

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010419\
 Data File : BG038939.D
 Acq On : 4 Jan 2019 21:06
 Operator : JU/SJ
 Sample : K1043-01
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 HB-1319-WS

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-BG121218.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	3.169	10	14	25	rVB	38187	62647	1.38%	0.399%
2	4.198	183	189	198	rBV	36748	59347	1.31%	0.378%
3	5.596	421	427	433	rVV	180909	304204	6.70%	1.940%
4	5.661	433	438	448	rVB	49721	113179	2.49%	0.722%
5	6.031	494	501	508	rBV	39585	64464	1.42%	0.411%
6	7.112	679	685	691	rBV	40132	65961	1.45%	0.421%
7	7.206	691	701	706	rVV	115268	249868	5.51%	1.593%
8	7.576	756	764	771	rBV	342180	617044	13.60%	3.934%
9	7.676	776	781	788	rVB	79878	133717	2.95%	0.853%
10	8.046	837	844	855	rBV2	93170	184339	4.06%	1.175%
11	8.152	855	862	869	rVB3	32741	65086	1.43%	0.415%
12	8.369	891	899	905	rBV	326136	595368	13.12%	3.796%
13	9.139	1023	1030	1037	rBV3	25790	47990	1.06%	0.306%
14	9.227	1037	1045	1052	rBV	278458	527453	11.62%	3.363%
15	10.243	1213	1218	1224	rVB2	42348	68001	1.50%	0.434%
16	10.484	1253	1259	1265	rBV2	27875	50010	1.10%	0.319%
17	10.860	1314	1323	1329	rBV	116708	237854	5.24%	1.517%
18	12.036	1516	1523	1531	rBV3	41972	79352	1.75%	0.506%
19	12.394	1576	1584	1591	rBV3	26603	48840	1.08%	0.311%
20	12.517	1599	1605	1611	rBV	32313	56321	1.24%	0.359%
21	12.746	1638	1644	1651	rBV3	28148	65231	1.44%	0.416%
22	13.175	1712	1717	1722	rBV2	29984	45556	1.00%	0.290%
23	13.305	1731	1739	1747	rVB	791173	1313664	28.95%	8.376%
24	14.121	1873	1878	1883	rVB2	52056	90220	1.99%	0.575%
25	14.685	1964	1974	1980	rBV2	198378	341617	7.53%	2.178%
26	16.178	2222	2228	2235	rBV	640751	1011145	22.28%	6.447%
27	16.348	2253	2257	2258	rBV	42339	50688	1.12%	0.323%
28	16.806	2331	2335	2343	rBV	128959	187755	4.14%	1.197%
29	17.435	2437	2442	2453	rVB2	242787	418940	9.23%	2.671%
30	18.334	2584	2595	2606	rBV	2159676	4537519	100.00%	28.931%
31	19.562	2801	2804	2811	rBV	129621	171765	3.79%	1.095%
32	20.038	2879	2885	2890	rVB	1380388	1846166	40.69%	11.771%
33	20.761	3005	3008	3012	rBV	70282	75713	1.67%	0.483%
34	21.712	3165	3170	3181	rVB	259241	437972	9.65%	2.792%

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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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Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	25.009	3724	3731	3747	rVB3	141859	453035	9.98%	2.889%
36	26.771	4019	4031	4045	rVB6	127611	551141	12.15%	3.514%
37	27.271	4109	4116	4126	rVB10	61769	200714	4.42%	1.280%
38	30.479	4651	4662	4674	rVB10	54308	254119	5.60%	1.620%

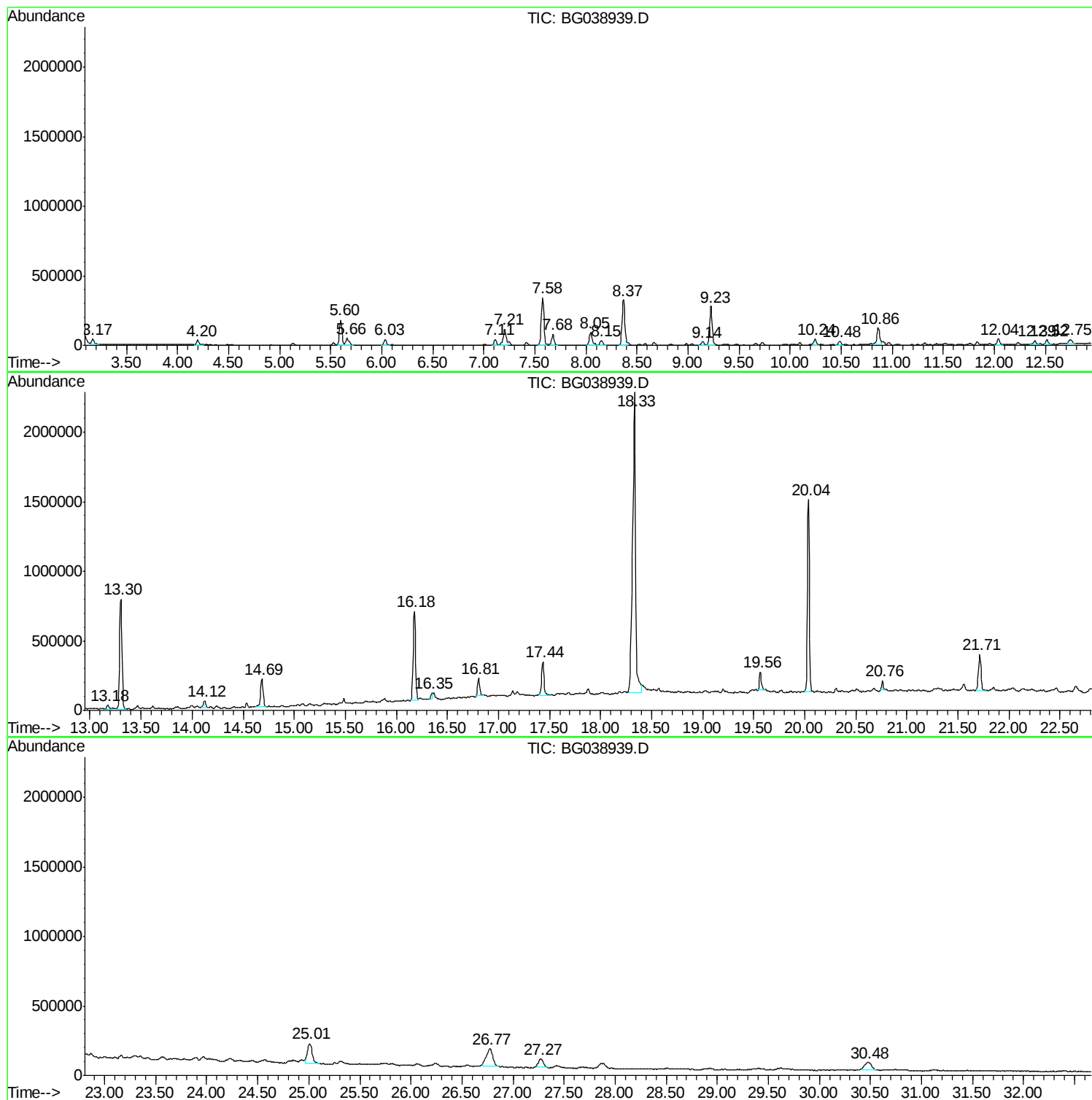
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG121218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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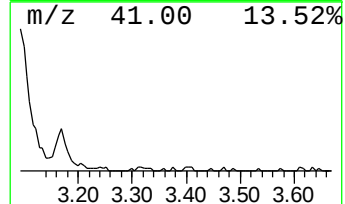
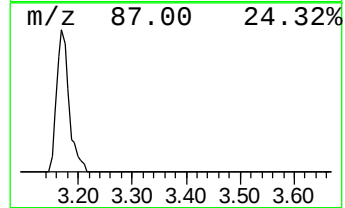
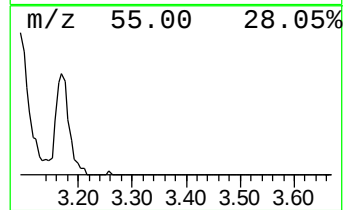
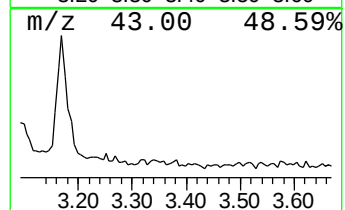
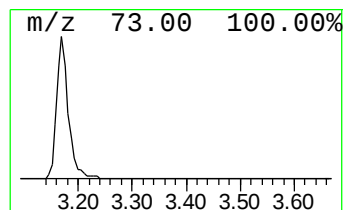
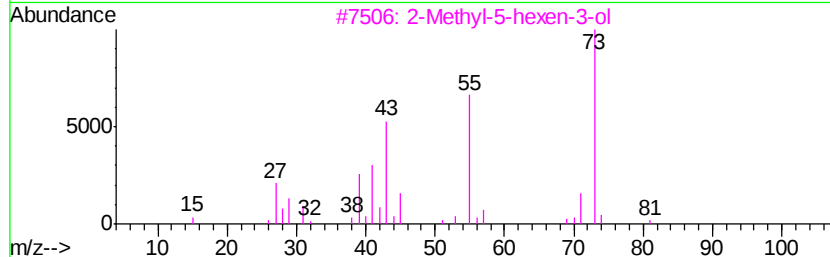
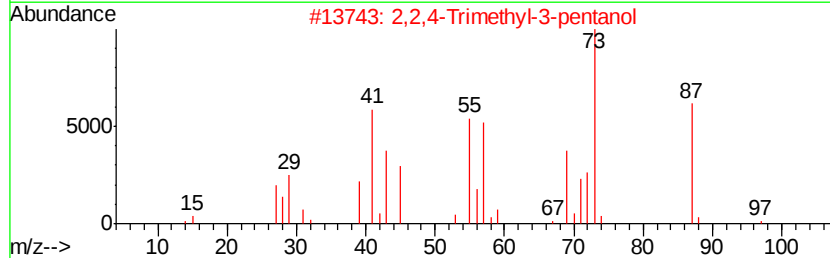
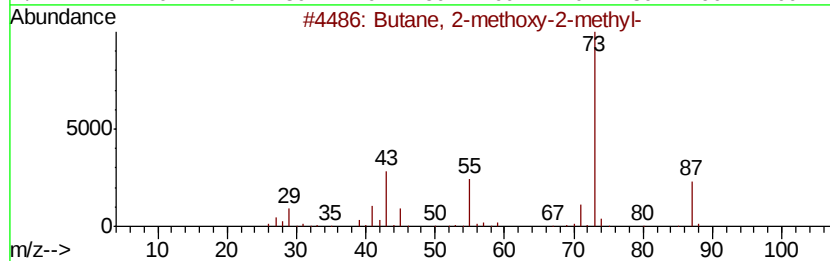
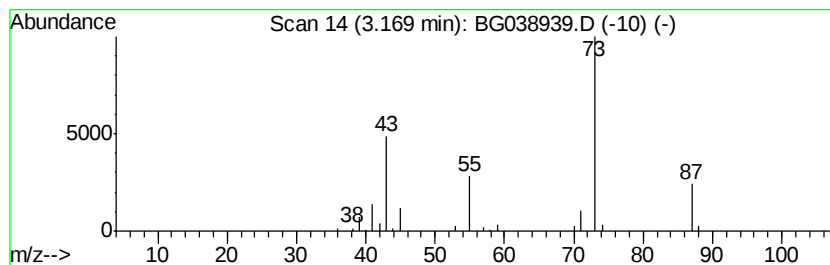
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG121218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.17	6.80 ng	62647	1,4-Dichlorobenzene-d4	8.05

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	42
3			2-Methyl-5-hexen-3-ol	114	C7H14O	032815-70-6	37
4			Silane, tetramethyl-	88	C4H12Si	000075-76-3	9
5			Silane, ethyltrimethyl-	102	C5H14Si	003439-38-1	9



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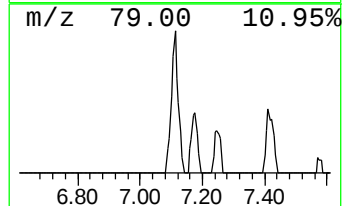
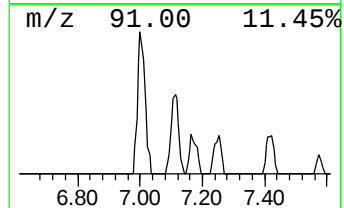
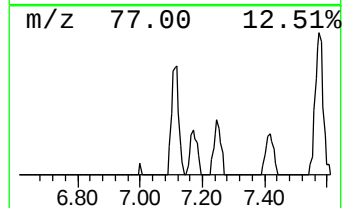
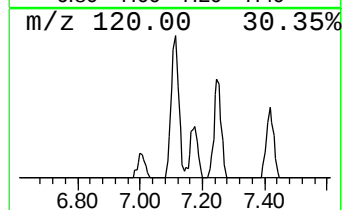
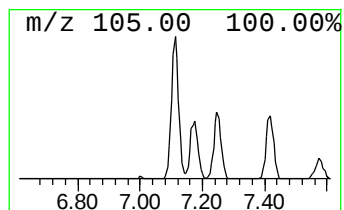
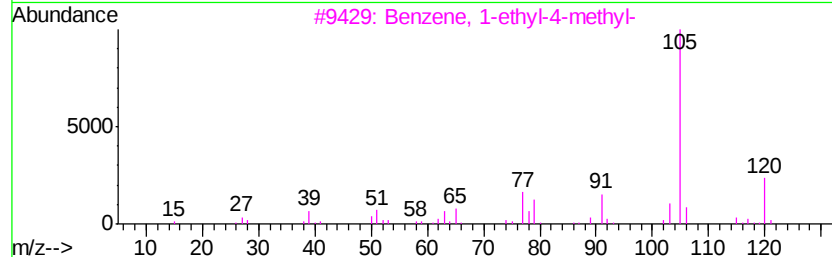
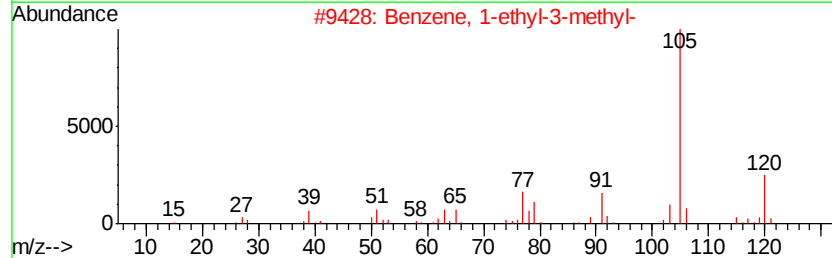
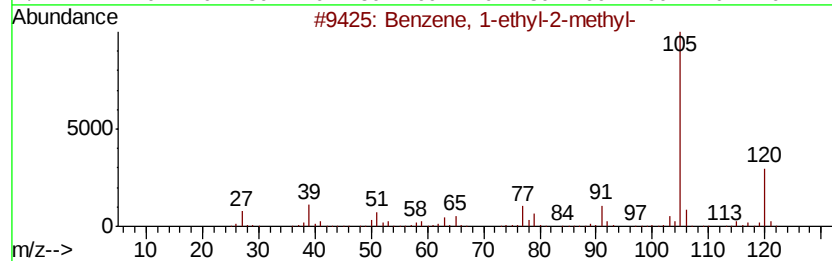
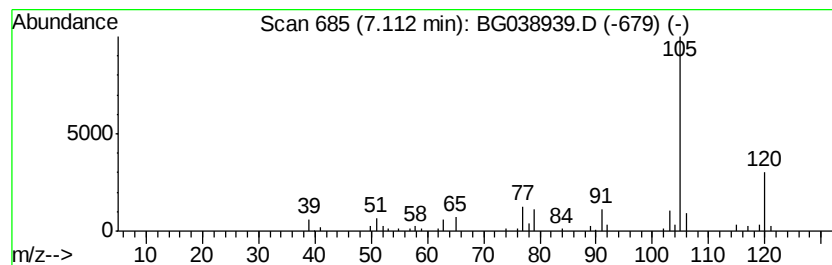
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TIC Library : C:\Database\NIST11.L
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Peak Number 5 Benzene, 1-ethyl-2-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.11	7.16 ng	65961	1,4-Dichlorobenzene-d4	8.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
3		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	94
4		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
5		Mesitylene	120	C9H12	000108-67-8	91



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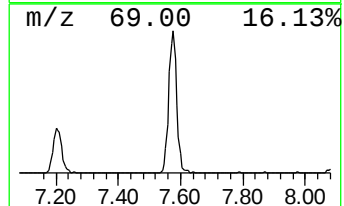
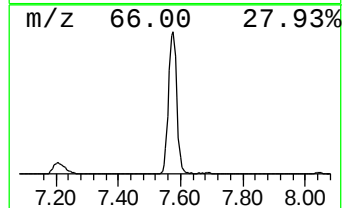
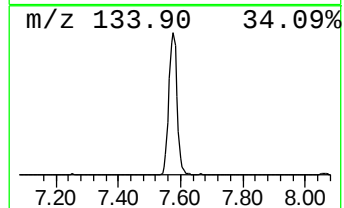
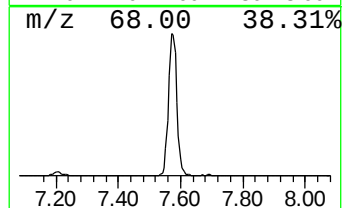
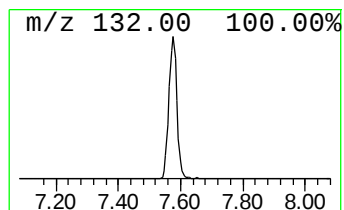
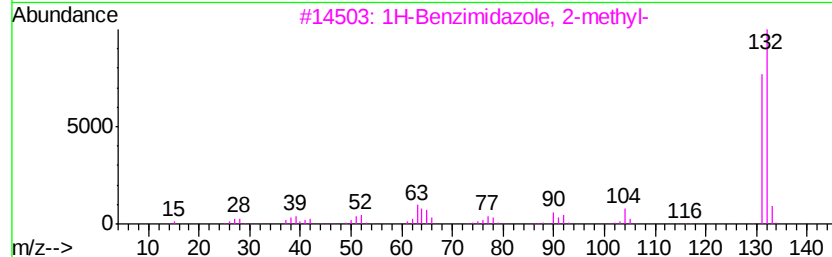
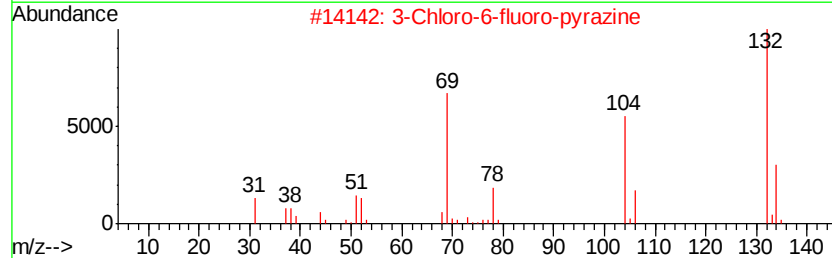
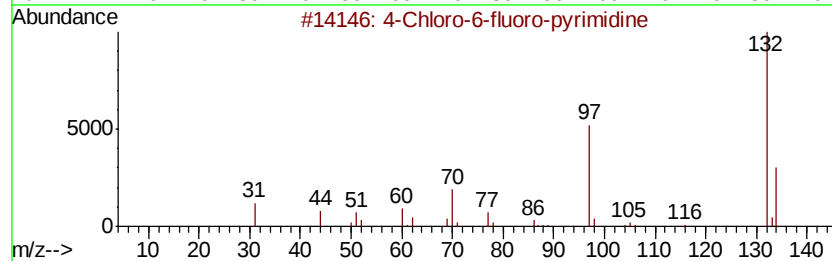
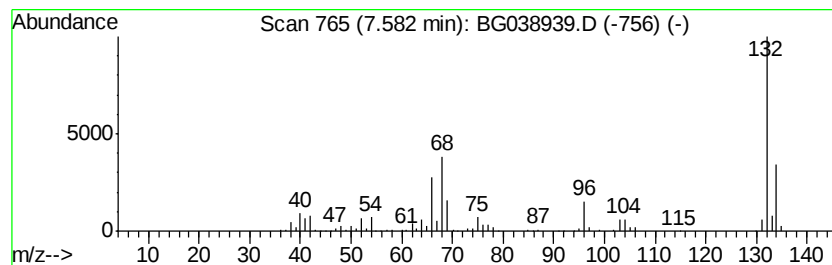
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TIC Library : C:\Database\NIST11.L
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Peak Number 6 unknown7.58 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.58	66.95 ng	617044	1,4-Dichlorobenzene-d4	8.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	35
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	27
4		Imidazole, 4,5-dicvano-1-methyl-	132	C6H4N4	019485-35-9	25
5		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	17



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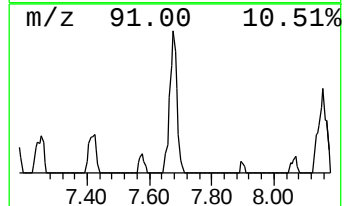
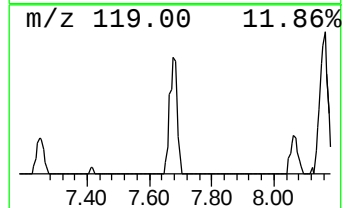
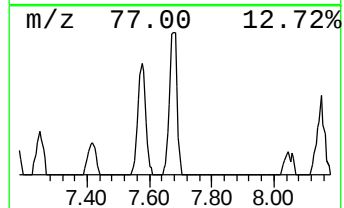
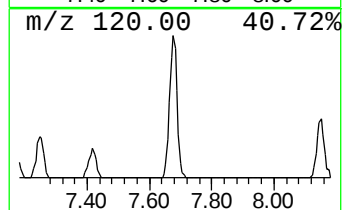
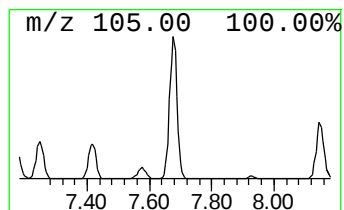
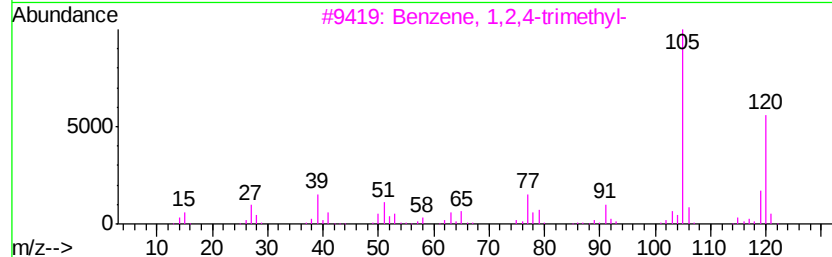
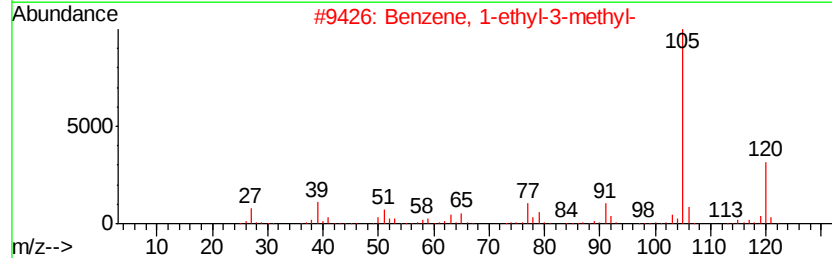
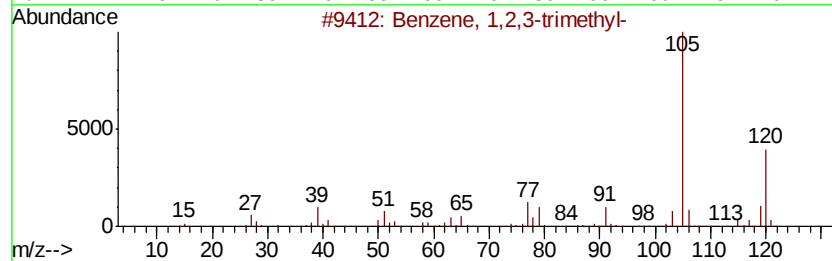
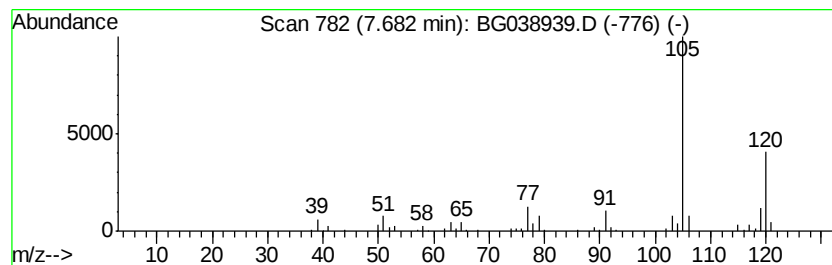
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Benzene, 1,2,3-trimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.68	14.51 ng	133717	1,4-Dichlorobenzene-d4	8.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
2		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
3		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
4		Mesitylene	120	C9H12	000108-67-8	91
5		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	90



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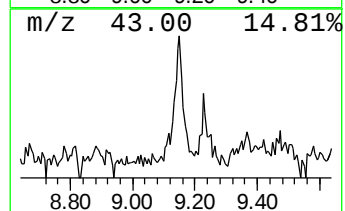
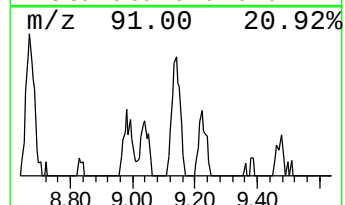
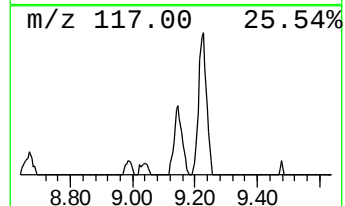
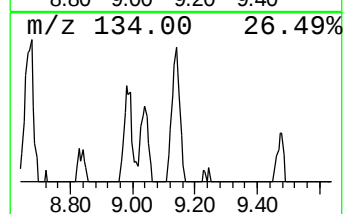
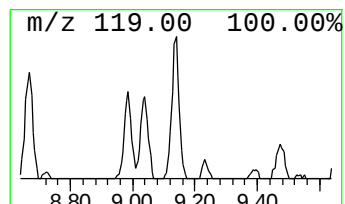
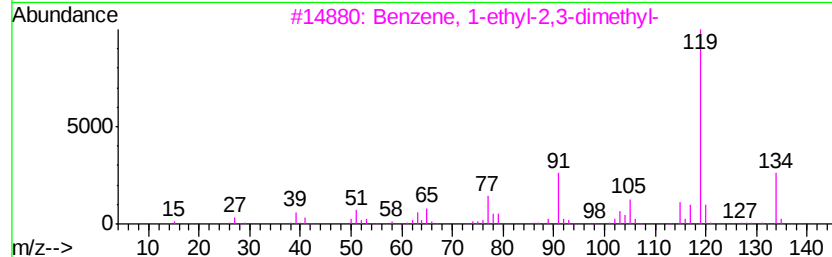
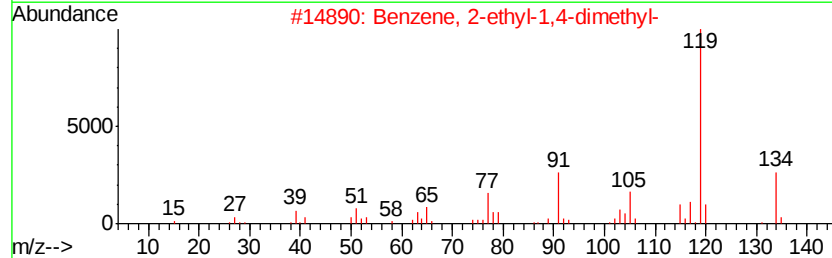
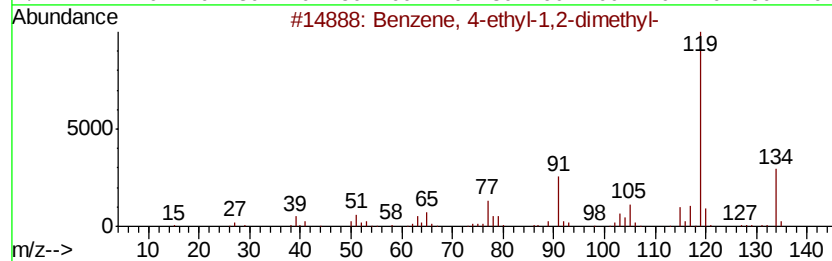
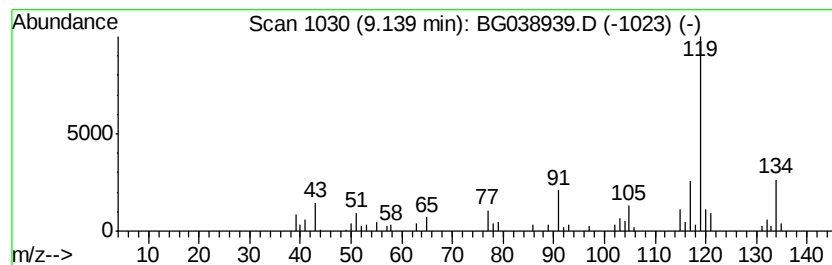
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.14	5.21 ng	47990	1,4-Dichlorobenzene-d4	8.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	93
2		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	93
3		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	93
4		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	87
5		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	81



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010419\
Data File : BG038939.D
Acq On : 4 Jan 2019 21:06
Operator : JU/SJ
Sample : K1043-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
HB-1319-WS

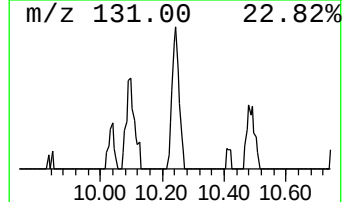
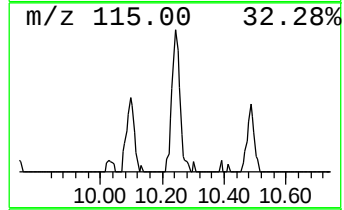
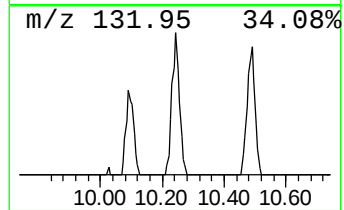
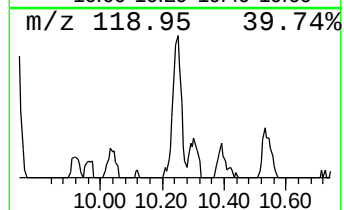
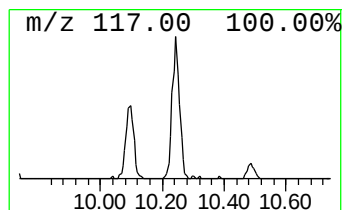
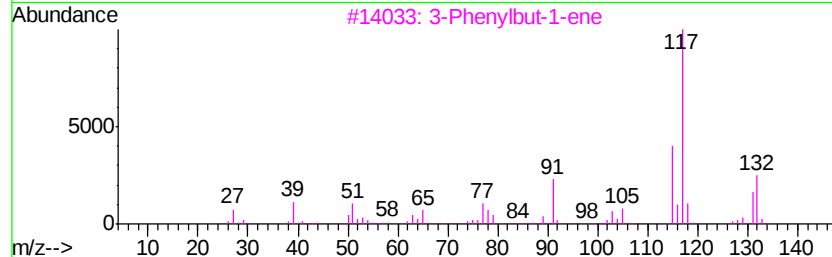
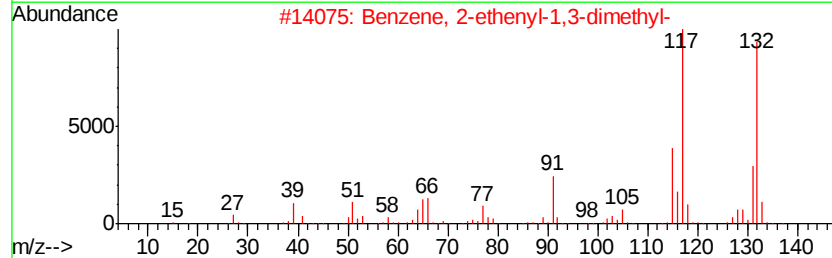
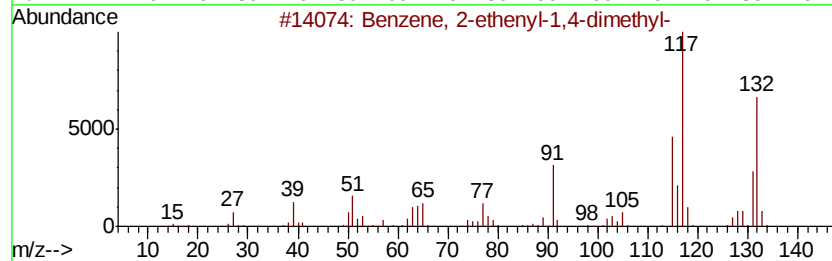
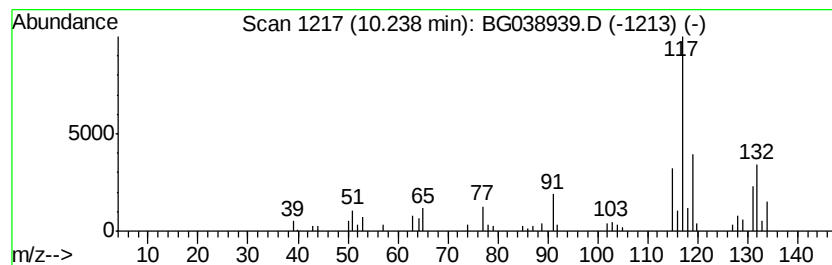
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.24	5.72 ng	68001	Naphthalene-d8	10.86

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	92
2			Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	87
3			3-Phenylbut-1-ene	132	C10H12	000934-10-1	87
4			1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	87
5			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010419\
Data File : BG038939.D
Acq On : 4 Jan 2019 21:06
Operator : JU/SJ
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Misc :
ALS Vial : 12 Sample Multiplier: 1

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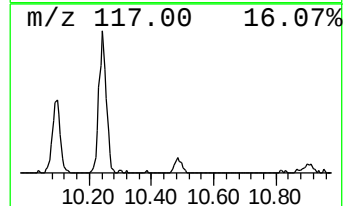
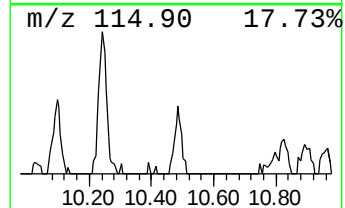
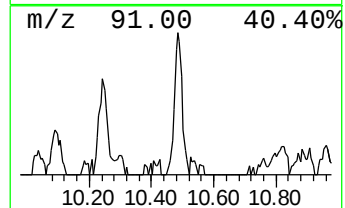
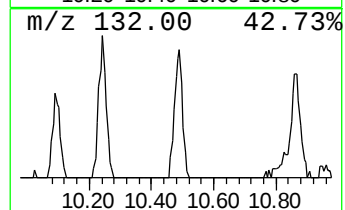
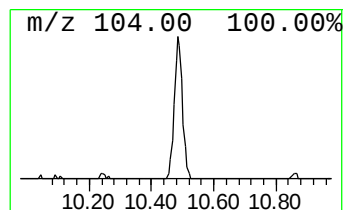
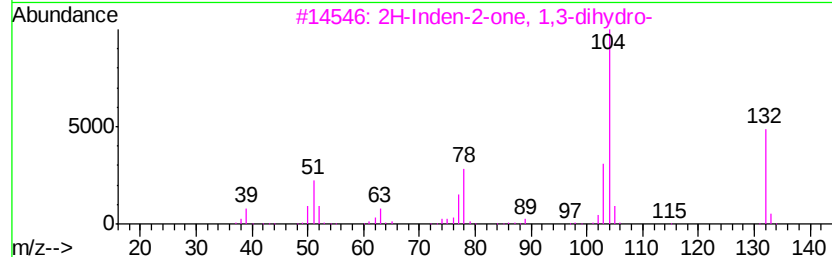
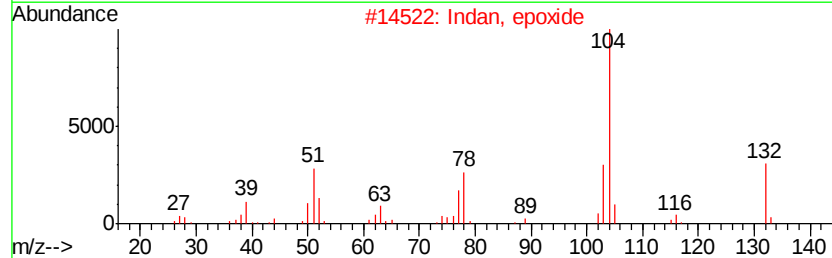
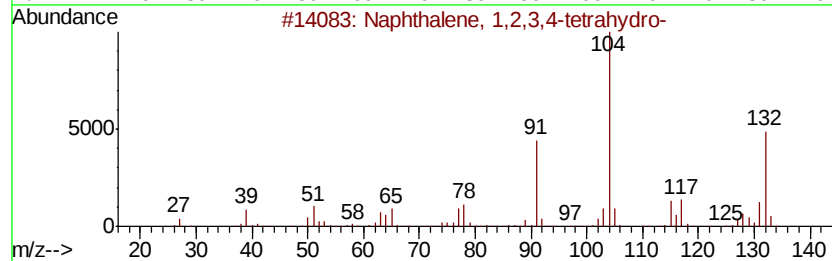
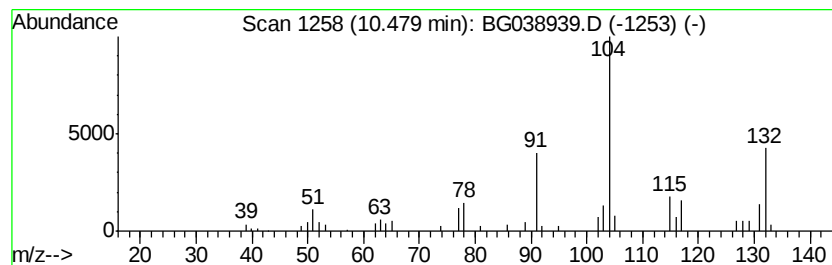
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Naphthalene, 1,2,3,4-tetrahydro- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.48	4.21 ng	50010	Naphthalene-d8	10.86

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	94
2			Indan, epoxide	132	C9H8O	000768-22-9	53
3			2H-Inden-2-one, 1,3-dihydro-	132	C9H8O	000615-13-4	50
4			2-Methylbenzyl cyanide	131	C9H9N	022364-68-7	43
5			1H-Inden-1-one, 2,3-dihydro-	132	C9H8O	000083-33-0	41



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010419\
Data File : BG038939.D
Acq On : 4 Jan 2019 21:06
Operator : JU/SJ
Sample : K1043-01
Misc :
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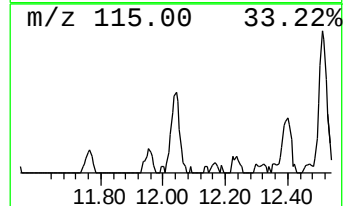
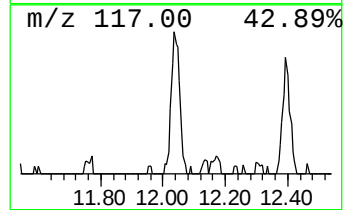
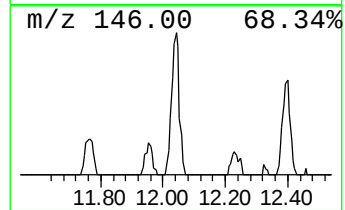
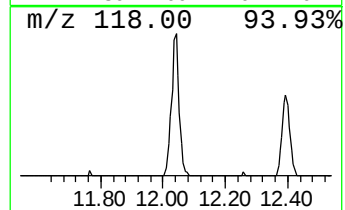
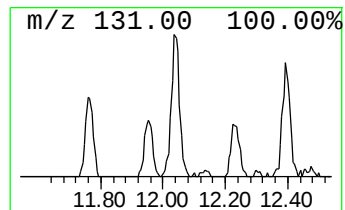
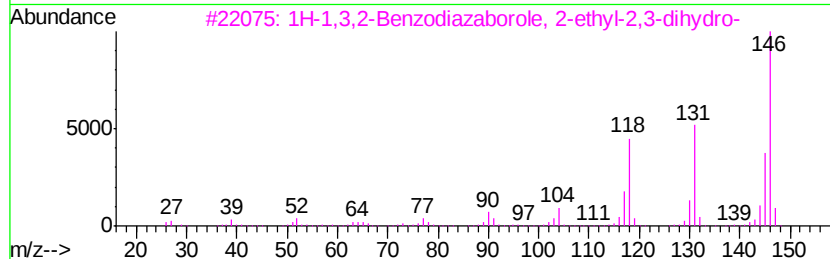
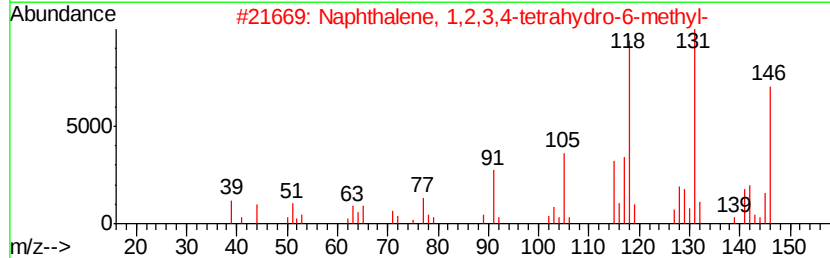
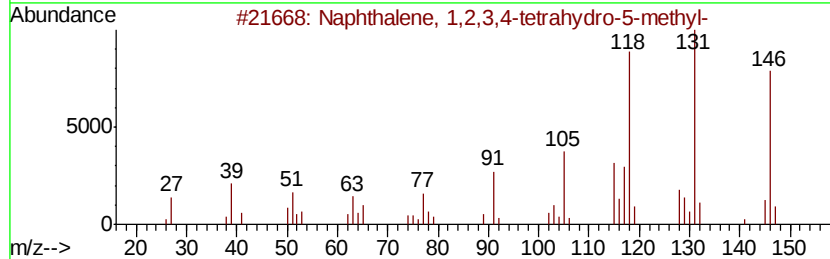
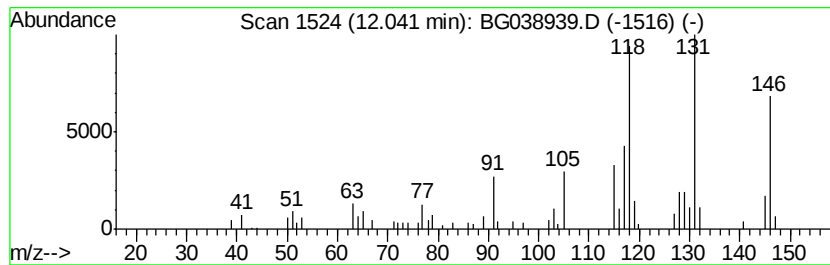
Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG121218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Naphthalene, 1,2,3,4-tetra... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.04	6.67 ng	79352	Naphthalene-d8	10.86	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	96
2	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	95
3	1H-1,3,2-Benzodiazaborole, 2-eth...	146	C8H11BN2	074663-81-3	87
4	1H-Inden-1-one, 2,3-dihydro-2-me...	146	C10H10O	017496-14-9	64
5	Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	60



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010419\
 Data File : BG038939.D
 Acq On : 4 Jan 2019 21:06
 Operator : JU/SJ
 Sample : K1043-01
 Misc :
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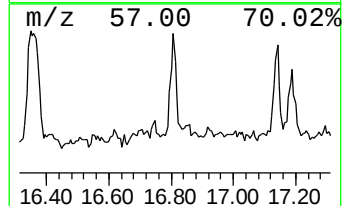
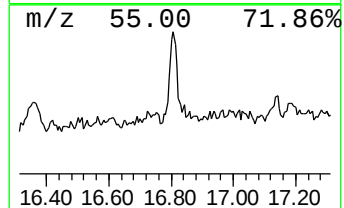
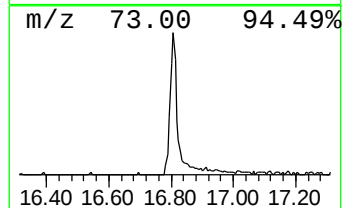
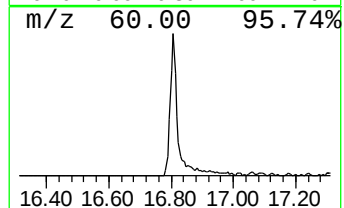
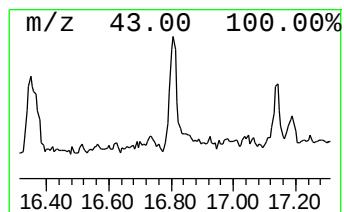
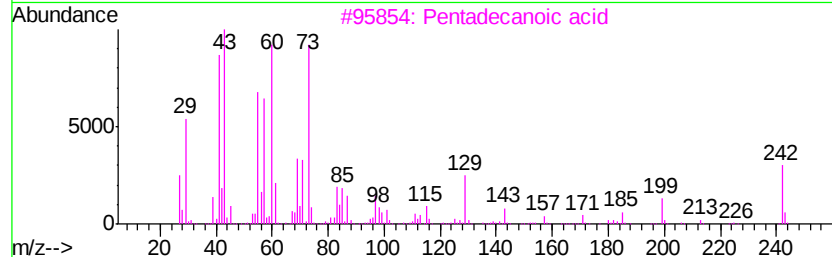
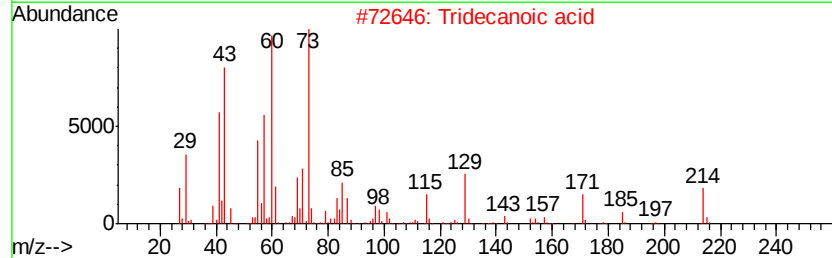
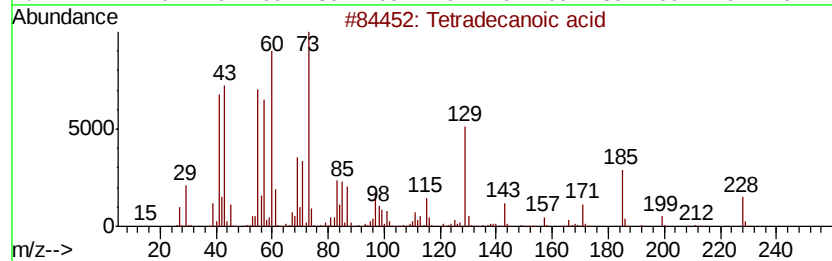
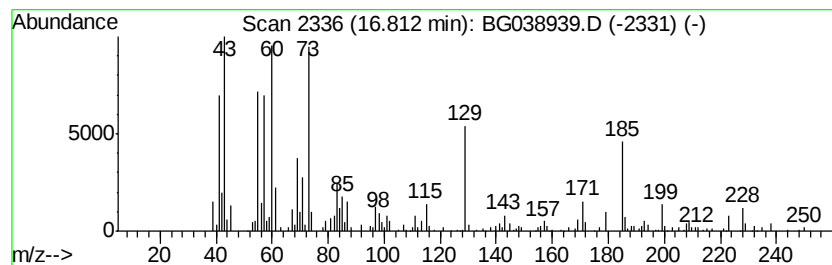
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG121218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 15 Tetradecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD		R.T.
16.81	8.96 ng	187755	Phenanthrene-d10		17.44
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetradecanoic acid	228	C14H28O2	000544-63-8	91
2	Tridecanoic acid	214	C13H26O2	000638-53-9	72
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	58
4	Dodecanoic acid	200	C12H24O2	000143-07-7	53
5	n-Decanoic acid	172	C10H20O2	000334-48-5	53



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010419\
Data File : BG038939.D
Acq On : 4 Jan 2019 21:06
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Sample : K1043-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
HB-1319-WS

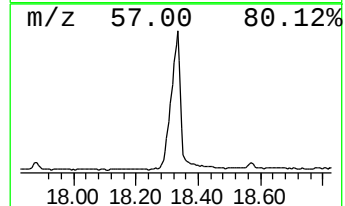
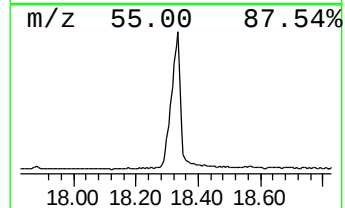
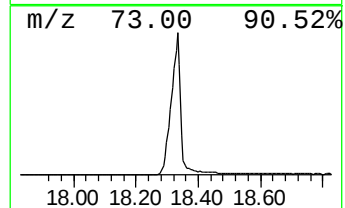
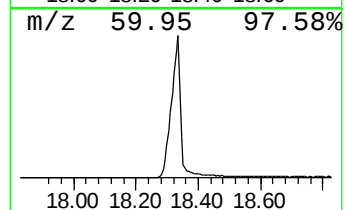
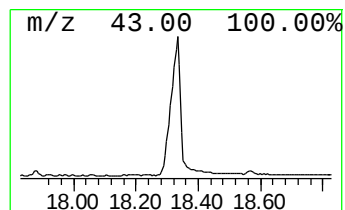
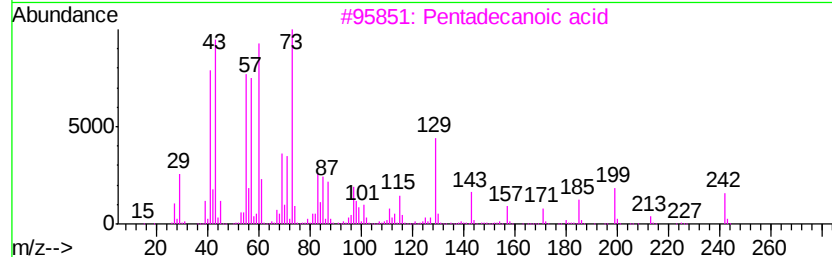
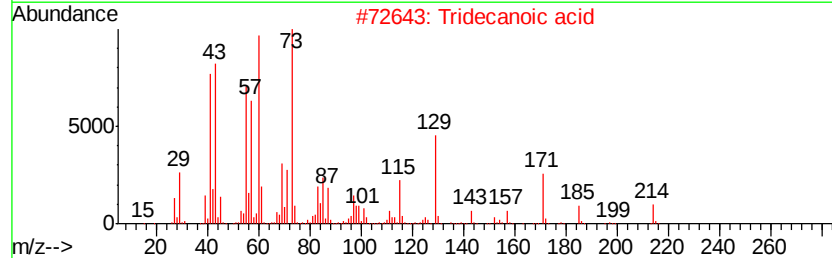
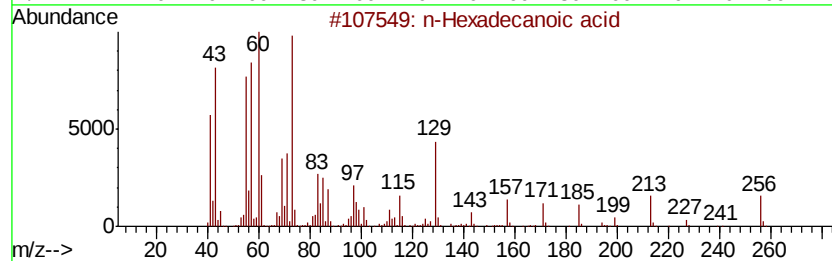
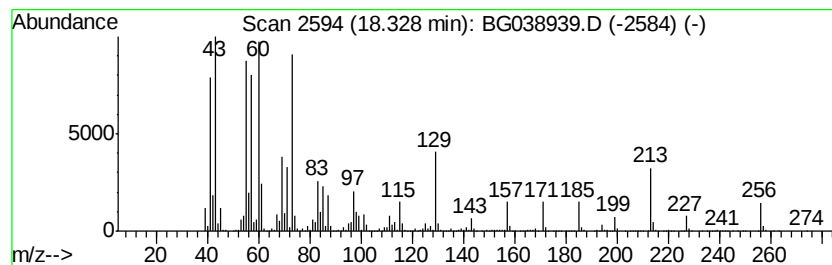
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 n-Hexadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.33	216.62 ng	4537520	Phenanthrene-d10	17.44

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	93
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	87
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	83
5		n-Decanoic acid	172	C10H20O2	000334-48-5	58



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010419\
Data File : BG038939.D
Acq On : 4 Jan 2019 21:06
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Misc :
ALS Vial : 12 Sample Multiplier: 1

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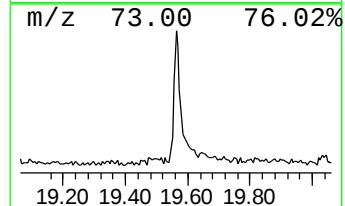
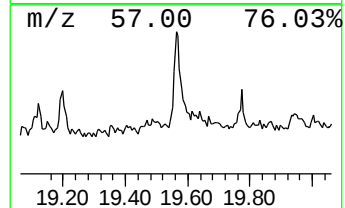
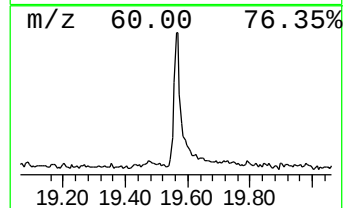
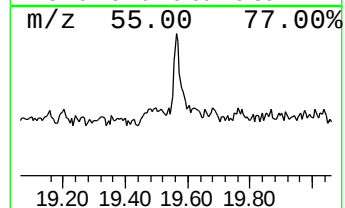
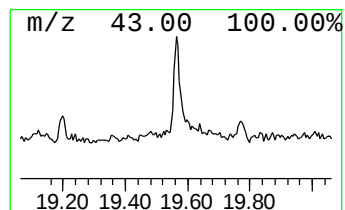
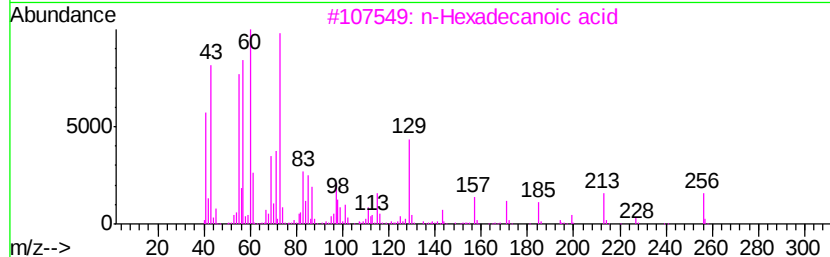
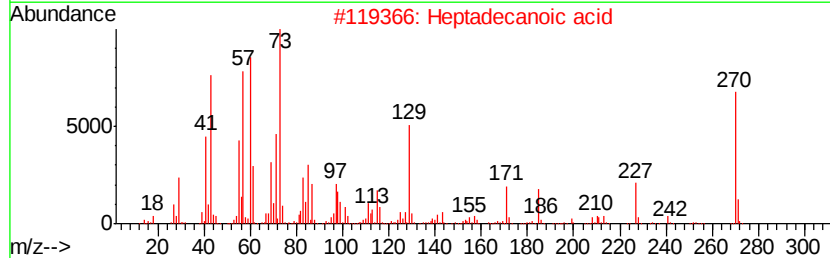
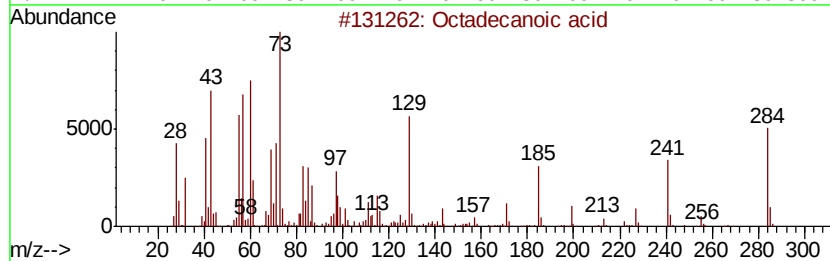
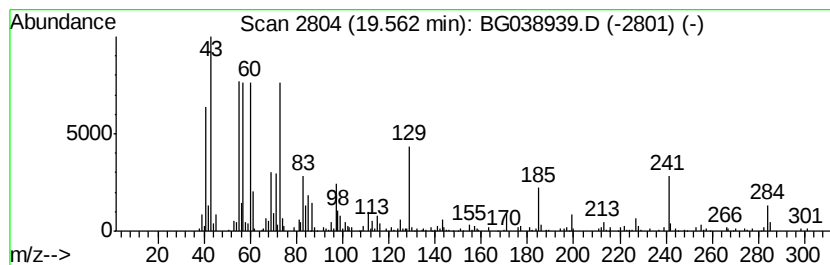
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Octadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.56	8.20 ng	171765	Phenanthrene-d10	17.44

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			Heptadecanoic acid	270	C17H34O2	000506-12-7	90
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	58
4			Octadecanoic acid, 2-(2-hydroxyve...	372	C22H44O4	000106-11-6	50
5			Nonadecanoic acid	298	C19H38O2	000646-30-0	49



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010419\
Data File : BG038939.D
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Operator : JU/SJ
Sample : K1043-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

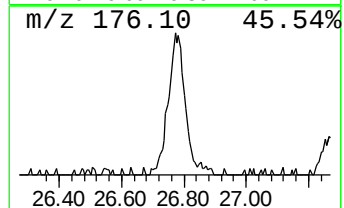
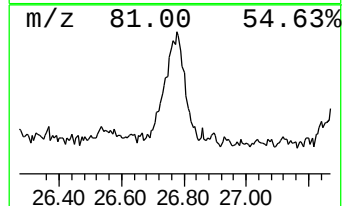
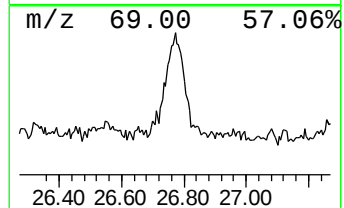
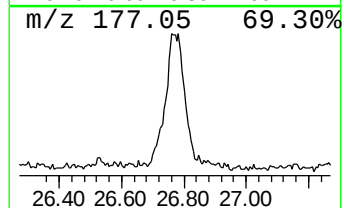
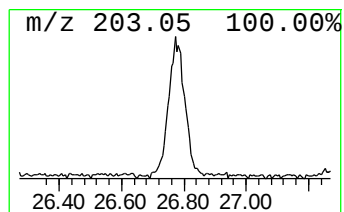
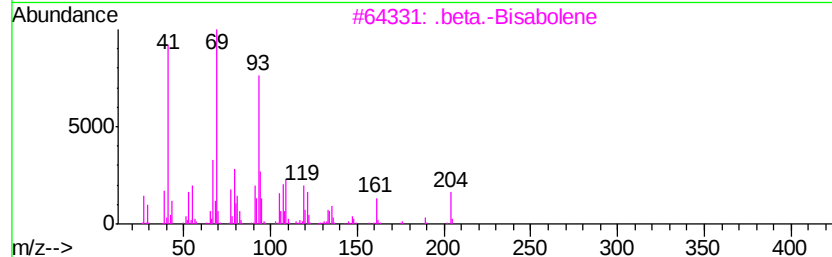
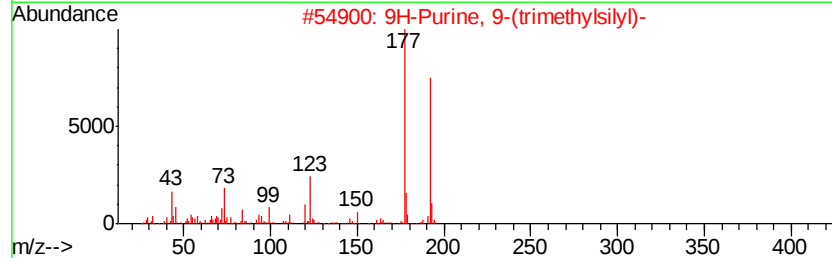
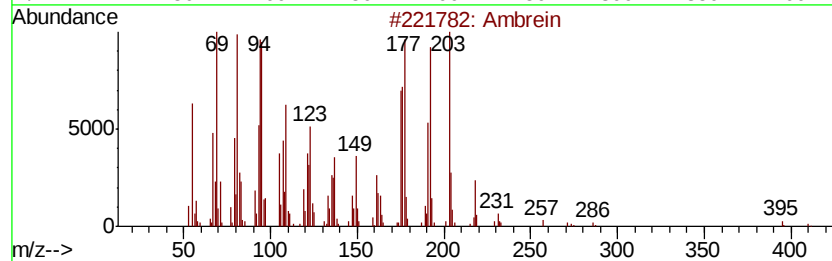
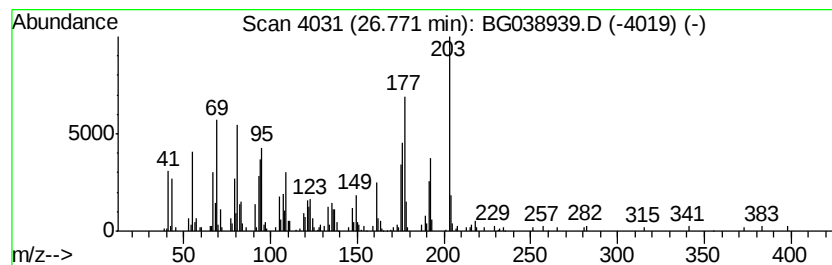
Instrument :
BNA_G
ClientSampleId :
HB-1319-WS

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG121218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 Ambrein Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.77	24.33 ng	551141	Perylene-d12	25.01
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Ambrein	428 C30H52O	1000292-95-3	50
2	9H-Purine, 9-(trimethylsilyl)-	192 C8H12N4Si	032865-85-3	15
3	.beta.-Bisabolene	204 C15H24	000495-61-4	15
4	Gitoxigenin	390 C23H34O5	000545-26-6	14
5	8-Amino-6-methoxy-4-methyl-5-[n-...	316 C19H28N2O2	108190-52-9	14



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010419\
Data File : BG038939.D
Acq On : 4 Jan 2019 21:06
Operator : JU/SJ
Sample : K1043-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
HB-1319-WS

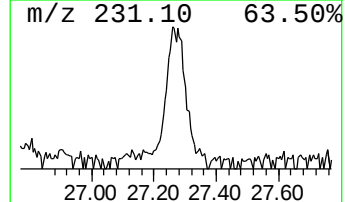
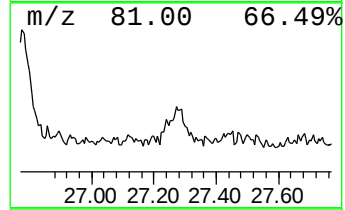
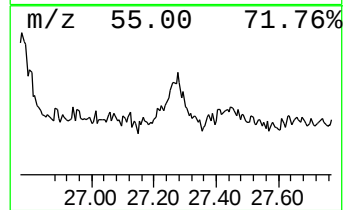
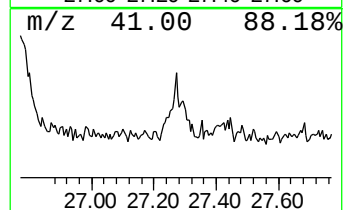
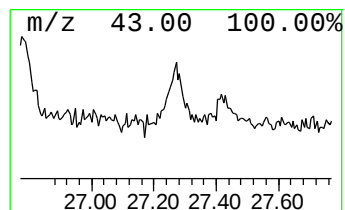
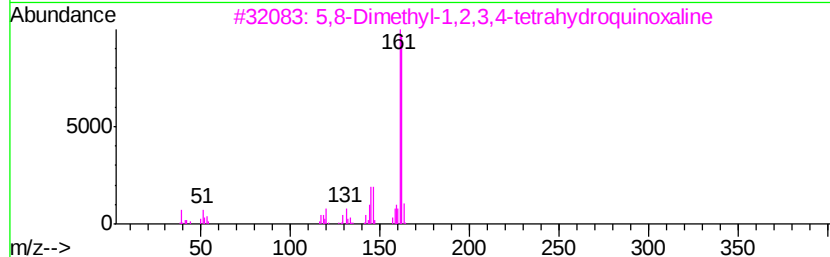
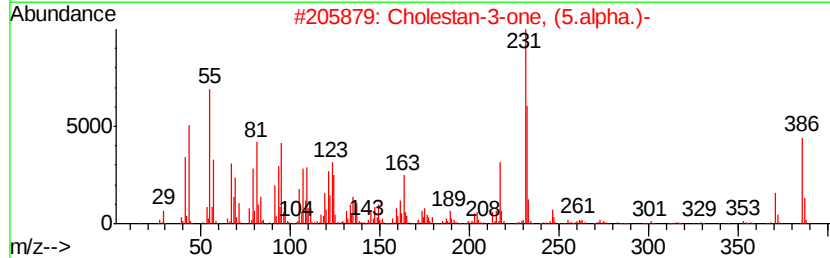
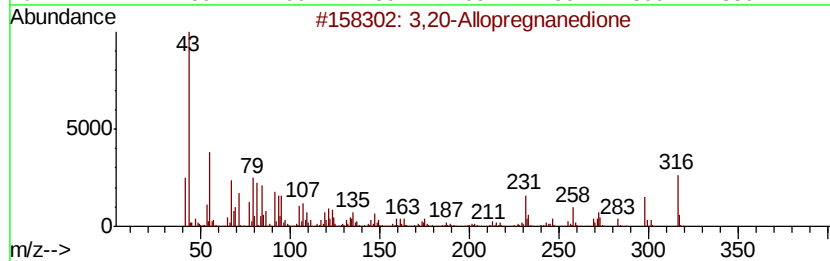
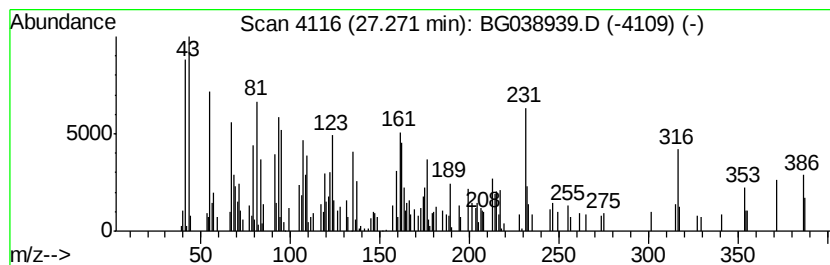
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 unknown27.27 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.27	8.86 ng	200714	Perylene-d12	25.01

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,20-Allopregnanedione			316	C21H32O2	000566-65-4	18
2	Cholestan-3-one, (5.alpha.)-			386	C27H46O	000566-88-1	15
3	5,8-Dimethyl-1,2,3,4-tetrahydroq...			162	C10H14N2	066102-39-4	15
4	Pregnane-3,20-dione, (5.beta.)-			316	C21H32O2	000128-23-4	15
5	Cholest-4-en-3-ol			386	C27H46O	014597-42-3	15



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010419\
Data File : BG038939.D
Acq On : 4 Jan 2019 21:06
Operator : JU/SJ
Sample : K1043-01
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
HB-1319-WS

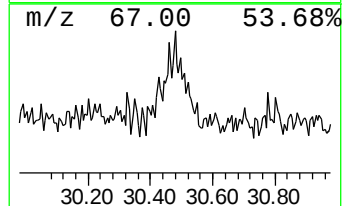
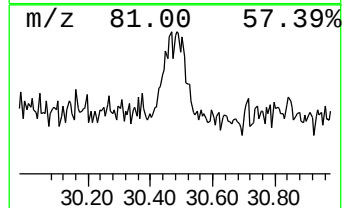
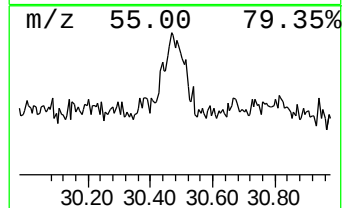
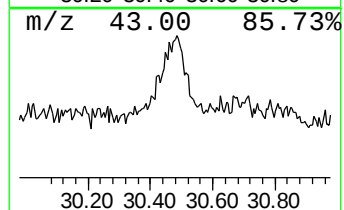
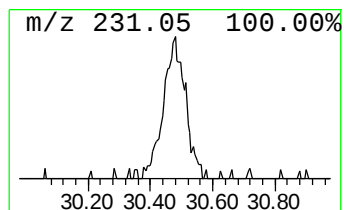
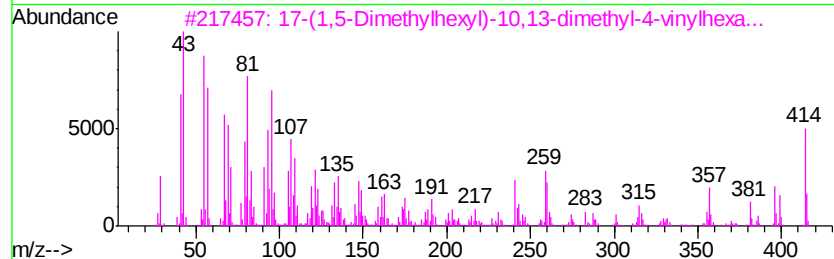
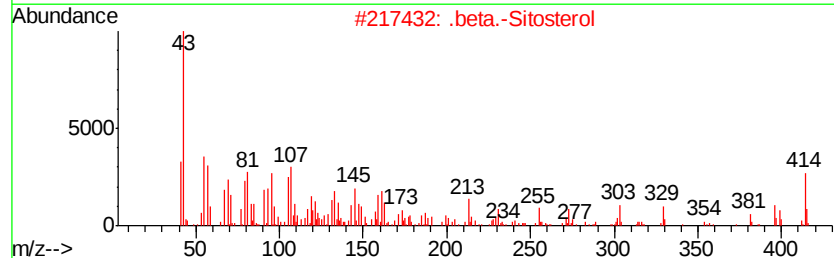
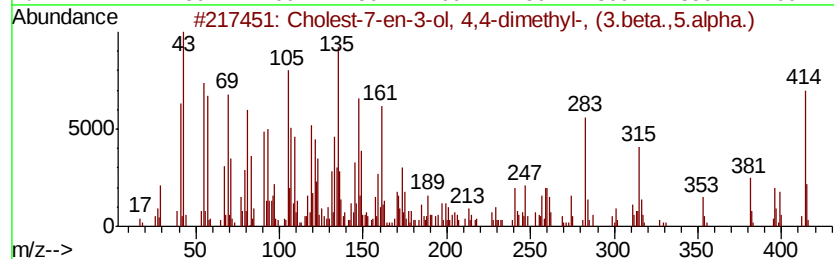
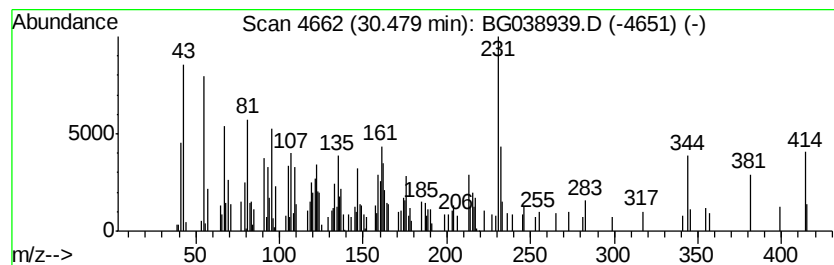
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG121218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 unknown30.48 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
30.48	11.22 ng	254119	Perylene-d12	25.01

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cholest-7-en-3-ol, 4,4-dimethyl-...	414	C29H50O	006384-28-7	43
2		.beta.-Sitosterol	414	C29H50O	000083-46-5	42
3		17-(1,5-Dimethylhexyl)-10,13-dim...	414	C29H50O	1000210-86-9	38
4		.gamma.-Sitosterol	414	C29H50O	000083-47-6	30
5		d-Norandrostane (5.alpha.,14.alp...	246	C18H30	1000281-13-8	25



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG010419\
 Data File : BG038939.D
 Acq On : 4 Jan 2019 21:06
 Operator : JU/SJ
 Sample : K1043-01
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 HB-1319-WS

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG121218.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...	3.17	6.8	ng	62647	1	8.05	184339	20.0
Benzene, 1-ethyl-...	7.11	7.2	ng	65961	1	8.05	184339	20.0
unknown7.58	7.58	67.0	ng	617044	1	8.05	184339	20.0
Benzene, 1,2,3-tr...	7.68	14.5	ng	133717	1	8.05	184339	20.0
Benzene, 4-ethyl-...	9.14	5.2	ng	47990	1	8.05	184339	20.0
Benzene, 2-etheny...	10.24	5.7	ng	68001	2	10.86	237854	20.0
Naphthalene, 1,2,...	10.48	4.2	ng	50010	2	10.86	237854	20.0
Naphthalene, 1,2,...	12.04	6.7	ng	79352	2	10.86	237854	20.0
Tetradecanoic acid	16.81	9.0	ng	187755	4	17.44	418940	20.0
n-Hexadecanoic acid	18.33	216.6	ng	4537520	4	17.44	418940	20.0
Octadecanoic acid	19.56	8.2	ng	171765	4	17.44	418940	20.0
Ambrein	26.77	24.3	ng	551141	6	25.01	453035	20.0
unknown27.27	27.27	8.9	ng	200714	6	25.01	453035	20.0
unknown30.48	30.48	11.2	ng	254119	6	25.01	453035	20.0