

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG080320\
 Data File : BG046156.D
 Acq On : 3 Aug 2020 19:23
 Operator : CG/JU
 Sample : L3536-01
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 JTY1-GW-01

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-BG073020.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	3.159	7	12	30	rVB	1550224	2329114	36.32%	7.100%
2	5.539	410	417	425	rBV	969337	1553540	24.23%	4.735%
3	7.113	678	685	696	rBV	663883	1130755	17.63%	3.447%
4	7.953	821	828	836	rBV	335145	563403	8.79%	1.717%
5	9.105	1016	1024	1037	rVB	976432	1741600	27.16%	5.309%
6	10.633	1279	1284	1296	rVB9	18228	66357	1.03%	0.202%
7	10.750	1296	1304	1314	rBV	456644	840586	13.11%	2.562%
8	11.097	1357	1363	1371	rVB	96371	170465	2.66%	0.520%
9	11.391	1404	1413	1422	rVB	81282	165166	2.58%	0.503%
10	13.206	1714	1722	1730	rBV	2551355	4044324	63.07%	12.328%
11	13.418	1753	1758	1765	rVB	77014	123617	1.93%	0.377%
12	14.029	1856	1862	1866	rBV	170910	245369	3.83%	0.748%
13	14.575	1948	1955	1970	rVB2	728237	1177664	18.36%	3.590%
14	15.321	2076	2082	2088	rVB	164041	241954	3.77%	0.738%
15	15.392	2090	2094	2100	rVB2	79620	113530	1.77%	0.346%
16	15.791	2155	2162	2167	rBV2	104916	161783	2.52%	0.493%
17	16.062	2199	2208	2216	rBV	2035406	3099534	48.33%	9.448%
18	16.150	2216	2223	2230	rVV	363512	545708	8.51%	1.663%
19	16.238	2233	2238	2247	rVB	61697	113928	1.78%	0.347%
20	17.248	2406	2410	2416	rBV7	38869	71311	1.11%	0.217%
21	17.319	2416	2422	2432	rVB	929276	1422368	22.18%	4.336%
22	17.577	2461	2466	2477	rBV	256774	414394	6.46%	1.263%
23	17.995	2527	2537	2543	rBV5	63998	118396	1.85%	0.361%
24	18.053	2543	2547	2552	rVV2	48973	66682	1.04%	0.203%
25	18.224	2570	2576	2583	rBV	432178	628921	9.81%	1.917%
26	18.982	2700	2705	2710	rBV3	76070	124323	1.94%	0.379%
27	19.493	2788	2792	2799	rBV	187790	267621	4.17%	0.816%
28	19.628	2811	2815	2822	rVB	181855	205416	3.20%	0.626%
29	19.740	2830	2834	2840	rVB	84467	123146	1.92%	0.375%
30	19.933	2861	2867	2873	rBV	5029968	6412748	100.00%	19.547%
31	20.027	2879	2883	2887	rVB	74398	98361	1.53%	0.300%
32	21.238	3084	3089	3100	rBV	395955	631132	9.84%	1.924%
33	21.561	3138	3144	3155	rBV2	1060939	1685703	26.29%	5.138%
34	24.669	3663	3673	3683	rBV2	636648	1748289	27.26%	5.329%

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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	25.944	3884	3890	3903	rvB10	38486	125783	1.96%	0.383%
36	28.753	4359	4368	4380	rvB10	59901	233588	3.64%	0.712%

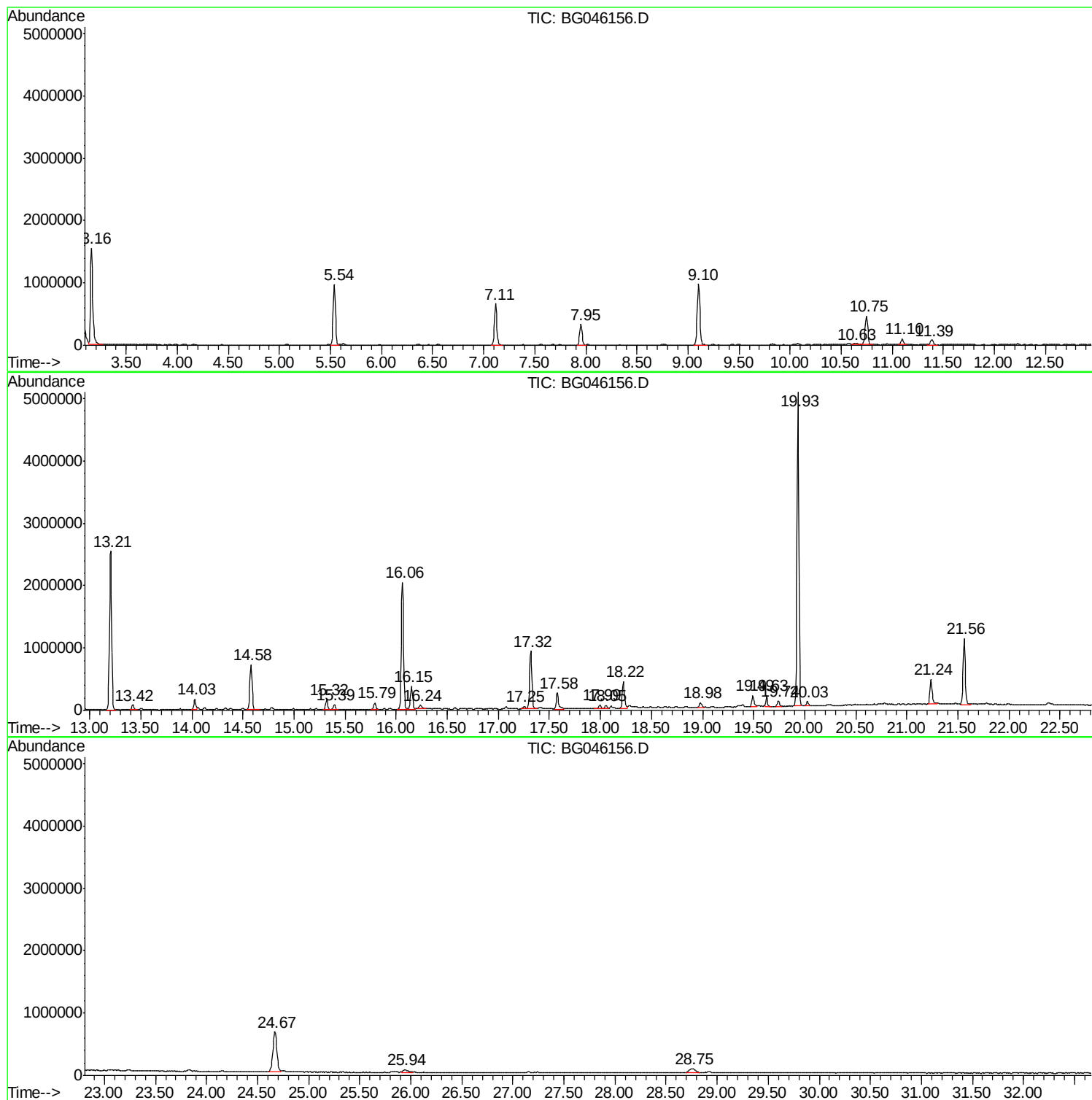
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG073020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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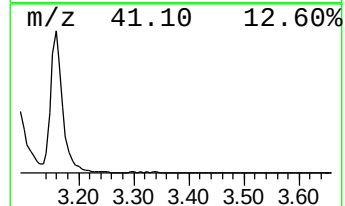
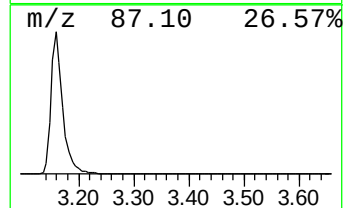
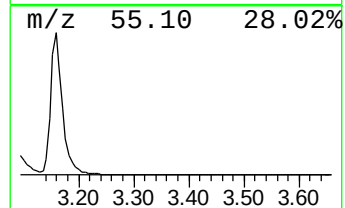
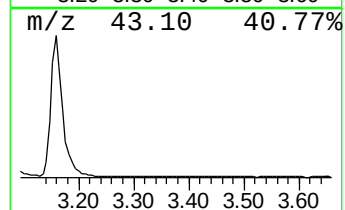
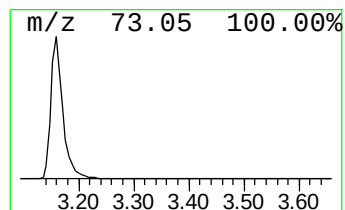
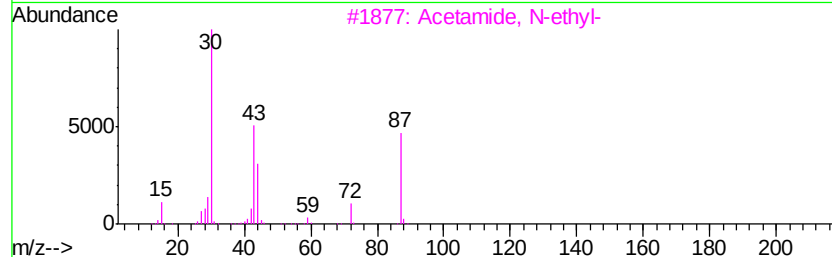
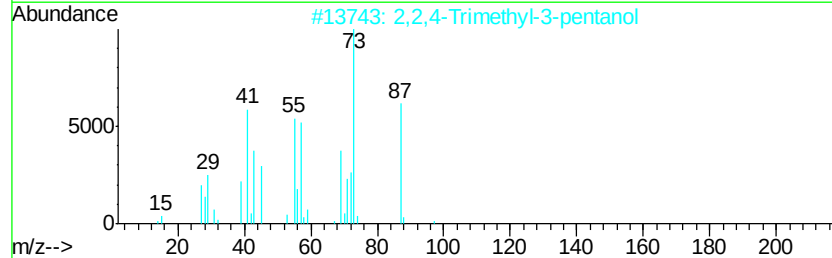
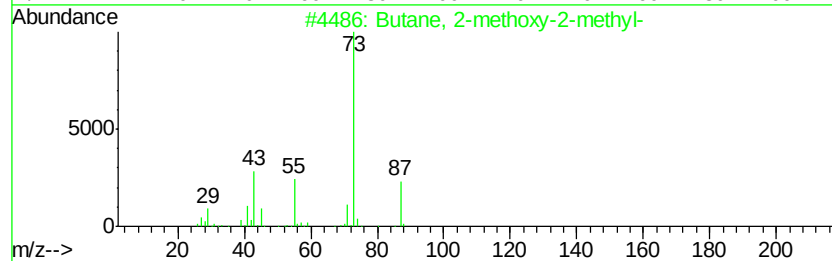
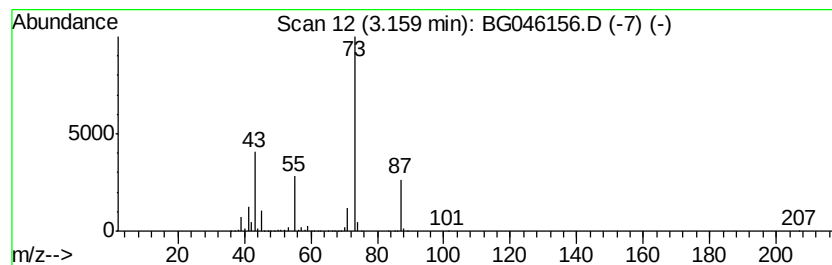
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG073020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.16	82.68 ng	2329110	1,4-Dichlorobenzene-d4	7.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	35
4		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	25
5		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12



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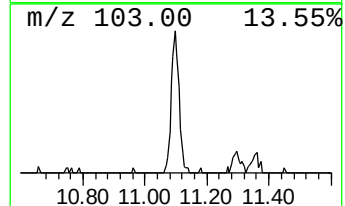
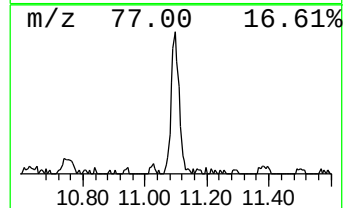
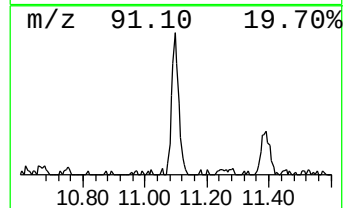
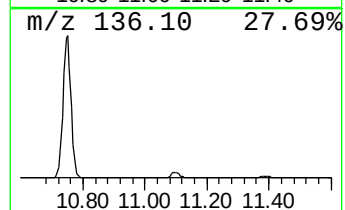
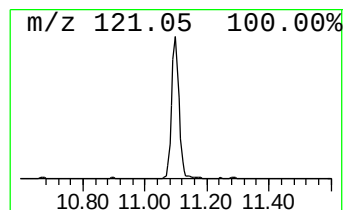
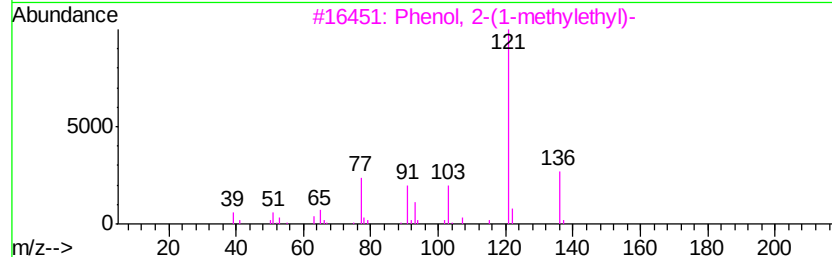
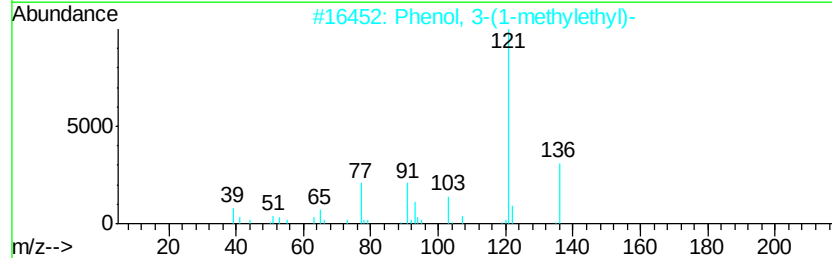
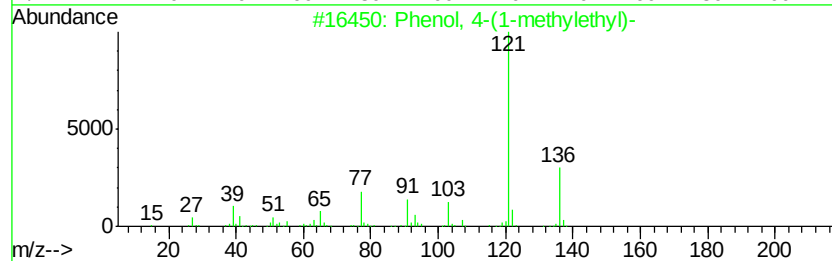
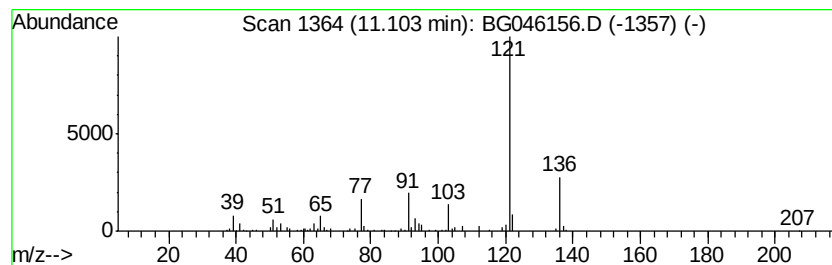
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG073020.M
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Phenol, 4-(1-methylethyl)- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.10	4.06 ng	170465	Napthalene-d8	10.75

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol,	4-(1-methylethyl)-	136	C9H12O	000099-89-8	94
2	Phenol,	3-(1-methylethyl)-	136	C9H12O	000618-45-1	94
3	Phenol,	2-(1-methylethyl)-	136	C9H12O	000088-69-7	91
4	Phenol,	2-ethyl-5-methyl-	136	C9H12O	001687-61-2	80
5	Phenol,	4-ethyl-3-methyl-	136	C9H12O	001123-94-0	80



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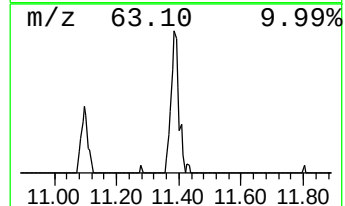
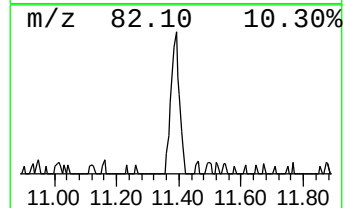
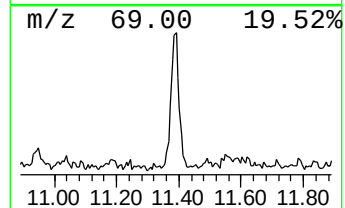
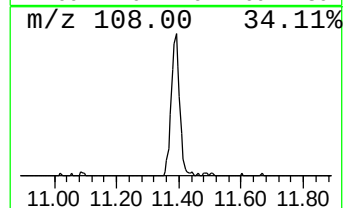
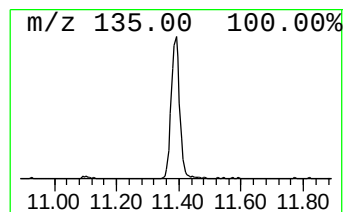
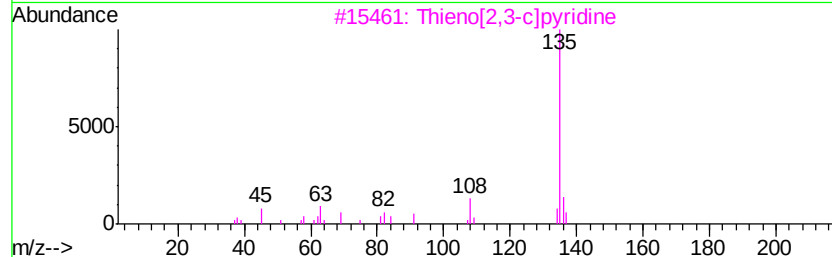
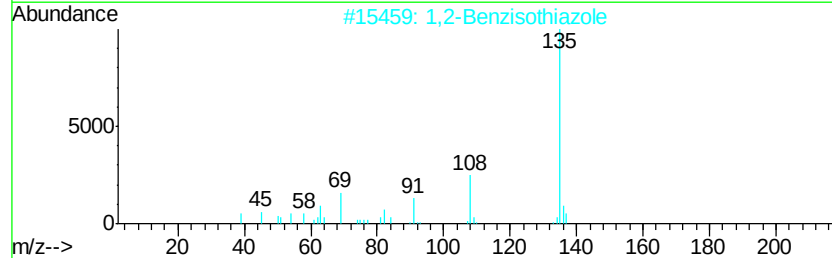
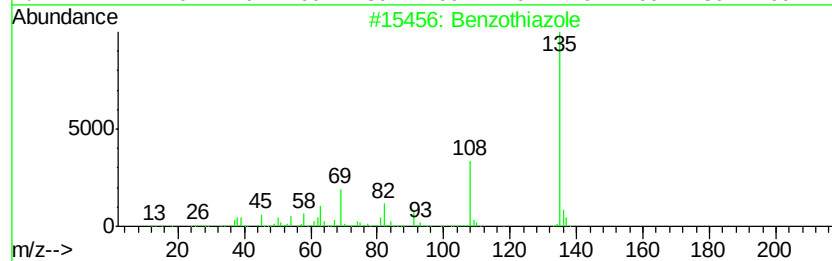
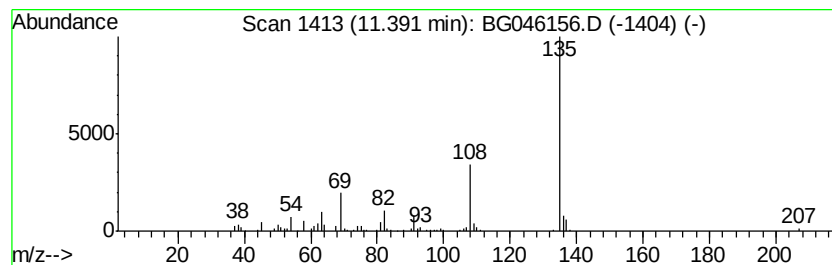
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG073020.M
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TIC Library : C:\Database\NIST11.L
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Peak Number 3 Benzothiazole Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.39	3.93 ng	165166	Napthalene-d8	10.75

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzothiazole	135	C7H5NS	000095-16-9	94
2		1,2-Benzisothiazole	135	C7H5NS	000272-16-2	91
3		Thieno[2,3-c]pyridine	135	C7H5NS	000272-12-8	80
4		3-Allylbenzothiazolium bromide	255	C10H10BrNS	016407-55-9	72
5		1H-Pyrazolo[3,4-d]pyrimidin-4-amine	135	C5H5N5	002380-63-4	59



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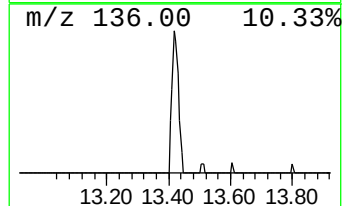
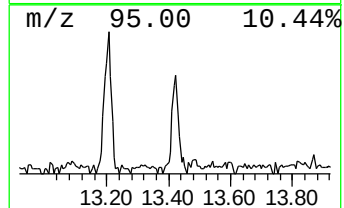
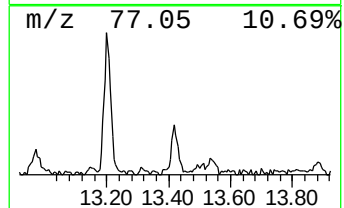
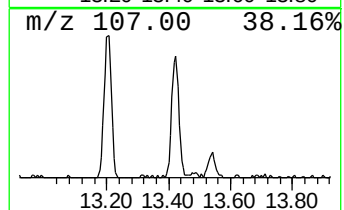
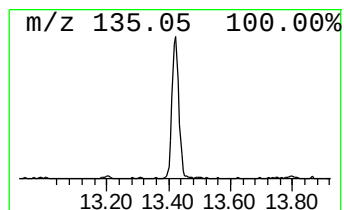
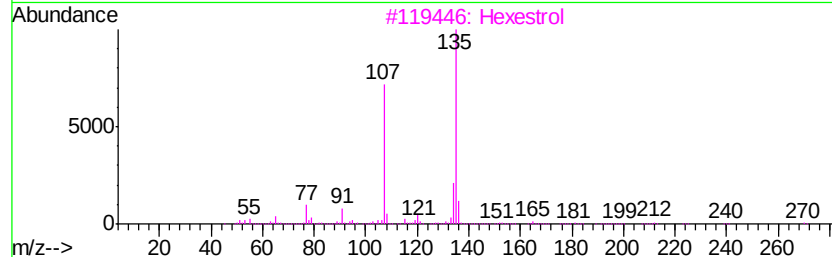
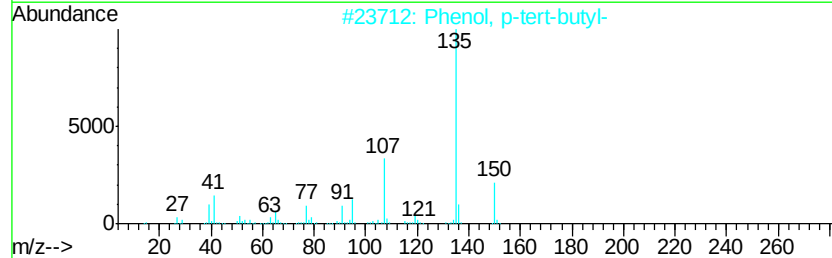
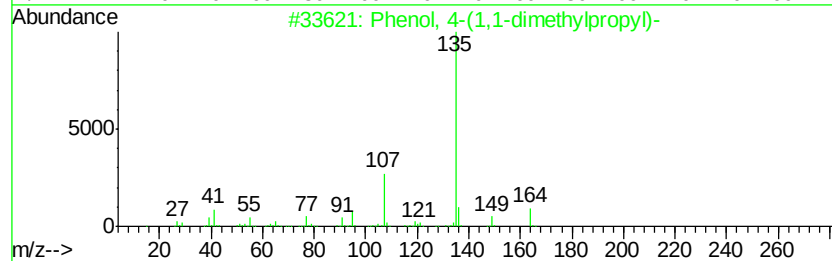
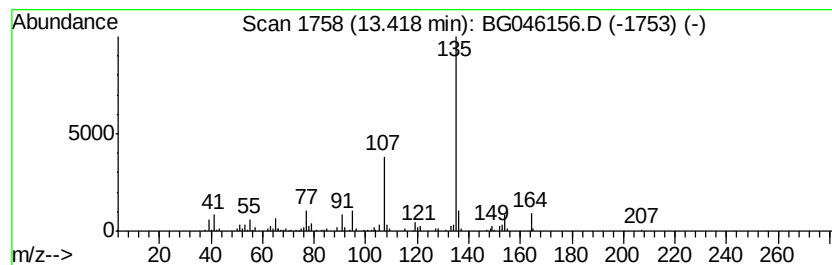
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Phenol, 4-(1,1-dimethylprop... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.42	2.10 ng	123617	Acenaphthene-d10	14.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	94
2		Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	64
3		Hexestrol	270	C18H22O2	000084-16-2	64
4		Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	64
5		Hexestrol	270	C18H22O2	005635-50-7	59



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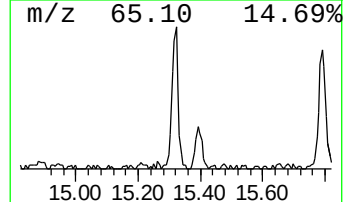
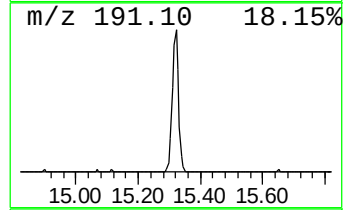
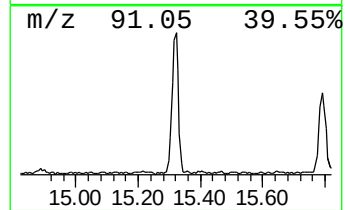
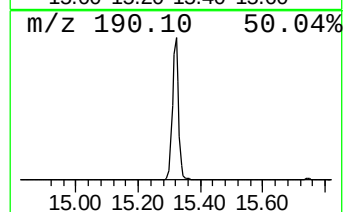
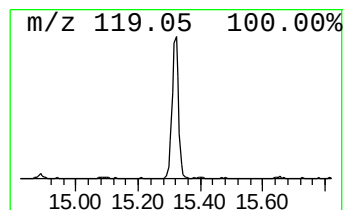
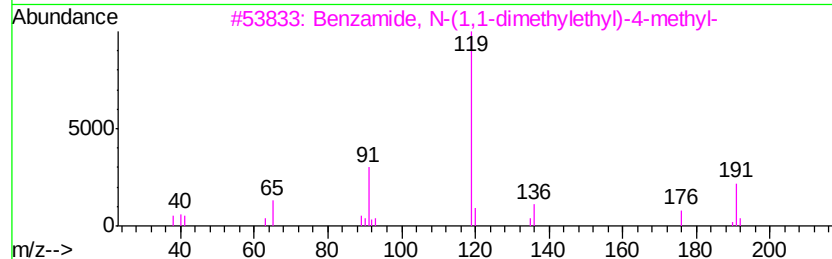
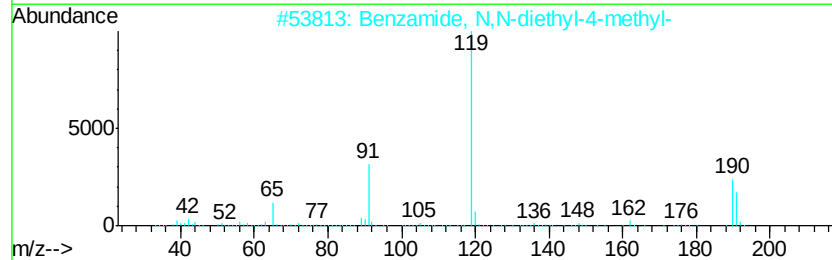
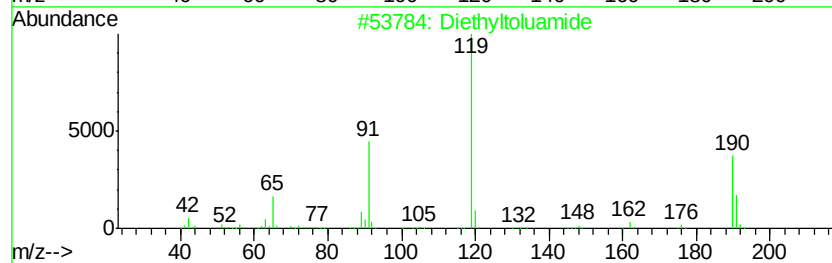
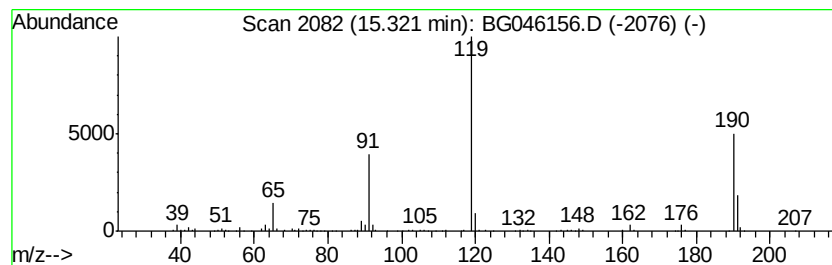
Instrument :
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ClientSampleId :
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TIC Library : C:\Database\NIST11.L
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Peak Number 5 Diethyltoluamide Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.32	4.11 ng	241954	Acenaphthene-d10	14.57	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Diethyltoluamide	191	C12H17NO	000134-62-3	96
2	Benzamide, N,N-diethyl-4-methyl-	191	C12H17NO	002728-05-4	93
3	Benzamide, N-(1,1-dimethylethyl)-4-methyl-	191	C12H17NO	042498-32-8	58
4	m-Toluic acid, 2-butyl ester	192	C12H16O2	005448-57-7	53
5	2,5-Pyrrolidinedione, 1-[(2-methoxy-2-oxoethyl)-	233	C12H11NO4	110920-14-4	50



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG080320\
Data File : BG046156.D
Acq On : 3 Aug 2020 19:23
Operator : CG/JU
Sample : L3536-01
Misc :
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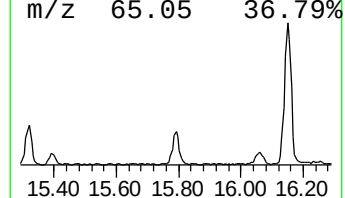
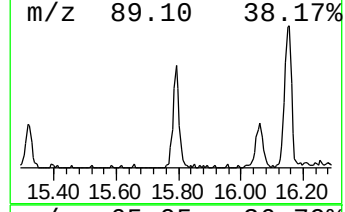
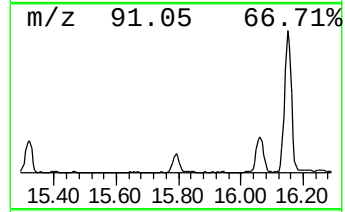
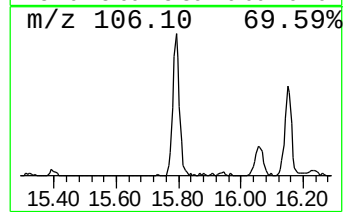
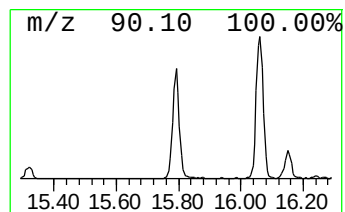
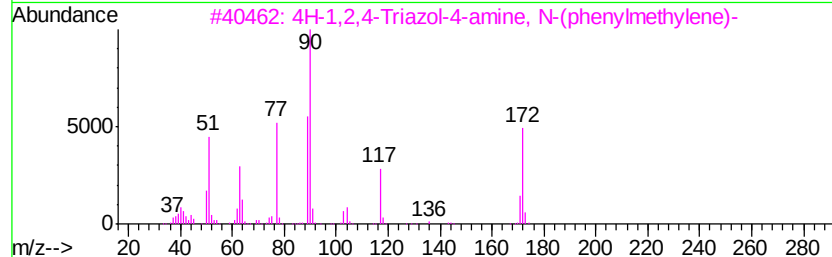
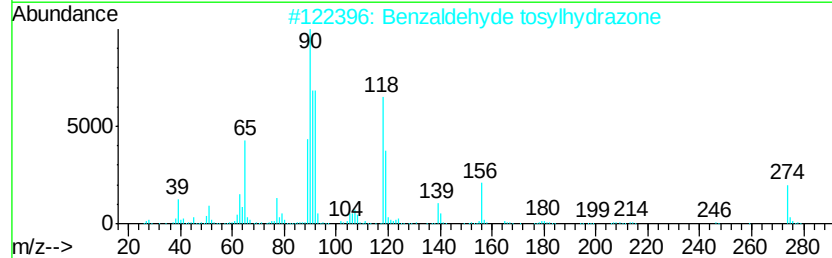
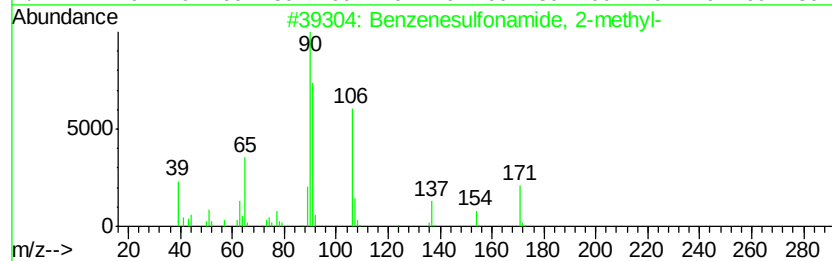
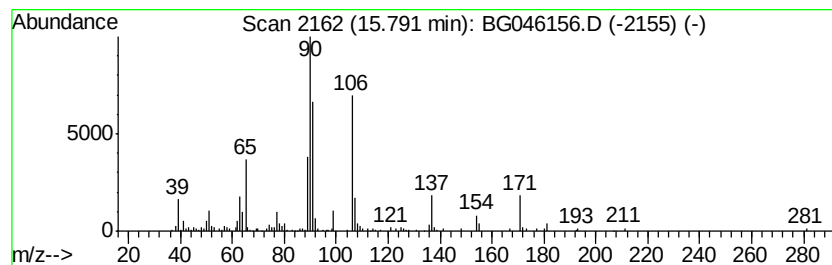
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Benzenesulfonamide, 2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD		R.T.	
15.79	2.75 ng	161783	Acenaphthene-d10		14.57	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzenesulfonamide, 2-methyl-	171	C7H9NO2S	000088-19-7	94
2		Benzaldehyde tosylhydrazone	274	C14H14N2O2S	001666-17-7	72
3		4H-1,2,4-Triazol-4-amine, N-(phe...	172	C9H8N4	018998-48-6	45
4		Bicyclo[4.1.0]hepta-1,3,5-triene	90	C7H6	004646-69-9	45
5		Thiourea, methyl-	90	C2H6N2S	000598-52-7	40



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG080320\
Data File : BG046156.D
Acq On : 3 Aug 2020 19:23
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Sample : L3536-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

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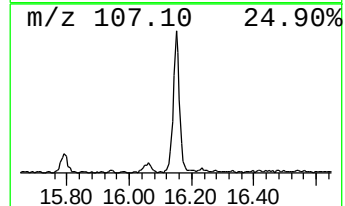
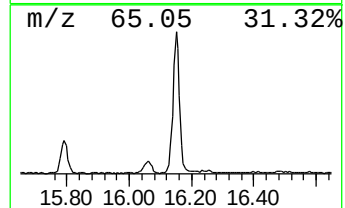
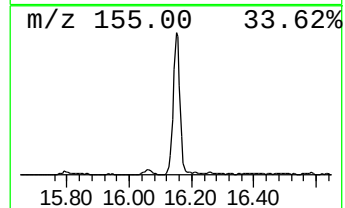
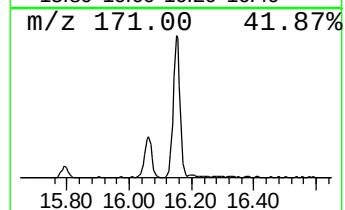
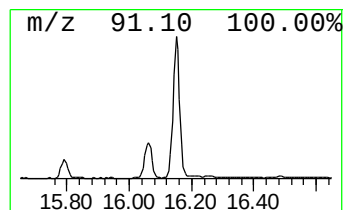
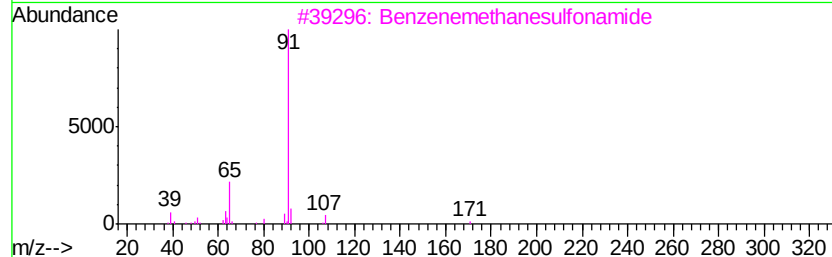
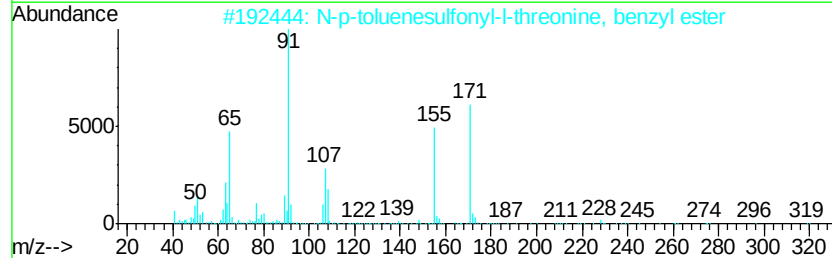
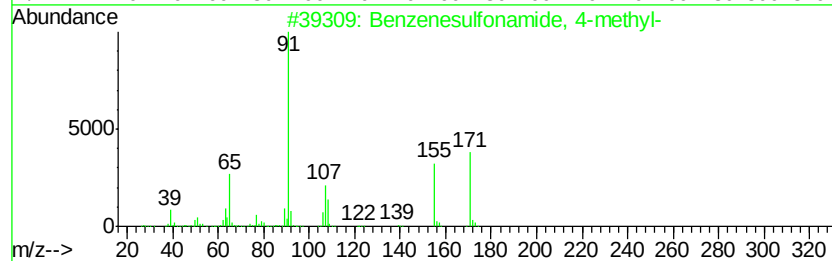
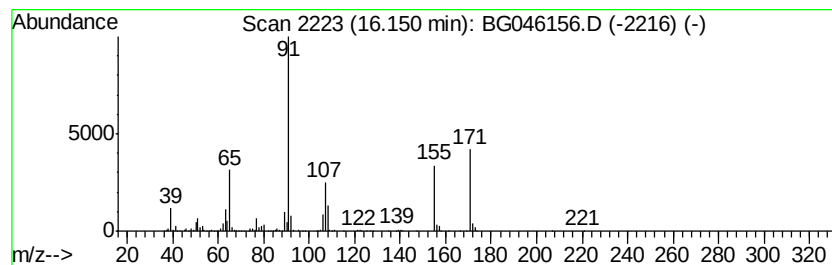
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Benzenesulfonamide, 4-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.15	7.67 ng	545708	Phenanthrene-d10	17.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzenesulfonamide, 4-methyl-	171	C7H9NO2S	000070-55-3	98
2		N-p-toluenesulfonyl-l-threonine,...	363	C18H21NO5S	1000130-34-4	90
3		Benzenemethanesulfonamide	171	C7H9NO2S	004563-33-1	47
4		4-Toluenesulfonylmethyl isocyanide	195	C9H9NO2S	036635-61-7	38
5		N-carbobenzyloxy-cysteine	255	C11H13NO4S	1000131-22-5	22



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG080320\
Data File : BG046156.D
Acq On : 3 Aug 2020 19:23
Operator : CG/JU
Sample : L3536-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
JTY1-GW-01

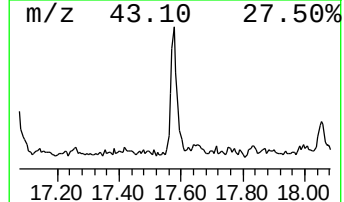
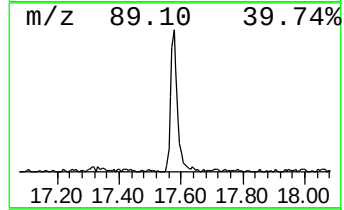
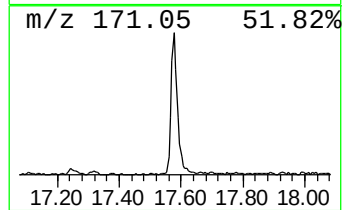
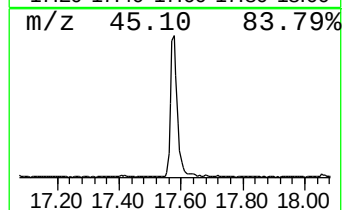
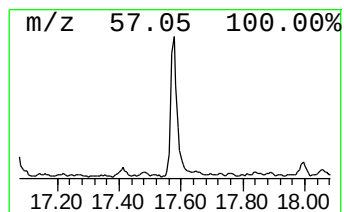
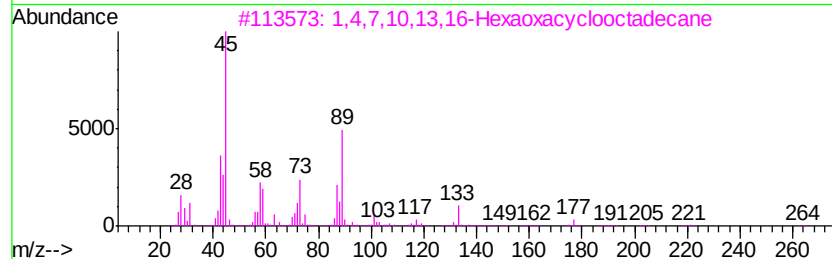
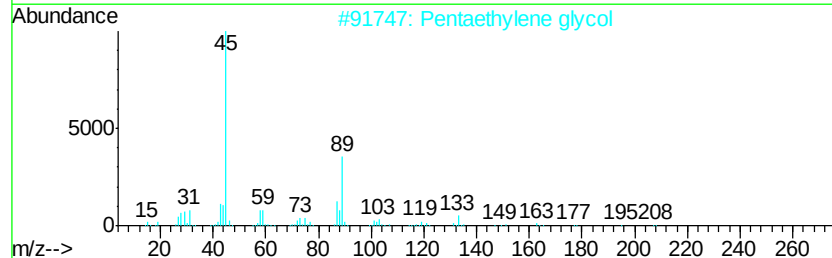
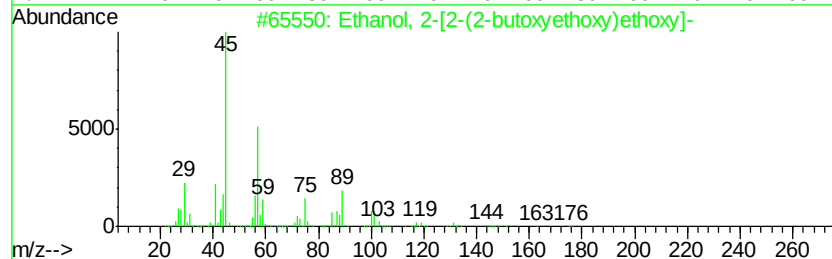
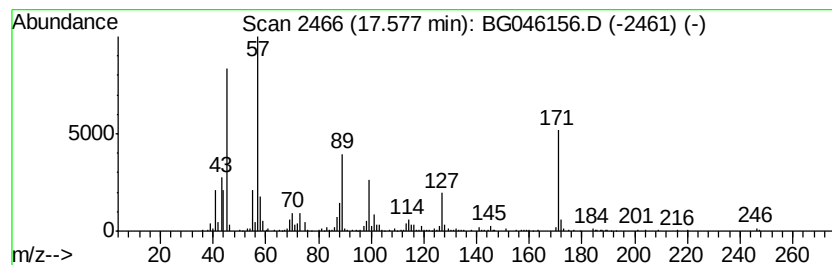
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 unknown17.58 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.58	5.83 ng	414394	Phenanthrene-d10	17.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethanol, 2-[2-(2-butoxyethoxy)et...	206	C10H22O4	000143-22-6	47
2		Pentaethylene glycol	238	C10H22O6	004792-15-8	43
3		1,4,7,10,13,16-Hexaoxacyclooctad...	264	C12H24O6	017455-13-9	38
4		Hexaethylene glycol	282	C12H26O7	002615-15-8	37
5		Tetraethylene glycol	194	C8H18O5	000112-60-7	32



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG080320\
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Acq On : 3 Aug 2020 19:23
Operator : CG/JU
Sample : L3536-01
Misc :
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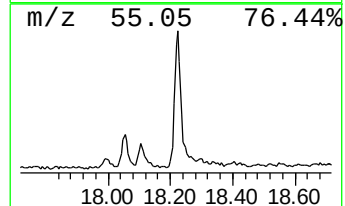
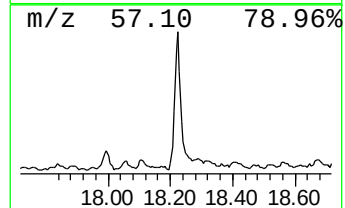
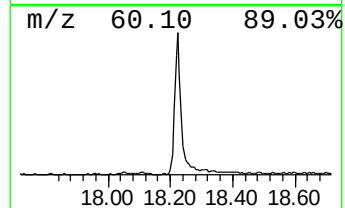
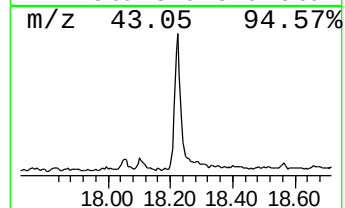
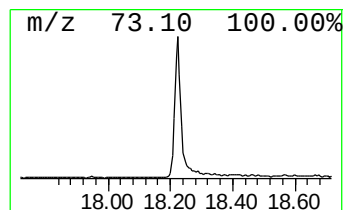
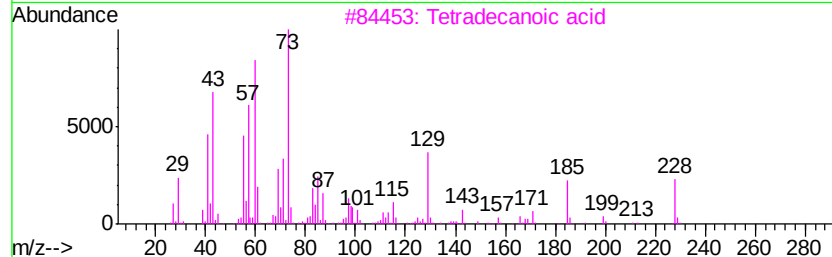
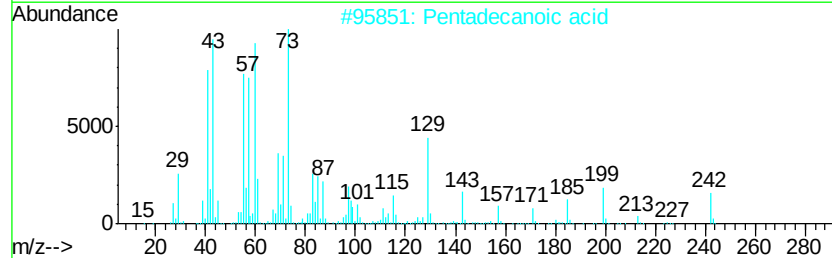
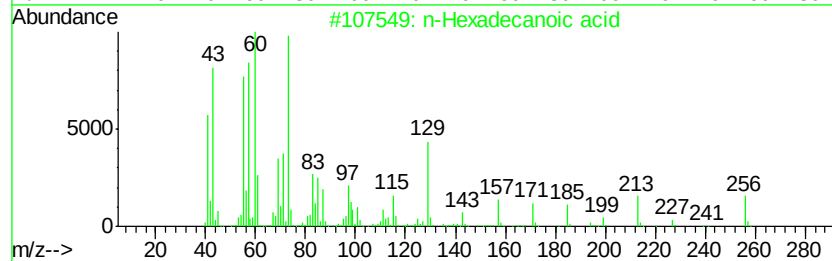
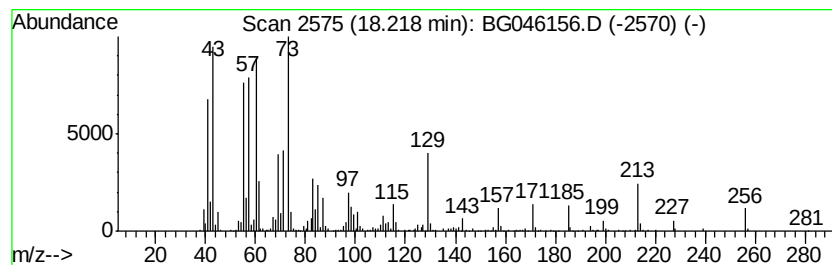
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.22	8.84 ng	628921	Phenanthrene-d10	17.32	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	87
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	80
4	Tridecanoic acid	214	C13H26O2	000638-53-9	76
5	Estra-1,3,5(10)-trien-17.beta.-ol	256	C18H24O	002529-64-8	30



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG080320\
Data File : BG046156.D
Acq On : 3 Aug 2020 19:23
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Sample : L3536-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
JTY1-GW-01

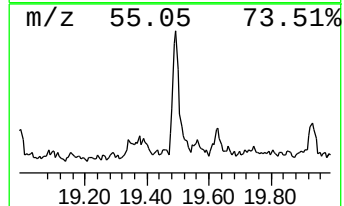
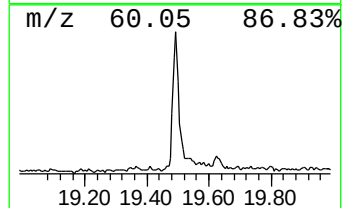
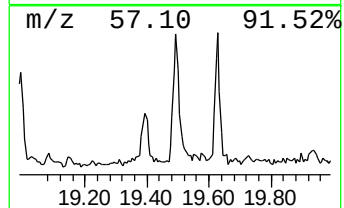
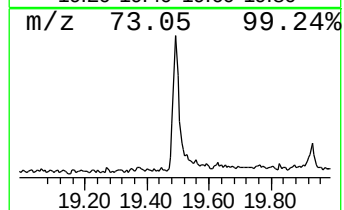
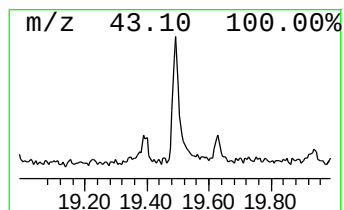
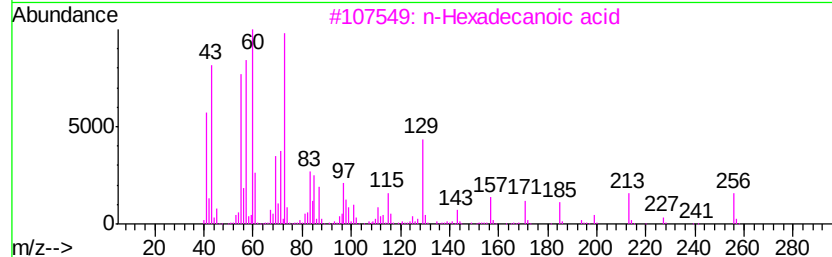
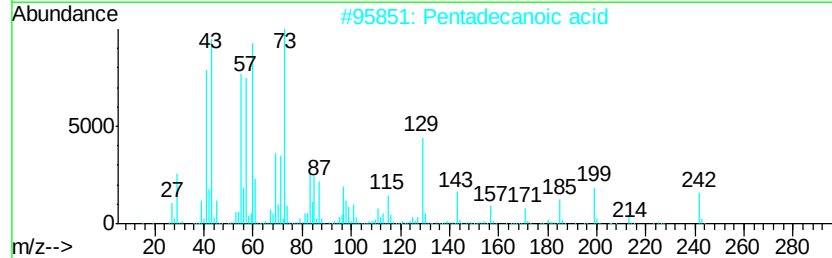
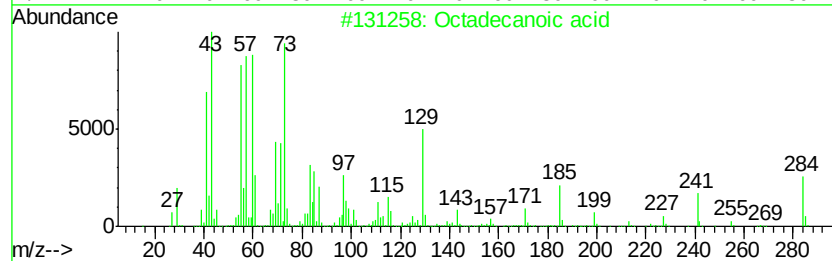
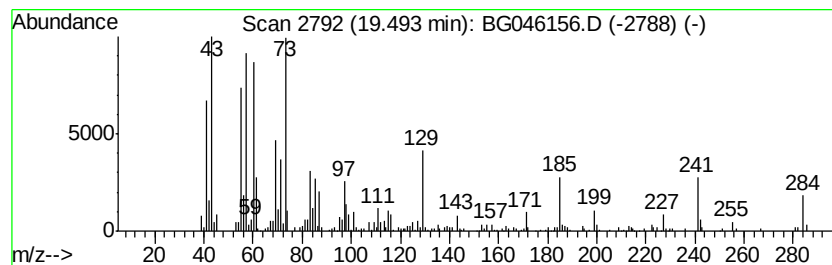
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Octadecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.49	3.18 ng	267621	Chrysene-d12	21.56

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octadecanoic acid	284	C18H36O2	000057-11-4	99
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	87
3		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	87
4		Octadecanoic acid, 2-(2-hydroxye...	372	C22H44O4	000106-11-6	80
5		Tetradecanoic acid	228	C14H28O2	000544-63-8	72



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG080320\
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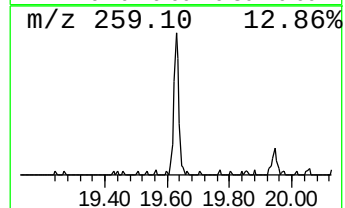
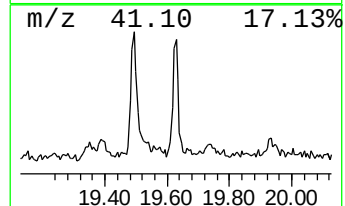
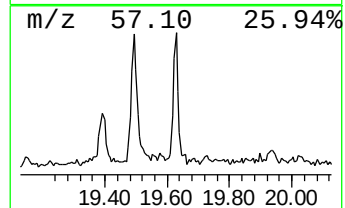
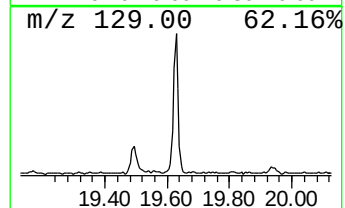
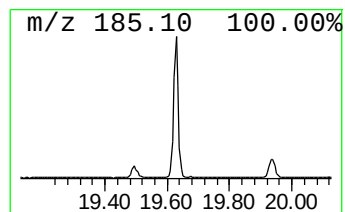
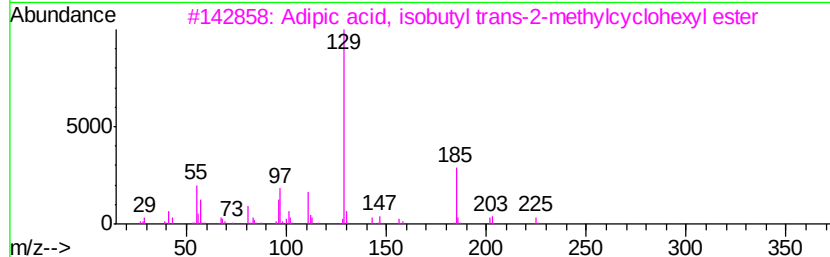
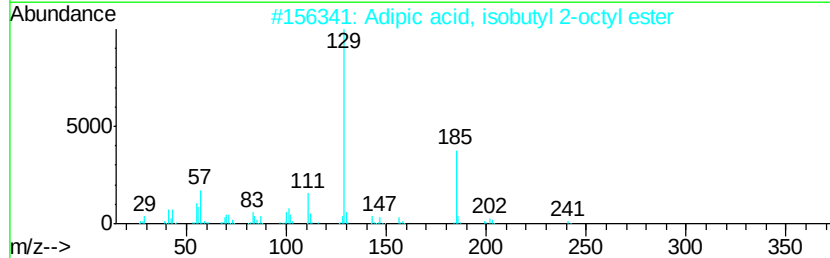
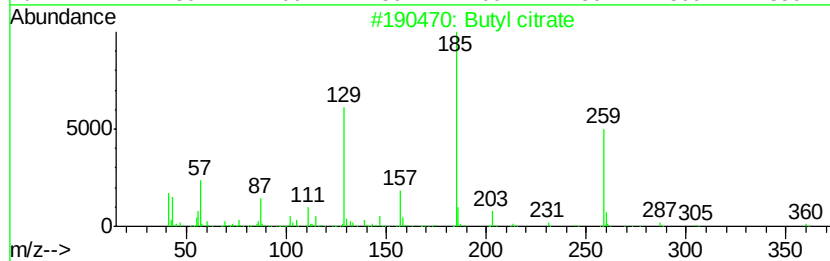
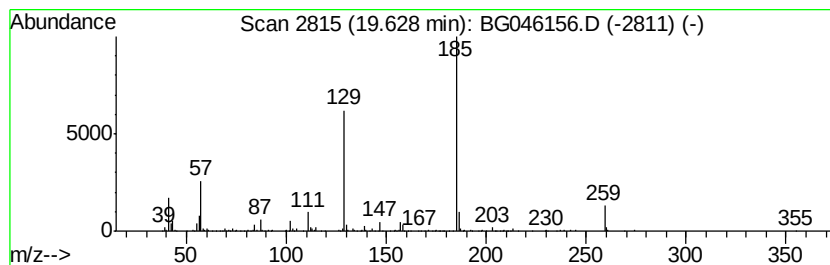
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Butyl citrate Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.63	2.44 ng	205416	Chrysene-d12	21.56
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Butyl citrate	360 C18H32O7	000077-94-1 78	
2	Adipic acid, isobutyl 2-octyl ester	314 C18H34O4	1000324-44-5 56	
3	Adipic acid, isobutyl trans-2-me...	298 C17H30O4	1000324-74-5 56	
4	Adipic acid, butyl 4-methylpent-...	286 C16H30O4	1000353-50-2 56	
5	Adipic acid, 3,3-dimethylbut-2-y...	286 C16H30O4	1000353-66-7 56	



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG080320\
Data File : BG046156.D
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Misc :
ALS Vial : 9 Sample Multiplier: 1

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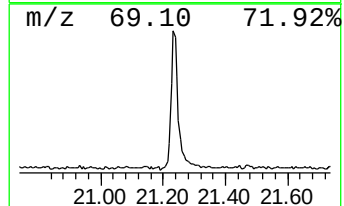
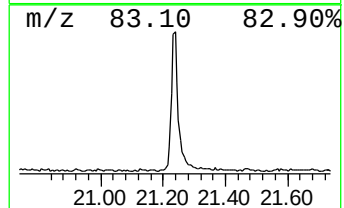
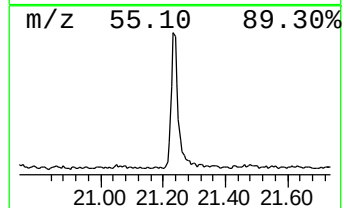
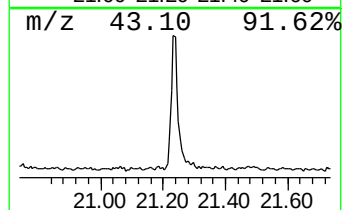
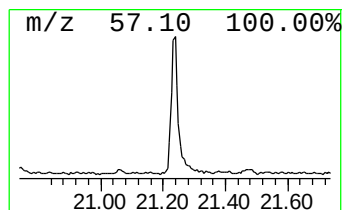
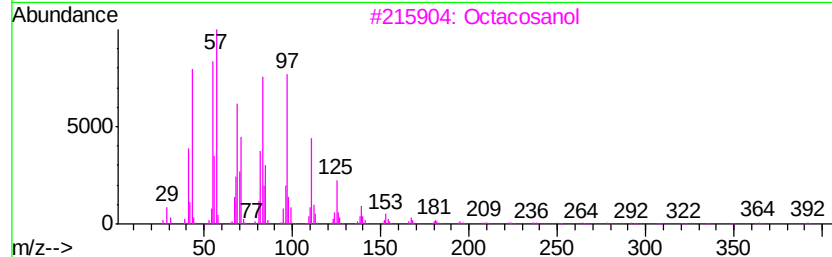
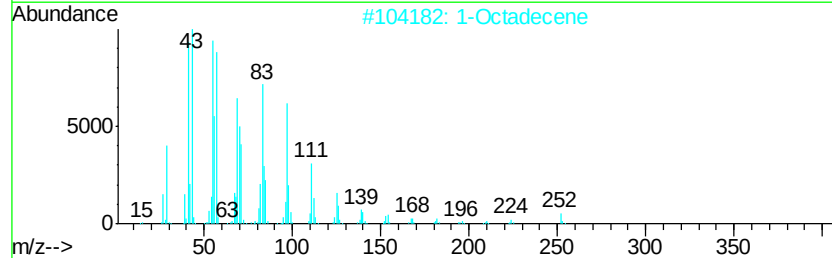
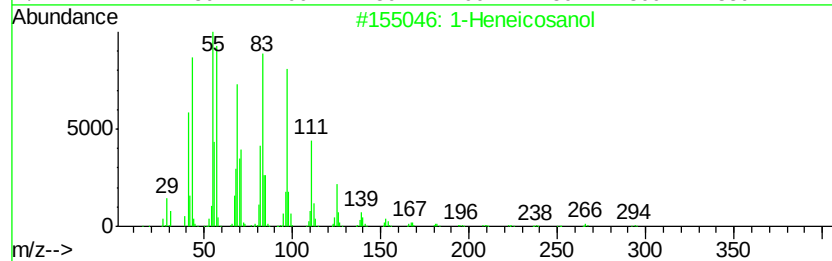
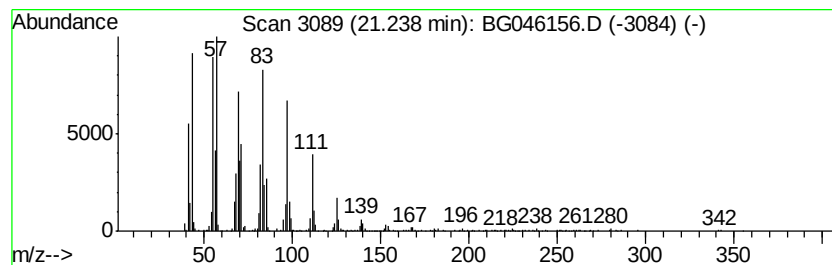
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 1-Heneicosanol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.24	7.49 ng	631132	Chrysene-d12	21.56

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Heneicosanol	312	C21H44O	015594-90-8	95
2			1-Octadecene	252	C18H36	000112-88-9	94
3			Octacosanol	410	C28H58O	000557-61-9	94
4			Pentafluoropropionic acid, hexad...	388	C19H33F5O2	006222-07-7	94
5			Behenic alcohol	326	C22H46O	000661-19-8	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG080320\
Data File : BG046156.D
Acq On : 3 Aug 2020 19:23
Operator : CG/JU
Sample : L3536-01
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
JTY1-GW-01

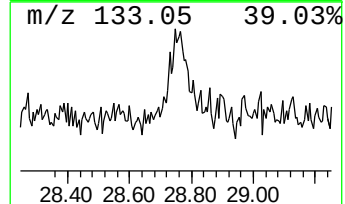
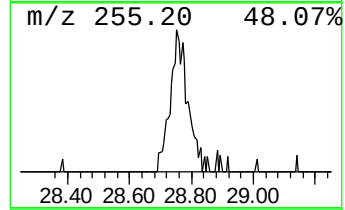
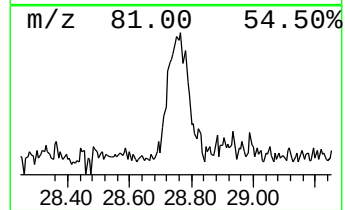
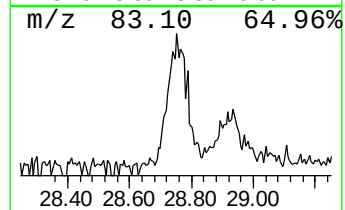
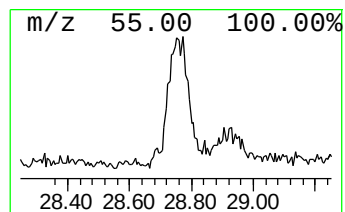
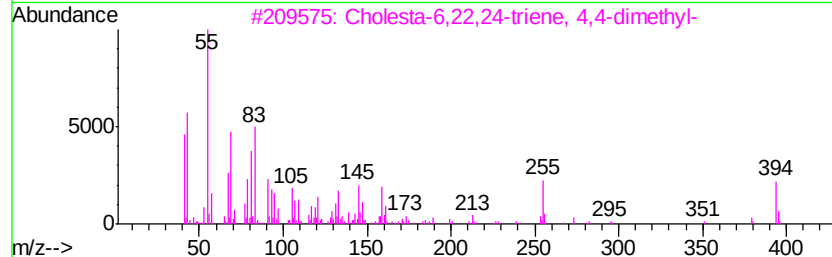
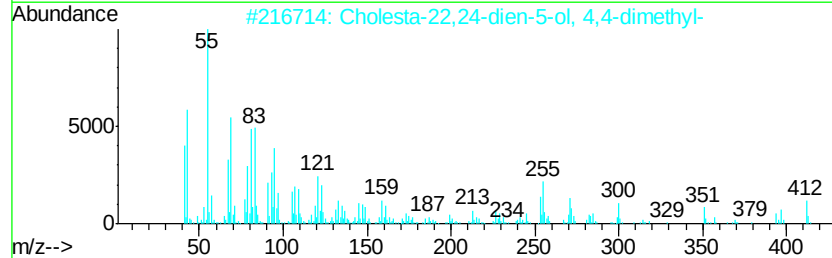
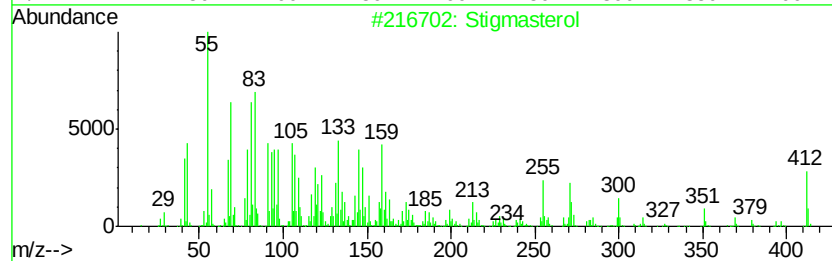
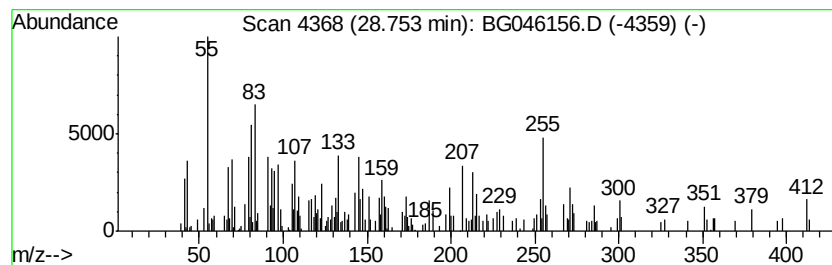
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Stigmasterol Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
28.75	2.67 ng	233588	Perylene-d12	24.67

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Stigmasterol	412	C29H48O	000083-48-7	95
2			Cholesta-22,24-dien-5-ol, 4,4-di...	412	C29H48O	1000128-66-1	89
3			Cholesta-6,22,24-triene, 4,4-dim...	394	C29H46	1000128-66-9	25
4			22-Stigmasten-3-one	412	C29H48O	004736-95-2	18
5			26,27-Dinorcholesta-5,22-dien-3-...	356	C25H40O	057597-10-1	14



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG080320\
 Data File : BG046156.D
 Acq On : 3 Aug 2020 19:23
 Operator : CG/JU
 Sample : L3536-01
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 JTY1-GW-01

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG073020.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...	3.16	82.7	ng	2329110	1	7.95	563403	20.0
Phenol, 4-(1-meth...	11.10	4.1	ng	170465	2	10.75	840586	20.0
Benzothiazole	11.39	3.9	ng	165166	2	10.75	840586	20.0
Phenol, 4-(1,1-di...	13.42	2.1	ng	123617	3	14.57	1177660	20.0
Diethyltoluamide	15.32	4.1	ng	241954	3	14.57	1177660	20.0
Benzenesulfonamid...	15.79	2.8	ng	161783	3	14.57	1177660	20.0
Benzenesulfonamid...	16.15	7.7	ng	545708	4	17.32	1422370	20.0
unknown17.58	17.58	5.8	ng	414394	4	17.32	1422370	20.0
n-Hexadecanoic acid	18.22	8.8	ng	628921	4	17.32	1422370	20.0
Octadecanoic acid	19.49	3.2	ng	267621	5	21.56	1685700	20.0
Butyl citrate	19.63	2.4	ng	205416	5	21.56	1685700	20.0
1-Heneicosanol	21.24	7.5	ng	631132	5	21.56	1685700	20.0
Stigmasterol	28.75	2.7	ng	233588	6	24.67	1748290	20.0