

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG022819\
 Data File : BG039671.D
 Acq On : 28 Feb 2019 22:14
 Operator : JU/SJ
 Sample : K1628-04
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 A0C6-2-0.5-1.0

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 3 % 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_G\METHODS\8270-BG012919.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.220	17	22	31	rBV2	77652	157489	4.37%	0.697%
2	3.291	31	34	44	rVB	48244	76501	2.12%	0.339%
3	5.711	437	446	457	rBV	1075062	1708923	47.42%	7.565%
4	7.315	710	719	732	rBV	1081297	1983727	55.05%	8.781%
5	7.691	776	783	800	rVB	1037122	1827107	50.70%	8.088%
6	8.166	856	864	875	rVB	180196	310718	8.62%	1.375%
7	8.490	911	919	927	rBV	741830	1301891	36.13%	5.763%
8	9.341	1056	1064	1075	rBV	649824	1168921	32.44%	5.174%
9	10.986	1336	1344	1352	rBV	247604	446506	12.39%	1.977%
10	13.412	1747	1757	1767	rBV	1670604	2609341	72.41%	11.551%
11	14.229	1890	1896	1903	rBV	93958	132390	3.67%	0.586%
12	14.787	1984	1991	1999	rBV2	394130	648211	17.99%	2.869%
13	16.273	2237	2244	2259	rBV	1384411	2131902	59.16%	9.437%
14	17.530	2450	2458	2469	rBV	534375	824908	22.89%	3.652%
15	18.159	2560	2565	2571	rVB	101753	127043	3.53%	0.562%
16	18.382	2597	2603	2615	rBV	141357	259854	7.21%	1.150%
17	19.451	2780	2785	2789	rBV	87523	103751	2.88%	0.459%
18	19.533	2792	2799	2810	rBV4	46477	104202	2.89%	0.461%
19	19.645	2815	2818	2827	rVV4	29443	51297	1.42%	0.227%
20	19.786	2839	2842	2849	rVB	144692	175288	4.86%	0.776%
21	20.127	2893	2900	2908	rBV	2625797	3603548	100.00%	15.952%
22	20.826	3014	3019	3025	rVB	238795	272686	7.57%	1.207%
23	21.413	3114	3119	3130	rBV2	163646	322363	8.95%	1.427%
24	21.819	3181	3188	3197	rVB	538809	907543	25.18%	4.017%
25	22.600	3315	3321	3325	rBV2	37406	64507	1.79%	0.286%
26	23.317	3436	3443	3449	rBV2	52737	101561	2.82%	0.450%
27	24.151	3579	3585	3594	rVB2	47457	98684	2.74%	0.437%
28	25.196	3746	3763	3774	rBV2	282793	956638	26.55%	4.235%
29	26.324	3948	3955	3964	rVB5	26486	75310	2.09%	0.333%
30	27.746	4193	4197	4207	rVB7	14162	37621	1.04%	0.167%

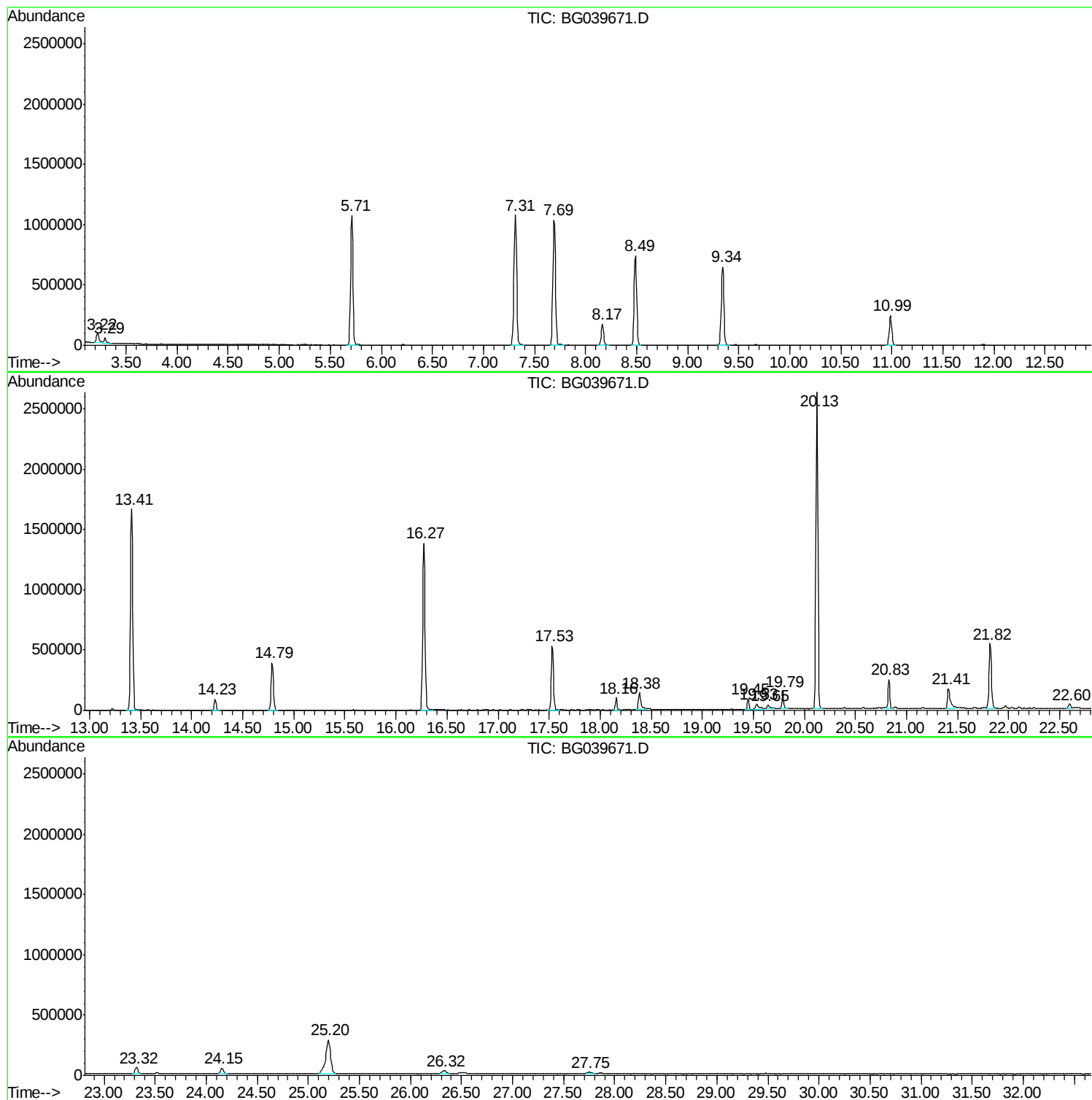
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG012919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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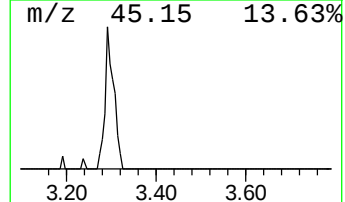
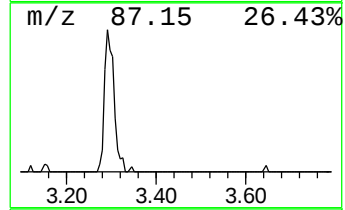
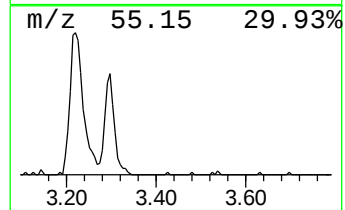
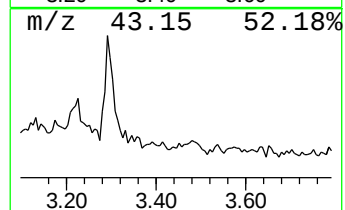
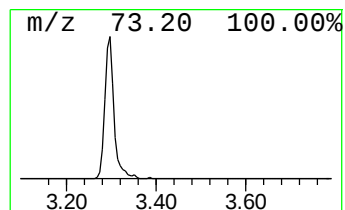
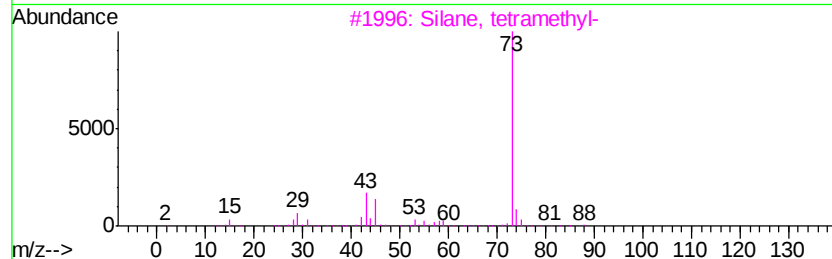
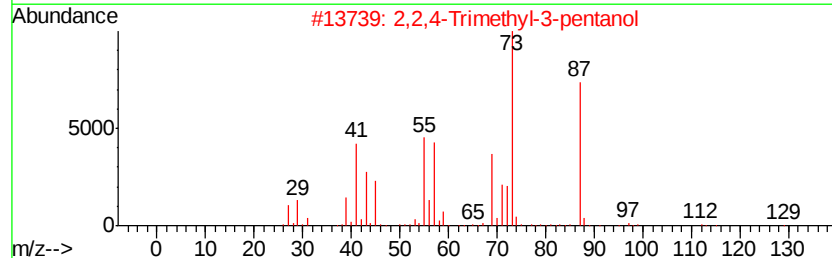
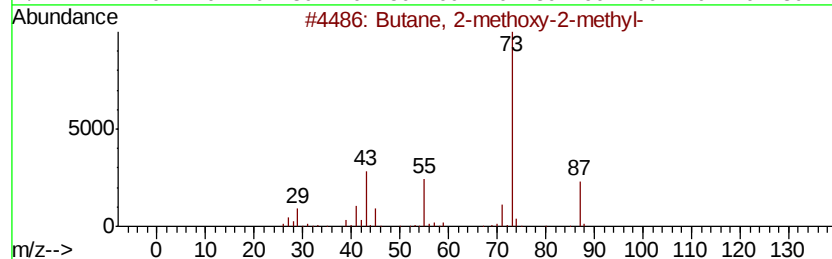
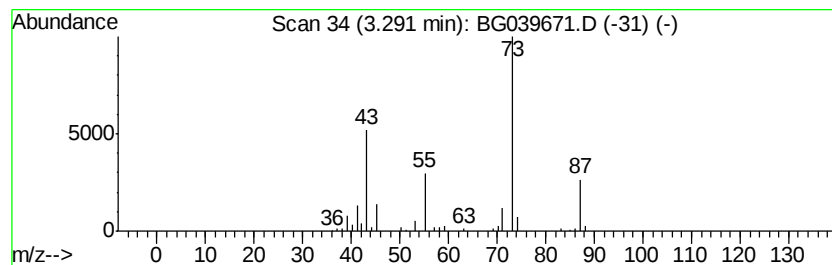
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Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.29	4.92 ng	76501	1,4-Dichlorobenzene-d4	8.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
3		Silane, tetramethyl-	88	C4H12Si	000075-76-3	37
4		1-Propene-1-thiol	74	C3H6S	000925-89-3	35
5		2-Methyl-5-hexen-3-ol	114	C7H14O	032815-70-6	28



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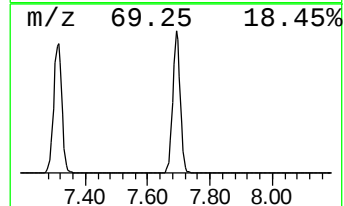
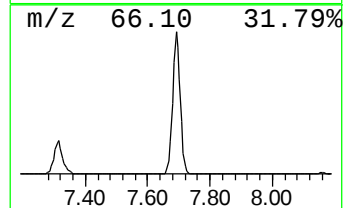
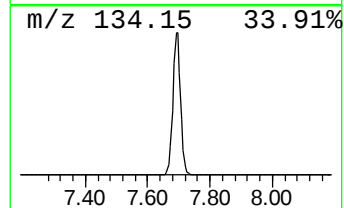
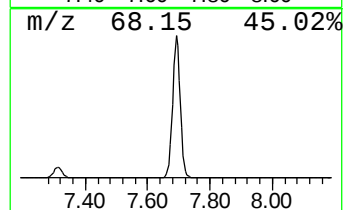
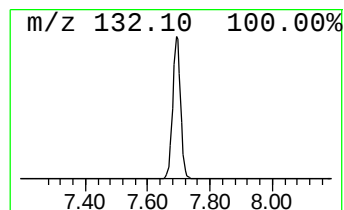
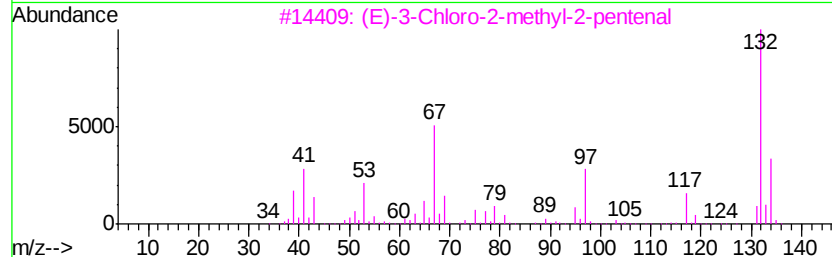
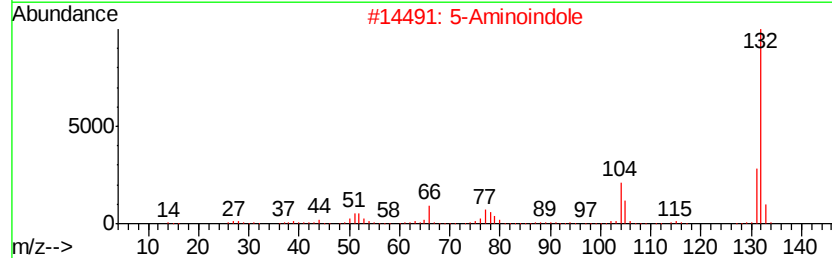
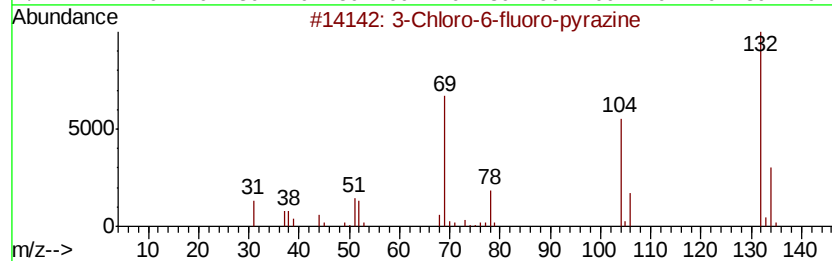
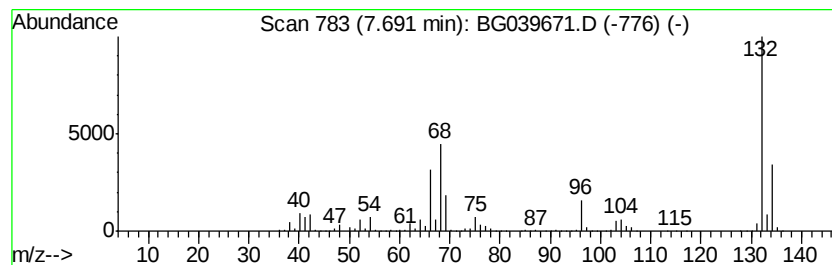
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Peak Number 3 unknown7.69 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.69	117.61 ng	1827110	1,4-Dichlorobenzene-d4	8.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Aminoindole	132	C8H8N2	005192-03-0	22
3		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
4		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	12
5		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12



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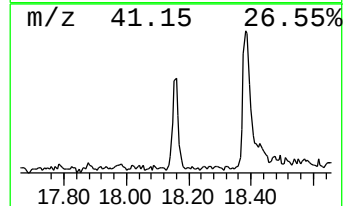
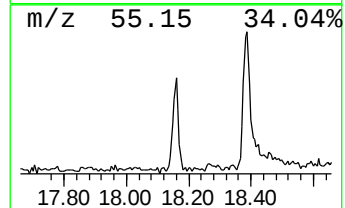
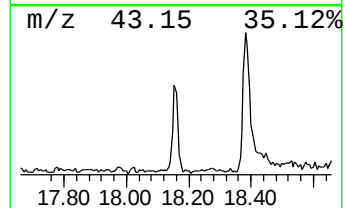
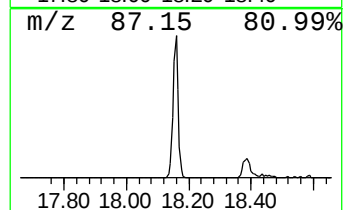
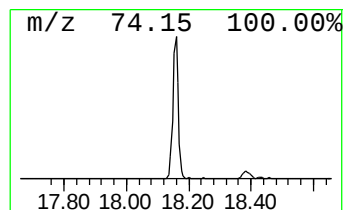
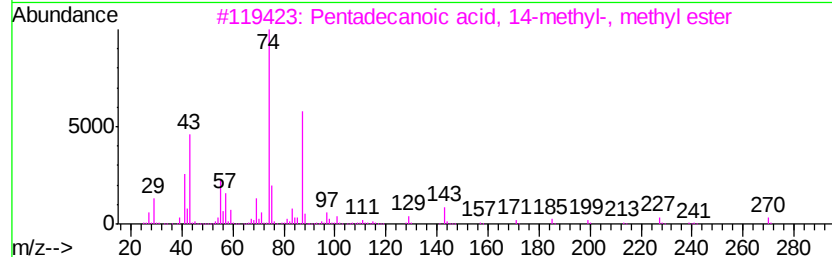
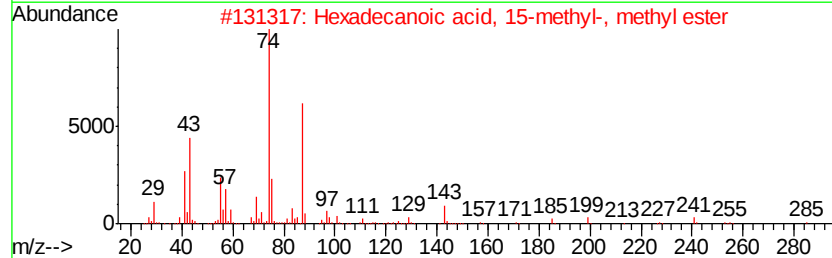
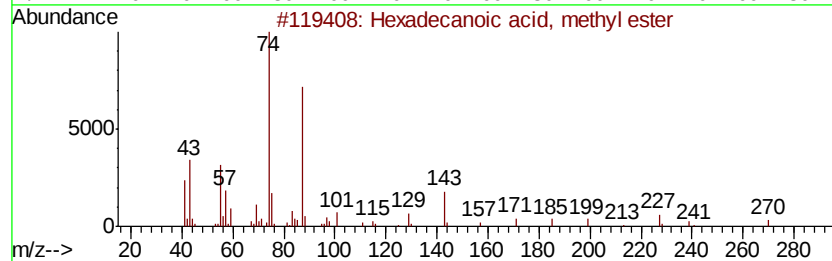
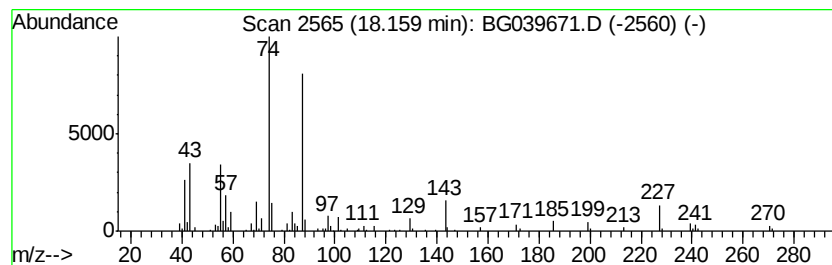
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Peak Number 5 Hexadecanoic acid, methyl e... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.16	3.08 ng	127043	Phenanthrene-d10	17.53

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	98
2		Hexadecanoic acid, 15-methyl-, m...	284	C18H36O2	006929-04-0	87
3		Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	86
4		Tridecanoic acid, 12-methyl-, me...	242	C15H30O2	005129-58-8	80
5		Methyl tetradecanoate	242	C15H30O2	000124-10-7	80



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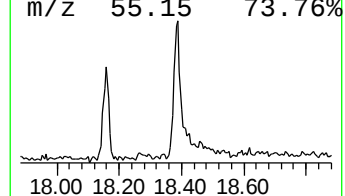
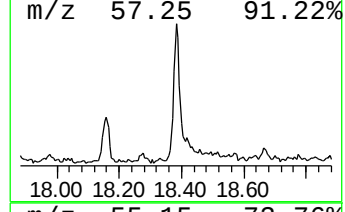
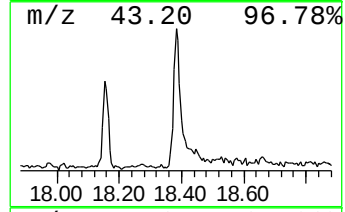
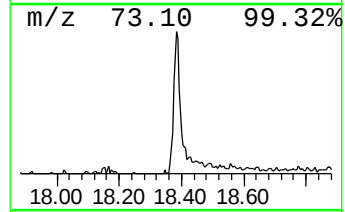
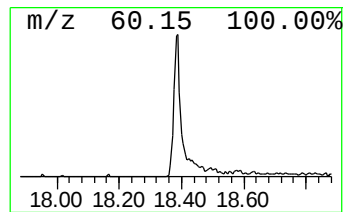
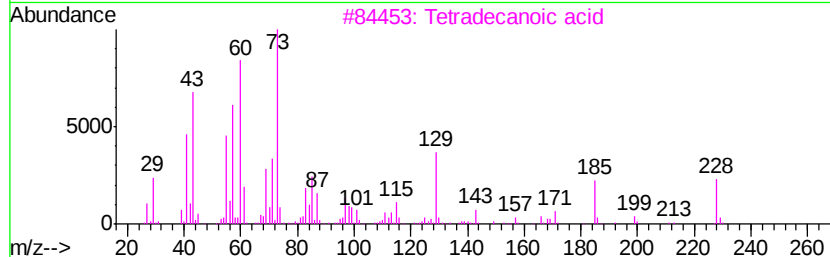
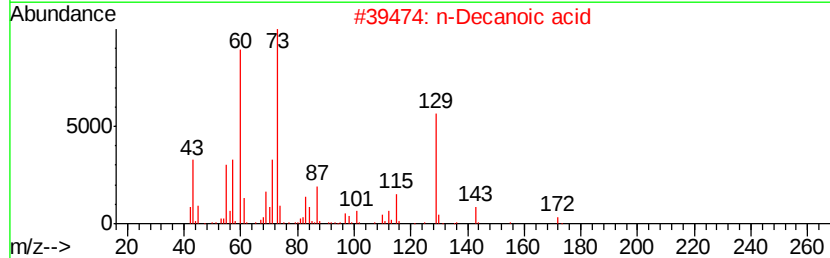
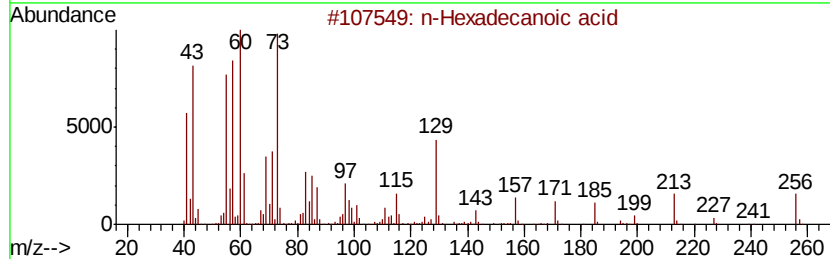
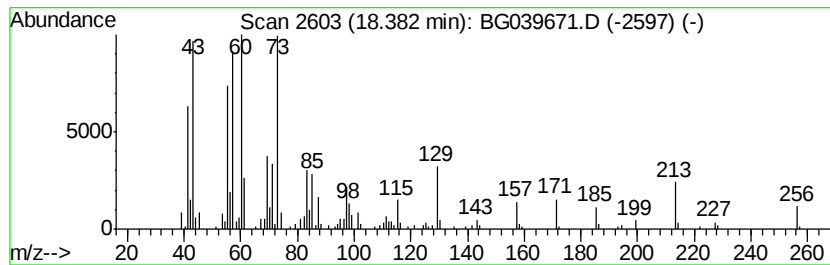
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Peak Number 6 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.38	6.30 ng	259854	Phenanthrene-d10	17.53

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		n-Decanoic acid	172	C10H20O2	000334-48-5	87
3		Tetradecanoic acid	228	C14H28O2	000544-63-8	72
4		Pentadecanoic acid	242	C15H30O2	001002-84-2	72
5		Tridecanoic acid	214	C13H26O2	000638-53-9	70



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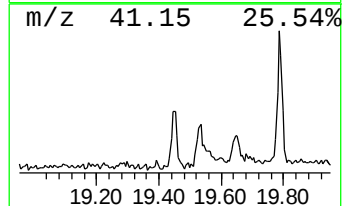
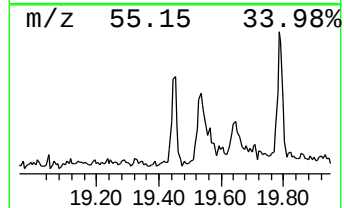
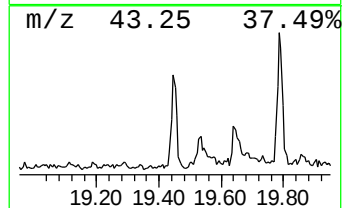
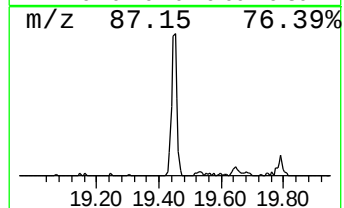
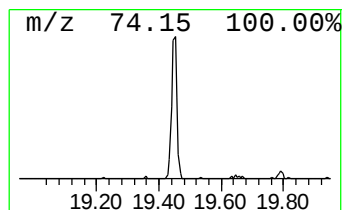
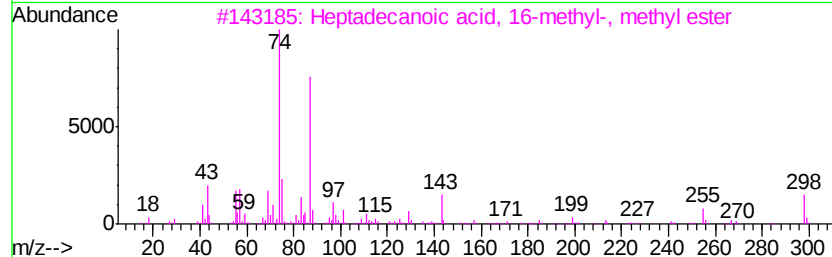
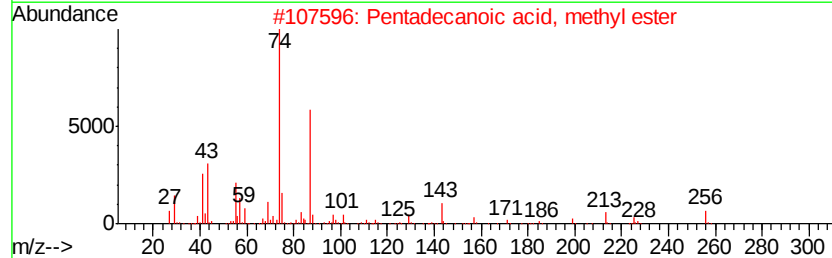
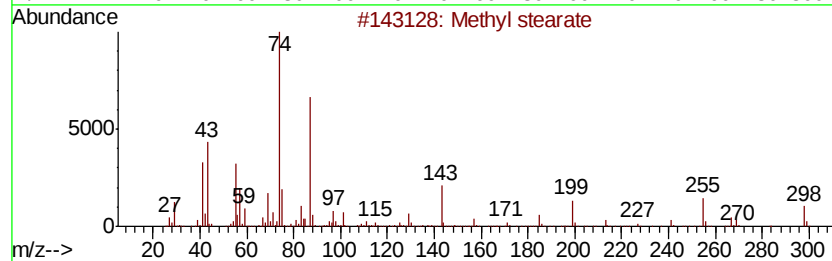
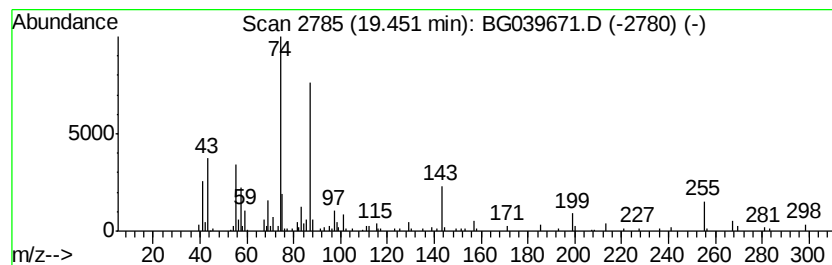
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Peak Number 7 Methyl stearate Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.45	2.52 ng	103751	Phenanthrene-d10	17.53
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Methyl stearate	298 C19H38O2	000112-61-8 90
2		Pentadecanoic acid, methyl ester	256 C16H32O2	007132-64-1 87
3		Heptadecanoic acid, 16-methyl-, ...	298 C19H38O2	005129-61-3 83
4		Hexadecanoic acid, 15-methyl-, m...	284 C18H36O2	006929-04-0 80
5		Methyl tetradecanoate	242 C15H30O2	000124-10-7 80



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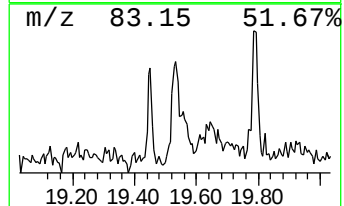
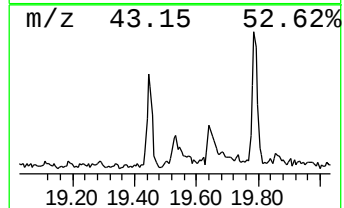
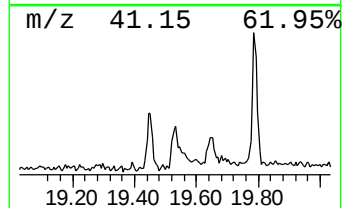
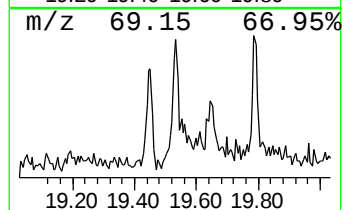
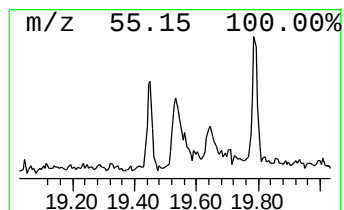
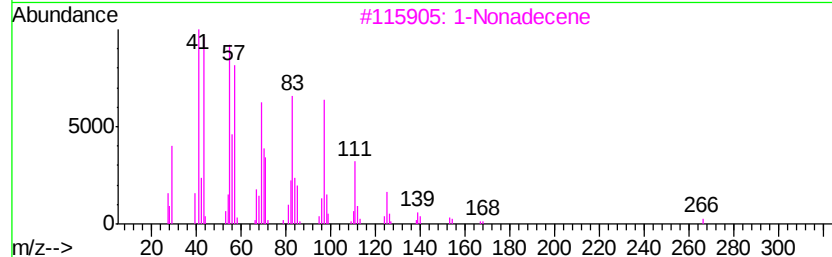
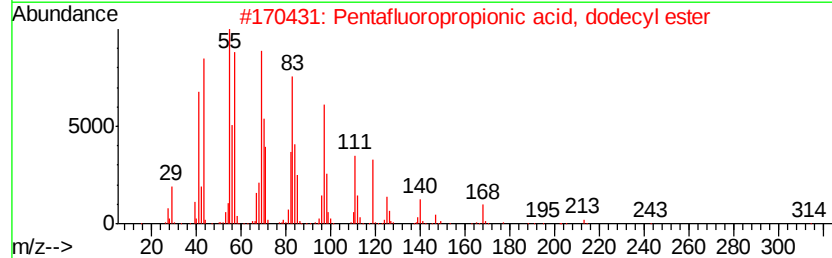
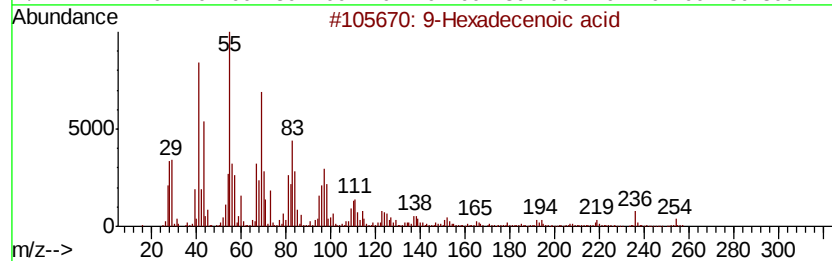
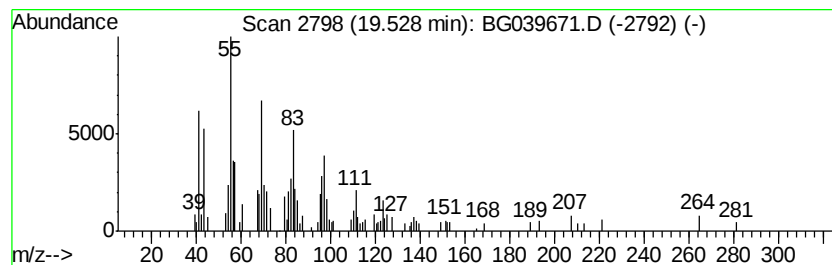
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 unknown19.53 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.53	2.53 ng	104202	Phenanthrene-d10	17.53

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9-Hexadecenoic acid	254	C16H30O2	002091-29-4	49
2		Pentafluoropropionic acid, dodec...	332	C15H25F5O2	006222-04-4	46
3		1-Nonadecene	266	C19H38	018435-45-5	45
4		Acetic acid n-octadecyl ester	312	C20H40O2	000822-23-1	45
5		Acetic acid, chloro-, hexadecyl ...	318	C18H35ClO2	052132-58-8	45



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG022819\
Data File : BG039671.D
Acq On : 28 Feb 2019 22:14
Operator : JU/SJ
Sample : K1628-04
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A0C6-2-0.5-1.0

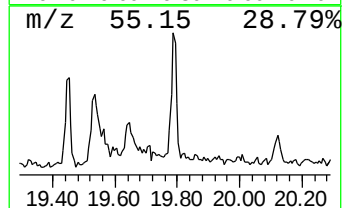
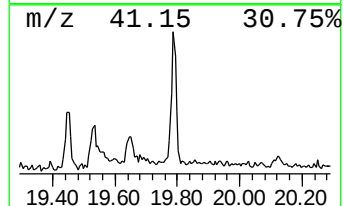
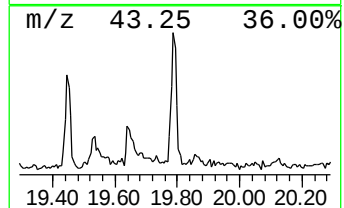
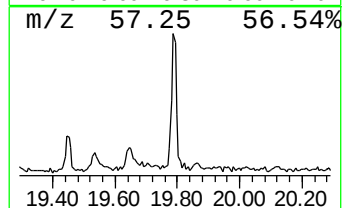
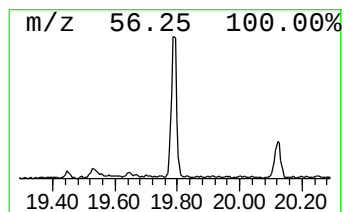
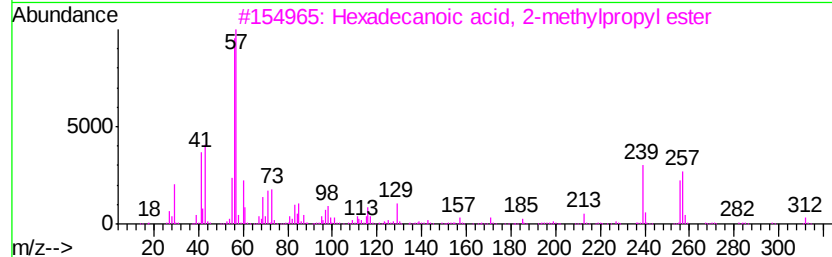
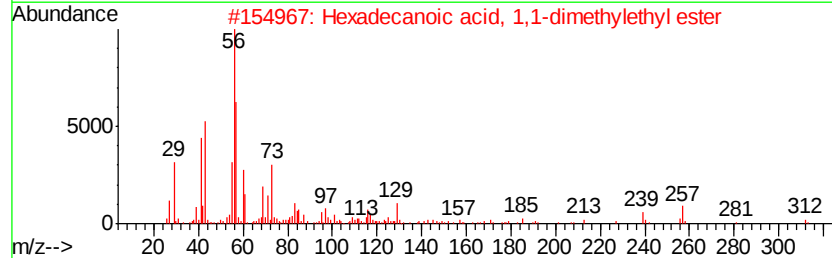
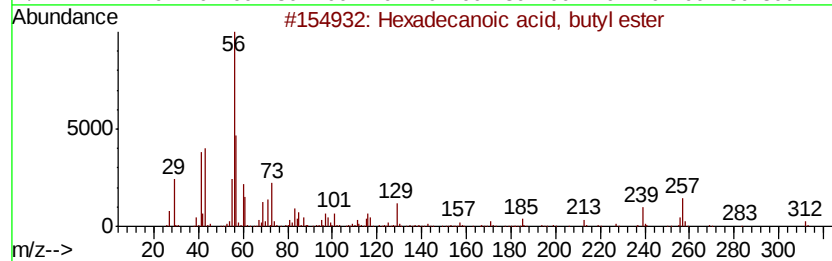
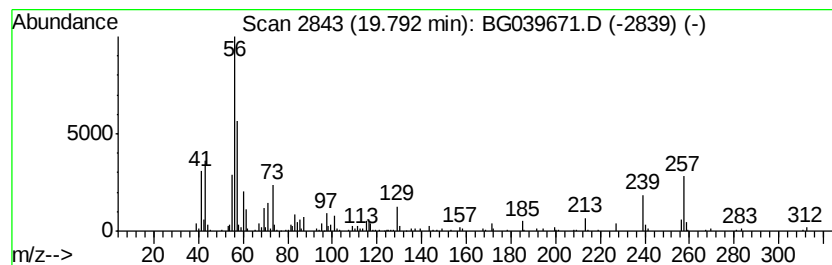
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Hexadecanoic acid, butyl ester Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.79	3.86 ng	175288	Chrysene-d12	21.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecanoic acid, butyl ester	312	C20H40O2	000111-06-8	97
2		Hexadecanoic acid, 1,1-dimethyle...	312	C20H40O2	031158-91-5	92
3		Hexadecanoic acid, 2-methylpropyl...	312	C20H40O2	000110-34-9	50
4		1-Hexene, 2,5-dimethyl-	112	C8H16	006975-92-4	38
5		Cyclobutane, 1-hexyl-2,3-dimethyl-	168	C12H24	055170-84-8	35



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG022819\
Data File : BG039671.D
Acq On : 28 Feb 2019 22:14
Operator : JU/SJ
Sample : K1628-04
Misc :
ALS Vial : 13 Sample Multiplier: 1

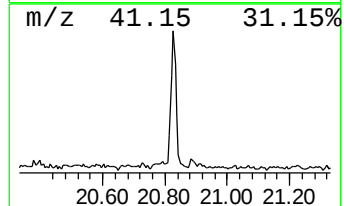
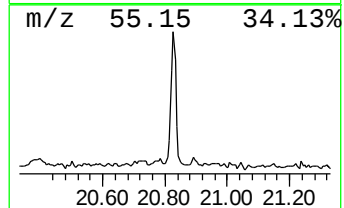
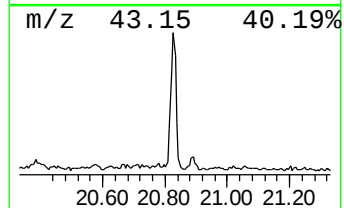
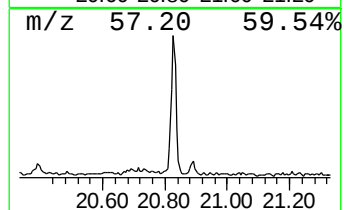
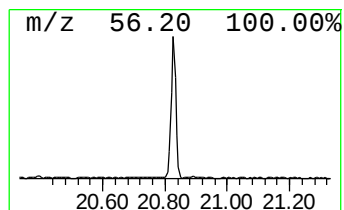
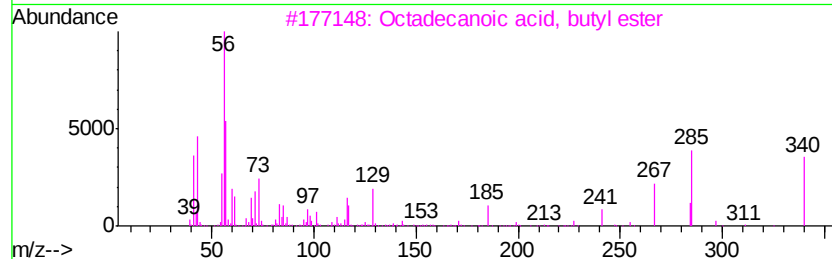
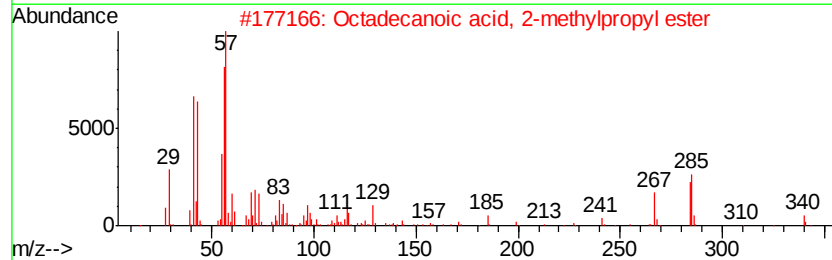
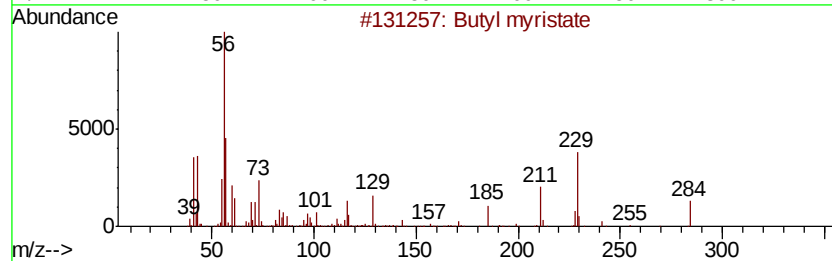
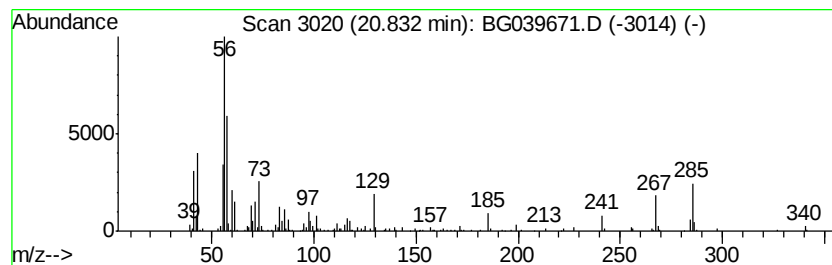
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Butyl myristate Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
20.83	6.01 ng	272686	Chrysene-d12	21.82	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butyl myristate	284	C18H36O2	000110-36-1	74
2	Octadecanoic acid, 2-methylpropyl ester	340	C22H44O2	000646-13-9	49
3	Octadecanoic acid, butyl ester	340	C22H44O2	000123-95-5	46
4	Cyclopentanecarboxylic acid, 2-allyl ester	129	C6H11NO2	037910-65-9	43
5	Cyclopentanecarboxylic acid, 2-allyl ester	129	C6H11NO2	040482-05-1	43



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG022819\
Data File : BG039671.D
Acq On : 28 Feb 2019 22:14
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Sample : K1628-04
Misc :
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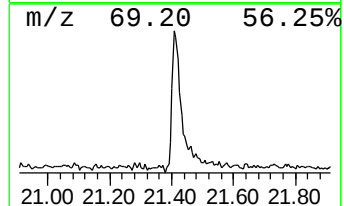
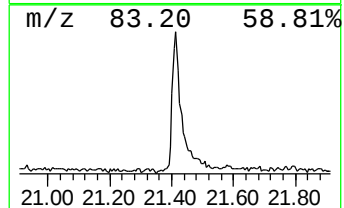
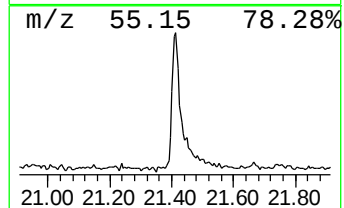
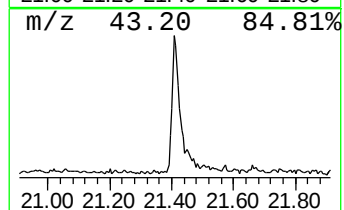
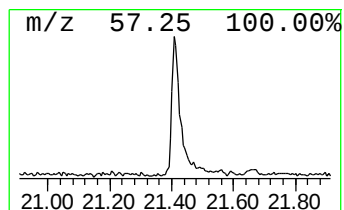
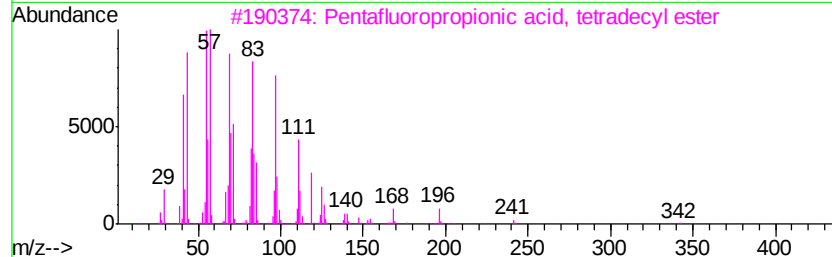
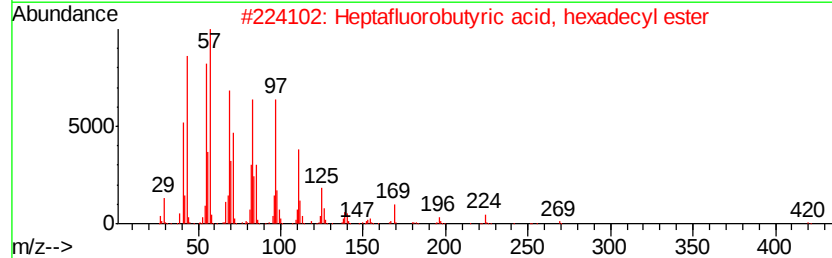
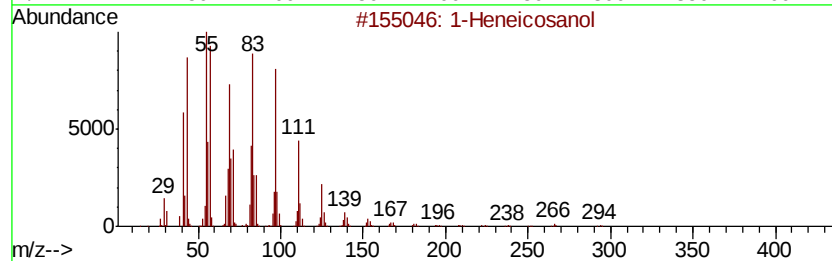
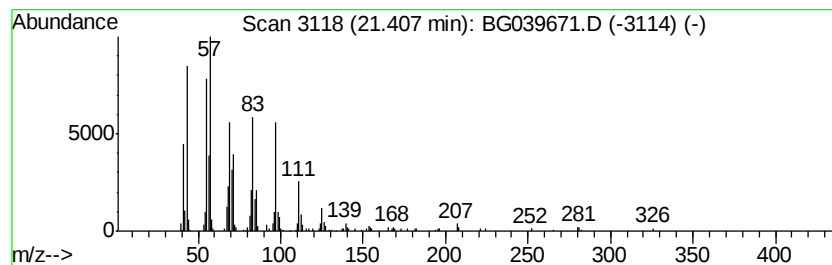
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 1-Heneicosanol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD		R.T.	
21.41	7.10 ng	322363	Chrysene-d12		21.82	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Heneicosanol	312	C21H44O	015594-90-8	93
2		Heptafluorobutyric acid, hexadec...	438	C20H33F7O2	006385-15-5	91
3		Pentafluoropropionic acid, tetra...	360	C17H29F5O2	006222-06-6	91
4		Heptacosyl acetate	438	C29H58O2	1000351-78-2	91
5		Octadecyl trifluoroacetate	366	C20H37F3O2	079392-43-1	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG022819\
Data File : BG039671.D
Acq On : 28 Feb 2019 22:14
Operator : JU/SJ
Sample : K1628-04
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
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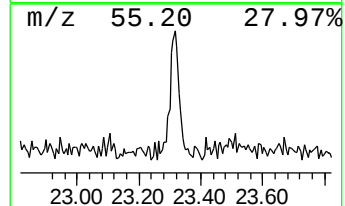
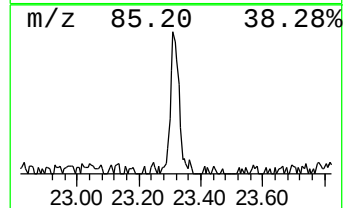
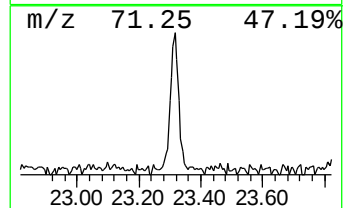
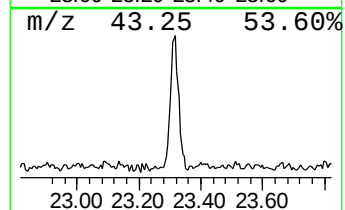
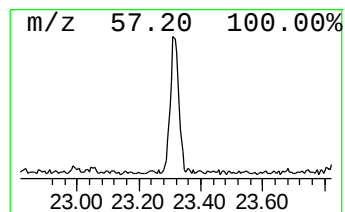
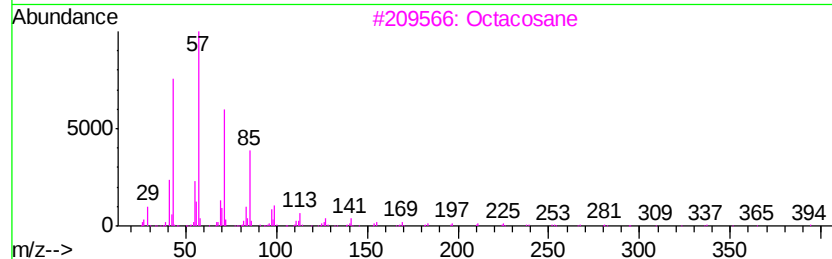
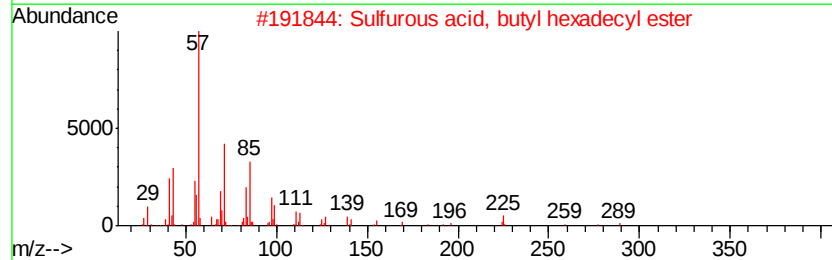
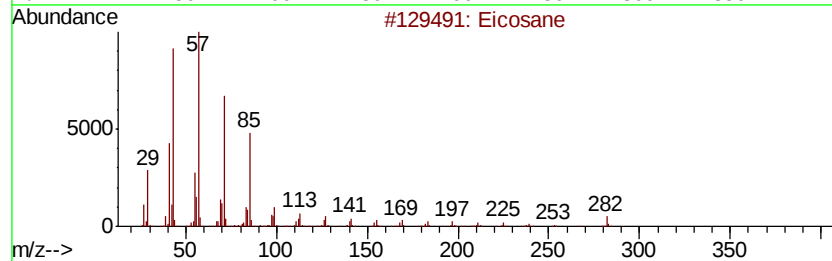
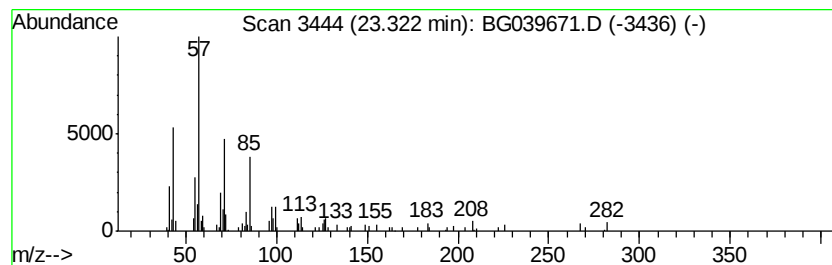
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Eicosane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.32	2.24 ng	101561	Chrysene-d12	21.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	94
2		Sulfurous acid, butyl hexadecyl ...	362	C20H42O3S	1000309-18-3	86
3		Octacosane	394	C28H58	000630-02-4	83
4		Tritetracontane	605	C43H88	007098-21-7	80
5		Sulfurous acid, butyl tetradecyl...	334	C18H38O3S	1000309-18-1	80



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG022819\
Data File : BG039671.D
Acq On : 28 Feb 2019 22:14
Operator : JU/SJ
Sample : K1628-04
Misc :
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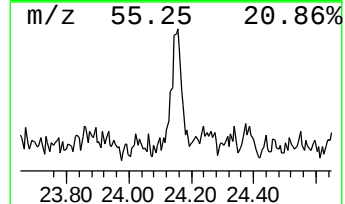
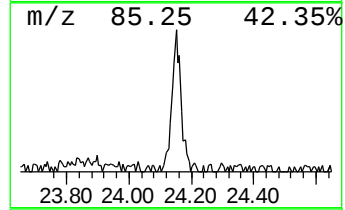
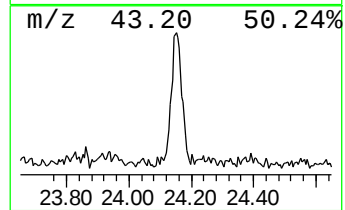
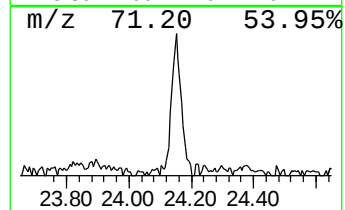
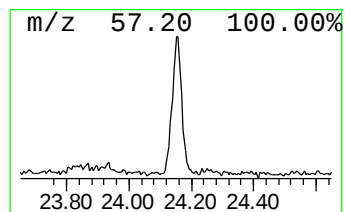
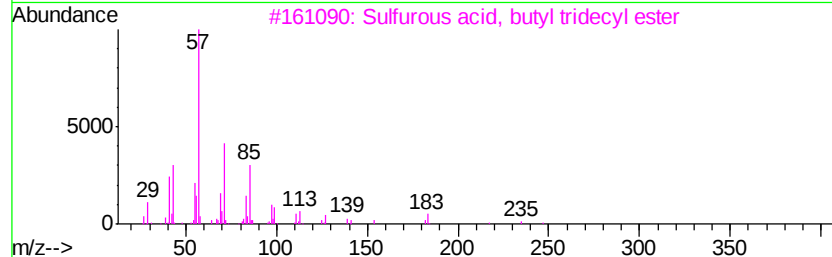
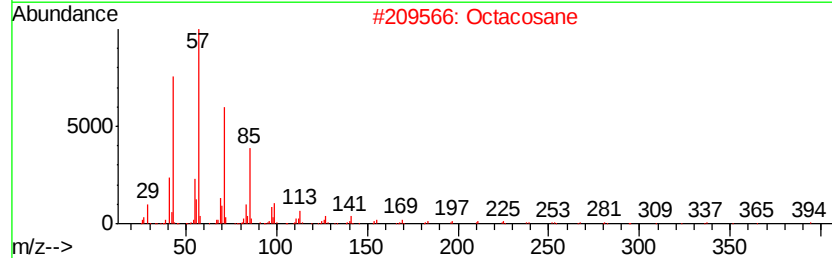
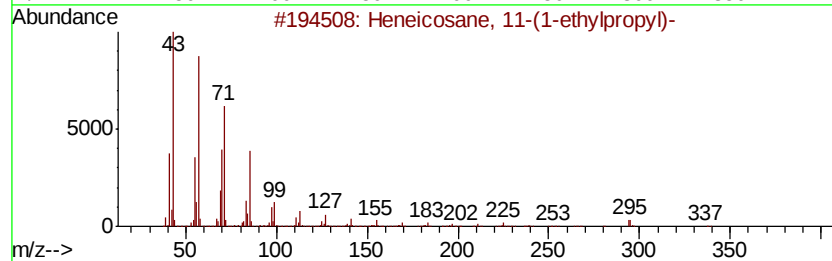
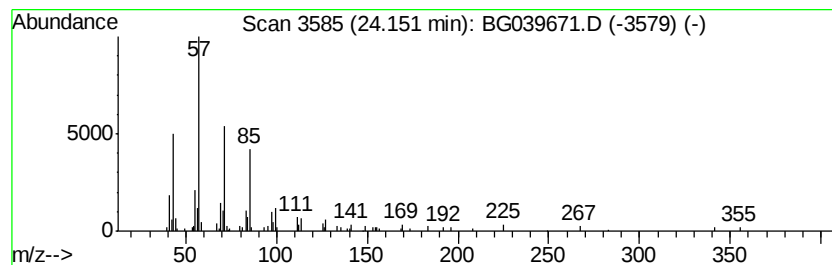
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Heneicosane, 11-(1-ethylpro... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.15	2.06 ng	98684	Perylene-d12	25.20
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Heneicosane, 11-(1-ethylpropyl)-	366 C26H54	055282-11-6	87
2	Octacosane	394 C28H58	000630-02-4	87
3	Sulfurous acid, butyl tridecyl e...	320 C17H36O3S	1000309-18-0	86
4	Tetracosane	338 C24H50	000646-31-1	86
5	Hexacosane	366 C26H54	000630-01-3	81



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG022819\
 Data File : BG039671.D
 Acq On : 28 Feb 2019 22:14
 Operator : JU/SJ
 Sample : K1628-04
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT 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
Butane, 2-methoxy...	3.29	4.9 ng		76501	1	8.17	310718	20.0
unknown7.69	7.69	117.6 ng		1827110	1	8.17	310718	20.0
Hexadecanoic acid...	18.16	3.1 ng		127043	4	17.53	824908	20.0
n-Hexadecanoic acid	18.38	6.3 ng		259854	4	17.53	824908	20.0
Methyl stearate	19.45	2.5 ng		103751	4	17.53	824908	20.0
unknown19.53	19.53	2.5 ng		104202	4	17.53	824908	20.0
Hexadecanoic acid...	19.79	3.9 ng		175288	5	21.82	907543	20.0
Butyl myristate	20.83	6.0 ng		272686	5	21.82	907543	20.0
1-Heneicosanol	21.41	7.1 ng		322363	5	21.82	907543	20.0
Eicosane	23.32	2.2 ng		101561	5	21.82	907543	20.0
Heneicosane, 11-(...	24.15	2.1 ng		98684	6	25.20	956638	20.0