

Data Path : Z:\HPCHEM1\BNA M\DATA\BM031218\
 Data File : BM014525.D
 Acq On : 12 Mar 2018 19:46
 Operator : SJ/JU
 Sample : J1813-02
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 F6F17

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-BM021418.M
 Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	6.687	640	646	662	rBV	172007	441942	18.97%	1.660%
2	6.834	664	671	685	rVB	148229	279176	11.98%	1.048%
3	7.022	697	703	715	rBV	233991	444363	19.08%	1.669%
4	7.475	772	780	795	rBV	208857	407310	17.48%	1.530%
5	8.210	899	905	925	rBV2	247683	540829	23.22%	2.031%
6	8.645	972	979	993	rBV	247385	480236	20.61%	1.803%
7	9.357	1094	1100	1119	rBV2	223952	456997	19.62%	1.716%
8	9.904	1185	1193	1211	rBV2	239700	585047	25.11%	2.197%
9	10.245	1243	1251	1259	rBV	312915	548238	23.53%	2.059%
10	10.422	1272	1281	1301	rBV2	192333	542248	23.28%	2.036%
11	13.010	1714	1721	1725	rBV5	18764	29100	1.25%	0.109%
12	13.116	1732	1739	1746	rVB4	33199	49844	2.14%	0.187%
13	13.257	1758	1763	1767	rBV4	44104	66651	2.86%	0.250%
14	13.374	1777	1783	1788	rBV2	298195	442425	18.99%	1.661%
15	13.427	1788	1792	1800	rVB2	121171	201123	8.63%	0.755%
16	13.569	1805	1816	1821	rVV	703274	1132535	48.62%	4.253%
17	13.616	1821	1824	1836	rVB	108135	195485	8.39%	0.734%
18	13.833	1853	1861	1870	rBV	769066	1174231	50.41%	4.409%
19	14.045	1888	1897	1900	rBV4	33713	68176	2.93%	0.256%
20	14.145	1907	1914	1922	rBV2	493812	779102	33.44%	2.926%
21	14.445	1960	1965	1988	rBV3	152379	556635	23.89%	2.090%
22	14.669	1998	2003	2008	rVB8	12760	28365	1.22%	0.107%
23	15.004	2056	2060	2066	rVV2	38989	52693	2.26%	0.198%
24	15.145	2079	2084	2093	rBV2	871821	1422057	61.04%	5.340%
25	15.327	2108	2115	2124	rBV2	273471	539372	23.15%	2.025%
26	15.427	2124	2132	2141	rVB2	255241	417112	17.91%	1.566%
27	15.627	2160	2166	2175	rVB4	58233	98551	4.23%	0.370%
28	15.868	2202	2207	2214	rBV4	29596	45136	1.94%	0.169%
29	16.904	2378	2383	2393	rVB	729764	1042445	44.75%	3.915%
30	17.010	2394	2401	2416	rBV2	1064976	1700788	73.01%	6.387%
31	17.757	2523	2528	2533	rBV5	19256	25645	1.10%	0.096%
32	17.815	2533	2538	2550	rBV2	225320	353745	15.19%	1.328%
33	18.956	2727	2732	2735	rBV4	53075	81788	3.51%	0.307%
34	18.986	2735	2737	2748	rVB9	36391	73876	3.17%	0.277%

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35	19.109	2753	2758	2769	rBV3	112063	193057	8.29%	0.725%
36	19.251	2778	2782	2788	rBV	364081	412948	17.73%	1.551%
37	19.321	2788	2794	2800	rVV	1475563	2060869	88.47%	7.739%
38	19.368	2800	2802	2806	rVB4	49518	54292	2.33%	0.204%
39	19.562	2832	2835	2840	rBV5	29443	36783	1.58%	0.138%
40	20.027	2909	2914	2922	rBV3	122969	168610	7.24%	0.633%
41	20.198	2936	2943	2947	rBV	1061065	1290811	55.41%	4.847%
42	20.245	2947	2951	2952	rBV2	186785	206893	8.88%	0.777%
43	20.309	2958	2962	2965	rVB	243152	274802	11.80%	1.032%
44	20.368	2969	2972	2976	rBV3	49596	71959	3.09%	0.270%
45	20.403	2976	2978	2983	rVV5	22566	27044	1.16%	0.102%
46	20.833	3047	3051	3060	rBV2	224440	295235	12.67%	1.109%
47	21.127	3096	3101	3114	rBV2	1028979	1481173	63.58%	5.562%
48	21.703	3194	3199	3205	rBV3	308351	485944	20.86%	1.825%
49	23.156	3440	3446	3452	rBV2	1331743	2329554	100.00%	8.748%
50	23.291	3463	3469	3479	rVB2	657086	1314868	56.44%	4.938%
51	23.744	3544	3546	3552	rVB4	90940	145200	6.23%	0.545%
52	25.544	3848	3852	3862	rVB10	210279	476465	20.45%	1.789%

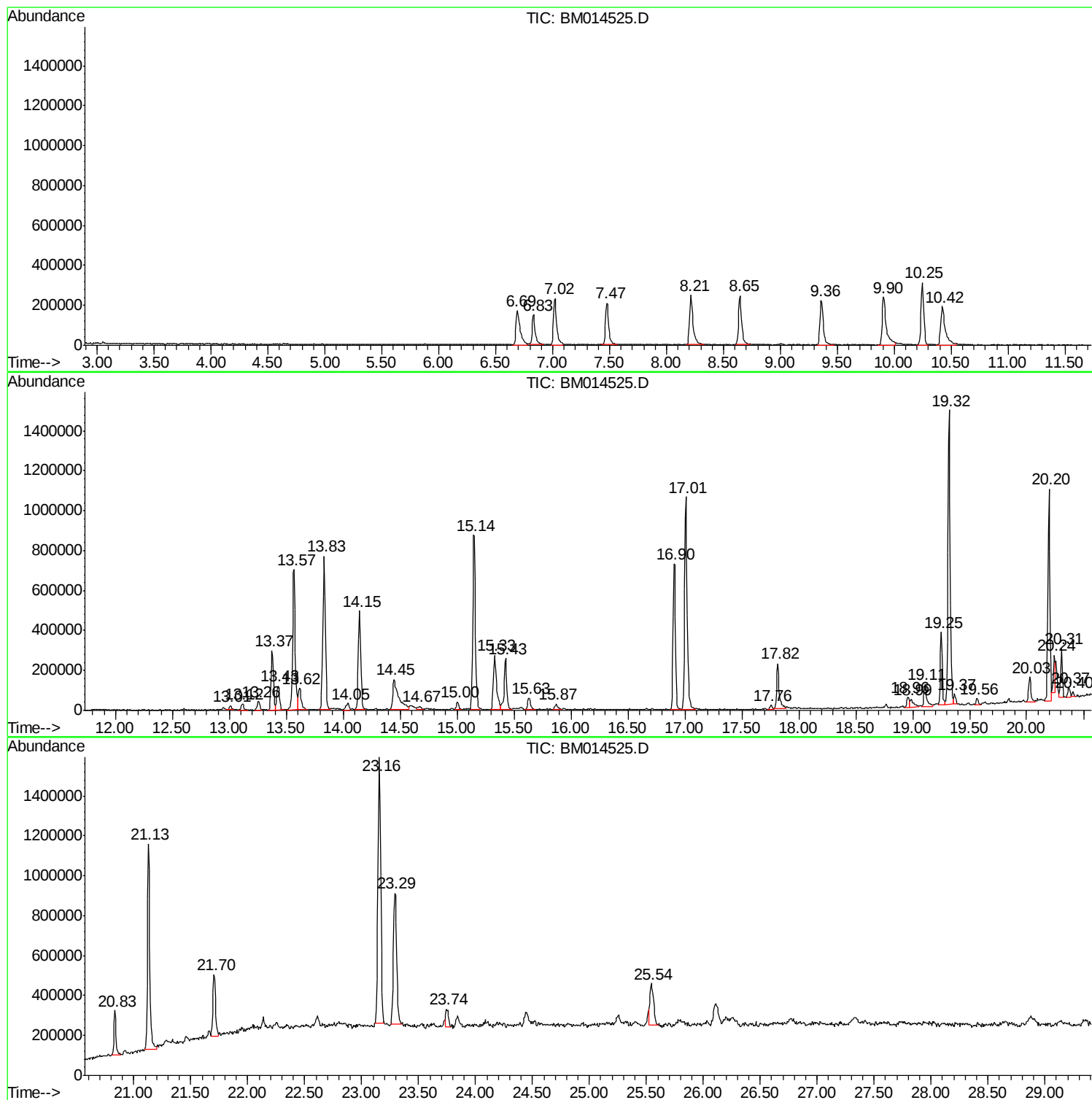
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Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM021418.M
Quant Title : SVOA CALIBRATION

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



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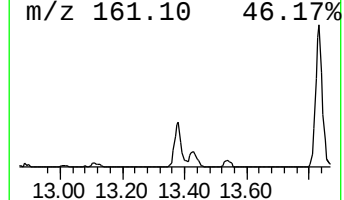
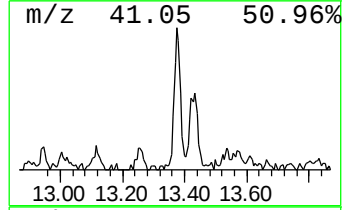
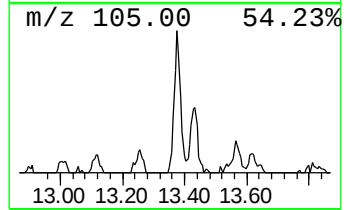
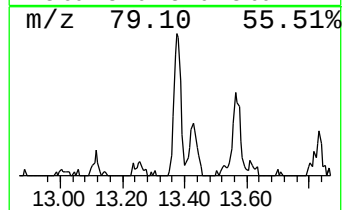
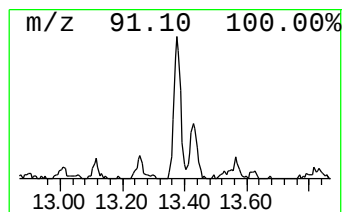
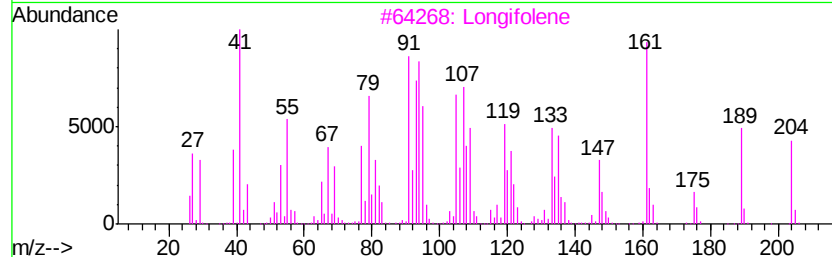
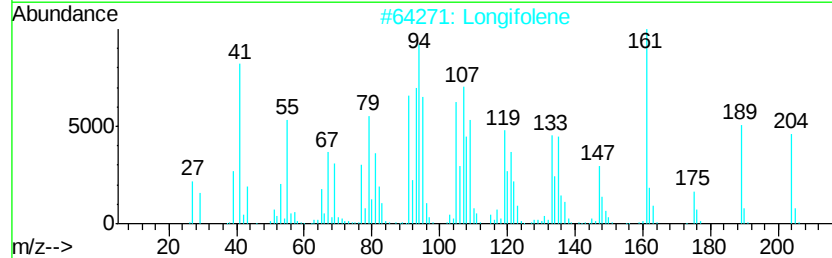
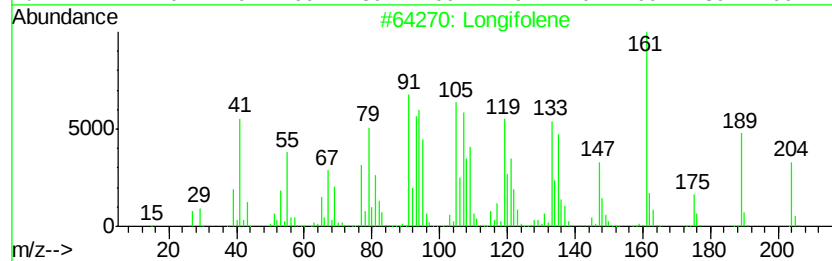
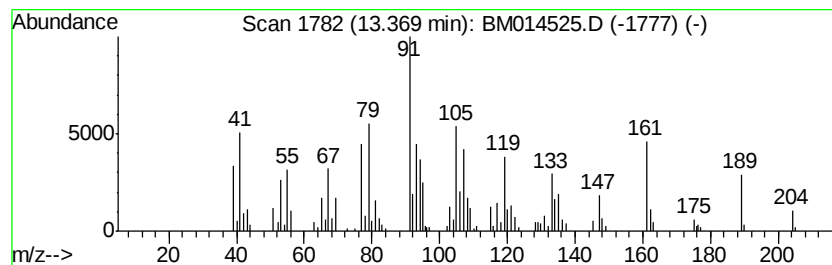
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Quant Title : SVOA CALIBRATION

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

Peak Number 1 Longifolene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.37	11.36 ng/ul	442425	Acenaphthene-d10	14.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Longifolene	204	C15H24	000475-20-7	99
2		Longifolene	204	C15H24	000475-20-7	96
3		Longifolene	204	C15H24	000475-20-7	95
4		Naphthalene, 1,2,3,5,6,7,8,8a-oc...	204	C15H24	004630-07-3	95
5		Naphthalene, 1,2,3,5,6,7,8,8a-oc...	204	C15H24	010219-75-7	95



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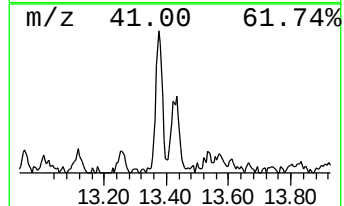
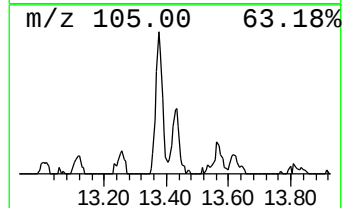
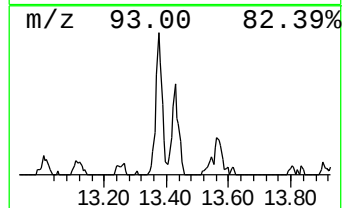
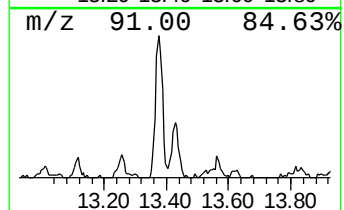
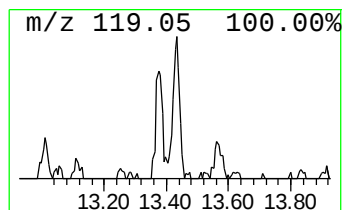
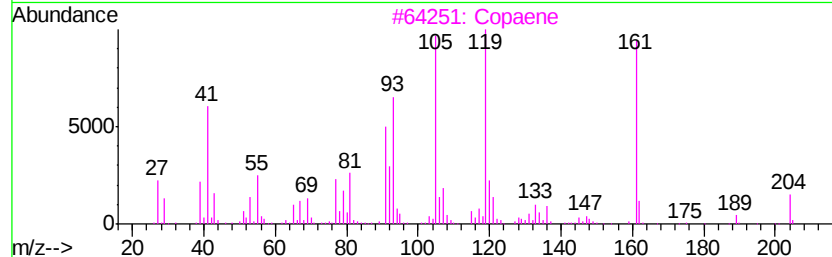
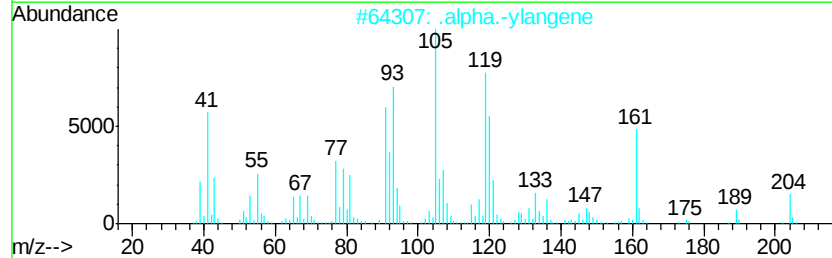
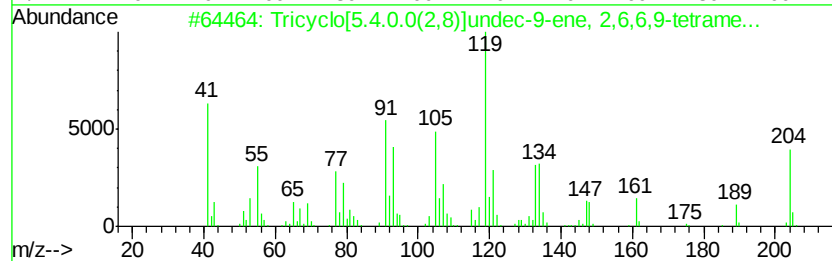
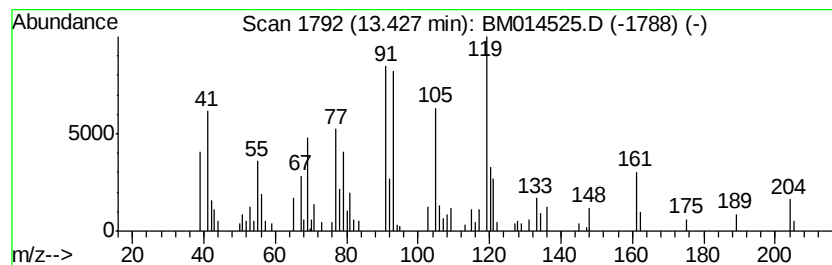
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 2 Tricyclo[5.4.0.0(2,8)]undec... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.43	5.16 ng/ul	201123	Acenaphthene-d10	14.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tricyclo[5.4.0.0(2,8)]undec-9-en...	204	C15H24	005989-08-2	58
2		.alpha.-ylangene	204	C15H24	1000374-19-0	55
3		Copaene	204	C15H24	003856-25-5	49
4		cis-.beta.-Farnesene	204	C15H24	028973-97-9	49
5		Di-epi-.alpha.-cedrene	204	C15H24	050894-66-1	46



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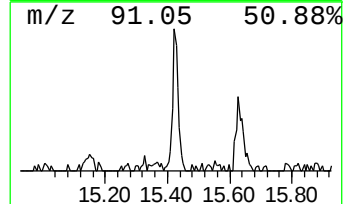
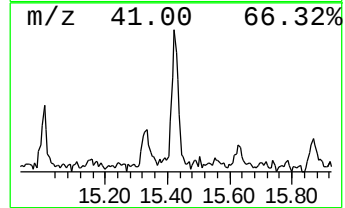
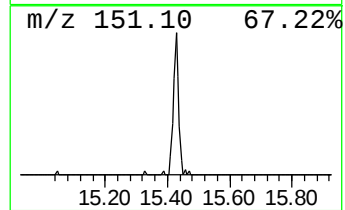
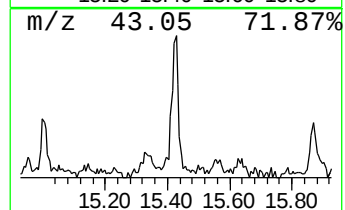
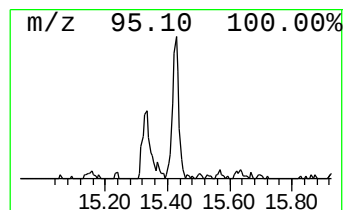
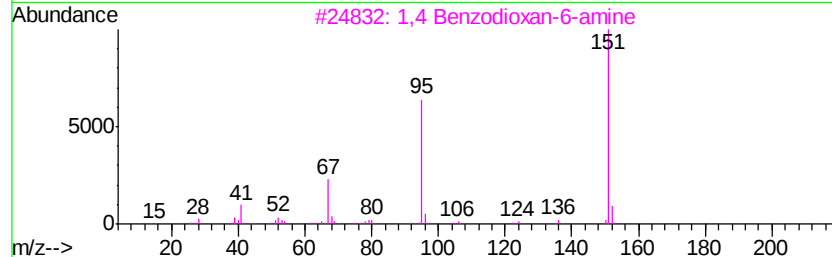
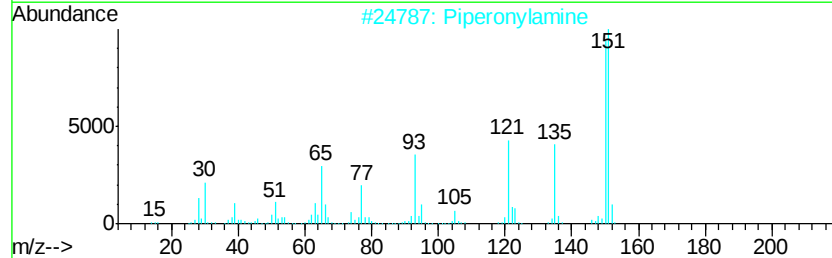
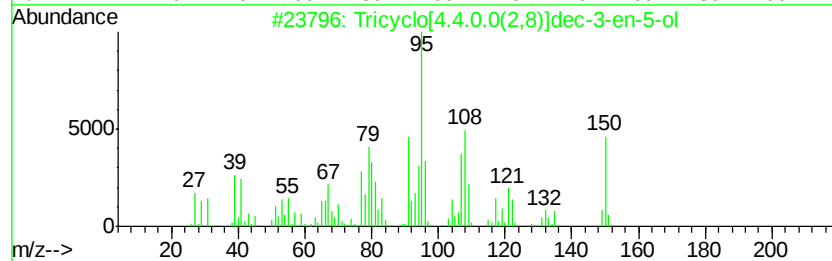
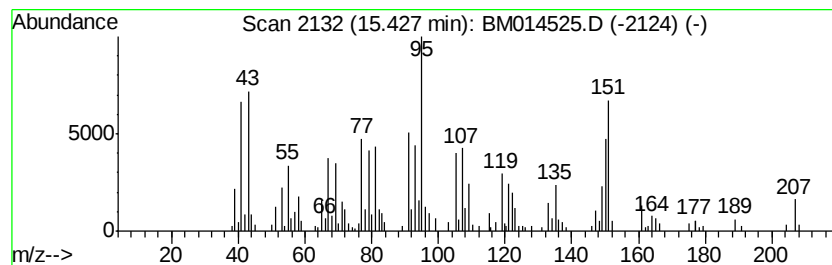
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Peak Number 3 unknown-01 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.43	10.71 ng/ul	417112	Acenaphthene-d10	14.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tricyclo[4.4.0.0(2,8)]dec-3-en-5-ol	150	C10H14O	1000193-38-7	46
2		Piperonylamine	151	C8H9NO2	002620-50-0	30
3		1,4 Benzodioxan-6-amine	151	C8H9NO2	022013-33-8	22
4		Naphthalene, decahydro-2,2-dimet...	166	C12H22	031124-85-3	22
5		1,4-Cyclohexanedimethanol	144	C8H16O2	000105-08-8	22



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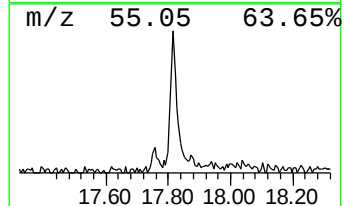
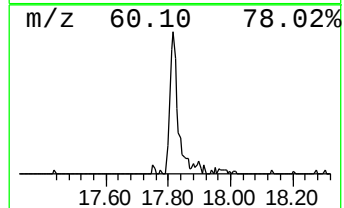
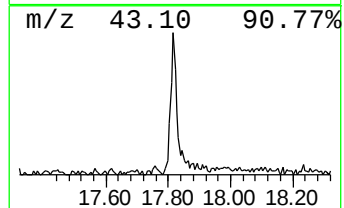
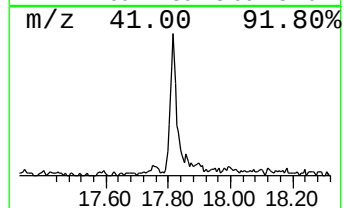
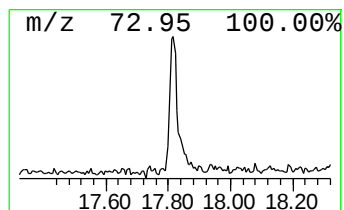
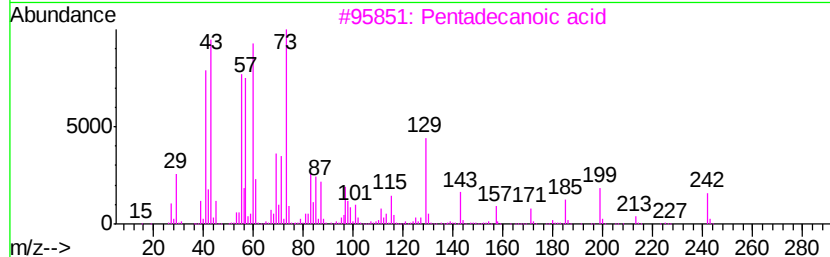
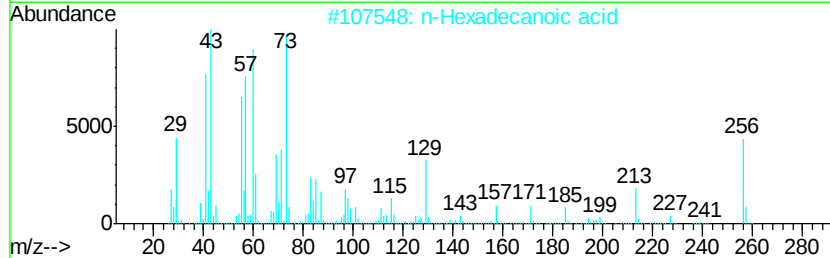
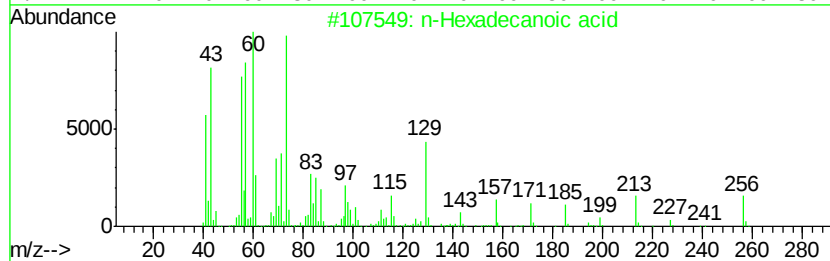
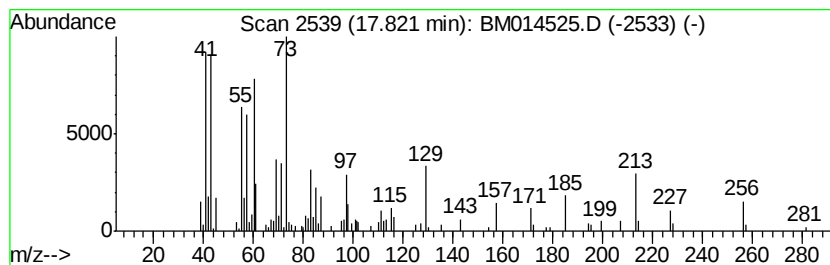
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Peak Number 4 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.82	6.79 ng/ul	353745	Phenanthrene-d10	16.91

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	76
2		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	70
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	62
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	62
5		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	60



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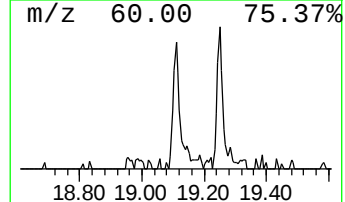
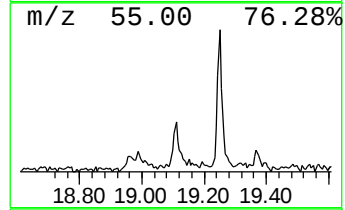
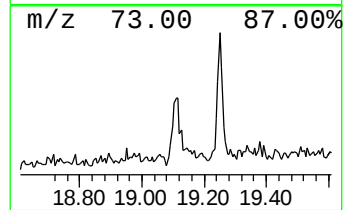
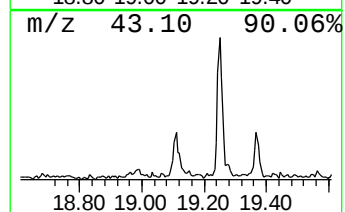
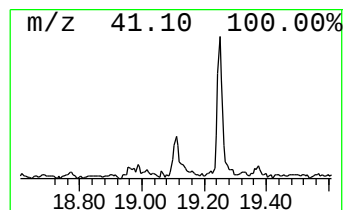
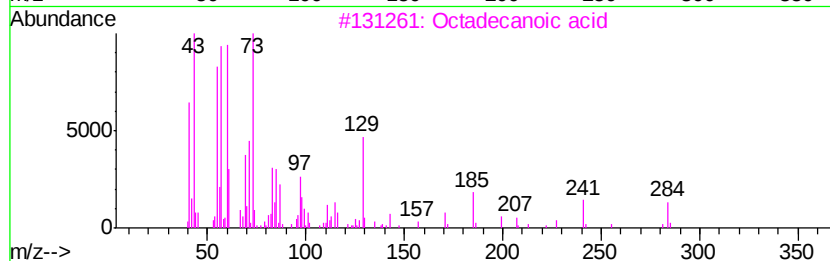
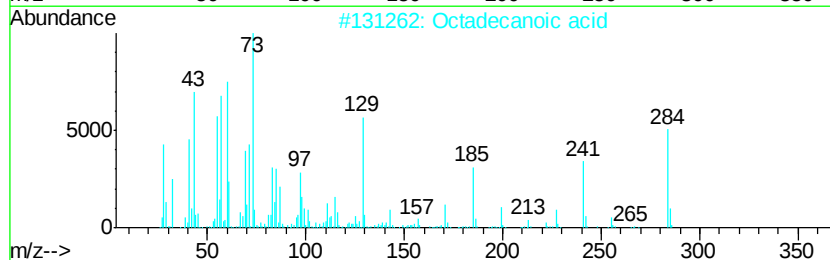
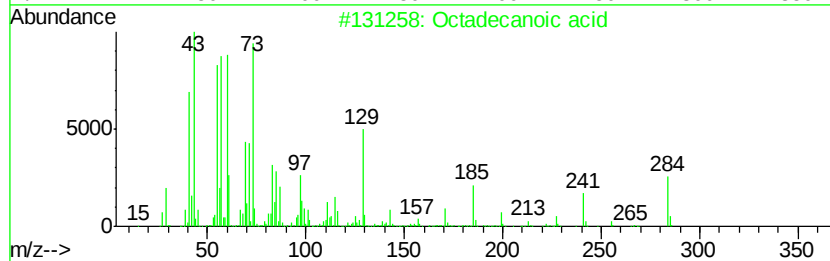
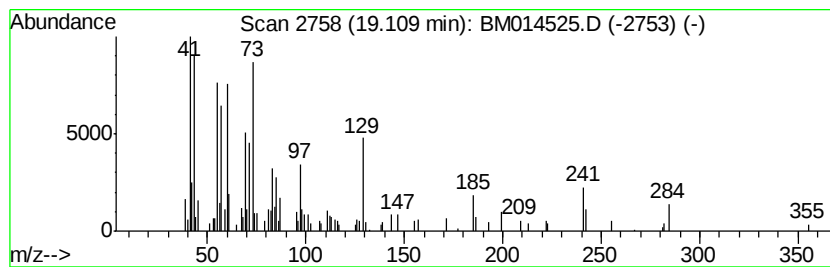
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Octadecanoic acid Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.11	2.61 ng/ul	193057	Chrysene-d12	21.13

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid		284	C18H36O2	000057-11-4	99
2	Octadecanoic acid		284	C18H36O2	000057-11-4	98
3	Octadecanoic acid		284	C18H36O2	000057-11-4	98
4	Octadecanoic acid		284	C18H36O2	000057-11-4	93
5	Pentadecanoic acid		242	C15H30O2	001002-84-2	76



Data Path : Z:\HPCHEM1\BNA M\DATA\BM031218\
Data File : BM014525.D
Acq On : 12 Mar 2018 19:46
Operator : SJ/JU
Sample : J1813-02
Misc :
ALS Vial : 9 Sample Multiplier: 1

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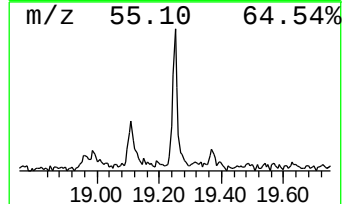
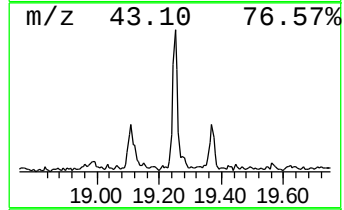
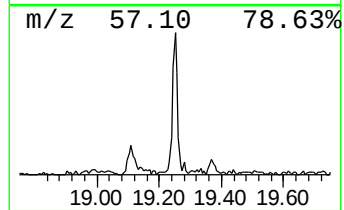
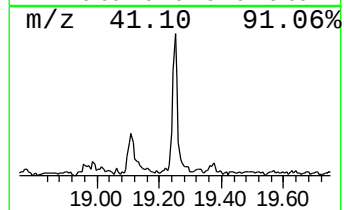
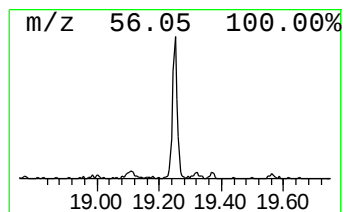
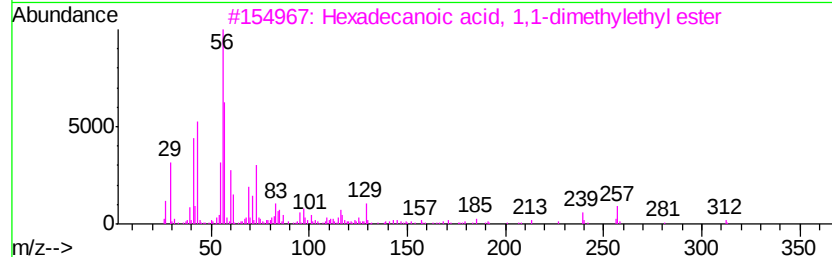
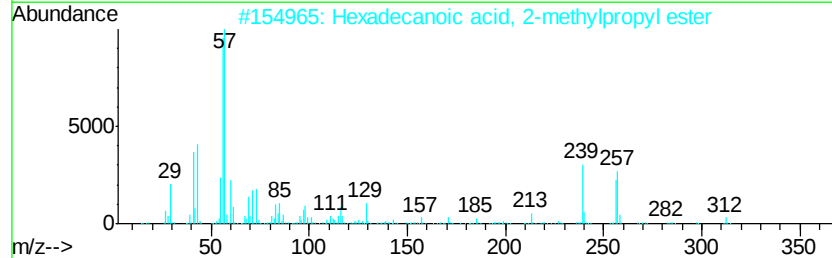
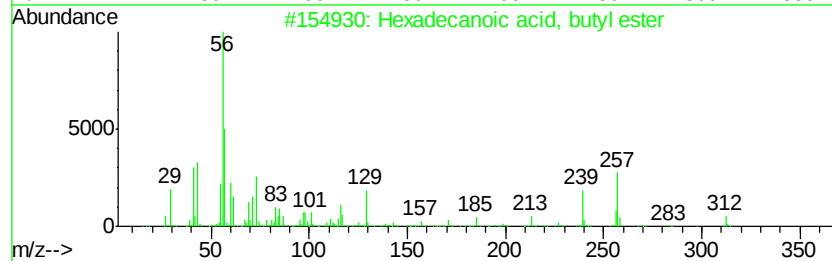
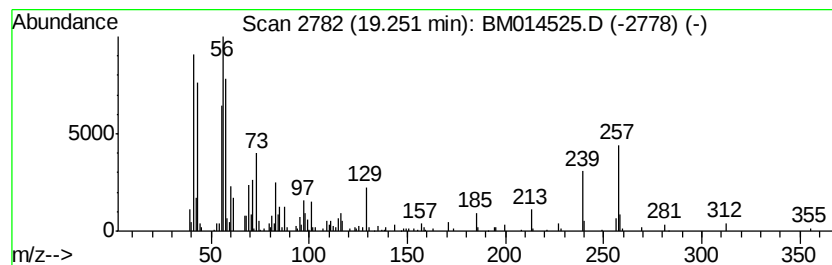
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Hexadecanoic acid, butyl ester Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.25	5.58 ng/ul	412948	Chrysene-d12	21.13

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecanoic acid, butyl ester	312	C20H40O2	000111-06-8	93
2		Hexadecanoic acid, 2-methylpropyl ester	312	C20H40O2	000110-34-9	93
3		Hexadecanoic acid, 1,1-dimethylethyl ester	312	C20H40O2	031158-91-5	55
4		Hexadecanoic acid, butyl ester	312	C20H40O2	000111-06-8	48
5		Butyl 4,8,12-trimethyl-tridecanoate	312	C20H40O2	1000336-63-2	38



Data Path : Z:\HPCHEM1\BNA M\DATA\BM031218\
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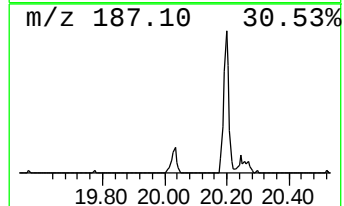
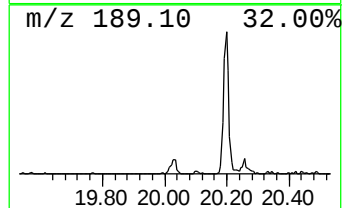
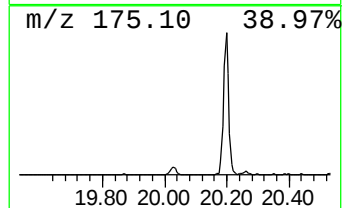
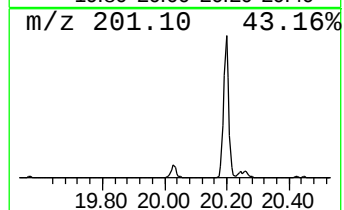
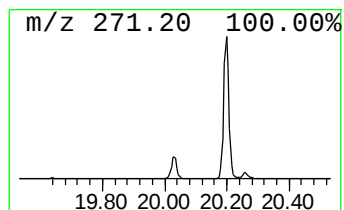
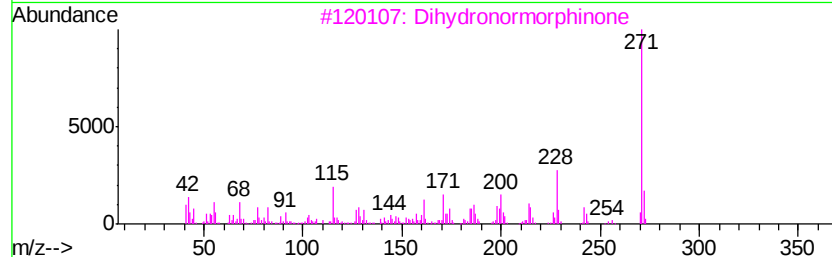
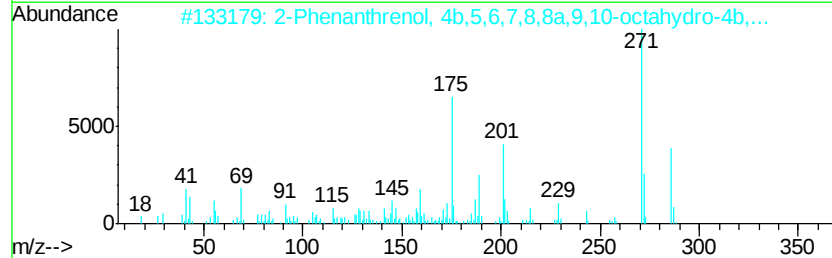
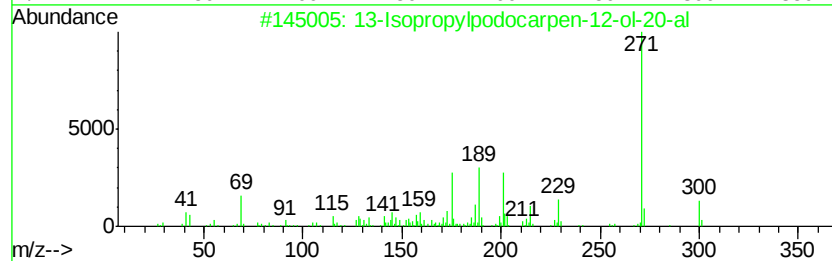
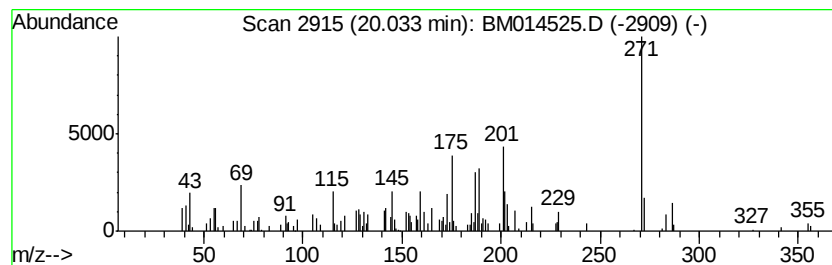
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 unknown-02 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.03	2.28 ng/ul	168610	Chrysene-d12	21.13

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	13-Isopropylpodocarpin-12-ol-20-al	300	C20H28O2	024035-37-8	47
2		2-Phenanthrenol, 4b,5,6,7,8,8a,9...	286	C20H30O	000511-15-9	35
3		Dihydronormorphinone	271	C16H17NO3	014696-23-2	35
4		Crinan-1-one	271	C16H17NO3	098301-98-5	35
5		Tridecanoic acid, tert-butyldime...	328	C19H40O2Si	104255-78-9	27



Data Path : Z:\HPCHEM1\BNA M\DATA\BM031218\
Data File : BM014525.D
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Misc :
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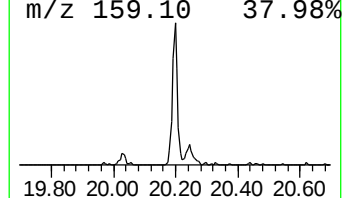
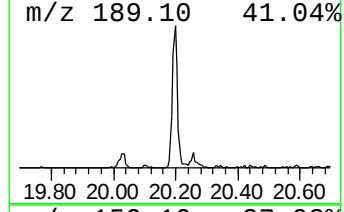
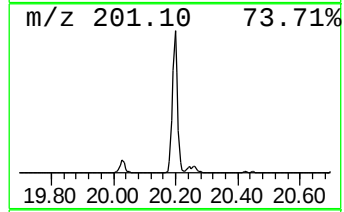
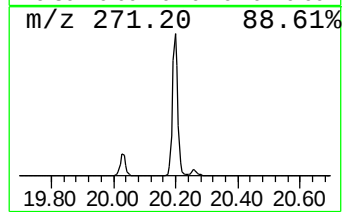
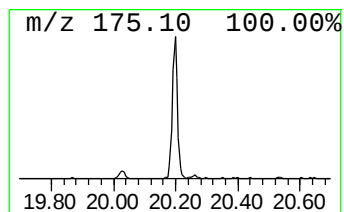
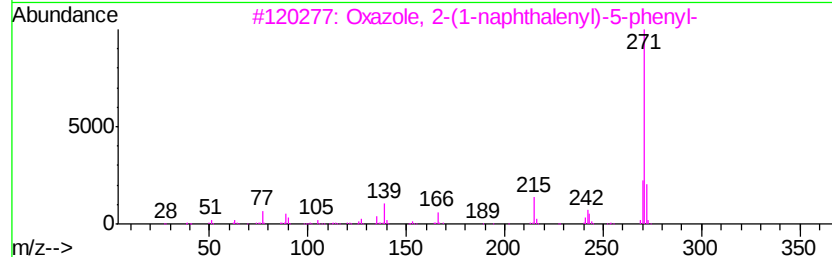
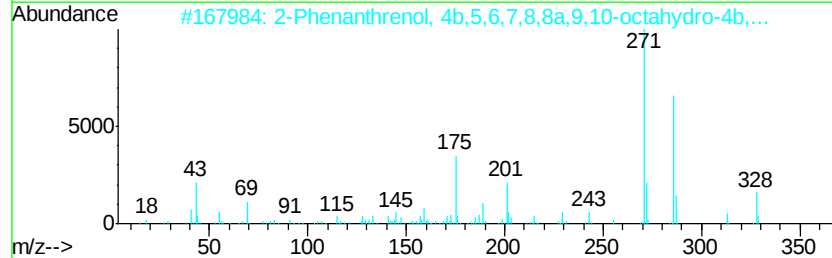
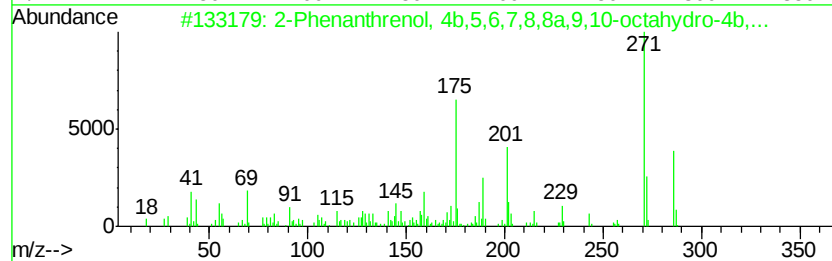
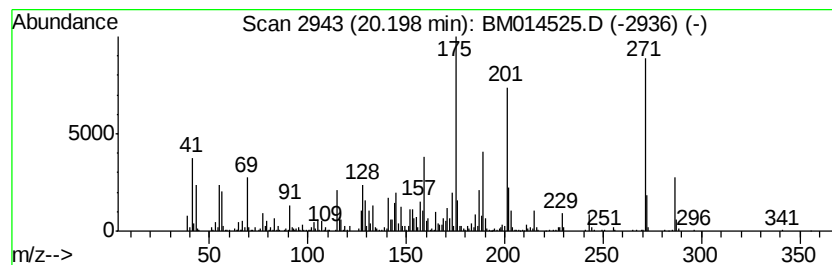
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 2-Phenanthrenol, 4b,5,6,7,8... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.20	17.43 ng/ul	1290810	Chrysene-d12	21.13

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Phenanthrenol, 4b,5,6,7,8,8a,9...	286	C20H30O	000511-15-9	94
2		2-Phenanthrenol, 4b,5,6,7,8,8a,9...	328	C22H32O2	015340-82-6	64
3		Oxazole, 2-(1-naphthalenyl)-5-ph...	271	C19H13NO	000846-63-9	15
4		Dihydroapohemanthamine	271	C16H17NO3	001153-01-1	14
5		7-Amino-4-methyl-1,8-naphthyridi...	175	C9H9N3O	001569-15-9	11



Data Path : Z:\HPCHEM1\BNA M\DATA\BM031218\
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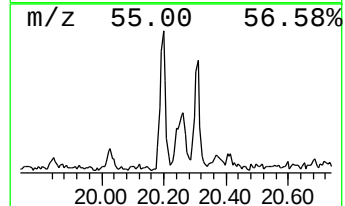
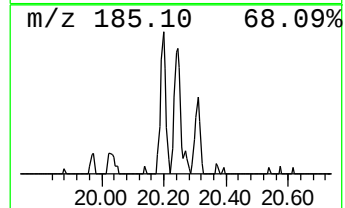
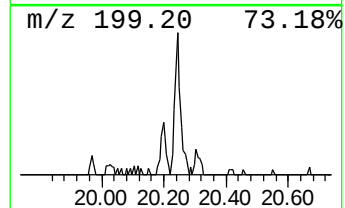
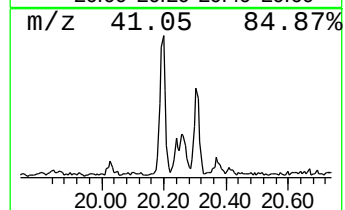
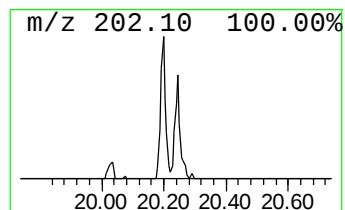
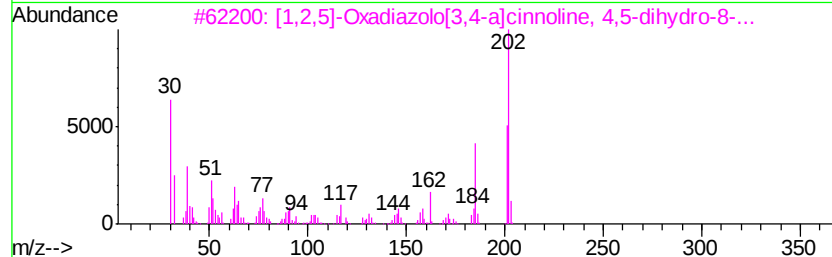
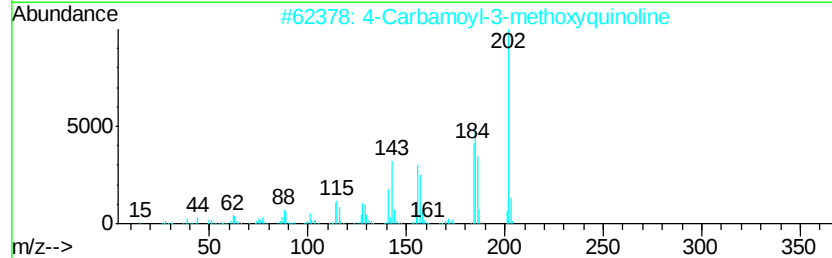
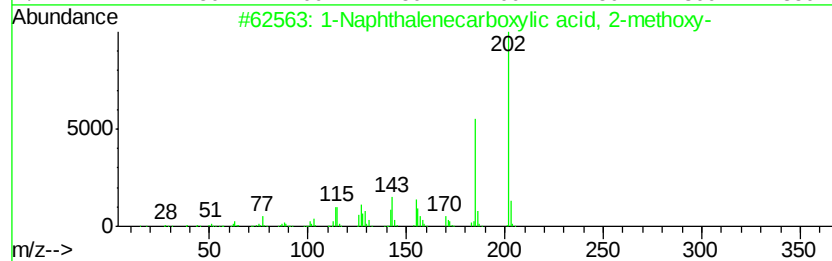
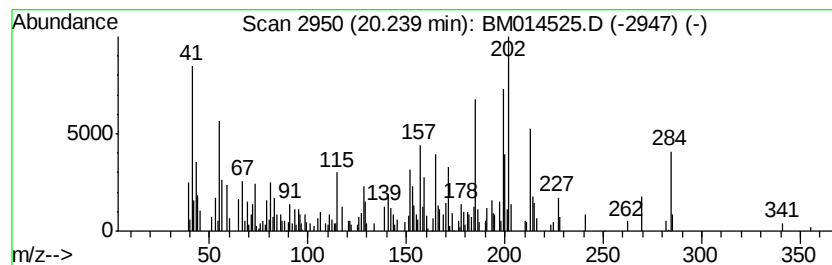
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 unknown-03 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.24	2.79 ng/ul	206893	Chrysene-d12	21.13

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Naphthalenecarboxylic acid, 2-...	202	C12H10O3	000947-62-6	46
2		4-Carbamoyl-3-methoxyquinoline	202	C11H10N2O2	020844-05-7	38
3		[1,2,5]-Oxadiazolo[3,4-a]cinnoli...	202	C10H10N4O	1000260-59-4	35
4		1-Naphthalenecarboxylic acid, 2-...	202	C12H10O3	000947-62-6	35
5		3,4-[Difluoromethylenedioxy]benz...	202	C8H4F2O4	000656-46-2	30



Data Path : Z:\HPCHEM1\BNA M\DATA\BM031218\
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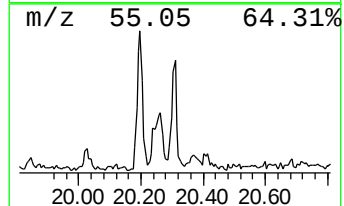
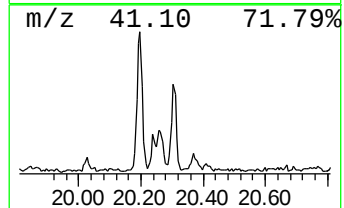
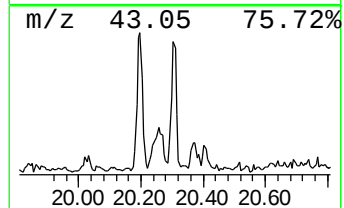
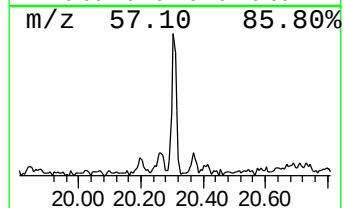
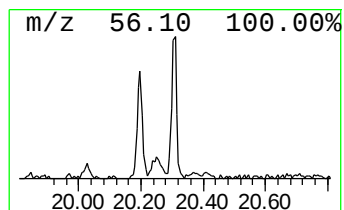
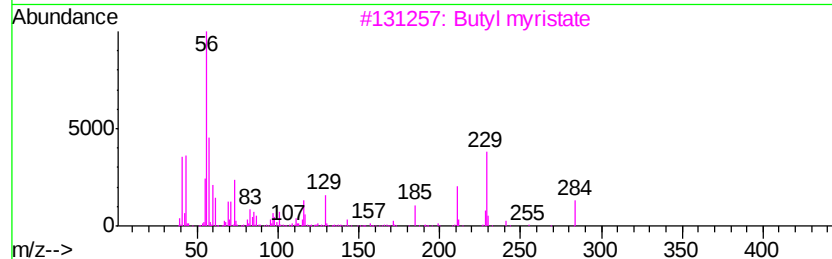
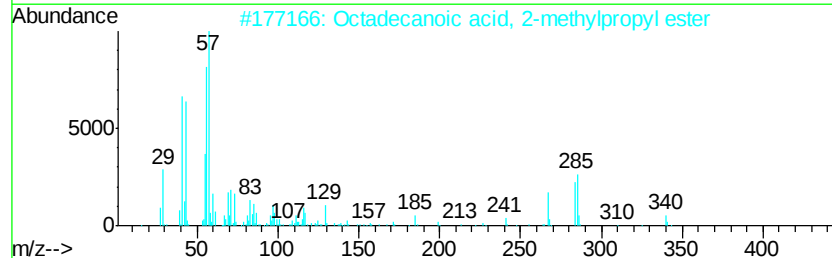
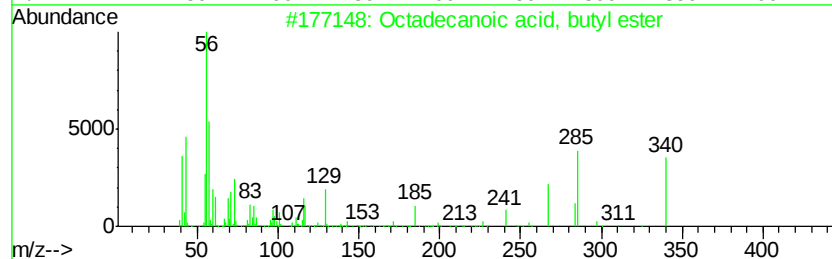
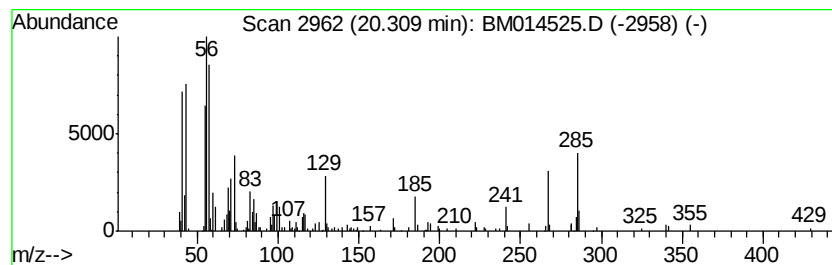
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Octadecanoic acid, butyl ester Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.31	3.71 ng/ul	274802	Chrysene-d12	21.13

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanoic acid, butyl ester	340	C22H44O2	000123-95-5	89
2		Octadecanoic acid, 2-methylpropyl ester	340	C22H44O2	000646-13-9	58
3		Butyl myristate	284	C18H36O2	000110-36-1	38
4		Piperidin-4-carboxylic acid	129	C6H11NO2	000498-94-2	35
5		Nipecotic acid	129	C6H11NO2	000498-95-3	35



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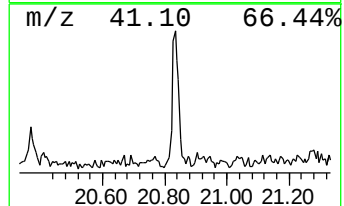
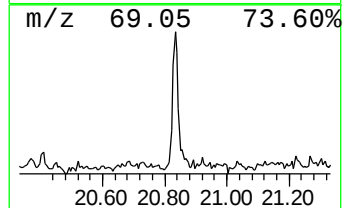
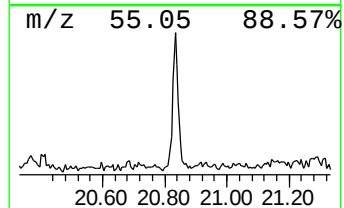
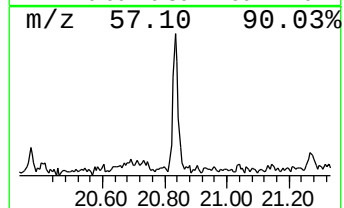
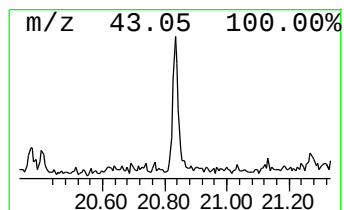
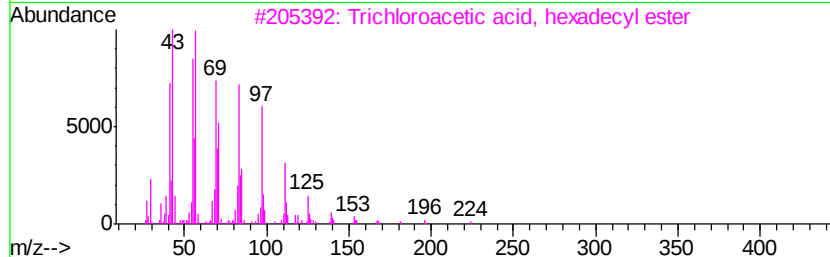
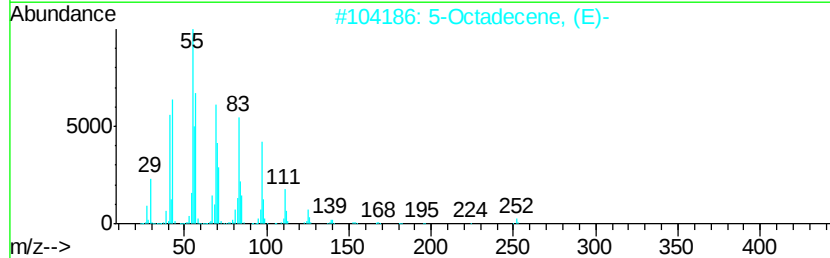
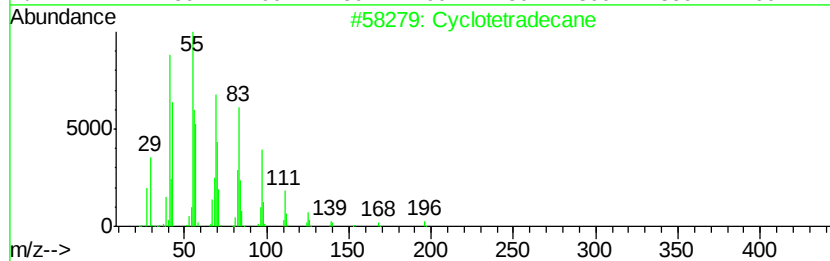
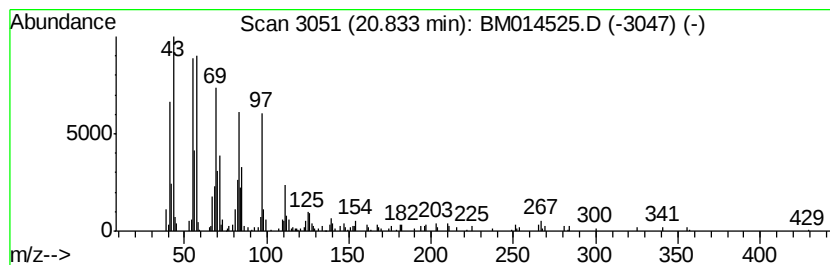
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 (DEL) Alkane: Cyclic20.83 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.83	3.99 ng/ul	295235	Chrysene-d12	21.13

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclotetradecane	196	C14H28	000295-17-0	93
2		5-Octadecene, (E)-	252	C18H36	007206-21-5	93
3		Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	91
4		Cyclopentadecane	210	C15H30	000295-48-7	89
5		9-Eicosene, (E)-	280	C20H40	074685-29-3	87



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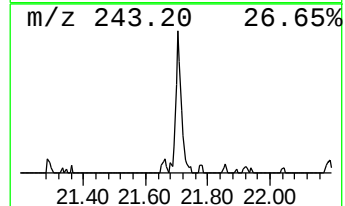
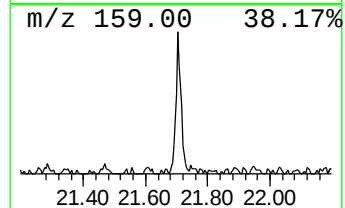
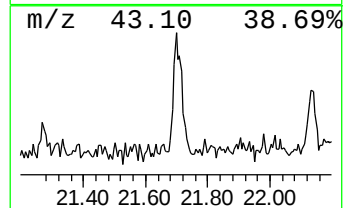
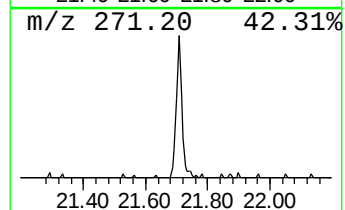
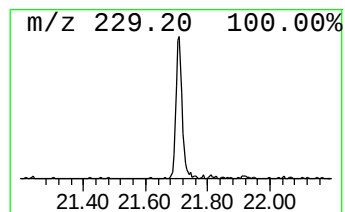
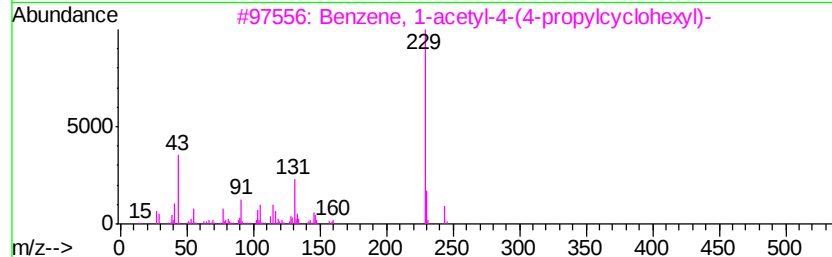
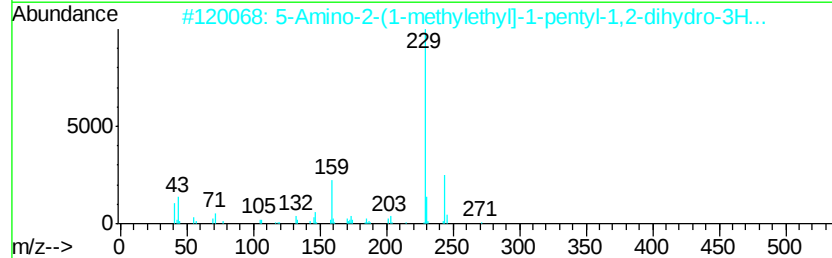
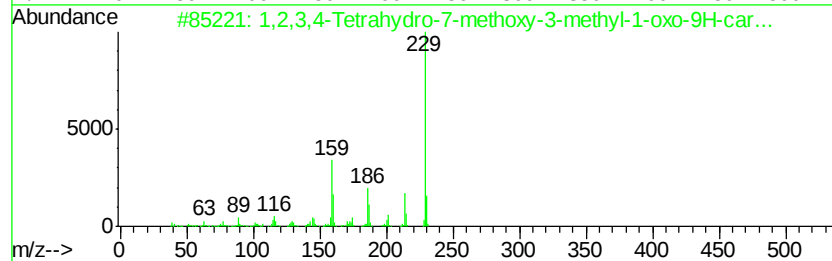
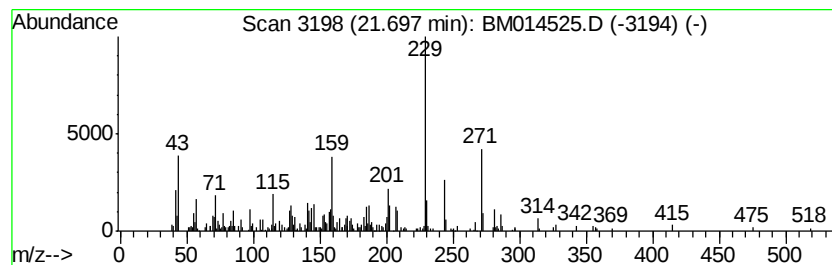
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 unknown-04 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.70	6.56 ng/ul	485944	Chrysene-d12	21.13

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,2,3,4-Tetrahydro-7-methoxy-3-m...	229	C14H15NO2	032550-51-9	38		
2	5-Amino-2-(1-methylethyl)-1-pent...	271	C15H21N5	1000326-95-8	32		
3	Benzene, 1-acetyl-4-(4-propylcyc...	244	C17H24O	078531-61-0	27		
4	1,3-Dihydro-1-(1-methyl-2-oxo-3-...	229	C13H11NO3	003988-53-2	27		
5	Cyclopent-2-ene-1-thioacetic aci...	286	C14H26O2SSi	068931-44-2	27		



Data Path : Z:\HPCHEM1\BNA M\DATA\BM031218\
Data File : BM014525.D
Acq On : 12 Mar 2018 19:46
Operator : SJ/JU
Sample : J1813-02
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampled :
F6F17

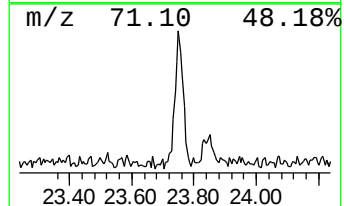
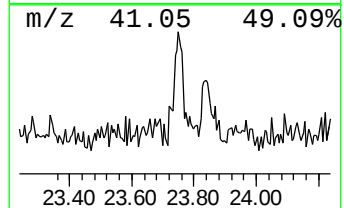
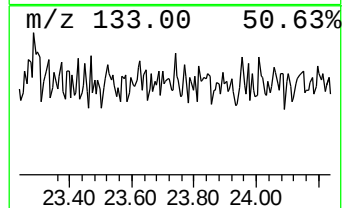
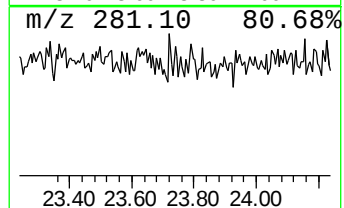
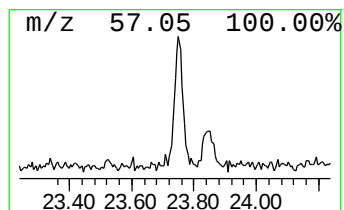
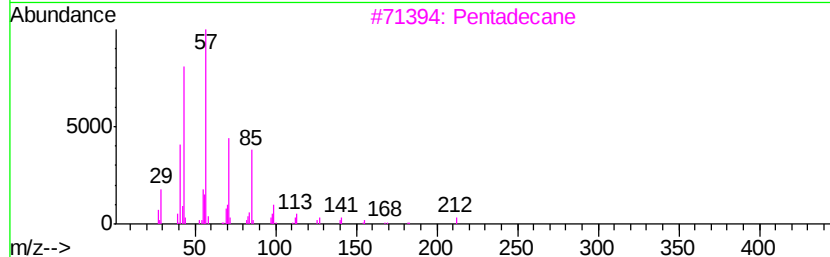
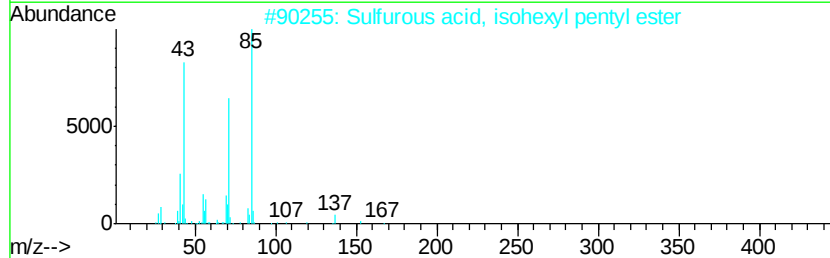
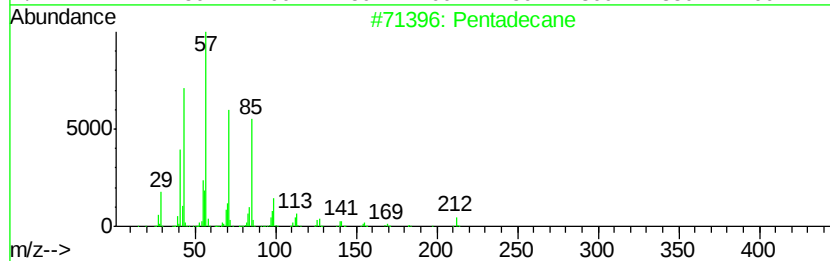
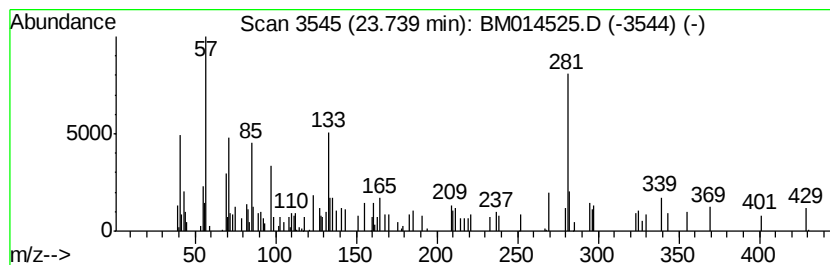
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 (DEL) Alkane: Straight-Chai... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.74	2.21 ng/ul	145200	Perylene-d12	23.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentadecane	212	C15H32	000629-62-9	18
2		Sulfurous acid, isohexyl pentyl ...	236	C11H24O3S	1000309-14-0	14
3		Pentadecane	212	C15H32	000629-62-9	14
4		3,5-Heptanedione, 4-ethyl-2,2,6,...	212	C13H24O2	167545-33-7	14
5		Tetracosane	338	C24H50	000646-31-1	14



Data Path : Z:\HPCHEM1\BNA M\DATA\BM031218\
Data File : BM014525.D
Acq On : 12 Mar 2018 19:46
Operator : SJ/JU
Sample : J1813-02
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
F6F17

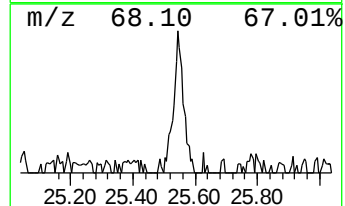
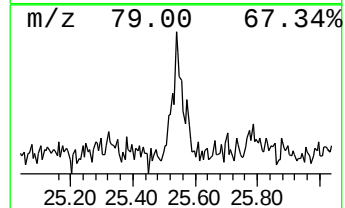
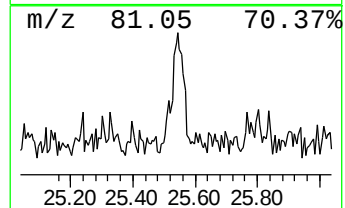
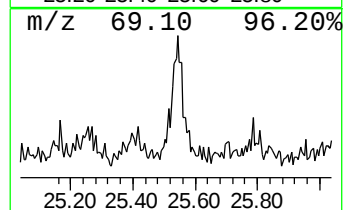
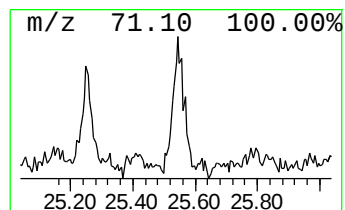
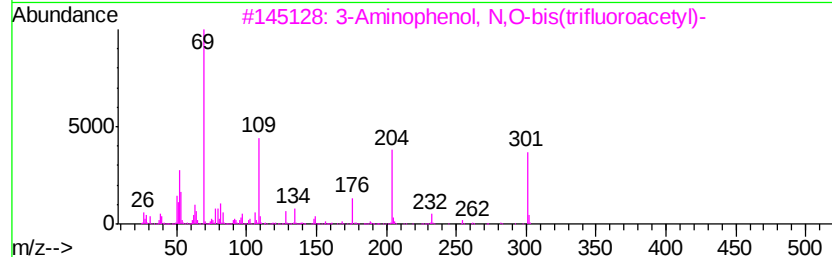
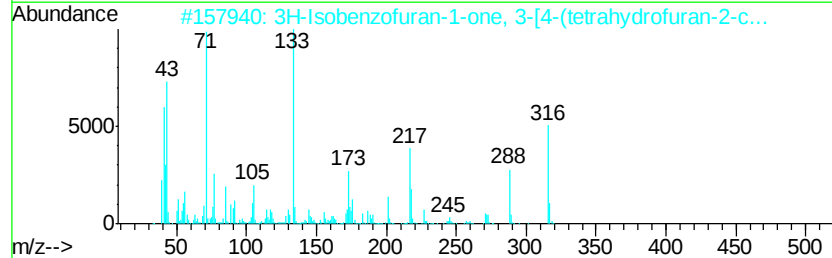
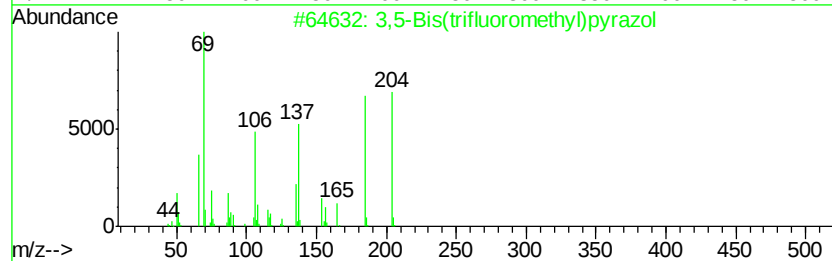
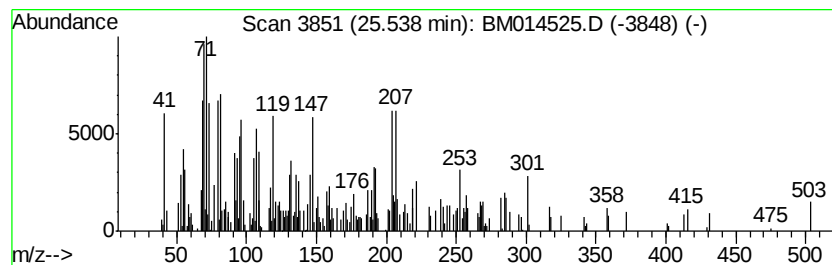
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 unknown-05 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.54	7.25 ng/ul	476465	Perylene-d12	23.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3,5-Bis(trifluoromethyl)pyrazol	204	C5H2F6N2	014704-41-7	25
2		3H-Isobenzofuran-1-one, 3-[4-(te...	316	C17H20N2O4	1000302-63-2	25
3		3-Aminophenol, N,O-bis(trifluoro...	301	C10H5F6N03	959257-40-0	22
4		3,7,11,15-Tetramethylhexadeca-1,...	290	C20H34O	068931-30-6	22
5		Himachala-2,4-diene	204	C15H24	060909-27-5	15



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM031218\
 Data File : BM014525.D
 Acq On : 12 Mar 2018 19:46
 Operator : SJ/JU
 Sample : J1813-02
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 F6F17

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM021418.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
Longifolene	13.37	11.4	ng/ul	442425	3	14.14	779102	20.0
Tricyclo[5.4.0.0(...	13.43	5.2	ng/ul	201123	3	14.14	779102	20.0
unknown-01	15.43	10.7	ng/ul	417112	3	14.14	779102	20.0
n-Hexadecanoic acid	17.82	6.8	ng/ul	353745	4	16.91	1042450	20.0
Octadecanoic acid	19.11	2.6	ng/ul	193057	5	21.13	1481170	20.0
Hexadecanoic acid...	19.25	5.6	ng/ul	412948	5	21.13	1481170	20.0
unknown-02	20.03	2.3	ng/ul	168610	5	21.13	1481170	20.0
2-Phenanthrenol, ...	20.20	17.4	ng/ul	1290810	5	21.13	1481170	20.0
unknown-03	20.24	2.8	ng/ul	206893	5	21.13	1481170	20.0
Octadecanoic acid...	20.31	3.7	ng/ul	274802	5	21.13	1481170	20.0
(DEL) Alkane: Cyc...	20.83	4.0	ng/ul	295235	5	21.13	1481170	20.0
unknown-04	21.70	6.6	ng/ul	485944	5	21.13	1481170	20.0
(DEL) Alkane: Str...	23.74	2.2	ng/ul	145200	6	23.30	1314870	20.0
unknown-05	25.54	7.3	ng/ul	476465	6	23.30	1314870	20.0