

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA M\DATA\BM120717\
 Data File : BM013254.D
 Acq On : 07 Dec 2017 21:39
 Operator : SJ/JU
 Sample : I6719-01
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 JHPH9

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-BM113017.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	3.240	23	26	33	rVB3	48170	62205	1.41%	0.102%
2	4.857	296	301	306	rVB2	42360	63362	1.43%	0.104%
3	6.875	638	644	654	rVB2	964560	1687748	38.15%	2.770%
4	7.039	661	672	679	rBV	746043	1116056	25.23%	1.832%
5	7.234	698	705	713	rVB	1045695	1630559	36.86%	2.676%
6	7.698	776	784	792	rBV	845189	1324193	29.94%	2.174%
7	8.404	896	904	910	rBV	1214820	1881636	42.54%	3.089%
8	8.463	910	914	922	rVB2	84184	140355	3.17%	0.230%
9	8.851	972	980	988	rBV	954627	1545923	34.95%	2.538%
10	9.263	1044	1050	1056	rVB4	24839	44888	1.01%	0.074%
11	9.569	1092	1102	1115	rBV	807913	1388449	31.39%	2.279%
12	10.104	1184	1193	1200	rBV	1219419	1985405	44.88%	3.259%
13	10.475	1249	1256	1264	rBV	983729	1687400	38.15%	2.770%
14	10.616	1269	1280	1293	rBV	839385	1517421	34.30%	2.491%
15	13.751	1807	1813	1818	rBV	1846909	2589651	58.54%	4.251%
16	13.798	1818	1821	1829	rVB	504307	671267	15.18%	1.102%
17	14.033	1851	1861	1867	rBV	2115281	3219002	72.77%	5.284%
18	14.186	1881	1887	1892	rBV4	37843	86297	1.95%	0.142%
19	14.245	1893	1897	1902	rVB	71396	97855	2.21%	0.161%
20	14.339	1905	1913	1922	rBV2	1282888	1981098	44.79%	3.252%
21	14.551	1941	1949	1965	rBV	811618	1239297	28.02%	2.034%
22	14.763	1978	1985	1990	rBV10	30633	60787	1.37%	0.100%
23	14.815	1990	1994	1999	rVB2	35285	56344	1.27%	0.092%
24	15.145	2046	2050	2056	rBV3	40180	68389	1.55%	0.112%
25	15.198	2056	2059	2064	rVB2	71965	93523	2.11%	0.154%
26	15.339	2076	2083	2089	rVB	2476750	3601603	81.42%	5.912%
27	15.462	2098	2104	2111	rBV	679079	929881	21.02%	1.526%
28	15.574	2119	2123	2127	rBV2	34247	59849	1.35%	0.098%
29	16.062	2201	2206	2214	rBV	73767	118769	2.68%	0.195%
30	16.857	2335	2341	2346	rBV2	50180	76776	1.74%	0.126%
31	17.086	2374	2380	2388	rBV2	1553399	2204739	49.84%	3.619%
32	17.186	2390	2397	2404	rBV2	2603122	3742775	84.61%	6.143%
33	17.992	2529	2534	2541	rBV2	462779	637582	14.41%	1.047%
34	19.127	2721	2727	2729	rBV3	60010	97739	2.21%	0.160%

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35	19.274	2748	2752	2763	rBV2	441624	590353	13.35%	0.969%
36	19.486	2782	2788	2794	rBV	3127509	4195151	94.84%	6.886%
37	19.539	2794	2797	2802	rVB6	52071	72099	1.63%	0.118%
38	20.003	2873	2876	2880	rBV6	93621	135171	3.06%	0.222%
39	20.368	2935	2938	2941	rBV5	35623	50636	1.14%	0.083%
40	20.997	3041	3045	3053	rVB2	425235	530927	12.00%	0.871%
41	21.274	3087	3092	3102	rBV	1864357	2419822	54.70%	3.972%
42	21.874	3189	3194	3203	rVB2	749280	1130661	25.56%	1.856%
43	21.986	3209	3213	3219	rBV	534405	722929	16.34%	1.187%
44	22.050	3221	3224	3231	rVB7	160363	227212	5.14%	0.373%
45	22.315	3265	3269	3275	rVB4	261241	461349	10.43%	0.757%
46	22.562	3308	3311	3316	rVB5	161286	195219	4.41%	0.320%
47	22.838	3355	3358	3361	rVV2	149861	186941	4.23%	0.307%
48	22.880	3361	3365	3373	rVB	714306	1096347	24.78%	1.800%
49	23.368	3442	3448	3460	rVB2	2315370	4423498	100.00%	7.261%
50	23.509	3466	3472	3483	rVB	1213514	2327019	52.61%	3.820%
51	23.709	3504	3506	3512	rVB5	154520	242080	5.47%	0.397%
52	23.956	3545	3548	3558	rVB8	156536	298109	6.74%	0.489%
53	24.133	3572	3578	3588	rVB2	369938	799879	18.08%	1.313%
54	24.780	3683	3688	3696	rVB7	334173	757330	17.12%	1.243%
55	26.938	4048	4055	4065	rVB5	424149	1188578	26.87%	1.951%
56	27.297	4109	4116	4125	rVB5	385760	1162809	26.29%	1.909%

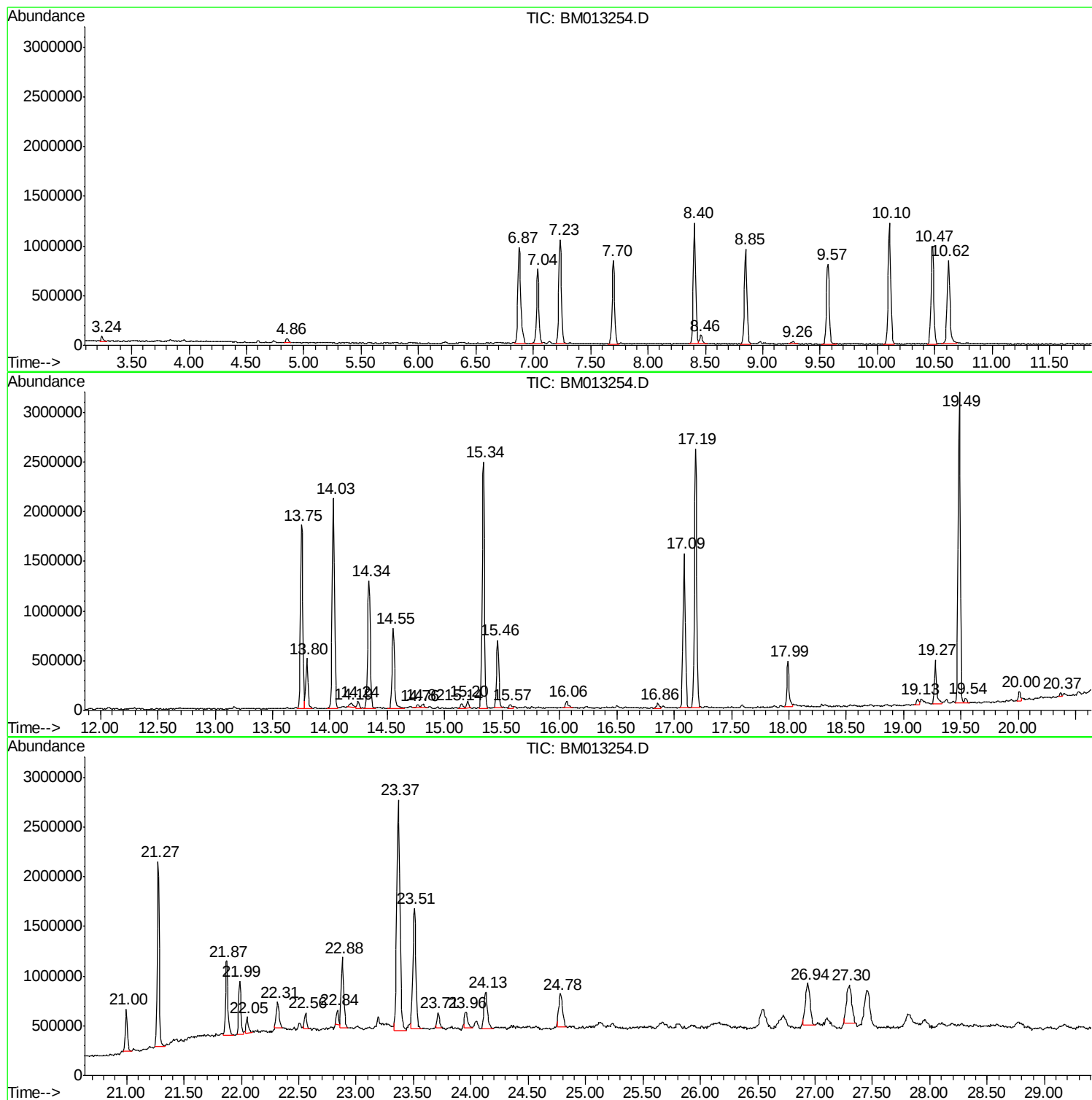
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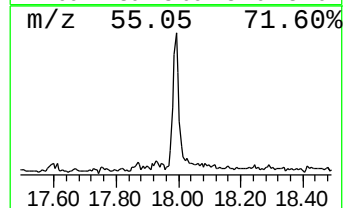
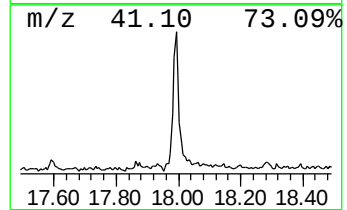
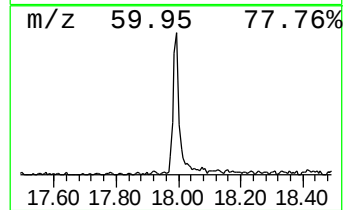
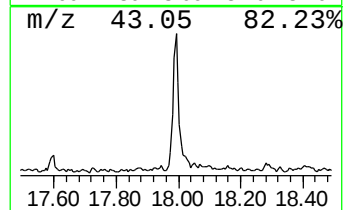
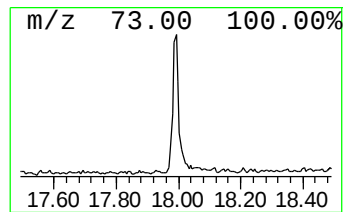
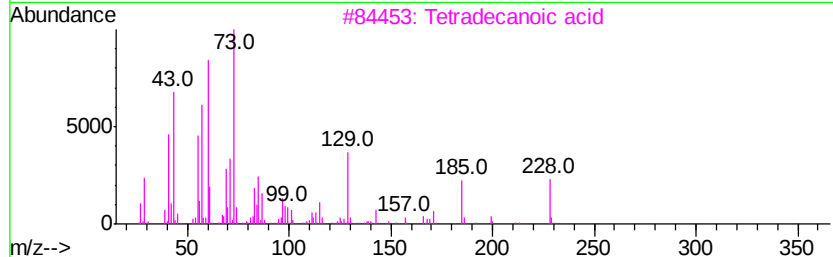
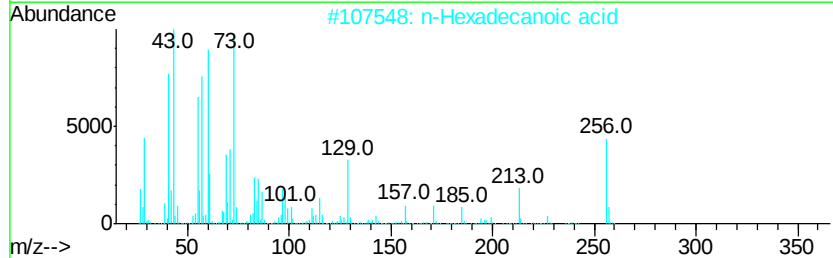
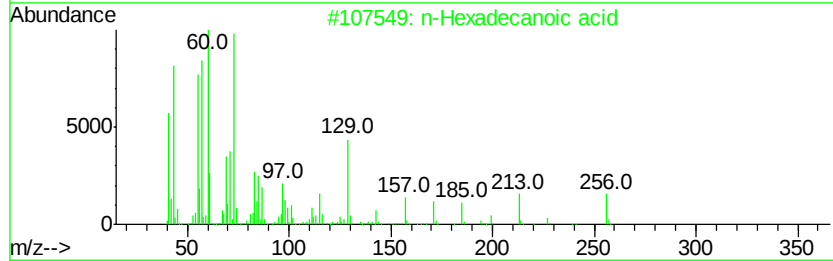
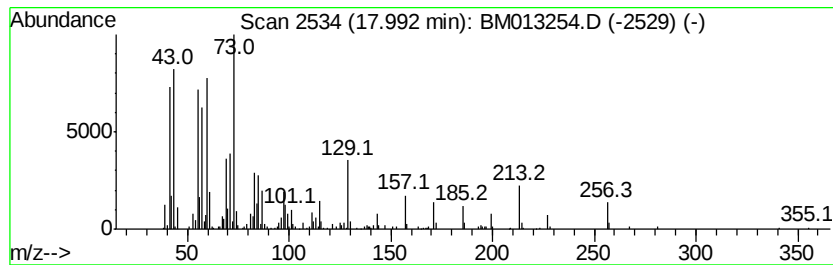
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Quant Title : SVOA CALIBRATION

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

Peak Number 1 n-Hexadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD		R.T.
17.99	5.78 ng/ul	637582	Phenanthrene-d10		17.09

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
3			Tetradecanoic acid	228	C14H28O2	000544-63-8	94
4			Octadecanoic acid	284	C18H36O2	000057-11-4	93
5			Tridecanoic acid	214	C13H26O2	000638-53-9	91



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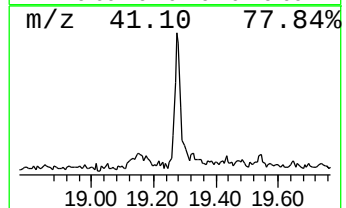
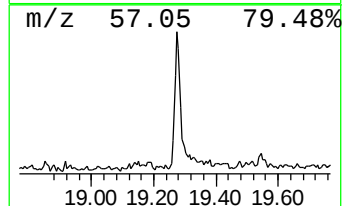
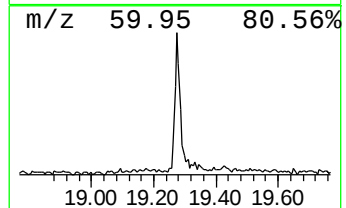
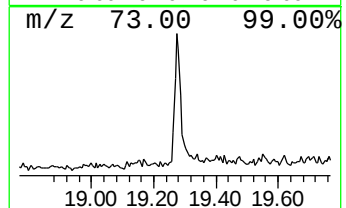
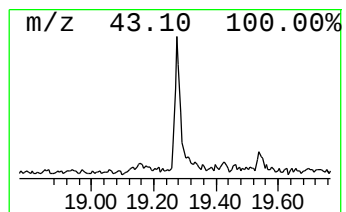
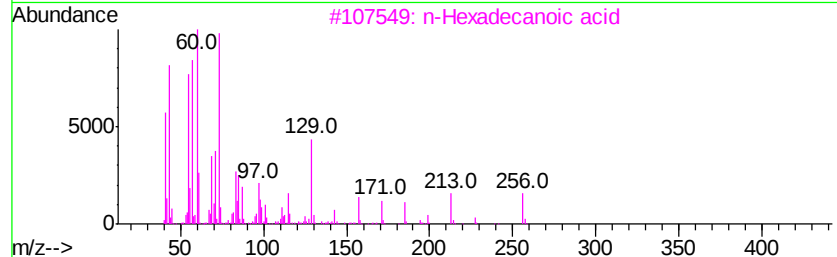
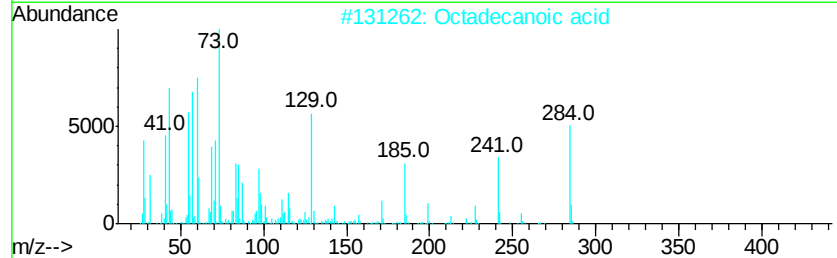
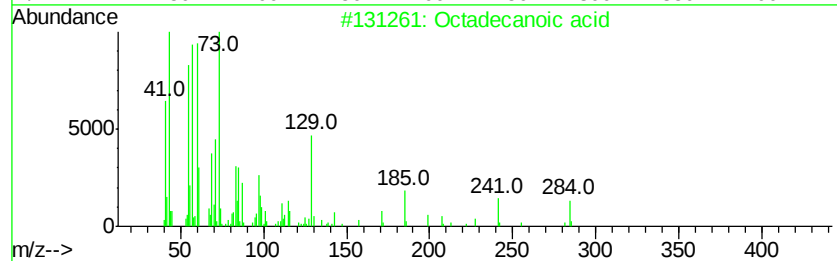
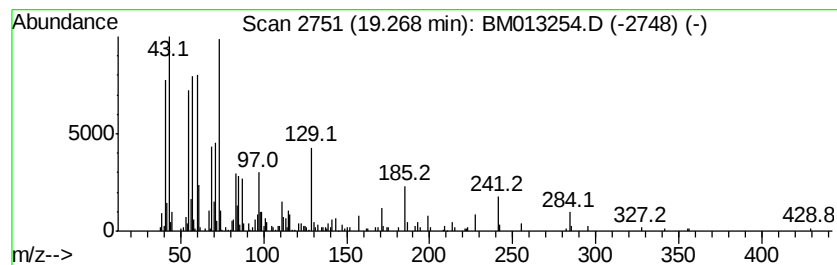
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM113017.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Octadecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.27	4.88 ng/ul	590353	Chrysene-d12	21.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2		Octadecanoic acid	284	C18H36O2	000057-11-4	99
3		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	94
4		Octadecanoic acid	284	C18H36O2	000057-11-4	94
5		Pentadecanoic acid	242	C15H30O2	001002-84-2	94



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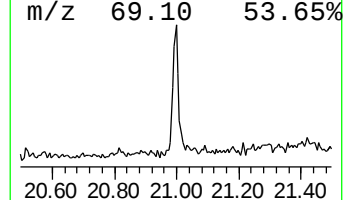
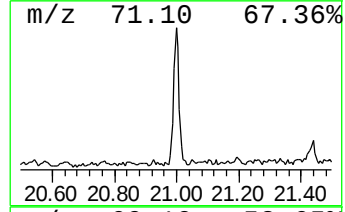
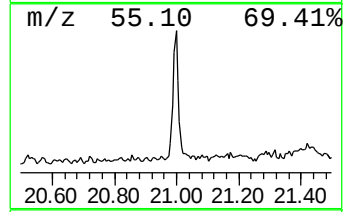
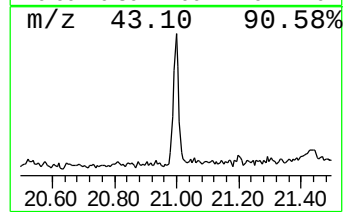
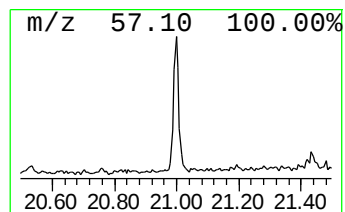
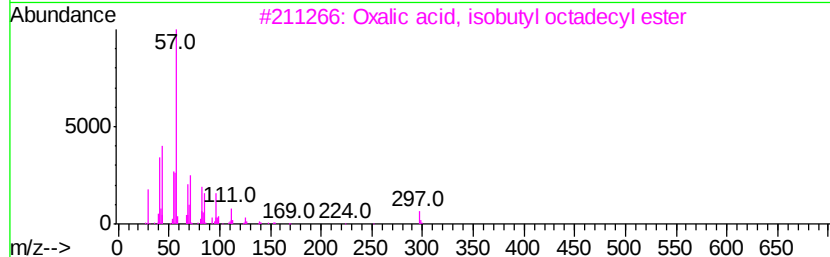
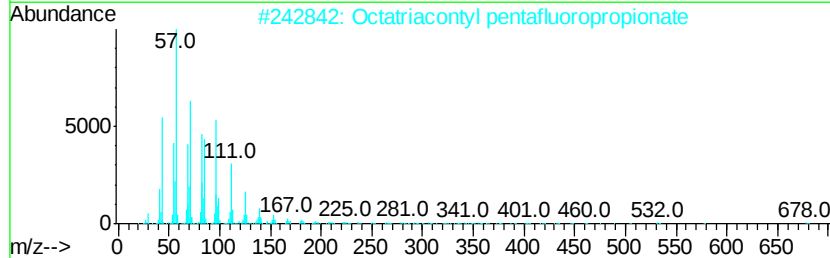
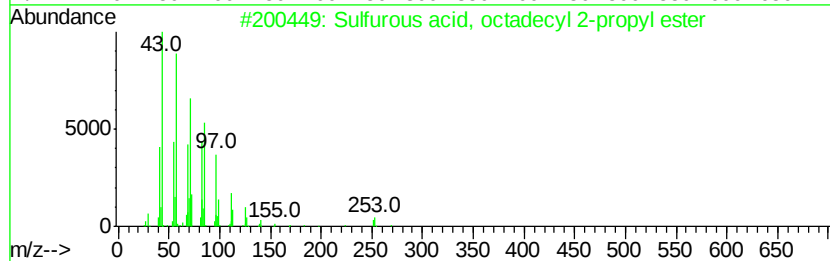
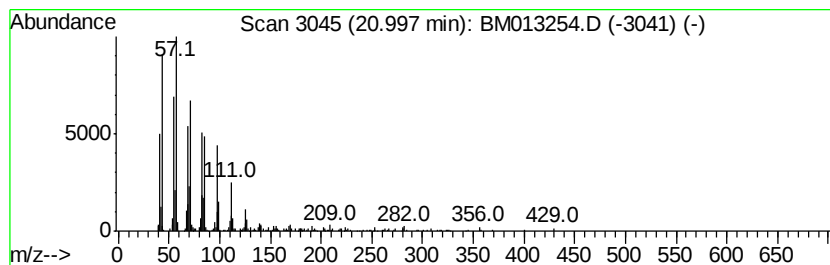
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Peak Number 3 Sulfurous acid, octadecyl 2... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.00	4.39 ng/ul	530927	Chrysene-d12	21.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Sulfurous acid, octadecyl 2-prop...	376	C21H44O3S	1000309-12-7	91
2		Octatriacontyl pentafluoropropio...	697	C41H77F5O2	1000351-89-1	91
3		Oxalic acid, isobutyl octadecyl ...	398	C24H46O4	1000309-38-3	91
4		Ethanol, 2-(hexadecyloxy)-	286	C18H38O2	002136-71-2	91
5		Oxalic acid, isobutyl hexadecyl ...	370	C22H42O4	1000309-38-1	91



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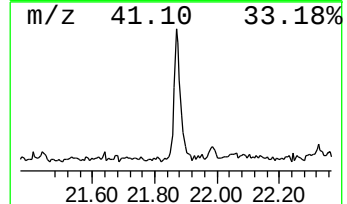
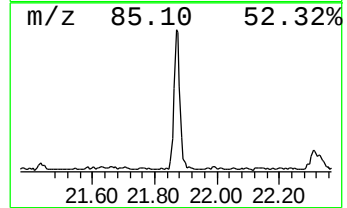
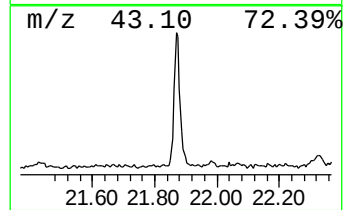
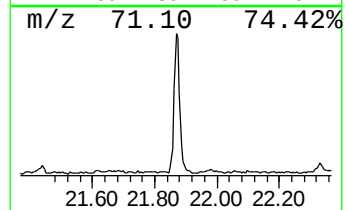
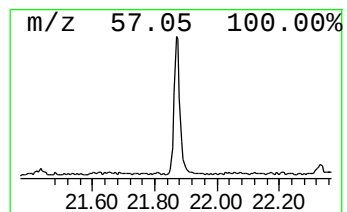
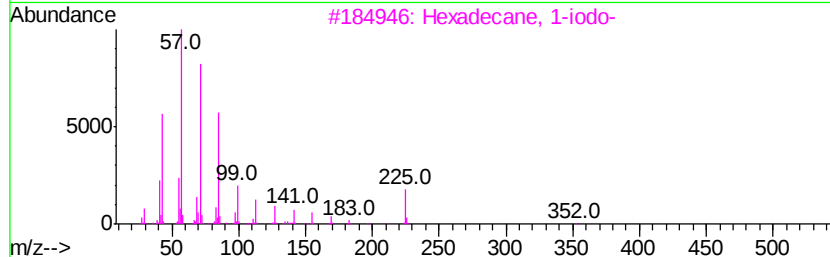
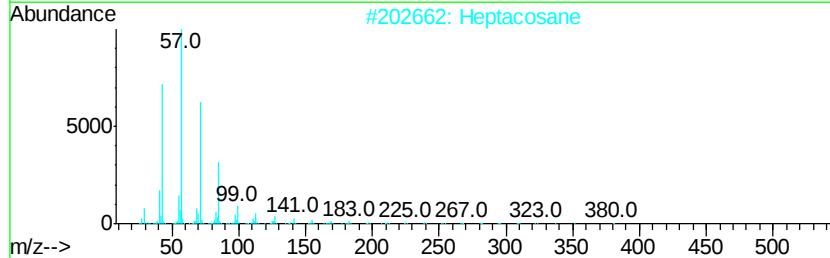
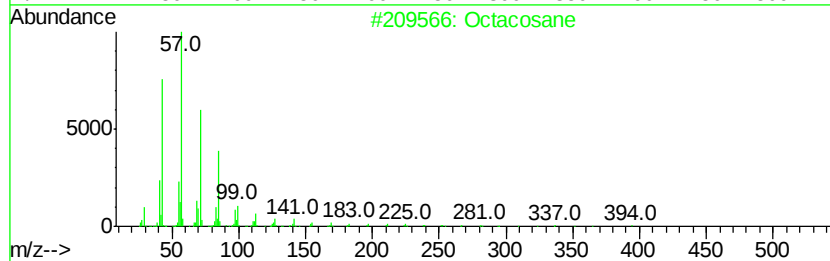
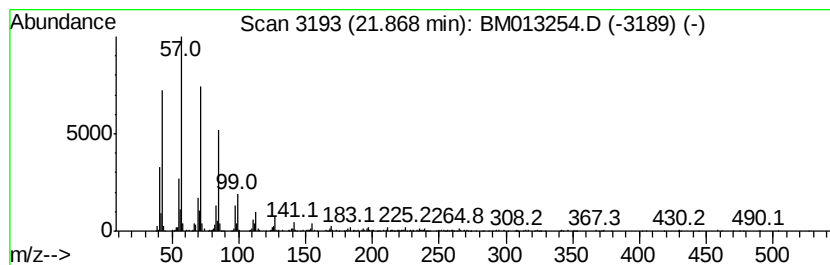
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Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 4

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21.87	9.34 ng/ul	1130660	Chrysene-d12	21.27

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octacosane	394	C28H58	000630-02-4	94
2			Heptacosane	380	C27H56	000593-49-7	93
3			Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	93
4			Heptadecane	240	C17H36	000629-78-7	93
5			Heptacosane	380	C27H56	000593-49-7	91



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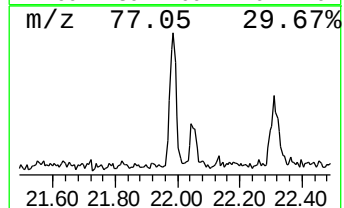
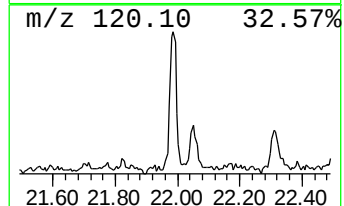
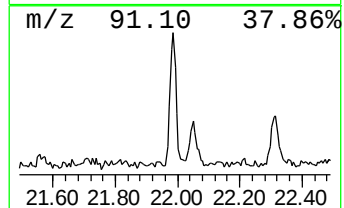
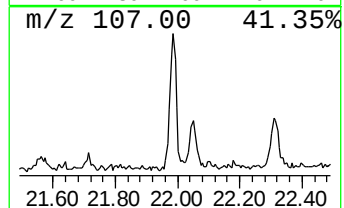
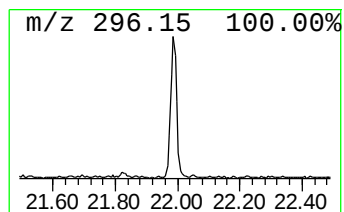
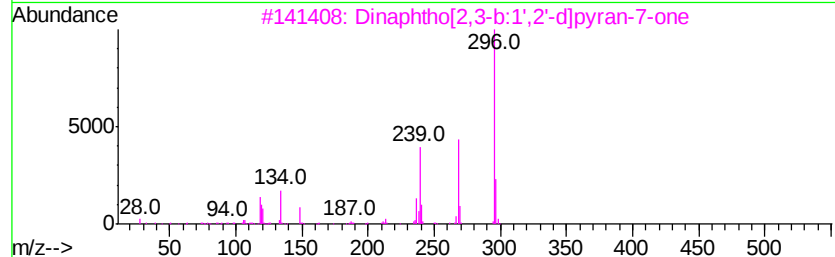
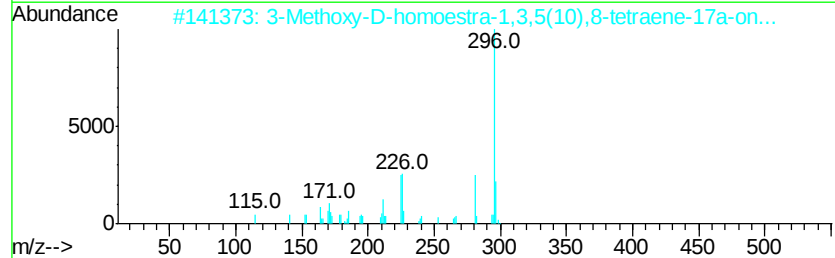
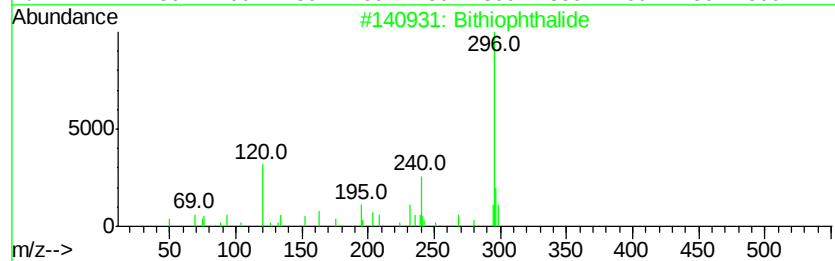
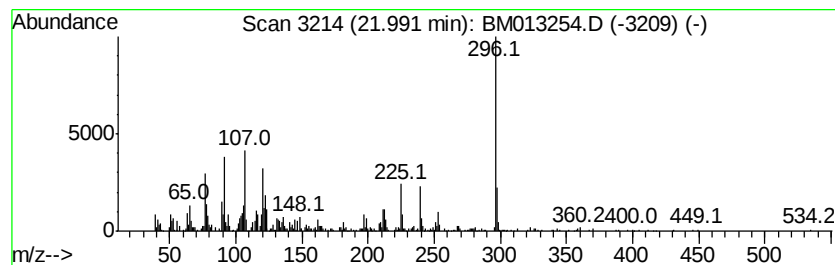
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 unknown-01 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.99	5.98 ng/ul	722929	Chrysene-d12	21.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Bithiophthalide	296	C16H8O2S2	032819-84-4	43
2		3-Methoxy-D-homoestra-1,3,5(10),...	296	C20H24O2	002000-72-8	40
3		Dinaphtho[2,3-b:1',2'-d]pyran-7-one	296	C21H12O2	154352-87-1	35
4		1H-Pyrazol-5-ol, 1,3-bis(4-metho...	296	C17H16N2O3	1000337-80-1	35
5		Thioindigo	296	C16H8O2S2	000522-75-8	32



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Data File : BM013254.D
Acq On : 07 Dec 2017 21:39
Operator : SJ/JU
Sample : I6719-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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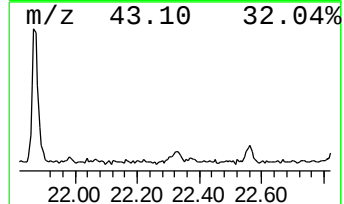
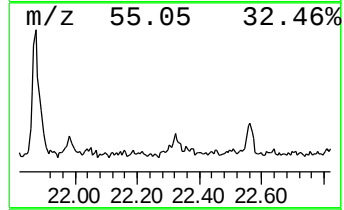
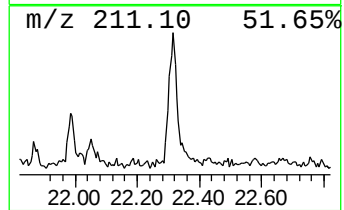
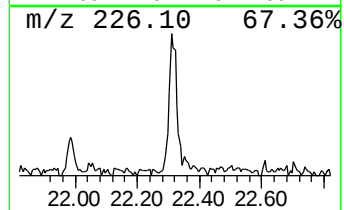
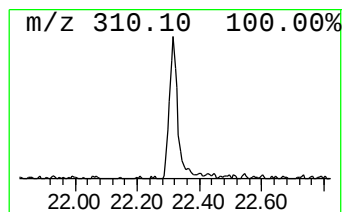
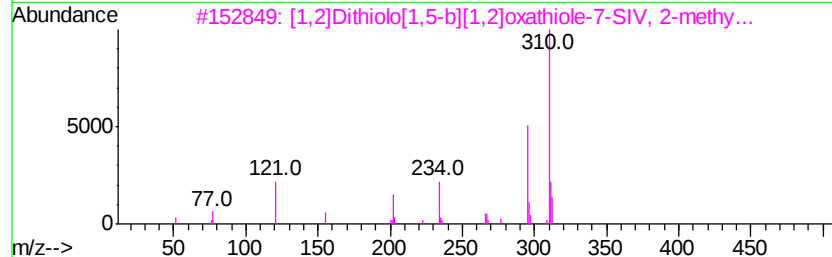
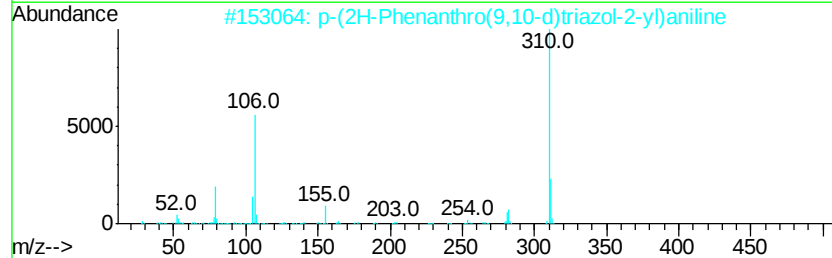
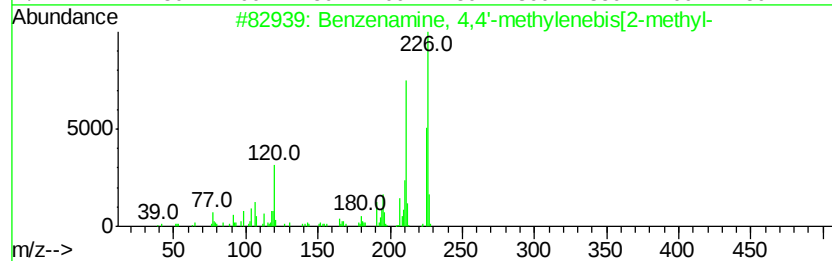
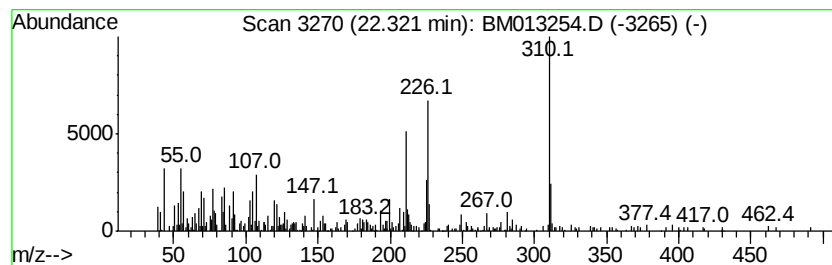
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Benzenamine, 4,4'-methylene... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.32	3.81 ng/ul	461349	Chrysene-d12	21.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzenamine, 4,4'-methylenebis[2...	226	C15H18N2	000838-88-0	50
2		p-(2H-Phenanthro(9,10-d)triazol-...	310	C20H14N4	000975-85-9	30
3		[1,2]Dithiolo[1,5-b][1,2]oxathio...	310	C18H14OS2	074810-20-1	25
4		Pvridine, 4-(1-methyl-2-pyrrolyl...	310	C22H18N2	005562-57-2	25
5		Ferrocene, 1,2-diacetyl-1',4-(1,...	310	C17H18FeO2	041558-93-4	25



Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA M\DATA\BM120717\
Data File : BM013254.D
Acq On : 07 Dec 2017 21:39
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Sample : I6719-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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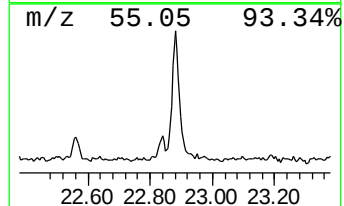
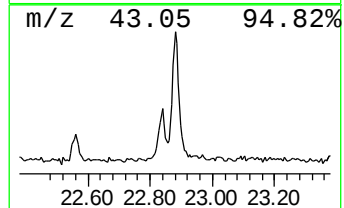
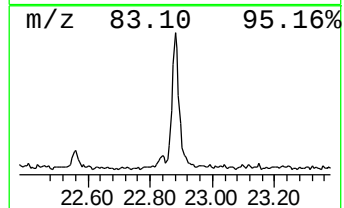
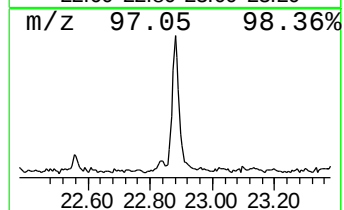
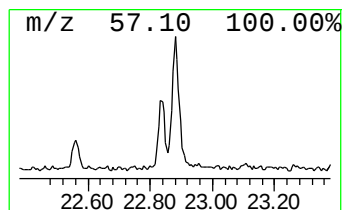
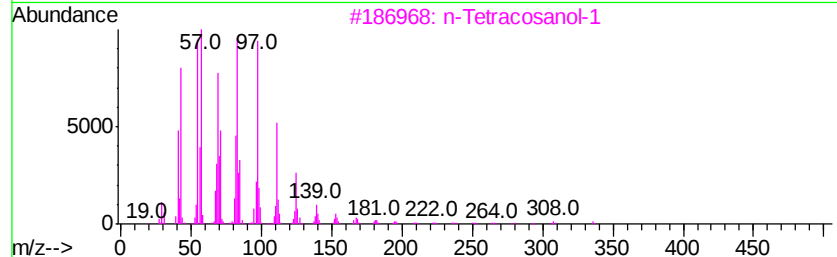
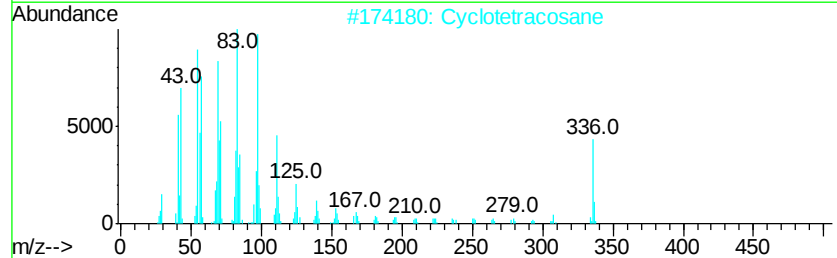
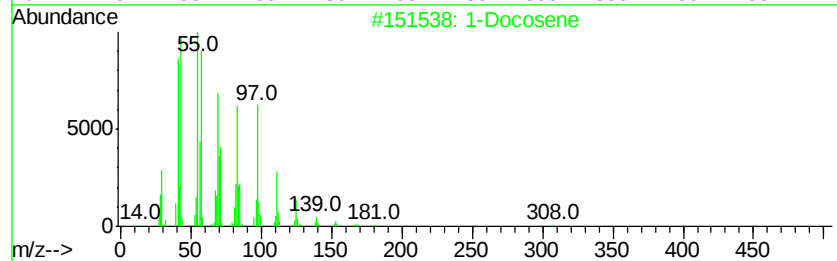
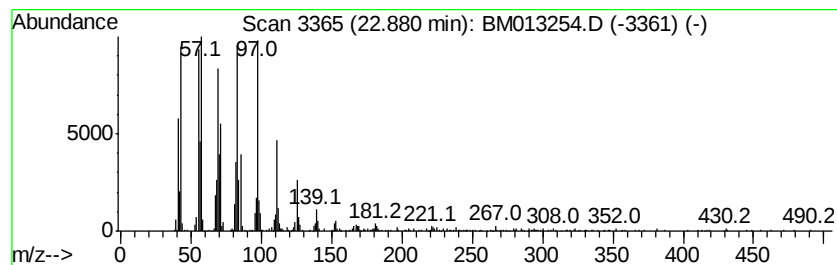
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Cyclic22.88 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.88	9.42 ng/ul	1096350	Perylene-d12	23.51

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Docosene	308	C22H44	001599-67-3	95
2			Cyclotetracosane	336	C24H48	000297-03-0	95
3			n-Tetracosanol-1	354	C24H50O	000506-51-4	94
4			1-Heptacosanol	396	C27H56O	002004-39-9	94
5			1-Heneicosanol	312	C21H44O	015594-90-8	93



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA M\DATA\BM120717\
Data File : BM013254.D
Acq On : 07 Dec 2017 21:39
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Sample : I6719-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

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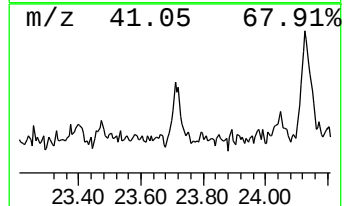
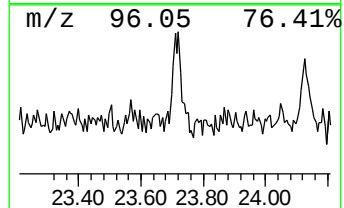
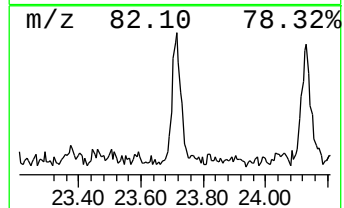
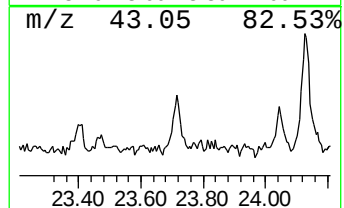
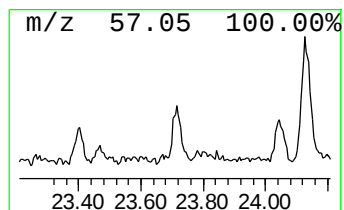
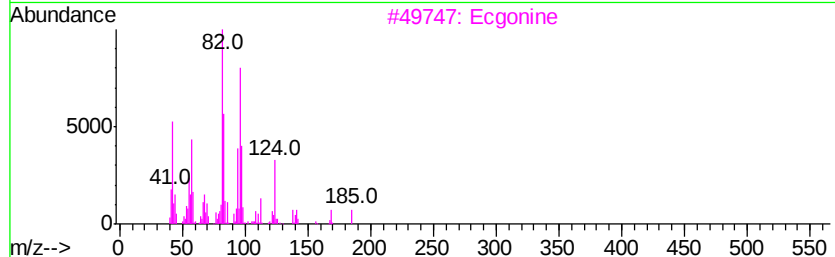
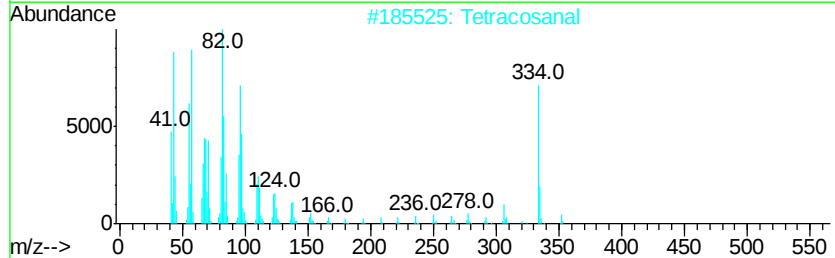
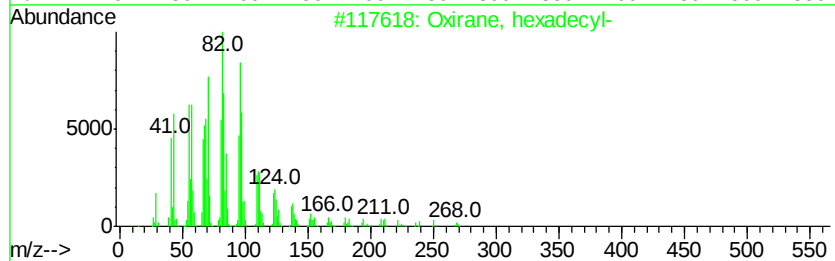
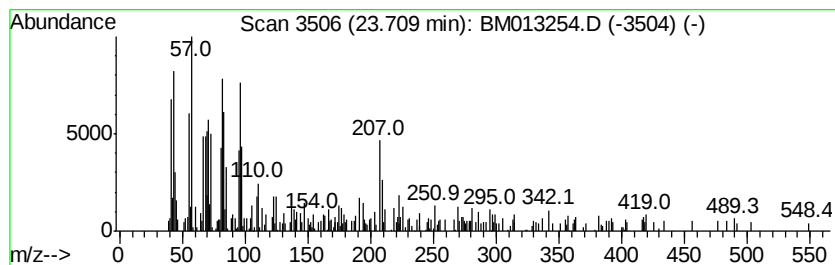
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Oxirane, hexadecyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.71	2.08 ng/ul	242080	Perylene-d12	23.51

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Oxirane, hexadecyl-	268	C18H36O	007390-81-0	68
2		Tetracosanal	352	C24H48O	057866-08-7	68
3		Ecgonine	185	C9H15NO3	000481-37-8	64
4		Oxirane, heptadecyl-	282	C19H38O	067860-04-2	64
5		1,19-Eicosadiene	278	C20H38	014811-95-1	64



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA M\DATA\BM120717\
Data File : BM013254.D
Acq On : 07 Dec 2017 21:39
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Sample : I6719-01
Misc :
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Instrument :
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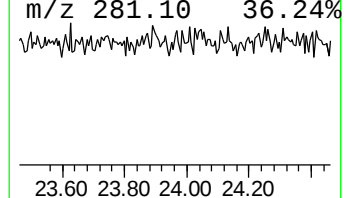
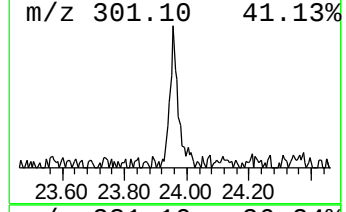
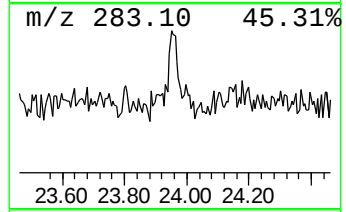
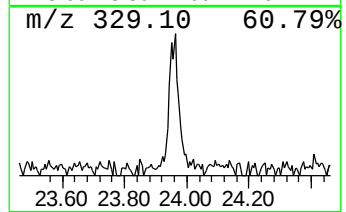
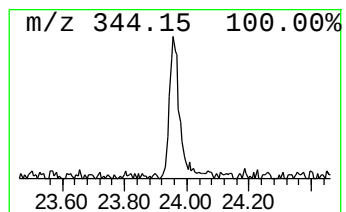
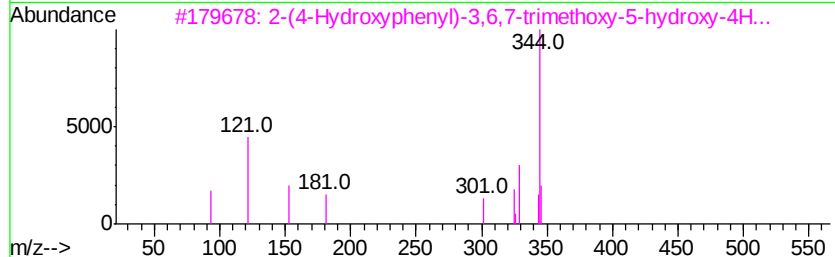
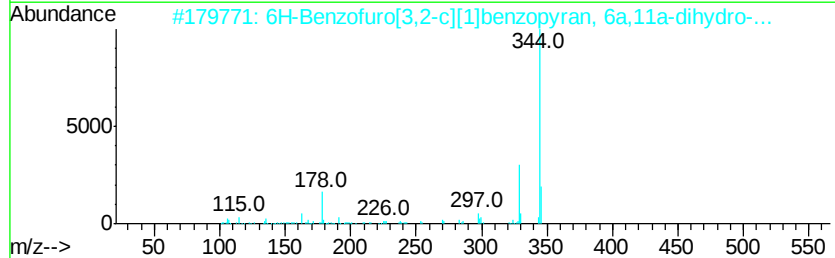
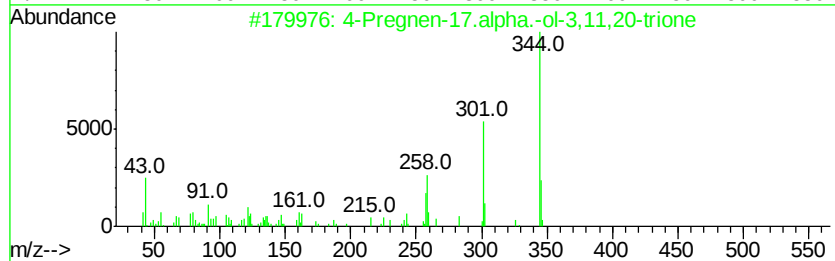
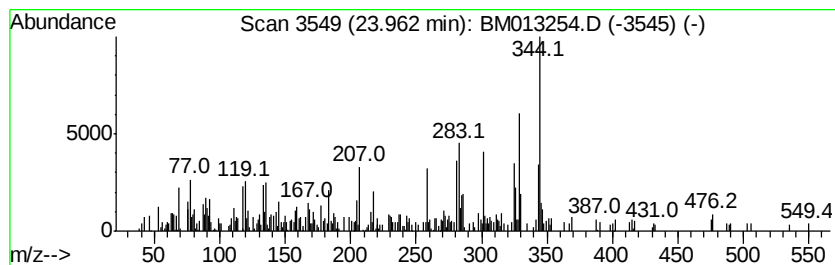
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 4-Pregnen-17.alpha.-ol-3,11... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.96	2.56 ng/ul	298109	Perylene-d12	23.51

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Pregnen-17.alpha.-ol-3,11,20-t...	344	C21H28O4	1000126-67-7	55
2			6H-Benzofuro[3,2-c][1]benzopyran...	344	C19H20O6	020390-16-3	38
3			2-(4-Hydroxyphenyl)-3,6,7-trimet...	344	C18H16O7	000569-80-2	37
4			Imidazo[4,5-a]indeno[1,2-b]quino...	344	C20H16N4O2	1000272-51-2	35
5			1H-Indazol-4-amine, N-(4-methylp...	344	C20H16N4O2	1000276-90-7	35



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA M\DATA\BM120717\
Data File : BM013254.D
Acq On : 07 Dec 2017 21:39
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ALS Vial : 16 Sample Multiplier: 1

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ClientSampleId :
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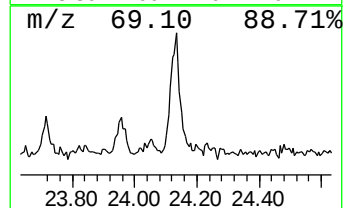
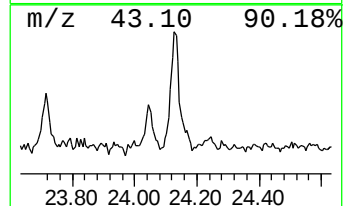
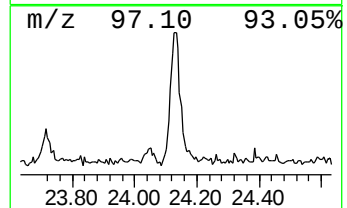
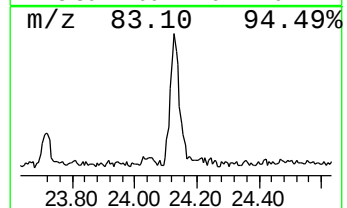
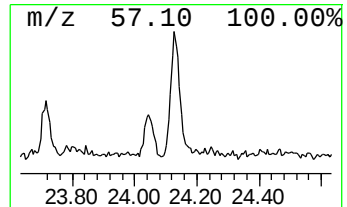
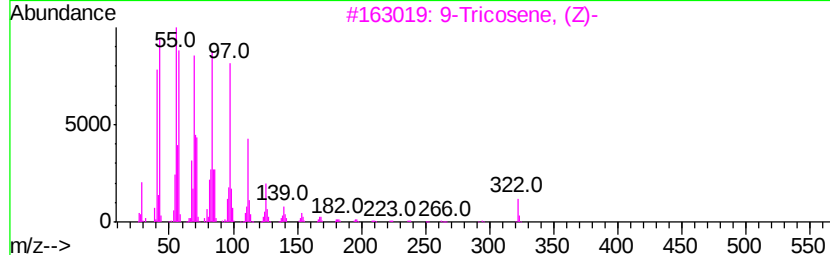
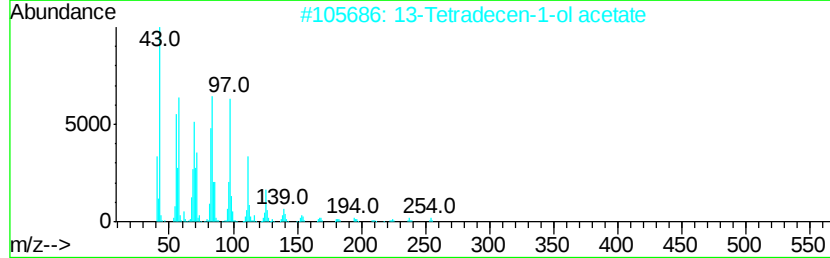
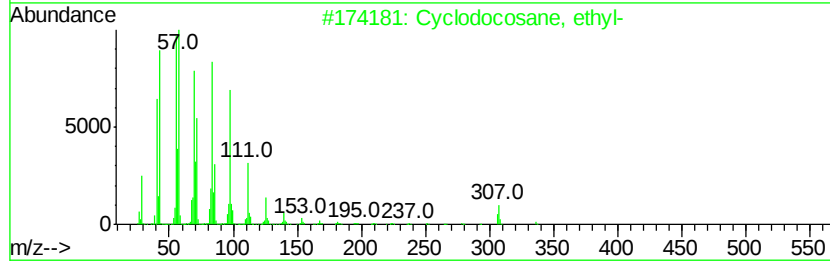
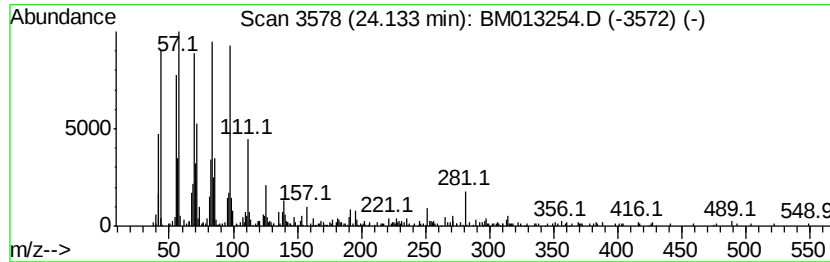
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 (DEL) Alkane: Cyclic24.13 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.13	6.87 ng/ul	799879	Perylene-d12	23.51

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclodocosane, ethyl-	336	C24H48	1000151-22-6	93
2		13-Tetradecen-1-ol acetate	254	C16H30O2	056221-91-1	91
3		9-Tricosene, (Z)-	322	C23H46	027519-02-4	90
4		Hexadecane-1,2-diol	258	C16H34O2	006920-24-7	90
5		Cyclotetracosane	336	C24H48	000297-03-0	87



Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA M\DATA\BM120717\
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Misc :
ALS Vial : 16 Sample Multiplier: 1

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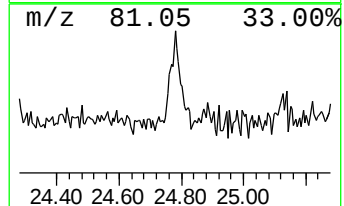
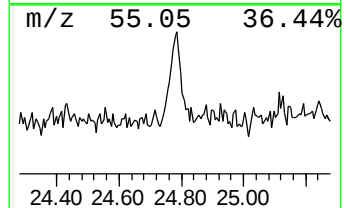
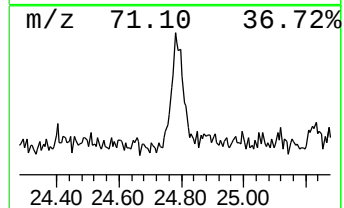
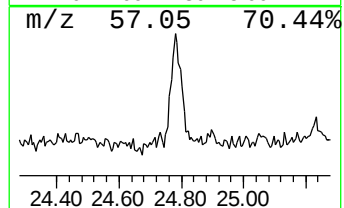
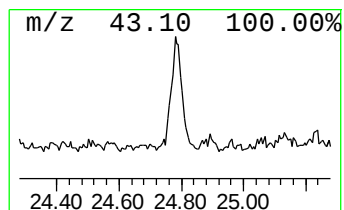
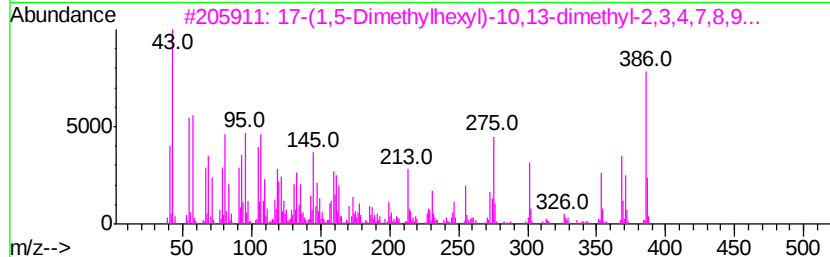
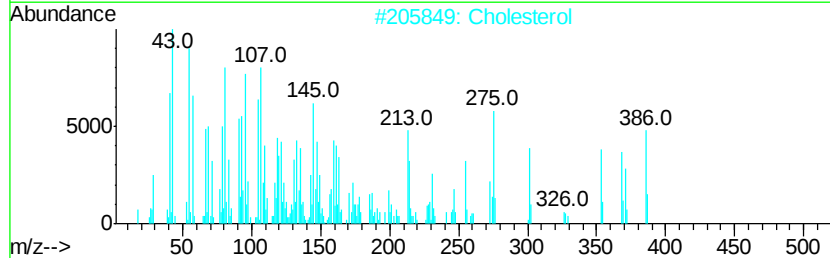
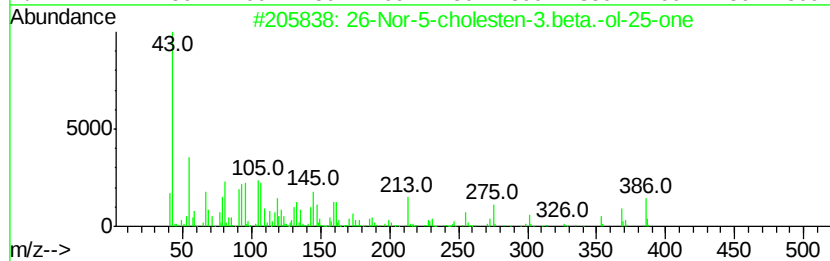
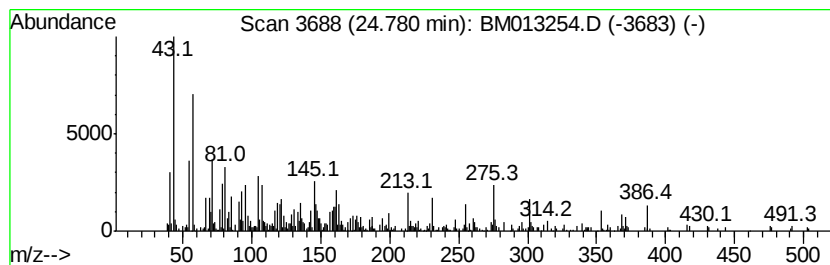
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 26-Nor-5-cholesten-3.beta.-... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.78	6.51 ng/ul	757330	Perylene-d12	23.51

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			26-Nor-5-cholesten-3.beta.-ol-25...	386	C26H42O2	007494-34-0	89
2			Cholesterol	386	C27H46O	000057-88-5	78
3			17-(1,5-Dimethylhexyl)-10,13-dim...	386	C27H46O	1000210-38-4	70
4			26-Nor-5-cholesten-3.beta.-ol-25...	386	C26H42O2	007494-34-0	64
5			Cholesterol	386	C27H46O	000057-88-5	50



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Acq On : 07 Dec 2017 21:39
Operator : SJ/JU
Sample : I6719-01
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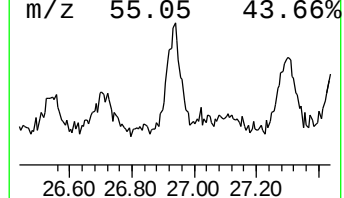
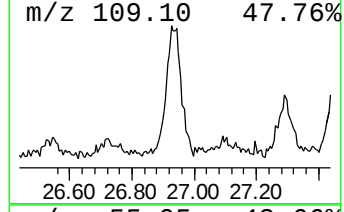
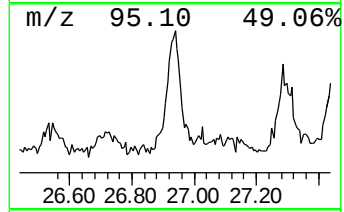
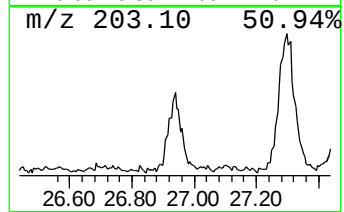
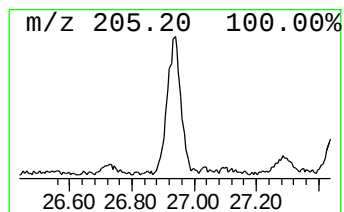
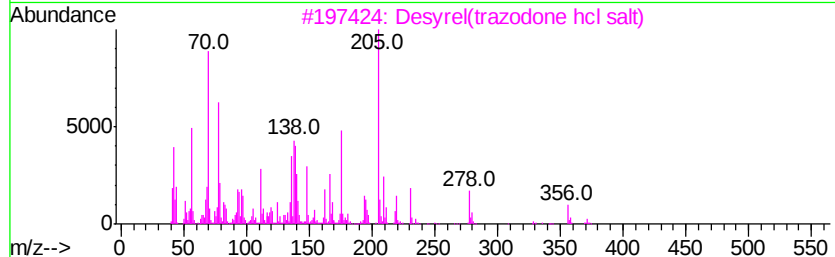
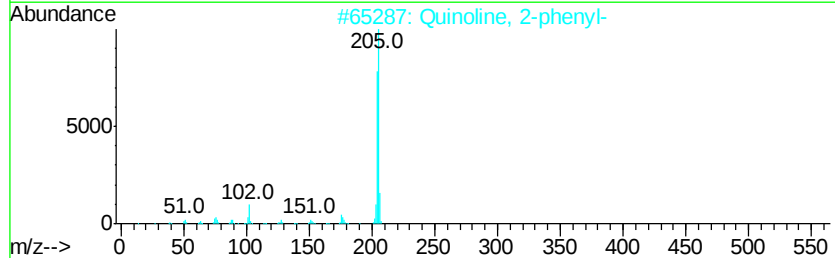
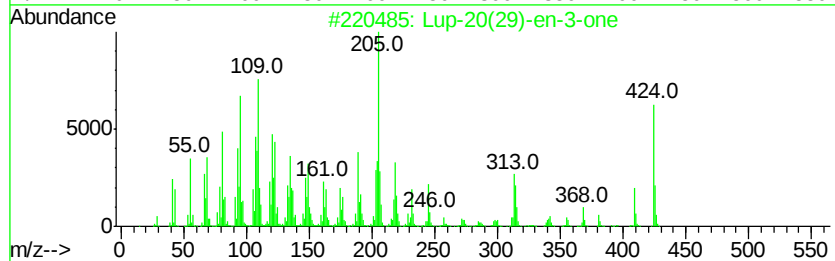
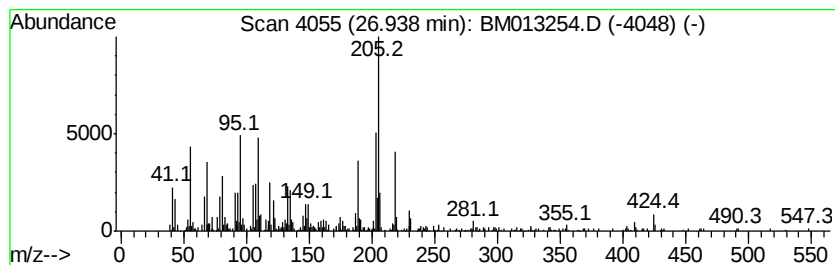
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM113017.M
Quant Title : SVOA CALIBRATION

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

Peak Number 12 Lup-20(29)-en-3-one Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.94	10.22 ng/ul	1188580	Perylene-d12	23.51
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Lup-20(29)-en-3-one	424 C30H48O	001617-70-5 64
2		Quinoline, 2-phenyl-	205 C15H11N	000612-96-4 30
3		Desvrel(trazodone hcl salt)	371 C19H22ClN5O	019794-93-5 30
4		1,3-Propanedione, 1,3-bis(tricyc...	340 C23H32O2	082094-53-9 27
5		6-Phenylisoquinoline	205 C15H11N	070125-61-0 25



Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA M\DATA\BM120717\
Data File : BM013254.D
Acq On : 07 Dec 2017 21:39
Operator : SJ/JU
Sample : I6719-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
JHPH9

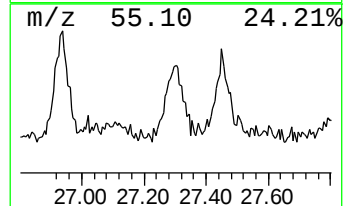
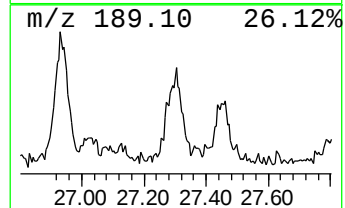
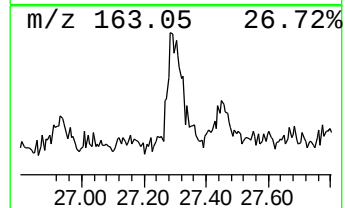
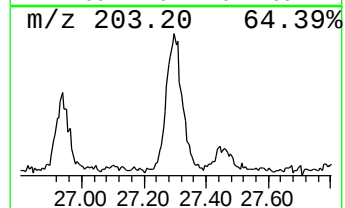
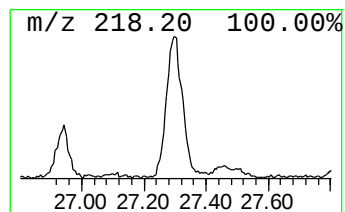
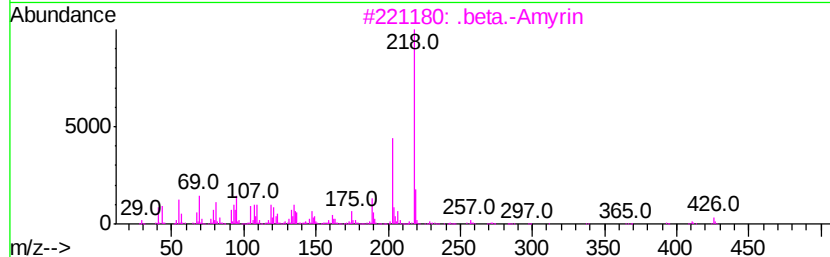
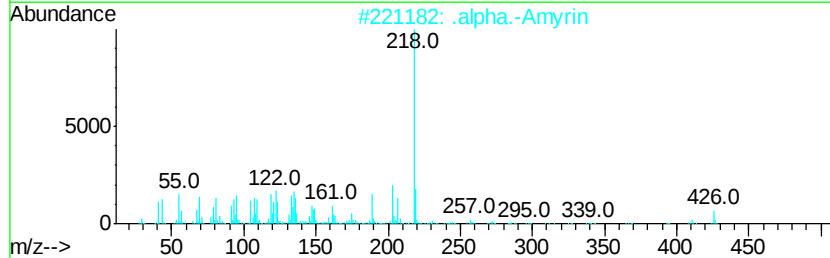
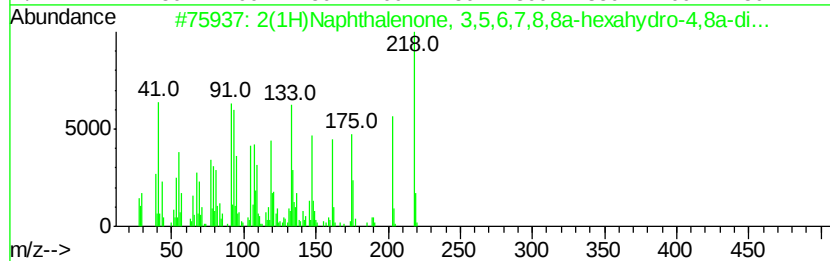
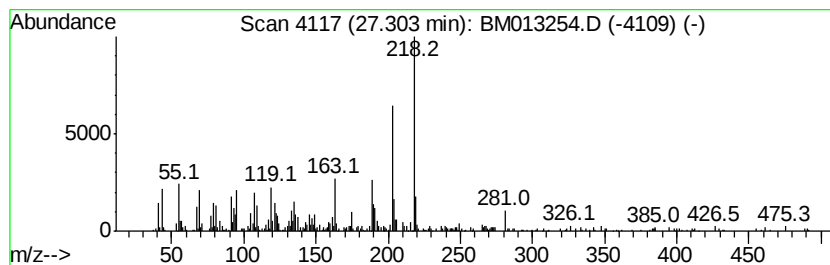
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM113017.M
Quant Title : SVOA CALIBRATION

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

Peak Number 13 2(1H)Naphthalenone, 3,5,6,7... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.30	9.99 ng/ul	1162810	Perylene-d12	23.51

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2(1H)Naphthalenone, 3,5,6,7,8,8a...	218	C15H22O	1000188-66-5	80
2		.alpha.-Amyrin	426	C30H50O	000638-95-9	47
3		.beta.-Amyrin	426	C30H50O	000559-70-6	43
4		5H-Oxazolo[3,2-a]pyridine-8-carb...	218	C12H14N2O2	1000337-58-3	38
5		Benzo[b]naphtho[2,3-d]furan	218	C16H10O	000243-42-5	38



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_M\DATA\BM120717\
 Data File : BM013254.D
 Acq On : 07 Dec 2017 21:39
 Operator : SJ/JU
 Sample : I6719-01
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 JHPH9

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM113017.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
n-Hexadecanoic acid	17.99	5.8	ng/ul	637582	4	17.09	2204740	20.0
Octadecanoic acid	19.27	4.9	ng/ul	590353	5	21.27	2419820	20.0
Sulfurous acid, o...	21.00	4.4	ng/ul	530927	5	21.27	2419820	20.0
(DEL) Alkane: Str...	21.87	9.3	ng/ul	1130660	5	21.27	2419820	20.0
unknown-01	21.99	6.0	ng/ul	722929	5	21.27	2419820	20.0
Benzenamine, 4,4'...	22.32	3.8	ng/ul	461349	5	21.27	2419820	20.0
(DEL) Alkane: Cyc...	22.88	9.4	ng/ul	1096350	6	23.51	2327020	20.0
Oxirane, hexadecyl-	23.71	2.1	ng/ul	242080	6	23.51	2327020	20.0
4-Pregnen-17.alph...	23.96	2.6	ng/ul	298109	6	23.51	2327020	20.0
(DEL) Alkane: Cyc...	24.13	6.9	ng/ul	799879	6	23.51	2327020	20.0
26-Nor-5-choleste...	24.78	6.5	ng/ul	757330	6	23.51	2327020	20.0
Lup-20(29)-en-3-one	26.94	10.2	ng/ul	1188580	6	23.51	2327020	20.0
2(1H)Naphthalenon...	27.30	10.0	ng/ul	1162810	6	23.51	2327020	20.0