

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN052219\
 Data File : BN005872.D
 Acq On : 23 May 2019 04:07
 Operator : JU/SJ
 Sample : K2927-19
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A42C0

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_N\METHODS\SOM-EPA-BN051819MA.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	6.757	636	641	654	rBV	188218	419169	9.83%	0.869%
2	6.910	662	667	678	rVB	180214	313752	7.36%	0.651%
3	7.104	694	700	712	rBV	245085	445403	10.45%	0.924%
4	7.563	772	778	786	rBV	273717	477069	11.19%	0.989%
5	8.275	894	899	914	rVB	273136	541376	12.70%	1.123%
6	8.710	967	973	980	rBV	276658	491471	11.53%	1.019%
7	9.122	1041	1043	1048	rVB2	8796	9984	0.23%	0.021%
8	9.428	1089	1095	1107	rVB	205015	384913	9.03%	0.798%
9	9.969	1181	1187	1205	rBV	263428	597113	14.01%	1.238%
10	10.322	1240	1247	1253	rBV	434153	754640	17.70%	1.565%
11	10.369	1253	1255	1262	rVB	34729	49114	1.15%	0.102%
12	10.475	1267	1273	1287	rBV	183911	450338	10.56%	0.934%
13	12.010	1529	1534	1545	rVB2	11434	26318	0.62%	0.055%
14	12.228	1564	1571	1579	rVB3	10077	20224	0.47%	0.042%
15	13.028	1702	1707	1717	rVB5	9396	16592	0.39%	0.034%
16	13.234	1736	1742	1752	rVB2	3884	9687	0.23%	0.020%
17	13.522	1786	1791	1796	rBV2	15823	25503	0.60%	0.053%
18	13.575	1796	1800	1802	rVV2	7803	11691	0.27%	0.024%
19	13.616	1802	1807	1812	rVV	814066	1132416	26.57%	2.349%
20	13.663	1812	1815	1824	rVB	286944	421180	9.88%	0.874%
21	13.757	1829	1831	1835	rVV2	8912	11182	0.26%	0.023%
22	13.887	1847	1853	1865	rBV	863689	1395067	32.73%	2.893%
23	14.110	1886	1891	1895	rBV	14884	20454	0.48%	0.042%
24	14.198	1899	1906	1913	rBV2	758199	1168884	27.42%	2.424%
25	14.263	1913	1917	1923	rVB	79886	115151	2.70%	0.239%
26	14.492	1951	1956	1966	rBV6	50314	163801	3.84%	0.340%
27	14.610	1972	1976	1983	rVB	52812	78456	1.84%	0.163%
28	14.675	1983	1987	1988	rBV4	7423	7850	0.18%	0.016%
29	14.834	2007	2014	2019	rVV3	8692	16601	0.39%	0.034%
30	14.886	2019	2023	2026	rVB2	8637	9936	0.23%	0.021%
31	15.004	2035	2043	2047	rBV6	11730	24157	0.57%	0.050%
32	15.075	2050	2055	2060	rVV4	20385	27288	0.64%	0.057%
33	15.204	2070	2077	2083	rVV	1188465	1730356	40.59%	3.589%
34	15.263	2083	2087	2099	rVB	93298	148758	3.49%	0.309%

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35	15.386	2099	2108	2112	rBV3	52010	104421	2.45%	0.217%
36	15.428	2112	2115	2121	rVV7	21116	49743	1.17%	0.103%
37	15.475	2121	2123	2126	rVV4	13753	18034	0.42%	0.037%
38	15.516	2126	2130	2133	rVV4	9936	17181	0.40%	0.036%
39	15.557	2133	2137	2145	rVV	28167	52897	1.24%	0.110%
40	15.710	2159	2163	2170	rVV	32230	64033	1.50%	0.133%
41	15.798	2171	2178	2182	rVB6	8859	19945	0.47%	0.041%
42	15.939	2197	2202	2209	rBV3	45593	80790	1.90%	0.168%
43	15.992	2209	2211	2214	rVB3	8975	8179	0.19%	0.017%
44	16.269	2249	2258	2263	rBV5	27321	58519	1.37%	0.121%
45	16.316	2263	2266	2272	rVV5	13593	21276	0.50%	0.044%
46	16.416	2277	2283	2287	rVB7	15101	29206	0.69%	0.061%
47	16.469	2289	2292	2294	rBV2	6526	7871	0.18%	0.016%
48	16.504	2294	2298	2301	rBV3	17608	26050	0.61%	0.054%
49	16.539	2301	2304	2307	rVB3	20352	24459	0.57%	0.051%
50	16.628	2314	2319	2325	rVB2	62201	92168	2.16%	0.191%
51	16.710	2326	2333	2335	rBV5	30810	68457	1.61%	0.142%
52	16.780	2340	2345	2356	rVB	85232	138451	3.25%	0.287%
53	16.963	2369	2376	2379	rBV2	1105939	1608205	37.73%	3.335%
54	17.004	2379	2383	2388	rVV	1213786	1655802	38.84%	3.434%
55	17.063	2388	2393	2397	rVV	1354078	2064845	48.44%	4.283%
56	17.098	2397	2399	2405	rVB	267852	293956	6.90%	0.610%
57	17.157	2405	2409	2410	rBV4	18141	23062	0.54%	0.048%
58	17.380	2440	2447	2456	rBV2	83106	201061	4.72%	0.417%
59	17.475	2459	2463	2468	rVB4	46103	64486	1.51%	0.134%
60	17.551	2471	2476	2478	rBV	25197	40826	0.96%	0.085%
61	17.686	2497	2499	2504	rVB2	23122	31285	0.73%	0.065%
62	17.839	2520	2525	2528	rBV	177419	238607	5.60%	0.495%
63	17.886	2528	2533	2541	rVV2	289921	542324	12.72%	1.125%
64	17.975	2544	2548	2552	rVV	92051	134474	3.15%	0.279%
65	18.033	2553	2558	2562	rVV2	226783	420088	9.85%	0.871%
66	18.075	2562	2565	2570	rVB	123230	165979	3.89%	0.344%
67	18.163	2576	2580	2584	rBV4	49305	84613	1.98%	0.175%
68	18.239	2591	2593	2598	rBV5	19010	25038	0.59%	0.052%
69	18.351	2607	2612	2617	rBV	134007	202927	4.76%	0.421%
70	18.410	2617	2622	2628	rVB	133938	208198	4.88%	0.432%
71	18.598	2651	2654	2658	rVB	52881	55494	1.30%	0.115%

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72	18.692	2667	2670	2674	rVB	49476	62646	1.47%	0.130%
73	18.774	2680	2684	2688	rBV	134873	202075	4.74%	0.419%
74	18.816	2688	2691	2692	rBV2	49065	53774	1.26%	0.112%
75	18.927	2705	2710	2715	rVB4	91796	171725	4.03%	0.356%
76	19.039	2724	2729	2737	rVB	1795392	2366283	55.51%	4.908%
77	19.380	2781	2787	2789	rBV	2043810	2963079	69.51%	6.146%
78	19.410	2789	2792	2801	rVV	1705333	2290810	53.74%	4.751%
79	19.486	2802	2805	2812	rVB2	138578	212571	4.99%	0.441%
80	19.592	2821	2823	2829	rVB	106592	135508	3.18%	0.281%
81	19.739	2844	2848	2849	rBV2	135753	189052	4.43%	0.392%
82	19.916	2875	2878	2881	rVB	217847	264612	6.21%	0.549%
83	20.016	2892	2895	2899	rBV	125019	156308	3.67%	0.324%
84	20.069	2901	2904	2910	rVB2	209895	330334	7.75%	0.685%
85	20.216	2926	2929	2932	rVB	144016	164065	3.85%	0.340%
86	20.263	2934	2937	2942	rVB2	146745	192676	4.52%	0.400%
87	20.680	3005	3008	3014	rVB	229683	275055	6.45%	0.570%
88	20.839	3032	3035	3038	rBV	221200	266669	6.26%	0.553%
89	20.874	3038	3041	3044	rVB	233850	258573	6.07%	0.536%
90	20.916	3044	3048	3057	rBV3	451992	790823	18.55%	1.640%
91	21.198	3089	3096	3100	rBV2	2034903	4262758	100.00%	8.841%
92	21.233	3100	3102	3107	rVB	1218500	1402590	32.90%	2.909%
93	21.357	3120	3123	3127	rVB3	230832	268407	6.30%	0.557%
94	21.792	3194	3197	3206	rVB2	295188	531159	12.46%	1.102%
95	22.245	3270	3274	3280	rBV3	848328	1234637	28.96%	2.561%
96	22.751	3354	3360	3366	rBV2	1005453	2768038	64.94%	5.741%
97	23.221	3435	3440	3443	rBV	518224	864201	20.27%	1.792%
98	23.274	3444	3449	3453	rVV	1212894	2377722	55.78%	4.931%
99	23.315	3453	3456	3465	rVB	724157	1185512	27.81%	2.459%
100	23.410	3466	3472	3477	rBV	838121	1514728	35.53%	3.142%

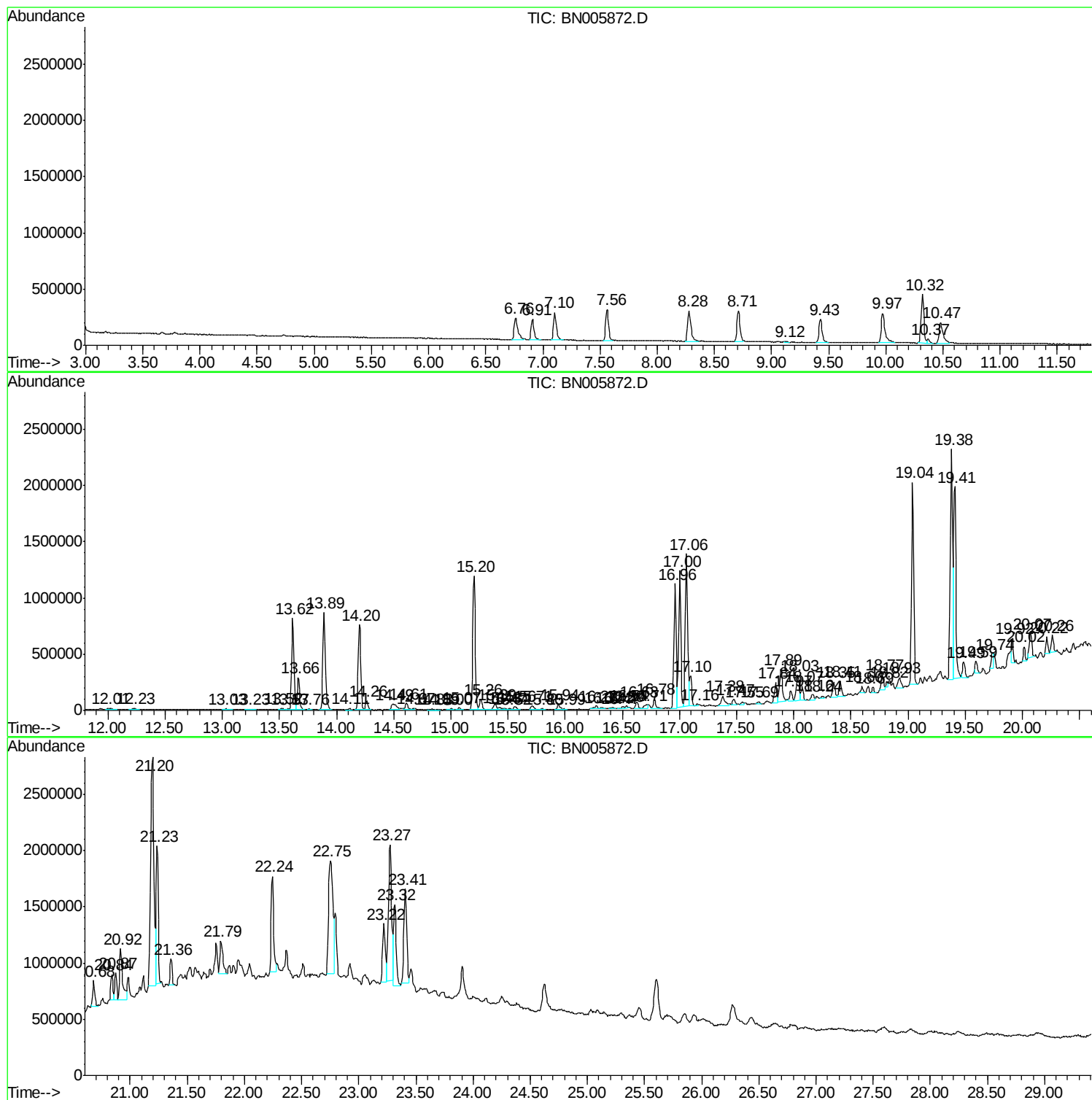
Sum of corrected areas: 48215234

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN051819MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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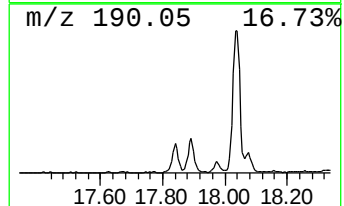
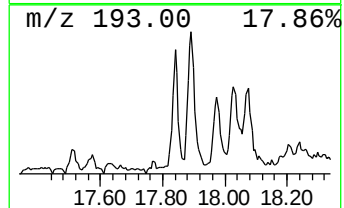
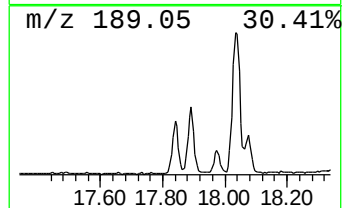
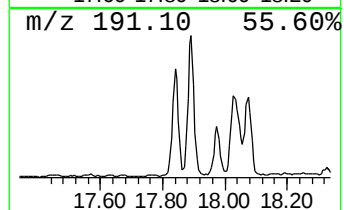
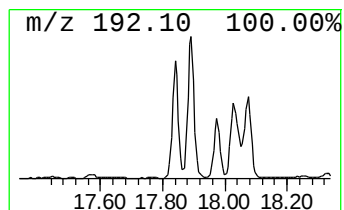
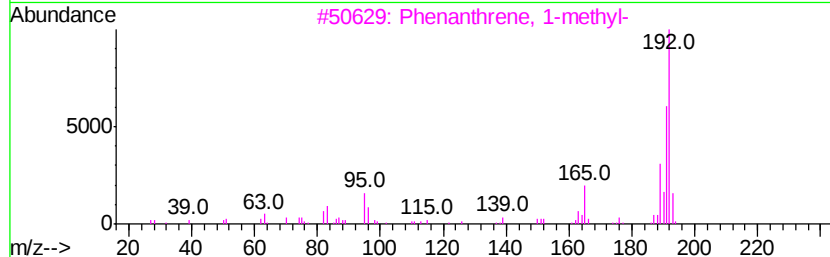
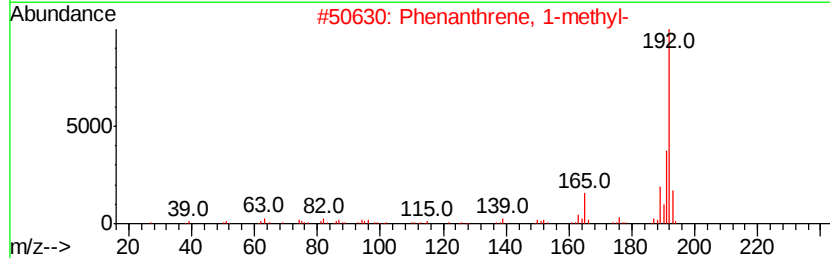
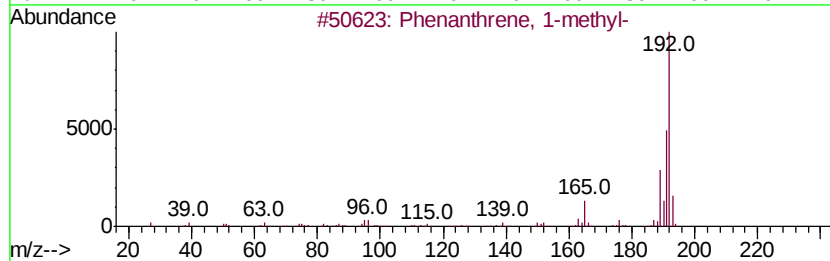
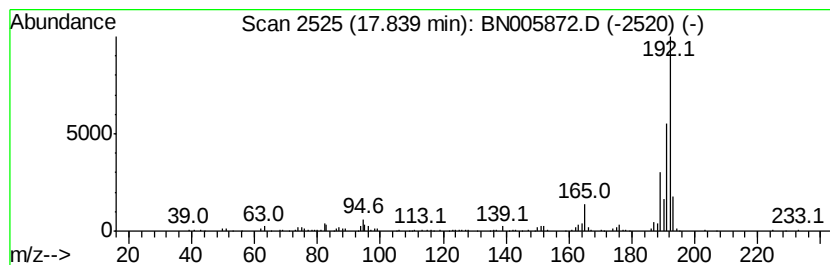
Instrument :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN051819MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.		
17.84	2.97 ng/ul	238607	Phenanthrene-d10	16.96		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95	
2	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95	
3	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	93	
4	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93	
5	Anthracene, 1-methyl-	192	C15H12	000610-48-0	91	



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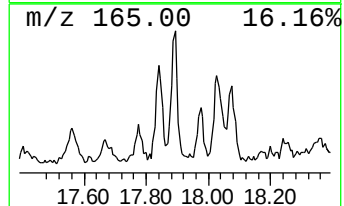
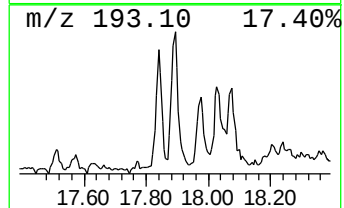
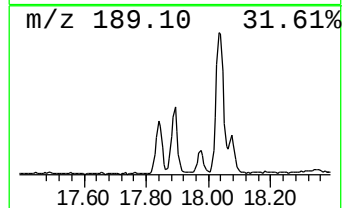
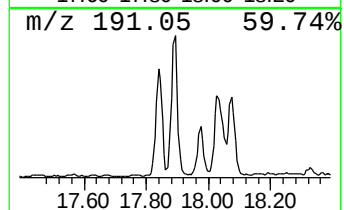
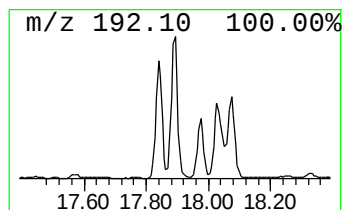
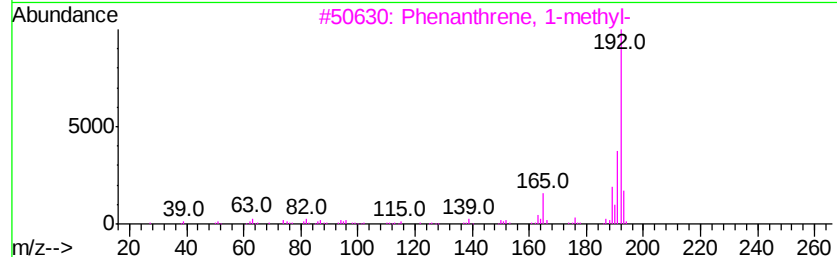
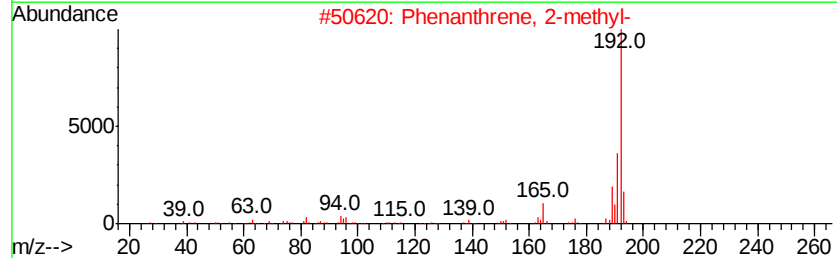
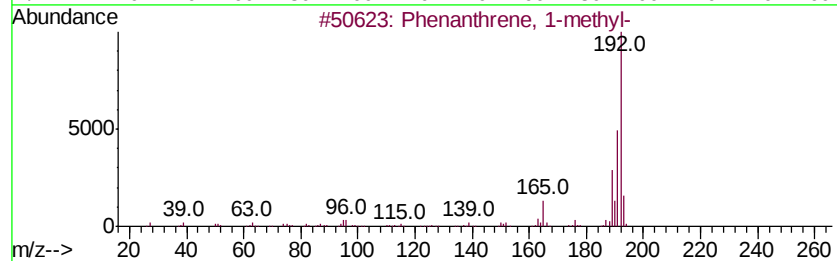
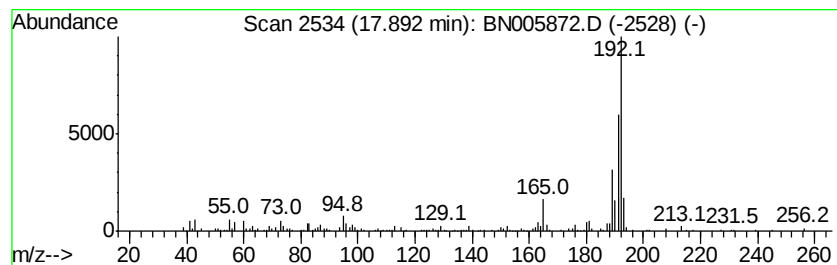
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.		
17.89	6.74 ng/ul	542324	Phenanthrene-d10	16.96		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	97	
2	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96	
3	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95	
4	Anthracene, 2-methyl-	192	C15H12	000613-12-7	94	
5	1H-Cyclopropa[1]phenanthrene, 1a, ...	192	C15H12	000949-41-7	93	



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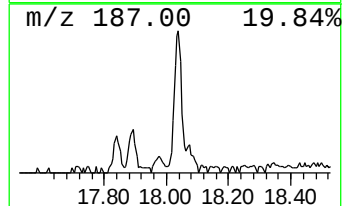
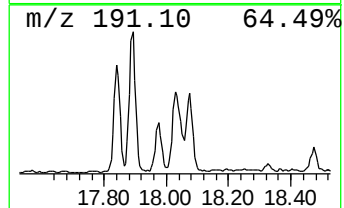
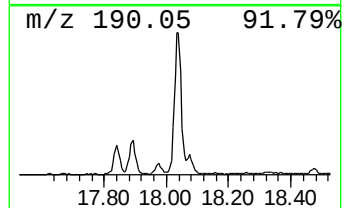
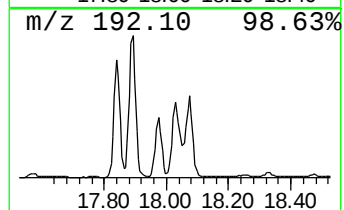
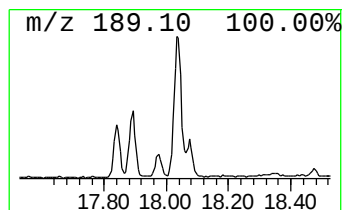
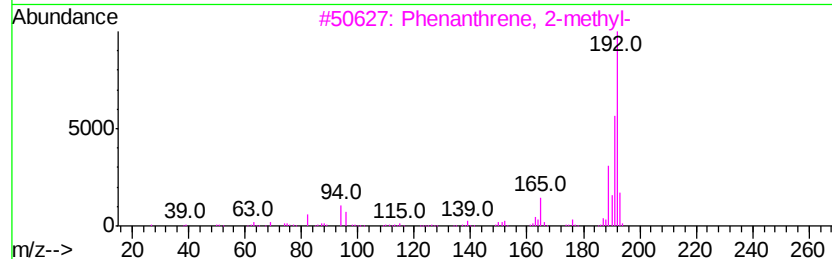
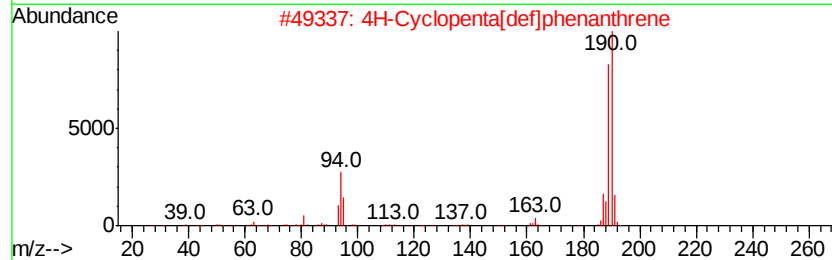
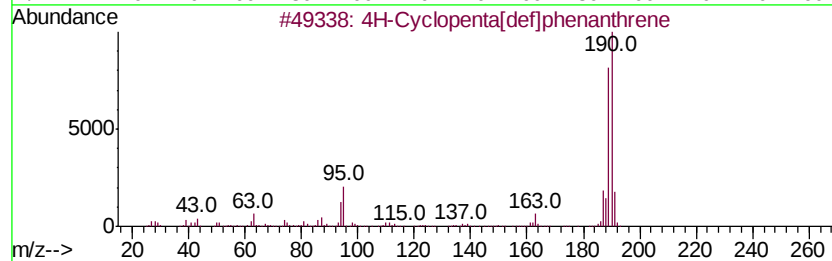
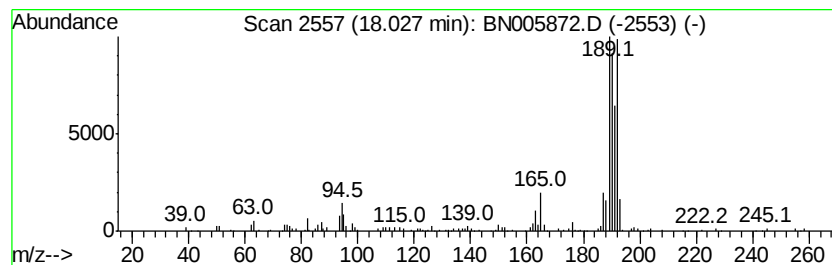
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 4H-Cyclopenta[def]phenanthrene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.03	5.22 ng/ul	420088	Phenanthrene-d10	16.96

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	64
2		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	58
3		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	42
4		Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	38
5		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	35



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Acq On : 23 May 2019 04:07
Operator : JU/SJ
Sample : K2927-19
Misc :
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Instrument :
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ClientSampleId :
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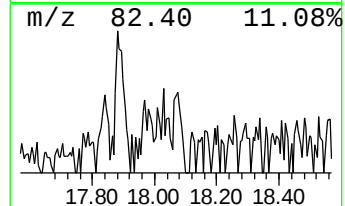
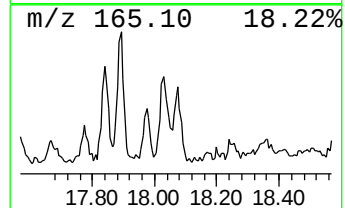
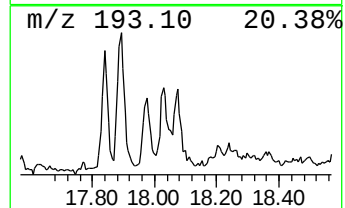
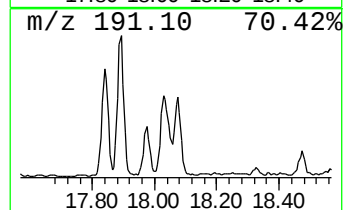
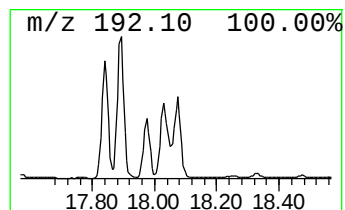
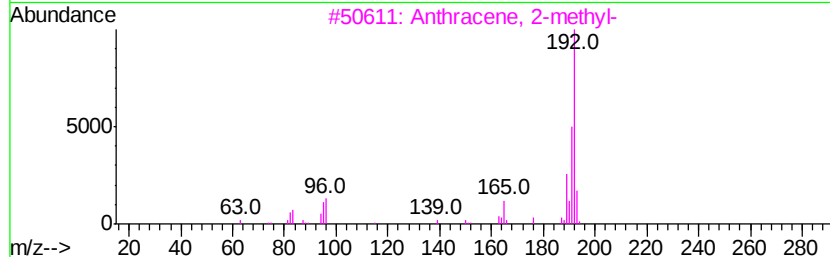
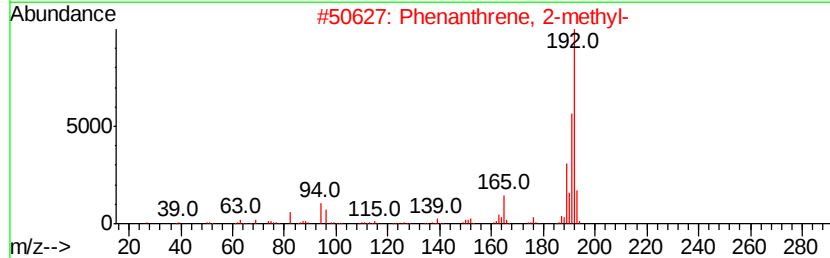
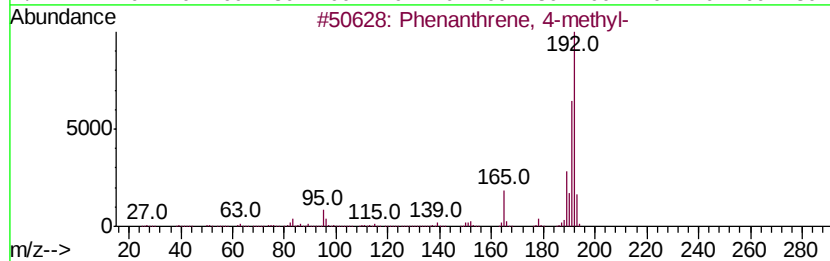
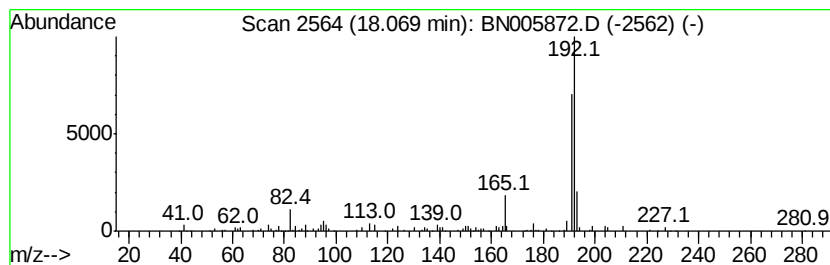
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Phenanthrene, 4-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.07	2.06 ng/ul	165979	Phenanthrene-d10	16.96

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	91
2		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	90
3		Anthracene, 2-methyl-	192	C15H12	000613-12-7	90
4		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	87
5		Propyne, 1,3-diphenyl-	192	C15H12	004980-70-5	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN052219\
Data File : BN005872.D
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Sample : K2927-19
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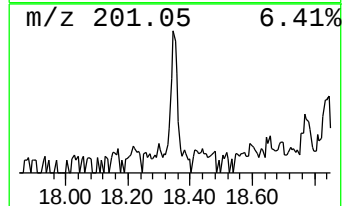
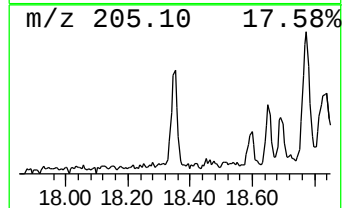
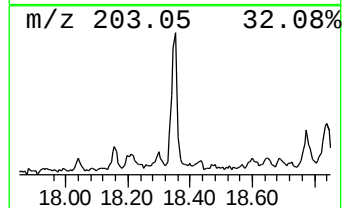
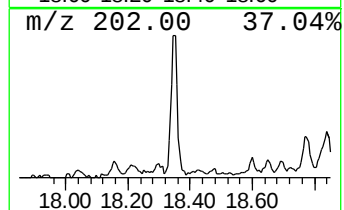
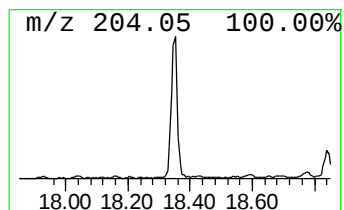
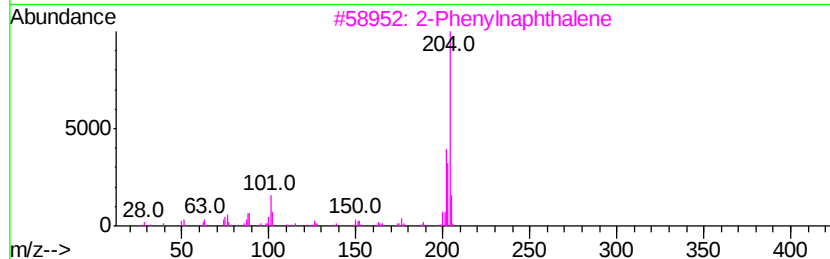
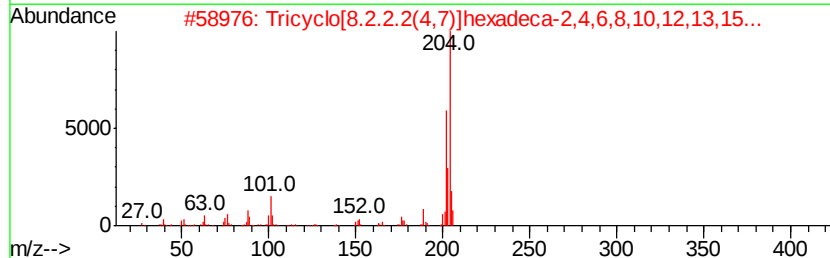
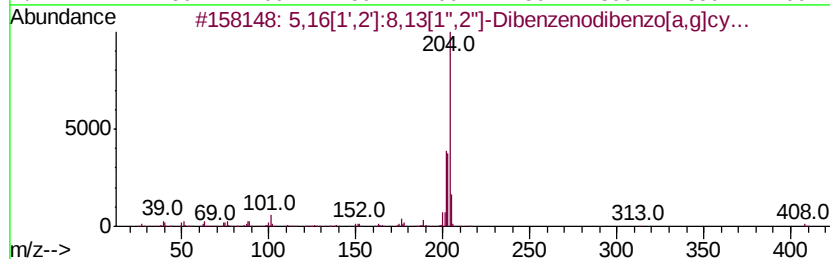
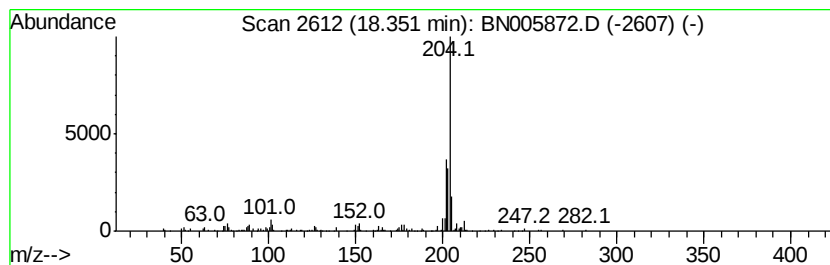
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 5,16[1',2']:8,13[1'',2'']-D... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.35	2.52 ng/ul	202927	Phenanthrene-d10	16.96

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	5,16	[1',2']	:8,13[1'',2'']-Dibenz...	408	C32H24	005672-97-9	86
2	Tricyclo	[8.2.2.2(4,7)]	hexadeca-2...	204	C16H12	006572-60-7	76
3	2-Phenyl	naphthalene		204	C16H12	035465-71-5	76
4	Naphthalene,	2-phenyl-		204	C16H12	000612-94-2	68
5	Naphthalene,	2-phenyl-		204	C16H12	000612-94-2	68



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN052219\
Data File : BN005872.D
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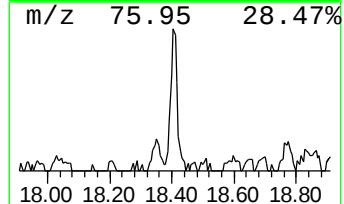
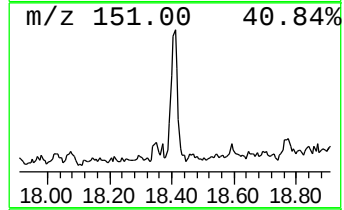
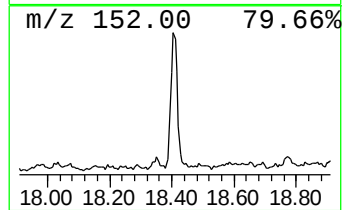
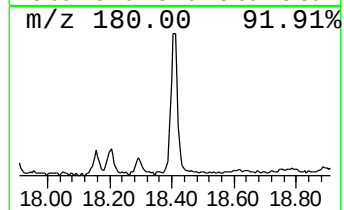
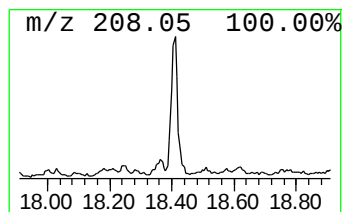
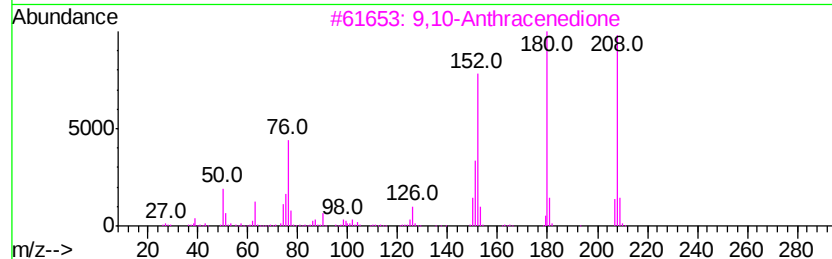
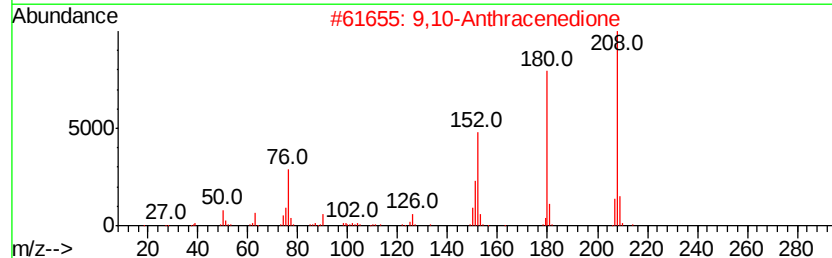
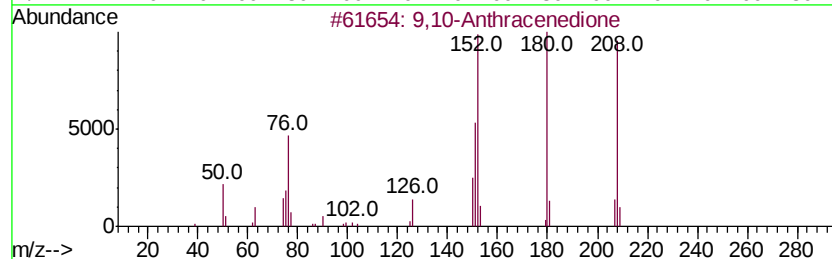
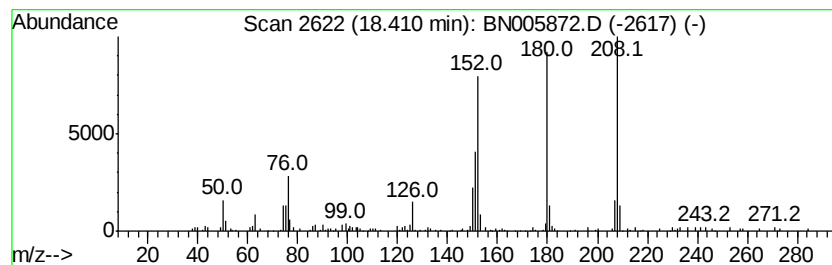
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 9,10-Anthracenedione Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.41	2.59 ng/ul	208198	Phenanthrene-d10	16.96

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9,10-Anthracenedione	208	C14H8O2	000084-65-1	98
2		9,10-Anthracenedione	208	C14H8O2	000084-65-1	94
3		9,10-Anthracenedione	208	C14H8O2	000084-65-1	87
4		1H-Phenalen-1-one	180	C13H8O	000548-39-0	78
5		9H-Fluoren-9-one	180	C13H8O	000486-25-9	70



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN052219\
Data File : BN005872.D
Acq On : 23 May 2019 04:07
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Sample : K2927-19
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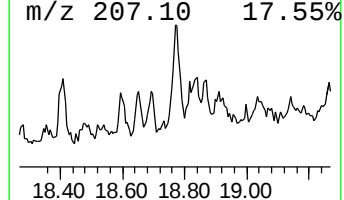
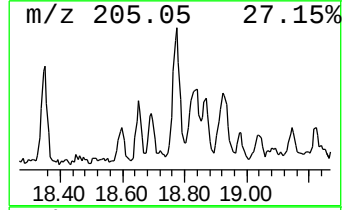
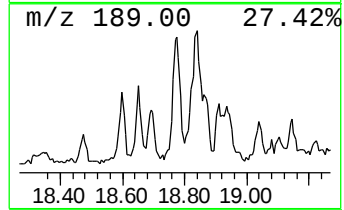
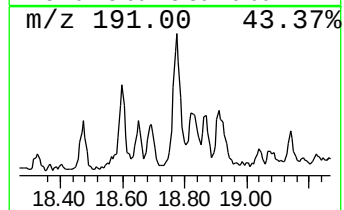
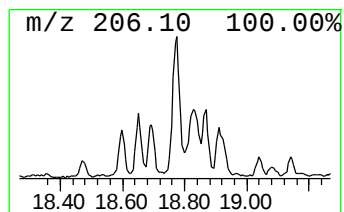
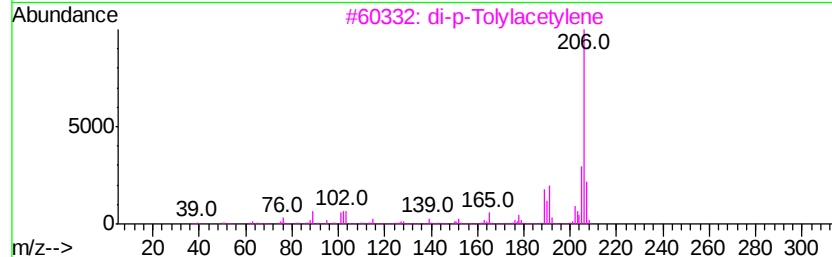
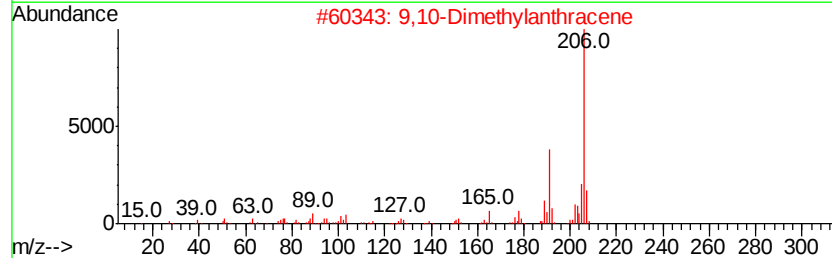
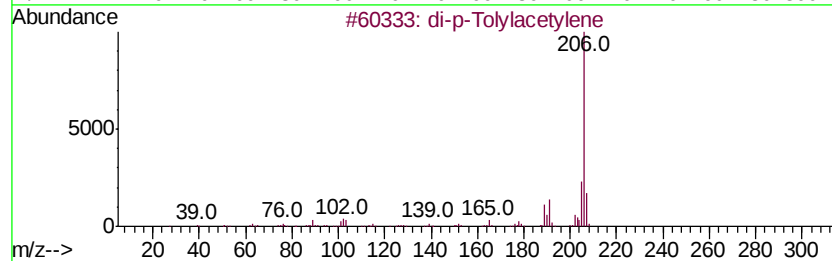
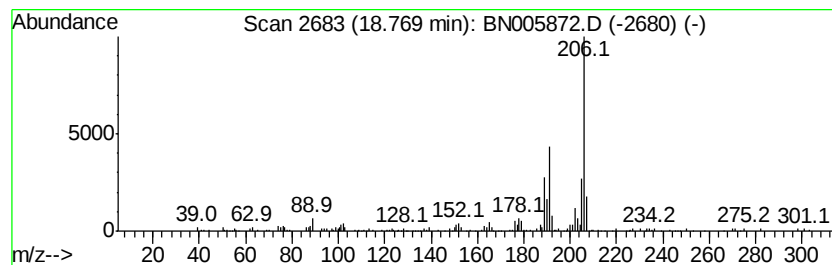
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 di-p-Tolylacetylene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.77	2.51 ng/ul	202075	Phenanthrene-d10	16.96

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	di-p-Tolylacetylene	206	C16H14	002789-88-0	93
2		9,10-Dimethylantracene	206	C16H14	000781-43-1	91
3		di-p-Tolylacetylene	206	C16H14	002789-88-0	91
4		Naphthalene, 1,2-dihydro-4-phenyl-	206	C16H14	007469-40-1	87
5		9,10-Dimethylantracene	206	C16H14	000781-43-1	83



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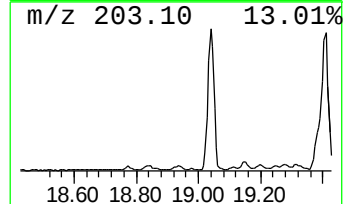
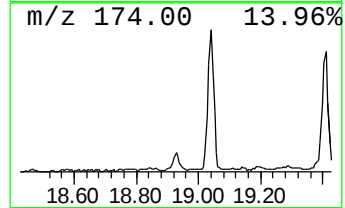
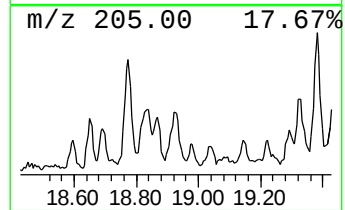
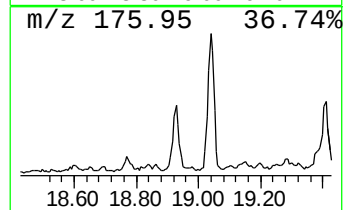
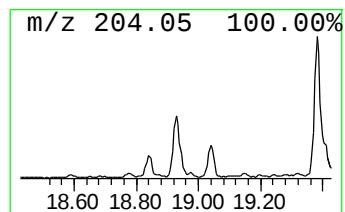
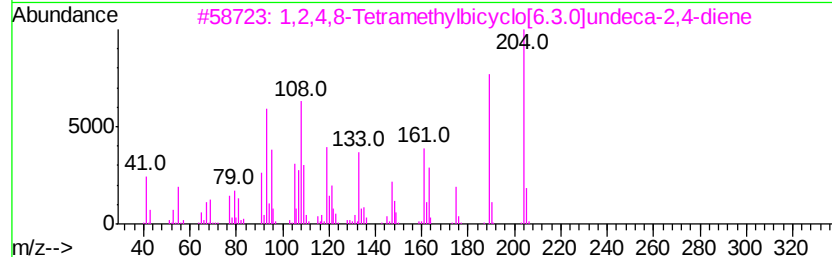
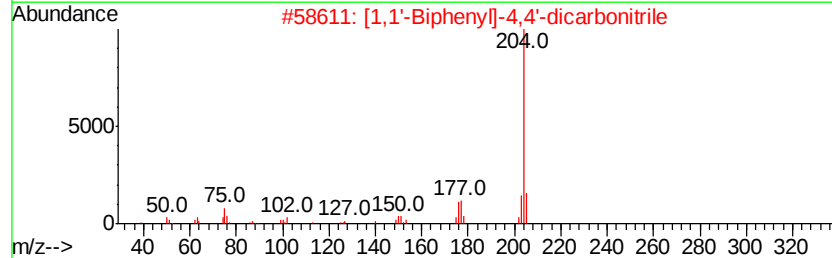
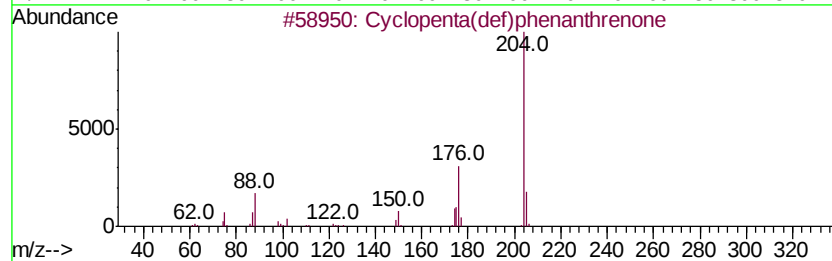
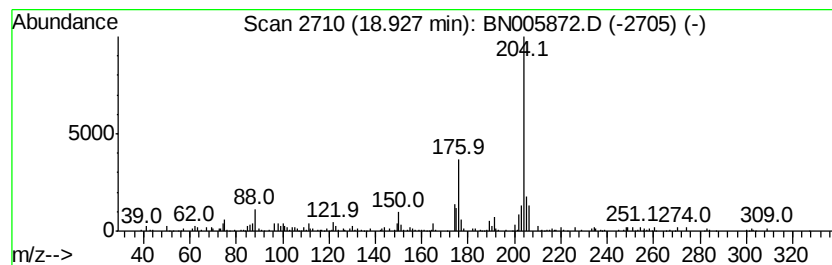
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Cyclopenta(def)phenanthrenone Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD		R.T.		
18.93	2.14 ng/ul	171725	Phenanthrene-d10		16.96		
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cvclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	87
2			[1,1'-Biphenyl]-4,4'-dicarbonitrile	204	C14H8N2	001591-30-6	59
3			1,2,4,8-Tetramethylbicyclo[6.3.0...	204	C15H24	137235-51-9	55
4			[1,1'-Biphenyl]-2,2'-dicarbonitrile	204	C14H8N2	004341-02-0	53
5			[1,1'-Biphenyl]-4,4'-dicarbonitrile	204	C14H8N2	001591-30-6	50



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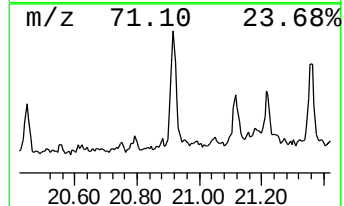
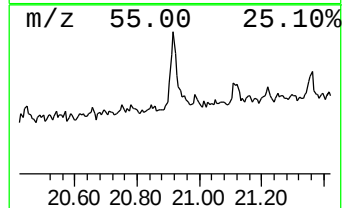
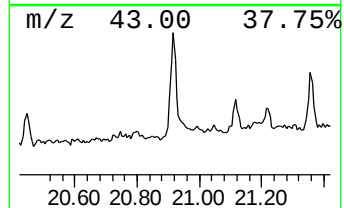
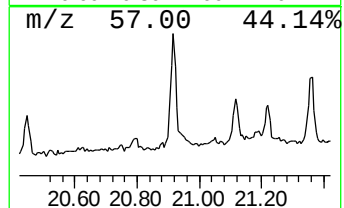
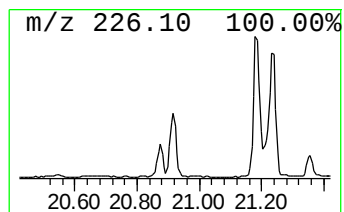
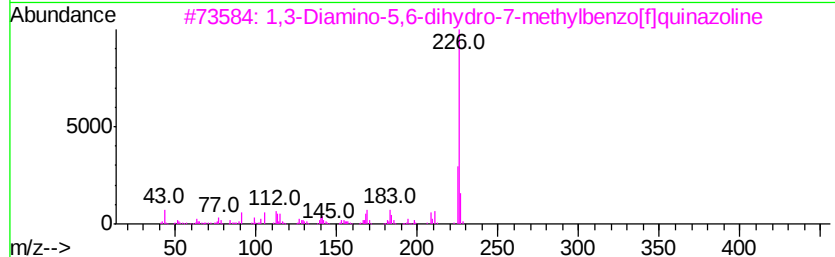
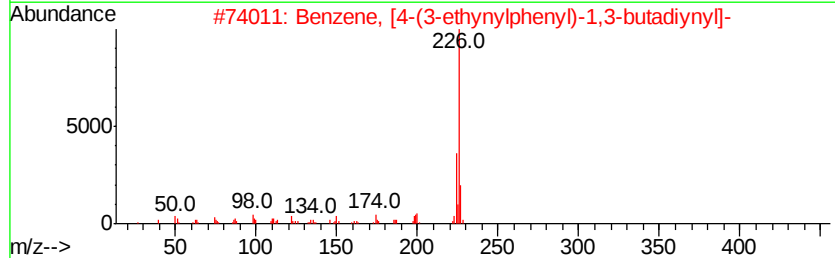
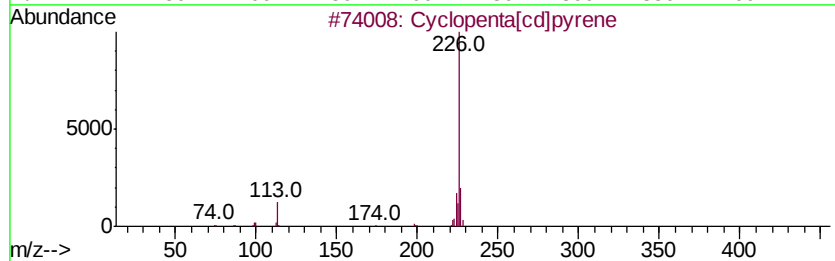
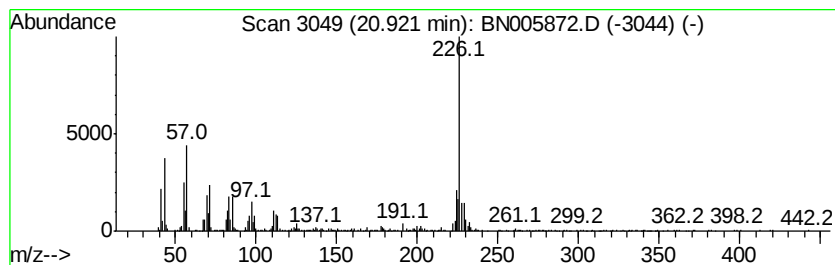
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Cyclopenta[cd]pyrene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.92	3.71 ng/ul	790823	Chrysene-d12	21.20

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopenta[cd]pyrene	226	C18H10	027208-37-3	62
2		Benzene, [4-(3-ethynylphenyl)-1,...	226	C18H10	1000115-87-3	43
3		1,3-Diamino-5,6-dihydro-7-methyl...	226	C13H14N4	037436-37-6	43
4		.beta.-Carboline, 7-methoxy-1,2-...	226	C14H14N2O	006519-18-2	43
5		p-Anisaldehyde phenylhydrazone	226	C14H14N2O	000622-73-1	43



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN052219\
Data File : BN005872.D
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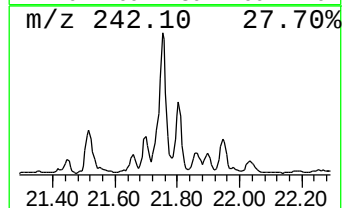
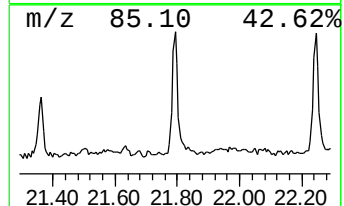
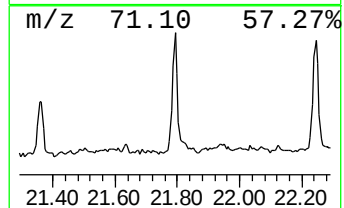
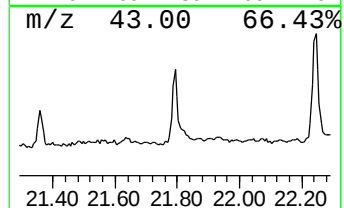
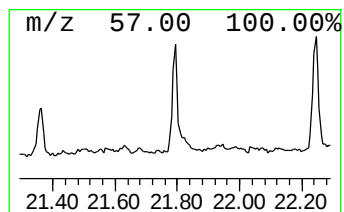
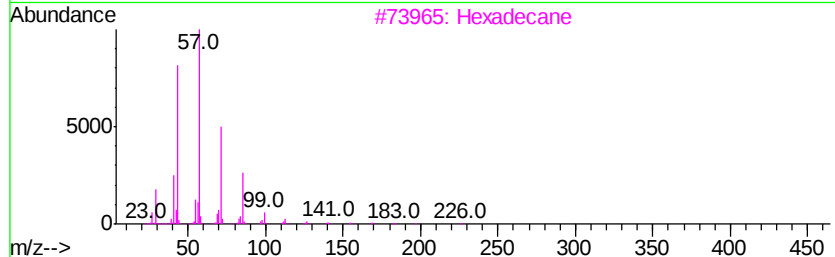
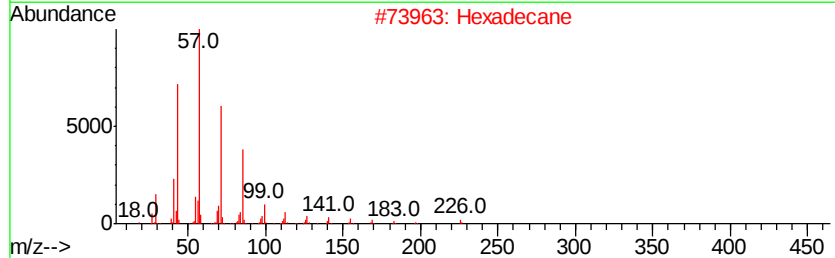
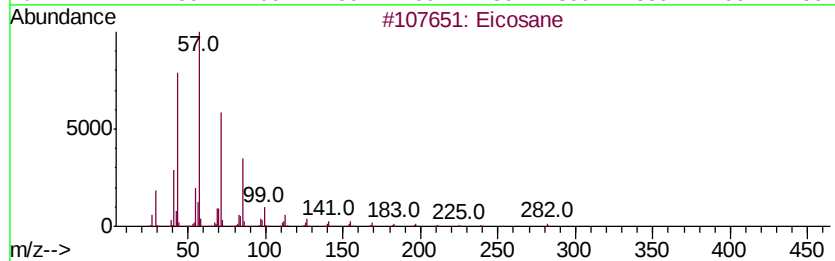
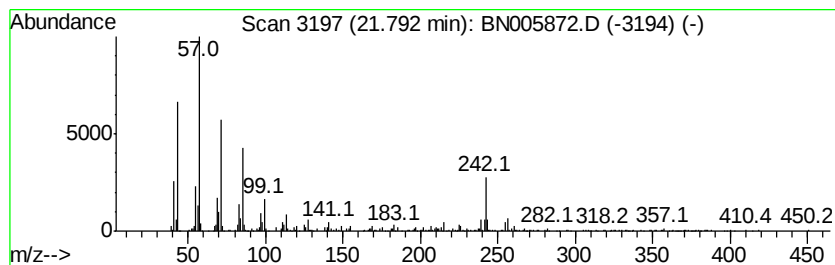
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.79	2.49 ng/ul	531159	Chrysene-d12	21.20

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane			282	C20H42	000112-95-8	96
2	Hexadecane			226	C16H34	000544-76-3	95
3	Hexadecane			226	C16H34	000544-76-3	93
4	Eicosane			282	C20H42	000112-95-8	93
5	Pentadecane			212	C15H32	000629-62-9	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN052219\
Data File : BN005872.D
Acq On : 23 May 2019 04:07
Operator : JU/SJ
Sample : K2927-19
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
A42C0

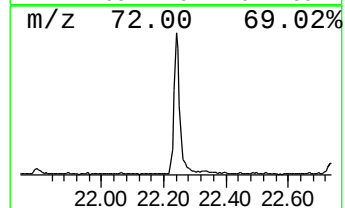
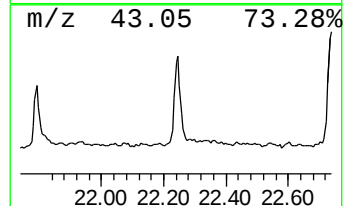
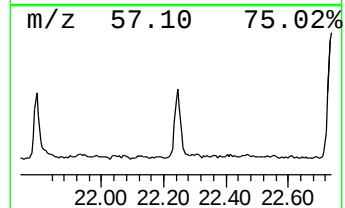
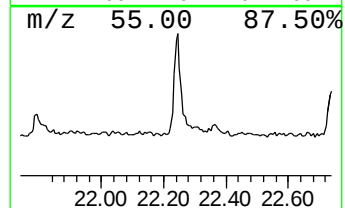
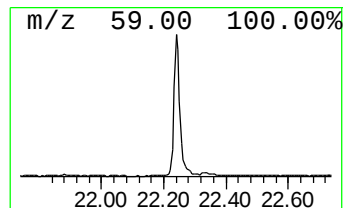
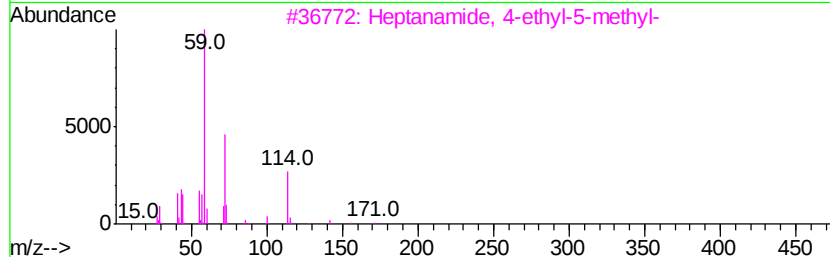
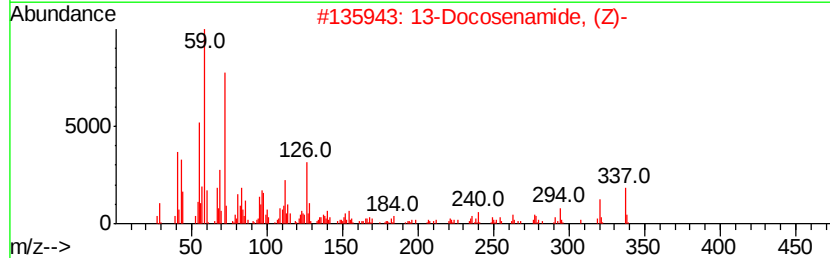
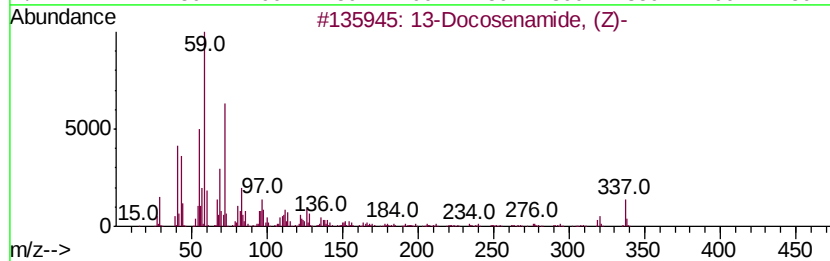
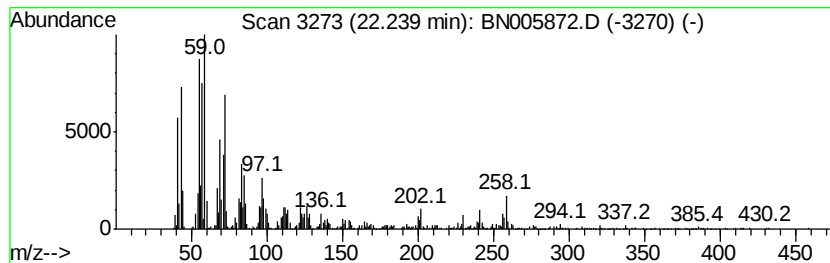
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 13-Docosenamide, (Z)- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.24	5.79 ng/ul	1234640	Chrysene-d12	21.20

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	70
2			13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	50
3			Heptanamide, 4-ethyl-5-methyl-	171	C10H21NO	054789-40-1	43
4			13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	38
5			Ethyl dodecyl ether	214	C14H30O	007289-37-4	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN052219\
Data File : BN005872.D
Acq On : 23 May 2019 04:07
Operator : JU/SJ
Sample : K2927-19
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
A42C0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN051819MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Phenanthrene, 1-m...	17.84	3.0	ng/ul	238607	4	16.96	1608210	20.0
Phenanthrene, 2-m...	17.89	6.7	ng/ul	542324	4	16.96	1608210	20.0
4H-Cyclopenta[def...	18.03	5.2	ng/ul	420088	4	16.96	1608210	20.0
Phenanthrene, 4-m...	18.07	2.1	ng/ul	165979	4	16.96	1608210	20.0
5,16[1',2']:8,13[...	18.35	2.5	ng/ul	202927	4	16.96	1608210	20.0
9,10-Anthracenedione	18.41	2.6	ng/ul	208198	4	16.96	1608210	20.0
di-p-Tolylacetylene	18.77	2.5	ng/ul	202075	4	16.96	1608210	20.0
Cyclopenta(def)ph...	18.93	2.1	ng/ul	171725	4	16.96	1608210	20.0
Cyclopenta[cd]pyrene	20.92	3.7	ng/ul	790823	5	21.20	4262760	20.0
(DEL) Alkane: Str...	21.79	2.5	ng/ul	531159	5	21.20	4262760	20.0
13-Docosenamide, ...	22.24	5.8	ng/ul	1234640	5	21.20	4262760	20.0