

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM041819\
 Data File : BM019967.D
 Acq On : 18 Apr 2019 23:22
 Operator : JU/SJ
 Sample : K2461-03
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PZ-15

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_M\METHODS\8270-BM041519.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	4.863	313	319	329	rBV	133652	212860	4.47%	0.788%
2	5.128	358	364	382	rBV	606419	1037972	21.77%	3.841%
3	6.716	628	634	649	rBV	273425	636095	13.34%	2.354%
4	7.051	683	691	709	rBV	1270622	2178700	45.70%	8.062%
5	7.510	762	769	784	rVB	335670	568322	11.92%	2.103%
6	7.822	813	822	837	rBV	1307048	2168806	45.49%	8.025%
7	8.586	947	952	960	rVB2	71075	113801	2.39%	0.421%
8	8.687	960	969	988	rBV	785099	1595319	33.46%	5.903%
9	9.004	1016	1023	1030	rVB2	158349	263411	5.53%	0.975%
10	9.110	1036	1041	1051	rVB	40967	80405	1.69%	0.298%
11	9.686	1129	1139	1147	rBV5	36401	85079	1.78%	0.315%
12	10.292	1236	1242	1255	rVB	329228	630697	13.23%	2.334%
13	11.016	1360	1365	1372	rBV2	25000	49228	1.03%	0.182%
14	11.651	1466	1473	1480	rBV2	38034	82507	1.73%	0.305%
15	12.780	1658	1665	1677	rBV	2308192	3435653	72.07%	12.713%
16	13.657	1810	1814	1825	rVB	40659	78105	1.64%	0.289%
17	14.180	1896	1903	1910	rVV2	538591	821449	17.23%	3.040%
18	15.692	2153	2160	2176	rBV	1723491	2820547	59.17%	10.437%
19	15.957	2200	2205	2219	rBV4	34403	101418	2.13%	0.375%
20	16.251	2250	2255	2264	rVB	95297	170255	3.57%	0.630%
21	16.945	2367	2373	2380	rBV2	663906	976392	20.48%	3.613%
22	17.833	2520	2524	2534	rBV	262034	427146	8.96%	1.581%
23	19.133	2741	2745	2752	rBV	147575	222367	4.66%	0.823%
24	19.362	2781	2784	2789	rBV	87651	108892	2.28%	0.403%
25	19.586	2817	2822	2831	rBV	3725508	4767167	100.00%	17.640%
26	20.856	3034	3038	3044	rVB	227399	274321	5.75%	1.015%
27	21.156	3085	3089	3095	rBV	850698	1215117	25.49%	4.496%
28	22.156	3256	3259	3261	rBV	116548	118770	2.49%	0.439%
29	22.639	3338	3341	3345	rVB2	152470	187137	3.93%	0.692%
30	23.180	3430	3433	3438	rVB2	162198	225747	4.74%	0.835%
31	23.350	3456	3462	3474	rVB	703584	1370788	28.75%	5.072%

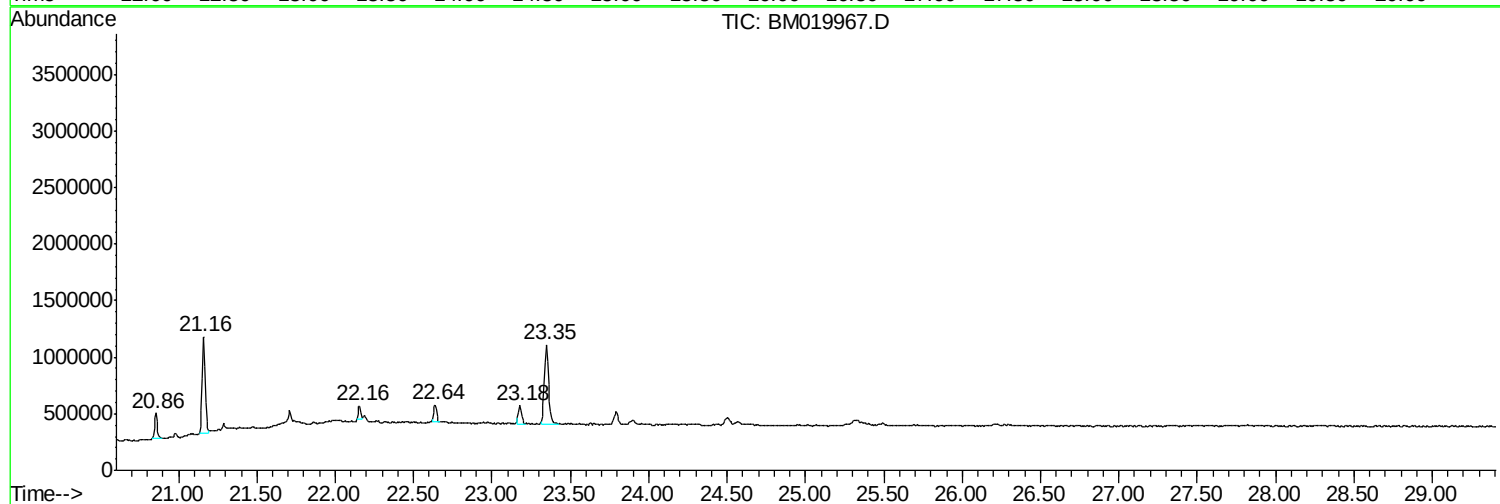
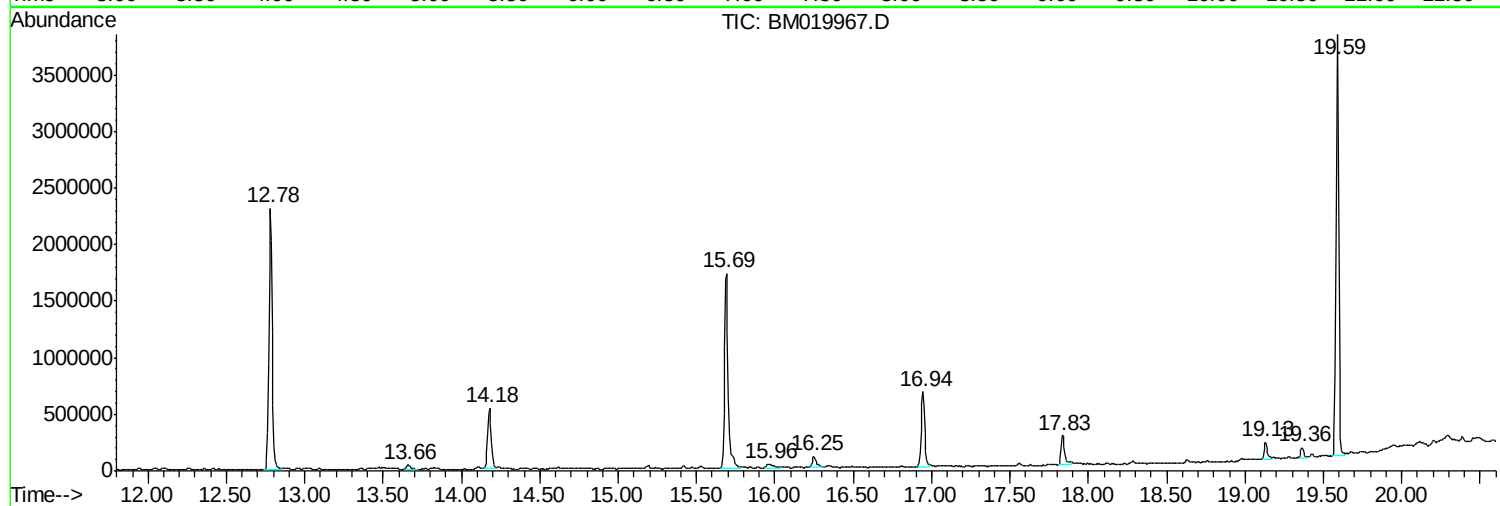
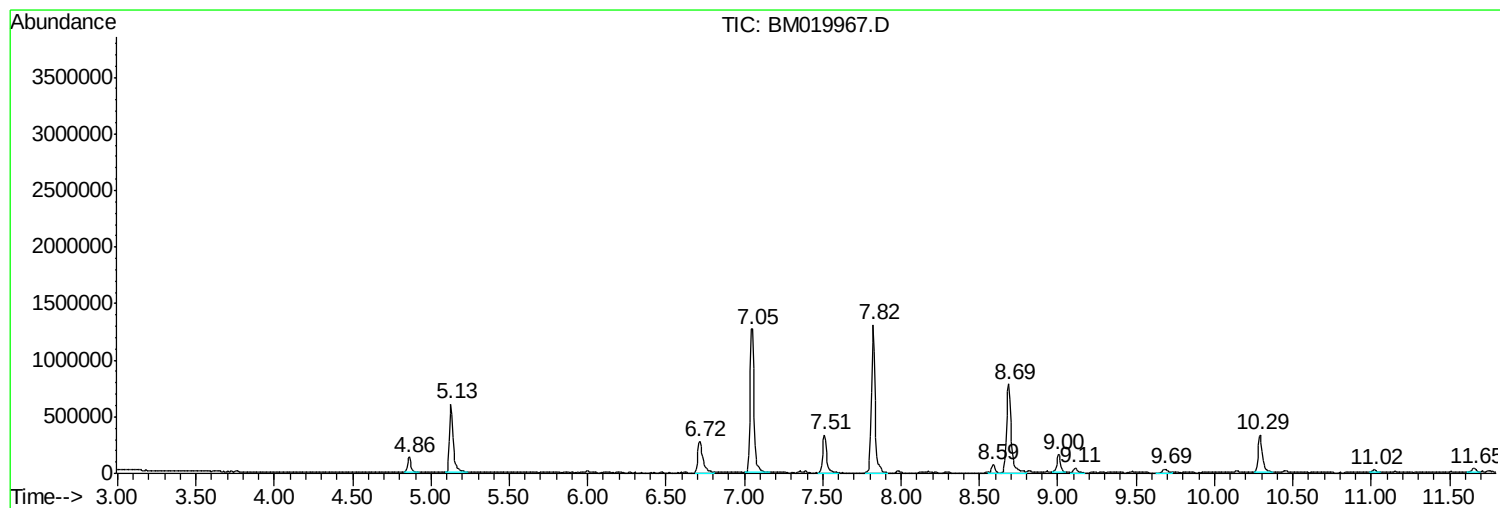
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Instrument :
BNA_M
ClientSampleId :
PZ-15

Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM041519.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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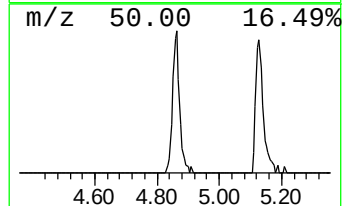
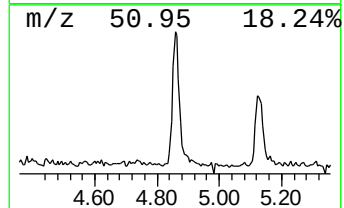
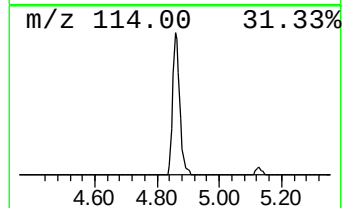
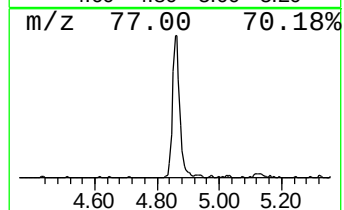
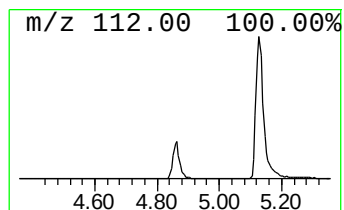
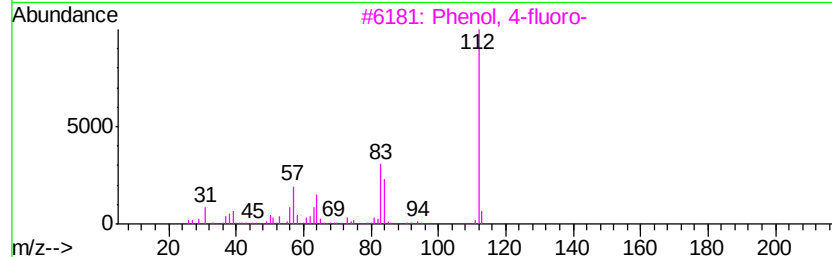
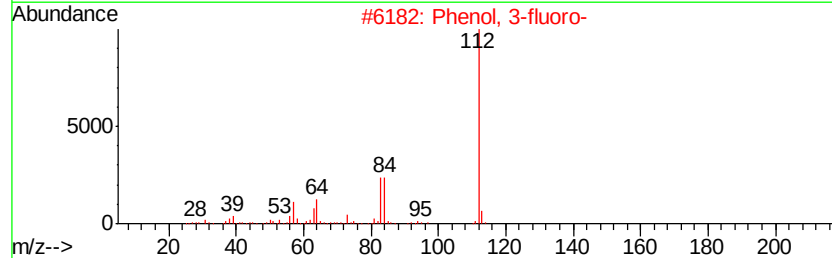
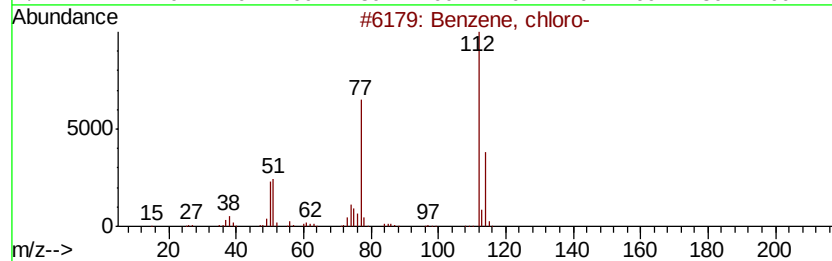
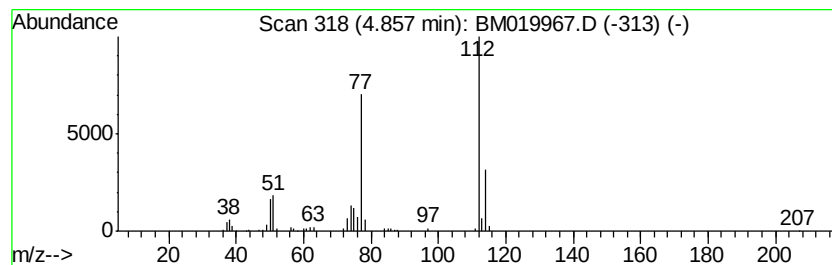
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown4.86 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.86	7.49 ng	212860	1,4-Dichlorobenzene-d4	7.51

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, chloro-	112	C6H5Cl	000108-90-7	97
2		Phenol, 3-fluoro-	112	C6H5FO	000372-20-3	38
3		Phenol, 4-fluoro-	112	C6H5FO	000371-41-5	25
4		3-Hydroxypyridazine 1-oxide	112	C4H4N2O2	018259-51-3	17
5		1,4-Cyclohexanedione	112	C6H8O2	000637-88-7	9



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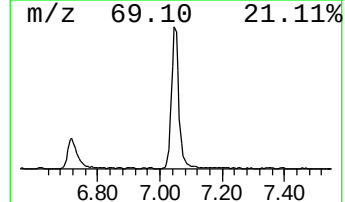
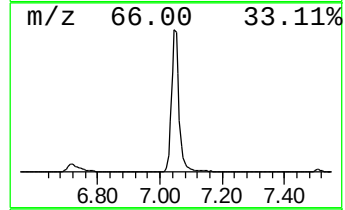
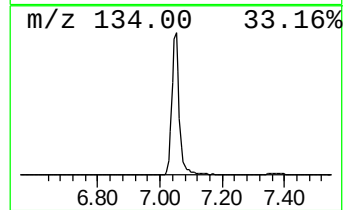
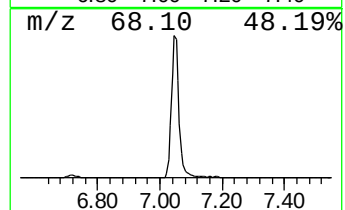
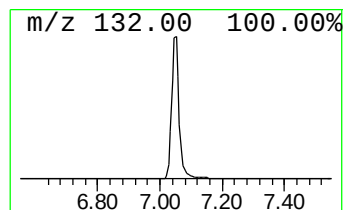
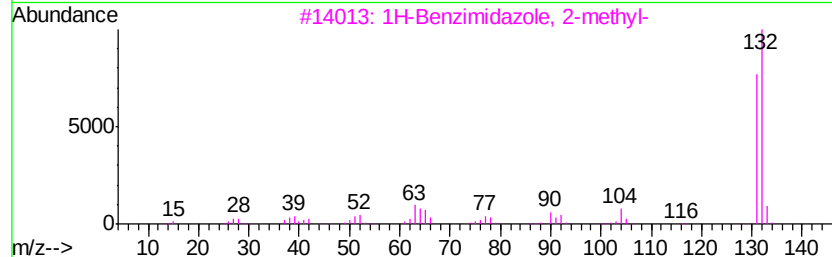
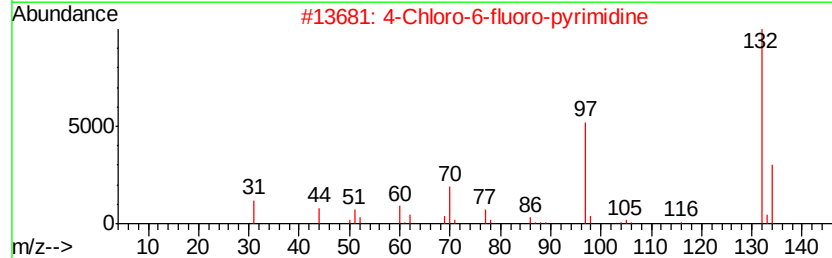
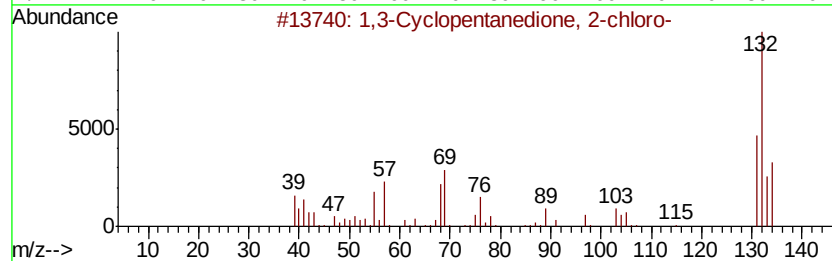
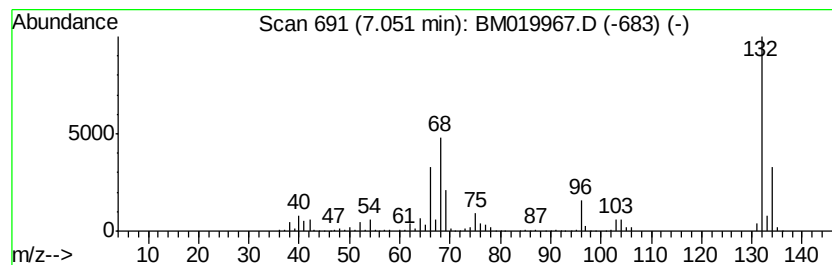
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown7.05 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.05	76.67 ng	2178700	1,4-Dichlorobenzene-d4	7.51

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	38
2		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
4		5-Aminoindole	132	C8H8N2	005192-03-0	22
5		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16



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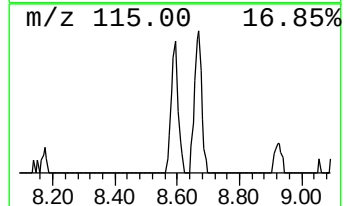
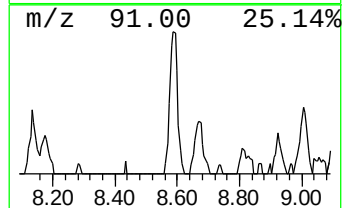
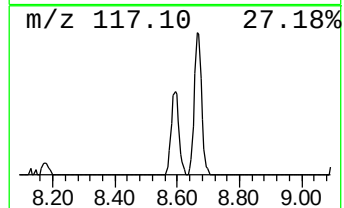
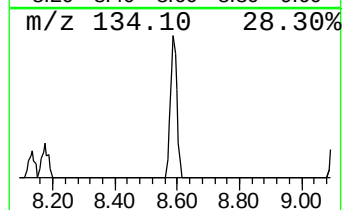
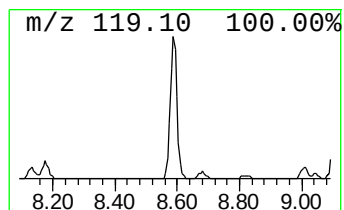
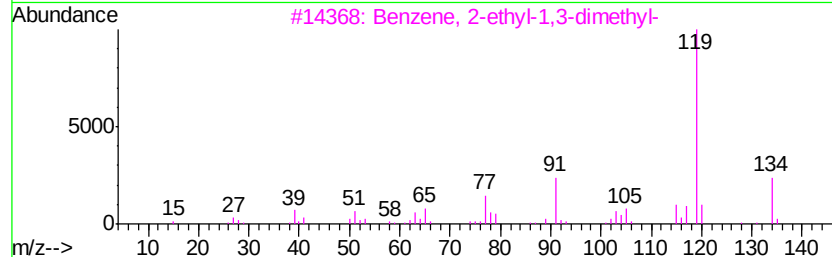
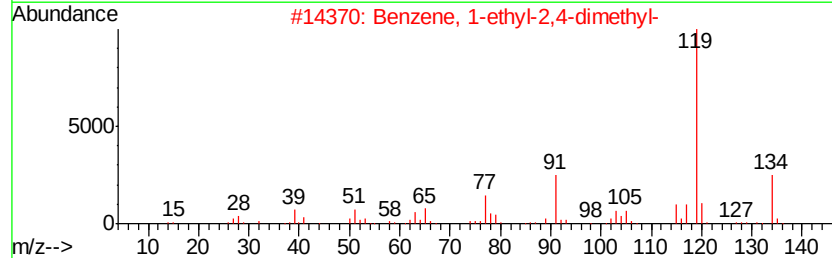
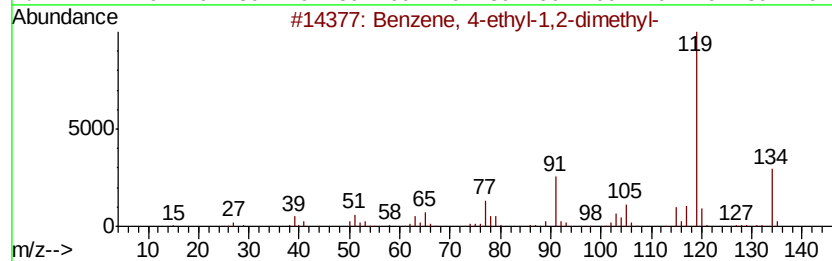
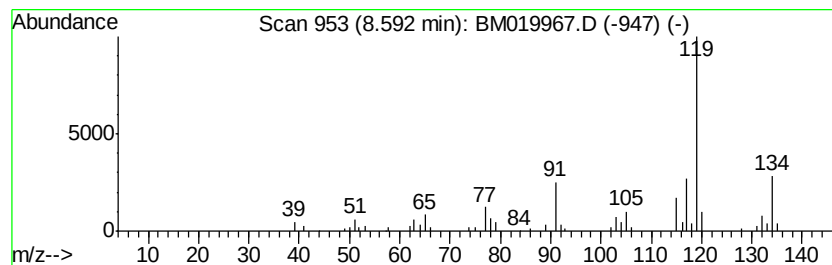
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.59	4.00 ng	113801	1,4-Dichlorobenzene-d4	7.51

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	94
2		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94
3		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	93
4		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	90
5		Benzene, 1-methyl-4-(1-methyleth...	134	C10H14	000099-87-6	81



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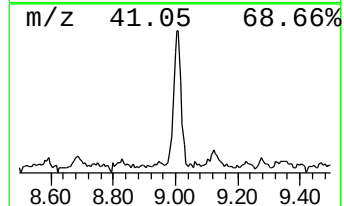
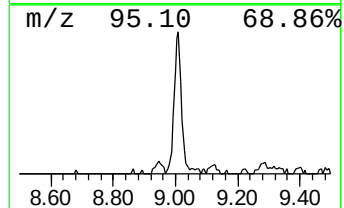
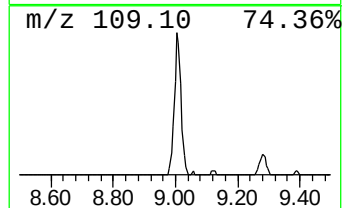
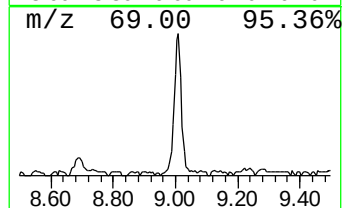
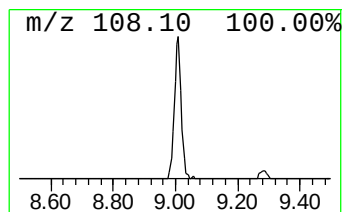
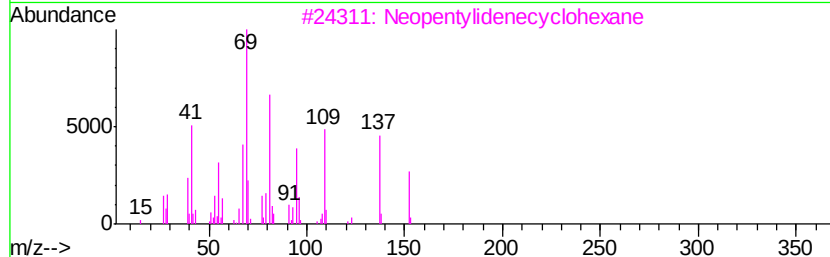
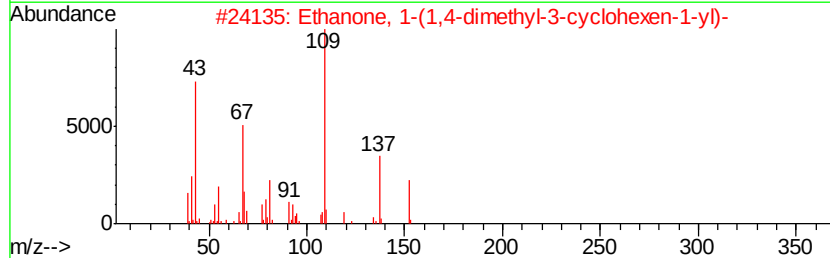
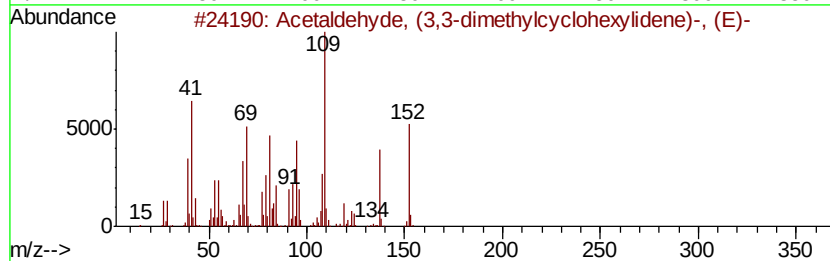
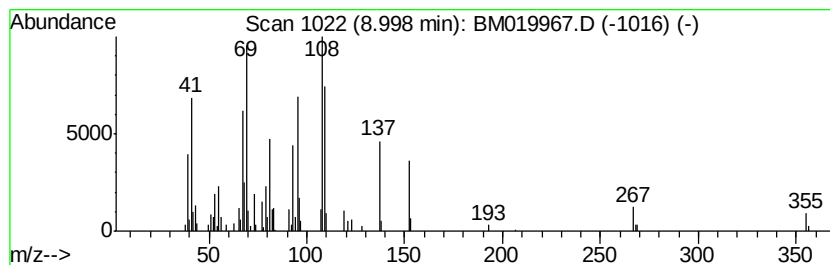
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 5 unknown9.00 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.00	8.35 ng	263411	Naphthalene-d8	10.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Acetaldehyde, (3,3-dimethylcyclo...	152	C10H16O	026532-25-2	42
2		Ethanone, 1-(1,4-dimethyl-3-cycl...	152	C10H16O	043219-68-7	41
3		Neopentylidenecyclohexane	152	C11H20	039546-80-0	38
4		trans-4a-Methyl-decahydronaphtha...	152	C11H20	002547-27-5	30
5		Cyclopentene,1-(2-propenyl)-	108	C8H12	037689-19-3	25



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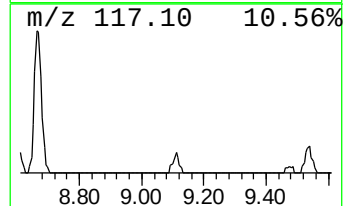
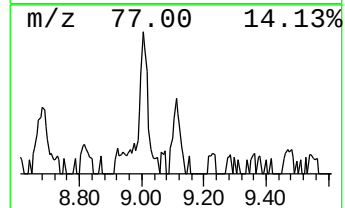
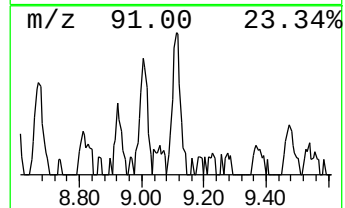
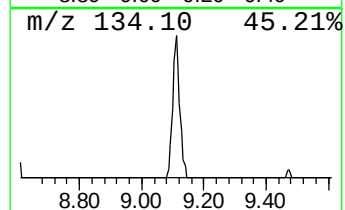
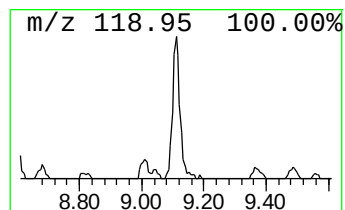
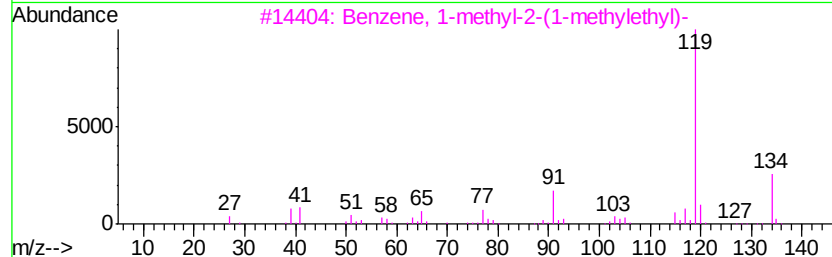
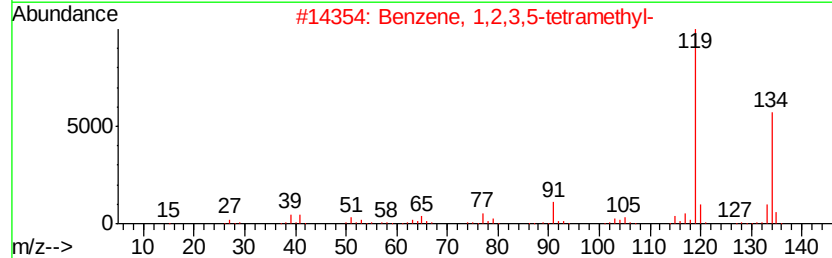
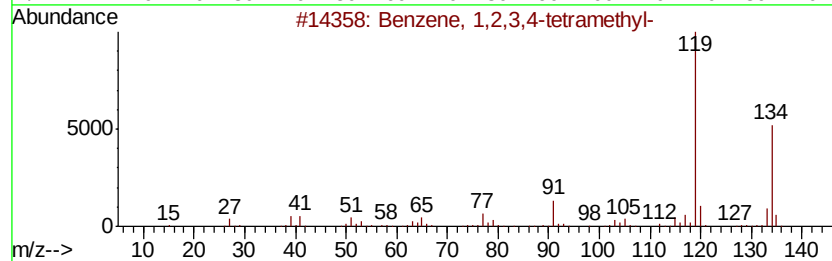
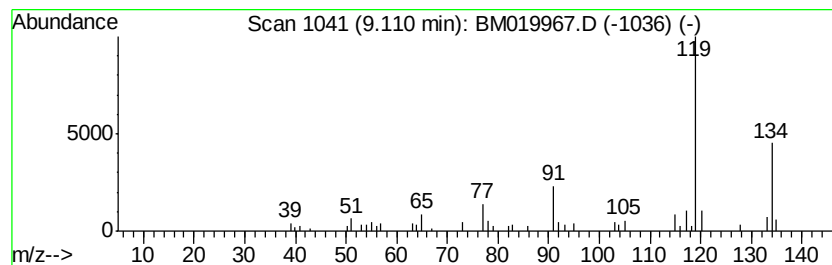
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.11	2.55 ng	80405	Naphthalene-d8	10.29

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	96
2	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	94
3	Benzene, 1-methyl-2-(1-methylethyl)-	134	C10H14	000527-84-4	94
4	Benzene, 1-methyl-4-(1-methylethyl)-	134	C10H14	000099-87-6	94
5	Benzene, 1-methyl-3-(1-methylethyl)-	134	C10H14	000535-77-3	93



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PZ-15

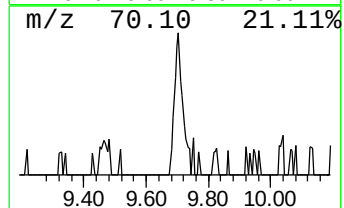
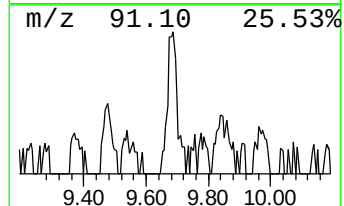
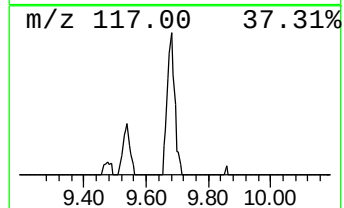
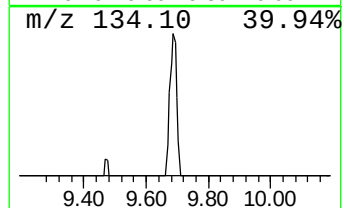
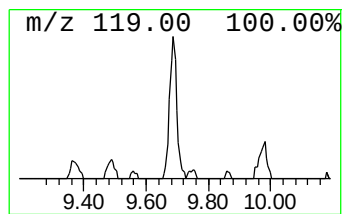
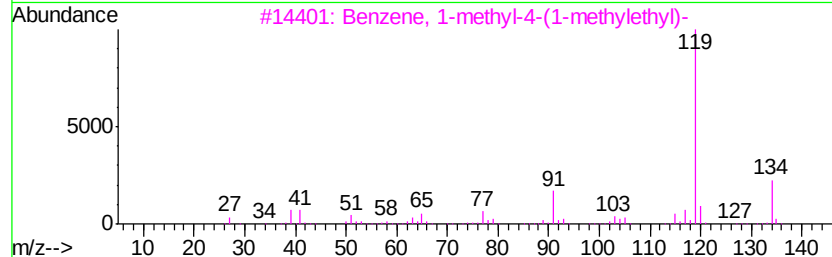
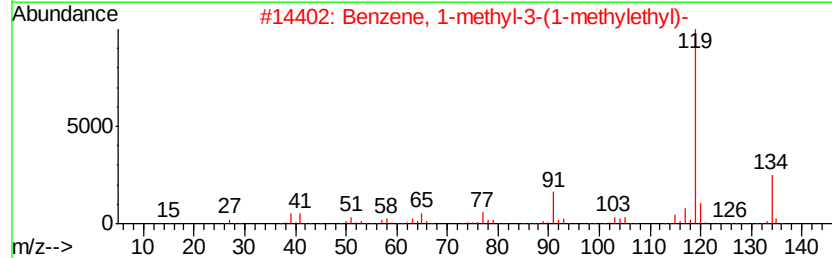
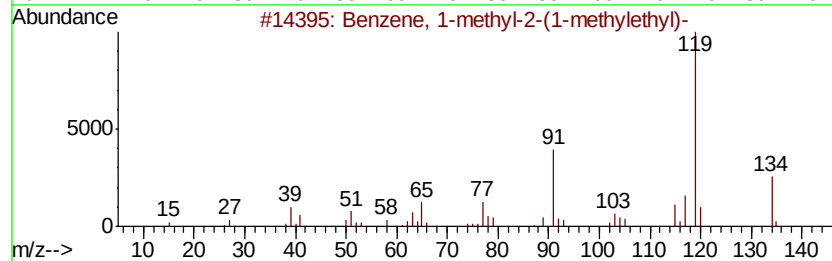
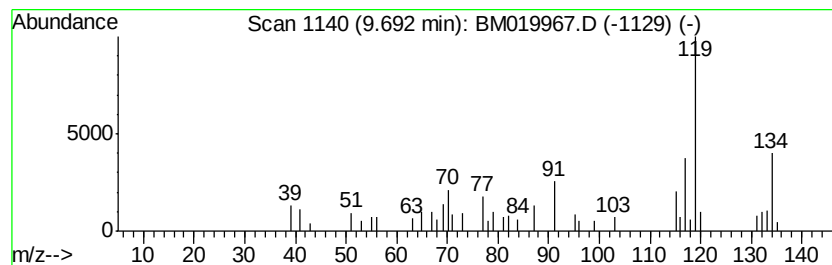
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Benzene, 1-methyl-2-(1-meth... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.69	2.70 ng	85079	Naphthalene-d8	10.29

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-2-(1-methylethyl)-	134	C10H14	000527-84-4	58		
2	Benzene, 1-methyl-3-(1-methylethyl)-	134	C10H14	000535-77-3	53		
3	Benzene, 1-methyl-4-(1-methylethyl)-	134	C10H14	000099-87-6	53		
4	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	52		
5	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	52		



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM041819\
Data File : BM019967.D
Acq On : 18 Apr 2019 23:22
Operator : JU/SJ
Sample : K2461-03
Misc :
ALS Vial : 15 Sample Multiplier: 1

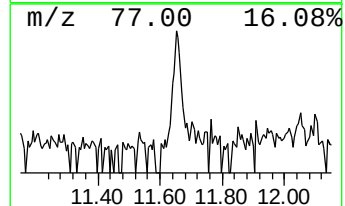
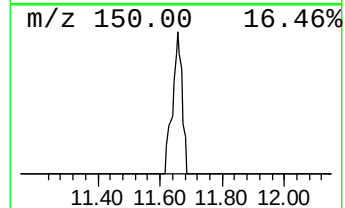
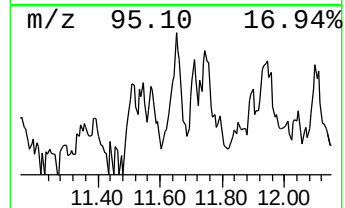
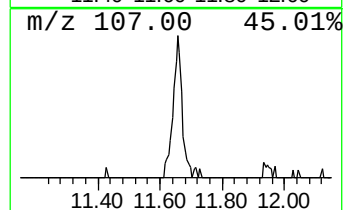
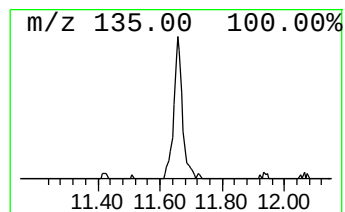
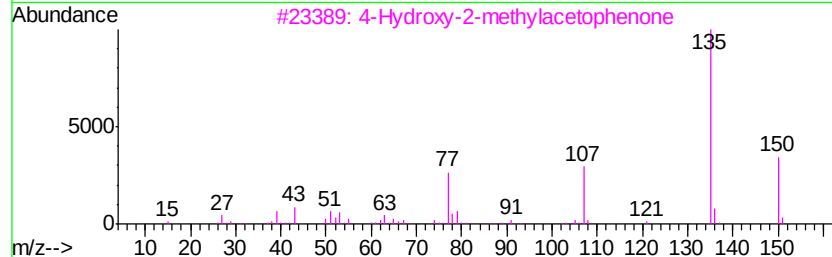
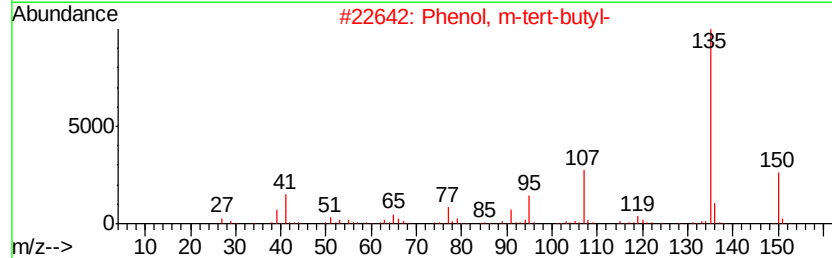
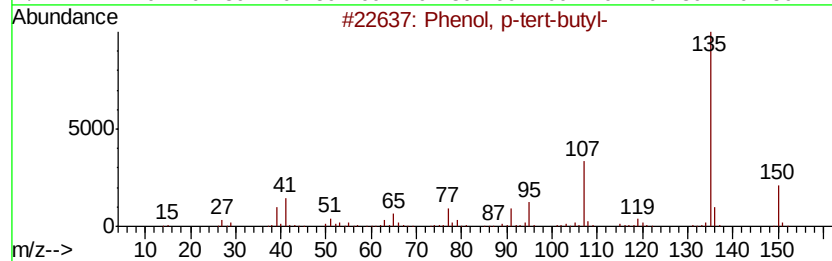
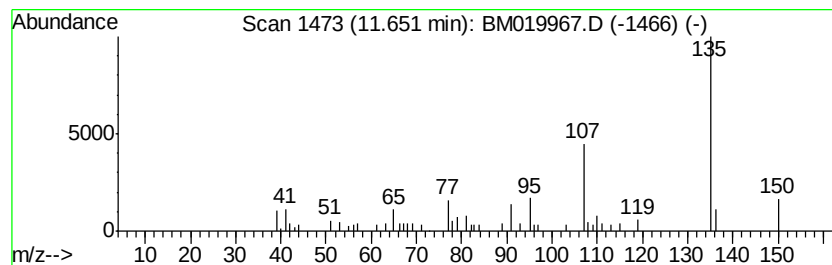
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Phenol, p-tert-butyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD		R.T.	
11.65	2.62 ng	82507	Naphthalene-d8		10.29	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, p-tert-butyl-		150	C10H14O	000098-54-4	90
2	Phenol, m-tert-butyl-		150	C10H14O	000585-34-2	86
3	4-Hydroxy-2-methylacetophenone		150	C9H10O2	000875-59-2	64
4	1H-Indole, 5-fluoro-		135	C8H6FN	000399-52-0	59
5	Hexestrol		270	C18H22O2	005635-50-7	59



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM041819\
Data File : BM019967.D
Acq On : 18 Apr 2019 23:22
Operator : JU/SJ
Sample : K2461-03
Misc :
ALS Vial : 15 Sample Multiplier: 1

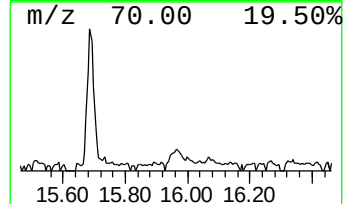
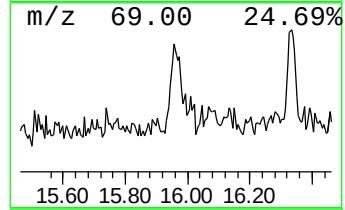
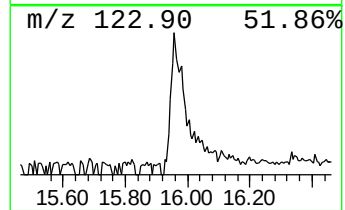
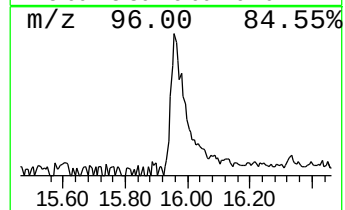
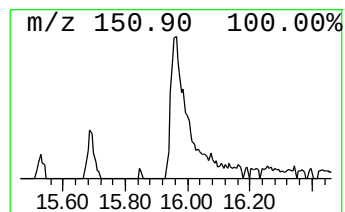
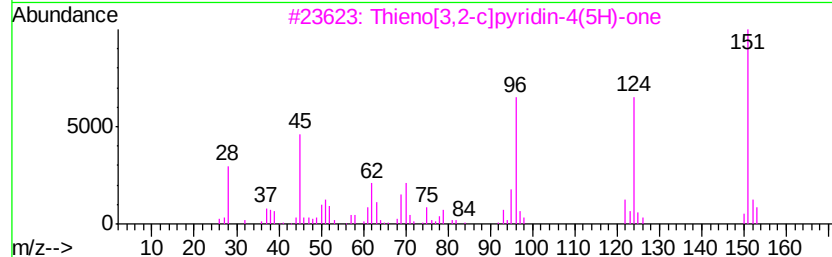
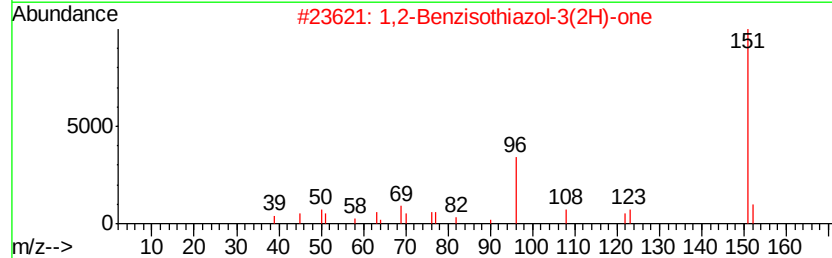
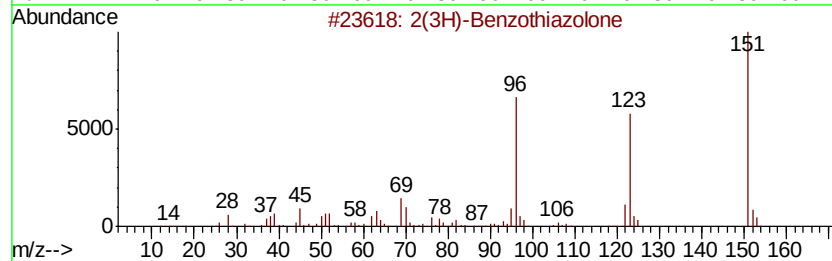
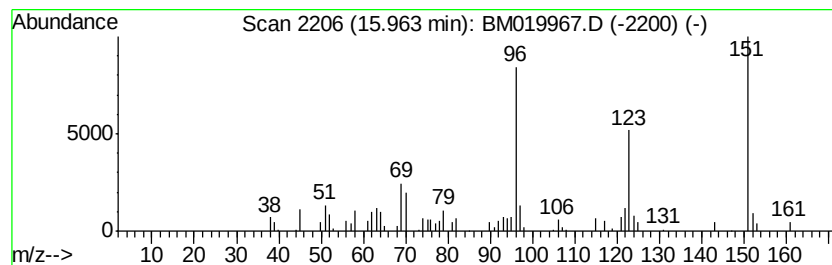
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM041519.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 2(3H)-Benzothiazolone Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD		R.T.		
15.96	2.08 ng	101418	Phenanthrene-d10		16.94		
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2(3H)-Benzothiazolone	151	C7H5NOS	000934-34-9	91
2			1,2-Benzisothiazol-3(2H)-one	151	C7H5NOS	002634-33-5	64
3			Thieno[3,2-c]pyridin-4(5H)-one	151	C7H5NOS	027685-92-3	47
4			2(3H)-Benzoxazolethione	151	C7H5NOS	002382-96-9	35
5			4-Methoxyformanilide	151	C8H9NO2	023896-88-0	35



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM041819\
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Sample : K2461-03
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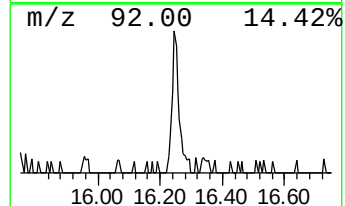
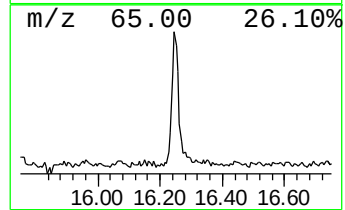
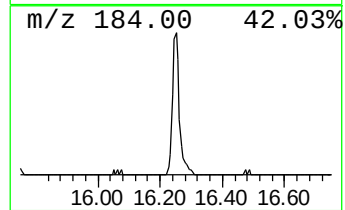
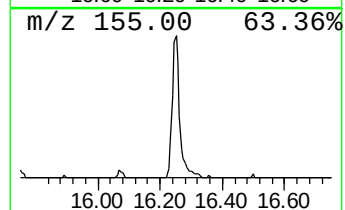
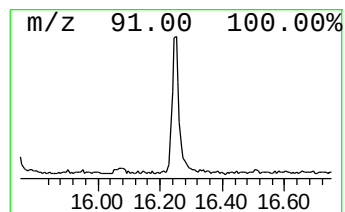
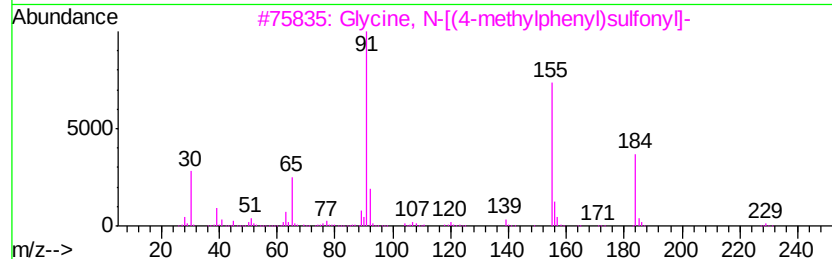
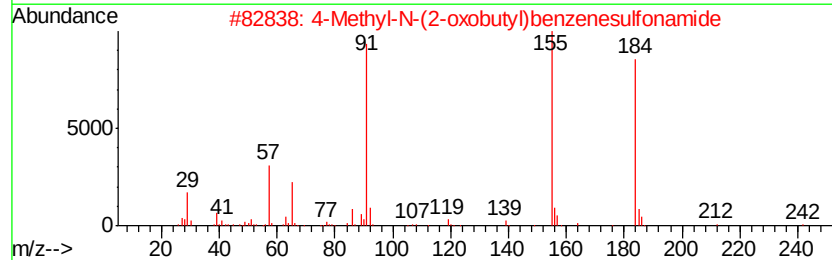
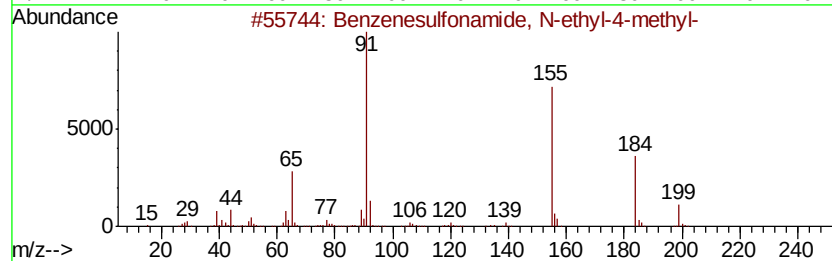
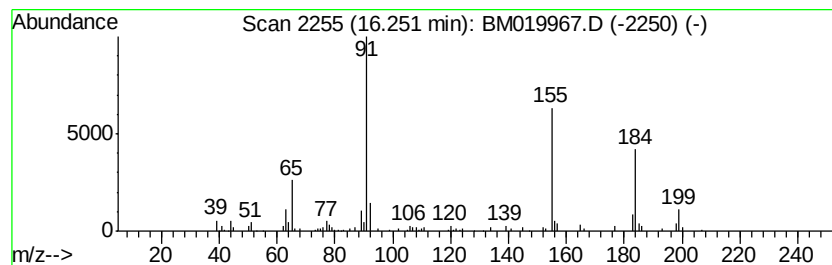
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Benzenesulfonamide, N-ethyl... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD		R.T.	
16.25	3.49 ng	170255	Phenanthrene-d10		16.94	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	98
2		4-Methyl-N-(2-oxobutyl)benzenesu...	241	C11H15NO3S	1000192-14-5	72
3		Glycine, N-[(4-methylphenyl)sulf...	229	C9H11NO4S	001080-44-0	64
4		(Toluene-4-sulfonylamino)-acetic...	283	C13H17NO4S	1000190-48-0	53
5		(O-p-Tosylisonitroso)malononitrile	249	C10H7N3O3S	020893-01-0	50



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM041819\
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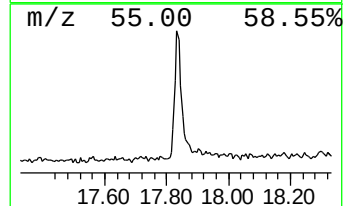
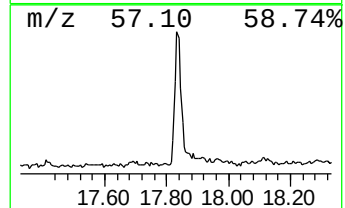
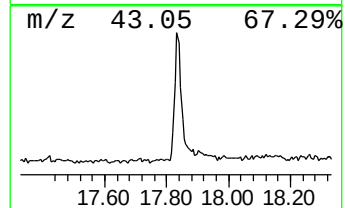
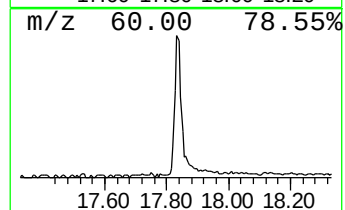
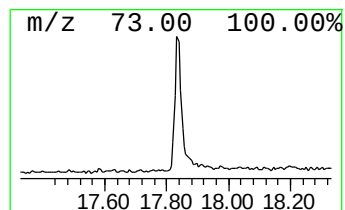
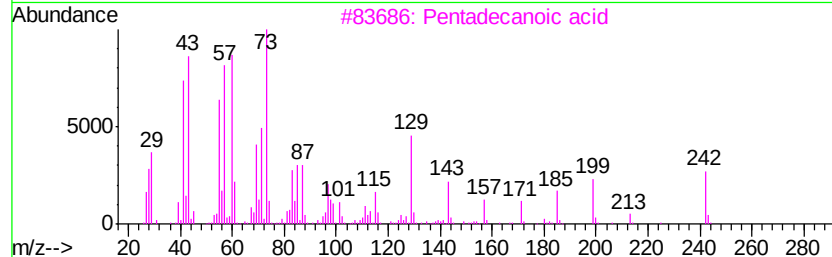
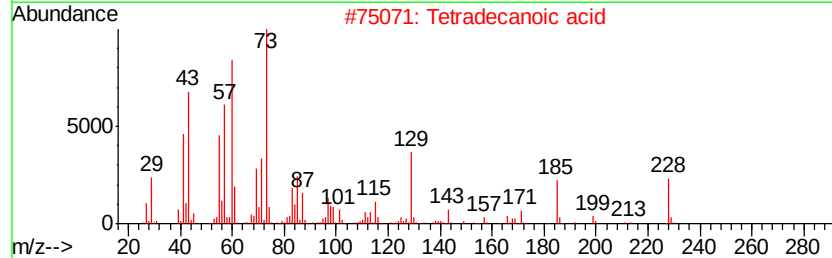
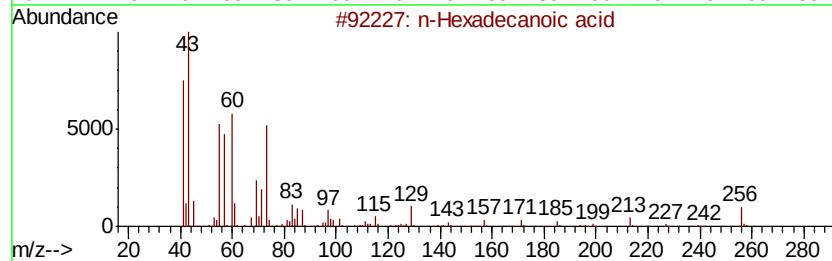
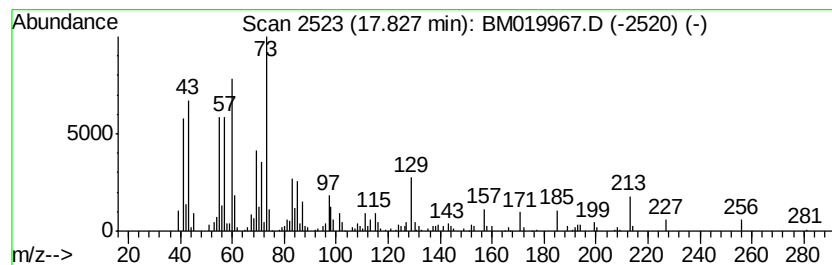
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD		R.T.
17.83	8.75 ng	427146	Phenanthrene-d10		16.94
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
2	Tetradecanoic acid	228	C14H28O2	000544-63-8	64
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	59
4	Tridecanoic acid	214	C13H26O2	000638-53-9	53
5	n-Decanoic acid	172	C10H20O2	000334-48-5	43



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM041819\
Data File : BM019967.D
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Sample : K2461-03
Misc :
ALS Vial : 15 Sample Multiplier: 1

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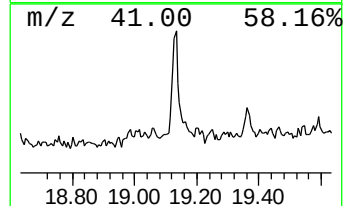
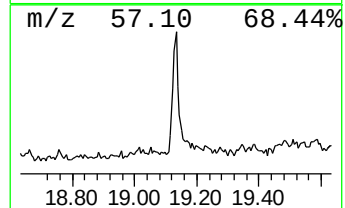
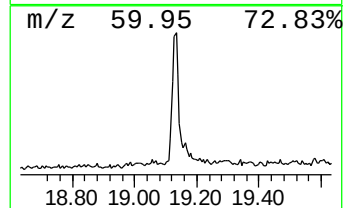
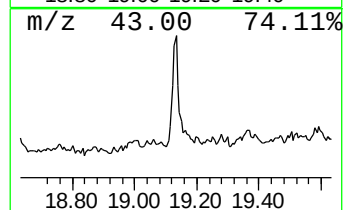
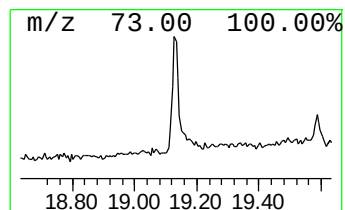
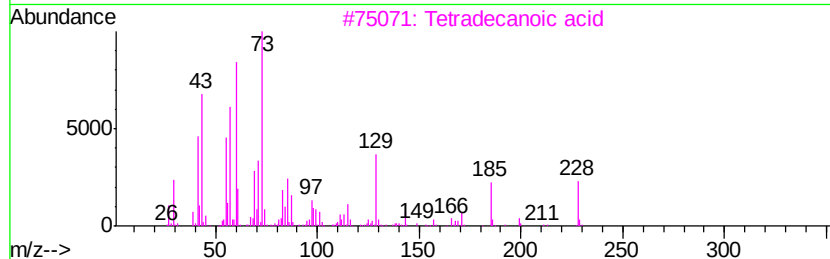
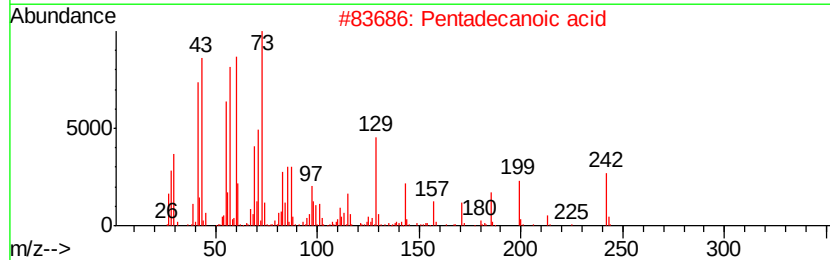
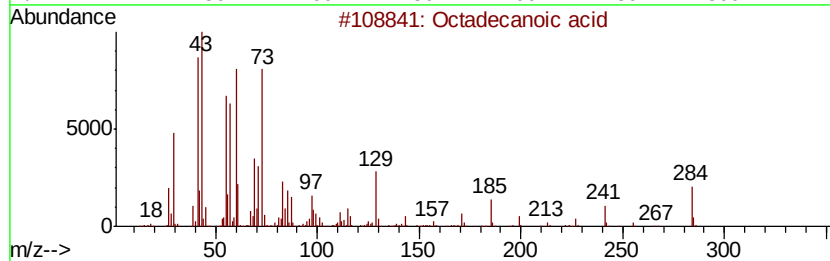
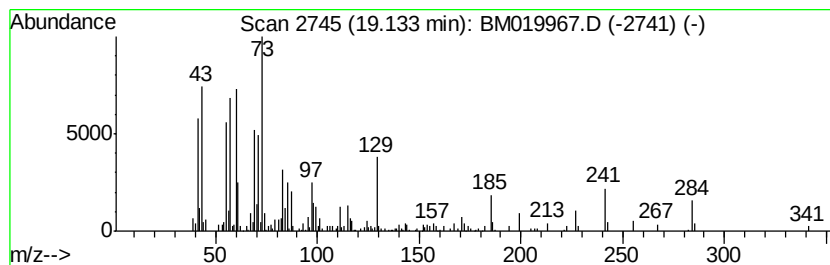
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 Octadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.13	3.66 ng	222367	Chrysene-d12	21.16

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octadecanoic acid	284	C18H36O2	000057-11-4	97
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	86
3		Tetradecanoic acid	228	C14H28O2	000544-63-8	62
4		Undecanoic acid	186	C11H22O2	000112-37-8	38
5		Glucitol, 6-O-nonyl-	308	C15H32O6	132696-05-0	35



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM041819\
Data File : BM019967.D
Acq On : 18 Apr 2019 23:22
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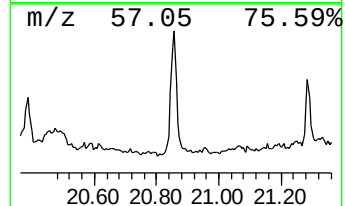
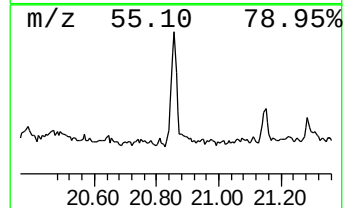
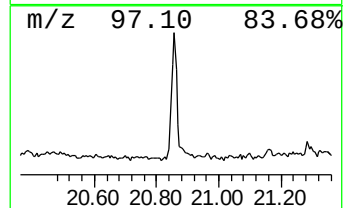
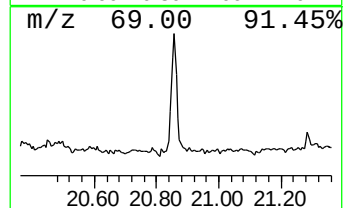
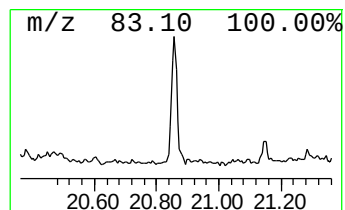
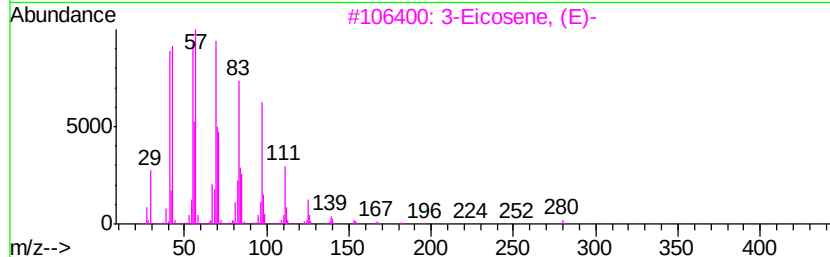
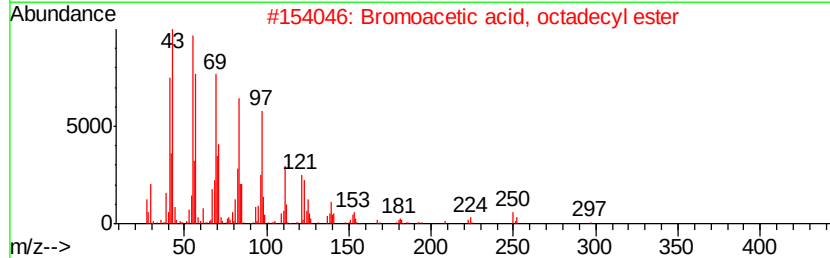
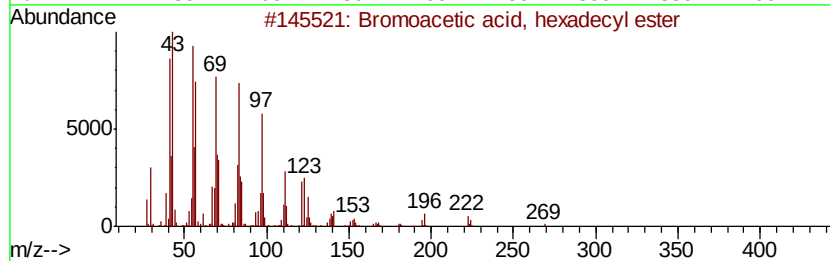
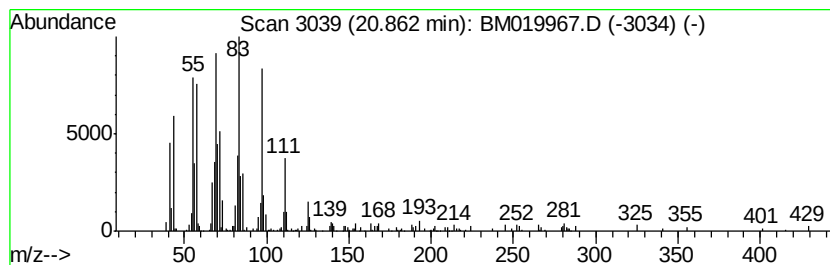
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 Bromoacetic acid, hexadecyl... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.86	4.52 ng	274321	Chrysene-d12	21.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Bromoacetic acid, hexadecyl ester	362	C18H35BrO2	005454-48-8	91
2		Bromoacetic acid, octadecyl ester	390	C20H39BrO2	018992-03-5	91
3		3-Eicosene, (E)-	280	C20H40	074685-33-9	91
4		Cyclotetradecane	196	C14H28	000295-17-0	89
5		2-Chloropropionic acid, octadec...	360	C21H41ClO2	088104-31-8	81



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM041819\
Data File : BM019967.D
Acq On : 18 Apr 2019 23:22
Operator : JU/SJ
Sample : K2461-03
Misc :
ALS Vial : 15 Sample Multiplier: 1

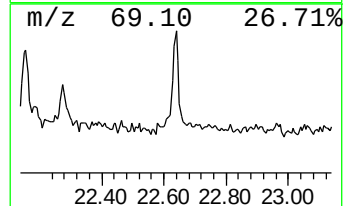
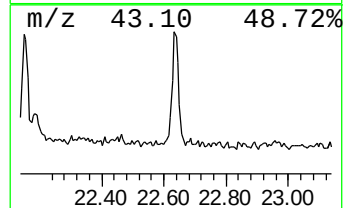
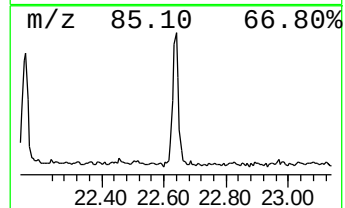
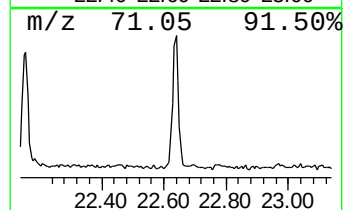
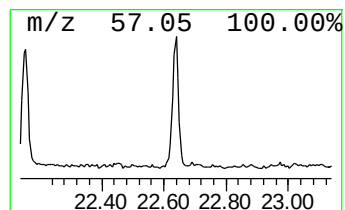
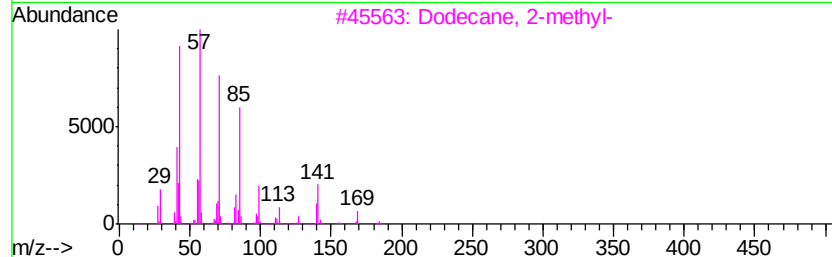
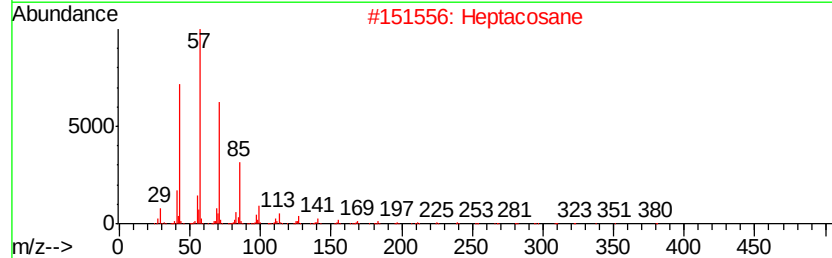
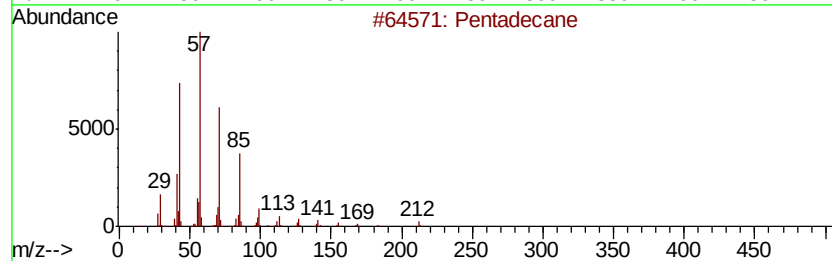
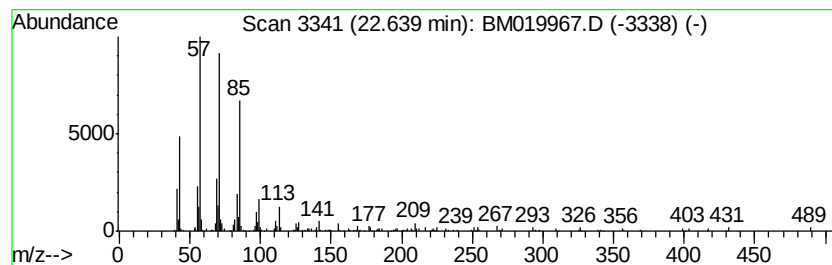
Instrument :
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ClientSampleId :
PZ-15

Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM041519.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 Pentadecane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.64	2.73 ng	187137	Perylene-d12	23.35
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Pentadecane	212 C15H32	000629-62-9	89
2	Heptacosane	380 C27H56	000593-49-7	83
3	Dodecane, 2-methyl-	184 C13H28	001560-97-0	80
4	Pentadecane, 8-heptyl-	310 C22H46	071005-15-7	72
5	Eicosane	282 C20H42	000112-95-8	72



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM041819\
Data File : BM019967.D
Acq On : 18 Apr 2019 23:22
Operator : JU/SJ
Sample : K2461-03
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
PZ-15

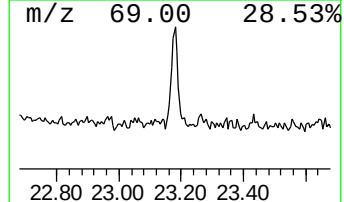
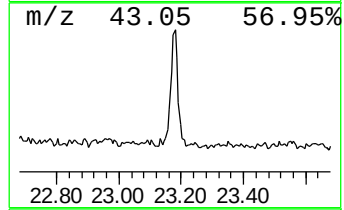
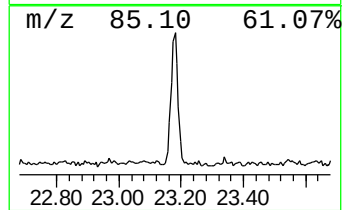
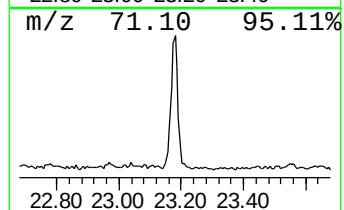
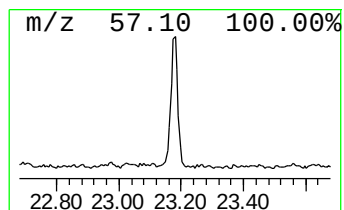
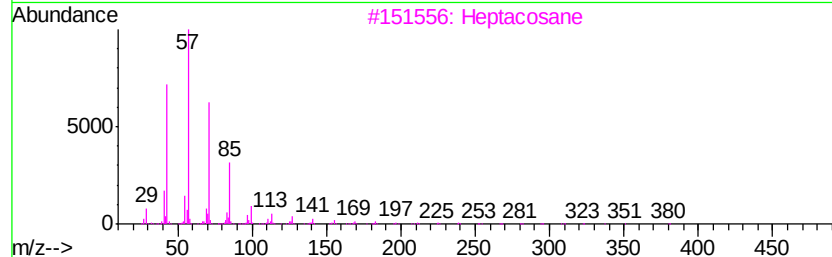
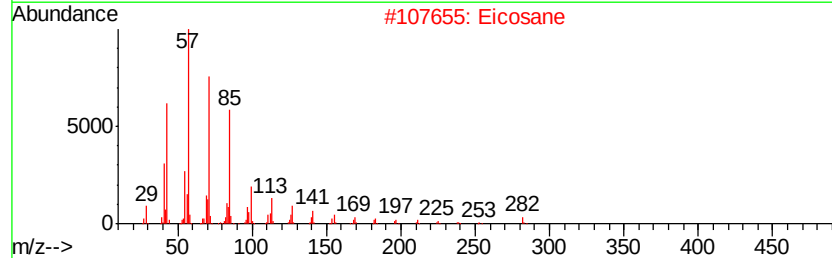
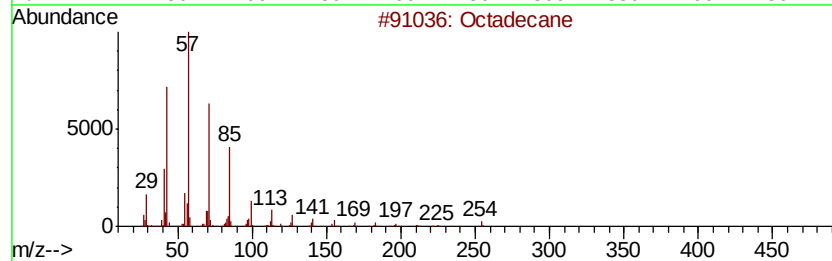
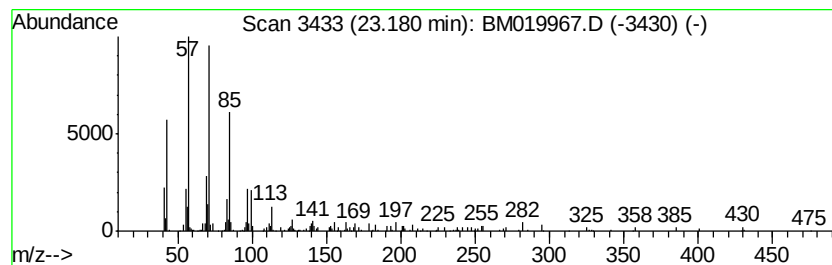
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM041519.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 15 Octadecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.18	3.29 ng	225747	Perylene-d12	23.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecane	254	C18H38	000593-45-3	90
2		Eicosane	282	C20H42	000112-95-8	83
3		Heptacosane	380	C27H56	000593-49-7	62
4		Nonane, 1-iodo-	254	C9H19I	004282-42-2	58
5		Heneicosane	296	C21H44	000629-94-7	52



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM041819\
 Data File : BM019967.D
 Acq On : 18 Apr 2019 23:22
 Operator : JU/SJ
 Sample : K2461-03
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PZ-15

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\8270-BM041519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown4.86	4.86	7.5	ng	212860	1	7.51	568322	20.0
unknown7.05	7.05	76.7	ng	2178700	1	7.51	568322	20.0
Benzene, 4-ethyl-...	8.59	4.0	ng	113801	1	7.51	568322	20.0
unknown9.00	9.00	8.3	ng	263411	2	10.29	630697	20.0
Benzene, 1,2,3,4-...	9.11	2.5	ng	80405	2	10.29	630697	20.0
Benzene, 1-methyl...	9.69	2.7	ng	85079	2	10.29	630697	20.0
Phenol, p-tert-bu...	11.65	2.6	ng	82507	2	10.29	630697	20.0
2(3H)-Benzothiazo...	15.96	2.1	ng	101418	4	16.94	976392	20.0
Benzenesulfonamid...	16.25	3.5	ng	170255	4	16.94	976392	20.0
n-Hexadecanoic acid	17.83	8.8	ng	427146	4	16.94	976392	20.0
Octadecanoic acid	19.13	3.7	ng	222367	5	21.16	1215120	20.0
Bromoacetic acid,...	20.86	4.5	ng	274321	5	21.16	1215120	20.0
Pentadecane	22.64	2.7	ng	187137	6	23.35	1370790	20.0
Octadecane	23.18	3.3	ng	225747	6	23.35	1370790	20.0