

Data Path : Z:\HPCHEM1\BNA M\DATA\BM042018\
 Data File : BM014771.D
 Acq On : 20 Apr 2018 19:54
 Operator : SJ/JU
 Sample : J2434-12
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AC7

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-BM041918.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	3.316	3	5	10	rVB	372609	427170	2.26%	0.203%
2	3.369	10	14	20	rVV	901199	1172915	6.21%	0.558%
3	3.422	20	23	34	rVB2	136133	199679	1.06%	0.095%
4	4.263	161	166	175	rVB	323999	452733	2.40%	0.215%
5	5.516	373	379	388	rVB2	140706	252487	1.34%	0.120%
6	7.663	737	744	758	rBV2	1028669	2028763	10.74%	0.965%
7	7.846	766	775	784	rBV	892898	1432692	7.59%	0.681%
8	8.063	803	812	819	rBV	1091513	1825045	9.67%	0.868%
9	8.545	887	894	906	rVB	1097919	1840360	9.75%	0.875%
10	9.240	1005	1012	1021	rBV	1028613	1685312	8.93%	0.802%
11	9.728	1087	1095	1102	rBV	1042229	1841783	9.75%	0.876%
12	10.457	1210	1219	1227	rBV	820965	1413808	7.49%	0.672%
13	10.992	1303	1310	1319	rBV	1304313	2194230	11.62%	1.044%
14	11.386	1370	1377	1384	rBV	1447106	2529407	13.40%	1.203%
15	11.510	1391	1398	1416	rBV	1017155	1894705	10.03%	0.901%
16	14.533	1905	1912	1916	rBV	2172473	3052110	16.16%	1.452%
17	14.580	1916	1920	1925	rVB	598079	825082	4.37%	0.392%
18	14.845	1958	1965	1977	rBV	2645792	3957383	20.96%	1.882%
19	14.992	1985	1990	1996	rVB	186984	233754	1.24%	0.111%
20	15.145	2005	2016	2022	rBV2	1985160	3183930	16.86%	1.514%
21	15.204	2022	2026	2035	rVB	171777	258987	1.37%	0.123%
22	15.298	2035	2042	2050	rBV	536193	827113	4.38%	0.393%
23	15.927	2145	2149	2155	rVB	224193	314172	1.66%	0.149%
24	16.121	2175	2182	2188	rBV	3175799	4615493	24.44%	2.195%
25	16.216	2194	2198	2204	rVB	265246	369310	1.96%	0.176%
26	16.333	2209	2218	2230	rVB4	84749	250106	1.32%	0.119%
27	16.463	2236	2240	2245	rVB2	148380	210507	1.11%	0.100%
28	16.780	2290	2294	2310	rVB	214533	436808	2.31%	0.208%
29	17.857	2471	2477	2481	rBV	2360519	3525845	18.67%	1.677%
30	17.898	2481	2484	2488	rVB	1305024	1695263	8.98%	0.806%
31	17.951	2488	2493	2498	rBV	3223279	4608150	24.41%	2.192%
32	18.245	2536	2543	2548	rBV2	285869	483663	2.56%	0.230%
33	18.568	2593	2598	2604	rBV3	131526	213166	1.13%	0.101%
34	18.680	2612	2617	2620	rBV	522493	713973	3.78%	0.340%

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35	18.710	2620	2622	2626	rVV	212596	242675	1.29%	0.115%
36	18.757	2626	2630	2636	rVB2	341644	504821	2.67%	0.240%
37	18.910	2649	2656	2659	rBV3	391407	731965	3.88%	0.348%
38	19.592	2768	2772	2778	rBV2	105314	216961	1.15%	0.103%
39	19.862	2812	2818	2823	rVB	6370468	8793770	46.57%	4.183%
40	19.915	2823	2827	2830	rBV2	181339	231619	1.23%	0.110%
41	20.192	2867	2874	2876	rBV2	4379961	7308895	38.71%	3.477%
42	20.215	2876	2878	2885	rVV	6818214	8606811	45.58%	4.094%
43	20.274	2885	2888	2893	rVB	248673	327297	1.73%	0.156%
44	20.386	2903	2907	2911	rBV	363656	463215	2.45%	0.220%
45	20.451	2914	2918	2924	rVB	536066	730129	3.87%	0.347%
46	20.521	2926	2930	2936	rBV3	456862	961371	5.09%	0.457%
47	20.580	2936	2940	2944	rBV	231843	291461	1.54%	0.139%
48	20.692	2946	2959	2963	rBV2	1645739	3641805	19.29%	1.732%
49	20.786	2971	2975	2979	rBV	1188903	1523625	8.07%	0.725%
50	20.851	2983	2986	2989	rVV	594474	718213	3.80%	0.342%
51	20.880	2989	2991	2995	rVB	242833	242427	1.28%	0.115%
52	20.986	3005	3009	3012	rBV	714400	939493	4.98%	0.447%
53	21.027	3013	3016	3023	rVB	699019	955544	5.06%	0.455%
54	21.421	3080	3083	3088	rVB3	416793	569317	3.02%	0.271%
55	21.592	3107	3112	3114	rBV2	1191789	1669188	8.84%	0.794%
56	21.621	3114	3117	3120	rVB	780538	950731	5.04%	0.452%
57	21.668	3120	3125	3128	rBV2	1178654	1730386	9.16%	0.823%
58	21.927	3164	3169	3175	rBV2	7772575	14539867	77.01%	6.916%
59	21.980	3175	3178	3183	rVB	6291353	8381534	44.39%	3.987%
60	22.109	3197	3200	3207	rVB	1097729	1555903	8.24%	0.740%
61	22.221	3216	3219	3223	rVB	831920	1087815	5.76%	0.517%
62	22.274	3224	3228	3234	rBV	1417752	2288516	12.12%	1.089%
63	22.533	3270	3272	3278	rVB	1082506	1411660	7.48%	0.671%
64	22.756	3307	3310	3314	rBV3	541421	915903	4.85%	0.436%
65	23.686	3460	3468	3473	rBV	7283664	18881315	100.00%	8.981%
66	23.874	3495	3500	3505	rVB	1005777	1718316	9.10%	0.817%
67	24.227	3554	3560	3565	rBV	4522615	9314097	49.33%	4.430%
68	24.286	3566	3570	3573	rVV2	2503172	5025671	26.62%	2.390%
69	24.339	3573	3579	3590	rVV	8172851	18605124	98.54%	8.850%
70	24.444	3593	3597	3601	rVV	1883659	3822067	20.24%	1.818%
71	24.497	3602	3606	3614	rVB2	2185249	4694343	24.86%	2.233%

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72	27.044	4030	4039	4050	rVB	4107631	13169133	69.75%	6.264%
73	27.850	4167	4176	4195	rVB	3209612	11507654	60.95%	5.474%

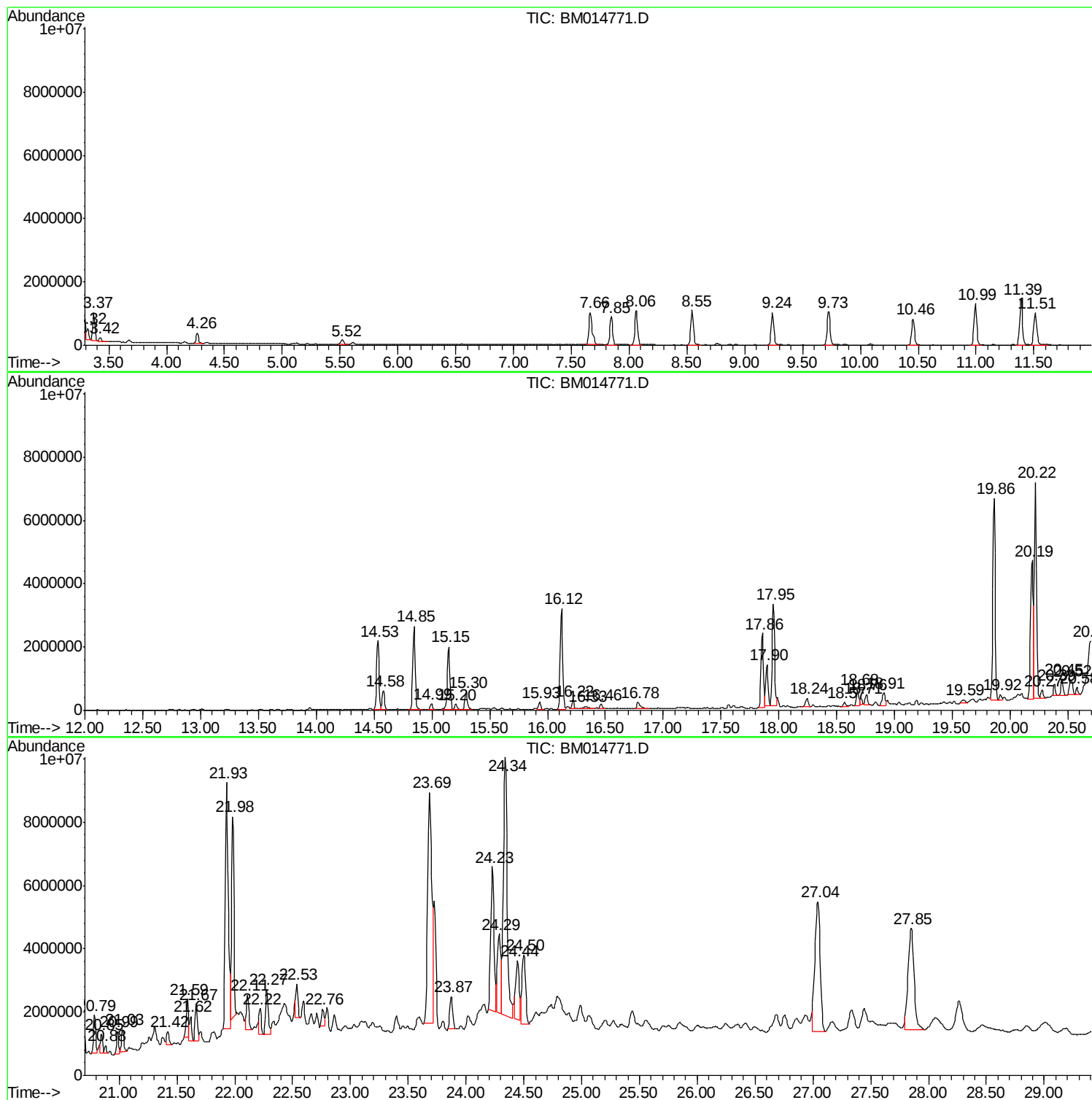
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Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM041918.M
Quant Title : SVOA CALIBRATION

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



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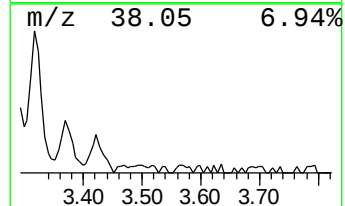
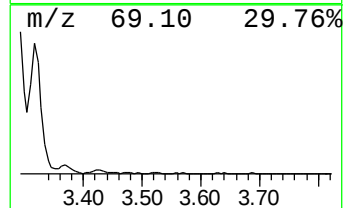
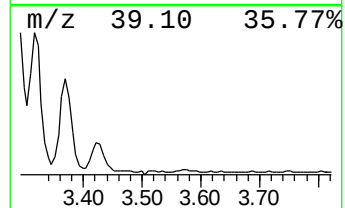
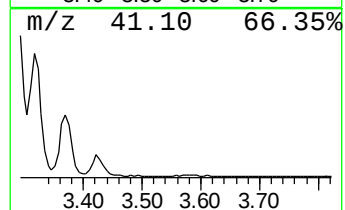
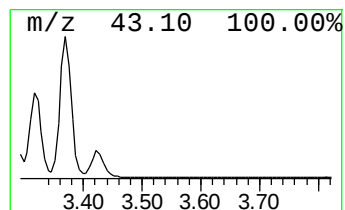
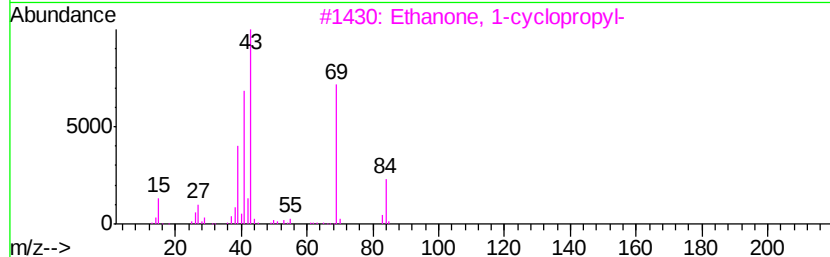
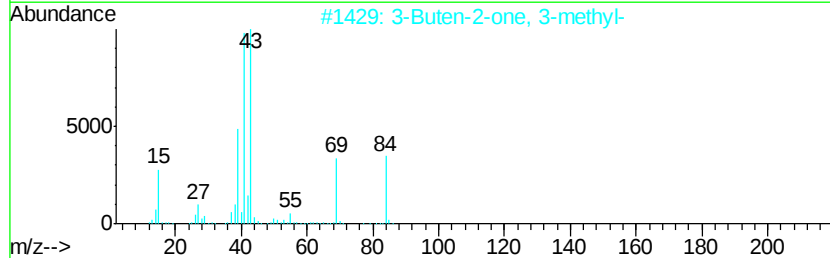
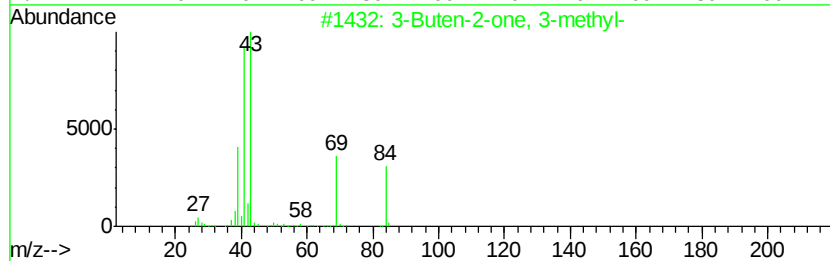
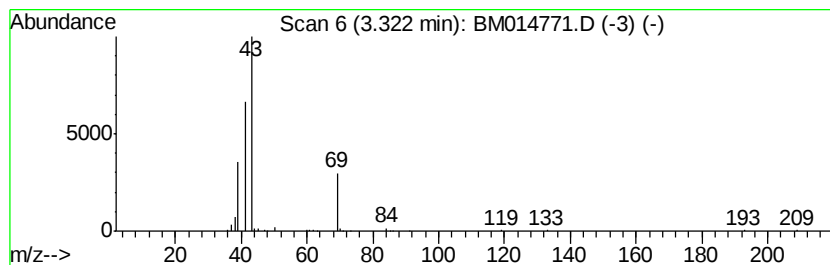
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Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM041918.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 3-Buten-2-one, 3-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
3.32	4.64 ng/ul	427170	1,4-Dichlorobenzene-d4	8.55	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Buten-2-one, 3-methyl-	84	C5H8O	000814-78-8	80
2	3-Buten-2-one, 3-methyl-	84	C5H8O	000814-78-8	38
3	Ethanone, 1-cvclopropyl-	84	C5H8O	000765-43-5	36
4	trans-Crotonamide	85	C4H7NO	000625-37-6	25
5	Furan, 2,3-dihydro-	70	C4H6O	001191-99-7	17



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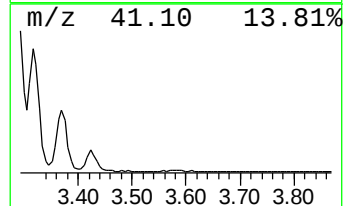
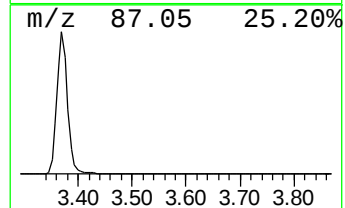
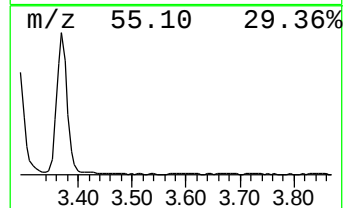
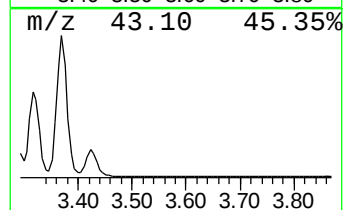
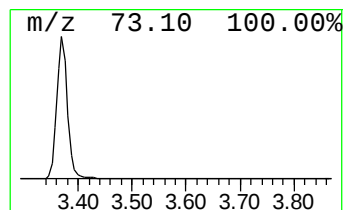
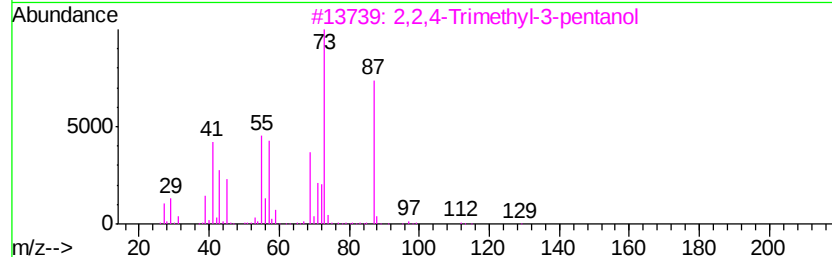
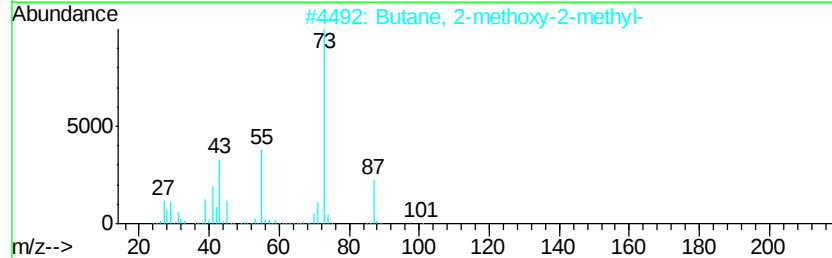
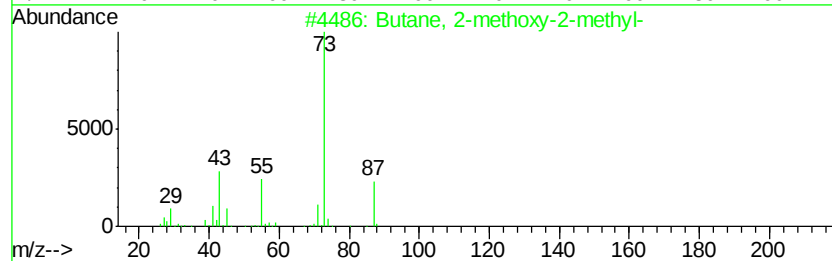
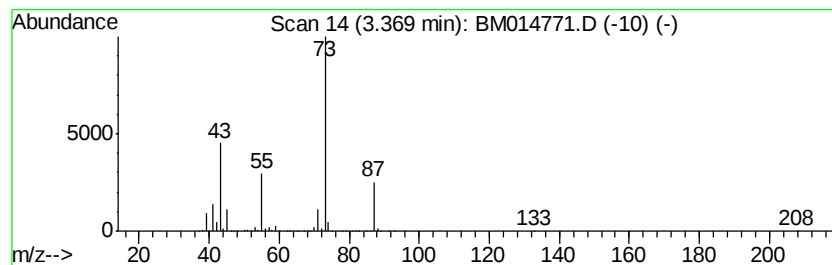
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM041918.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.37	12.75 ng/ul	1172920	1,4-Dichlorobenzene-d4	8.55

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		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
4		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
5		3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	37



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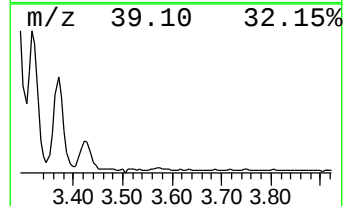
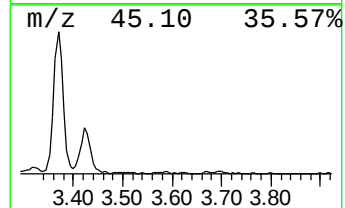
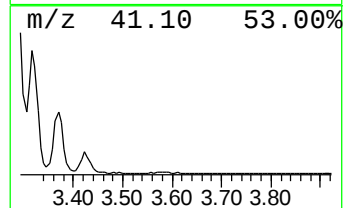
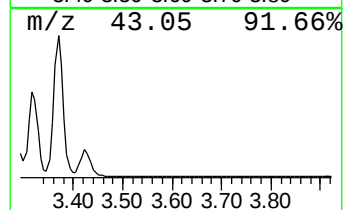
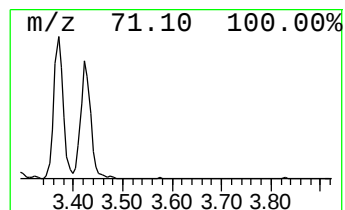
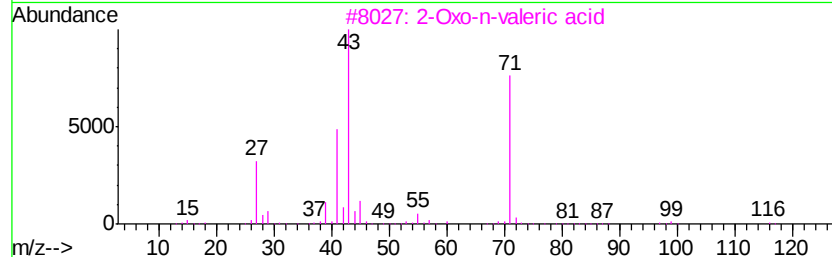
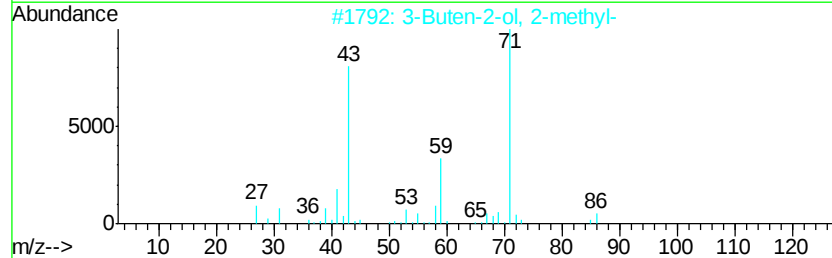
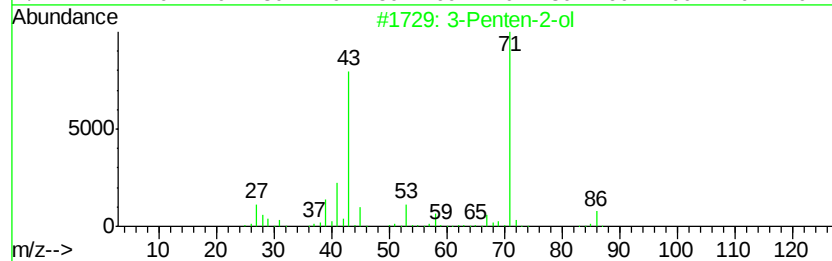
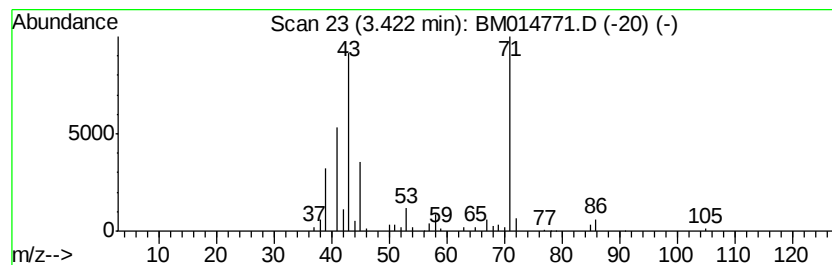
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 3 3-Penten-2-ol Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.42	2.17 ng/ul	199679	1,4-Dichlorobenzene-d4	8.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Penten-2-ol	86	C5H10O	001569-50-2	72
2		3-Buten-2-ol, 2-methyl-	86	C5H10O	000115-18-4	64
3		2-Oxo-n-valeric acid	116	C5H8O3	001821-02-9	64
4		3-Buten-2-ol, 2-methyl-	86	C5H10O	000115-18-4	64
5		Butane, 2-bromo-2-methyl-	150	C5H11Br	000507-36-8	59



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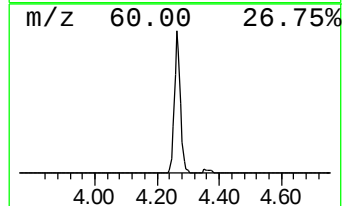
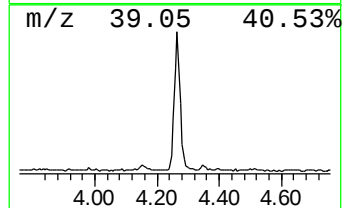
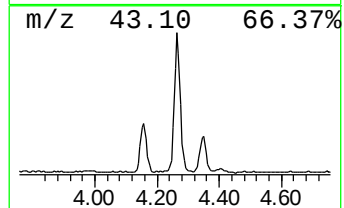
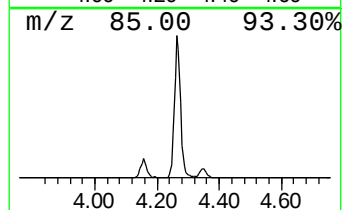
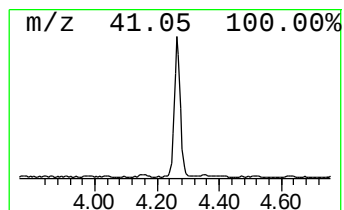
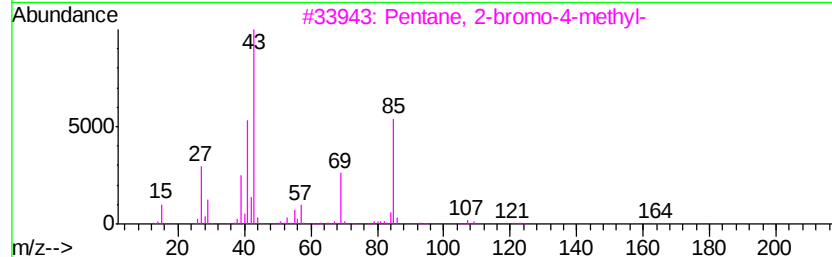
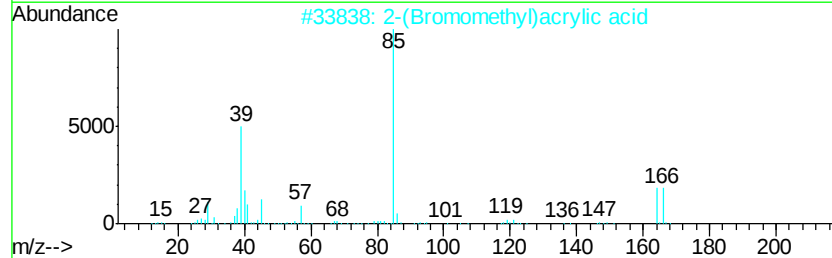
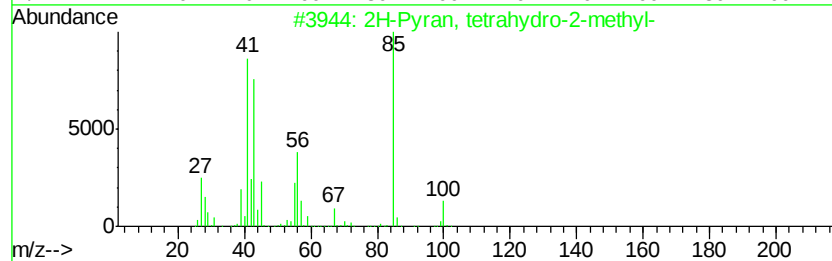
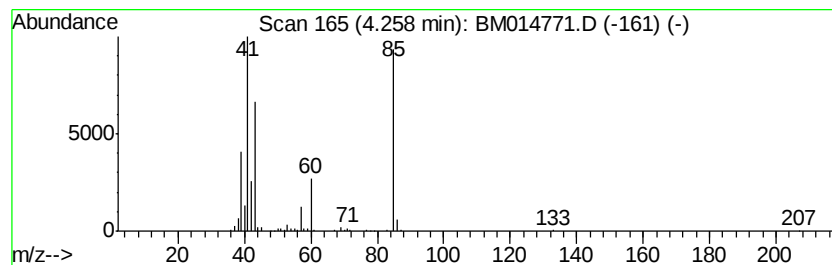
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 unknown-01 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
4.26	4.92 ng/ul	452733	1,4-Dichlorobenzene-d4	8.55	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2H-Pyran, tetrahydro-2-methyl-	100	C6H12O	010141-72-7	42
2	2-(Bromomethyl)acrylic acid	164	C4H5BrO2	072707-66-5	33
3	Pentane, 2-bromo-4-methyl-	164	C6H13Br	030310-22-6	33
4	2H-Pyran, tetrahydro-2-methyl-	100	C6H12O	010141-72-7	33
5	Methacrylamide	85	C4H7NO	000079-39-0	28



Data Path : Z:\HPCHEM1\BNA M\DATA\BM042018\
Data File : BM014771.D
Acq On : 20 Apr 2018 19:54
Operator : SJ/JU
Sample : J2434-12
Misc :
ALS Vial : 18 Sample Multiplier: 1

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ClientSampleId :
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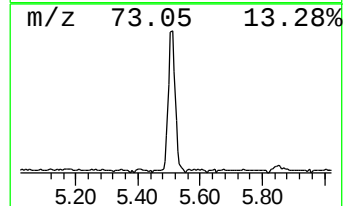
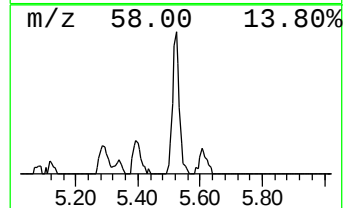
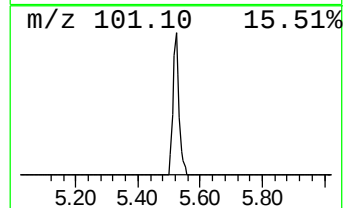
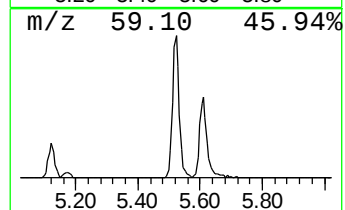
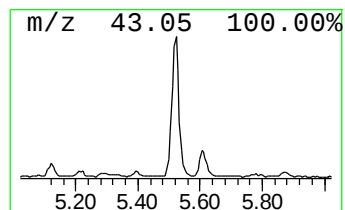
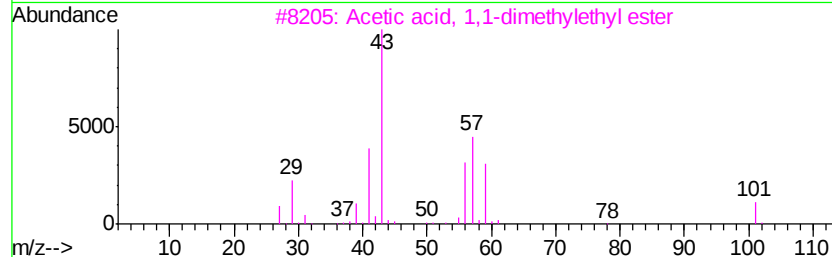
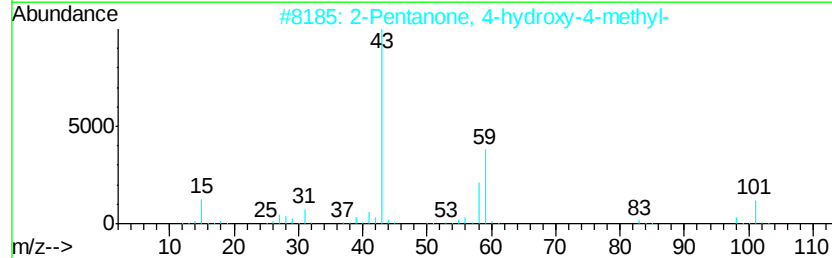
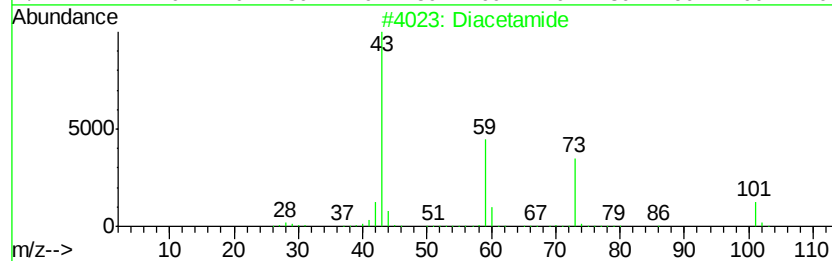
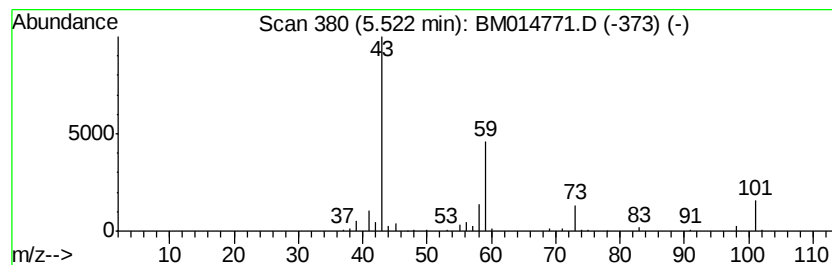
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Diacetamide Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.52	2.74 ng/ul	252487	1,4-Dichlorobenzene-d4	8.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Diacetamide	101	C4H7NO2	000625-77-4	59
2		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	42
3		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	40
4		2-Butanol, 2-methyl-, acetate	130	C7H14O2	000625-16-1	38
5		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	36



Data Path : Z:\HPCHEM1\BNA M\DATA\BM042018\
Data File : BM014771.D
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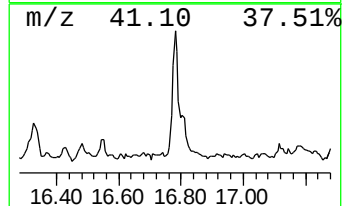
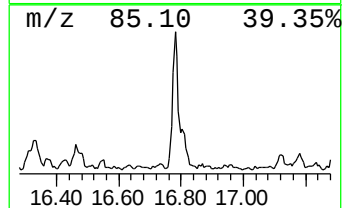
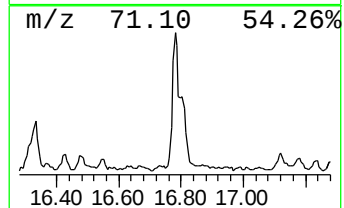
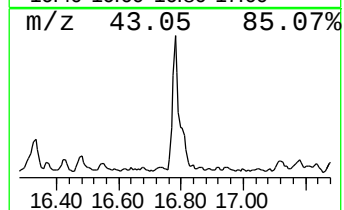
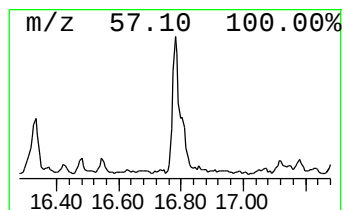
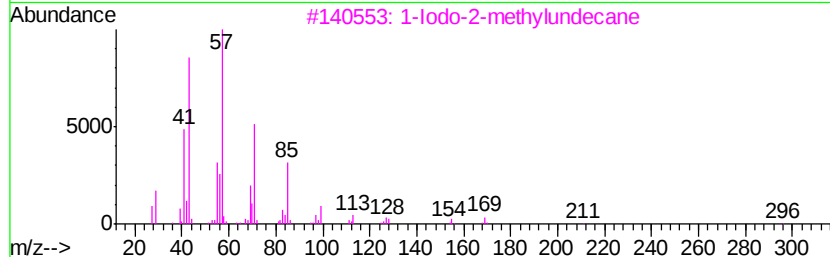
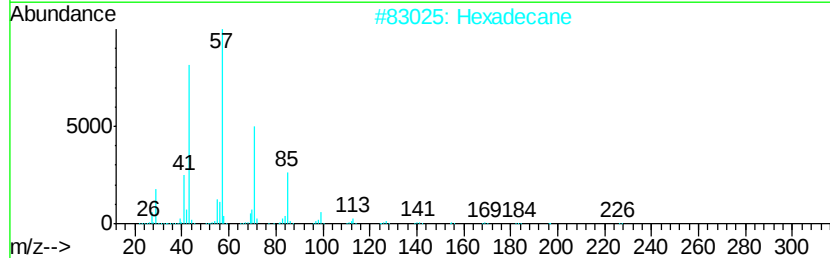
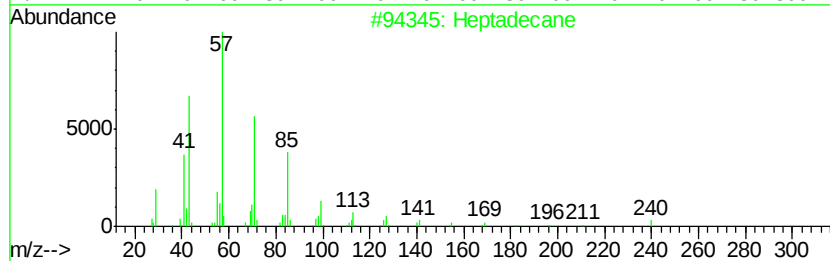
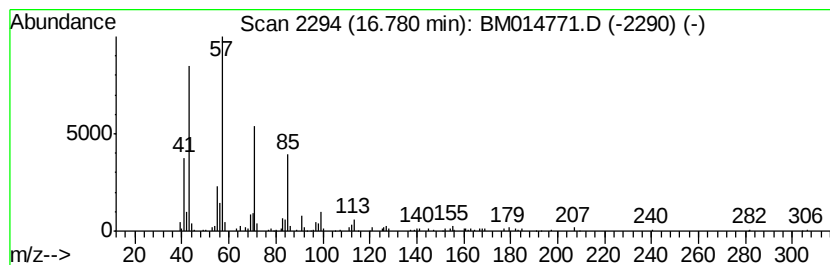
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.78	2.48 ng/ul	436808	Phenanthrene-d10	17.86

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecane	240	C17H36	000629-78-7	94
2			Hexadecane	226	C16H34	000544-76-3	91
3			1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	90
4			Pentadecane	212	C15H32	000629-62-9	90
5			Eicosane, 10-methyl-	296	C21H44	054833-23-7	90



Data Path : Z:\HPCHEM1\BNA M\DATA\BM042018\
Data File : BM014771.D
Acq On : 20 Apr 2018 19:54
Operator : SJ/JU
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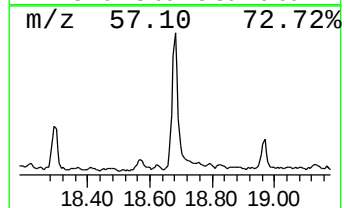
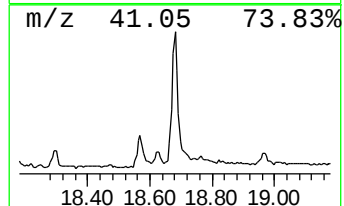
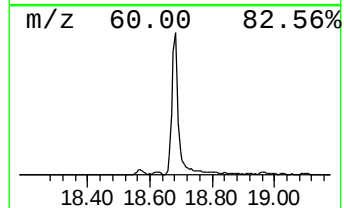
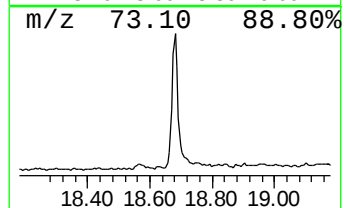
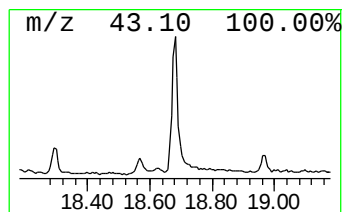
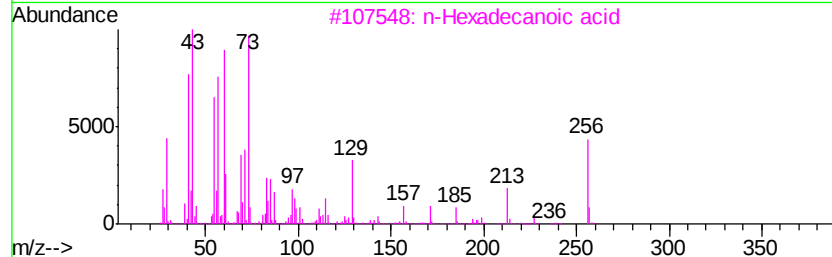
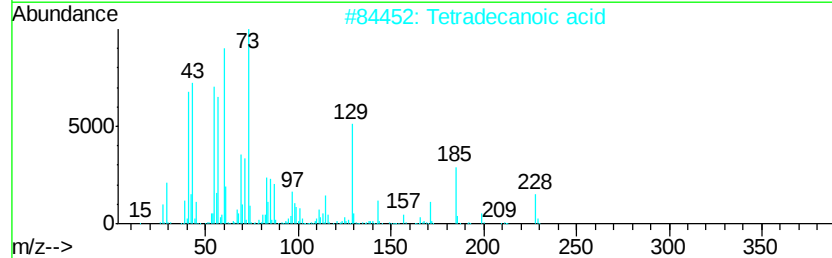
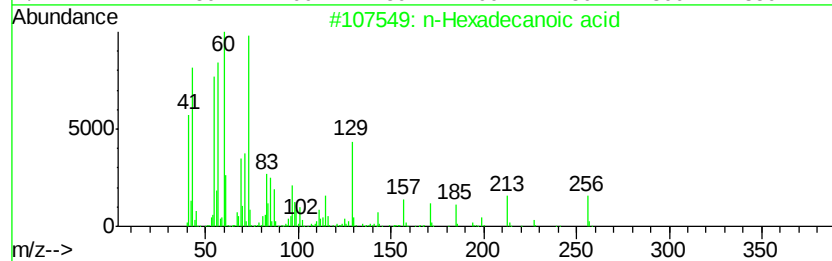
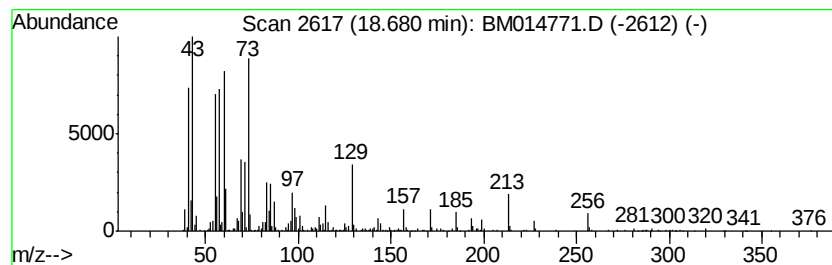
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 n-Hexadecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.68	4.05 ng/ul	713973	Phenanthrene-d10	17.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		Tetradecanoic acid	228	C14H28O2	000544-63-8	97
3		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
4		Tridecanoic acid	214	C13H26O2	000638-53-9	95
5		Octadecanoic acid	284	C18H36O2	000057-11-4	87



Data Path : Z:\HPCHEM1\BNA M\DATA\BM042018\
Data File : BM014771.D
Acq On : 20 Apr 2018 19:54
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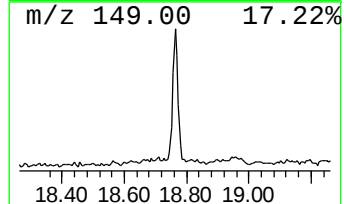
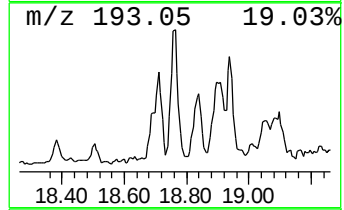
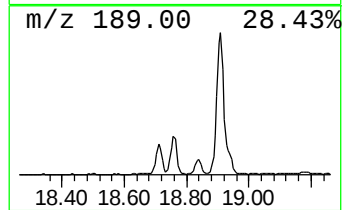
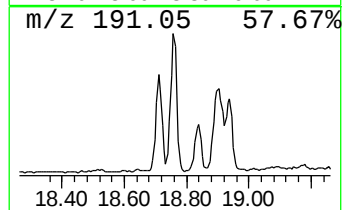
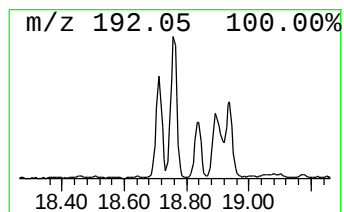
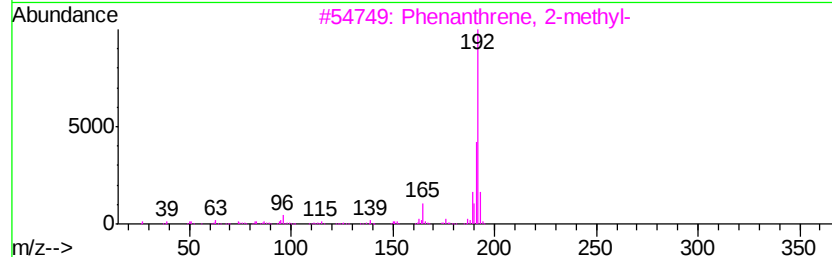
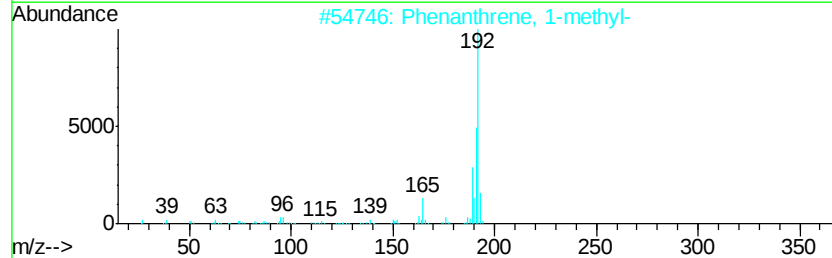
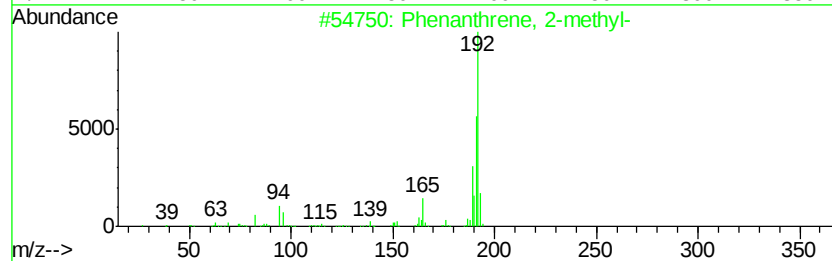
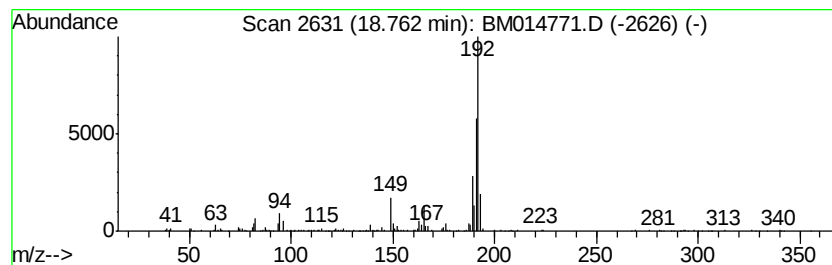
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Phenanthrene, 2-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.76	2.86 ng/ul	504821	Phenanthrene-d10	17.86

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	98
2		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	97
3		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
4		Anthracene, 2-methyl-	192	C15H12	000613-12-7	95
5		Anthracene, 2-methyl-	192	C15H12	000613-12-7	94



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Acq On : 20 Apr 2018 19:54
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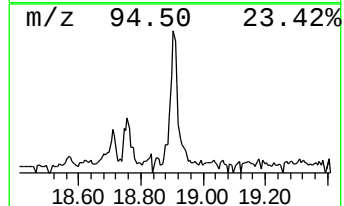
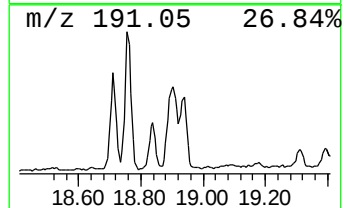
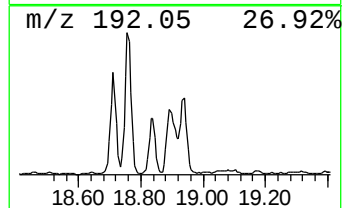
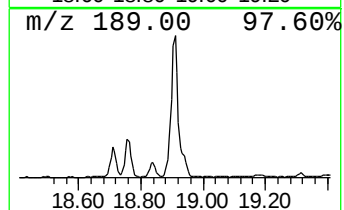
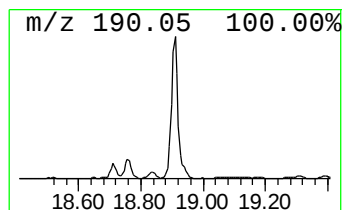
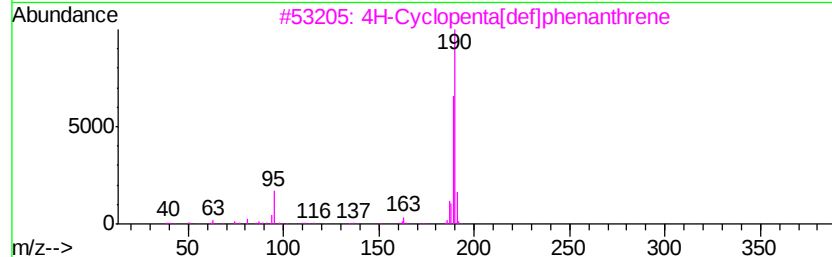
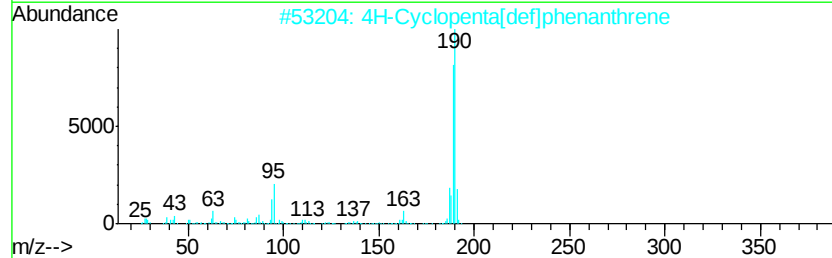
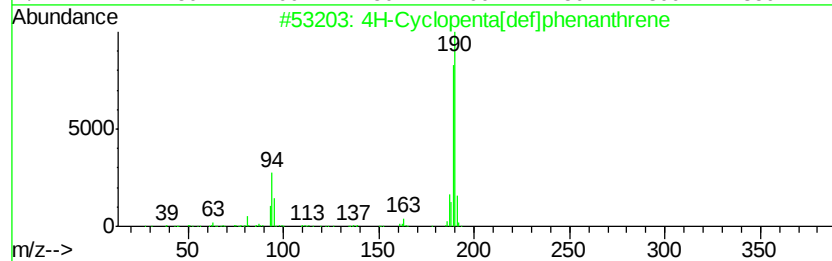
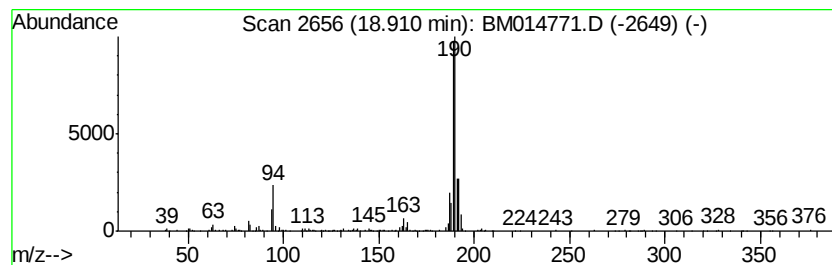
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 4H-Cyclopenta[def]phenanthrene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.91	4.15 ng/ul	731965	Phenanthrene-d10	17.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	87
2		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	76
3		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	64
4		2,3,5,6-Tetramethylterephthalald...	190	C12H14O2	007072-01-7	58
5		6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	45



Data Path : Z:\HPCHEM1\BNA M\DATA\BM042018\
Data File : BM014771.D
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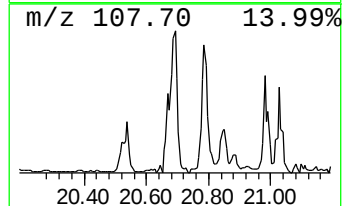
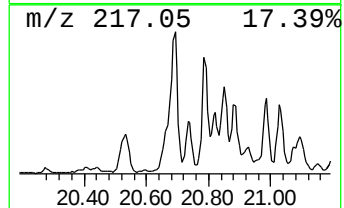
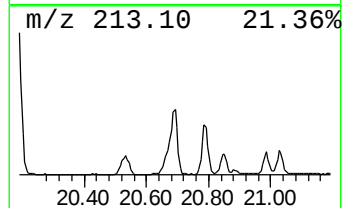
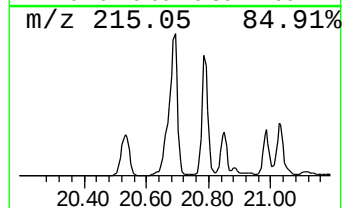
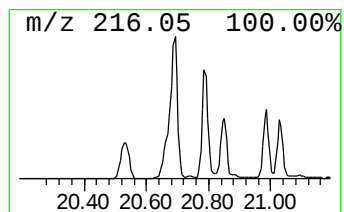
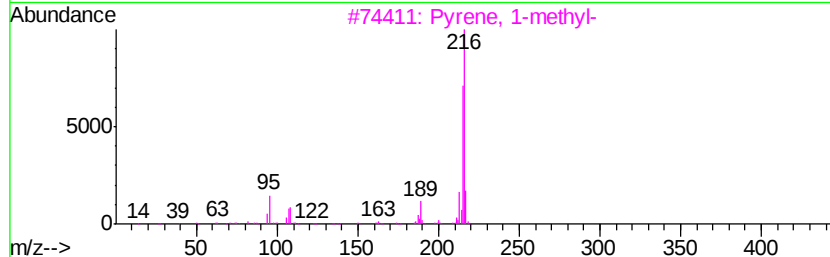
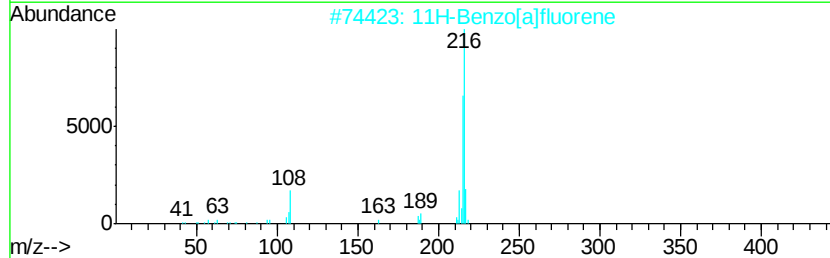
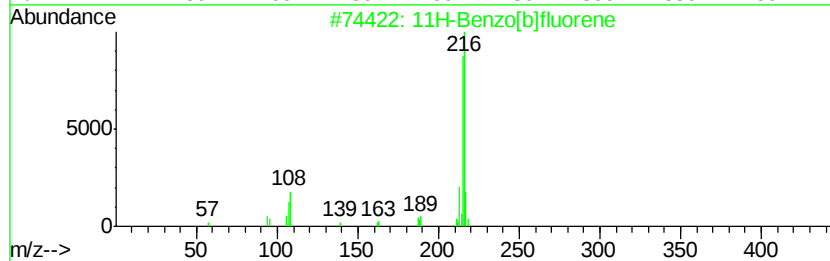
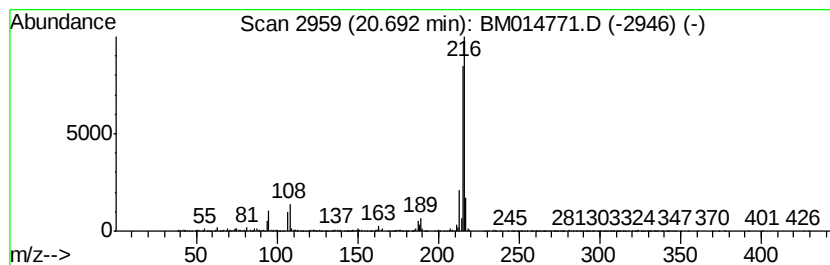
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 11H-Benzo[b]fluorene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.69	5.01 ng/ul	3641810	Chrysene-d12	21.94

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	94
2			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	93
3			Pvrene, 1-methyl-	216	C17H12	002381-21-7	93
4			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	93
5			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	91



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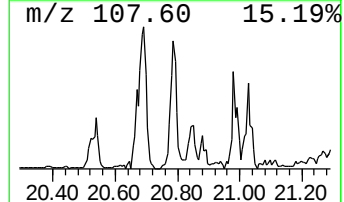
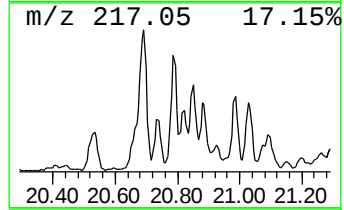
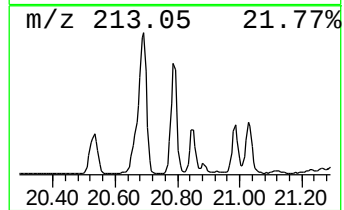
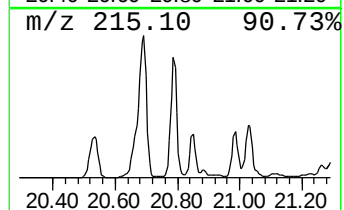
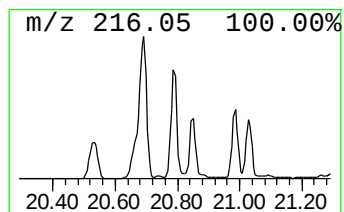
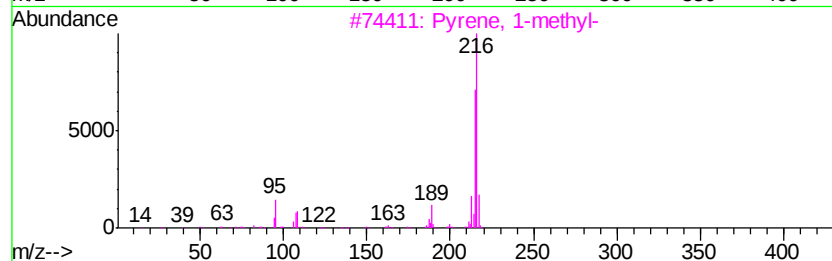
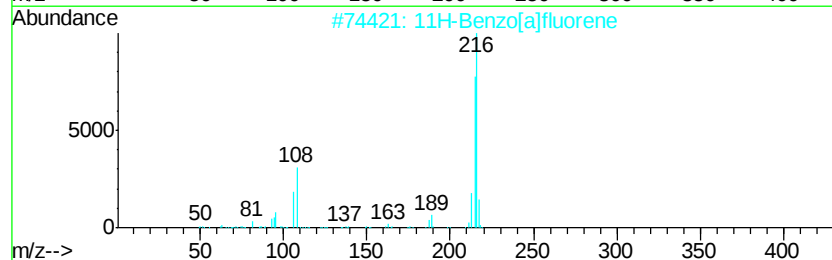
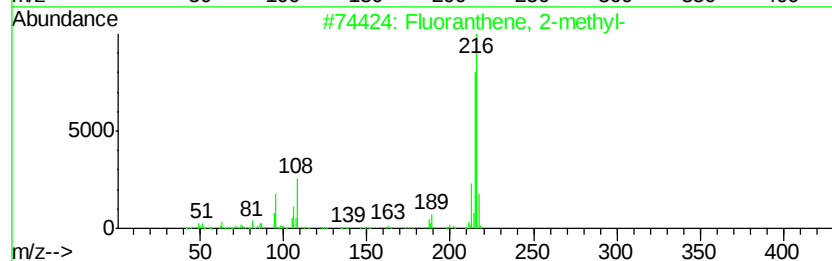
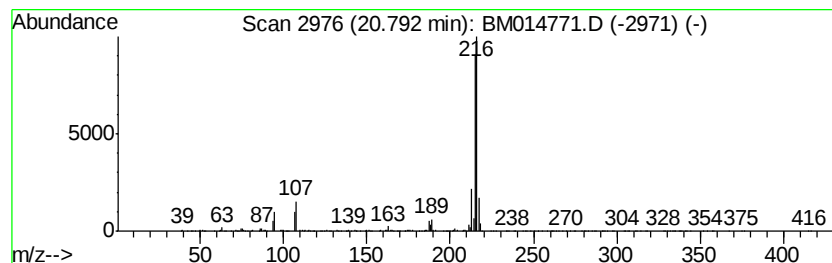
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 Fluoranthene, 2-methyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.79	2.10 ng/ul	1523630	Chrysene-d12	21.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	95
2		11H-Benzo[a]fluorene	216	C17H12	000238-84-6	91
3		Pyrene, 1-methyl-	216	C17H12	002381-21-7	90
4		11H-Benzo[b]fluorene	216	C17H12	000243-17-4	90
5		11H-Benzo[b]fluorene	216	C17H12	000243-17-4	87



Data Path : Z:\HPCHEM1\BNA M\DATA\BM042018\
Data File : BM014771.D
Acq On : 20 Apr 2018 19:54
Operator : SJ/JU
Sample : J2434-12
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0AC7

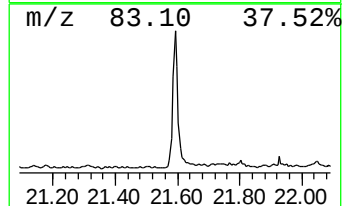
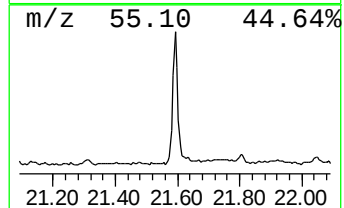
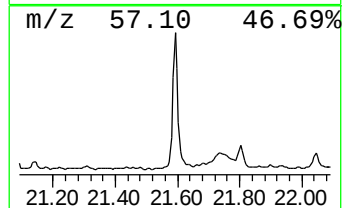
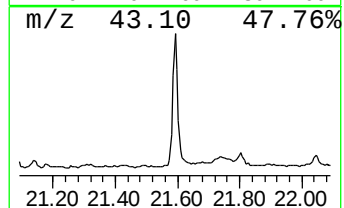
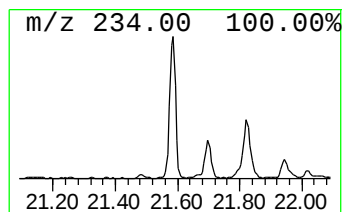
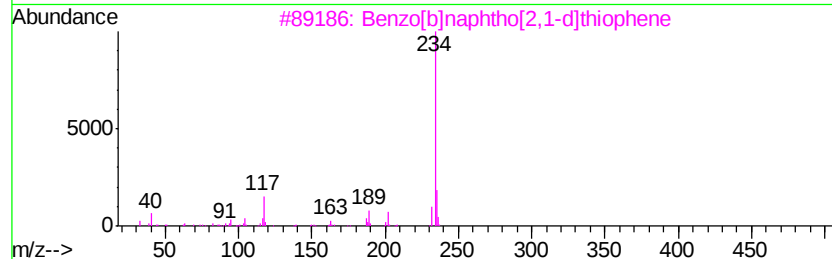
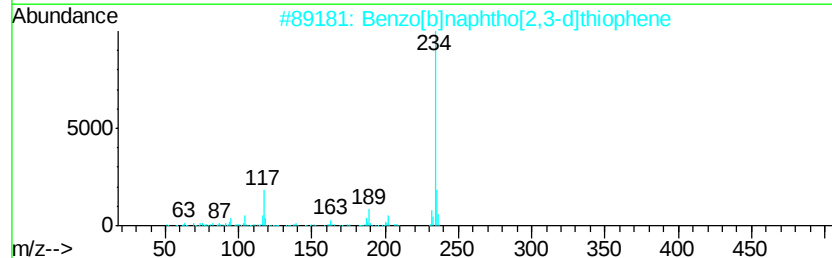
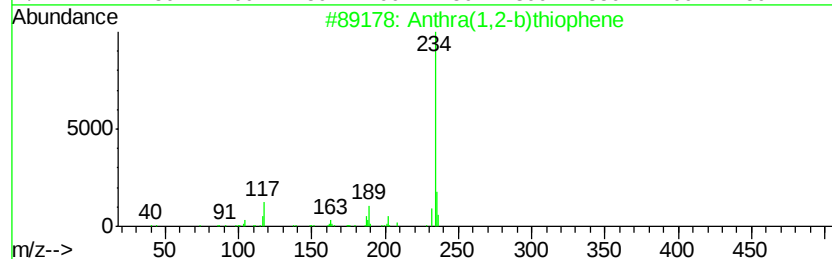
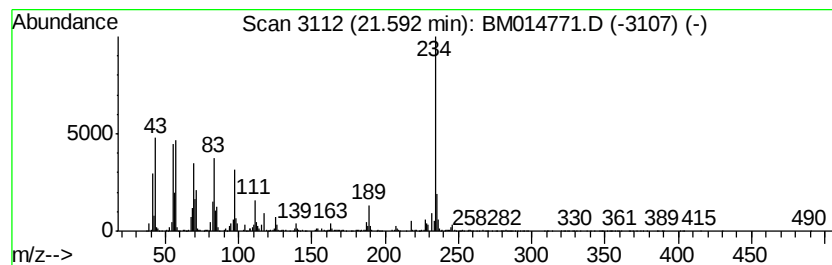
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 Anthra(1,2-b)thiophene Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.59	2.30 ng/ul	1669190	Chrysene-d12	21.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Anthra(1,2-b)thiophene	234	C16H10S	000227-86-1	86
2		Benzo[b]naphtho[2,3-d]thiophene	234	C16H10S	000243-46-9	70
3		Benzo[b]naphtho[2,1-d]thiophene	234	C16H10S	000239-35-0	70
4		Benzo[b]naphtho[2,1-d]thiophene	234	C16H10S	000239-35-0	70
5		Benzo[b]naphtho[2,1-d]thiophene	234	C16H10S	000239-35-0	70



Data Path : Z:\HPCHEM1\BNA M\DATA\BM042018\
Data File : BM014771.D
Acq On : 20 Apr 2018 19:54
Operator : SJ/JU
Sample : J2434-12
Misc :
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ClientSampleId :
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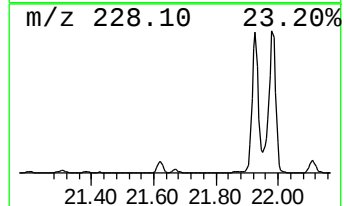
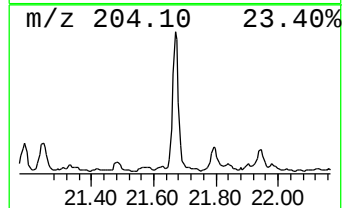
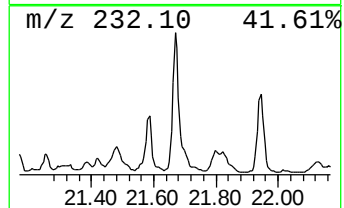
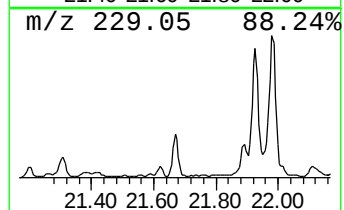
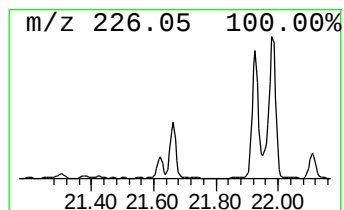
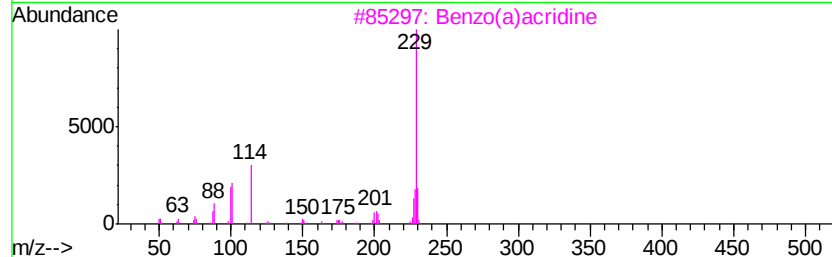
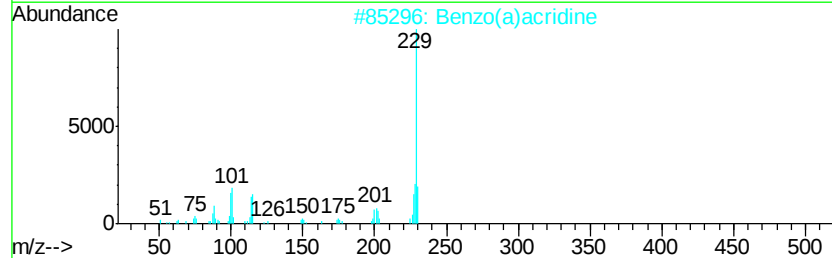
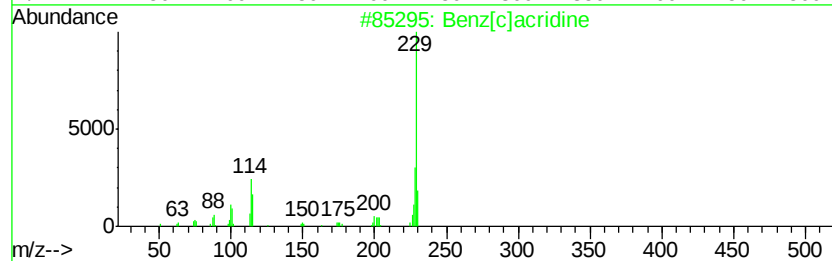
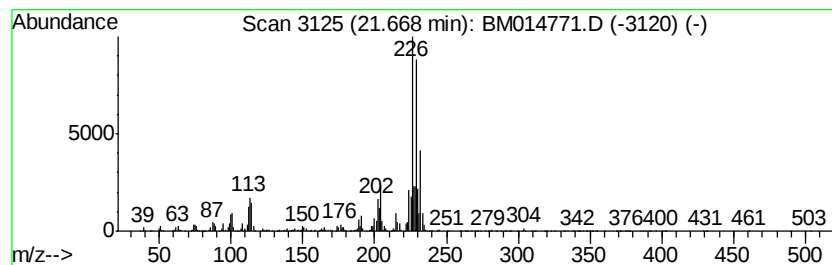
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 unknown-02 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.67	2.38 ng/ul	1730390	Chrysene-d12	21.94

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benz[c]acridine	229	C17H11N	000225-51-4	42
2			Benzo(a)acridine	229	C17H11N	000225-11-6	27
3			Benzo(a)acridine	229	C17H11N	000225-11-6	22
4			Ferrocene, [(hydroxvimino)methyl...	229	C11H11FeNO	032679-07-5	22
5			Naphtho(2,3-h)quinoline	229	C17H11N	000084-56-0	22



Data Path : Z:\HPCHEM1\BNA M\DATA\BM042018\
Data File : BM014771.D
Acq On : 20 Apr 2018 19:54
Operator : SJ/JU
Sample : J2434-12
Misc :
ALS Vial : 18 Sample Multiplier: 1

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ClientSampleId :
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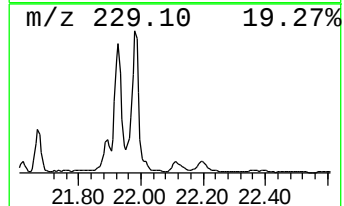
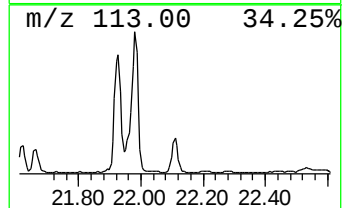
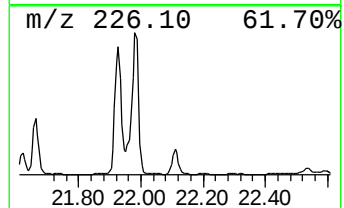
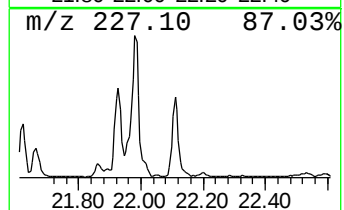
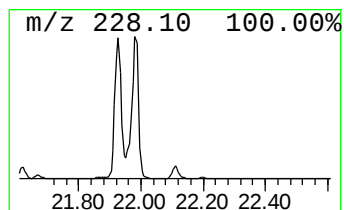
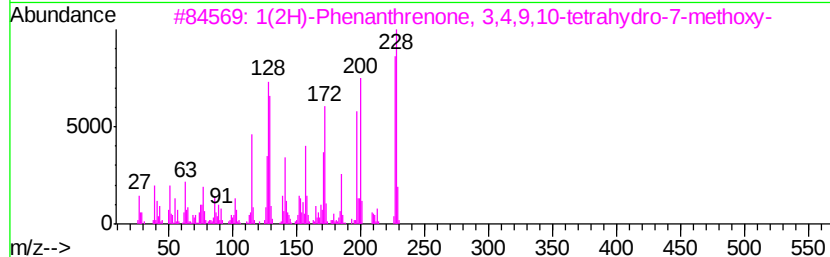
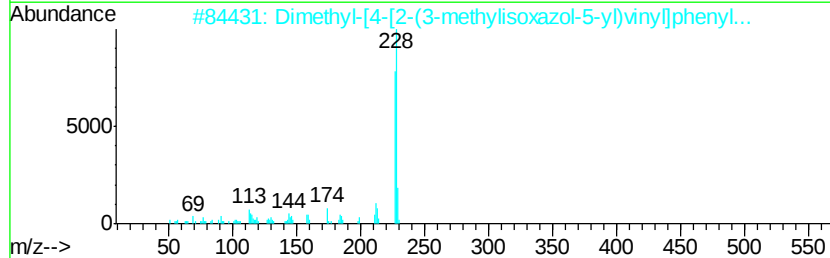
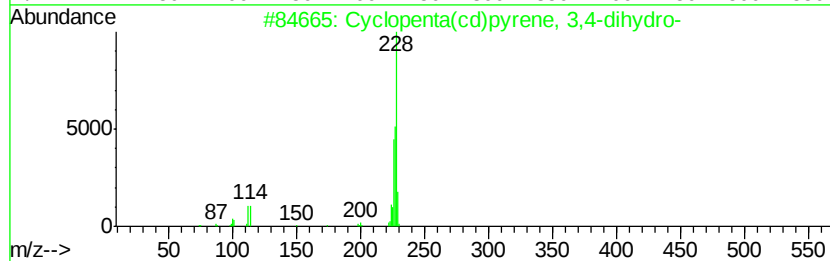
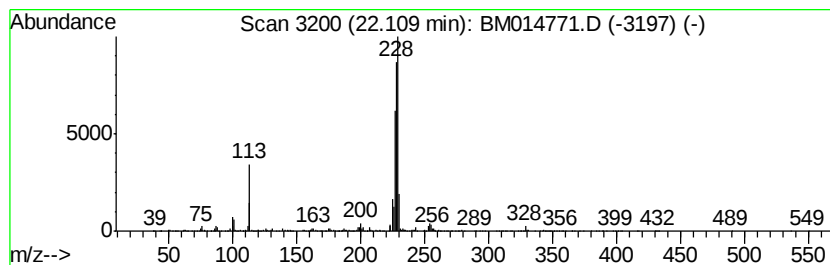
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 Cyclopenta(cd)pyrene, 3,4-d... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.11	2.14 ng/ul	1555900	Chrysene-d12	21.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopenta(cd)pyrene, 3,4-dihydro-	228	C18H12	025732-74-5	55
2		Dimethyl-[4-[2-(3-methylisoxazol...	228	C14H16N2O	1000306-39-6	50
3		1(2H)-Phenanthrenone, 3,4,9,10-t...	228	C15H16O2	014427-61-3	50
4		Benzo[c]phenanthrene	228	C18H12	000195-19-7	49
5		Benzo[c]phenanthrene	228	C18H12	000195-19-7	47



Data Path : Z:\HPCHEM1\BNA M\DATA\BM042018\
Data File : BM014771.D
Acq On : 20 Apr 2018 19:54
Operator : SJ/JU
Sample : J2434-12
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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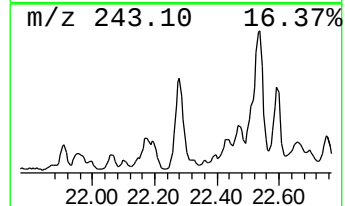
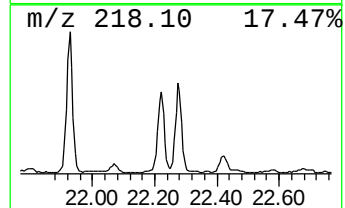
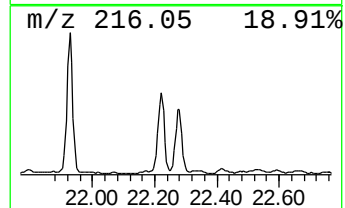
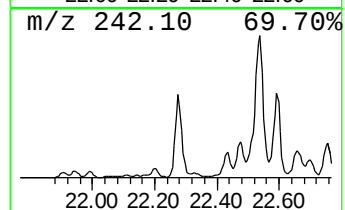
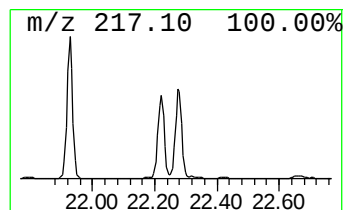
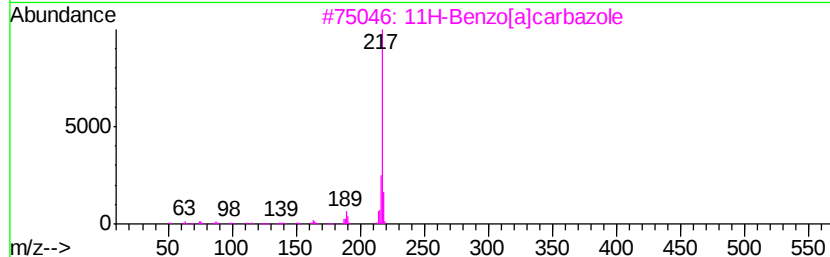
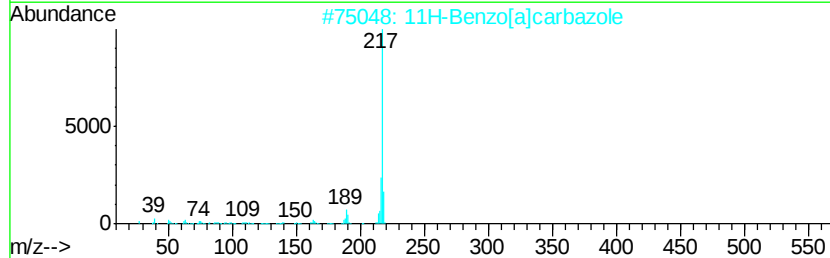
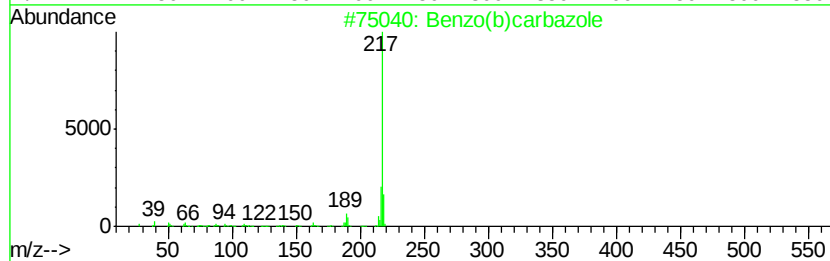
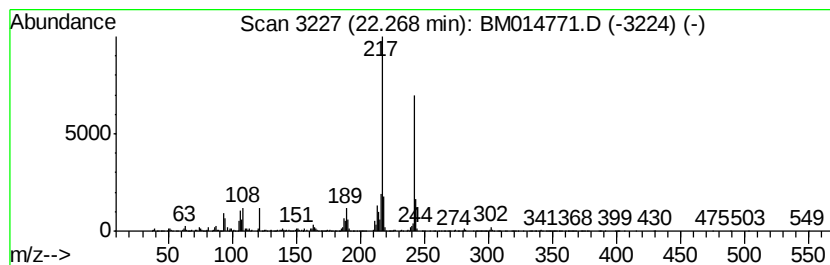
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 Benzo(b)carbazole Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.27	3.15 ng/ul	2288520	Chrysene-d12	21.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo(b)carbazole	217	C16H11N	000243-28-7	90
2		11H-Benzo[a]carbazole	217	C16H11N	000239-01-0	78
3		11H-Benzo[a]carbazole	217	C16H11N	000239-01-0	64
4		Benzo(c)carbazole	217	C16H11N	034777-33-8	60
5		7H-Benzo[c]carbazole	217	C16H11N	000205-25-4	60



Data Path : Z:\HPCHEM1\BNA M\DATA\BM042018\
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Sample : J2434-12
Misc :
ALS Vial : 18 Sample Multiplier: 1

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ClientSampleId :
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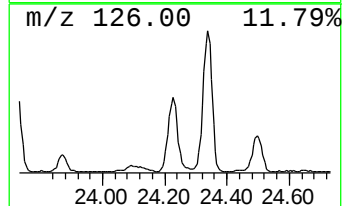
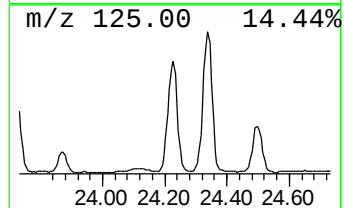
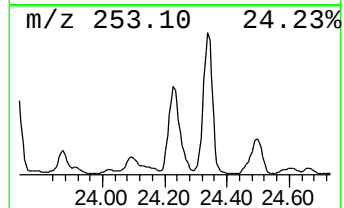
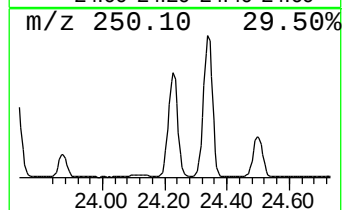
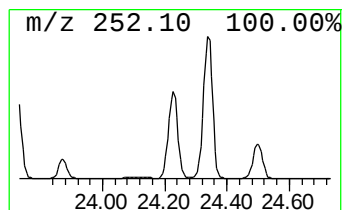
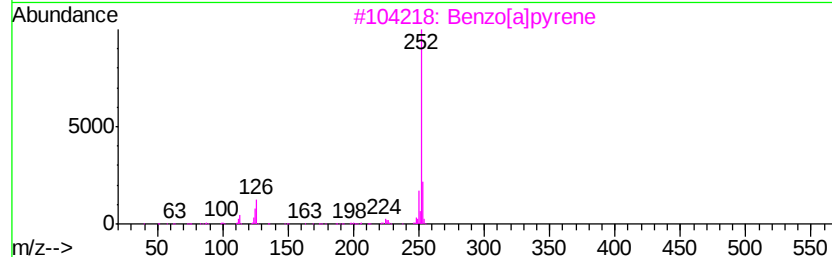
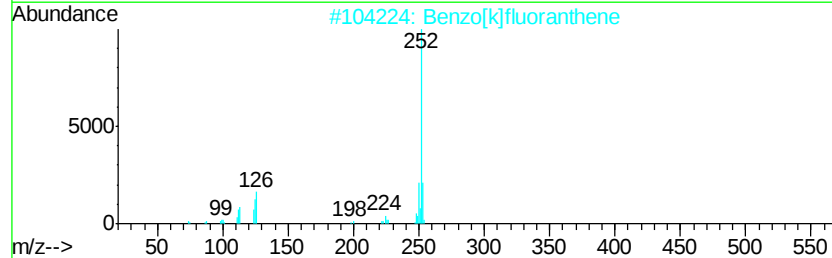
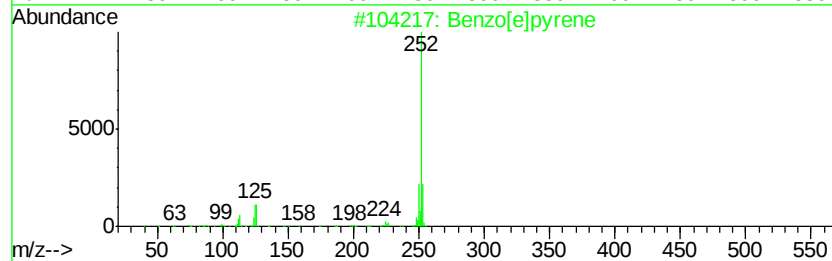
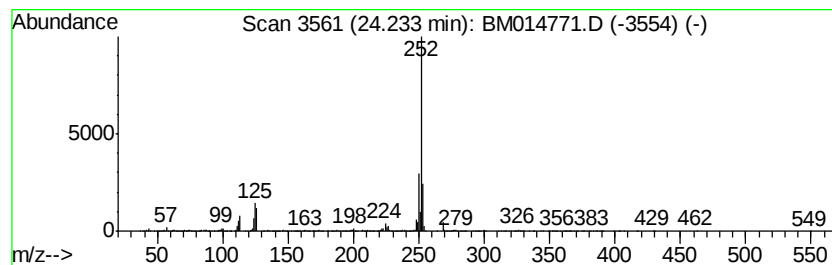
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Quant Title : SVOA CALIBRATION

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

Peak Number 17 Benzo[e]pyrene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.23	48.74 ng/ul	9314100	Perylene-d12	24.44

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[e]pyrene	252	C20H12	000192-97-2	99
2			Benzo[k]fluoranthene	252	C20H12	000207-08-9	98
3			Benzo[a]pyrene	252	C20H12	000050-32-8	98
4			Benzo[e]acephenanthrylene	252	C20H12	000205-99-2	97
5			Perylene	252	C20H12	000198-55-0	96



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM042018\
 Data File : BM014771.D
 Acq On : 20 Apr 2018 19:54
 Operator : SJ/JU
 Sample : J2434-12
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM041918.M
 Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
3-Buten-2-one, 3-...	3.32	4.6	ng/ul	427170	1	8.55	1840360	20.0
Butane, 2-methoxy...	3.37	12.8	ng/ul	1172920	1	8.55	1840360	20.0
3-Penten-2-ol	3.42	2.2	ng/ul	199679	1	8.55	1840360	20.0
unknown-01	4.26	4.9	ng/ul	452733	1	8.55	1840360	20.0
Diacetamide	5.52	2.7	ng/ul	252487	1	8.55	1840360	20.0
(DEL) Alkane: Str...	16.78	2.5	ng/ul	436808	4	17.86	3525850	20.0
n-Hexadecanoic acid	18.68	4.0	ng/ul	713973	4	17.86	3525850	20.0
Phenanthrene, 2-m...	18.76	2.9	ng/ul	504821	4	17.86	3525850	20.0
4H-Cyclopenta[def...	18.91	4.2	ng/ul	731965	4	17.86	3525850	20.0
11H-Benzo[b]fluorene	20.69	5.0	ng/ul	3641810	5	21.94	14539900	20.0
Fluoranthene, 2-m...	20.79	2.1	ng/ul	1523630	5	21.94	14539900	20.0
Anthra(1,2-b)thio...	21.59	2.3	ng/ul	1669190	5	21.94	14539900	20.0
unknown-02	21.67	2.4	ng/ul	1730390	5	21.94	14539900	20.0
Cyclopenta(cd)pyr...	22.11	2.1	ng/ul	1555900	5	21.94	14539900	20.0
Benzo(b)carbazole	22.27	3.1	ng/ul	2288520	5	21.94	14539900	20.0
Benzo[e]pyrene	24.23	48.7	ng/ul	9314100	6	24.44	3822070	20.0