

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM073018\
 Data File : BM016131.D
 Acq On : 30 Jul 2018 23:50
 Operator : SJ/JU
 Sample : J4149-19
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 JJ1H8

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:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM071318MA.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	3.458	26	29	39	rVB2	13540	23388	1.01%	0.105%
2	7.357	688	692	710	rBB	68606	118699	5.14%	0.531%
3	7.522	714	720	729	rBV	190678	306350	13.26%	1.369%
4	7.722	749	754	770	rBV	262561	435315	18.84%	1.946%
5	8.198	829	835	848	rBB	185604	313602	13.57%	1.402%
6	8.916	952	957	973	rBV	191338	343889	14.88%	1.537%
7	9.381	1030	1036	1048	rBV	299146	503746	21.80%	2.251%
8	10.104	1153	1159	1174	rBB	280372	481478	20.84%	2.152%
9	10.639	1244	1250	1267	rBV	360670	662408	28.67%	2.961%
10	11.016	1307	1314	1326	rBV	283762	482572	20.89%	2.157%
11	12.386	1538	1547	1554	rBV6	8506	24459	1.06%	0.109%
12	12.528	1565	1571	1573	rBV3	15164	23717	1.03%	0.106%
13	12.710	1598	1602	1613	rVB6	17268	37106	1.61%	0.166%
14	12.816	1613	1620	1626	rBV5	11295	24771	1.07%	0.111%
15	12.939	1633	1641	1644	rBV5	15837	39441	1.71%	0.176%
16	12.969	1644	1646	1653	rVB6	14684	25408	1.10%	0.114%
17	13.357	1707	1712	1719	rBV4	18354	34757	1.50%	0.155%
18	13.486	1730	1734	1744	rVB7	24626	57185	2.48%	0.256%
19	13.598	1747	1753	1760	rBV5	31056	74049	3.20%	0.331%
20	13.692	1763	1769	1774	rBV5	69232	132216	5.72%	0.591%
21	13.780	1774	1784	1789	rVV5	53397	156804	6.79%	0.701%
22	13.833	1789	1793	1797	rVV2	63451	101663	4.40%	0.454%
23	13.875	1797	1800	1805	rVB5	24024	40154	1.74%	0.179%
24	14.033	1821	1827	1828	rBV4	27717	41984	1.82%	0.188%
25	14.110	1836	1840	1843	rBV	61351	85428	3.70%	0.382%
26	14.145	1843	1846	1851	rVB4	43235	63378	2.74%	0.283%
27	14.233	1855	1861	1868	rVV	902257	1282022	55.49%	5.730%
28	14.304	1868	1873	1883	rVB5	83273	204825	8.87%	0.915%
29	14.386	1883	1887	1889	rBV4	37252	52304	2.26%	0.234%
30	14.422	1890	1893	1899	rVB8	25191	39401	1.71%	0.176%
31	14.522	1904	1910	1916	rBV	968635	1367307	59.18%	6.111%
32	14.604	1921	1924	1929	rVB3	49888	87099	3.77%	0.389%
33	14.686	1934	1938	1943	rVB3	42680	73367	3.18%	0.328%
34	14.739	1944	1947	1951	rBV6	19344	33159	1.44%	0.148%

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35	14.822	1956	1961	1969	rVB2	496923	777367	33.65%	3.474%
36	14.892	1969	1973	1974	rBV3	18333	24177	1.05%	0.108%
37	15.027	1989	1996	1997	rBV4	42833	79576	3.44%	0.356%
38	15.057	1997	2001	2008	rVV3	100321	199697	8.64%	0.893%
39	15.351	2048	2051	2052	rBV	29294	29603	1.28%	0.132%
40	15.574	2086	2089	2095	rVB5	61045	100951	4.37%	0.451%
41	15.663	2101	2104	2109	rBV4	35672	67175	2.91%	0.300%
42	15.810	2123	2129	2135	rBV	1262927	1743480	75.46%	7.792%
43	15.951	2148	2153	2158	rBV	529044	813172	35.20%	3.634%
44	16.163	2187	2189	2195	rVB4	82452	115386	4.99%	0.516%
45	16.257	2202	2205	2211	rBV5	58786	78085	3.38%	0.349%
46	16.804	2294	2298	2304	rVB5	46756	73022	3.16%	0.326%
47	17.168	2357	2360	2366	rVB8	34540	49374	2.14%	0.221%
48	17.245	2370	2373	2378	rBV3	80636	115442	5.00%	0.516%
49	17.304	2380	2383	2389	rVB2	85085	119625	5.18%	0.535%
50	17.557	2420	2426	2432	rBV	719113	1083943	46.91%	4.845%
51	17.657	2437	2443	2449	rVB	1474184	2088958	90.41%	9.336%
52	18.404	2567	2570	2579	rVB	121015	210787	9.12%	0.942%
53	18.886	2648	2652	2659	rVB	62534	93343	4.04%	0.417%
54	19.662	2780	2784	2791	rVB2	88782	139224	6.03%	0.622%
55	19.915	2821	2827	2835	rBV	1787341	2310441	100.00%	10.326%
56	21.674	3121	3126	3135	rBV	924531	1153670	49.93%	5.156%
57	23.886	3494	3502	3510	rVB	1118312	2123740	91.92%	9.492%
58	24.033	3520	3527	3537	rVB	521286	1010913	43.75%	4.518%

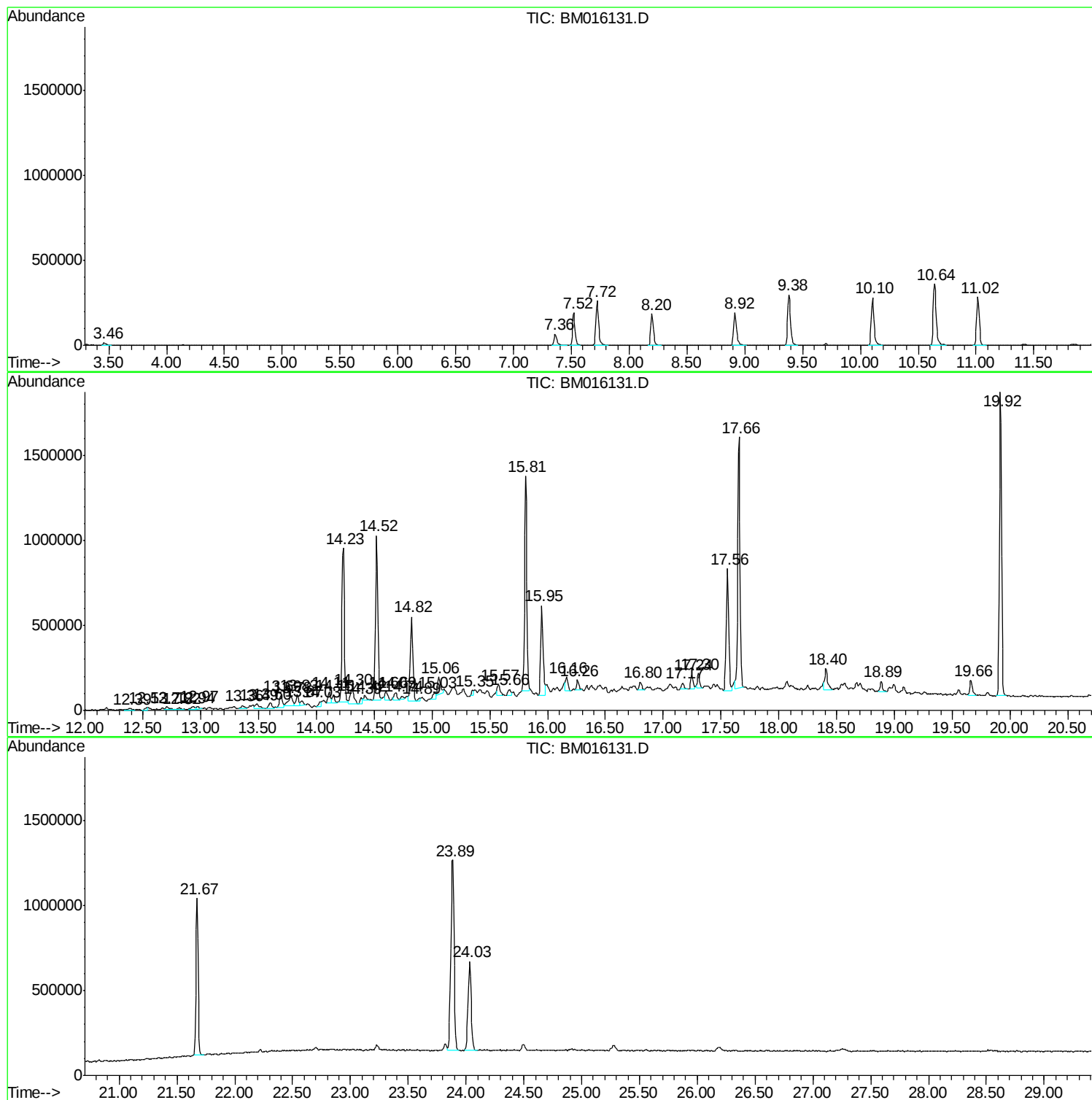
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Quant Title : SVOA CALIBRATION

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



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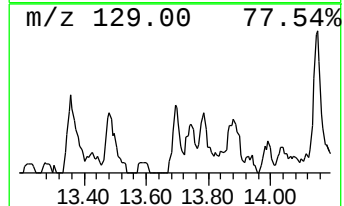
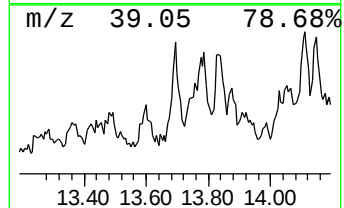
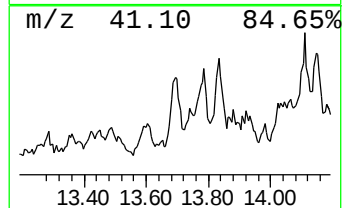
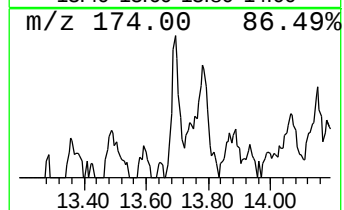
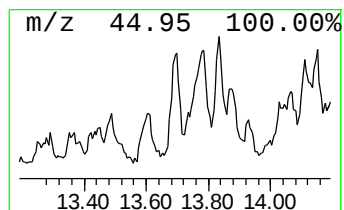
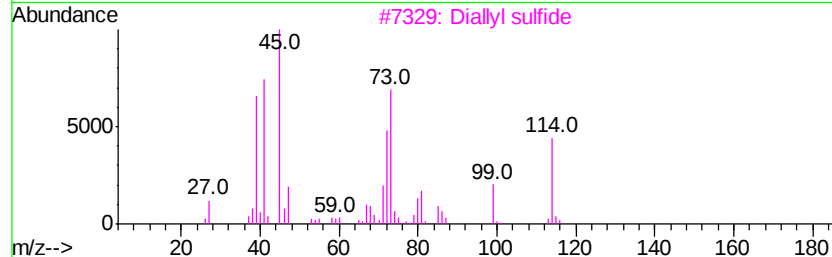
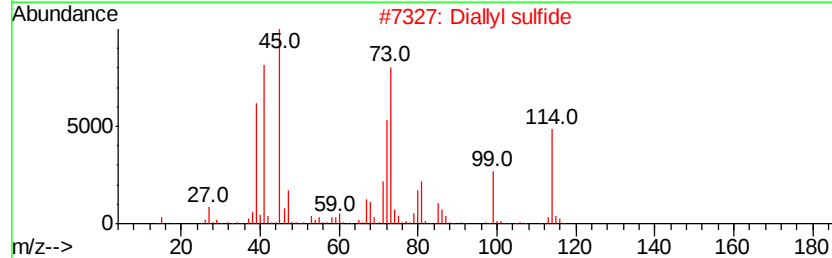
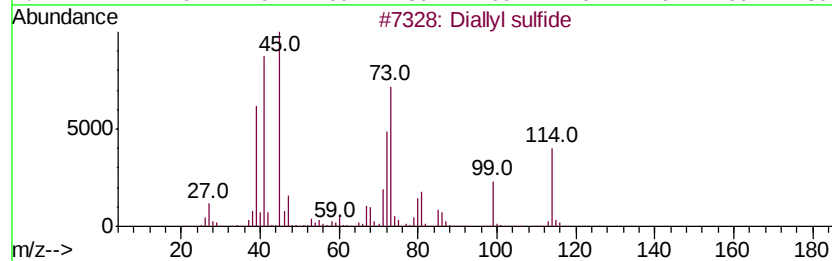
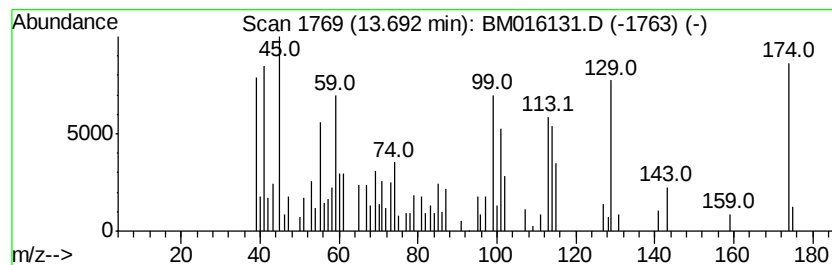
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Quant Title : SVOA CALIBRATION

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

Peak Number 1 unknown-01 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.69	3.40 ng/ul	132216	Acenaphthene-d10	14.82

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Diallyl sulfide	114	C6H10S	000592-88-1	49
2			Diallyl sulfide	114	C6H10S	000592-88-1	46
3			Diallyl sulfide	114	C6H10S	000592-88-1	46
4			Cyclopropanecarboxylic acid, met...	100	C5H8O2	002868-37-3	45
5			5-Hexen-2-ol, 5-methyl-	114	C7H14O	050551-88-7	38



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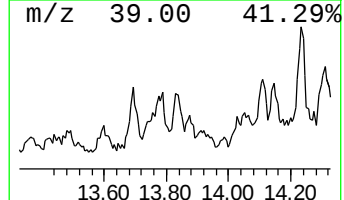
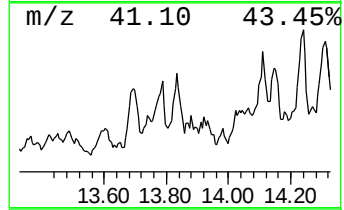
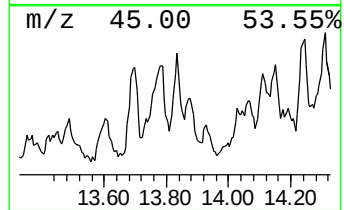
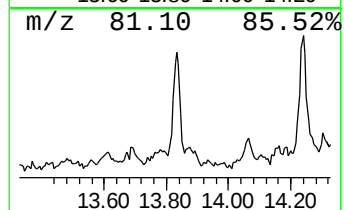
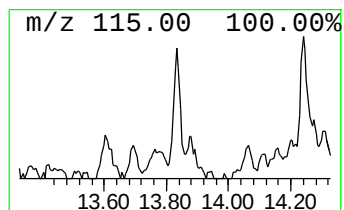
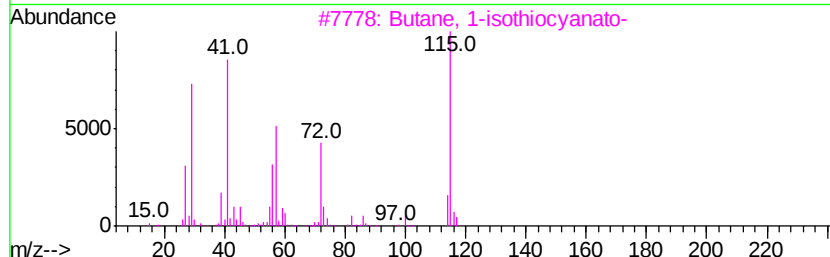
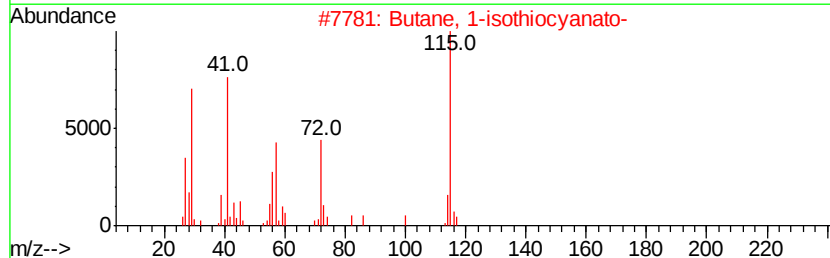
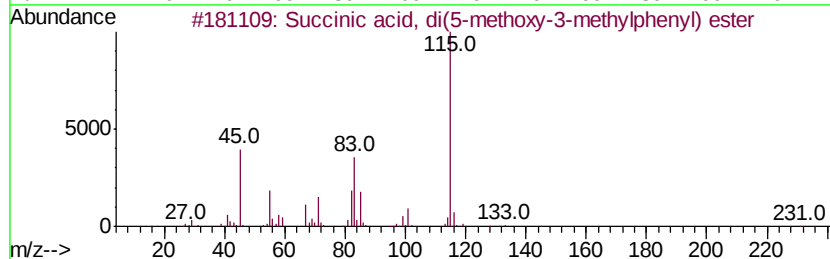
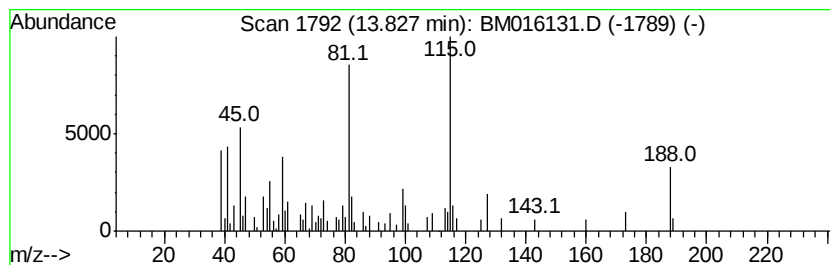
Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM071318MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 unknown-02 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.83	2.62 ng/ul	101663	Acenaphthene-d10	14.82
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Succinic acid, di(5-methoxy-3-me...	346 C18H34O6	1000330-26-5	27
2	Butane, 1-isothiocyanato-	115 C5H9NS	000592-82-5	18
3	Butane, 1-isothiocyanato-	115 C5H9NS	000592-82-5	18
4	Cyclopropane, 1-chloro-2-(1-prop...	116 C6H9Cl	074752-94-6	15
5	Allyloxy-t-butyltrimethylsilane	172 C9H20OSi	085807-85-8	14



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ALS Vial : 12 Sample Multiplier: 1

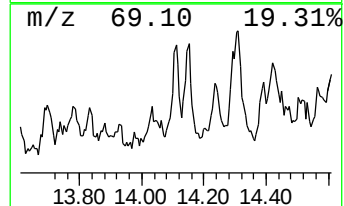
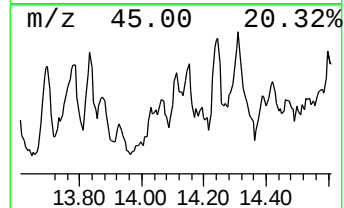
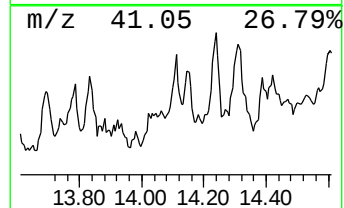
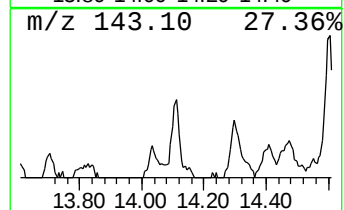
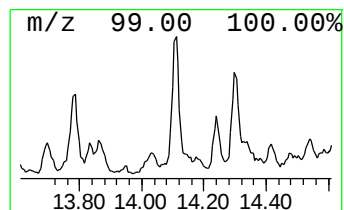
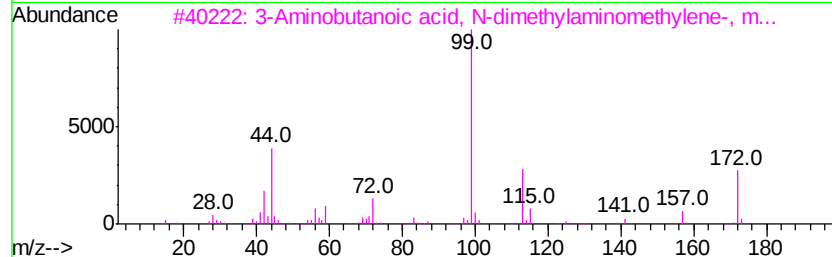
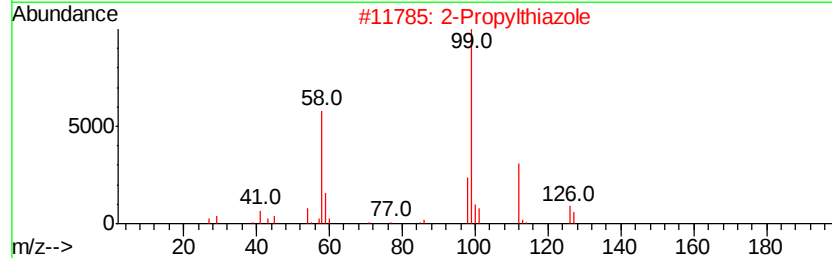
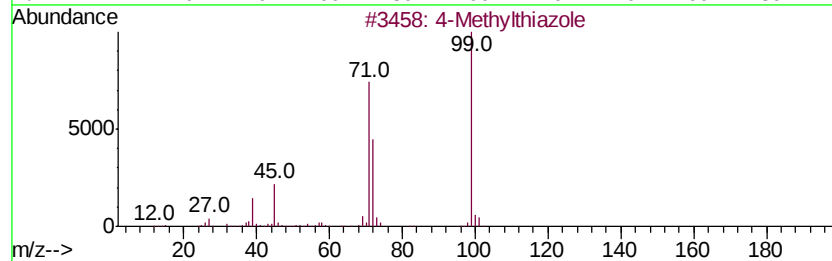
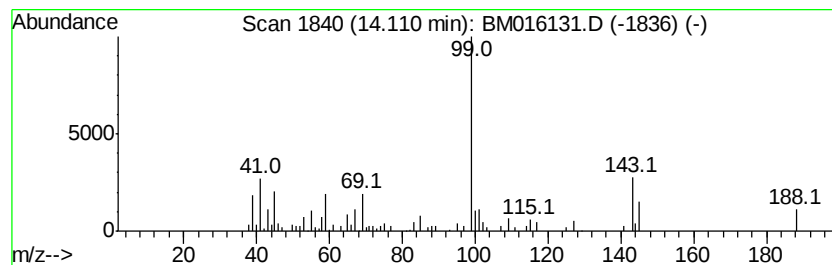
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TIC Library : C:\Database\NIST11.L
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Peak Number 4 4-Methylthiazole Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.11	2.20 ng/ul	85428	Acenaphthene-d10	14.82
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	4-Methylthiazole	99	C4H5NS	000693-95-8 50
2	2-Propylthiazole	127	C6H9NS	017626-75-4 47
3	3-Aminobutanoic acid, N-dimethyl...	172	C8H16N2O2	1000375-77-6 43
4	Thiazole, 5-methyl-	99	C4H5NS	003581-89-3 40
5	1,4-Dioxaspiro[4.5]decane, 8-(me...	188	C9H16O2S	055103-51-0 38



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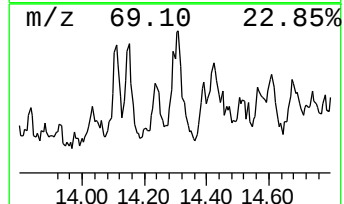
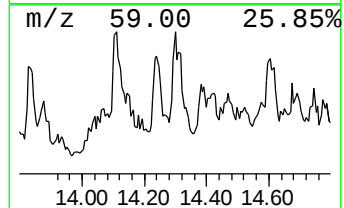
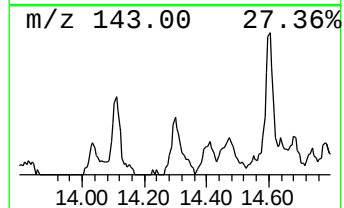
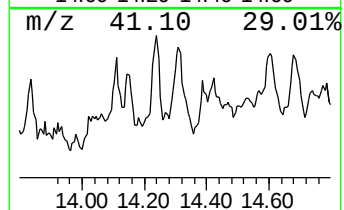
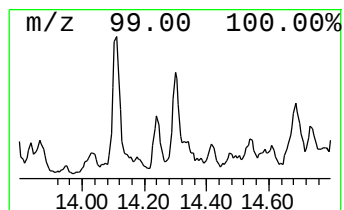
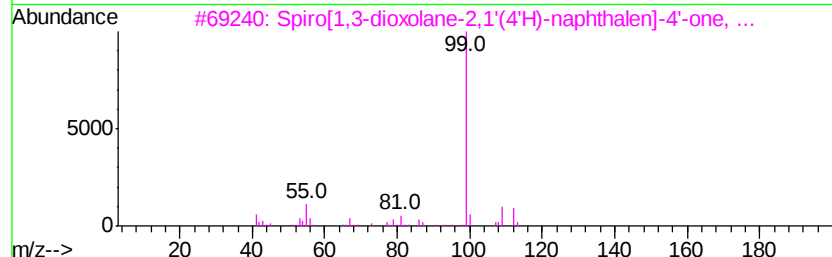
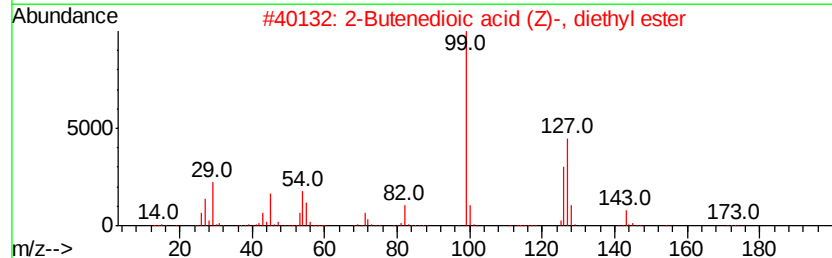
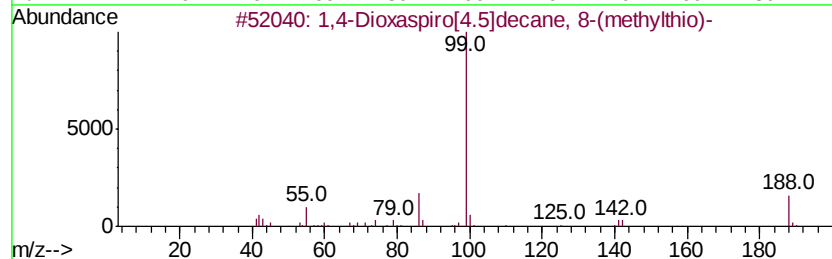
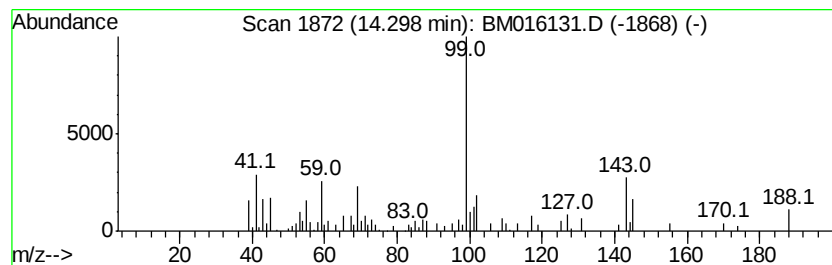
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST11.L
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Peak Number 5 unknown-03 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.30	5.27 ng/ul	204825	Acenaphthene-d10	14.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,4-Dioxaspiro[4.5]decane, 8-(me...	188	C9H16O2S	055103-51-0	38
2		2-Butenedioic acid (Z)-, diethyl...	172	C8H12O4	000141-05-9	38
3		Spiro[1.3-dioxolane-2.1'(4'H)-na...	210	C12H18O3	021727-93-5	38
4		Phosphoric acid, diethyl pentyl ...	224	C9H21O4P	020195-08-8	37
5		4,8,12,16-Tetramethylheptadecan...	324	C21H40O2	096168-15-9	37



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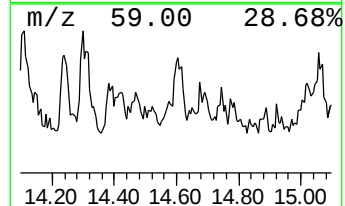
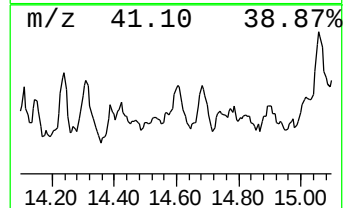
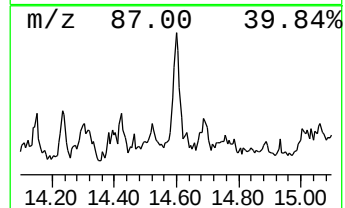
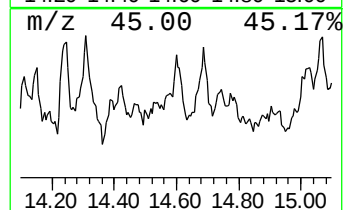
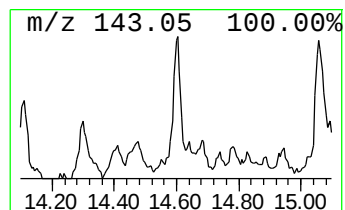
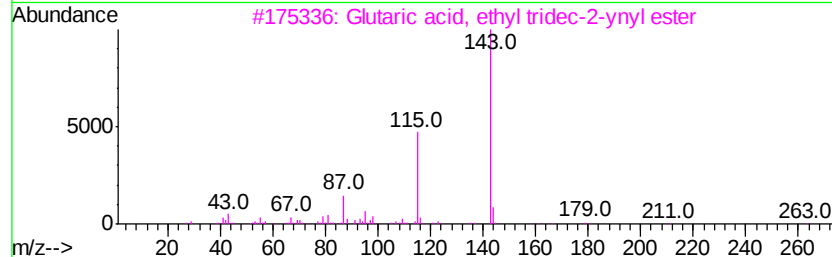
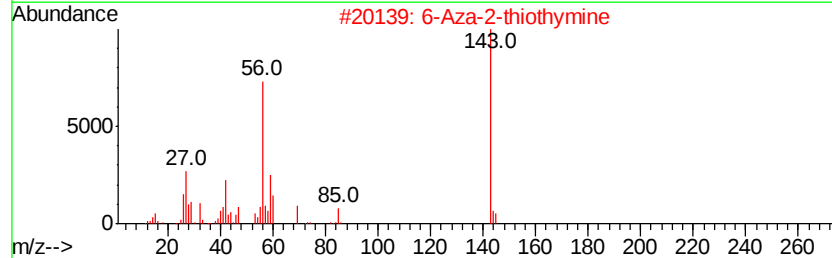
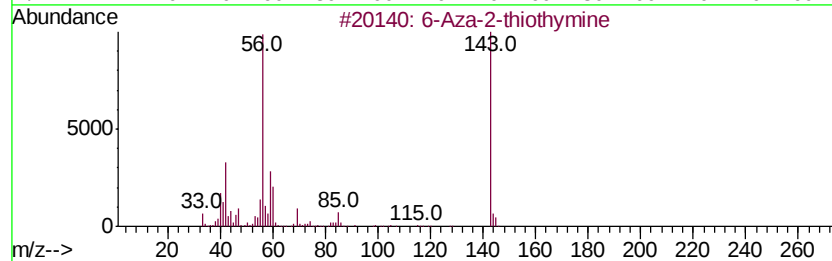
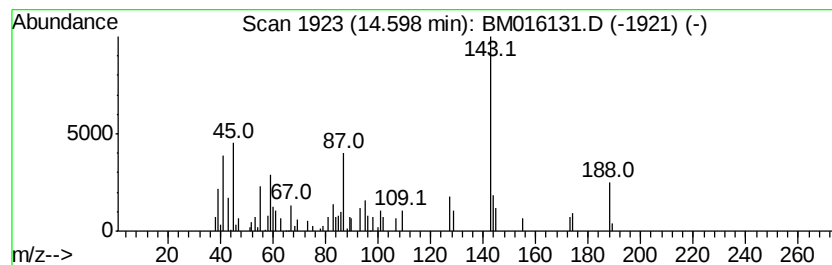
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 unknown-04 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.60	2.24 ng/ul	87099	Acenaphthene-d10	14.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	6-Aza-2-thiothymine	143	C4H5N3OS	000615-76-9	38
2		6-Aza-2-thiothymine	143	C4H5N3OS	000615-76-9	38
3		Glutaric acid, ethyl tridec-2-yn...	338	C20H34O4	1000360-11-9	32
4		Malonic acid, butyl 3-hexyl ester	244	C13H24O4	1000349-12-0	32
5		4-Chloro-6-(methylamino)pyrimidine	143	C5H6ClN3	065766-32-7	32



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM073018\
Data File : BM016131.D
Acq On : 30 Jul 2018 23:50
Operator : SJ/JU
Sample : J4149-19
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
JJ1H8

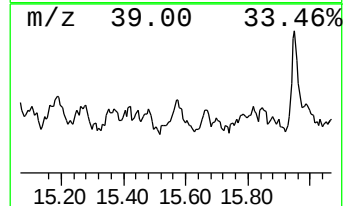
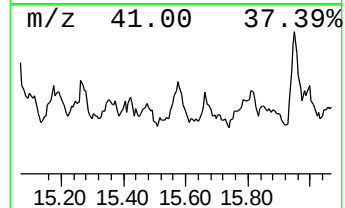
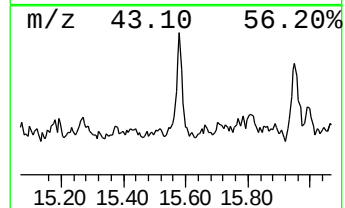
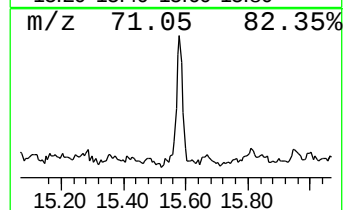
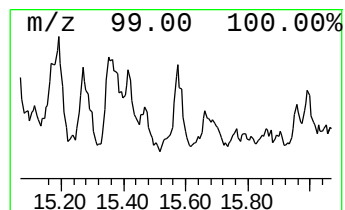
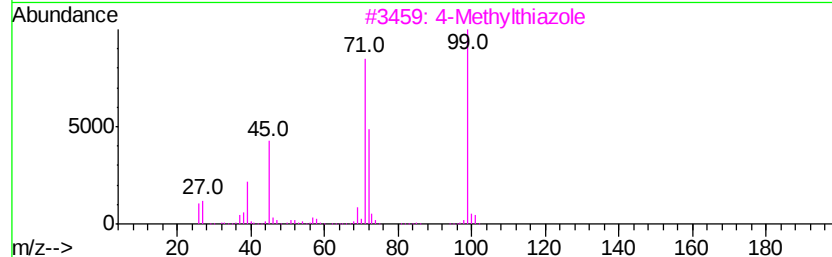
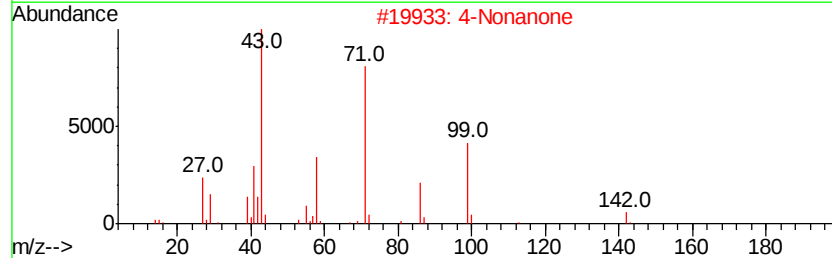
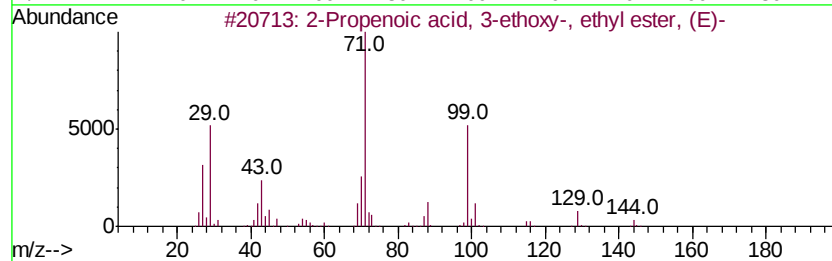
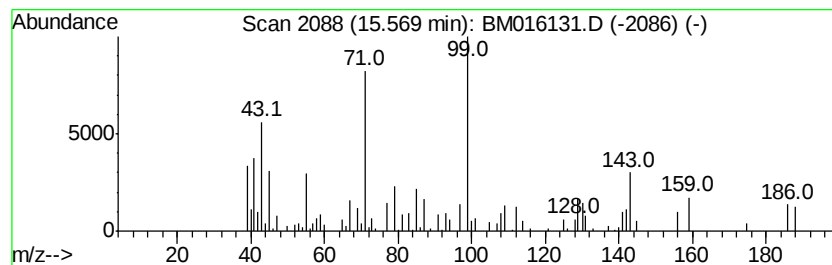
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 2-Propenoic acid, 3-ethoxy-... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.57	2.60 ng/ul	100951	Acenaphthene-d10	14.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Propenoic acid, 3-ethoxy-, eth...	144	C7H12O3	005941-55-9	52
2		4-Nonanone	142	C9H18O	004485-09-0	47
3		4-Methylthiazole	99	C4H5NS	000693-95-8	47
4		Silane, ethenyldiethylmethyl-	128	C7H16Si	018292-29-0	38
5		Pantolactone	130	C6H10O3	000599-04-2	35



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM073018\
Data File : BM016131.D
Acq On : 30 Jul 2018 23:50
Operator : SJ/JU
Sample : J4149-19
Misc :
ALS Vial : 12 Sample Multiplier: 1

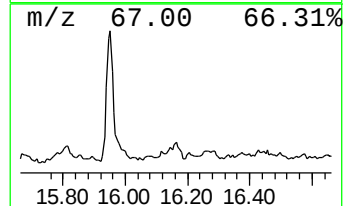
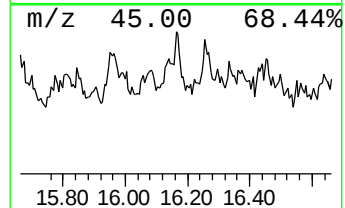
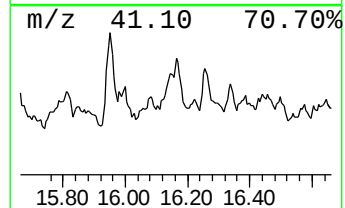
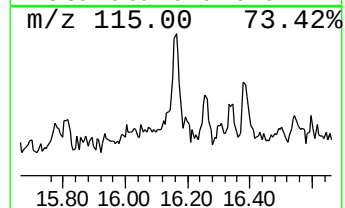
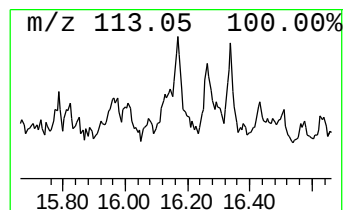
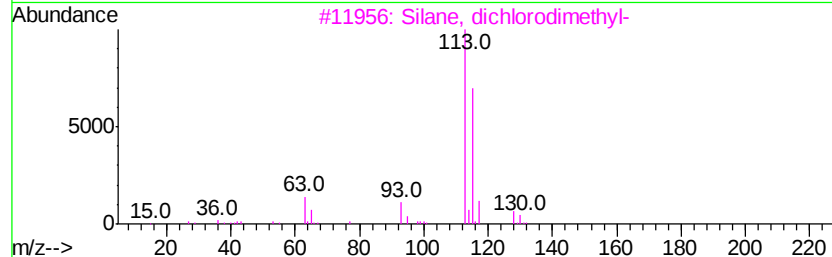
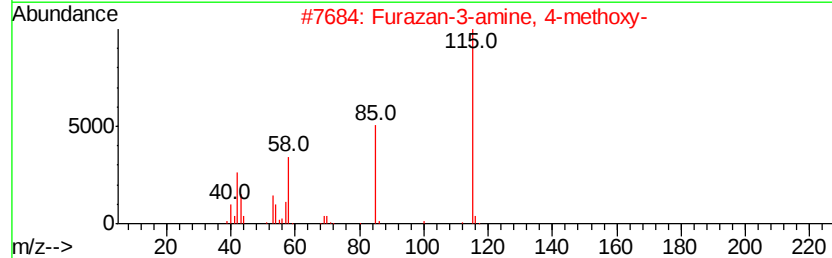
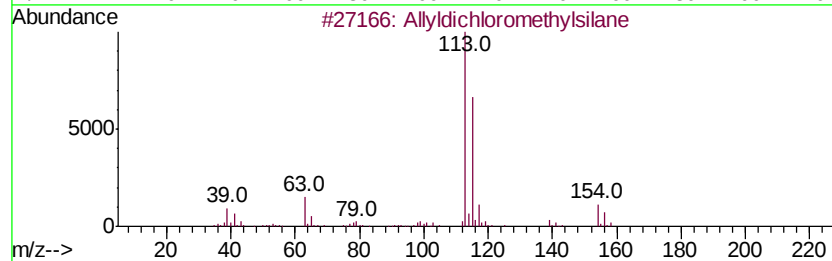
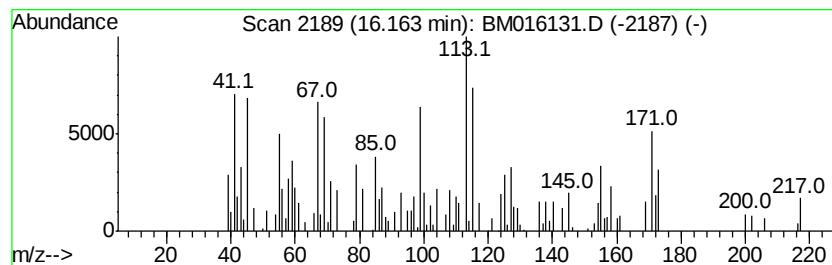
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 unknown-05 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.16	2.97 ng/ul	115386	Acenaphthene-d10	14.82
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Allvldichloromethylsilane	154 C4H8Cl2Si	001873-92-3 38
2		Furazan-3-amine, 4-methoxy-	115 C3H5N3O2	078350-48-8 18
3		Silane, dichlorodimethyl-	128 C2H6Cl2Si	000075-78-5 16
4		Silane, dichlorodimethyl-	128 C2H6Cl2Si	000075-78-5 16
5		Glutaric acid, 3-methylbut-2-eny...	242 C13H22O4	1000360-08-5 14



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM073018\
Data File : BM016131.D
Acq On : 30 Jul 2018 23:50
Operator : SJ/JU
Sample : J4149-19
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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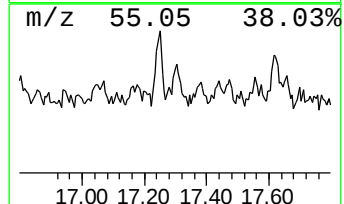
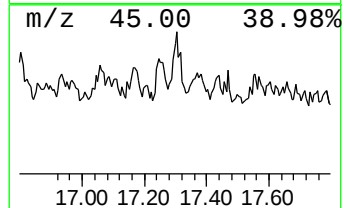
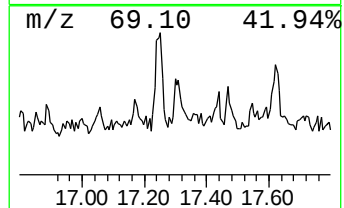
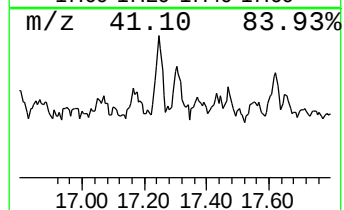
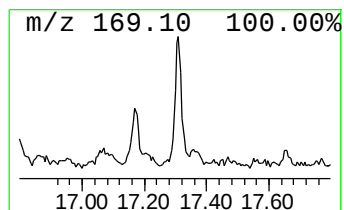
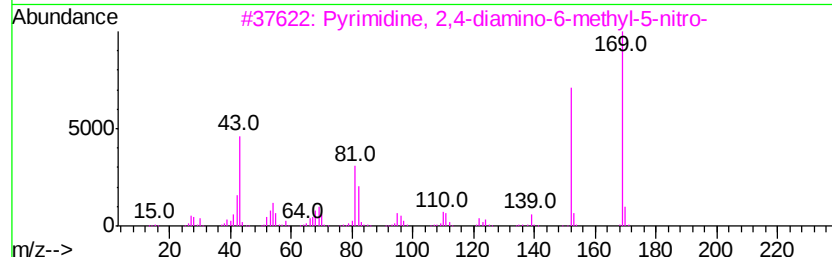
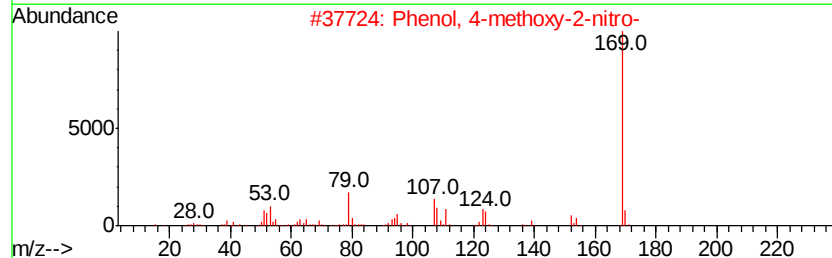
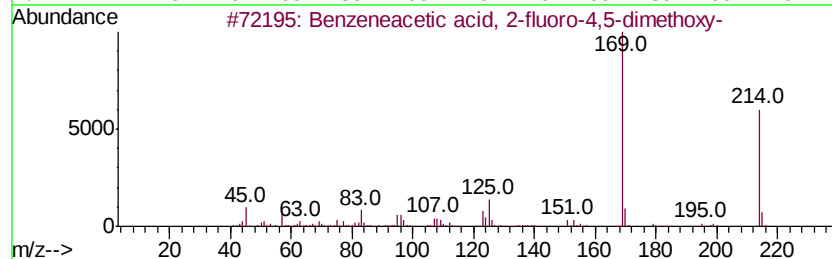
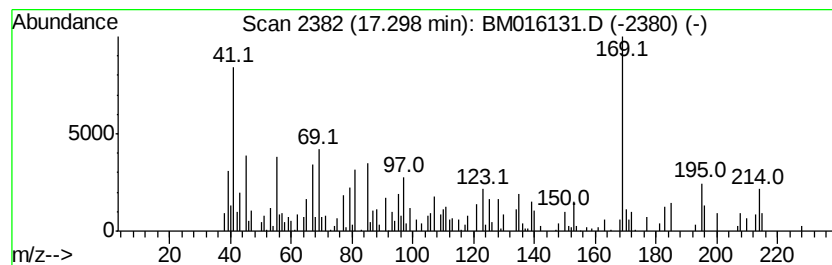
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Benzeneacetic acid, 2-fluor... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.30	2.21 ng/ul	119625	Phenanthrene-d10	17.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzeneacetic acid, 2-fluoro-4,5...	214	C10H11F04	091407-50-0	50
2		Phenol, 4-methoxy-2-nitro-	169	C7H7N04	001568-70-3	49
3		Pvrimidine, 2,4-diamino-6-methyl...	169	C5H7N5O2	002829-59-6	49
4		2-Chlorophenyl isothiocyanate	169	C7H4ClNS	002740-81-0	47
5		2-chloro-4,5-methylenedioxypheny...	212	C10H9ClO3	1000378-96-4	47



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM073018\
Data File : BM016131.D
Acq On : 30 Jul 2018 23:50
Operator : SJ/JU
Sample : J4149-19
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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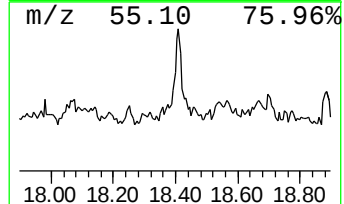
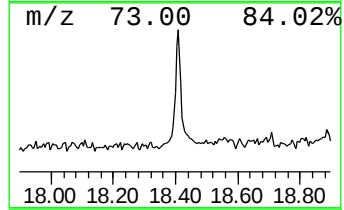
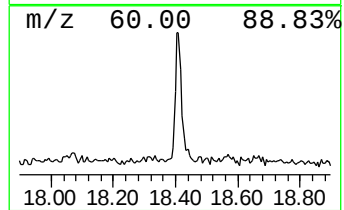
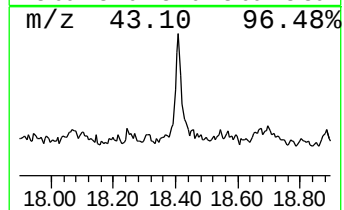
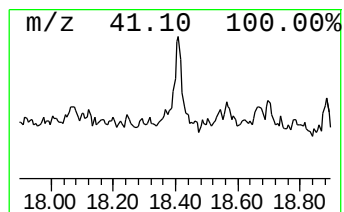
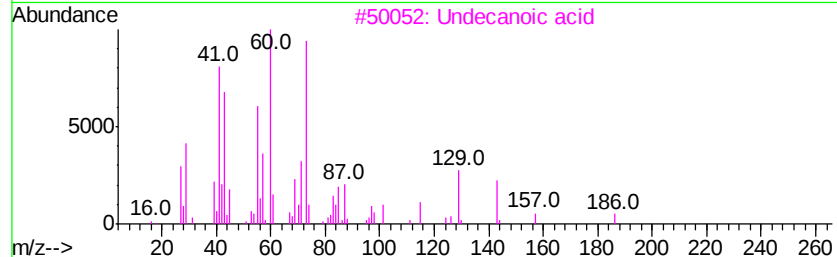
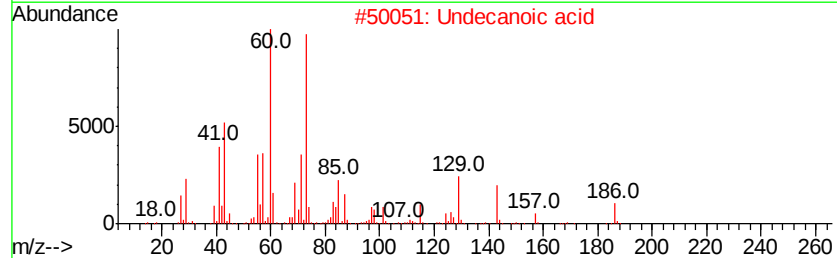
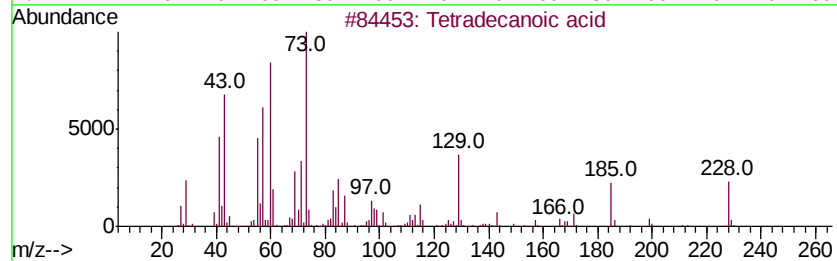
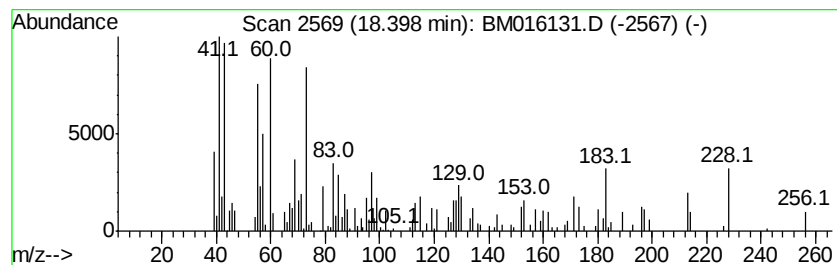
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 Tetradecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.40	3.89 ng/ul	210787	Phenanthrene-d10	17.56

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecanoic acid	228	C14H28O2	000544-63-8	70
2			Undecanoic acid	186	C11H22O2	000112-37-8	58
3			Undecanoic acid	186	C11H22O2	000112-37-8	58
4			n-Decanoic acid	172	C10H20O2	000334-48-5	55
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	53



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM073018\
Data File : BM016131.D
Acq On : 30 Jul 2018 23:50
Operator : SJ/JU
Sample : J4149-19
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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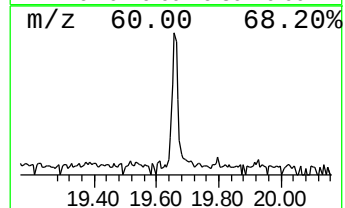
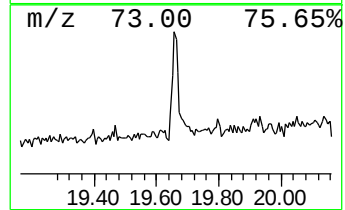
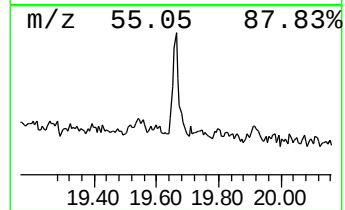
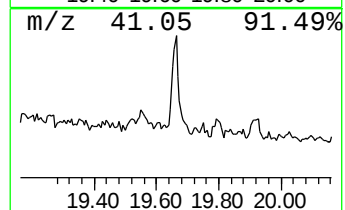
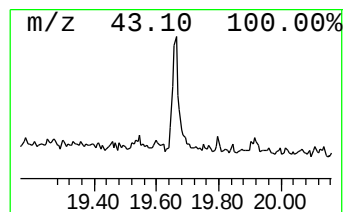
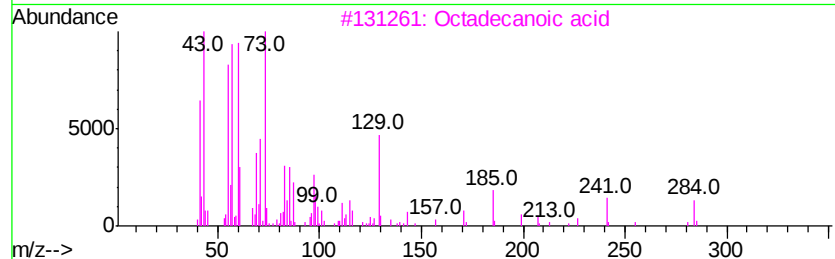
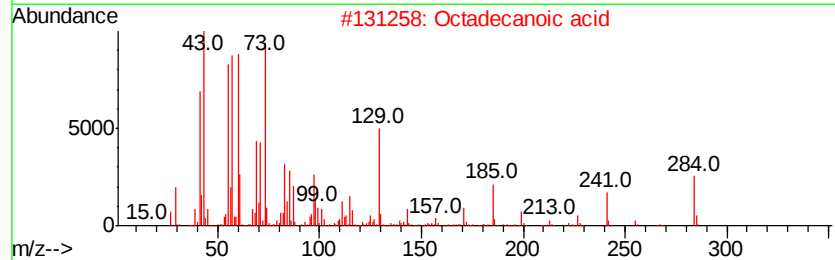
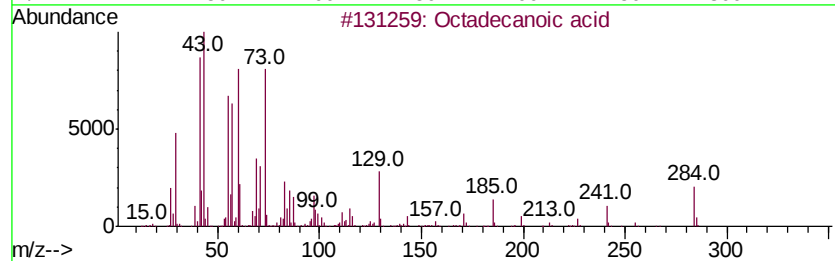
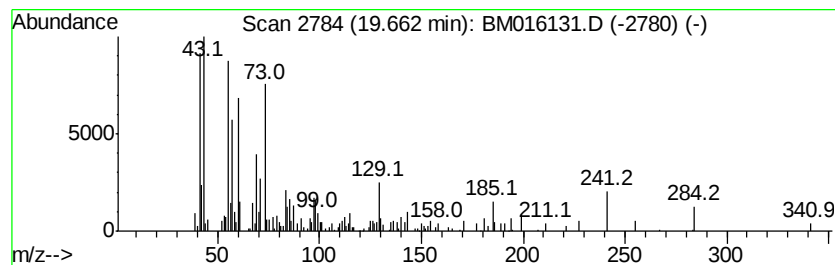
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 Octadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.66	2.41 ng/ul	139224	Chrysene-d12	21.67

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	93
2			Octadecanoic acid	284	C18H36O2	000057-11-4	90
3			Octadecanoic acid	284	C18H36O2	000057-11-4	83
4			Octadecanoic acid	284	C18H36O2	000057-11-4	83
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	62



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM073018\
Data File : BM016131.D
Acq On : 30 Jul 2018 23:50
Operator : SJ/JU
Sample : J4149-19
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM071318MA.M
Quant Title : SVOA CALIBRATION

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

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					#	RT	Resp	Conc
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unknown-02	13.83	2.6	ng/ul	101663	3	14.82	777367	20.0
4-Methylthiazole	14.11	2.2	ng/ul	85428	3	14.82	777367	20.0
unknown-03	14.30	5.3	ng/ul	204825	3	14.82	777367	20.0
unknown-04	14.60	2.2	ng/ul	87099	3	14.82	777367	20.0
2-Propenoic acid,...	15.57	2.6	ng/ul	100951	3	14.82	777367	20.0
unknown-05	16.16	3.0	ng/ul	115386	3	14.82	777367	20.0
Cyclohexane, 1-me...	17.24	2.1	ng/ul	115442	4	17.56	1083940	20.0
Benzeneacetic aci...	17.30	2.2	ng/ul	119625	4	17.56	1083940	20.0
Tetradecanoic acid	18.40	3.9	ng/ul	210787	4	17.56	1083940	20.0
Octadecanoic acid	19.66	2.4	ng/ul	139224	5	21.67	1153670	20.0