

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG022819\
 Data File : BG039672.D
 Acq On : 28 Feb 2019 22:53
 Operator : JU/SJ
 Sample : K1643-01
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 INTERIOR-WALL

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	3.219	15	22	31	rBV	76927	159262	5.98%	0.736%
2	3.295	31	35	45	rVB	34892	55236	2.08%	0.255%
3	5.709	439	446	456	rBV	779474	1244379	46.76%	5.753%
4	7.143	683	690	697	rBV3	13063	33427	1.26%	0.155%
5	7.307	710	718	732	rBV	820645	1434814	53.92%	6.633%
6	7.695	776	784	801	rVB	752757	1343142	50.47%	6.209%
7	8.165	857	864	873	rBV	179778	306502	11.52%	1.417%
8	8.488	911	919	928	rVB	572445	977581	36.73%	4.519%
9	9.340	1056	1064	1073	rBV	468100	848299	31.88%	3.922%
10	9.428	1073	1079	1097	rVB2	32373	83874	3.15%	0.388%
11	10.268	1215	1222	1234	rBV4	29127	70034	2.63%	0.324%
12	10.985	1337	1344	1352	rBV	258522	454373	17.07%	2.100%
13	11.748	1467	1474	1486	rBV4	40038	89935	3.38%	0.416%
14	12.753	1638	1645	1653	rBV2	33022	71900	2.70%	0.332%
15	13.411	1747	1757	1766	rBV	1357567	2095141	78.73%	9.686%
16	14.227	1891	1896	1903	rBV2	46548	70452	2.65%	0.326%
17	14.691	1970	1975	1982	rVB	20481	32542	1.22%	0.150%
18	14.791	1982	1992	2005	rBV3	394319	696439	26.17%	3.220%
19	15.161	2048	2055	2058	rBV5	17561	34799	1.31%	0.161%
20	15.578	2122	2126	2132	rVB	102892	146714	5.51%	0.678%
21	15.702	2141	2147	2151	rBV4	17185	31753	1.19%	0.147%
22	16.272	2237	2244	2258	rBV	998078	1585166	59.57%	7.328%
23	16.894	2344	2350	2357	rBV5	42222	87961	3.31%	0.407%
24	17.535	2452	2459	2470	rBV2	505174	811376	30.49%	3.751%
25	18.028	2540	2543	2549	rVB3	24831	39904	1.50%	0.184%
26	18.398	2599	2606	2614	rBV	1159632	1882210	70.73%	8.701%
27	18.469	2614	2618	2623	rVB	218634	315782	11.87%	1.460%
28	19.291	2755	2758	2764	rVB2	60054	75344	2.83%	0.348%
29	19.438	2780	2783	2788	rVB2	50550	61768	2.32%	0.286%
30	19.532	2796	2799	2805	rBV4	86155	136540	5.13%	0.631%
31	19.649	2815	2819	2832	rBV2	379554	608405	22.86%	2.813%
32	19.867	2852	2856	2859	rVB	80919	95298	3.58%	0.441%
33	20.125	2894	2900	2906	rBV	2070716	2661213	100.00%	12.302%
34	20.395	2943	2946	2950	rVB2	75819	79738	3.00%	0.369%

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Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % 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

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35	20.889	3028	3030	3034	rVB2	64778	65748	2.47%	0.304%
36	21.412	3115	3119	3131	rVB	195285	349267	13.12%	1.615%
37	21.664	3158	3162	3166	rBV	127580	163755	6.15%	0.757%
38	21.823	3183	3189	3196	rVB	486923	850102	31.94%	3.930%
39	21.970	3211	3214	3220	rVB3	42638	54427	2.05%	0.252%
40	22.598	3317	3321	3328	rVB	43826	62694	2.36%	0.290%
41	23.315	3438	3443	3452	rVB2	64619	122603	4.61%	0.567%
42	24.155	3580	3586	3594	rVB2	58597	130734	4.91%	0.604%
43	25.201	3750	3764	3775	rVB3	285019	983058	36.94%	4.545%
44	26.328	3950	3956	3966	rVB3	35915	98215	3.69%	0.454%
45	26.487	3979	3983	3987	rBV5	14737	29762	1.12%	0.138%

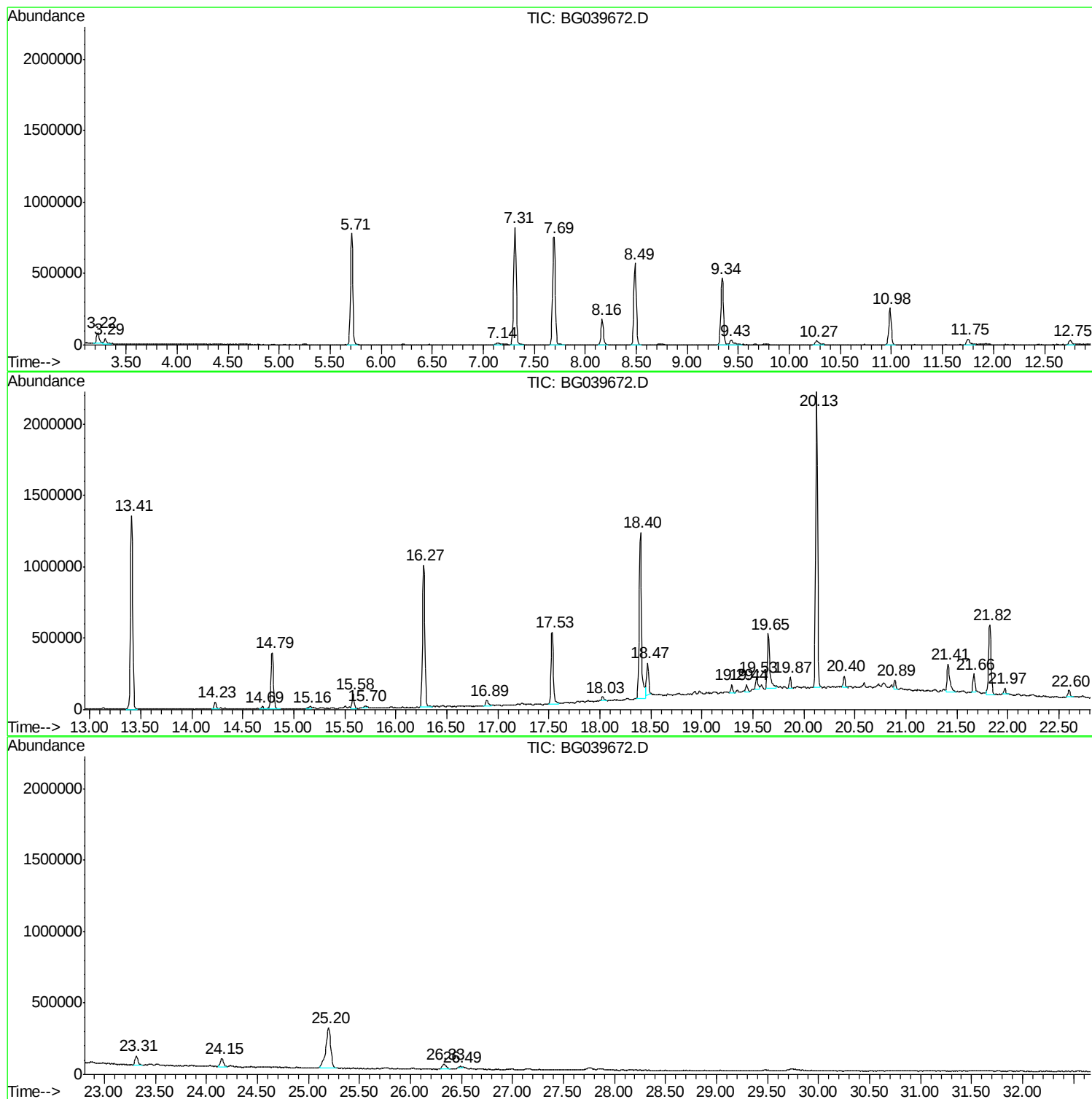
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG012919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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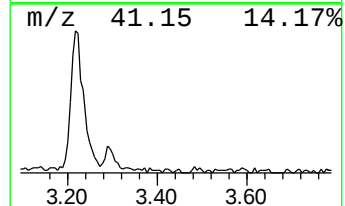
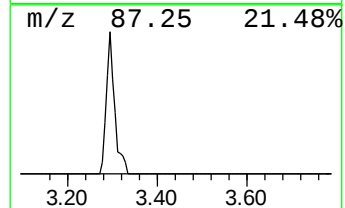
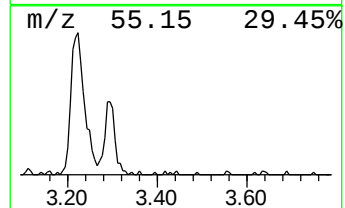
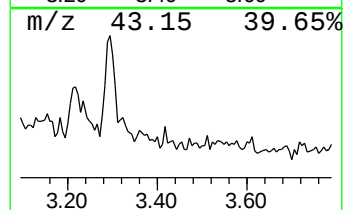
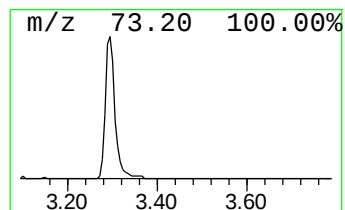
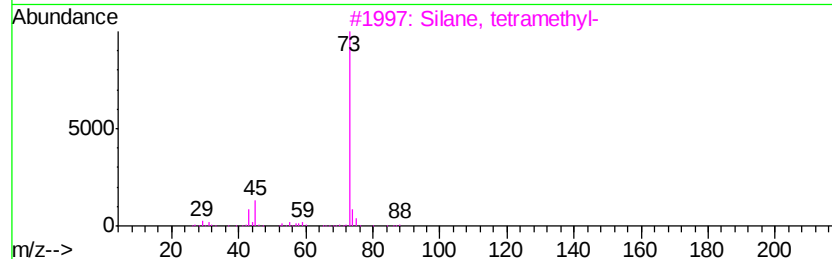
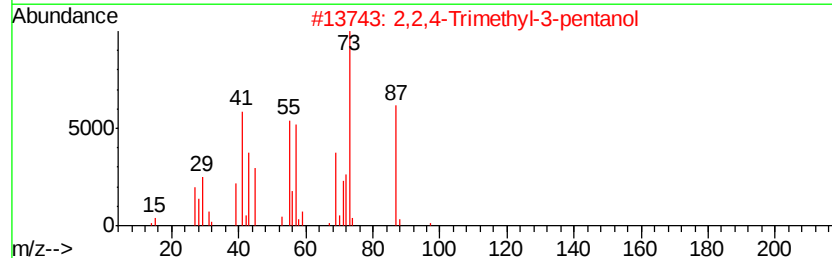
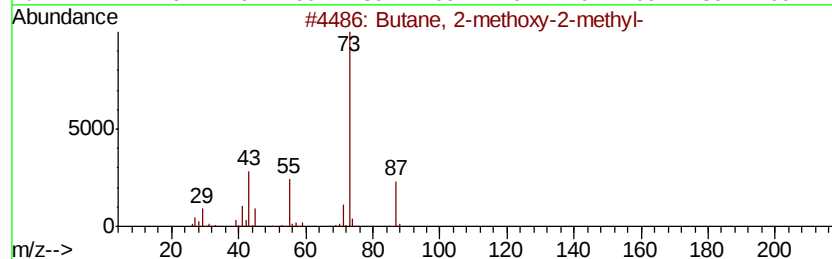
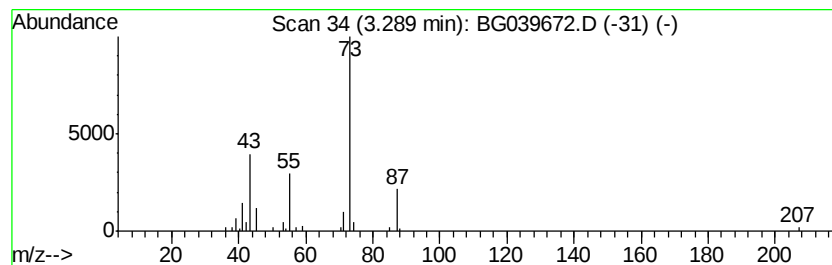
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.29	3.60 ng	55236	1,4-Dichlorobenzene-d4	8.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	64
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	38
3		Silane, tetramethyl-	88	C4H12Si	000075-76-3	35
4		2-Methyl-5-hexen-3-ol	114	C7H14O	032815-70-6	25
5		Boronic acid, ethyl-, dimethyl e...	102	C4H11BO2	007318-82-3	9



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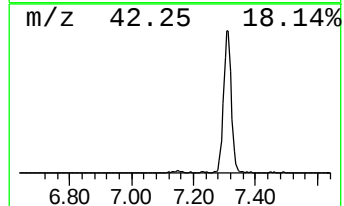
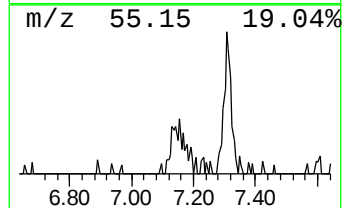
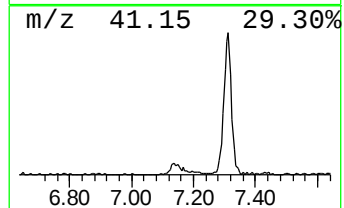
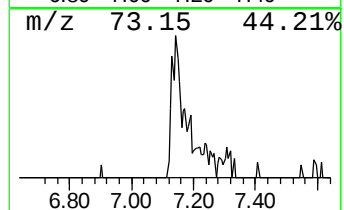
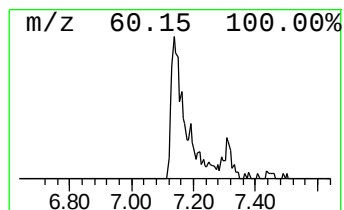
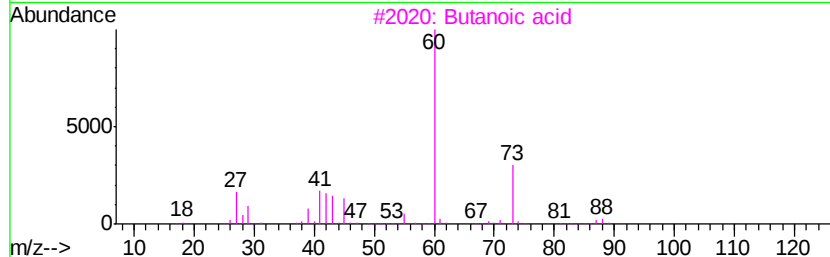
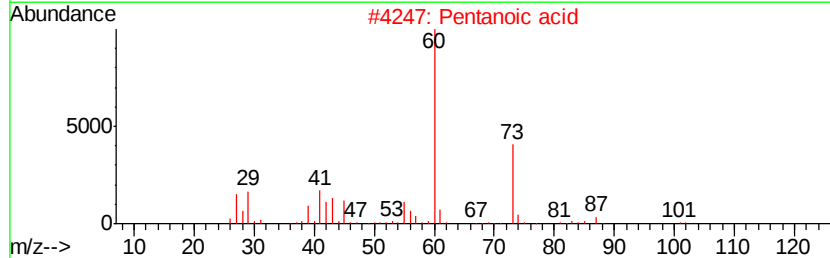
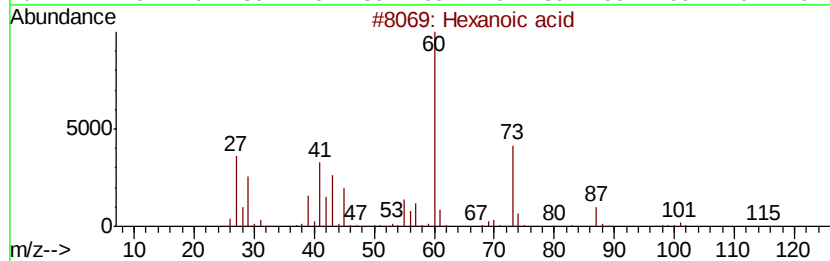
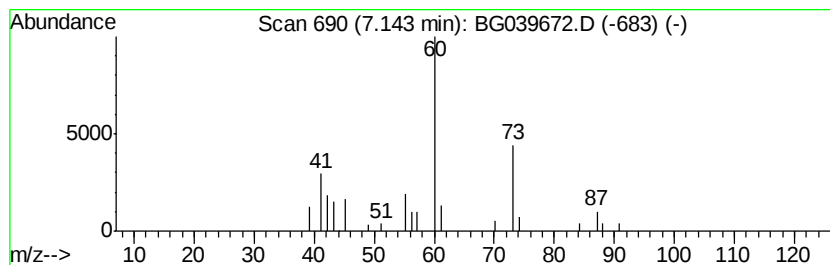
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Hexanoic acid Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.14	2.18 ng	33427	1,4-Dichlorobenzene-d4	8.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexanoic acid	116	C6H12O2	000142-62-1	78
2		Pentanoic acid	102	C5H10O2	000109-52-4	56
3		Butanoic acid	88	C4H8O2	000107-92-6	53
4		Propanedioic acid, propyl-	146	C6H10O4	000616-62-6	50
5		Butanoic acid, 4-chloro-	122	C4H7ClO2	000627-00-9	9



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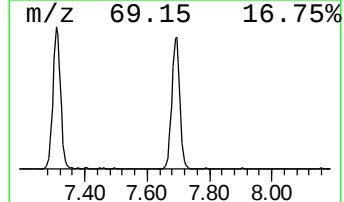
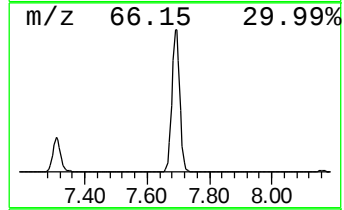
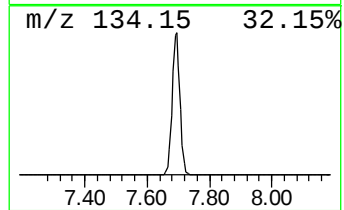
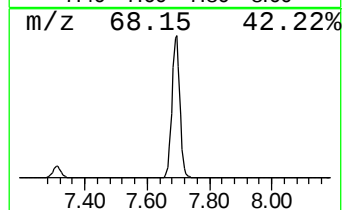
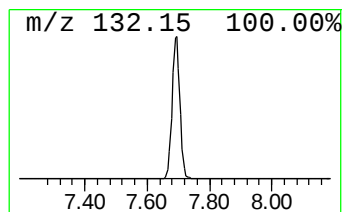
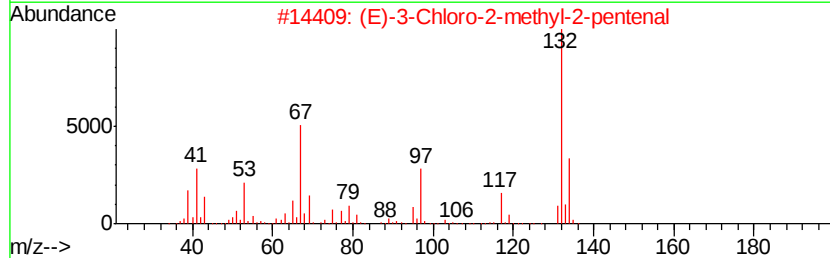
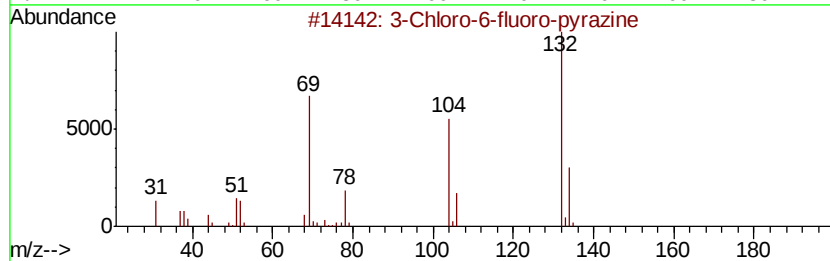
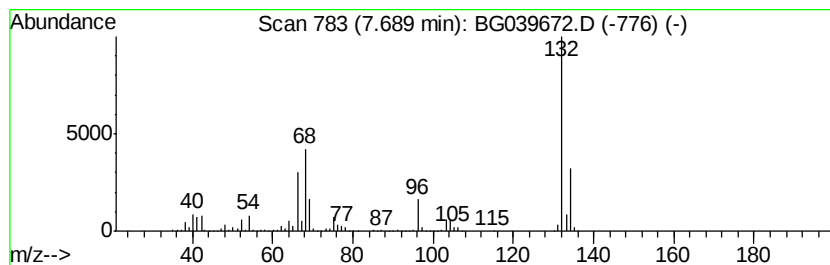
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG012919.M
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 unknown7.69 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.69	87.64 ng	1343140	1,4-Dichlorobenzene-d4	8.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
3		Tranvlcyproline-propionyl	189	C12H15NO	1000123-86-3	12
4		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	12
5		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12



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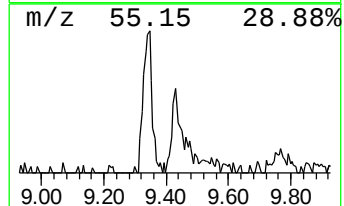
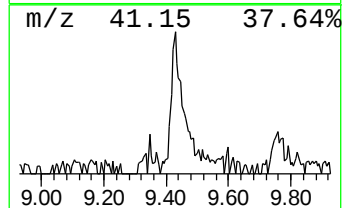
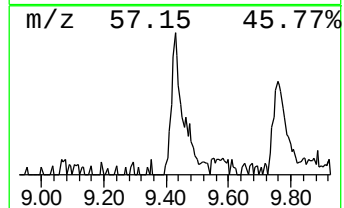
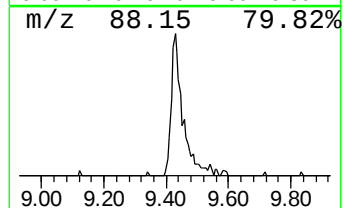
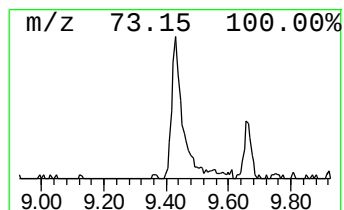
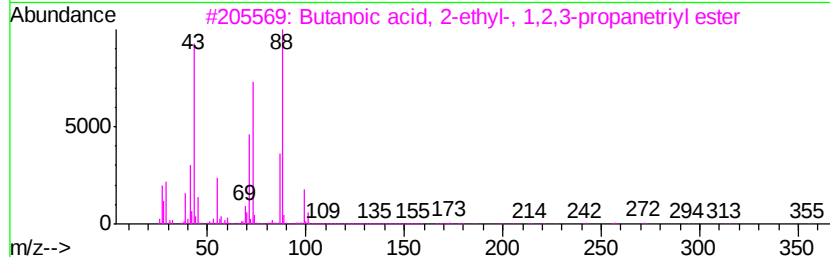
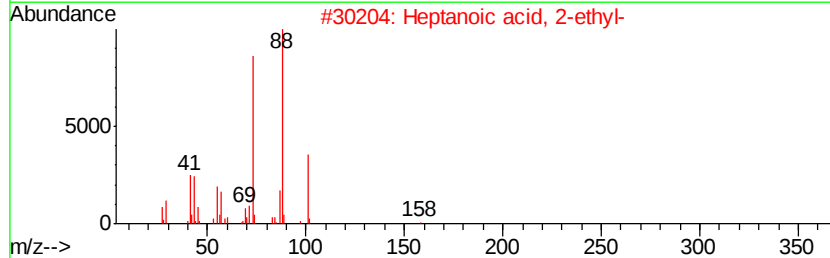
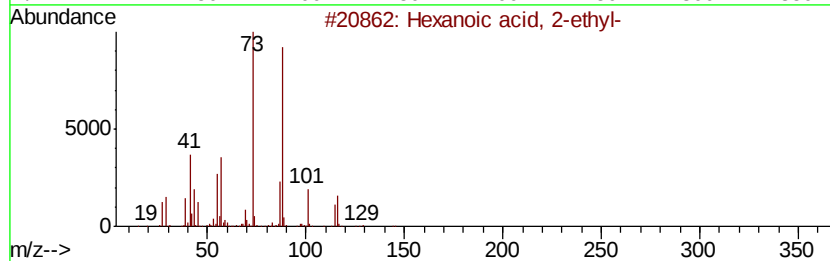
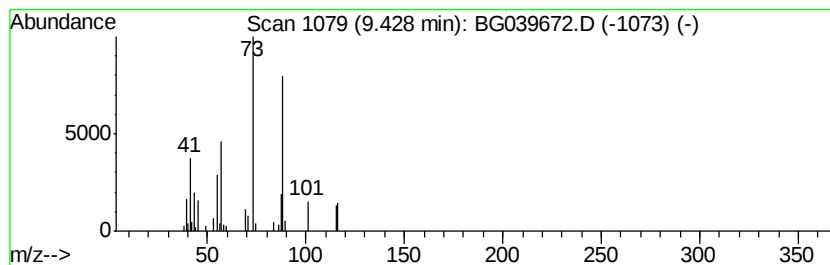
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Hexanoic acid, 2-ethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.43	5.47 ng	83874	1,4-Dichlorobenzene-d4	8.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	80
2		Heptanoic acid, 2-ethyl-	158	C9H18O2	003274-29-1	45
3		Butanoic acid, 2-ethyl-, 1,2,3-p...	386	C21H38O6	056554-54-2	40
4		Heptanoic acid, ethyl ester	158	C9H18O2	000106-30-9	38
5		1,4-Dioxane	88	C4H8O2	000123-91-1	38



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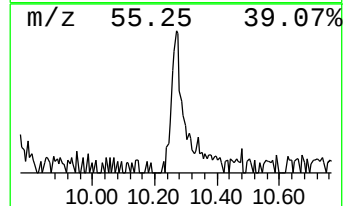
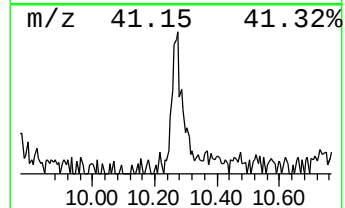
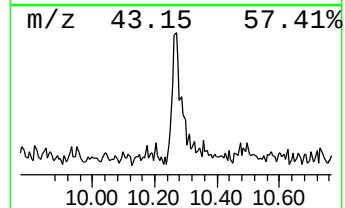
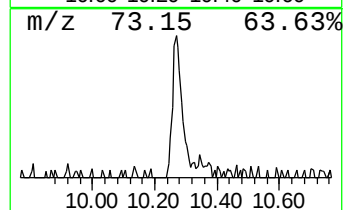
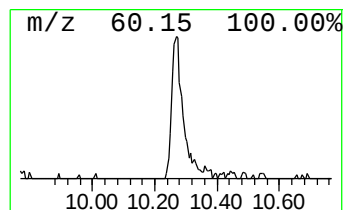
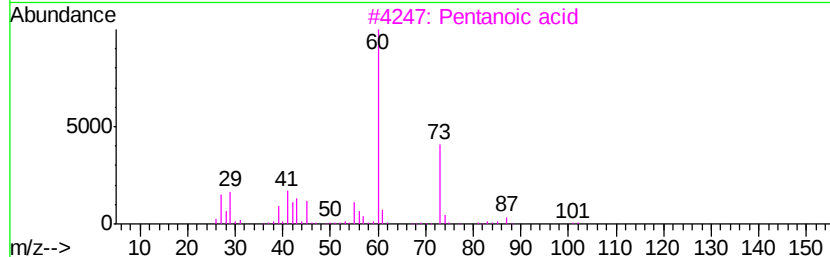
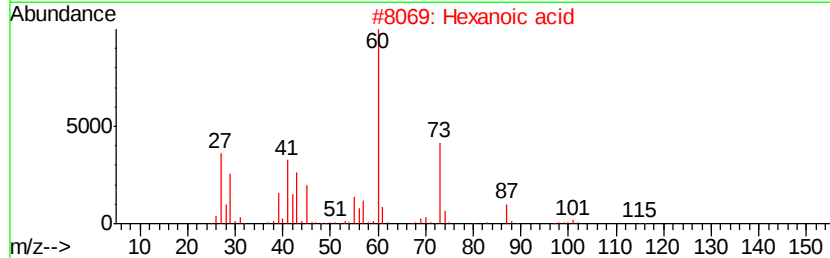
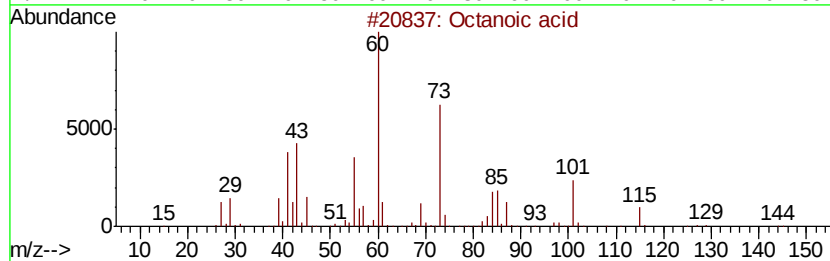
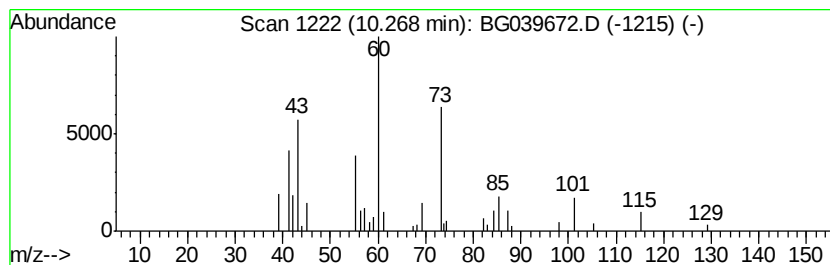
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Octanoic acid Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.27	3.08 ng	70034	Naphthalene-d8	10.98

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octanoic acid		144	C8H16O2	000124-07-2	78
2	Hexanoic acid		116	C6H12O2	000142-62-1	64
3	Pentanoic acid		102	C5H10O2	000109-52-4	59
4	Nonanoic acid		158	C9H18O2	000112-05-0	45
5	Octanoic acid, silver(1+) salt		250	C8H15AgO2	024927-67-1	37



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG022819\
Data File : BG039672.D
Acq On : 28 Feb 2019 22:53
Operator : JU/SJ
Sample : K1643-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
INTERIOR-WALL

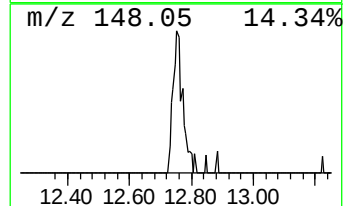
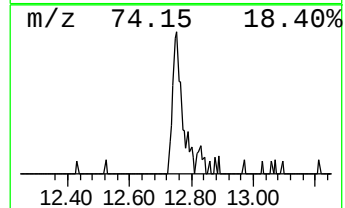
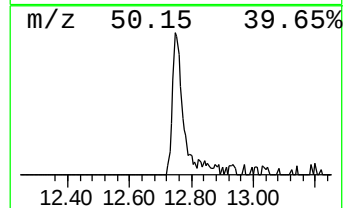
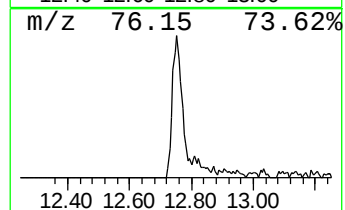
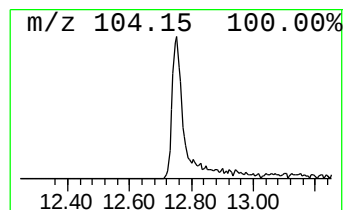
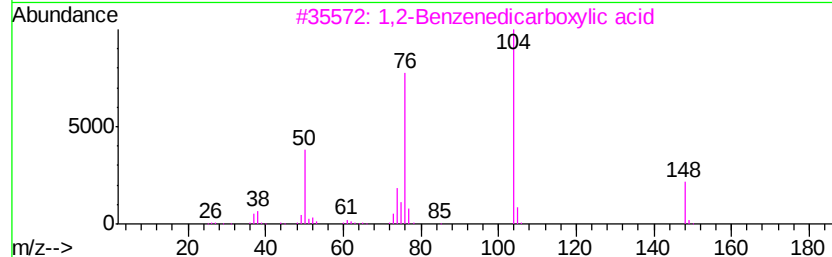
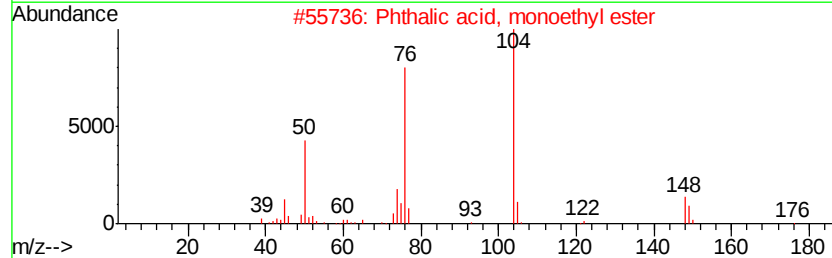
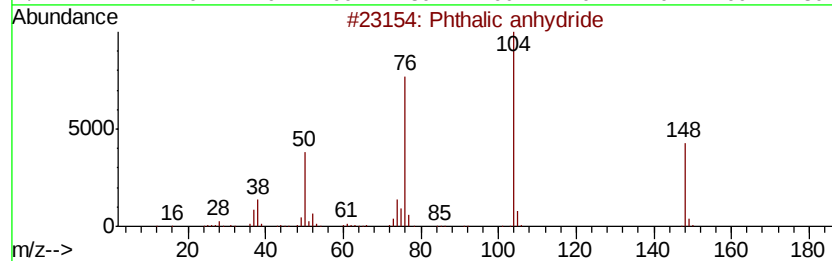
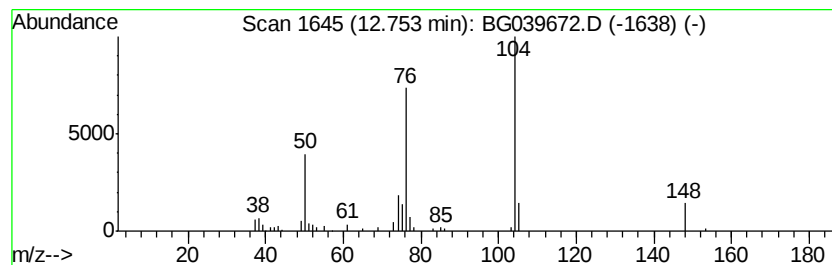
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Phthalic anhydride Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.75	3.16 ng	71900	Naphthalene-d8	10.98

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phthalic anhydride	148	C8H4O3	000085-44-9	91
2		Phthalic acid, monoethyl ester	194	C10H10O4	002306-33-4	90
3		1,2-Benzenedicarboxylic acid	166	C8H6O4	000088-99-3	83
4		Methyl hydrogen phthalate	180	C9H8O4	004376-18-5	83
5		1H-Isoindole-1,3(2H)-dione, 2-(h...	177	C9H7N03	000118-29-6	78



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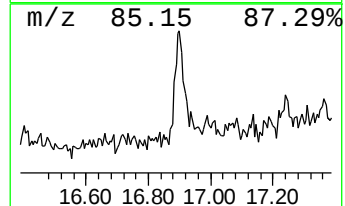
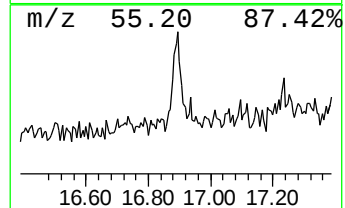
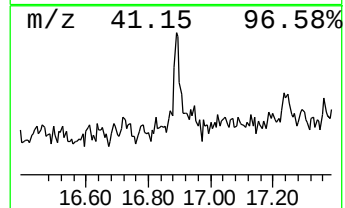
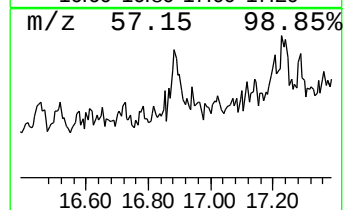
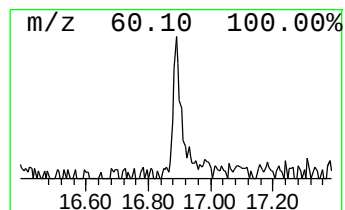
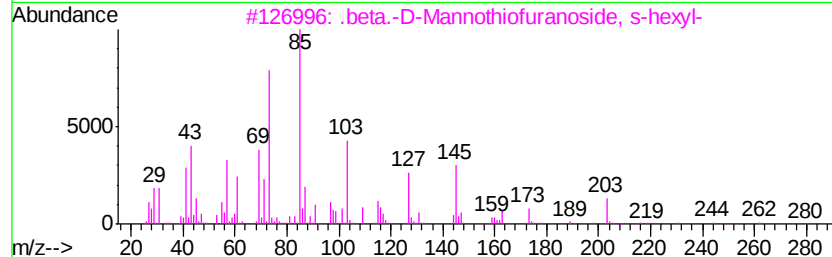
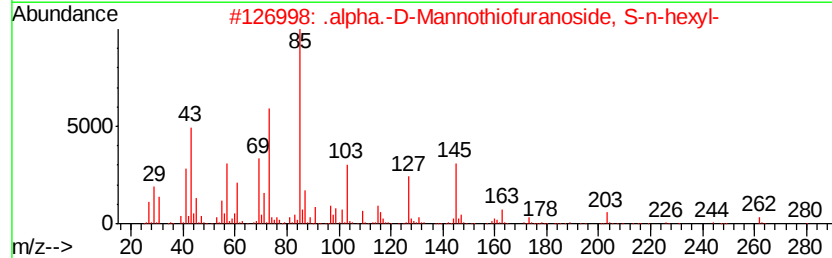
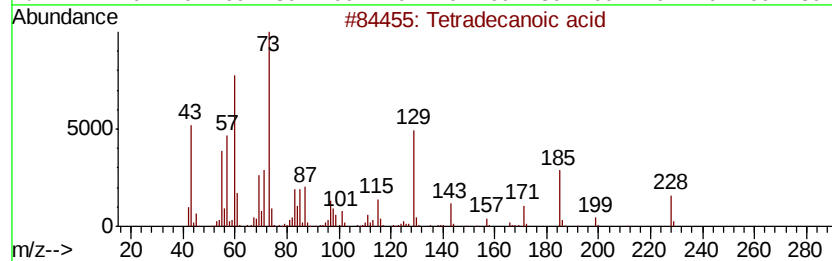
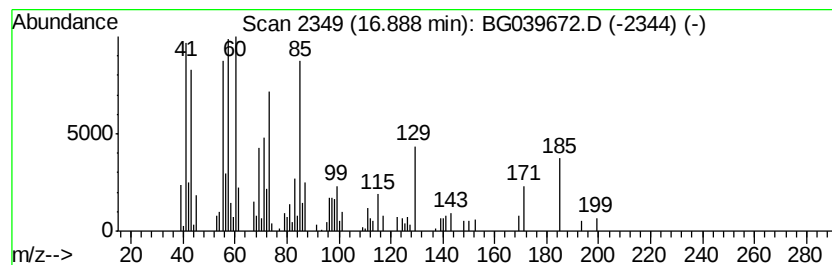
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Tetradecanoic acid Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.		
16.89	2.17 ng	87961	Phenanthrene-d10	17.53		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetradecanoic acid	228	C14H28O2	000544-63-8	87
2		.alpha.-D-Mannothiofuranoside, S...	280	C12H24O5S	1000159-73-9	46
3		.beta.-D-Mannothiofuranoside, s-...	280	C12H24O5S	1000158-98-0	43
4		Pentanoic acid, 2-octyl ester	214	C13H26O2	006938-48-3	38
5		(4S,5S)-(-)-2-Methyl-5-phenyl-2-...	191	C11H13NO2	053732-41-5	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG022819\
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Instrument :
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ClientSampleId :
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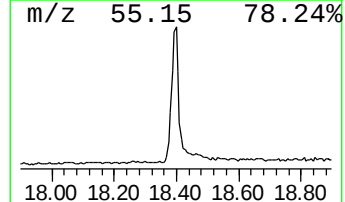
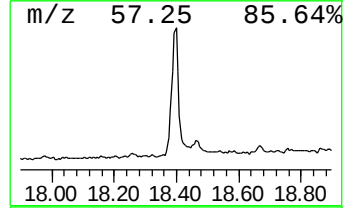
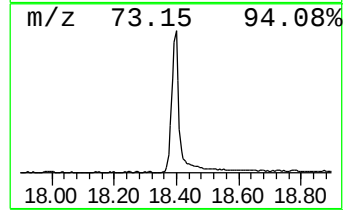
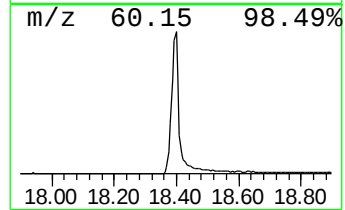
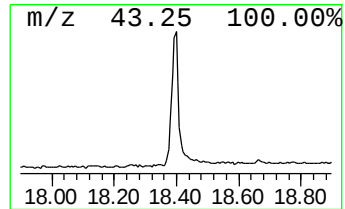
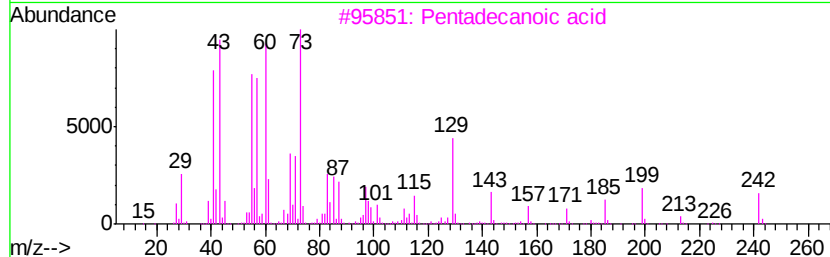
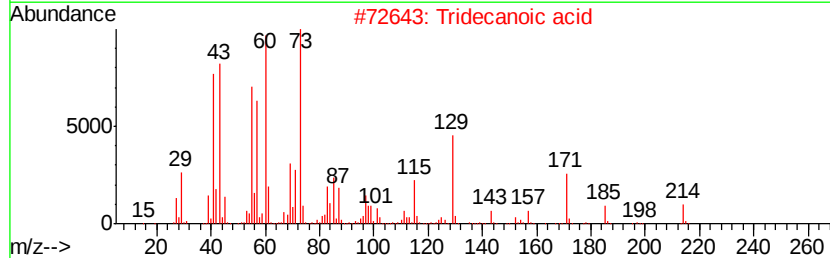
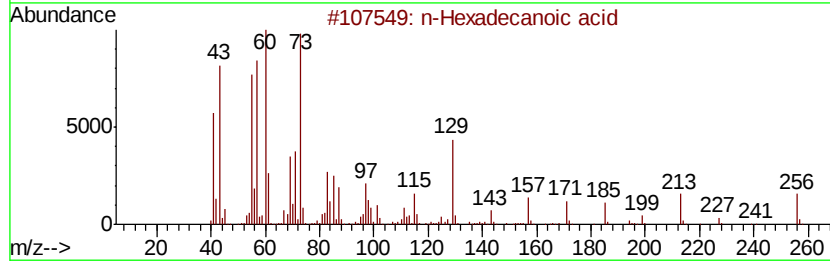
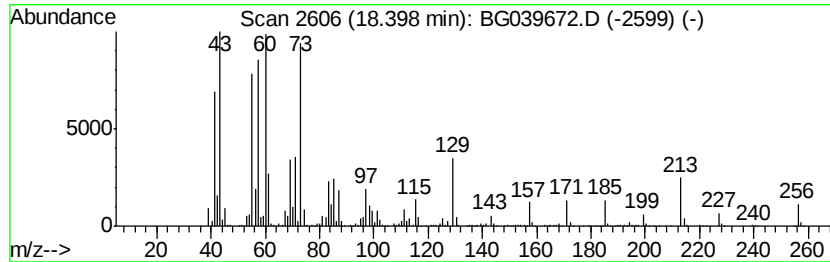
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
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Peak Number 11 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.40	46.40 ng	1882210	Phenanthrene-d10	17.53

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		Tridecanoic acid	214	C13H26O2	000638-53-9	89
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	81
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	72
5		n-Decanoic acid	172	C10H20O2	000334-48-5	47



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG022819\
Data File : BG039672.D
Acq On : 28 Feb 2019 22:53
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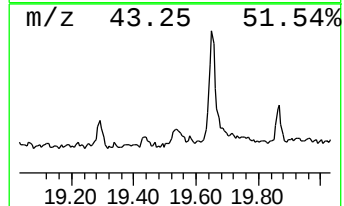
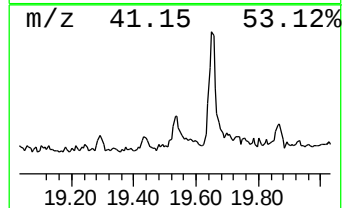
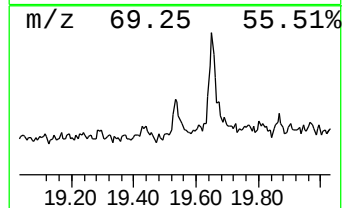
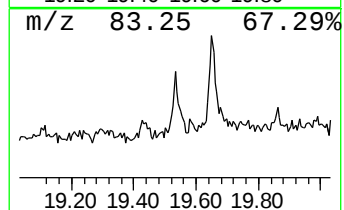
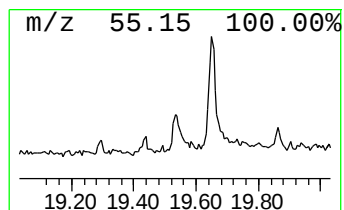
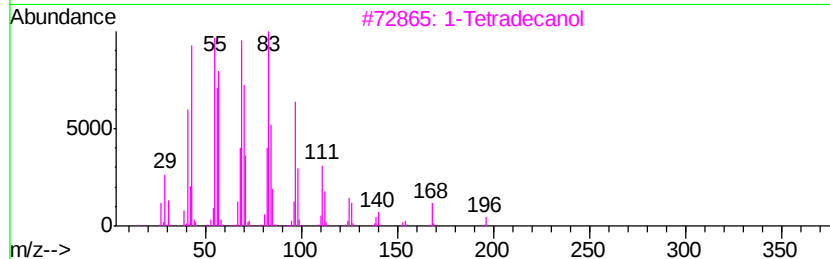
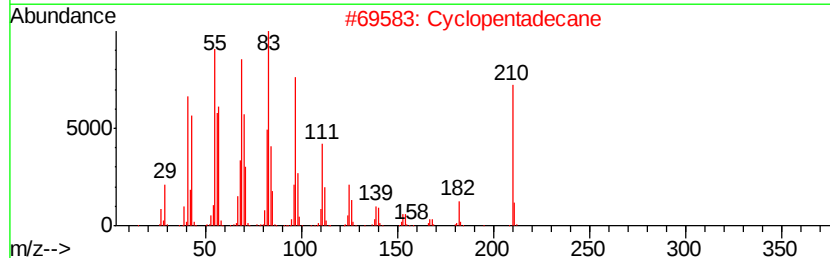
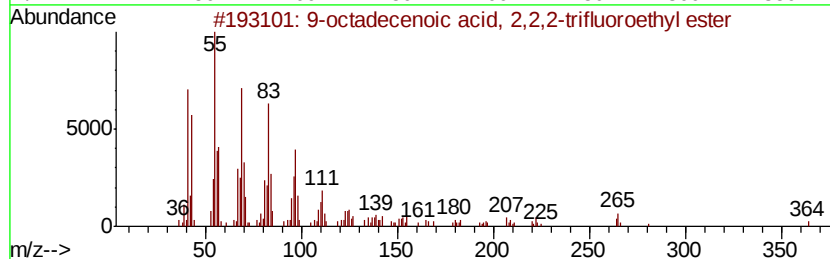
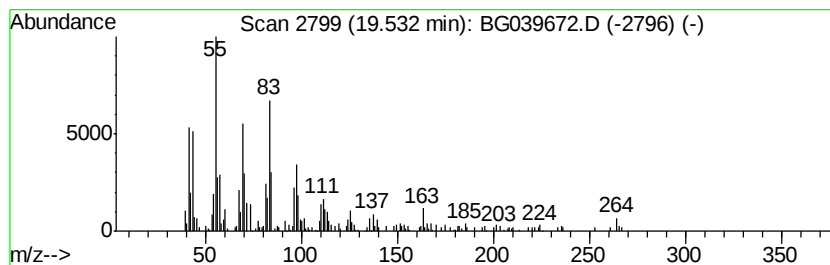
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 9-octadecenoic acid, 2,2,2-... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.53	3.37 ng	136540	Phenanthrene-d10	17.53

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9-octadecenoic acid, 2,2,2-trifl...	364	C20H35F3O2	1000376-54-3	58
2		Cyclopentadecane	210	C15H30	000295-48-7	53
3		1-Tetradecanol	214	C14H30O	000112-72-1	52
4		Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	46
5		Cetene	224	C16H32	000629-73-2	46



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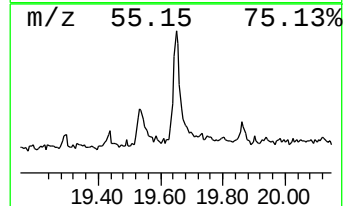
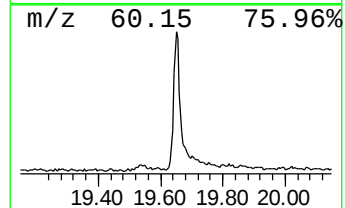
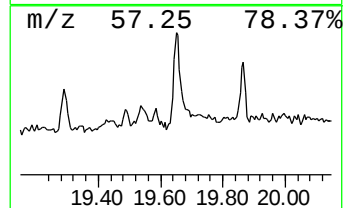
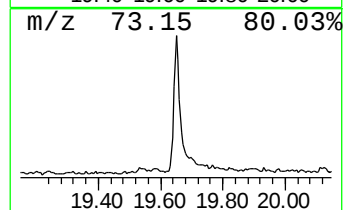
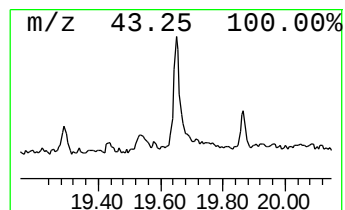
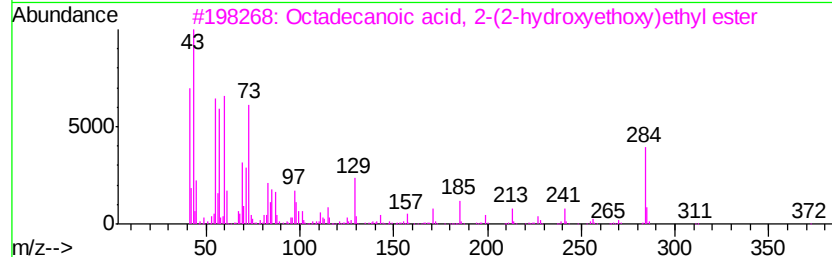
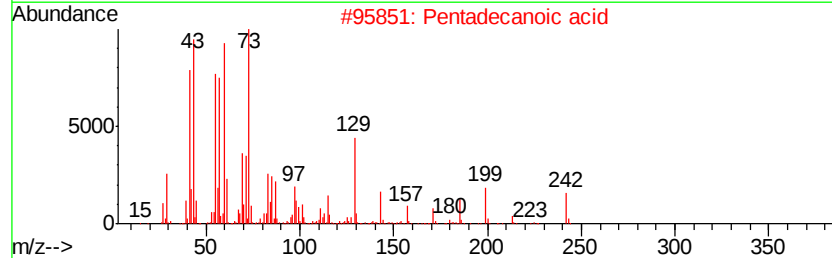
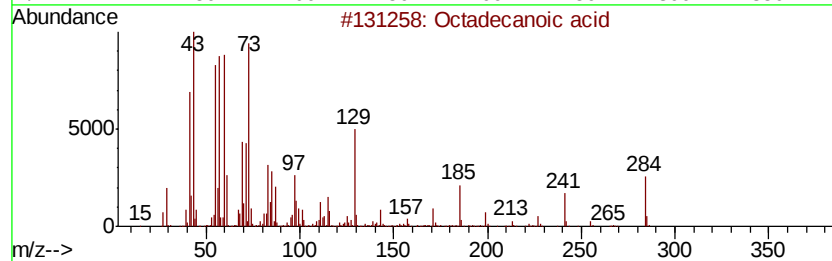
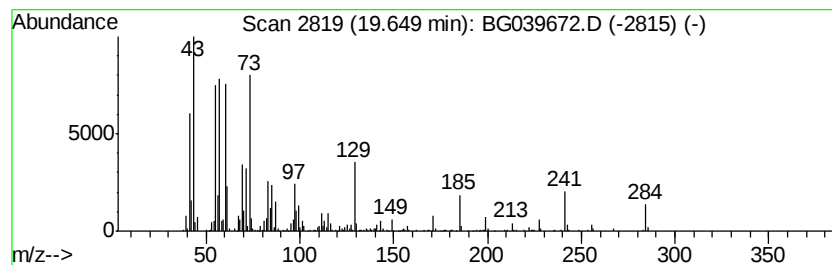
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TIC Library : C:\Database\NIST11.L
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Peak Number 13 Octadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.65	15.00 ng	608405	Phenanthrene-d10	17.53

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	86
3		Octadecanoic acid, 2-(2-hydroxyve...	372	C22H44O4	000106-11-6	80
4		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	70
5		Tetradecanoic acid	228	C14H28O2	000544-63-8	62



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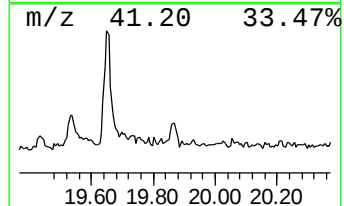
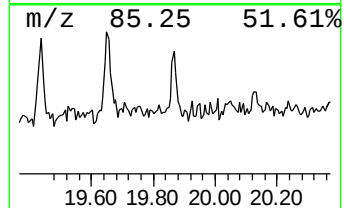
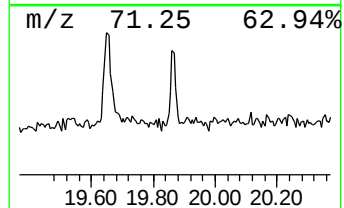
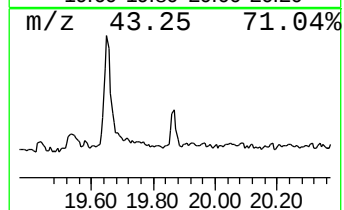
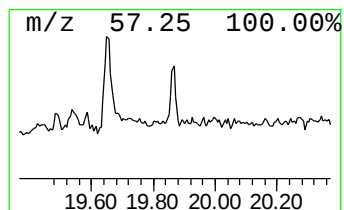
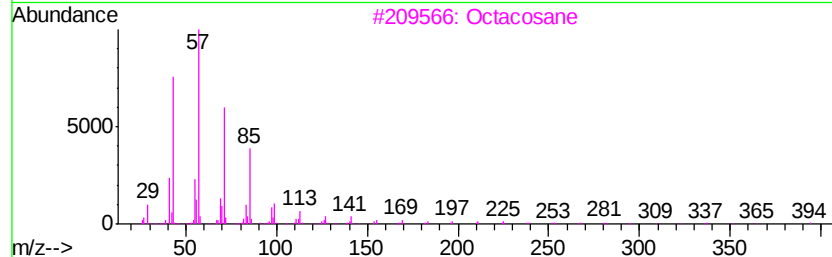
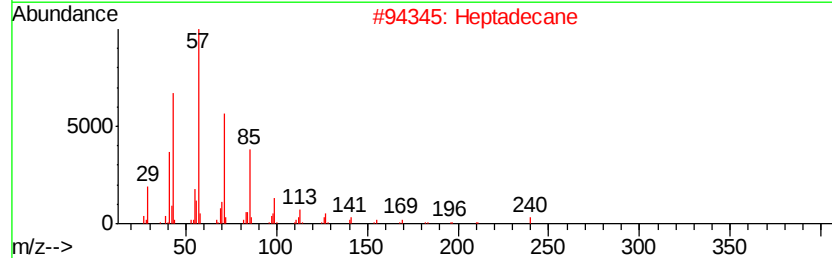
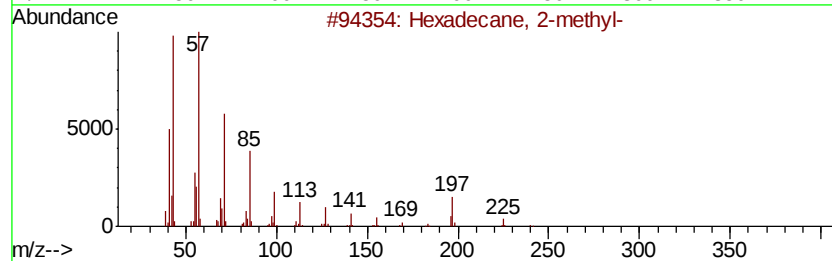
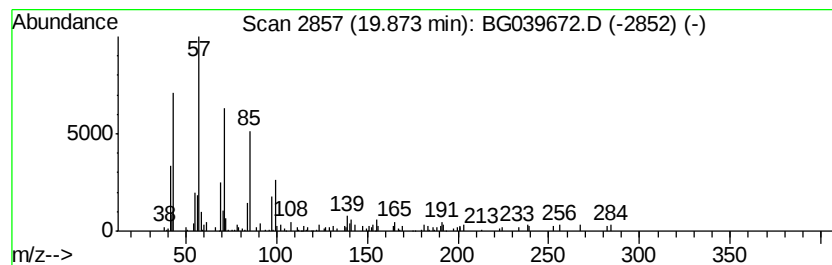
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 Hexadecane, 2-methyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.87	2.24 ng	95298	Chrysene-d12	21.82

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane, 2-methyl-	240	C17H36	001560-92-5	80
2			Heptadecane	240	C17H36	000629-78-7	80
3			Octacosane	394	C28H58	000630-02-4	72
4			Hentriacontane	437	C31H64	000630-04-6	72
5			3,5-Dimethyldodecane	198	C14H30	107770-99-0	72



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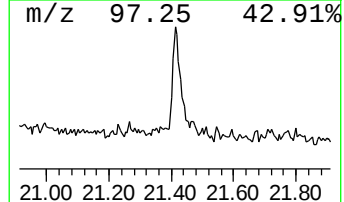
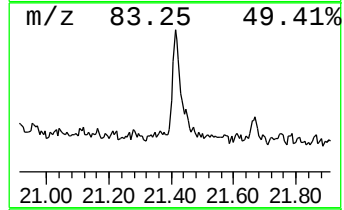
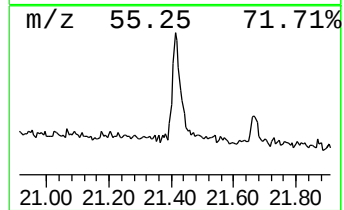
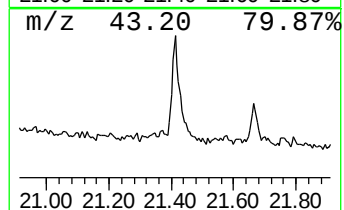
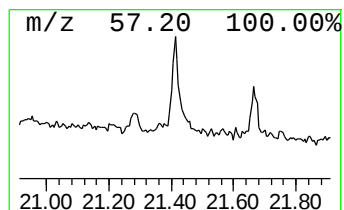
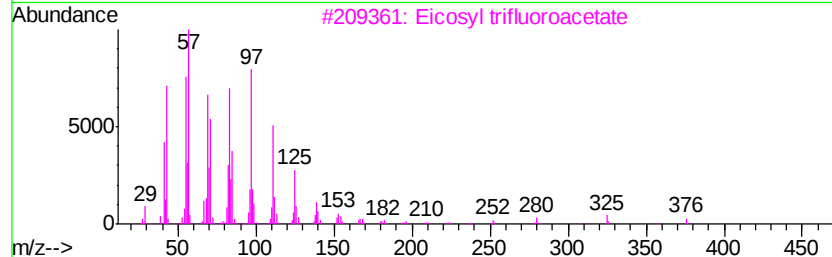
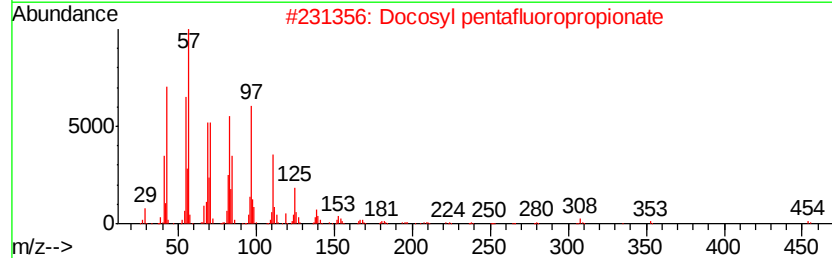
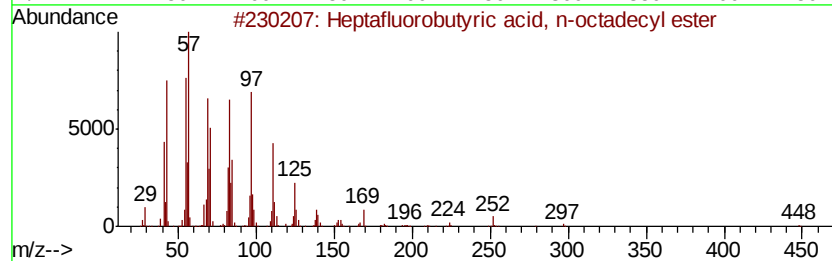
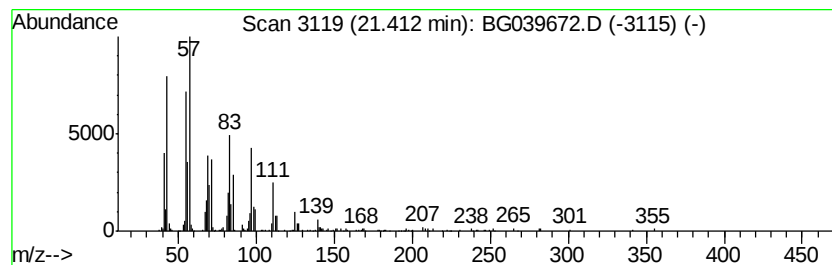
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Heptafluorobutyric acid, n-... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.41	8.22 ng	349267	Chrysene-d12	21.82

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptafluorobutyric acid, n-octad...	466	C22H37F7O2	000400-57-7	91
2			Docosyl pentafluoropropionate	472	C25H45F5O2	1000351-80-9	91
3			Eicosyl trifluoroacetate	394	C22H41F3O2	1000351-75-2	91
4			Heptadecyl heptafluorobutyrte	452	C21H35F7O2	959085-66-6	91
5			Pentafluoropropionic acid, octad...	416	C21H37F5O2	959261-25-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG022819\
Data File : BG039672.D
Acq On : 28 Feb 2019 22:53
Operator : JU/SJ
Sample : K1643-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
INTERIOR-WALL

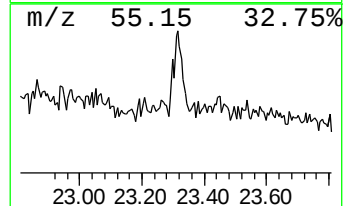
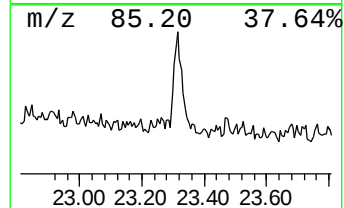
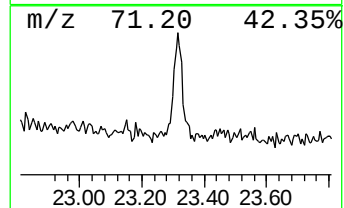
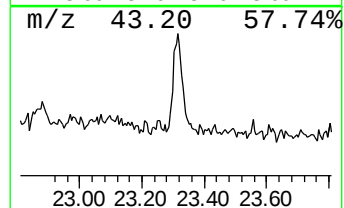
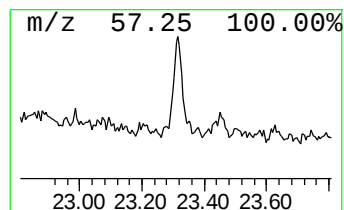
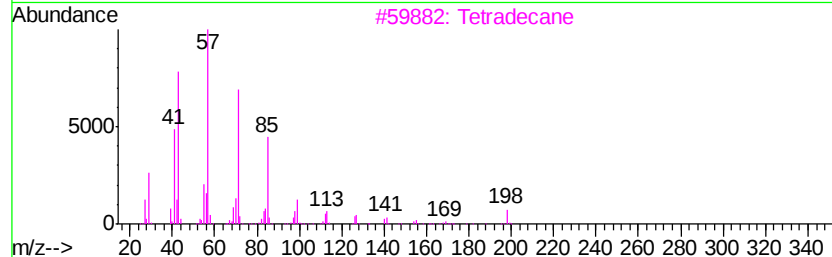
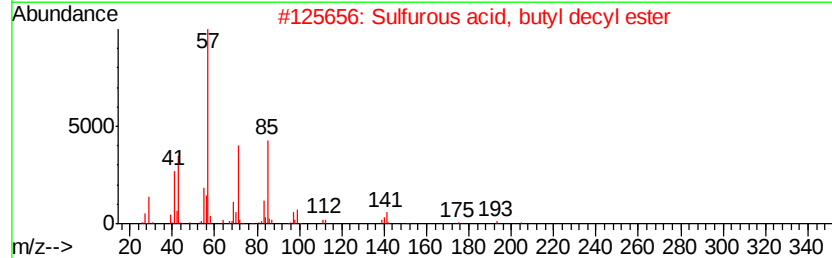
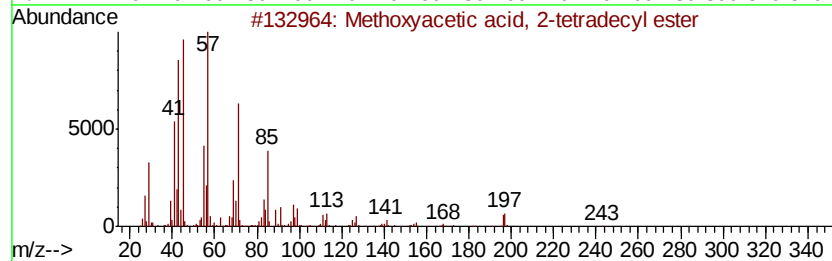
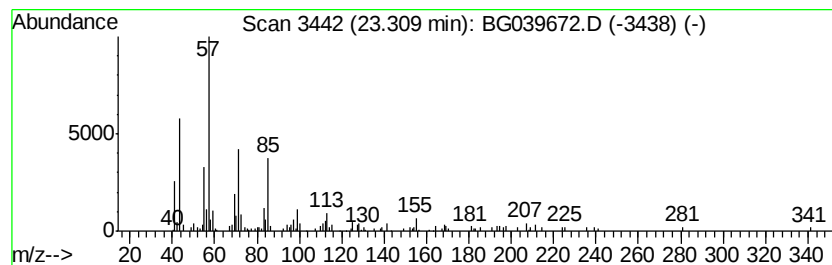
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Methoxyacetic acid, 2-tetra... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.31	2.88 ng	122603	Chrysene-d12	21.82

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methoxvacetic acid, 2-tetradecyl...	286	C17H34O3	1000282-04-8	72
2			Sulfurous acid, butyl decyl ester	278	C14H30O3S	1000309-17-7	72
3			Tetradecane	198	C14H30	000629-59-4	72
4			Hentriacontane	437	C31H64	000630-04-6	68
5			Hexadecane	226	C16H34	000544-76-3	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG022819\
Data File : BG039672.D
Acq On : 28 Feb 2019 22:53
Operator : JU/SJ
Sample : K1643-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
INTERIOR-WALL

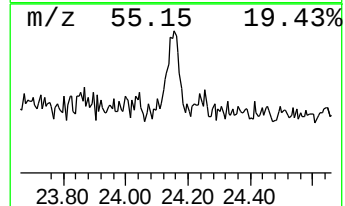
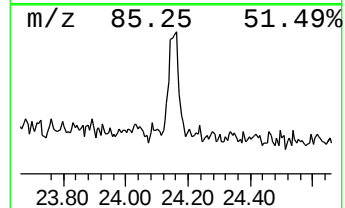
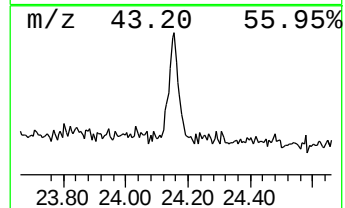
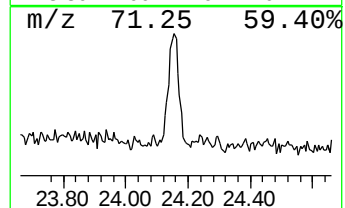
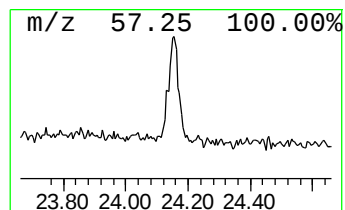
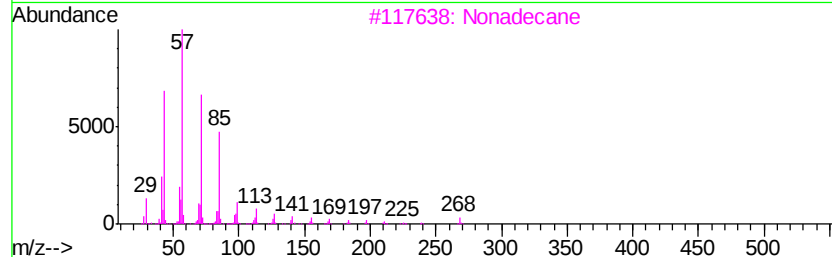
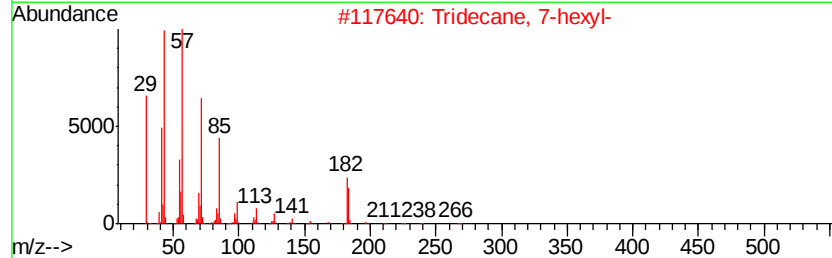
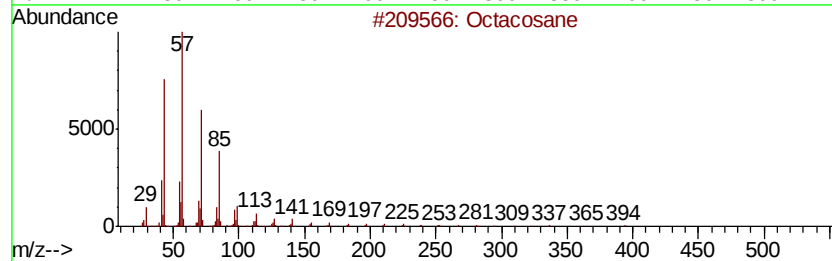
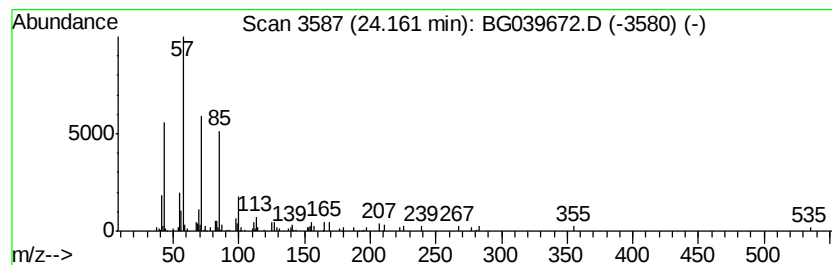
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Octacosane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.16	2.66 ng	130734	Perylene-d12	25.20

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosane	394	C28H58	000630-02-4	80
2		Tridecane, 7-hexyl-	268	C19H40	007225-66-3	80
3		Nonadecane	268	C19H40	000629-92-5	80
4		Tridecane, 1-iodo-	310	C13H27I	035599-77-0	80
5		Heptadecane, 2-methyl-	254	C18H38	001560-89-0	72



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG022819\
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 Sample : K1643-01
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 INTERIOR-WALL

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	3.29	3.6 ng		55236	1	8.16	306502	20.0
Hexanoic acid	7.14	2.2 ng		33427	1	8.16	306502	20.0
unknown7.69	7.69	87.6 ng		1343140	1	8.16	306502	20.0
Hexanoic acid, 2-...	9.43	5.5 ng		83874	1	8.16	306502	20.0
Octanoic acid	10.27	3.1 ng		70034	2	10.98	454373	20.0
Phthalic anhydride	12.75	3.2 ng		71900	2	10.98	454373	20.0
Tetradecanoic acid	16.89	2.2 ng		87961	4	17.53	811376	20.0
n-Hexadecanoic acid	18.40	46.4 ng		1882210	4	17.53	811376	20.0
9-octadecenoic ac...	19.53	3.4 ng		136540	4	17.53	811376	20.0
Octadecanoic acid	19.65	15.0 ng		608405	4	17.53	811376	20.0
Hexadecane, 2-met...	19.87	2.2 ng		95298	5	21.82	850102	20.0
Heptafluorobutyri...	21.41	8.2 ng		349267	5	21.82	850102	20.0
Methoxyacetic aci...	23.31	2.9 ng		122603	5	21.82	850102	20.0
Octacosane	24.16	2.7 ng		130734	6	25.20	983058	20.0