

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG040621\
 Data File : BG048604.D
 Acq On : 6 Apr 2021 15:05
 Operator : CG/JU
 Sample : M1933-06
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BENNETT_BH_01_2-2.5

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

Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG033021.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BG048604.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.638	385	391	398	rBV	362982	564595	11.00%	3.079%
2	7.248	658	665	674	rBV	367843	628724	12.25%	3.429%
3	8.094	802	809	817	rBV2	93392	159532	3.11%	0.870%
4	9.263	1000	1008	1016	rBV	234934	407237	7.94%	2.221%
5	10.920	1283	1290	1298	rVB	112426	211842	4.13%	1.155%
6	13.370	1699	1707	1713	rBV	500924	792398	15.44%	4.321%
7	14.751	1935	1942	1948	rBV	156089	276614	5.39%	1.508%
8	16.238	2188	2195	2209	rVB	507422	823457	16.05%	4.491%
9	17.501	2404	2410	2415	rBV	219838	341734	6.66%	1.864%
10	18.376	2551	2559	2569	rVB6	64748	139435	2.72%	0.760%
11	18.605	2583	2598	2612	rBV2	429982	1780178	34.70%	9.708%
12	18.764	2613	2625	2635	rVV	631338	2101713	40.96%	11.461%
13	18.911	2635	2650	2661	rVB	1836878	5130725	100.00%	27.979%
14	19.058	2668	2675	2687	rVV2	131095	348466	6.79%	1.900%
15	19.275	2707	2712	2717	rVB3	30543	63070	1.23%	0.344%
16	19.645	2768	2775	2788	rVB3	242829	655867	12.78%	3.577%
17	19.845	2804	2809	2817	rVB5	64766	150540	2.93%	0.821%
18	20.110	2849	2854	2863	rVB	1030177	1334752	26.01%	7.279%
19	20.691	2946	2953	2960	rBV2	119271	226038	4.41%	1.233%
20	21.426	3074	3078	3084	rBV7	67878	148944	2.90%	0.812%
21	21.490	3084	3089	3096	rVB2	125638	236124	4.60%	1.288%
22	21.684	3117	3122	3133	rVB2	103528	230893	4.50%	1.259%
23	21.802	3135	3142	3149	rVB2	239686	433364	8.45%	2.363%
24	22.407	3237	3245	3255	rVB4	99908	272359	5.31%	1.485%
25	23.476	3419	3427	3434	rVB2	96097	233204	4.55%	1.272%
26	24.716	3629	3638	3648	rVB5	85912	243763	4.75%	1.329%
27	25.151	3703	3712	3721	rBV2	141843	402012	7.84%	2.192%

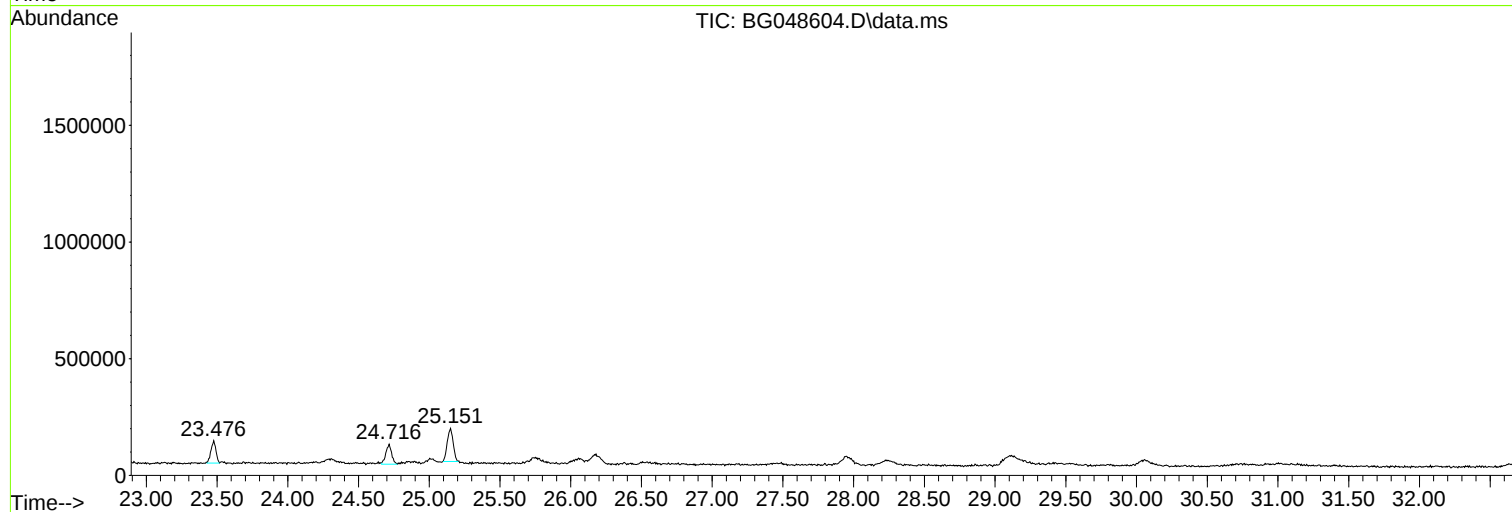
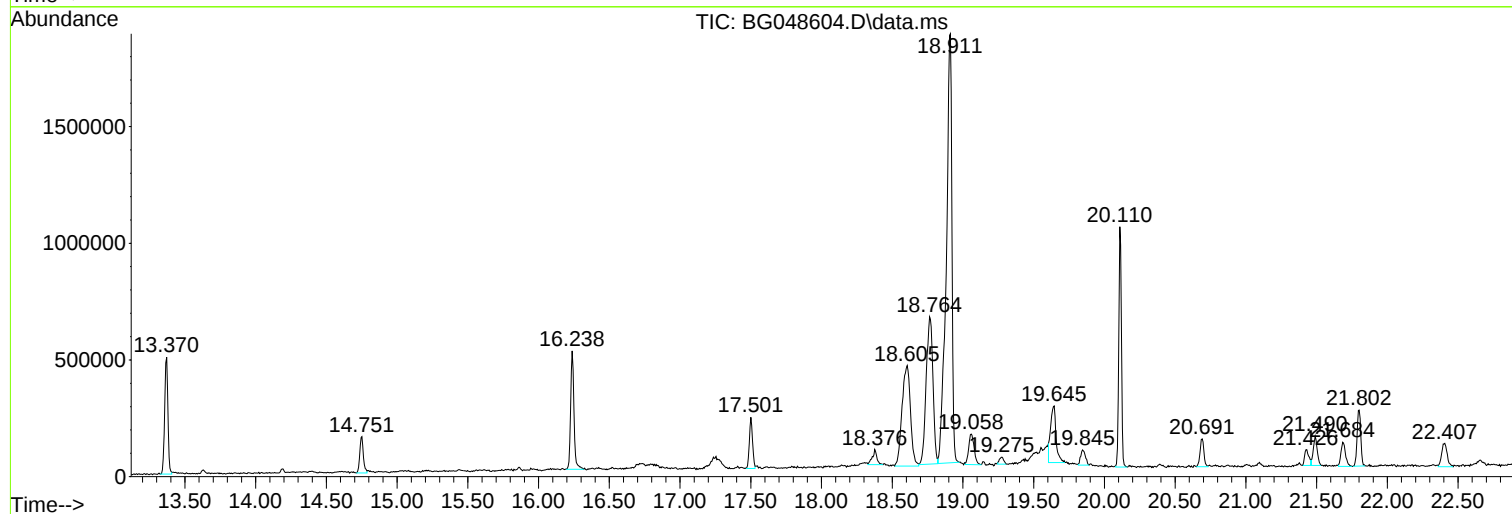
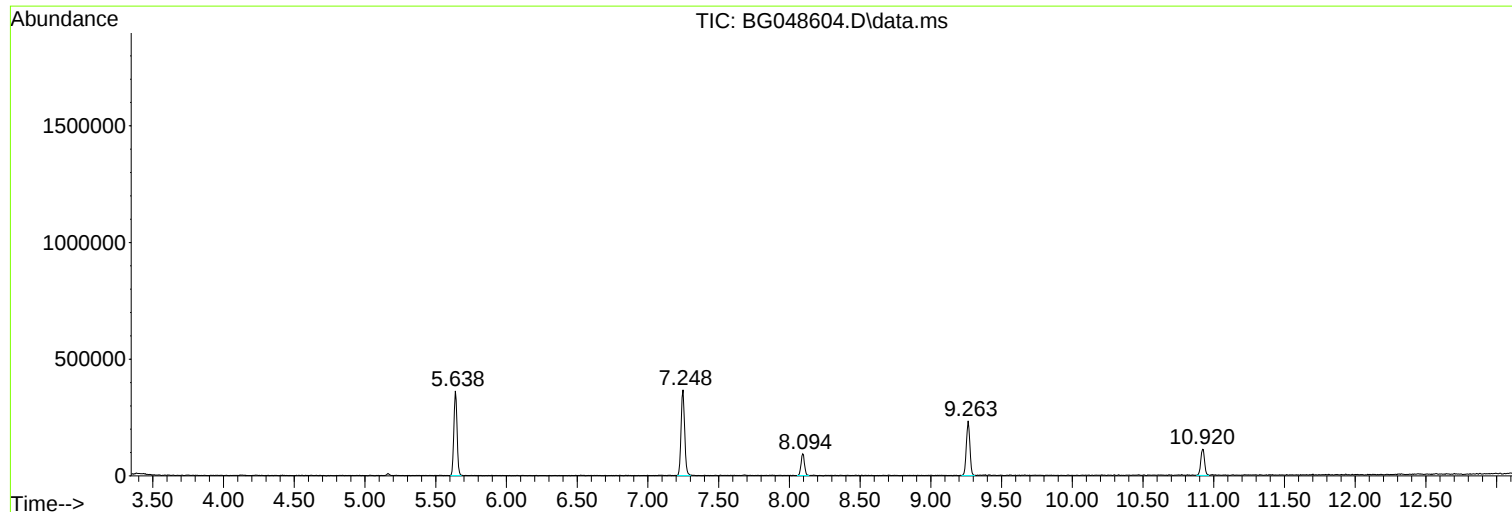
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG033021.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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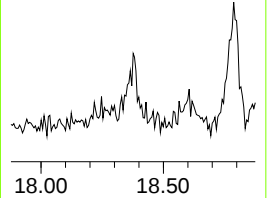
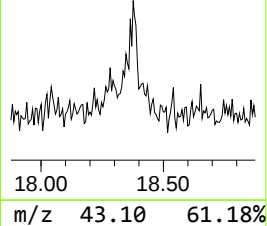
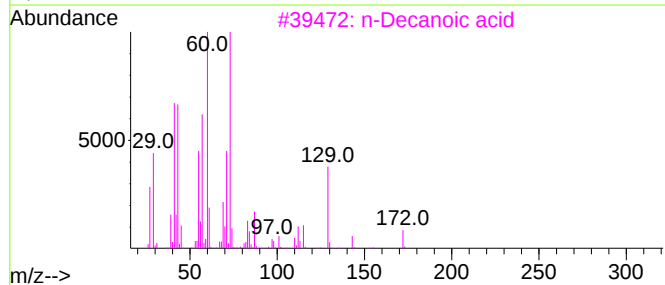
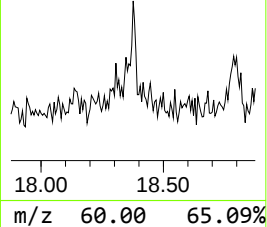
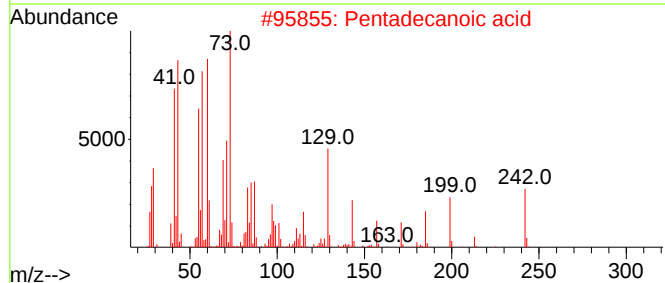
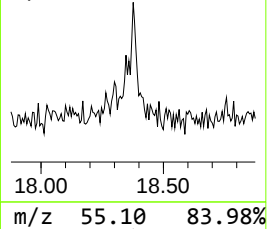
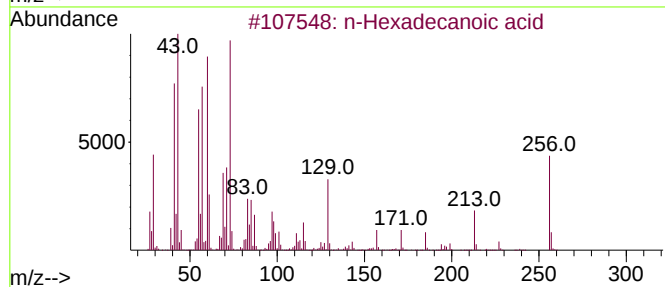
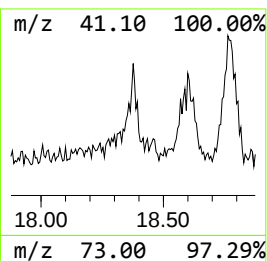
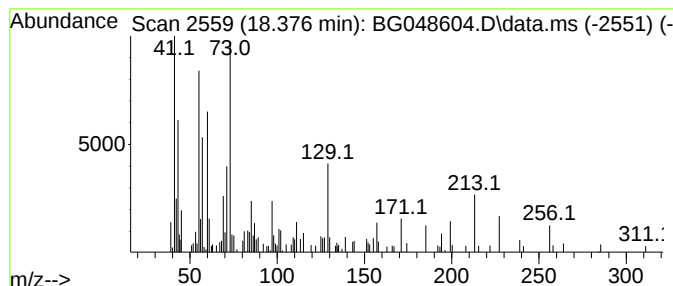
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TIC Library : C:\Database\NIST11.L
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Peak Number 1 n-Hexadecanoic acid Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.376	8.16 ng	139435	Phenanthrene-d10	17.501

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	91
2			Pentadecanoic acid	242	C15H30O2	001002-84-2	53
3			n-Decanoic acid	172	C10H20O2	000334-48-5	43
4			Dodecanoic acid	200	C12H24O2	000143-07-7	38
5			Ethanone, 1-(4,5-dihydro-2-thiaz...	129	C5H7NO5	029926-41-8	25



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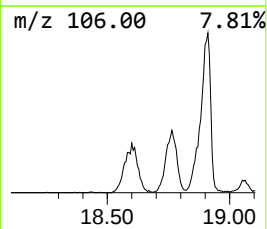
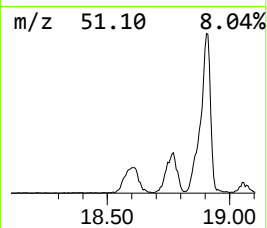
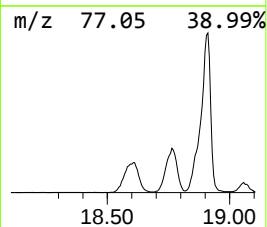
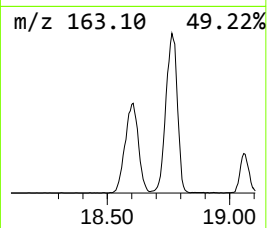
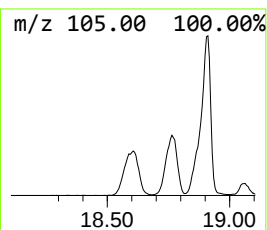
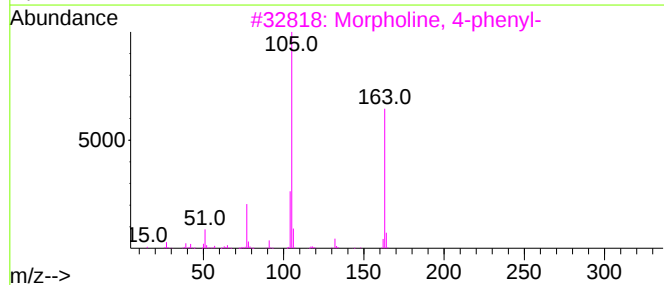
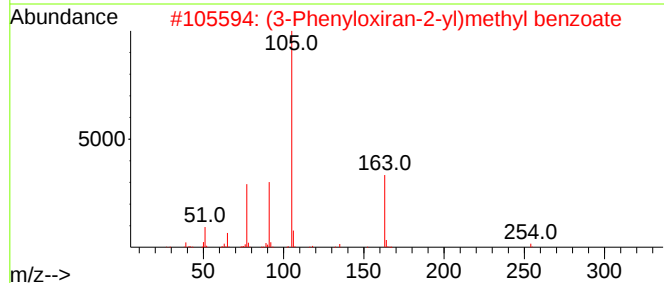
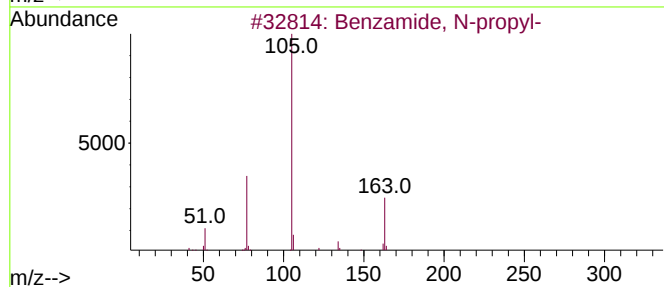
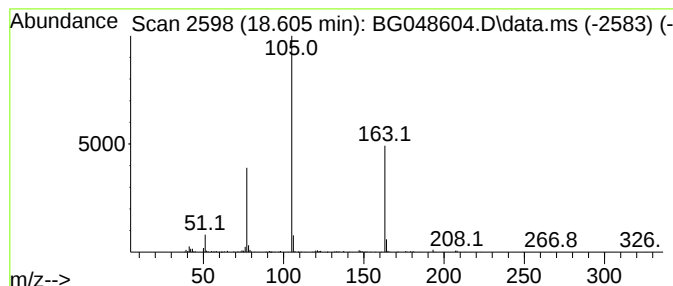
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Peak Number 2 Benzamide, N-propyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.605	104.19 ng	1780180	Phenanthrene-d10	17.501

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzamide, N-propyl-	163	C10H13NO	010546-70-0	78
2			(3-Phenyloxiran-2-yl)methyl benz...	254	C16H14O3	1000366-96-1	64
3			Morpholine, 4-phenyl-	163	C10H13NO	000092-53-5	50
4			Vinyl benzoate	148	C9H8O2	000769-78-8	43
5			Benzoic acid, phenyl ester	198	C13H10O2	000093-99-2	43



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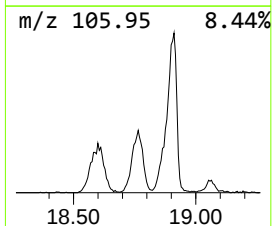
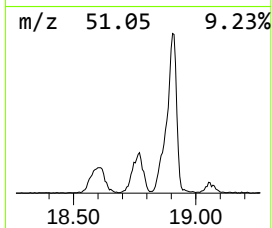
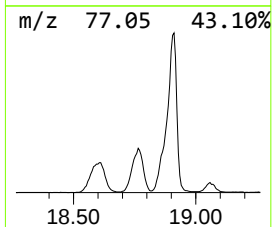
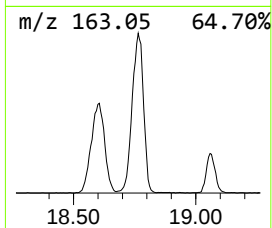
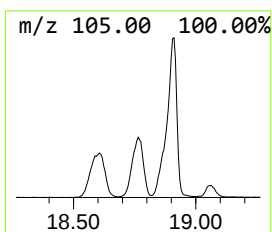
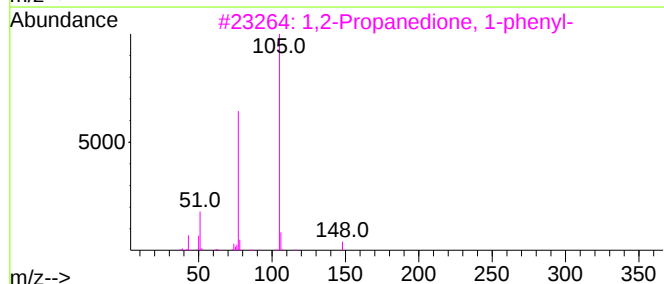
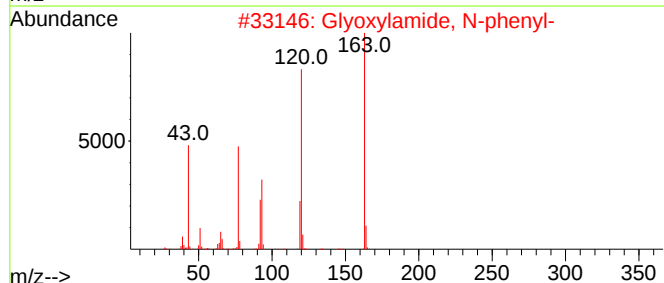
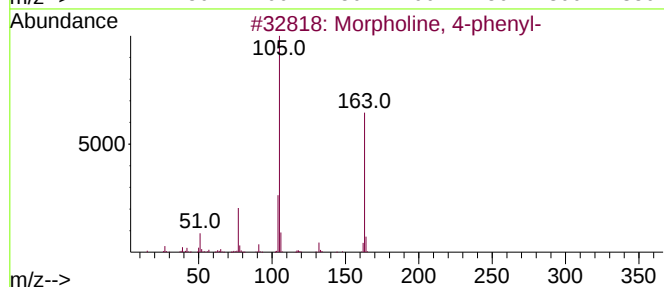
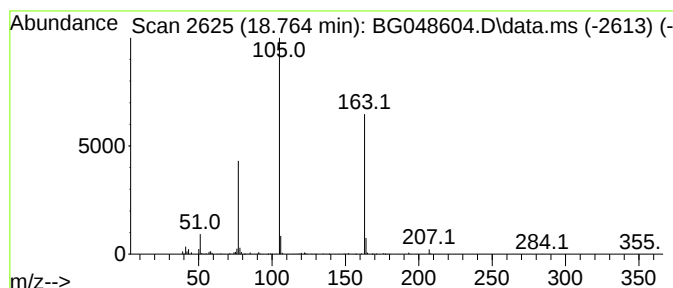
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TIC Library : C:\Database\NIST11.L
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Peak Number 3 Morpholine, 4-phenyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.764	123.00 ng	2101710	Phenanthrene-d10	17.501

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Morpholine, 4-phenyl-	163	C10H13NO	000092-53-5	72
2			Glyoxylamide, N-phenyl-	163	C9H9NO2	032331-52-5	53
3			1,2-Propanedione, 1-phenyl-	148	C9H8O2	000579-07-7	47
4			Vinyl benzoate	148	C9H8O2	000769-78-8	47
5			Benzeneacetic acid, .alpha.-oxo-...	178	C10H10O3	001603-79-8	47



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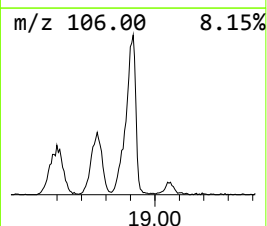
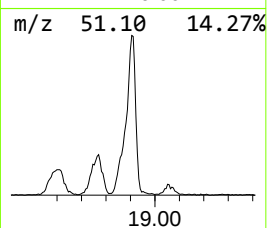
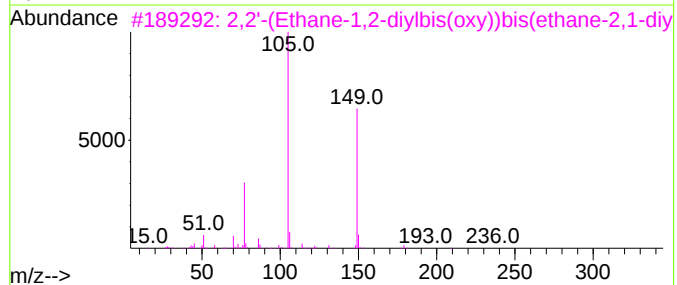
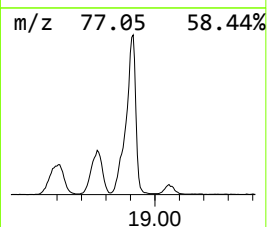
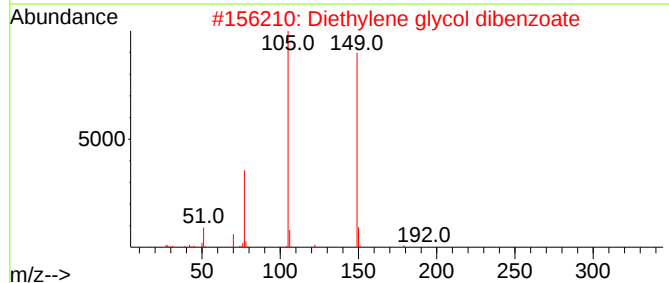
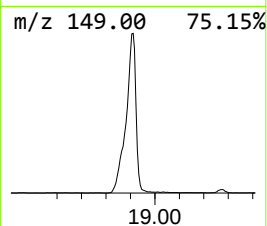
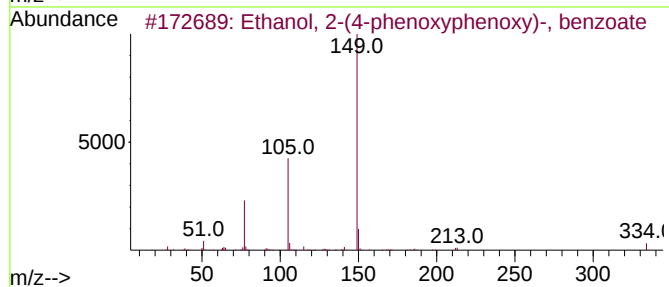
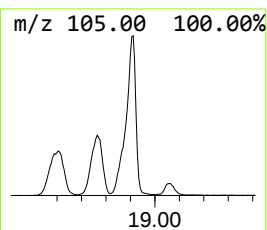
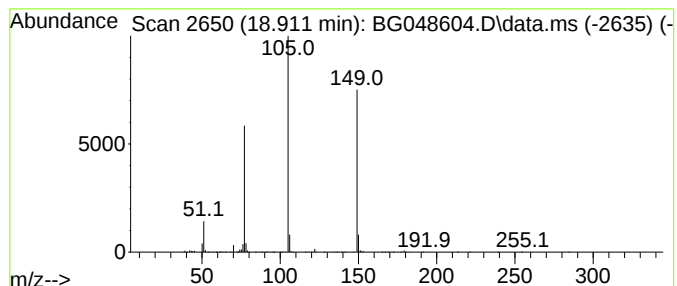
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Peak Number 4 Ethanol, 2-(4-phenoxypheno... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.911	300.28 ng	5130730	Phenanthrene-d10	17.501

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(4-phenoxyphenoxy)-, ...	334	C21H18O4	055191-59-8	83
2			Diethylene glycol dibenzoate	314	C18H18O5	000120-55-8	72
3			2,2'-(Ethane-1,2-diylbis(oxy))bi...	358	C20H22O6	1000366-99-4	72
4			1,3-Dioxolane, 2-(methoxymethyl)...	194	C11H14O3	1000156-71-2	72
5			2-(2-Carboxyvinyl)pyridine, trans	149	C8H7NO2	007340-22-9	50



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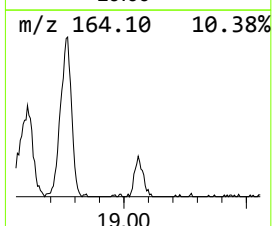
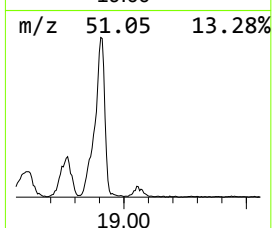
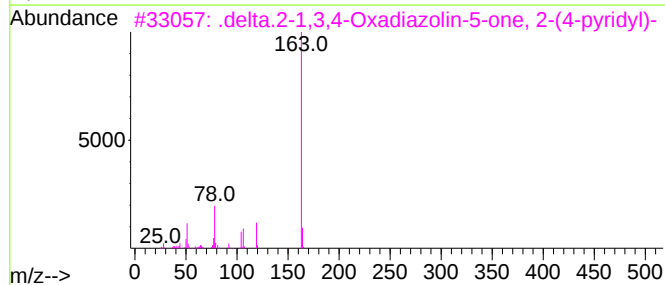
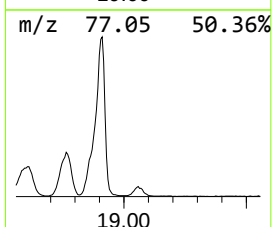
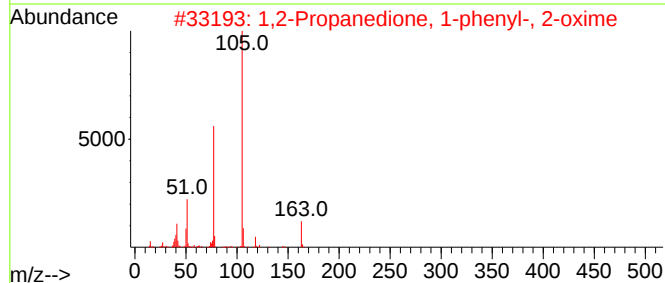
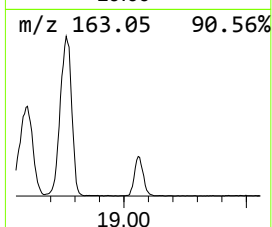
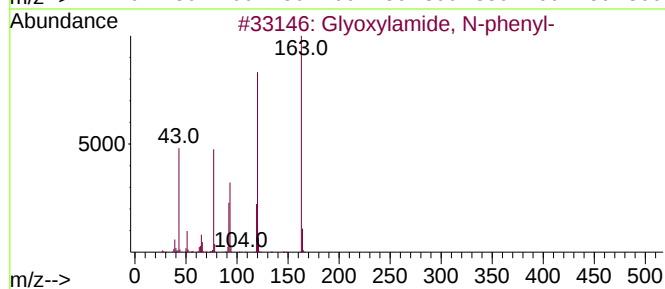
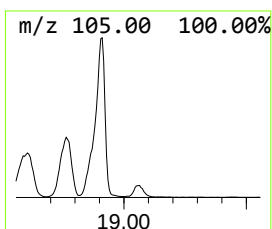
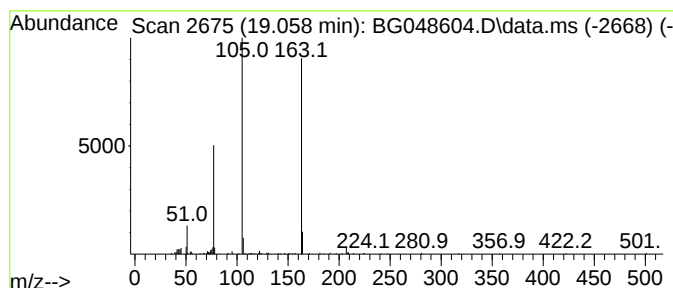
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Peak Number 5 Glyoxylamide, N-phenyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.058	20.39 ng	348466	Phenanthrene-d10	17.501	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Glyoxylamide, N-phenyl-	163	C9H9NO2	032331-52-5	64
2	1,2-Propanedione, 1-phenyl-, 2-o...	163	C9H9NO2	000119-51-7	59
3	.delta.2-1,3,4-Oxadiazolin-5-one...	163	C7H5N3O2	002845-82-1	47
4	2,2-Bis[4-(benzoyloxy)phenyl]pro...	436	C29H24O4	002297-14-5	35
5	Benzenepropanenitrile, .beta.-oxo-	145	C9H7NO	000614-16-4	27



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BENNETT_BH_01_2-2.5

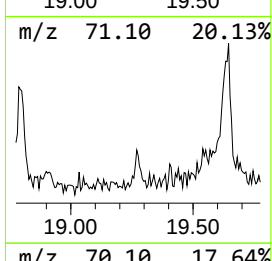
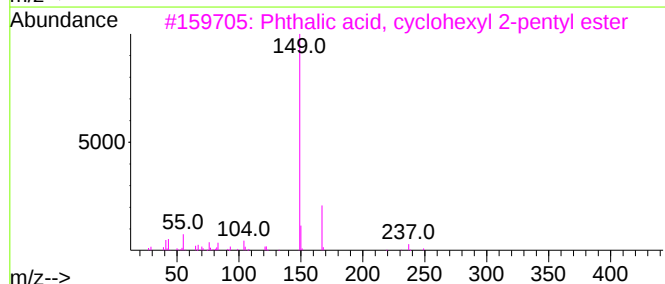
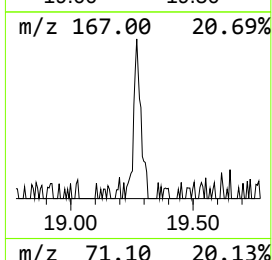
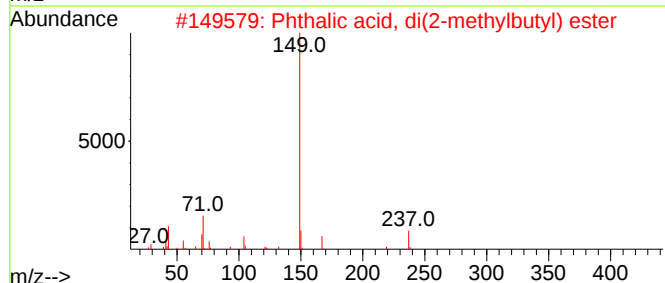
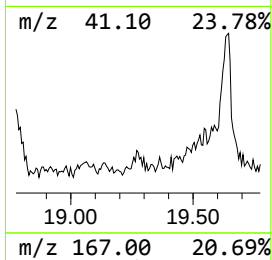
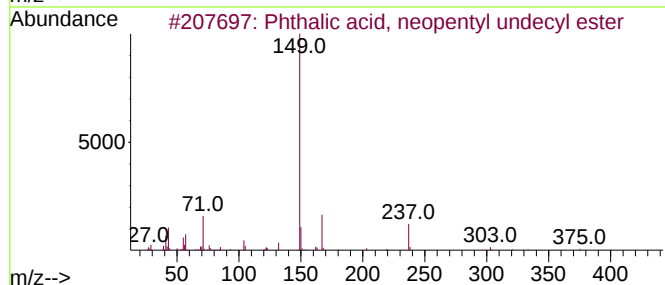
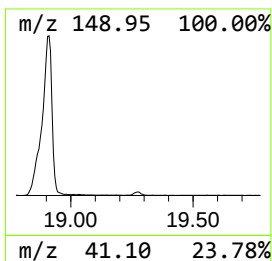
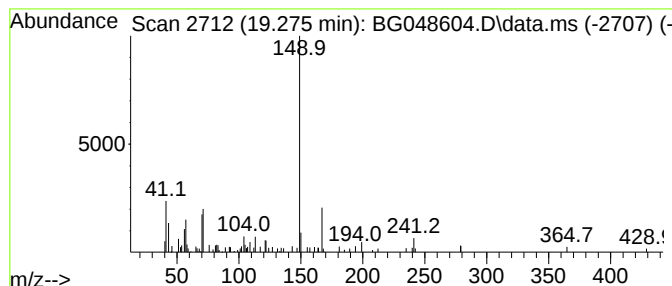
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Phthalic acid, neopentyl un... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.275	3.69 ng	63070	Phenanthrene-d10	17.501

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phthalic acid, neopentyl undecyl...	390	C24H38O4	1000309-07-4	64
2			Phthalic acid, di(2-methylbutyl)...	306	C18H26O4	1000322-82-5	59
3			Phthalic acid, cyclohexyl 2-pent...	318	C19H26O4	1000315-55-3	53
4			2-(Nonyloxycarbonyl)benzoic acid	292	C17H24O4	1000373-89-9	53
5			Phthalic acid, 6-methylhept-2-yl...	530	C34H58O4	1000377-97-2	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG040621\
Data File : BG048604.D
Acq On : 6 Apr 2021 15:05
Operator : CG/JU
Sample : M1933-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BENNETT_BH_01_2-2.5

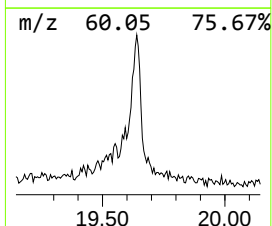
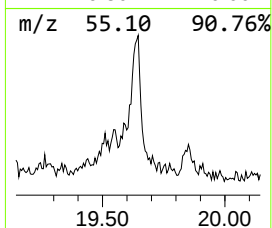
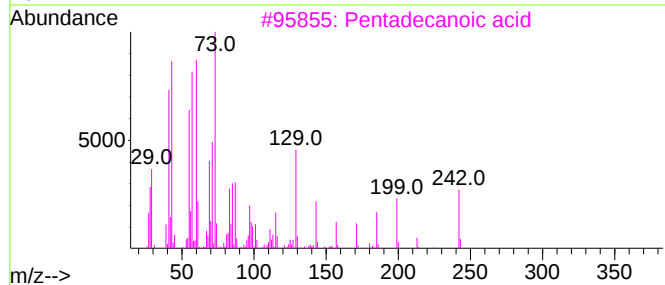
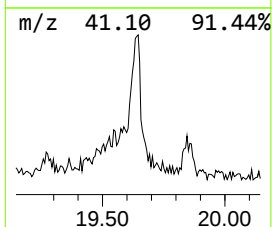
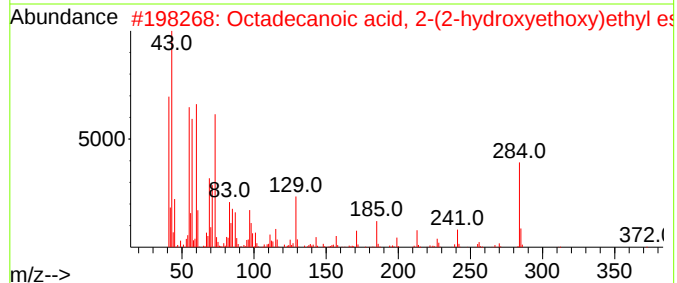
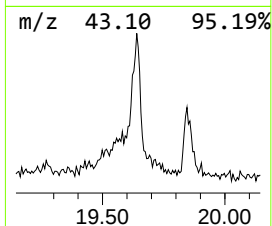
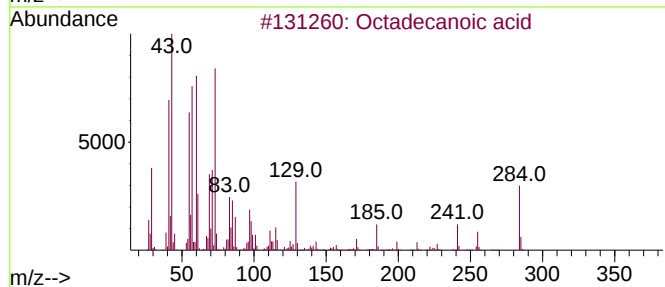
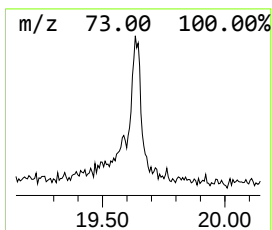
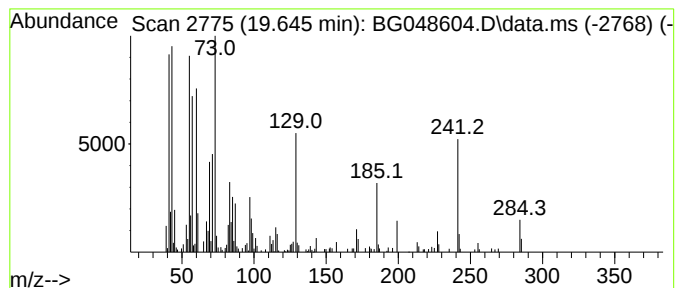
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Octadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.645	38.38 ng	655867	Phenanthrene-d10	17.501

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	91
2			Octadecanoic acid, 2-(2-hydroxye...	372	C22H44O4	000106-11-6	87
3			Pentadecanoic acid	242	C15H30O2	001002-84-2	78
4			Tridecanoic acid	214	C13H26O2	000638-53-9	70
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	68



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG040621\
Data File : BG048604.D
Acq On : 6 Apr 2021 15:05
Operator : CG/JU
Sample : M1933-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

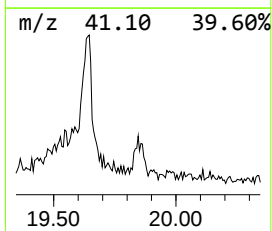
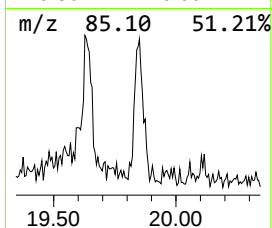
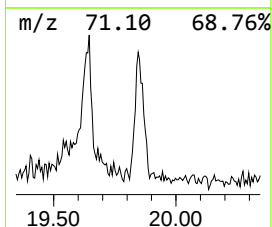
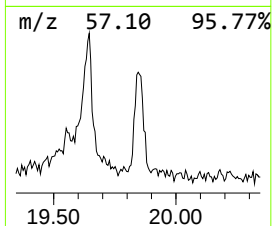
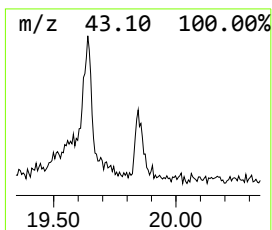
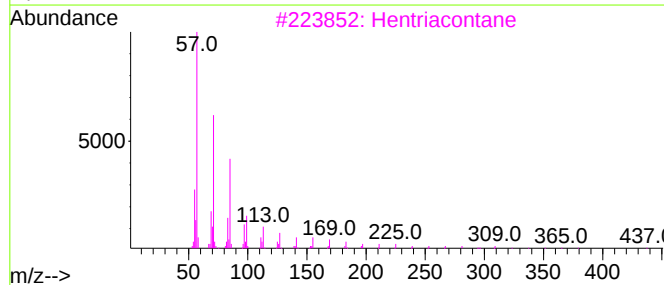
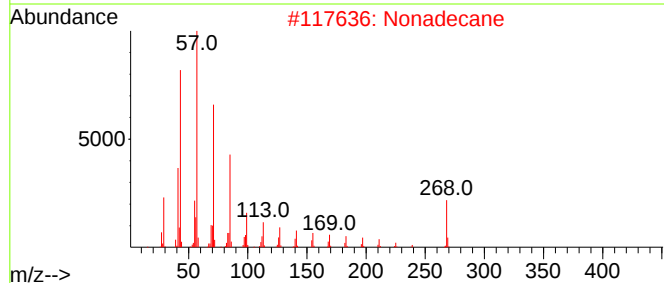
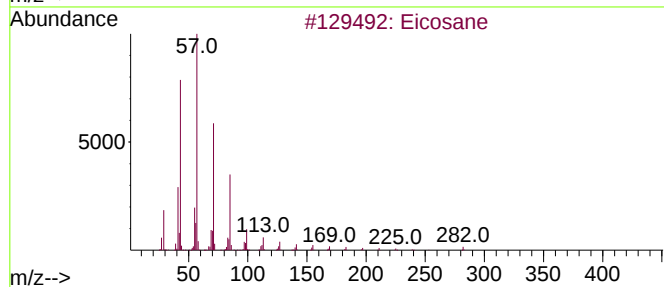
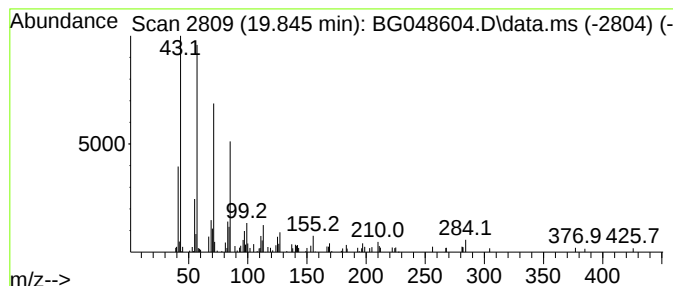
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG033021.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Eicosane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.845	6.95 ng	150540	Chrysene-d12	21.802	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane	282	C20H42	000112-95-8	91
2	Nonadecane	268	C19H40	000629-92-5	91
3	Hentriacontane	437	C31H64	000630-04-6	87
4	Octadecane, 2-methyl-	268	C19H40	001560-88-9	87
5	Docosane	310	C22H46	000629-97-0	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG040621\
Data File : BG048604.D
Acq On : 6 Apr 2021 15:05
Operator : CG/JU
Sample : M1933-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

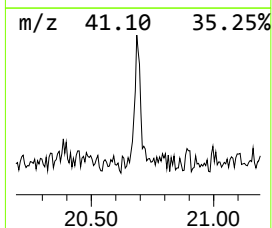
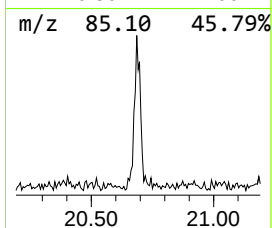
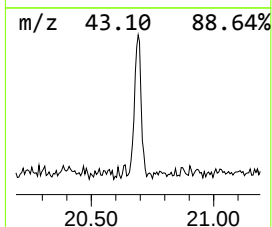
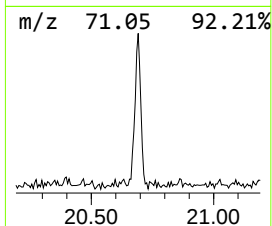
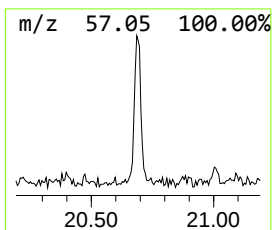
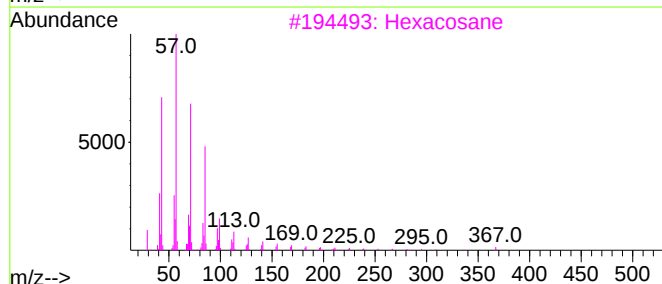
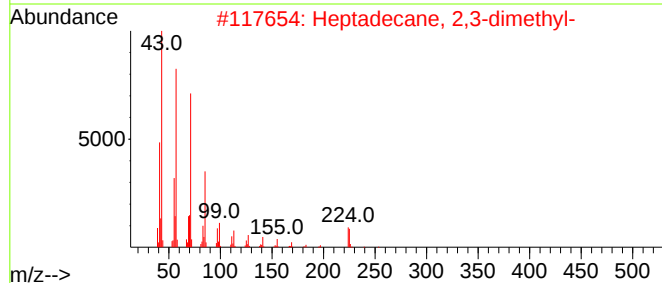
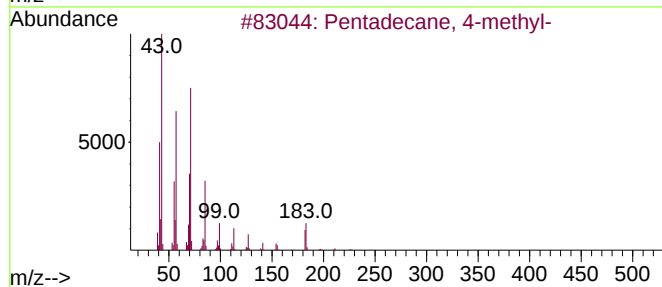
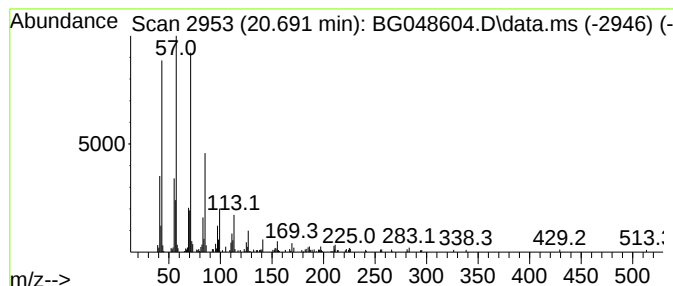
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Pentadecane, 4-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD		R.T.	
20.691	10.43 ng	226038	Chrysene-d12		21.802	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Pentadecane, 4-methyl-		226	C16H34	002801-87-8	91
2	Heptadecane, 2,3-dimethyl-		268	C19H40	061868-03-9	91
3	Hexacosane		366	C26H54	000630-01-3	90
4	2-methyloctacosane		408	C29H60	1000376-72-8	87
5	Tetracosane		338	C24H50	000646-31-1	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG040621\
Data File : BG048604.D
Acq On : 6 Apr 2021 15:05
Operator : CG/JU
Sample : M1933-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

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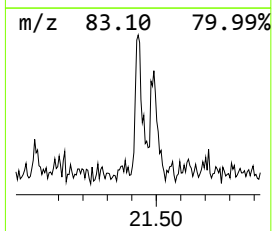
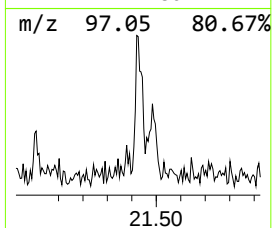
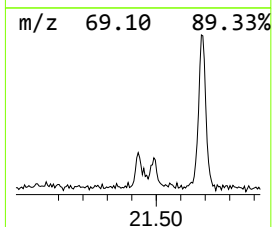
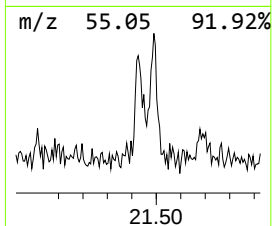
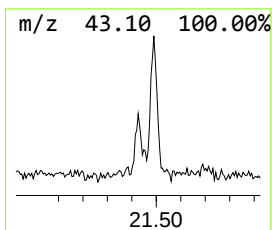
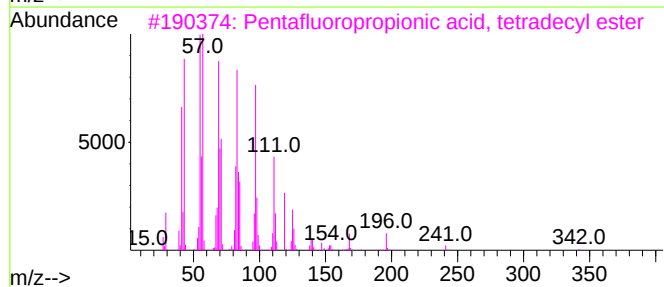
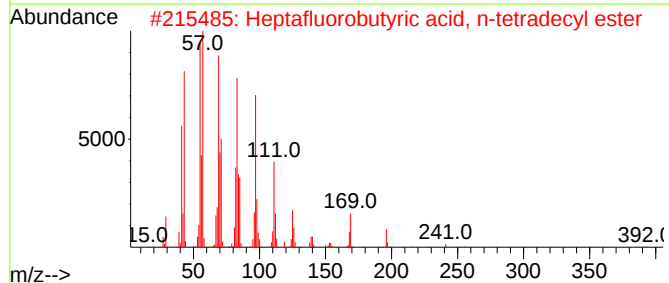
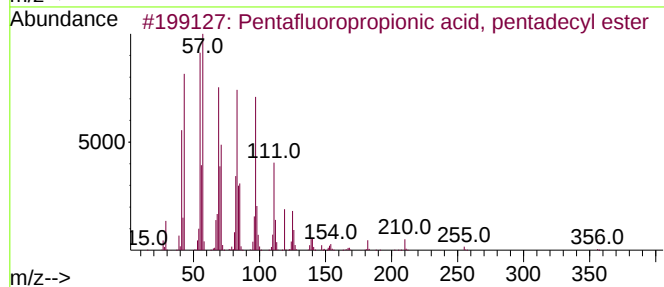
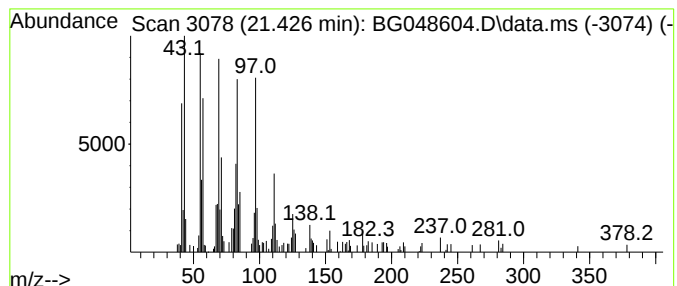
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Pentafluoropropionic acid, ... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.426	6.87 ng	148944	Chrysene-d12	21.802

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	87
2			Heptafluorobutyric acid, n-tetra...	410	C18H29F7O2	007365-36-8	87
3			Pentafluoropropionic acid, tetra...	360	C17H29F5O2	006222-06-6	86
4			Trifluoroacetic acid, pentadecyl...	324	C17H31F3O2	959010-23-2	86
5			Bromoacetic acid, hexadecyl ester	362	C18H35BrO2	005454-48-8	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG040621\
Data File : BG048604.D
Acq On : 6 Apr 2021 15:05
Operator : CG/JU
Sample : M1933-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
BENNETT_BH_01_2-2.5

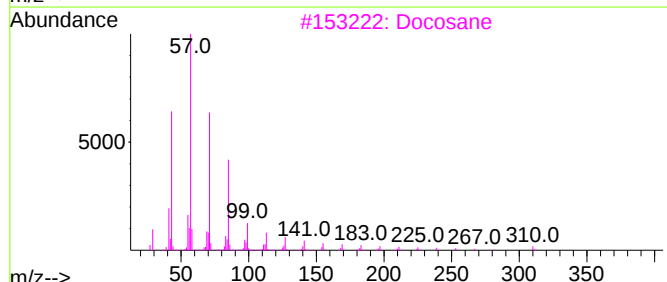
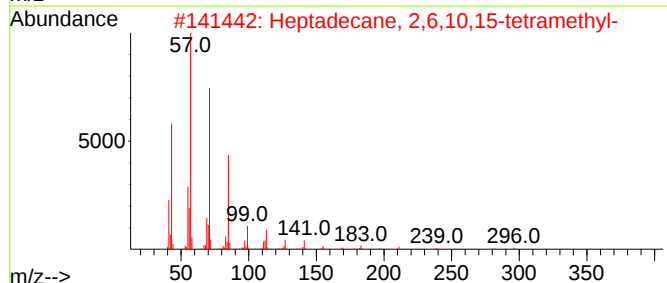
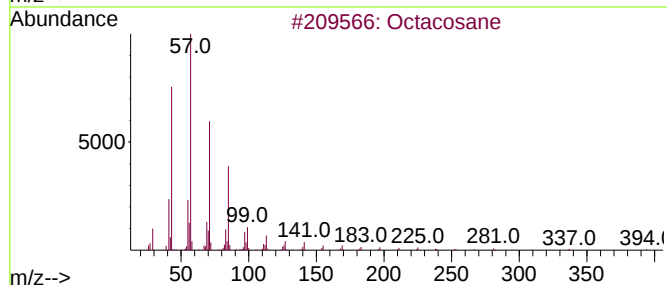
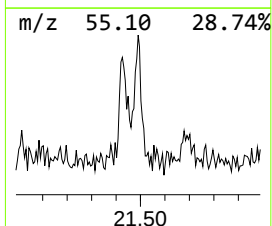
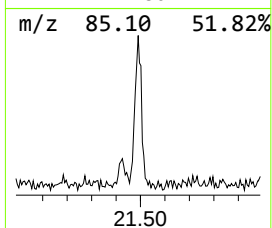
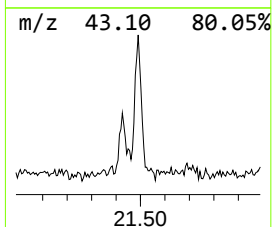
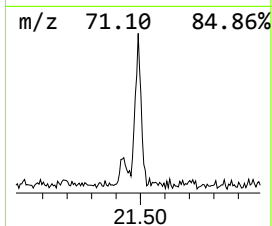
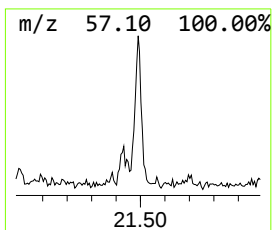
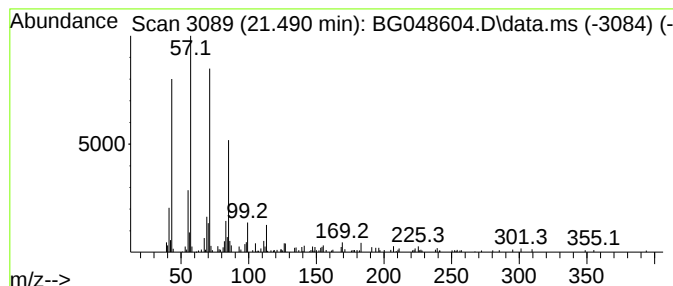
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Octacosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.490	10.90 ng	236124	Chrysene-d12	21.802

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octacosane	394	C28H58	000630-02-4	90
2			Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	78
3			Docosane	310	C22H46	000629-97-0	78
4			Decane, 2,4,6-trimethyl-	184	C13H28	062108-27-4	72
5			Eicosane	282	C20H42	000112-95-8	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG040621\
Data File : BG048604.D
Acq On : 6 Apr 2021 15:05
Operator : CG/JU
Sample : M1933-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

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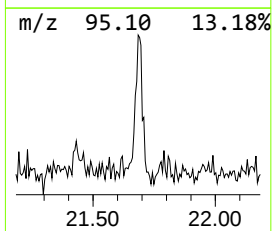
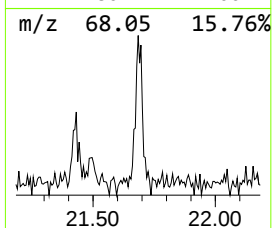
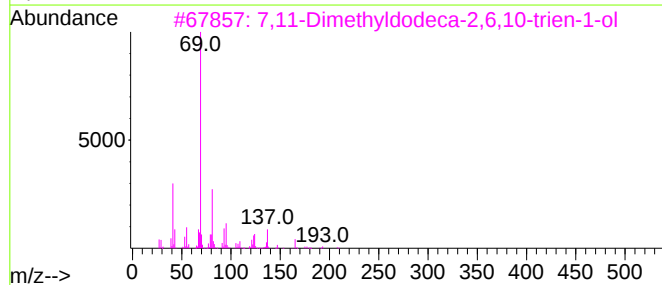
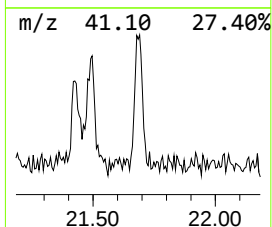
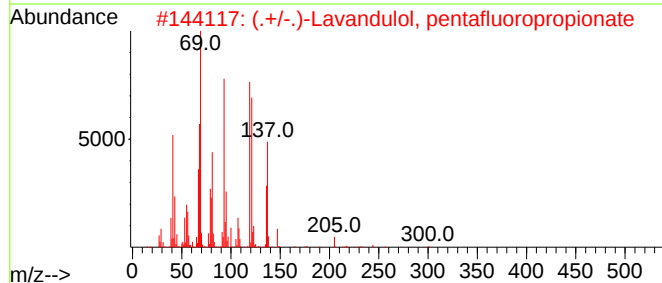
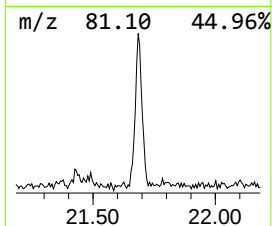
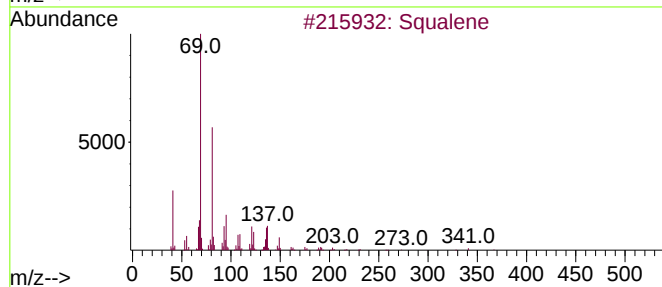
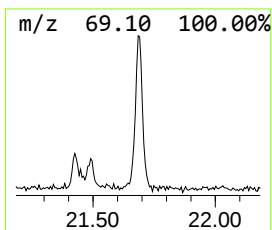
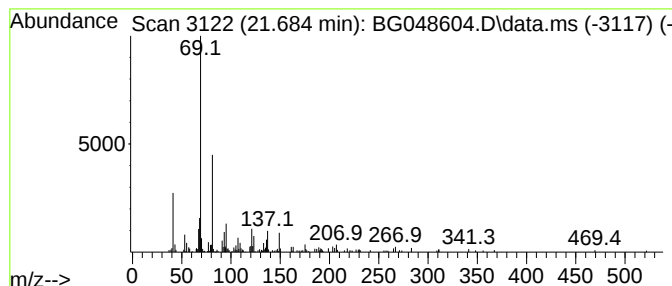
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Squalene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.684	10.66 ng	230893	Chrysene-d12	21.802

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Squalene	410	C30H50	000111-02-4	86
2			(.+/-)-Lavandulol, pentafluorop...	300	C13H17F5O2	1000352-63-1	72
3			7,11-Dimethyldodeca-2,6,10-trien...	208	C14H24O	125529-10-4	64
4			5,9,13-Pentadecatrien-2-one, 6,1...	262	C18H30O	001117-52-8	59
5			Bis(3-methylbut-3-enyl) phthalate	302	C18H22O4	022711-53-1	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG040621\
Data File : BG048604.D
Acq On : 6 Apr 2021 15:05
Operator : CG/JU
Sample : M1933-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

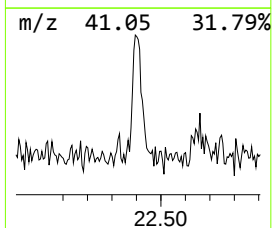
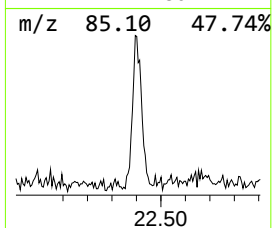
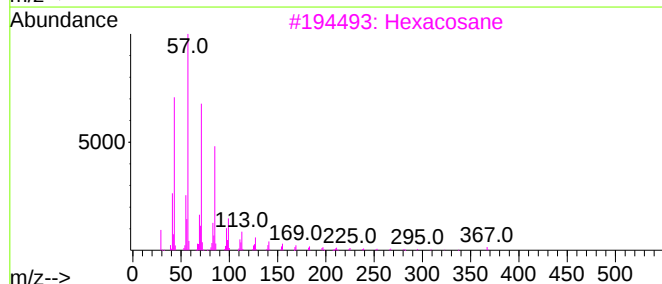
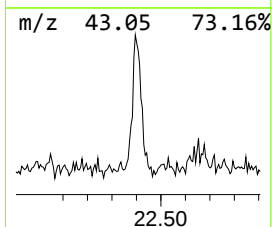
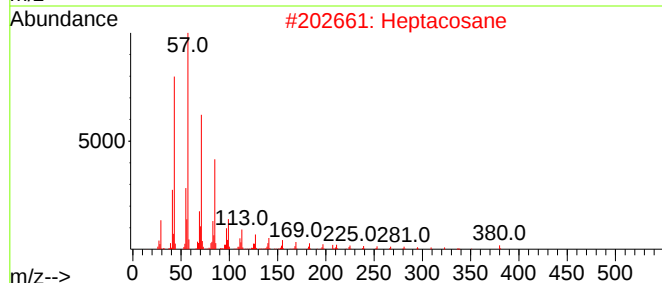
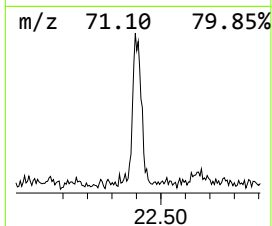
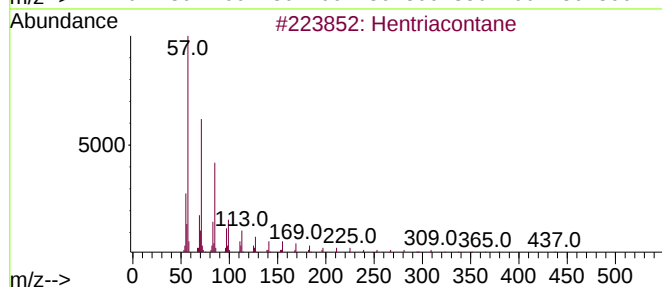
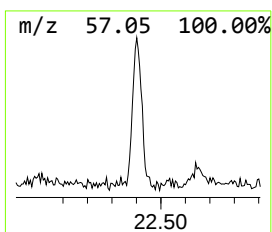
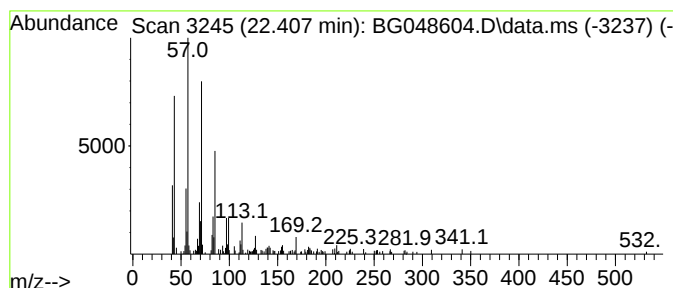
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG033021.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Hentriacontane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
22.407	12.57 ng	272359	Chrysene-d12	21.802	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hentriacontane	437	C31H64	000630-04-6	91
2	Heptacosane	380	C27H56	000593-49-7	91
3	Hexacosane	366	C26H54	000630-01-3	91
4	Pentacosane	352	C25H52	000629-99-2	91
5	Nonadecane	268	C19H40	000629-92-5	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG040621\
Data File : BG048604.D
Acq On : 6 Apr 2021 15:05
Operator : CG/JU
Sample : M1933-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BENNETT_BH_01_2-2.5

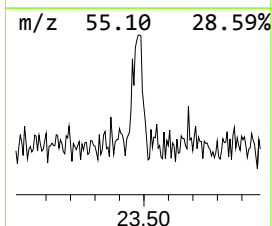
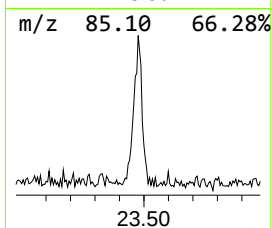
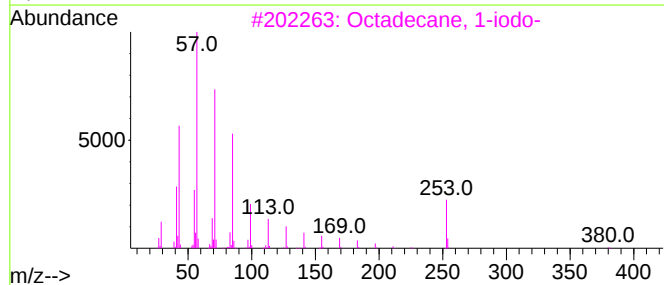
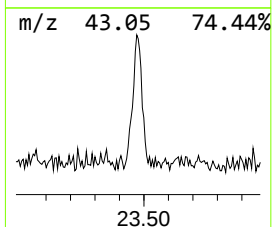
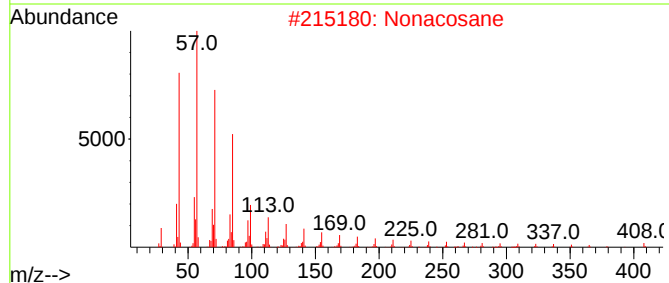
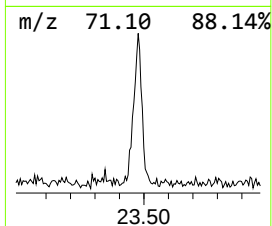
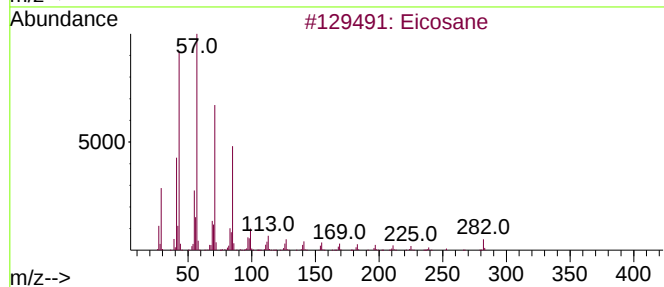
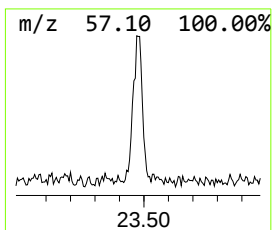
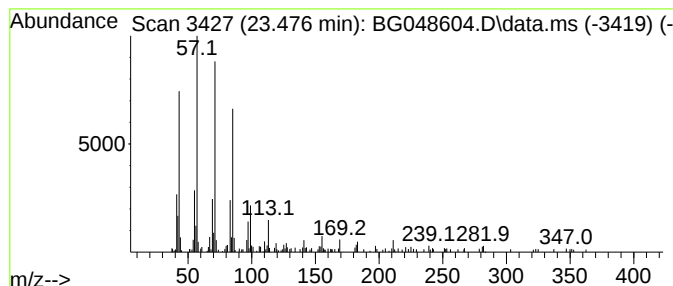
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Nonacosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.476	10.76 ng	233204	Chrysene-d12	21.802

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Eicosane	282	C20H42	000112-95-8	87
2		Nonacosane	408	C29H60	000630-03-5	83
3		Octadecane, 1-iodo-	380	C18H37I	000629-93-6	83
4		2-methyloctacosane	408	C29H60	1000376-72-8	80
5		Docosane	310	C22H46	000629-97-0	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG040621\
Data File : BG048604.D
Acq On : 6 Apr 2021 15:05
Operator : CG/JU
Sample : M1933-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BENNETT_BH_01_2-2.5

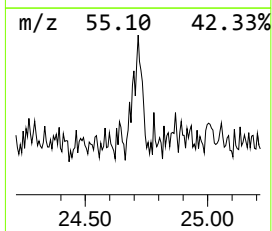
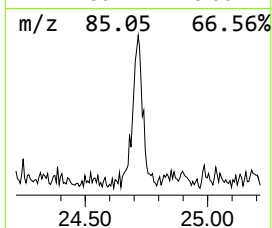
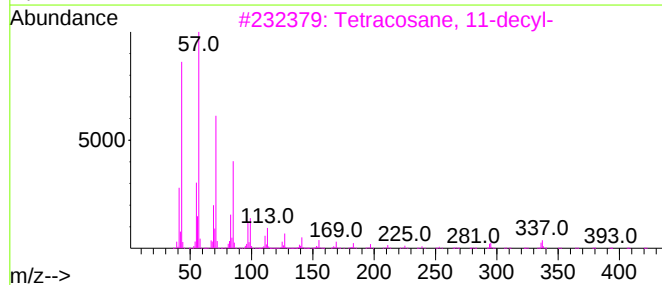
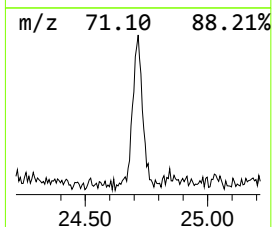
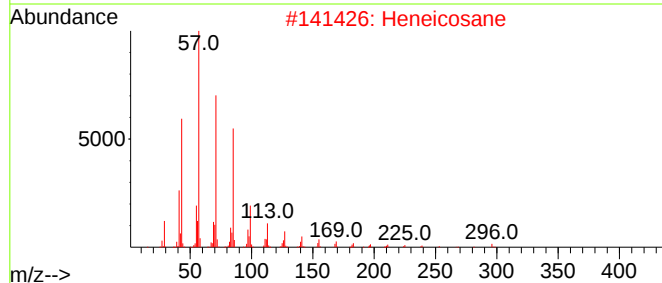
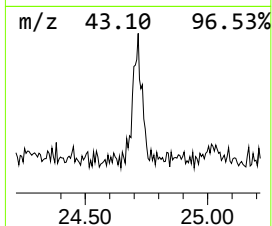
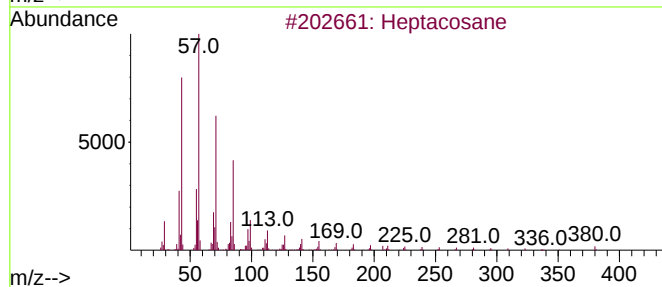
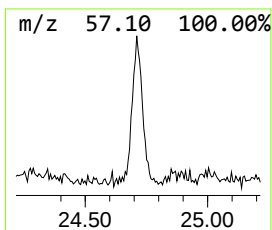
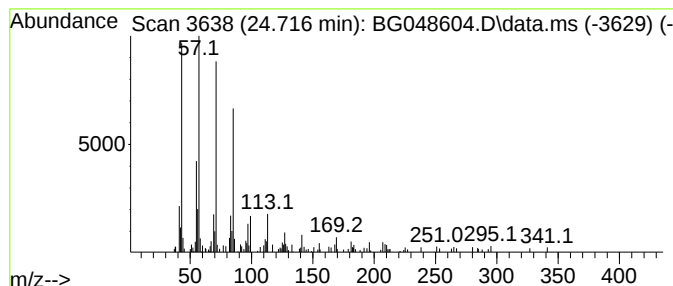
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Heptacosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.716	12.13 ng	243763	Perylene-d12	25.151

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptacosane	380	C27H56	000593-49-7	91
2			Heneicosane	296	C21H44	000629-94-7	91
3			Tetracosane, 11-decyl-	479	C34H70	055429-84-0	91
4			2-methyloctacosane	408	C29H60	1000376-72-8	91
5			Docosane, 11-butyl-	366	C26H54	013475-76-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG040621\
 Data File : BG048604.D
 Acq On : 6 Apr 2021 15:05
 Operator : CG/JU
 Sample : M1933-06
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BENNETT_BH_01_2-2.5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG033021.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
n-Hexadecanoic ...	18.376	8.2	ng	139435	4	17.501	341734	20.0
Benzamide, N-pr...	18.605	104.2	ng	1780180	4	17.501	341734	20.0
Morpholine, 4-p...	18.764	123.0	ng	2101710	4	17.501	341734	20.0
Ethanol, 2-(4-p...	18.911	300.3	ng	5130730	4	17.501	341734	20.0
Glyoxylamide, N...	19.058	20.4	ng	348466	4	17.501	341734	20.0
Phthalic acid, ...	19.275	3.7	ng	63070	4	17.501	341734	20.0
Octadecanoic acid	19.645	38.4	ng	655867	4	17.501	341734	20.0
Eicosane	19.845	7.0	ng	150540	5	21.802	433364	20.0
Pentadecane, 4-...	20.691	10.4	ng	226038	5	21.802	433364	20.0
Pentafluoroprop...	21.426	6.9	ng	148944	5	21.802	433364	20.0
Octacosane	21.490	10.9	ng	236124	5	21.802	433364	20.0
Squalene	21.684	10.7	ng	230893	5	21.802	433364	20.0
Hentriacontane	22.407	12.6	ng	272359	5	21.802	433364	20.0
Nonacosane	23.476	10.8	ng	233204	5	21.802	433364	20.0
Heptacosane	24.716	12.1	ng	243763	6	25.151	402012	20.0