

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG101821\
 Data File : BG050617.D
 Acq On : 18 Oct 2021 14:45
 Operator : CG/JU
 Sample : M4259-07
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 WC-1-BD

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: BG050617.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.797	385	393	405	rBV	645542	1072261	14.73%	3.886%
2	7.395	657	665	674	rBV	626224	1137501	15.63%	4.122%
3	8.258	805	812	820	rBV	160407	280308	3.85%	1.016%
4	9.428	1002	1011	1023	rBV	404318	731970	10.06%	2.652%
5	11.079	1284	1292	1300	rBV	213163	398346	5.47%	1.444%
6	13.499	1697	1704	1711	rVB	923163	1421963	19.54%	5.153%
7	14.874	1931	1938	1948	rBV	332047	533934	7.34%	1.935%
8	16.355	2184	2190	2204	rBV	689911	1089874	14.97%	3.949%
9	17.618	2399	2405	2415	rVB	413303	656756	9.02%	2.380%
10	18.423	2532	2542	2547	rBV4	30149	73496	1.01%	0.266%
11	19.345	2686	2699	2709	rBV	854685	2711378	37.25%	9.825%
12	19.463	2709	2719	2727	rVV	1190241	3241468	44.53%	11.746%
13	19.563	2727	2736	2752	rVV	2807034	7278869	100.00%	26.377%
14	19.698	2753	2759	2772	rVB	283173	621825	8.54%	2.253%
15	20.203	2840	2845	2852	rVB	1559254	2111878	29.01%	7.653%
16	20.327	2859	2866	2877	rVB	146548	290375	3.99%	1.052%
17	21.067	2986	2992	2999	rBV	244188	394374	5.42%	1.429%
18	21.519	3065	3069	3072	rBV3	48049	82399	1.13%	0.299%
19	21.613	3080	3085	3091	rVB	170799	283040	3.89%	1.026%
20	21.837	3116	3123	3127	rBV	137845	303383	4.17%	1.099%
21	21.919	3133	3137	3143	rVB	400235	704884	9.68%	2.554%
22	22.800	3279	3287	3299	rVB	189772	443322	6.09%	1.606%
23	23.664	3429	3434	3445	rVB	51309	107953	1.48%	0.391%
24	23.858	3458	3467	3477	rBV2	136575	378069	5.19%	1.370%
25	25.109	3672	3680	3692	rVB2	108230	331294	4.55%	1.201%
26	25.333	3709	3718	3730	rVB2	242188	751843	10.33%	2.725%
27	26.608	3932	3935	3948	rVB3	54626	162794	2.24%	0.590%

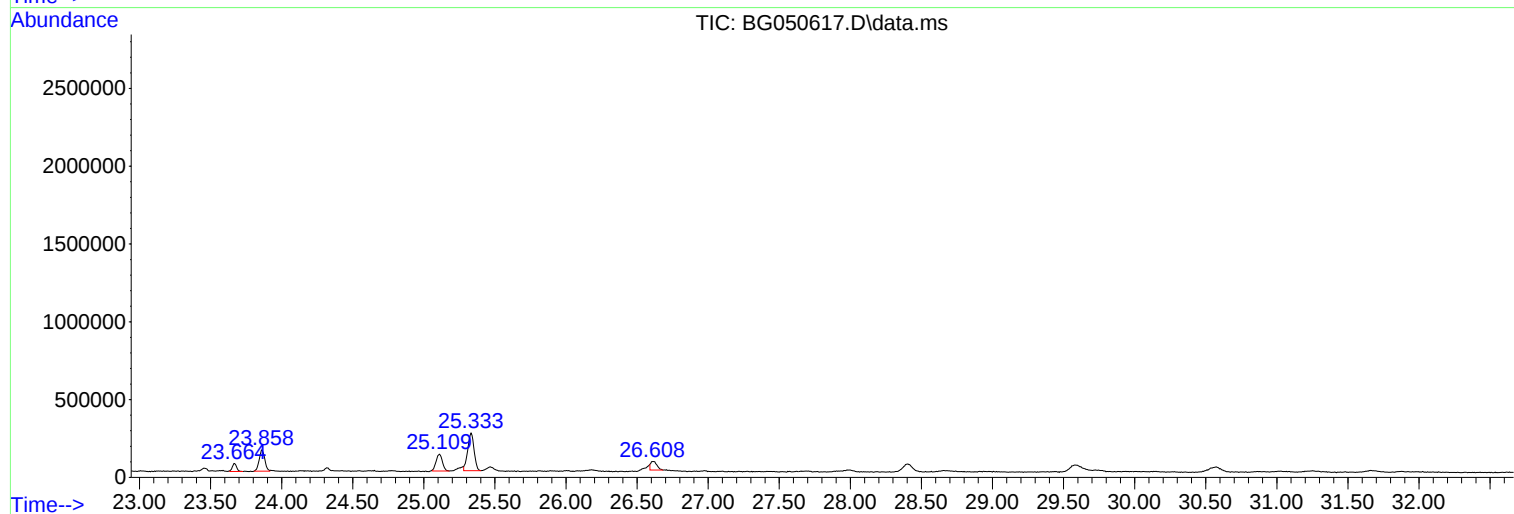
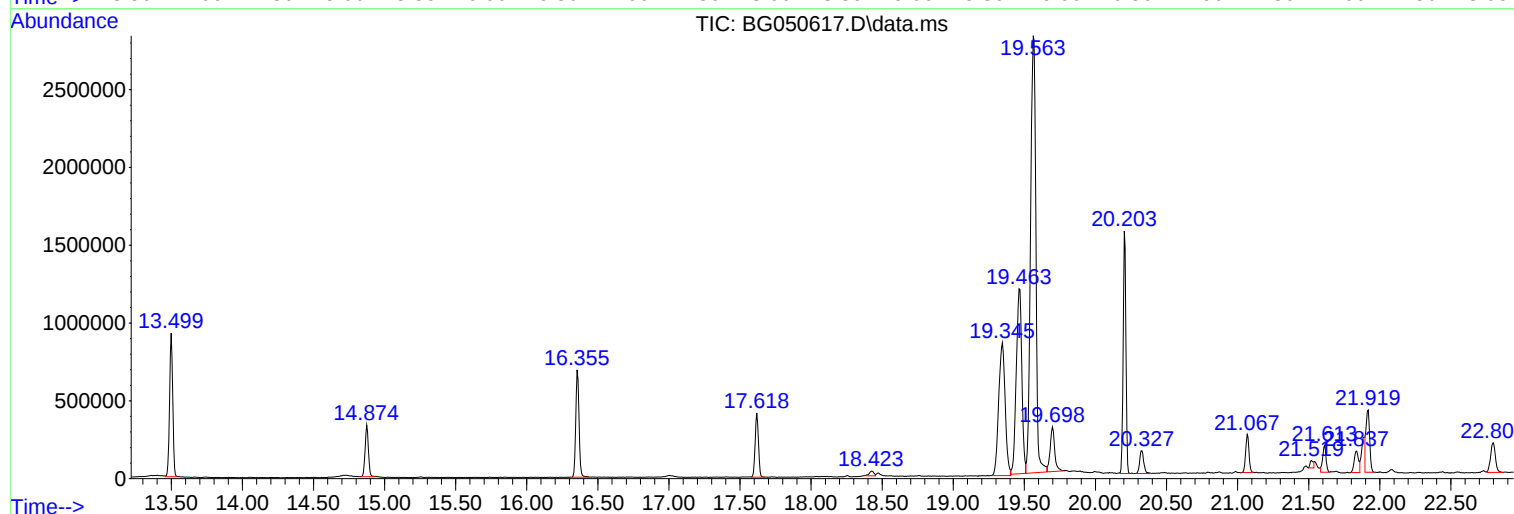
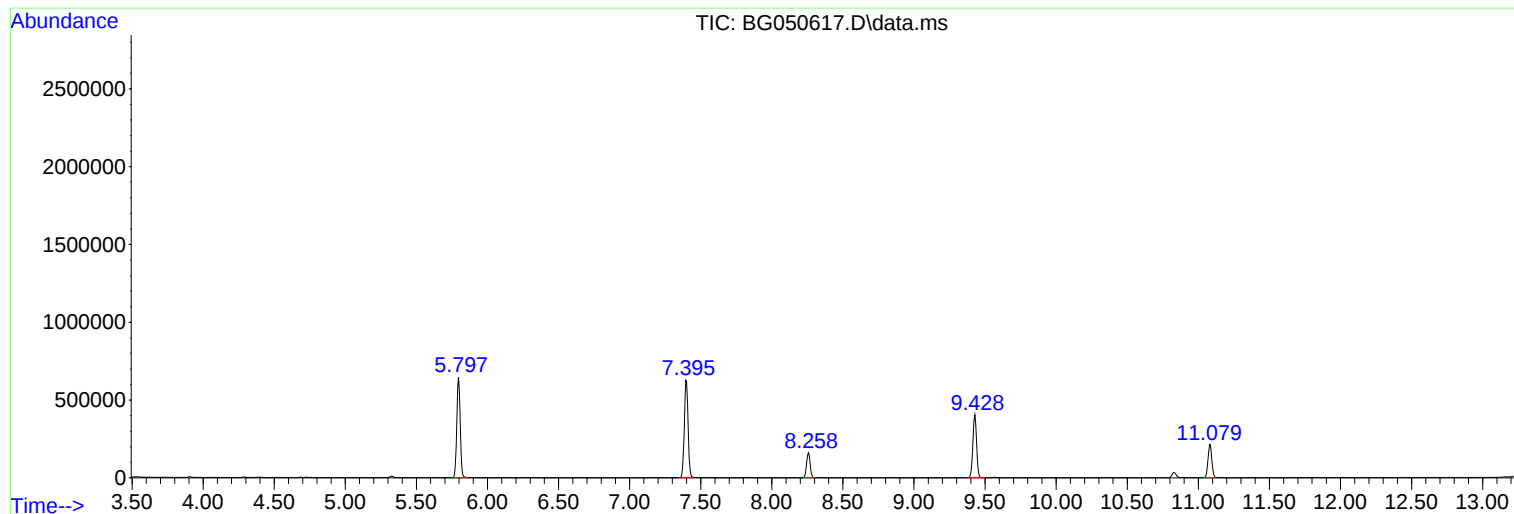
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG100621.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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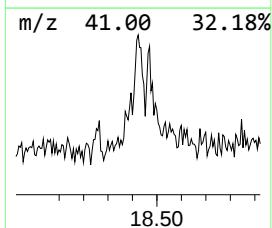
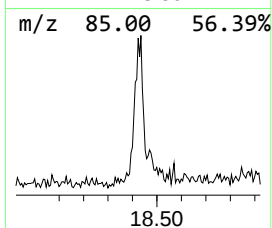
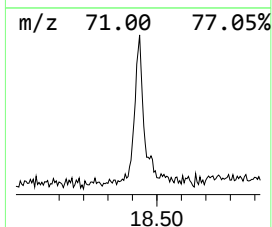
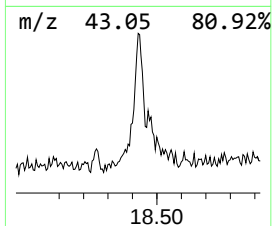
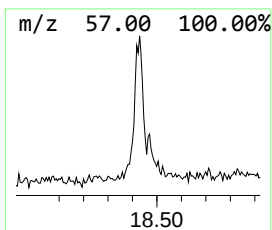
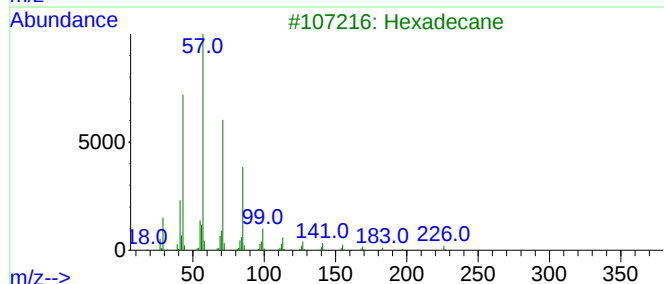
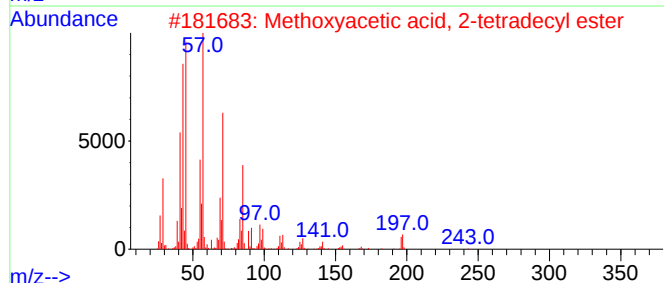
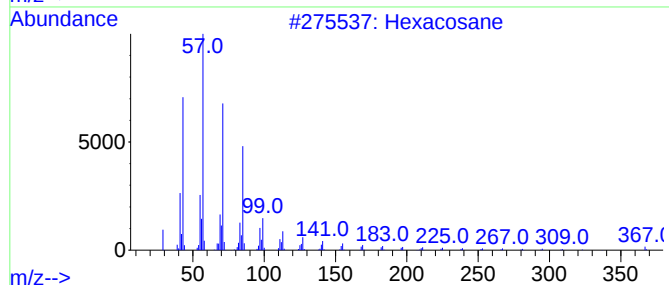
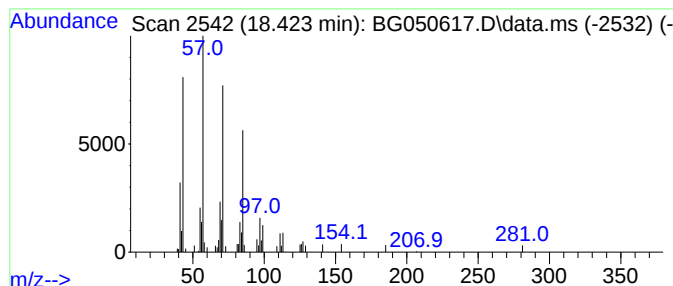
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Hexacosane Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.423	2.24 ng	73496	Phenanthrene-d10	17.618

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexacosane	366	C26H54	000630-01-3	91
2			Methoxyacetic acid, 2-tetradecyl...	286	C17H34O3	1000282-04-8	90
3			Hexadecane	226	C16H34	000544-76-3	86
4			Sulfurous acid, pentadecyl 2-pro...	334	C18H38O3S	1000309-12-6	86
5			Tetradecane	198	C14H30	000629-59-4	86



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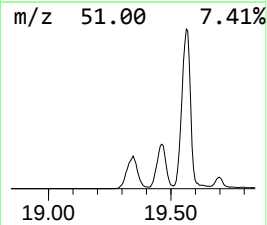
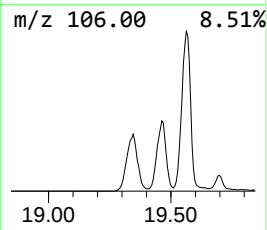
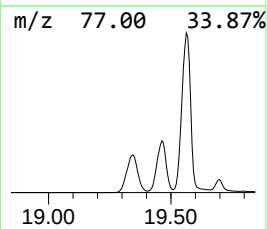
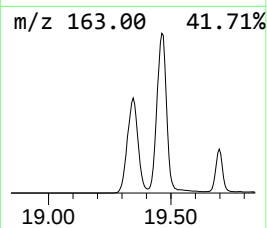
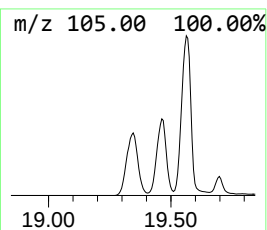
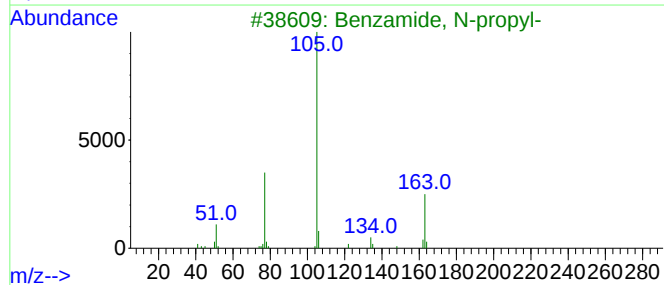
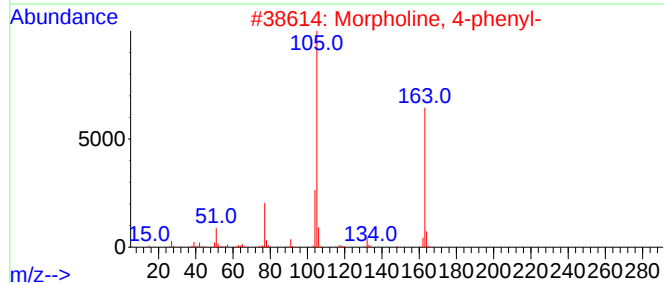
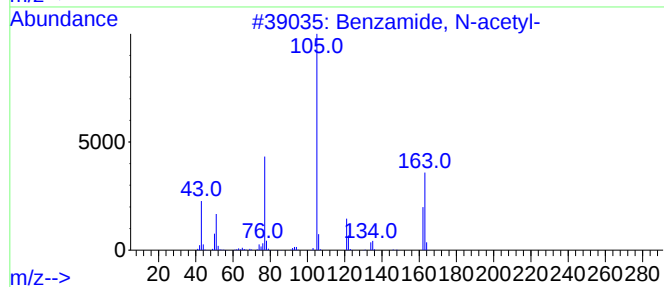
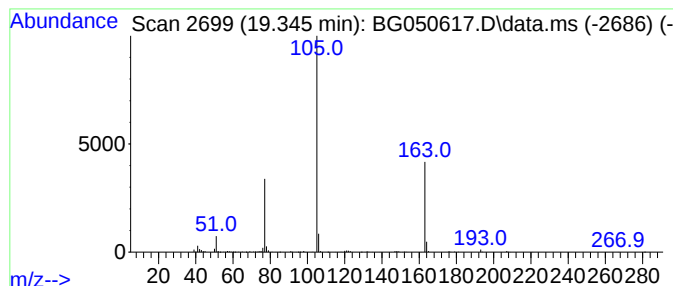
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Peak Number 2 Benzamide, N-acetyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.345	82.57 ng	2711380	Phenanthrene-d10	17.618

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzamide, N-acetyl-	163	C9H9NO2	001575-95-7	64
2			Morpholine, 4-phenyl-	163	C10H13NO	000092-53-5	56
3			Benzamide, N-propyl-	163	C10H13NO	010546-70-0	56
4			Oxybis(propane-1,2-diyl) dibenzoate	342	C20H22O5	000094-03-1	53
5			N-(5-Pyrimidinyl)benzamide	199	C11H9N3O	103854-65-5	38



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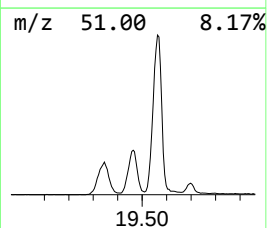
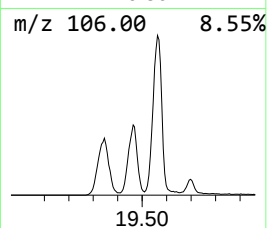
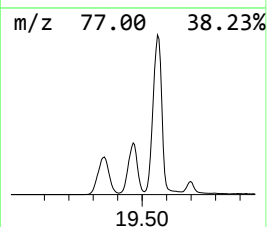
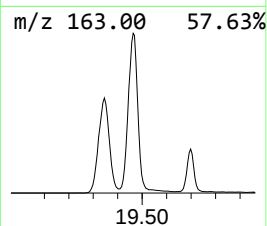
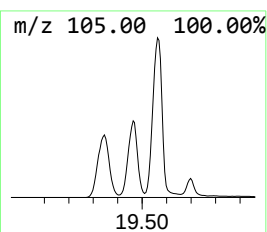
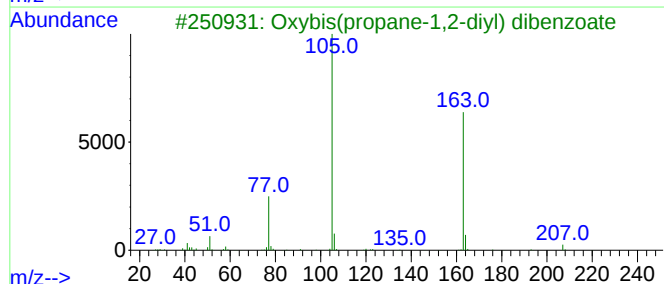
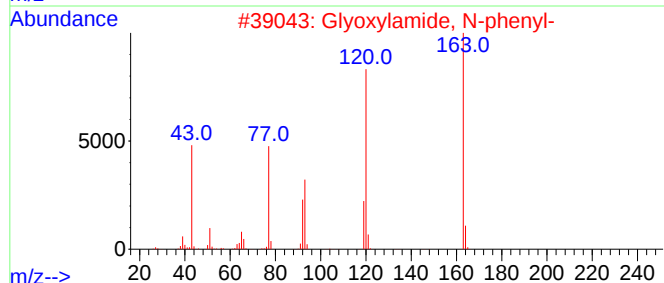
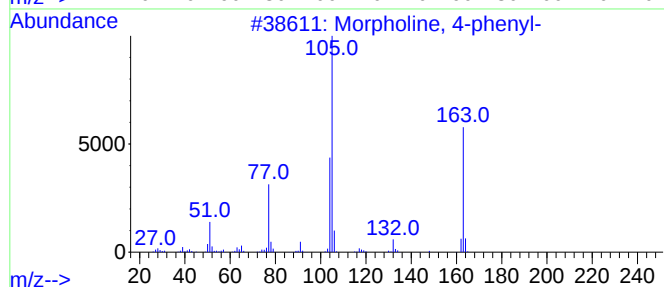
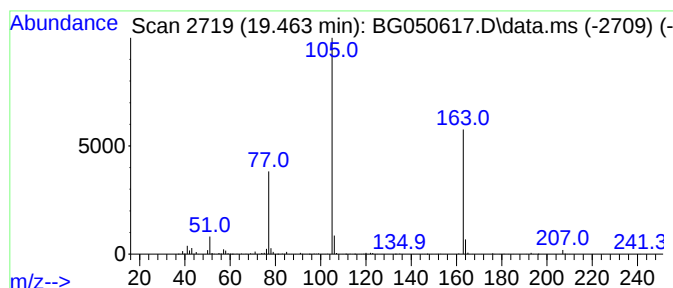
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Peak Number 3 Morpholine, 4-phenyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.463	98.71 ng	3241470	Phenanthrene-d10	17.618

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			Oxybis(propane-1,2-diyl) dibenzoate	342	C20H22O5	000094-03-1	47
4			1,3-Benzenediol, monobenzoate	214	C13H10O3	000136-36-7	43
5			N-(5-Nitrothiazol-2-yl)-benzamide	249	C10H7N3O3S	064398-84-1	43



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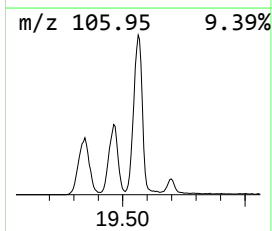
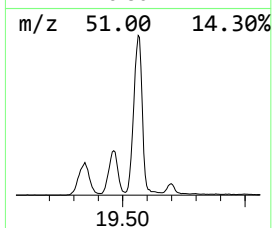
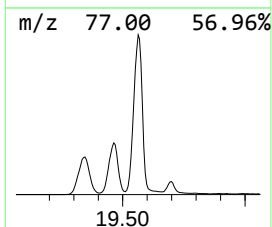
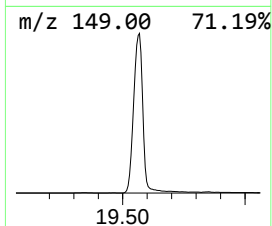
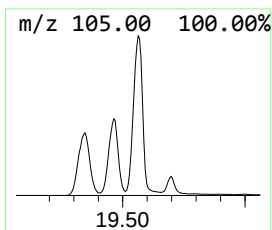
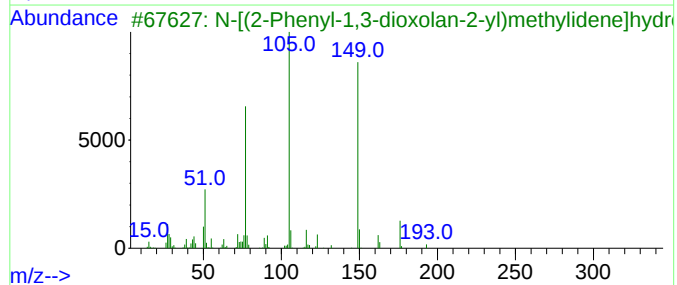
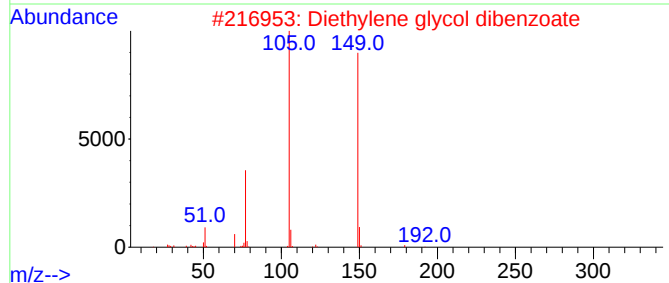
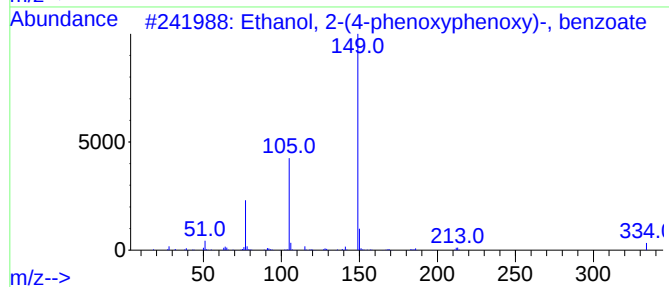
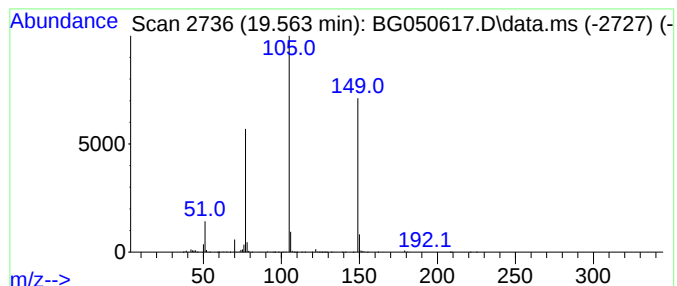
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Peak Number 4 Ethanol, 2-(4-phenoxypheno... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.563	221.66 ng	7278870	Phenanthrene-d10	17.618

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			N-[(2-Phenyl-1,3-dioxolan-2-yl)m...	193	C10H11NO3	1000411-48-9	78
4			1-Triazene, 3,3-dimethyl-1-phenyl-	149	C8H11N3	007227-91-0	64
5			1,2-Propanedione, 1-phenyl-	148	C9H8O2	000579-07-7	50



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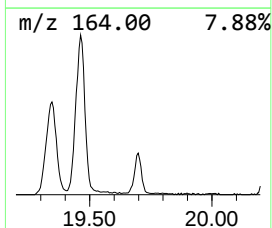
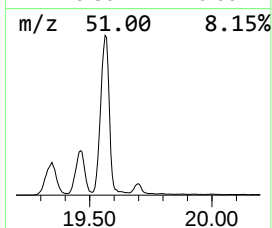
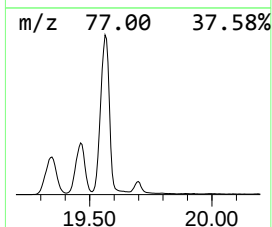
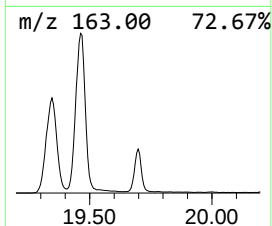
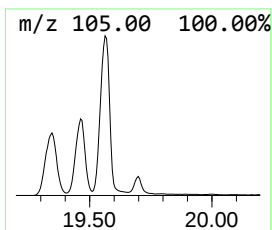
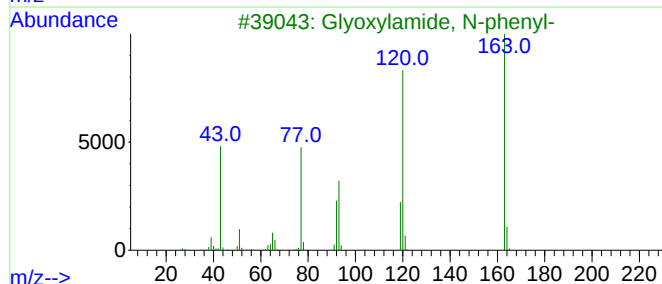
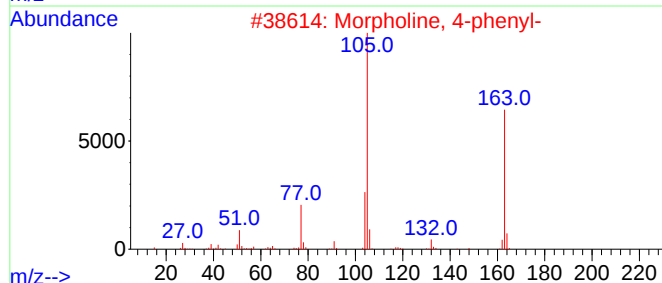
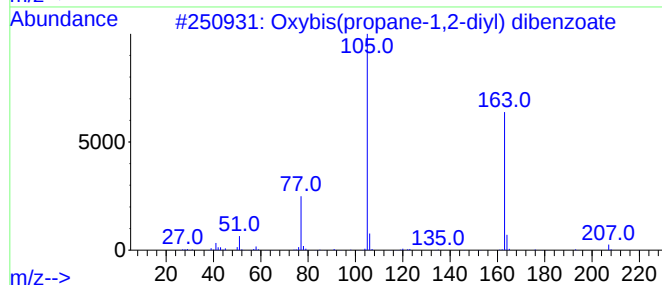
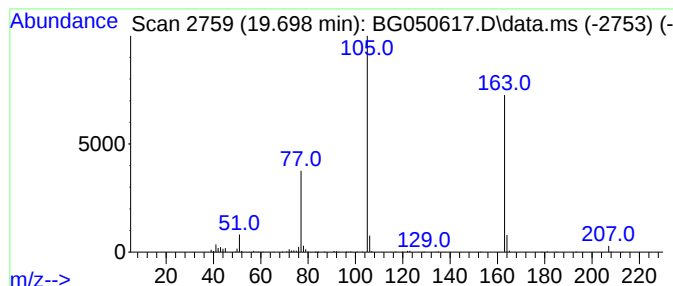
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Peak Number 5 Oxybis(propene-1,2-diyl) di... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.698	18.94 ng	621825	Phenanthrene-d10	17.618

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Oxybis(propene-1,2-diyl) dibenzoate	342	C20H22O5	000094-03-1	74
2			Morpholine, 4-phenyl-	163	C10H13NO	000092-53-5	64
3			Glyoxylamide, N-phenyl-	163	C9H9NO2	032331-52-5	50
4			.delta.2-1,3,4-Oxadiazolin-5-one...	163	C7H5N3O2	002845-82-1	43
5			Benzoic acid, phenyl ester	198	C13H10O2	000093-99-2	38



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WC-1-BD

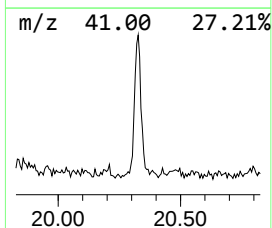
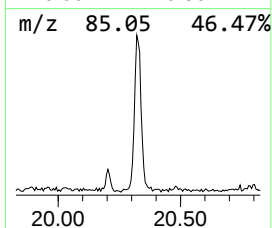
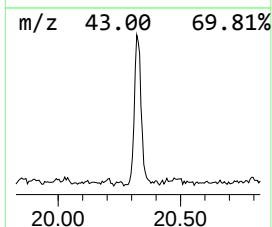
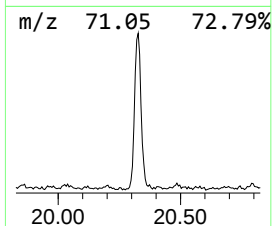
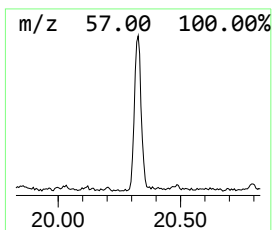
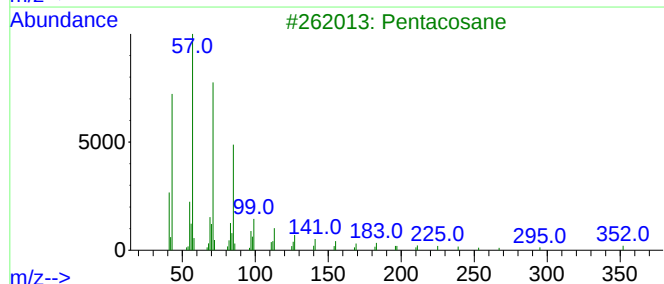
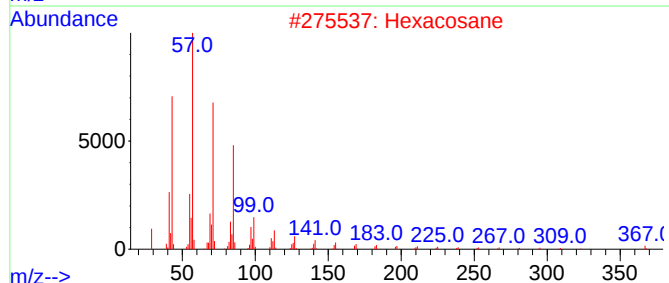
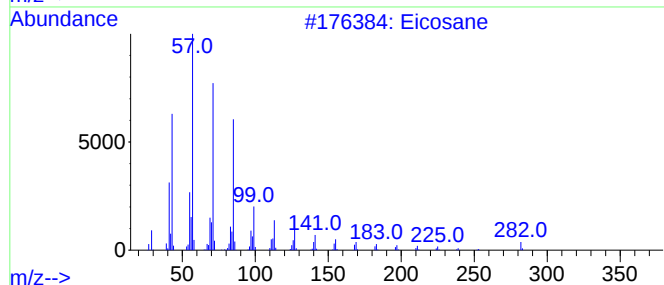
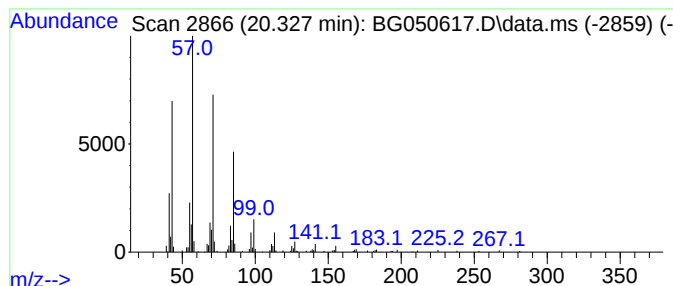
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Eicosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.327	8.24 ng	290375	Chrysene-d12	21.919

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane	282	C20H42	000112-95-8	94
2			Hexacosane	366	C26H54	000630-01-3	91
3			Pentacosane	352	C25H52	000629-99-2	91
4			Heneicosane	296	C21H44	000629-94-7	91
5			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG101821\
Data File : BG050617.D
Acq On : 18 Oct 2021 14:45
Operator : CG/JU
Sample : M4259-07
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
WC-1-BD

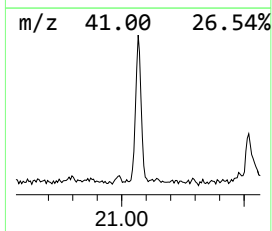
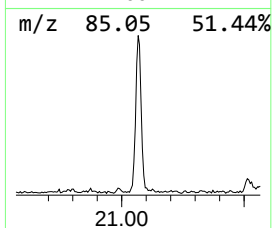
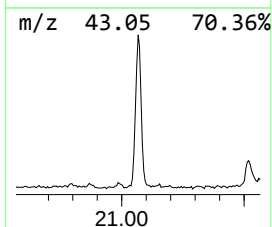
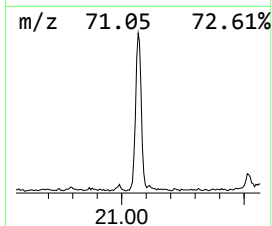
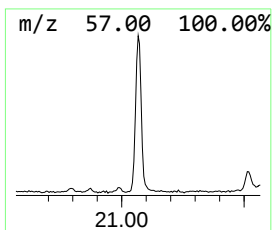
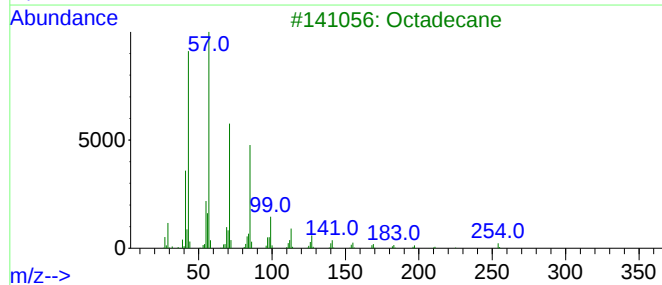
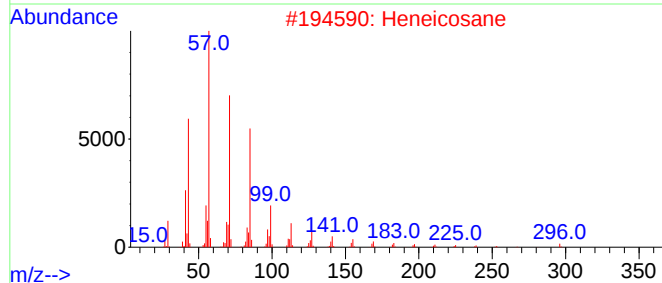
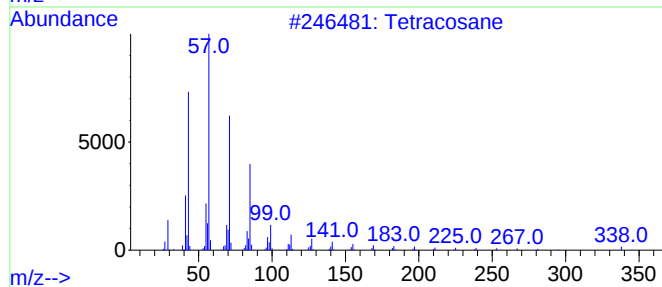
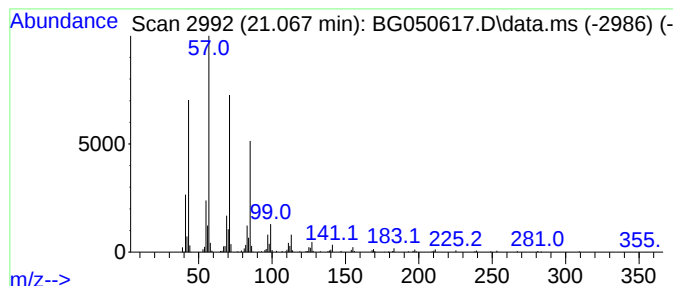
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Tetracosane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.067	11.19 ng	394374	Chrysene-d12	21.919

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetracosane	338	C24H50	000646-31-1	91
2			Heneicosane	296	C21H44	000629-94-7	91
3			Octadecane	254	C18H38	000593-45-3	91
4			Heptacosane	380	C27H56	000593-49-7	87
5			Pentadecane	212	C15H32	000629-62-9	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG101821\
Data File : BG050617.D
Acq On : 18 Oct 2021 14:45
Operator : CG/JU
Sample : M4259-07
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
WC-1-BD

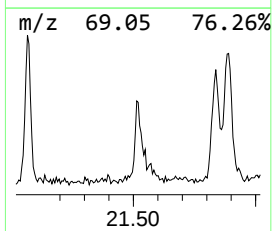
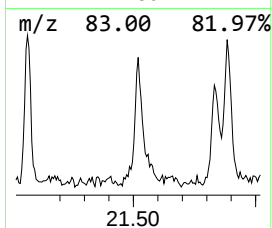
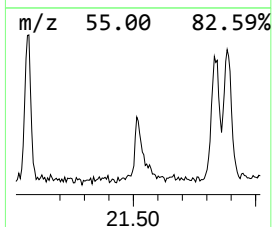
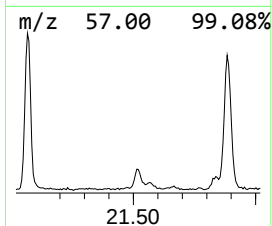
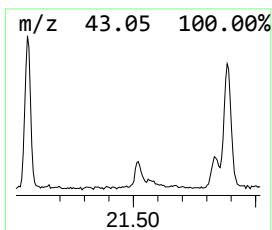
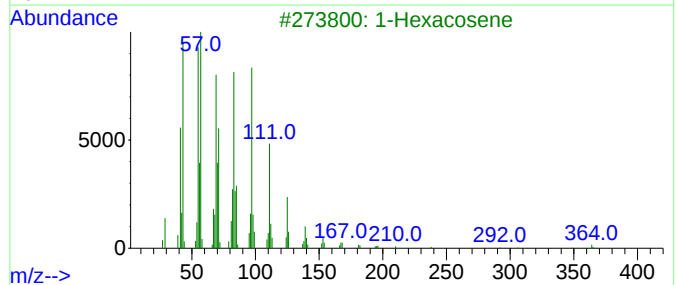
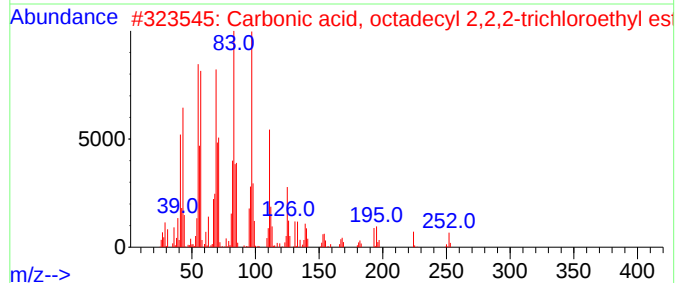
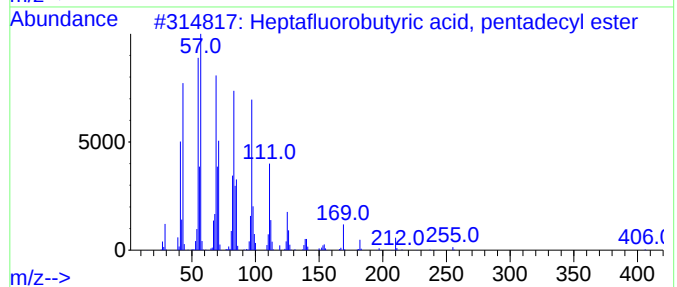
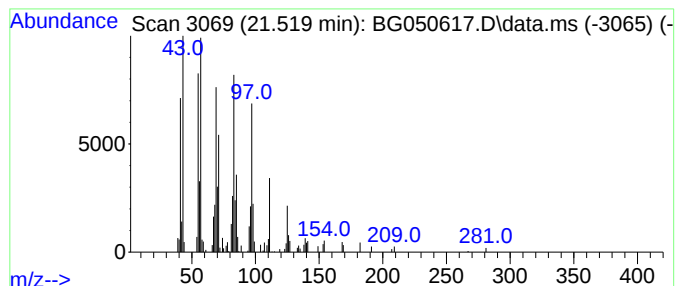
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Heptafluorobutyric acid, pe... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.519	2.34 ng	82399	Chrysene-d12	21.919

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	94
2			Carbonic acid, octadecyl 2,2,2-t...	444	C21H39Cl3O3	1000314-56-3	91
3			1-Hexacosene	364	C26H52	018835-33-1	91
4			Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	91
5			Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG101821\
Data File : BG050617.D
Acq On : 18 Oct 2021 14:45
Operator : CG/JU
Sample : M4259-07
Misc :
ALS Vial : 8 Sample Multiplier: 1

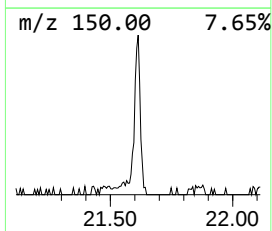
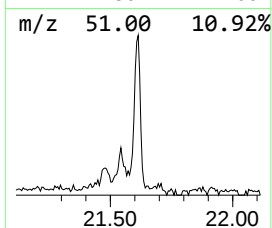
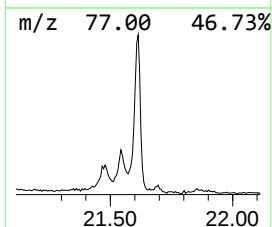
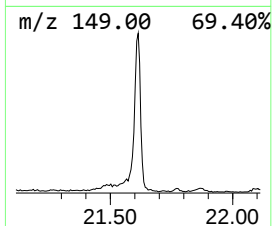
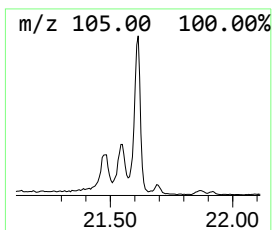
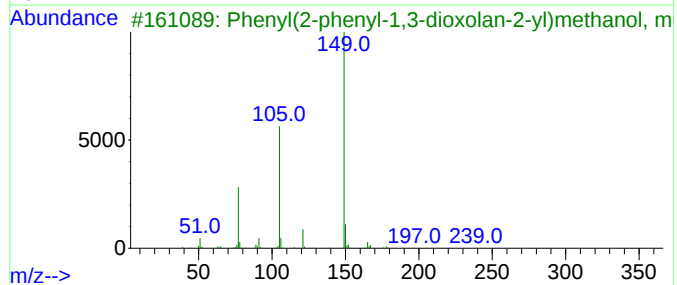
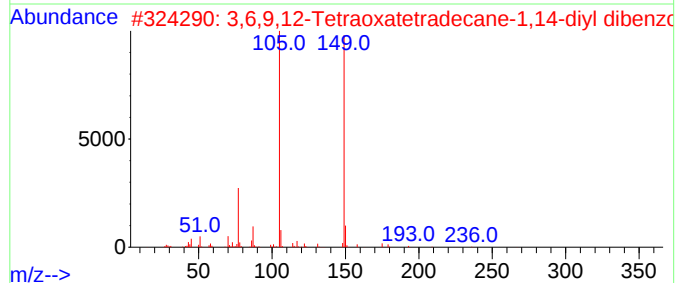
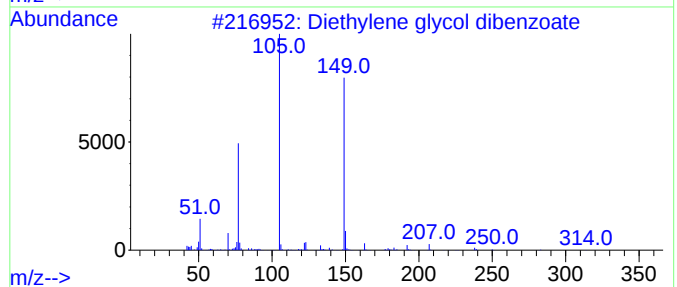
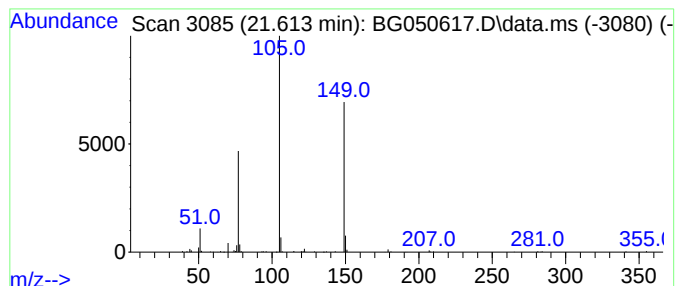
Instrument :
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ClientSampleId :
WC-1-BD

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 Diethylene glycol dibenzoate Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.613	8.03 ng	283040	Chrysene-d12	21.919	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Diethylene glycol dibenzoate	314	C18H18O5	000120-55-8	78
2	3,6,9,12-Tetraoxatetradecane-1,1...	446	C24H30O8	1000366-97-0	78
3	Phenyl(2-phenyl-1,3-dioxolan-2-y...	270	C17H18O3	1000507-07-6	72
4	N-[(2-Phenyl-1,3-dioxolan-2-yl)m...	193	C10H11NO3	1000411-48-9	64
5	2-(2-(2-Methoxyethoxy)ethoxy)eth...	268	C14H20O5	1000366-99-1	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG101821\
Data File : BG050617.D
Acq On : 18 Oct 2021 14:45
Operator : CG/JU
Sample : M4259-07
Misc :
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Instrument :
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ClientSampleId :
WC-1-BD

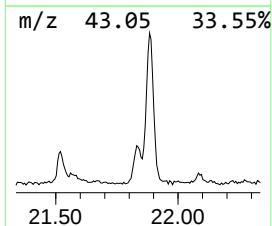
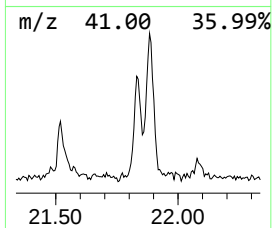
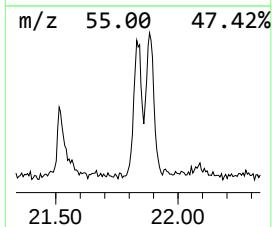
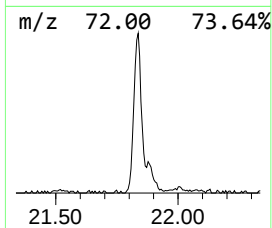
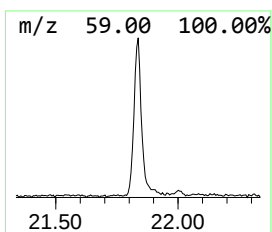
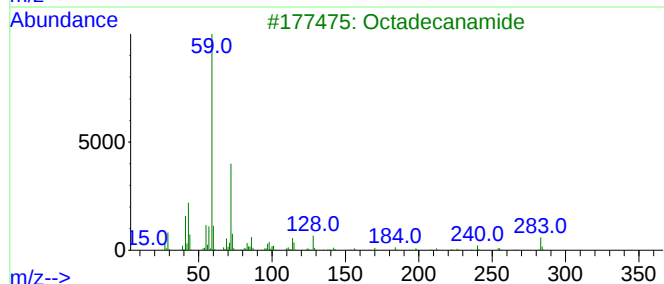
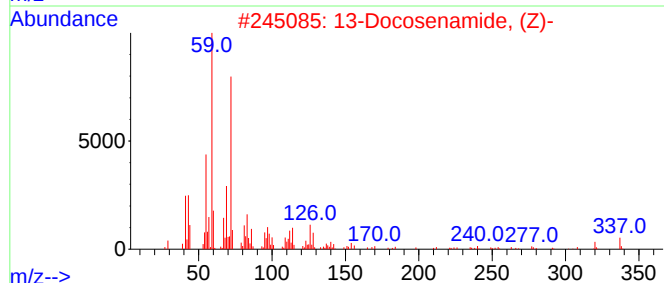
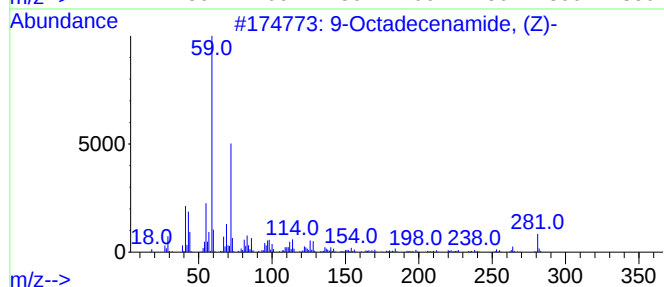
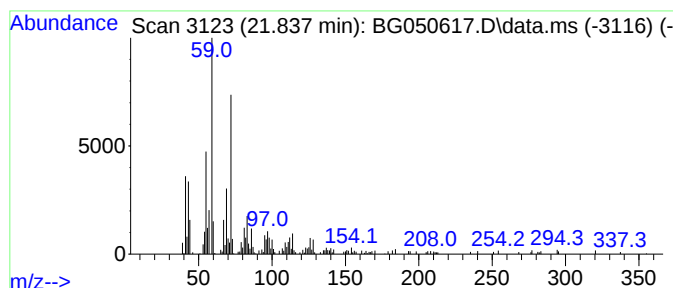
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 9-Octadecenamide, (Z)- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.837	8.61 ng	303383	Chrysene-d12	21.919

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	91
2		13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	91
3		Octadecanamide	283	C18H37NO	000124-26-5	47
4		trans-2,7-Dimethyl-4,6-octadien-...	154	C10H18O	1010281-69-5	38
5		N-Ethyl-3-methyl-2-butanamine, N...	157	C9H19NO	1849351-93-4	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG101821\
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Acq On : 18 Oct 2021 14:45
Operator : CG/JU
Sample : M4259-07
Misc :
ALS Vial : 8 Sample Multiplier: 1

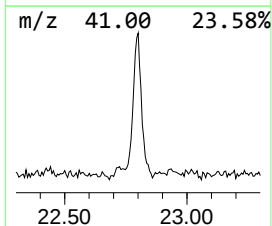
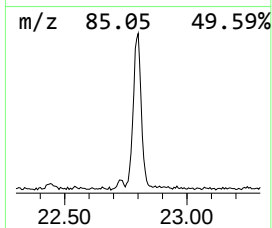
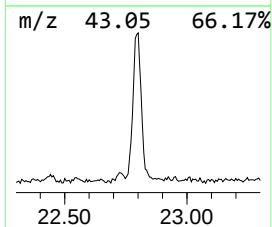
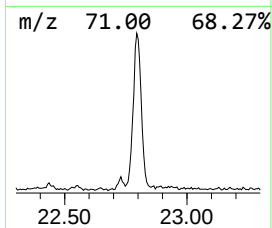
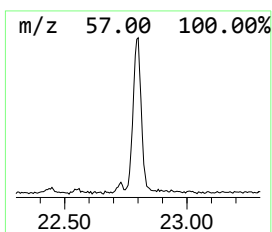
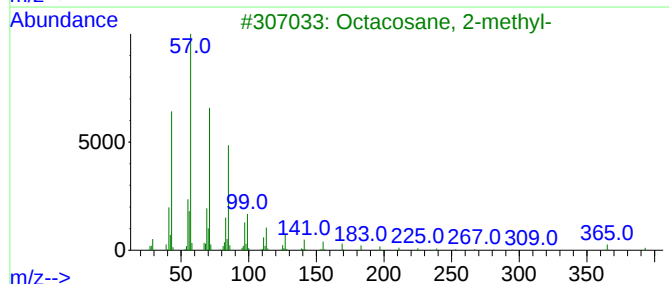
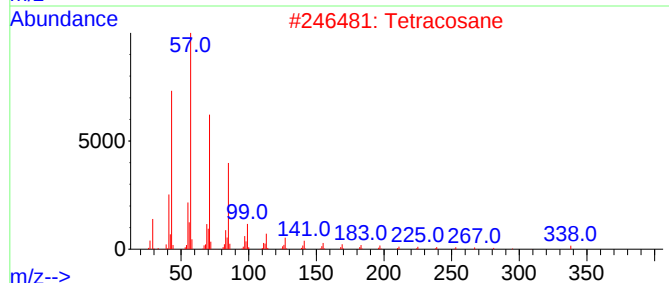
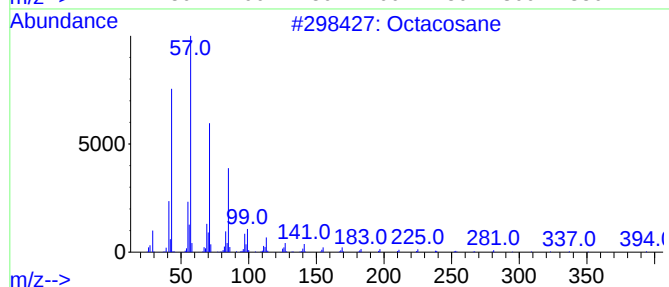
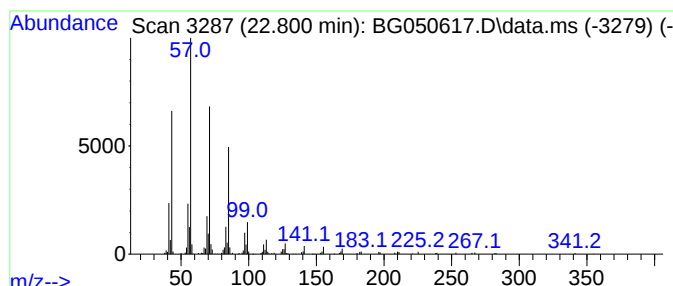
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Octacosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
22.800	12.58 ng	443322	Chrysene-d12	21.919	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octacosane	394	C28H58	000630-02-4	91
2	Tetracosane	338	C24H50	000646-31-1	91
3	Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
4	Eicosane	282	C20H42	000112-95-8	90
5	Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG101821\
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Acq On : 18 Oct 2021 14:45
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Sample : M4259-07
Misc :
ALS Vial : 8 Sample Multiplier: 1

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ClientSampleId :
WC-1-BD

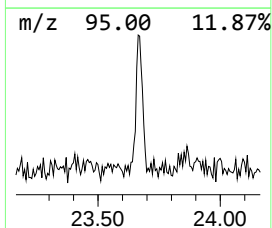
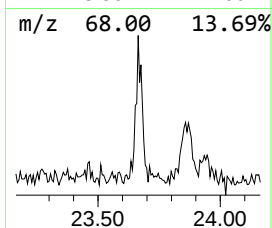
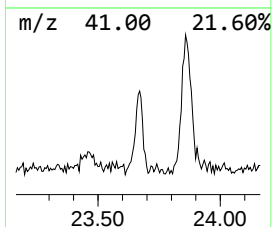
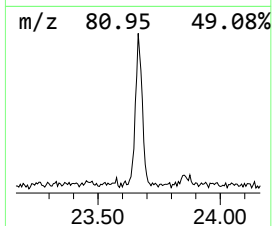
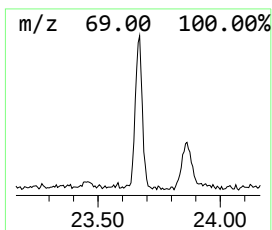
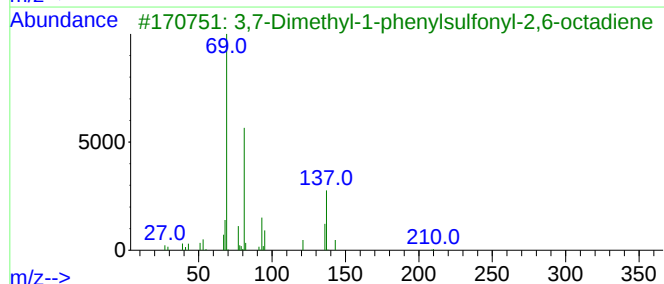
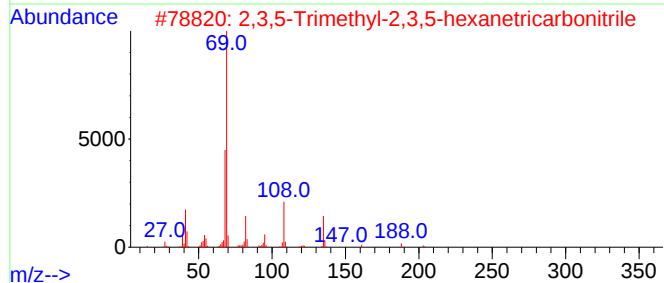
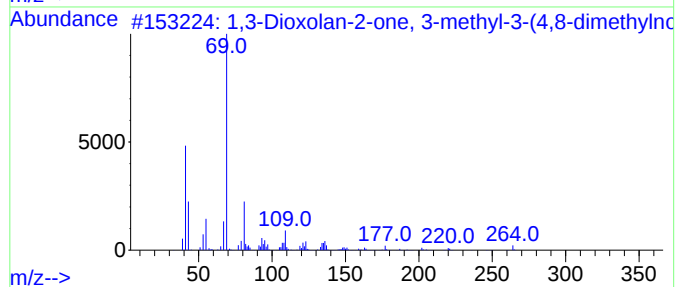
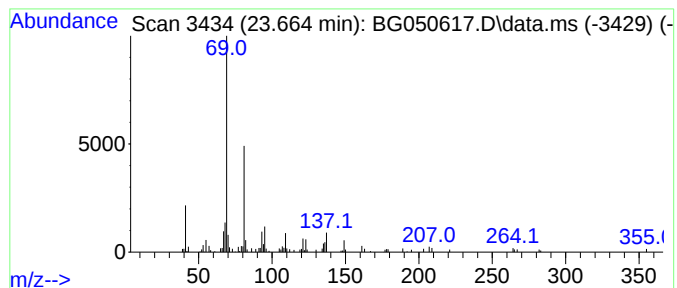
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 1,3-Dioxolan-2-one, 3-methy... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.664	2.87 ng	107953	Perylene-d12	25.333

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,3-Dioxolan-2-one, 3-methyl-3-(...	264	C16H24O3	336879-76-6	72
2			2,3,5-Trimethyl-2,3,5-hexanetric...	203	C12H17N3	003974-79-6	64
3			3,7-Dimethyl-1-phenylsulfonyl-2,...	278	C16H22O2S	1000432-27-5	64
4			3,7,11-Trimethyl-2,6,10-dodecatr...	217	C15H23N	029789-67-1	64
5			Geranyl bromide	216	C10H17Br	006138-90-5	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG101821\
Data File : BG050617.D
Acq On : 18 Oct 2021 14:45
Operator : CG/JU
Sample : M4259-07
Misc :
ALS Vial : 8 Sample Multiplier: 1

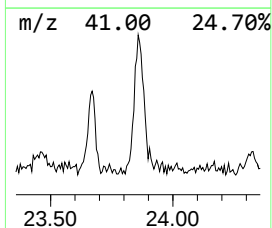
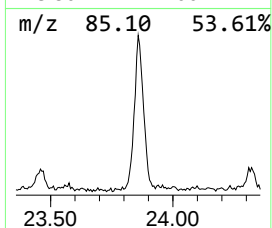
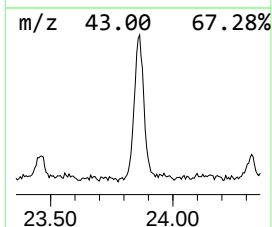
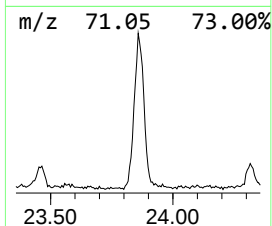
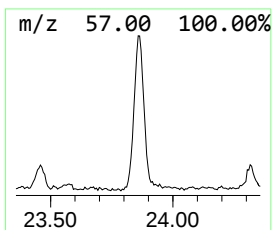
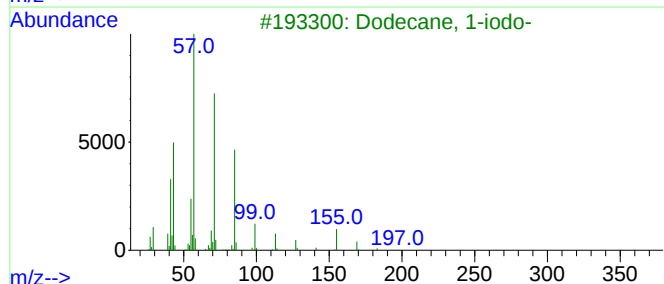
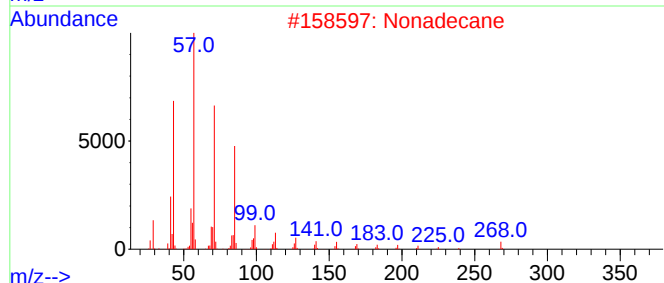
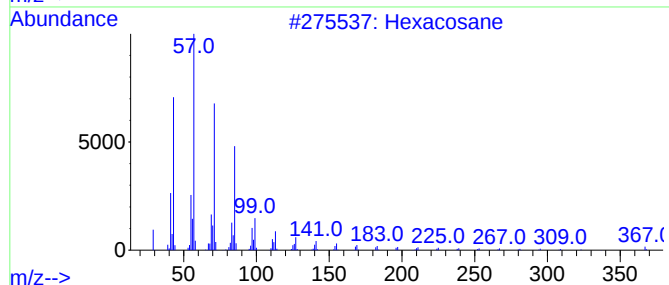
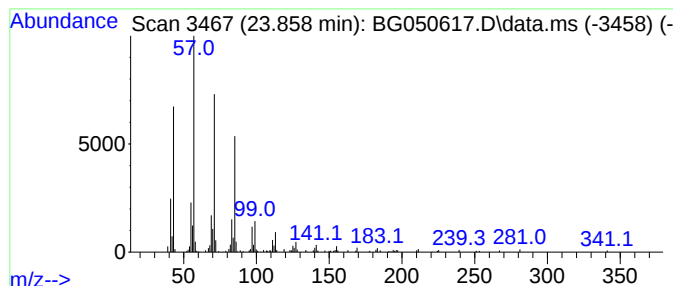
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ClientSampleId :
WC-1-BD

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 Nonadecane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
23.858	10.06 ng	378069	Perylene-d12	25.333	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexacosane	366	C26H54	000630-01-3	87
2	Nonadecane	268	C19H40	000629-92-5	86
3	Dodecane, 1-iodo-	296	C12H25I	004292-19-7	83
4	Octacosane	394	C28H58	000630-02-4	81
5	Tetratetracontane	619	C44H90	007098-22-8	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG101821\
Data File : BG050617.D
Acq On : 18 Oct 2021 14:45
Operator : CG/JU
Sample : M4259-07
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
WC-1-BD

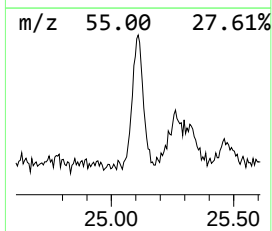
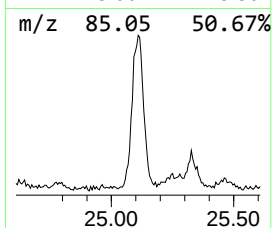
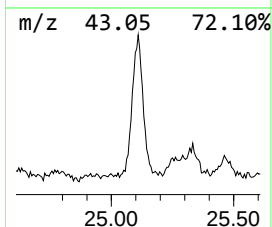
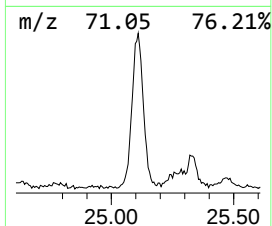
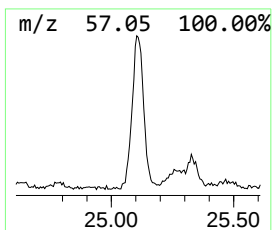
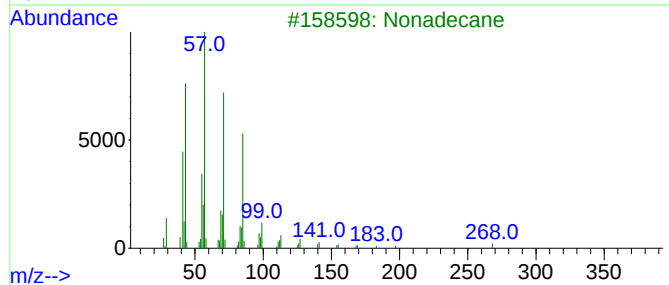
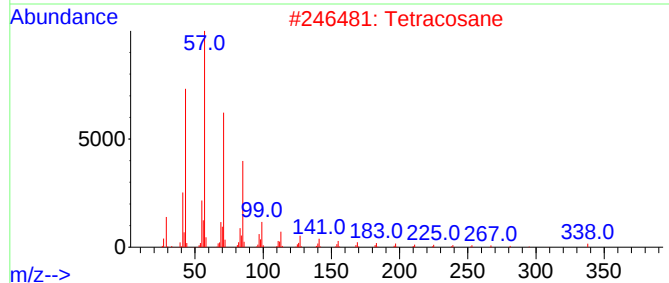
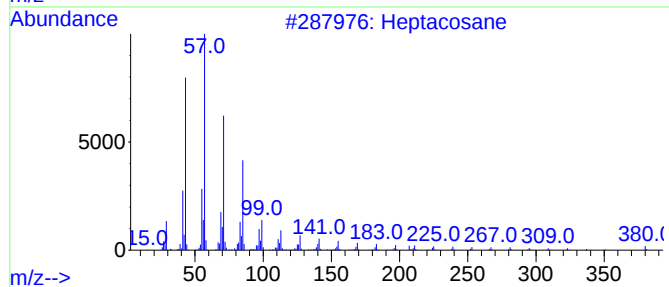
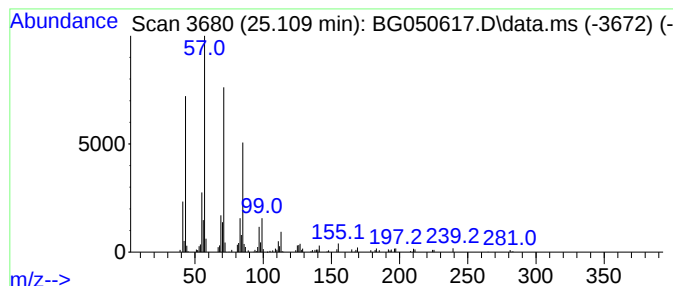
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Heptacosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.109	8.81 ng	331294	Perylene-d12	25.333

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptacosane	380	C27H56	000593-49-7	91
2		Tetracosane	338	C24H50	000646-31-1	91
3		Nonadecane	268	C19H40	000629-92-5	90
4		5,5,7,7-Tetraethylundecane	268	C19H40	1000360-42-0	90
5		Octacosane, 2-methyl-	408	C29H60	001560-98-1	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG101821\
Data File : BG050617.D
Acq On : 18 Oct 2021 14:45
Operator : CG/JU
Sample : M4259-07
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
WC-1-BD

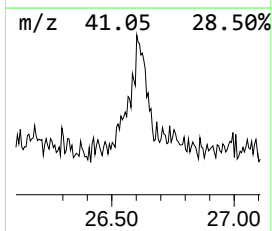
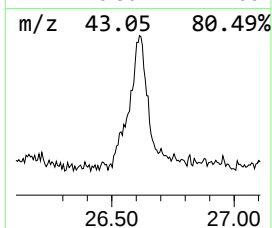
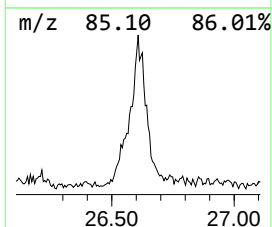
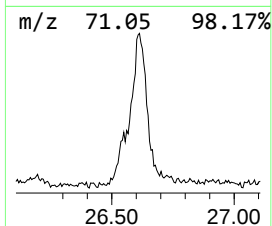
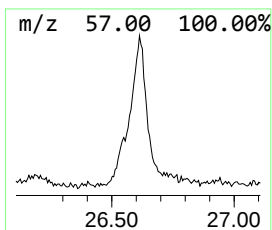
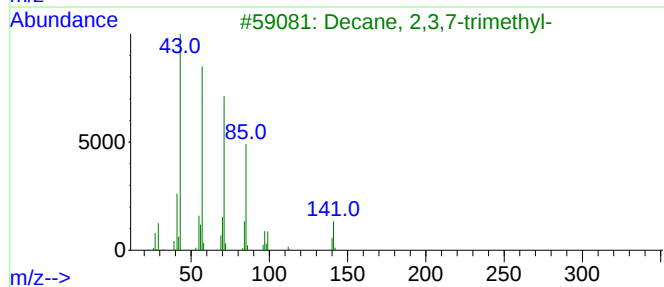
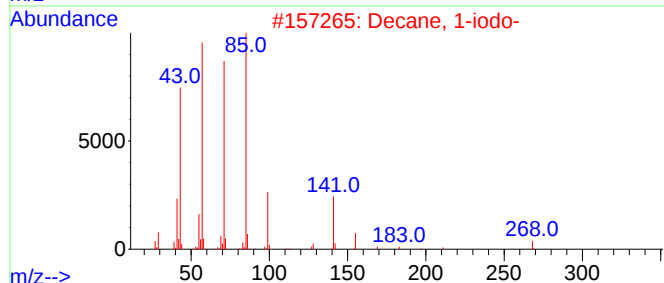
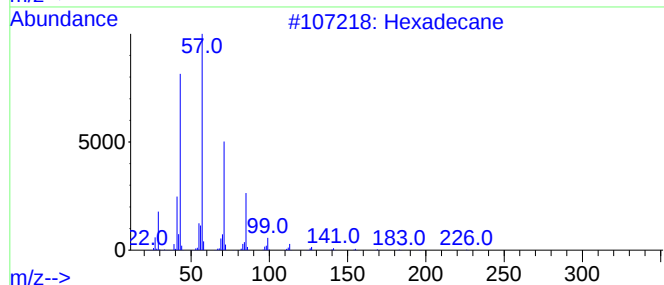
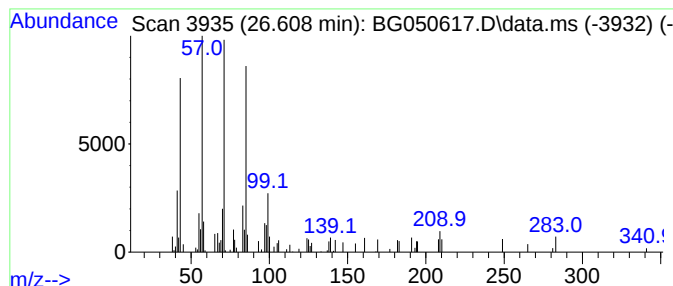
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Hexadecane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.608	4.33 ng	162794	Perylene-d12	25.333

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane	226	C16H34	000544-76-3	64
2			Decane, 1-iodo-	268	C10H21I	002050-77-3	59
3			Decane, 2,3,7-trimethyl-	184	C13H28	062238-13-5	59
4			Sulfurous acid, 2-ethylhexyl hex...	278	C14H30O3S	1000309-20-2	53
5			1-Iodoundecane	282	C11H23I	004282-44-4	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG101821\
 Data File : BG050617.D
 Acq On : 18 Oct 2021 14:45
 Operator : CG/JU
 Sample : M4259-07
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 WC-1-BD

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
Hexacosane	18.423	2.2	ng	73496	4	17.618	656756	20.0
Benzamide, N-ac...	19.345	82.6	ng	2711380	4	17.618	656756	20.0
Morpholine, 4-p...	19.463	98.7	ng	3241470	4	17.618	656756	20.0
Ethanol, 2-(4-p...	19.563	221.7	ng	7278870	4	17.618	656756	20.0
Oxybis(propane-...	19.698	18.9	ng	621825	4	17.618	656756	20.0
Eicosane	20.327	8.2	ng	290375	5	21.919	704884	20.0
Tetracosane	21.067	11.2	ng	394374	5	21.919	704884	20.0
Heptafluorobuty...	21.519	2.3	ng	82399	5	21.919	704884	20.0
Diethylene glyc...	21.613	8.0	ng	283040	5	21.919	704884	20.0
9-Octadecenamid...	21.837	8.6	ng	303383	5	21.919	704884	20.0
Octacosane	22.800	12.6	ng	443322	5	21.919	704884	20.0
1,3-Dioxolan-2-...	23.664	2.9	ng	107953	6	25.333	751843	20.0
Nonadecane	23.858	10.1	ng	378069	6	25.333	751843	20.0
Heptacosane	25.109	8.8	ng	331294	6	25.333	751843	20.0
Hexadecane	26.608	4.3	ng	162794	6	25.333	751843	20.0