

Data Path : Z:\HPCHEM1\BNA M\DATA\BM103017\
 Data File : BM012587.D
 Acq On : 30 Oct 2017 23:26
 Operator : SJ/JU
 Sample : I5886-05
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 ESNE5

Integration Parameters: LSCINT.P

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

Filtering: 5
 Min Area: 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:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102417.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	3.140	6	9	14	rVB	137727	174521	1.55%	0.136%
2	3.863	127	132	138	rVV	315056	425646	3.77%	0.332%
3	4.993	318	324	331	rVB2	80181	135855	1.20%	0.106%
4	7.022	662	669	680	rBV2	778732	1415386	12.55%	1.105%
5	7.187	691	697	705	rBV	644679	966437	8.57%	0.754%
6	7.392	724	732	739	rBV	820986	1315647	11.67%	1.027%
7	7.857	803	811	818	rBV	1314814	2104390	18.66%	1.643%
8	8.557	924	930	941	rBV	646288	1028954	9.13%	0.803%
9	9.010	1000	1007	1015	rBV	761188	1248496	11.07%	0.975%
10	9.728	1121	1129	1136	rBV	622161	1068568	9.48%	0.834%
11	10.269	1212	1221	1231	rBV	988172	1629594	14.45%	1.272%
12	10.645	1276	1285	1292	rBV	1659387	2792220	24.76%	2.180%
13	10.780	1302	1308	1317	rBV	443929	843751	7.48%	0.659%
14	11.963	1498	1509	1518	rBV	92497	157035	1.39%	0.123%
15	13.892	1829	1837	1842	rBV	1639539	2222331	19.71%	1.735%
16	13.939	1842	1845	1849	rVB	103772	137773	1.22%	0.108%
17	14.180	1879	1886	1897	rBV	1756695	2800776	24.84%	2.186%
18	14.486	1930	1938	1945	rBV2	2317418	3566130	31.63%	2.784%
19	14.545	1945	1948	1957	rVB	70001	113153	1.00%	0.088%
20	14.680	1966	1971	1980	rBV	479172	814716	7.23%	0.636%
21	14.892	2002	2007	2015	rVB3	58905	132958	1.18%	0.104%
22	15.474	2100	2106	2112	rBV	2243212	3329039	29.52%	2.599%
23	15.533	2112	2116	2122	rVB2	150443	206705	1.83%	0.161%
24	15.598	2122	2127	2135	rBV	508085	695716	6.17%	0.543%
25	15.839	2162	2168	2176	rVB	419252	595023	5.28%	0.464%
26	16.221	2225	2233	2237	rBV3	75568	200394	1.78%	0.156%
27	16.886	2341	2346	2353	rVB	84260	130660	1.16%	0.102%
28	17.045	2369	2373	2381	rVB	109200	166066	1.47%	0.130%
29	17.227	2399	2404	2408	rBV2	3021968	4242060	37.62%	3.311%
30	17.268	2408	2411	2416	rVB	2287884	2981489	26.44%	2.327%
31	17.327	2416	2421	2425	rBV	2403666	3367636	29.87%	2.629%
32	17.421	2432	2437	2443	rVB	89593	157425	1.40%	0.123%
33	17.627	2465	2472	2477	rBV	242040	395286	3.51%	0.309%
34	17.745	2486	2492	2498	rVB3	58278	126239	1.12%	0.099%

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35	18.115	2548	2555	2558	rVV2	438916	835852	7.41%	0.652%
36	18.151	2558	2561	2566	rVB	337045	461329	4.09%	0.360%
37	18.298	2580	2586	2590	rBV3	417178	698184	6.19%	0.545%
38	18.333	2590	2592	2599	rVB2	146100	186107	1.65%	0.145%
39	18.603	2633	2638	2641	rBV	236869	338244	3.00%	0.264%
40	18.645	2641	2645	2649	rVB	446522	586633	5.20%	0.458%
41	19.021	2704	2709	2713	rBV2	123084	210711	1.87%	0.164%
42	19.156	2728	2732	2741	rBV	196553	327789	2.91%	0.256%
43	19.280	2745	2753	2760	rVV	7253711	9935790	88.12%	7.756%
44	19.392	2767	2772	2776	rBV4	235586	407654	3.62%	0.318%
45	19.527	2791	2795	2798	rBV	121222	136767	1.21%	0.107%
46	19.615	2804	2810	2812	rBV2	3954279	5580197	49.49%	4.356%
47	19.645	2812	2815	2822	rVV	5897518	7581441	67.24%	5.918%
48	19.715	2823	2827	2832	rVB3	199764	298155	2.64%	0.233%
49	19.821	2841	2845	2850	rVB	276532	341257	3.03%	0.266%
50	19.886	2852	2856	2861	rVB	143068	208208	1.85%	0.163%
51	19.962	2864	2869	2876	rBV3	268817	684702	6.07%	0.534%
52	20.021	2876	2879	2882	rVB	117096	134724	1.19%	0.105%
53	20.139	2895	2899	2908	rVB	578616	883294	7.83%	0.690%
54	20.239	2912	2916	2919	rBV	374166	481519	4.27%	0.376%
55	20.297	2920	2926	2929	rVV2	327818	552118	4.90%	0.431%
56	20.333	2929	2932	2935	rVB	177751	196824	1.75%	0.154%
57	20.433	2947	2949	2953	rVB	154237	159289	1.41%	0.124%
58	20.480	2953	2957	2961	rBV3	218451	324972	2.88%	0.254%
59	20.880	3022	3025	3030	rVB	524722	656081	5.82%	0.512%
60	21.044	3048	3053	3057	rBV2	648962	1013202	8.99%	0.791%
61	21.086	3057	3060	3062	rVV	531366	641157	5.69%	0.501%
62	21.121	3062	3066	3071	rVV3	757581	1481275	13.14%	1.156%
63	21.174	3071	3075	3085	rVB2	830391	1264377	11.21%	0.987%
64	21.397	3106	3113	3117	rBV3	5508645	11275749	100.00%	8.802%
65	21.433	3117	3119	3130	rVB	3717971	4736189	42.00%	3.697%
66	21.556	3136	3140	3145	rBV	717458	960133	8.52%	0.750%
67	21.715	3163	3167	3172	rBV2	489047	754515	6.69%	0.589%
68	23.009	3380	3387	3391	rBV	3927308	8769559	77.77%	6.846%
69	23.497	3465	3470	3475	rBV	1935511	3585950	31.80%	2.799%
70	23.550	3475	3479	3483	rVV2	2162489	3942413	34.96%	3.078%
71	23.597	3483	3487	3495	rVB	2885197	5241005	46.48%	4.091%

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72	23.697	3498	3504	3509	rBV	2944205	5393896	47.84%	4.211%
73	26.026	3894	3900	3913	rVB2	1818151	5147589	45.65%	4.018%

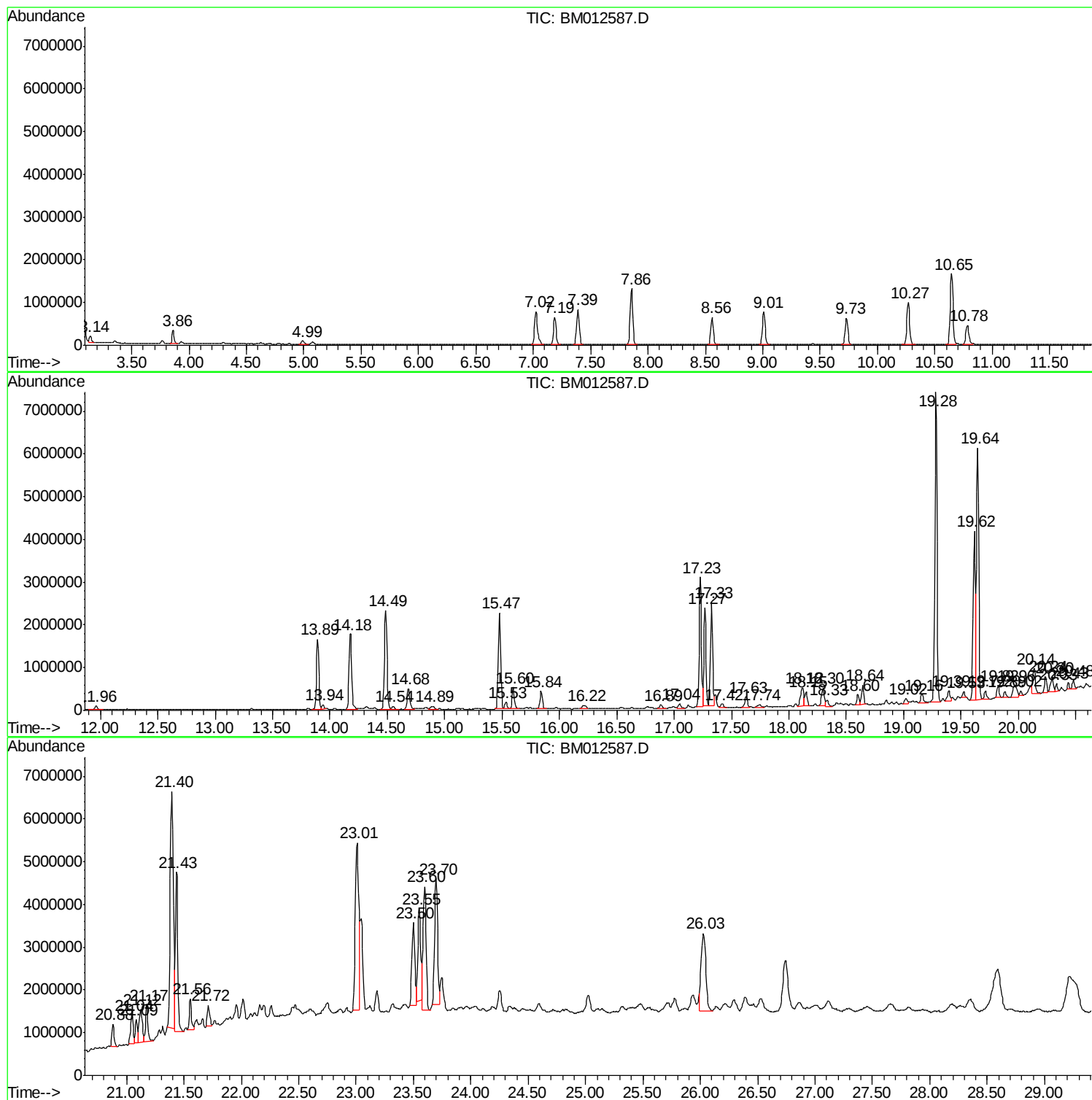
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Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM102417.M
Quant Title : SVOA CALIBRATION

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



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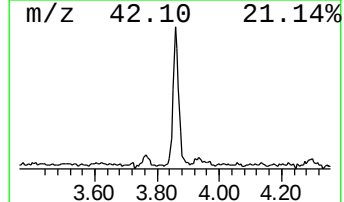
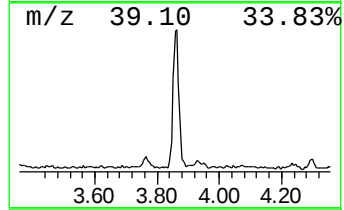
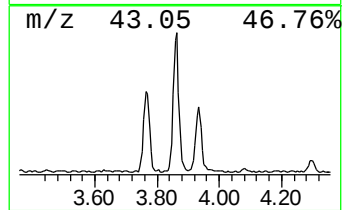
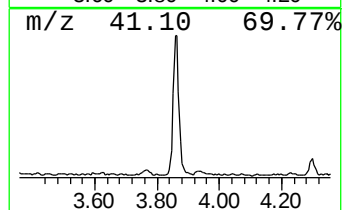
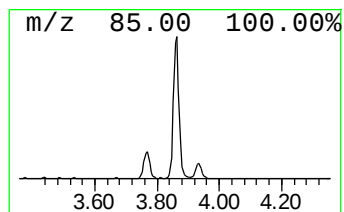
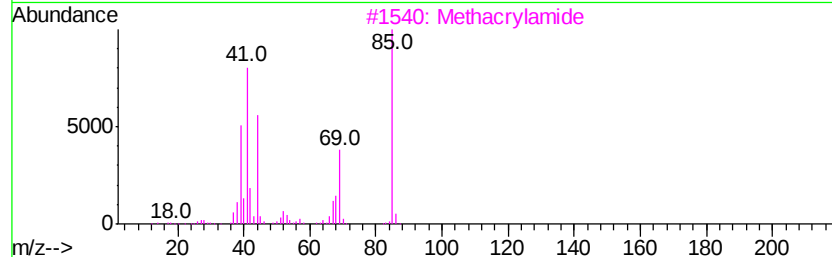
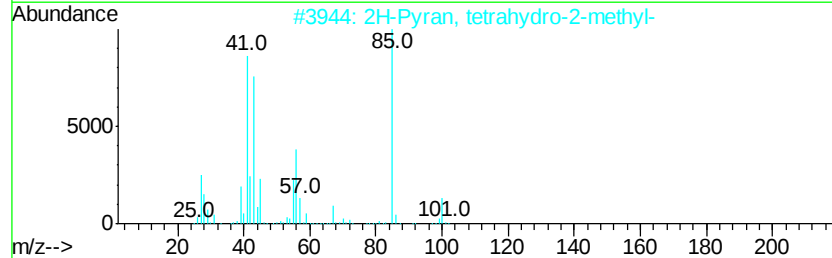
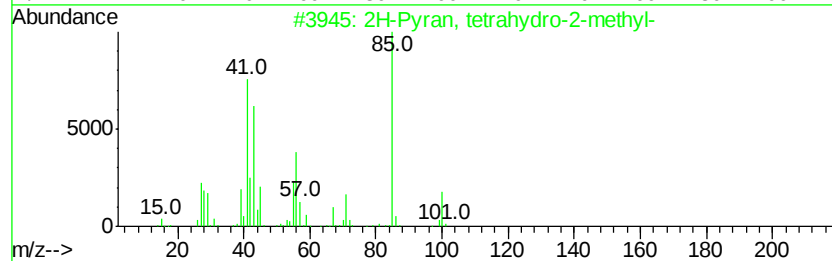
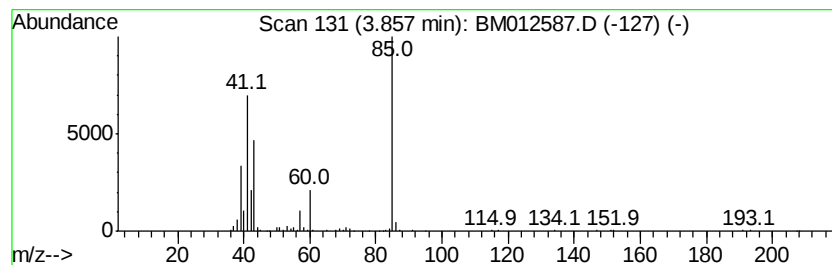
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown-01 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.86	4.05 ng/ul	425646	1,4-Dichlorobenzene-d4	7.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2H-Pyran, tetrahydro-2-methyl-	100	C6H12O	010141-72-7	38
2		2H-Pyran, tetrahydro-2-methyl-	100	C6H12O	010141-72-7	38
3		Methacrylamide	85	C4H7NO	000079-39-0	36
4		Methacrylamide	85	C4H7NO	000079-39-0	12
5		2(3H)-Furanone, dihydro-5-propyl-	128	C7H12O2	000105-21-5	9



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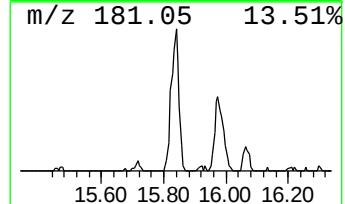
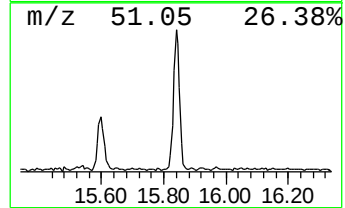
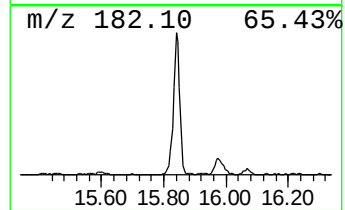
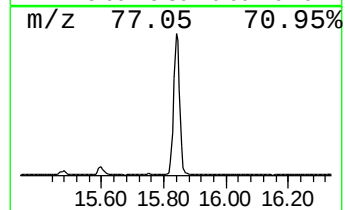
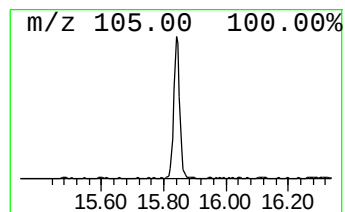
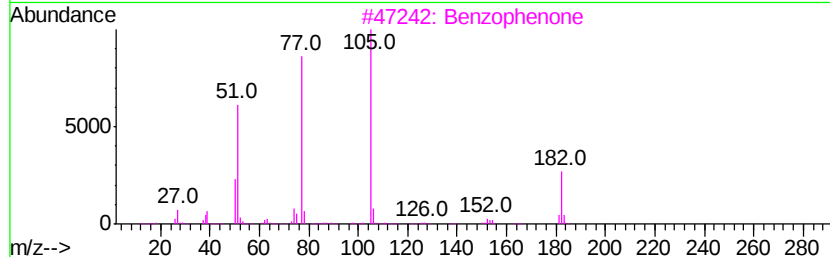
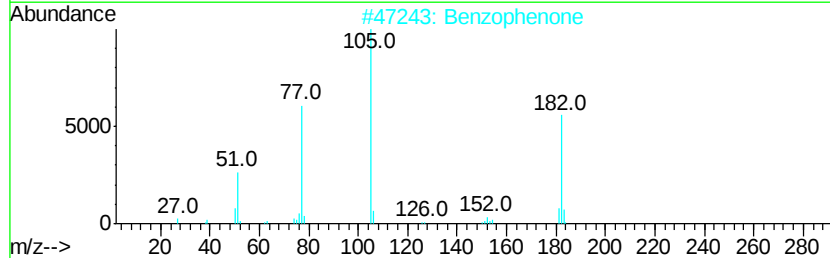
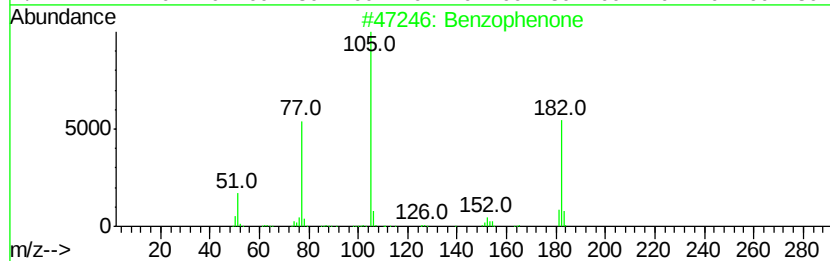
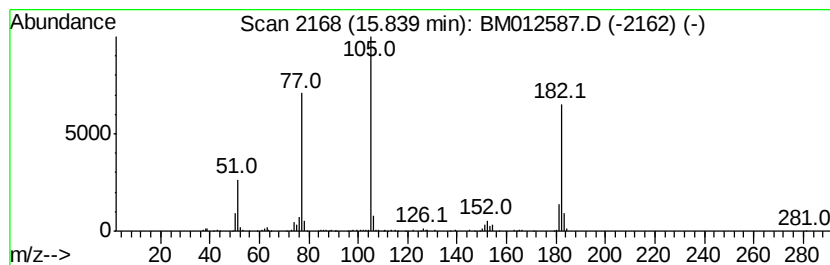
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM102417.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzophenone Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.84	3.34 ng/ul	595023	Acenaphthene-d10	14.49

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzophenone		182	C13H10O	000119-61-9	97
2	Benzophenone		182	C13H10O	000119-61-9	96
3	Benzophenone		182	C13H10O	000119-61-9	95
4	Benzophenone		182	C13H10O	000119-61-9	95
5	Benzophenone		182	C13H10O	000119-61-9	87



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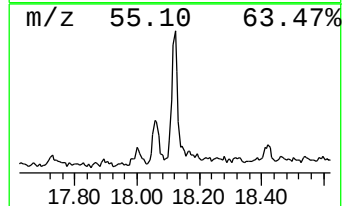
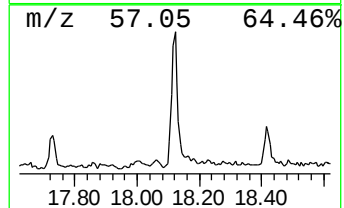
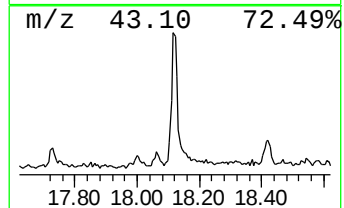
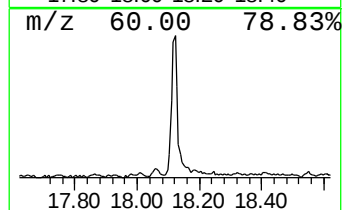
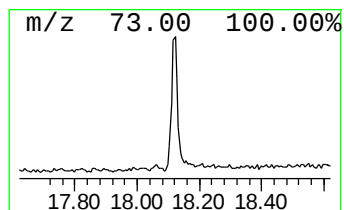
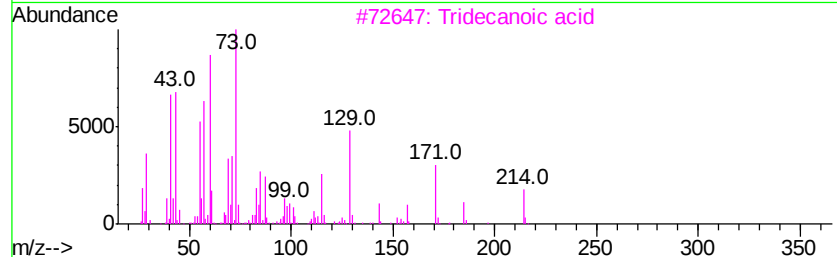
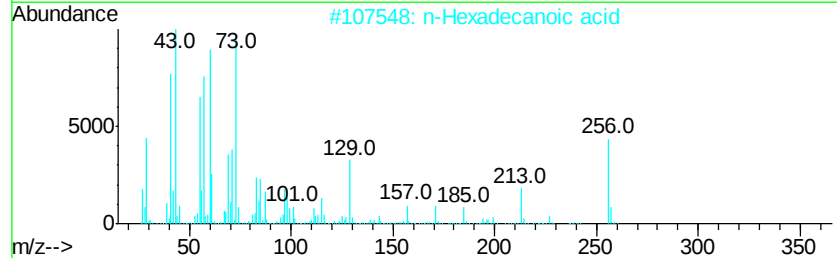
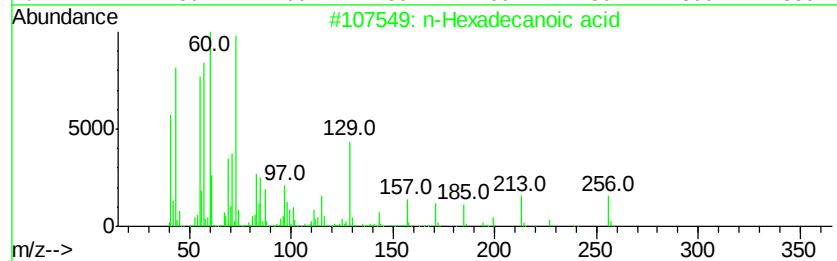
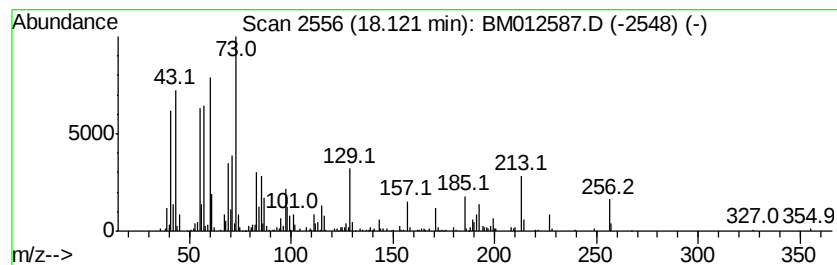
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Peak Number 3 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.12	3.94 ng/ul	835852	Phenanthrene-d10	17.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
3		Tridecanoic acid	214	C13H26O2	000638-53-9	91
4		Tridecanoic acid	214	C13H26O2	000638-53-9	83
5		Pentadecanoic acid	242	C15H30O2	001002-84-2	60



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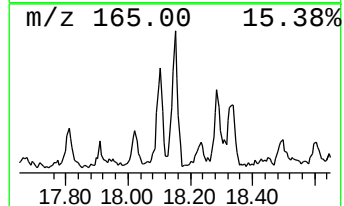
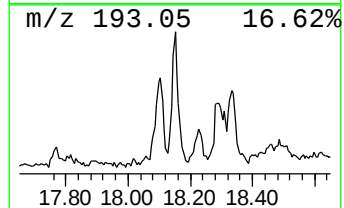
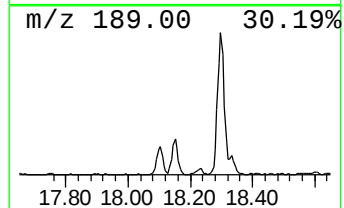
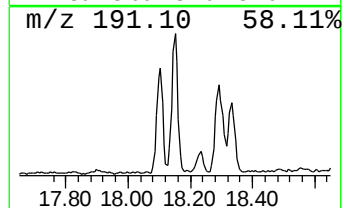
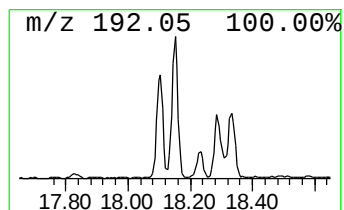
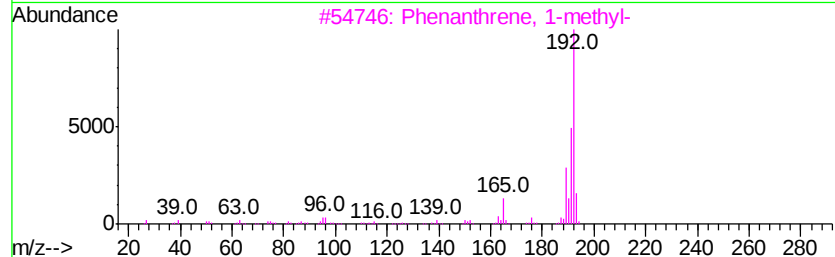
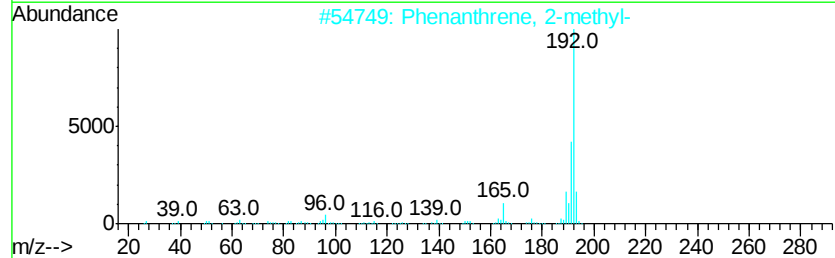
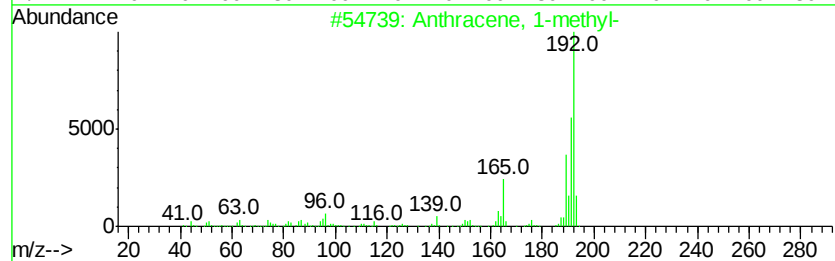
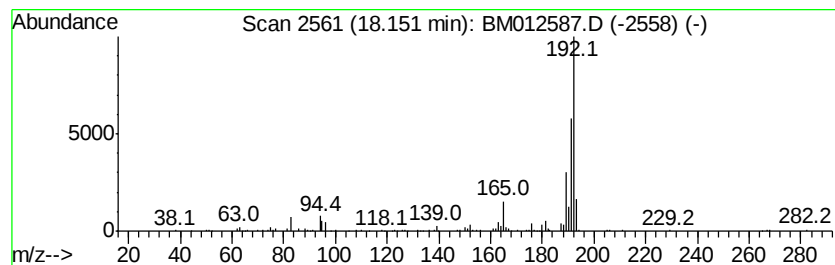
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Anthracene, 1-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.15	2.18 ng/ul	461329	Phenanthrene-d10	17.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Anthracene, 1-methyl-	192	C15H12	000610-48-0	93
2		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93
3		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	93
4		1H-Cyclopropa[1,1]phenanthrene, 1a,...	192	C15H12	000949-41-7	93
5		1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	93



Data Path : Z:\HPCHEM1\BNA M\DATA\BM103017\
Data File : BM012587.D
Acq On : 30 Oct 2017 23:26
Operator : SJ/JU
Sample : I5886-05
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ESNE5

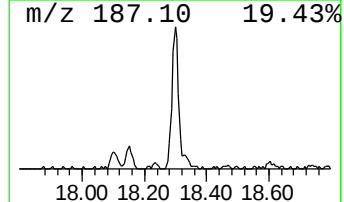
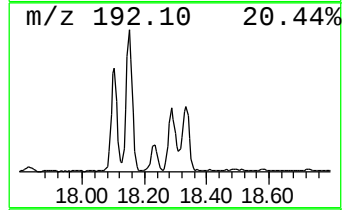
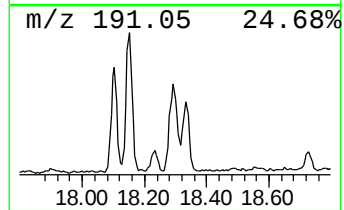
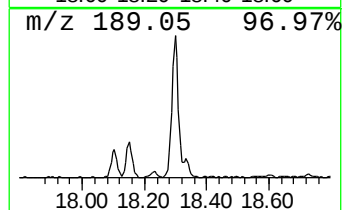
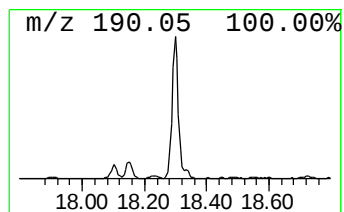
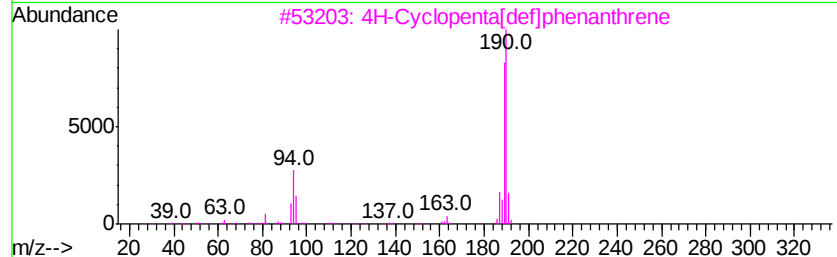
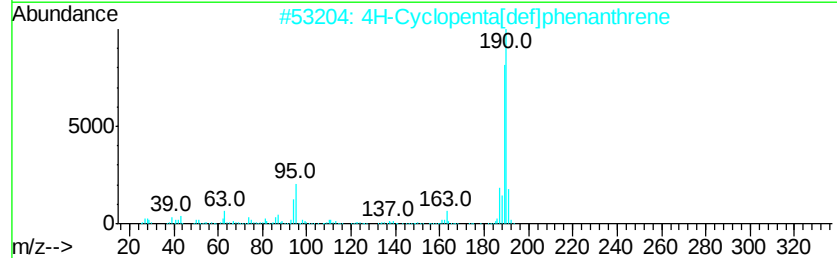
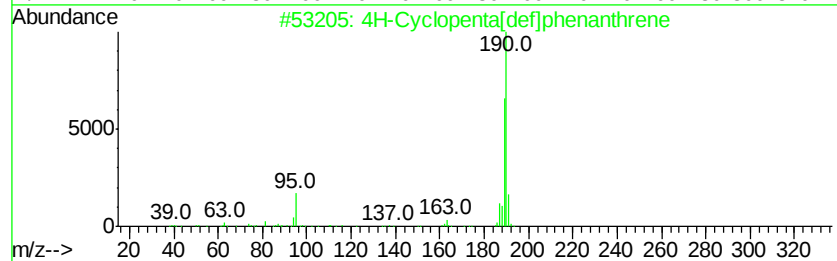
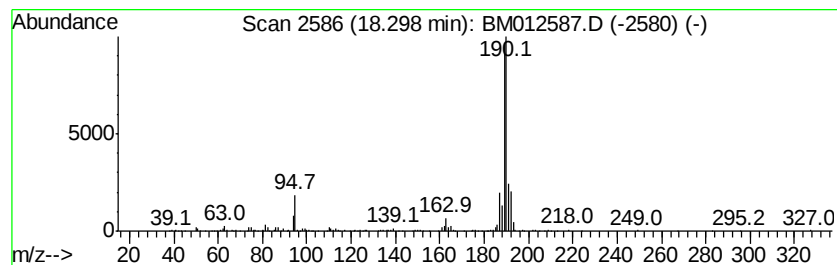
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 4H-Cyclopenta[def]phenanthrene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.30	3.29 ng/ul	698184	Phenanthrene-d10	17.23

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	94
2		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	90
3		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	87
4		2,3,5,6-Tetramethylterephthalald...	190	C12H14O2	007072-01-7	50
5		Pyrazole-4-carboxaldehyde, 3-(4-...	190	C10H7FN2O	306936-57-2	50



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ClientSampleId :
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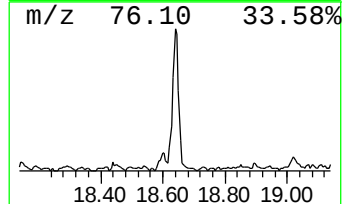
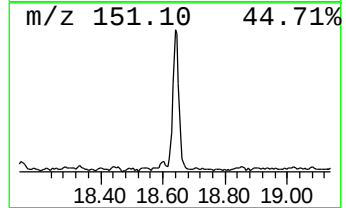
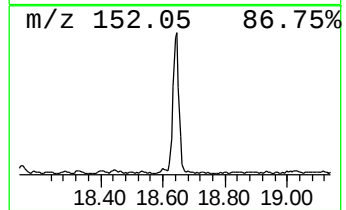
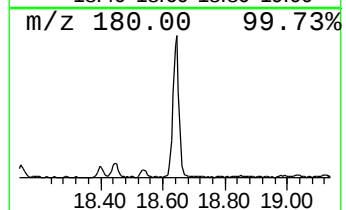
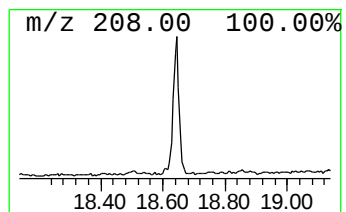
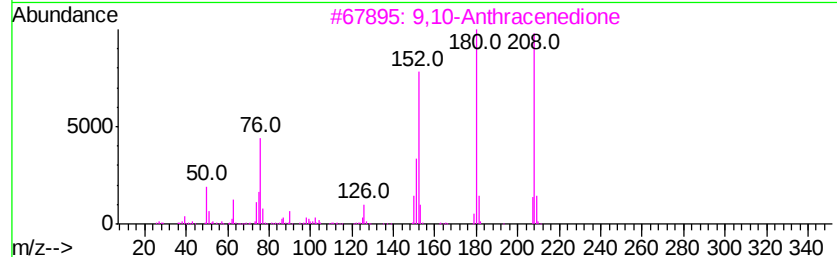
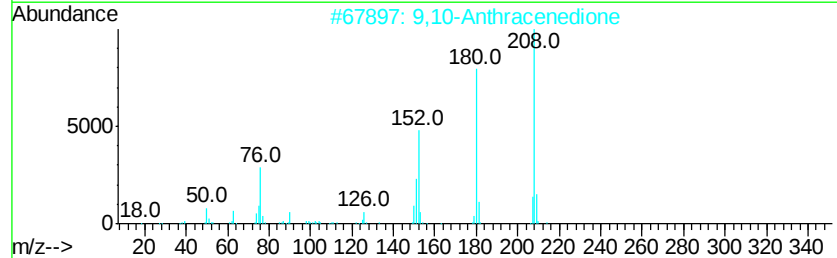
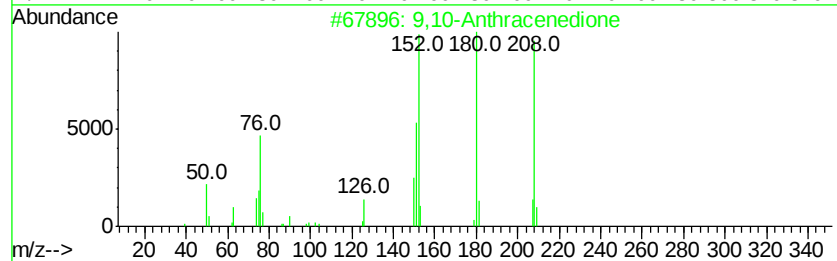
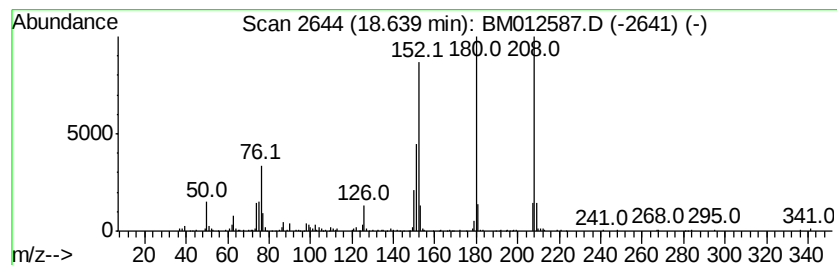
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 9,10-Anthracenedione Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.64	2.77 ng/ul	586633	Phenanthrene-d10	17.23

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	94
4		Benzo[cl]cinoline	180	C12H8N2	000230-17-1	91
5		1H-Phenalen-1-one	180	C13H8O	000548-39-0	64



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Sample : I5886-05
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ALS Vial : 19 Sample Multiplier: 1

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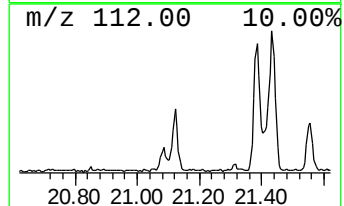
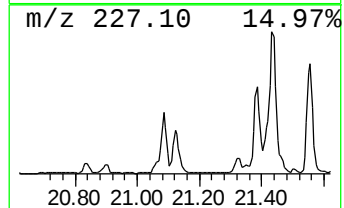
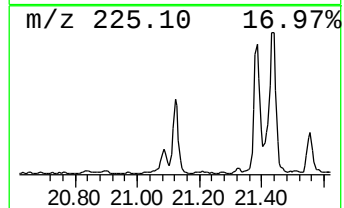
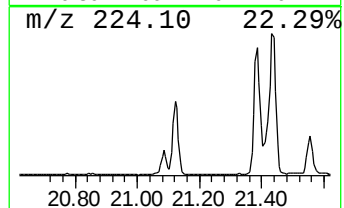
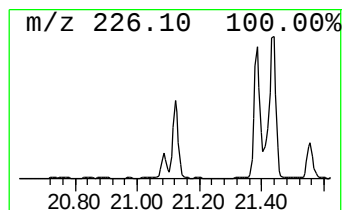
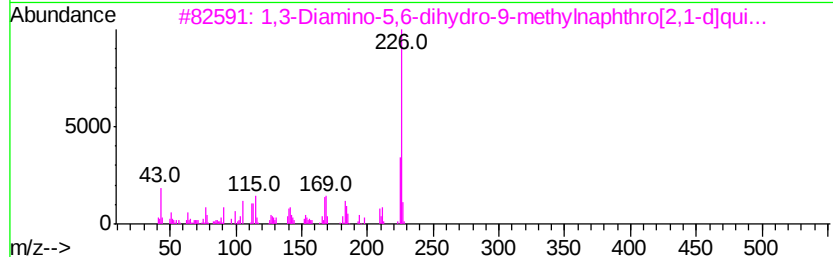
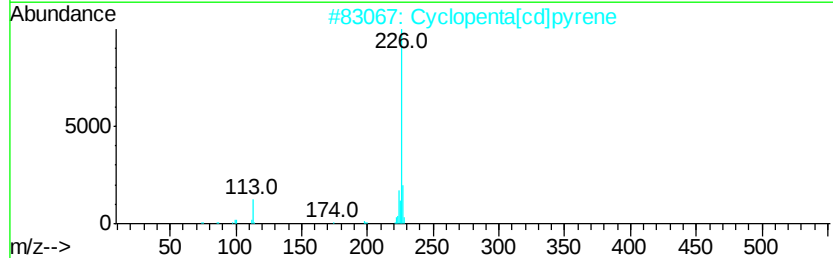
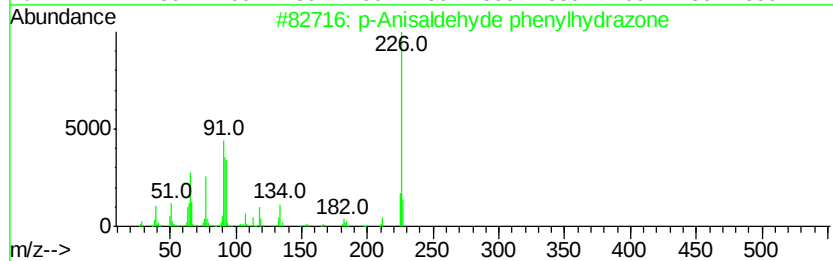
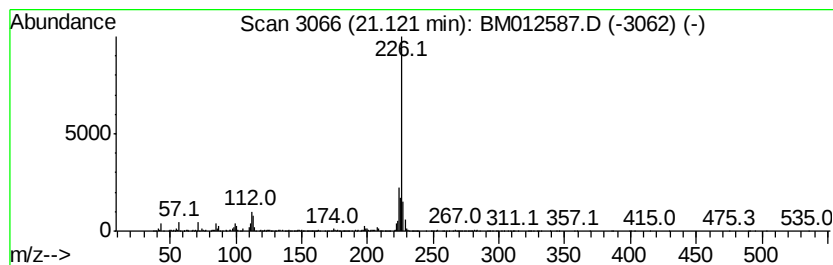
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 p-Anisaldehyde phenylhydrazone Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.12	2.63 ng/ul	1481280	Chrysene-d12	21.40

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	p-Anisaldehydve phenylhvdrazone	226	C14H14N2O	000622-73-1	52
2		Cyclopenta[cd]pyrene	226	C18H10	027208-37-3	52
3		1,3-Diamino-5,6-dihydro-9-methyl...	226	C13H14N4	037436-39-8	50
4		1,8-Diazacyclotetradecane-2,7-dione	226	C12H22N2O2	004266-66-4	50
5		2-Furfurylidene-7-hydroxy-1-inda...	226	C14H10O3	1000244-80-4	47



Data Path : Z:\HPCHEM1\BNA M\DATA\BM103017\
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ClientSampleId :
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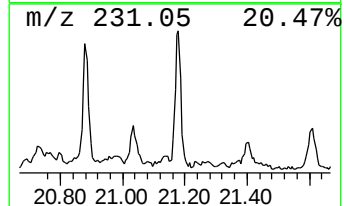
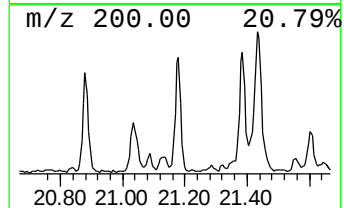
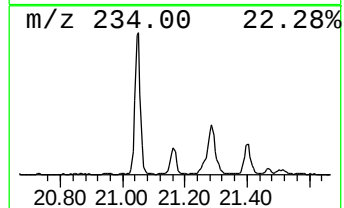
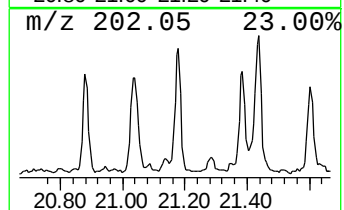
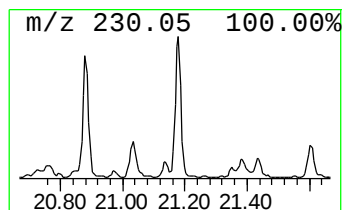
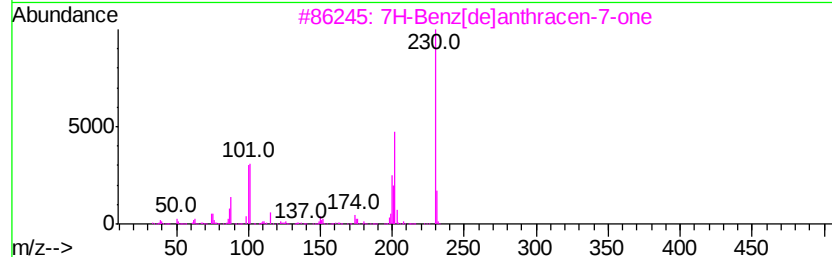
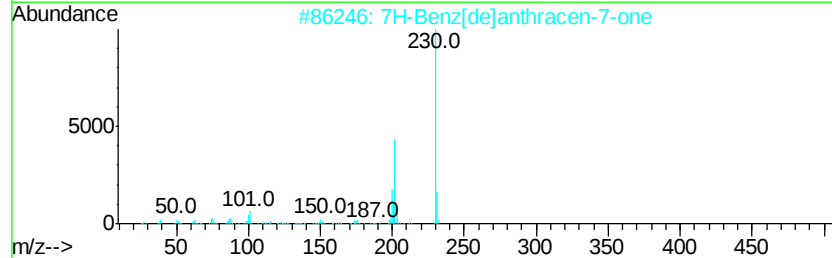
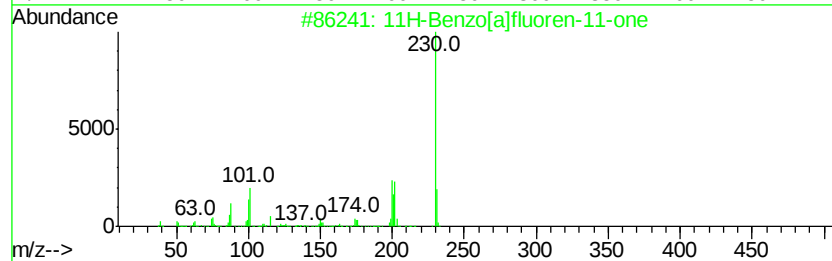
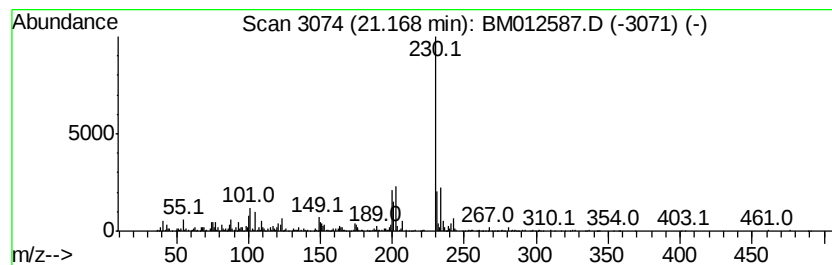
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 11H-Benzo[a]fluoren-11-one Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.17	2.24 ng/ul	1264380	Chrysene-d12	21.40

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	11H-Benzo[a]fluoren-11-one	230	C17H10O	000479-79-8	96
2		7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	90
3		7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	90
4		7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	83
5		(E)-.alpha.,.alpha.'-Dicyanostil...	230	C16H10N2	002450-55-7	80



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Data File : BM012587.D
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Operator : SJ/JU
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Misc :
ALS Vial : 19 Sample Multiplier: 1

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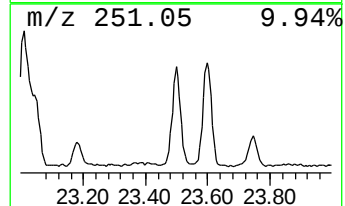
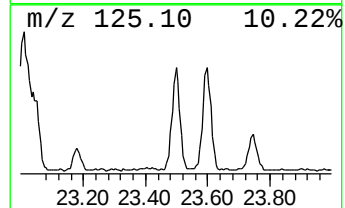
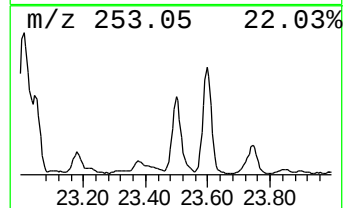
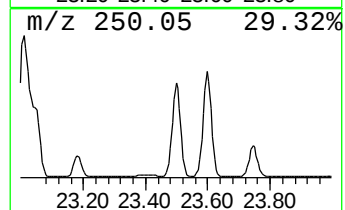
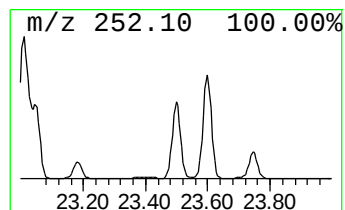
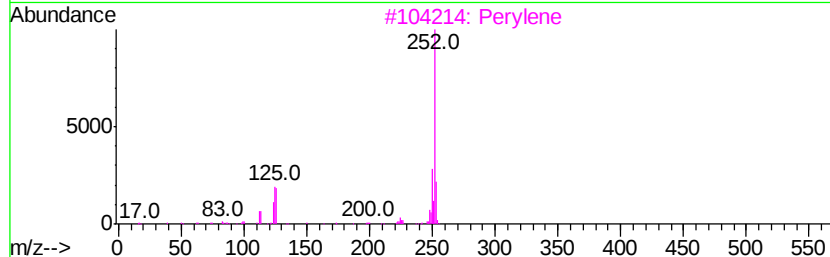
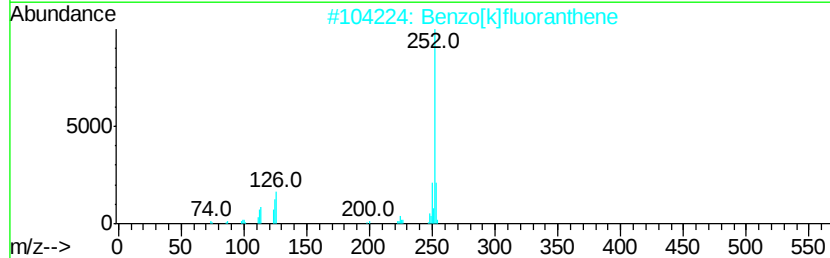
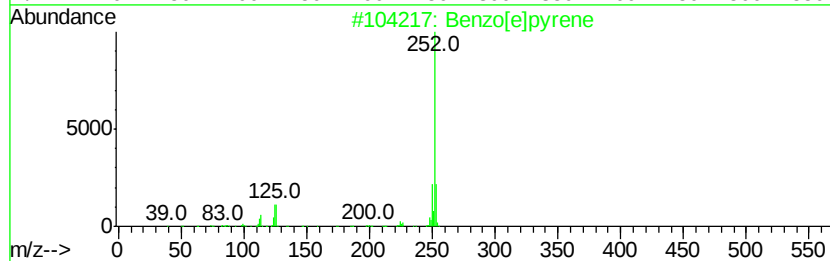
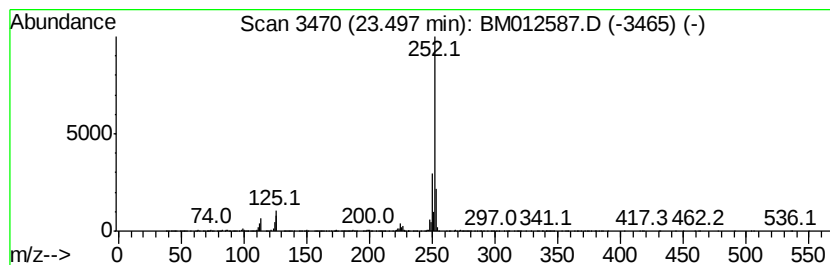
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Benzo[e]pyrene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.50	13.30 ng/ul	3585950	Perylene-d12	23.70

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	97
3		Perylene	252	C20H12	000198-55-0	94
4		Benzo[e]acephenanthrylene	252	C20H12	000205-99-2	93
5		Benzo[a]pyrene	252	C20H12	000050-32-8	91



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Quant Title : SVOA CALIBRATION

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

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					#	RT	Resp	Conc
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Benzophenone	15.84	3.3	ng/ul	595023	3	14.49	3566130	20.0
n-Hexadecanoic acid	18.12	3.9	ng/ul	835852	4	17.23	4242060	20.0
Anthracene, 1-met...	18.15	2.2	ng/ul	461329	4	17.23	4242060	20.0
4H-Cyclopenta[def...	18.30	3.3	ng/ul	698184	4	17.23	4242060	20.0
9,10-Anthracenedione	18.64	2.8	ng/ul	586633	4	17.23	4242060	20.0
p-Anisaldehyde ph...	21.12	2.6	ng/ul	1481280	5	21.40	11275700	20.0
11H-Benzo[a]fluor...	21.17	2.2	ng/ul	1264380	5	21.40	11275700	20.0
Benzo[e]pyrene	23.50	13.3	ng/ul	3585950	6	23.70	5393900	20.0