

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM101817\
 Data File : BM012306.D
 Acq On : 19 Oct 2017 07:25
 Operator : SJ/JU
 Sample : I5747-06 2X
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 B-4(1-3)-100617

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-BM101217.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	7.181	688	696	703	rVB	11945393	18148217	8.44%	4.012%
2	7.381	725	730	733	rVV	1916603	2827810	1.32%	0.625%
3	7.416	733	736	743	rVB2	1409314	2175215	1.01%	0.481%
4	7.734	784	790	794	rVB	1554970	2325073	1.08%	0.514%
5	7.904	811	819	823	rBV3	1754569	3710838	1.73%	0.820%
6	8.410	899	905	917	rBV2	1007148	2250144	1.05%	0.497%
7	8.986	993	1003	1007	rBV2	2296759	4389397	2.04%	0.970%
8	10.533	1260	1266	1270	rBV	1976379	3157934	1.47%	0.698%
9	10.580	1270	1274	1280	rVB	1367130	2208727	1.03%	0.488%
10	10.716	1292	1297	1303	rVB	1441493	2201134	1.02%	0.487%
11	11.980	1507	1512	1519	rBV	1666966	2579596	1.20%	0.570%
12	14.439	1924	1930	1935	rBV2	1776603	2656115	1.24%	0.587%
13	16.586	2290	2295	2301	rBV	1493257	2247563	1.05%	0.497%
14	17.186	2392	2397	2401	rBV2	2149729	3105038	1.44%	0.686%
15	17.233	2401	2405	2410	rVV	1615399	2277542	1.06%	0.503%
16	18.109	2542	2554	2565	rBV	29859102	61983991	28.83%	13.702%
17	18.980	2697	2702	2706	rBV3	1364407	2364872	1.10%	0.523%
18	19.298	2737	2756	2767	rBV4	52652350	215021142	100.00%	47.533%
19	19.392	2767	2772	2780	rVB2	22103381	31156463	14.49%	6.888%
20	19.492	2785	2789	2791	rVV3	1668432	2648303	1.23%	0.585%
21	19.521	2791	2794	2799	rVV	1493753	2292835	1.07%	0.507%
22	19.586	2800	2805	2807	rVV2	5706323	7792398	3.62%	1.723%
23	19.615	2807	2810	2815	rVB	6746156	8814567	4.10%	1.949%
24	20.109	2890	2894	2899	rVB	2136287	3071000	1.43%	0.679%
25	21.056	3052	3055	3060	rVB	3596138	4124011	1.92%	0.912%
26	21.362	3102	3107	3113	rBV2	8072207	16199538	7.53%	3.581%
27	21.415	3113	3116	3125	rVB	5569652	7208077	3.35%	1.593%
28	22.991	3378	3384	3388	rBV	5985641	13185574	6.13%	2.915%
29	23.162	3409	3413	3419	rVB	1750654	2760875	1.28%	0.610%
30	23.480	3462	3467	3472	rBV	2854045	4942514	2.30%	1.093%
31	23.580	3479	3484	3493	rVB	5152650	10124655	4.71%	2.238%
32	23.680	3497	3501	3505	rVV	1471900	2409532	1.12%	0.533%

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Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
B-4(1-3)-100617

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 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-BM101217.M
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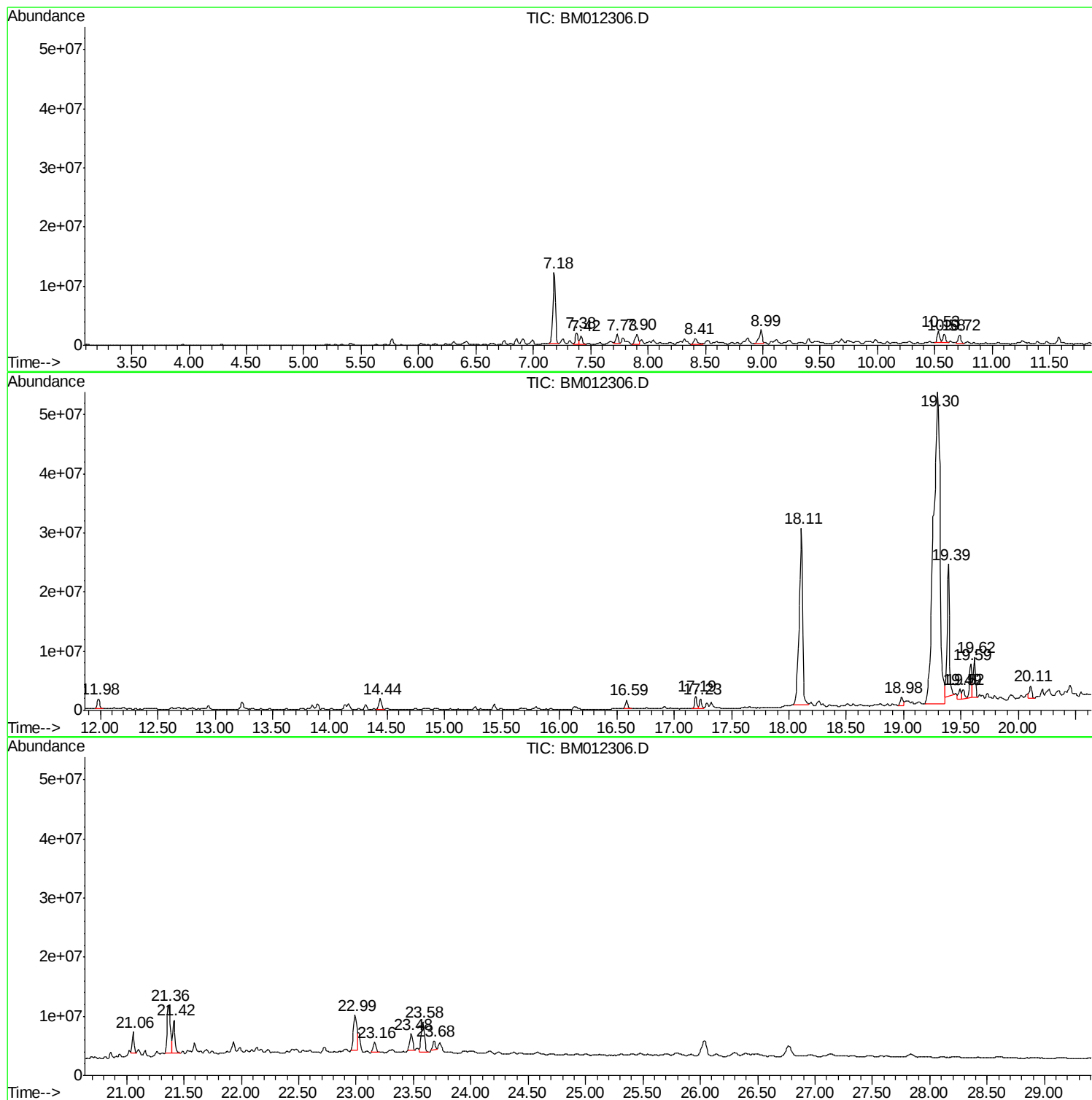
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM101217.M
Quant Title : SVOA CALIBRATION

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



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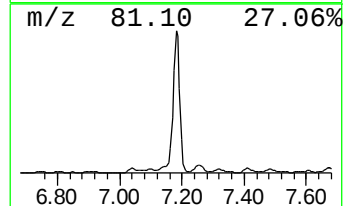
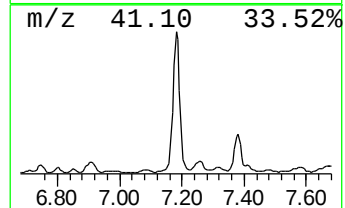
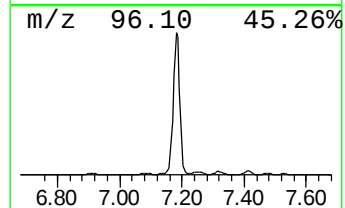
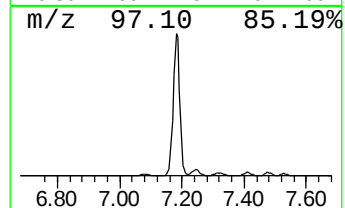
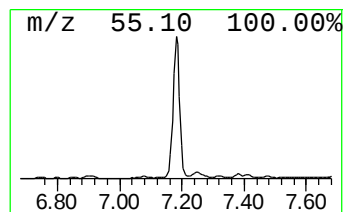
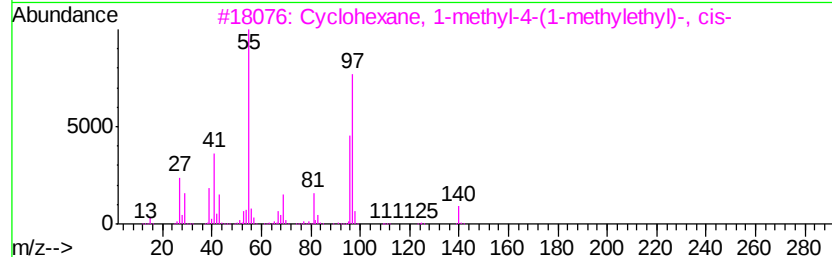
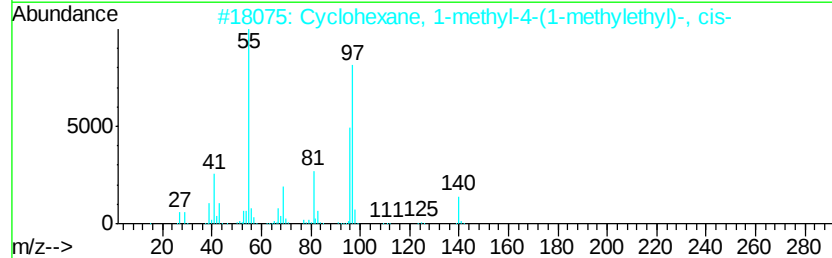
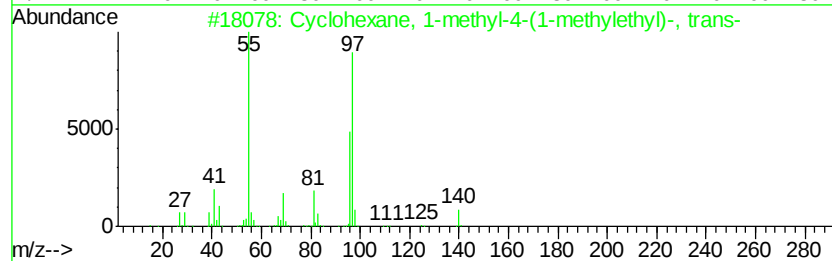
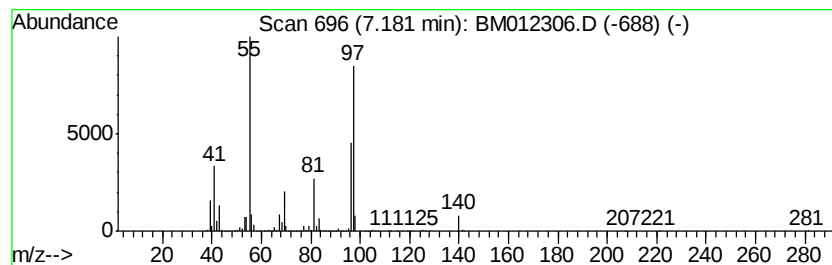
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM101217.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Cyclohexane, 1-methyl-4-(1-... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.18	156.11 ng/ul	18148200	1,4-Dichlorobenzene-d4	7.78

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1-methyl-4-(1-methy...	140	C10H20	001678-82-6	94
2			Cyclohexane, 1-methyl-4-(1-methy...	140	C10H20	006069-98-3	94
3			Cyclohexane, 1-methyl-4-(1-methy...	140	C10H20	006069-98-3	91
4			m-Menthane, (1S,3S)-(+)-	140	C10H20	013837-67-7	91
5			Cyclohexane, 1-methyl-3-(1-methy...	140	C10H20	016580-24-8	87



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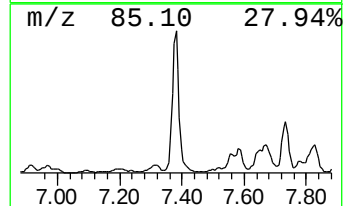
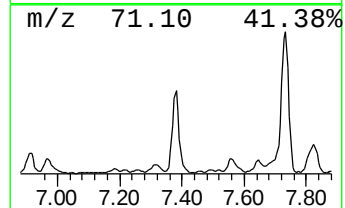
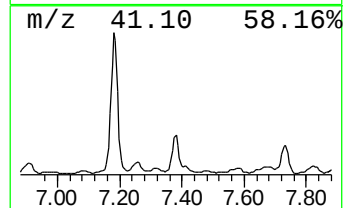
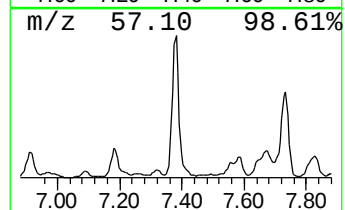
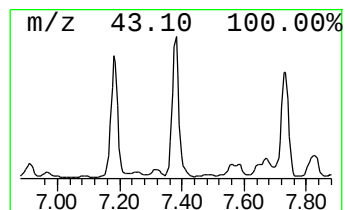
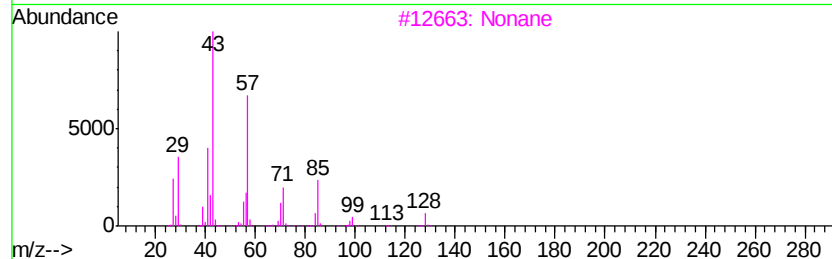
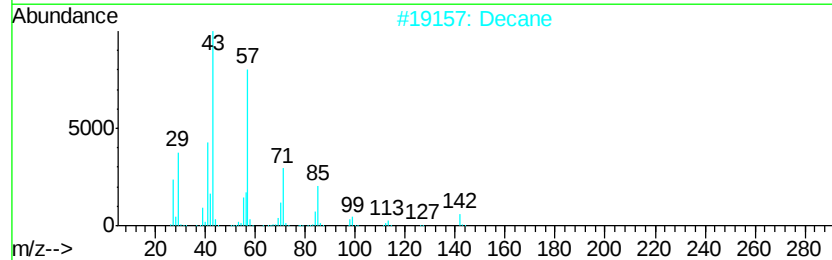
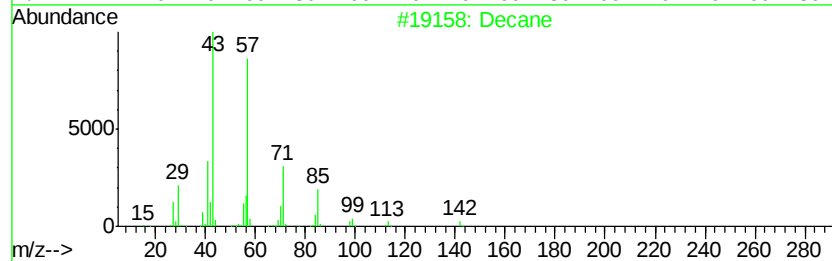
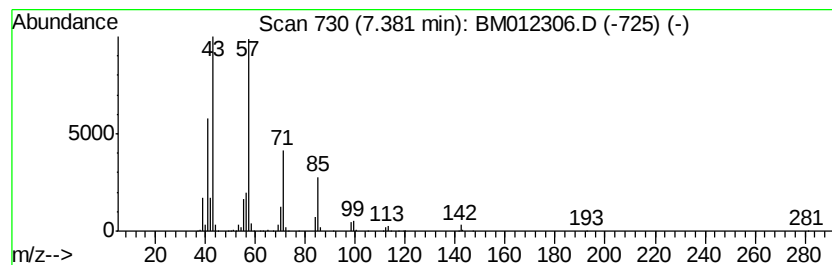
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Quant Title : SVOA CALIBRATION

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

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.38	24.32 ng/ul	2827810	1,4-Dichlorobenzene-d4	7.78

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Decane	142	C10H22	000124-18-5	95
2		Decane	142	C10H22	000124-18-5	90
3		Nonane	128	C9H20	000111-84-2	80
4		Undecane	156	C11H24	001120-21-4	78
5		Sulfurous acid, 2-propyl tetradec...	320	C17H36O3S	1000309-12-5	78



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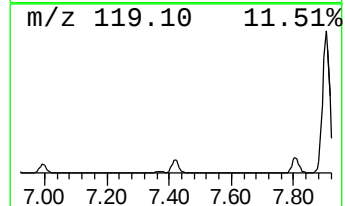
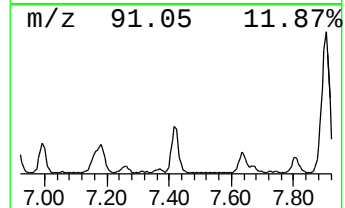
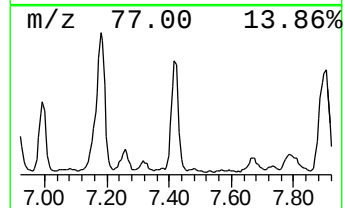
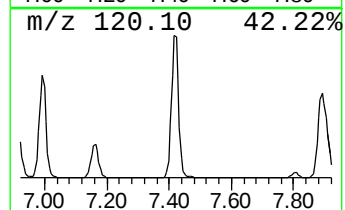
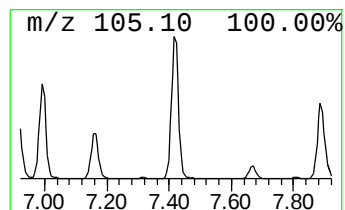
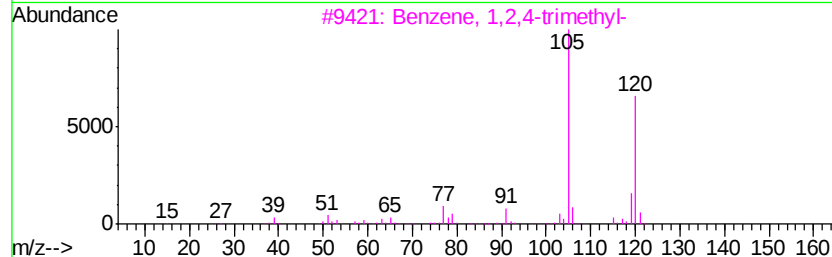
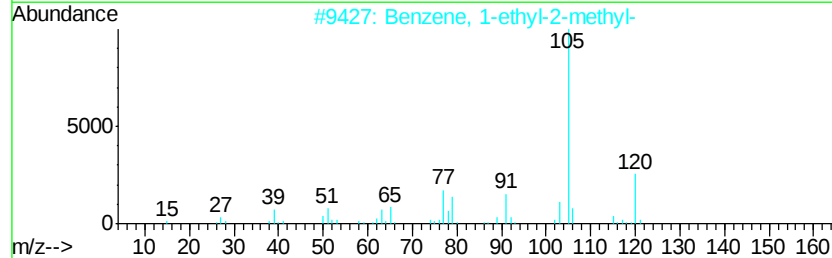
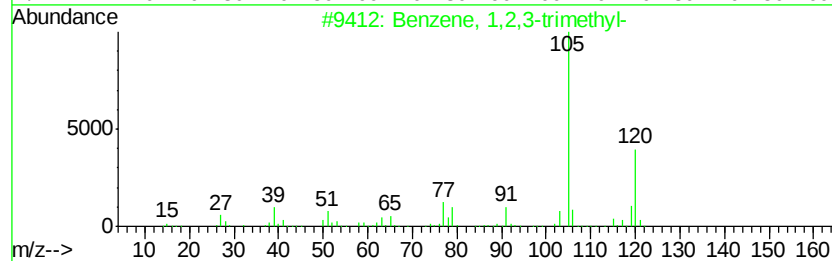
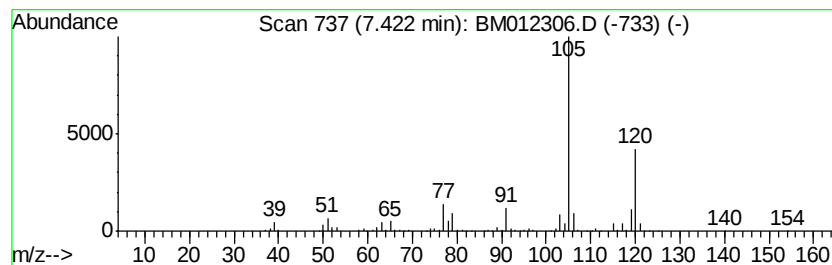
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Benzene, 1,2,3-trimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.42	18.71 ng/ul	2175220	1,4-Dichlorobenzene-d4	7.78

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
2		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	93
3		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	93
4		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	93
5		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91



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ALS Vial : 25 Sample Multiplier: 1

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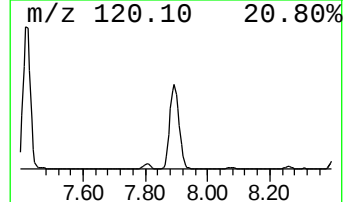
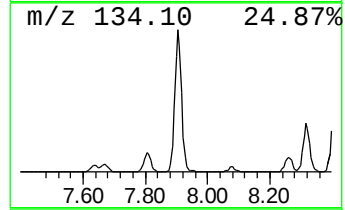
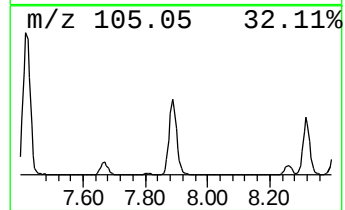
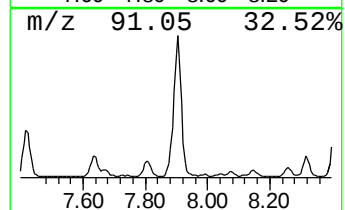
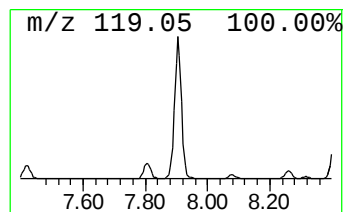
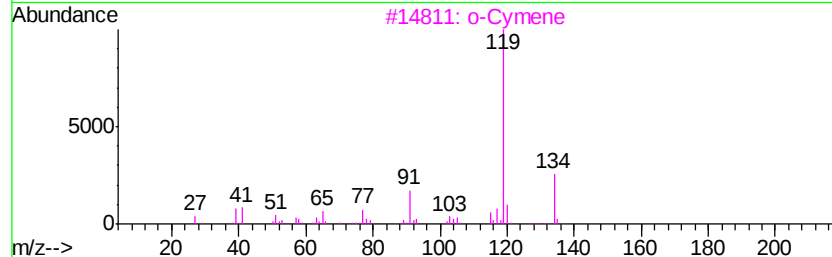
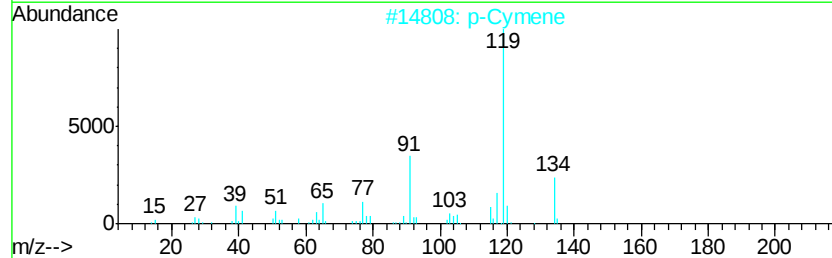
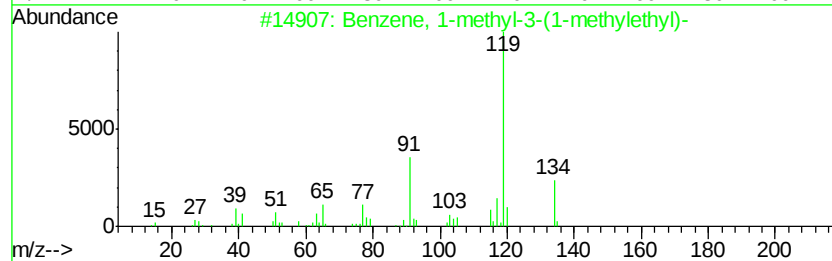
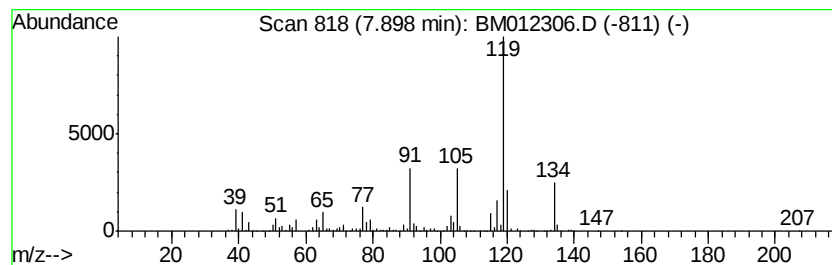
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Benzene, 1-methyl-3-(1-meth... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.90	31.92 ng/ul	3710840	1,4-Dichlorobenzene-d4	7.78

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene,	1-methyl-3-(1-methyleth...	134	C10H14		000535-77-3	97
2	p-Cymene		134	C10H14		000099-87-6	97
3	o-Cymene		134	C10H14		000527-84-4	93
4	o-Cymene		134	C10H14		000527-84-4	92
5	Benzene,	1-ethyl-2,3-dimethyl-	134	C10H14		000933-98-2	91



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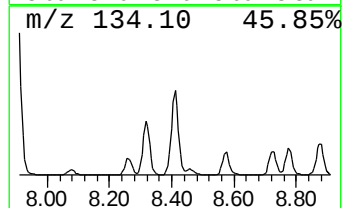
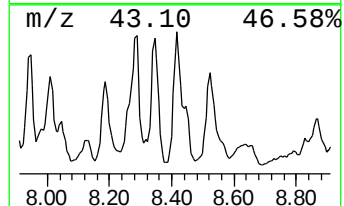
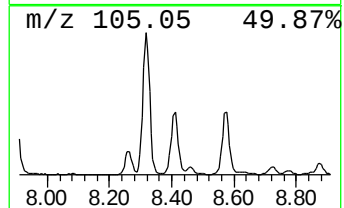
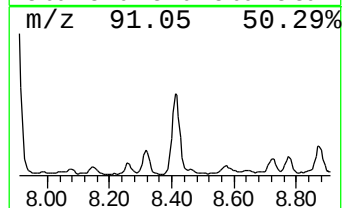
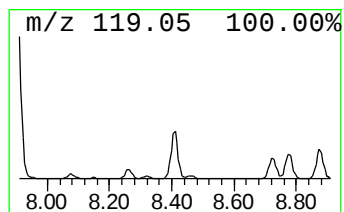
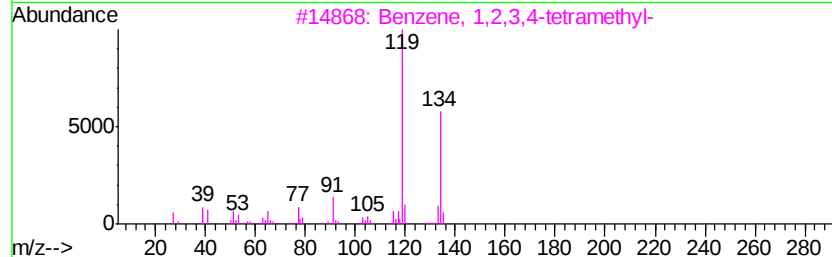
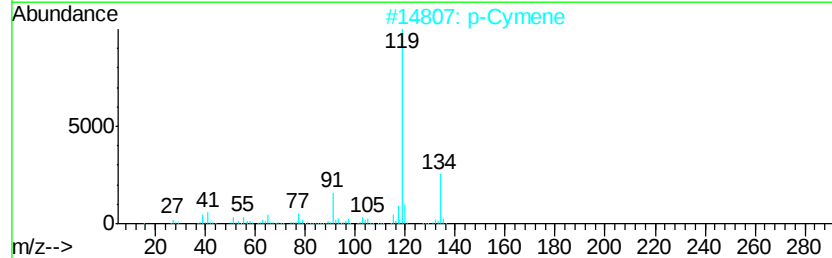
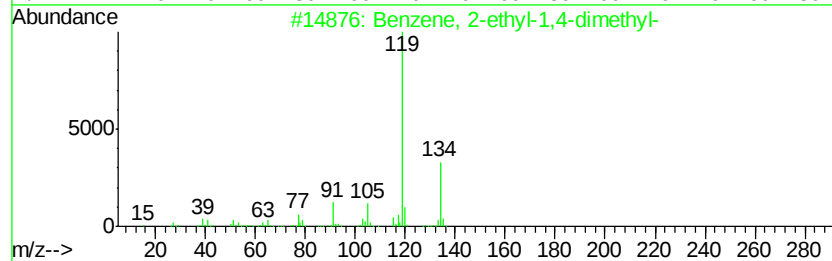
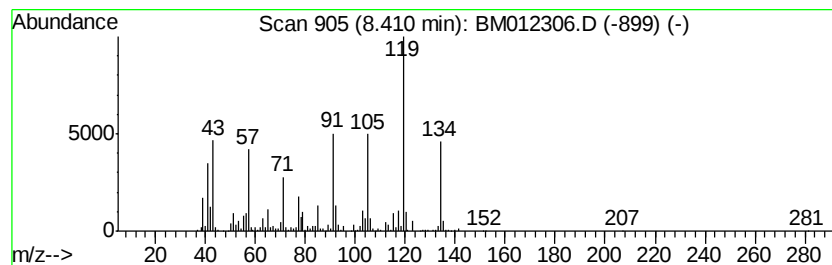
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.41	19.36 ng/ul	2250140	1,4-Dichlorobenzene-d4	7.78

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	93
2		p-Cymene	134	C10H14	000099-87-6	93
3		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	93
4		o-Cymene	134	C10H14	000527-84-4	90
5		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	89



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ClientSampleId :
B-4(1-3)-100617

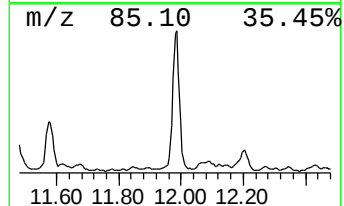
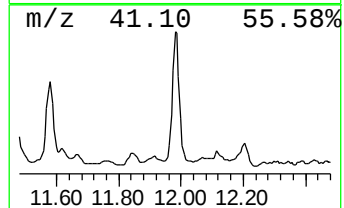
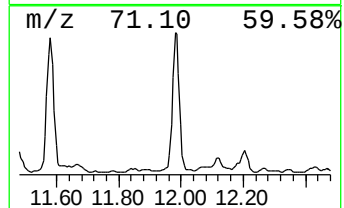
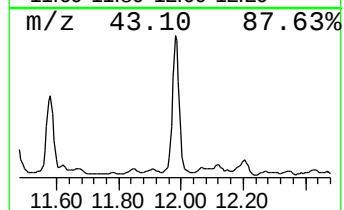
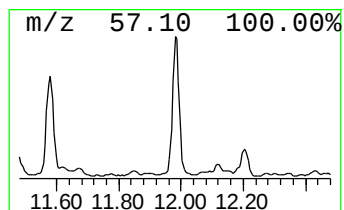
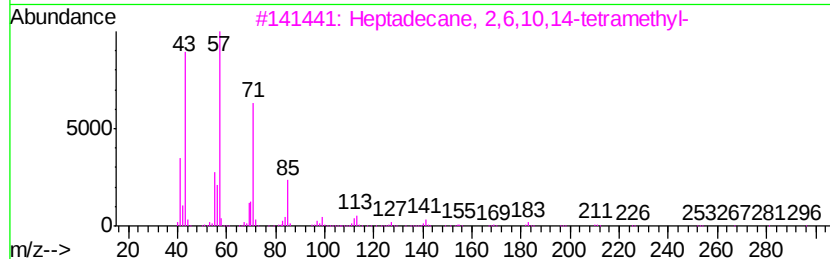
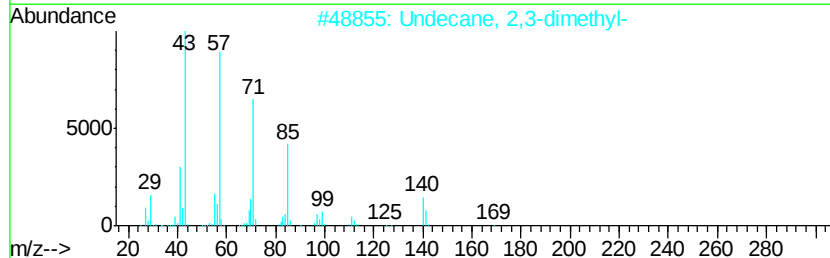
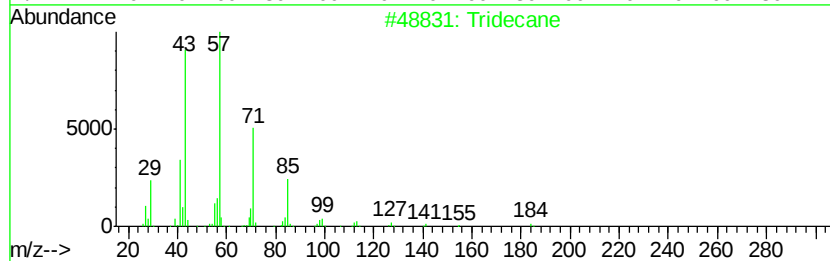
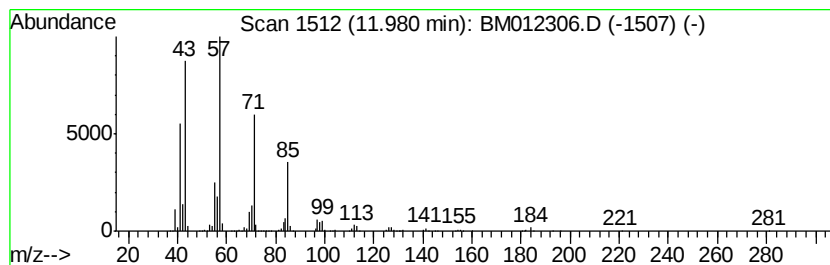
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.98	23.36 ng/ul	2579600	Napthalene-d8	10.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tridecane	184	C13H28	000629-50-5	93
2		Undecane, 2,3-dimethyl-	184	C13H28	017312-77-5	86
3		Heptadecane, 2,6,10,14-tetramethyl-	296	C21H44	018344-37-1	86
4		Octane, 2,4,6-trimethyl-	156	C11H24	062016-37-9	86
5		Hexadecane	226	C16H34	000544-76-3	86



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM101817\
Data File : BM012306.D
Acq On : 19 Oct 2017 07:25
Operator : SJ/JU
Sample : I5747-06 2X
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
B-4(1-3)-100617

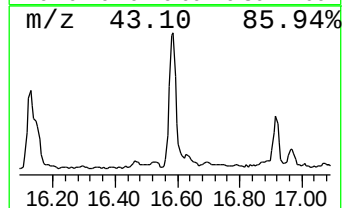
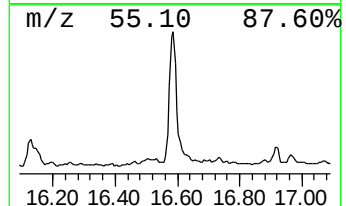
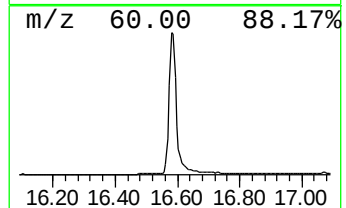
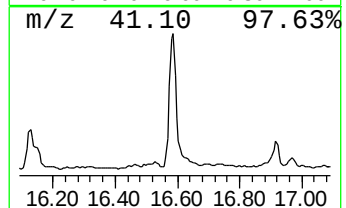
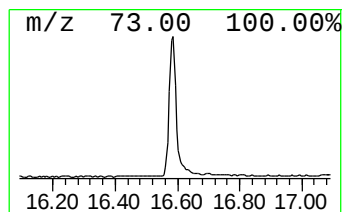
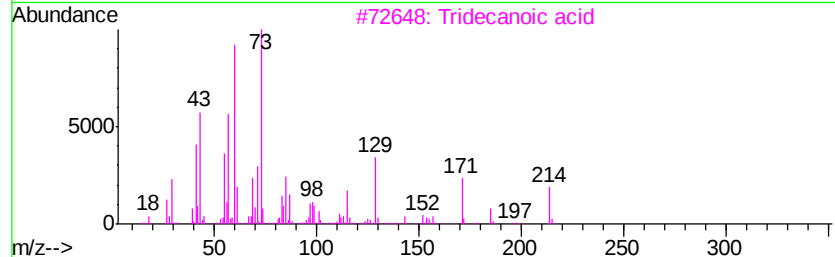
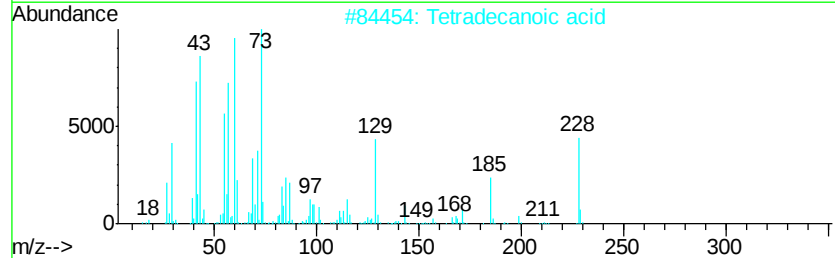
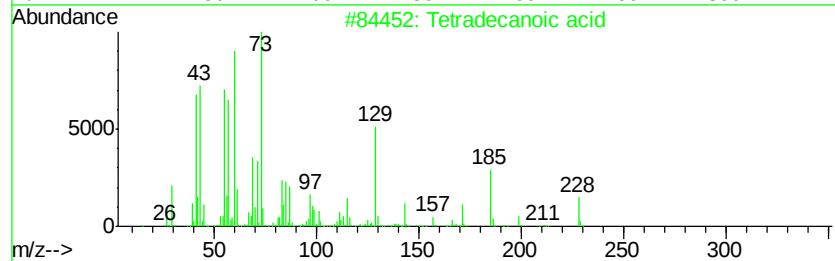
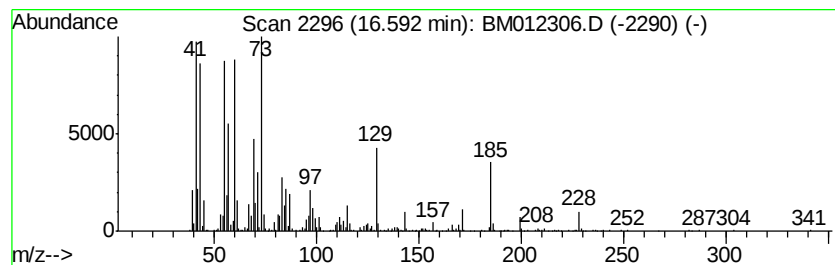
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Tetradecanoic acid Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.59	14.48 ng/ul	2247560	Phenanthrene-d10	17.19

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetradecanoic acid	228	C14H28O2	000544-63-8	96
2		Tetradecanoic acid	228	C14H28O2	000544-63-8	94
3		Tridecanoic acid	214	C13H26O2	000638-53-9	80
4		Tridecanoic acid	214	C13H26O2	000638-53-9	80
5		n-Decanoic acid	172	C10H20O2	000334-48-5	70



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM101817\
Data File : BM012306.D
Acq On : 19 Oct 2017 07:25
Operator : SJ/JU
Sample : I5747-06 2X
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
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ClientSampleId :
B-4(1-3)-100617

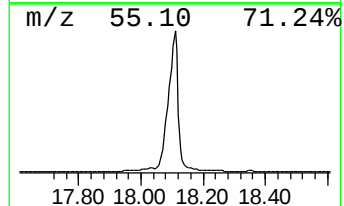
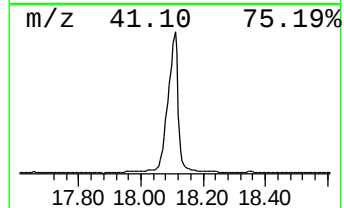
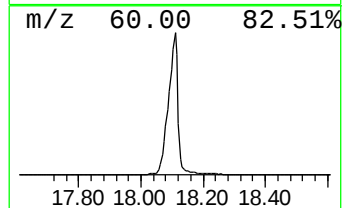
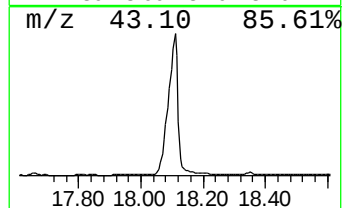
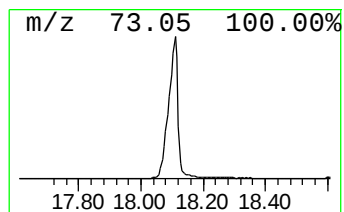
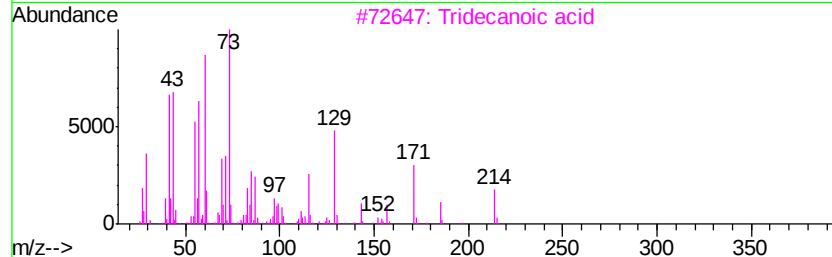
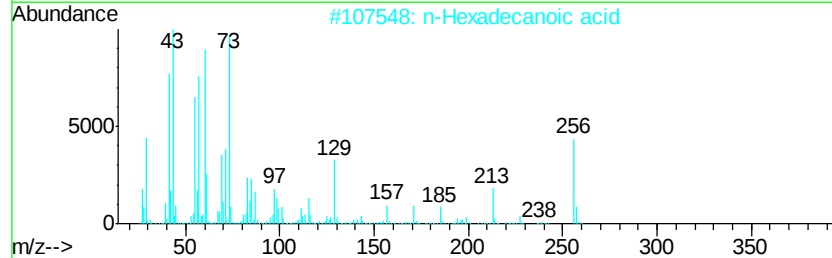
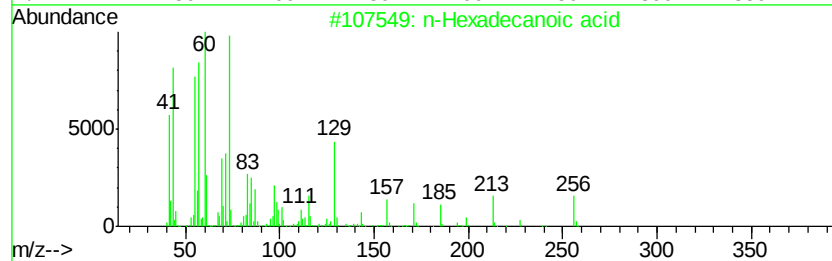
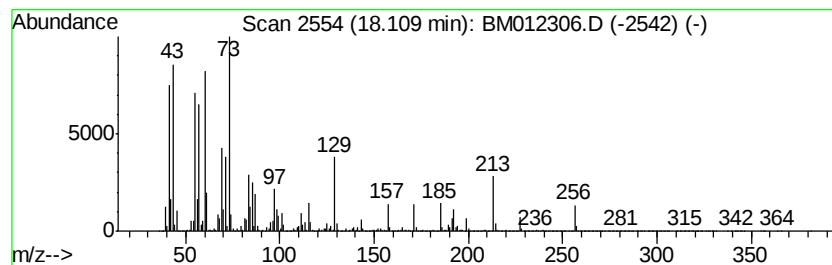
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 n-Hexadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.11	399.25 ng/ul	61984000	Phenanthrene-d10	17.19

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
3		Tridecanoic acid	214	C13H26O2	000638-53-9	92
4		Pentadecanoic acid	242	C15H30O2	001002-84-2	91
5		Tridecanoic acid	214	C13H26O2	000638-53-9	90



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM101817\
Data File : BM012306.D
Acq On : 19 Oct 2017 07:25
Operator : SJ/JU
Sample : I5747-06 2X
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
B-4(1-3)-100617

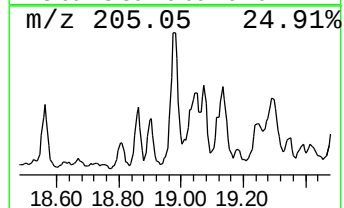
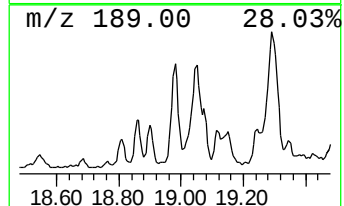
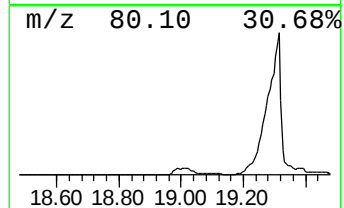
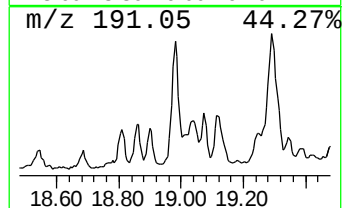
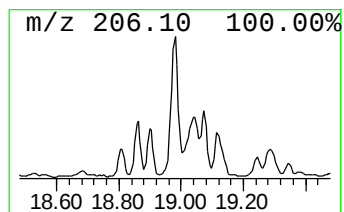
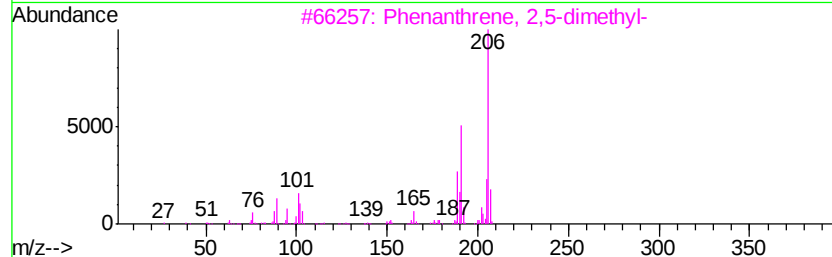
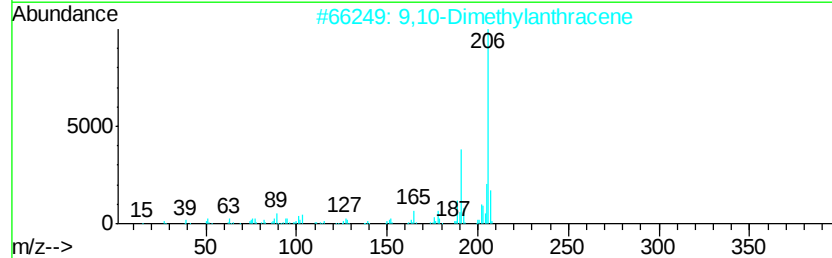
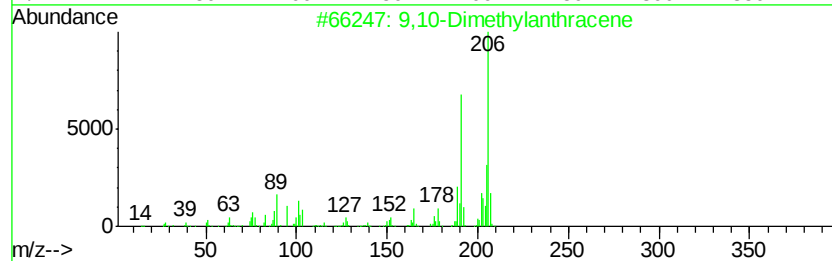
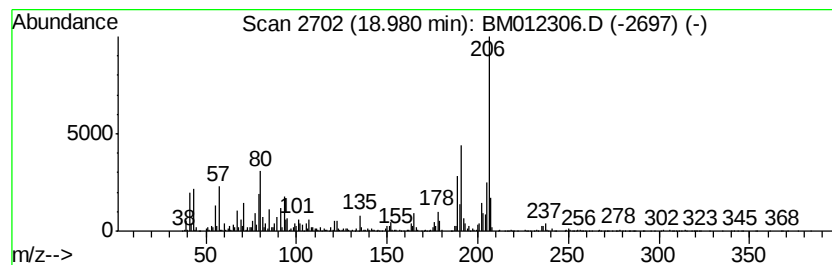
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 9,10-Dimethylantracene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.98	15.23 ng/ul	2364870	Phenanthrene-d10	17.19

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9,10-Dimethylantracene	206	C16H14	000781-43-1	95
2		9,10-Dimethylantracene	206	C16H14	000781-43-1	94
3		Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	93
4		di-p-Tolylacetylene	206	C16H14	002789-88-0	93
5		Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	91



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM101817\
Data File : BM012306.D
Acq On : 19 Oct 2017 07:25
Operator : SJ/JU
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Misc :
ALS Vial : 25 Sample Multiplier: 1

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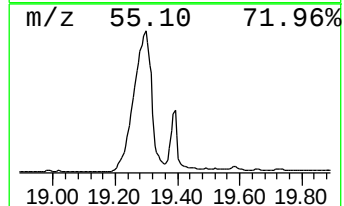
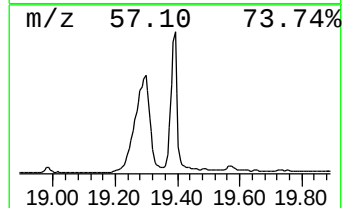
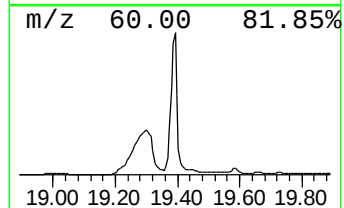
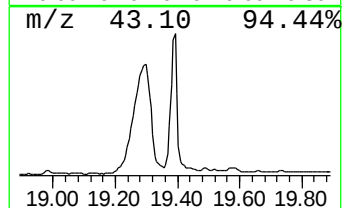
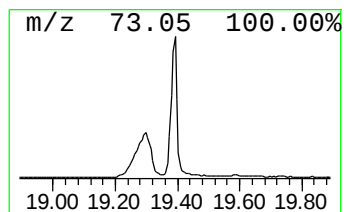
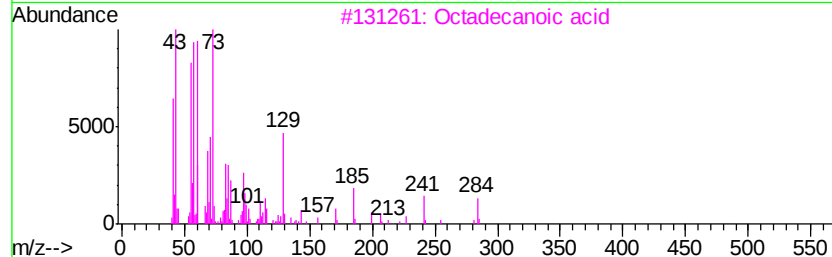
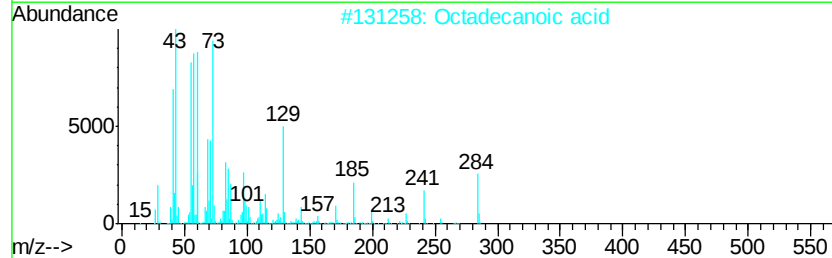
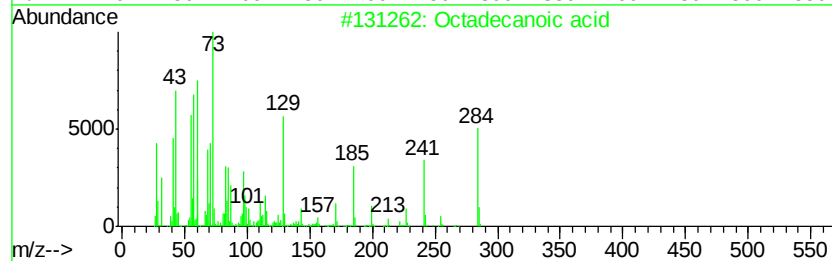
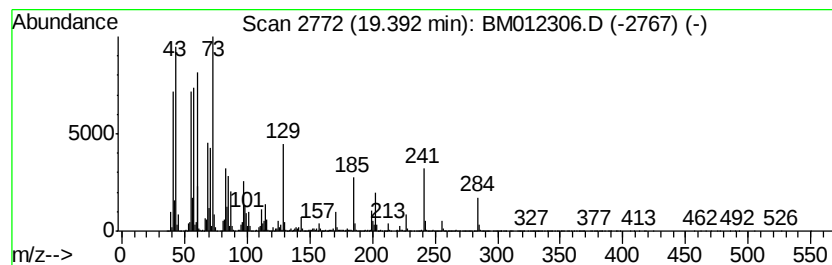
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Octadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.39	38.47 ng/ul	31156500	Chrysene-d12	21.37

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		Octadecanoic acid	284	C18H36O2	000057-11-4	99
4		Octadecanoic acid	284	C18H36O2	000057-11-4	98
5		Pentadecanoic acid	242	C15H30O2	001002-84-2	86



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM101817\
Data File : BM012306.D
Acq On : 19 Oct 2017 07:25
Operator : SJ/JU
Sample : I5747-06 2X
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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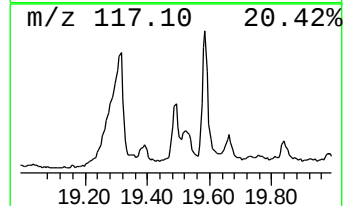
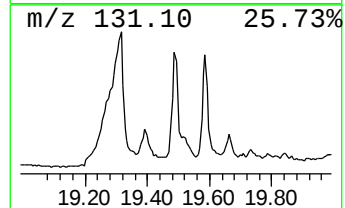
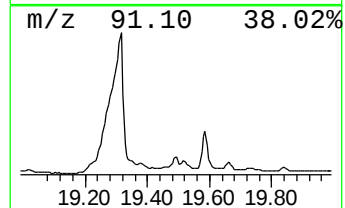
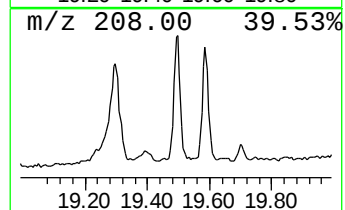
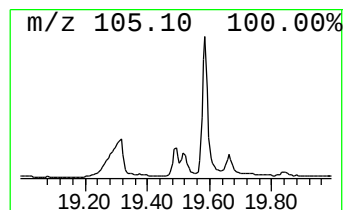
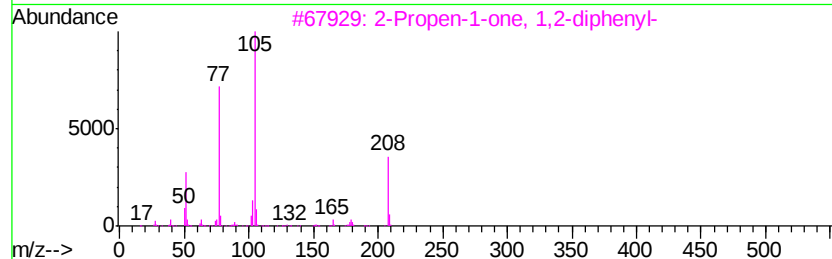
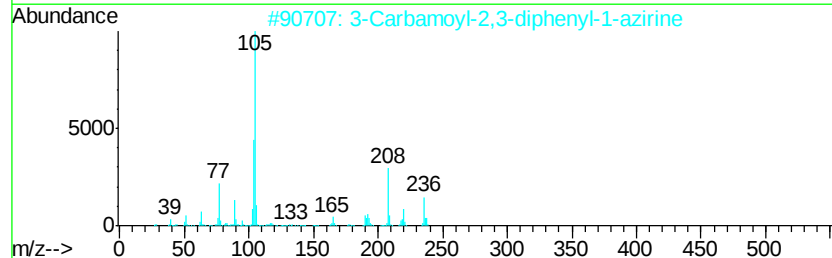
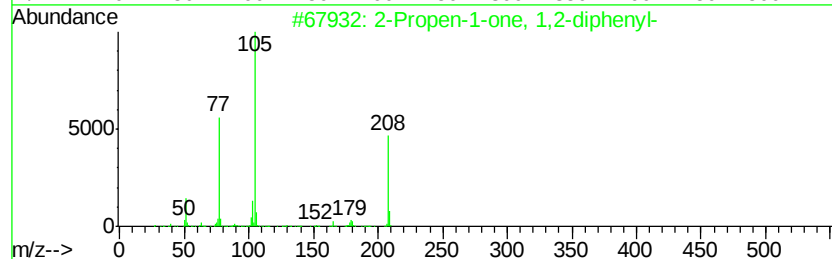
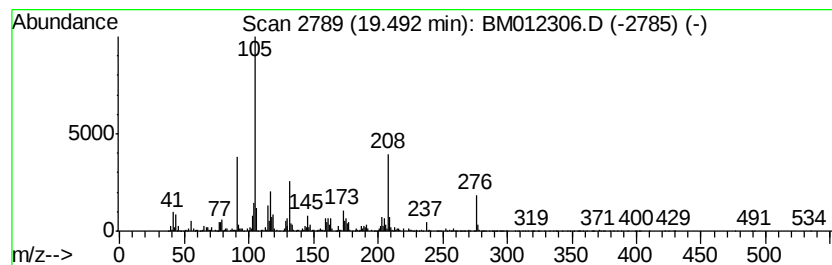
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 unknown-01 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.49	3.27 ng/ul	2648300	Chrysene-d12	21.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Propen-1-one, 1,2-diphenyl-	208	C15H12O	004452-11-3	25
2		3-Carbamoyl-2,3-diphenyl-1-azirine	236	C15H12N2O	033859-59-5	23
3		2-Propen-1-one, 1,2-diphenyl-	208	C15H12O	004452-11-3	17
4		Benzene, 1,1',1''-[1-(3-methyl-3...	410	C31H38	074685-55-5	16
5		4,4,6-Trimethyl-2-phenyl-5,6-dih...	203	C13H17NO	026939-21-9	11



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM101817\
Data File : BM012306.D
Acq On : 19 Oct 2017 07:25
Operator : SJ/JU
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Misc :
ALS Vial : 25 Sample Multiplier: 1

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ClientSampleId :
B-4(1-3)-100617

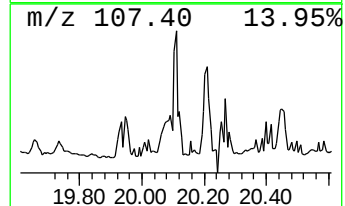
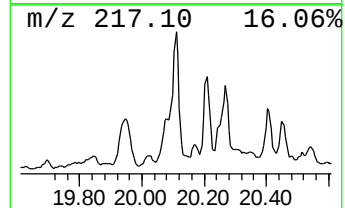
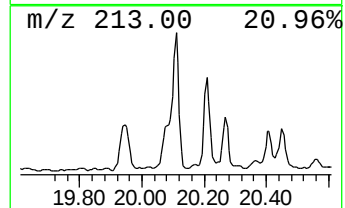
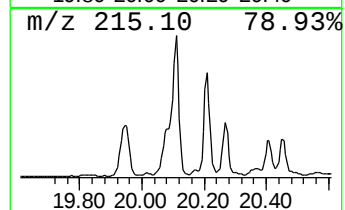
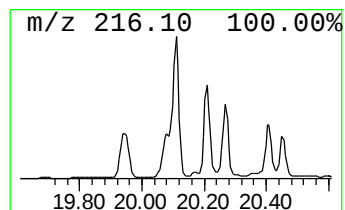
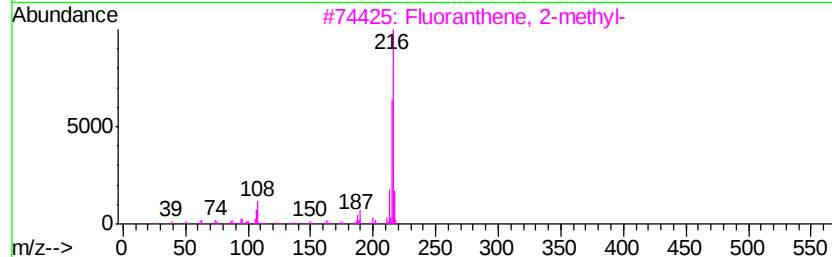
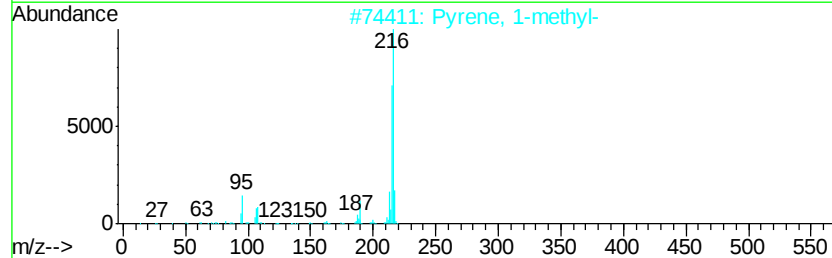
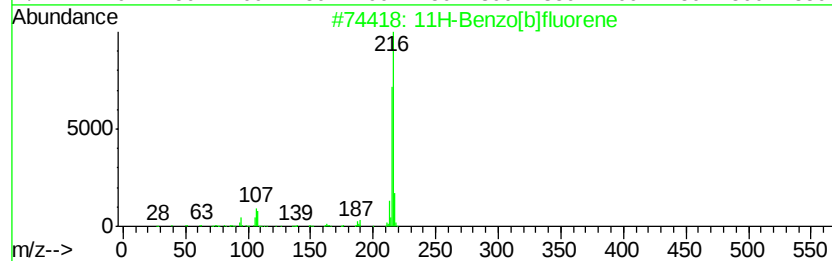
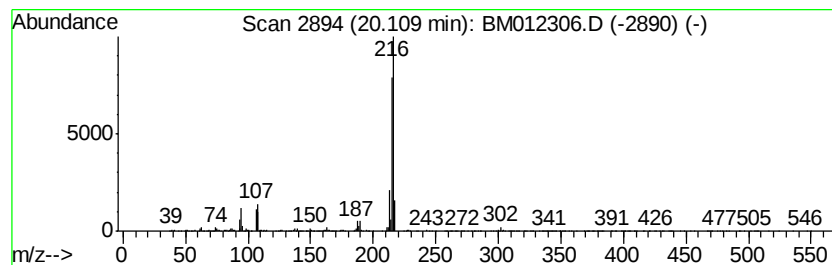
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 11H-Benzo[b]fluorene Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.11	3.79 ng/ul	3071000	Chrysene-d12	21.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	87
2		Pyrene, 1-methyl-	216	C17H12	002381-21-7	83
3		Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	81
4		11H-Benzo[a]fluorene	216	C17H12	000238-84-6	72
5		Pyrene, 1-methyl-	216	C17H12	002381-21-7	72



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM101817\
Data File : BM012306.D
Acq On : 19 Oct 2017 07:25
Operator : SJ/JU
Sample : I5747-06 2X
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
B-4(1-3)-100617

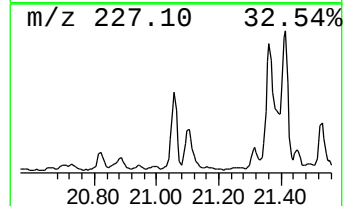
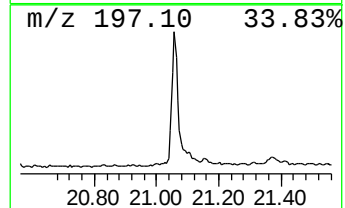
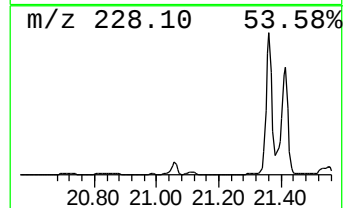
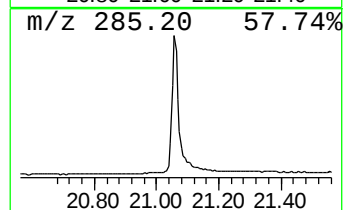
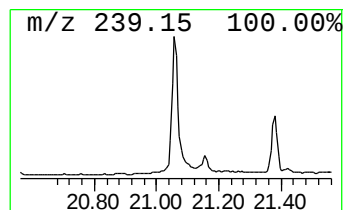
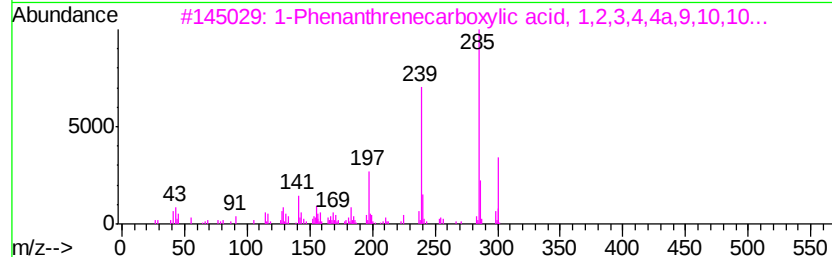
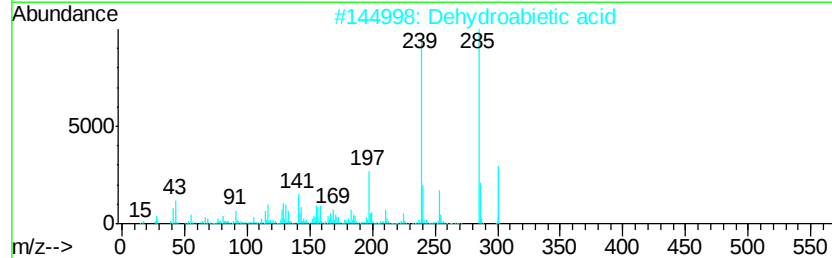
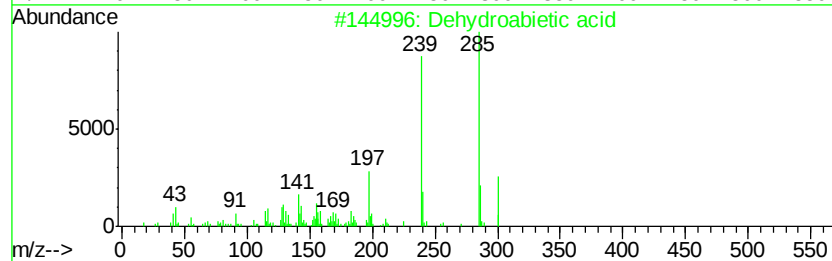
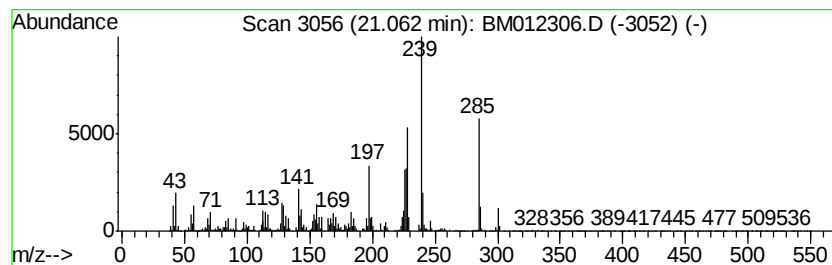
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 Dehydroabietic acid Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.06	5.09 ng/ul	4124010	Chrysene-d12	21.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dehydroabietic acid	300	C20H28O2	001740-19-8	99
2		Dehydroabietic acid	300	C20H28O2	001740-19-8	83
3		1-Phenanthrenecarboxylic acid, 1...	300	C20H28O2	005155-70-4	83
4		Dehydroabietic acid	300	C20H28O2	001740-19-8	64
5		trans-1,2-Bis(methyldichlorosilyl...	252	C4H8Cl4Si2	065899-10-7	56



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM101817\
Data File : BM012306.D
Acq On : 19 Oct 2017 07:25
Operator : SJ/JU
Sample : I5747-06 2X
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
B-4(1-3)-100617

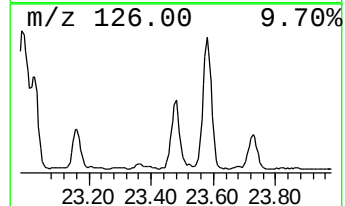
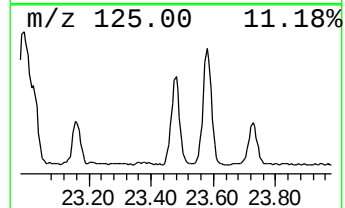
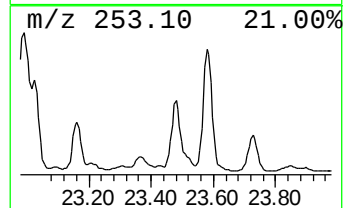
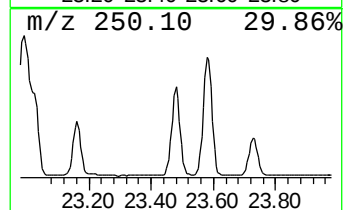
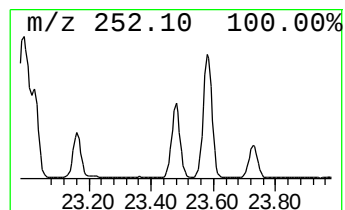
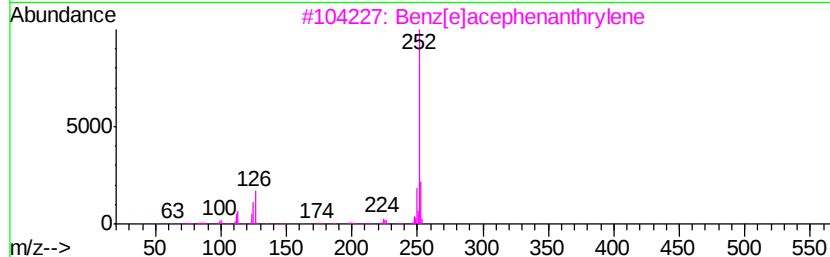
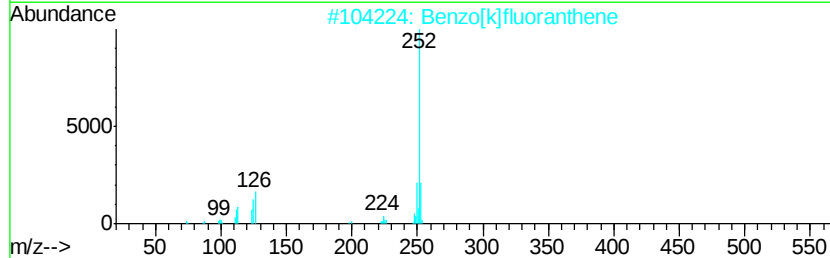
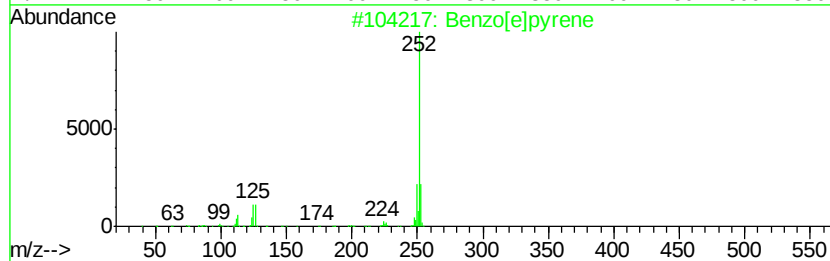
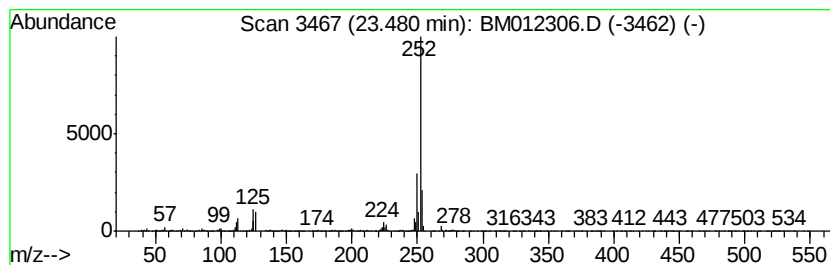
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM101217.M
Quant Title : SVOA CALIBRATION

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

Peak Number 16 Benzo[e]pyrene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.48	41.02 ng/ul	4942510	Perylene-d12	23.68

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[e]pyrene	252	C20H12	000192-97-2	99
2		Benzo[k]fluoranthene	252	C20H12	000207-08-9	97
3		Benz[e]acephenanthrylene	252	C20H12	000205-99-2	95
4		Benz[e]acephenanthrylene	252	C20H12	000205-99-2	94
5		Benzo[a]pyrene	252	C20H12	000050-32-8	93



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM101817\
 Data File : BM012306.D
 Acq On : 19 Oct 2017 07:25
 Operator : SJ/JU
 Sample : I5747-06 2X
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 B-4(1-3)-100617

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM101217.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
Cyclohexane, 1-me...	7.18	156.1	ng/ul	18148200	1	7.78	2325070	20.0
(DEL) Alkane: Str...	7.38	24.3	ng/ul	2827810	1	7.78	2325070	20.0
Benzene, 1,2,3-tr...	7.42	18.7	ng/ul	2175220	1	7.78	2325070	20.0
Benzene, 1-methyl...	7.90	31.9	ng/ul	3710840	1	7.78	2325070	20.0
Benzene, 2-ethyl-...	8.41	19.4	ng/ul	2250140	1	7.78	2325070	20.0
(DEL) Alkane: Str...	11.98	23.4	ng/ul	2579600	2	10.58	2208730	20.0
Tetradecanoic acid	16.59	14.5	ng/ul	2247560	4	17.19	3105040	20.0
n-Hexadecanoic acid	18.11	399.3	ng/ul	61984000	4	17.19	3105040	20.0
9,10-Dimethylanthe...	18.98	15.2	ng/ul	2364870	4	17.19	3105040	20.0
Octadecanoic acid	19.39	38.5	ng/ul	31156500	5	21.37	16199500	20.0
unknown-01	19.49	3.3	ng/ul	2648300	5	21.37	16199500	20.0
11H-Benzo[b]fluorene	20.11	3.8	ng/ul	3071000	5	21.37	16199500	20.0
Dehydroabietic acid	21.06	5.1	ng/ul	4124010	5	21.37	16199500	20.0
Benzo[e]pyrene	23.48	41.0	ng/ul	4942510	6	23.68	2409530	20.0