

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG101321\
 Data File : BG050572.D
 Acq On : 13 Oct 2021 16:08
 Operator : CG/JU
 Sample : M4047-03
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 149-ENV-1-COMPOSITE

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-BG100621.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BG050572.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.794	385	392	404	rBV	620263	1017140	12.76%	2.976%
2	7.398	657	665	675	rBV	620574	1100465	13.81%	3.220%
3	8.256	804	811	818	rBV	146383	257719	3.23%	0.754%
4	9.431	1001	1011	1018	rBV	377632	701168	8.80%	2.052%
5	10.829	1243	1249	1261	rVB	132370	230220	2.89%	0.674%
6	11.082	1285	1292	1301	rVB	189233	351542	4.41%	1.029%
7	13.503	1696	1704	1710	rBV	793933	1273255	15.97%	3.725%
8	14.878	1932	1938	1947	rBV	291532	464019	5.82%	1.358%
9	16.358	2184	2190	2199	rBV	668933	1052510	13.20%	3.080%
10	17.621	2399	2405	2410	rBV	390063	609488	7.65%	1.783%
11	18.138	2483	2493	2504	rVB5	120376	464243	5.82%	1.358%
12	19.173	2657	2669	2680	rBV2	973402	3271009	41.04%	9.571%
13	19.296	2680	2690	2700	rBV	1313691	4300958	53.96%	12.584%
14	19.419	2700	2711	2721	rVB	3045703	7971189	100.00%	23.323%
15	19.554	2727	2734	2742	rBV	273798	685590	8.60%	2.006%
16	19.725	2755	2763	2786	rVB2	526321	1626529	20.41%	4.759%
17	20.207	2839	2845	2857	rVB2	1452925	2543724	31.91%	7.443%
18	21.000	2974	2980	2988	rVB	531317	880829	11.05%	2.577%
19	21.617	3079	3085	3093	rVB	286904	456398	5.73%	1.335%
20	21.828	3114	3121	3129	rVB	367295	752857	9.44%	2.203%
21	21.916	3130	3136	3143	rVV2	362415	637385	8.00%	1.865%
22	22.751	3270	3278	3287	rBV2	371313	847193	10.63%	2.479%
23	23.820	3453	3460	3471	rVB	236539	645193	8.09%	1.888%
24	25.083	3668	3675	3686	rVB2	207686	621770	7.80%	1.819%
25	25.330	3710	3717	3726	rVB2	208594	570918	7.16%	1.670%
26	26.593	3922	3932	3943	rVB3	105509	428151	5.37%	1.253%
27	28.409	4231	4241	4256	rVB4	85517	416215	5.22%	1.218%

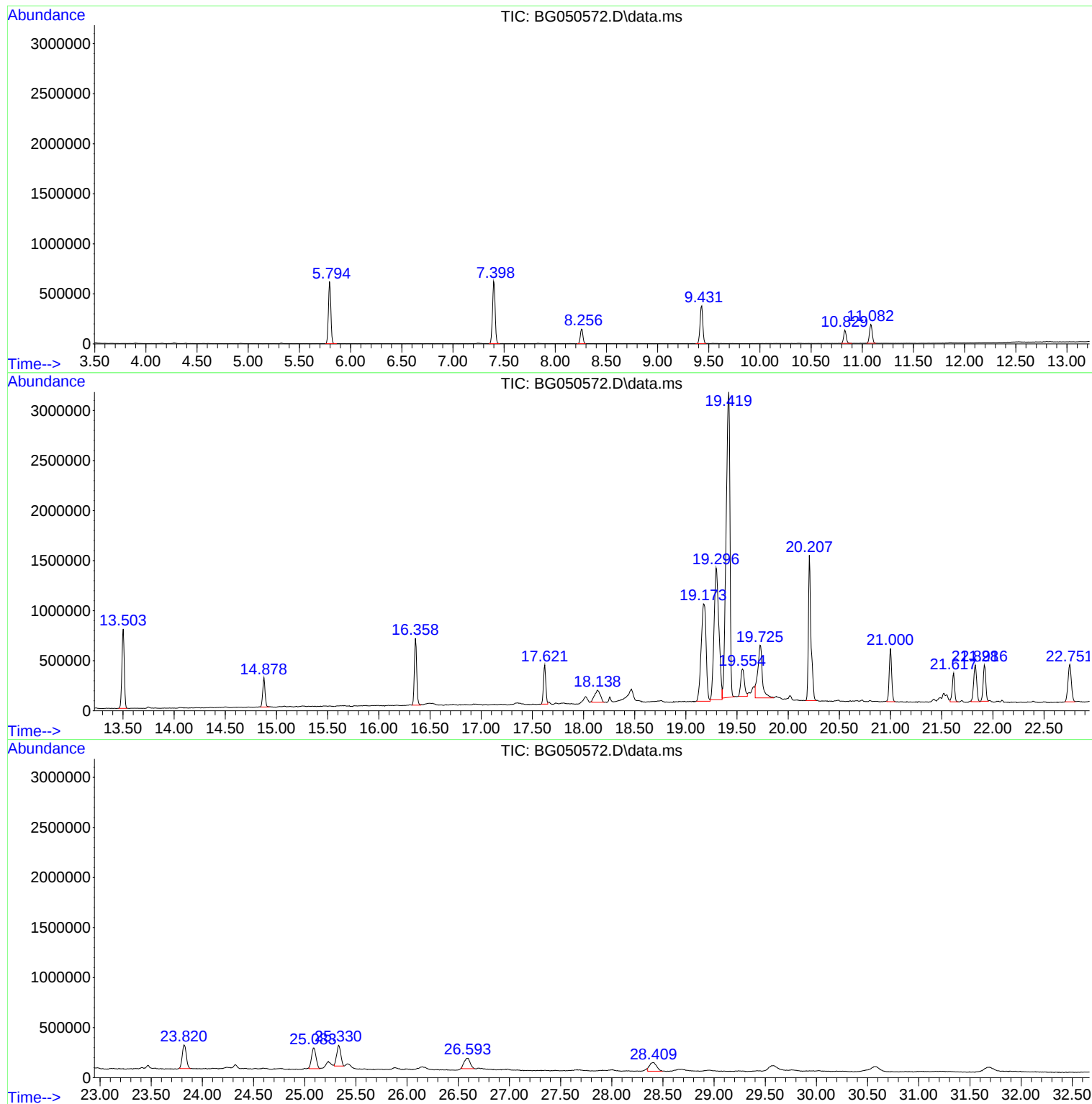
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG100621.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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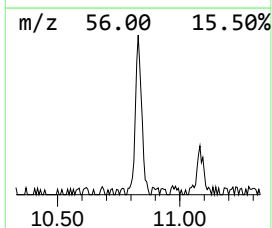
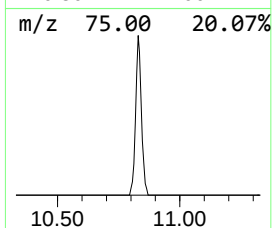
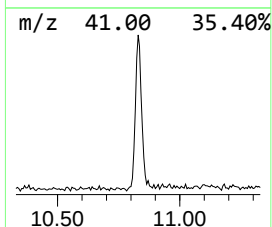
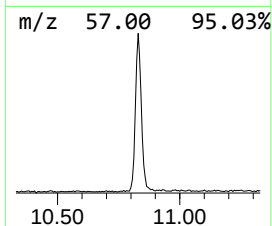
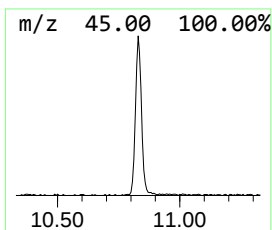
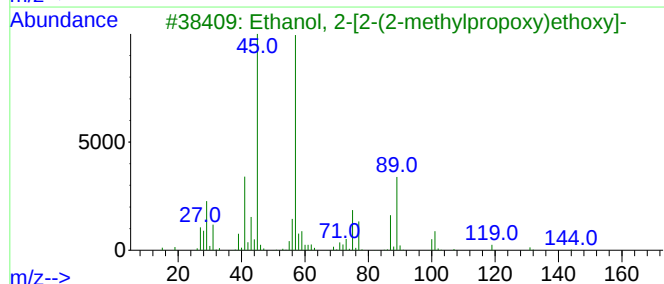
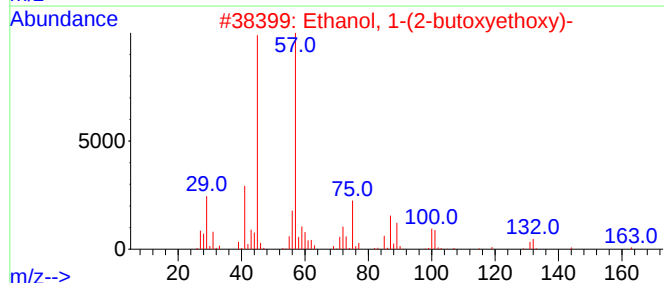
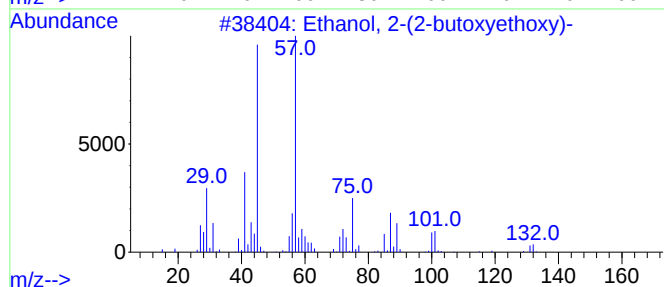
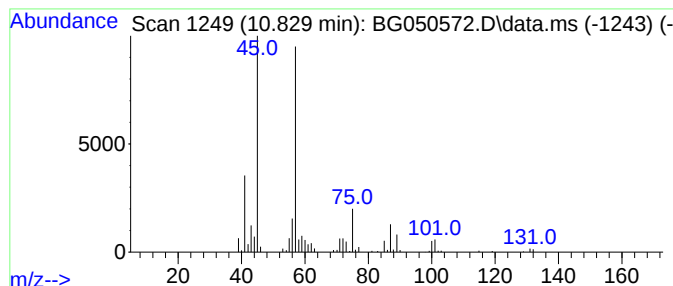
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG100621.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.829	13.10 ng	230220	Naphthalene-d8	11.082	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
2	Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	90
3	Ethanol, 2-[2-(2-methylpropoxy)e...	162	C8H18O3	018912-80-6	59
4	Di-sec-Butyl ether	130	C8H18O	006863-58-7	53
5	3,3-Dimethylbutane-2-ol	102	C6H14O	000464-07-3	50



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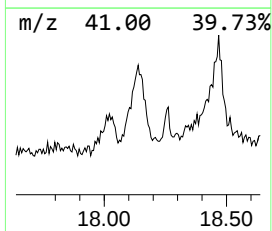
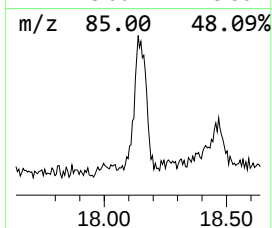
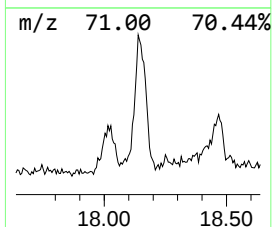
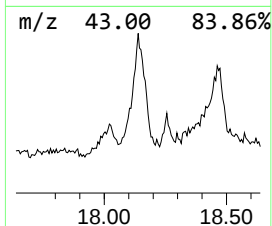
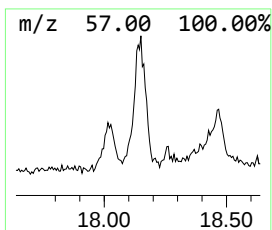
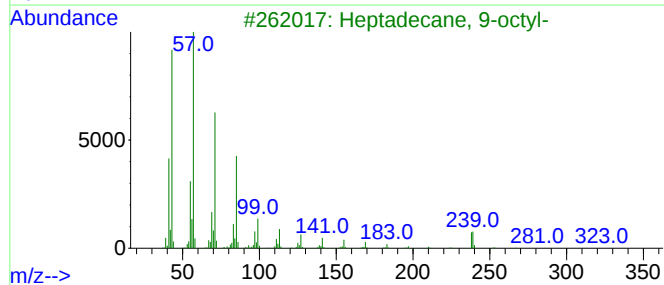
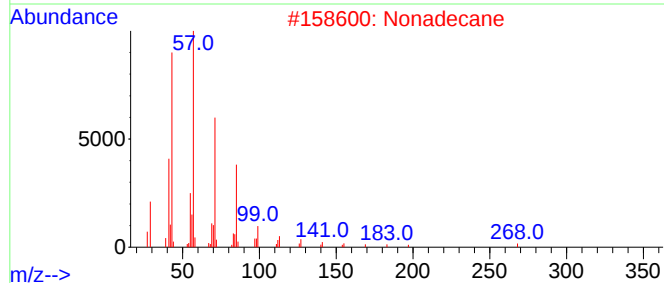
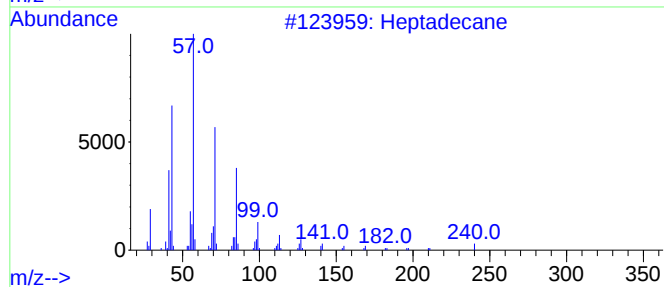
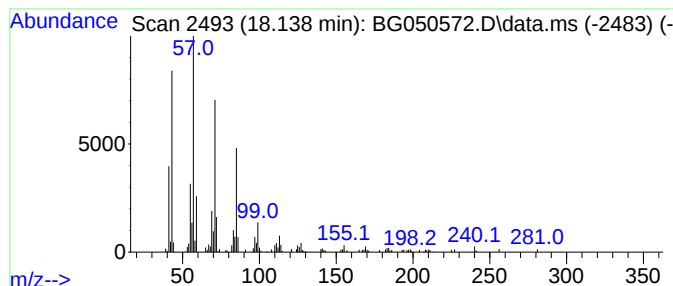
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TIC Library : C:\Database\NIST20.L
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Peak Number 2 Heptadecane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.139	15.23 ng	464243	Phenanthrene-d10	17.621

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecane	240	C17H36	000629-78-7	93
2			Nonadecane	268	C19H40	000629-92-5	68
3			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	68
4			Hexacosane	366	C26H54	000630-01-3	68
5			Octadecane, 1-iodo-	380	C18H37I	000629-93-6	64



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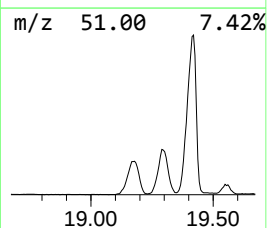
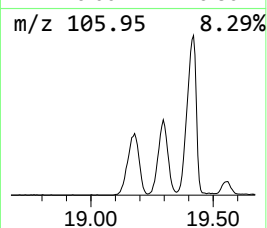
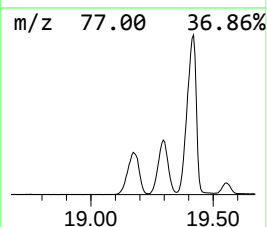
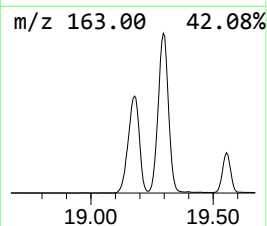
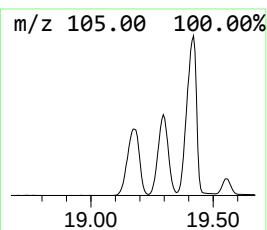
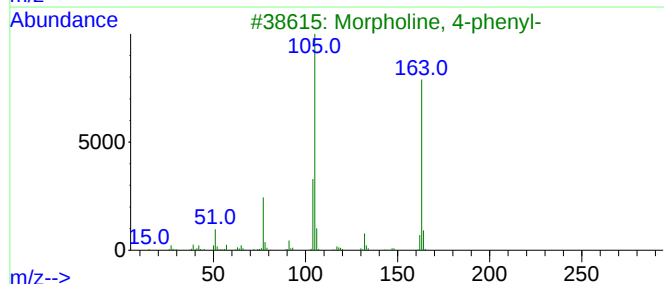
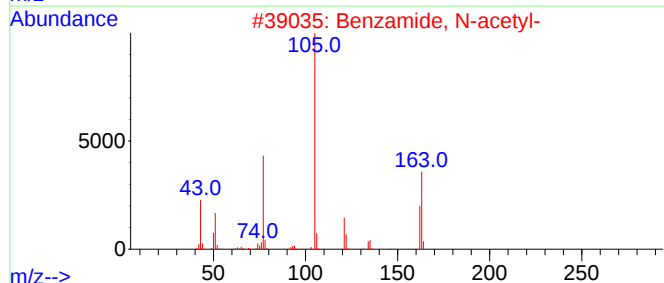
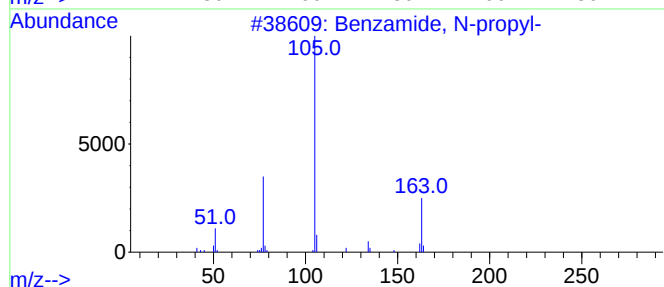
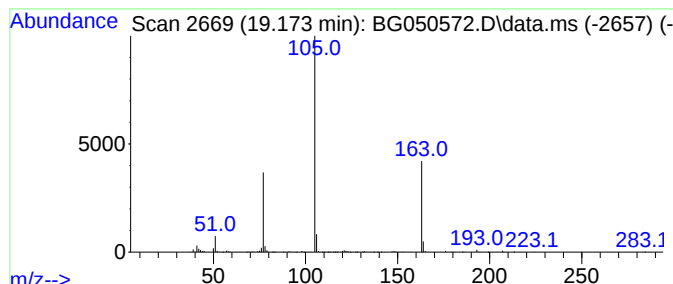
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TIC Library : C:\Database\NIST20.L
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Peak Number 3 Benzamide, N-propyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.172	107.34 ng	3271010	Phenanthrene-d10	17.621

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzamide, N-propyl-	163	C10H13NO	010546-70-0	72
2			Benzamide, N-acetyl-	163	C9H9NO2	001575-95-7	56
3			Morpholine, 4-phenyl-	163	C10H13NO	000092-53-5	53
4			Phenylglyoxylic Acid, 3-methylbu...	220	C13H16O3	1000453-46-1	38
5			1-(Benzoyloxy)propan-2-yl benzoate	284	C17H16O4	019224-26-1	38



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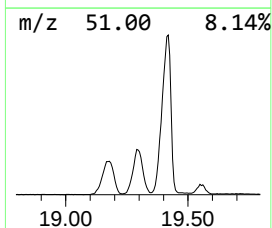
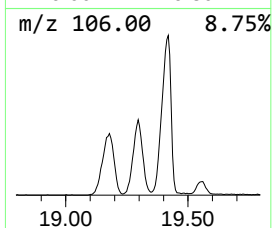
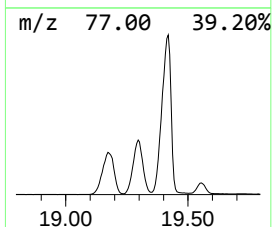
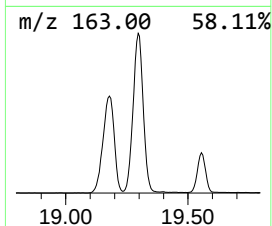
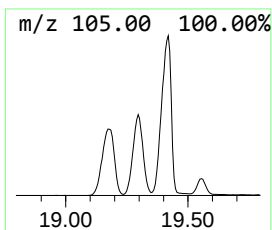
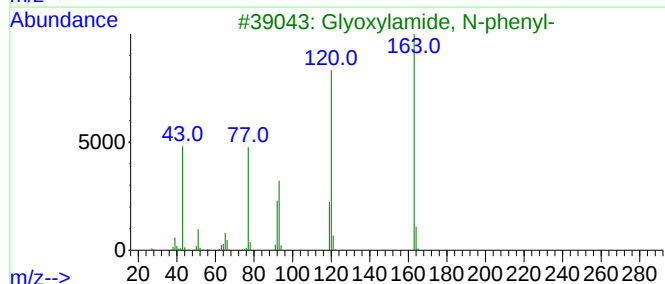
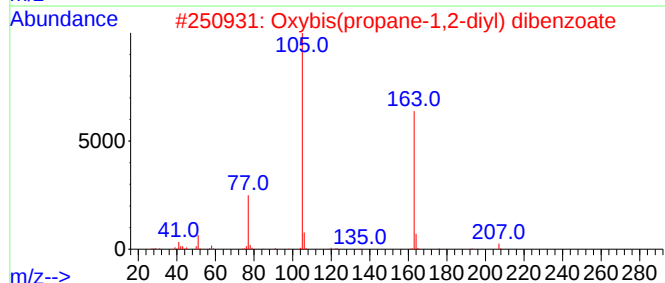
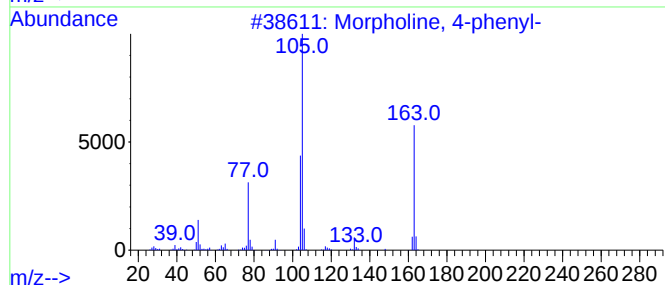
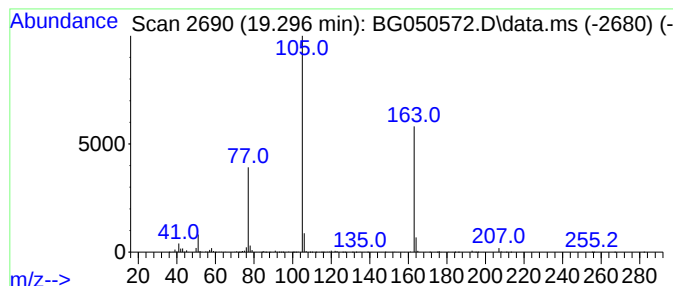
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TIC Library : C:\Database\NIST20.L
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Peak Number 4 Morpholine, 4-phenyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.296	141.13 ng	4300960	Phenanthrene-d10	17.621

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Morpholine, 4-phenyl-	163	C10H13NO	000092-53-5	72
2			Oxybis(propane-1,2-diyl) dibenzoate	342	C20H22O5	000094-03-1	58
3			Glyoxylamide, N-phenyl-	163	C9H9NO2	032331-52-5	53
4			Benzoic acid, phenyl ester	198	C13H10O2	000093-99-2	43
5			N-(2-Fluoro-5-nitrophenyl)benzamide	260	C13H9FN2O3	144205-37-8	43



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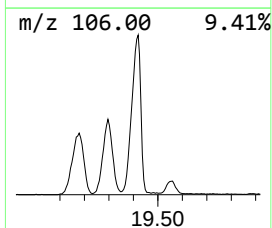
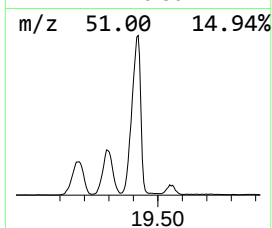
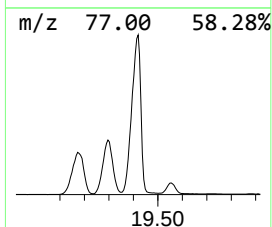
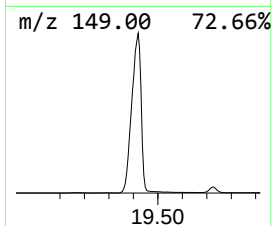
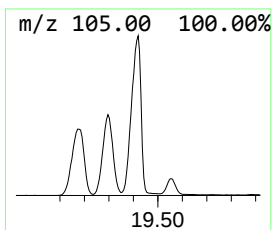
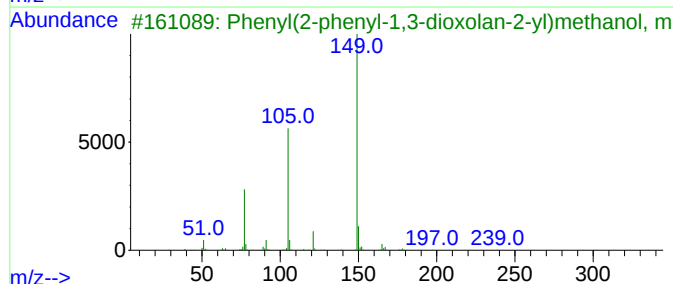
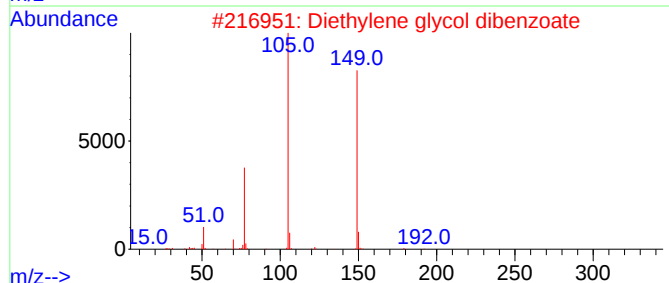
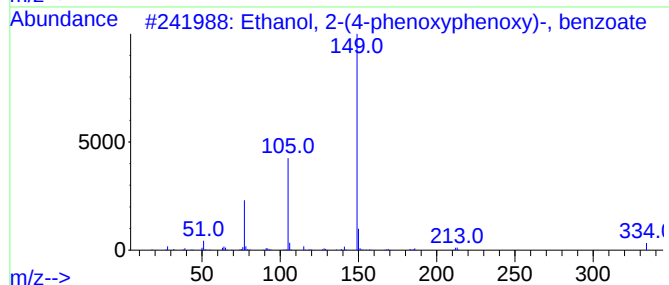
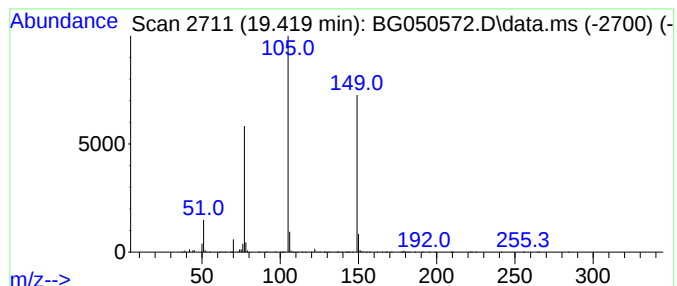
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Peak Number 5 Ethanol, 2-(4-phenoxypheno... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.419	261.57 ng	7971190	Phenanthrene-d10	17.621

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	80
3			Phenyl(2-phenyl-1,3-dioxolan-2-y...	270	C17H18O3	1000507-07-6	78
4			Benzamide, N,N-dimethyl-	149	C9H11NO	000611-74-5	78
5			N-[(2-Phenyl-1,3-dioxolan-2-yl)m...	193	C10H11NO3	1000411-48-9	78



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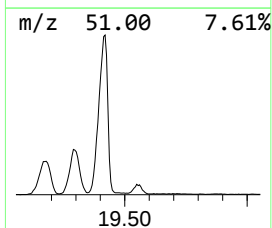
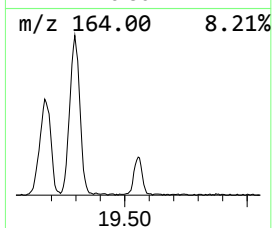
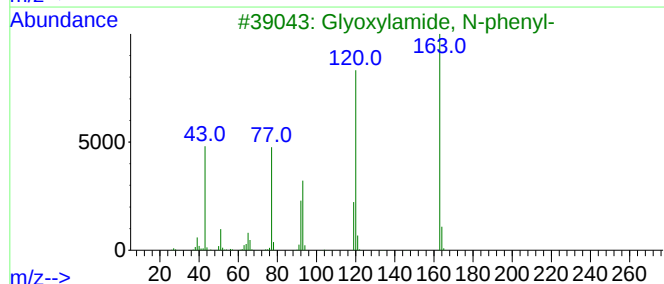
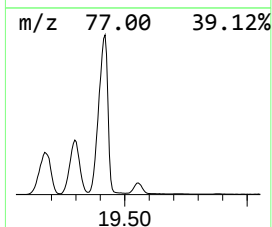
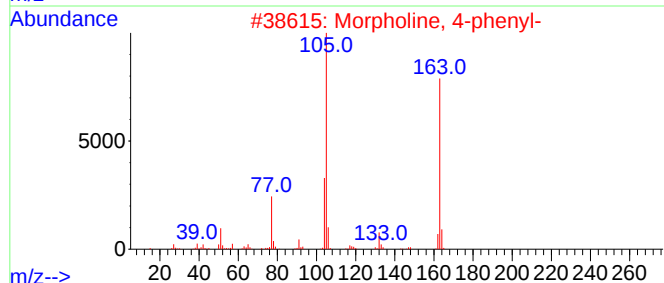
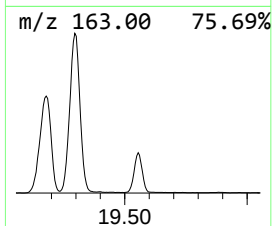
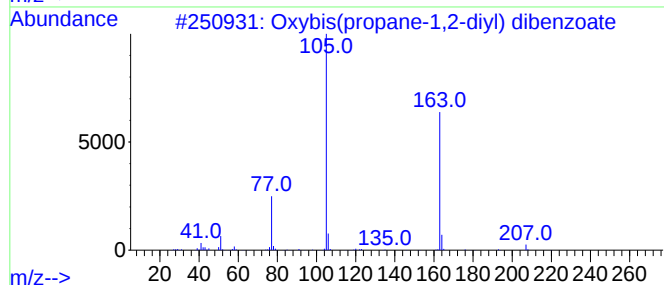
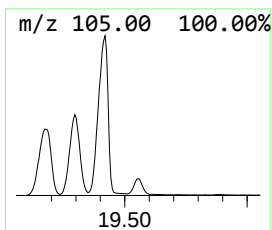
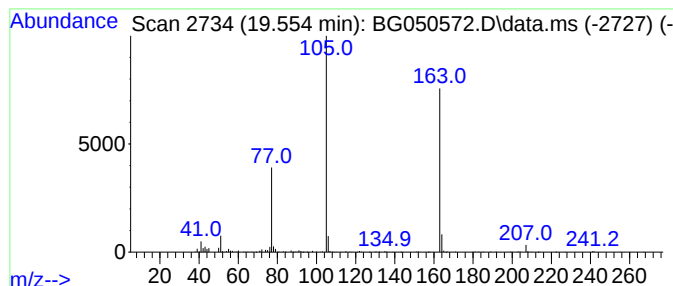
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Oxybis(propene-1,2-diyl) di... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.554	22.50 ng	685590	Phenanthrene-d10	17.621

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Oxybis(propene-1,2-diyl) dibenzoate	342	C20H22O5	000094-03-1	72
2			Morpholine, 4-phenyl-	163	C10H13NO	000092-53-5	59
3			Glyoxylamide, N-phenyl-	163	C9H9NO2	032331-52-5	59
4			N-Methoxy-N-methylbenzamide	165	C9H11NO2	006919-61-5	27
5			1-(Benzoyloxy)propan-2-yl benzoate	284	C17H16O4	019224-26-1	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG101321\
Data File : BG050572.D
Acq On : 13 Oct 2021 16:08
Operator : CG/JU
Sample : M4047-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
149-ENV-1-COMPOSITE

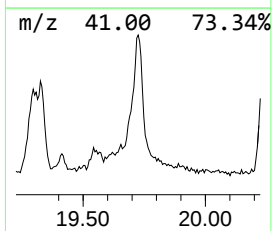
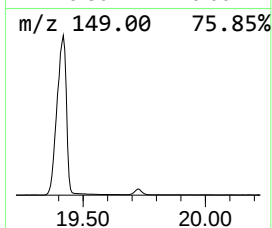
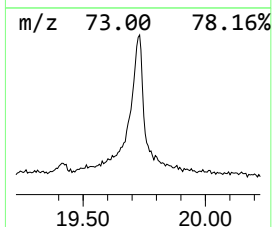
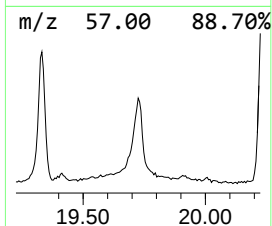
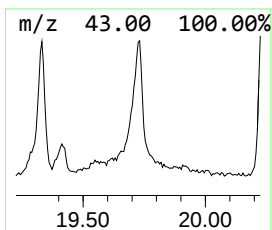
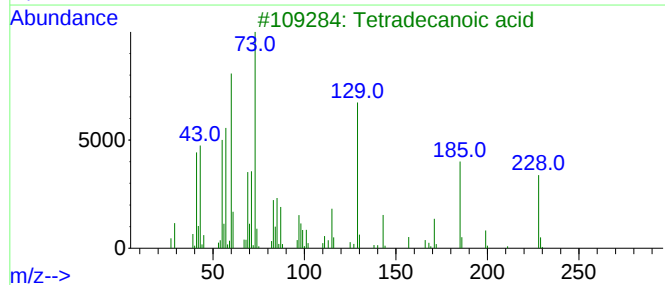
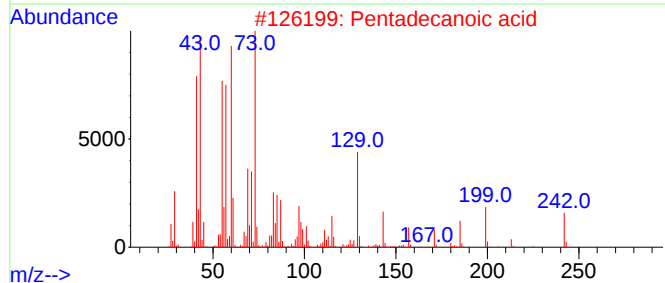
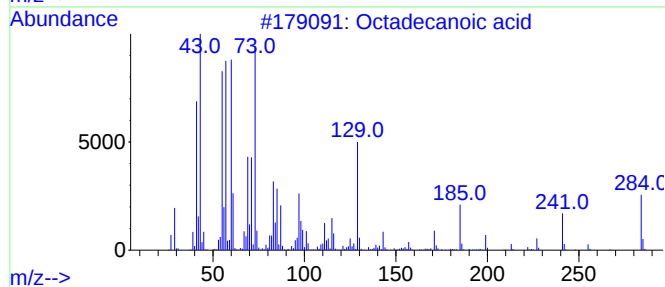
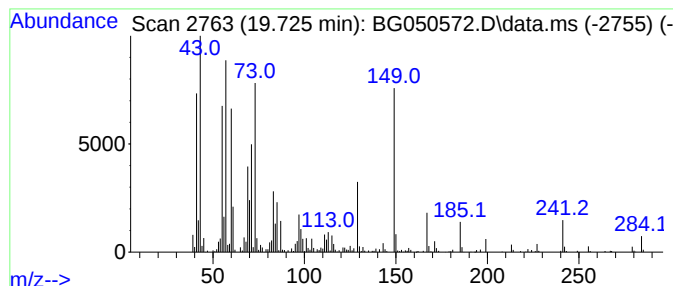
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Octadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.725	53.37 ng	1626530	Phenanthrene-d10	17.621

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	95
2			Pentadecanoic acid	242	C15H30O2	001002-84-2	64
3			Tetradecanoic acid	228	C14H28O2	000544-63-8	50
4			Octadecanoic acid, 2-(2-hydroxye...	372	C22H44O4	000106-11-6	49
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	45



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG101321\
Data File : BG050572.D
Acq On : 13 Oct 2021 16:08
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Sample : M4047-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
149-ENV-1-COMPOSITE

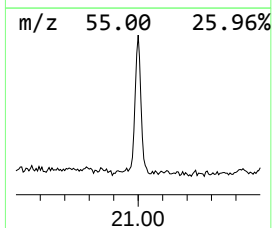
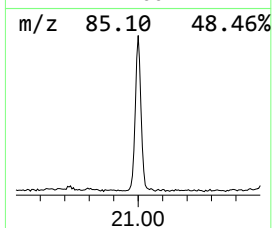
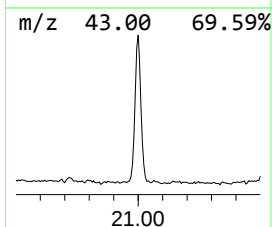
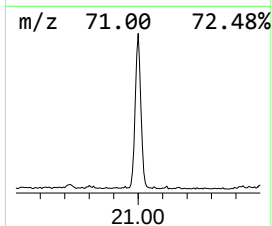
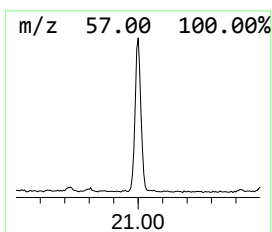
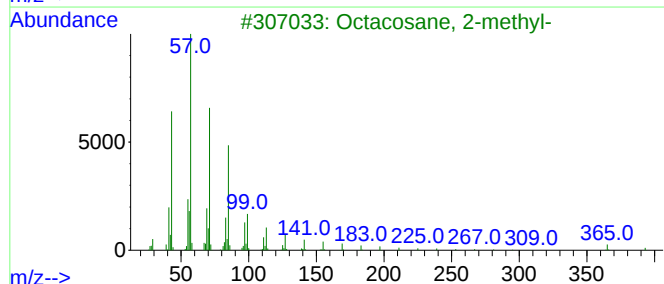
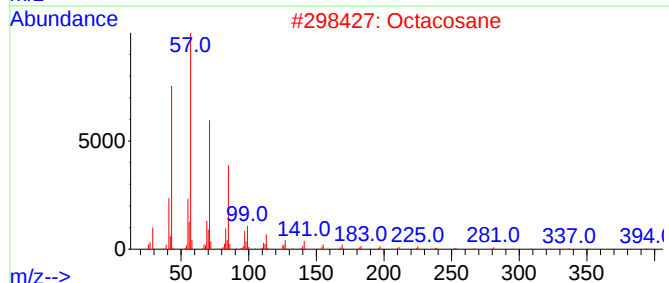
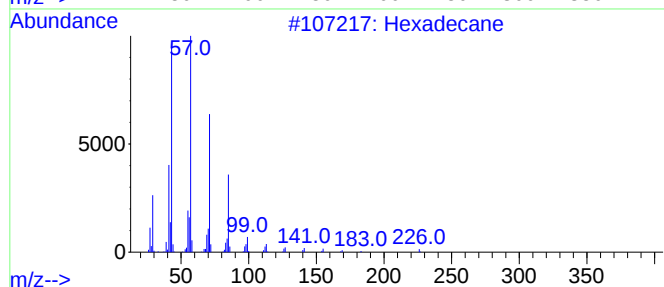
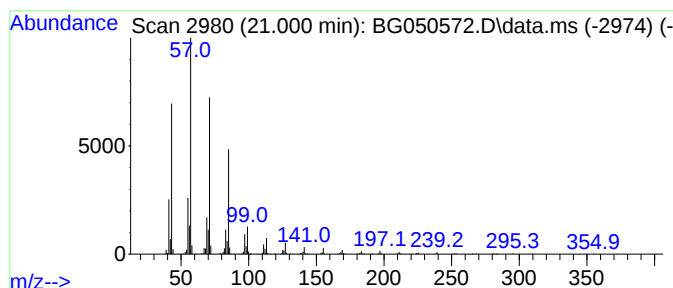
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Hexadecane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.000	27.64 ng	880829	Chrysene-d12	21.916

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hexadecane	226	C16H34	000544-76-3	94
2		Octacosane	394	C28H58	000630-02-4	91
3		Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
4		Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	90
5		Heneicosane	296	C21H44	000629-94-7	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG101321\
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Acq On : 13 Oct 2021 16:08
Operator : CG/JU
Sample : M4047-03
Misc :
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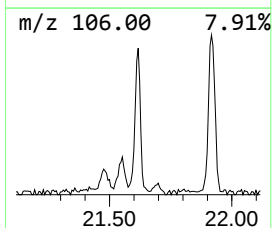
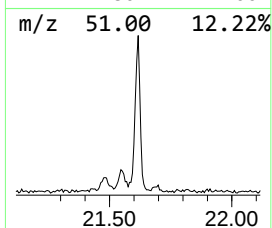
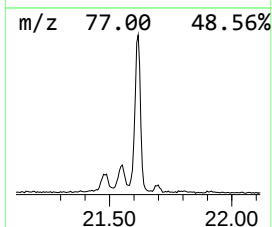
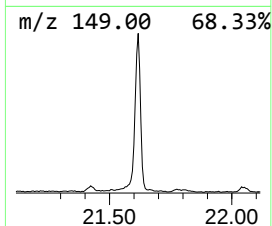
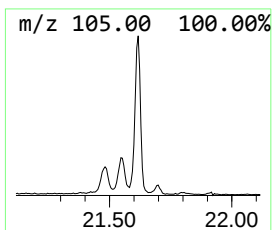
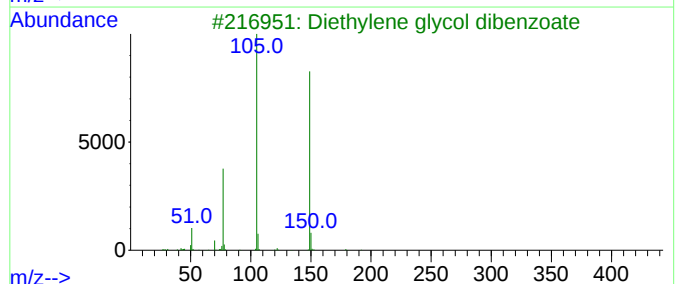
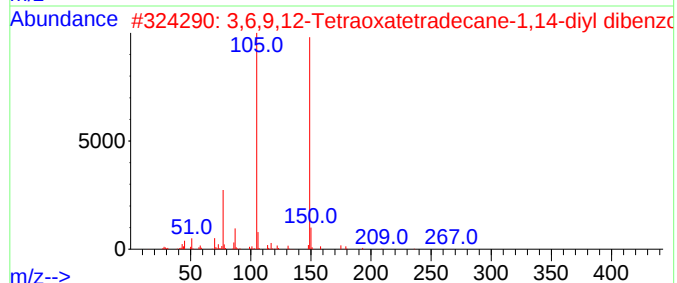
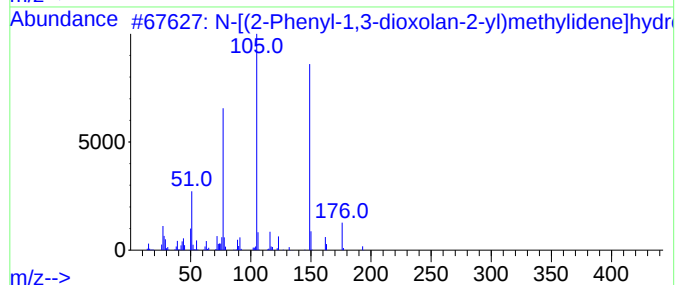
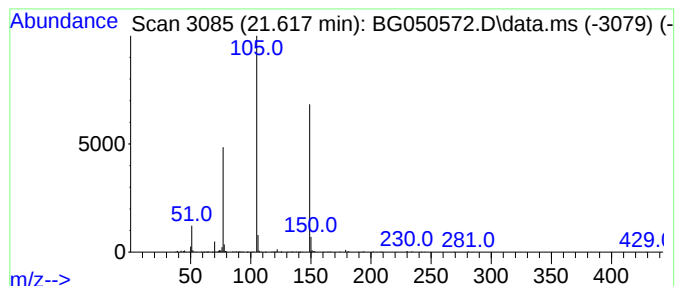
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149-ENV-1-COMPOSITE

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG100621.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 N-[(2-Phenyl-1,3-dioxolan-2-yl)methylidene]hydrazine Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD			R.T.
21.617	14.32 ng	456398	Chrysene-d12			21.916
Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		N-[(2-Phenyl-1,3-dioxolan-2-yl)methylidene]hydrazine	193	C10H11NO3	1000411-48-9	78
2		3,6,9,12-Tetraoxatetradecane-1,14-diyl dibenzoate	446	C24H30O8	1000366-97-0	78
3		Diethylene glycol dibenzoate	314	C18H18O5	000120-55-8	53
4		N-(2-Fluoro-5-nitrophenyl)benzamide	260	C13H9FN2O3	144205-37-8	47
5		N-(2-Methyl-4-nitrophenyl)benzamide	256	C14H12N2O3	104478-92-4	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG101321\
Data File : BG050572.D
Acq On : 13 Oct 2021 16:08
Operator : CG/JU
Sample : M4047-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
149-ENV-1-COMPOSITE

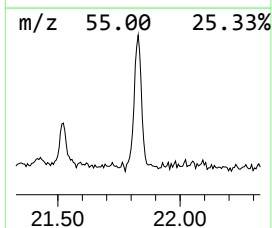
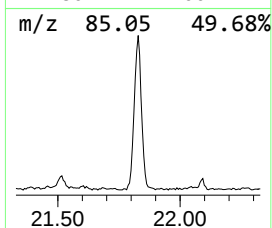
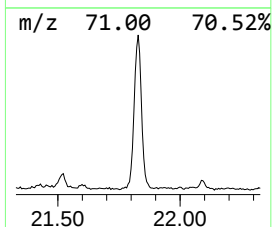
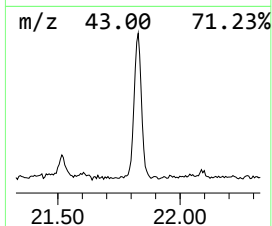
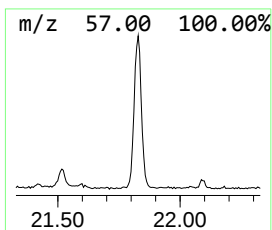
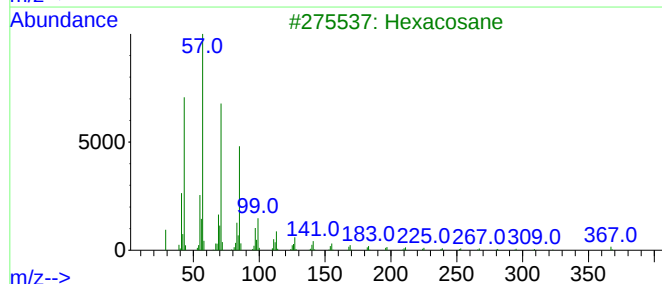
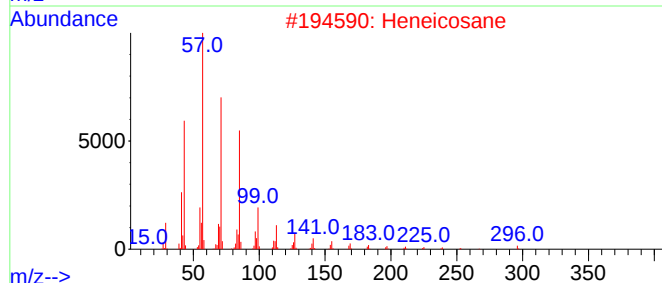
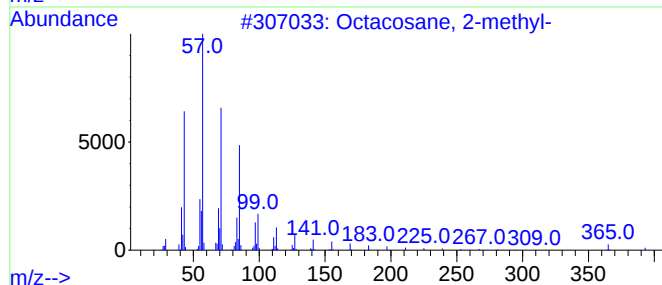
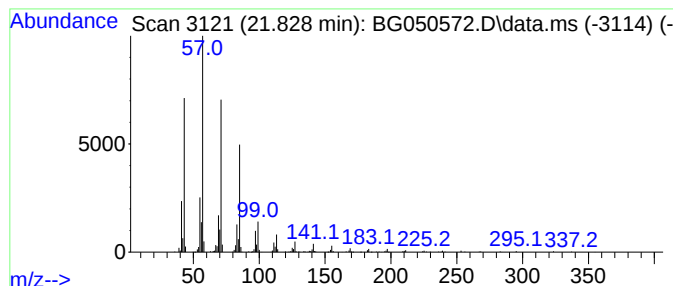
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Octacosane, 2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.828	23.62 ng	752857	Chrysene-d12	21.916

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
2		Heneicosane	296	C21H44	000629-94-7	91
3		Hexacosane	366	C26H54	000630-01-3	90
4		Docosane, 1-iodo-	436	C22H45I	1000406-31-9	87
5		Heptacosane	380	C27H56	000593-49-7	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG101321\
Data File : BG050572.D
Acq On : 13 Oct 2021 16:08
Operator : CG/JU
Sample : M4047-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
149-ENV-1-COMPOSITE

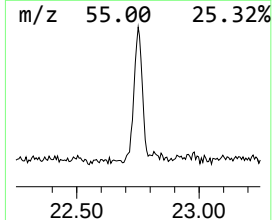
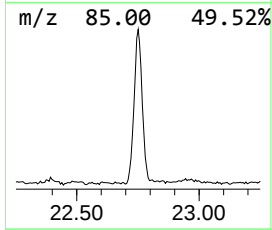
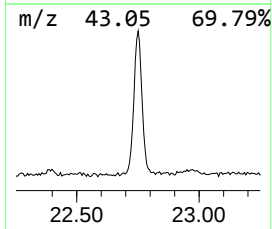
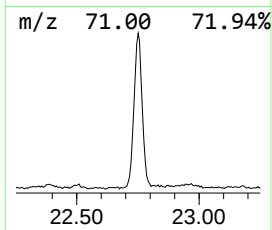
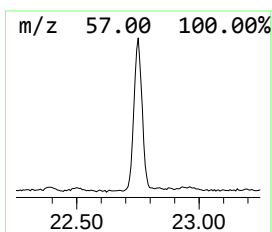
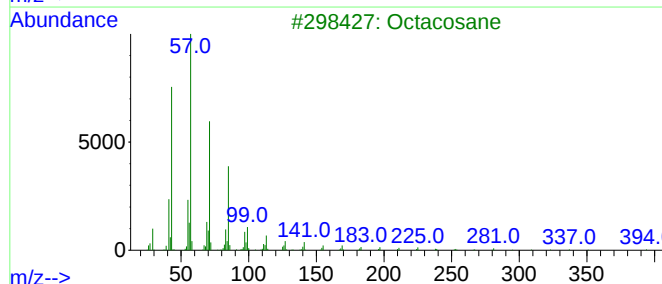
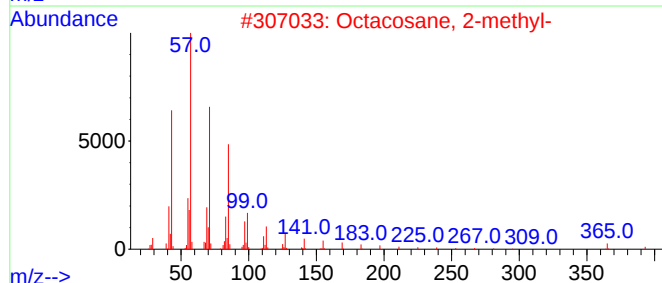
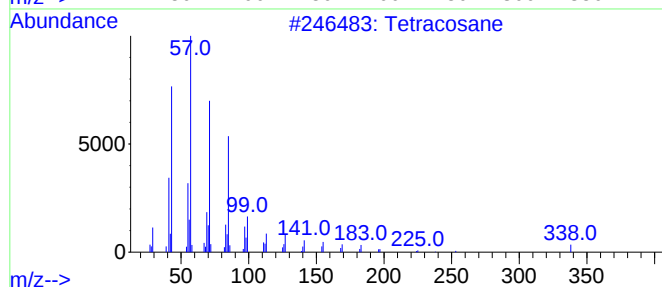
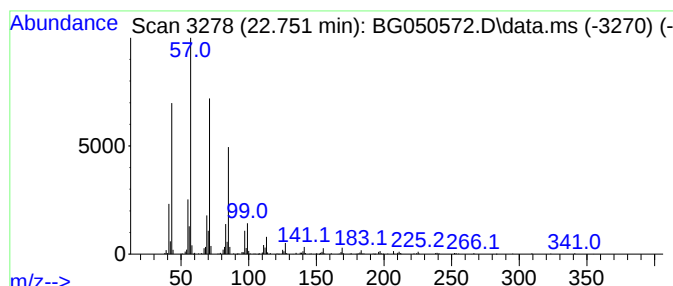
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 Tetracosane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.751	26.58 ng	847193	Chrysene-d12	21.916

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetracosane	338	C24H50	000646-31-1	91
2		Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
3		Octacosane	394	C28H58	000630-02-4	91
4		Heptadecane	240	C17H36	000629-78-7	91
5		Pentacosane	352	C25H52	000629-99-2	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG101321\
Data File : BG050572.D
Acq On : 13 Oct 2021 16:08
Operator : CG/JU
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Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
149-ENV-1-COMPOSITE

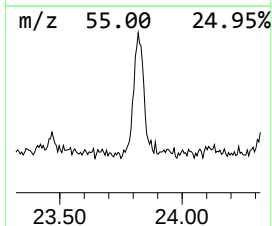
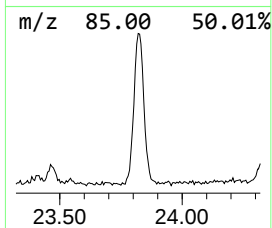
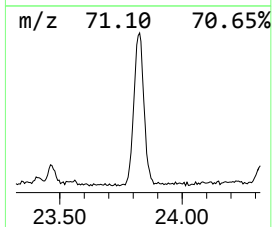
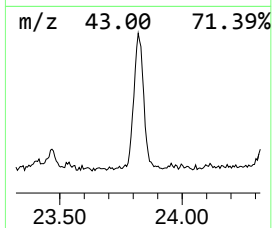
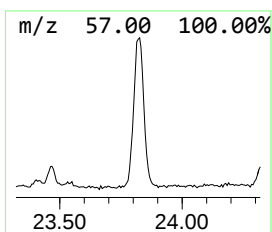
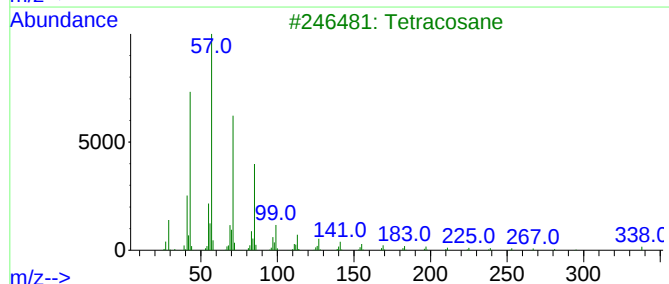
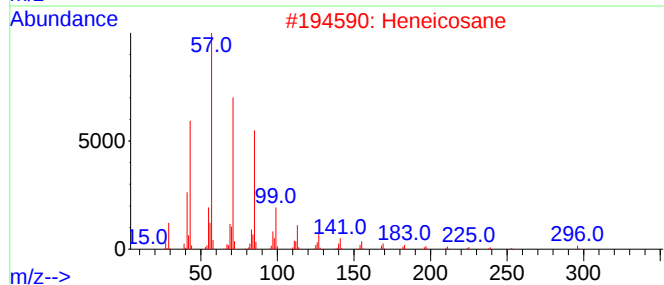
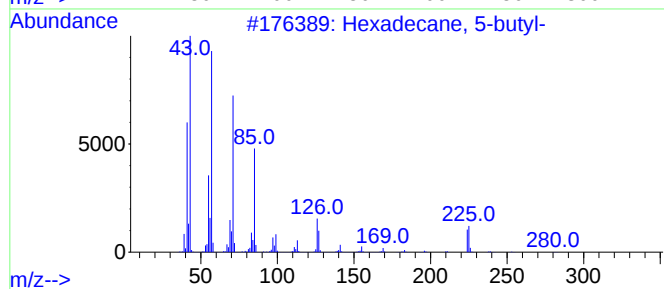
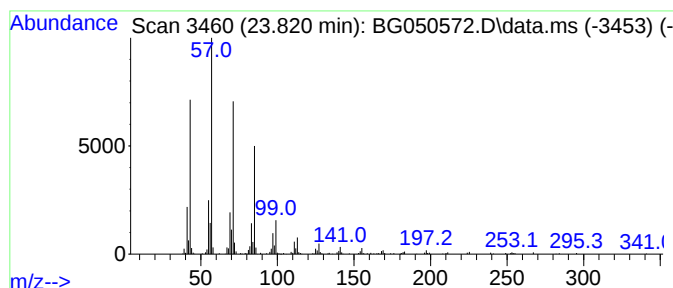
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 Hexadecane, 5-butyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.820	22.60 ng	645193	Perylene-d12	25.330

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane, 5-butyl-	282	C20H42	006912-07-8	93	
2	Heneicosane	296	C21H44	000629-94-7	91	
3	Tetracosane	338	C24H50	000646-31-1	91	
4	Heptadecane, 3-methyl-	254	C18H38	006418-44-6	90	
5	Hexacosane	366	C26H54	000630-01-3	90	



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG101321\
Data File : BG050572.D
Acq On : 13 Oct 2021 16:08
Operator : CG/JU
Sample : M4047-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

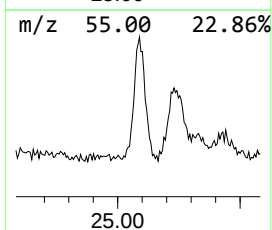
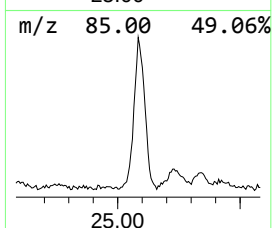
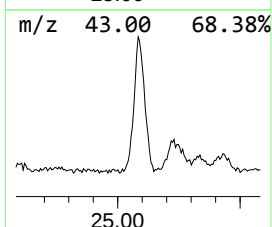
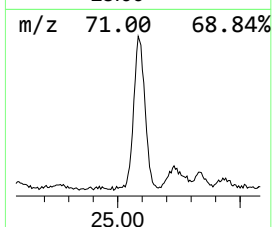
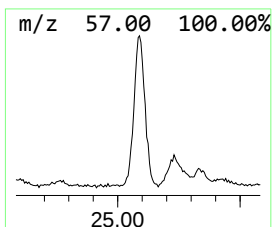
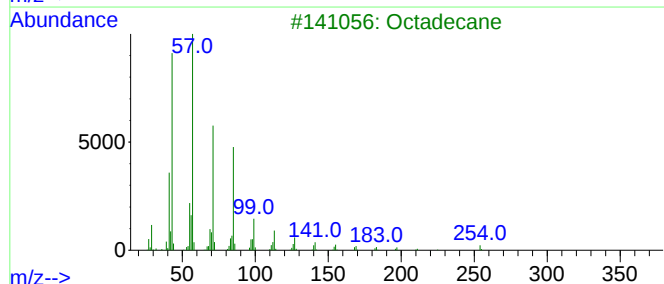
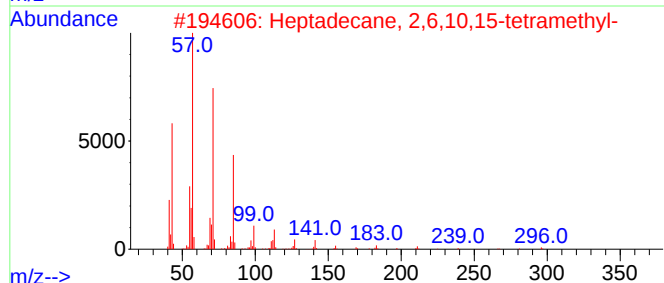
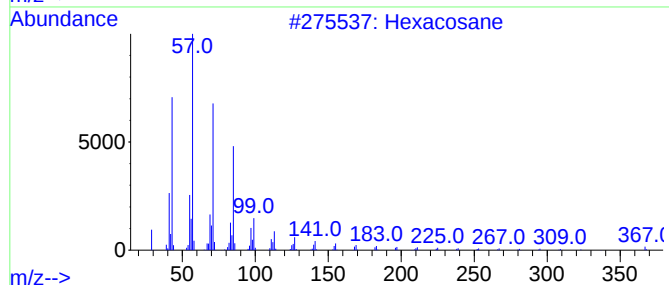
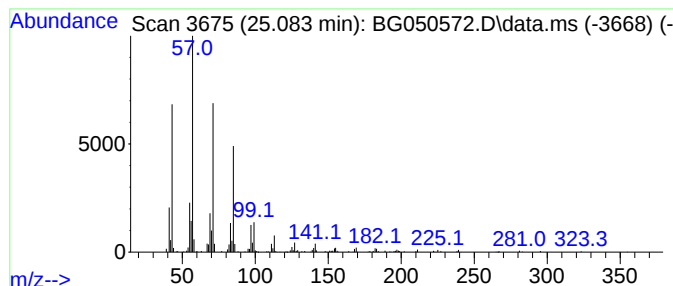
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BNA_G
ClientSampleId :
149-ENV-1-COMPOSITE

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG100621.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 Heptadecane, 2,6,10,15-tetr... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD		R.T.	
25.083	21.78 ng	621770	Perylene-d12		25.330	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Hexacosane		366	C26H54	000630-01-3	91
2	Heptadecane, 2,6,10,15-tetramethyl-		296	C21H44	054833-48-6	87
3	Octadecane		254	C18H38	000593-45-3	87
4	Tetracosane		338	C24H50	000646-31-1	87
5	Tridecane, 1-iodo-		310	C13H27I	035599-77-0	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG101321\
Data File : BG050572.D
Acq On : 13 Oct 2021 16:08
Operator : CG/JU
Sample : M4047-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
149-ENV-1-COMPOSITE

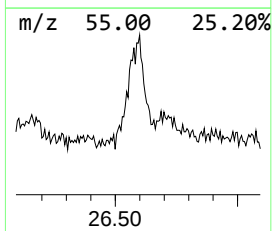
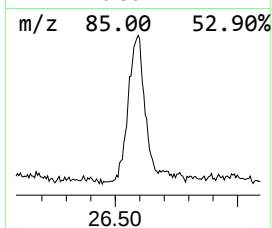
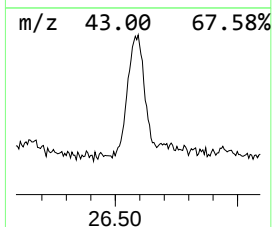
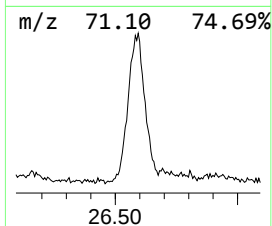
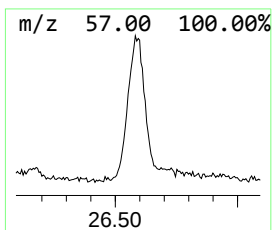
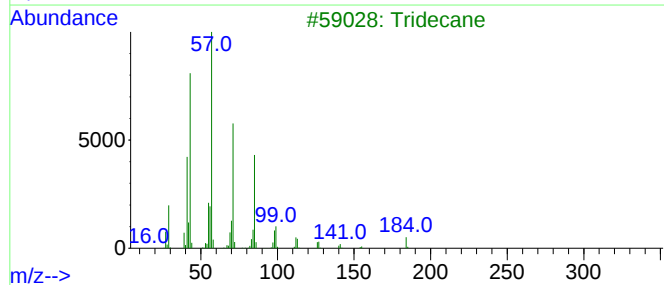
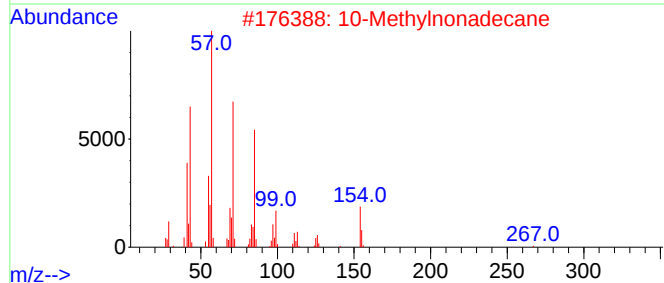
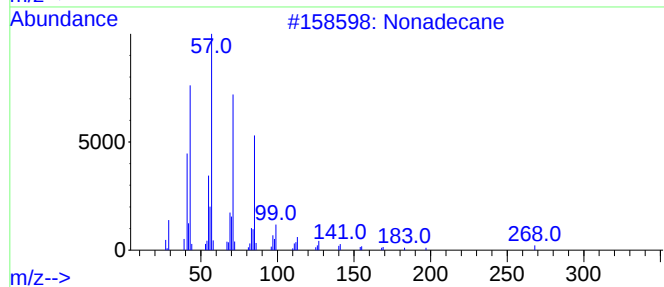
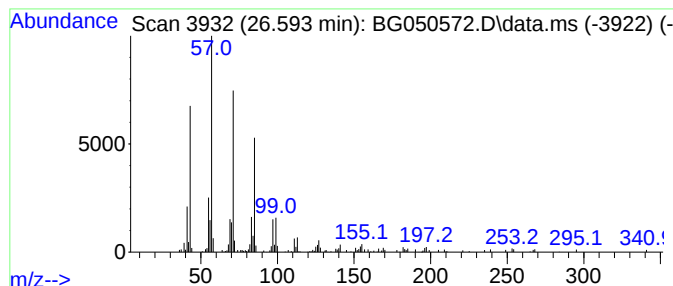
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 Nonadecane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.593	15.00 ng	428151	Perylene-d12	25.330

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonadecane	268	C19H40	000629-92-5	93
2			10-Methylnonadecane	282	C20H42	056862-62-5	87
3			Tridecane	184	C13H28	000629-50-5	87
4			Tetracosane	338	C24H50	000646-31-1	87
5			Dodecane, 2-methyl-	184	C13H28	001560-97-0	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG101321\
Data File : BG050572.D
Acq On : 13 Oct 2021 16:08
Operator : CG/JU
Sample : M4047-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
149-ENV-1-COMPOSITE

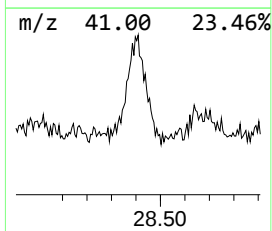
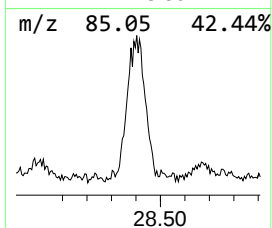
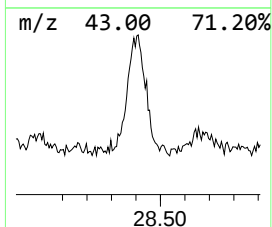
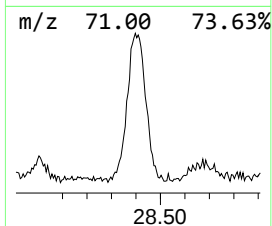
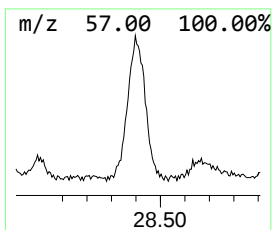
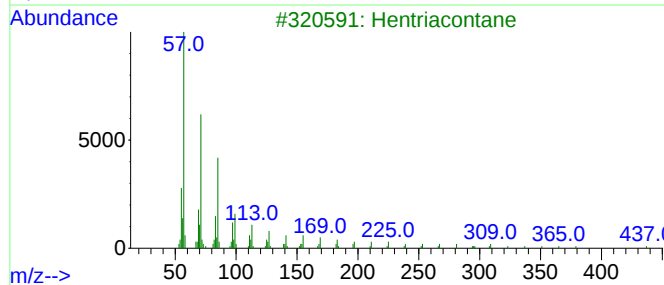
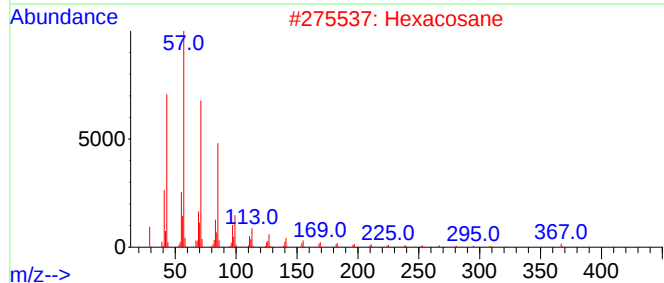
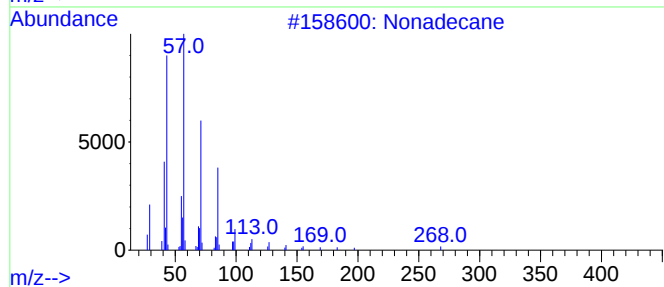
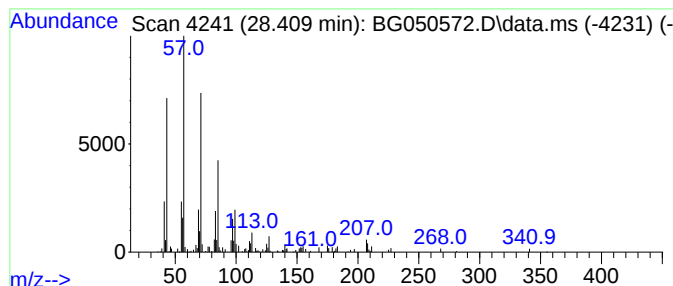
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG100621.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 15 Hexacosane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
28.409	14.58 ng	416215	Perylene-d12	25.330

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Nonadecane	268	C19H40	000629-92-5	91
2		Hexacosane	366	C26H54	000630-01-3	87
3		Hentriacontane	437	C31H64	000630-04-6	87
4		Heneicosane	296	C21H44	000629-94-7	87
5		Octacosane, 2-methyl-	408	C29H60	001560-98-1	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG101321\
 Data File : BG050572.D
 Acq On : 13 Oct 2021 16:08
 Operator : CG/JU
 Sample : M4047-03
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 149-ENV-1-COMPOSITE

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG100621.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST0.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Ethanol, 2-(2-b...	10.829	13.1	ng	230220	2	11.082	351542	20.0
Heptadecane	18.139	15.2	ng	464243	4	17.621	609488	20.0
Benzamide, N-pr...	19.172	107.3	ng	3271010	4	17.621	609488	20.0
Morpholine, 4-p...	19.296	141.1	ng	4300960	4	17.621	609488	20.0
Ethanol, 2-(4-p...	19.419	261.6	ng	7971190	4	17.621	609488	20.0
Oxybis(propane-...	19.554	22.5	ng	685590	4	17.621	609488	20.0
Octadecanoic acid	19.725	53.4	ng	1626530	4	17.621	609488	20.0
Hexadecane	21.000	27.6	ng	880829	5	21.916	637385	20.0
N-[(2-Phenyl-1,...	21.617	14.3	ng	456398	5	21.916	637385	20.0
Octacosane, 2-m...	21.828	23.6	ng	752857	5	21.916	637385	20.0
Tetracosane	22.751	26.6	ng	847193	5	21.916	637385	20.0
Hexadecane, 5-b...	23.820	22.6	ng	645193	6	25.330	570918	20.0
Heptadecane, 2,...	25.083	21.8	ng	621770	6	25.330	570918	20.0
Nonadecane	26.593	15.0	ng	428151	6	25.330	570918	20.0
Hexacosane	28.409	14.6	ng	416215	6	25.330	570918	20.0