

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN051119\
 Data File : BN005624.D
 Acq On : 11 May 2019 13:42
 Operator : JU/SJ
 Sample : K2603-06MEDL 30X
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 GAJ58MEDL

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-BN041919MA.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	7.581	775	781	795	rBV	734256	1269743	29.12%	2.958%
2	10.340	1243	1250	1269	rBV	1020640	1925343	44.16%	4.485%
3	12.246	1565	1574	1584	rBV2	49309	100878	2.31%	0.235%
4	13.045	1705	1710	1721	rBV	83549	141331	3.24%	0.329%
5	13.393	1760	1769	1776	rBV4	94657	252333	5.79%	0.588%
6	13.540	1785	1794	1798	rBV	205466	363839	8.35%	0.848%
7	13.593	1798	1803	1807	rBV	110819	146500	3.36%	0.341%
8	13.640	1807	1811	1814	rBV	43558	62358	1.43%	0.145%
9	13.675	1814	1817	1828	rVB3	75228	130307	2.99%	0.304%
10	13.775	1828	1834	1838	rBV2	113588	196010	4.50%	0.457%
11	13.934	1849	1861	1870	rBV2	546711	954599	21.90%	2.224%
12	14.216	1900	1909	1916	rBV2	1782994	2846245	65.28%	6.630%
13	14.281	1916	1920	1923	rVV2	78467	117171	2.69%	0.273%
14	14.316	1923	1926	1931	rVB2	54423	69597	1.60%	0.162%
15	14.440	1941	1947	1955	rVB3	28157	64989	1.49%	0.151%
16	14.628	1968	1979	1986	rBV	133292	251334	5.76%	0.585%
17	14.704	1986	1992	1997	rBV	87113	116688	2.68%	0.272%
18	14.840	2008	2015	2021	rBV3	93587	199500	4.58%	0.465%
19	14.898	2021	2025	2029	rVB	64122	83854	1.92%	0.195%
20	14.998	2035	2042	2043	rBV2	39948	57143	1.31%	0.133%
21	15.098	2054	2059	2064	rVB	149639	211967	4.86%	0.494%
22	15.175	2066	2072	2075	rBV6	35095	78468	1.80%	0.183%
23	15.216	2075	2079	2084	rVB3	212365	309453	7.10%	0.721%
24	15.275	2084	2089	2102	rVB	663946	1086105	24.91%	2.530%
25	15.398	2107	2110	2114	rBV4	42073	54850	1.26%	0.128%
26	15.498	2122	2127	2134	rBV6	32677	78502	1.80%	0.183%
27	15.575	2135	2140	2148	rVB2	75319	129682	2.97%	0.302%
28	15.728	2158	2166	2173	rBV2	45443	103122	2.37%	0.240%
29	15.957	2200	2205	2213	rBV	89665	166002	3.81%	0.387%
30	16.281	2251	2260	2265	rVV2	136193	330663	7.58%	0.770%
31	16.334	2265	2269	2274	rVV	115026	173449	3.98%	0.404%
32	16.434	2281	2286	2291	rVV2	70328	121892	2.80%	0.284%
33	16.516	2297	2300	2303	rVV4	30908	50991	1.17%	0.119%
34	16.557	2303	2307	2310	rVV2	45751	69793	1.60%	0.163%

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Integration Parameters: LSCINT.P

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35	16.634	2316	2320	2328	rVV	154229	238927	5.48%	0.557%
36	16.710	2328	2333	2337	rVV5	26265	48699	1.12%	0.113%
37	16.757	2337	2341	2343	rVV2	56696	77631	1.78%	0.181%
38	16.792	2343	2347	2361	rVB	245147	378275	8.68%	0.881%
39	16.975	2370	2378	2381	rBV2	2254935	3208342	73.59%	7.473%
40	17.016	2381	2385	2391	rVV	3144537	4359833	100.00%	10.156%
41	17.075	2391	2395	2398	rVV2	125056	207952	4.77%	0.484%
42	17.110	2398	2401	2407	rVV	311573	450854	10.34%	1.050%
43	17.169	2407	2411	2417	rVV4	45826	99637	2.29%	0.232%
44	17.257	2423	2426	2430	rVV4	24716	48464	1.11%	0.113%
45	17.298	2430	2433	2437	rVV4	28021	46955	1.08%	0.109%
46	17.498	2463	2467	2472	rVV2	133221	183791	4.22%	0.428%
47	17.557	2472	2477	2486	rVV4	150289	265156	6.08%	0.618%
48	17.704	2492	2502	2508	rVB3	80117	198086	4.54%	0.461%
49	17.781	2510	2515	2521	rBV4	41935	71877	1.65%	0.167%
50	17.851	2522	2527	2532	rVV	513807	750566	17.22%	1.748%
51	17.904	2532	2536	2543	rVV	544007	772802	17.73%	1.800%
52	17.986	2546	2550	2554	rVV	72914	106150	2.43%	0.247%
53	18.045	2554	2560	2564	rVV2	625757	1021353	23.43%	2.379%
54	18.086	2564	2567	2574	rVB	324417	455868	10.46%	1.062%
55	18.186	2580	2584	2587	rBV3	44933	56219	1.29%	0.131%
56	18.263	2593	2597	2600	rVB2	36224	46671	1.07%	0.109%
57	18.363	2609	2614	2618	rVV	279286	394736	9.05%	0.919%
58	18.410	2618	2622	2632	rVB2	229259	365896	8.39%	0.852%
59	18.616	2653	2657	2662	rVV4	54297	86650	1.99%	0.202%
60	18.663	2662	2665	2669	rVV	64219	83285	1.91%	0.194%
61	18.704	2669	2672	2678	rVB2	47959	66358	1.52%	0.155%
62	18.786	2681	2686	2689	rBV2	208450	341399	7.83%	0.795%
63	18.828	2689	2693	2700	rVV4	166861	449189	10.30%	1.046%
64	18.880	2700	2702	2706	rVB	100216	105655	2.42%	0.246%
65	18.933	2706	2711	2718	rBV3	112051	237196	5.44%	0.553%
66	19.051	2723	2731	2737	rBV	1189642	1632915	37.45%	3.804%
67	19.122	2739	2743	2746	rBV2	42746	56054	1.29%	0.131%
68	19.204	2753	2757	2762	rVB	338222	441147	10.12%	1.028%
69	19.269	2762	2768	2770	rBV3	54508	80801	1.85%	0.188%
70	19.328	2775	2778	2783	rVB	53352	76762	1.76%	0.179%
71	19.416	2783	2793	2804	rBV	1573229	2415106	55.39%	5.626%

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72	19.504	2804	2808	2813	rVB3	46374	76097	1.75%	0.177%
73	19.598	2819	2824	2827	rBV5	25913	41641	0.96%	0.097%
74	19.663	2832	2835	2842	rVB5	28135	43526	1.00%	0.101%
75	19.757	2846	2851	2859	rBV2	127416	260934	5.98%	0.608%
76	19.892	2869	2874	2876	rVV3	117294	180198	4.13%	0.420%
77	19.922	2876	2879	2884	rVB2	201883	285707	6.55%	0.666%
78	20.027	2894	2897	2903	rVV2	107498	126500	2.90%	0.295%
79	20.086	2903	2907	2911	rVB	172973	206448	4.74%	0.481%
80	20.227	2927	2931	2935	rVV	119086	149824	3.44%	0.349%
81	20.269	2935	2938	2944	rVB	132910	183193	4.20%	0.427%
82	20.686	3005	3009	3014	rVB	90068	142063	3.26%	0.331%
83	20.851	3033	3037	3040	rBV3	92505	118971	2.73%	0.277%
84	20.892	3040	3044	3047	rVV	103961	127095	2.92%	0.296%
85	20.927	3047	3050	3055	rVV	66234	98698	2.26%	0.230%
86	20.992	3055	3061	3067	rVB2	97454	191469	4.39%	0.446%
87	21.098	3075	3079	3081	rBV	33820	42333	0.97%	0.099%
88	21.210	3091	3098	3102	rBV2	2398286	3618558	83.00%	8.429%
89	21.251	3102	3105	3112	rVB	422558	576467	13.22%	1.343%
90	21.433	3132	3136	3144	rBV2	110779	176321	4.04%	0.411%
91	21.769	3190	3193	3197	rVB	90836	100918	2.31%	0.235%
92	21.822	3199	3202	3207	rVB2	83416	100049	2.29%	0.233%
93	21.963	3223	3226	3229	rBV3	50053	67568	1.55%	0.157%
94	22.274	3277	3279	3284	rVB3	43047	46853	1.07%	0.109%
95	22.774	3359	3364	3369	rBV	236723	482348	11.06%	1.124%
96	22.939	3389	3392	3397	rVB2	53262	82961	1.90%	0.193%
97	23.233	3438	3442	3447	rBV	119686	189368	4.34%	0.441%
98	23.327	3454	3458	3465	rVB2	223427	377502	8.66%	0.879%
99	23.421	3467	3474	3493	rVB	1279108	2499586	57.33%	5.822%
100	23.951	3562	3564	3572	rVB3	51420	94961	2.18%	0.221%

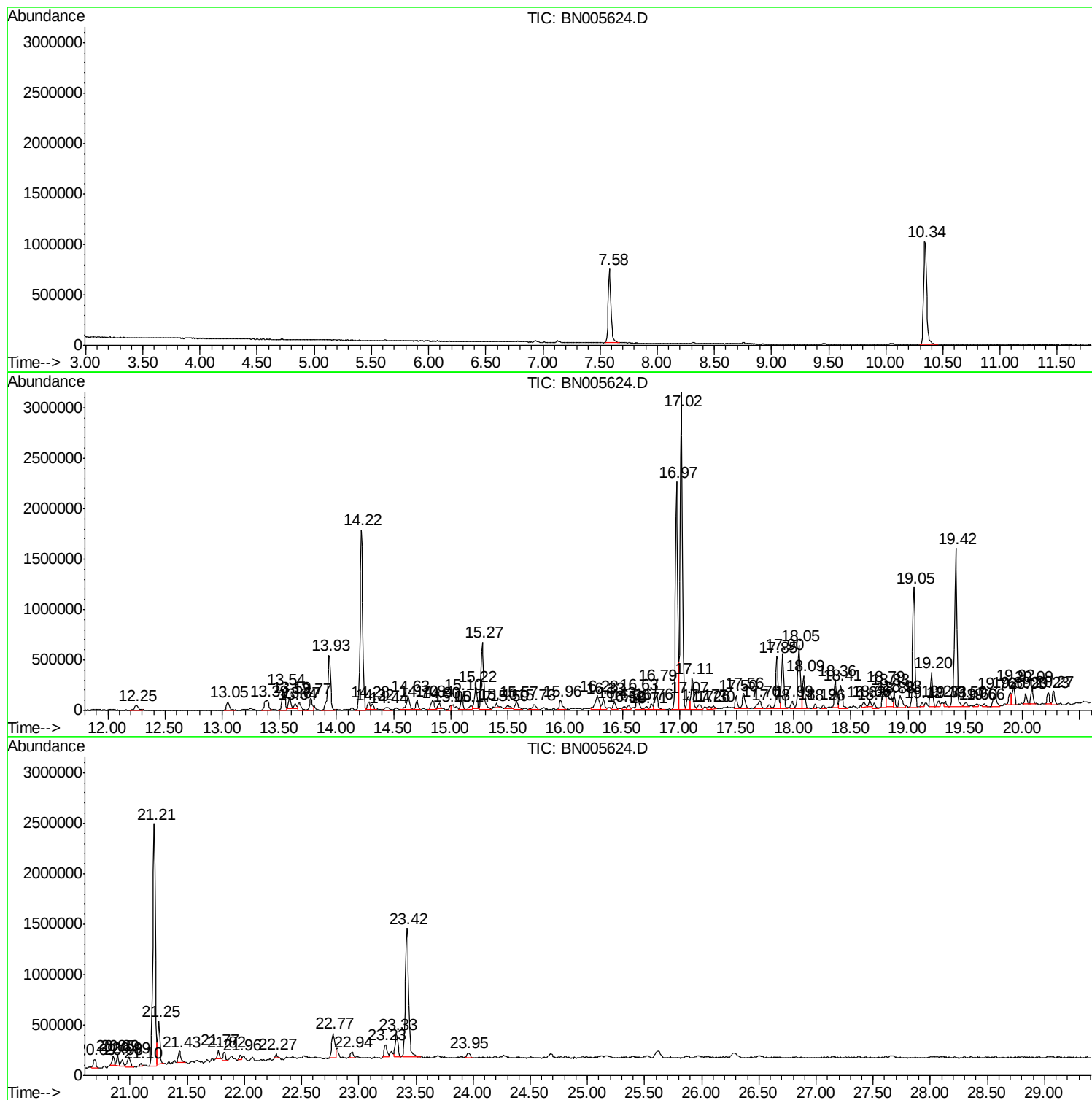
Sum of corrected areas: 42930115

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN041919MA.M
Quant Title : SVOA CALIBRATION

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



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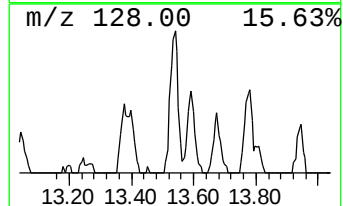
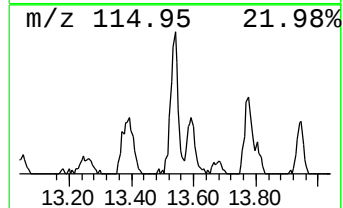
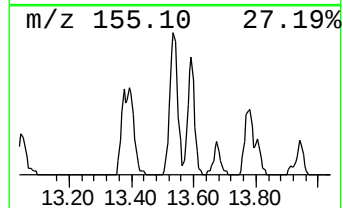
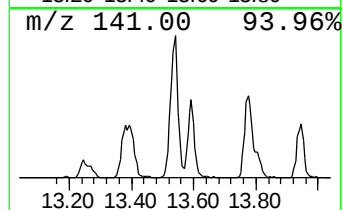
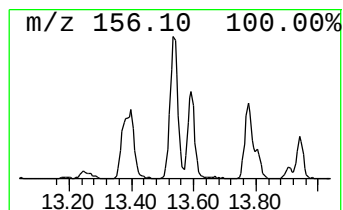
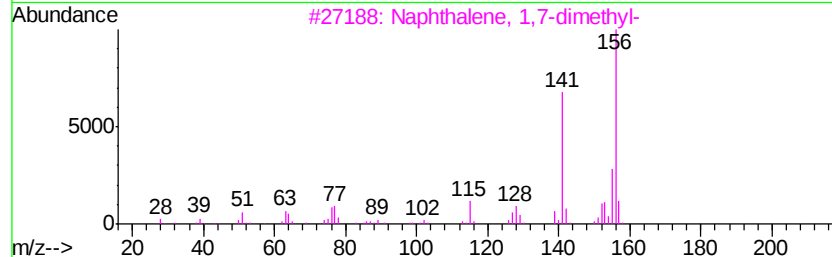
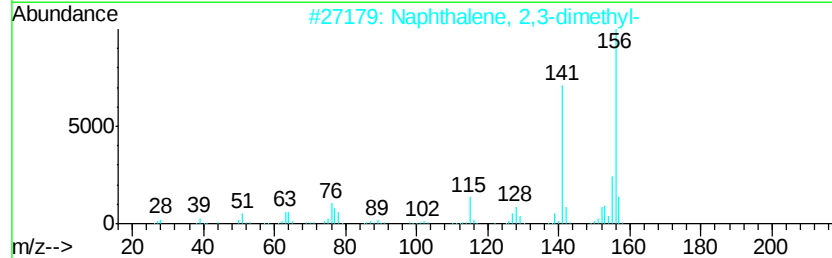
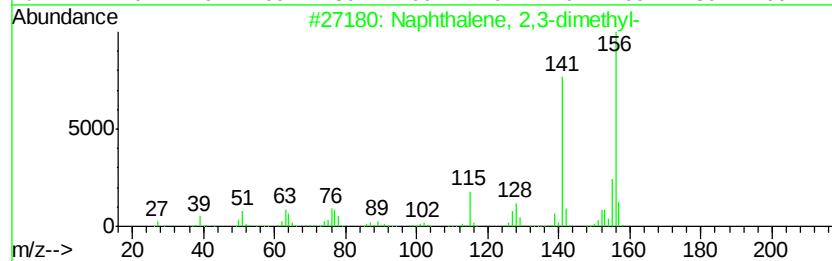
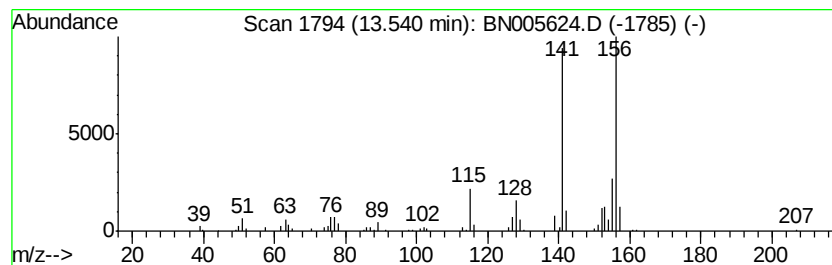
Instrument :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN041919MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Naphthalene, 2,3-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.54	2.56 ng/ul	363839	Acenaphthene-d10	14.22
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 2,3-dimethyl-	156 C12H12	000581-40-8 97
2		Naphthalene, 2,3-dimethyl-	156 C12H12	000581-40-8 97
3		Naphthalene, 1,7-dimethyl-	156 C12H12	000575-37-1 97
4		Naphthalene, 1,4-dimethyl-	156 C12H12	000571-58-4 97
5		Naphthalene, 2,3-dimethyl-	156 C12H12	000581-40-8 96



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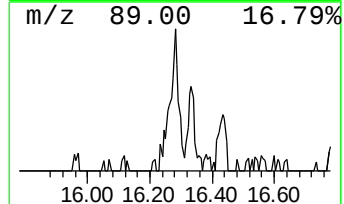
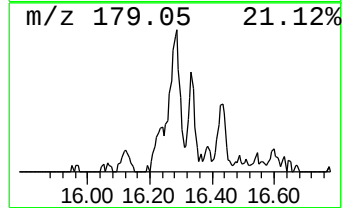
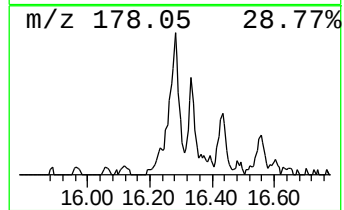
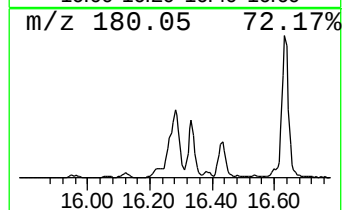
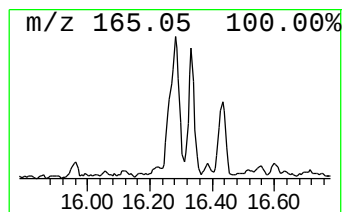
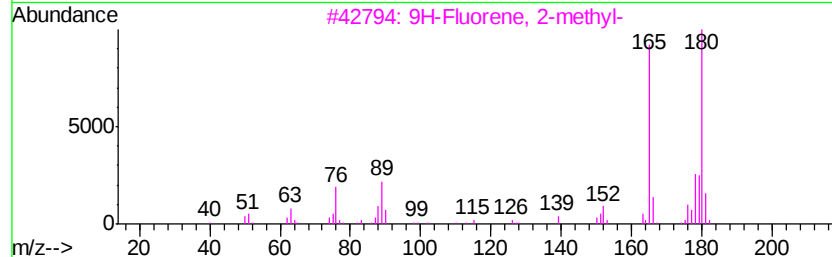
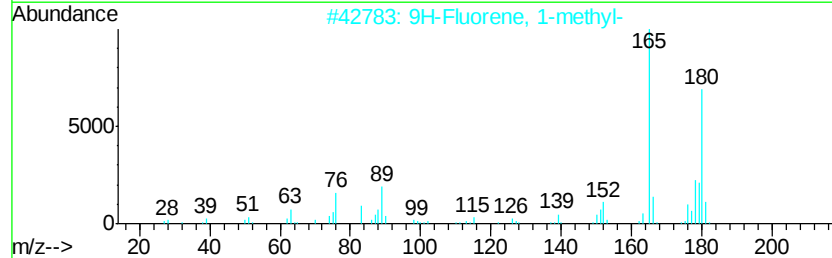
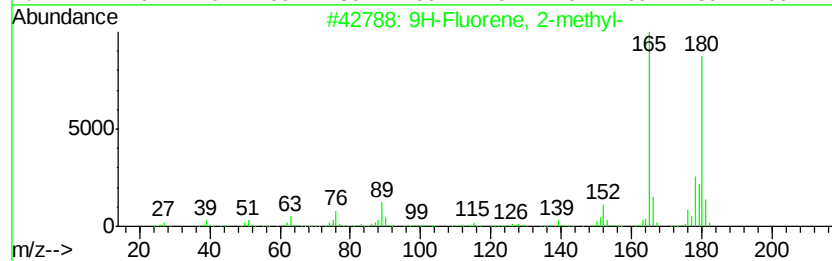
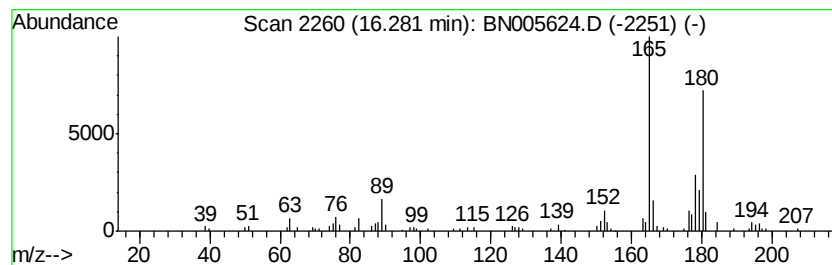
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 9H-Fluorene, 2-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.28	2.06 ng/ul	330663	Phenanthrene-d10	16.97

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9H-Fluorene, 2-methyl-	180	C14H12	001430-97-3	94
2			9H-Fluorene, 1-methyl-	180	C14H12	001730-37-6	93
3			9H-Fluorene, 2-methyl-	180	C14H12	001430-97-3	89
4			9H-Fluorene, 1-methyl-	180	C14H12	001730-37-6	89
5			9H-Fluorene, 1-methyl-	180	C14H12	001730-37-6	89



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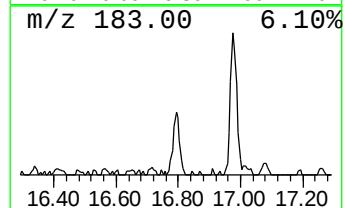
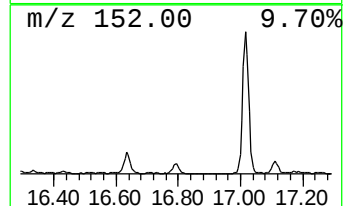
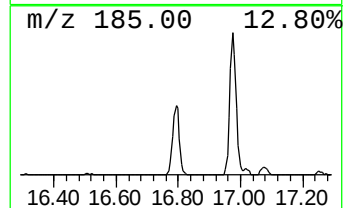
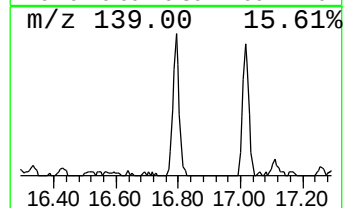
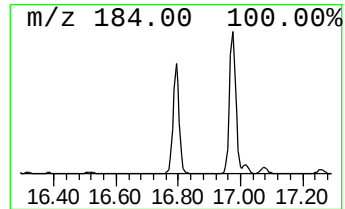
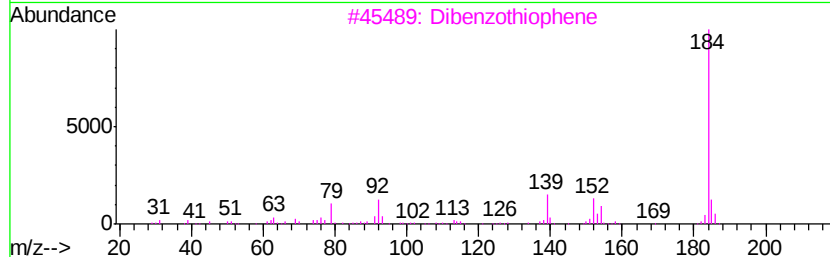
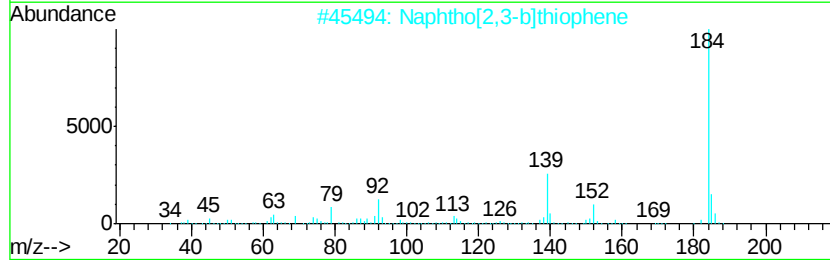
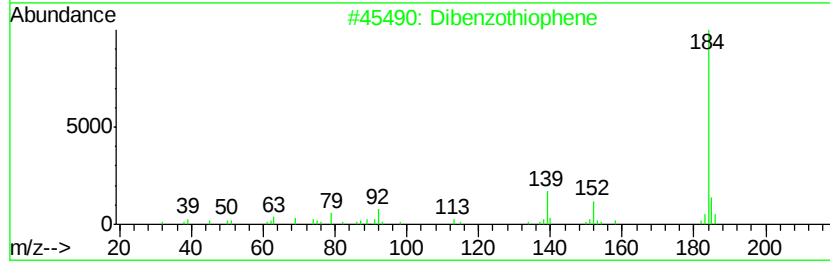
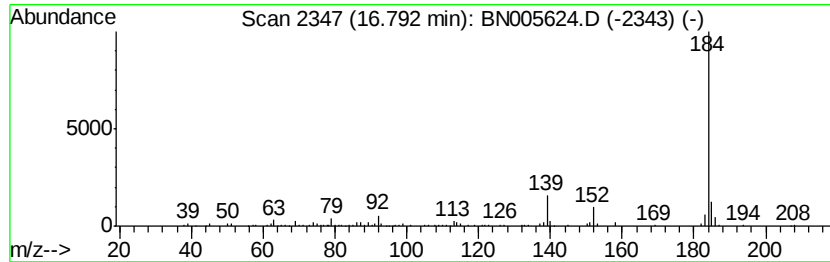
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TIC Integration Parameters: LSCINT.P

Peak Number 3 Dibenzothiophene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.79	2.36 ng/ul	378275	Phenanthrene-d10	16.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dibenzothiophene	184	C12H8S	000132-65-0	96
2		Naphtho[2,3-b]thiophene	184	C12H8S	000268-77-9	94
3		Dibenzothiophene	184	C12H8S	000132-65-0	94
4		Dibenzothiophene	184	C12H8S	000132-65-0	94
5		Dibenzothiophene	184	C12H8S	000132-65-0	94



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Data File : BN005624.D
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Operator : JU/SJ
Sample : K2603-06MEDL 30X
Misc :
ALS Vial : 4 Sample Multiplier: 1

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ClientSampleId :
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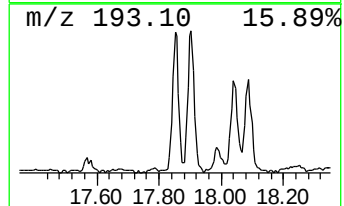
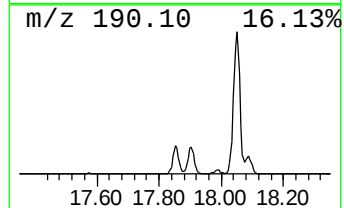
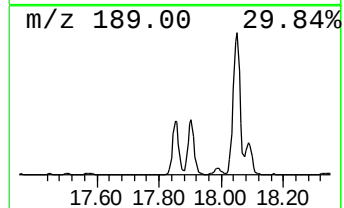
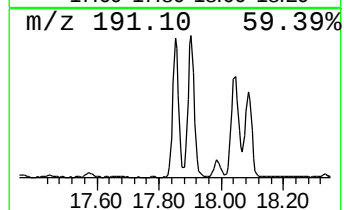
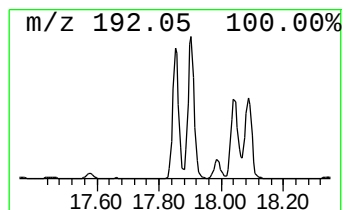
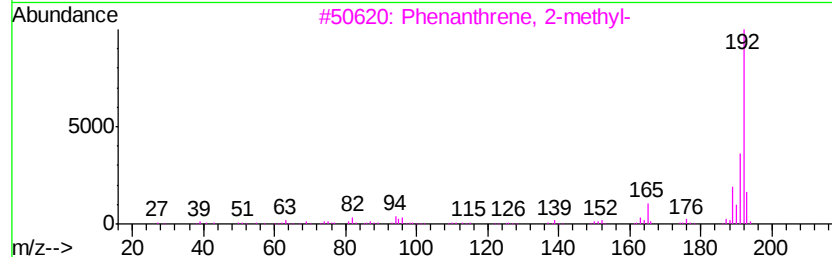
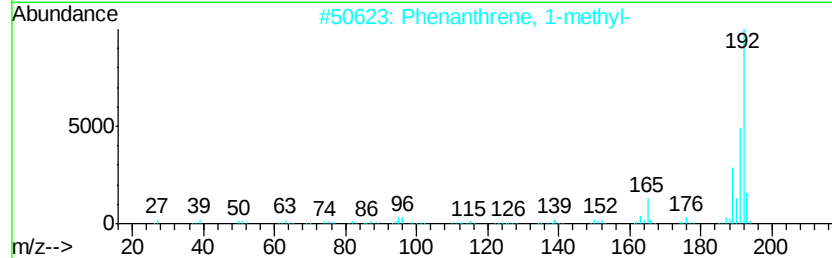
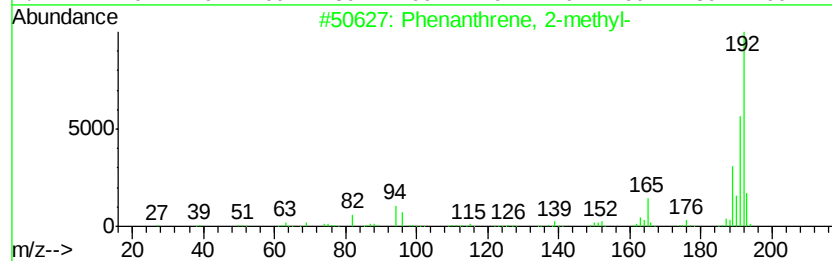
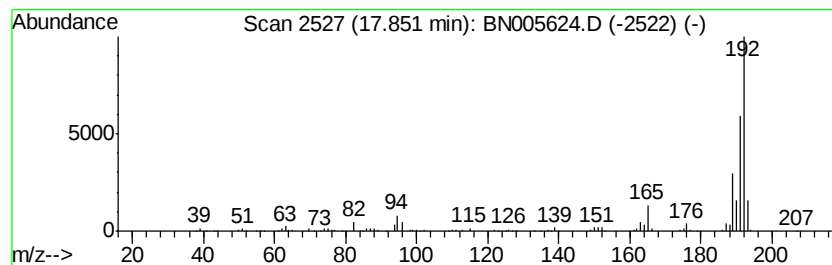
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Phenanthrene, 2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.85	4.68 ng/ul	750566	Phenanthrene-d10	16.97

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	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	Anthracene, 2-methyl-	192	C15H12	000613-12-7	92	



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Operator : JU/SJ
Sample : K2603-06MEDL 30X
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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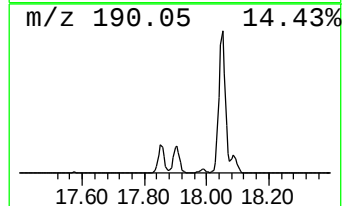
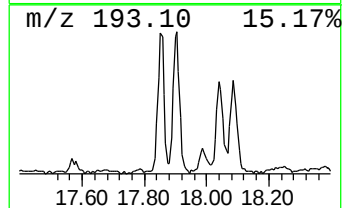
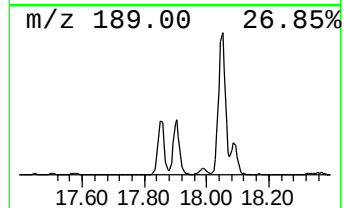
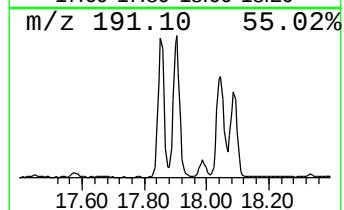
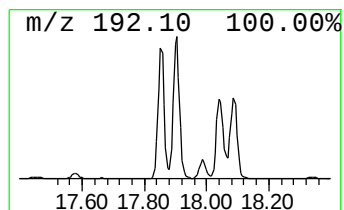
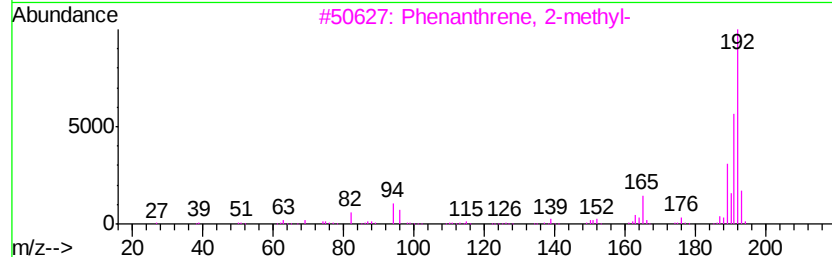
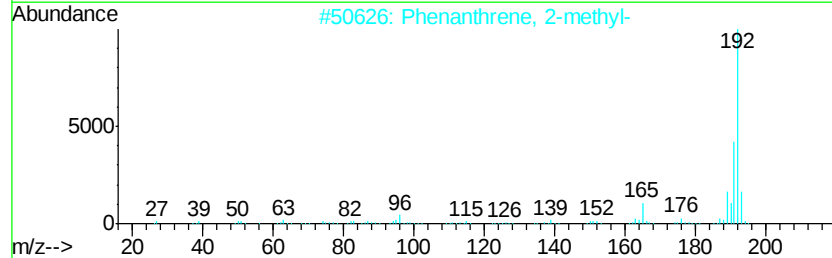
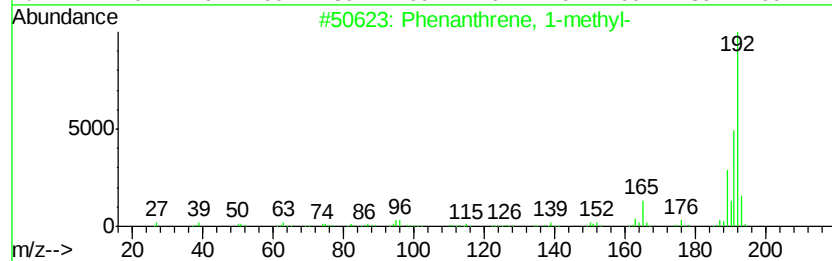
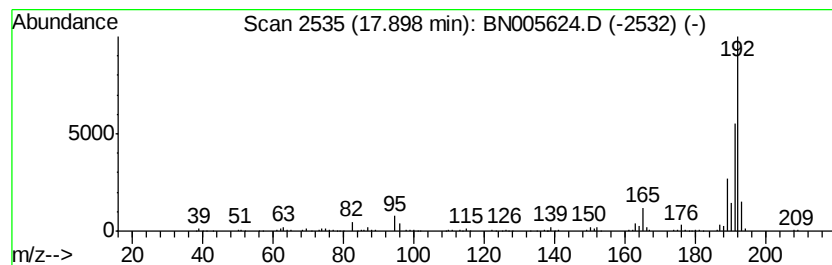
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.90	4.82 ng/ul	772802	Phenanthrene-d10	16.97

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	97	
3	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95	
4	1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	95	
5	Anthracene, 2-methyl-	192	C15H12	000613-12-7	95	



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN051119\
Data File : BN005624.D
Acq On : 11 May 2019 13:42
Operator : JU/SJ
Sample : K2603-06MEDL 30X
Misc :
ALS Vial : 4 Sample Multiplier: 1

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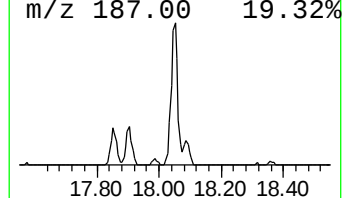
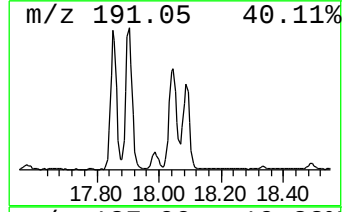
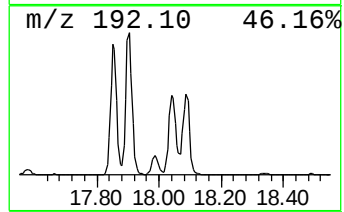
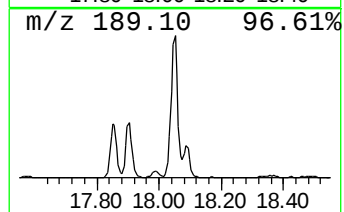
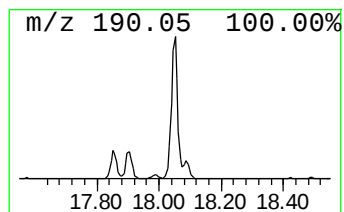
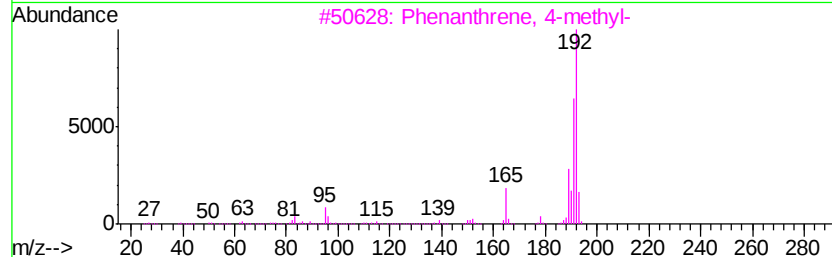
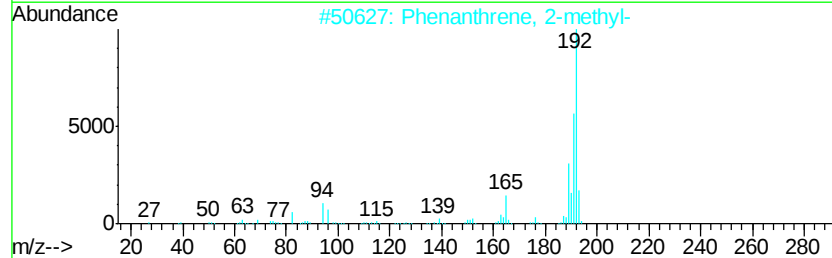
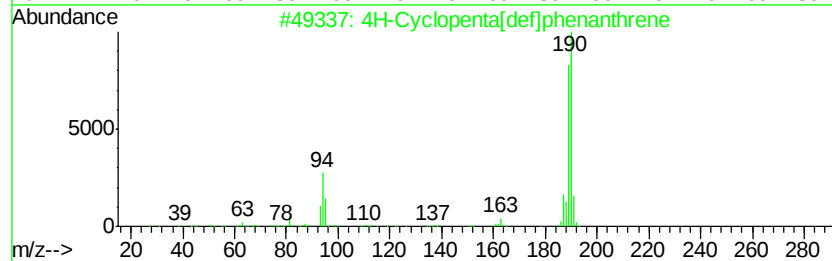
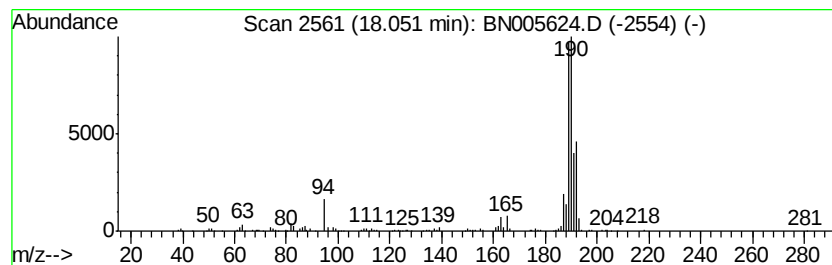
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 4H-Cyclopenta[def]phenanthrene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.05	6.37 ng/ul	1021350	Phenanthrene-d10	16.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	90
2		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	46
3		Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	42
4		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	42
5		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	42



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN051119\
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Acq On : 11 May 2019 13:42
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Misc :
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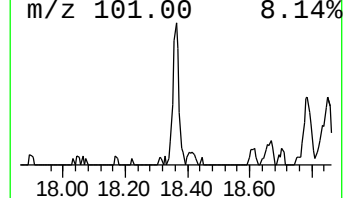
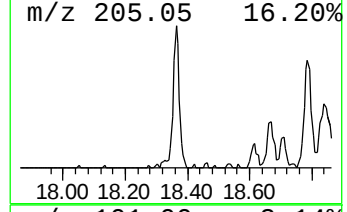
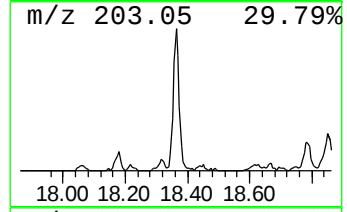
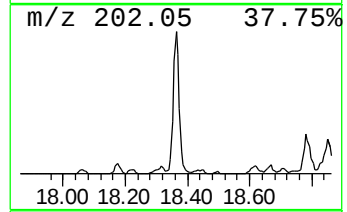
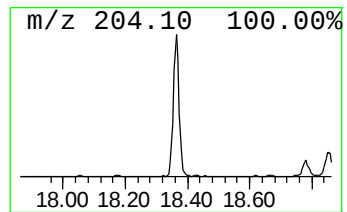
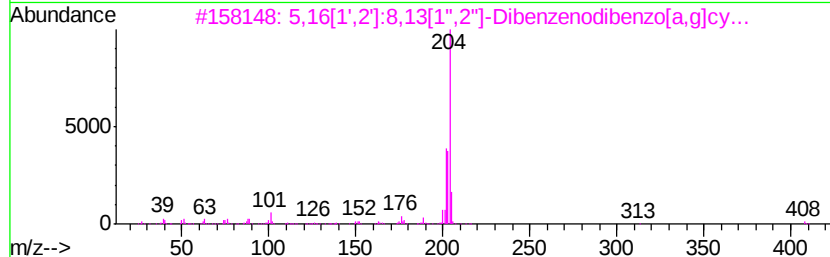
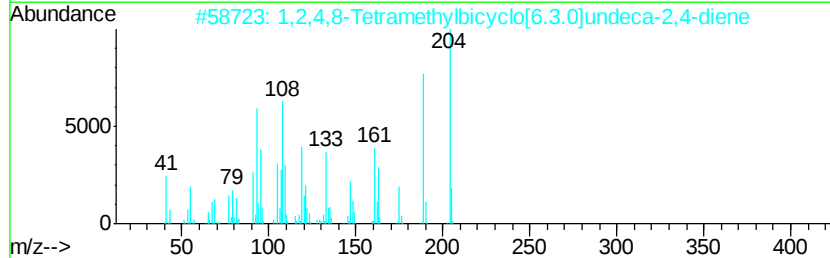
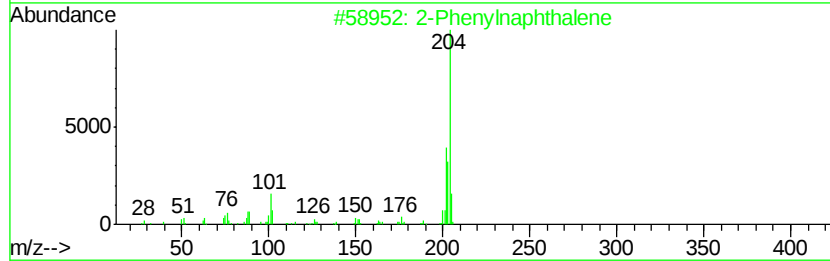
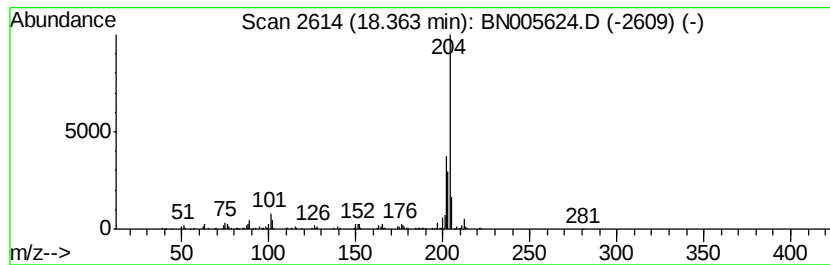
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 2-Phenylanthralene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.36	2.46 ng/ul	394736	Phenanthrene-d10	16.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Phenylanthralene	204	C16H12	035465-71-5	93
2		1,2,4,8-Tetramethylbicyclo[6.3.0...]	204	C15H24	137235-51-9	90
3		5,16[1',2']:8,13[1'',2'']-Dibenz...	408	C32H24	005672-97-9	86
4		Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	81
5		Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	76



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN051119\
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Misc :
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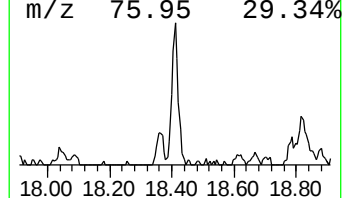
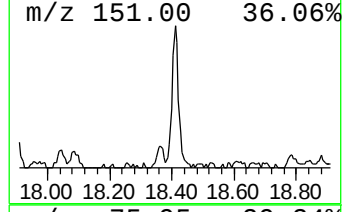
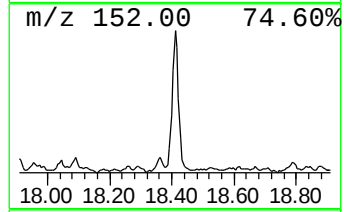
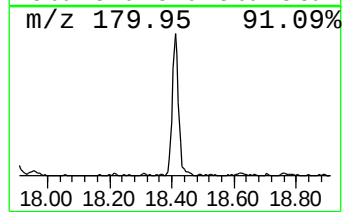
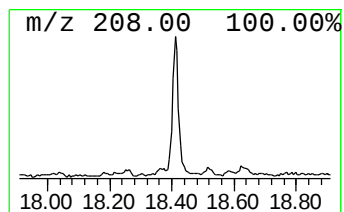
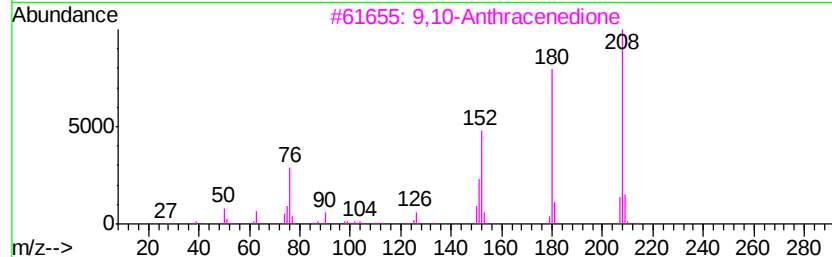
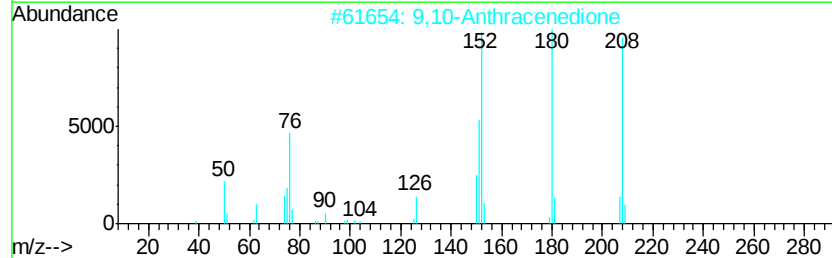
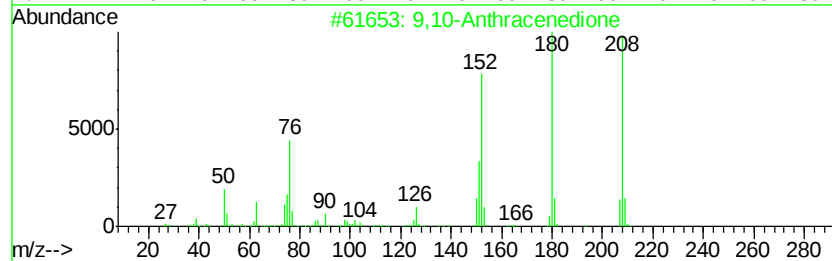
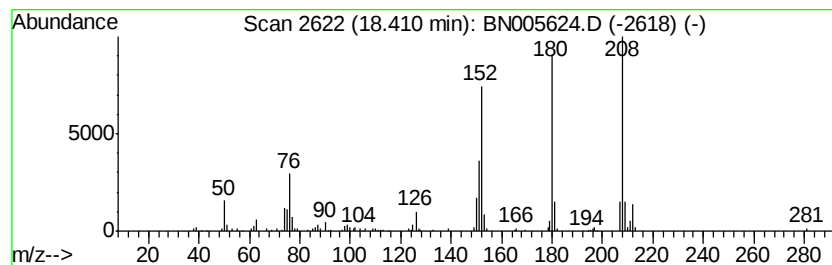
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 9,10-Anthracenedione Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.41	2.28 ng/ul	365896	Phenanthrene-d10	16.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9,10-Anthracenedione	208	C14H8O2	000084-65-1	97
2		9,10-Anthracenedione	208	C14H8O2	000084-65-1	97
3		9,10-Anthracenedione	208	C14H8O2	000084-65-1	90
4		Benzo[c]cinnoline	180	C12H8N2	000230-17-1	83
5		Benzo[c]cinnoline	180	C12H8N2	000230-17-1	78



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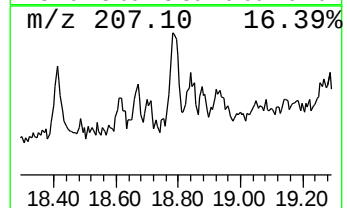
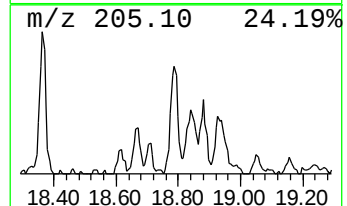
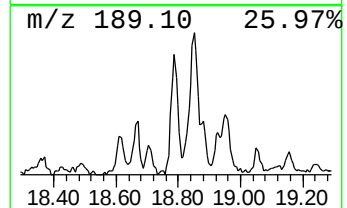
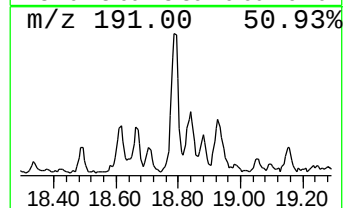
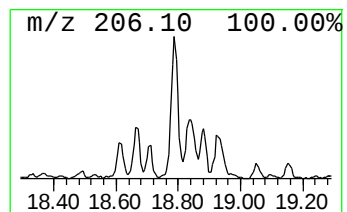
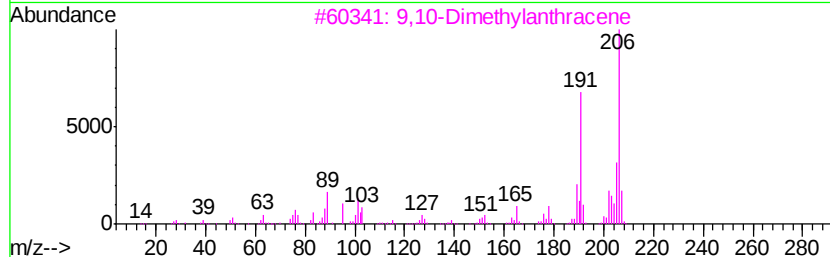
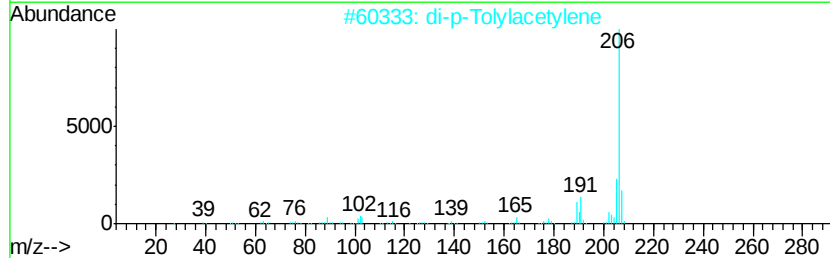
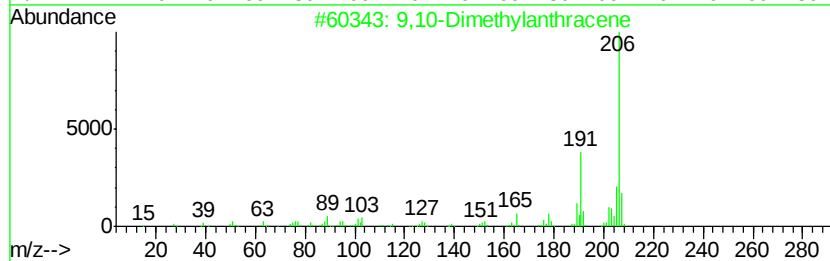
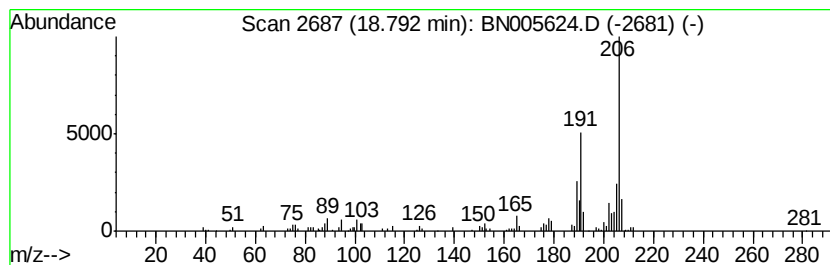
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 9,10-Dimethylantracene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.79	2.13 ng/ul	341399	Phenanthrene-d10	16.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9,10-Dimethylantracene	206	C16H14	000781-43-1	94
2		di-p-Tolylacetylene	206	C16H14	002789-88-0	94
3		9,10-Dimethylantracene	206	C16H14	000781-43-1	93
4		Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	93
5		Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN051119\
Data File : BN005624.D
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Misc :
ALS Vial : 4 Sample Multiplier: 1

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ClientSampleId :
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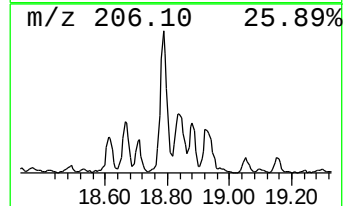
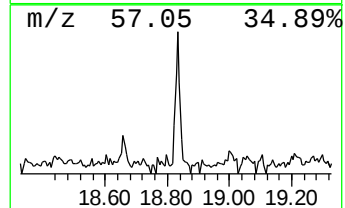
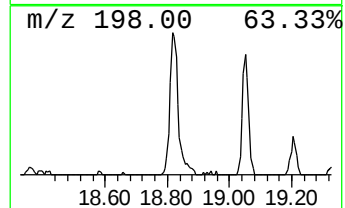
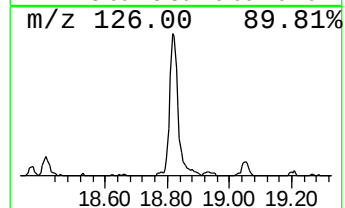
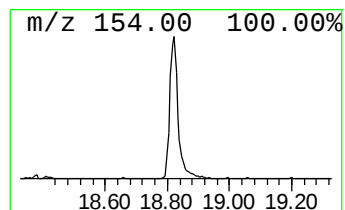
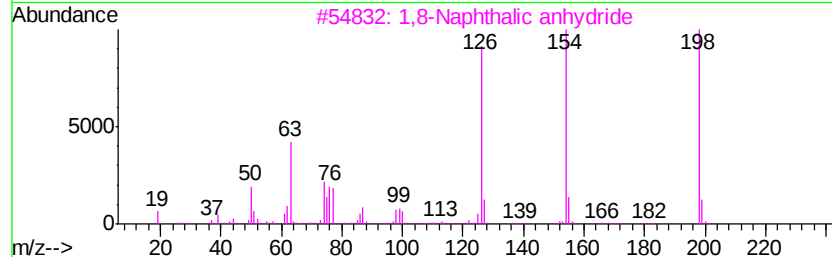
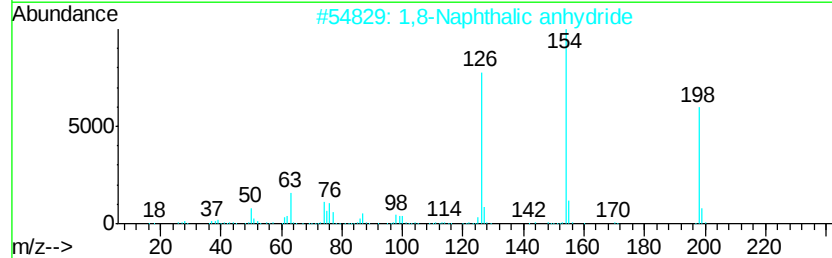
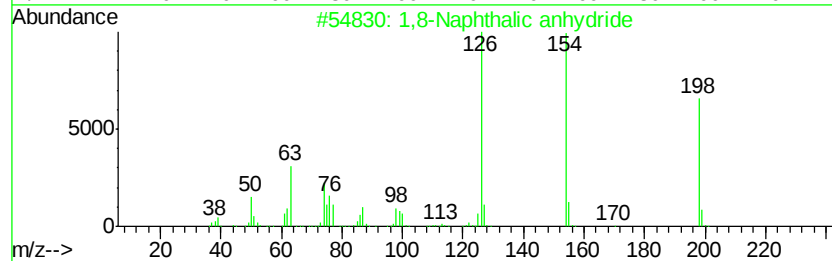
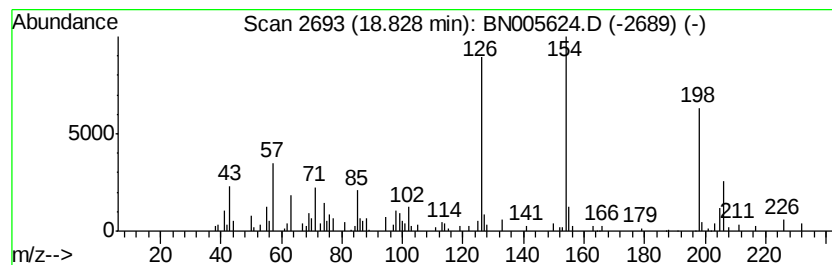
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 1,8-Naphthalic anhydride Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.83	2.80 ng/ul	449189	Phenanthrene-d10	16.97

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,8-Naphthalic anhydride	198	C12H6O3	000081-84-5	89
2			1,8-Naphthalic anhydride	198	C12H6O3	000081-84-5	89
3			1,8-Naphthalic anhydride	198	C12H6O3	000081-84-5	64
4			1,8-Naphthalic anhydride	198	C12H6O3	000081-84-5	60
5			1H-Benz[f]indene-1,2,3-trione	210	C13H6O3	020927-01-9	35



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN051119\
Data File : BN005624.D
Acq On : 11 May 2019 13:42
Operator : JU/SJ
Sample : K2603-06MEDL 30X
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
GAJ58MEDL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN041919MA.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
Naphthalene, 2,3-...	13.54	2.6	ng/ul	363839	3	14.22	2846250	20.0
9H-Fluorene, 2-me...	16.28	2.1	ng/ul	330663	4	16.97	3208340	20.0
Dibenzothiophene	16.79	2.4	ng/ul	378275	4	16.97	3208340	20.0
Phenanthrene, 2-m...	17.85	4.7	ng/ul	750566	4	16.97	3208340	20.0
Phenanthrene, 1-m...	17.90	4.8	ng/ul	772802	4	16.97	3208340	20.0
4H-Cyclopenta[def...	18.05	6.4	ng/ul	1021350	4	16.97	3208340	20.0
2-Phenylnaphthalene	18.36	2.5	ng/ul	394736	4	16.97	3208340	20.0
9,10-Anthracenedione	18.41	2.3	ng/ul	365896	4	16.97	3208340	20.0
9,10-Dimethylanth...	18.79	2.1	ng/ul	341399	4	16.97	3208340	20.0
1,8-Naphthalic an...	18.83	2.8	ng/ul	449189	4	16.97	3208340	20.0