	1430	rBV9	9539	22634	0.41%	0.039%
25	12.239	1567	1573	1579	rBV2	23140	46639	0.84%	0.080%
26	12.457	1605	1610	1614	rBV2	23878	34254	0.61%	0.059%
27	13.251	1739	1745	1751	rBV3	21928	37185	0.67%	0.064%
28	13.592	1797	1803	1811	rVB9	13956	38480	0.69%	0.066%
29	13.745	1818	1829	1833	rBV2	28327	58362	1.04%	0.100%
30	13.827	1833	1843	1854	rVV	649844	990623	17.74%	1.699%
31	14.110	1880	1891	1905	rBV	863493	1435881	25.71%	2.462%
32	14.262	1911	1917	1922	rBV9	13851	26257	0.47%	0.045%
33	14.415	1933	1943	1950	rBV	1518575	2246635	40.23%	3.853%
34	14.480	1950	1954	1961	rVB	108237	169036	3.03%	0.290%
35	14.539	1961	1964	1970	rVB7	16806	28199	0.50%	0.048%
36	14.627	1973	1979	1985	rBV	160753	315219	5.64%	0.541%

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Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
 Title : SVOA CALIBRATION

37	14.815	2007	2011	2019	rVV	61058	114094	2.04%	0.196%
38	14.892	2019	2024	2028	rVB5	20916	35978	0.64%	0.062%
39	15.033	2042	2048	2053	rBV5	13661	30057	0.54%	0.052%
40	15.086	2053	2057	2062	rVB3	17312	26634	0.48%	0.046%
41	15.209	2074	2078	2085	rBV8	13955	31215	0.56%	0.054%
42	15.274	2085	2089	2096	rVB2	20656	35537	0.64%	0.061%
43	15.409	2105	2112	2117	rBV	1111834	1562157	27.97%	2.679%
44	15.462	2117	2121	2126	rVV	101683	148305	2.66%	0.254%
45	15.527	2127	2132	2139	rVV	139270	201994	3.62%	0.346%
46	15.586	2139	2142	2146	rVB3	16690	23687	0.42%	0.041%
47	15.633	2146	2150	2152	rBV5	22764	37543	0.67%	0.064%
48	15.668	2153	2156	2161	rVB4	36947	53846	0.96%	0.092%
49	15.774	2167	2174	2181	rVB	507799	778527	13.94%	1.335%
50	15.904	2192	2196	2204	rVB2	36591	72353	1.30%	0.124%
51	15.998	2206	2212	2218	rVB7	24509	56384	1.01%	0.097%
52	16.133	2230	2235	2243	rBV5	49636	104168	1.87%	0.179%
53	16.333	2266	2269	2274	rBV	41819	53956	0.97%	0.093%
54	16.392	2276	2279	2282	rVB	19248	25083	0.45%	0.043%
55	16.468	2289	2292	2296	rVB4	29137	40778	0.73%	0.070%
56	16.515	2297	2300	2303	rVB3	21071	23010	0.41%	0.039%
57	16.562	2305	2308	2314	rBV7	21555	28788	0.52%	0.049%
58	16.621	2314	2318	2323	rBV7	15273	23200	0.42%	0.040%
59	16.692	2326	2330	2334	rVV5	46242	69279	1.24%	0.119%
60	16.739	2334	2338	2341	rVB4	24407	30681	0.55%	0.053%
61	16.809	2346	2350	2358	rVB	67669	105010	1.88%	0.180%
62	16.915	2365	2368	2374	rVB5	29764	43560	0.78%	0.075%
63	16.974	2374	2378	2385	rVB	83640	118876	2.13%	0.204%
64	17.039	2385	2389	2392	rBV3	42837	72674	1.30%	0.125%
65	17.068	2392	2394	2401	rVB3	54054	72980	1.31%	0.125%
66	17.156	2403	2409	2413	rBV	1886993	2806412	50.25%	4.813%
67	17.198	2413	2416	2421	rVV	1509181	2010549	36.00%	3.448%
68	17.256	2421	2426	2429	rVV	1200249	1748339	31.30%	2.998%
69	17.286	2429	2431	2436	rVV	328777	422782	7.57%	0.725%
70	17.339	2436	2440	2447	rVB4	116681	200225	3.59%	0.343%
71	17.556	2474	2477	2482	rVB	122646	164053	2.94%	0.281%
72	17.609	2482	2486	2489	rBV3	57272	77199	1.38%	0.132%
73	17.756	2504	2511	2516	rBV2	263640	480276	8.60%	0.824%
74	17.874	2527	2531	2538	rBV5	76503	120369	2.16%	0.206%
75	17.992	2548	2551	2554	rBV3	71911	120218	2.15%	0.206%
76	18.050	2554	2561	2564	rBV2	769799	1212730	21.71%	2.080%

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77	18.162	2576	2580	2585	rVB	83075	120426	2.16%	0.207%
78	18.227	2585	2591	2595	rBV2	562835	899446	16.10%	1.542%
79	18.286	2598	2601	2605	rVB	198747	238857	4.28%	0.410%
80	18.327	2605	2608	2614	rBV2	152865	233587	4.18%	0.401%
81	18.515	2635	2640	2641	rBV2	183985	266256	4.77%	0.457%
82	18.568	2646	2649	2654	rVB2	196449	293183	5.25%	0.503%
83	18.686	2666	2669	2676	rVB2	130235	161805	2.90%	0.277%
84	18.950	2710	2714	2717	rBV2	112437	174243	3.12%	0.299%
85	19.044	2727	2730	2732	rBV2	157509	161326	2.89%	0.277%
86	19.086	2733	2737	2745	rVB2	295582	482400	8.64%	0.827%
87	19.209	2753	2758	2765	rVB	2956297	4014990	71.89%	6.885%
88	19.327	2775	2778	2780	rBV	208557	218756	3.92%	0.375%
89	19.544	2810	2815	2817	rBV	1878577	2646500	47.39%	4.538%
90	19.574	2817	2820	2829	rVB	2166275	2866823	51.33%	4.916%
91	20.062	2901	2903	2909	rVB2	358467	481213	8.62%	0.825%
92	20.168	2918	2921	2924	rVB	225932	246209	4.41%	0.422%
93	20.468	2968	2972	2976	rBV	532064	604385	10.82%	1.036%
94	20.539	2980	2984	2989	rVV	2823868	3043084	54.49%	5.218%
95	21.062	3070	3073	3080	rVB3	553330	801767	14.36%	1.375%
96	21.286	3108	3111	3115	rVB	672366	764452	13.69%	1.311%
97	21.380	3120	3127	3132	rVV2	2394908	5584968	100.00%	9.577%
98	21.427	3132	3135	3147	rVB	1321999	2047741	36.67%	3.512%
99	23.432	3470	3476	3482	rBV	905989	2315886	41.47%	3.971%
100	24.374	3630	3636	3649	rVB	1253769	3197878	57.26%	5.484%

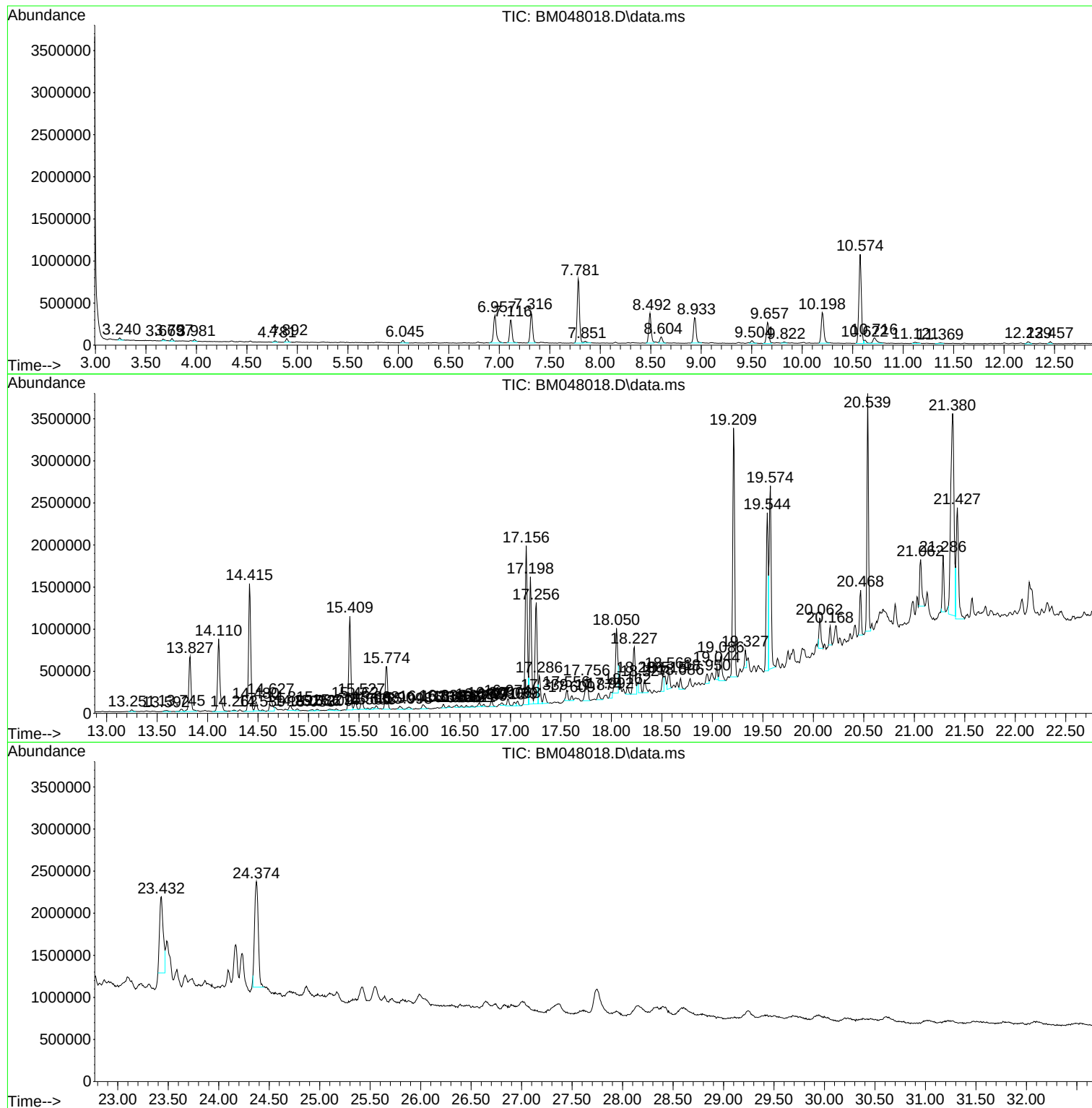
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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



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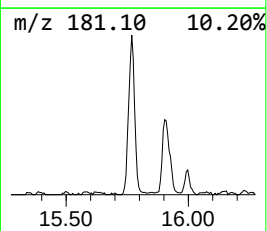
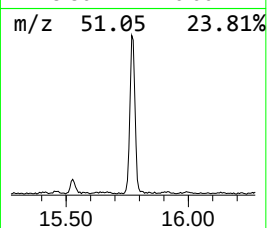
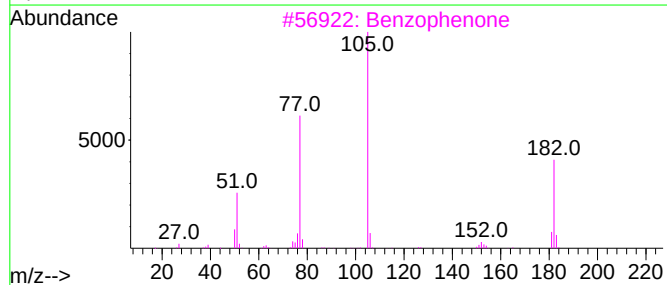
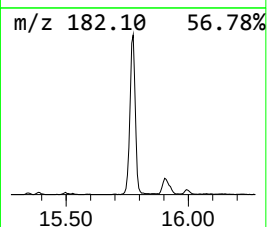
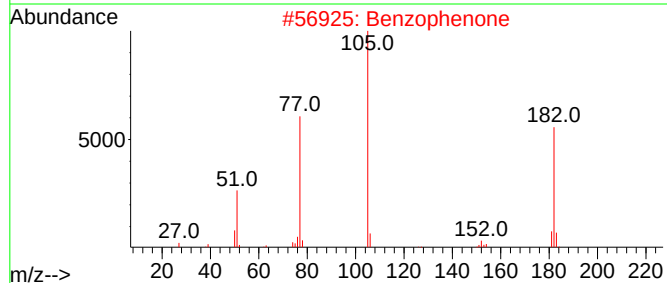
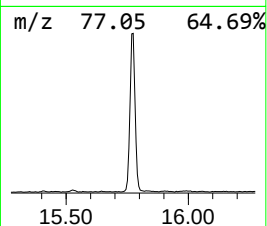
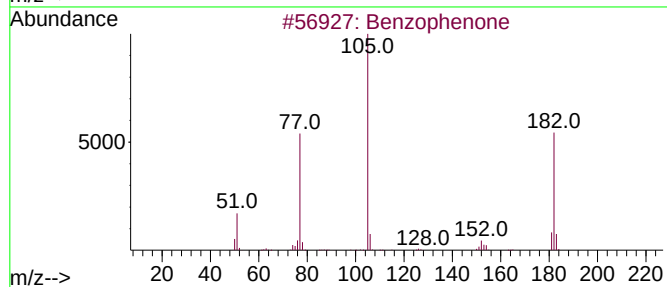
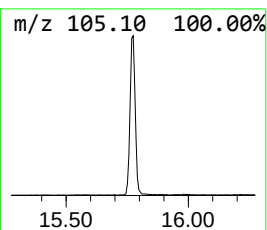
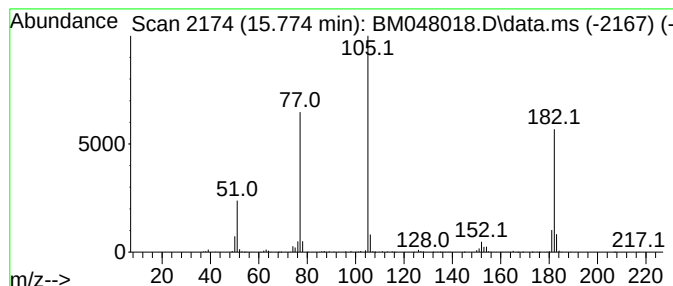
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzophenone Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.774	6.93 ng/ul	778527	Acenaphthene-d10	14.415

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			Benzophenone	182	C13H10O	000119-61-9	96
3			Benzophenone	182	C13H10O	000119-61-9	95
4			Benzophenone	182	C13H10O	000119-61-9	90
5			N-Methylbenzamide, N-chlorodiflu...	247	C10H8ClF2NO2	1000446-98-2	47



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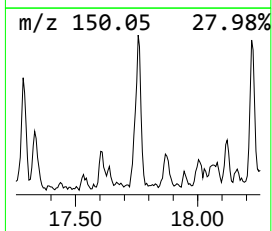
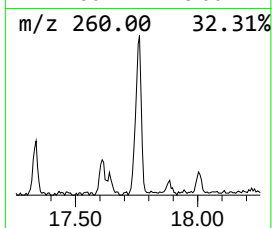
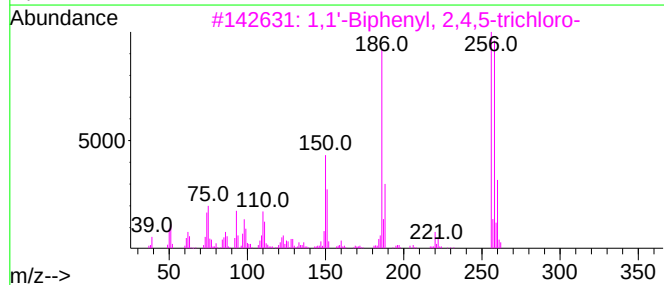
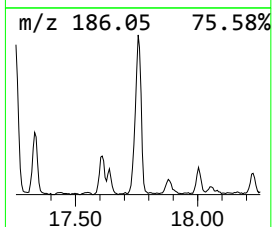
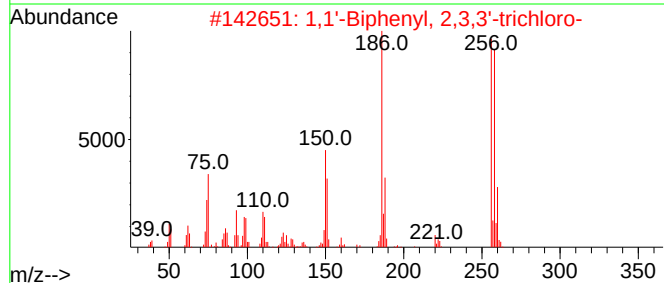
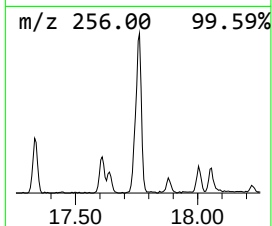
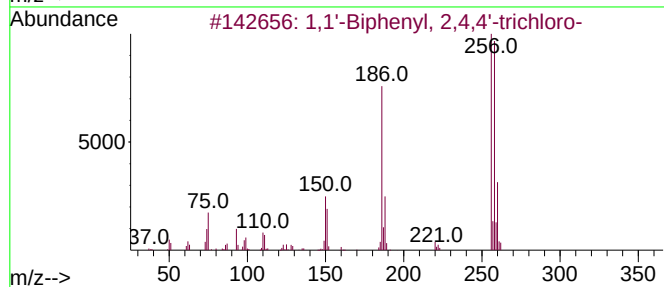
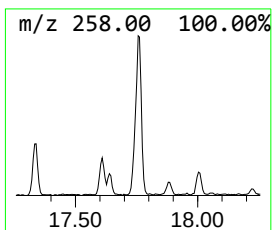
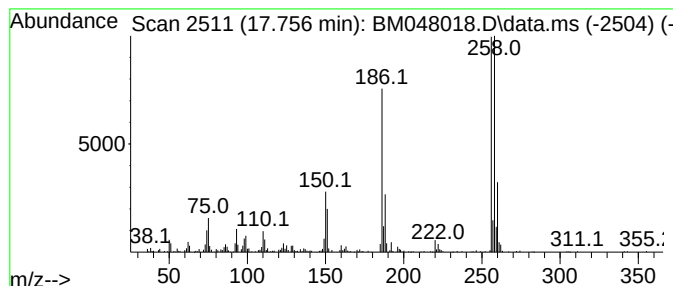
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.756	3.42 ng/ul	480276	Phenanthrene-d10	17.156	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,4,4'-trichloro-	256	C12H7Cl3	007012-37-5	99
2	1,1'-Biphenyl, 2,3,3'-trichloro-	256	C12H7Cl3	038444-84-7	99
3	1,1'-Biphenyl, 2,4,5-trichloro-	256	C12H7Cl3	015862-07-4	99
4	1,1'-Biphenyl, 2,3,4'-Trichloro-	256	C12H7Cl3	038444-85-8	98
5	1,1'-Biphenyl, 2,3',5-trichloro-	256	C12H7Cl3	038444-81-4	98



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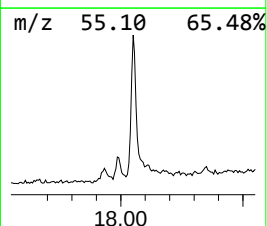
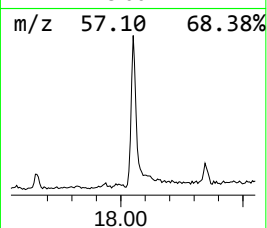
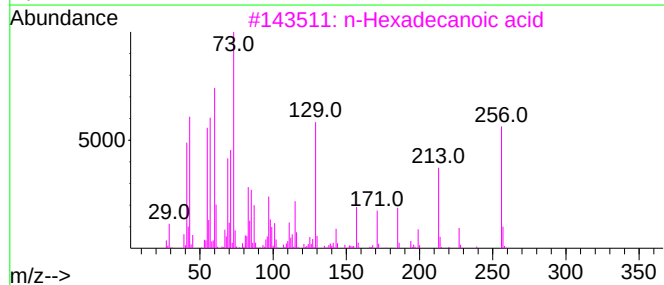
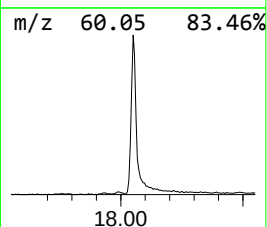
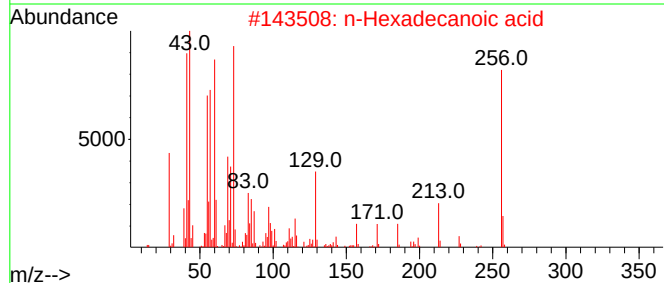
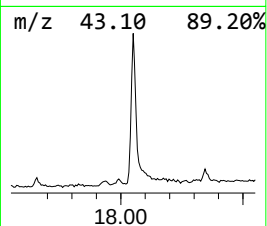
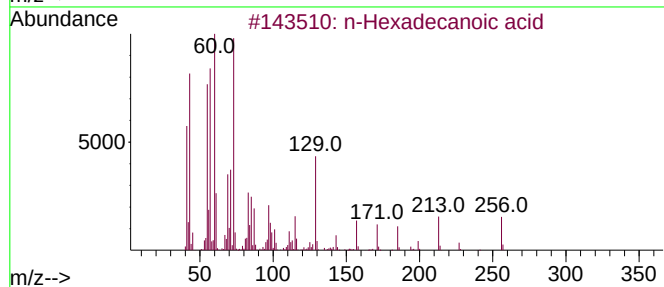
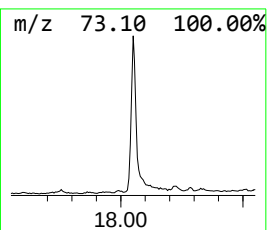
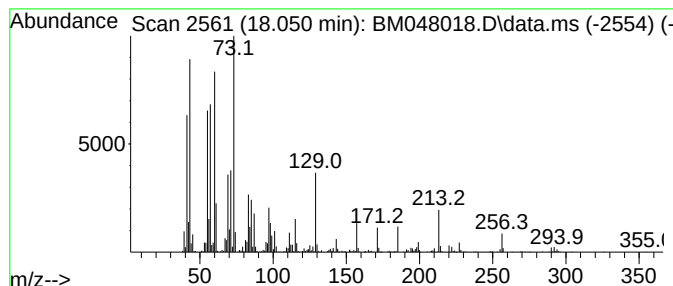
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.050	8.64 ng/ul	1212730	Phenanthrene-d10	17.156

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
5			Tridecanoic acid	214	C13H26O2	000638-53-9	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101124\
Data File : BM048018.D
Acq On : 11 Oct 2024 19:55
Operator : RC/JU
Sample : P4237-15
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
A4EN7

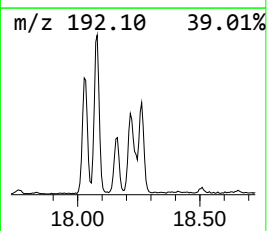
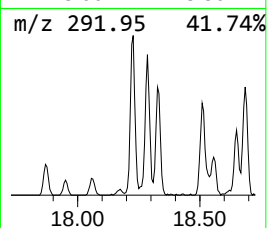
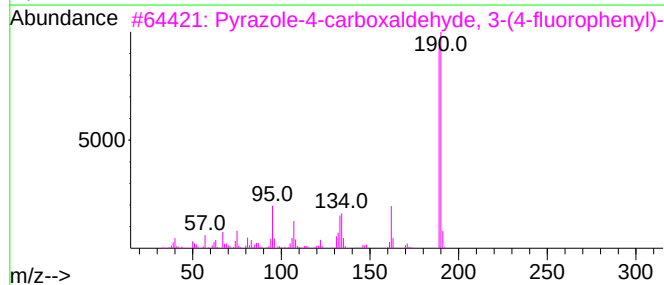
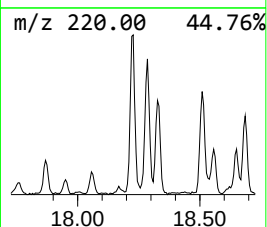
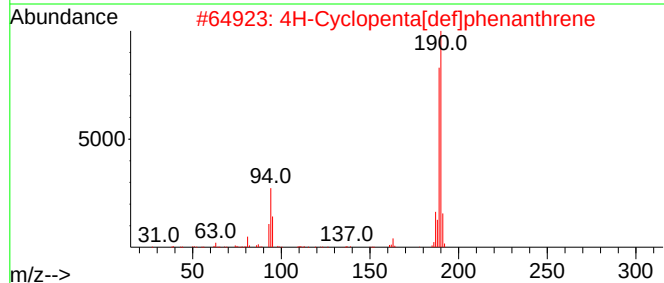
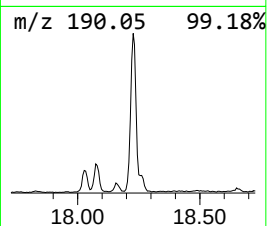
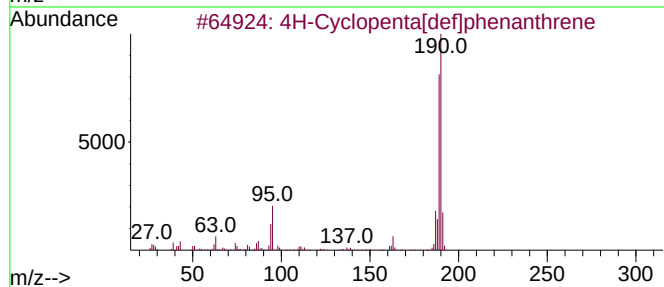
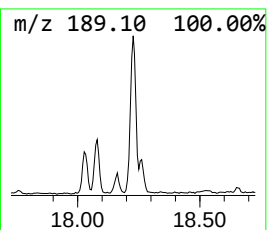
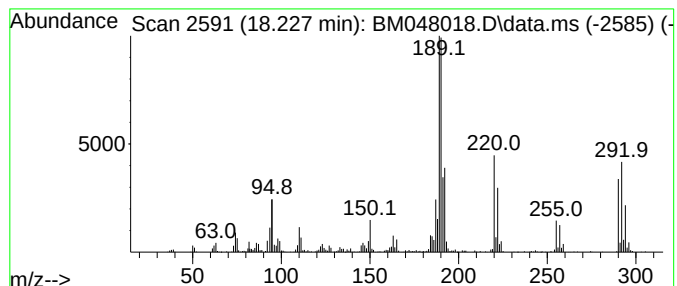
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 4H-Cyclopenta[def]phenanthrene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.227	6.41 ng/ul	899446	Phenanthrene-d10	17.156

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	50
2			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	43
3			Pyrazole-4-carboxaldehyde, 3-(4-...	190	C10H7FN2O	306936-57-2	25
4			6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	25
5			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101124\
Data File : BM048018.D
Acq On : 11 Oct 2024 19:55
Operator : RC/JU
Sample : P4237-15
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
A4EN7

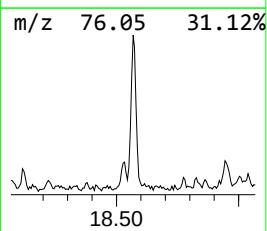
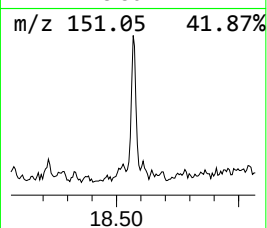
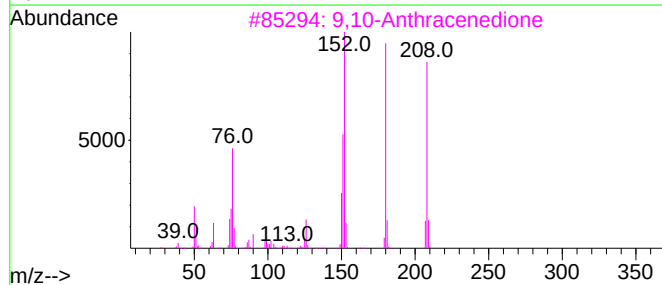
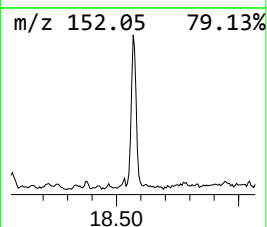
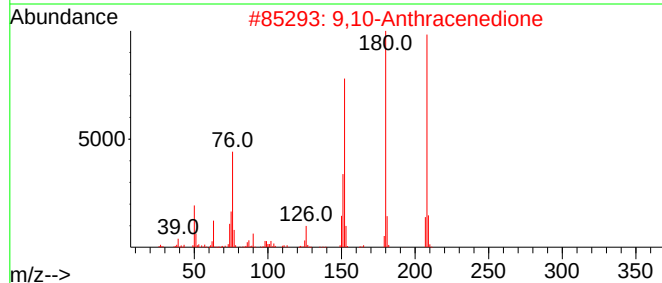
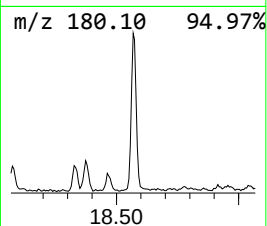
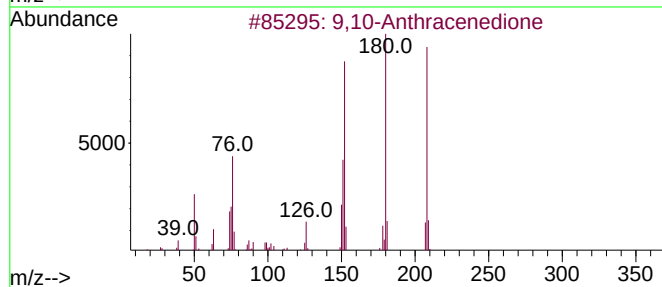
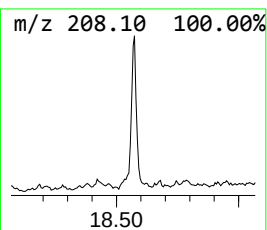
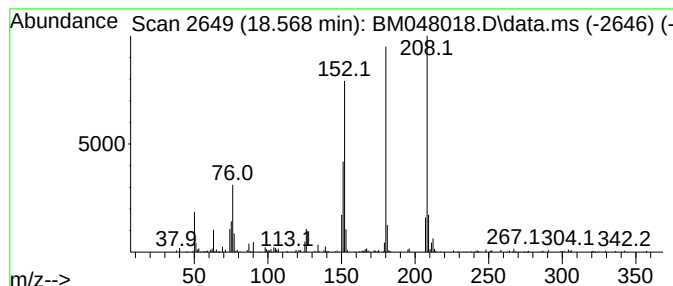
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 9,10-Anthracenedione Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.568	2.09 ng/ul	293183	Phenanthrene-d10	17.156

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	95
4	Benzo[c]cinnoline			180	C12H8N2	000230-17-1	64
5	Benzo[h]cinnoline			180	C12H8N2	000230-31-9	62



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101124\
Data File : BM048018.D
Acq On : 11 Oct 2024 19:55
Operator : RC/JU
Sample : P4237-15
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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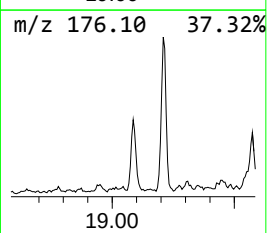
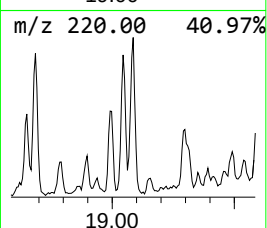
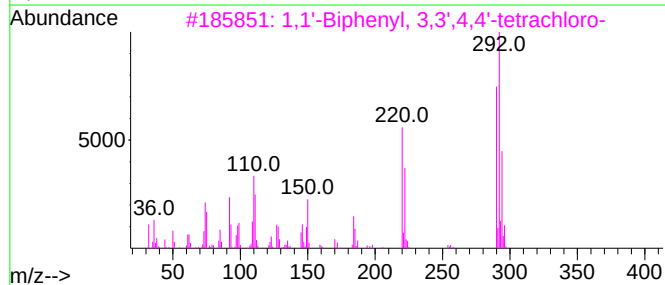
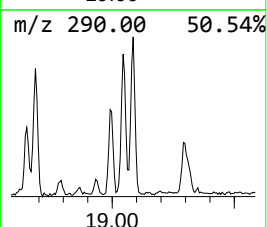
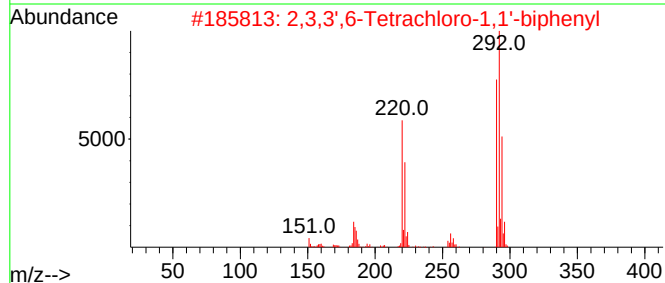
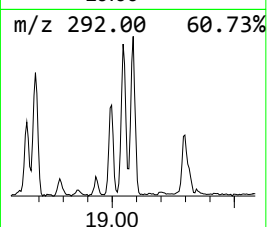
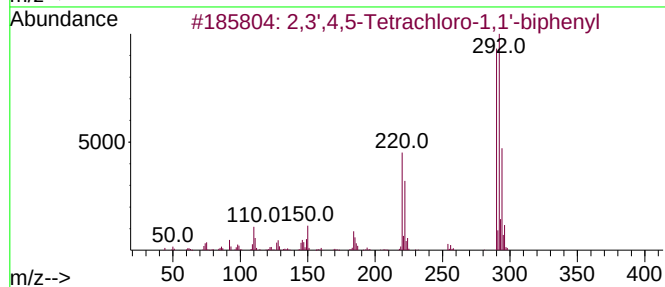
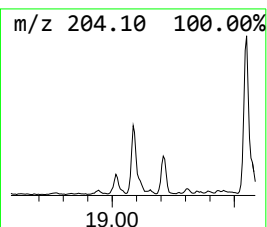
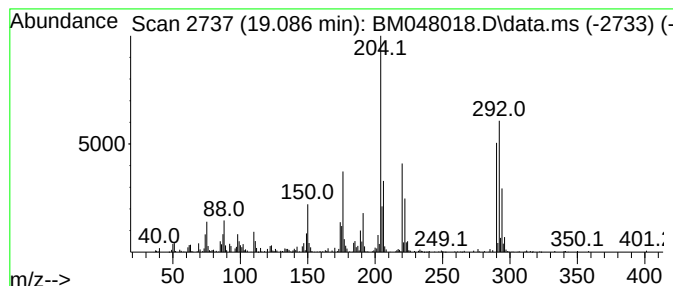
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 2,3',4,5-Tetrachloro-1,1'-b... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.086	3.44 ng/ul	482400	Phenanthrene-d10	17.156

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,3',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	073575-53-8	81		
2	2,3,3',6-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	074472-33-6	64		
3	1,1'-Biphenyl, 3,3',4,4'-tetrach...	290	C12H6Cl4	032598-13-3	64		
4	1,1'-Biphenyl, 2,3,3',4'-tetrach...	290	C12H6Cl4	041464-43-1	62		
5	1,1'-Biphenyl, 2,3,4',6-tetrachl...	290	C12H6Cl4	052663-58-8	56		



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101124\
Data File : BM048018.D
Acq On : 11 Oct 2024 19:55
Operator : RC/JU
Sample : P4237-15
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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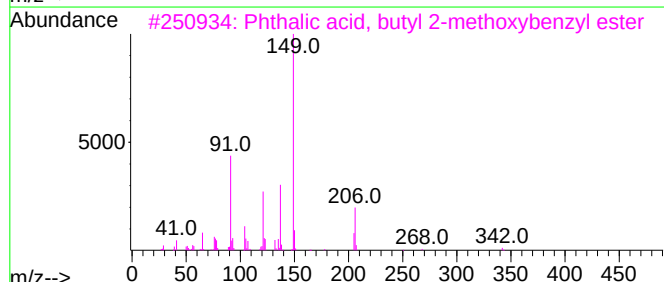
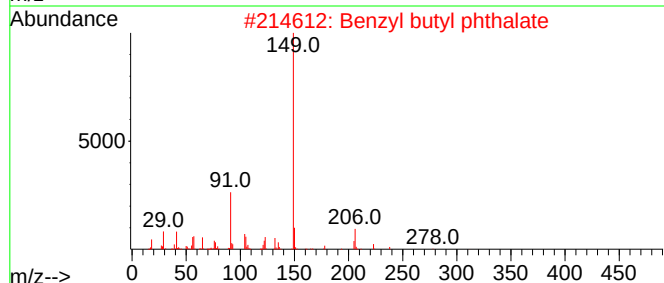
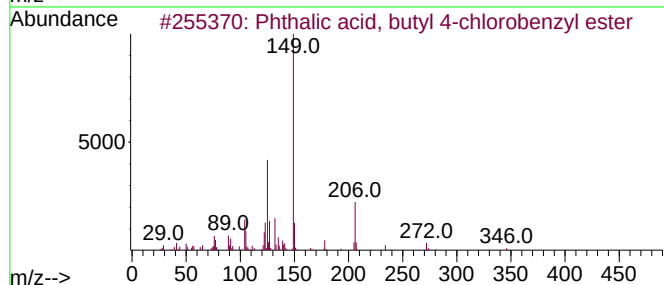
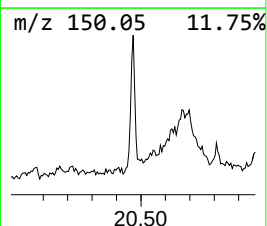
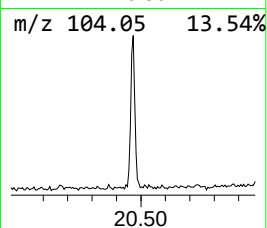
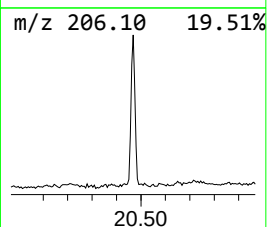
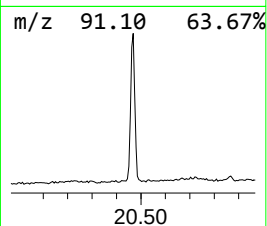
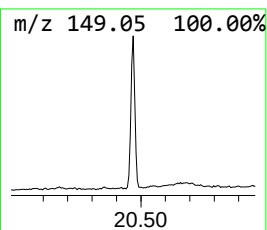
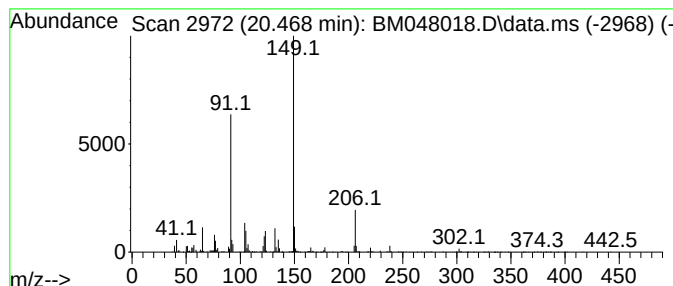
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Phthalic acid, butyl 4-chlo... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.468	2.16 ng/ul	604385	Chrysene-d12	21.386

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phthalic acid, butyl 4-chloroben...	346	C19H19ClO4	1000309-03-0	90
2			Benzyl butyl phthalate	312	C19H20O4	000085-68-7	86
3			Phthalic acid, butyl 2-methoxybe...	342	C20H22O5	1000382-49-3	72
4			Benzyl butyl phthalate	312	C19H20O4	000085-68-7	70
5			Benzyl butyl phthalate	312	C19H20O4	000085-68-7	62



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101124\
Data File : BM048018.D
Acq On : 11 Oct 2024 19:55
Operator : RC/JU
Sample : P4237-15
Misc :
ALS Vial : 9 Sample Multiplier: 1

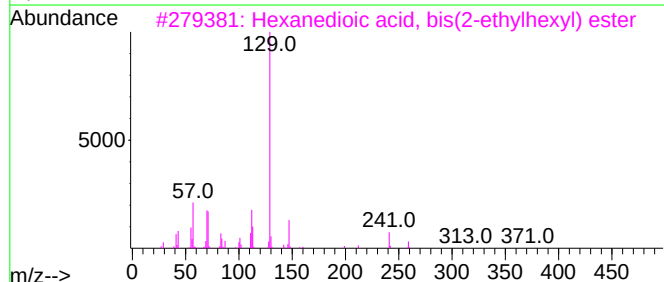
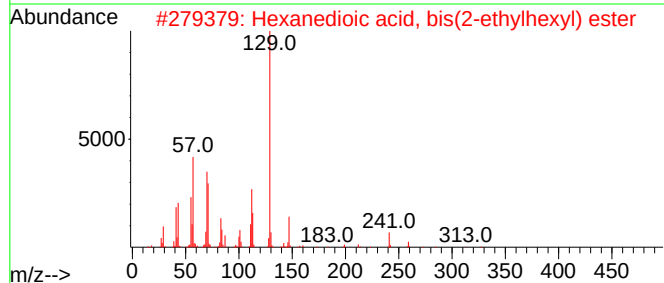
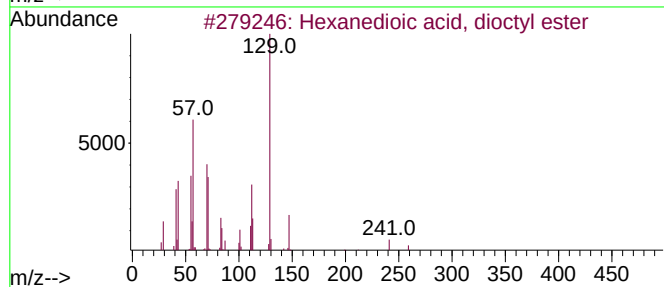
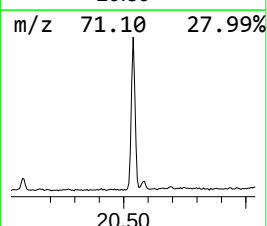
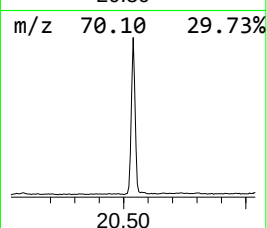
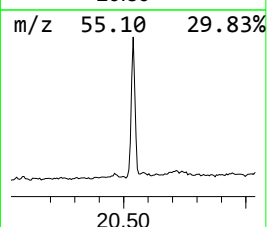
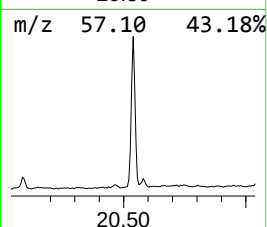
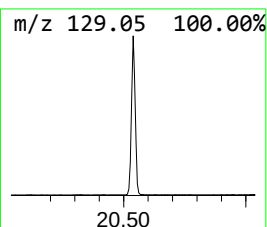
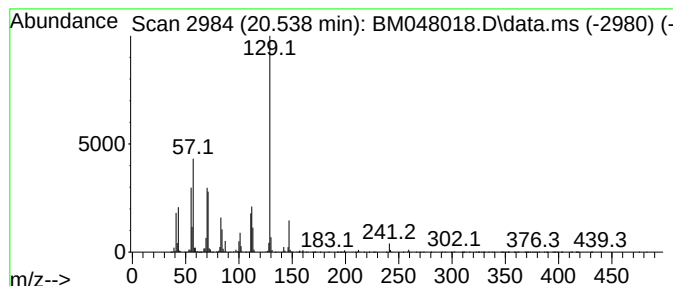
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 8 Hexanedioic acid, dioctyl e... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
20.538	10.90 ng/ul	3043080	Chrysene-d12	21.386	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexanedioic acid, dioctyl ester	370	C22H42O4	000123-79-5	95
2	Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	91
3	Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	91
4	Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	89
5	Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101124\
Data File : BM048018.D
Acq On : 11 Oct 2024 19:55
Operator : RC/JU
Sample : P4237-15
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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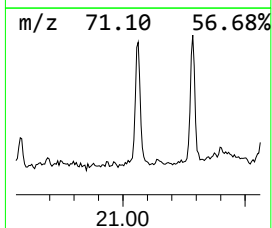
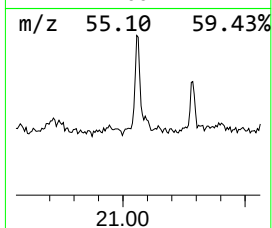
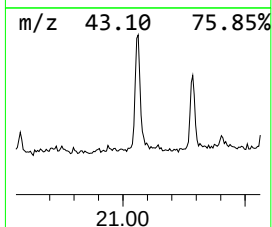
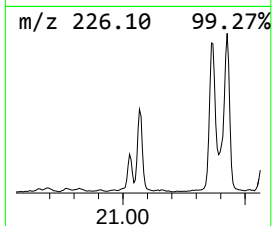
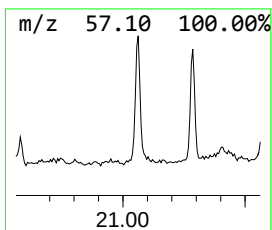
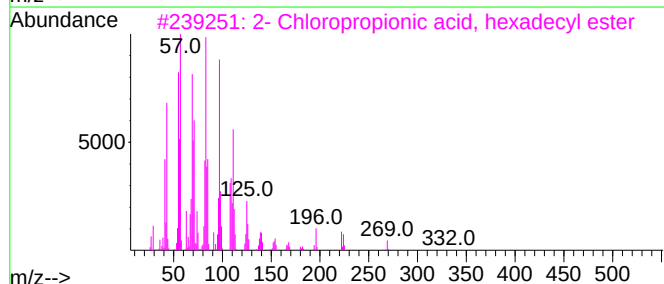
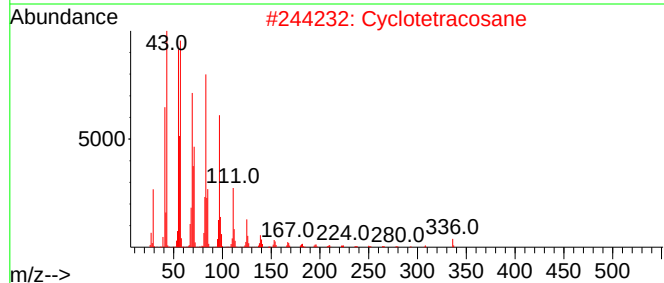
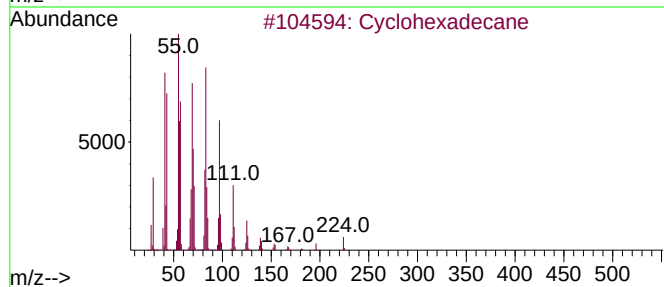
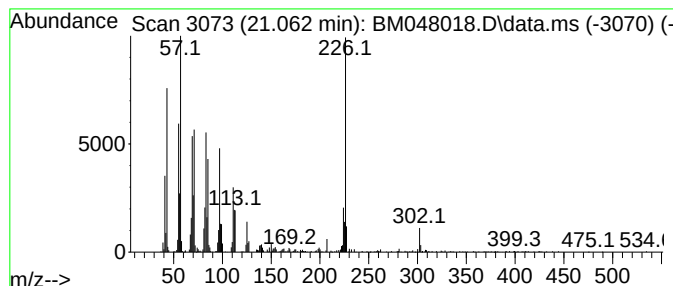
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Cyclic21.062 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.062	2.87 ng/ul	801767	Chrysene-d12	21.386

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexadecane	224	C16H32	000295-65-8	86
2			Cyclotetracosane	336	C24H48	000297-03-0	76
3			2- Chloropropionic acid, hexadec...	332	C19H37ClO2	086711-81-1	51
4			Nonadecyl trifluoroacetate	380	C21H39F3O2	1000351-76-3	50
5			Octacosyl trifluoroacetate	506	C30H57F3O2	1000351-74-9	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101124\
Data File : BM048018.D
Acq On : 11 Oct 2024 19:55
Operator : RC/JU
Sample : P4237-15
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
A4EN7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzophenone	15.774	6.9	ng/ul	778527	3	14.415	2246640	20.0
1,1'-Biphenyl, ...	17.756	3.4	ng/ul	480276	4	17.156	2806410	20.0
n-Hexadecanoic ...	18.050	8.6	ng/ul	1212730	4	17.156	2806410	20.0
4H-Cyclopenta[d...]	18.227	6.4	ng/ul	899446	4	17.156	2806410	20.0
9,10-Anthracene...	18.568	2.1	ng/ul	293183	4	17.156	2806410	20.0
2,3',4,5-Tetrac...	19.086	3.4	ng/ul	482400	4	17.156	2806410	20.0
Phthalic acid, ...	20.468	2.2	ng/ul	604385	5	21.386	5584970	20.0
Hexanedioic aci...	20.538	10.9	ng/ul	3043080	5	21.386	5584970	20.0
(DEL) Alkane: C...	21.062	2.9	ng/ul	801767	5	21.386	5584970	20.0