

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121720\
 Data File : BP004348.D
 Acq On : 17 Dec 2020 19:22
 Operator : CG/JU
 Sample : L5067-02
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A54E2

Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 0.1 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP121020MA.M
 Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.834	107	110	116	rVB2	51247	74759	1.15%	0.096%
2	6.993	641	647	663	rVB2	555941	1240530	19.16%	1.588%
3	7.164	670	676	686	rBV	446490	704132	10.88%	0.901%
4	7.363	704	710	717	rBV	586859	937220	14.48%	1.200%
5	7.422	717	720	734	rVB	64222	122381	1.89%	0.157%
6	7.834	783	790	798	rBV	860717	1382080	21.35%	1.769%
7	8.534	902	909	918	rBV	662206	1073616	16.59%	1.374%
8	8.987	979	986	996	rVB	539141	890647	13.76%	1.140%
9	9.422	1054	1060	1065	rBV3	26908	43526	0.67%	0.056%
10	9.710	1101	1109	1116	rBV	438753	750007	11.59%	0.960%
11	10.246	1193	1200	1210	rBV	707942	1193710	18.44%	1.528%
12	10.628	1257	1265	1271	rBV	1221731	2056845	31.77%	2.633%
13	10.675	1271	1273	1280	rVV	38908	60282	0.93%	0.077%
14	10.763	1280	1288	1307	rVB	493227	920613	14.22%	1.179%
15	11.634	1430	1436	1439	rBV	29027	51606	0.80%	0.066%
16	12.122	1514	1519	1526	rVV10	10588	29896	0.46%	0.038%
17	12.222	1530	1536	1542	rVV2	14263	29073	0.45%	0.037%
18	12.293	1542	1548	1555	rVB	31387	51933	0.80%	0.066%
19	12.510	1580	1585	1594	rVB	25339	42108	0.65%	0.054%
20	12.793	1628	1633	1637	rVV	15893	28002	0.43%	0.036%
21	13.081	1677	1682	1686	rBV2	20456	28956	0.45%	0.037%
22	13.310	1716	1721	1726	rBV3	26151	40867	0.63%	0.052%
23	13.375	1726	1732	1739	rVV3	31227	56484	0.87%	0.072%
24	13.798	1799	1804	1809	rVB2	29072	48509	0.75%	0.062%
25	13.887	1809	1819	1823	rBV	1180187	1741552	26.90%	2.230%
26	13.928	1823	1826	1831	rVB	133583	189085	2.92%	0.242%
27	14.034	1837	1844	1848	rBV4	13909	30870	0.48%	0.040%
28	14.169	1855	1867	1880	rBV	1542136	2614290	40.39%	3.347%
29	14.387	1899	1904	1909	rVB2	16715	24033	0.37%	0.031%
30	14.481	1909	1920	1926	rBV	1871866	2832961	43.76%	3.627%
31	14.545	1926	1931	1939	rVB	47800	81722	1.26%	0.105%
32	14.675	1947	1953	1961	rBV	425150	677041	10.46%	0.867%
33	14.881	1983	1988	1995	rVB	39945	74502	1.15%	0.095%
34	14.957	1995	2001	2006	rVB	19793	32416	0.50%	0.041%

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Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP121020MA.M
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35	15.104	2018	2026	2031	rBV3	26737	67365	1.04%	0.086%
36	15.287	2047	2057	2063	rBV3	33697	93352	1.44%	0.120%
37	15.345	2063	2067	2072	rVB3	23769	36609	0.57%	0.047%
38	15.481	2082	2090	2095	rBV	1871957	2810974	43.42%	3.599%
39	15.534	2095	2099	2105	rVB4	61700	86679	1.34%	0.111%
40	15.604	2105	2111	2118	rBV	270877	404111	6.24%	0.517%
41	15.722	2123	2131	2133	rBV2	19072	40507	0.63%	0.052%
42	15.834	2145	2150	2159	rVB2	32444	68903	1.06%	0.088%
43	15.975	2170	2174	2182	rVB	31142	67157	1.04%	0.086%
44	16.063	2182	2189	2195	rBV2	15222	35191	0.54%	0.045%
45	16.169	2202	2207	2210	rBV3	17389	26964	0.42%	0.035%
46	16.210	2210	2214	2221	rVV7	44129	88172	1.36%	0.113%
47	16.539	2268	2270	2275	rVV5	26892	44634	0.69%	0.057%
48	16.592	2275	2279	2285	rVB2	27042	41773	0.65%	0.053%
49	16.775	2307	2310	2314	rVV2	22338	34930	0.54%	0.045%
50	16.816	2314	2317	2320	rVB	29944	35773	0.55%	0.046%
51	16.892	2326	2330	2337	rVB2	66733	103298	1.60%	0.132%
52	17.057	2355	2358	2366	rVB	72382	109001	1.68%	0.140%
53	17.239	2383	2389	2393	rBV	2380352	3434864	53.06%	4.397%
54	17.281	2393	2396	2401	rVV	1039254	1517210	23.44%	1.942%
55	17.339	2401	2406	2411	rVV	2127662	3188399	49.26%	4.082%
56	17.375	2411	2412	2417	rVV	240372	213871	3.30%	0.274%
57	17.434	2417	2422	2428	rVB3	47300	85200	1.32%	0.109%
58	17.528	2435	2438	2440	rBV2	18668	24196	0.37%	0.031%
59	17.645	2450	2458	2462	rBV2	92836	163862	2.53%	0.210%
60	17.763	2474	2478	2482	rVB2	31660	46574	0.72%	0.060%
61	17.822	2485	2488	2494	rVB	73327	88182	1.36%	0.113%
62	17.963	2509	2512	2518	rVB	38549	63820	0.99%	0.082%
63	18.110	2533	2537	2541	rBV	251952	377487	5.83%	0.483%
64	18.157	2541	2545	2550	rVB	340883	496738	7.67%	0.636%
65	18.239	2555	2559	2563	rVV	93196	127598	1.97%	0.163%
66	18.298	2564	2569	2573	rVV2	342759	609967	9.42%	0.781%
67	18.334	2573	2575	2582	rVB	202472	262392	4.05%	0.336%
68	18.598	2616	2620	2623	rBV	187150	252547	3.90%	0.323%
69	18.633	2623	2626	2632	rVB3	186329	261216	4.04%	0.334%
70	18.716	2638	2640	2645	rBV3	35044	42189	0.65%	0.054%
71	18.833	2658	2660	2664	rVV2	71365	72885	1.13%	0.093%

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72	18.881	2664	2668	2672	rVV2	138770	206310	3.19%	0.264%
73	18.922	2672	2675	2679	rVB2	72753	89088	1.38%	0.114%
74	19.004	2684	2689	2693	rBV	220652	360778	5.57%	0.462%
75	19.139	2708	2712	2718	rBV	169224	283762	4.38%	0.363%
76	19.257	2727	2732	2738	rVB	2721333	3415711	52.77%	4.373%
77	19.351	2746	2748	2752	rVB2	72983	88749	1.37%	0.114%
78	19.404	2754	2757	2760	rVV	108699	123033	1.90%	0.158%
79	19.580	2782	2787	2790	rBV2	3095864	4639801	71.68%	5.940%
80	19.610	2790	2792	2800	rVV	2470868	3073334	47.48%	3.935%
81	19.680	2801	2804	2811	rVB2	146920	235250	3.63%	0.301%
82	19.786	2819	2822	2827	rVB	106716	139351	2.15%	0.178%
83	19.928	2841	2846	2852	rBV3	224256	545111	8.42%	0.698%
84	20.092	2871	2874	2880	rVB	355962	472075	7.29%	0.604%
85	20.192	2888	2891	2894	rVB	150103	164706	2.54%	0.211%
86	20.251	2897	2901	2904	rVB2	272324	394599	6.10%	0.505%
87	20.386	2921	2924	2927	rBV	175305	208560	3.22%	0.267%
88	20.427	2929	2931	2935	rVB	195951	222206	3.43%	0.284%
89	20.822	2995	2998	3005	rVB	309225	452108	6.98%	0.579%
90	21.027	3030	3033	3035	rBV	212900	252056	3.89%	0.323%
91	21.057	3035	3038	3044	rVB3	359827	659224	10.18%	0.844%
92	21.245	3067	3070	3074	rVB	661709	772336	11.93%	0.989%
93	21.333	3079	3085	3089	rVV2	3542342	6473207	100.00%	8.287%
94	21.375	3089	3092	3102	rVB	1448614	1932639	29.86%	2.474%
95	22.980	3360	3365	3371	rBV2	2315262	4564325	70.51%	5.843%
96	23.498	3448	3453	3457	rBV	631909	1206396	18.64%	1.544%
97	23.551	3457	3462	3466	rVV	1848011	3407417	52.64%	4.362%
98	23.598	3466	3470	3477	rVB	1365261	2477785	38.28%	3.172%
99	23.704	3481	3488	3494	rBV	1920915	3978637	61.46%	5.094%
100	24.298	3584	3589	3599	rVB2	739597	1591475	24.59%	2.037%

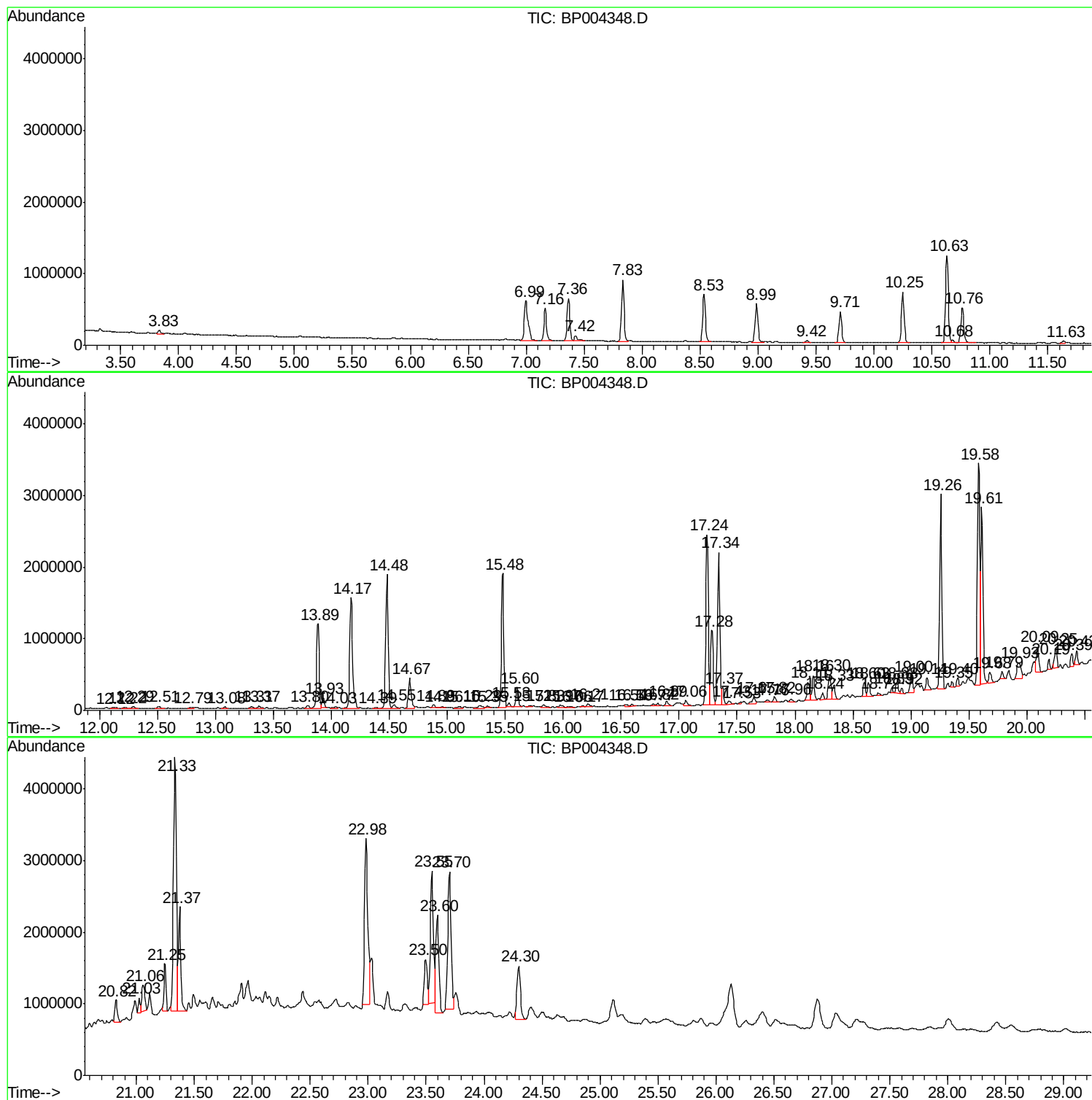
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP121020MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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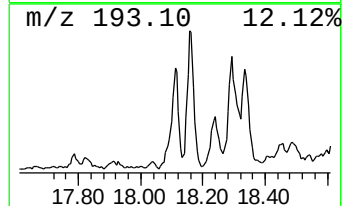
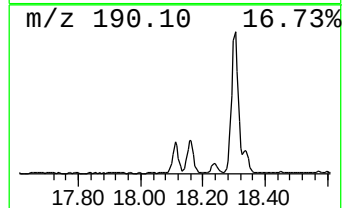
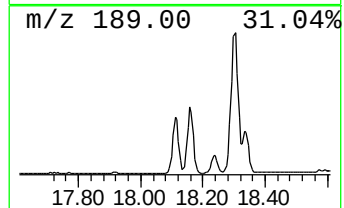
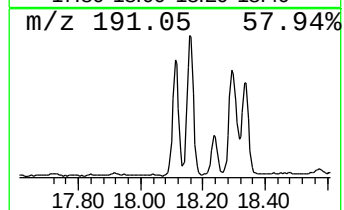
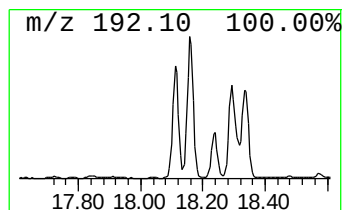
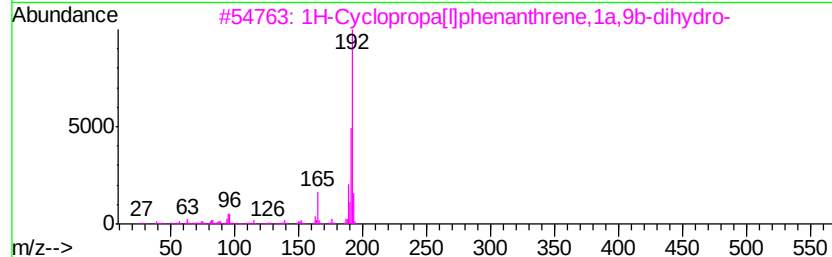
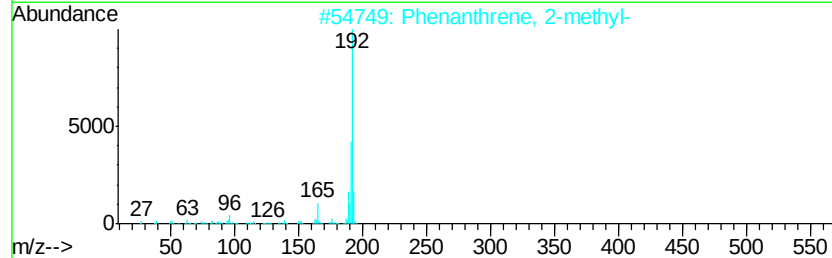
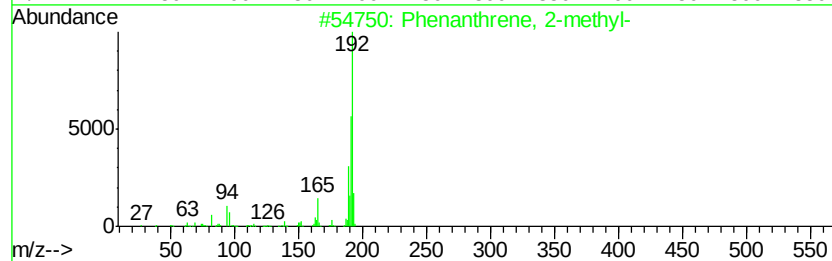
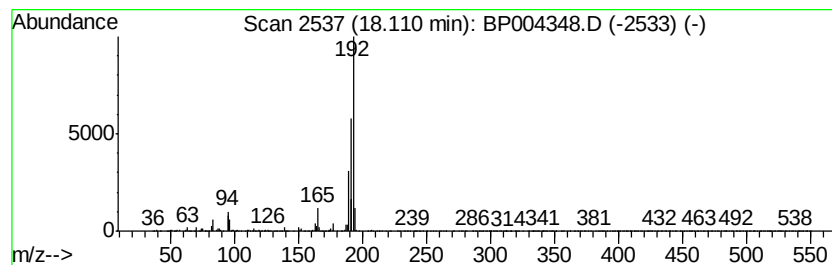
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP121020MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Phenanthrene, 2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.11	2.20 ng/ul	377487	Phenanthrene-d10	17.24	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93
2	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	90
3	1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	89
4	Anthracene, 2-methyl-	192	C15H12	000613-12-7	89
5	Anthracene, 9-methyl-	192	C15H12	000779-02-2	83



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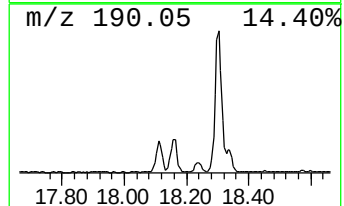
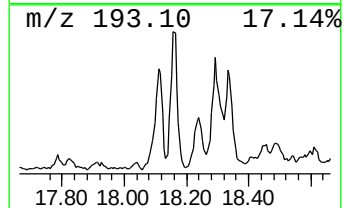
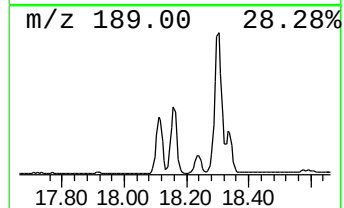
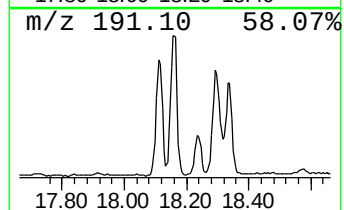
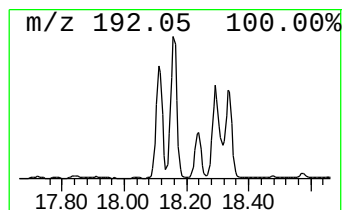
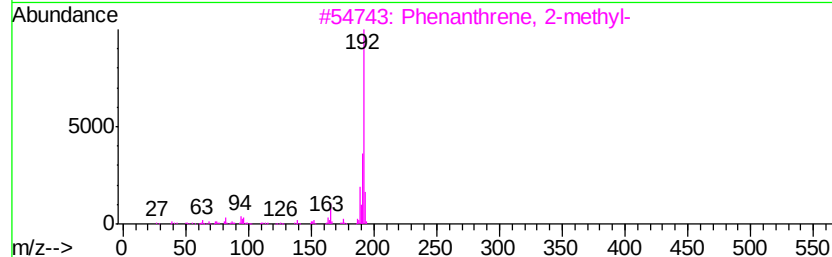
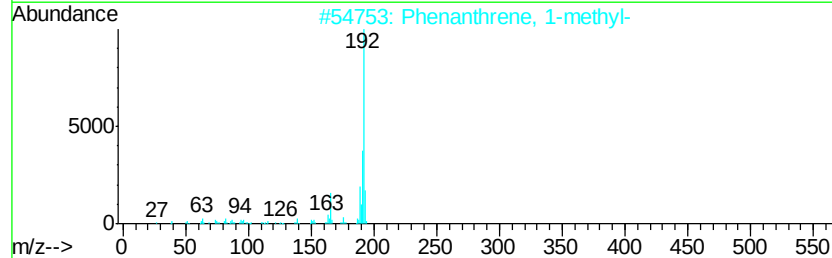
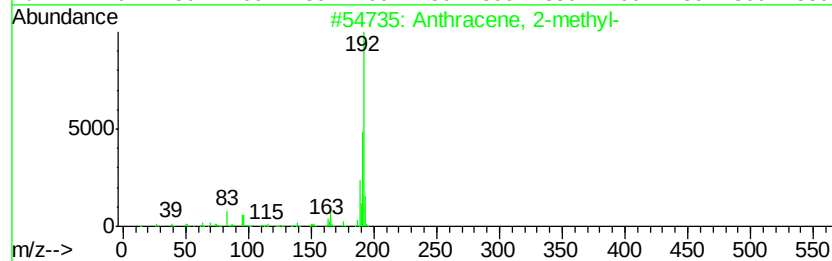
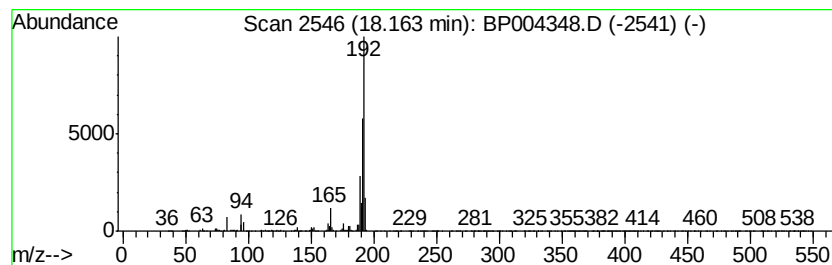
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP121020MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Anthracene, 2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.16	2.89 ng/ul	496738	Phenanthrene-d10	17.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
2		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95
3		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
4		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
5		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93



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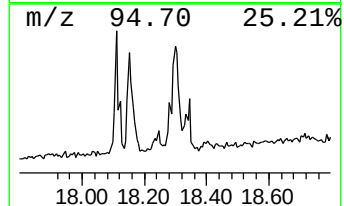
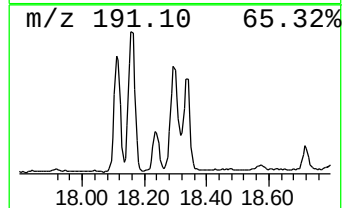
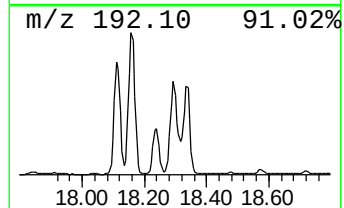
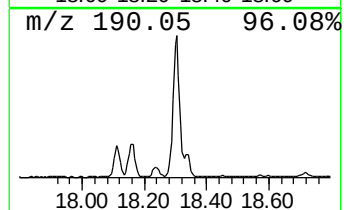
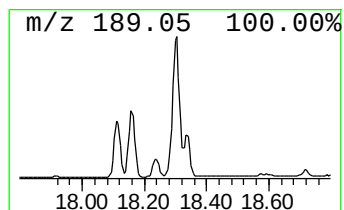
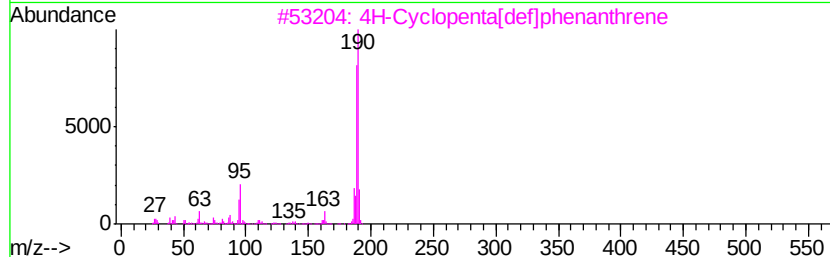
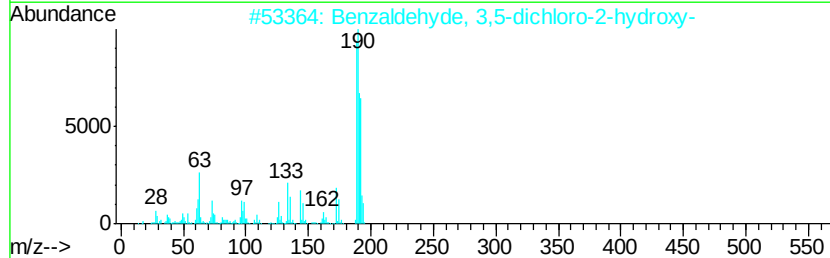
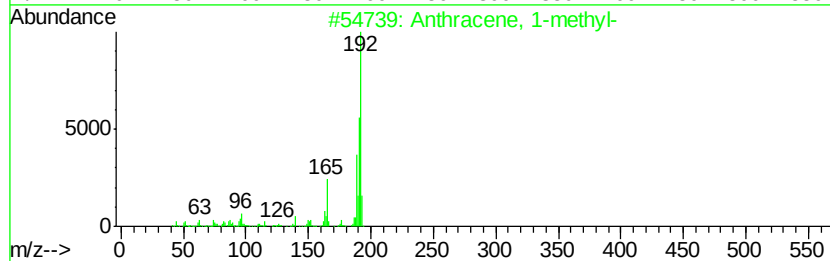
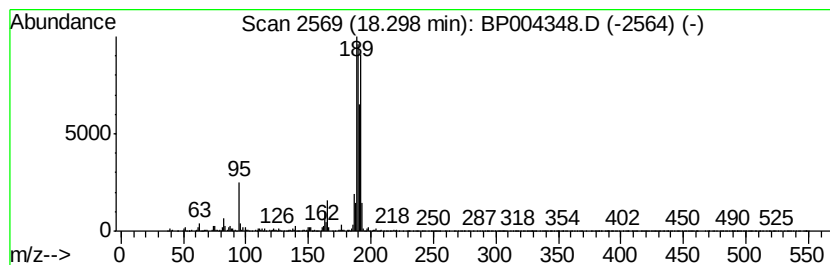
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Peak Number 3 Anthracene, 1-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.		
18.30	3.55 ng/ul	609967	Phenanthrene-d10	17.24		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Anthracene, 1-methyl-	192	C15H12	000610-48-0	81
2		Benzaldehyde, 3,5-dichloro-2-hyd...	190	C7H4Cl2O2	000090-60-8	72
3		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	60
4		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	53
5		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	46



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121720\
Data File : BP004348.D
Acq On : 17 Dec 2020 19:22
Operator : CG/JU
Sample : L5067-02
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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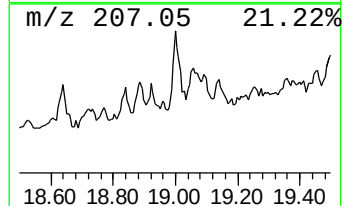
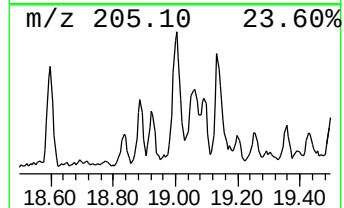
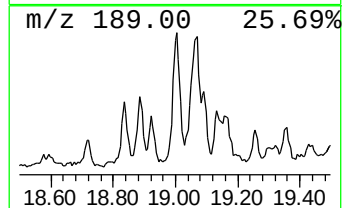
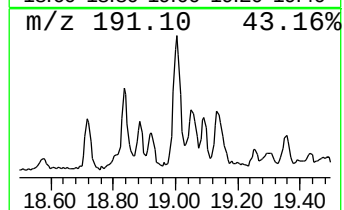
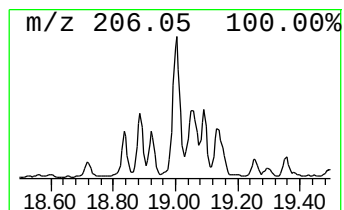
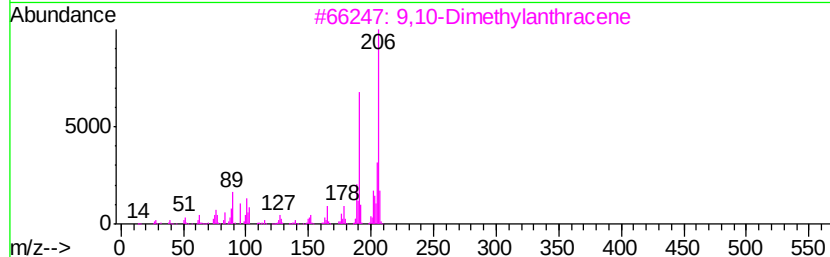
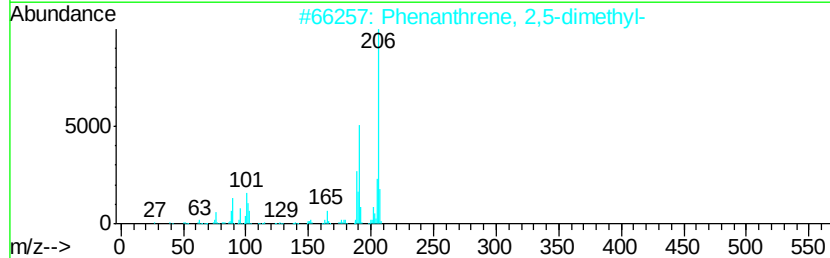
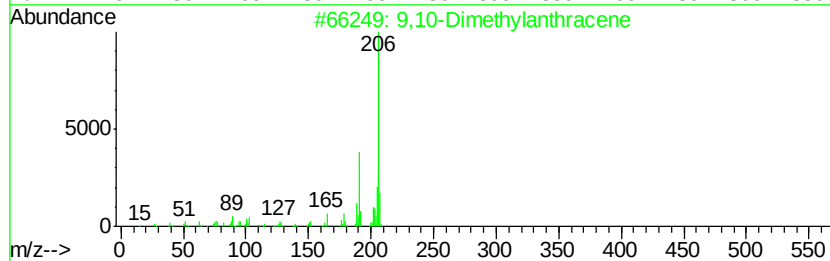
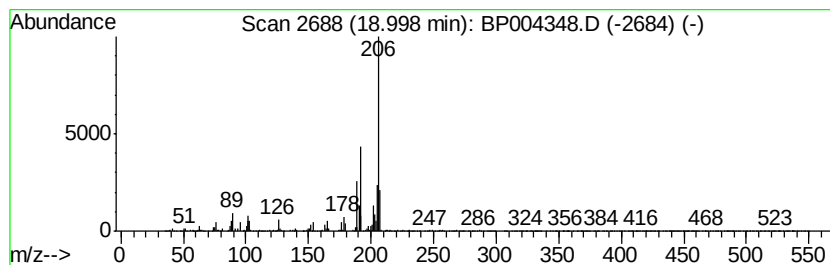
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 9,10-Dimethylantracene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.00	2.10 ng/ul	360778	Phenanthrene-d10	17.24

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9,10-Dimethylantracene	206	C16H14	000781-43-1	96
2			Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	93
3			9,10-Dimethylantracene	206	C16H14	000781-43-1	91
4			di-p-Tolylacetylene	206	C16H14	002789-88-0	91
5			Phenanthrene, 2,7-dimethyl-	206	C16H14	001576-69-8	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121720\
Data File : BP004348.D
Acq On : 17 Dec 2020 19:22
Operator : CG/JU
Sample : L5067-02
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A54E2

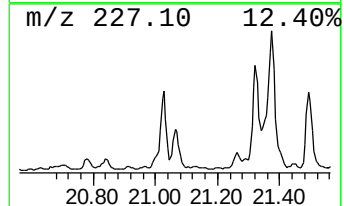
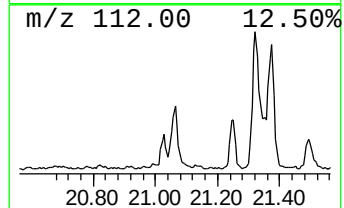
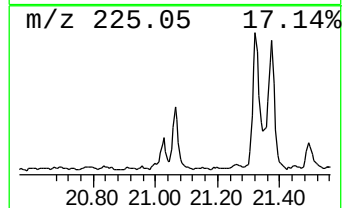
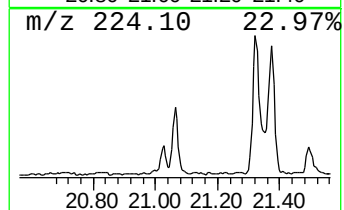
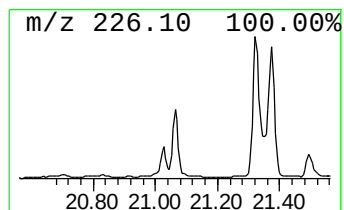
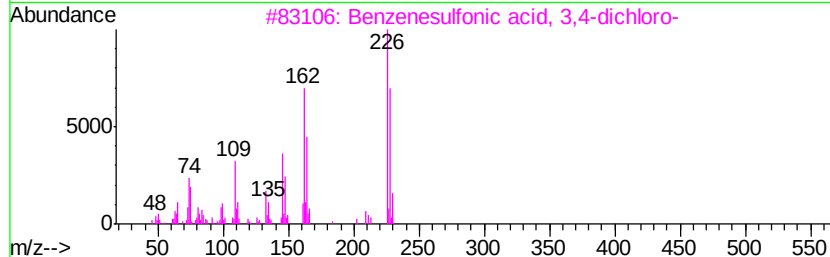
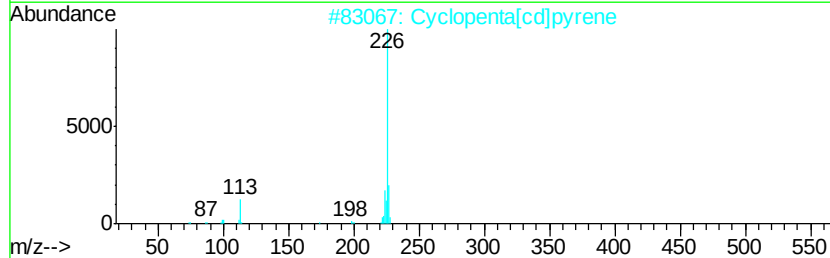
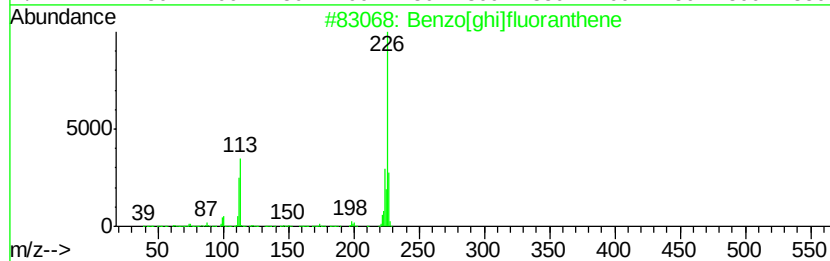
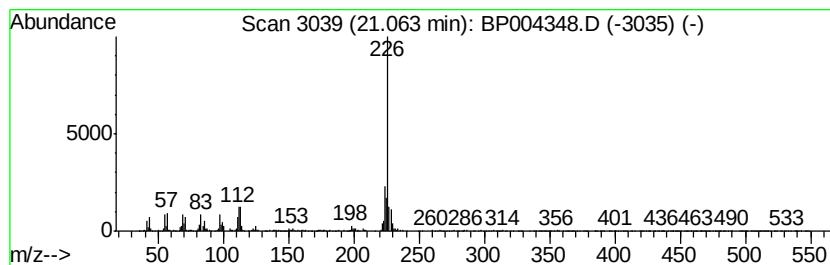
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 unknown-01 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.06	2.04 ng/ul	659224	Chrysene-d12	21.34

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[ghi]fluoranthene	226	C18H10	000203-12-3	47
2		Cyclopenta[cd]pyrene	226	C18H10	027208-37-3	46
3		Benzenesulfonic acid, 3,4-dichloro-	226	C6H4Cl2O3S	000939-95-7	35
4		1,3-Diamino-5,6-dihydro-9-methyl...	226	C13H14N4	037436-39-8	23
5		Benzonitrile, 2-(5-methyl-2-thie...	226	C13H10N2S	315670-19-0	23



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121720\
Data File : BP004348.D
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Operator : CG/JU
Sample : L5067-02
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A54E2

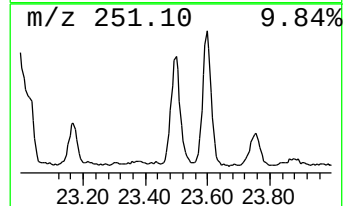
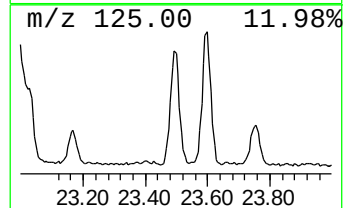
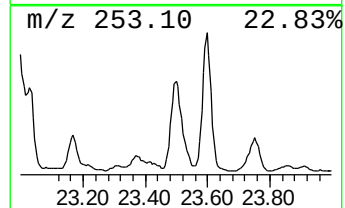
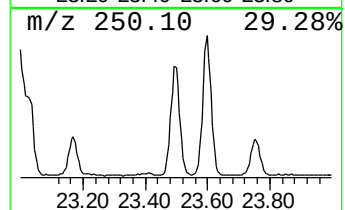
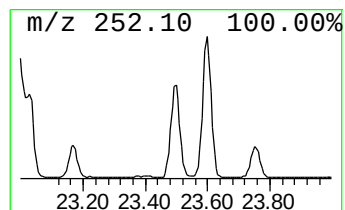
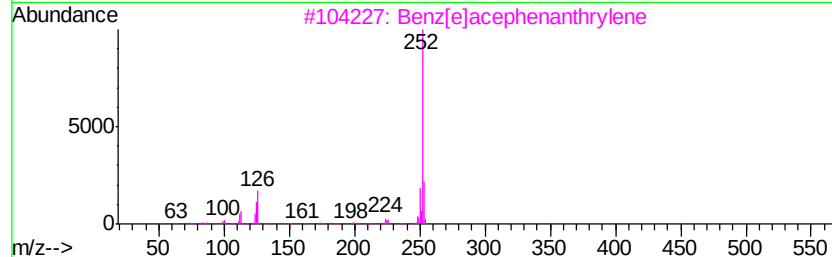
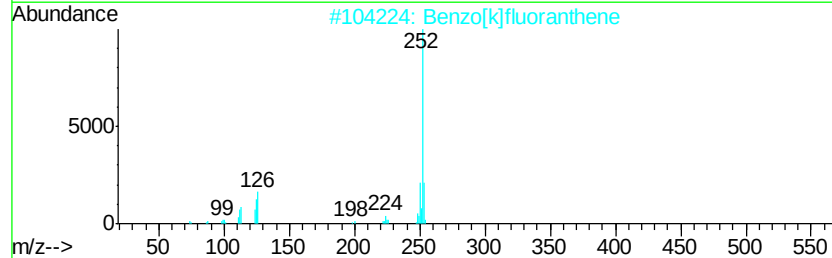
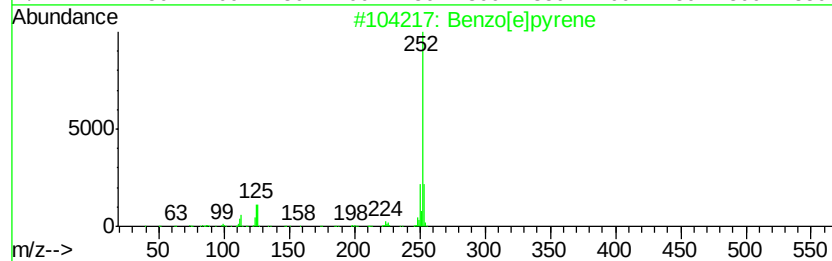
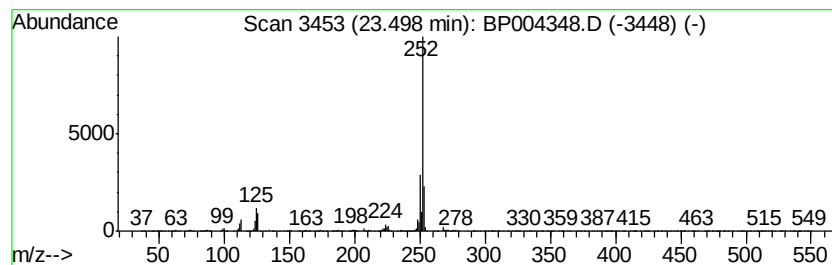
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Benzo[e]pyrene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.50	6.06 ng/ul	1206400	Perylene-d12	23.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[e]pyrene	252	C20H12	000192-97-2	99
2		Benzo[k]fluoranthene	252	C20H12	000207-08-9	98
3		Benzo[e]acephenanthrylene	252	C20H12	000205-99-2	96
4		Benzo[i]fluoranthene	252	C20H12	000205-82-3	95
5		Benzo[a]pyrene	252	C20H12	000050-32-8	94



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121720\
Data File : BP004348.D
Acq On : 17 Dec 2020 19:22
Operator : CG/JU
Sample : L5067-02
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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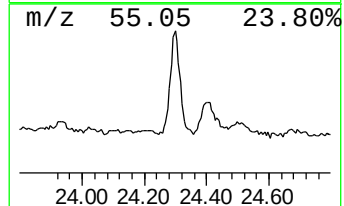
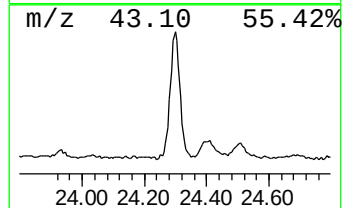
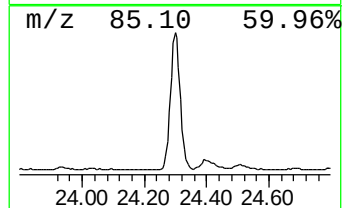
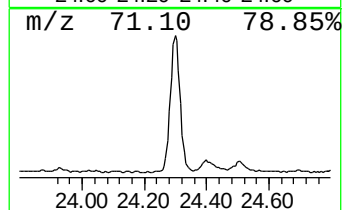
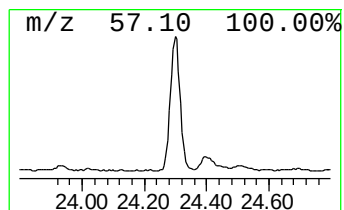
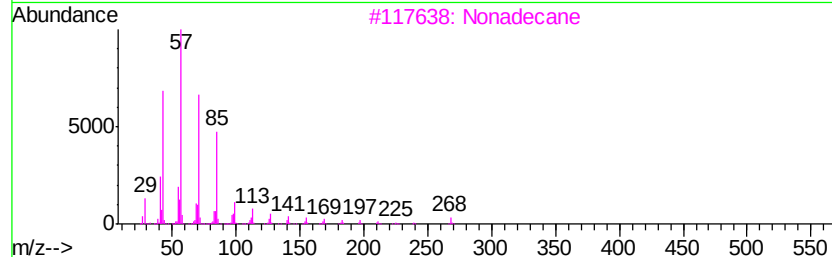
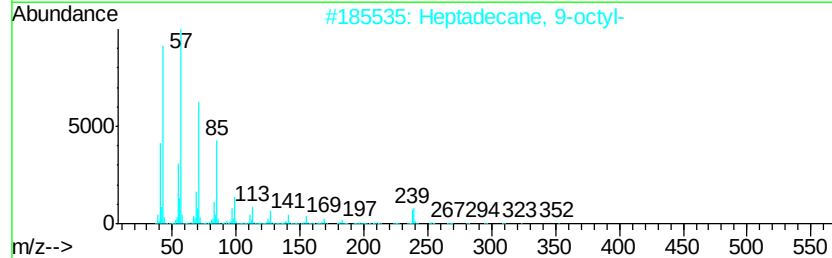
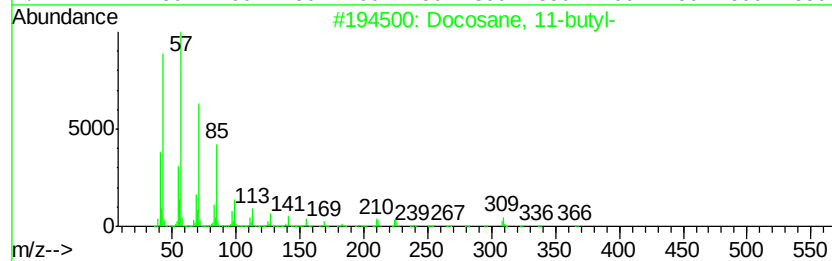
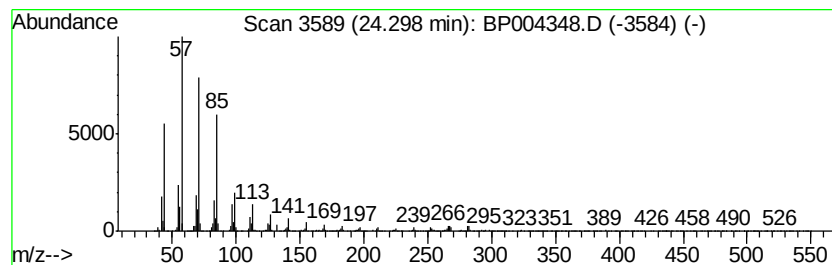
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP121020MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.30	8.00 ng/ul	1591480	Perylene-d12	23.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Docosane, 11-butyl-	366	C26H54	013475-76-8	94
2		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	93
3		Nonadecane	268	C19H40	000629-92-5	93
4		2-Bromo dodecane	248	C12H25Br	013187-99-0	93
5		Docosane, 11-butyl-	366	C26H54	013475-76-8	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP121720\
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Sample : L5067-02
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP121020MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Phenanthrene, 2-m...	18.11	2.2	ng/ul	377487	4	17.24	3434860	20.0
Anthracene, 2-met...	18.16	2.9	ng/ul	496738	4	17.24	3434860	20.0
Anthracene, 1-met...	18.30	3.5	ng/ul	609967	4	17.24	3434860	20.0
9,10-Dimethylanthe...	19.00	2.1	ng/ul	360778	4	17.24	3434860	20.0
unknown-01	21.06	2.0	ng/ul	659224	5	21.34	6473210	20.0
Benzo[e]pyrene	23.50	6.1	ng/ul	1206400	6	23.70	3978640	20.0
(DEL) Alkane: Str...	24.30	8.0	ng/ul	1591480	6	23.70	3978640	20.0