

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121523\
 Data File : BM043371.D
 Acq On : 15 Dec 2023 23:13
 Operator : MA/JU
 Sample : 05794-12
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BH3Q7

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-BM121523.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BM043371.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.822	105	108	118	rBV	89670	148382	2.21%	0.144%
2	3.928	121	126	133	rVV	172065	235595	3.50%	0.228%
3	5.069	315	320	331	rVB	73225	120395	1.79%	0.117%
4	5.963	468	472	482	rVB3	32535	69239	1.03%	0.067%
5	7.151	664	674	689	rVB	732464	1346805	20.03%	1.305%
6	7.334	695	705	715	rBV	564775	937286	13.94%	0.908%
7	7.522	730	737	747	rVV	681907	1118209	16.63%	1.084%
8	7.992	807	817	825	rVV	1907616	3032438	45.11%	2.939%
9	8.704	931	938	946	rBV	552157	916446	13.63%	0.888%
10	8.828	952	959	975	rVB	447868	778437	11.58%	0.754%
11	9.163	1009	1016	1022	rBV	675919	1163025	17.30%	1.127%
12	9.886	1129	1139	1149	rBV	537682	932557	13.87%	0.904%
13	10.033	1156	1164	1174	rVV	87659	178553	2.66%	0.173%
14	10.245	1188	1200	1207	rBV2	45808	105311	1.57%	0.102%
15	10.422	1223	1230	1240	rVV	869217	1427716	21.24%	1.384%
16	10.804	1285	1295	1301	rBV	2709726	4577282	68.09%	4.436%
17	10.851	1301	1303	1310	rVV2	64045	92422	1.37%	0.090%
18	10.945	1310	1319	1338	rVB	428026	793474	11.80%	0.769%
19	11.345	1380	1387	1396	rBV5	39331	77042	1.15%	0.075%
20	12.457	1570	1576	1586	rVB	36750	76832	1.14%	0.074%
21	12.680	1607	1614	1620	rVB	37689	67954	1.01%	0.066%
22	13.463	1735	1747	1752	rVV2	70701	146676	2.18%	0.142%
23	13.522	1752	1757	1765	rVV	64992	124804	1.86%	0.121%
24	13.686	1778	1785	1790	rVB	143046	222485	3.31%	0.216%
25	13.945	1818	1829	1834	rVV2	61670	122486	1.82%	0.119%
26	14.039	1840	1845	1854	rVB	1393662	1954131	29.07%	1.894%
27	14.316	1880	1892	1904	rBV	1644219	2624847	39.04%	2.544%
28	14.621	1936	1944	1951	rBV	3729940	5535466	82.34%	5.365%
29	14.686	1951	1955	1963	rVV	180272	281786	4.19%	0.273%
30	14.821	1972	1978	1988	rBV	358809	593751	8.83%	0.575%
31	15.021	2008	2012	2020	rVV	137681	245118	3.65%	0.238%
32	15.410	2071	2078	2081	rBV7	31471	67194	1.00%	0.065%
33	15.615	2106	2113	2118	rVV	2065348	2960912	44.04%	2.870%
34	15.668	2118	2122	2127	rVV2	186013	284864	4.24%	0.276%
35	15.733	2128	2133	2140	rVV	325622	499253	7.43%	0.484%
36	15.833	2146	2150	2154	rVV6	35486	71062	1.06%	0.069%

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Smoothing : OFF

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Min Area: 0.1 % of largest Peak

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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-BM121523.MA.M
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37	15.963	2167	2172	2180	rVB	110553	177913	2.65%	0.172%
38	16.110	2192	2197	2205	rVB	69991	137158	2.04%	0.133%
39	16.333	2229	2235	2242	rVV5	67729	156106	2.32%	0.151%
40	16.557	2269	2273	2281	rVB2	62978	115461	1.72%	0.112%
41	16.662	2282	2291	2297	rBV4	155817	311790	4.64%	0.302%
42	16.898	2323	2331	2334	rBV3	47975	109442	1.63%	0.106%
43	16.939	2335	2338	2340	rVV3	54969	80514	1.20%	0.078%
44	16.968	2340	2343	2347	rVV	97332	140653	2.09%	0.136%
45	17.021	2347	2352	2358	rVV	143396	222923	3.32%	0.216%
46	17.092	2358	2364	2366	rVV2	39476	73159	1.09%	0.071%
47	17.127	2367	2370	2375	rVV5	68023	120027	1.79%	0.116%
48	17.186	2375	2380	2387	rVB	140308	223246	3.32%	0.216%
49	17.368	2404	2411	2414	rBV	3727472	5397749	80.29%	5.232%
50	17.409	2414	2418	2423	rVV	2701283	3717121	55.29%	3.603%
51	17.462	2423	2427	2431	rVV2	1918414	2791838	41.53%	2.706%
52	17.498	2431	2433	2438	rVV	440301	525759	7.82%	0.510%
53	17.557	2439	2443	2447	rVB3	50033	85279	1.27%	0.083%
54	17.774	2472	2480	2489	rBV	210752	388361	5.78%	0.376%
55	17.886	2493	2499	2505	rVV3	64435	152862	2.27%	0.148%
56	17.945	2506	2509	2517	rVB2	66748	118714	1.77%	0.115%
57	18.245	2555	2560	2561	rBV	293759	384574	5.72%	0.373%
58	18.292	2565	2568	2573	rVB	421148	556553	8.28%	0.539%
59	18.368	2579	2581	2586	rVB	93525	115537	1.72%	0.112%
60	18.433	2587	2592	2596	rBV2	382272	685418	10.20%	0.664%
61	18.468	2596	2598	2606	rVB	202602	268643	4.00%	0.260%
62	18.556	2609	2613	2616	rBV4	58931	95039	1.41%	0.092%
63	18.745	2640	2645	2648	rBV	209154	278959	4.15%	0.270%
64	18.786	2648	2652	2656	rVV3	162592	230965	3.44%	0.224%
65	18.839	2657	2661	2664	rVB2	125352	166841	2.48%	0.162%
66	18.986	2683	2686	2689	rVB2	71572	79423	1.18%	0.077%
67	19.068	2697	2700	2704	rVB2	88942	134682	2.00%	0.131%
68	19.156	2710	2715	2719	rBV2	193706	332603	4.95%	0.322%
69	19.198	2719	2722	2724	rBV2	65321	93072	1.38%	0.090%
70	19.292	2734	2738	2746	rBV2	211934	460600	6.85%	0.446%
71	19.415	2754	2759	2765	rVB	3763000	5031119	74.84%	4.876%
72	19.751	2810	2816	2818	rBV2	2356327	3397333	50.53%	3.293%
73	19.780	2818	2821	2829	rVB	3240631	4155978	61.82%	4.028%
74	19.845	2829	2832	2836	rBV2	140860	233807	3.48%	0.227%
75	19.892	2836	2840	2844	rVB	1389600	1606428	23.89%	1.557%
76	19.933	2845	2847	2849	rBV3	86290	103004	1.53%	0.100%

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Integrator: RTE
 Smoothing : OFF Filtering: 5
 Sampling : 1 Min Area: 0.1 % of largest Peak
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 Stop Thrs : 0 Peak Location: TOP

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77	19.992	2854	2857	2867	rVB2	299096	533377	7.93%	0.517%
78	20.274	2901	2905	2908	rBV2	194581	246159	3.66%	0.239%
79	20.374	2919	2922	2923	rBV	123709	138993	2.07%	0.135%
80	20.398	2923	2926	2929	rVV	449677	602922	8.97%	0.584%
81	20.450	2930	2935	2942	rVB2	1275331	2033392	30.25%	1.971%
82	20.586	2953	2958	2961	rBV2	762330	958451	14.26%	0.929%
83	20.639	2965	2967	2974	rVB	552657	663886	9.88%	0.643%
84	20.839	2996	3001	3004	rBV	4536585	5184446	77.12%	5.025%
85	20.868	3004	3006	3009	rVV	442354	501295	7.46%	0.486%
86	20.909	3010	3013	3020	rVB3	564153	749520	11.15%	0.726%
87	21.015	3027	3031	3037	rBV2	419001	634768	9.44%	0.615%
88	21.227	3063	3067	3069	rBV	300853	430638	6.41%	0.417%
89	21.262	3070	3073	3078	rVB2	483646	691400	10.28%	0.670%
90	21.539	3113	3120	3124	rBV2	3664646	6722877	100.00%	6.516%
91	21.580	3124	3127	3137	rVB	1377952	1900825	28.27%	1.842%
92	21.821	3164	3168	3172	rBV	1335021	1589967	23.65%	1.541%
93	22.062	3206	3209	3212	rBV	301327	451590	6.72%	0.438%
94	22.250	3237	3241	3250	rVB5	490101	795488	11.83%	0.771%
95	22.415	3266	3269	3272	rBV3	292767	448633	6.67%	0.435%
96	23.227	3401	3407	3412	rBV	1108016	2843710	42.30%	2.756%
97	23.744	3491	3495	3500	rBV	607614	1002874	14.92%	0.972%
98	23.803	3500	3505	3509	rVV2	883189	1637349	24.35%	1.587%
99	23.850	3509	3513	3519	rVB	811140	1444476	21.49%	1.400%
100	23.962	3526	3532	3538	rBV	1844133	3580563	53.26%	3.470%

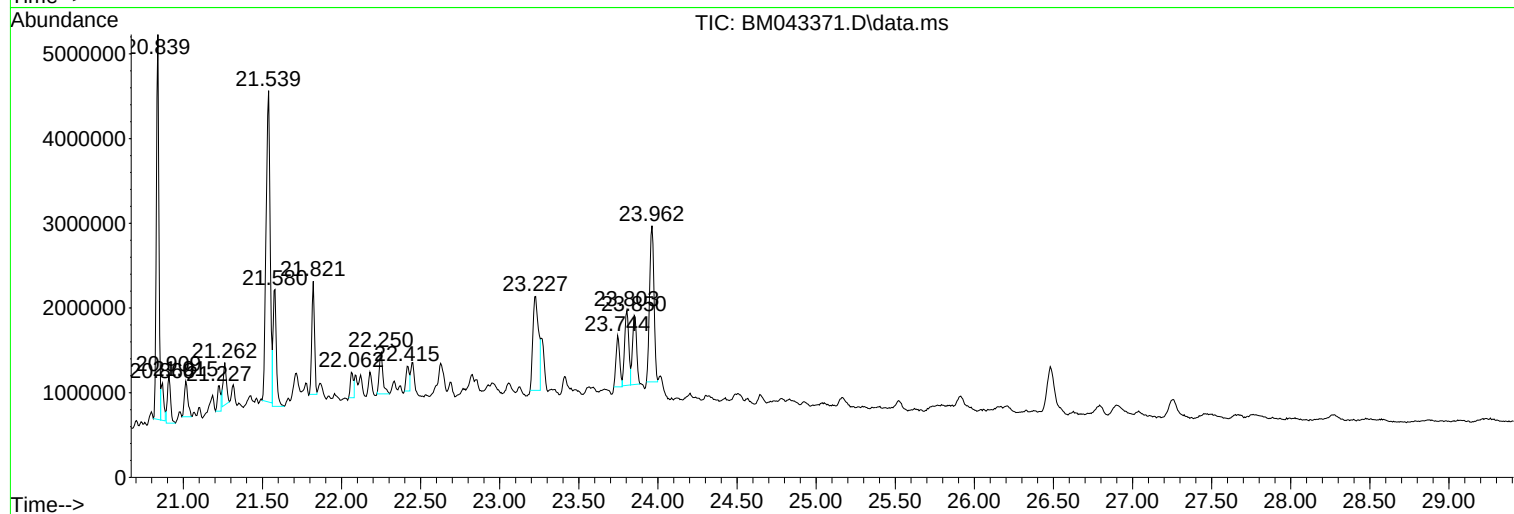
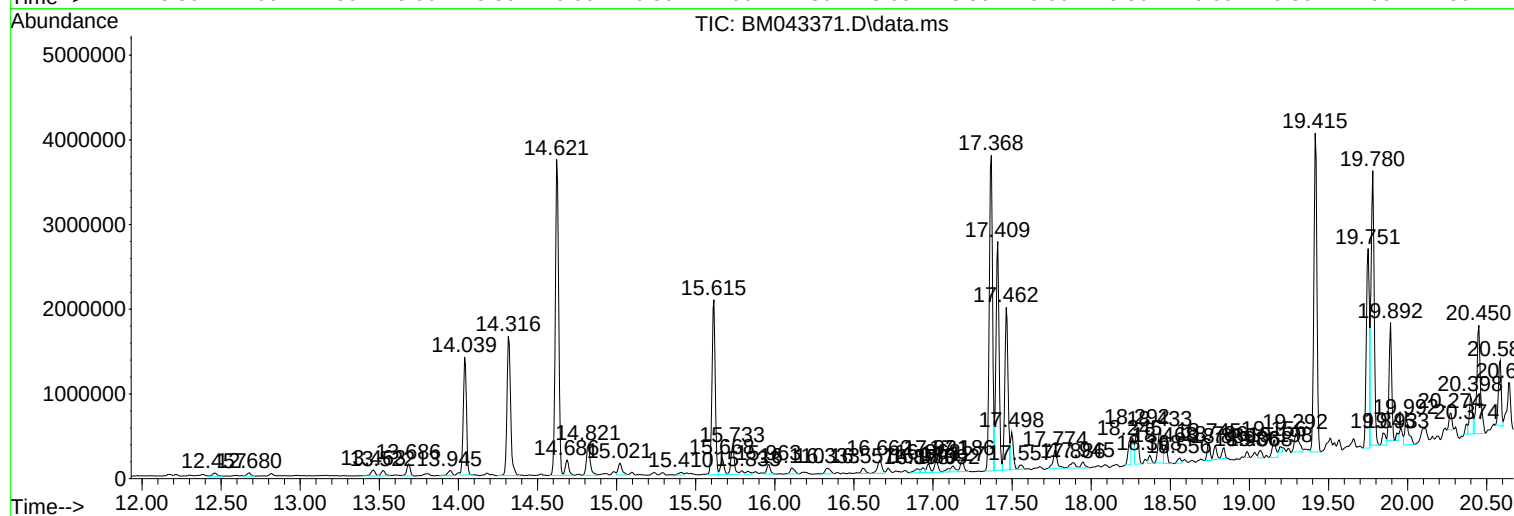
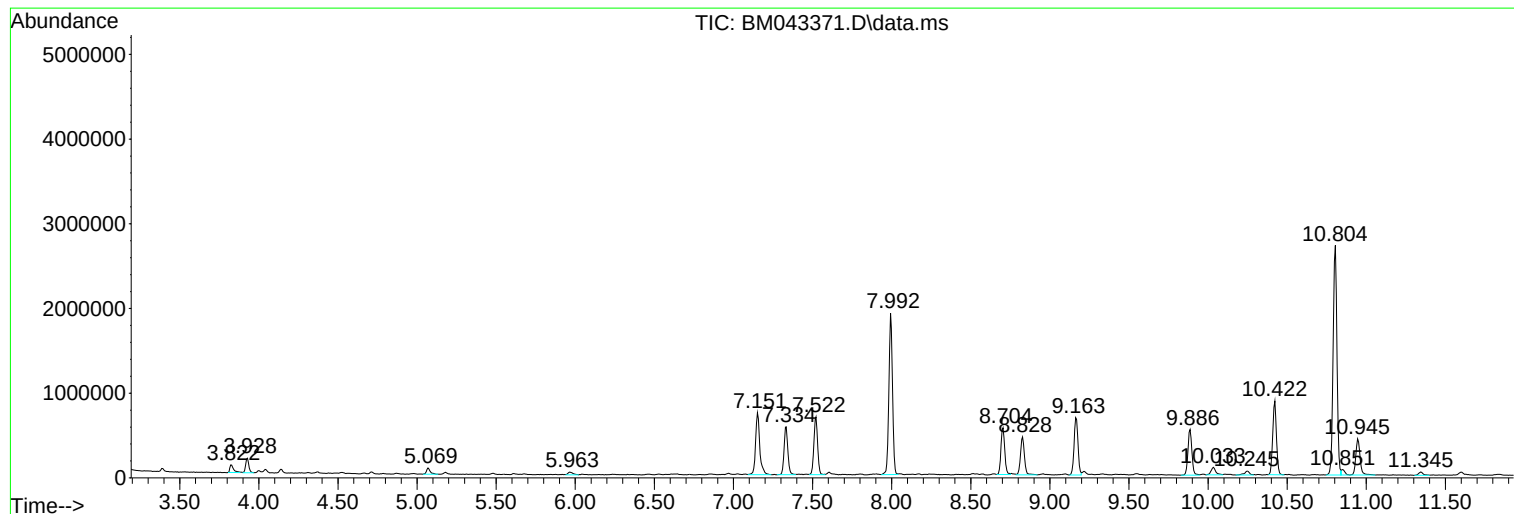
Sum of corrected areas: 103174487

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.MA.M
Quant Title : SVOA CALIBRATION

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



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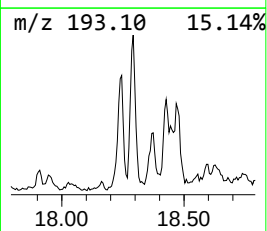
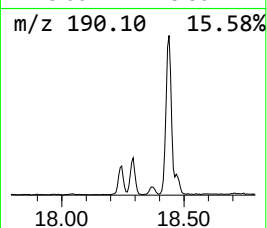
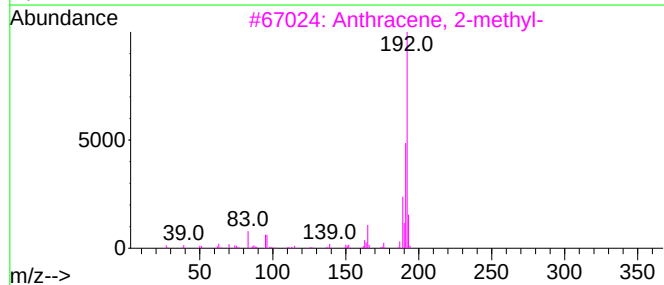
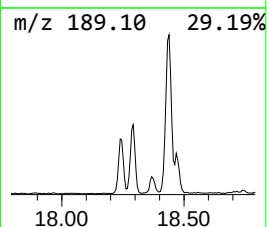
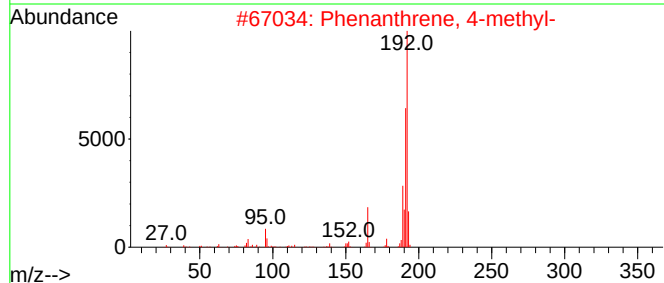
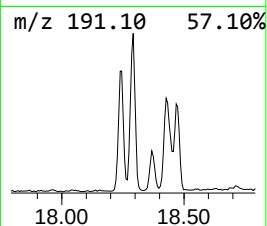
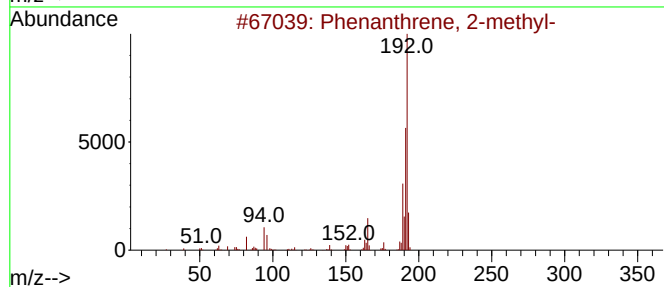
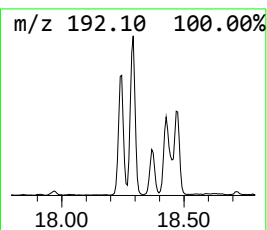
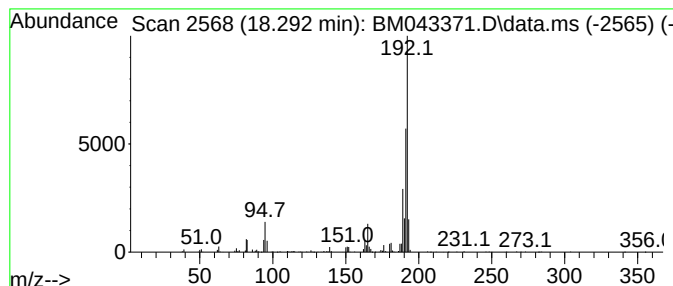
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Phenanthrene, 2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.292	2.06 ng/ul	556553	Phenanthrene-d10	17.368

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
2		Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	95
3		Anthracene, 2-methyl-	192	C15H12	000613-12-7	95
4		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	94
5		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	94



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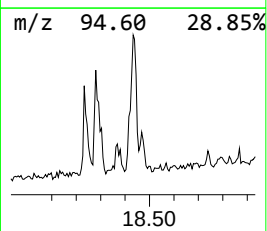
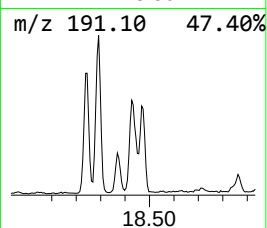
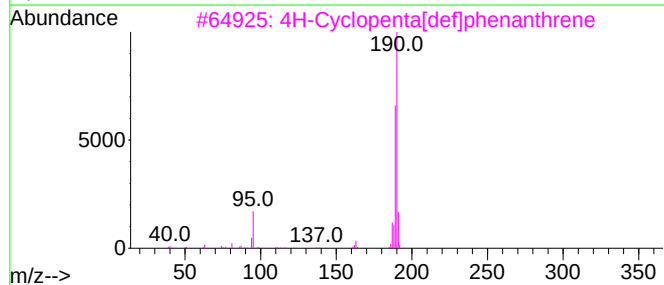
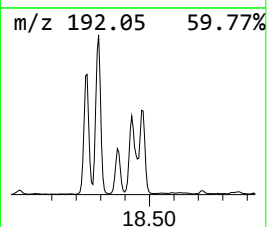
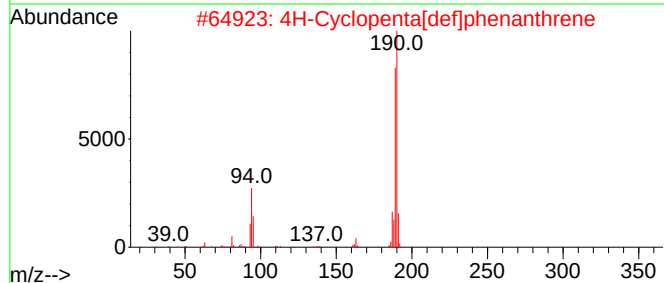
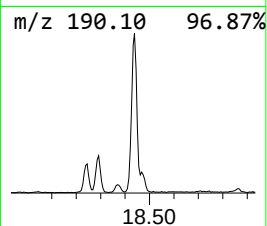
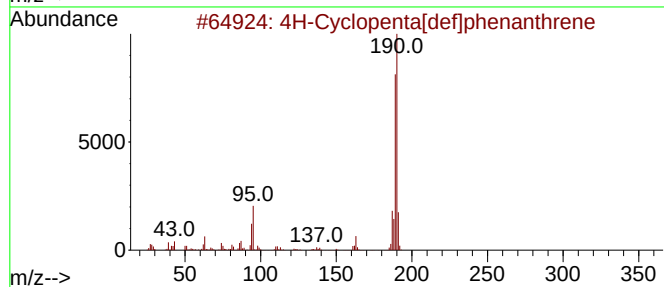
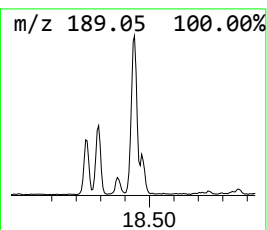
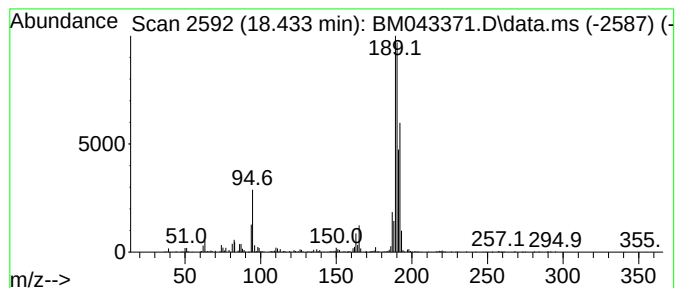
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 4H-Cyclopenta[def]phenanthrene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.433	2.54 ng/ul	685418	Phenanthrene-d10	17.368

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	60
2			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	58
3			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	35
4			1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	30
5			Benzene, 1,1'-(1,2-propadienyld...	192	C15H12	014251-57-1	25



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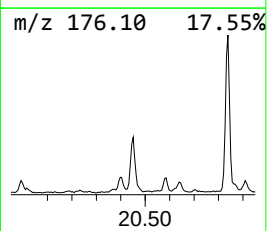
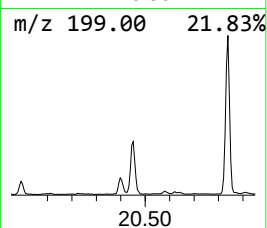
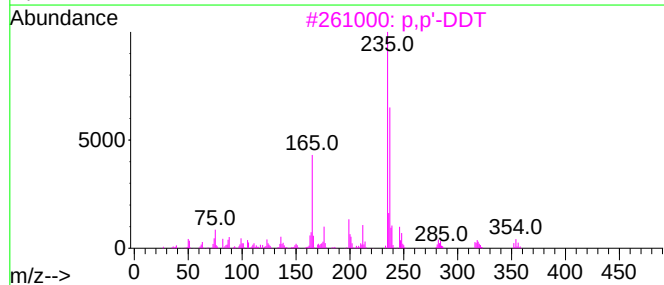
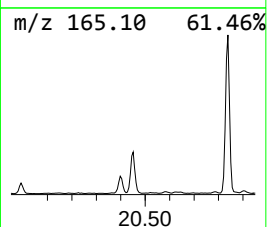
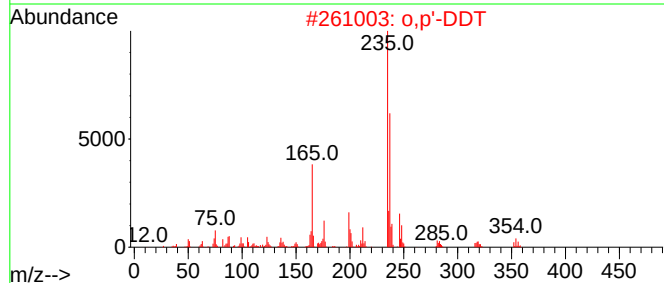
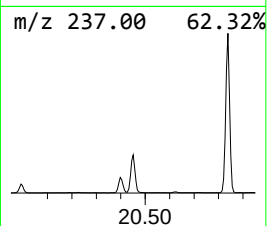
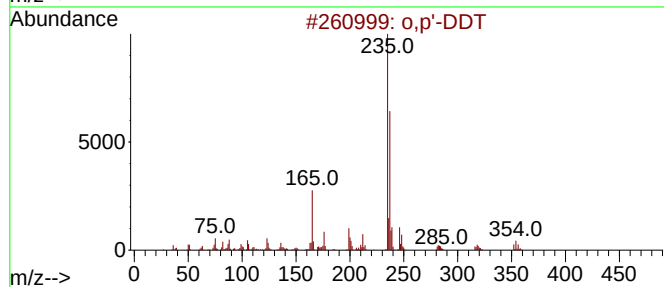
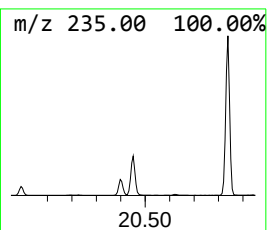
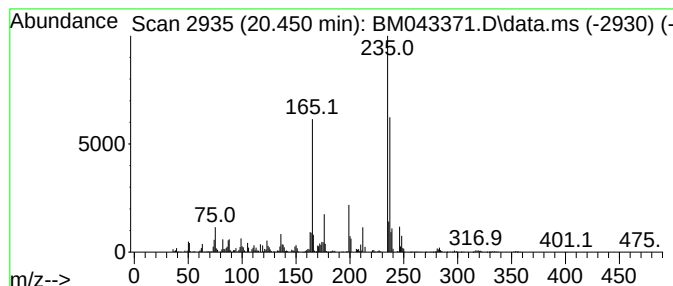
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 o,p'-DDT Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.451	6.05 ng/ul	2033390	Chrysene-d12	21.539

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			o,p'-DDT	352	C14H9Cl5	000789-02-6	97
2			o,p'-DDT	352	C14H9Cl5	000789-02-6	91
3			p,p'-DDT	352	C14H9Cl5	000050-29-3	90
4			3-Butanone, 1,1-bis(4-chlorophen...	320	C18H18Cl2O	1000158-20-4	89
5			p,p'-DDT	352	C14H9Cl5	000050-29-3	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121523\
Data File : BM043371.D
Acq On : 15 Dec 2023 23:13
Operator : MA/JU
Sample : 05794-12
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BH3Q7

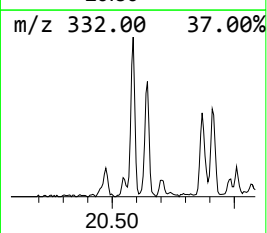
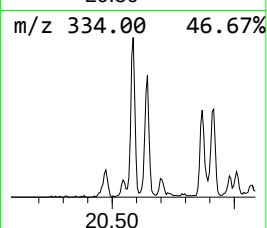
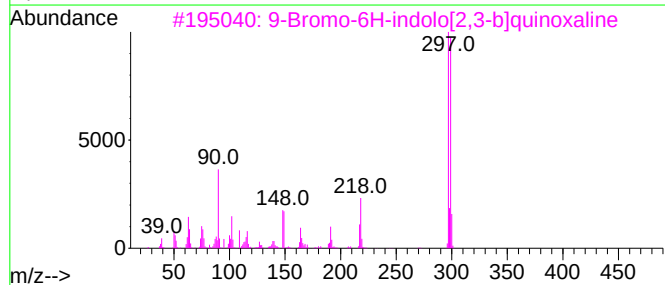
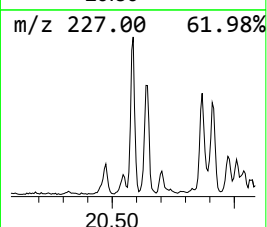
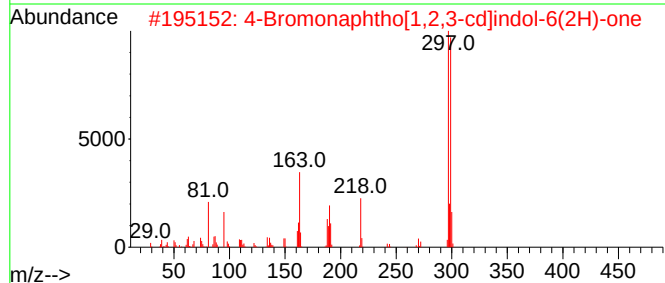
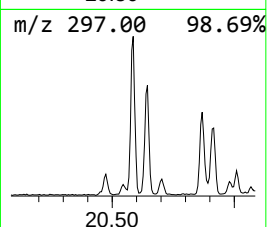
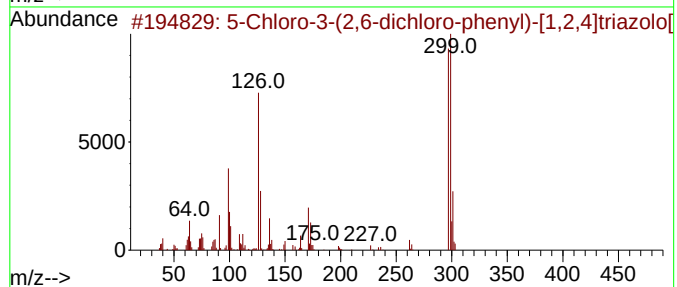
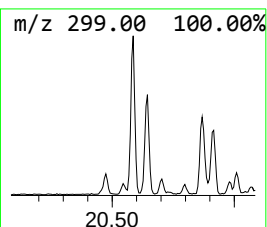
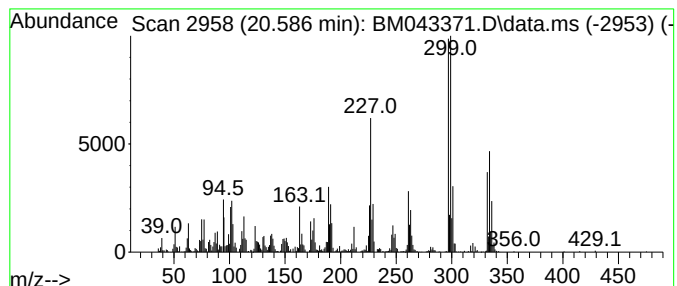
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 unknown-01 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.586	2.85 ng/ul	958451	Chrysene-d12	21.539

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			5-Chloro-3-(2,6-dichloro-phenyl)...	297	C12H6Cl3N3	1000274-53-0	46
2			4-Bromonaphtho[1,2,3-cd]indol-6(...	297	C15H8BrNO	1000459-95-4	45
3			9-Bromo-6H-indolo[2,3-b]quinoxaline	297	C14H8BrN3	057743-36-9	38
4			1,3,5-Triazine, 2-chloro-4-(4-ch...	332	C15H10Cl2N4O	333772-26-2	22
5			6-Chlorobenzo(a)phenothiazin-5-one	297	C16H8ClNOS	025947-05-1	18



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121523\
Data File : BM043371.D
Acq On : 15 Dec 2023 23:13
Operator : MA/JU
Sample : 05794-12
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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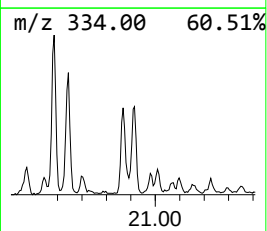
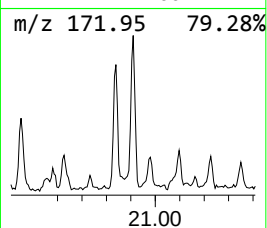
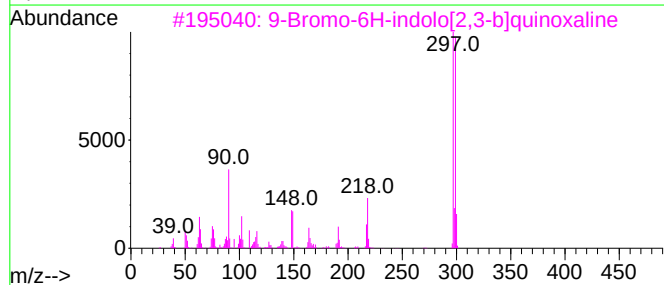
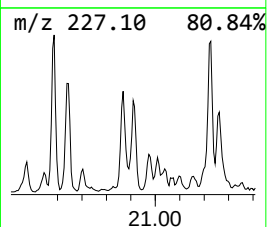
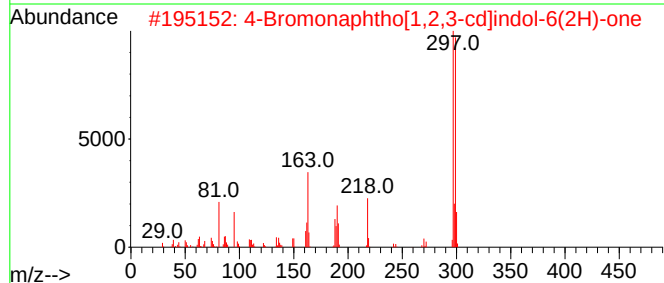
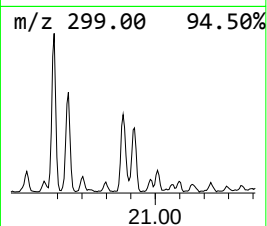
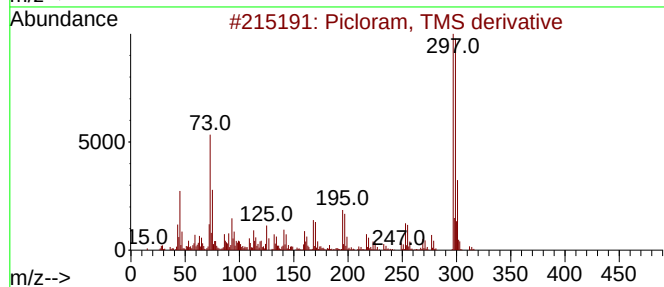
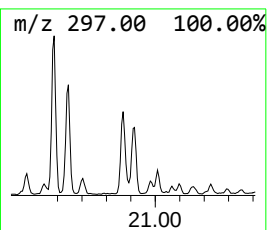
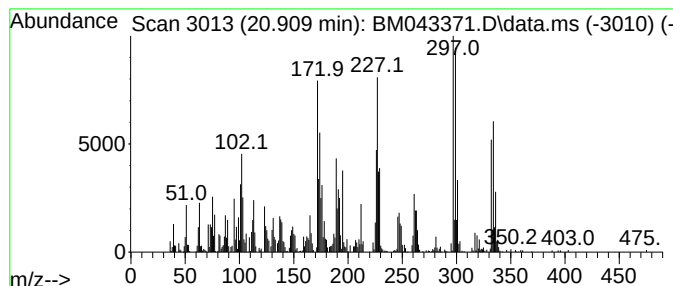
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Picloram, TMS derivative Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.909	2.23 ng/ul	749520	Chrysene-d12	21.539

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Picloram, TMS derivative	312	C9H11Cl3N2O2Si	1000492-53-1	51
2			4-Bromonaphtho[1,2,3-cd]indol-6(...	297	C15H8BrNO	1000459-95-4	20
3			9-Bromo-6H-indolo[2,3-b]quinoxaline	297	C14H8BrN3	057743-36-9	14
4			2,6-Lutidine, 3,5-dichloro-4-[p-...	297	C14H13Cl2NS	1000252-28-1	12
5			1-[5-Amino-6-uracilyl]-2-[2,6-di...	297	C12H9Cl2N3O2	1000214-68-4	10



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121523\
Data File : BM043371.D
Acq On : 15 Dec 2023 23:13
Operator : MA/JU
Sample : 05794-12
Misc :
ALS Vial : 20 Sample Multiplier: 1

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ClientSampleId :
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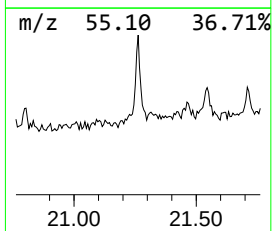
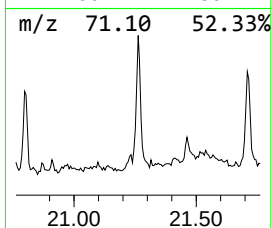
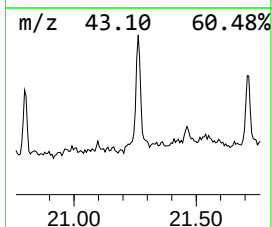
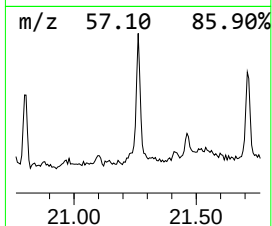
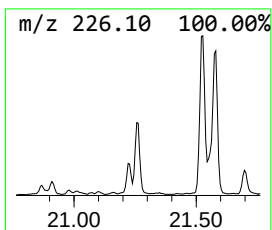
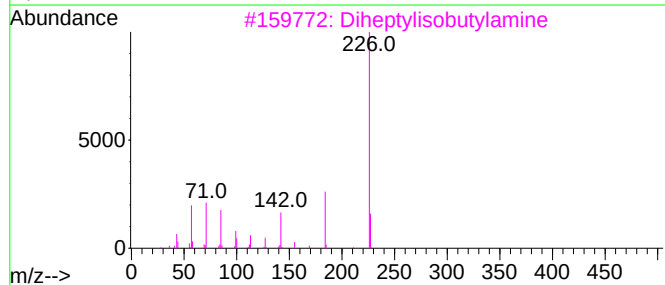
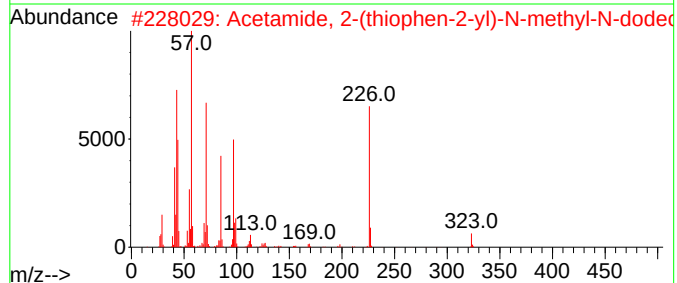
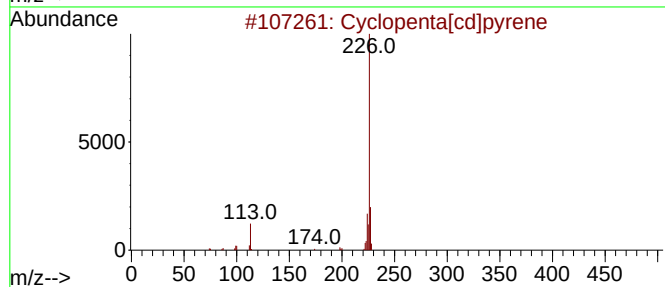
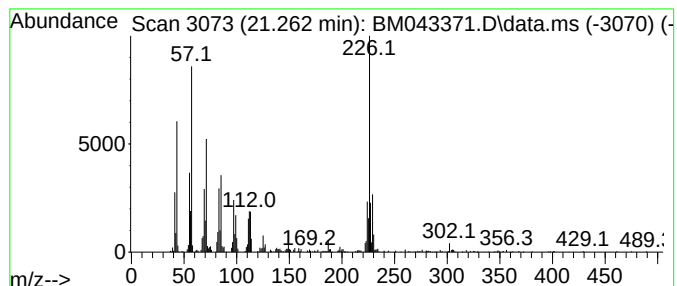
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Cyclopenta[cd]pyrene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.262	2.06 ng/ul	691400	Chrysene-d12	21.539

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopenta[cd]pyrene	226	C18H10	027208-37-3	55
2			Acetamide, 2-(thiophen-2-yl)-N-m...	323	C19H33NOS	1000421-20-9	38
3			Diheptylisobutylamine	269	C18H39N	1000310-32-0	37
4			benzenamine, 4-(2-benzothiazolyl)-	226	C13H10N2S	1000400-93-4	35
5			.beta.-Carboline, 7-methoxy-1,2-...	226	C14H14N2O	006519-18-2	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121523\
Data File : BM043371.D
Acq On : 15 Dec 2023 23:13
Operator : MA/JU
Sample : 05794-12
Misc :
ALS Vial : 20 Sample Multiplier: 1

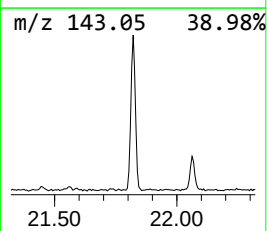
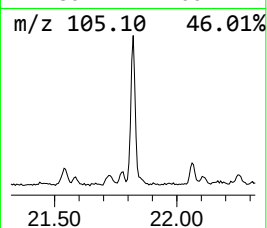
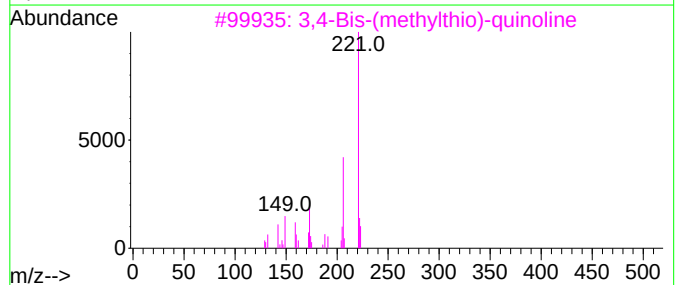
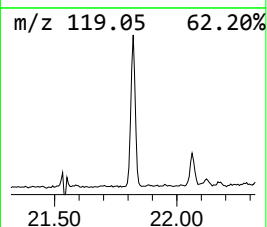
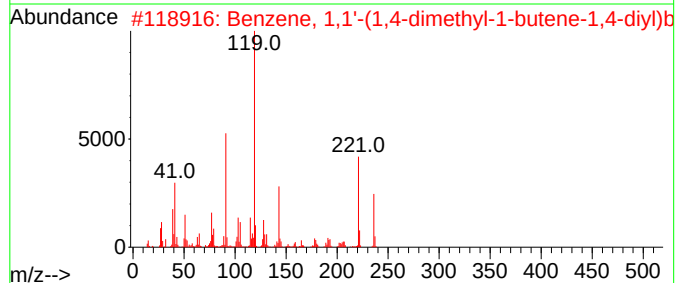
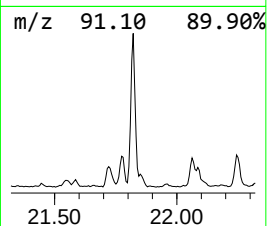
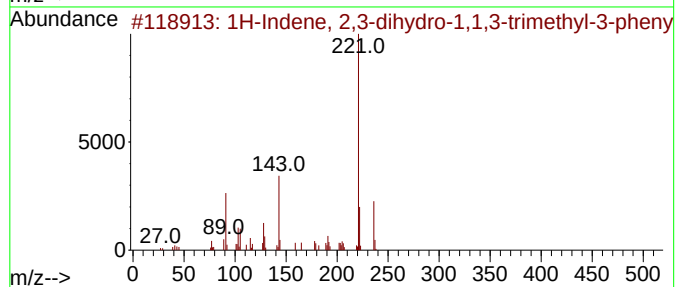
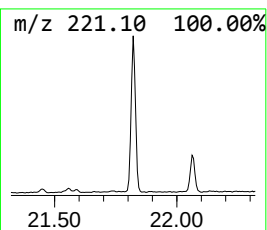
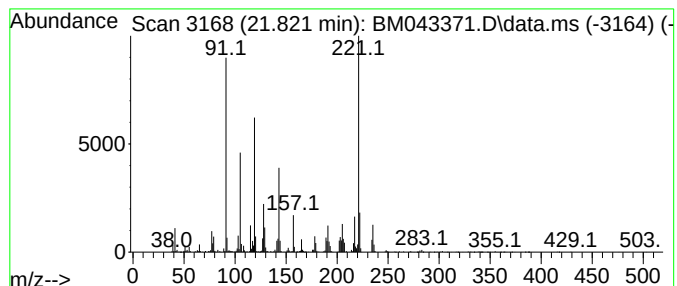
Instrument :
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 1H-Indene, 2,3-dihydro-1,1,... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD		R.T.	
21.821	4.73 ng/ul	1589970	Chrysene-d12		21.539	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-1,1,3-tri...		236	C18H20	003910-35-8	50
2	Benzene, 1,1'-(1,4-dimethyl-1-bu...		236	C18H20	052161-54-3	45
3	3,4-Bis-(methylthio)-quinoline		221	C11H11NS2	074579-34-3	43
4	1H-Indene, 2,3-dihydro-1,1,3-tri...		236	C18H20	003910-35-8	38
5	1H-Indole, 1-ethyl-2-phenyl-		221	C16H15N	013228-39-2	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121523\
Data File : BM043371.D
Acq On : 15 Dec 2023 23:13
Operator : MA/JU
Sample : 05794-12
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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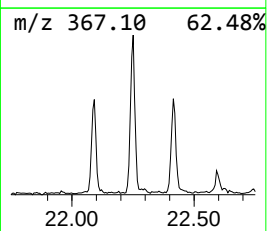
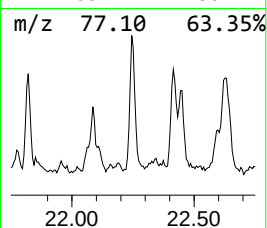
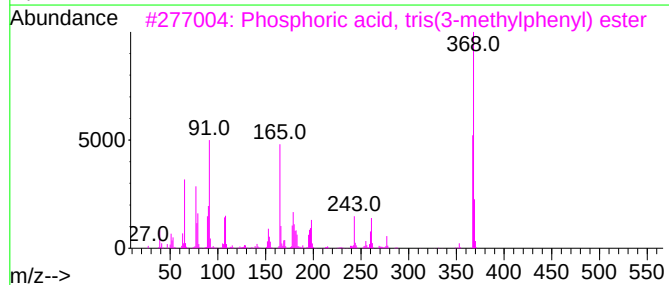
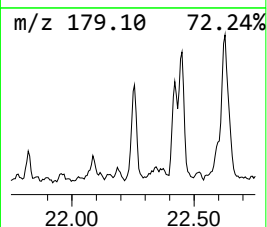
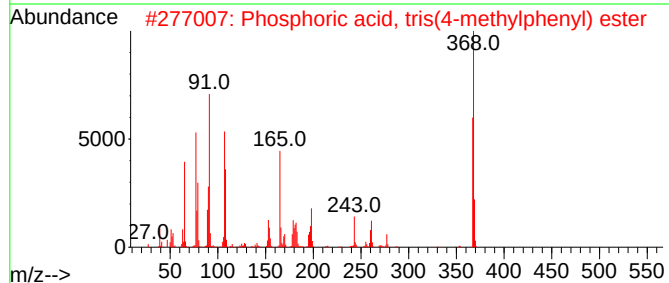
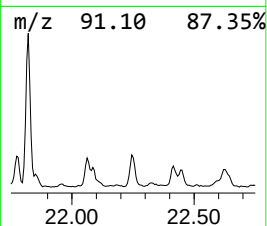
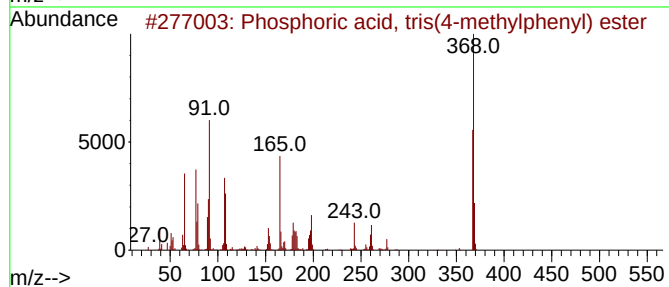
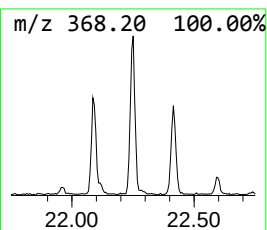
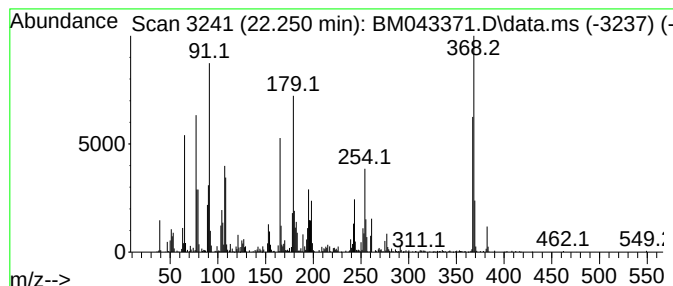
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Phosphoric acid, tris(4-met... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.250	2.37 ng/ul	795488	Chrysene-d12	21.539

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phosphoric acid, tris(4-methylph...	368	C21H21O4P	000078-32-0	97
2			Phosphoric acid, tris(4-methylph...	368	C21H21O4P	000078-32-0	95
3			Phosphoric acid, tris(3-methylph...	368	C21H21O4P	000563-04-2	89
4			Phosphoric acid, tris(4-methylph...	368	C21H21O4P	000078-32-0	86
5			Phosphoric acid, tris(4-methylph...	368	C21H21O4P	000078-32-0	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121523\
Data File : BM043371.D
Acq On : 15 Dec 2023 23:13
Operator : MA/JU
Sample : 05794-12
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
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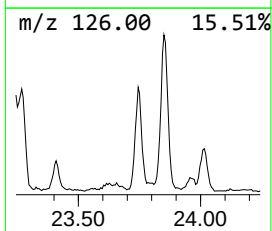
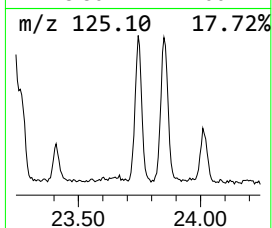
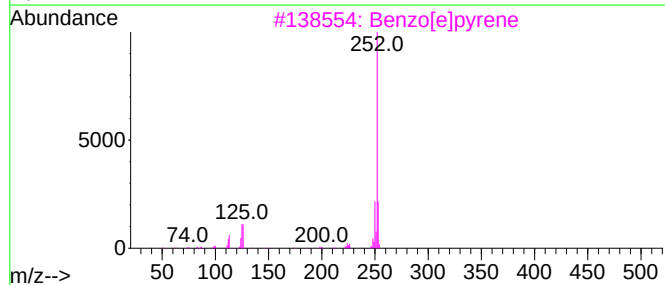
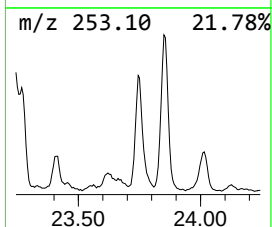
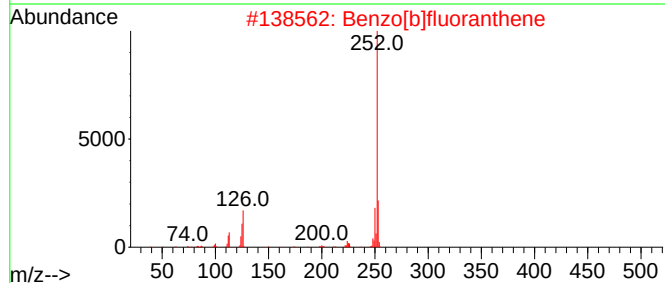
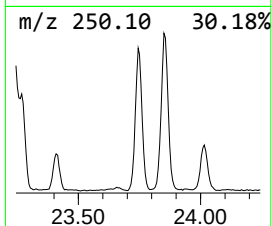
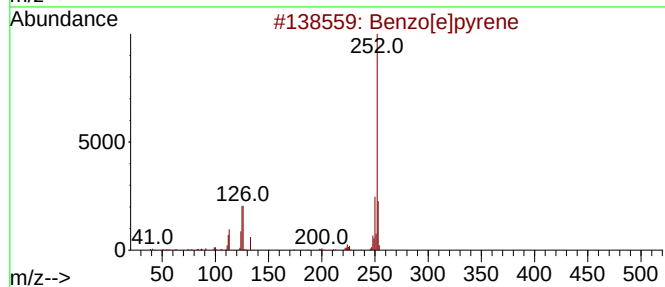
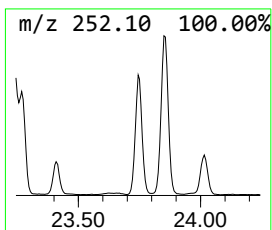
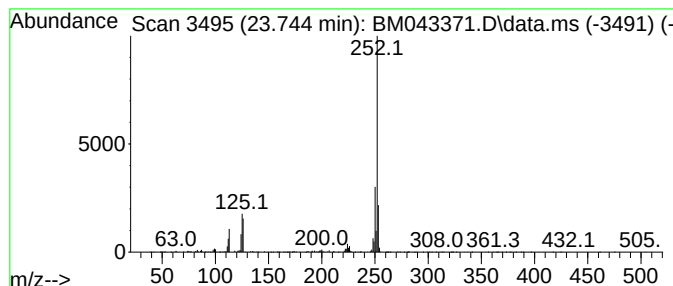
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 Benzo[e]pyrene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.744	5.60 ng/ul	1002870	Perylene-d12	23.962

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[e]pyrene	252	C20H12	000192-97-2	99
2			Benzo[b]fluoranthene	252	C20H12	000205-99-2	98
3			Benzo[e]pyrene	252	C20H12	000192-97-2	98
4			Benzo[k]fluoranthene	252	C20H12	000207-08-9	98
5			Perylene	252	C20H12	000198-55-0	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121523\
Data File : BM043371.D
Acq On : 15 Dec 2023 23:13
Operator : MA/JU
Sample : 05794-12
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BH3Q7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.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
Phenanthrene, 2...	18.292	2.1	ng/u1	556553	4	17.368	5397750	20.0
4H-Cyclopenta[d...	18.433	2.5	ng/u1	685418	4	17.368	5397750	20.0
o,p'-DDT	20.451	6.0	ng/u1	2033390	5	21.539	6722880	20.0
unknown-01	20.586	2.9	ng/u1	958451	5	21.539	6722880	20.0
Picloram, TMS d...	20.909	2.2	ng/u1	749520	5	21.539	6722880	20.0
Cyclopenta[cd]p...	21.262	2.1	ng/u1	691400	5	21.539	6722880	20.0
1H-Indene, 2,3-...	21.821	4.7	ng/u1	1589970	5	21.539	6722880	20.0
Phosphoric acid...	22.250	2.4	ng/u1	795488	5	21.539	6722880	20.0
Benzo[e]pyrene	23.744	5.6	ng/u1	1002870	6	23.962	3580560	20.0