

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121824\
 Data File : BP023542.D
 Acq On : 20 Dec 2024 01:25
 Operator : RC/JU
 Sample : P5265-06
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 E28Z1

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-BP112624.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP023542.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.093	14	18	25	rVB	51739	75265	0.87%	0.065%
2	3.505	84	88	111	rBV	603619	1041273	12.01%	0.893%
3	3.805	136	139	149	rVB	18473	34523	0.40%	0.030%
4	6.222	542	550	556	rVB10	11090	26581	0.31%	0.023%
5	6.664	621	625	632	rVV5	13702	27551	0.32%	0.024%
6	6.740	632	638	651	rVV	823646	1648667	19.02%	1.413%
7	6.875	655	661	675	rVB	743168	1176891	13.58%	1.009%
8	7.075	688	695	707	rBV	889062	1467556	16.93%	1.258%
9	7.516	759	770	779	rBV	1033807	1629033	18.79%	1.396%
10	8.240	886	893	911	rBV	1008104	1956440	22.57%	1.677%
11	8.669	959	966	984	rBV	820288	1500943	17.31%	1.287%
12	8.822	987	992	1000	rVB8	16163	34018	0.39%	0.029%
13	9.052	1026	1031	1036	rBV2	24767	41987	0.48%	0.036%
14	9.228	1055	1061	1068	rBV	56885	100139	1.16%	0.086%
15	9.381	1079	1087	1098	rBV	663187	1226968	14.15%	1.052%
16	9.669	1129	1136	1144	rBV	13943	42102	0.49%	0.036%
17	9.922	1172	1179	1197	rBV	983828	1999384	23.06%	1.714%
18	10.263	1227	1237	1243	rBV	1392495	2334599	26.93%	2.001%
19	10.310	1243	1245	1254	rVB	48037	79239	0.91%	0.068%
20	10.428	1257	1265	1288	rBV	774280	1737483	20.04%	1.489%
21	10.834	1326	1334	1343	rVB7	20426	44357	0.51%	0.038%
22	11.675	1472	1477	1484	rVV2	24105	36212	0.42%	0.031%
23	11.787	1489	1496	1501	rBV3	14311	36164	0.42%	0.031%
24	12.646	1638	1642	1648	rVB2	24497	35003	0.40%	0.030%
25	12.793	1661	1667	1674	rBV	207655	348366	4.02%	0.299%
26	12.946	1685	1693	1699	rBV	67027	95941	1.11%	0.082%
27	13.157	1724	1729	1742	rVB2	30553	77736	0.90%	0.067%
28	13.557	1789	1797	1804	rBV	1902401	3086503	35.61%	2.646%
29	13.616	1804	1807	1810	rVB	34651	44795	0.52%	0.038%
30	13.822	1835	1842	1857	rVB2	1934296	3320822	38.31%	2.847%
31	14.034	1873	1878	1882	rBV	69945	80575	0.93%	0.069%
32	14.134	1888	1895	1903	rBV	2191392	3494136	40.31%	2.995%
33	14.204	1903	1907	1913	rVB	30463	44402	0.51%	0.038%
34	14.387	1930	1938	1941	rBV9	15197	32678	0.38%	0.028%
35	14.434	1941	1946	1962	rVV2	400353	1172517	13.53%	1.005%
36	14.546	1962	1965	1969	rVV2	46478	90475	1.04%	0.078%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
 Title : SVOA CALIBRATION

37	14.593	1969	1973	1982	rVV10	39305	111984	1.29%	0.096%
38	14.669	1983	1986	1994	rVB7	22738	35649	0.41%	0.031%
39	15.004	2039	2043	2048	rVB	45470	61369	0.71%	0.053%
40	15.145	2061	2067	2074	rBV	2798102	3894870	44.93%	3.339%
41	15.322	2091	2097	2105	rBV	243545	419287	4.84%	0.359%
42	15.416	2109	2113	2117	rVB3	42114	70485	0.81%	0.060%
43	15.522	2124	2131	2147	rVB	656413	1092422	12.60%	0.936%
44	15.881	2188	2192	2194	rBV	68360	96881	1.12%	0.083%
45	16.351	2267	2272	2277	rBV2	58998	109411	1.26%	0.094%
46	16.598	2310	2314	2320	rVB	77399	110780	1.28%	0.095%
47	16.698	2327	2331	2336	rBV2	47815	75290	0.87%	0.065%
48	16.757	2337	2341	2348	rVB2	86176	125420	1.45%	0.108%
49	16.822	2348	2352	2354	rBV	72246	87225	1.01%	0.075%
50	16.851	2354	2357	2364	rVB7	61166	97798	1.13%	0.084%
51	16.940	2365	2372	2376	rBV	3382809	4679147	53.98%	4.011%
52	16.981	2376	2379	2384	rVV	1383543	2050730	23.66%	1.758%
53	17.045	2384	2390	2400	rVB	2846250	4335556	50.01%	3.716%
54	17.134	2401	2405	2409	rBV4	47234	74720	0.86%	0.064%
55	17.375	2443	2446	2453	rVB	84834	139313	1.61%	0.119%
56	17.463	2458	2461	2465	rBV2	48239	81839	0.94%	0.070%
57	17.563	2473	2478	2482	rBV6	63192	154559	1.78%	0.132%
58	17.757	2507	2511	2516	rVB7	68665	117488	1.36%	0.101%
59	17.822	2517	2522	2528	rBV	1080569	1858010	21.43%	1.593%
60	17.887	2528	2533	2553	rVV	3405934	5980595	68.99%	5.127%
61	18.039	2555	2559	2564	rVB2	252892	396324	4.57%	0.340%
62	18.339	2607	2610	2613	rBV	121141	143187	1.65%	0.123%
63	18.422	2620	2624	2629	rBV	348216	559013	6.45%	0.479%
64	18.528	2640	2642	2648	rVB4	89953	113694	1.31%	0.097%
65	18.598	2649	2654	2657	rBV6	84326	172793	1.99%	0.148%
66	18.904	2702	2706	2709	rBV2	185606	265548	3.06%	0.228%
67	18.969	2714	2717	2722	rVB	154132	213716	2.47%	0.183%
68	19.075	2728	2735	2747	rBV	4474816	8668626	100.00%	7.431%
69	19.222	2755	2760	2766	rBV2	985658	1473168	16.99%	1.263%
70	19.428	2790	2795	2798	rBV	3681273	5479814	63.21%	4.697%
71	19.457	2798	2800	2804	rVV	3682165	3879923	44.76%	3.326%
72	19.492	2804	2806	2819	rVB	923837	1428692	16.48%	1.225%
73	19.657	2832	2834	2836	rBV	98513	98974	1.14%	0.085%
74	19.704	2839	2842	2844	rVV2	248199	299208	3.45%	0.256%
75	19.728	2844	2846	2849	rVV2	217428	236850	2.73%	0.203%
76	19.792	2854	2857	2860	rVV5	200375	267471	3.09%	0.229%

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 Title : SVOA CALIBRATION

77	20.098	2904	2909	2911	rBV4	211148	431717	4.98%	0.370%
78	20.457	2966	2970	2972	rBV4	160771	238698	2.75%	0.205%
79	20.480	2972	2974	2976	rBV2	221199	174737	2.02%	0.150%
80	20.533	2978	2983	2988	rBV3	832981	1645865	18.99%	1.411%
81	20.851	3034	3037	3039	rBV2	223025	246946	2.85%	0.212%
82	20.939	3048	3052	3058	rVB2	441389	847870	9.78%	0.727%
83	20.998	3059	3062	3064	rBV2	230995	255590	2.95%	0.219%
84	21.039	3064	3069	3075	rVV3	1323023	2381847	27.48%	2.042%
85	21.192	3092	3095	3104	rVB2	482719	881092	10.16%	0.755%
86	21.369	3119	3125	3130	rBV2	3529928	6000324	69.22%	5.143%
87	21.422	3130	3134	3139	rVB	979400	1488850	17.18%	1.276%
88	21.769	3191	3193	3194	rBV	184266	128249	1.48%	0.110%
89	22.027	3229	3237	3246	rVB	2606702	6751562	77.89%	5.787%
90	22.592	3331	3333	3336	rVB	471663	401205	4.63%	0.344%
91	22.939	3390	3392	3394	rBV	422051	247313	2.85%	0.212%
92	23.151	3426	3428	3430	rBV3	185615	215737	2.49%	0.185%
93	23.533	3489	3493	3502	rBV2	1058782	3095322	35.71%	2.653%
94	24.233	3609	3612	3619	rVB	442482	837880	9.67%	0.718%
95	24.316	3620	3626	3634	rBV	1294295	3193430	36.84%	2.737%
96	24.539	3657	3664	3672	rBV	1774045	5163099	59.56%	4.426%
97	25.021	3743	3746	3755	rVB2	663034	1287974	14.86%	1.104%
98	25.645	3850	3852	3854	rBV	182082	182099	2.10%	0.156%
99	27.063	4091	4093	4095	rBV	407460	374811	4.32%	0.321%
100	27.098	4097	4099	4103	rVB	386781	407825	4.70%	0.350%

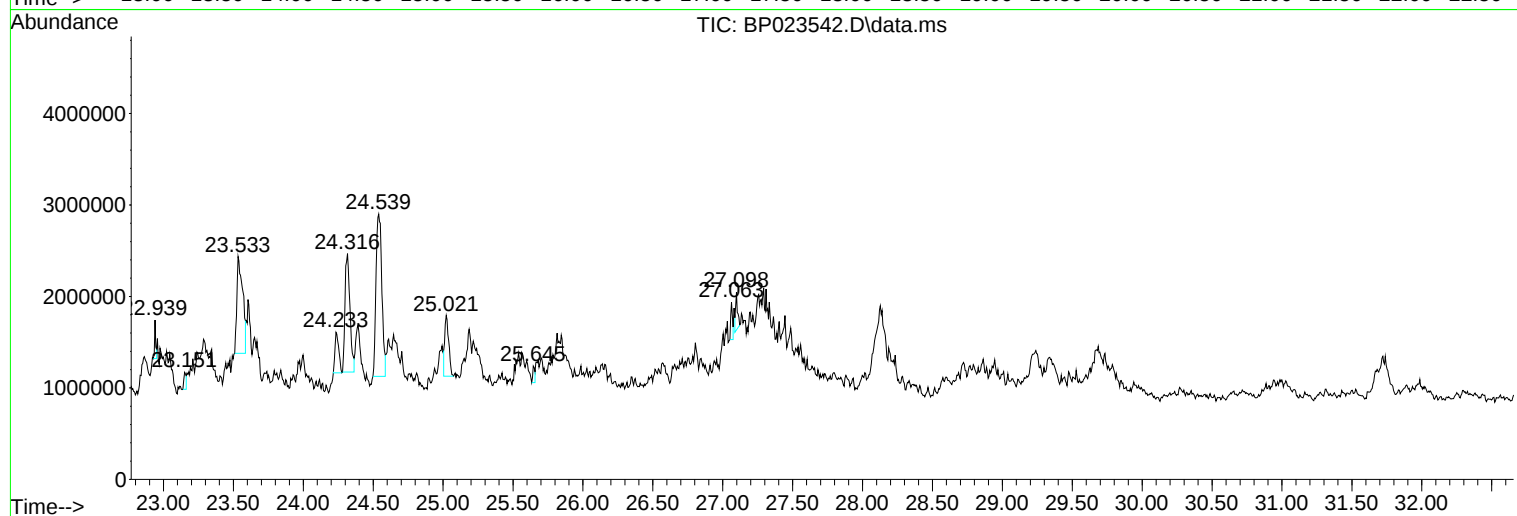
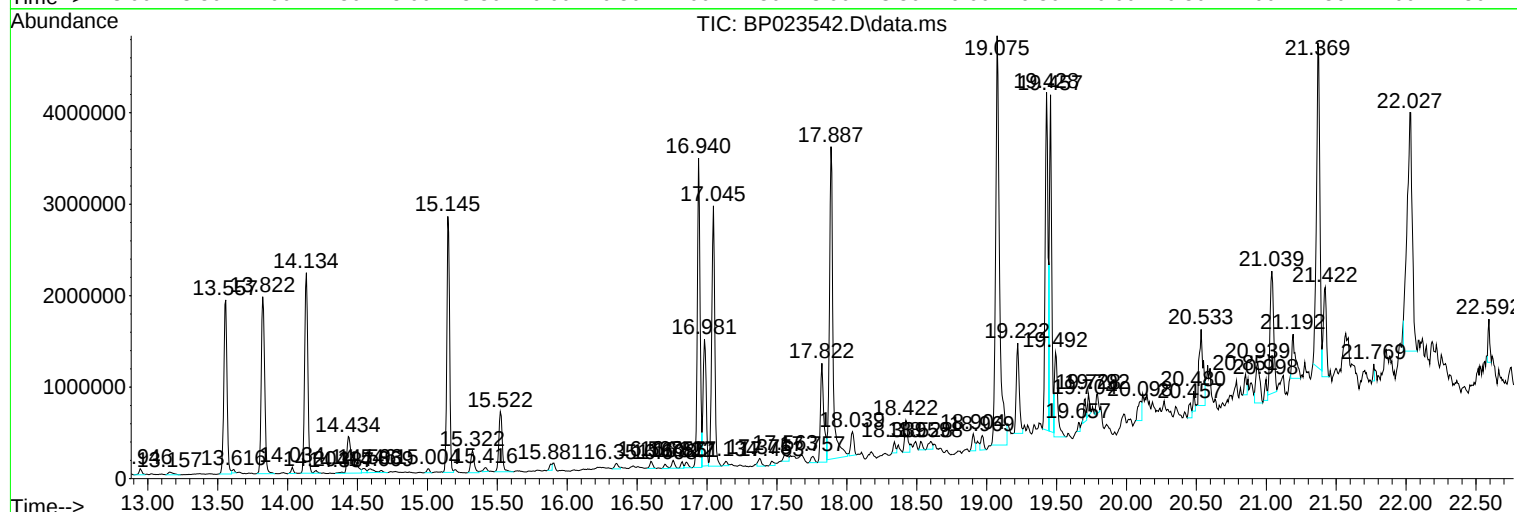
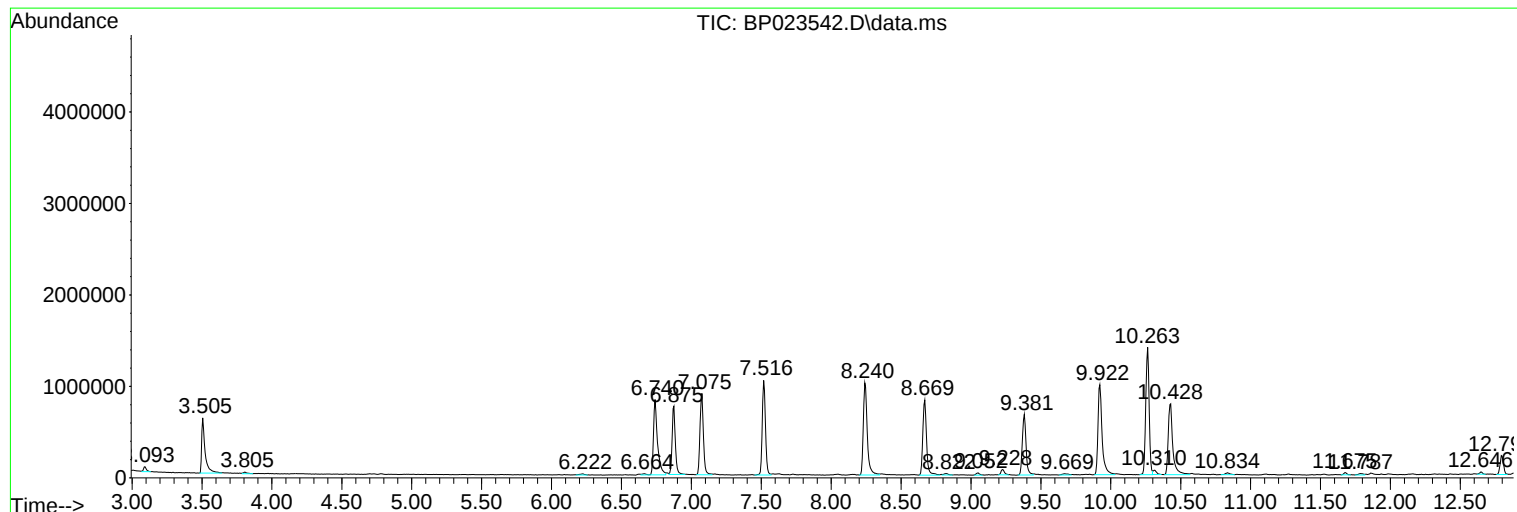
Sum of corrected areas: 116659195

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Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

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



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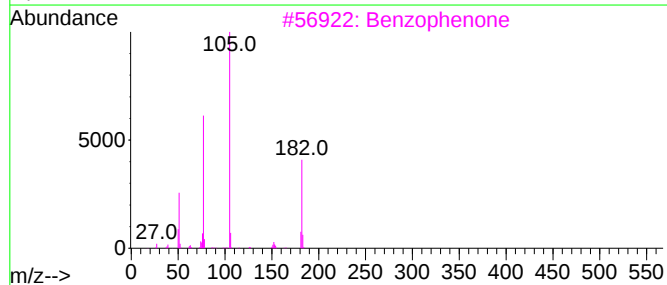
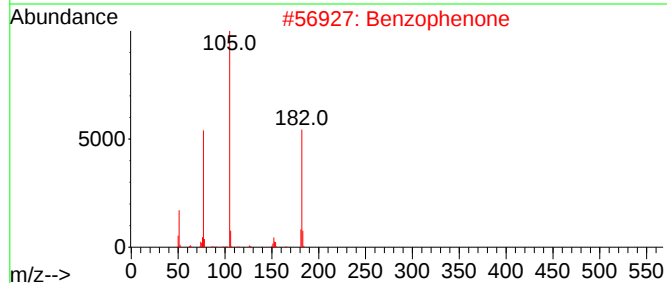
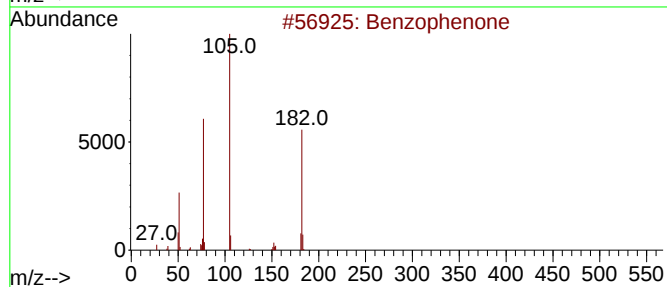
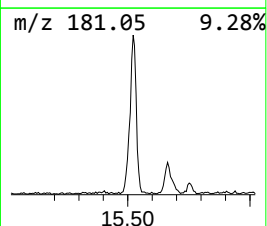
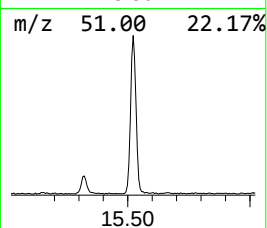
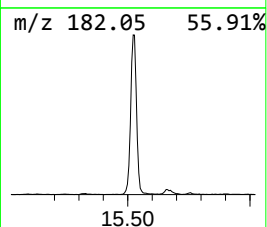
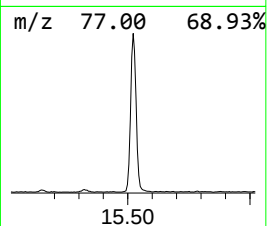
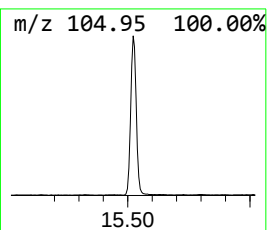
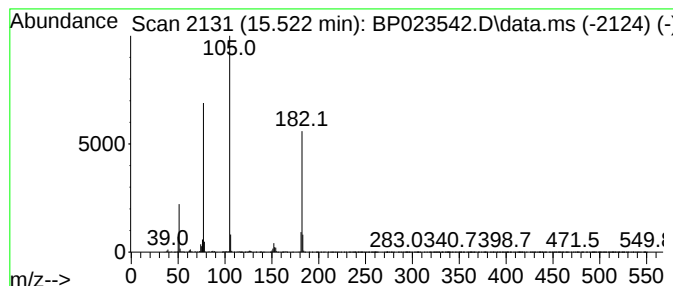
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzophenone Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.522	6.25 ng/ul	1092420	Acenaphthene-d10	14.134

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			Benzophenone	182	C13H10O	000119-61-9	95
3			Benzophenone	182	C13H10O	000119-61-9	94
4			Benzophenone	182	C13H10O	000119-61-9	93
5			Benzophenone	182	C13H10O	000119-61-9	91



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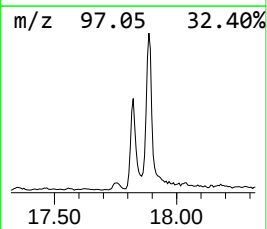
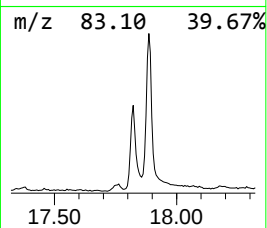
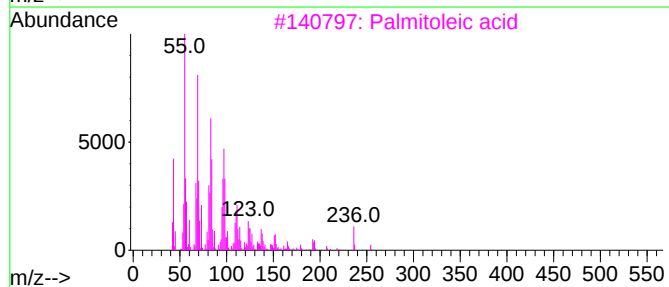
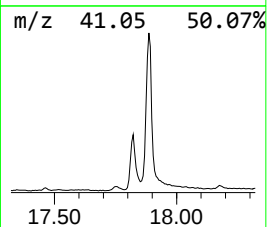
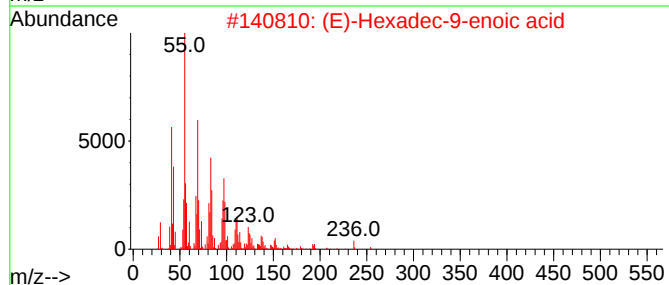
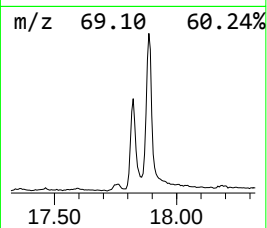
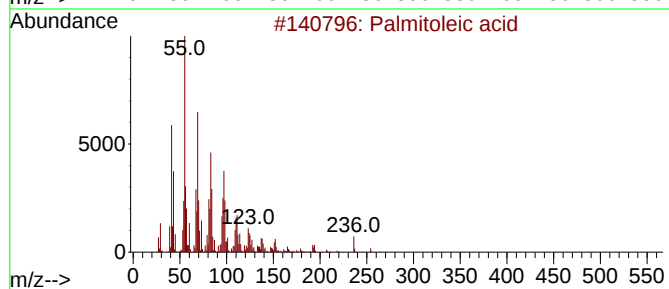
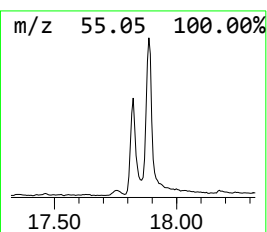
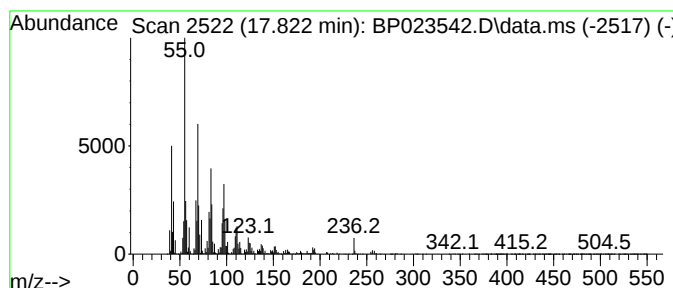
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Palmitoleic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.822	7.94 ng/ul	1858010	Phenanthrene-d10	16.939

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Palmitoleic acid	254	C16H30O2	000373-49-9	99
2			(E)-Hexadec-9-enoic acid	254	C16H30O2	010030-73-6	98
3			Palmitoleic acid	254	C16H30O2	000373-49-9	96
4			cis-7-Hexadecenoic acid	254	C16H30O2	002416-19-5	94
5			cis-10-Nonadecenoic acid	296	C19H36O2	073033-09-7	91



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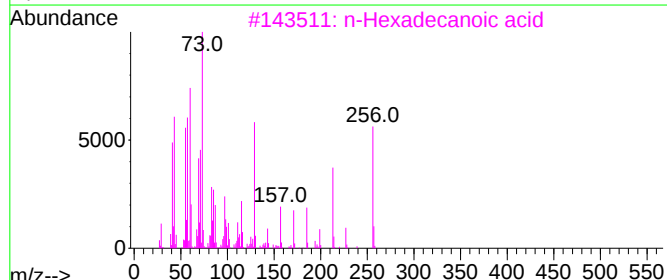
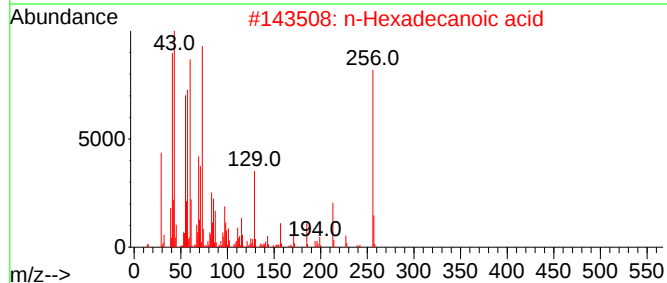
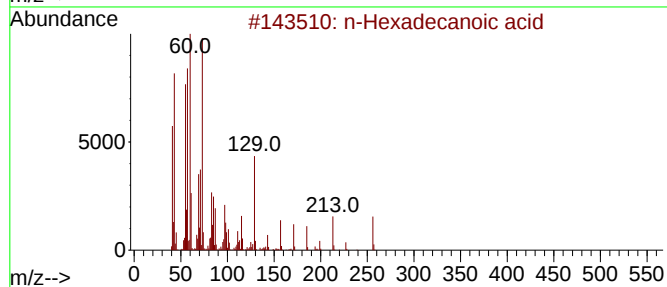
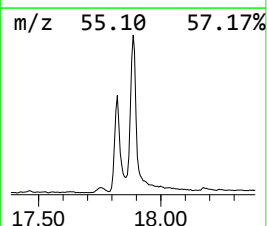
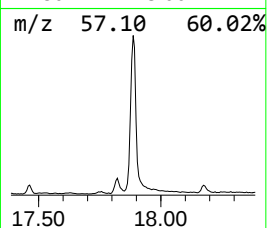
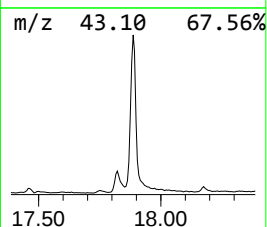
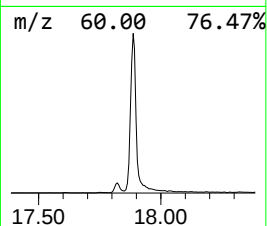
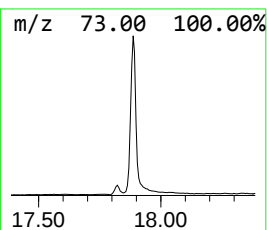
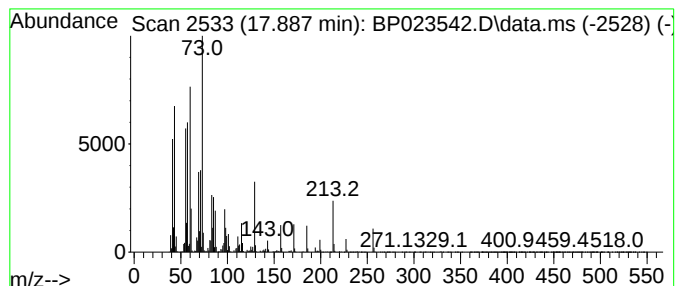
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 n-Hexadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.887	25.56 ng/ul	5980600	Phenanthrene-d10	16.939

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	98
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121824\
Data File : BP023542.D
Acq On : 20 Dec 2024 01:25
Operator : RC/JU
Sample : P5265-06
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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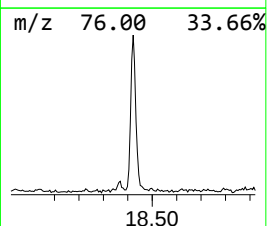
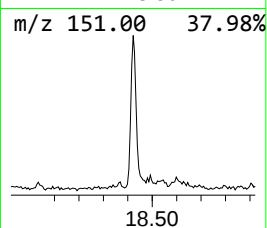
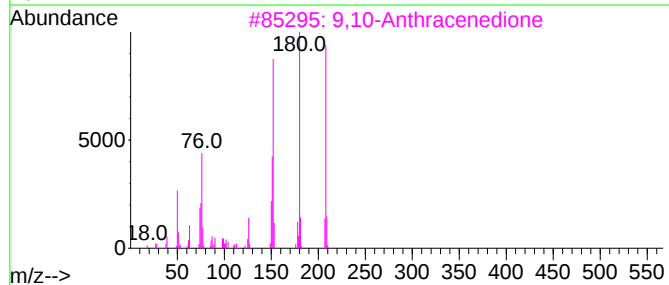
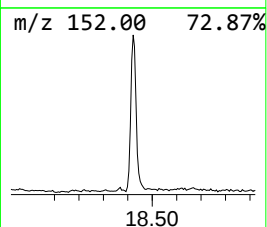
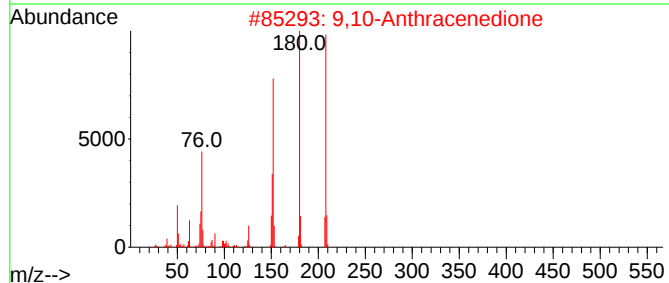
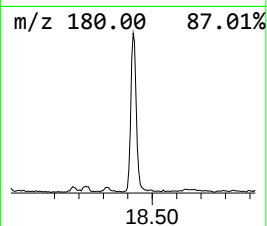
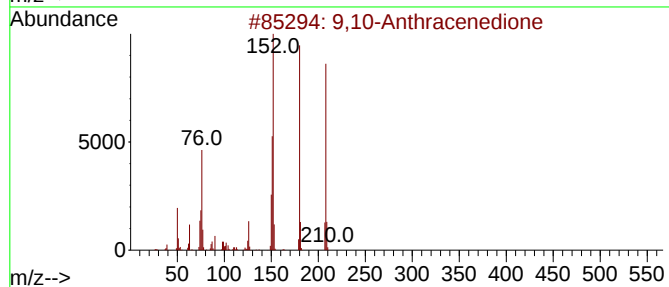
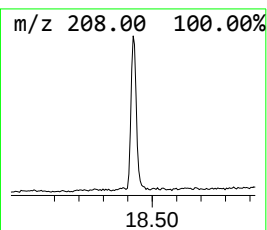
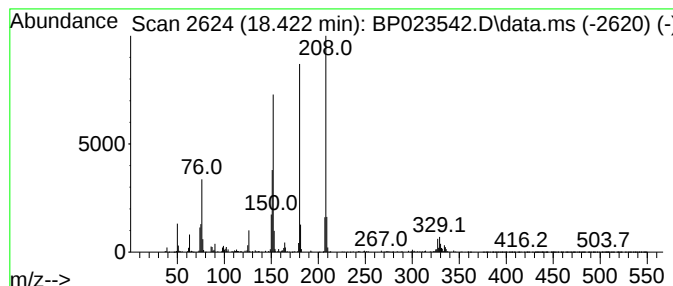
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 9,10-Anthracenedione Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.422	2.39 ng/ul	559013	Phenanthrene-d10	16.939

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9,10-Anthracenedione	208	C14H8O2	000084-65-1	98
2			9,10-Anthracenedione	208	C14H8O2	000084-65-1	97
3			9,10-Anthracenedione	208	C14H8O2	000084-65-1	96
4			9,10-Anthracenedione	208	C14H8O2	000084-65-1	94
5			Benzo[c]cinnoline	180	C12H8N2	000230-17-1	55



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121824\
Data File : BP023542.D
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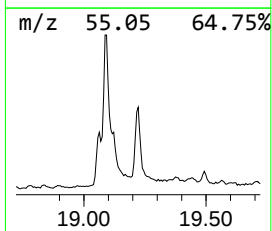
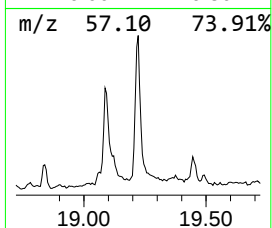
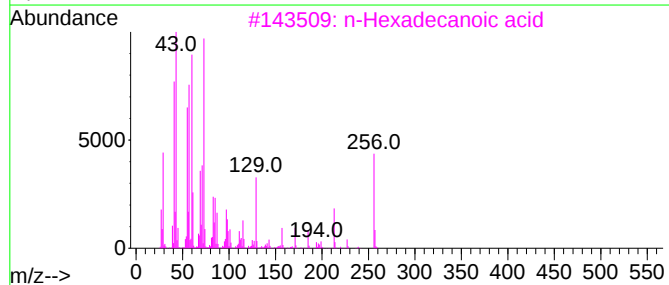
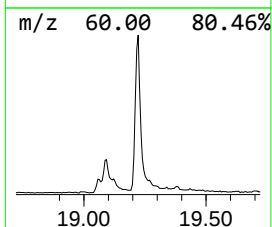
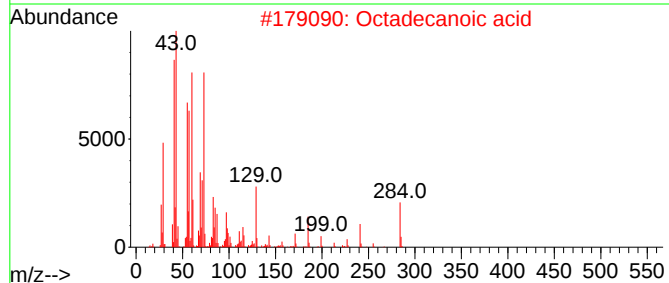
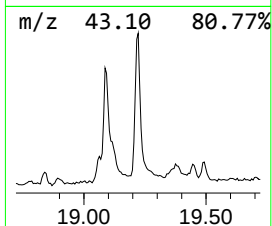
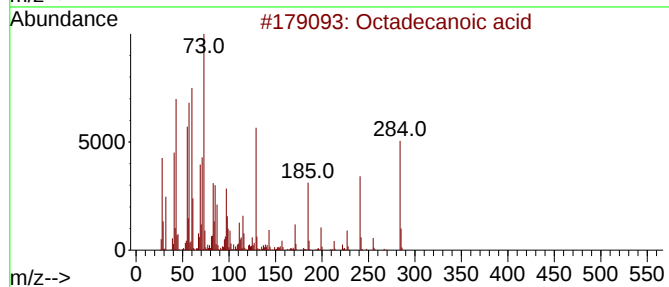
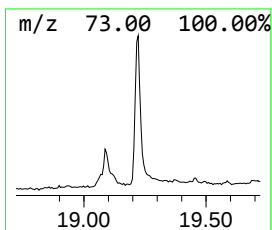
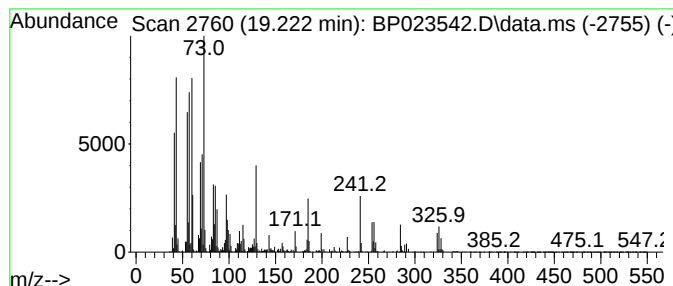
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Octadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.222	4.91 ng/ul	1473170	Chrysene-d12	21.375

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			Octadecanoic acid	284	C18H36O2	000057-11-4	98
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
4			Octadecanoic acid	284	C18H36O2	000057-11-4	95
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121824\
Data File : BP023542.D
Acq On : 20 Dec 2024 01:25
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Sample : P5265-06
Misc :
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Instrument :
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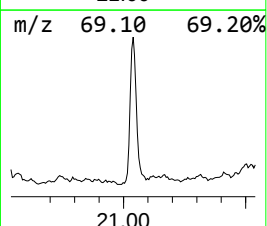
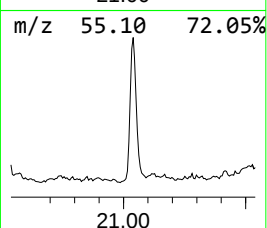
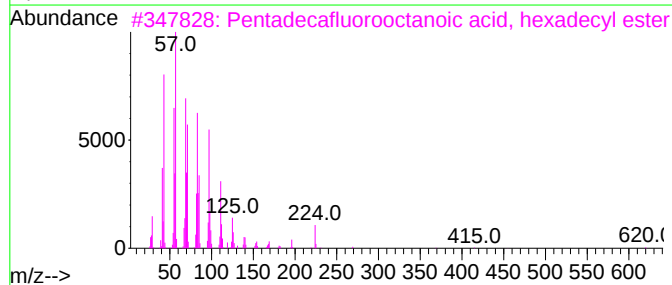
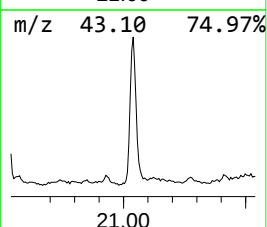
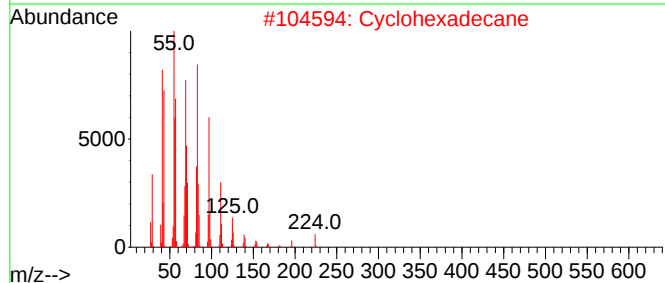
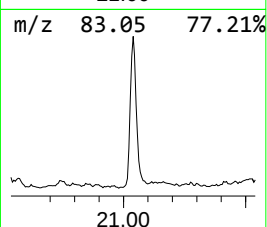
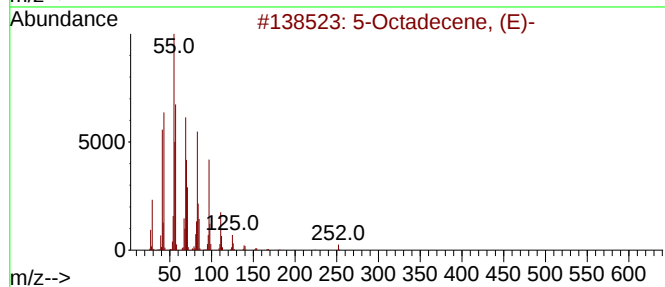
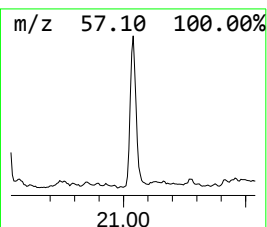
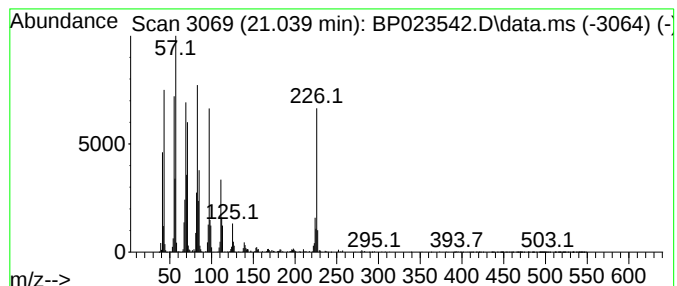
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.039	7.94 ng/ul	2381850	Chrysene-d12	21.375

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			5-Octadecene, (E)-	252	C18H36	007206-21-5	93
2			Cyclohexadecane	224	C16H32	000295-65-8	91
3			Pentadecafluorooctanoic acid, he...	638	C24H33F15O2	1000406-04-6	90
4			1-Octadecene	252	C18H36	000112-88-9	90
5			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121824\
Data File : BP023542.D
Acq On : 20 Dec 2024 01:25
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Misc :
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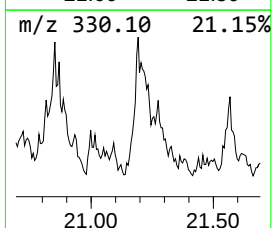
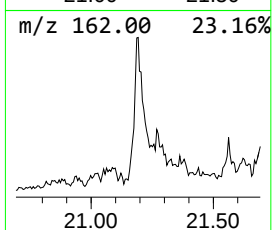
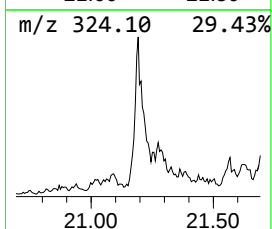
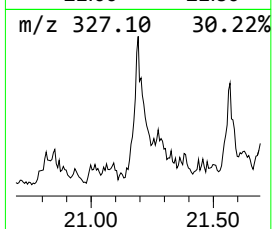
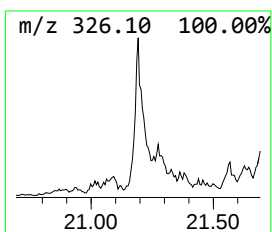
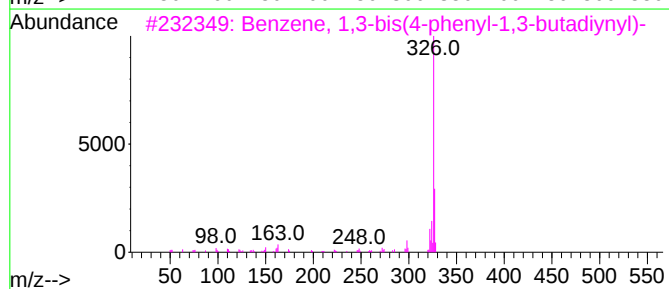
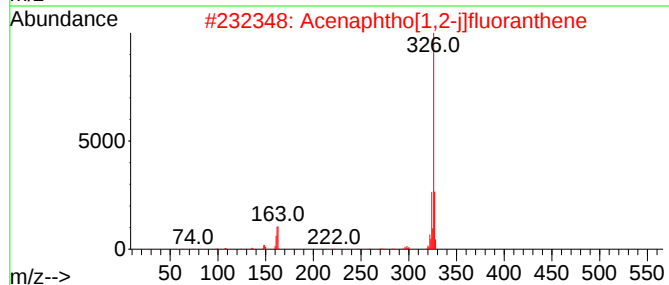
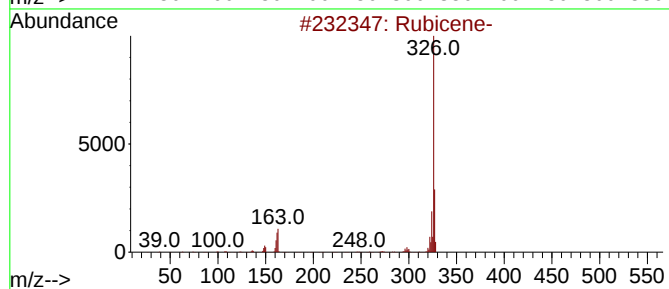
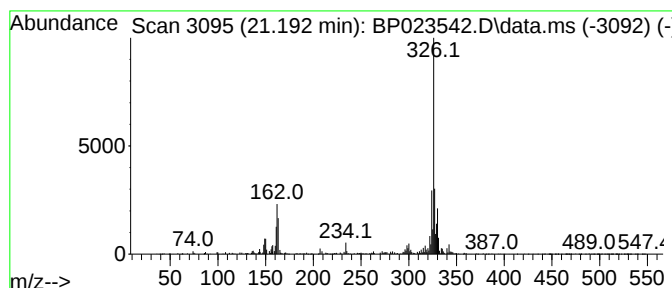
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Rubicene- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.192	2.94 ng/ul	881092	Chrysene-d12	21.375

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Rubicene-	326	C26H14	000197-61-5	92
2			Acenaphtho[1,2-j]fluoranthene	326	C26H14	000193-21-5	83
3			Benzene, 1,3-bis(4-phenyl-1,3-bu...	326	C26H14	037902-13-9	55
4			1,1'-Biphenyl, 2,2',3',4,5-Penta...	324	C12H5C15	041464-51-1	53
5			2,3,3',4,5'-Pentachloro-1,1'-bip...	324	C12H5C15	070362-41-3	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121824\
Data File : BP023542.D
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Instrument :
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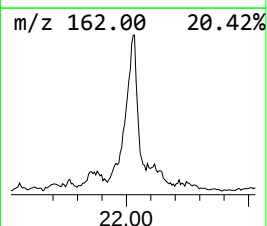
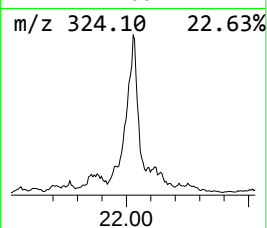
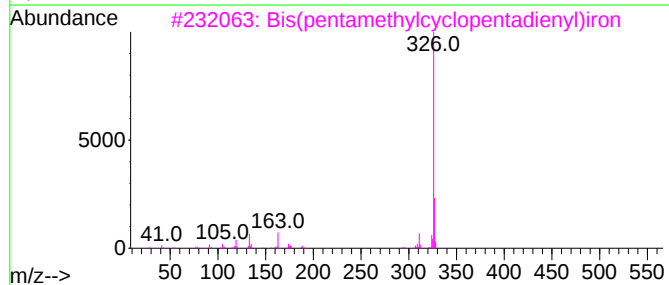
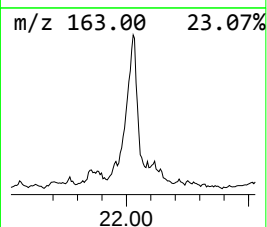
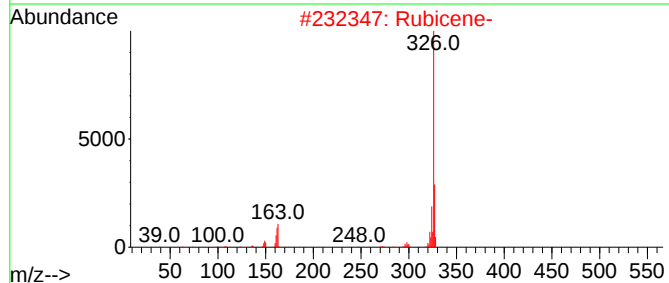
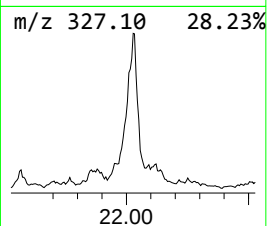
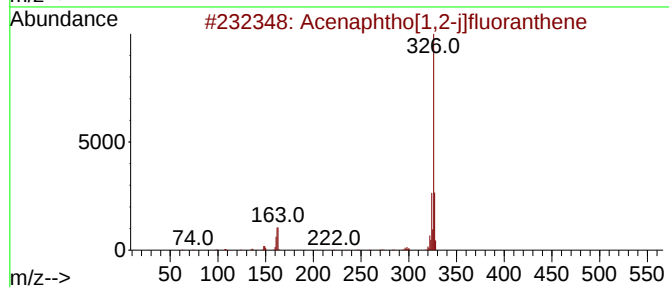
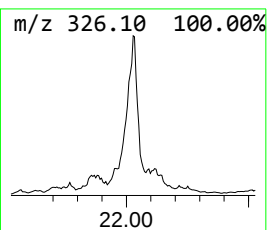
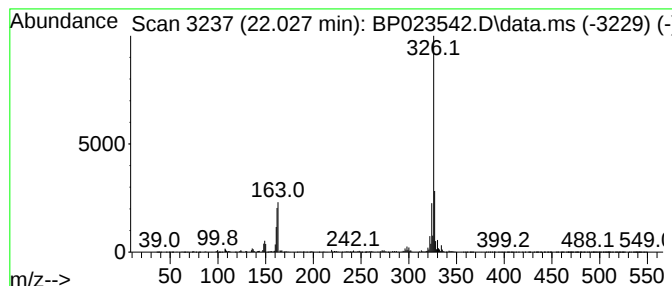
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Acenaphtho[1,2-j]fluoranthene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.028	22.50 ng/ul	6751560	Chrysene-d12	21.375

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acenaphtho[1,2-j]fluoranthene	326	C26H14	000193-21-5	97
2			Rubicene-	326	C26H14	000197-61-5	96
3			Bis(pentamethylcyclopentadienyl)...	326	C20H30Fe	012126-50-0	50
4			quino[2,3-b]acridine-7,14-dione,...	326	C21H14N2O2	1000397-96-1	50
5			Benzene, 1,3-bis(4-phenyl-1,3-bu...	326	C26H14	037902-13-9	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121824\
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Acq On : 20 Dec 2024 01:25
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Misc :
ALS Vial : 19 Sample Multiplier: 1

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ClientSampleId :
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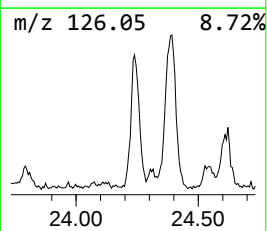
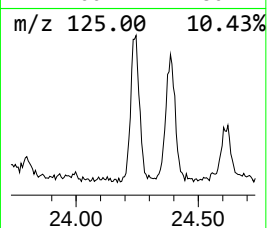
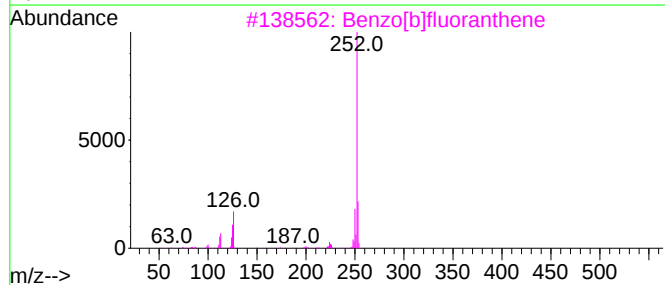
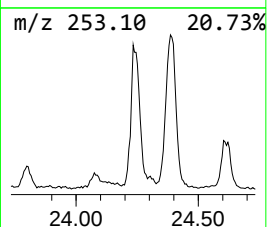
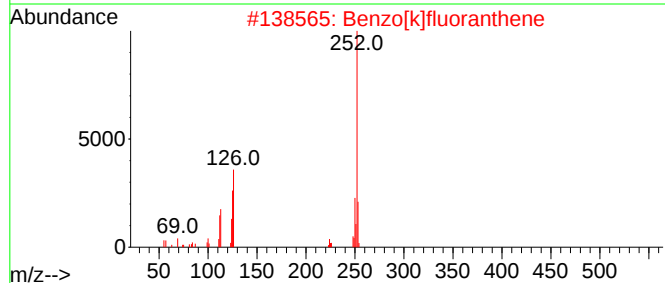
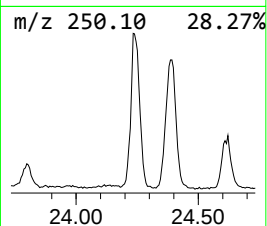
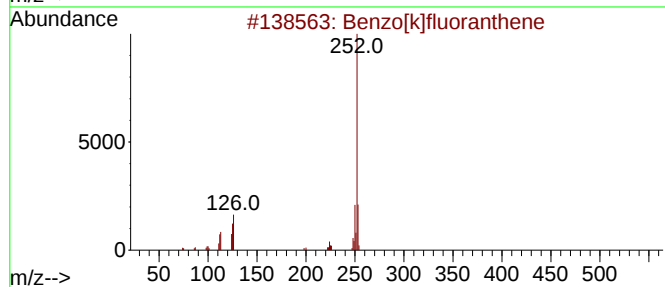
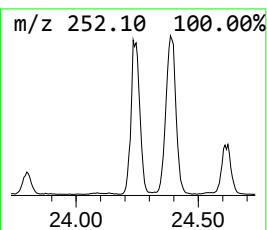
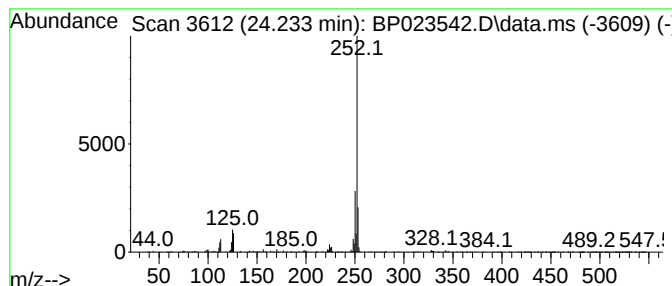
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 Benzo[e]pyrene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.233	3.25 ng/ul	837880	Perylene-d12	24.539

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[k]fluoranthene	252	C20H12	000207-08-9	97
2			Benzo[k]fluoranthene	252	C20H12	000207-08-9	95
3			Benzo[b]fluoranthene	252	C20H12	000205-99-2	94
4			Benzo[e]pyrene	252	C20H12	000192-97-2	94
5			Benzo[b]fluoranthene	252	C20H12	000205-99-2	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121824\
Data File : BP023542.D
Acq On : 20 Dec 2024 01:25
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Sample : P5265-06
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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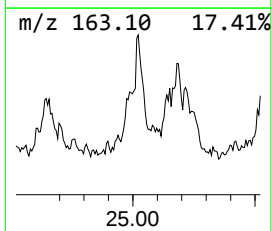
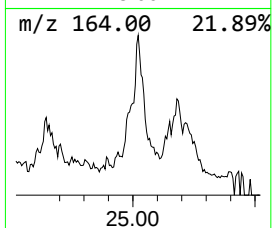
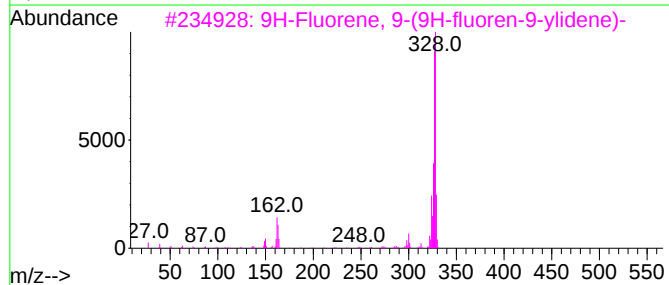
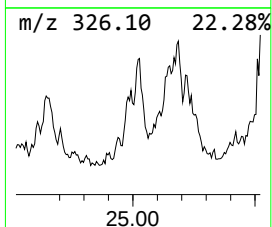
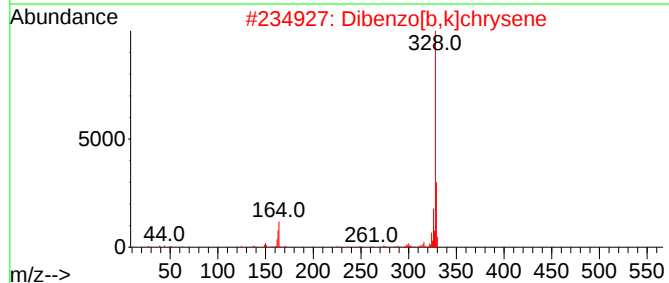
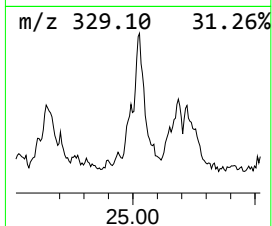
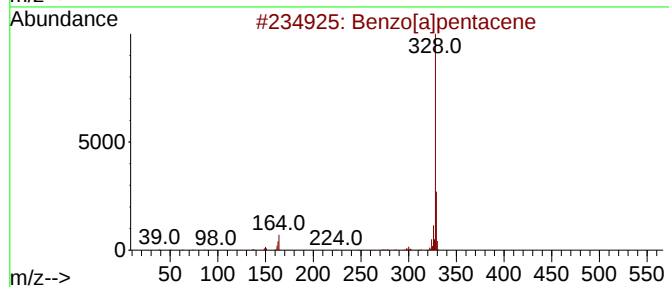
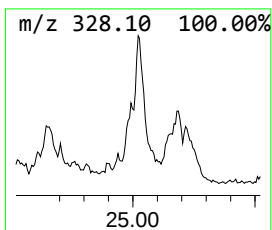
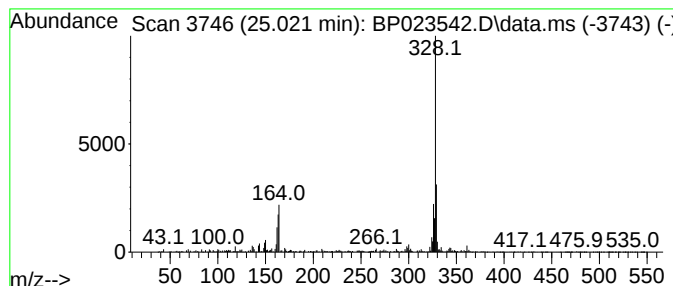
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 Benzo[a]pentacene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.021	4.99 ng/ul	1287970	Perylene-d12	24.539

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[a]pentacene	328	C26H16	000239-98-5	91
2			Dibenzo[b,k]chrysene	328	C26H16	000217-54-9	74
3			9H-Fluorene, 9-(9H-fluoren-9-ylidene)-	328	C26H16	000746-47-4	52
4			1-Amino-4-(4-toluidino)anthra-9,10-dione	328	C21H16N2O2	039774-73-7	50
5			[1,1'-Biphenyl]-4,4'-diamine, 2,2'-bis(4-aminophenyl)-	328	C12H4F8N2	001038-66-0	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121824\
Data File : BP023542.D
Acq On : 20 Dec 2024 01:25
Operator : RC/JU
Sample : P5265-06
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E28Z1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.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	15.522	6.3	ng/ul	1092420	3	14.134	3494140	20.0
Palmitoleic acid	17.822	7.9	ng/ul	1858010	4	16.939	4679150	20.0
n-Hexadecanoic ...	17.887	25.6	ng/ul	5980600	4	16.939	4679150	20.0
9,10-Anthracene...	18.422	2.4	ng/ul	559013	4	16.939	4679150	20.0
Octadecanoic acid	19.222	4.9	ng/ul	1473170	5	21.375	6000320	20.0
(DEL) Alkane: S...	21.039	7.9	ng/ul	2381850	5	21.375	6000320	20.0
Rubicene-	21.192	2.9	ng/ul	881092	5	21.375	6000320	20.0
Acenaphtho[1,2-...	22.028	22.5	ng/ul	6751560	5	21.375	6000320	20.0
Benzo[e]pyrene	24.233	3.3	ng/ul	837880	6	24.539	5163100	20.0
Benzo[a]pentacene	25.021	5.0	ng/ul	1287970	6	24.539	5163100	20.0