

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062415\
 Data File : BM001785.D
 Acq On : 24 Jun 2015 16:56
 Operator : TP/IZ
 Sample : G2740-05
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 H0FA7

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:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM061515.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	2.810	16	21	25	rVV	80709	115090	6.15%	0.489%
2	2.881	29	33	53	rVB	1400763	1655615	88.51%	7.031%
3	6.828	697	704	714	rBB	132883	220397	11.78%	0.936%
4	6.998	727	733	740	rBV	96710	143443	7.67%	0.609%
5	7.193	760	766	772	rBB	130009	195724	10.46%	0.831%
6	7.663	840	846	853	rVB	229240	351903	18.81%	1.494%
7	8.363	960	965	973	rVB	145498	224417	12.00%	0.953%
8	8.481	980	985	993	rVB	64576	101849	5.45%	0.433%
9	8.822	1036	1043	1050	rBV	120176	188804	10.09%	0.802%
10	9.539	1159	1165	1171	rBV	96172	155074	8.29%	0.659%
11	9.686	1183	1190	1198	rVB2	12642	23807	1.27%	0.101%
12	10.069	1249	1255	1266	rVB	160478	261227	13.97%	1.109%
13	10.451	1311	1320	1326	rBV	313275	514633	27.51%	2.186%
14	10.504	1326	1329	1337	rVB	21945	31226	1.67%	0.133%
15	10.592	1339	1344	1353	rBV	94074	154780	8.27%	0.657%
16	12.122	1599	1604	1610	rVB	12203	18861	1.01%	0.080%
17	12.345	1637	1642	1650	rVB2	12208	18867	1.01%	0.080%
18	12.916	1735	1739	1744	rVB	15985	18928	1.01%	0.080%
19	13.633	1854	1861	1866	rBV2	11286	20432	1.09%	0.087%
20	13.727	1872	1877	1882	rVV	263276	361788	19.34%	1.536%
21	13.774	1882	1885	1891	rVB	88579	113784	6.08%	0.483%
22	14.010	1919	1925	1937	rVB	296504	468536	25.05%	1.990%
23	14.321	1970	1978	1984	rBV2	447634	675938	36.14%	2.871%
24	14.380	1984	1988	1994	rVV	47283	65821	3.52%	0.280%
25	14.516	2005	2011	2020	rBV	115865	167149	8.94%	0.710%
26	14.716	2041	2045	2053	rVB	33048	46615	2.49%	0.198%
27	15.310	2141	2146	2151	rBV	364541	506131	27.06%	2.149%
28	15.368	2151	2156	2161	rVV2	49373	77295	4.13%	0.328%
29	15.433	2163	2167	2174	rBV	70220	95329	5.10%	0.405%
30	15.663	2202	2206	2211	rVB	14828	20369	1.09%	0.087%
31	15.810	2226	2231	2239	rVB	16147	29583	1.58%	0.126%
32	16.051	2265	2272	2277	rVB3	23896	46233	2.47%	0.196%
33	16.715	2381	2385	2392	rVB	30245	43725	2.34%	0.186%
34	16.886	2409	2414	2423	rVB	24870	39399	2.11%	0.167%

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Integration Parameters: LSCINT.P

Integrator: RTE
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35	17.062	2438	2444	2448	rBV	574573	834074	44.59%	3.542%
36	17.104	2448	2451	2456	rVV	573446	801294	42.84%	3.403%
37	17.162	2456	2461	2464	rVV	413975	564890	30.20%	2.399%
38	17.198	2464	2467	2472	rVB	142989	184457	9.86%	0.783%
39	17.257	2472	2477	2482	rBV3	13538	21459	1.15%	0.091%
40	17.445	2502	2509	2510	rBV2	21952	31853	1.70%	0.135%
41	17.468	2510	2513	2519	rVB	50714	62829	3.36%	0.267%
42	17.674	2546	2548	2553	rVB2	25176	29018	1.55%	0.123%
43	17.945	2584	2594	2598	rBV2	102196	231368	12.37%	0.983%
44	17.986	2598	2601	2607	rVB	125885	164308	8.78%	0.698%
45	18.062	2609	2614	2619	rVB	47076	65851	3.52%	0.280%
46	18.133	2619	2626	2630	rBV2	110521	207496	11.09%	0.881%
47	18.168	2630	2632	2637	rVB	71694	79771	4.26%	0.339%
48	18.445	2673	2679	2682	rBV	63230	82366	4.40%	0.350%
49	18.486	2682	2686	2691	rVB	36384	53716	2.87%	0.228%
50	18.539	2691	2695	2698	rBV2	18594	24993	1.34%	0.106%
51	18.692	2716	2721	2724	rBV2	24684	33548	1.79%	0.142%
52	18.739	2724	2729	2732	rBV3	11403	21857	1.17%	0.093%
53	18.774	2732	2735	2739	rVB3	24084	32266	1.73%	0.137%
54	18.862	2744	2750	2754	rBV3	52149	91446	4.89%	0.388%
55	18.927	2756	2761	2763	rVV3	29030	47109	2.52%	0.200%
56	18.956	2763	2766	2769	rVB	22853	28275	1.51%	0.120%
57	19.004	2769	2774	2782	rBV2	60236	112717	6.03%	0.479%
58	19.127	2789	2795	2800	rVB	917830	1173817	62.76%	4.985%
59	19.239	2808	2814	2818	rBV3	45740	84542	4.52%	0.359%
60	19.274	2818	2820	2824	rVB	21115	23736	1.27%	0.101%
61	19.321	2824	2828	2830	rBV5	20106	28144	1.50%	0.120%
62	19.462	2846	2852	2854	rBV2	573142	821362	43.91%	3.488%
63	19.492	2854	2857	2865	rVB	910730	1203108	64.32%	5.109%
64	19.562	2865	2869	2875	rVB2	68027	99018	5.29%	0.421%
65	19.668	2882	2887	2893	rVB	57302	80846	4.32%	0.343%
66	19.727	2893	2897	2902	rVB2	37426	52002	2.78%	0.221%
67	19.815	2906	2912	2918	rBV4	69745	166535	8.90%	0.707%
68	19.951	2930	2935	2938	rBV5	51064	96530	5.16%	0.410%
69	19.986	2938	2941	2948	rVB3	92653	149123	7.97%	0.633%
70	20.092	2955	2959	2962	rBV	51237	64378	3.44%	0.273%
71	20.145	2962	2968	2973	rVV2	108252	186520	9.97%	0.792%

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72	20.192	2973	2976	2979	rVB3	24709	26960	1.44%	0.114%
73	20.227	2979	2982	2986	rBV2	29807	42718	2.28%	0.181%
74	20.286	2989	2992	2995	rVV	57142	68332	3.65%	0.290%
75	20.333	2995	3000	3004	rVB	64776	88898	4.75%	0.378%
76	20.509	3027	3030	3032	rVB3	23228	23804	1.27%	0.101%
77	20.574	3039	3041	3045	rVB4	20587	28905	1.55%	0.123%
78	20.650	3052	3054	3059	rVB6	19799	21433	1.15%	0.091%
79	20.692	3059	3061	3065	rBV5	13500	23825	1.27%	0.101%
80	20.739	3065	3069	3075	rVV	102903	137512	7.35%	0.584%
81	20.892	3091	3095	3096	rBV2	45353	57290	3.06%	0.243%
82	20.945	3100	3104	3106	rVV	85877	128672	6.88%	0.546%
83	20.974	3106	3109	3115	rVV3	161523	301627	16.13%	1.281%
84	21.033	3115	3119	3128	rVB2	92837	168733	9.02%	0.717%
85	21.174	3140	3143	3146	rVB	51293	54011	2.89%	0.229%
86	21.256	3146	3157	3161	rBV2	976857	1870461	100.00%	7.943%
87	21.297	3161	3164	3174	rVB	455671	605407	32.37%	2.571%
88	21.409	3180	3183	3189	rBV	101235	129002	6.90%	0.548%
89	21.574	3206	3211	3215	rBV2	66210	112410	6.01%	0.477%
90	21.803	3247	3250	3253	rVB	97452	106163	5.68%	0.451%
91	21.997	3280	3283	3286	rBV2	51455	64713	3.46%	0.275%
92	22.815	3415	3422	3434	rBV	304603	985102	52.67%	4.183%
93	22.986	3447	3451	3456	rVB	64360	96593	5.16%	0.410%
94	23.291	3498	3503	3507	rBV	172549	308628	16.50%	1.311%
95	23.339	3507	3511	3515	rVV2	250886	439460	23.49%	1.866%
96	23.386	3515	3519	3526	rVB	318884	545117	29.14%	2.315%
97	23.480	3529	3535	3541	rBV	522743	998341	53.37%	4.240%
98	25.715	3909	3915	3925	rVB2	152042	438307	23.43%	1.861%
99	26.068	3971	3975	3984	rVB2	78399	178927	9.57%	0.760%
100	26.403	4024	4032	4042	rVB2	93583	286787	15.33%	1.218%

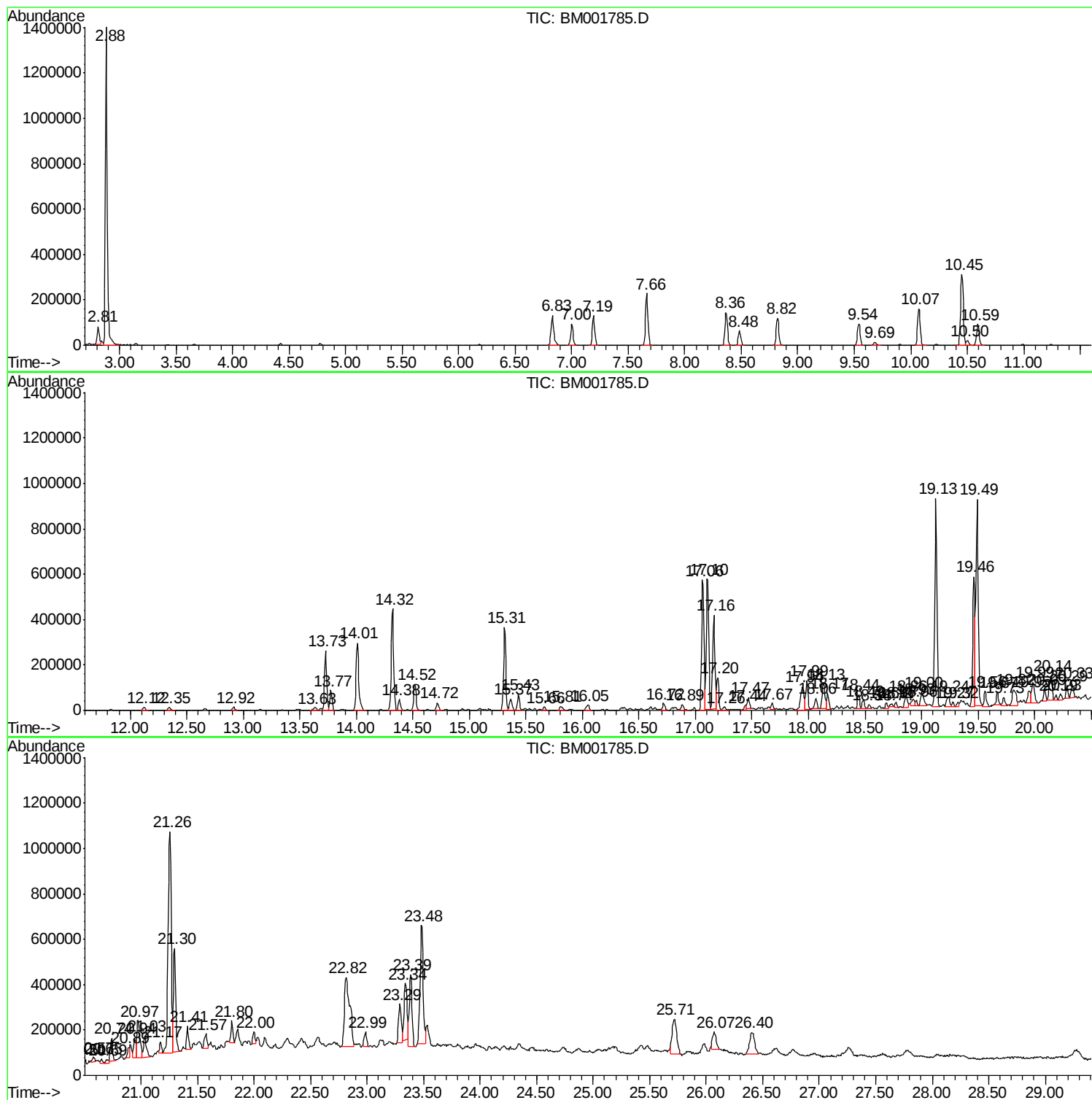
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM061515.M
Quant Title : SVOA CALIBRATION

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



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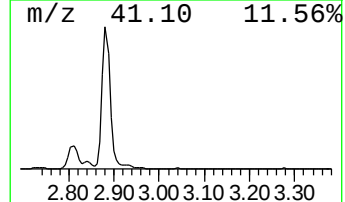
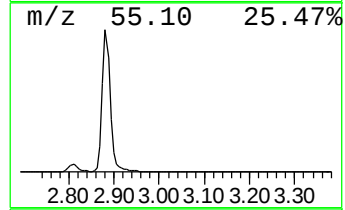
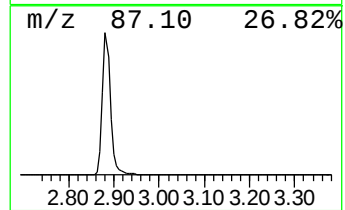
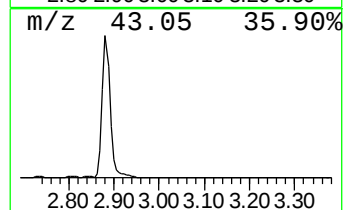
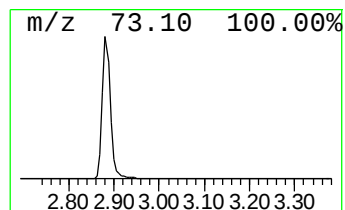
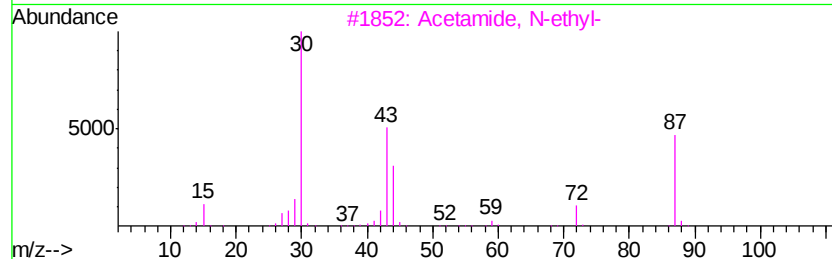
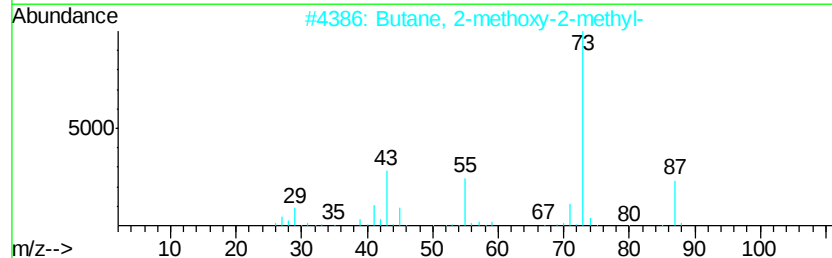
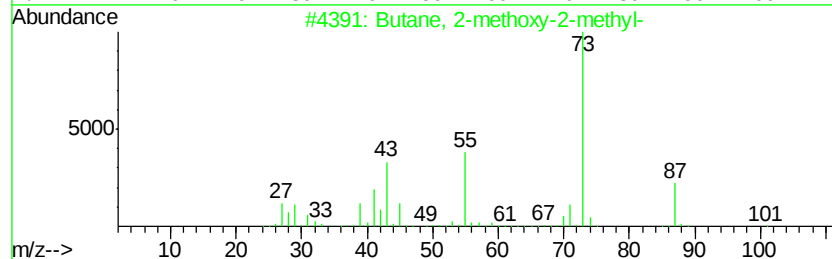
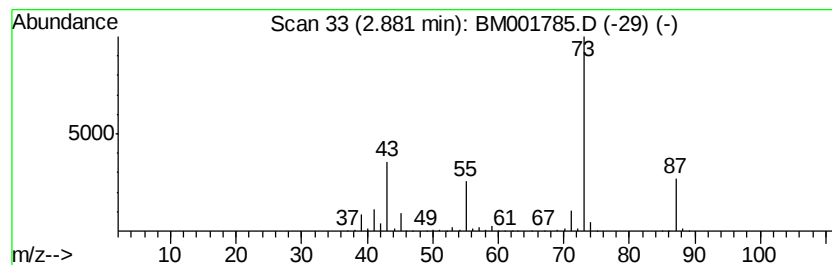
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM061515.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.88	94.10 ng/ul	1655620	1,4-Dichlorobenzene-d4	7.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
3		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	22
4		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17
5		N-Ethylformamide	73	C3H7NO	000627-45-2	9



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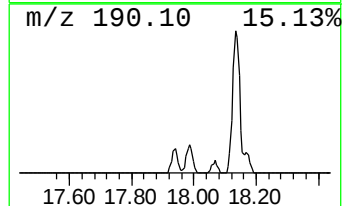
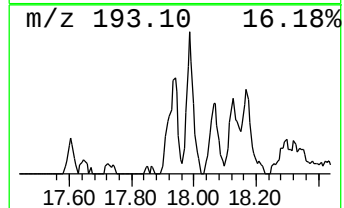
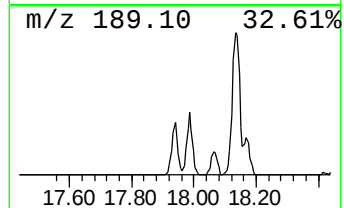
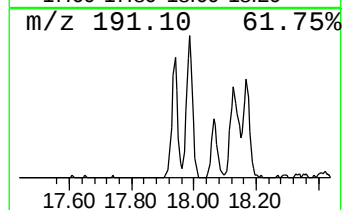
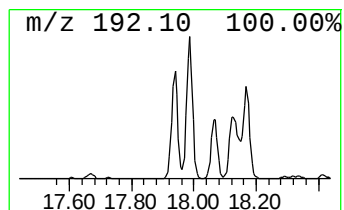
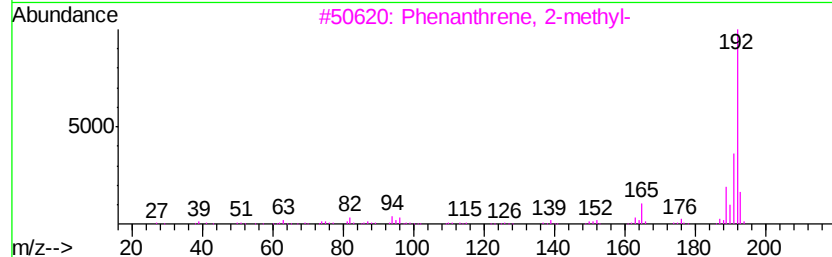
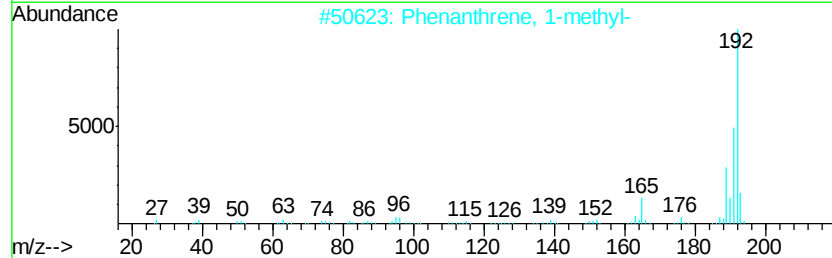
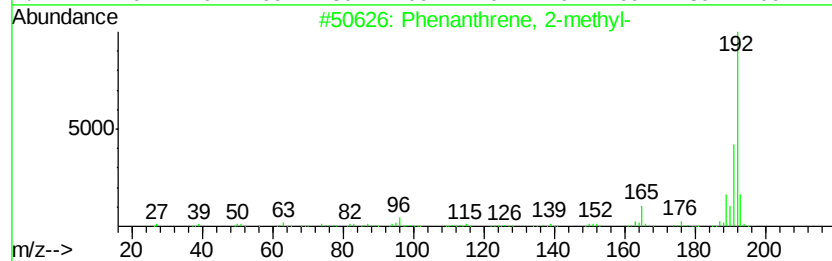
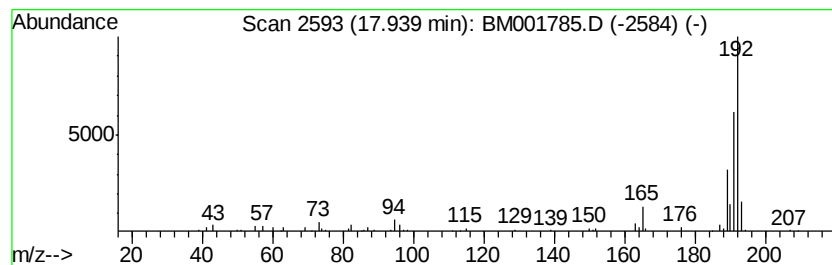
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.94	5.55 ng/ul	231368	Phenanthrene-d10	17.06	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	97
2	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	97
3	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
4	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
5	Anthracene, 2-methyl-	192	C15H12	000613-12-7	95



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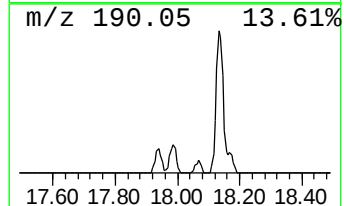
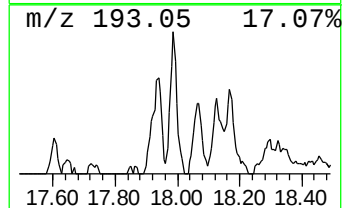
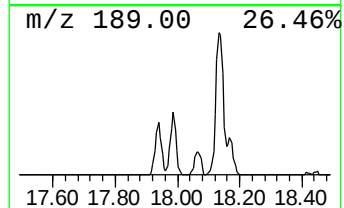
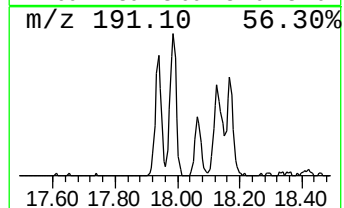
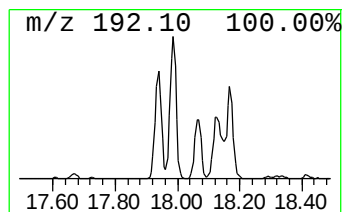
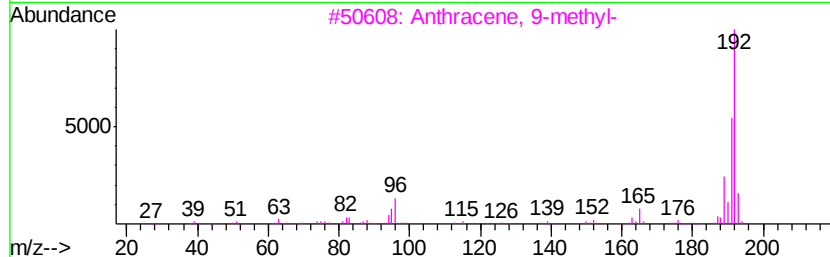
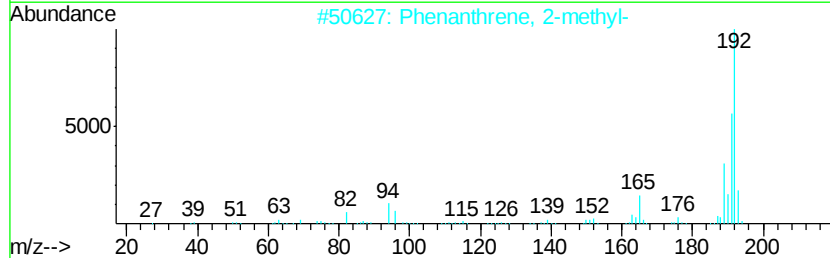
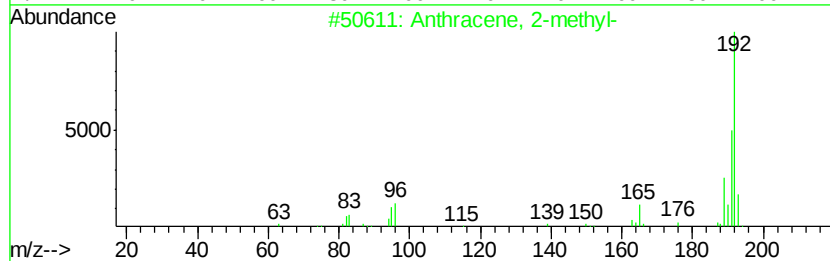
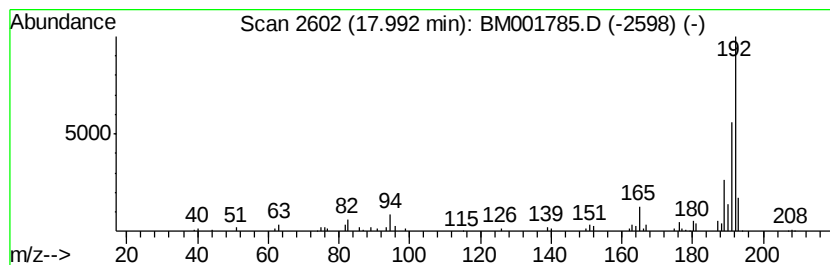
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TIC Integration Parameters: LSCINT.P

Peak Number 4 Anthracene, 2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.99	3.94 ng/ul	164308	Phenanthrene-d10	17.06

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Anthracene, 2-methyl-	192	C15H12	000613-12-7	94
2		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93
3		Anthracene, 9-methyl-	192	C15H12	000779-02-2	91
4		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	91
5		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	91



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062415\
Data File : BM001785.D
Acq On : 24 Jun 2015 16:56
Operator : TP/IZ
Sample : G2740-05
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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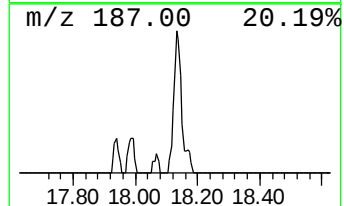
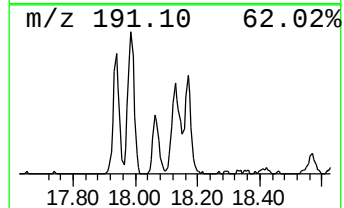
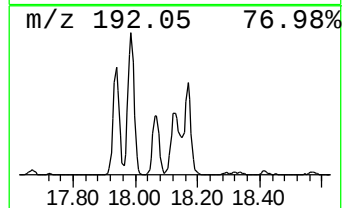
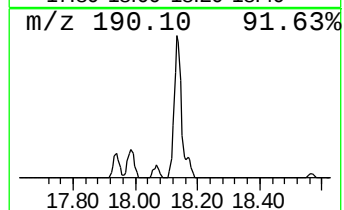
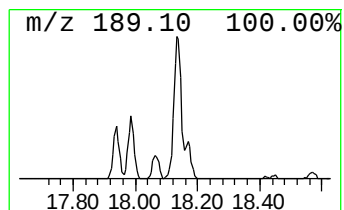
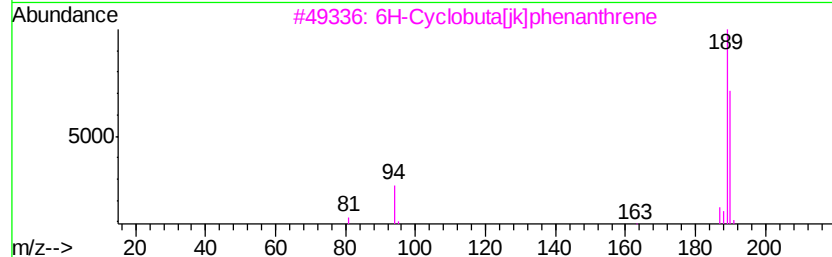
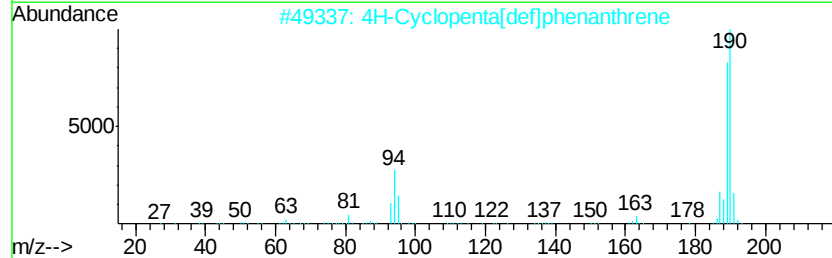
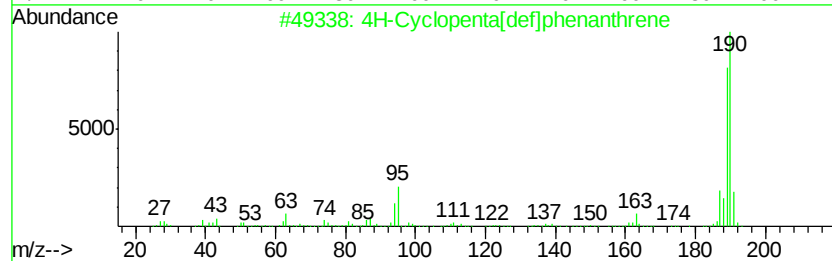
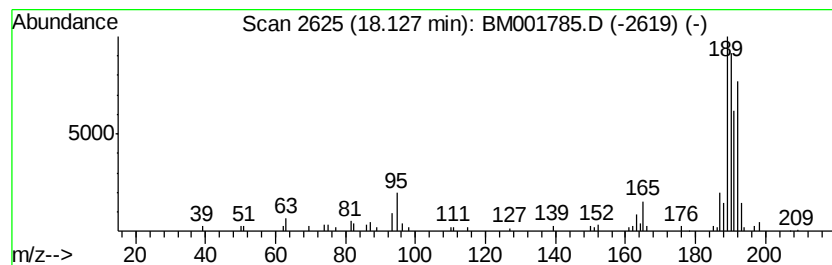
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.13	4.98 ng/ul	207496	Phenanthrene-d10	17.06

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	58
2		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	52
3		6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	52
4		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	50
5		Thiophene-3-carboxaldehyde, 5-ch...	190	C5H4BClO3S	036155-87-0	50



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062415\
Data File : BM001785.D
Acq On : 24 Jun 2015 16:56
Operator : TP/IZ
Sample : G2740-05
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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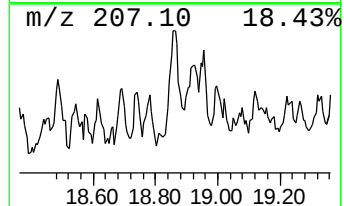
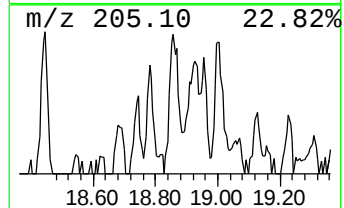
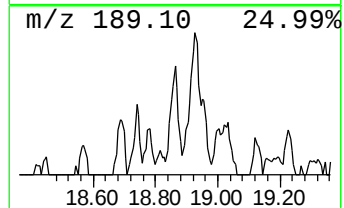
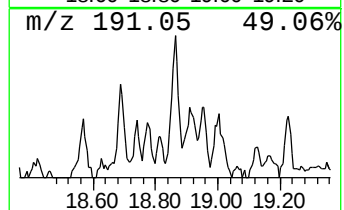
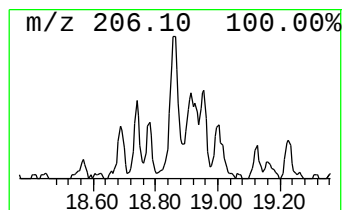
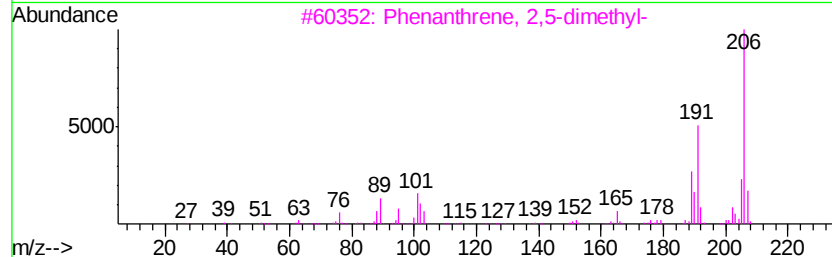
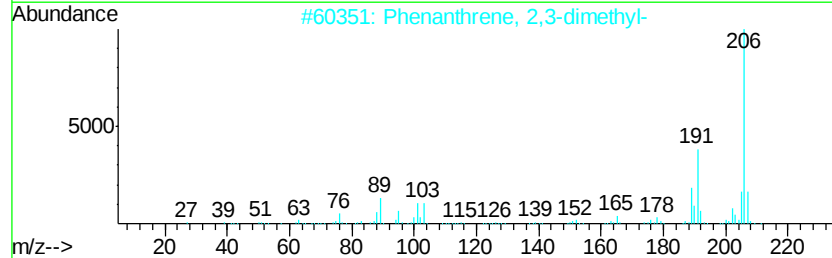
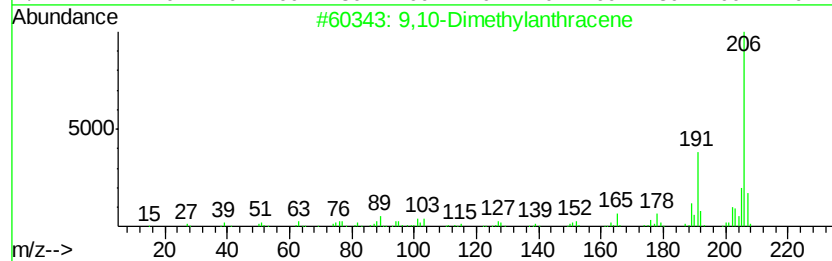
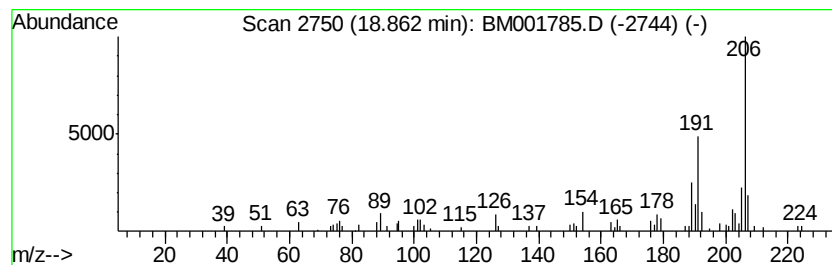
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 9,10-Dimethylantracene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.86	2.19 ng/ul	91446	Phenanthrene-d10	17.06

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,10-Dimethylantracene	206	C16H14	000781-43-1	93	
2	Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	93	
3	Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	91	
4	9,10-Dimethylantracene	206	C16H14	000781-43-1	90	
5	di-p-Tolylacetylene	206	C16H14	002789-88-0	90	



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062415\
Data File : BM001785.D
Acq On : 24 Jun 2015 16:56
Operator : TP/IZ
Sample : G2740-05
Misc :
ALS Vial : 8 Sample Multiplier: 1

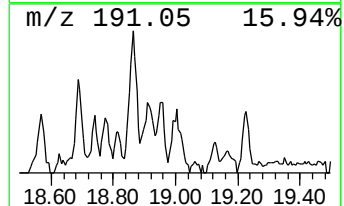
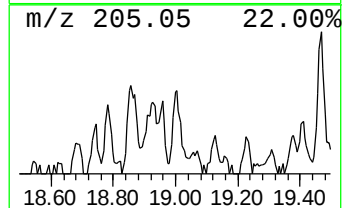
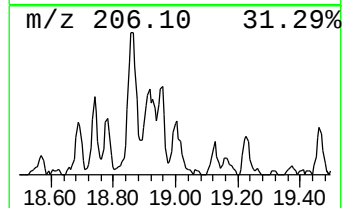
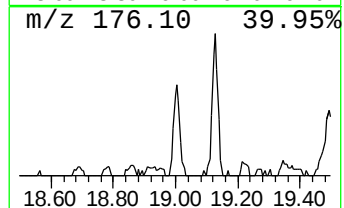
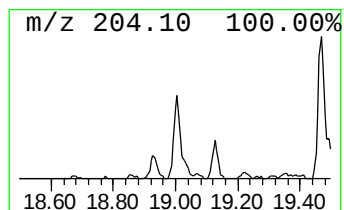
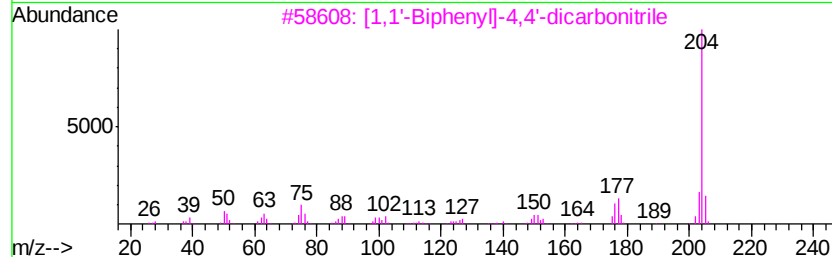
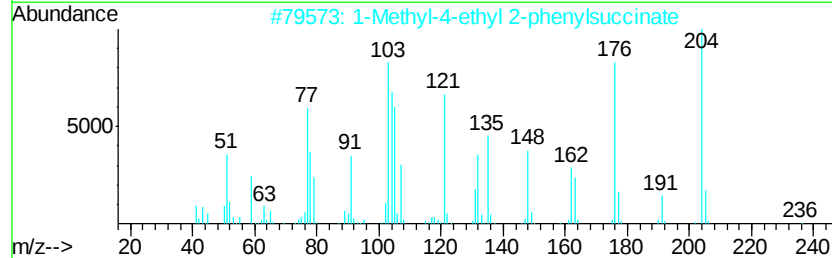
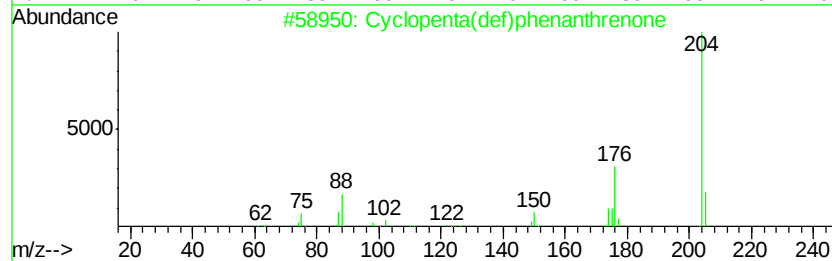
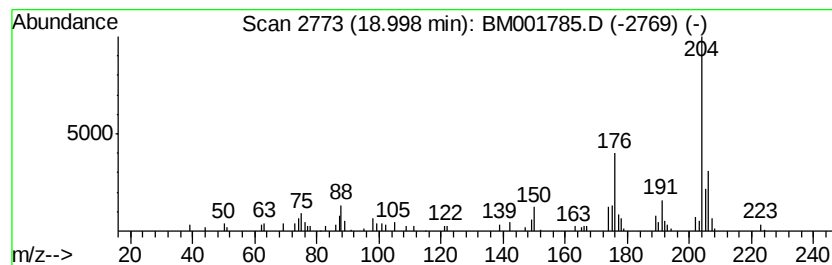
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM061515.M
Quant Title : SVOA CALIBRATION

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

Peak Number 7 Cyclopenta(def)phenanthrenone Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.		
19.00	2.70 ng/ul	112717	Phenanthrene-d10	17.06		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	81
2		1-Methyl-4-ethyl 2-phenylsuccinate	236	C13H16O4	132545-36-9	50
3		[1,1'-Biphenyl]-4,4'-dicarbonitrile	204	C14H8N2	001591-30-6	46
4		Benzene, 1-chloro-4-phenoxy-	204	C12H9ClO	007005-72-3	46
5		[1,1'-Biphenyl-4-ol], 2-chloro-	204	C12H9ClO	023719-22-4	43



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062415\
Data File : BM001785.D
Acq On : 24 Jun 2015 16:56
Operator : TP/IZ
Sample : G2740-05
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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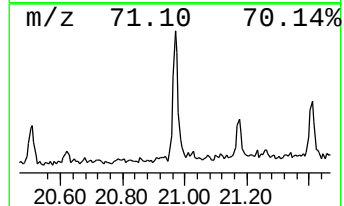
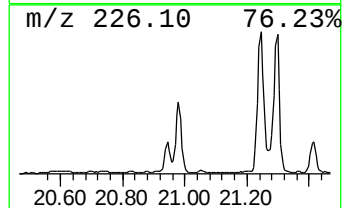
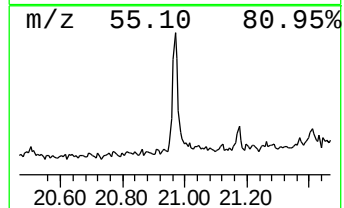
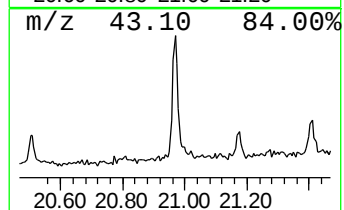
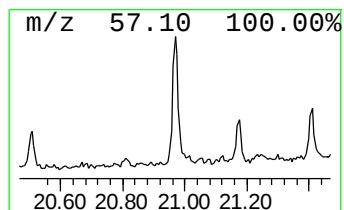
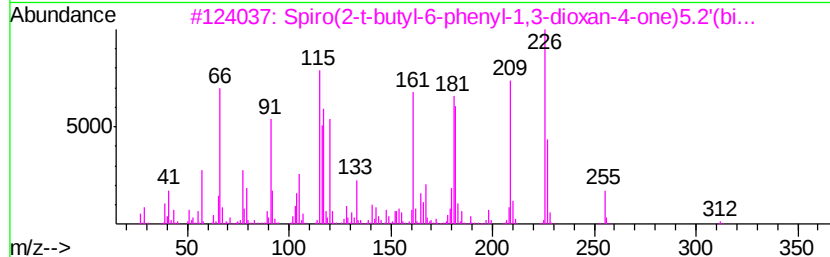
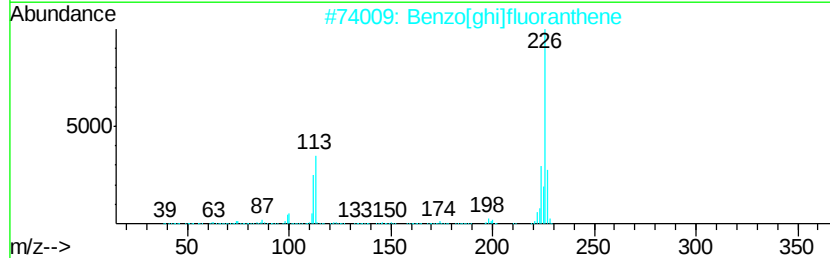
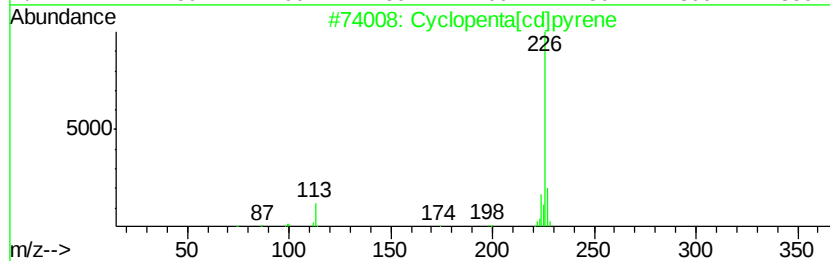
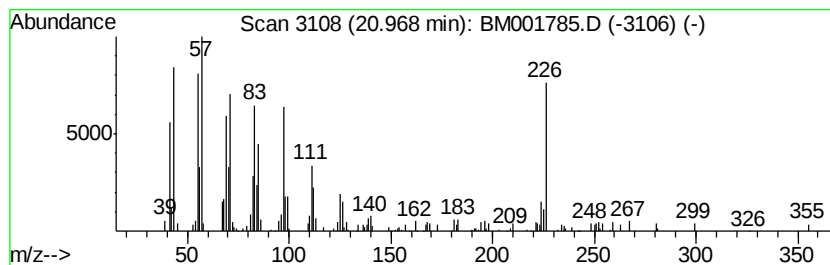
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 unknown-01 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.97	3.23 ng/ul	301627	Chrysene-d12	21.26

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopenta[cd]pyrene	226	C18H10	027208-37-3	47
2		Benzo[ghi]fluoranthene	226	C18H10	000203-12-3	28
3		Spiro(2-t-butyl-6-phenyl-1,3-dio...	312	C20H24O3	1000191-40-7	25
4		1,3-Diamino-5,6-dihydro-9-methyl...	226	C13H14N4	037436-39-8	23
5		5(10H)-Pyrido[3,4-b]quinolone, 7...	226	C13H10N2O2	1000256-99-5	23



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062415\
Data File : BM001785.D
Acq On : 24 Jun 2015 16:56
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Sample : G2740-05
Misc :
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ClientSampleId :
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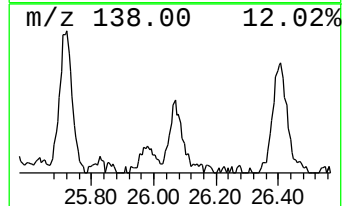
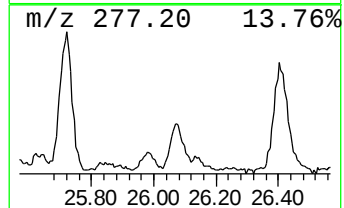
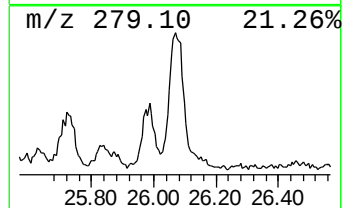
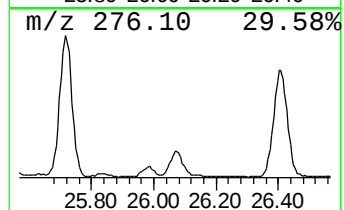
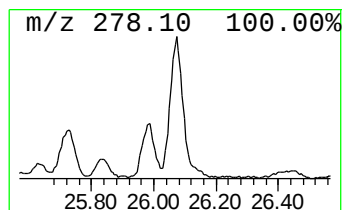
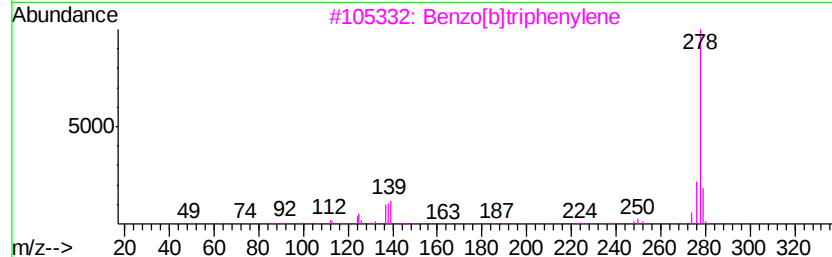
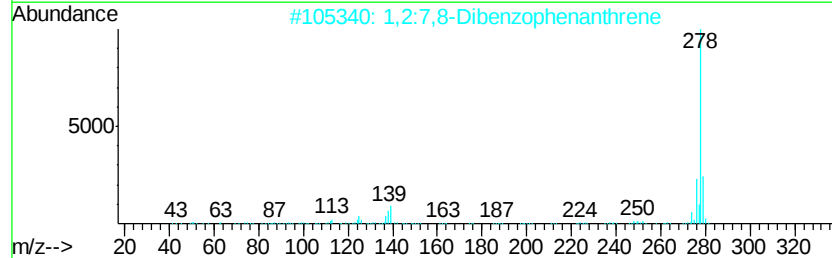
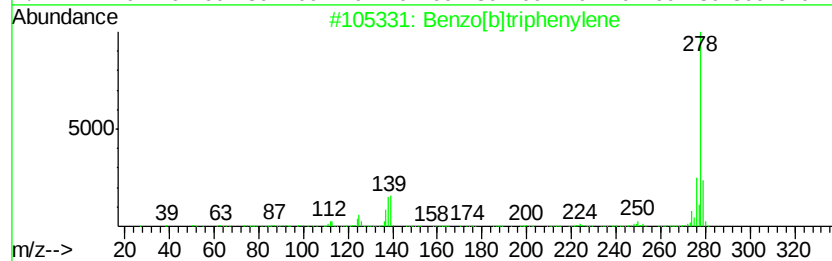
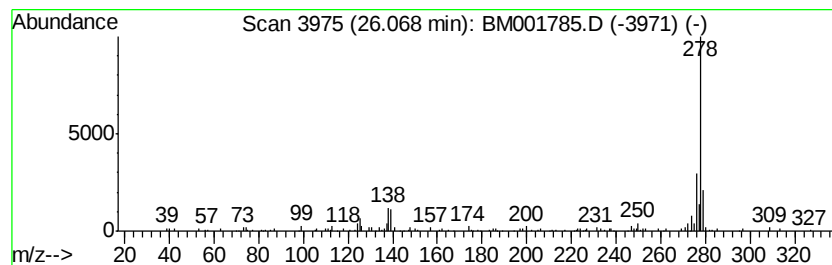
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Cyclic-01 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.07	3.58 ng/ul	178927	Perylene-d12	23.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[b]triphenylene	278	C22H14	000215-58-7	94
2		1,2:7,8-Dibenzophenanthrene	278	C22H14	000213-46-7	90
3		Benzo[b]triphenylene	278	C22H14	000215-58-7	86
4		Dibenz[a,h]anthracene	278	C22H14	000053-70-3	86
5		Benzo[b]triphenylene	278	C22H14	000215-58-7	81



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062415\
Data File : BM001785.D
Acq On : 24 Jun 2015 16:56
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Sample : G2740-05
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM061515.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
Butane, 2-methoxy...	2.88	94.1	ng/ul	1655620	1	7.66	351903	20.0
Phenanthrene, 2-m...	17.94	5.5	ng/ul	231368	4	17.06	834074	20.0
Anthracene, 2-met...	17.99	3.9	ng/ul	164308	4	17.06	834074	20.0
4H-Cyclopenta[def...	18.13	5.0	ng/ul	207496	4	17.06	834074	20.0
9,10-Dimethylanthe...	18.86	2.2	ng/ul	91446	4	17.06	834074	20.0
Cyclopenta(def)ph...	19.00	2.7	ng/ul	112717	4	17.06	834074	20.0
unknown-01	20.97	3.2	ng/ul	301627	5	21.26	1870460	20.0
(DEL) Alkane: Cyc...	26.07	3.6	ng/ul	178927	6	23.49	998341	20.0