

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021119\
 Data File : BG039437.D
 Acq On : 11 Feb 2019 16:54
 Operator : JU/SJ
 Sample : K1424-01
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SS2-RB200035

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	15	22	31	rBV2	100853	206760	6.46%	0.833%
2	3.296	31	35	44	rVB	48129	72986	2.28%	0.294%
3	3.842	122	128	135	rBV	38781	69487	2.17%	0.280%
4	4.236	179	195	201	rBV	102251	304602	9.51%	1.228%
5	4.988	315	323	330	rBV	54952	118762	3.71%	0.479%
6	5.141	341	349	356	rBV	43342	87007	2.72%	0.351%
7	5.252	362	368	375	rVB	58198	91198	2.85%	0.368%
8	5.710	437	446	460	rVB	1002204	1575820	49.22%	6.352%
9	7.308	710	718	729	rBV	1063280	1842097	57.53%	7.425%
10	7.696	775	784	791	rBV	978541	1669213	52.13%	6.729%
11	8.172	858	865	874	rBV	199969	327341	10.22%	1.319%
12	8.495	912	920	928	rBV	671422	1144643	35.75%	4.614%
13	9.347	1057	1065	1074	rBV	605038	1063352	33.21%	4.286%
14	10.991	1338	1345	1352	rBV	283925	514516	16.07%	2.074%
15	13.418	1750	1758	1767	rBV	1567479	2432344	75.97%	9.805%
16	14.234	1892	1897	1903	rVB	101078	147100	4.59%	0.593%
17	14.792	1985	1992	2001	rBV2	470538	786198	24.56%	3.169%
18	16.278	2238	2245	2256	rBV	1363180	2068426	64.60%	8.338%
19	17.536	2451	2459	2466	rBV	628822	1005381	31.40%	4.053%
20	17.724	2484	2491	2500	rBV9	13419	35365	1.10%	0.143%
21	18.393	2599	2605	2616	rBV	193933	327358	10.22%	1.320%
22	19.539	2791	2800	2803	rBV8	18134	55558	1.74%	0.224%
23	19.656	2816	2820	2828	rBV2	61816	102067	3.19%	0.411%
24	20.132	2895	2901	2907	rVB	2471291	3201731	100.00%	12.906%
25	20.584	2975	2978	2984	rBV5	19575	35949	1.12%	0.145%
26	20.902	3029	3032	3036	rBV3	29699	37455	1.17%	0.151%
27	21.424	3116	3121	3134	rVB	339870	504373	15.75%	2.033%
28	21.830	3183	3190	3197	rVB2	619390	1065812	33.29%	4.296%
29	21.988	3213	3217	3228	rVB2	105513	183318	5.73%	0.739%
30	22.623	3319	3325	3337	rBV	193347	363871	11.36%	1.467%
31	23.345	3442	3448	3457	rVB2	176819	334952	10.46%	1.350%
32	24.197	3585	3593	3600	rBV	230555	520304	16.25%	2.097%
33	25.208	3754	3765	3778	rVB3	443147	1400317	43.74%	5.645%
34	25.754	3851	3858	3868	rVB4	38101	104983	3.28%	0.423%

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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	26.388	3956	3966	3980	rVB2	175397	575341	17.97%	2.319%
36	26.529	3980	3990	4002	rBV6	47396	165027	5.15%	0.665%
37	27.828	4202	4211	4218	rBV3	34333	111632	3.49%	0.450%
38	28.679	4350	4356	4357	rBV4	24038	38057	1.19%	0.153%
39	29.560	4497	4506	4507	rBV8	31262	70856	2.21%	0.286%
40	29.807	4537	4548	4550	rBV8	17011	46445	1.45%	0.187%

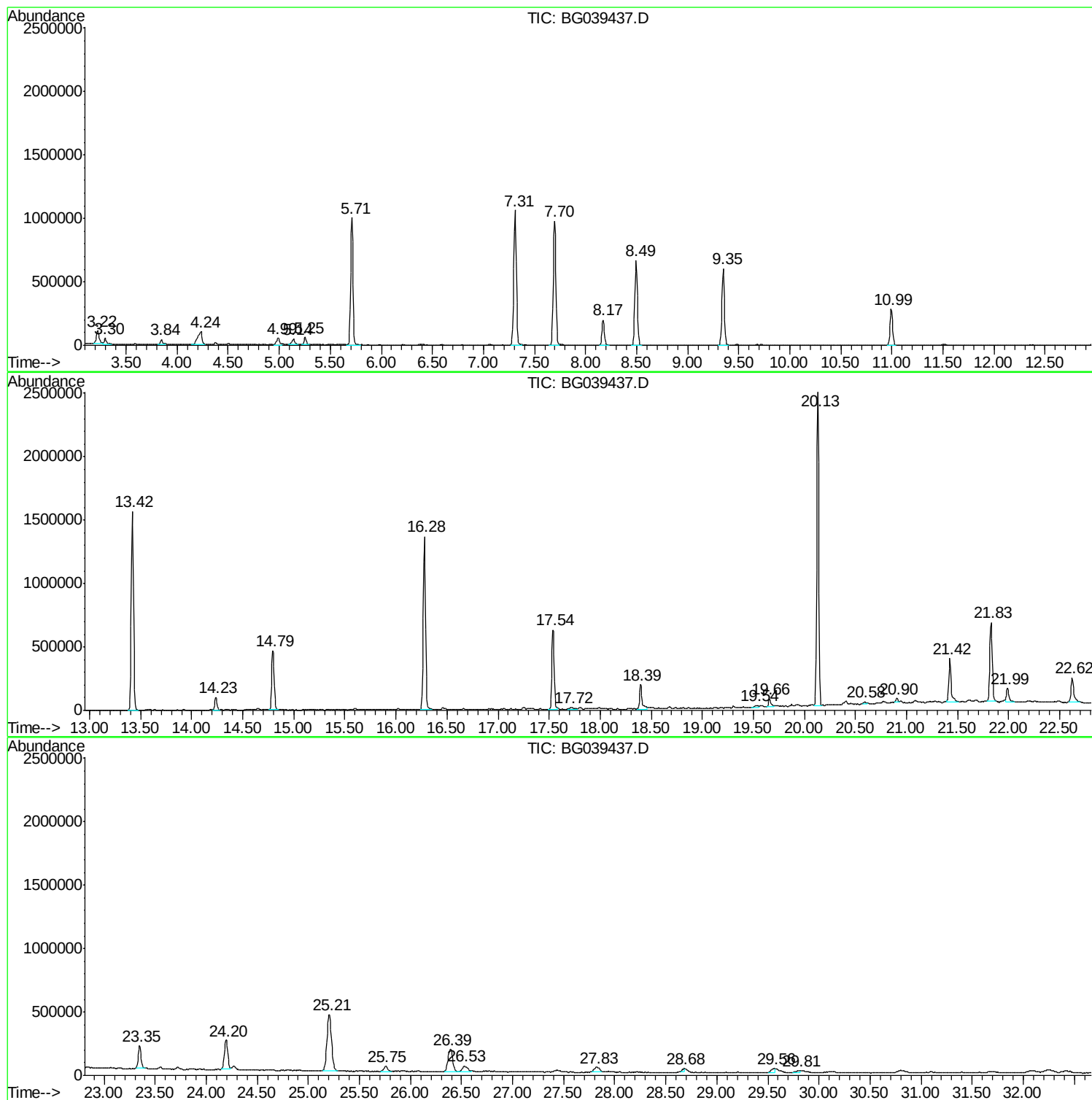
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG012919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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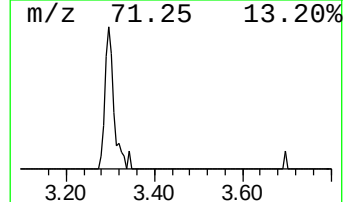
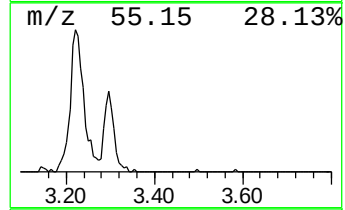
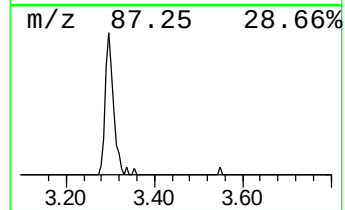
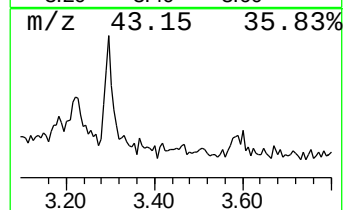
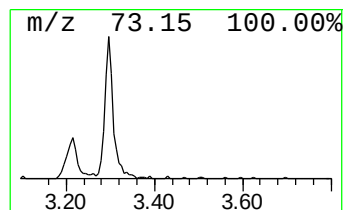
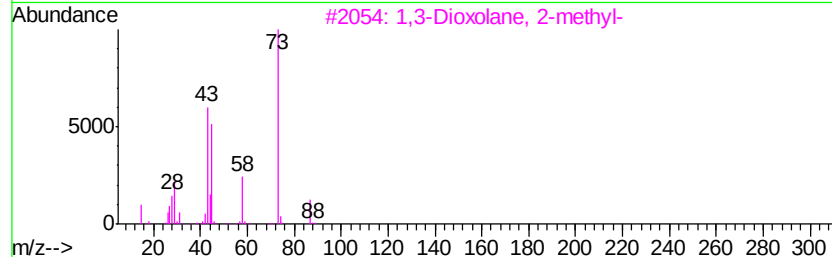
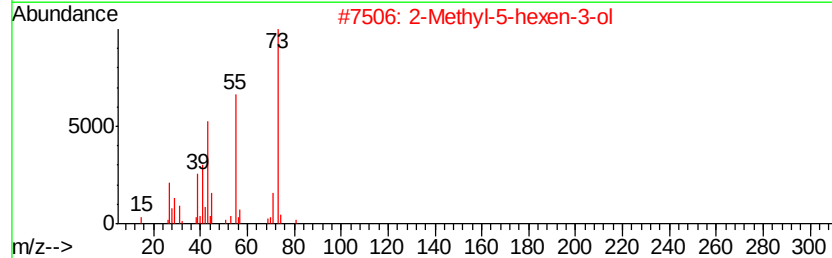
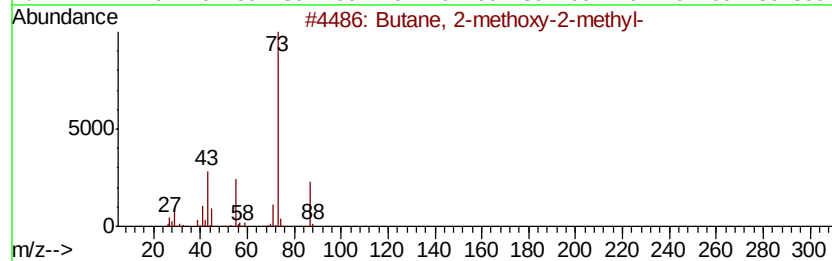
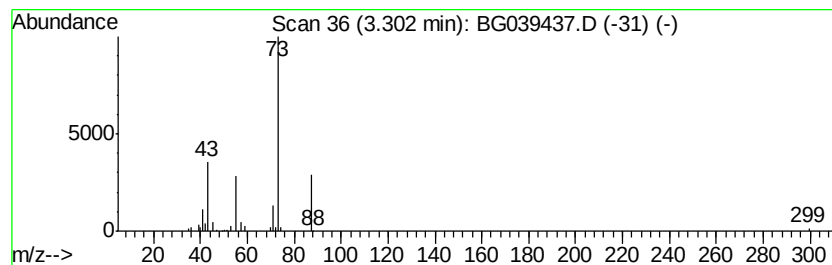
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG012919.M
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown3.30 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.30	4.46 ng	72986	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	39
2		2-Methyl-5-hexen-3-ol	114	C7H14O	032815-70-6	25
3		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	23
4		N-Ethylformamide	73	C3H7NO	000627-45-2	9
5		Silane, tetramethyl-	88	C4H12Si	000075-76-3	9



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Instrument :
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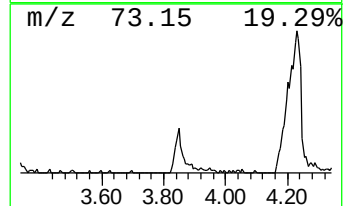
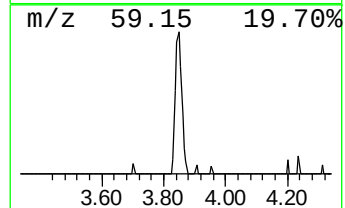
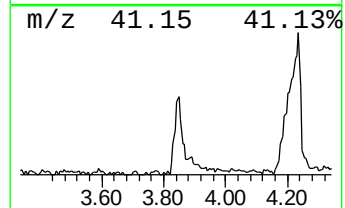
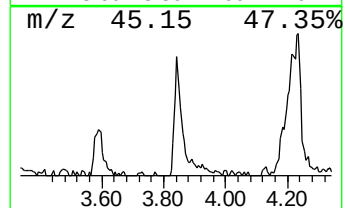
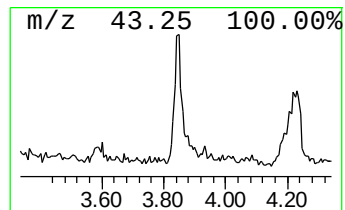
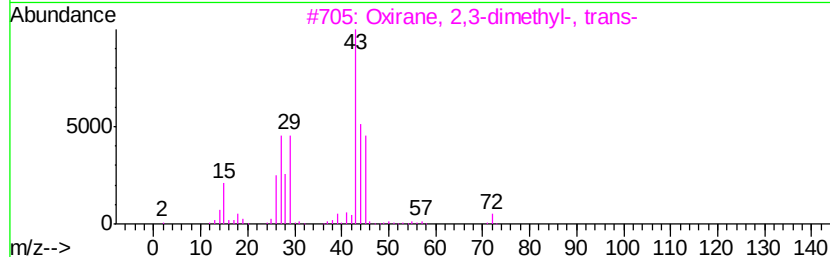
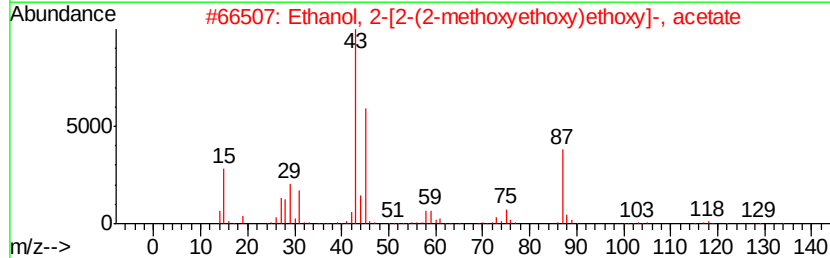
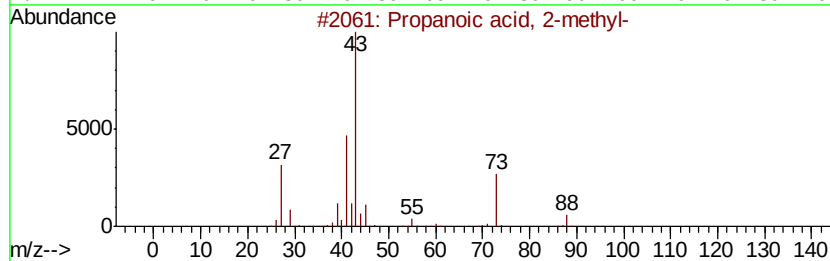
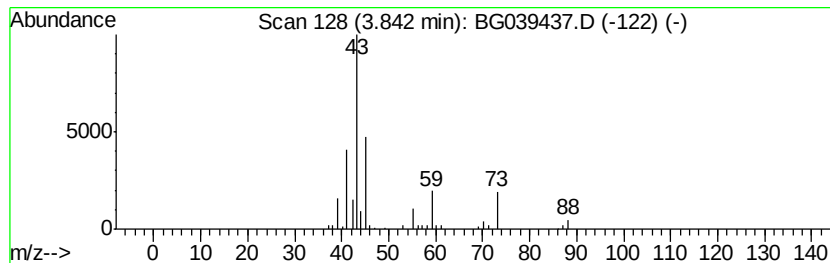
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TIC Library : C:\Database\NIST11.L
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Peak Number 3 unknown3.84 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.84	4.25 ng	69487	1,4-Dichlorobenzene-d4	8.17

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propanoic acid, 2-methyl-	88	C4H8O2	000079-31-2	43
2			Ethanol, 2-[2-(2-methoxyethoxy)ethoxy]-	206	C9H18O5	003610-27-3	25
3			Oxirane, 2,3-dimethyl-, trans-	72	C4H8O	021490-63-1	23
4			Oxirane, 2,3-dimethyl-, cis-	72	C4H8O	001758-33-4	23
5			Acetamide, N-(2-hydroxyethyl)-	103	C4H9NO2	000142-26-7	17



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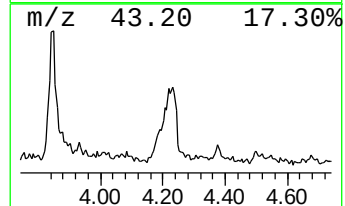
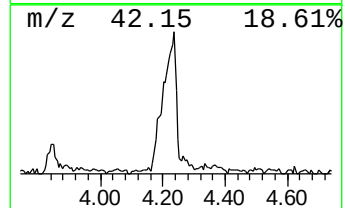
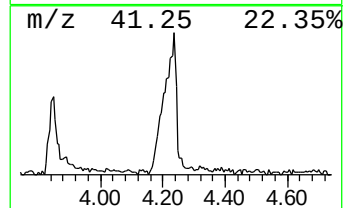
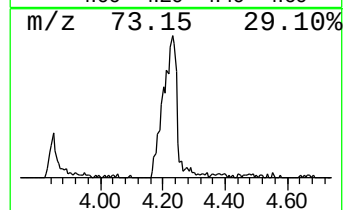
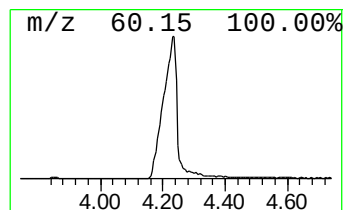
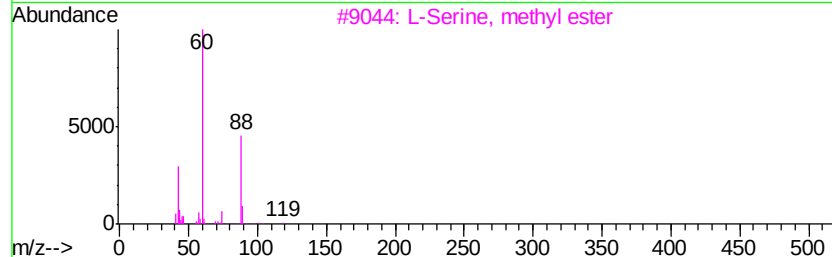
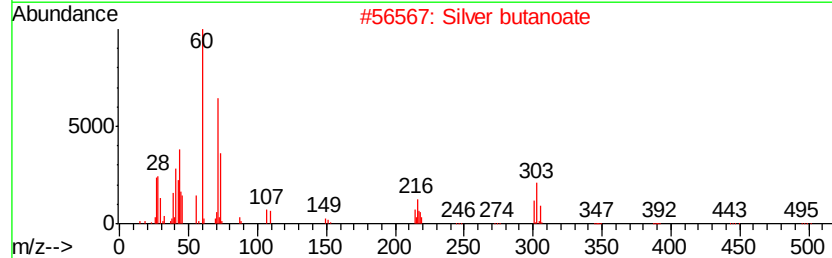
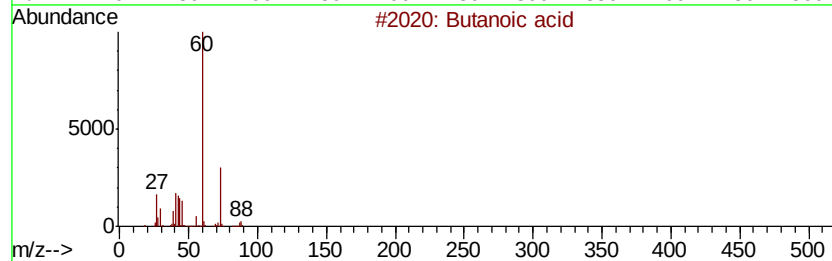
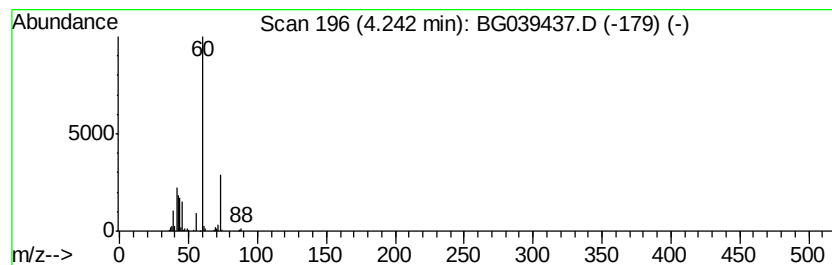
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TIC Library : C:\Database\NIST11.L
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Peak Number 4 Butanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.24	18.61 ng	304602	1,4-Dichlorobenzene-d4	8.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butanoic acid	88	C4H8O2	000107-92-6	56
2		Silver butanoate	194	C4H7AqO2	005076-24-4	43
3		L-Serine, methyl ester	119	C4H9NO3	002788-84-3	36
4		Butanoic acid, 4-chloro-	122	C4H7ClO2	000627-00-9	9
5		4-Bromobutyric acid	166	C4H7BrO2	002623-87-2	9



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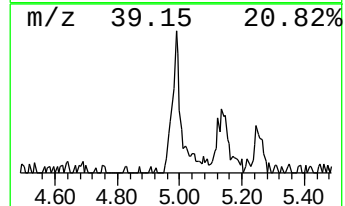
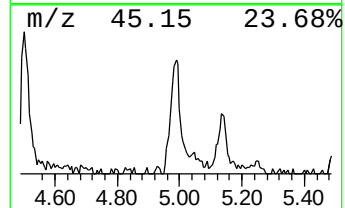
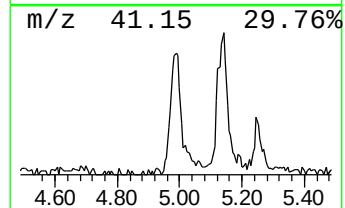
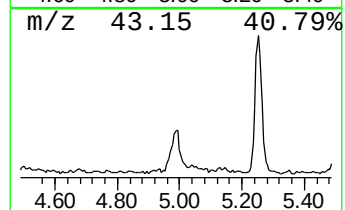
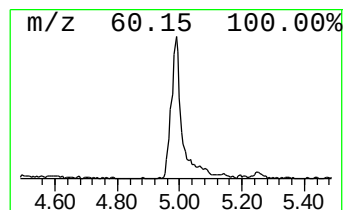
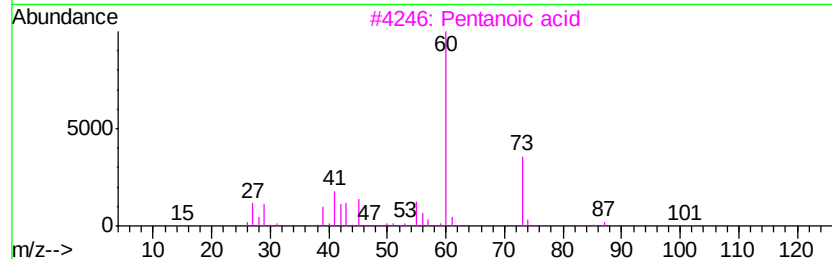
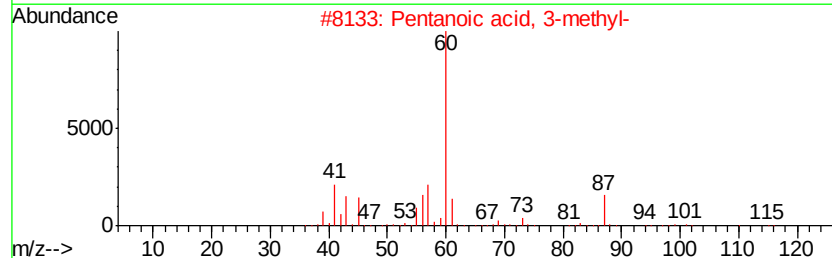
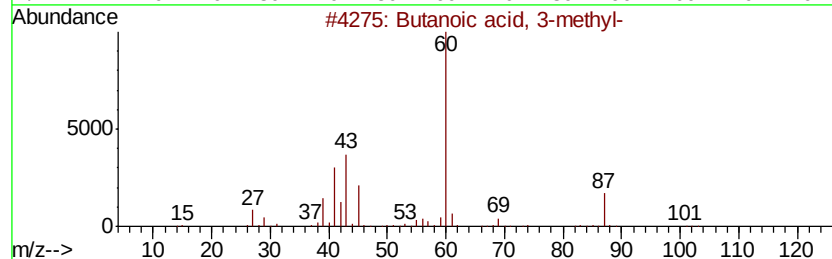
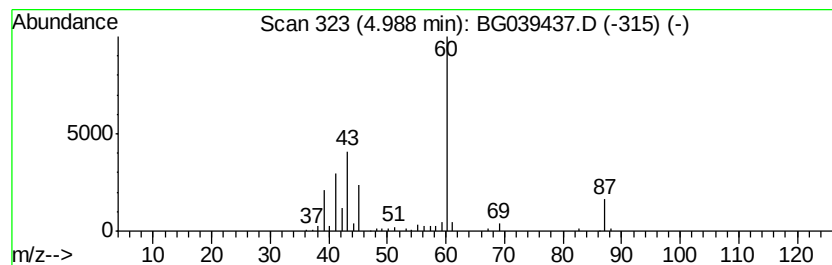
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TIC Library : C:\Database\NIST11.L
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Peak Number 5 Butanoic acid, 3-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.99	7.26 ng	118762	1,4-Dichlorobenzene-d4	8.17

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Butanoic acid, 3-methyl-	102	C5H10O2	000503-74-2	78
2		Pentanoic acid, 3-methyl-	116	C6H12O2	000105-43-1	72
3		Pentanoic acid	102	C5H10O2	000109-52-4	38
4		Formic acid hydrazide	60	CH4N2O	000624-84-0	9
5		Benserazide	257	C10H15N3O5	000322-35-0	9



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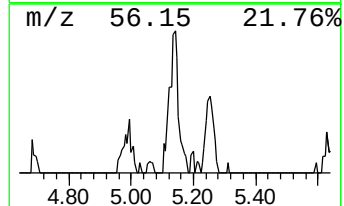
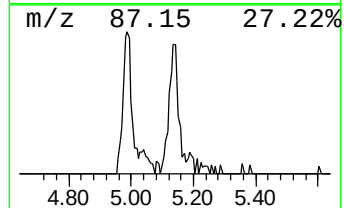
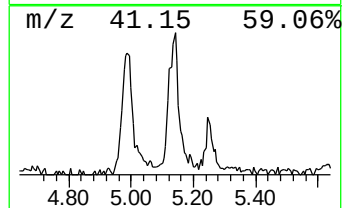
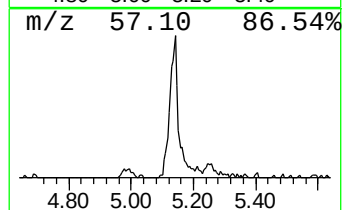
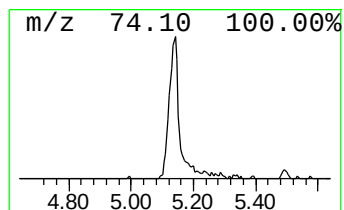
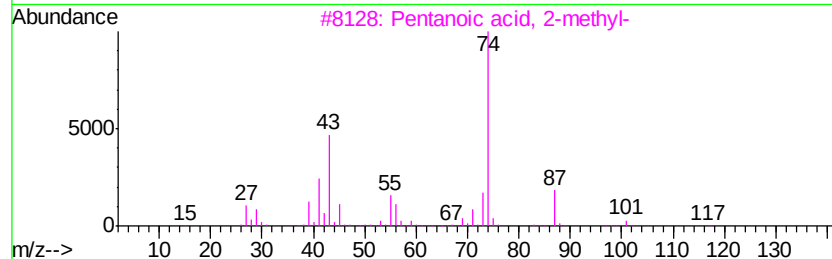
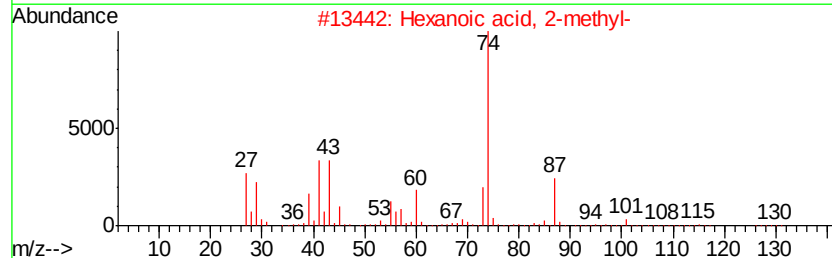
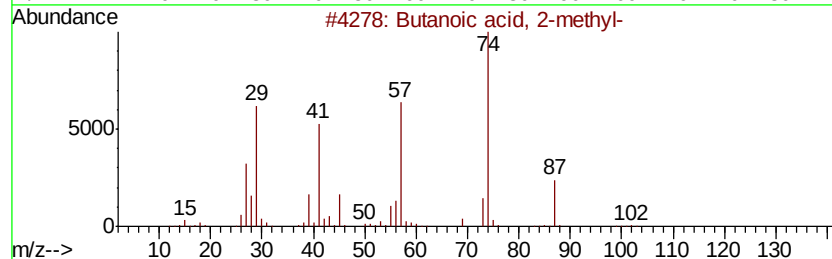
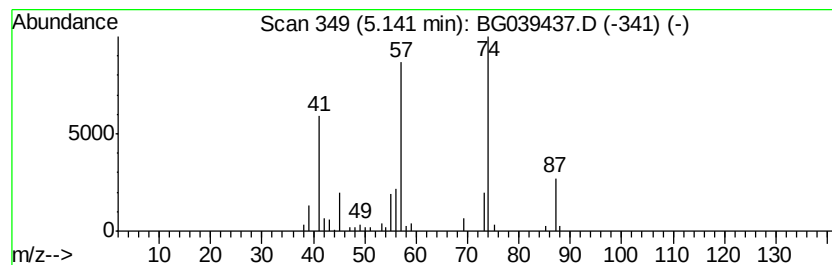
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Peak Number 6 Butanoic acid, 2-methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.14	5.32 ng	87007	1,4-Dichlorobenzene-d4	8.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butanoic acid, 2-methyl-	102	C5H10O2	000116-53-0	78
2		Hexanoic acid, 2-methyl-	130	C7H14O2	004536-23-6	42
3		Pentanoic acid, 2-methyl-	116	C6H12O2	000097-61-0	33
4		Propanoic acid	74	C3H6O2	000079-09-4	25
5		Oxirane, (butoxymethyl)-	130	C7H14O2	002426-08-6	17



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Data File : BG039437.D
Acq On : 11 Feb 2019 16:54
Operator : JU/SJ
Sample : K1424-01
Misc :
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Instrument :
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ClientSampleId :
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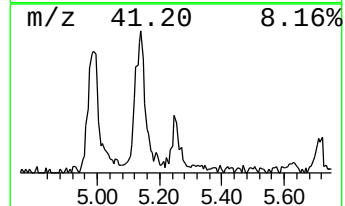
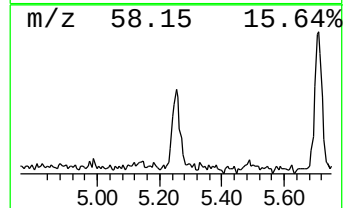
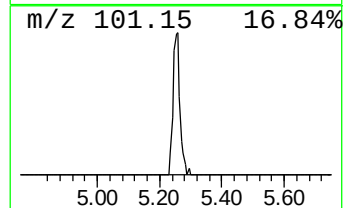
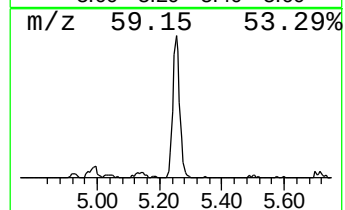
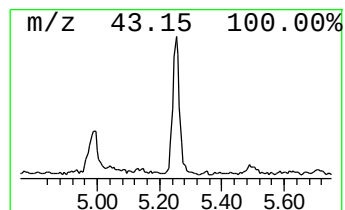
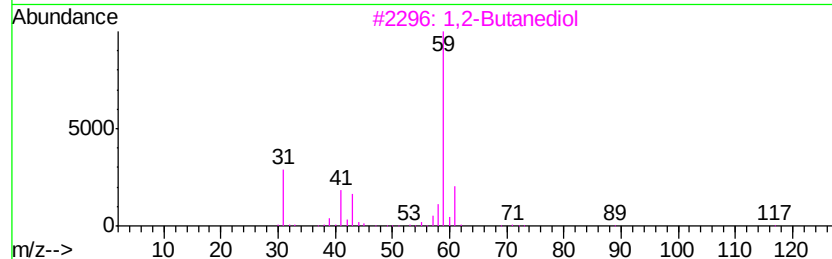
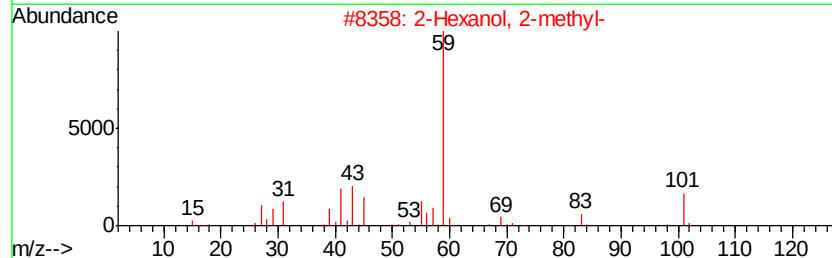
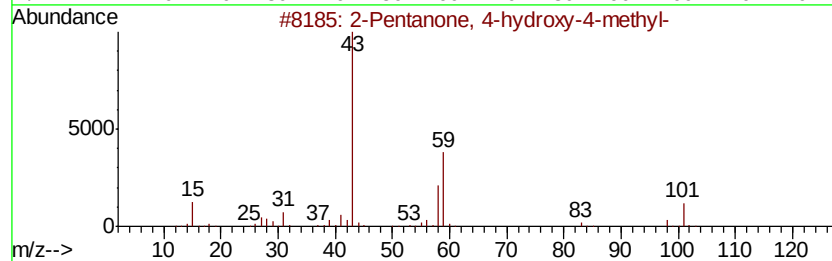
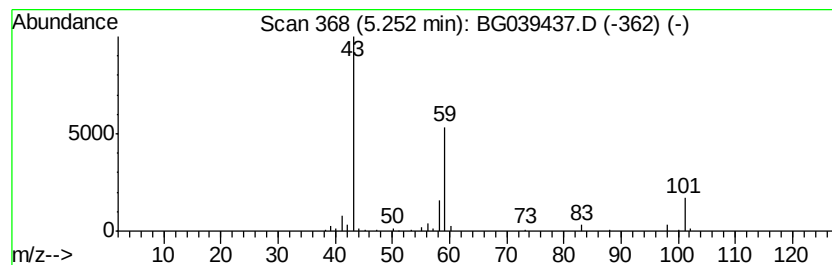
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.25	5.57 ng	91198	1,4-Dichlorobenzene-d4	8.17

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2	2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28
3	1,2-Butanediol	90	C4H10O2	000584-03-2	23
4	5-Hexen-2-one	98	C6H10O	000109-49-9	9
5	Formamide, N,N-diethyl-	101	C5H11NO	000617-84-5	9



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021119\
Data File : BG039437.D
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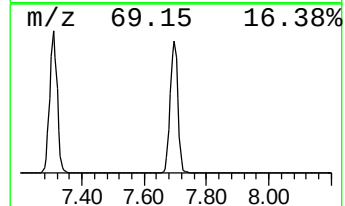
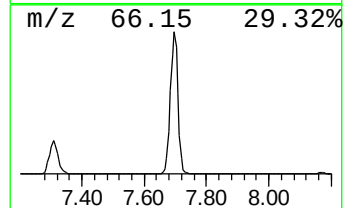
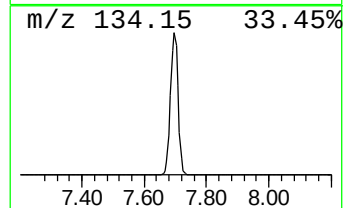
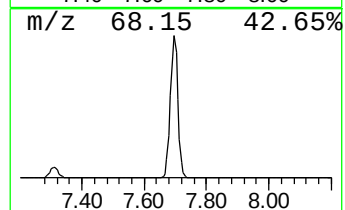
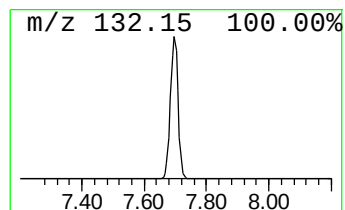
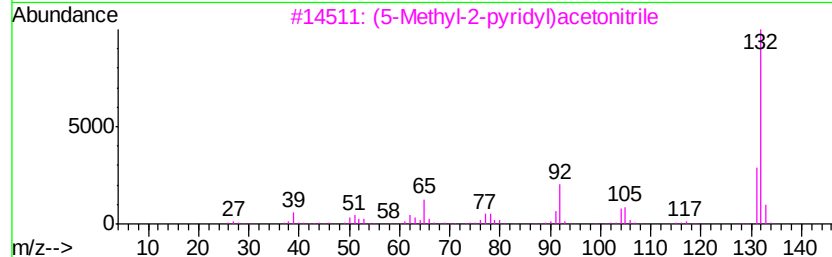
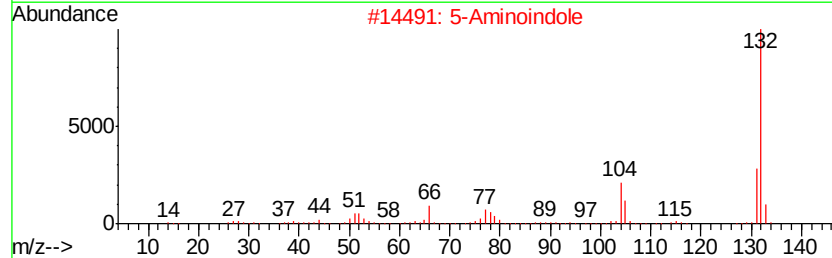
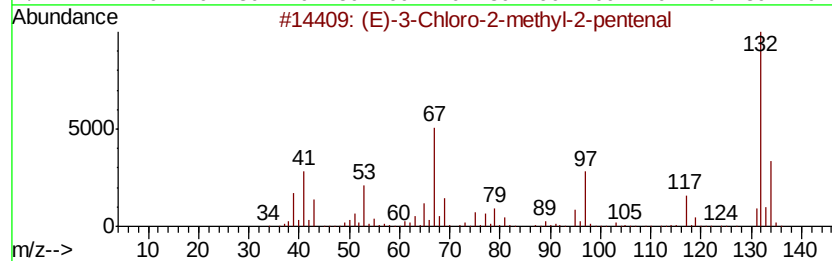
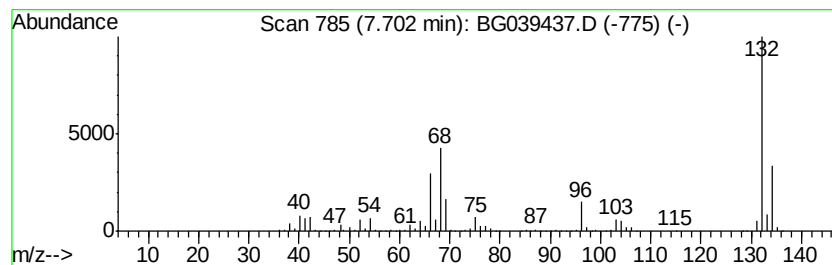
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 unknown7.70 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.70	101.99 ng	1669210	1,4-Dichlorobenzene-d4	8.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
2		5-Aminoindole	132	C8H8N2	005192-03-0	12
3		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	10
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021119\
Data File : BG039437.D
Acq On : 11 Feb 2019 16:54
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Sample : K1424-01
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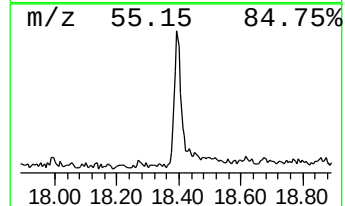
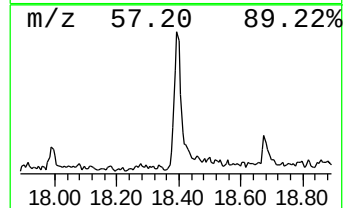
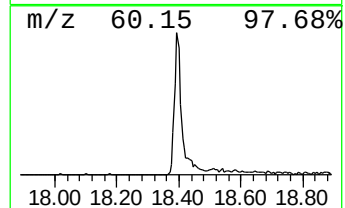
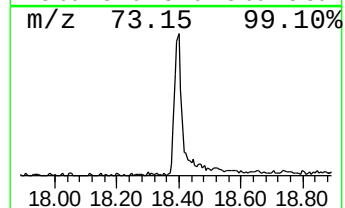
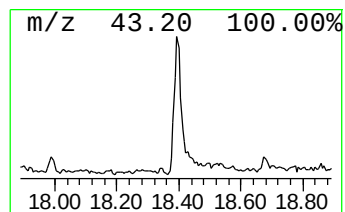
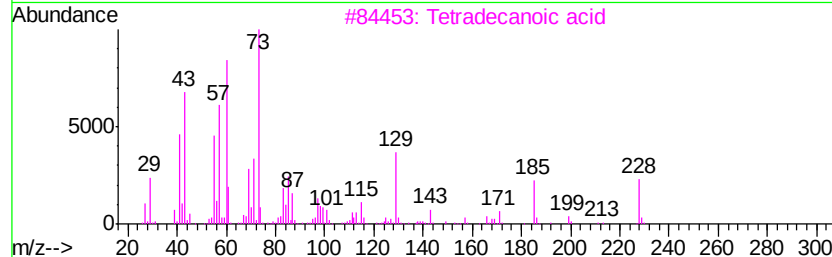
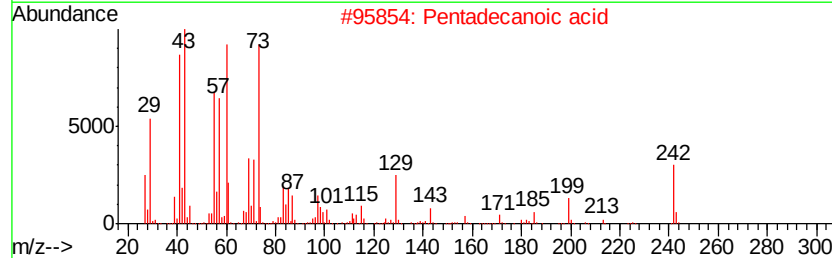
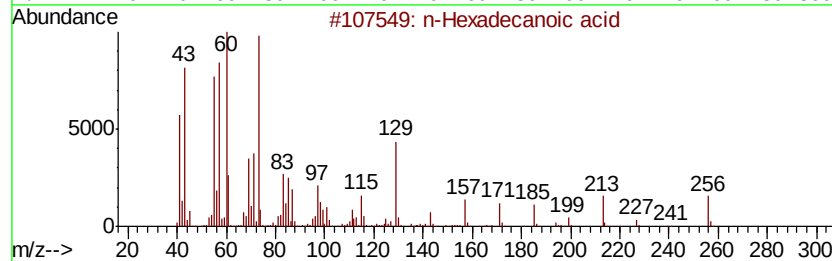
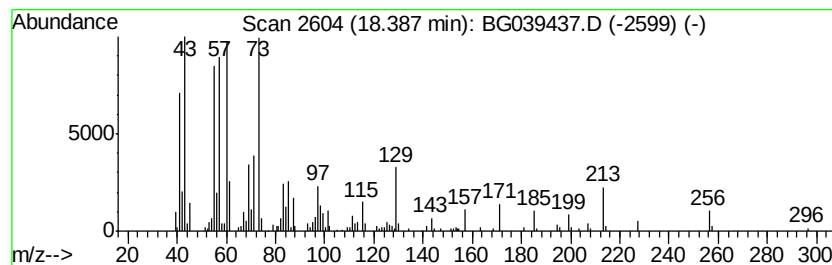
Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG012919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 n-Hexadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.39	6.51 ng	327358	Phenanthrene-d10	17.54	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	80
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	72
4	Nonanoic acid	158	C9H18O2	000112-05-0	43
5	n-Decanoic acid	172	C10H20O2	000334-48-5	25



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021119\
Data File : BG039437.D
Acq On : 11 Feb 2019 16:54
Operator : JU/SJ
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Misc :
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ClientSampleId :
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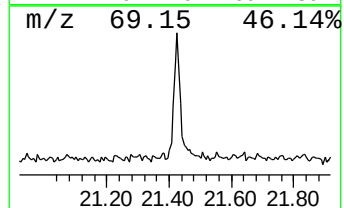
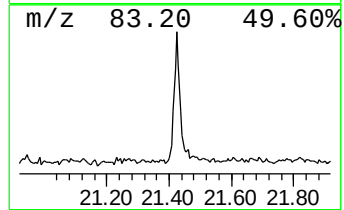
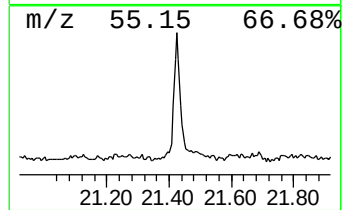
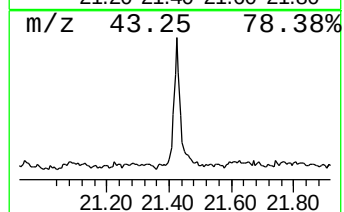
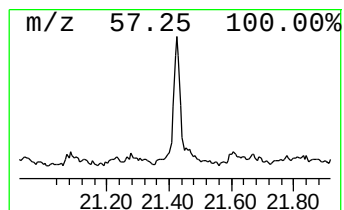
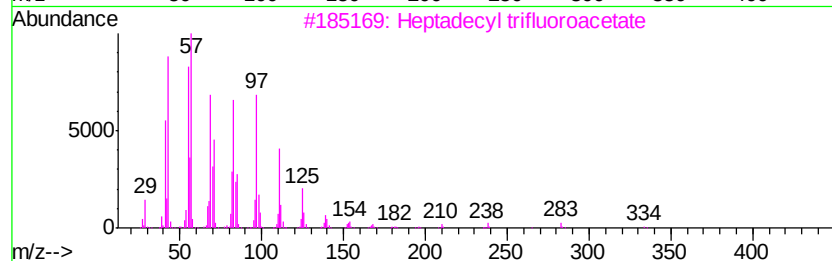
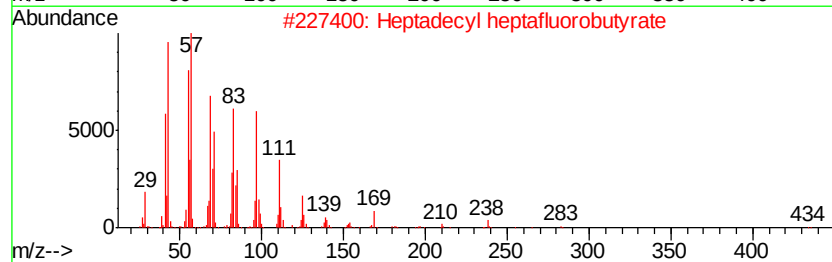
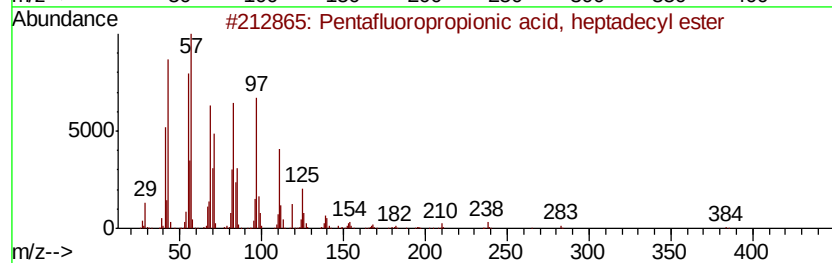
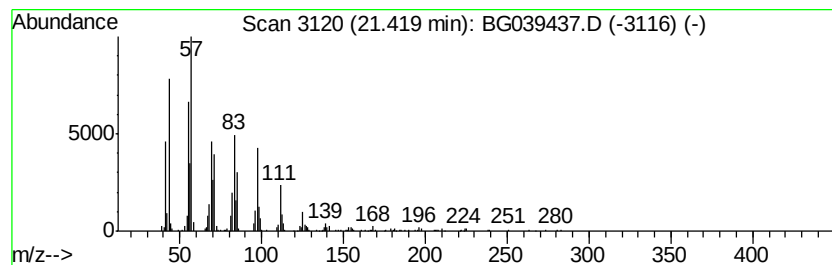
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Pentafluoropropionic acid, ... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.42	9.46 ng	504373	Chrysene-d12	21.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentafluoropropionic acid, hepta...	402	C20H35F5O2	959218-78-1	91
2		Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	91
3		Heptadecyl trifluoroacetate	352	C19H35F3O2	1000351-87-0	91
4		Octadecyl trifluoroacetate	366	C20H37F3O2	079392-43-1	91
5		Heptafluorobutyric acid, hexadec...	438	C20H33F7O2	006385-15-5	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021119\
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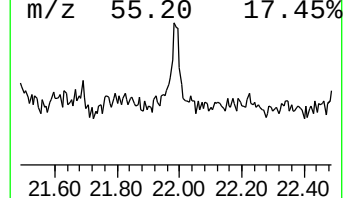
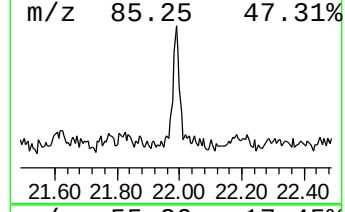
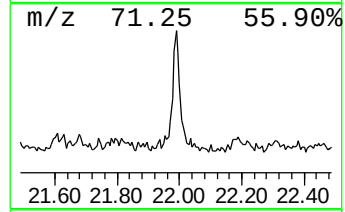
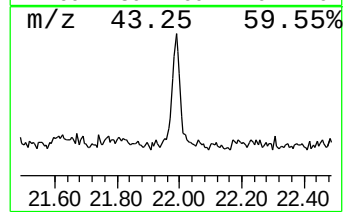
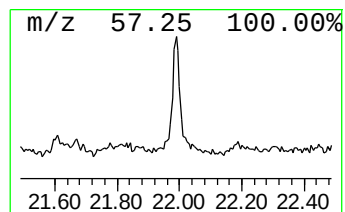
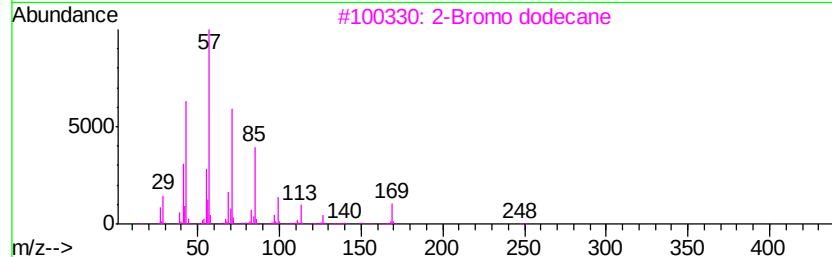
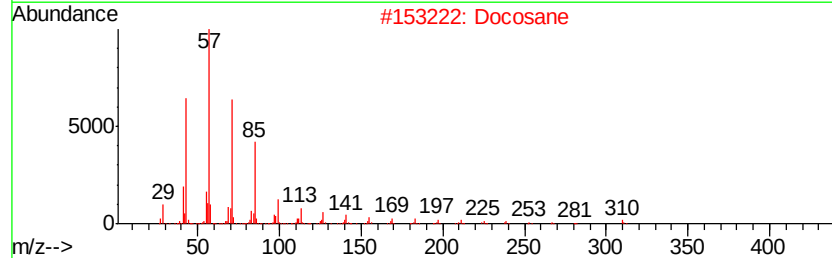
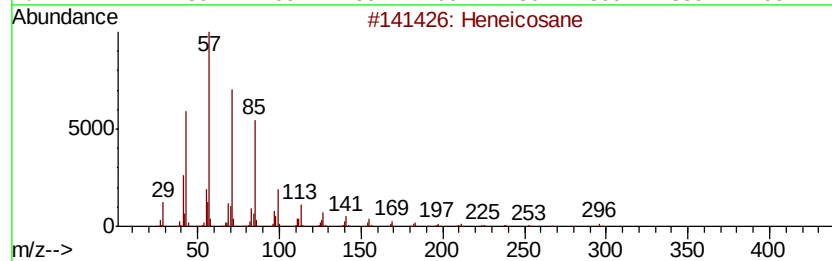
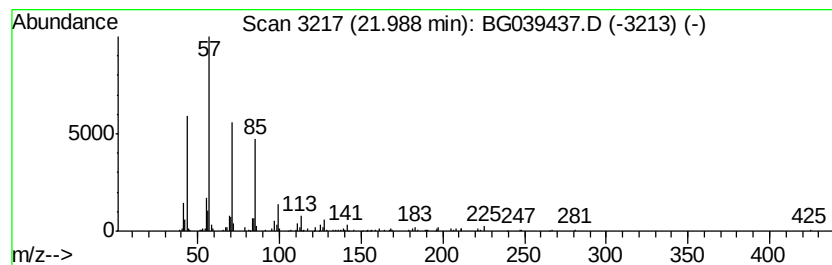
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 13 Heneicosane Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD		R.T.
21.99	3.44 ng	183318	Chrysene-d12		21.83
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	90
2	Docosane	310	C22H46	000629-97-0	87
3	2-Bromo dodecane	248	C12H25Br	013187-99-0	86
4	Heptadecane	240	C17H36	000629-78-7	86
5	Hexadecane	226	C16H34	000544-76-3	78



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021119\
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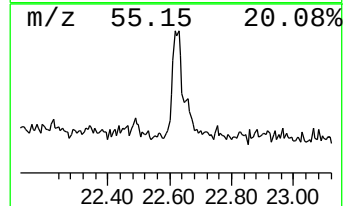
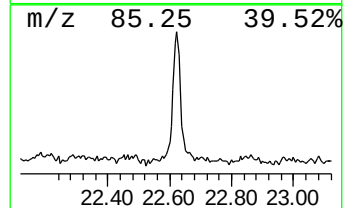
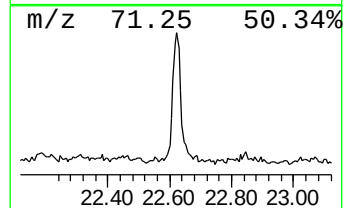
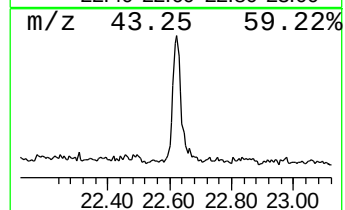
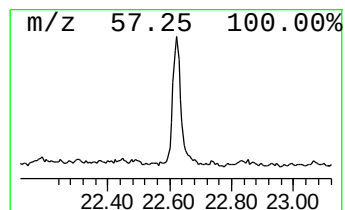
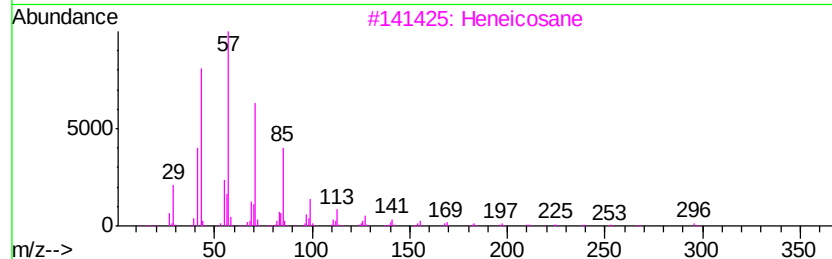
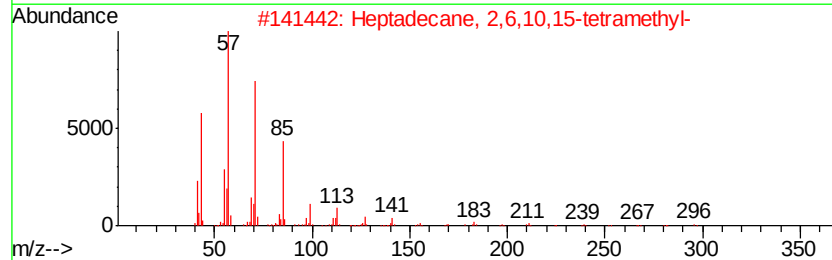
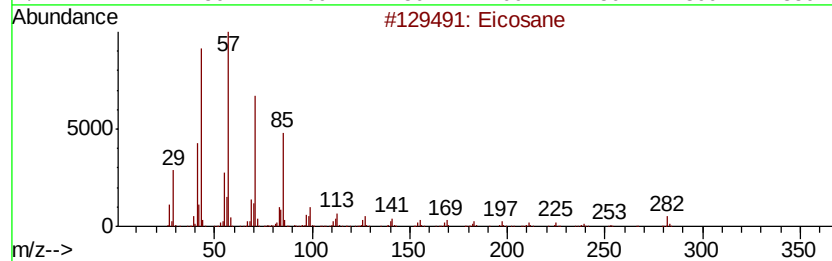
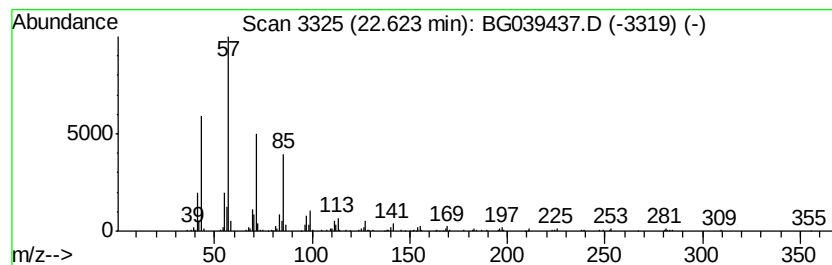
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Eicosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.62	6.83 ng	363871	Chrysene-d12	21.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	93
2		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	91
3		Heneicosane	296	C21H44	000629-94-7	90
4		Tetracosane	338	C24H50	000646-31-1	90
5		Octacosane	394	C28H58	000630-02-4	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021119\
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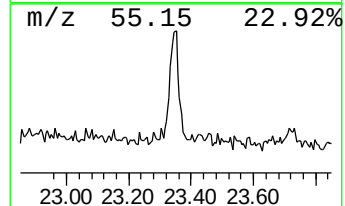
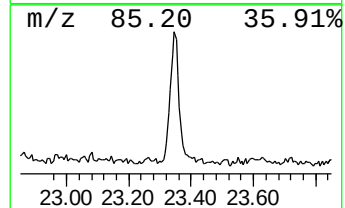
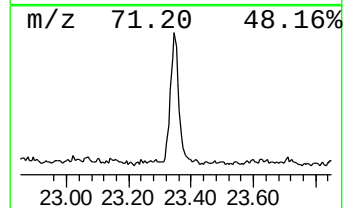
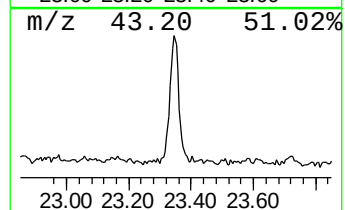
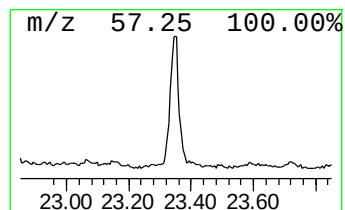
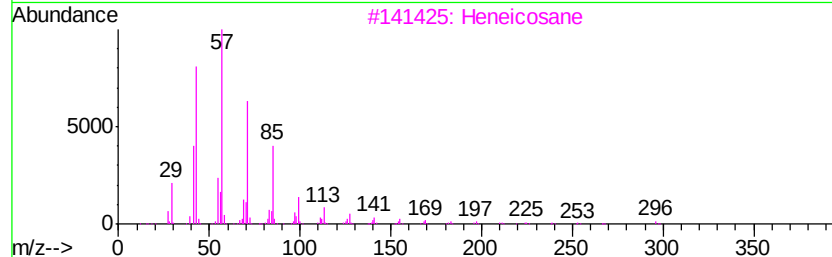
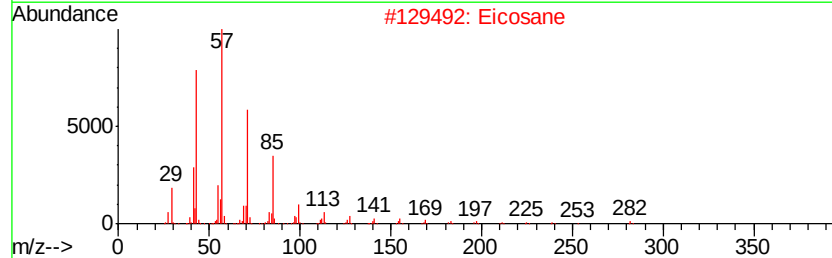
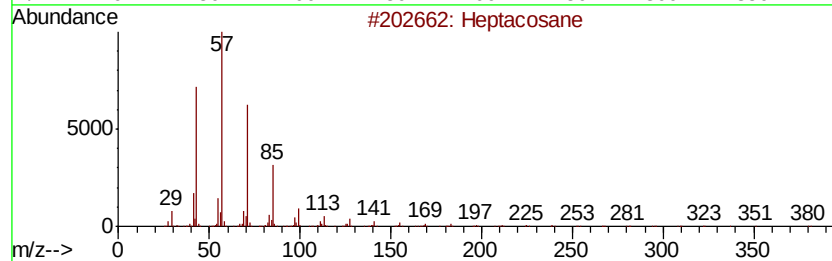
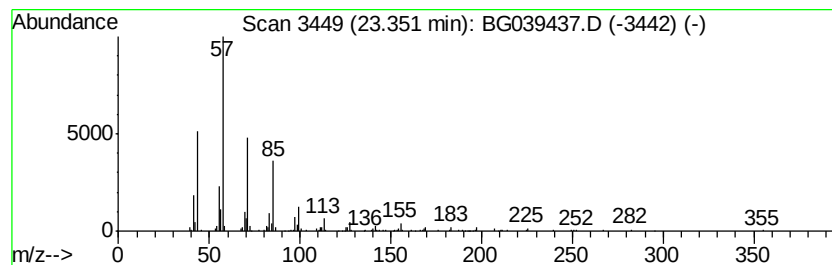
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 15 Heptacosane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.35	6.29 ng	334952	Chrysene-d12	21.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptacosane	380	C27H56	000593-49-7	91
2		Eicosane	282	C20H42	000112-95-8	87
3		Heneicosane	296	C21H44	000629-94-7	86
4		Nonadecane	268	C19H40	000629-92-5	83
5		Docosane	310	C22H46	000629-97-0	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021119\
Data File : BG039437.D
Acq On : 11 Feb 2019 16:54
Operator : JU/SJ
Sample : K1424-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

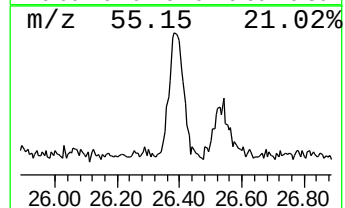
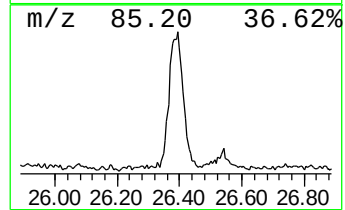
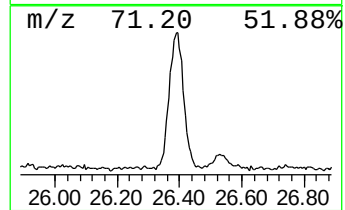
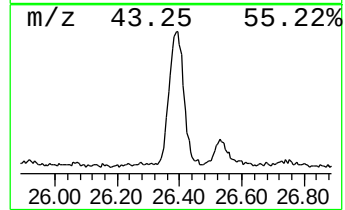
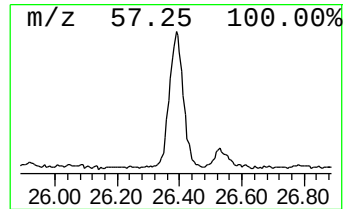
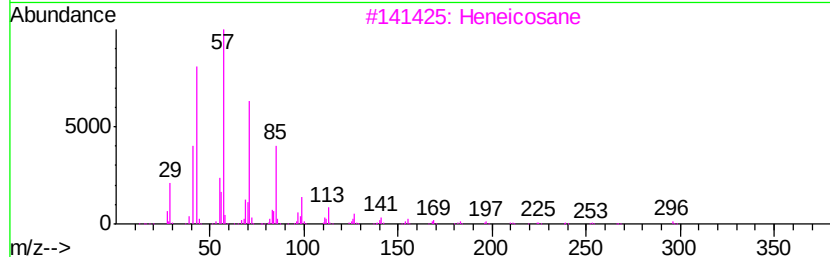
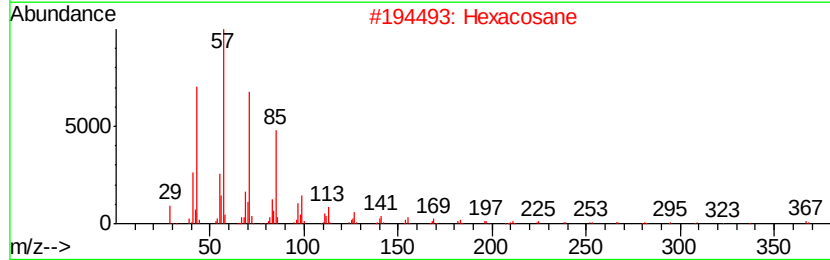
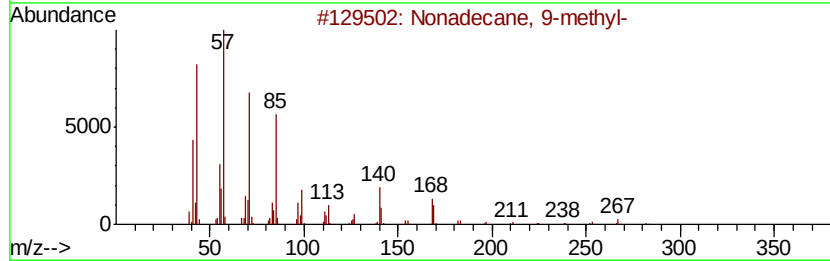
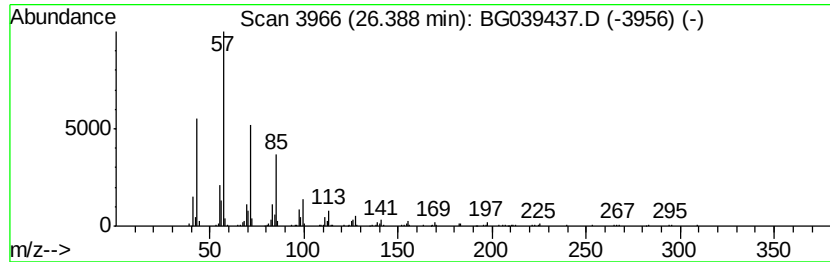
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Nonadecane, 9-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD		R.T.
26.39	8.22 ng	575341	Perylene-d12		25.21
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonadecane, 9-methyl-	282	C20H42	013287-24-6	94
2	Hexacosane	366	C26H54	000630-01-3	91
3	Heneicosane	296	C21H44	000629-94-7	91
4	Octacosane	394	C28H58	000630-02-4	87
5	Pentadecane	212	C15H32	000629-62-9	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG021119\
Data File : BG039437.D
Acq On : 11 Feb 2019 16:54
Operator : JU/SJ
Sample : K1424-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SS2-RB200035

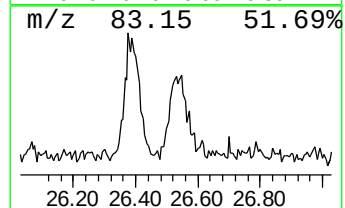
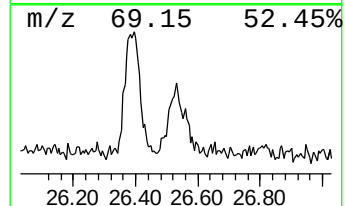
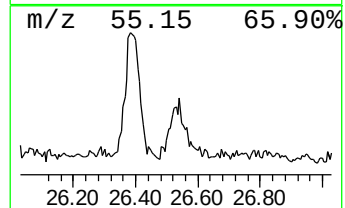
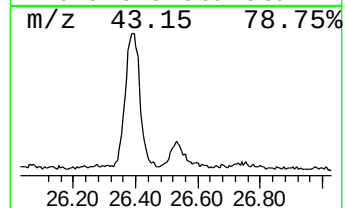
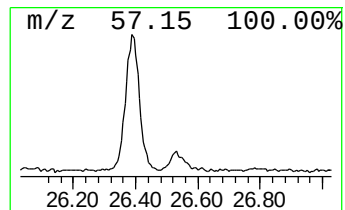
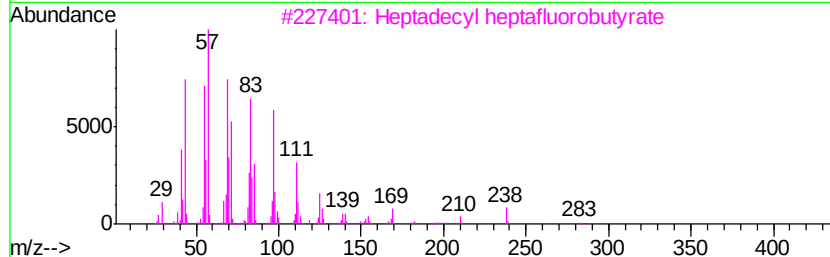
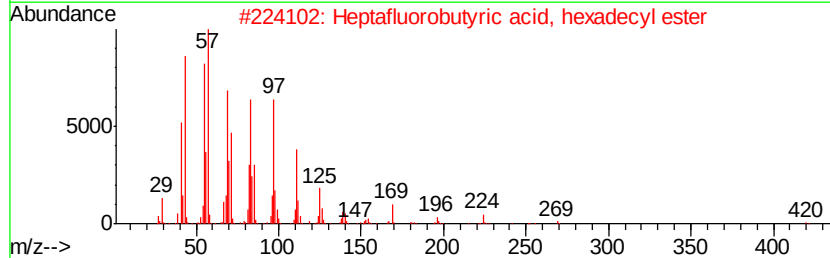
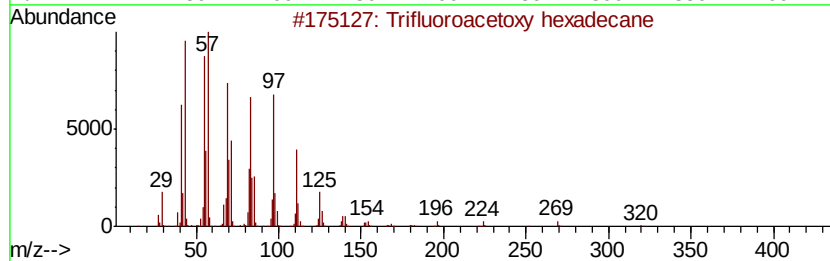
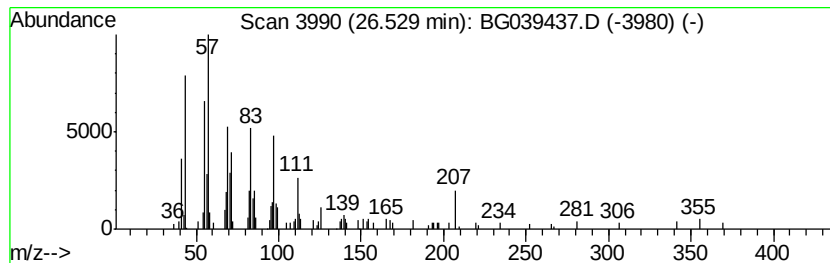
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 Trifluoroacetoxy hexadecane Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.53	2.36 ng	165027	Perylene-d12	25.21

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	93
2		Heptafluorobutyric acid, hexadec...	438	C20H33F7O2	006385-15-5	90
3		Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	90
4		Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	90
5		E-15-Heptadecenal	252	C17H32O	1000130-97-9	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG021119\
 Data File : BG039437.D
 Acq On : 11 Feb 2019 16:54
 Operator : JU/SJ
 Sample : K1424-01
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SS2-RB200035

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
unknown3.30	3.30	4.5	ng	72986	1	8.17	327341	20.0
unknown3.84	3.84	4.3	ng	69487	1	8.17	327341	20.0
Butanoic acid	4.24	18.6	ng	304602	1	8.17	327341	20.0
Butanoic acid, 3-...	4.99	7.3	ng	118762	1	8.17	327341	20.0
Butanoic acid, 2-...	5.14	5.3	ng	87007	1	8.17	327341	20.0
2-Pentanone, 4-hy...	5.25	5.6	ng	91198	1	8.17	327341	20.0
unknown7.70	7.70	102.0	ng	1669210	1	8.17	327341	20.0
n-Hexadecanoic acid	18.39	6.5	ng	327358	4	17.54	1005380	20.0
Pentafluoropropio...	21.42	9.5	ng	504373	5	21.83	1065810	20.0
Heneicosane	21.99	3.4	ng	183318	5	21.83	1065810	20.0
Eicosane	22.62	6.8	ng	363871	5	21.83	1065810	20.0
Heptacosane	23.35	6.3	ng	334952	5	21.83	1065810	20.0
Nonadecane, 9-met...	26.39	8.2	ng	575341	6	25.21	1400320	20.0
Trifluoroacetoxy ...	26.53	2.4	ng	165027	6	25.21	1400320	20.0