

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080320\
 Data File : BM026986.D
 Acq On : 03 Aug 2020 19:50
 Operator : JU/CG
 Sample : L3424-05DL 25X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0S80DL

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_M\METHODS\SOM-EPA-BM072720MA.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	3.940	158	162	166	rVB3	7546	10596	0.30%	0.043%
2	5.687	454	459	464	rBV5	4301	6360	0.18%	0.026%
3	6.834	650	654	663	rVB	14045	23926	0.69%	0.096%
4	7.004	677	683	688	rBV2	11195	17100	0.49%	0.069%
5	7.198	711	716	724	rVB2	15199	24342	0.70%	0.098%
6	7.663	788	795	802	rBV	428508	652349	18.78%	2.621%
7	8.198	880	886	896	rBV	27550	48573	1.40%	0.195%
8	8.363	907	914	919	rVV2	17178	25981	0.75%	0.104%
9	8.804	984	989	994	rVB2	14894	21954	0.63%	0.088%
10	9.522	1104	1111	1116	rBV2	10556	15257	0.44%	0.061%
11	10.057	1196	1202	1209	rVB2	16650	28334	0.82%	0.114%
12	10.433	1258	1266	1271	rBV	551035	898177	25.85%	3.609%
13	10.481	1271	1274	1284	rVB	65285	107039	3.08%	0.430%
14	10.569	1284	1289	1292	rBV3	4768	7652	0.22%	0.031%
15	11.816	1496	1501	1505	rVB6	4923	7598	0.22%	0.031%
16	11.939	1515	1522	1529	rBV3	5378	12129	0.35%	0.049%
17	12.104	1541	1550	1558	rVB	25752	42272	1.22%	0.170%
18	12.327	1581	1588	1592	rVB2	9525	16520	0.48%	0.066%
19	13.127	1717	1724	1732	rVB2	7953	12924	0.37%	0.052%
20	13.704	1818	1822	1826	rVV	28738	40154	1.16%	0.161%
21	13.745	1826	1829	1837	rVB3	12567	20689	0.60%	0.083%
22	14.010	1864	1874	1889	rBV2	158885	309537	8.91%	1.244%
23	14.292	1916	1922	1929	rVV	797430	1145391	32.97%	4.602%
24	14.357	1929	1933	1937	rVB2	16919	25614	0.74%	0.103%
25	14.486	1950	1955	1962	rVB5	7217	12815	0.37%	0.051%
26	14.692	1983	1990	1999	rBV	33742	50566	1.46%	0.203%
27	15.168	2066	2071	2077	rBV5	8541	16813	0.48%	0.068%
28	15.286	2085	2091	2096	rBV	47774	64013	1.84%	0.257%
29	15.339	2096	2100	2106	rVV3	11887	22160	0.64%	0.089%
30	15.398	2108	2110	2114	rVB4	7211	8538	0.25%	0.034%
31	15.545	2128	2135	2138	rBV6	7518	15056	0.43%	0.060%
32	15.645	2148	2152	2155	rBV4	8639	12346	0.36%	0.050%
33	15.680	2155	2158	2162	rVB5	6429	9192	0.26%	0.037%
34	15.751	2167	2170	2173	rVV3	5843	8115	0.23%	0.033%

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Integration Parameters: LSCINT.P

Integrator: RTE
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Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM072720MA.M
 Title : SVOA CALIBRATION

35	15.786	2173	2176	2183	rVB3	9646	18871	0.54%	0.076%
36	16.010	2208	2214	2221	rBV3	31265	62282	1.79%	0.250%
37	16.068	2221	2224	2229	rVB5	9188	13594	0.39%	0.055%
38	16.304	2258	2264	2266	rBV6	3409	6231	0.18%	0.025%
39	16.615	2312	2317	2323	rBV	17804	27063	0.78%	0.109%
40	16.686	2323	2329	2339	rVB2	40080	67715	1.95%	0.272%
41	16.833	2348	2354	2356	rBV5	11709	19620	0.56%	0.079%
42	16.886	2360	2363	2366	rVB4	7139	8819	0.25%	0.035%
43	17.033	2382	2388	2393	rBV	939614	1317614	37.92%	5.294%
44	17.074	2393	2395	2400	rVB	122954	140156	4.03%	0.563%
45	17.133	2400	2405	2406	rBV2	48856	63814	1.84%	0.256%
46	17.168	2406	2411	2419	rVB	334199	500886	14.42%	2.013%
47	17.433	2449	2456	2463	rBV	33586	60032	1.73%	0.241%
48	17.480	2463	2464	2466	rVV2	8960	6866	0.20%	0.028%
49	17.527	2469	2472	2475	rVV5	8532	15511	0.45%	0.062%
50	17.568	2475	2479	2481	rVV4	15672	24807	0.71%	0.100%
51	17.615	2485	2487	2492	rVV6	9482	11770	0.34%	0.047%
52	17.692	2496	2500	2503	rVV6	11956	18549	0.53%	0.075%
53	17.733	2504	2507	2516	rVB7	14674	26204	0.75%	0.105%
54	17.821	2519	2522	2525	rBV5	4239	6029	0.17%	0.024%
55	17.915	2534	2538	2541	rBV3	8854	11897	0.34%	0.048%
56	17.956	2541	2545	2551	rVB4	34312	56983	1.64%	0.229%
57	18.033	2554	2558	2563	rBV2	48300	72301	2.08%	0.291%
58	18.109	2563	2571	2580	rVB3	38914	95053	2.74%	0.382%
59	18.233	2587	2592	2594	rBV6	8169	15434	0.44%	0.062%
60	18.298	2599	2603	2610	rVB	31962	49168	1.42%	0.198%
61	18.386	2615	2618	2621	rVV2	22200	34270	0.99%	0.138%
62	18.445	2625	2628	2633	rVB2	33067	44004	1.27%	0.177%
63	18.662	2663	2665	2669	rVB5	6729	5992	0.17%	0.024%
64	18.751	2678	2680	2684	rVB4	10455	9735	0.28%	0.039%
65	18.827	2689	2693	2698	rBV3	40429	65440	1.88%	0.263%
66	18.909	2699	2707	2711	rBV5	60780	102805	2.96%	0.413%
67	18.968	2713	2717	2726	rVB	103914	141474	4.07%	0.568%
68	19.045	2727	2730	2733	rBV2	22768	26027	0.75%	0.105%
69	19.092	2733	2738	2745	rBV	640741	832887	23.97%	3.347%
70	19.186	2752	2754	2759	rVB	21616	26007	0.75%	0.105%
71	19.245	2759	2764	2768	rBV	81169	116804	3.36%	0.469%

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 Title : SVOA CALIBRATION

72	19.456	2790	2800	2809	rBV	837662	1353882	38.97%	5.440%
73	19.639	2828	2831	2837	rBV2	35311	49196	1.42%	0.198%
74	19.692	2837	2840	2846	rVB2	40538	65588	1.89%	0.264%
75	19.798	2851	2858	2862	rBV3	89712	180905	5.21%	0.727%
76	19.927	2875	2880	2889	rVB4	119249	307550	8.85%	1.236%
77	20.009	2892	2894	2895	rBV2	15996	13084	0.38%	0.053%
78	20.056	2899	2902	2905	rVB2	26497	32316	0.93%	0.130%
79	20.115	2906	2912	2915	rBV2	108573	180386	5.19%	0.725%
80	20.250	2932	2935	2939	rBV	108752	143447	4.13%	0.576%
81	20.298	2940	2943	2951	rVV5	47902	106461	3.06%	0.428%
82	20.515	2976	2980	2984	rBV5	50483	91829	2.64%	0.369%
83	20.656	3001	3004	3008	rBV3	33460	57393	1.65%	0.231%
84	20.703	3008	3012	3018	rVB2	110946	196398	5.65%	0.789%
85	20.868	3034	3040	3044	rBV3	94135	218818	6.30%	0.879%
86	20.909	3044	3047	3049	rVV2	93449	104533	3.01%	0.420%
87	20.945	3049	3053	3058	rVV	255966	347173	9.99%	1.395%
88	21.221	3093	3100	3104	rBV2	1382099	2549241	73.37%	10.243%
89	21.256	3104	3106	3114	rVB	580790	805752	23.19%	3.238%
90	21.392	3126	3129	3136	rBV2	117439	189166	5.44%	0.760%
91	21.527	3148	3152	3155	rBV	138511	185400	5.34%	0.745%
92	21.821	3199	3202	3209	rVB3	130123	172039	4.95%	0.691%
93	22.509	3315	3319	3325	rBV2	117780	260982	7.51%	1.049%
94	22.633	3336	3340	3347	rVB2	215193	339165	9.76%	1.363%
95	22.774	3358	3364	3369	rBV2	1460323	3474516	100.00%	13.961%
96	22.939	3388	3392	3398	rVB	232768	354181	10.19%	1.423%
97	23.244	3438	3444	3450	rVB	648379	1177861	33.90%	4.733%
98	23.333	3454	3459	3467	rVB	560184	1030104	29.65%	4.139%
99	23.433	3469	3476	3481	rBV	685895	1330644	38.30%	5.347%
100	25.638	3843	3851	3860	rVB	622209	1705913	49.10%	6.855%

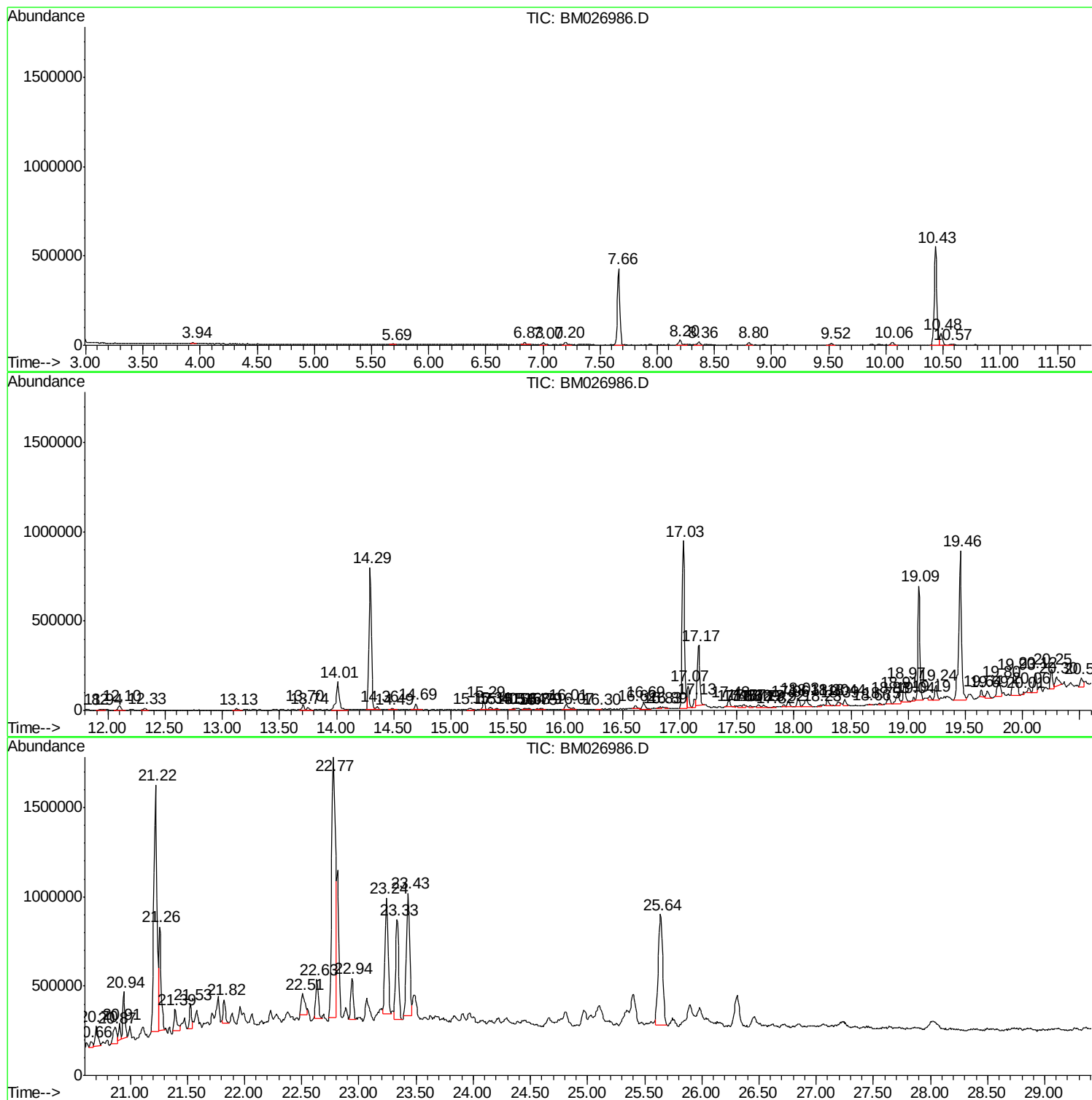
Sum of corrected areas: 24886719

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Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM072720MA.M
Quant Title : SVOA CALIBRATION

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



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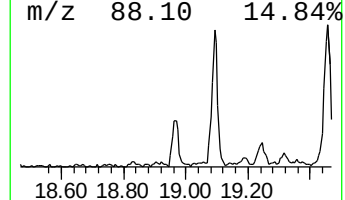
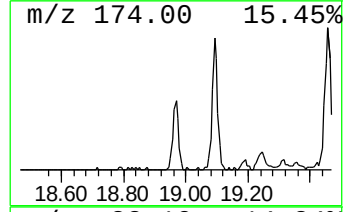
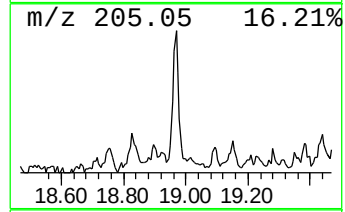
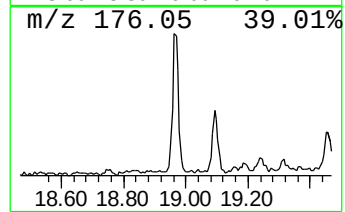
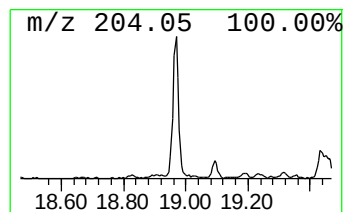
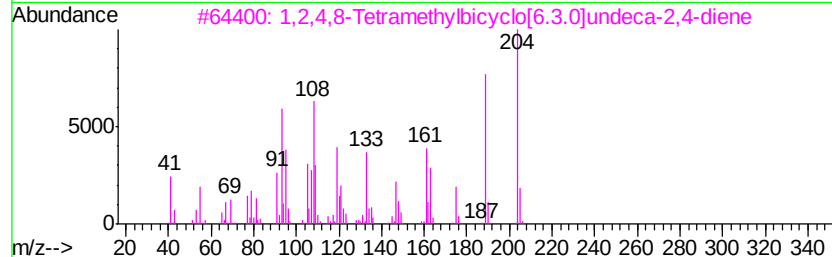
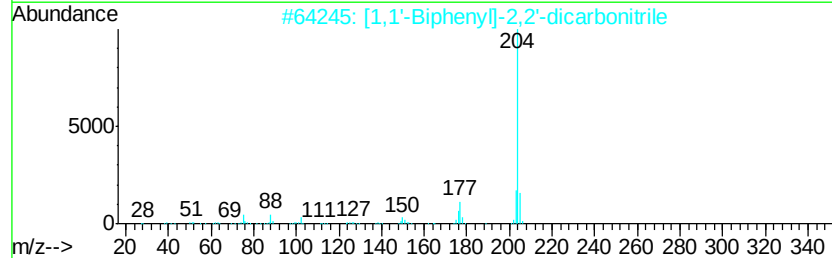
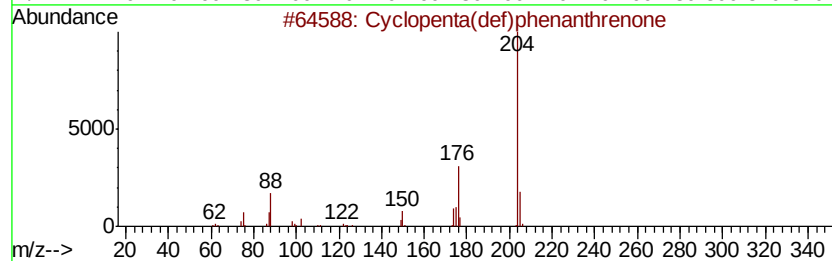
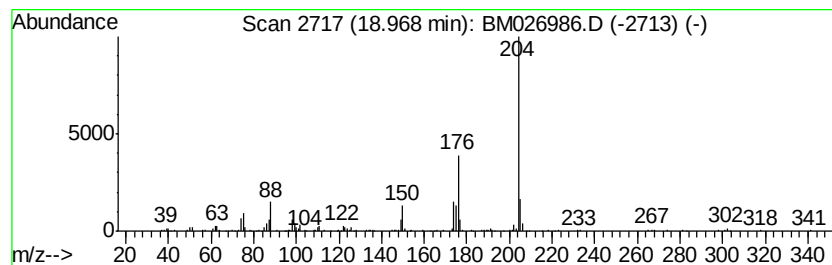
Instrument :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM072720MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Cyclopenta(def)phenanthrenone Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD		R.T.	
18.97	2.15 ng/ul	141474	Phenanthrene-d10		17.03	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cvclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	91
2		[1,1'-Biphenyl]-2,2'-dicarbonitrile	204	C14H8N2	004341-02-0	50
3		1,2,4,8-Tetramethylbicyclo[6.3.0]undeca-2,4-diene	204	C15H24	137235-51-9	49
4		[1,1'-Biphenyl]-4,4'-dicarbonitrile	204	C14H8N2	001591-30-6	47
5		9-Acridinecarbonitrile	204	C14H8N2	005326-19-2	47



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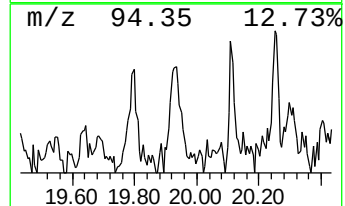
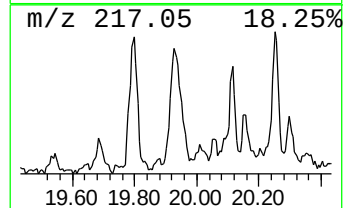
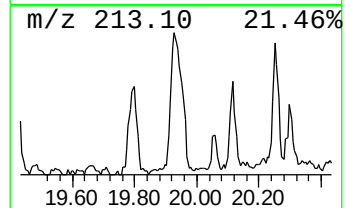
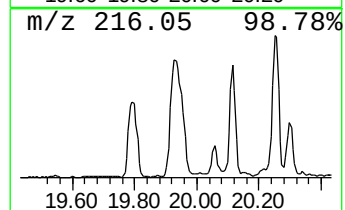
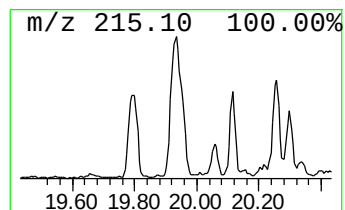
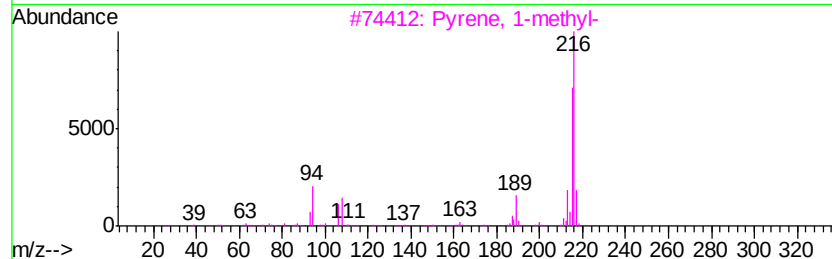
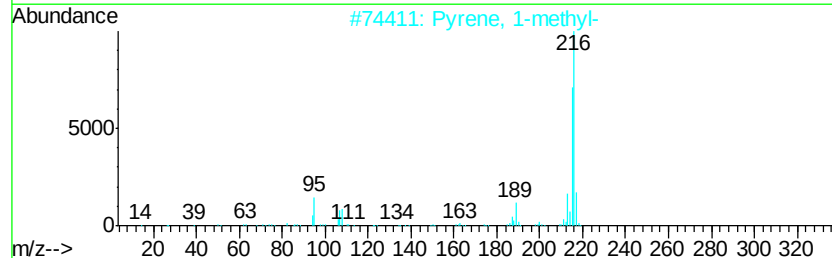
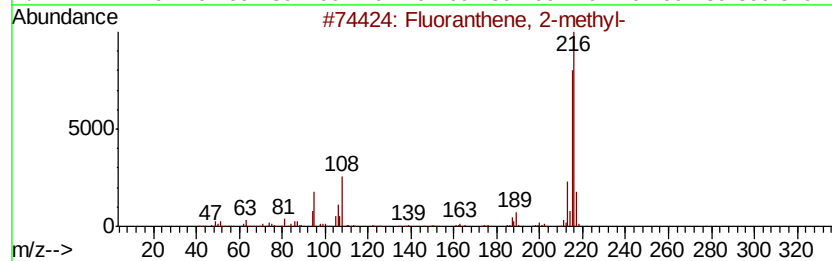
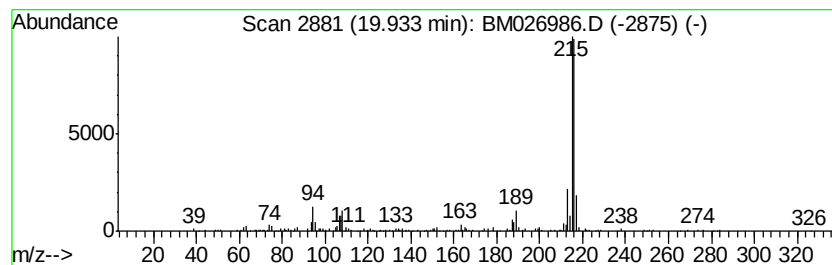
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM072720MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Fluoranthene, 2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.93	2.41 ng/ul	307550	Chrysene-d12	21.22

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	95
2		Pyrene, 1-methyl-	216	C17H12	002381-21-7	95
3		Pyrene, 1-methyl-	216	C17H12	002381-21-7	92
4		Pyrene, 1-methyl-	216	C17H12	002381-21-7	90
5		11H-Benzo[b]fluorene	216	C17H12	000243-17-4	87



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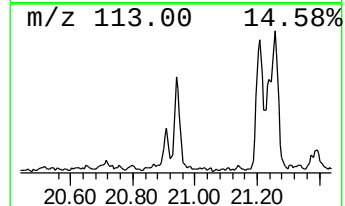
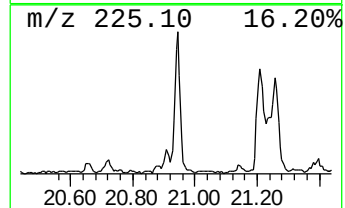
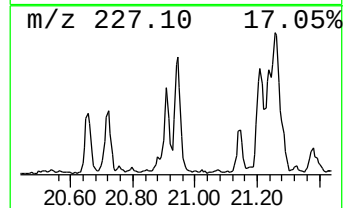
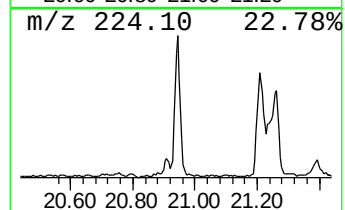
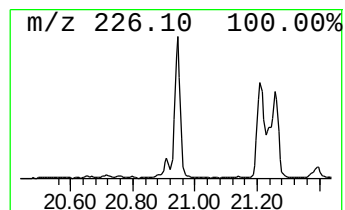
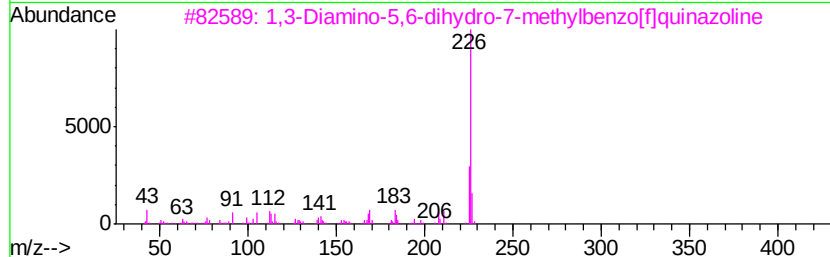
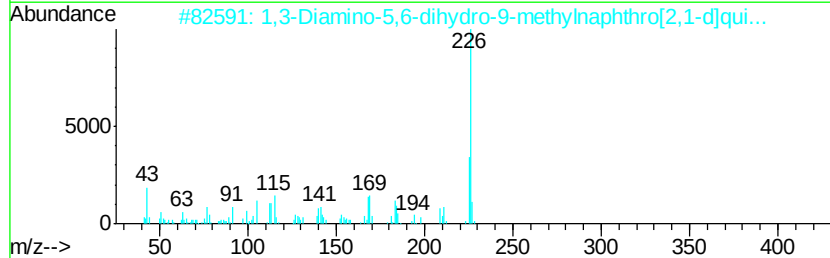
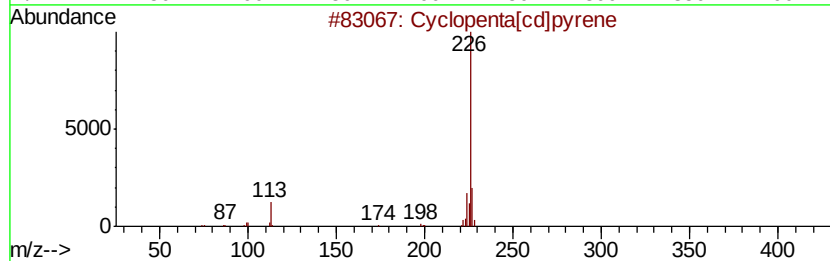
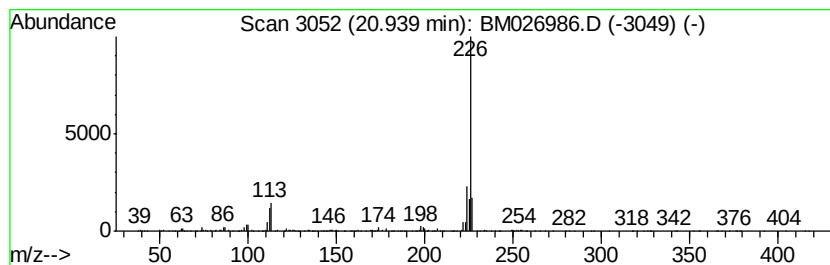
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Cyclopenta[cd]pyrene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.94	2.72 ng/ul	347173	Chrysene-d12	21.22

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopenta[cd]pyrene	226	C18H10	027208-37-3	72
2		1,3-Diamino-5,6-dihydro-9-methyl...	226	C13H14N4	037436-39-8	53
3		1,3-Diamino-5,6-dihydro-7-methyl...	226	C13H14N4	037436-37-6	53
4		Benzo[ghi]fluoranthene	226	C18H10	000203-12-3	47
5		9H-Thioxanthen-9-one, 2-methyl-	226	C14H10OS	015774-82-0	42



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080320\
Data File : BM026986.D
Acq On : 03 Aug 2020 19:50
Operator : JU/CG
Sample : L3424-05DL 25X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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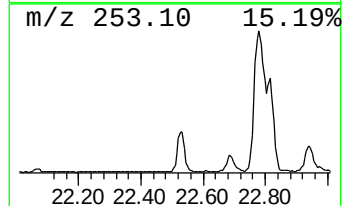
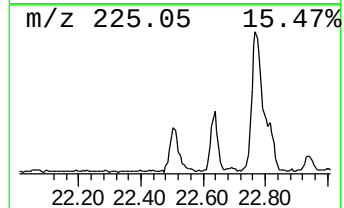
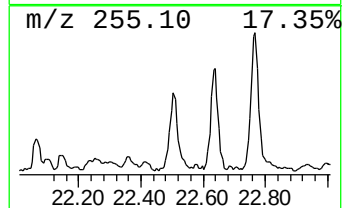
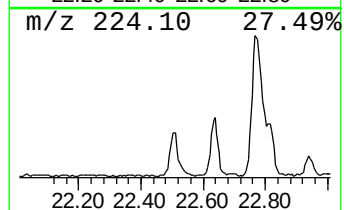
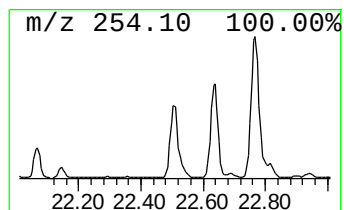
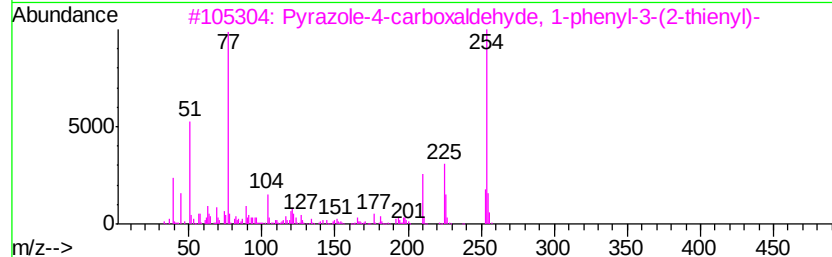
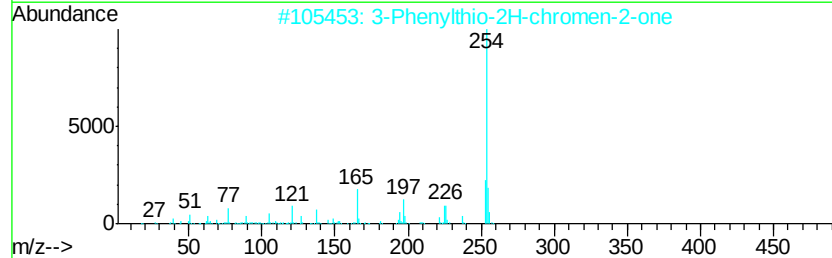
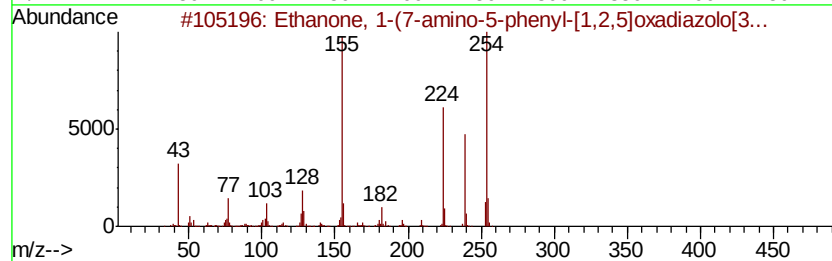
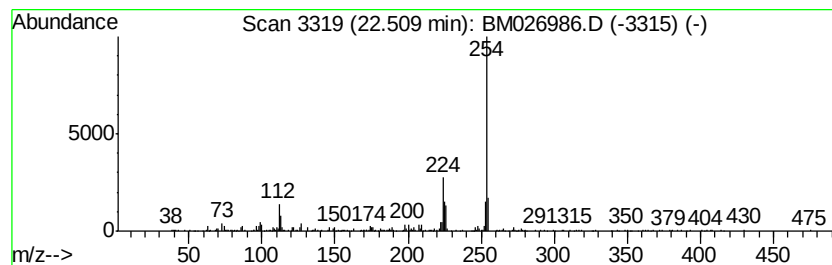
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Ethanone, 1-(7-amino-5-phen... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.51	3.92 ng/ul	260982	Perylene-d12	23.43

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethanone, 1-(7-amino-5-phenyl-[1...	254	C13H10N4O2	1000316-41-1	59
2		3-Phenylthio-2H-chromen-2-one	254	C15H10O2S	072167-95-4	58
3		Pyrazole-4-carboxaldehyde, 1-phe...	254	C14H10N2OS	210825-11-9	53
4		3-Hydroxy-1-methoxyvanthraquinone	254	C15H10O4	028504-24-7	53
5		Phenanthrene, 2-phenyl-	254	C20H14	004325-77-3	52



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080320\
Data File : BM026986.D
Acq On : 03 Aug 2020 19:50
Operator : JU/CG
Sample : L3424-05DL 25X
Misc :
ALS Vial : 13 Sample Multiplier: 1

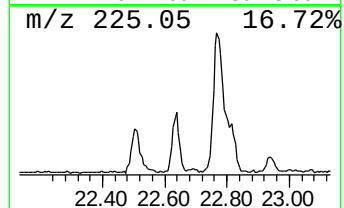
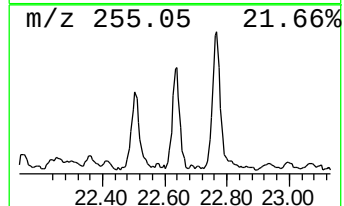
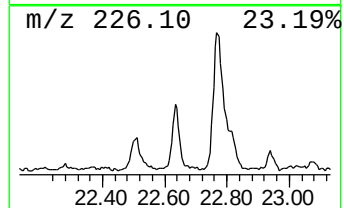
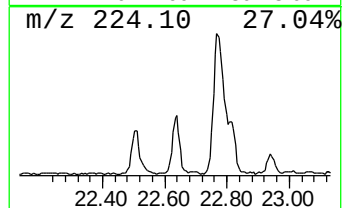
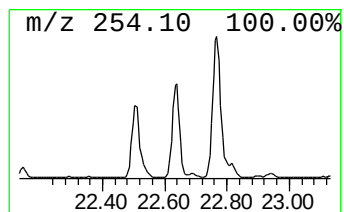
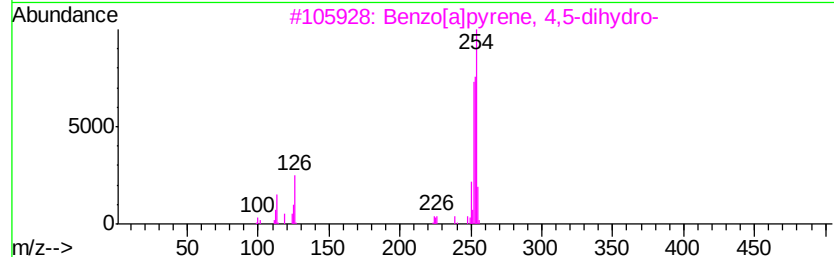
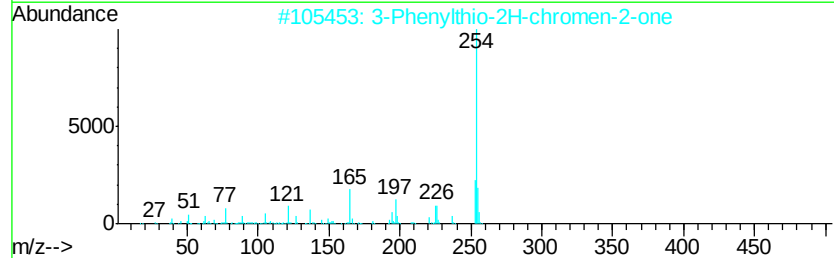
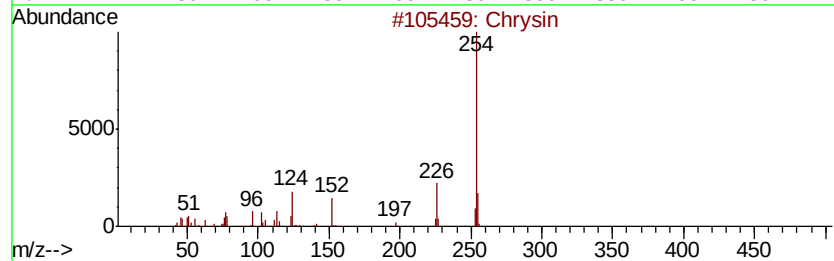
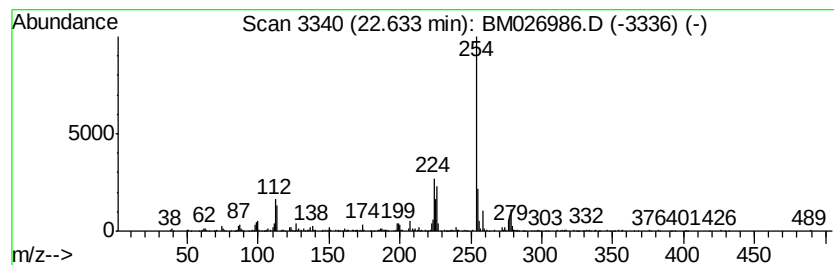
Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM072720MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 5 Chrysin Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.63	5.10 ng/ul	339165	Perylene-d12	23.43
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Chrysin		254 C15H10O4	000480-40-0 50
2	3-Phenylthio-2H-chromen-2-one		254 C15H10O2S	072167-95-4 47
3	Benzo[a]pyrene, 4,5-dihydro-		254 C20H14	057652-66-1 47
4	Chrysin		254 C15H10O4	000480-40-0 47
5	Phenanthrene, 2-phenyl-		254 C20H14	004325-77-3 46



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080320\
Data File : BM026986.D
Acq On : 03 Aug 2020 19:50
Operator : JU/CG
Sample : L3424-05DL 25X
Misc :
ALS Vial : 13 Sample Multiplier: 1

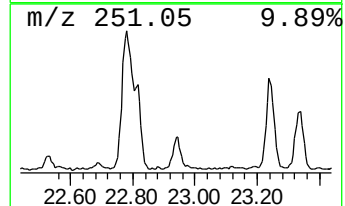
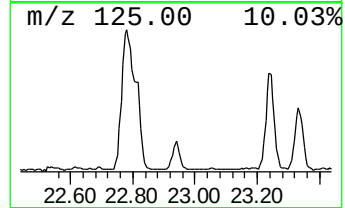
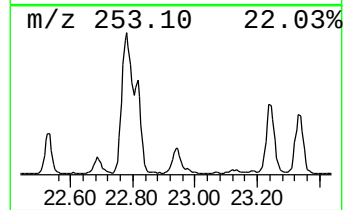
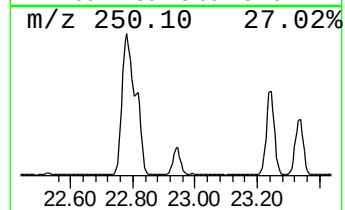
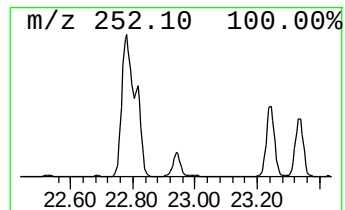
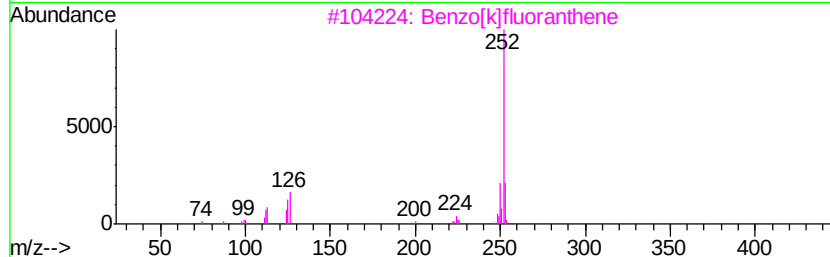
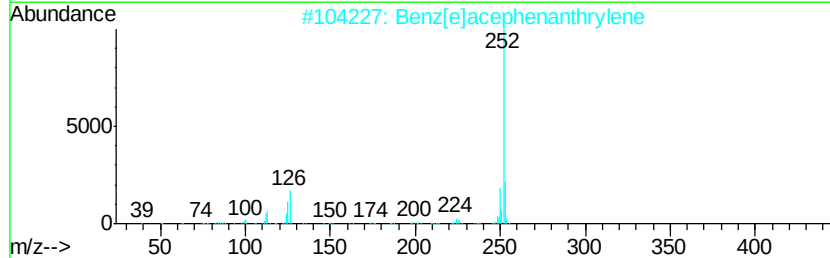
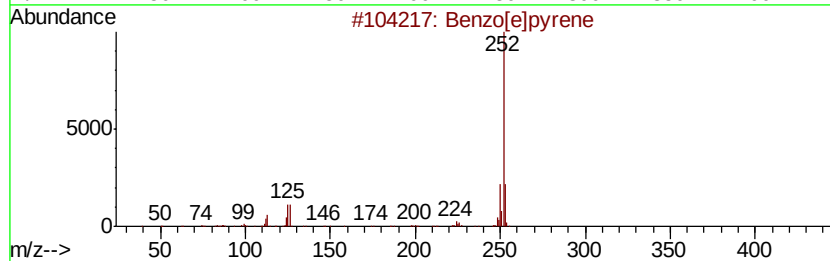
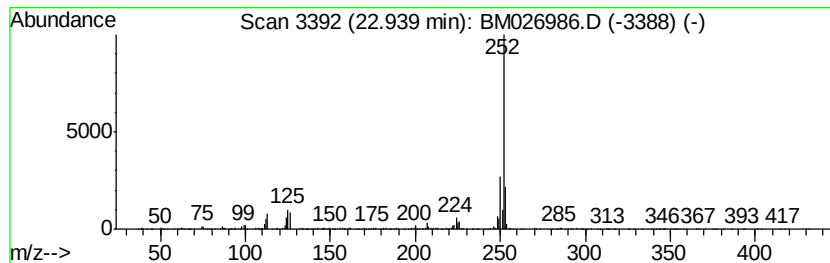
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM072720MA.M
Quant Title : SVOA CALIBRATION

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.
22.94	5.32 ng/ul	354181	Perylene-d12	23.43
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzo[e]pyrene	252 C20H12	000192-97-2 95
2		Benzo[e]acephenanthrylene	252 C20H12	000205-99-2 93
3		Benzo[k]fluoranthene	252 C20H12	000207-08-9 93
4		Perylene	252 C20H12	000198-55-0 93
5		Benzo[a]pyrene	252 C20H12	000050-32-8 91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080320\
Data File : BM026986.D
Acq On : 03 Aug 2020 19:50
Operator : JU/CG
Sample : L3424-05DL 25X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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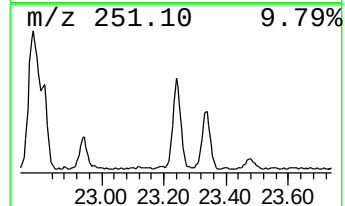
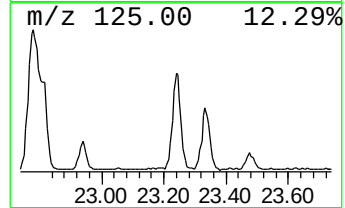
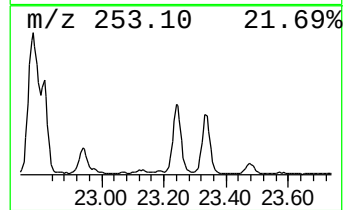
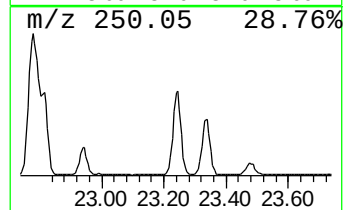
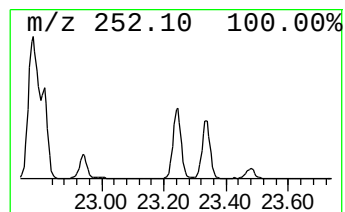
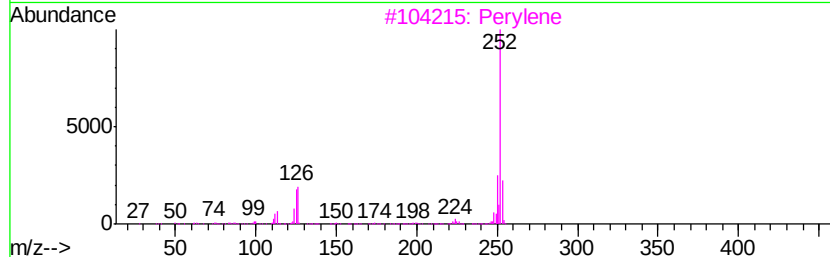
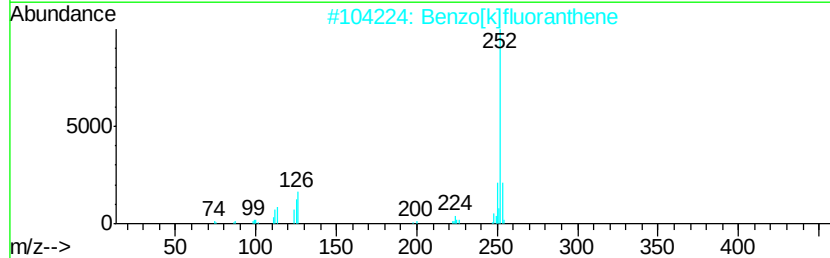
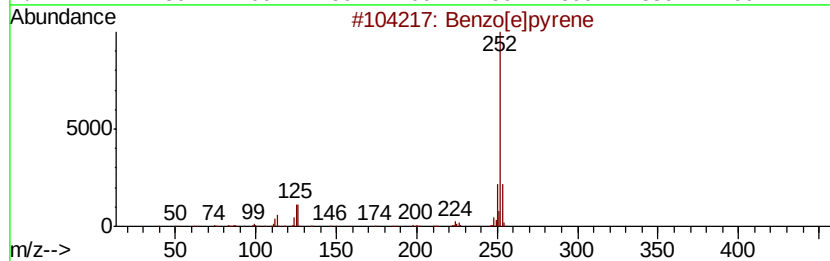
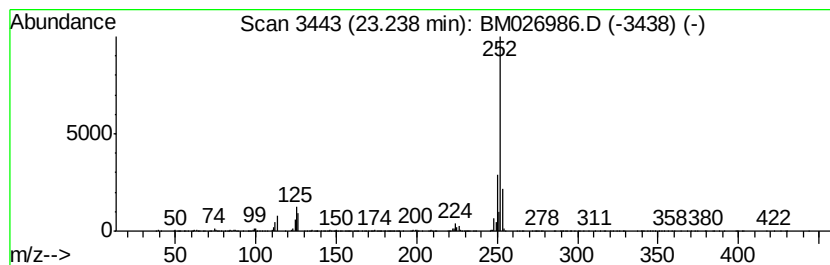
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 unknown-01 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.24	17.70 ng/ul	1177860	Perylene-d12	23.43

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[e]pyrene	252	C20H12	000192-97-2	98
2		Benzo[k]fluoranthene	252	C20H12	000207-08-9	98
3		Perylene	252	C20H12	000198-55-0	94
4		Perylene	252	C20H12	000198-55-0	94
5		Benz[e]acephenanthrylene	252	C20H12	000205-99-2	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM080320\
Data File : BM026986.D
Acq On : 03 Aug 2020 19:50
Operator : JU/CG
Sample : L3424-05DL 25X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM072720MA.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
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Fluoranthene, 2-m...	19.93	2.4	ng/ul	307550	5	21.22	2549240	20.0
Cyclopenta[cd]pyrene	20.94	2.7	ng/ul	347173	5	21.22	2549240	20.0
Ethanone, 1-(7-am...	22.51	3.9	ng/ul	260982	6	23.43	1330640	20.0
Chrysin	22.63	5.1	ng/ul	339165	6	23.43	1330640	20.0
Benzo[e]pyrene	22.94	5.3	ng/ul	354181	6	23.43	1330640	20.0
unknown-01	23.24	17.7	ng/ul	1177860	6	23.43	1330640	20.0