

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG061019\
 Data File : BG041161.D
 Acq On : 10 Jun 2019 12:18
 Operator : HP/JU
 Sample : K3223-02
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SS043MW014-190605

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-BG052919.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	3.187	12	17	26	rVB	173420	271168	3.88%	1.036%
2	3.434	53	59	74	rBV	2367067	3914317	56.05%	14.948%
3	5.202	349	360	364	rBV	3424818	6983733	100.00%	26.670%
4	5.649	429	436	445	rBV2	73672	159349	2.28%	0.609%
5	7.247	702	708	715	rBV	75432	135493	1.94%	0.517%
6	7.635	767	774	793	rBV	215543	527209	7.55%	2.013%
7	8.105	847	854	864	rBV	152137	262853	3.76%	1.004%
8	8.428	902	909	918	rVB	585012	1016030	14.55%	3.880%
9	8.922	986	993	1001	rVB	104849	169418	2.43%	0.647%
10	9.280	1045	1054	1068	rBV	478715	880933	12.61%	3.364%
11	10.925	1327	1334	1347	rBV	207748	421769	6.04%	1.611%
12	11.166	1368	1375	1387	rVB2	58814	109448	1.57%	0.418%
13	11.295	1390	1397	1407	rBV	115105	211983	3.04%	0.810%
14	13.358	1739	1748	1757	rBV	1549774	2381451	34.10%	9.095%
15	13.587	1780	1787	1799	rBV	394391	621090	8.89%	2.372%
16	14.738	1976	1983	1995	rVB2	337743	573323	8.21%	2.189%
17	15.649	2132	2138	2147	rBV	154670	239335	3.43%	0.914%
18	17.482	2441	2450	2460	rBV	485224	757533	10.85%	2.893%
19	20.073	2885	2891	2897	rBV	2919989	3890909	55.71%	14.859%
20	21.348	3104	3108	3115	rBV2	98928	155427	2.23%	0.594%
21	21.754	3171	3177	3187	rBV	510880	870342	12.46%	3.324%
22	21.901	3197	3202	3208	rVB	82683	112366	1.61%	0.429%
23	22.523	3302	3308	3315	rVB	85388	143355	2.05%	0.547%
24	23.223	3422	3427	3439	rVB	86130	163421	2.34%	0.624%
25	24.045	3561	3567	3575	rVB2	76652	151916	2.18%	0.580%
26	25.026	3725	3734	3735	rBV2	64713	134118	1.92%	0.512%
27	25.079	3736	3743	3759	rVB2	263902	804966	11.53%	3.074%
28	26.178	3923	3930	3941	rVB4	44574	122321	1.75%	0.467%

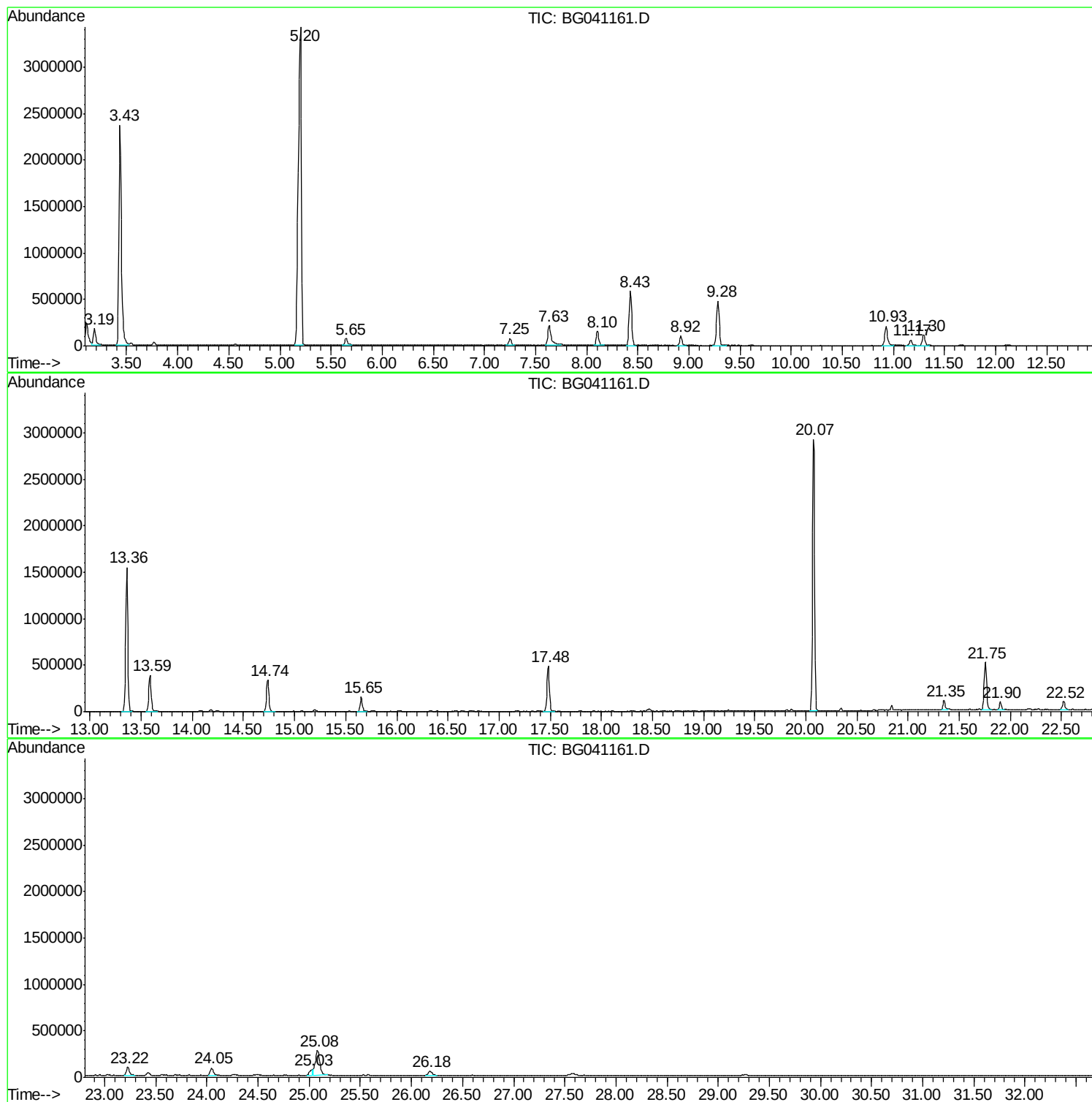
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Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SS043MW014-190605

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG052919.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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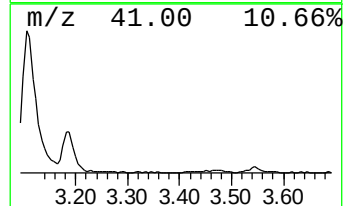
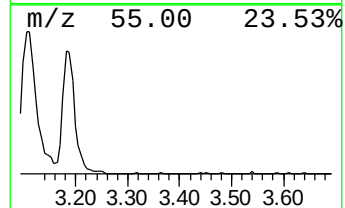
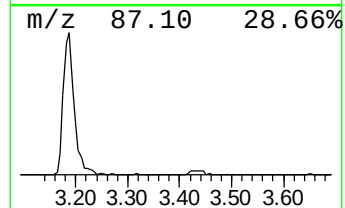
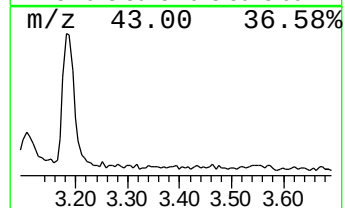
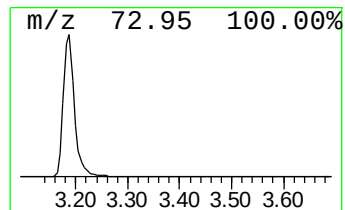
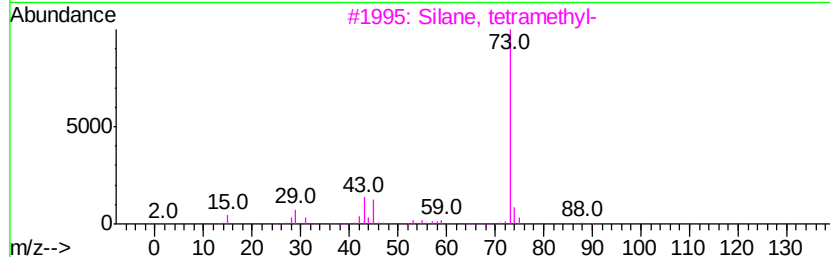
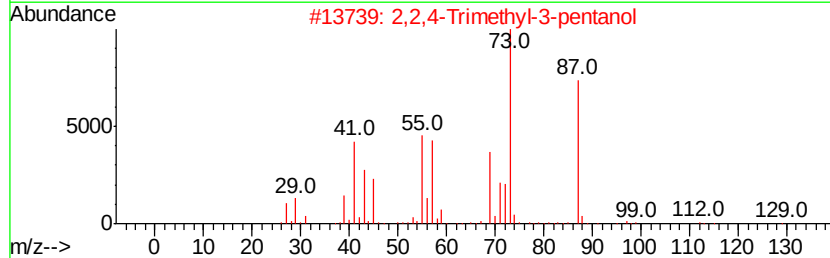
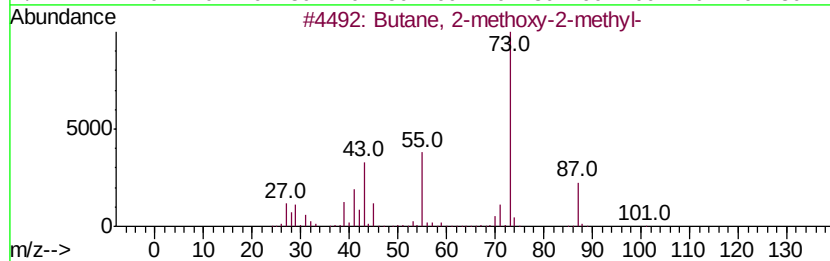
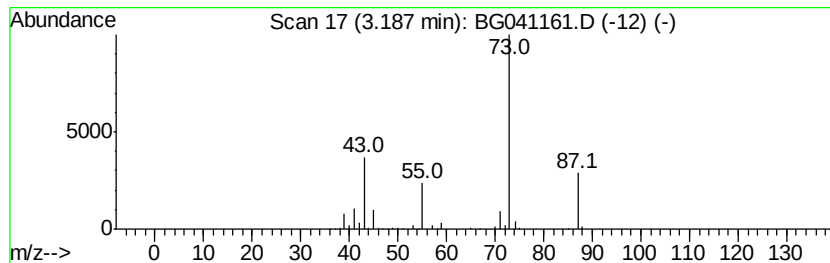
Instrument :
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ClientSampleId :
SS043MW014-190605

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG052919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.19	20.63 ng	271168	1,4-Dichlorobenzene-d4	8.10
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Butane, 2-methoxy-2-methyl-	102 C6H14O	000994-05-8 72
2		2,2,4-Trimethyl-3-pentanol	130 C8H18O	005162-48-1 40
3		Silane, tetramethyl-	88 C4H12Si	000075-76-3 9
4		1-Propene-1-thiol	74 C3H6S	000925-89-3 9
5		N-Ethylformamide	73 C3H7NO	000627-45-2 9



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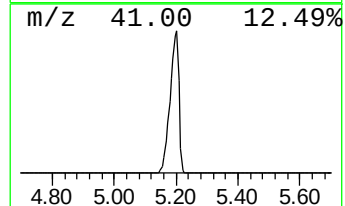
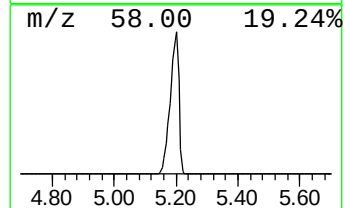
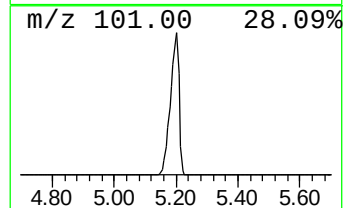
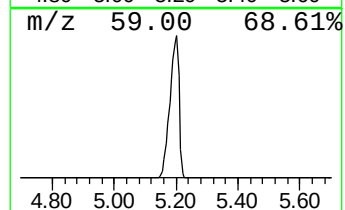
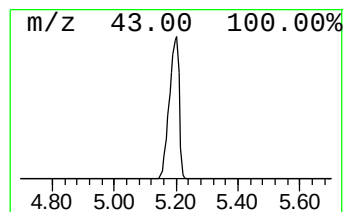
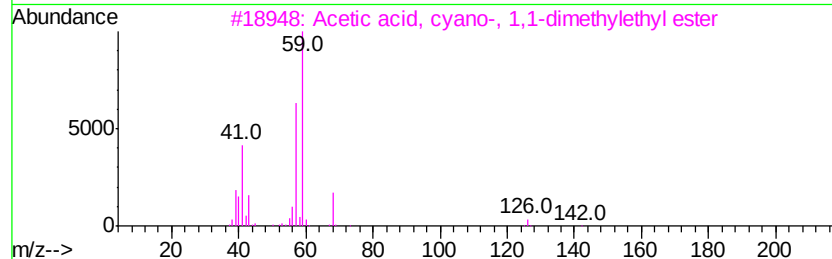
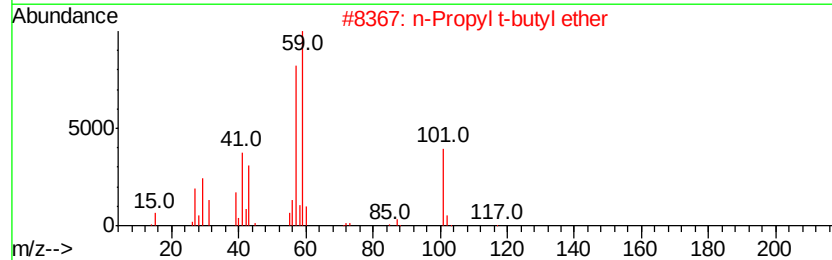
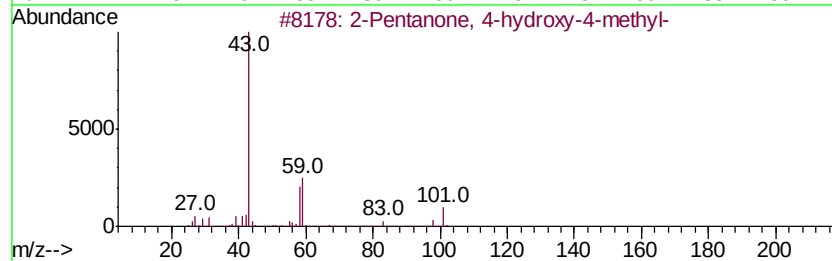
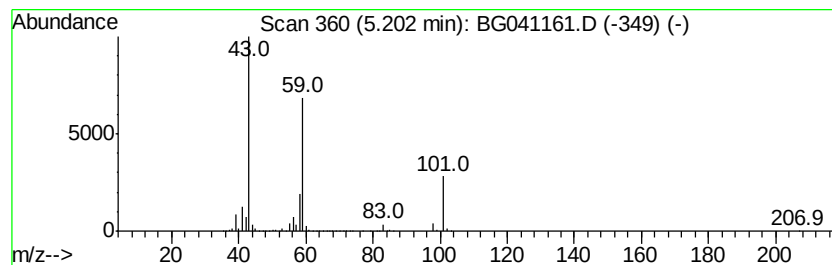
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.20	531.38 ng	6983730	1,4-Dichlorobenzene-d4	8.10

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2		000123-42-2	53
2	n-Propyl t-butyl ether	116	C7H16O		1000334-81-9	28
3	Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2		001116-98-9	25
4	2-Nonanone, 9-hydroxy-	158	C9H18O2		025368-56-3	25
5	3-Hydroxy-3-methyl-2-butanone	102	C5H10O2		000115-22-0	16



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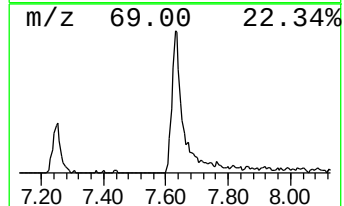
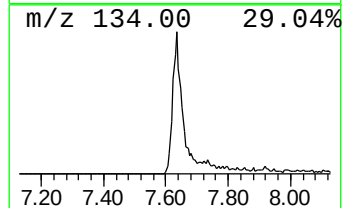
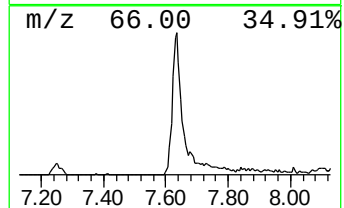
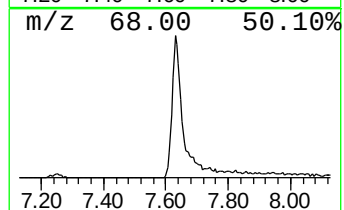
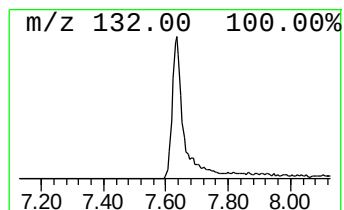
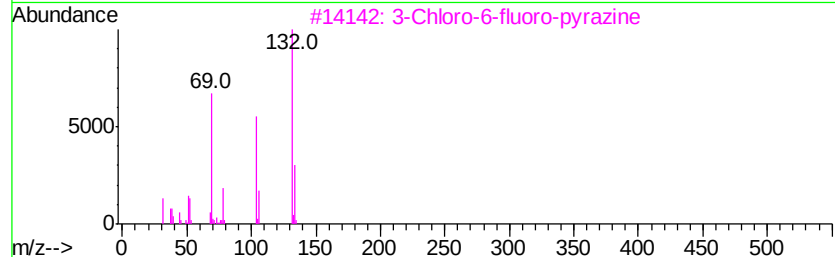
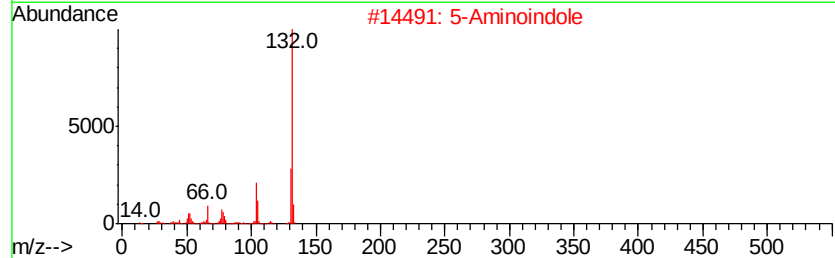
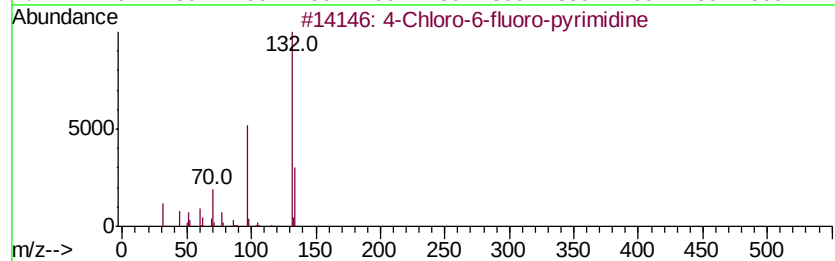
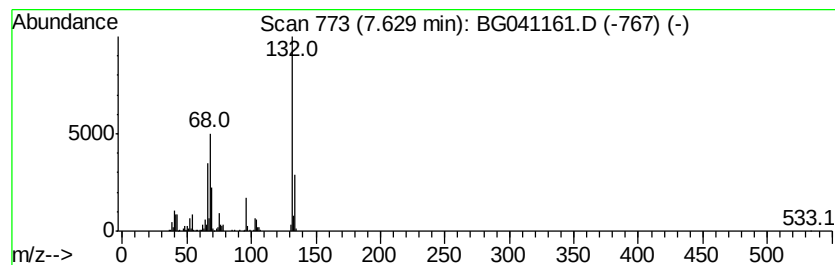
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TIC Library : C:\Database\NIST11.L
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Peak Number 4 unknown7.63 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.63	40.11 ng	527209	1,4-Dichlorobenzene-d4	8.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
2		5-Aminoindole	132	C8H8N2	005192-03-0	27
3		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
4		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
5		3H-Benzo[4,5]imidazo[1,2-c]oxazo...	218	C12H14N2O2	1000316-99-9	12



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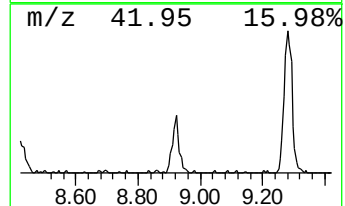
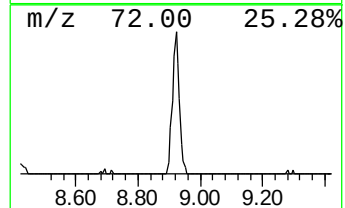
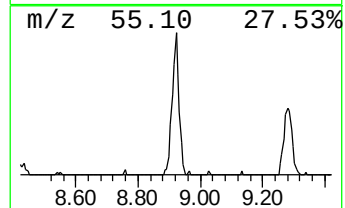
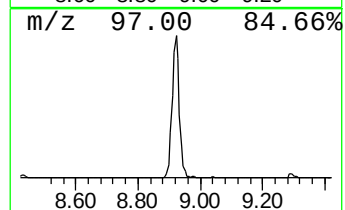
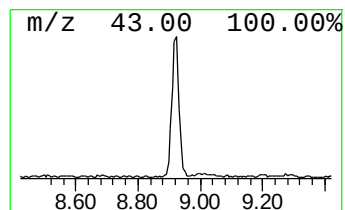
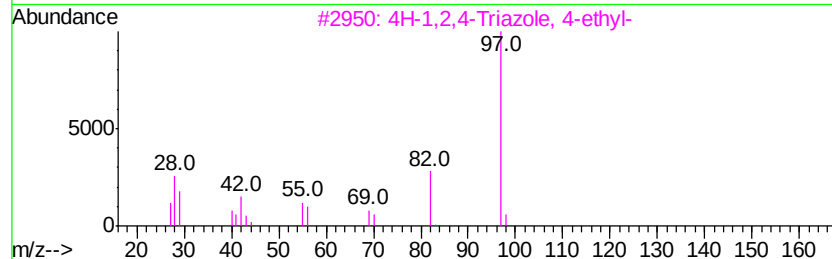
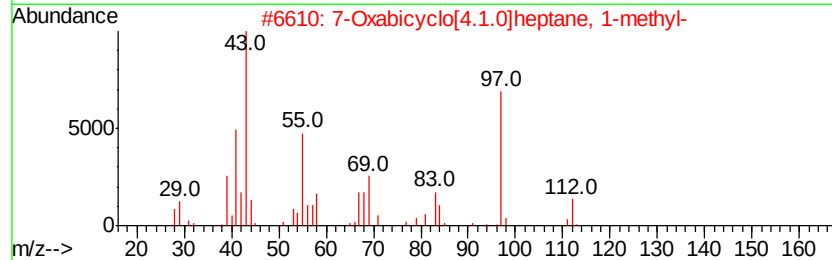
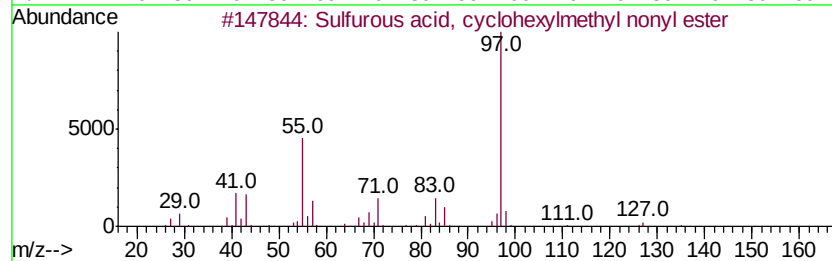
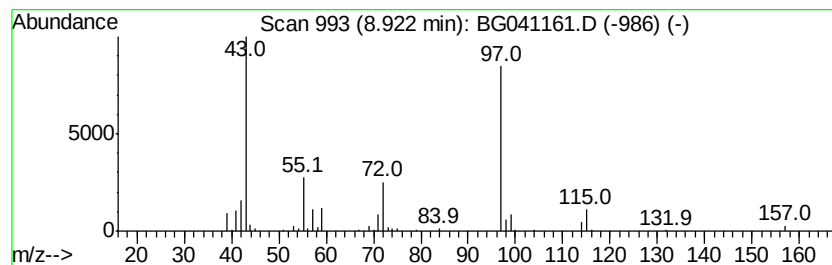
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Peak Number 6 unknown8.92 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.92	12.89 ng	169418	1,4-Dichlorobenzene-d4	8.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Sulfurous acid, cyclohexylmethyl...	304	C16H32O3S	1000309-21-8	43
2		7-Oxabicyclo[4.1.0]heptane, 1-me...	112	C7H12O	001713-33-3	42
3		4H-1,2,4-Triazole, 4-ethyl-	97	C4H7N3	043183-55-7	37
4		2,4-Pentanedione, 3-methyl-	114	C6H10O2	000815-57-6	25
5		Sulfurous acid, cyclohexylmethyl...	388	C22H44O3S	1000309-22-3	25



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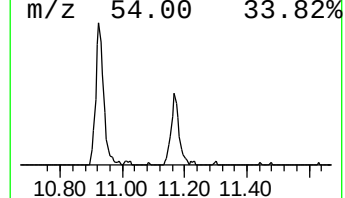
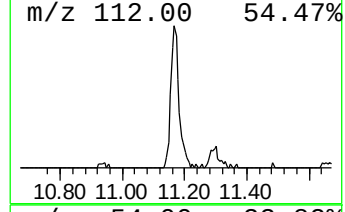
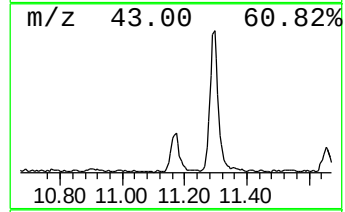
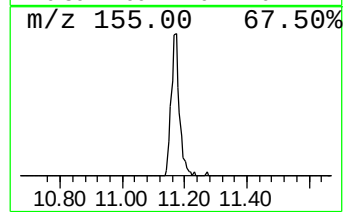
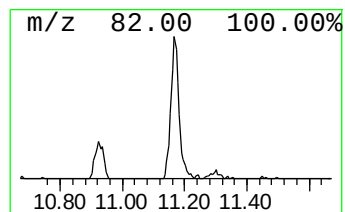
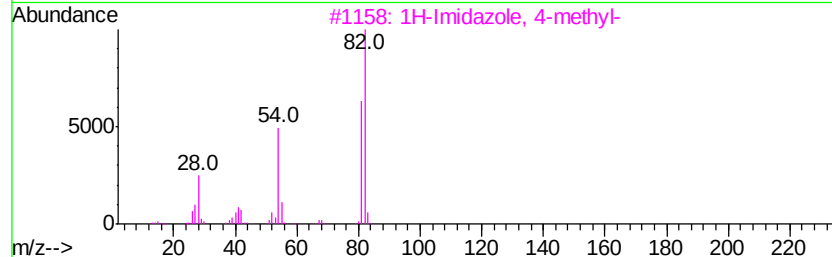
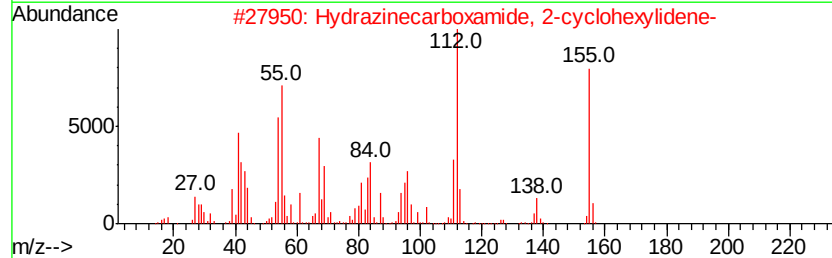
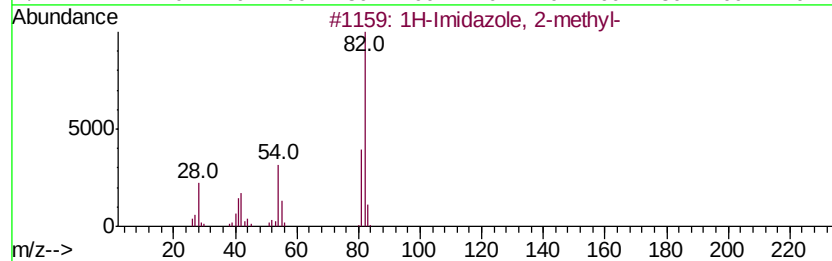
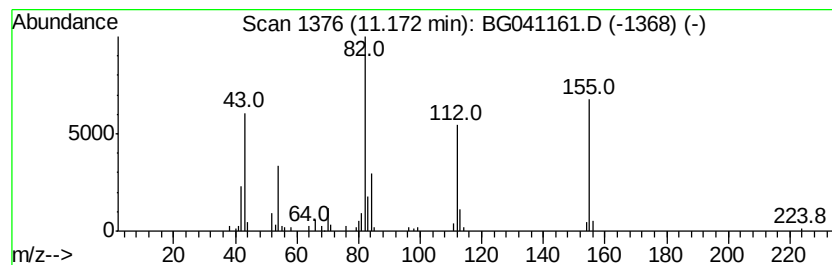
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 unknown11.17 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.17	5.19 ng	109448	Naphthalene-d8	10.93

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Imidazole, 2-methyl-	82	C4H6N2	000693-98-1	38
2		Hydrazinecarboxamide, 2-cyclohex...	155	C7H13N3O	001589-61-3	32
3		1H-Imidazole, 4-methyl-	82	C4H6N2	000822-36-6	16
4		3-Aminocrotononitrile	82	C4H6N2	001118-61-2	12
5		2,2-Dimethyl-3(2H)-furanone	112	C6H8O2	035298-48-7	11



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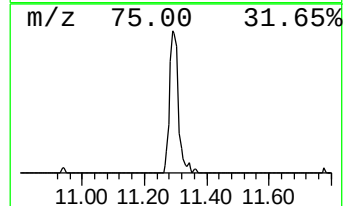
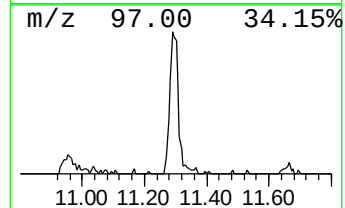
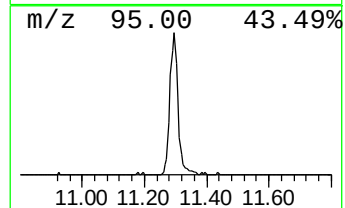
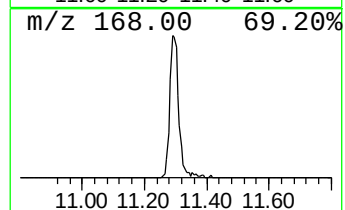
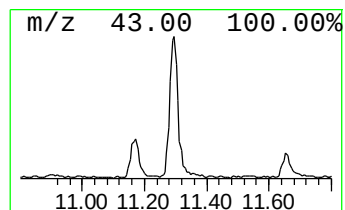
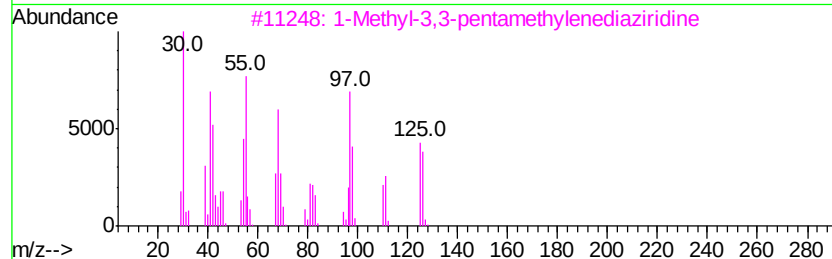
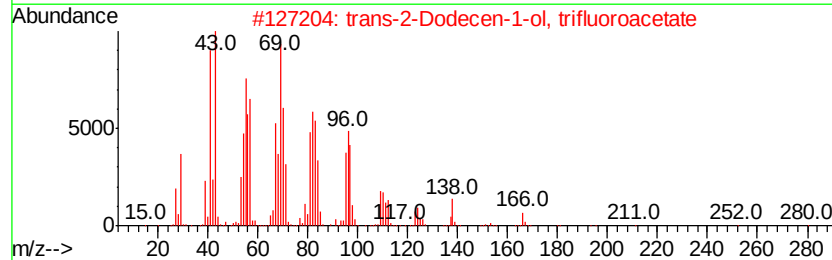
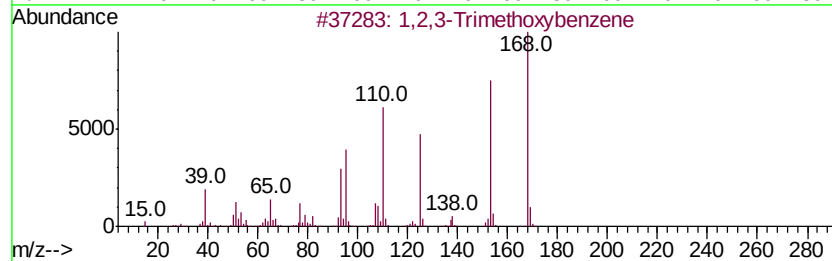
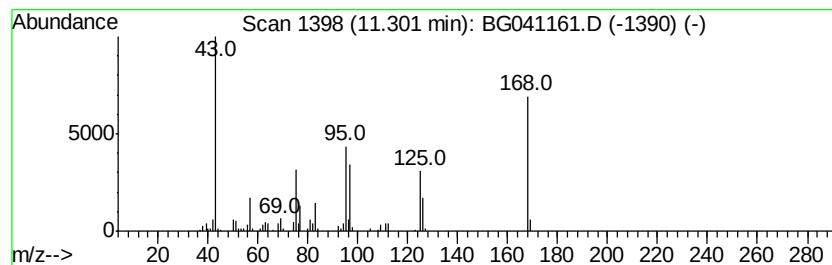
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 unknown11.30 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.30	10.05 ng	211983	Naphthalene-d8	10.93

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,2,3-Trimethoxybenzene	168	C9H12O3	000634-36-6	37
2		trans-2-Dodecen-1-ol, trifluoroa...	280	C14H23F3O2	1000352-26-2	11
3		1-Methyl-3,3-pentamethylenediazi...	126	C7H14N2	026177-34-4	10
4		3,5-Dimethoxybenzyl alcohol	168	C9H12O3	000705-76-0	10
5		6-Dodecanol acetate	228	C14H28O2	060826-23-5	10



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG061019\
Data File : BG041161.D
Acq On : 10 Jun 2019 12:18
Operator : HP/JU
Sample : K3223-02
Misc :
ALS Vial : 7 Sample Multiplier: 1

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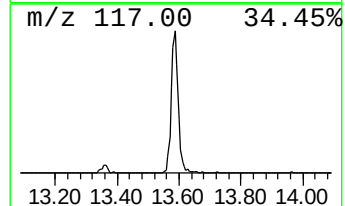
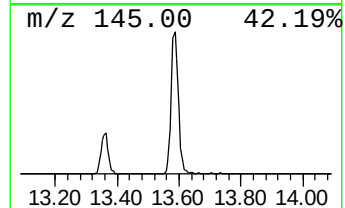
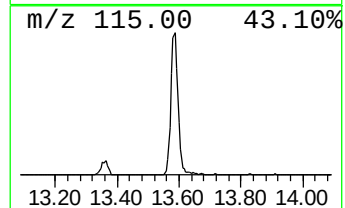
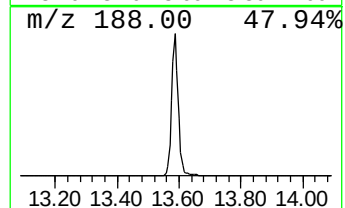
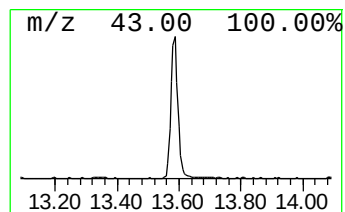
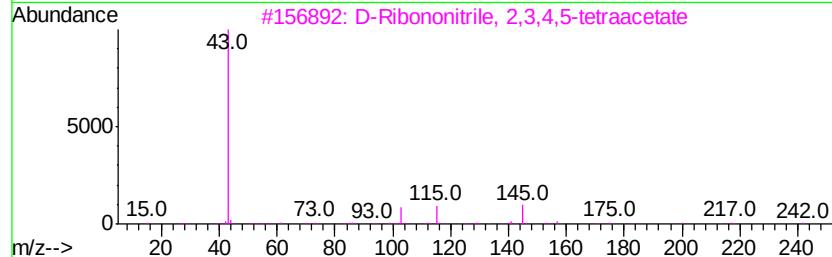
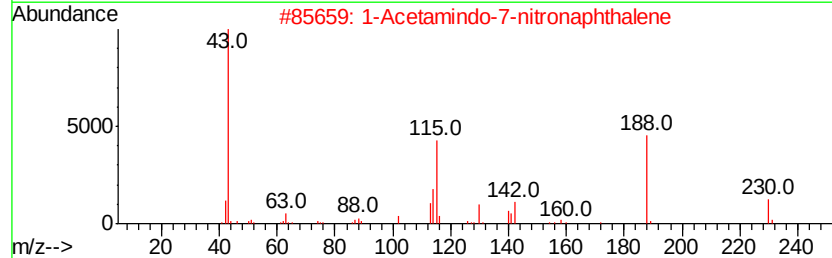
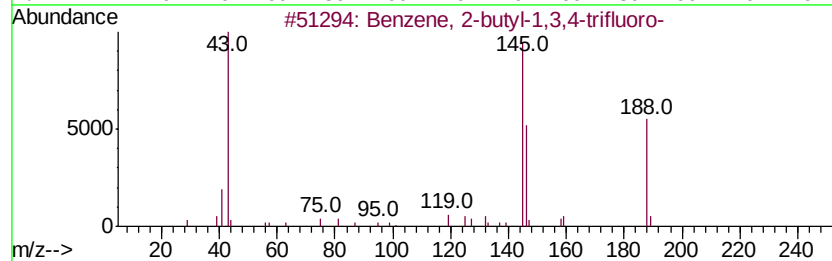
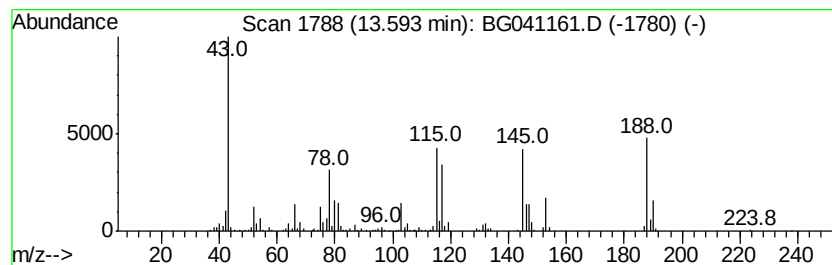
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 unknown13.59 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.59	21.67 ng	621090	Acenaphthene-d10	14.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-butyl-1,3,4-trifluoro-	188	C10H11F3	022872-43-1	22
2		1-Acetamido-7-nitronaphthalene	230	C12H10N2O3	102877-10-1	17
3		D-Ribonitrile, 2,3,4,5-tetraac...	315	C13H17N08	025546-50-3	10
4		Lyxonitrile, 2,3,4,5-tetraacet...	315	C13H17N08	034360-56-0	10
5		Cyclopropanecarboxylic acid, 2-p...	190	C12H14O2	1000221-86-7	9



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG061019\
Data File : BG041161.D
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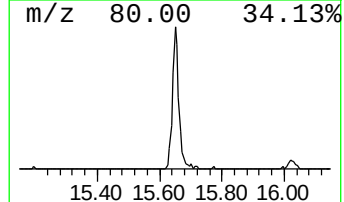
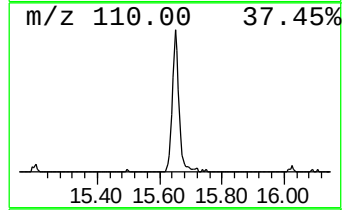
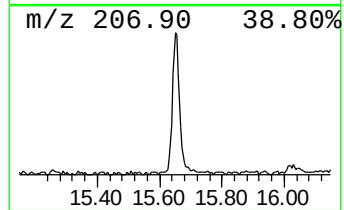
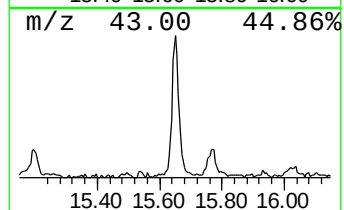
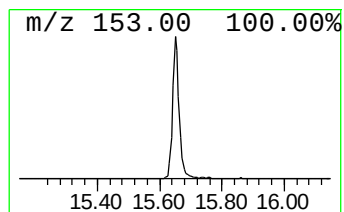
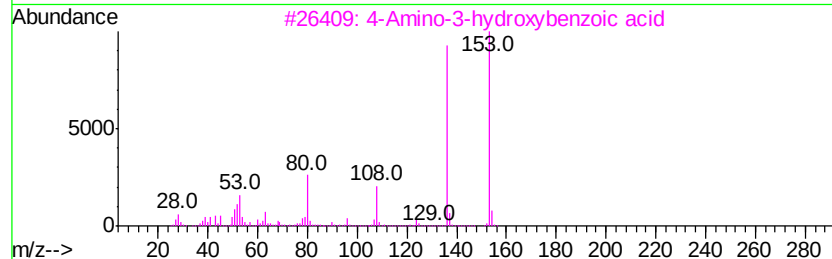
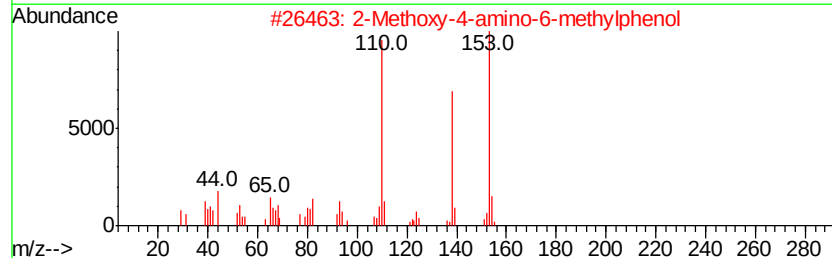
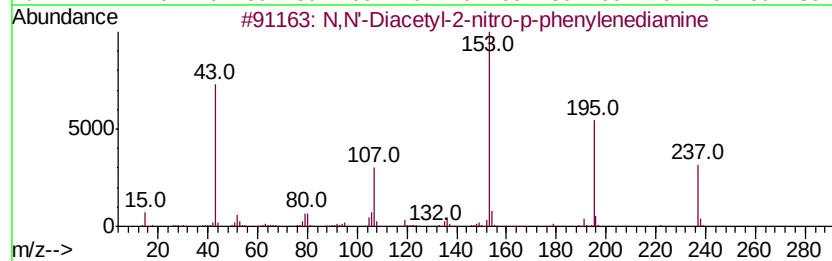
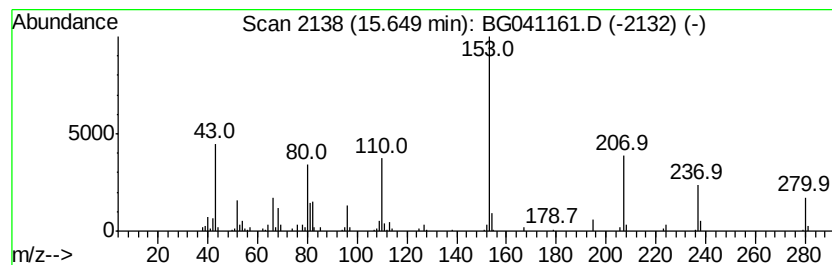
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 unknown15.65 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.65	8.35 ng	239335	Acenaphthene-d10	14.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	N,N'-Diacetyl-2-nitro-p-phenylen...	237	C10H11N3O4	005345-53-9	38
2		2-Methoxy-4-amino-6-methylphenol	153	C8H11NO2	1000289-26-0	38
3		4-Amino-3-hydroxybenzoic acid	153	C7H7NO3	002374-03-0	32
4		1,4-Benzenediamine, 2-nitro-	153	C6H7N3O2	005307-14-2	25
5		1,3-Benzenediamine, 5-nitro-	153	C6H7N3O2	005042-55-7	23



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG061019\
Data File : BG041161.D
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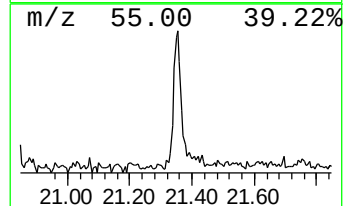
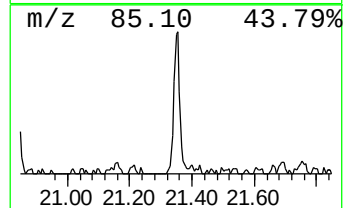
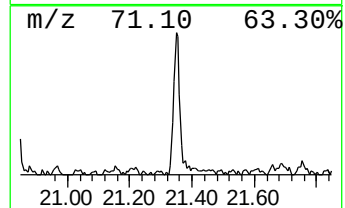
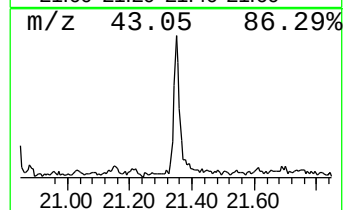
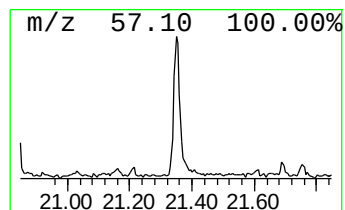
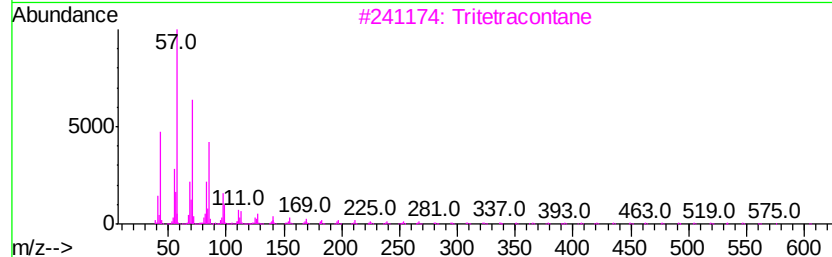
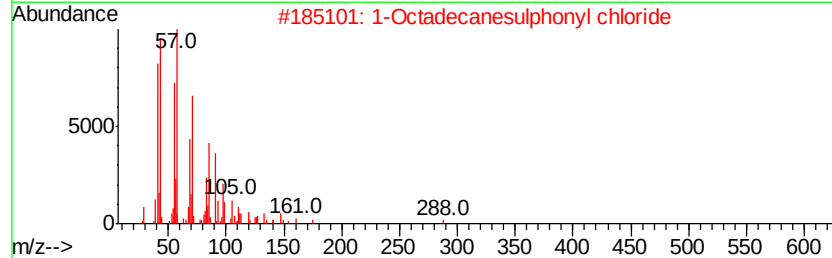
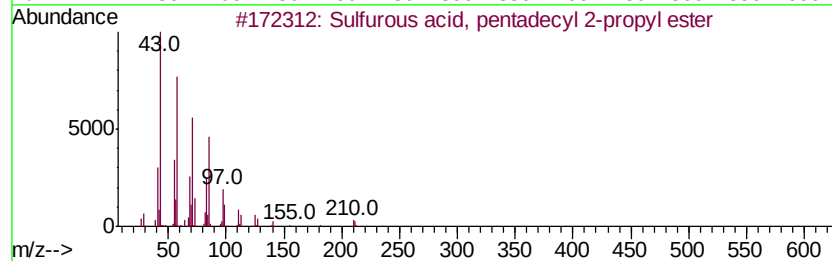
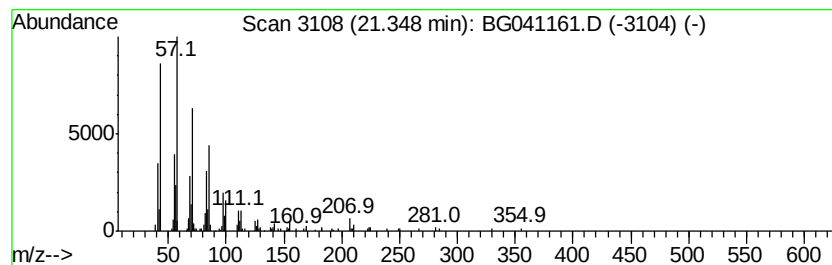
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Sulfurous acid, pentadecyl ... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.35	3.57 ng	155427	Chrysene-d12	21.76

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Sulfurous acid, pentadecyl 2-pro...	334	C18H38O3S	1000309-12-6	91
2		1-Octadecanesulphonyl chloride	352	C18H37ClO2S	1000342-70-4	91
3		Tritetracontane	605	C43H88	007098-21-7	91
4		Sulfurous acid, butyl octadecyl ...	390	C22H46O3S	1000309-18-5	91
5		Tetracosane, 1-bromo-	416	C24H49Br	006946-24-3	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG061019\
Data File : BG041161.D
Acq On : 10 Jun 2019 12:18
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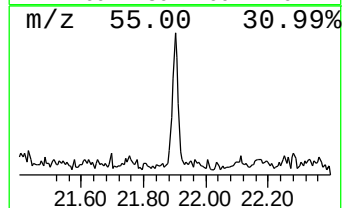
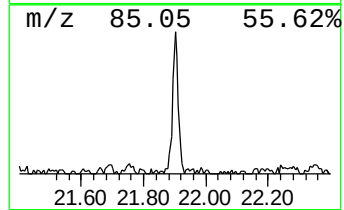
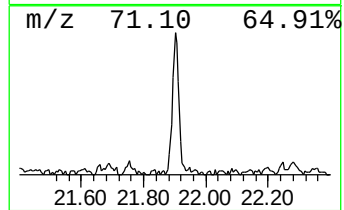
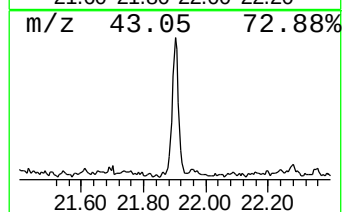
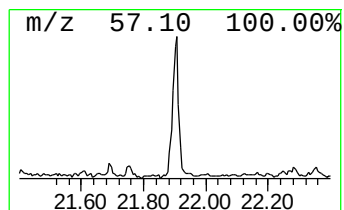
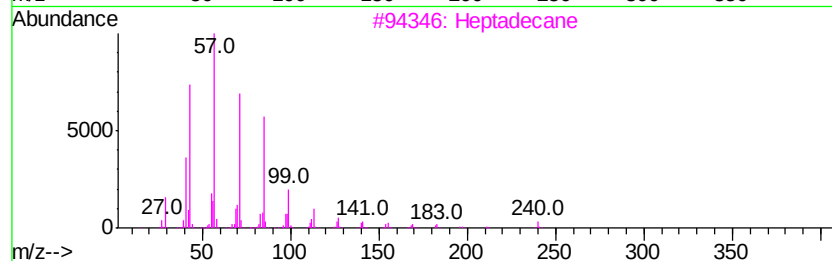
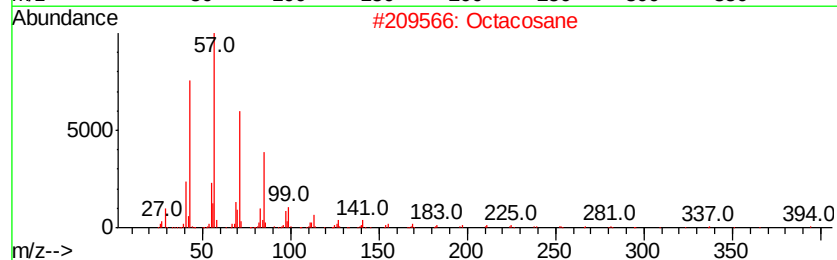
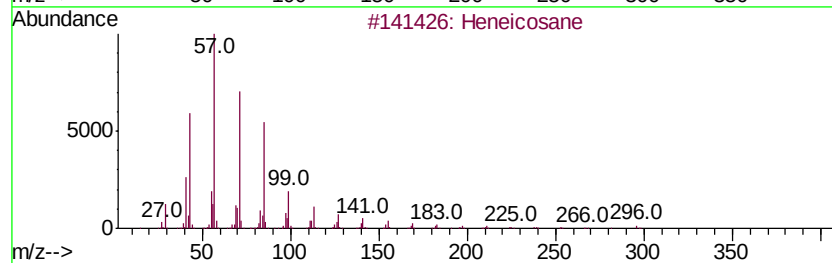
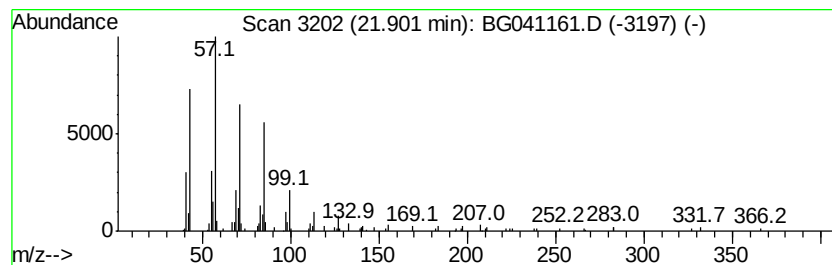
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Heneicosane Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.90	2.58 ng	112366	Chrysene-d12	21.76

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	87
2		Octacosane	394	C28H58	000630-02-4	87
3		Heptadecane	240	C17H36	000629-78-7	86
4		Heptadecane, 4-propyl-	282	C20H42	055044-10-5	86
5		Tetracosane	338	C24H50	000646-31-1	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG061019\
Data File : BG041161.D
Acq On : 10 Jun 2019 12:18
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Misc :
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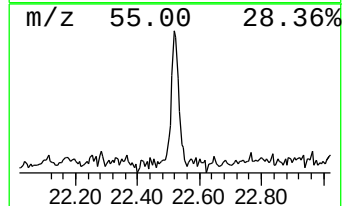
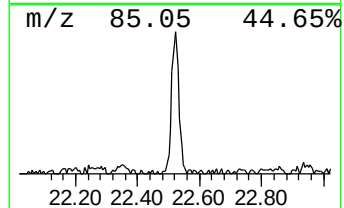
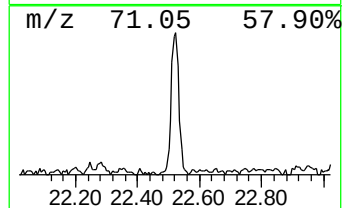
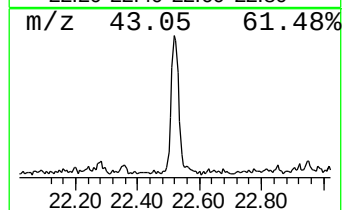
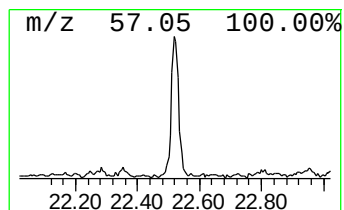
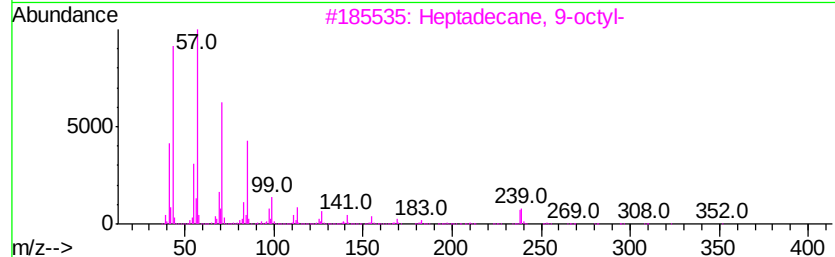
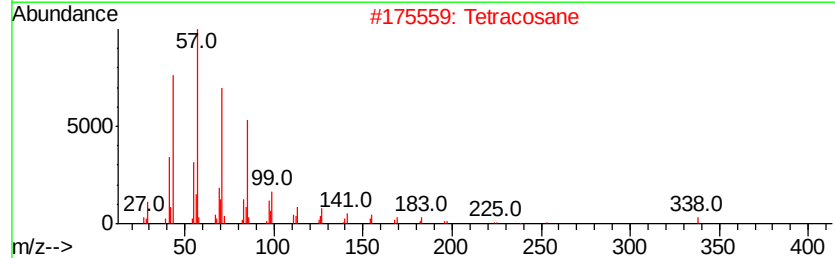
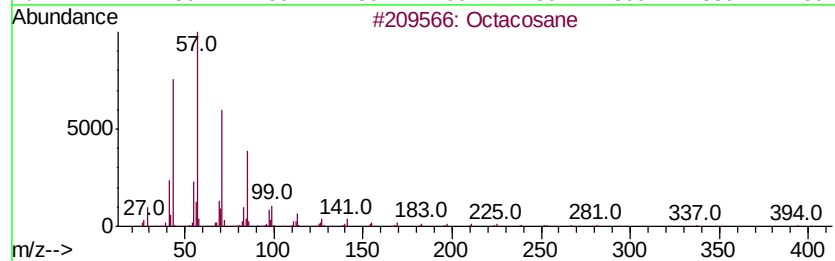
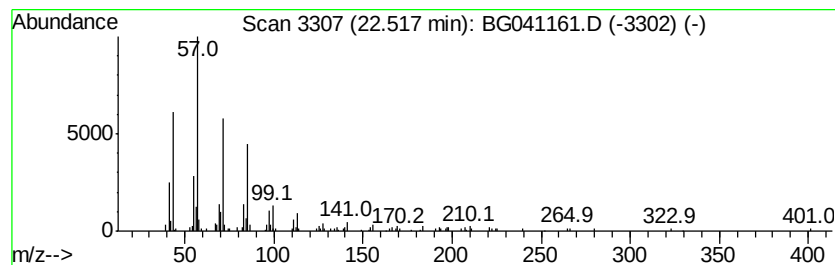
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Octacosane Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.52	3.29 ng	143355	Chrysene-d12	21.76

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosane	394	C28H58	000630-02-4	90
2		Tetracosane	338	C24H50	000646-31-1	90
3		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	87
4		Hexacosane	366	C26H54	000630-01-3	87
5		Eicosane, 7-hexyl-	366	C26H54	055333-99-8	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG061019\
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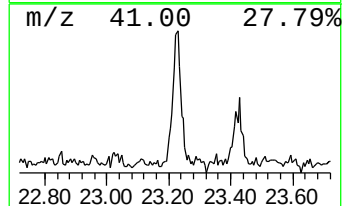
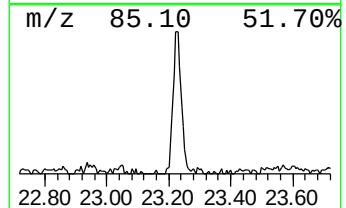
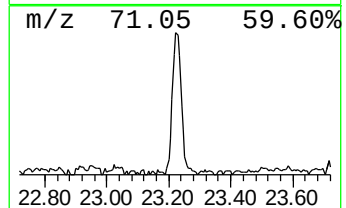
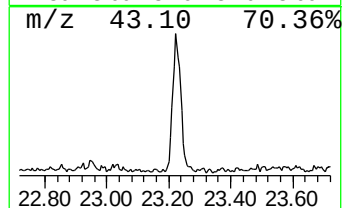
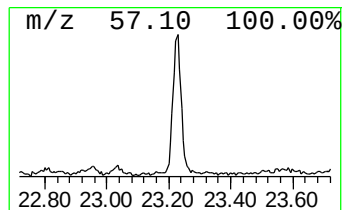
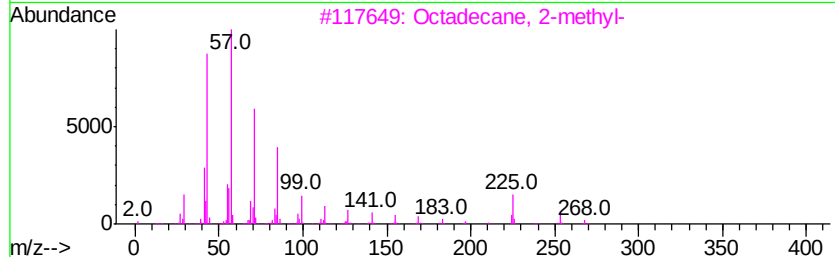
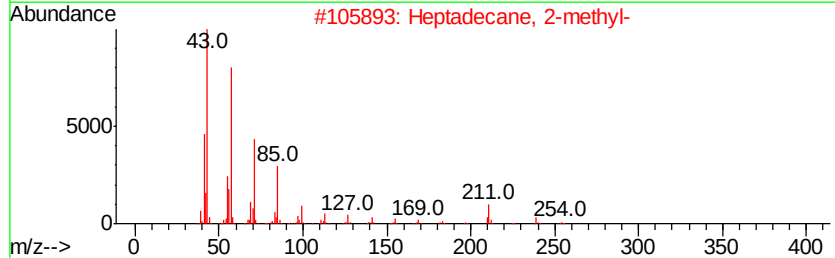
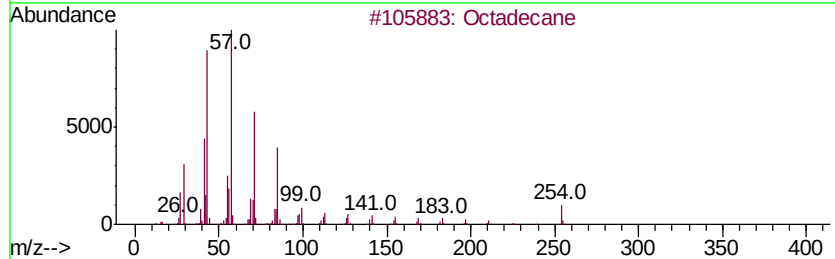
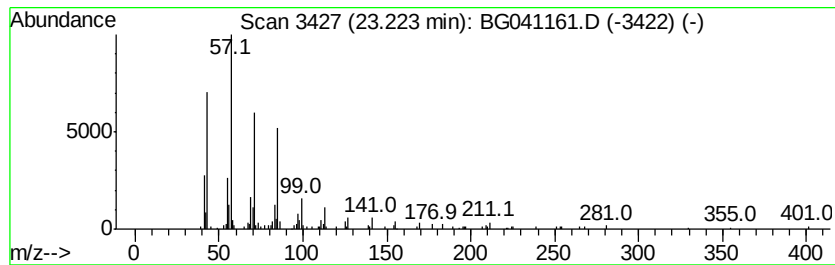
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 Octadecane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.22	3.76 ng	163421	Chrysene-d12	21.76

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecane	254	C18H38	000593-45-3	93
2			Heptadecane, 2-methyl-	254	C18H38	001560-89-0	93
3			Octadecane, 2-methyl-	268	C19H40	001560-88-9	89
4			Eicosane, 7-hexyl-	366	C26H54	055333-99-8	87
5			Hexacosane	366	C26H54	000630-01-3	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG061019\
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Misc :
ALS Vial : 7 Sample Multiplier: 1

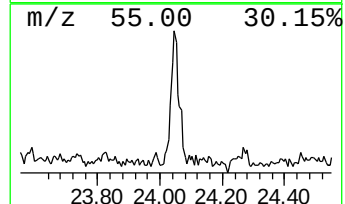
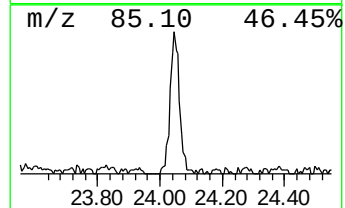
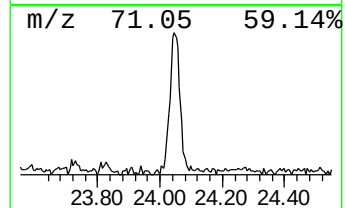
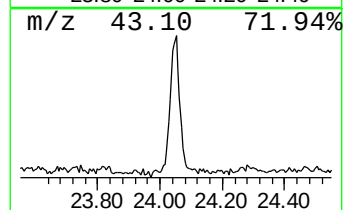
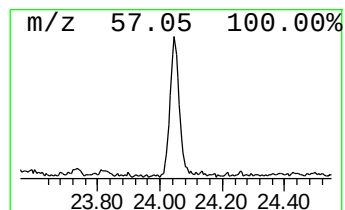
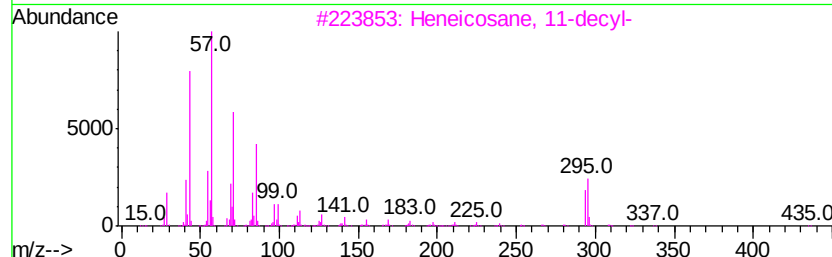
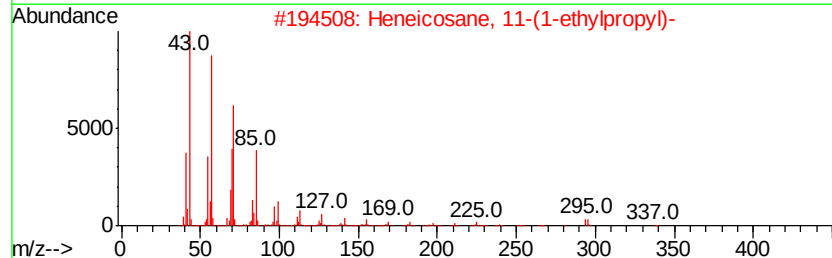
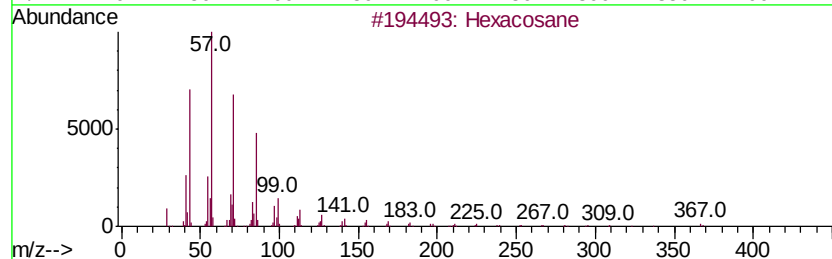
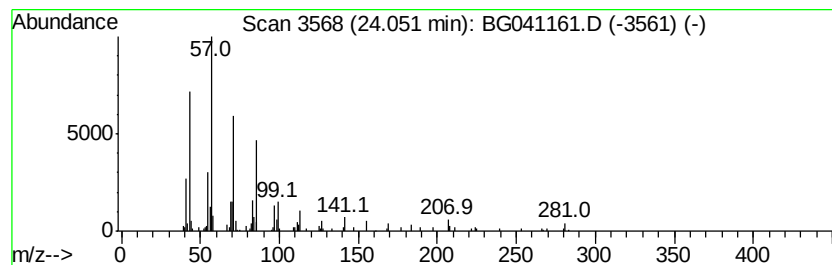
Instrument :
BNA_G
ClientSampleId :
SS043MW014-190605

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG052919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Hexacosane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.05	3.77 ng	151916	Perylene-d12	25.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Hexacosane	366 C26H54	000630-01-3	87
2	Heneicosane, 11-(1-ethylpropyl)-	366 C26H54	055282-11-6	86
3	Heneicosane, 11-decyl-	437 C31H64	055320-06-4	86
4	Heptadecane	240 C17H36	000629-78-7	83
5	Eicosane	282 C20H42	000112-95-8	80



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG061019\
 Data File : BG041161.D
 Acq On : 10 Jun 2019 12:18
 Operator : HP/JU
 Sample : K3223-02
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SS043MW014-190605

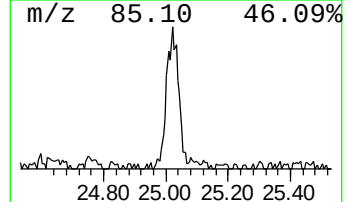
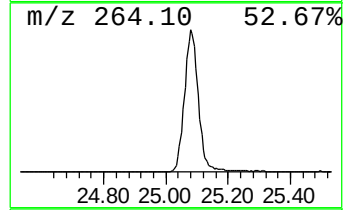
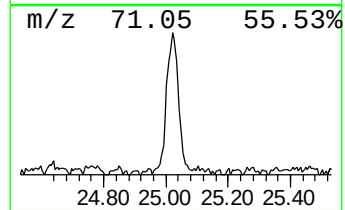
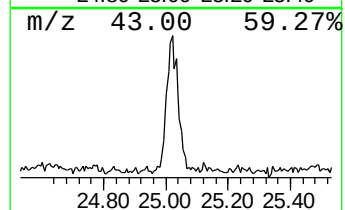
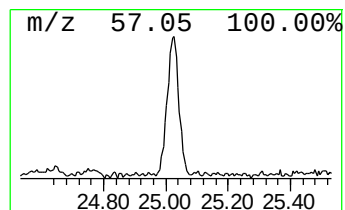
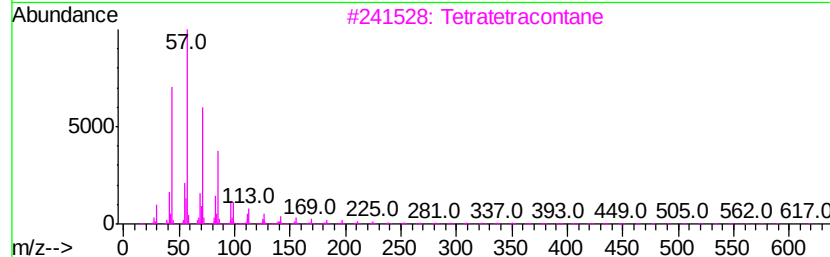
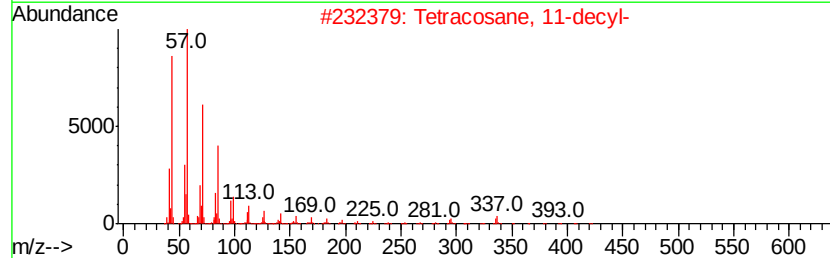
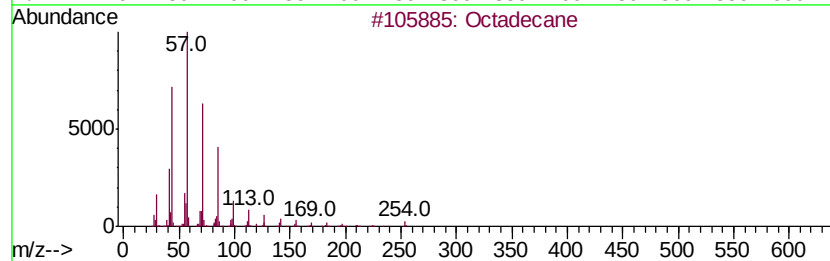
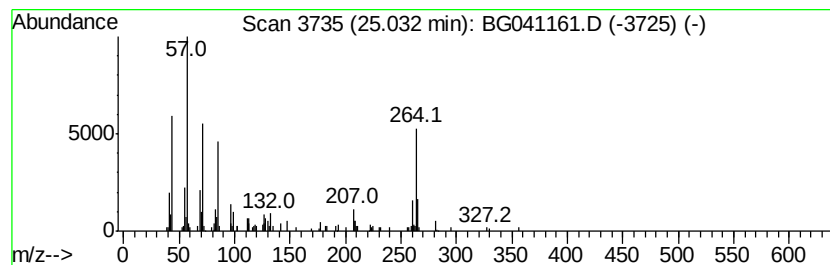
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 16 Tetracosane, 11-decyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.03	3.33 ng	134118	Perylene-d12	25.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecane	254	C18H38	000593-45-3	91
2		Tetracosane, 11-decyl-	479	C34H70	055429-84-0	60
3		Tetratetracontane	619	C44H90	007098-22-8	60
4		2-methyloctacosane	408	C29H60	1000376-72-8	53
5		Sulfurous acid, butyl tetradecyl...	334	C18H38O3S	1000309-18-1	53



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG061019\
Data File : BG041161.D
Acq On : 10 Jun 2019 12:18
Operator : HP/JU
Sample : K3223-02
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SS043MW014-190605

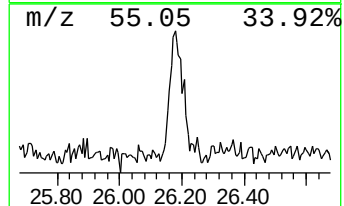
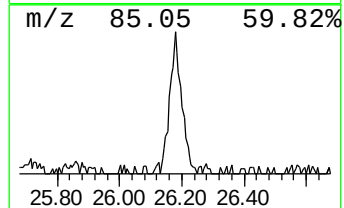
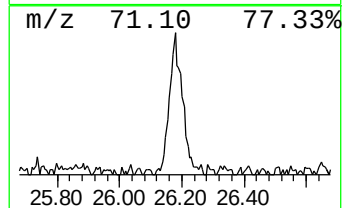
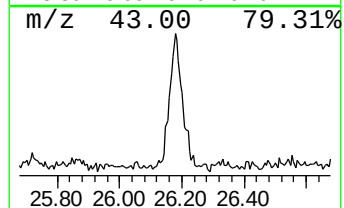
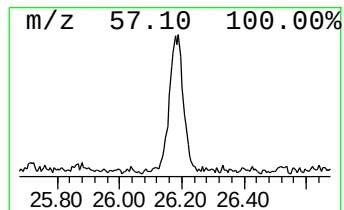
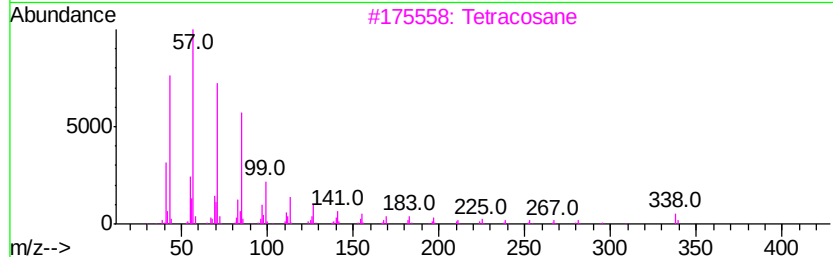
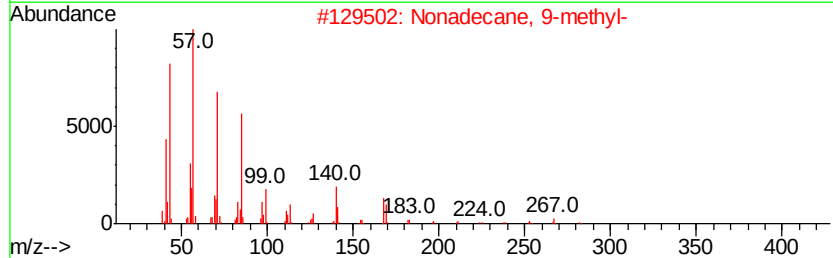
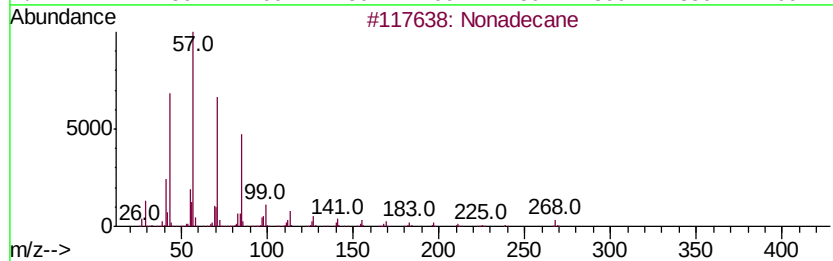
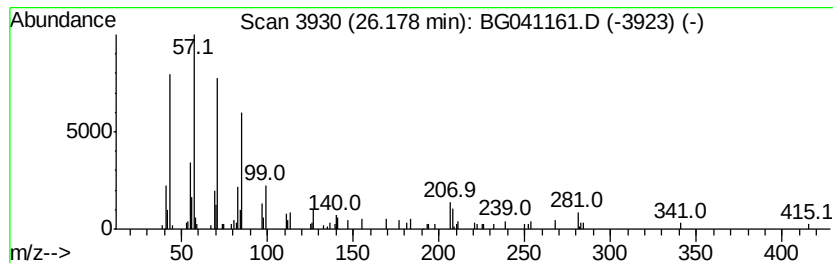
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG052919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Nonadecane Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.18	3.04 ng	122321	Perylene-d12	25.08

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonadecane	268	C19H40	000629-92-5	93
2			Nonadecane, 9-methyl-	282	C20H42	013287-24-6	90
3			Tetracosane	338	C24H50	000646-31-1	83
4			Octacosane	394	C28H58	000630-02-4	83
5			Hexacosane	366	C26H54	000630-01-3	74



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG061019\
 Data File : BG041161.D
 Acq On : 10 Jun 2019 12:18
 Operator : HP/JU
 Sample : K3223-02
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SS043MW014-190605

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG052919.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...	3.19	20.6	ng	271168	1	8.10	262853	20.0
2-Pentanone, 4-hy...	5.20	531.4	ng	6983730	1	8.10	262853	20.0
unknown7.63	7.63	40.1	ng	527209	1	8.10	262853	20.0
unknown8.92	8.92	12.9	ng	169418	1	8.10	262853	20.0
unknown11.17	11.17	5.2	ng	109448	2	10.93	421769	20.0
unknown11.30	11.30	10.1	ng	211983	2	10.93	421769	20.0
unknown13.59	13.59	21.7	ng	621090	3	14.74	573323	20.0
unknown15.65	15.65	8.3	ng	239335	3	14.74	573323	20.0
Sulfurous acid, p...	21.35	3.6	ng	155427	5	21.76	870342	20.0
Heneicosane	21.90	2.6	ng	112366	5	21.76	870342	20.0
Octacosane	22.52	3.3	ng	143355	5	21.76	870342	20.0
Octadecane	23.22	3.8	ng	163421	5	21.76	870342	20.0
Hexacosane	24.05	3.8	ng	151916	6	25.08	804966	20.0
Tetracosane, 11-d...	25.03	3.3	ng	134118	6	25.08	804966	20.0
Nonadecane	26.18	3.0	ng	122321	6	25.08	804966	20.0