

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG060719\
 Data File : BG041122.D
 Acq On : 8 Jun 2019 6:39
 Operator : HP/JU
 Sample : K3191-07
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 CULVER-CBTC-TP1-060419

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0
 Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

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

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.183	12	16	23	rVB	153529	213175	5.73%	0.953%
2	5.639	427	434	444	rBV	305842	498766	13.40%	2.229%
3	7.243	700	707	715	rBV	211162	379690	10.20%	1.697%
4	7.631	765	773	787	rBV	585672	1014718	27.26%	4.535%
5	8.107	847	854	859	rBV	144694	255028	6.85%	1.140%
6	8.430	901	909	919	rVB	534495	922995	24.80%	4.125%
7	9.282	1045	1054	1062	rBV	490345	876915	23.56%	3.919%
8	9.382	1063	1071	1082	rVB2	70285	165328	4.44%	0.739%
9	10.927	1327	1334	1347	rBV	190576	372557	10.01%	1.665%
10	11.697	1458	1465	1479	rBV	62064	129879	3.49%	0.580%
11	12.702	1630	1636	1643	rBV5	16870	38978	1.05%	0.174%
12	12.990	1680	1685	1695	rBV4	20011	39122	1.05%	0.175%
13	13.360	1738	1748	1757	rBV	1467971	2267127	60.90%	10.133%
14	14.176	1882	1887	1894	rBV2	51891	87086	2.34%	0.389%
15	14.735	1974	1982	1989	rBV2	356024	573706	15.41%	2.564%
16	16.221	2228	2235	2251	rBV	1333520	2001812	53.78%	8.947%
17	17.478	2442	2449	2455	rBV2	506125	775397	20.83%	3.466%
18	17.655	2474	2479	2483	rVB	25330	37727	1.01%	0.169%
19	18.330	2589	2594	2604	rBV2	256850	389470	10.46%	1.741%
20	19.458	2782	2786	2788	rBV2	40790	51285	1.38%	0.229%
21	19.488	2788	2791	2801	rVV2	153196	222319	5.97%	0.994%
22	19.599	2805	2810	2822	rVV2	175317	261766	7.03%	1.170%
23	19.740	2822	2834	2842	rVV	3309757	3722487	100.00%	16.637%
24	20.075	2881	2891	2896	rBV	2916902	3714464	99.78%	16.602%
25	20.346	2932	2937	2942	rVB	34202	44726	1.20%	0.200%
26	20.722	2997	3001	3005	rBV6	33774	48360	1.30%	0.216%
27	20.839	3017	3021	3025	rBV2	65292	70600	1.90%	0.316%
28	21.350	3103	3108	3123	rVB2	193480	311369	8.36%	1.392%
29	21.756	3169	3177	3186	rBV2	532379	901666	24.22%	4.030%
30	21.903	3197	3202	3208	rVB	93003	129850	3.49%	0.580%
31	22.520	3302	3307	3318	rBV	103380	167385	4.50%	0.748%
32	23.225	3421	3427	3436	rBV2	103167	196515	5.28%	0.878%
33	23.424	3455	3461	3467	rBV3	24931	46063	1.24%	0.206%
34	24.053	3560	3568	3584	rVB3	91843	213185	5.73%	0.953%

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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 >
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35	25.075	3725	3742	3757	rVB2	293460	1071803	28.79%	4.790%
36	26.180	3922	3930	3939	rVB4	56098	160798	4.32%	0.719%

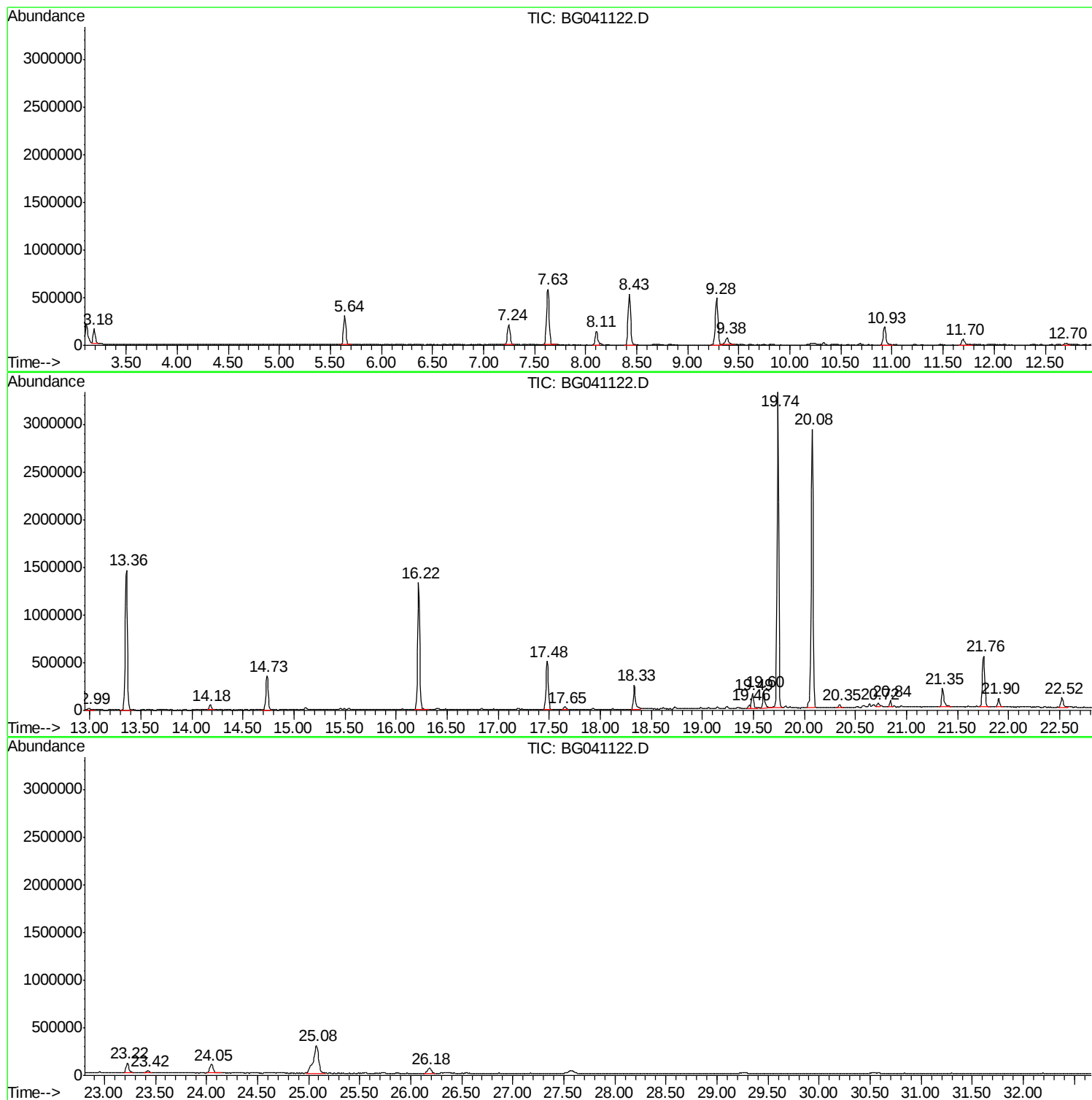
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P



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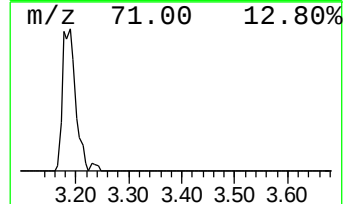
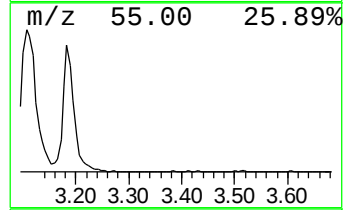
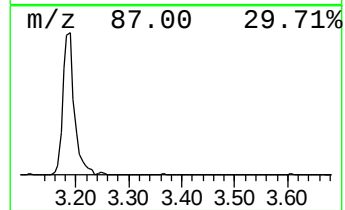
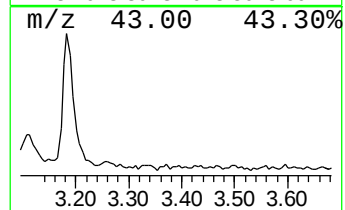
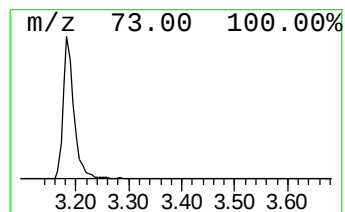
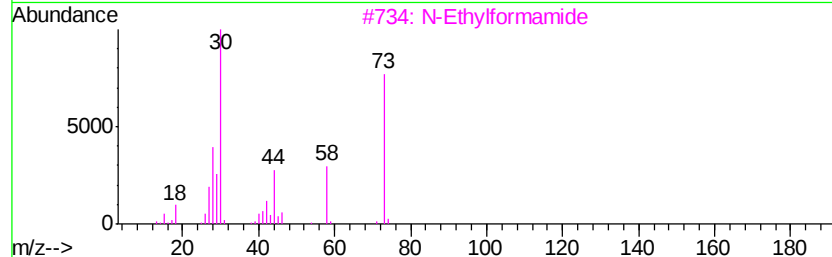
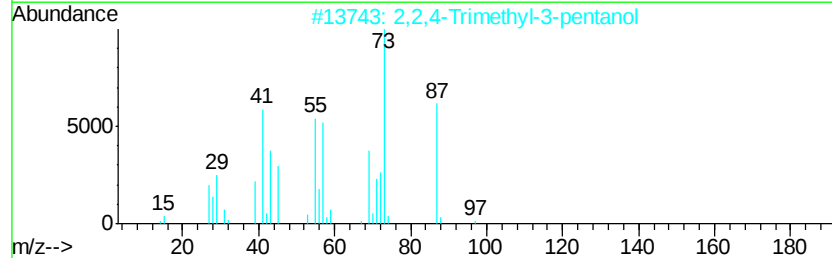
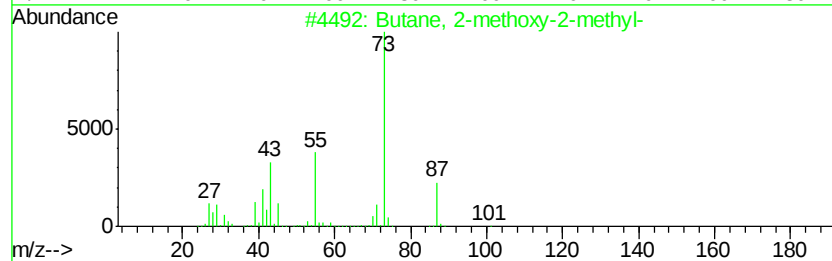
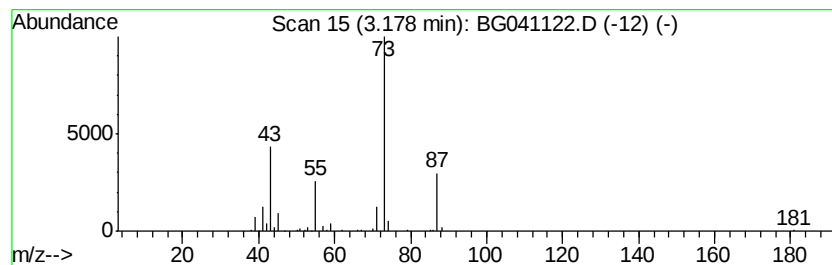
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.18	16.72 ng	213175	1,4-Dichlorobenzene-d4	8.11

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	72
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	64
3		N-Ethylformamide	73	C3H7NO	000627-45-2	9
4		Propane, 2-methoxy-2-methyl-	88	C5H12O	001634-04-4	9
5		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	9



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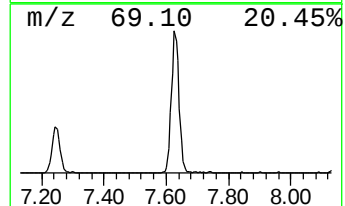
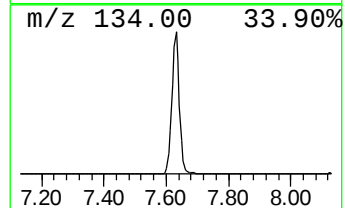
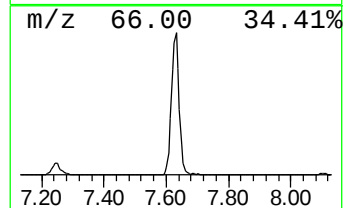
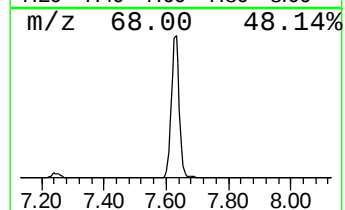
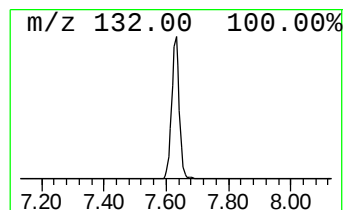
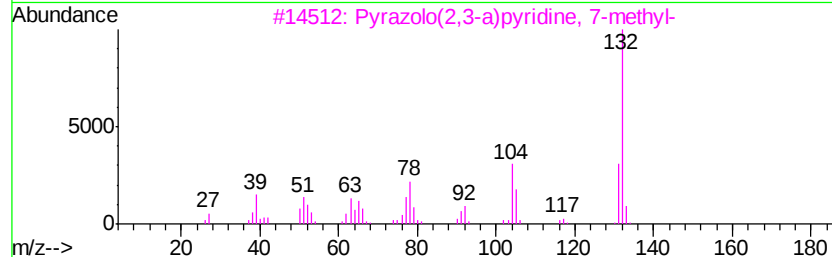
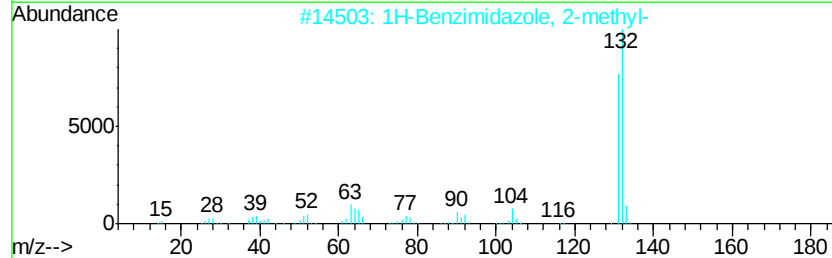
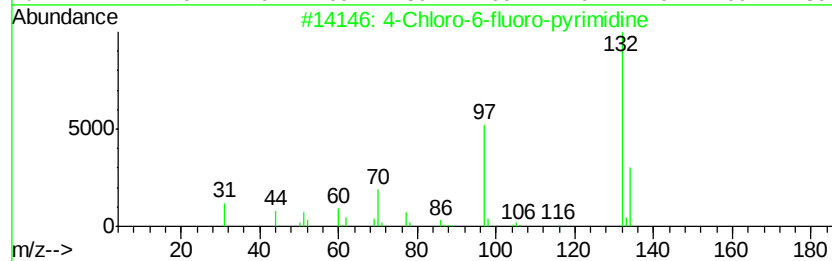
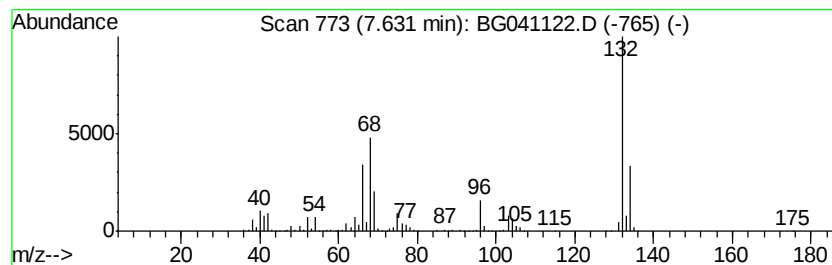
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown7.63 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.63	79.58 ng	1014720	1,4-Dichlorobenzene-d4	8.11

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
2		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
3		Pvrazolo(2,3-a)pvridine, 7-methyl-	132	C8H8N2	034760-58-2	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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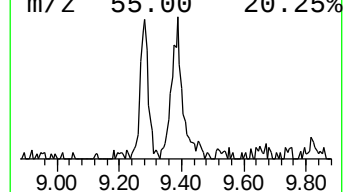
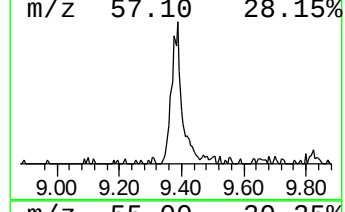
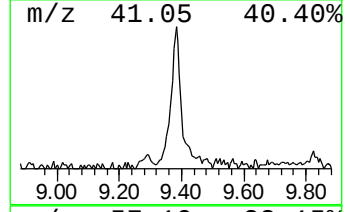
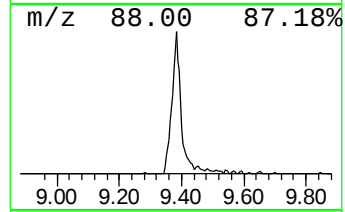
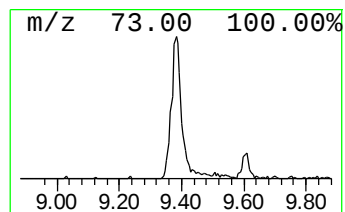
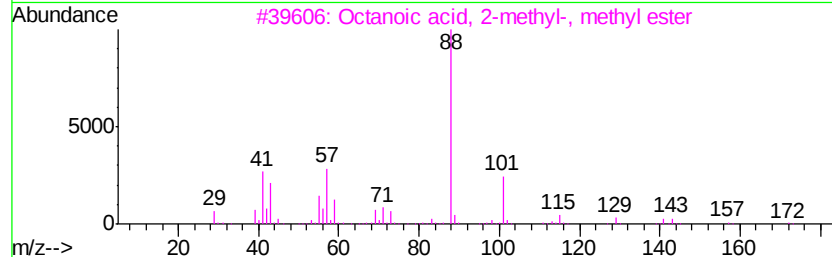
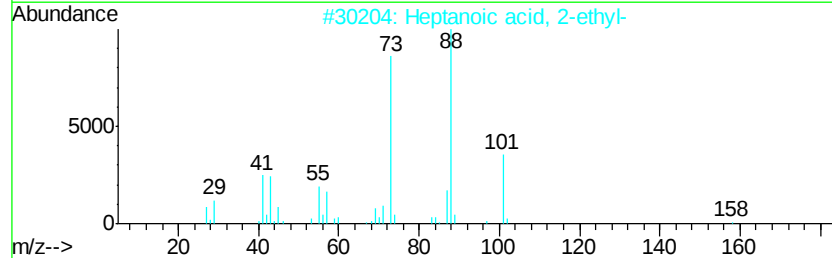
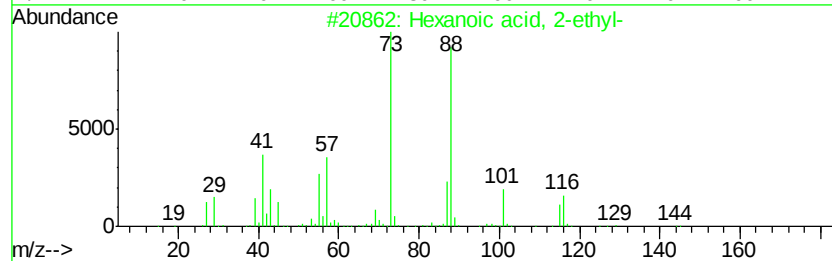
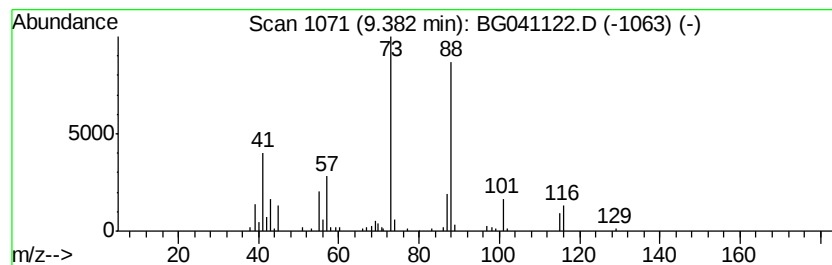
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Peak Number 4 Hexanoic acid, 2-ethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.38	12.97 ng	165328	1,4-Dichlorobenzene-d4	8.11

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	83	
2	Heptanoic acid, 2-ethyl-	158	C9H18O2	003274-29-1	53	
3	Octanoic acid, 2-methyl-, methyl...	172	C10H20O2	002177-86-8	38	
4	Pentanoic acid, 2,4-dimethyl-, m...	144	C8H16O2	071672-33-8	28	
5	Pentanoic acid, 2-methyl-, methy...	130	C7H14O2	002177-77-7	28	



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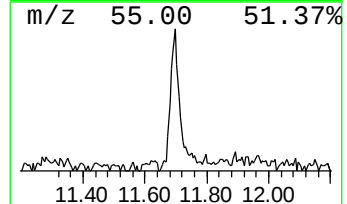
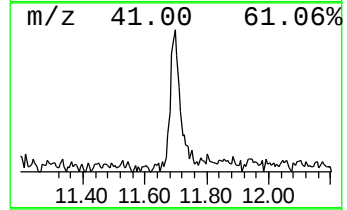
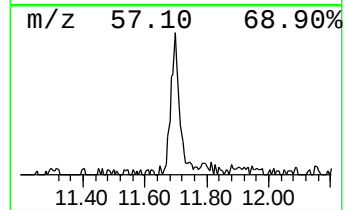
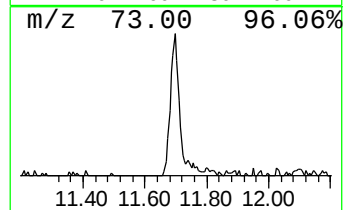
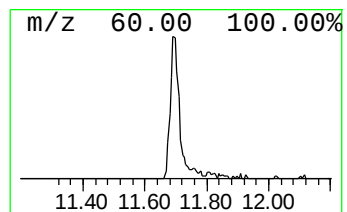
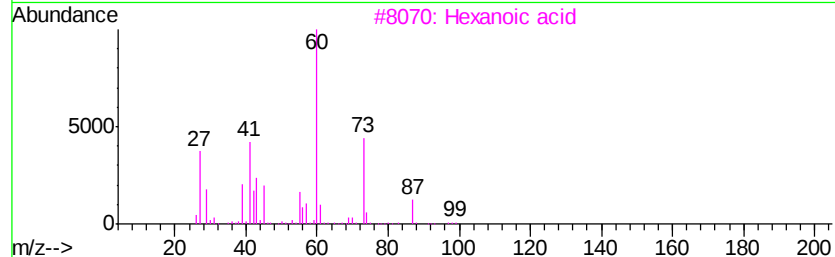
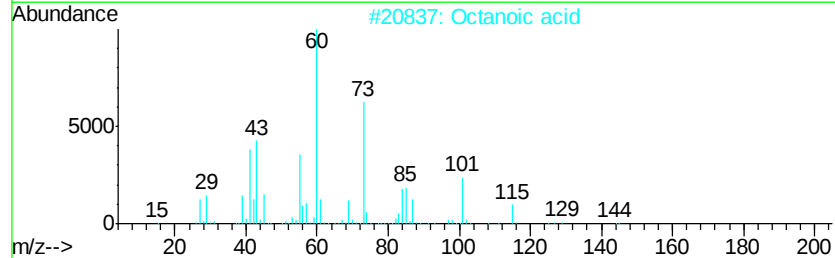
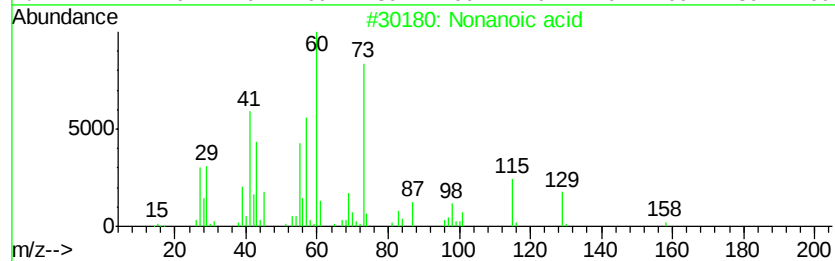
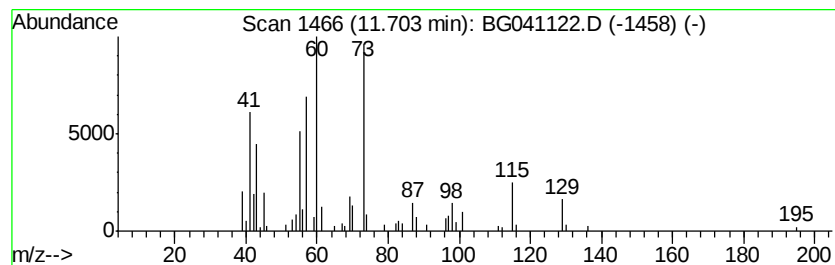
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Peak Number 5 Nonanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.70	6.97 ng	129879	Naphthalene-d8	10.93	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonanoic acid	158	C9H18O2	000112-05-0	72
2	Octanoic acid	144	C8H16O2	000124-07-2	59
3	Hexanoic acid	116	C6H12O2	000142-62-1	50
4	Pentanoic acid, 3-methyl-	116	C6H12O2	000105-43-1	47
5	.alpha.-D-Glucose	180	C6H12O6	000492-62-6	38



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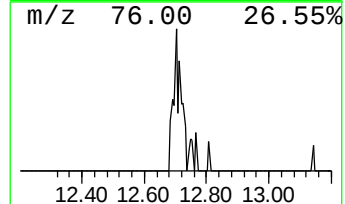
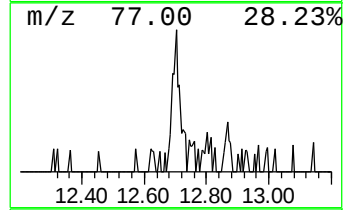
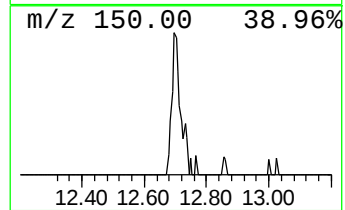
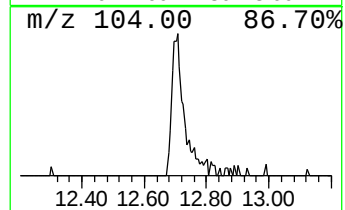
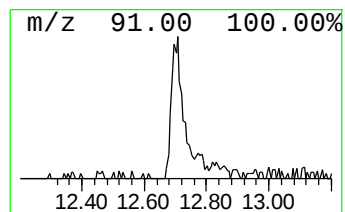
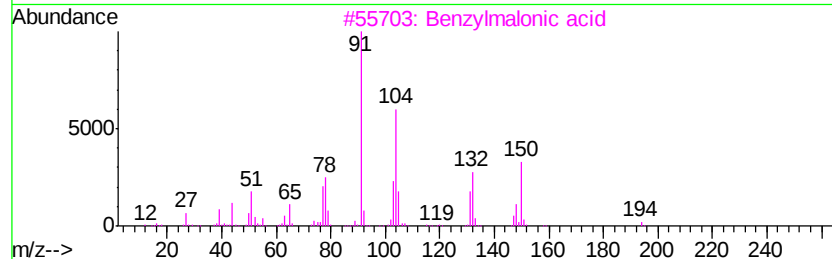
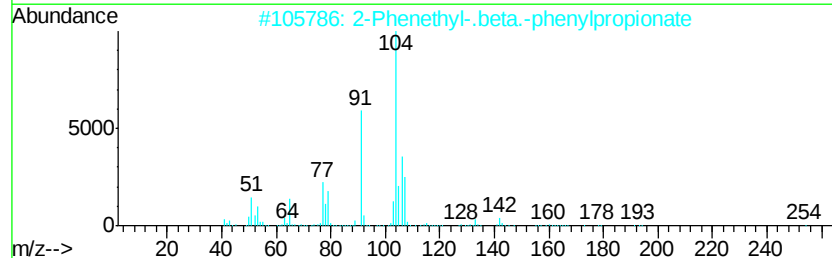
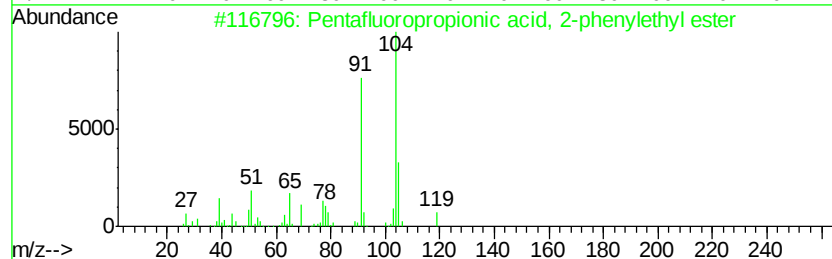
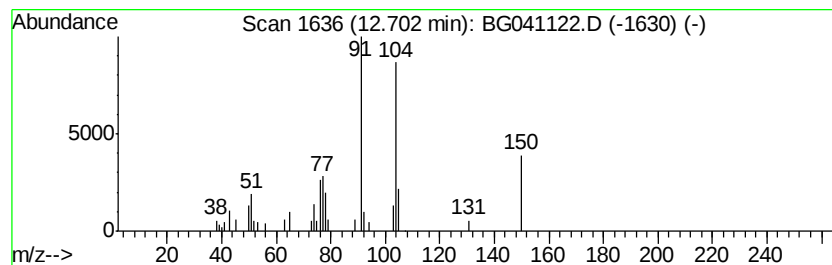
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Peak Number 6 Pentafluoropropionic acid, ... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.70	2.09 ng	38978	Naphthalene-d8	10.93

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Pentafluoropropionic acid, 2-phe...	268	C11H9F5O2	081089-48-7	50
2		2-Phenethyl-.beta.-phenylpropionate	254	C17H18O2	028049-10-7	42
3		Benzylmalonic acid	194	C10H10O4	000616-75-1	42
4		4-Nitrosophenyl-.beta.-phenylpro...	255	C15H13NO3	1000129-40-0	42
5		Styrene	104	C8H8	000100-42-5	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG060719\
Data File : BG041122.D
Acq On : 8 Jun 2019 6:39
Operator : HP/JU
Sample : K3191-07
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
CULVER-CBTC-TP1-060419

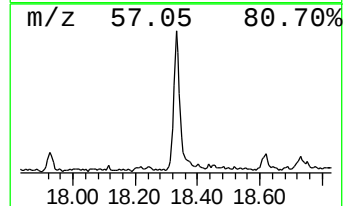
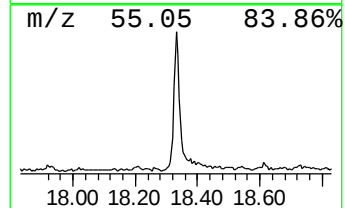
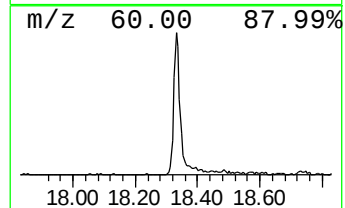
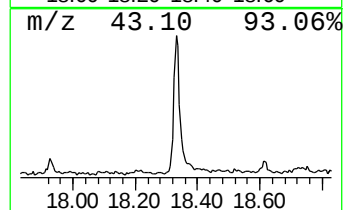
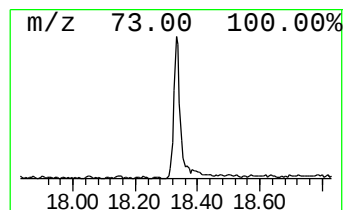
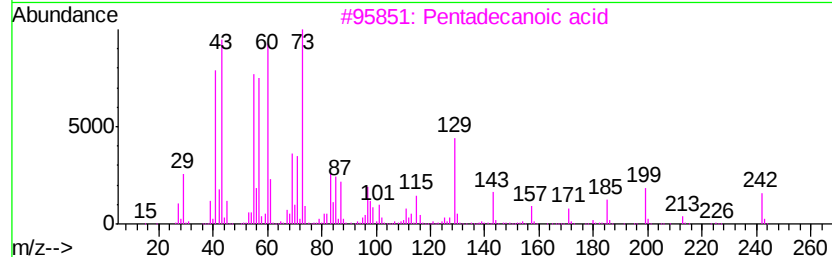
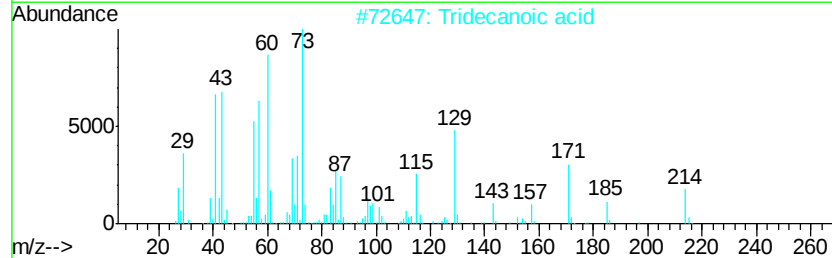
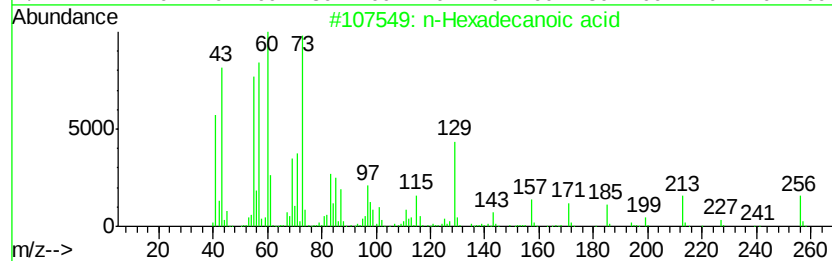
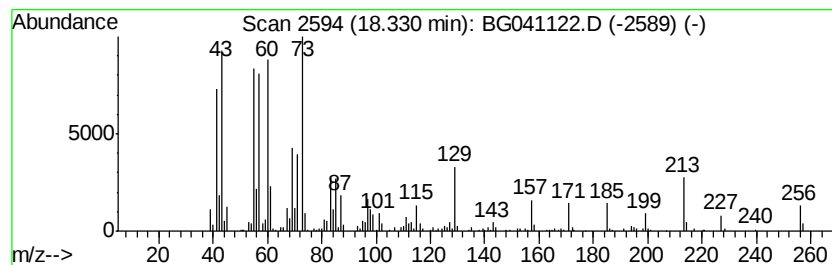
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.33	10.05 ng	389470	Phenanthrene-d10	17.48

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	91
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	70
4		n-Decanoic acid	172	C10H20O2	000334-48-5	25
5		Oxalic acid, propyl tridecyl ester	314	C18H34O4	1000309-26-6	15



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG060719\
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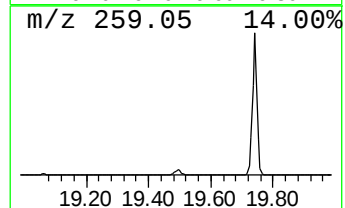
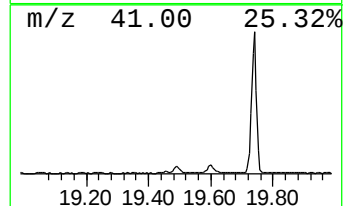
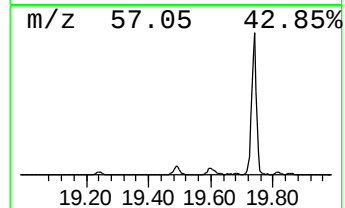
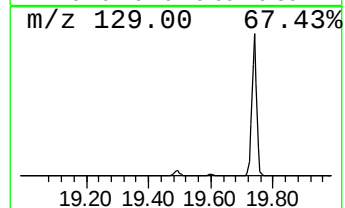
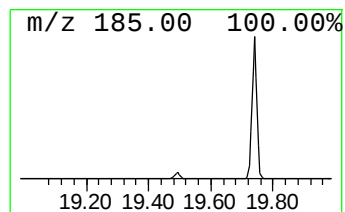
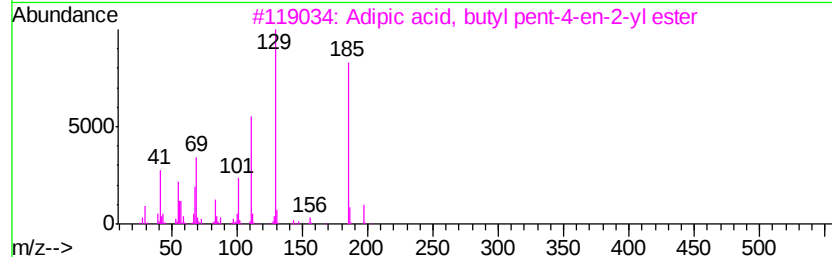
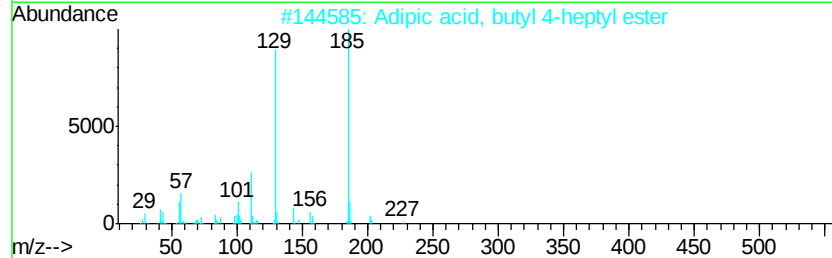
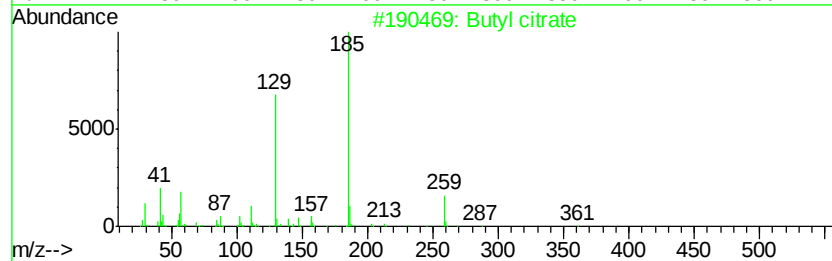
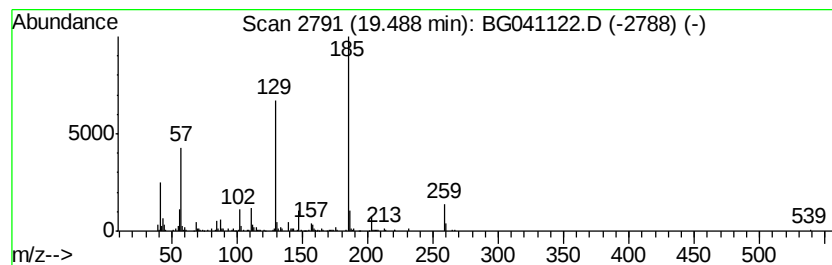
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Adipic acid, butyl 4-heptyl... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.49	5.73 ng	222319	Phenanthrene-d10	17.48	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butyl citrate	360	C18H32O7	000077-94-1	90
2	Adipic acid, butyl 4-heptyl ester	300	C17H32O4	1000349-73-7	59
3	Adipic acid, butyl pent-4-en-2-yl ester	270	C15H26O4	1000354-11-9	53
4	Adipic acid, 3,3-dimethylbut-2-yl ester	286	C16H30O4	1000353-66-7	50
5	Adipic acid, butyl 3-hexyl ester	286	C16H30O4	1000353-62-3	50



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG060719\
Data File : BG041122.D
Acq On : 8 Jun 2019 6:39
Operator : HP/JU
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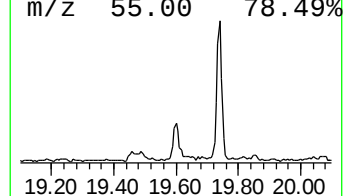
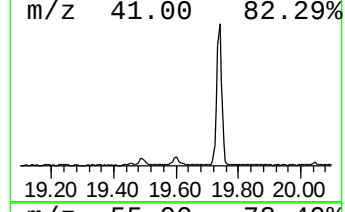
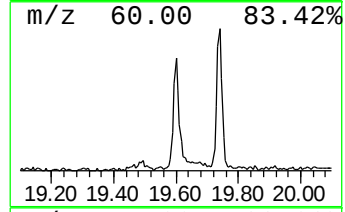
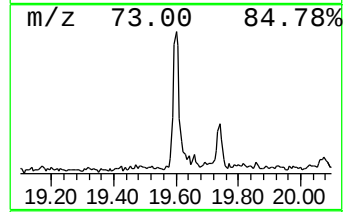
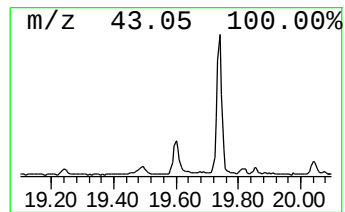
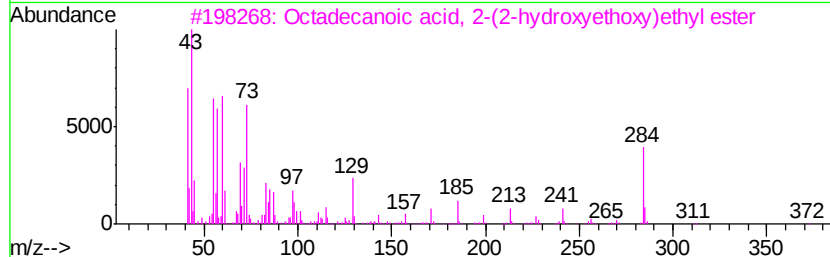
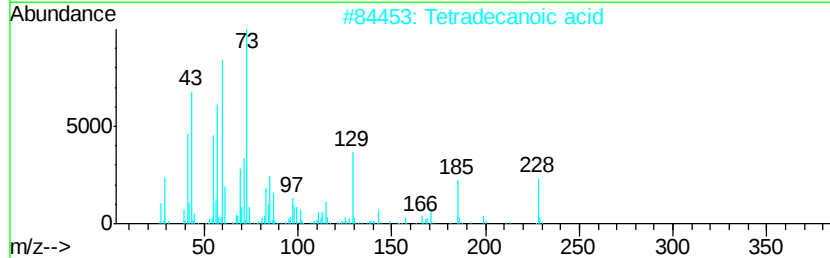
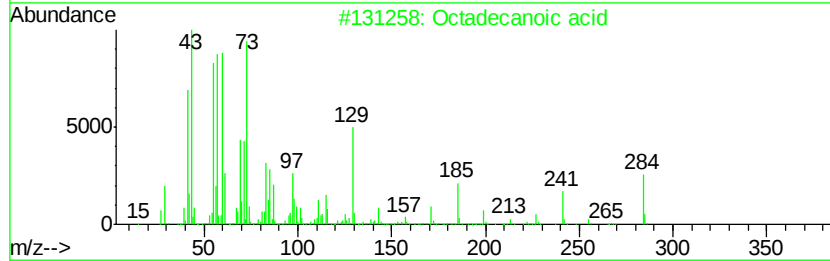
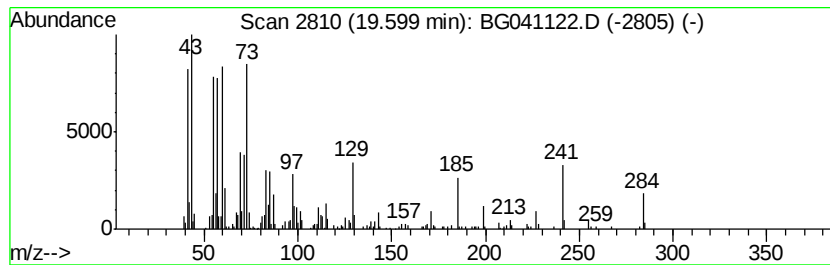
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Octadecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.60	6.75 ng	261766	Phenanthrene-d10	17.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2		Tetradecanoic acid	228	C14H28O2	000544-63-8	83
3		Octadecanoic acid, 2-(2-hydroxyve...	372	C22H44O4	000106-11-6	72
4		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	68
5		n-Decanoic acid	172	C10H20O2	000334-48-5	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG060719\
Data File : BG041122.D
Acq On : 8 Jun 2019 6:39
Operator : HP/JU
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Instrument :
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ClientSampleId :
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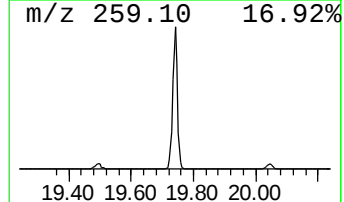
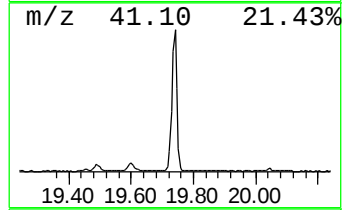
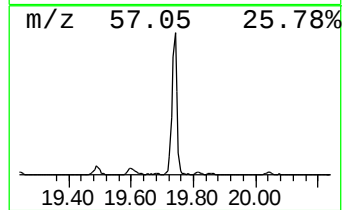
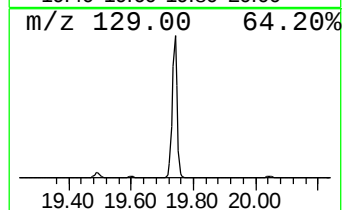
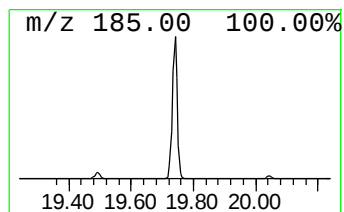
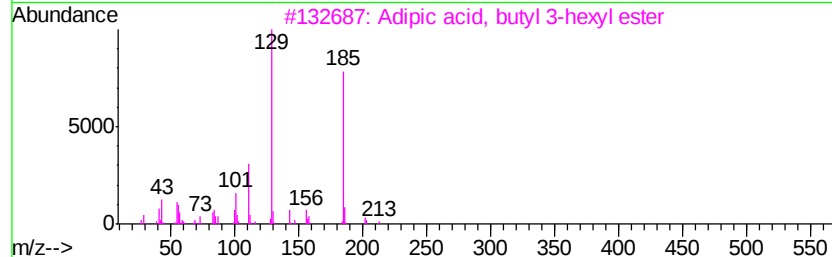
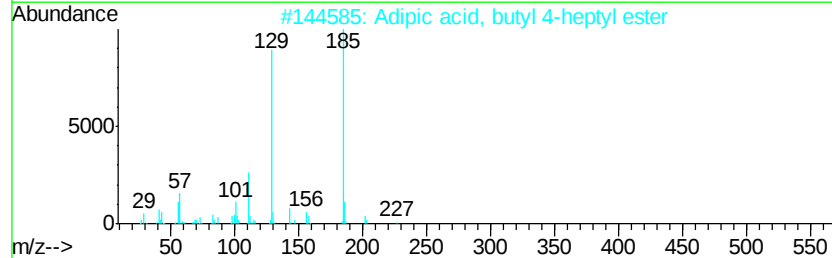
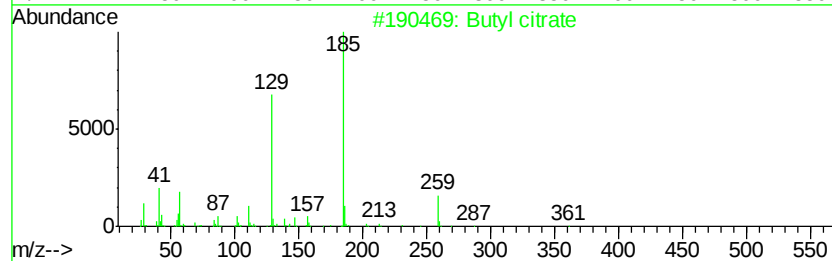
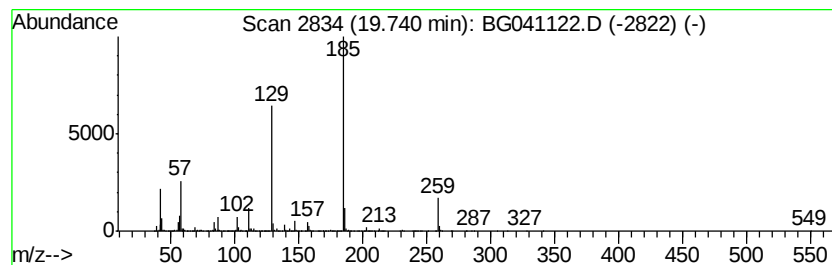
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Butyl citrate Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.74	82.57 ng	3722490	Chrysene-d12	21.76

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butyl citrate	360	C18H32O7	000077-94-1	91
2		Adipic acid, butyl 4-heptyl ester	300	C17H32O4	1000349-73-7	64
3		Adipic acid, butyl 3-hexyl ester	286	C16H30O4	1000353-62-3	56
4		Adipic acid, 4-heptyl isobutyl e...	300	C17H32O4	1000349-73-6	56
5		Hexanedioic acid, bis(2-methylpr...	258	C14H26O4	000141-04-8	56



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG060719\
Data File : BG041122.D
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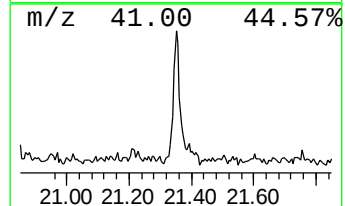
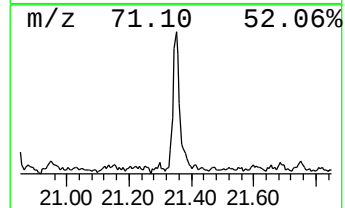
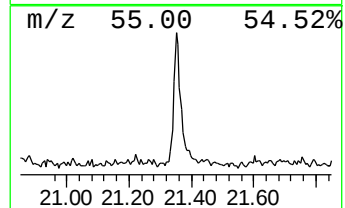
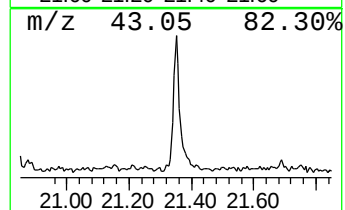
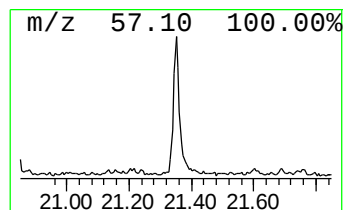
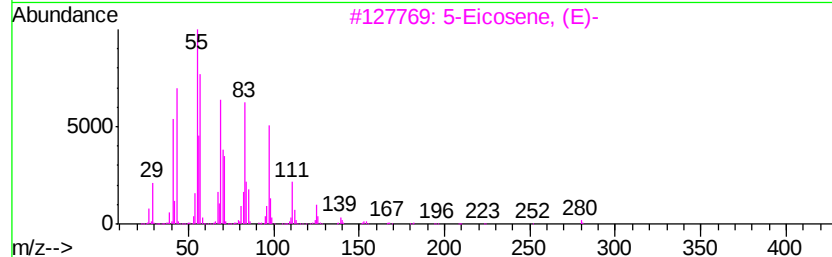
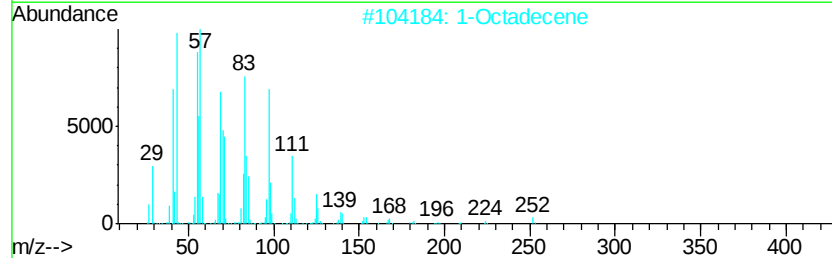
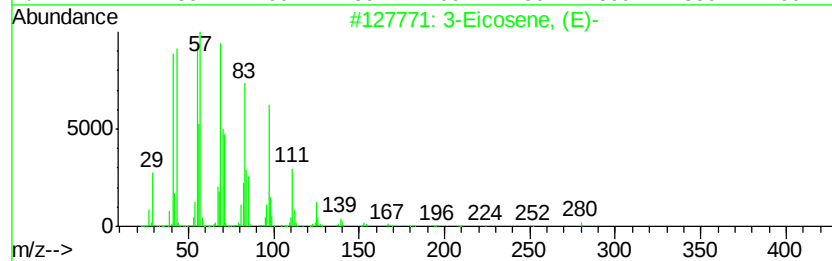
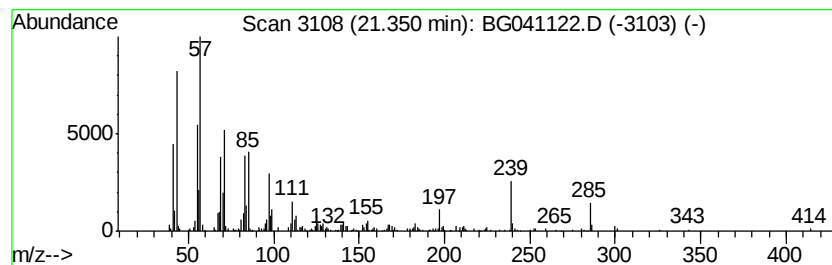
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 3-Eicosene, (E)- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.35	6.91 ng	311369	Chrysene-d12	21.76

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Eicosene, (E)-	280	C20H40	074685-33-9	95
2		1-Octadecene	252	C18H36	000112-88-9	94
3		5-Eicosene, (E)-	280	C20H40	074685-30-6	86
4		1-Heptadecene	238	C17H34	006765-39-5	86
5		9-Eicosene, (E)-	280	C20H40	074685-29-3	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG060719\
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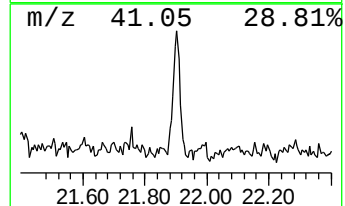
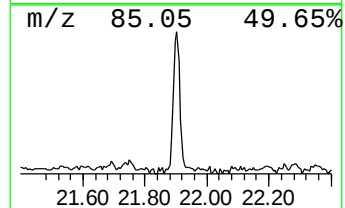
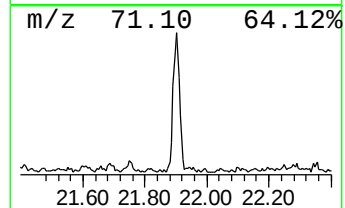
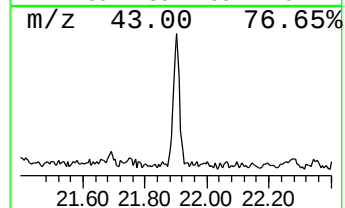
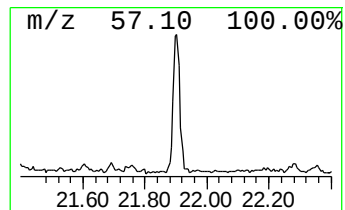
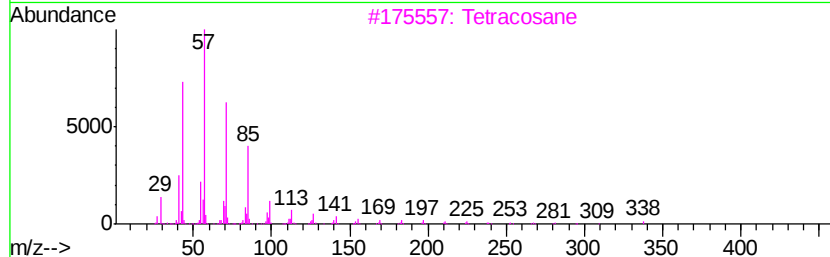
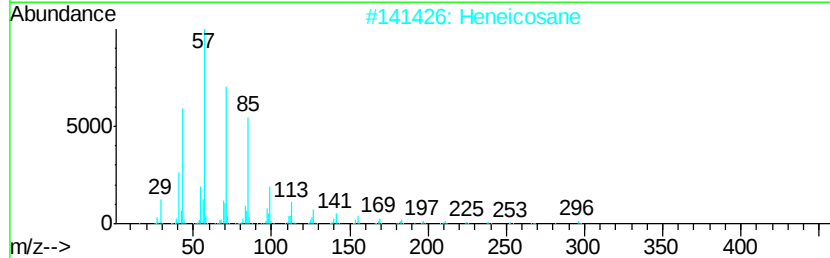
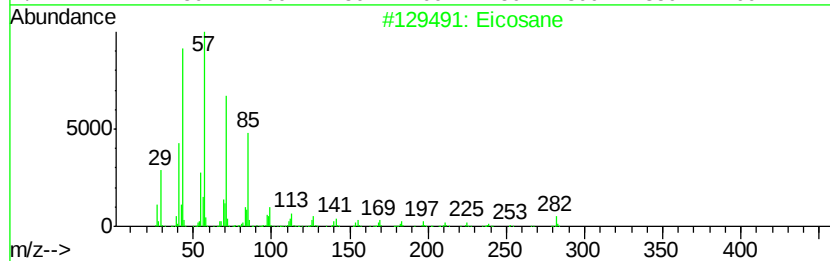
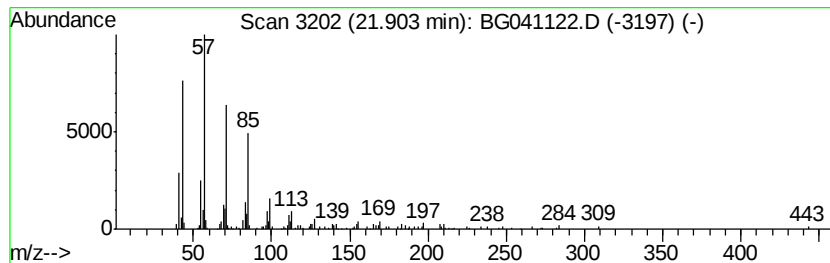
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Eicosane Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.90	2.88 ng	129850	Chrysene-d12	21.76

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	93
2		Heneicosane	296	C21H44	000629-94-7	87
3		Tetracosane	338	C24H50	000646-31-1	87
4		2-methyloctacosane	408	C29H60	1000376-72-8	87
5		Docosane	310	C22H46	000629-97-0	86



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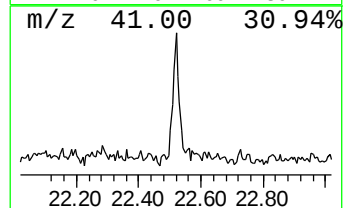
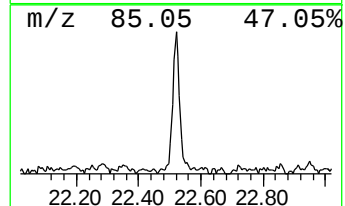
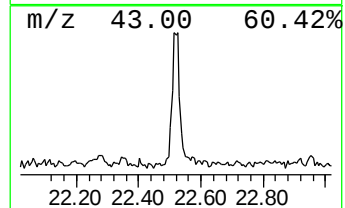
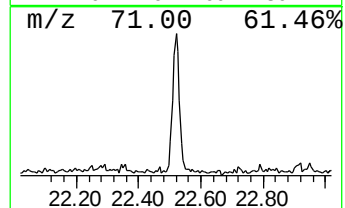
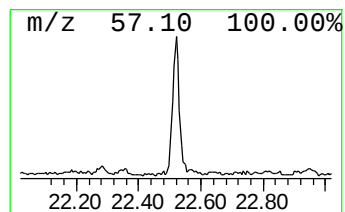
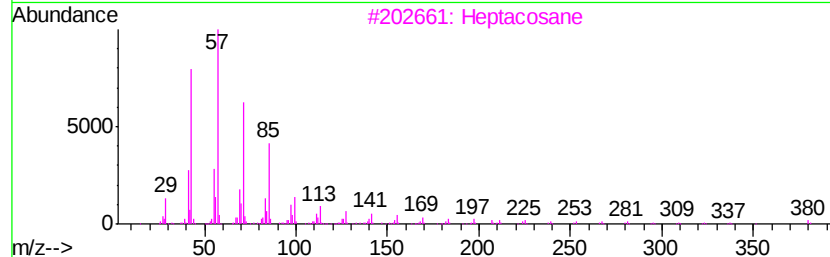
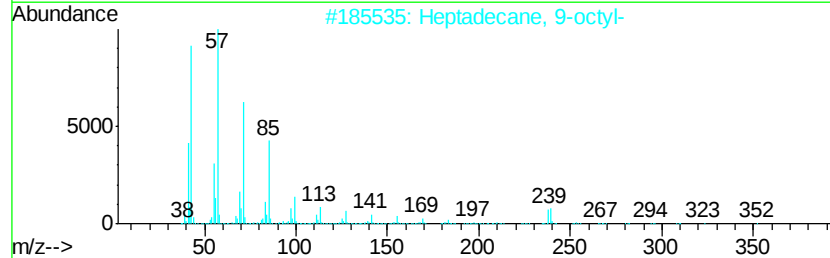
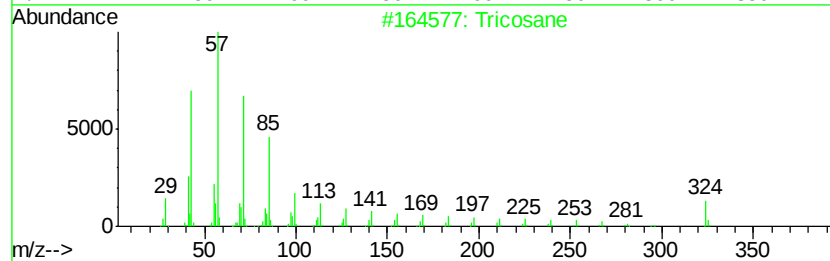
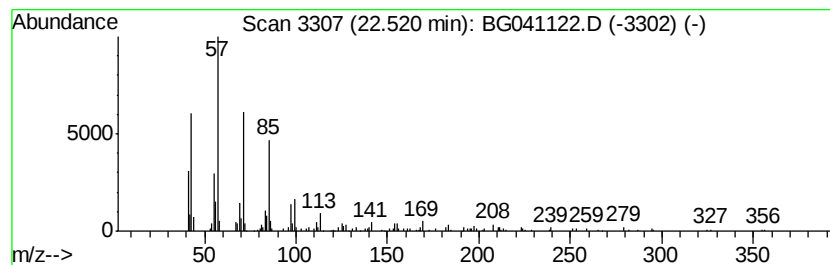
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Tricosane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.52	3.71 ng	167385	Chrysene-d12	21.76

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tricosane	324	C23H48	000638-67-5	91
2			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	87
3			Heptacosane	380	C27H56	000593-49-7	87
4			Heneicosane	296	C21H44	000629-94-7	86
5			2-methyloctacosane	408	C29H60	1000376-72-8	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG060719\
Data File : BG041122.D
Acq On : 8 Jun 2019 6:39
Operator : HP/JU
Sample : K3191-07
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
CULVER-CBTC-TP1-060419

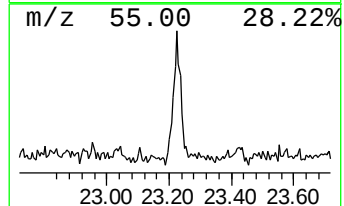
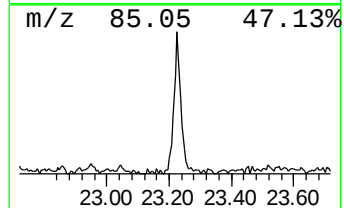
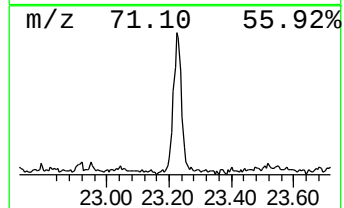
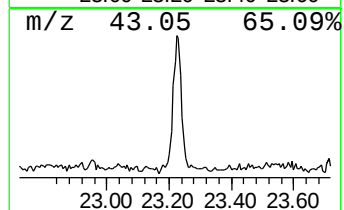
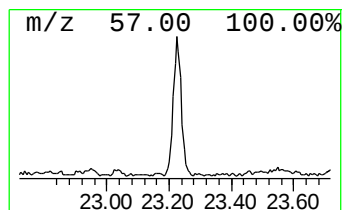
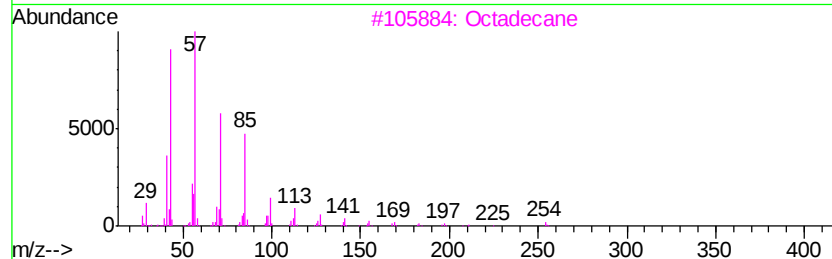
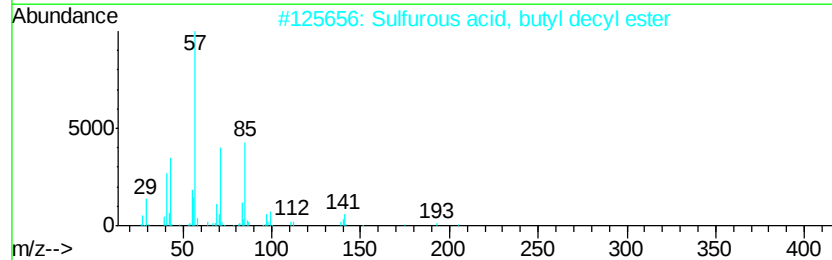
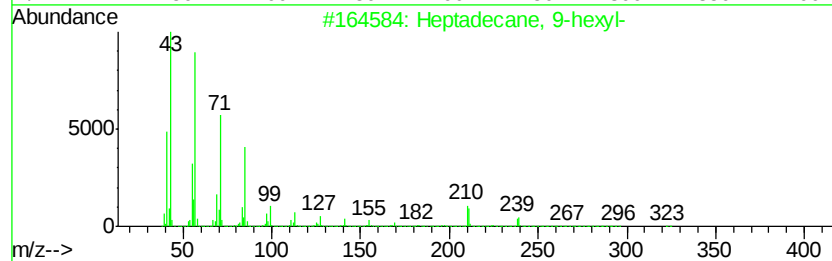
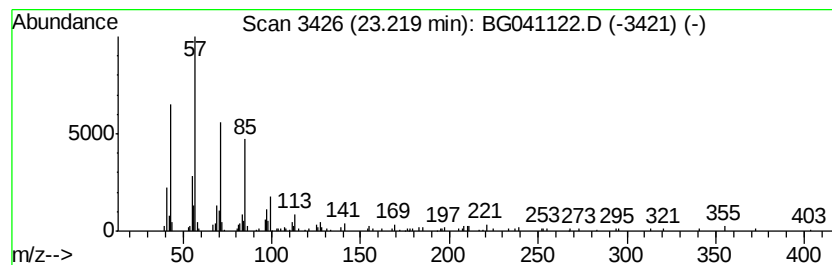
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Heptadecane, 9-hexyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.22	4.36 ng	196515	Chrysene-d12	21.76

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecane, 9-hexyl-	324	C23H48	055124-79-3	86
2			Sulfurous acid, butyl decyl ester	278	C14H30O3S	1000309-17-7	86
3			Octadecane	254	C18H38	000593-45-3	80
4			Tetratetracontane	619	C44H90	007098-22-8	76
5			Octacosane	394	C28H58	000630-02-4	76



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG060719\
Data File : BG041122.D
Acq On : 8 Jun 2019 6:39
Operator : HP/JU
Sample : K3191-07
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
CULVER-CBTC-TP1-060419

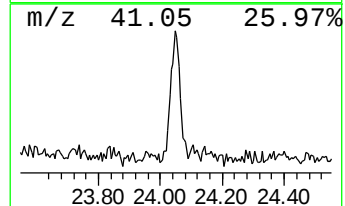
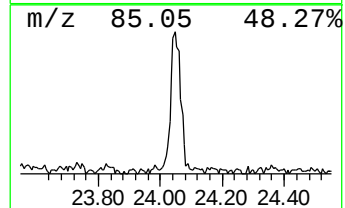
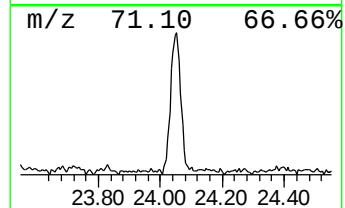
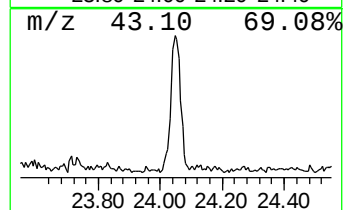
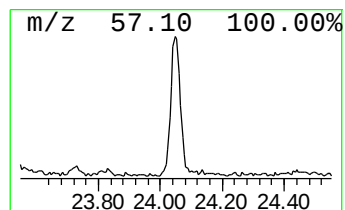
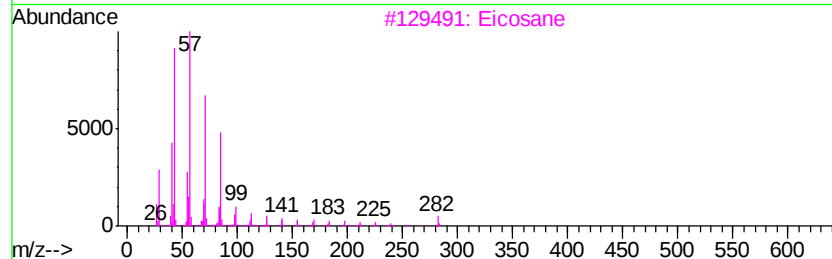
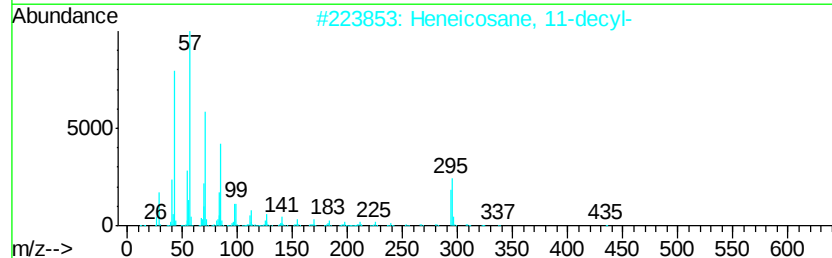
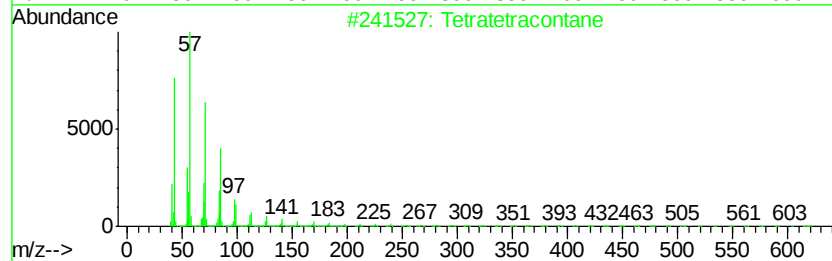
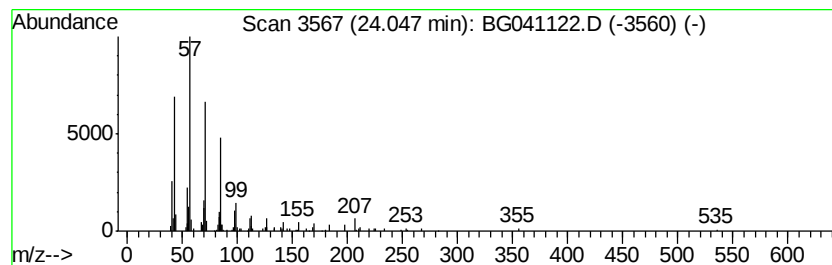
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Tetratetracontane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.05	3.98 ng	213185	Perylene-d12	25.08

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetratetracontane	619	C44H90	007098-22-8	90
2			Heneicosane, 11-decyl-	437	C31H64	055320-06-4	90
3			Eicosane	282	C20H42	000112-95-8	90
4			Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	89
5			Octacosane	394	C28H58	000630-02-4	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG060719\
 Data File : BG041122.D
 Acq On : 8 Jun 2019 6:39
 Operator : HP/JU
 Sample : K3191-07
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 CULVER-CBTC-TP1-060419

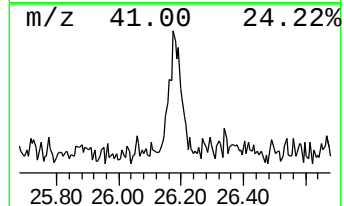
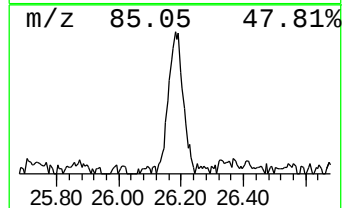
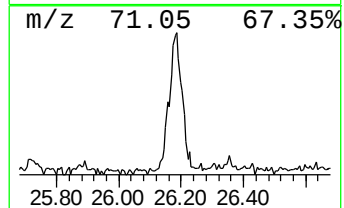
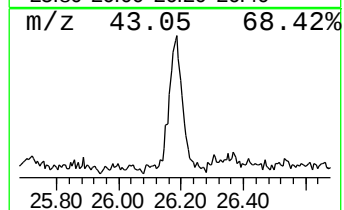
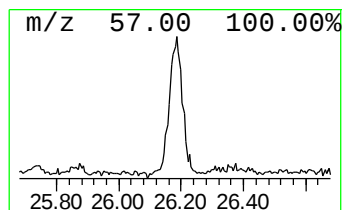
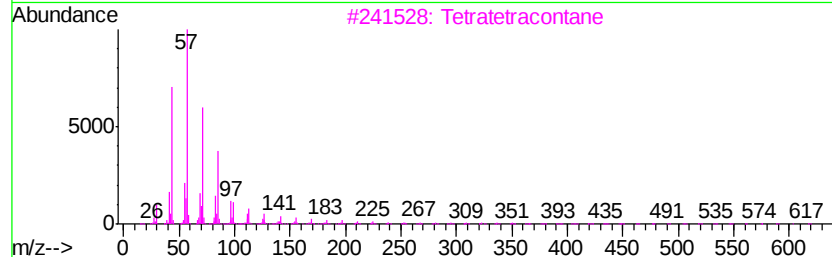
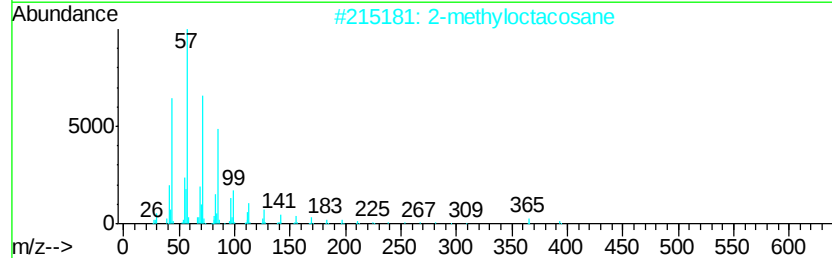
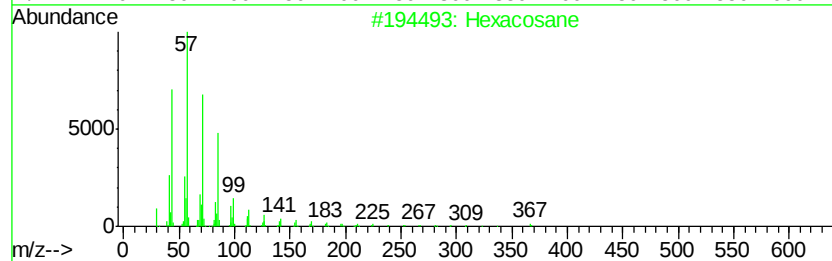
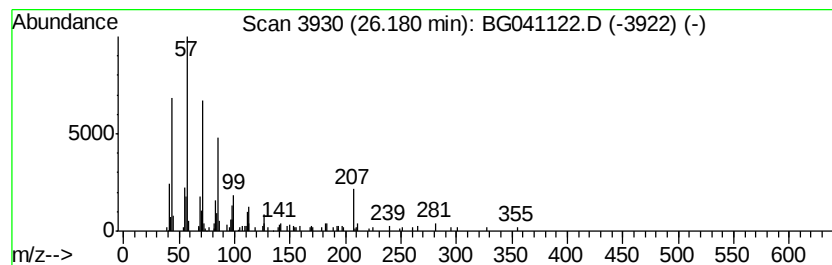
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 16 Hexacosane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.18	3.00 ng	160798	Perylene-d12	25.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexacosane	366	C26H54	000630-01-3	87
2		2-methyloctacosane	408	C29H60	1000376-72-8	87
3		Tetratetracontane	619	C44H90	007098-22-8	87
4		Heneicosane, 11-decyl-	437	C31H64	055320-06-4	83
5		Heptacosane	380	C27H56	000593-49-7	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG060719\
 Data File : BG041122.D
 Acq On : 8 Jun 2019 6:39
 Operator : HP/JU
 Sample : K3191-07
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

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

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	3.18	16.7	ng	213175	1	8.11	255028	20.0
unknown7.63	7.63	79.6	ng	1014720	1	8.11	255028	20.0
Hexanoic acid, 2-...	9.38	13.0	ng	165328	1	8.11	255028	20.0
Nonanoic acid	11.70	7.0	ng	129879	2	10.93	372557	20.0
Pentafluoropropio...	12.70	2.1	ng	38978	2	10.93	372557	20.0
n-Hexadecanoic acid	18.33	10.1	ng	389470	4	17.48	775397	20.0
Adipic acid, buty...	19.49	5.7	ng	222319	4	17.48	775397	20.0
Octadecanoic acid	19.60	6.8	ng	261766	4	17.48	775397	20.0
Butyl citrate	19.74	82.6	ng	3722490	5	21.76	901666	20.0
3-Eicosene, (E)-	21.35	6.9	ng	311369	5	21.76	901666	20.0
Eicosane	21.90	2.9	ng	129850	5	21.76	901666	20.0
Tricosane	22.52	3.7	ng	167385	5	21.76	901666	20.0
Heptadecane, 9-he...	23.22	4.4	ng	196515	5	21.76	901666	20.0
Tetratetracontane	24.05	4.0	ng	213185	6	25.08	1071800	20.0
Hexacosane	26.18	3.0	ng	160798	6	25.08	1071800	20.0