

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081721\
 Data File : BG049874.D
 Acq On : 17 Aug 2021 12:24
 Operator : CG/JU
 Sample : M3412-02
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PG-2

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-BG080321.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BG049874.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.132	10	16	30	rVB	145467	265642	21.21%	2.681%
2	5.100	343	351	357	rBV2	20870	33821	2.70%	0.341%
3	5.582	426	433	442	rBV	450342	739599	59.06%	7.465%
4	7.198	700	708	720	rBV	474141	853806	68.18%	8.618%
5	8.026	842	849	856	rBV	106132	185272	14.79%	1.870%
6	9.195	1039	1048	1058	rBV	315878	574648	45.89%	5.800%
7	10.611	1284	1289	1296	rBV2	24497	43763	3.49%	0.442%
8	10.846	1322	1329	1338	rVB2	141806	256198	20.46%	2.586%
9	13.295	1738	1746	1754	rBV	683937	1090163	87.05%	11.003%
10	14.682	1976	1982	1988	rVB	198151	326789	26.09%	3.298%
11	16.174	2230	2236	2245	rVB	678456	1035079	82.65%	10.447%
12	17.431	2445	2450	2459	rVB	240271	376249	30.04%	3.798%
13	18.083	2559	2561	2568	rVB8	11678	16080	1.28%	0.162%
14	18.312	2596	2600	2606	rBV5	39014	54786	4.37%	0.553%
15	18.923	2696	2704	2705	rBV4	58395	126879	10.13%	1.281%
16	20.039	2889	2894	2900	rBV	950725	1252357	100.00%	12.640%
17	20.121	2900	2908	2918	rVB2	127348	303107	24.20%	3.059%
18	20.344	2940	2946	2954	rVB7	36531	94262	7.53%	0.951%
19	20.585	2983	2987	2991	rBV5	34636	48934	3.91%	0.494%
20	21.038	3056	3064	3076	rVB2	149377	370194	29.56%	3.736%
21	21.343	3112	3116	3123	rVB5	55996	101660	8.12%	1.026%
22	21.425	3126	3130	3134	rVB	42391	58061	4.64%	0.586%
23	21.713	3174	3179	3185	rVB2	236678	395068	31.55%	3.988%
24	22.019	3225	3231	3242	rVB2	116251	280370	22.39%	2.830%
25	22.500	3309	3313	3318	rBV4	30086	53157	4.24%	0.537%
26	23.182	3423	3429	3439	rVB4	113251	333989	26.67%	3.371%
27	24.586	3661	3668	3675	rVB3	71115	218157	17.42%	2.202%
28	24.985	3728	3736	3745	rVB	146729	419450	33.49%	4.234%

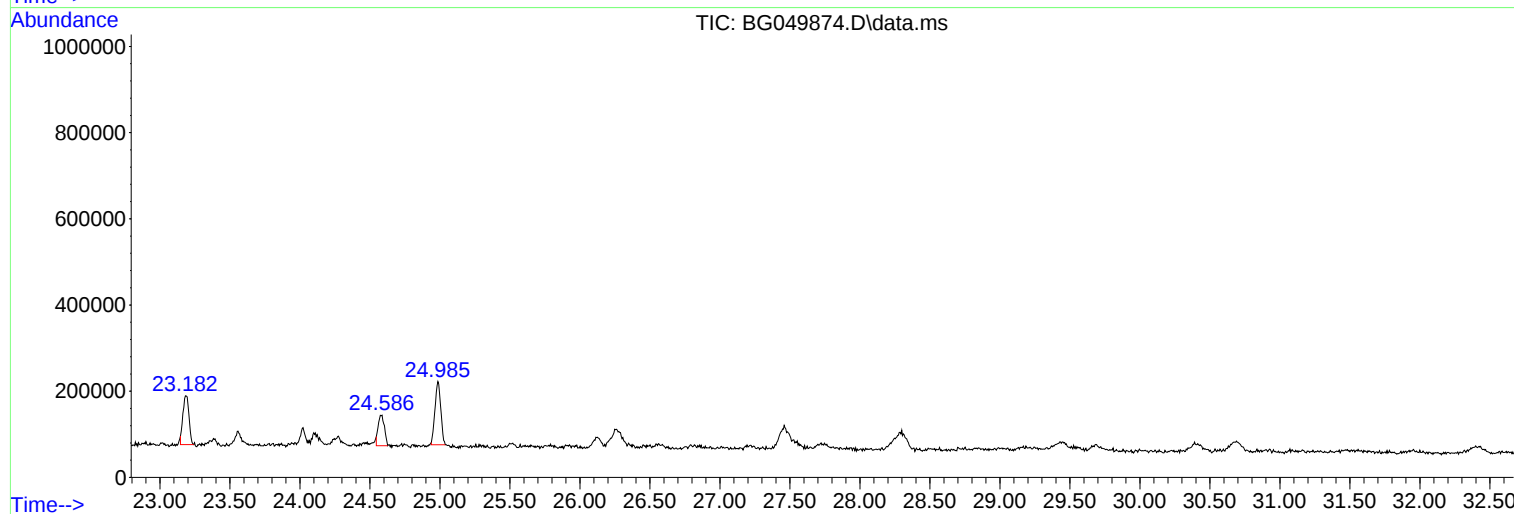
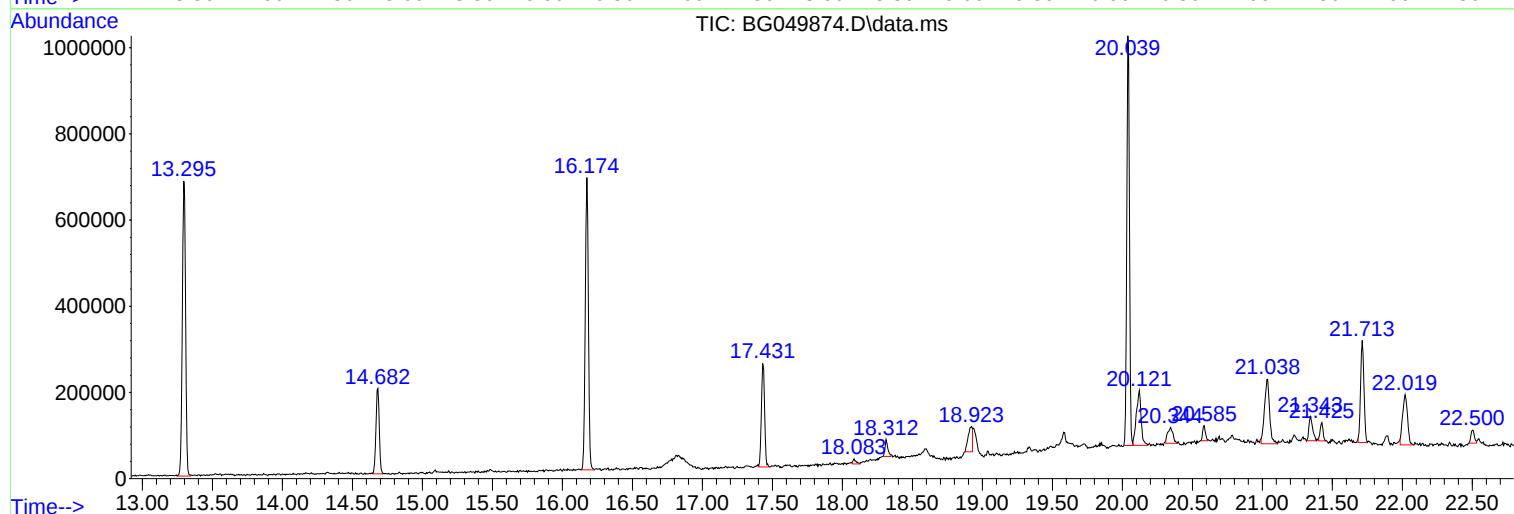
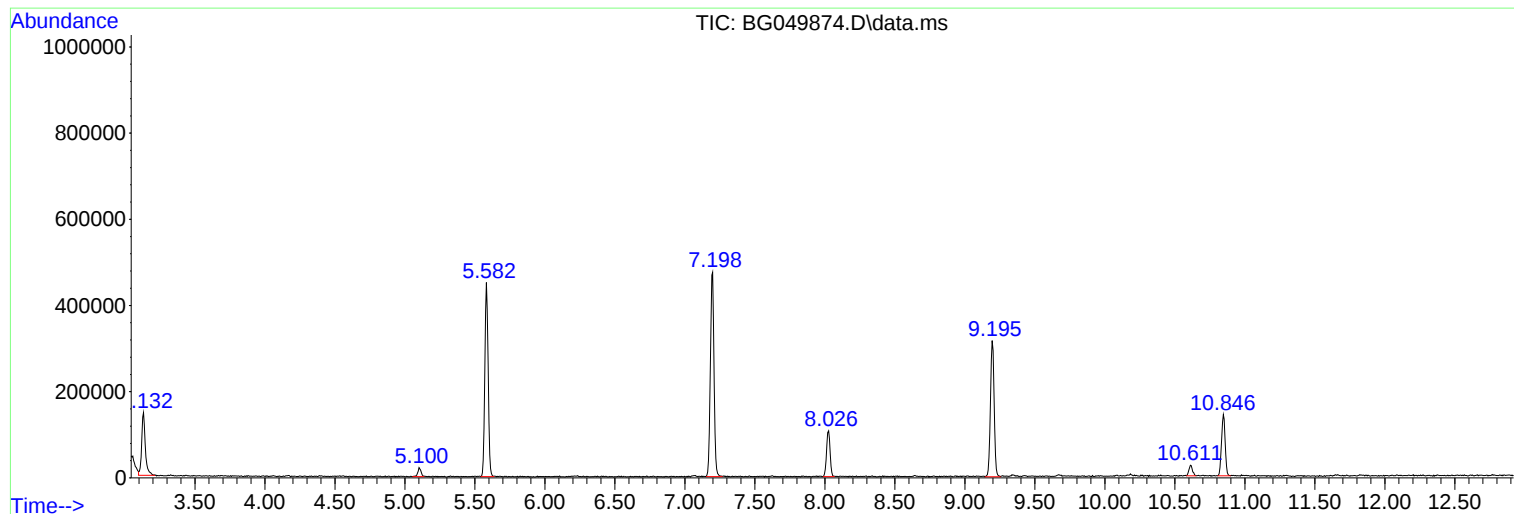
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG080321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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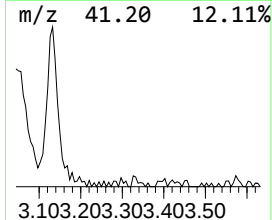
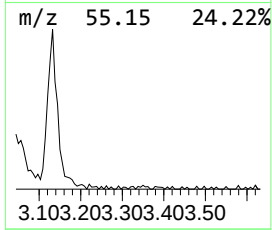
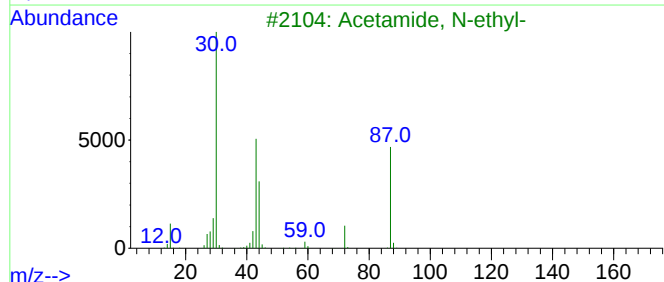
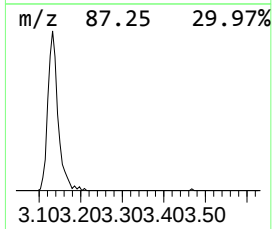
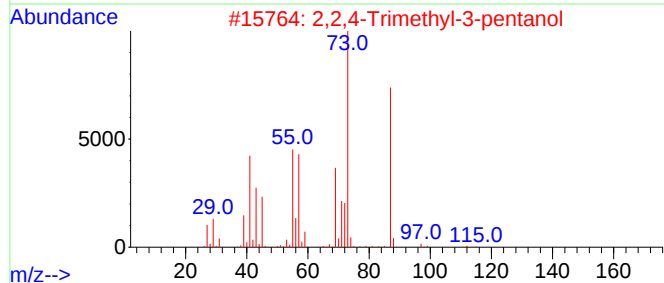
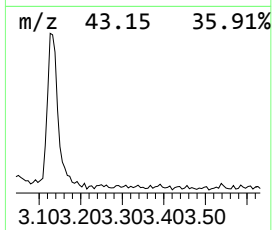
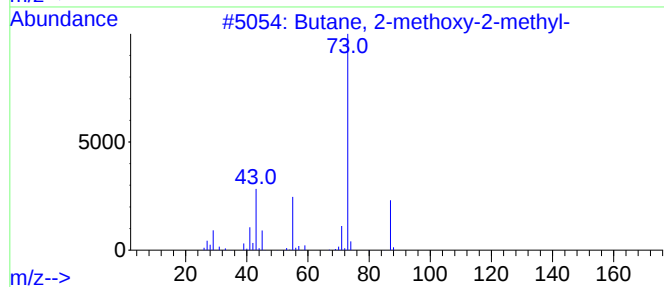
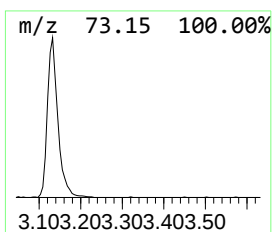
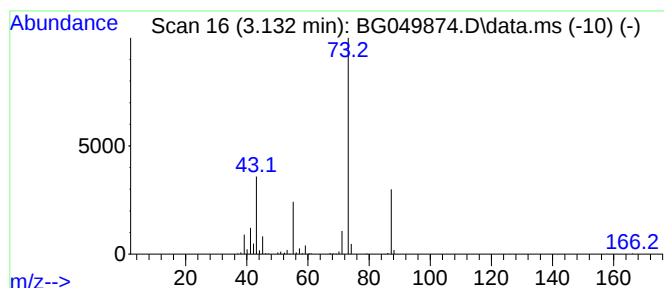
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.132	28.68 ng	265642	1,4-Dichlorobenzene-d4	8.026

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
3			Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	27
4			1,3-Dioxolane, 2-propyl-	116	C6H12O2	003390-13-4	23
5			1-Propene-1-thiol	74	C3H6S	000925-89-3	9



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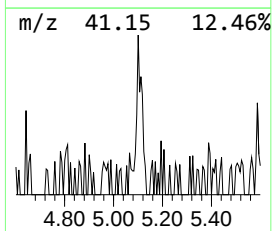
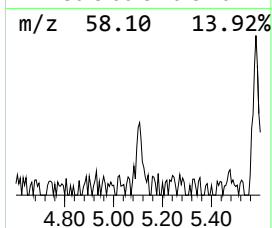
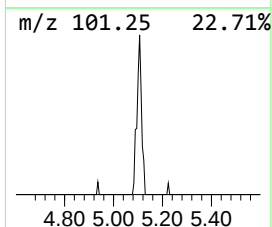
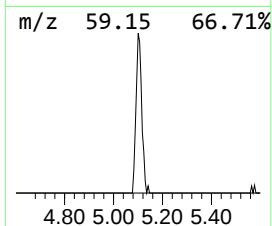
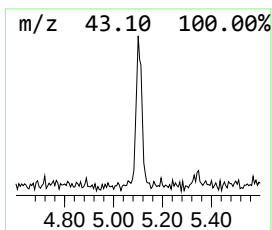
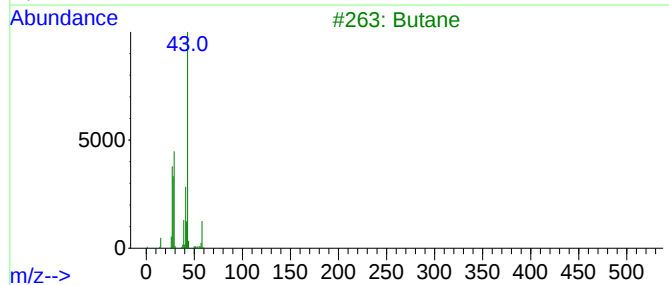
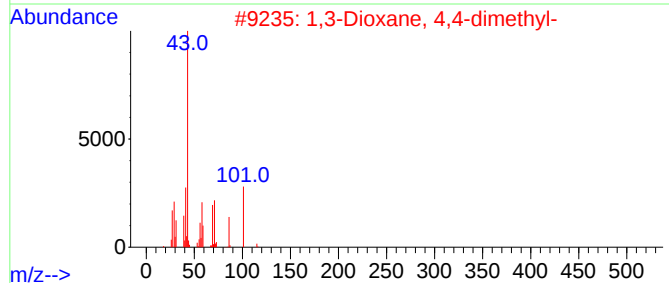
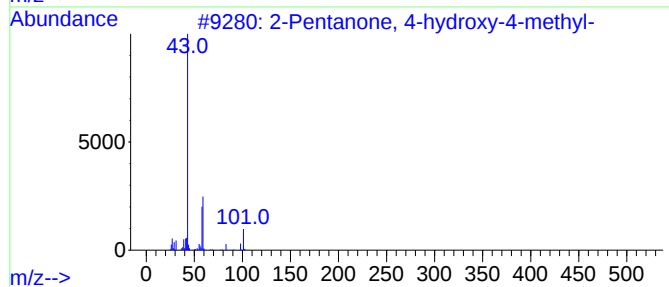
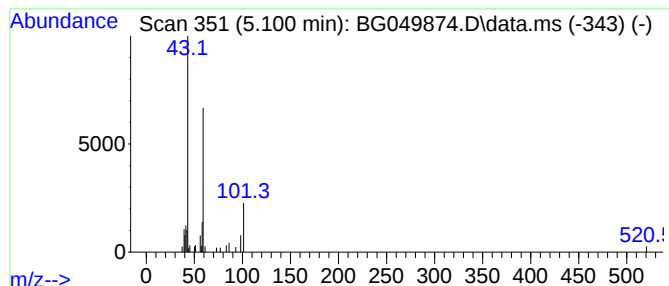
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TIC Library : C:\Database\NIST20.L
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.100	3.65 ng	33821	1,4-Dichlorobenzene-d4	8.026

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			1,3-Dioxane, 4,4-dimethyl-	116	C6H12O2	000766-15-4	17
3			Butane	58	C4H10	000106-97-8	9
4			4-Penten-2-one, 3-methyl-	98	C6H10O	000758-87-2	9
5			Methyl 1-methylcyclopropyl ketone	98	C6H10O	001567-75-5	9



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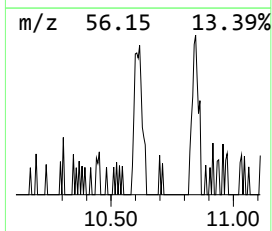
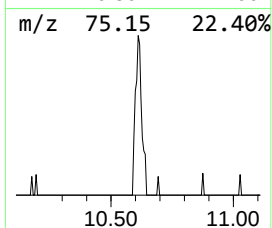
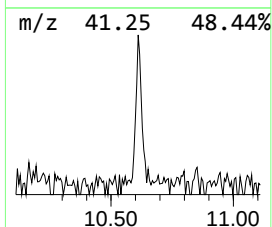
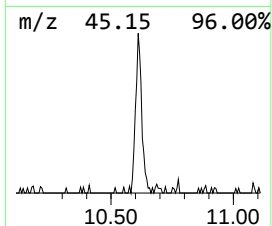
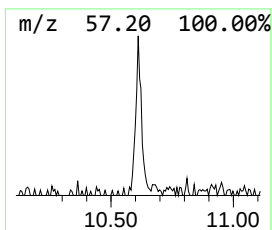
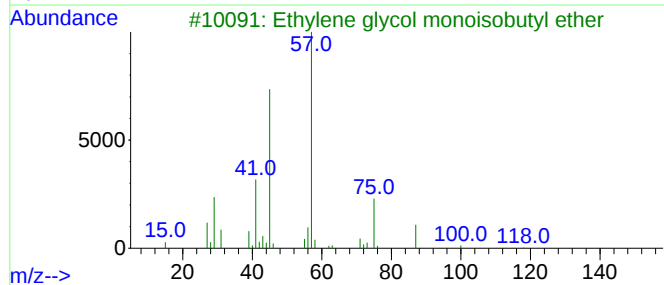
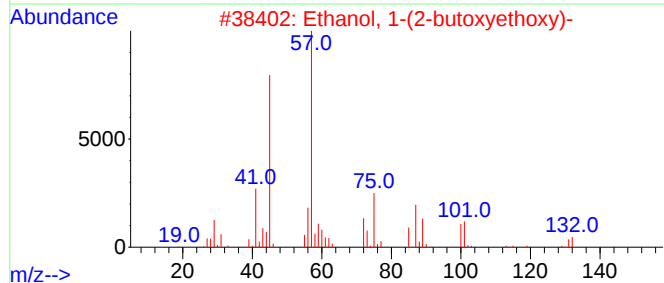
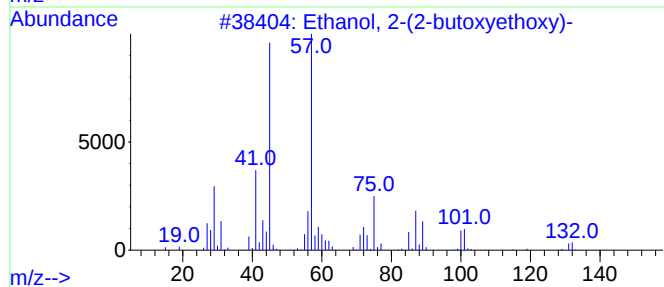
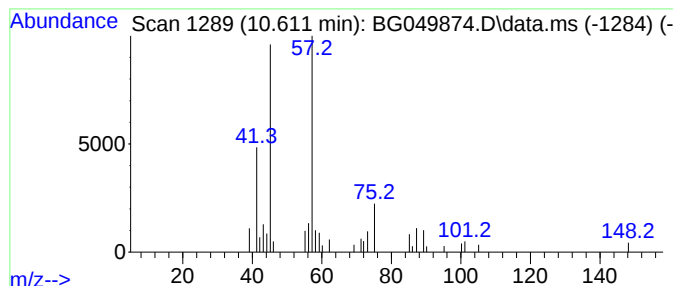
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.611	3.42 ng	43763	Naphthalene-d8	10.846

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	83
2			Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	74
3			Ethylene glycol monoisobutyl ether	118	C6H14O2	004439-24-1	64
4			15-Crown-5	220	C10H20O5	033100-27-5	50
5			DL-Lactamide, butyl ether	145	C7H15NO2	1000452-56-5	50



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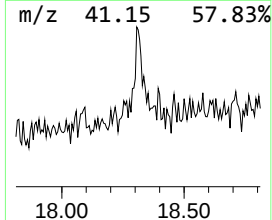
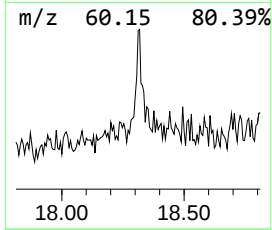
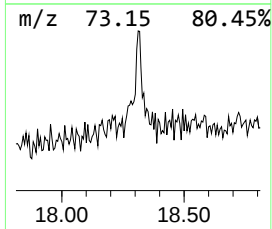
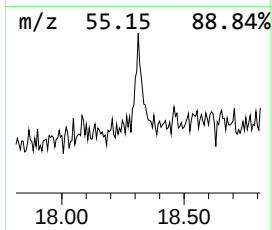
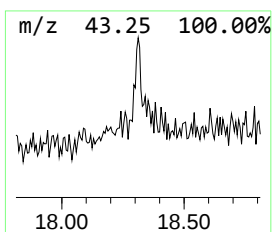
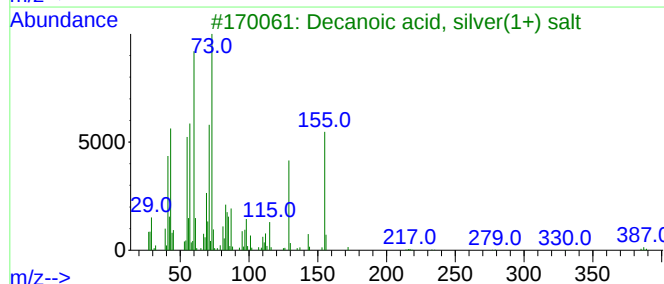
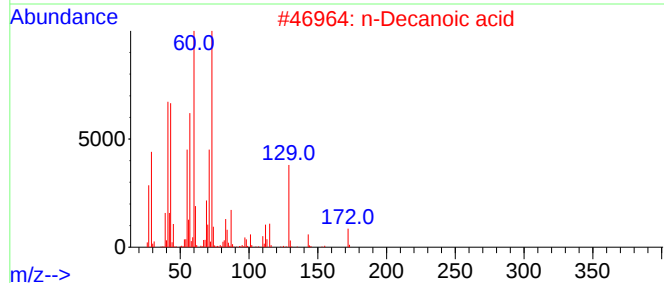
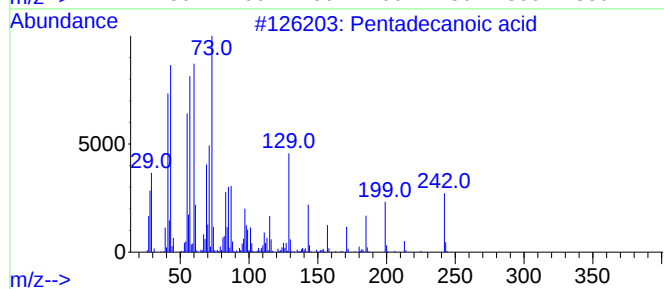
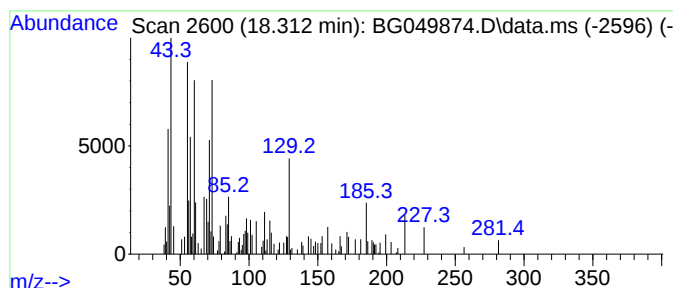
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TIC Library : C:\Database\NIST20.L
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Peak Number 4 Pentadecanoic acid Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD		R.T.	
18.312	2.91 ng	54786	Phenanthrene-d10		17.431	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Pentadecanoic acid		242	C15H30O2	001002-84-2	80
2	n-Decanoic acid		172	C10H20O2	000334-48-5	62
3	Decanoic acid, silver(1+) salt		278	C10H19AgO2	013126-67-5	58
4	Dodecanoic acid		200	C12H24O2	000143-07-7	53
5	Tetradecanoic acid		228	C14H28O2	000544-63-8	53



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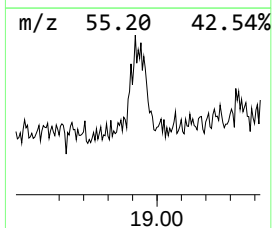
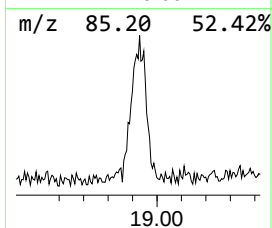
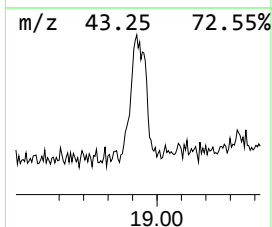
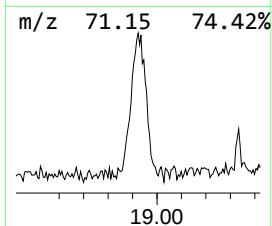
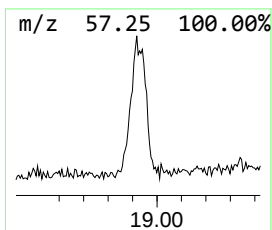
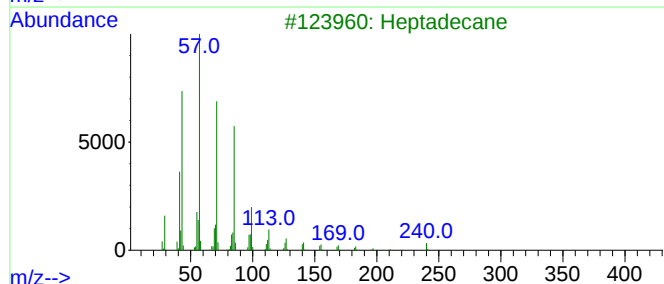
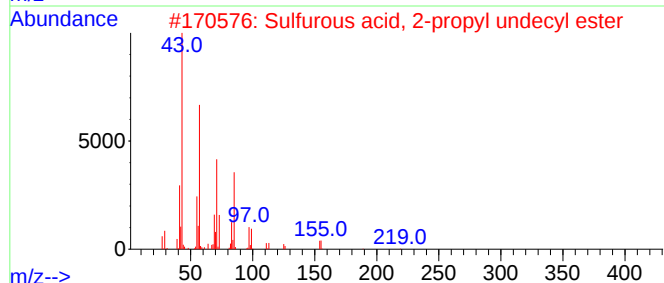
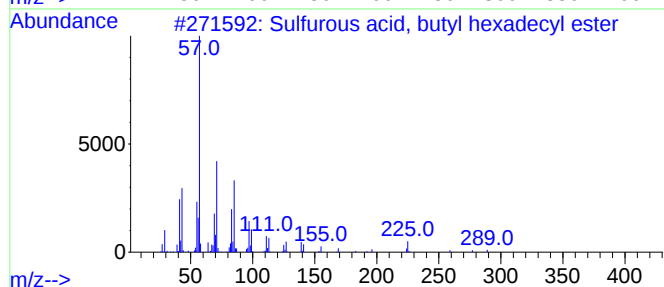
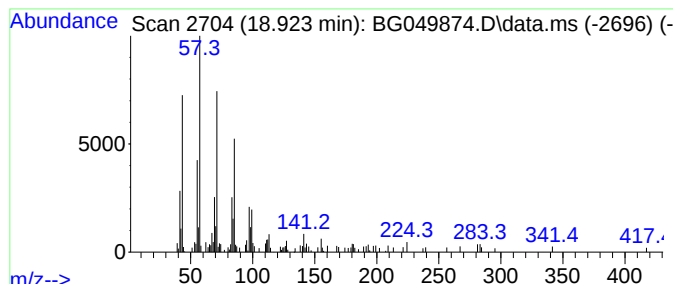
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TIC Library : C:\Database\NIST20.L
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Peak Number 5 Sulfurous acid, butyl hexad... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.923	6.74 ng	126879	Phenanthrene-d10	17.431

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Sulfurous acid, butyl hexadecyl ...	362	C20H42O3S	1000309-18-3	78
2			Sulfurous acid, 2-propyl undecyl...	278	C14H30O3S	1000309-12-2	72
3			Heptadecane	240	C17H36	000629-78-7	59
4			Pentadecane	212	C15H32	000629-62-9	59
5			Tridecane	184	C13H28	000629-50-5	58



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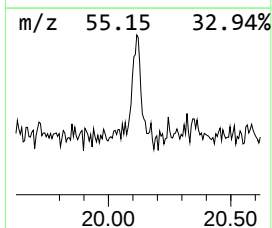
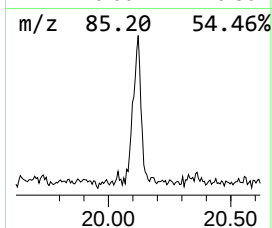
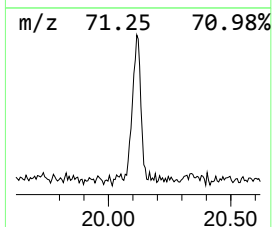
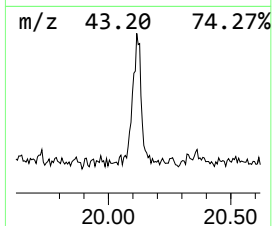
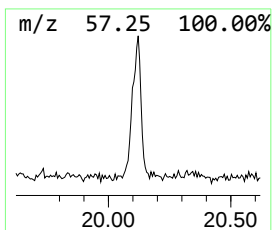
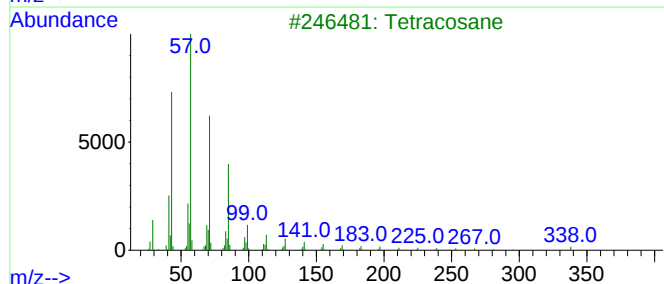
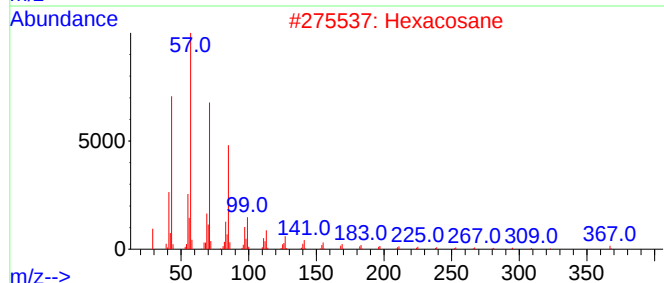
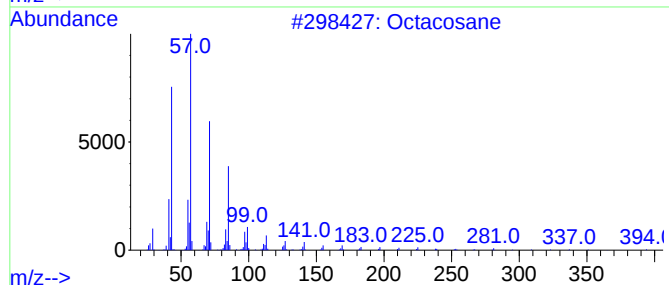
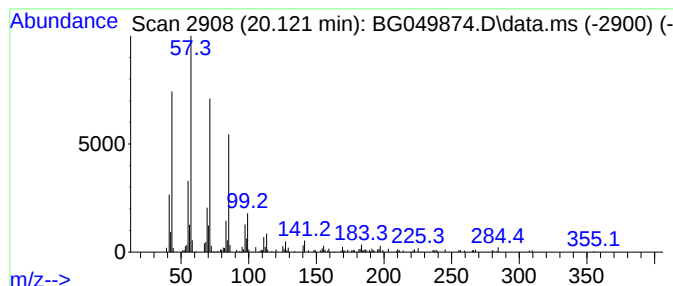
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Octacosane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.121	15.34 ng	303107	Chrysene-d12	21.713

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octacosane	394	C28H58	000630-02-4	87
2			Hexacosane	366	C26H54	000630-01-3	87
3			Tetracosane	338	C24H50	000646-31-1	80
4			Heptadecane	240	C17H36	000629-78-7	80
5			Octadecane	254	C18H38	000593-45-3	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081721\
Data File : BG049874.D
Acq On : 17 Aug 2021 12:24
Operator : CG/JU
Sample : M3412-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
PG-2

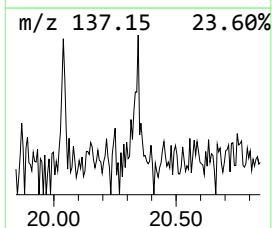
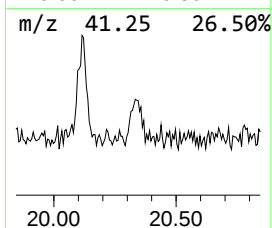
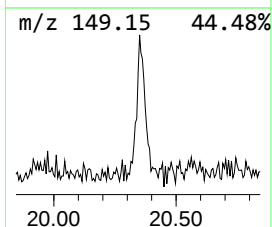
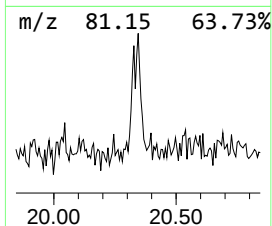
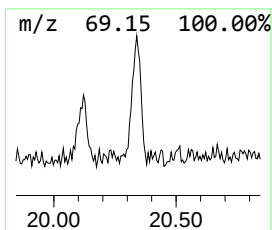
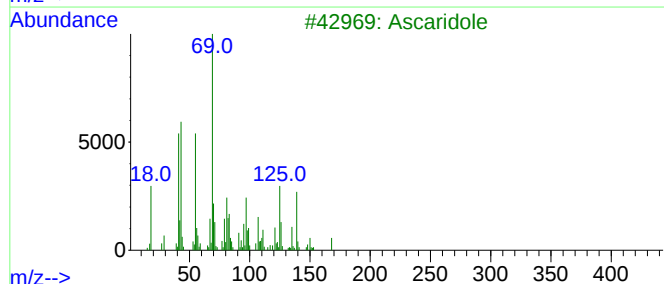
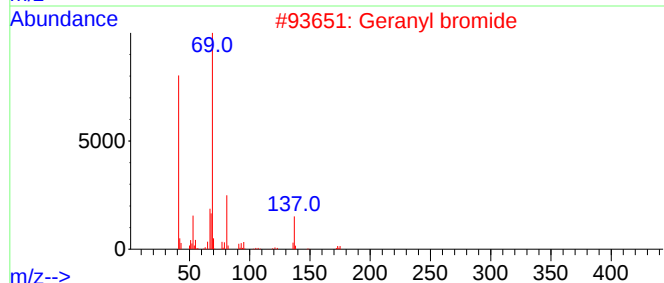
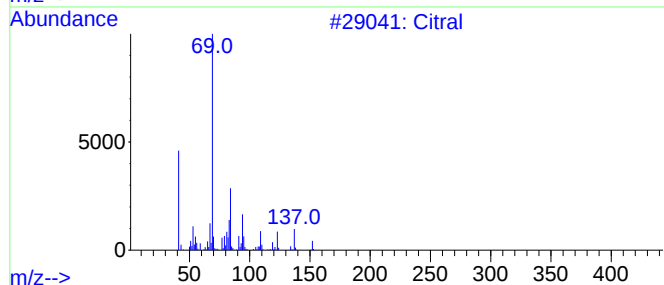
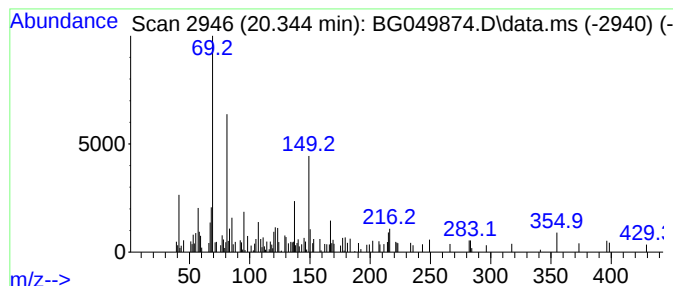
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 unknown20.345 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.345	4.77 ng	94262	Chrysene-d12	21.713

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Citral	152	C10H16O	005392-40-5	43
2			Geranyl bromide	216	C10H17Br	006138-90-5	43
3			Ascaridole	168	C10H16O2	000512-85-6	38
4			Docosa-2,6,10,14,18-pentaen-22-a...	384	C27H44O	1000163-04-7	38
5			2,6,10,14,18-Pentamethyl-2,6,10,...	342	C25H42	075581-03-2	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081721\
Data File : BG049874.D
Acq On : 17 Aug 2021 12:24
Operator : CG/JU
Sample : M3412-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

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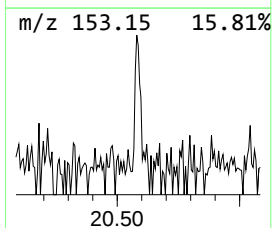
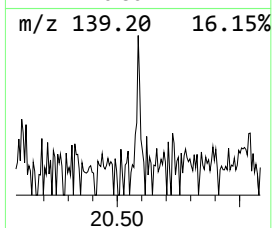
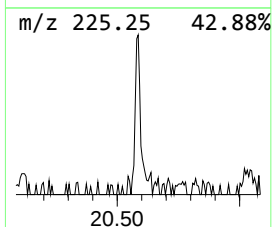
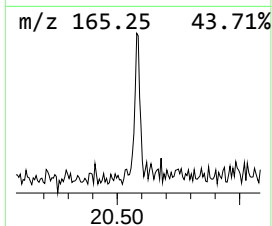
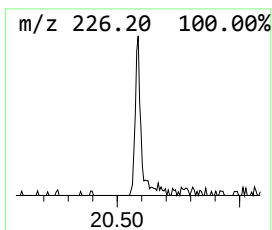
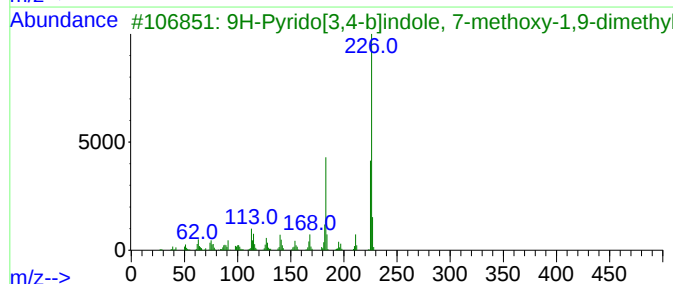
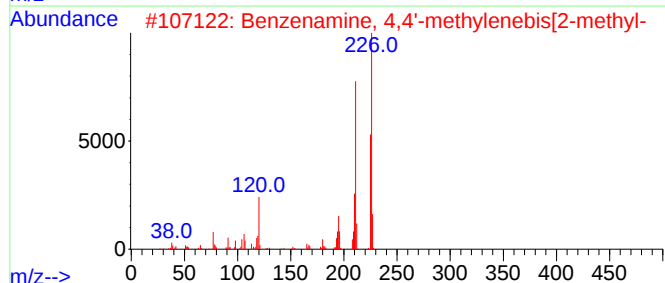
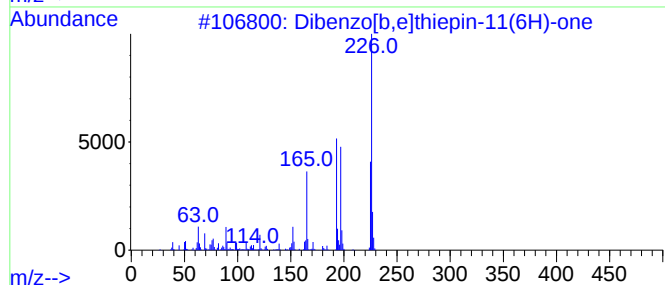
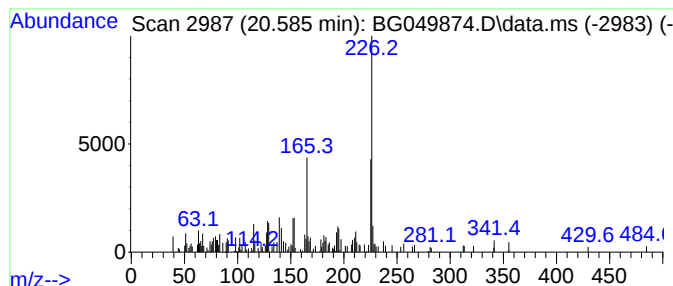
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Dibenzo[b,e]thiepin-11(6H)-one Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.585	2.48 ng	48934	Chrysene-d12	21.713

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dibenzo[b,e]thiepin-11(6H)-one	226	C14H10O5	001531-77-7	83
2			Benzenamine, 4,4'-methylenebis[2...	226	C15H18N2	000838-88-0	58
3			9H-Pyrido[3,4-b]indole, 7-methox...	226	C14H14N2O	143502-37-8	52
4			Acridin-9-amine, 1,2,3,4-tetrahy...	226	C15H18N2	297758-19-1	52
5			1,3-Diamino-5,6-dihydro-9-methyl...	226	C13H14N4	037436-39-8	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081721\
Data File : BG049874.D
Acq On : 17 Aug 2021 12:24
Operator : CG/JU
Sample : M3412-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
PG-2

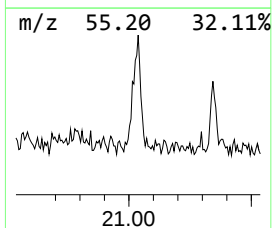
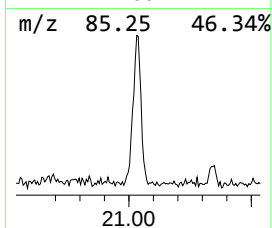
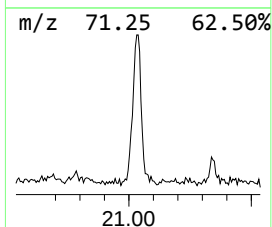
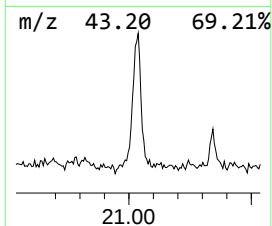
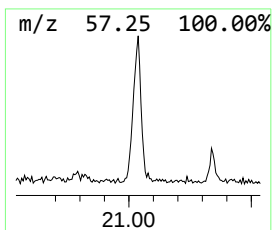
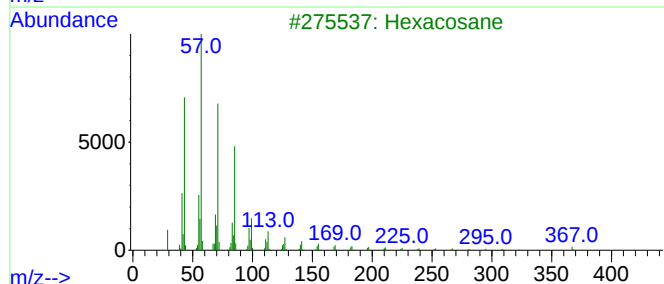
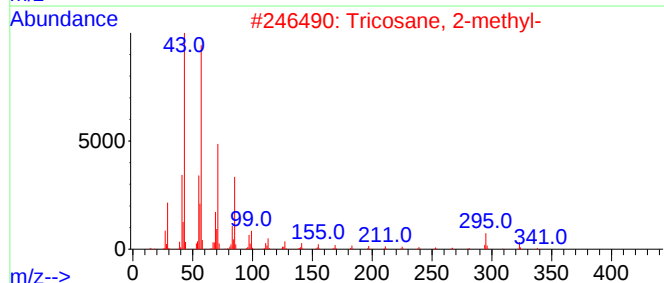
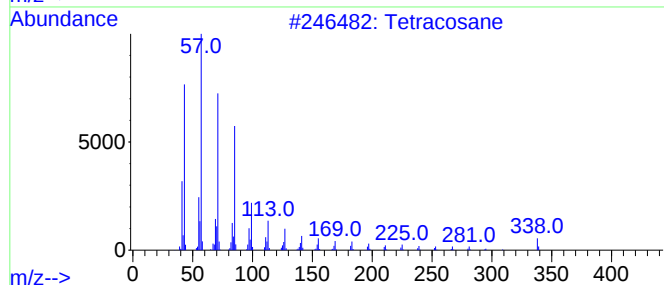
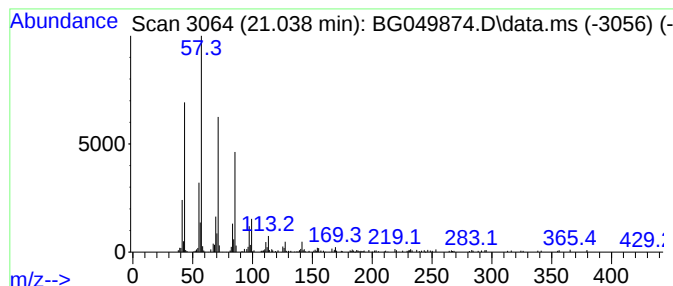
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Tetracosane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.038	18.74 ng	370194	Chrysene-d12	21.713

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetracosane	338	C24H50	000646-31-1	94
2			Tricosane, 2-methyl-	338	C24H50	001928-30-9	93
3			Hexacosane	366	C26H54	000630-01-3	90
4			Tricosane	324	C23H48	000638-67-5	90
5			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081721\
Data File : BG049874.D
Acq On : 17 Aug 2021 12:24
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Sample : M3412-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampled :
PG-2

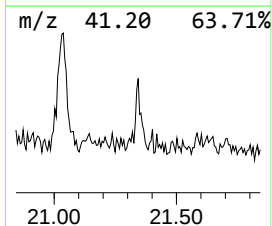
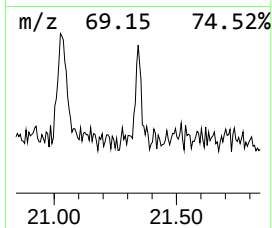
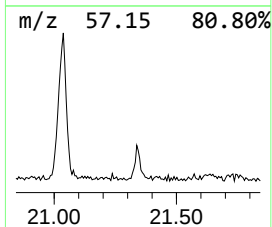
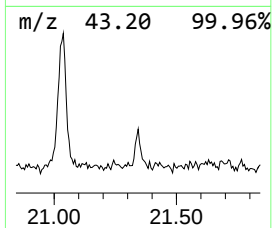
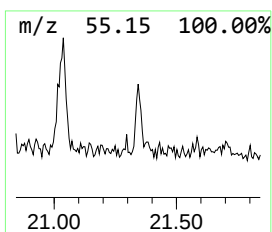
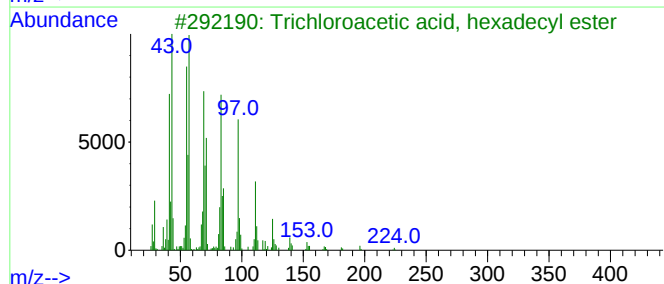
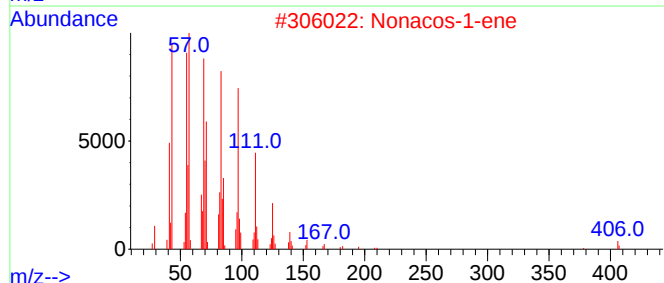
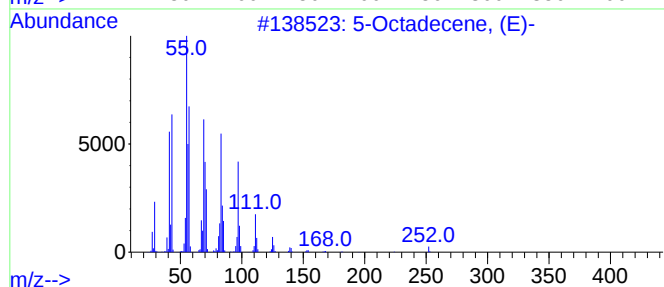
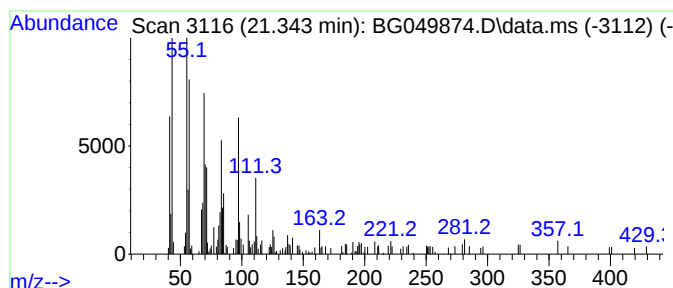
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 5-Octadecene, (E)- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.343	5.15 ng	101660	Chrysene-d12	21.713

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Octadecene, (E)-		252	C18H36	007206-21-5	93
2	Nonacos-1-ene		406	C29H58	018835-35-3	91
3	Trichloroacetic acid, hexadecyl ...		386	C18H33Cl3O2	074339-54-1	90
4	9-Octadecene, (E)-		252	C18H36	007206-25-9	89
5	1-Octadecene		252	C18H36	000112-88-9	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081721\
Data File : BG049874.D
Acq On : 17 Aug 2021 12:24
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Sample : M3412-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampled :
PG-2

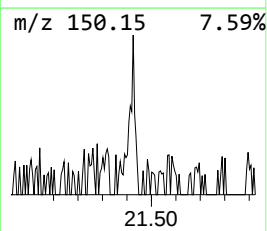
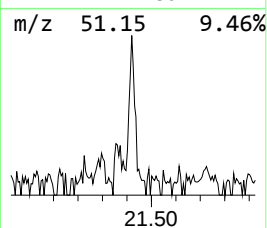
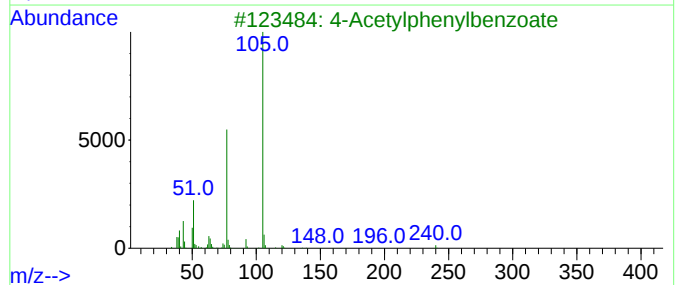
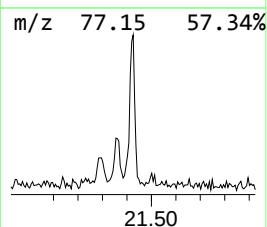
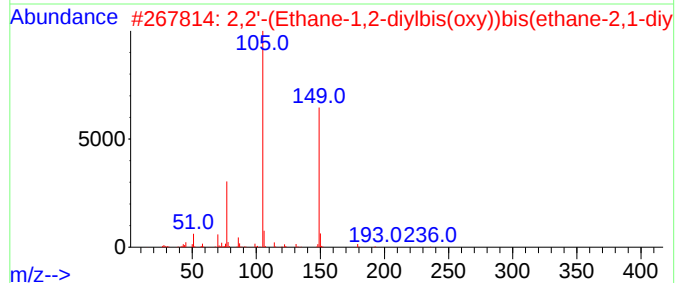
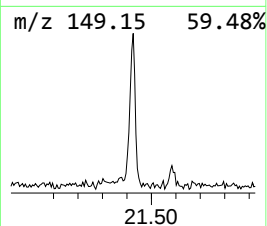
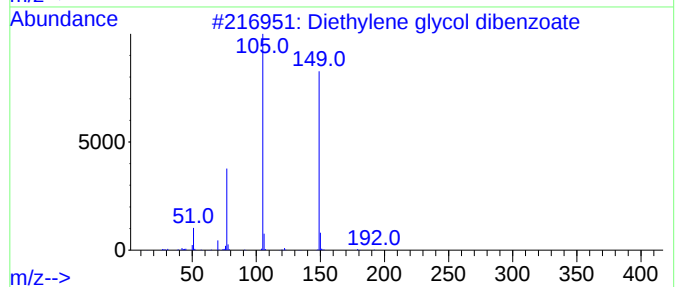
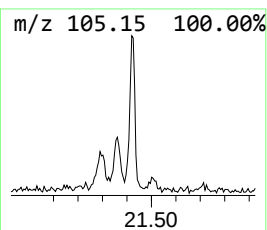
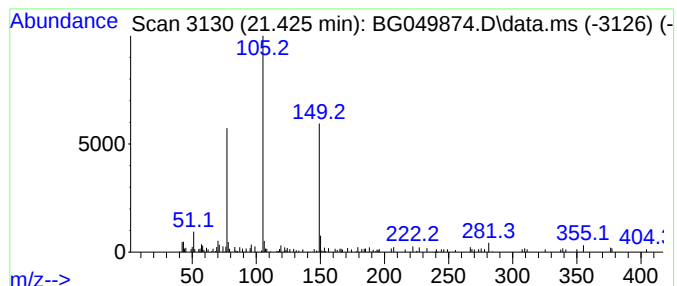
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Diethylene glycol dibenzoate Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.425	2.94 ng	58061	Chrysene-d12	21.713

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Diethylene glycol dibenzoate	314	C18H18O5	000120-55-8	50
2			2,2'-(Ethane-1,2-diylbis(oxy))bi...	358	C20H22O6	1000366-99-4	50
3			4-Acetylphenylbenzoate	240	C15H12O3	001523-18-8	43
4			Disulfide, dibenzoyl	274	C14H10O2S2	000644-32-6	38
5			Phenacyl thiocyanate	177	C9H7NOS	005399-30-4	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081721\
Data File : BG049874.D
Acq On : 17 Aug 2021 12:24
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Sample : M3412-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

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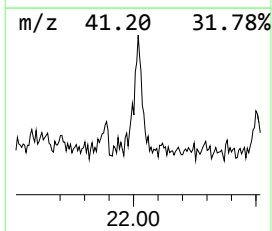
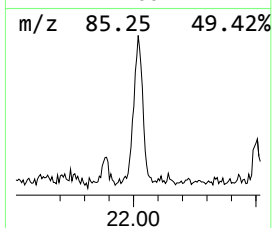
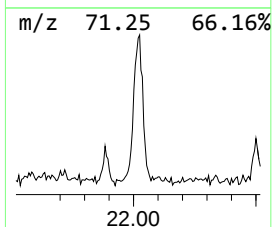
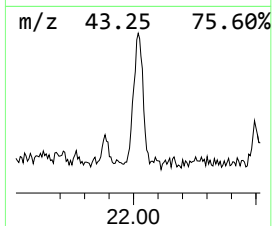
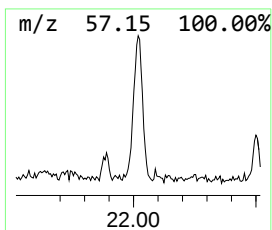
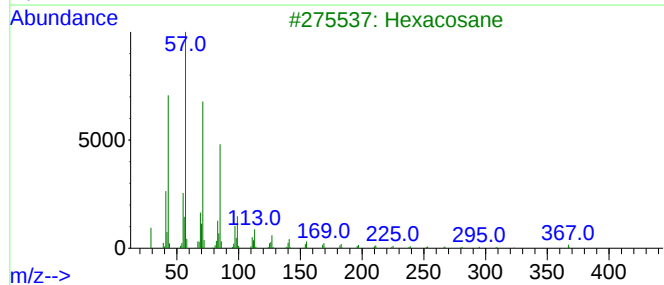
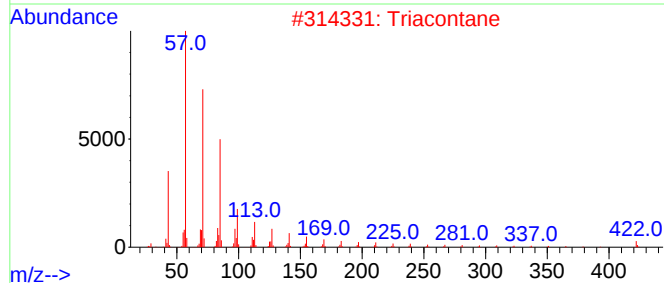
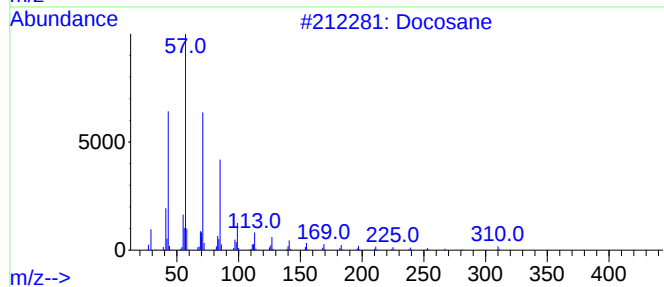
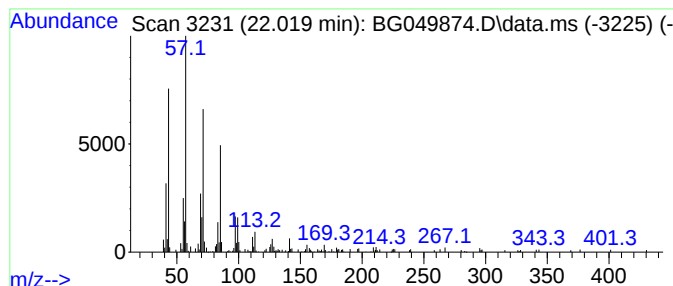
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Docosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.019	14.19 ng	280370	Chrysene-d12	21.713

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Docosane	310	C22H46	000629-97-0	87
2			Triacontane	422	C30H62	000638-68-6	87
3			Hexacosane	366	C26H54	000630-01-3	87
4			Dodecane, 2-methyl-	184	C13H28	001560-97-0	87
5			Tetratetracontane	619	C44H90	007098-22-8	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081721\
Data File : BG049874.D
Acq On : 17 Aug 2021 12:24
Operator : CG/JU
Sample : M3412-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
PG-2

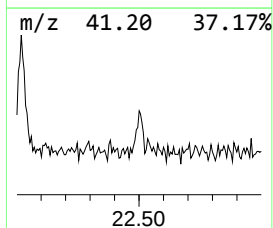
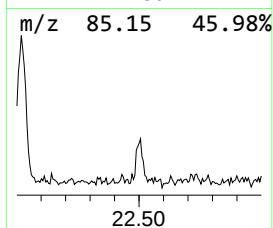
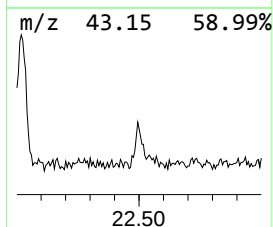
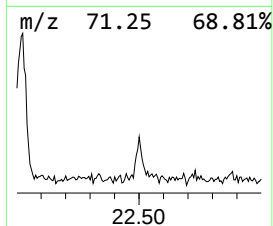
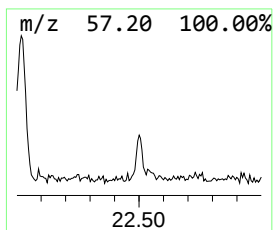
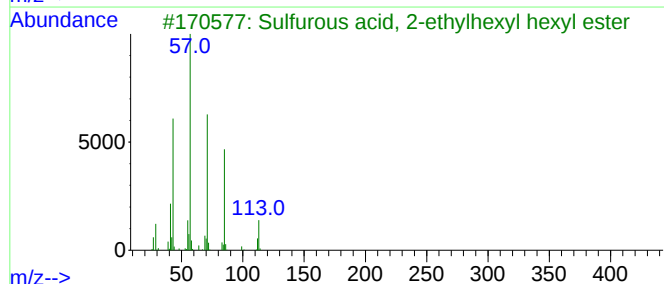
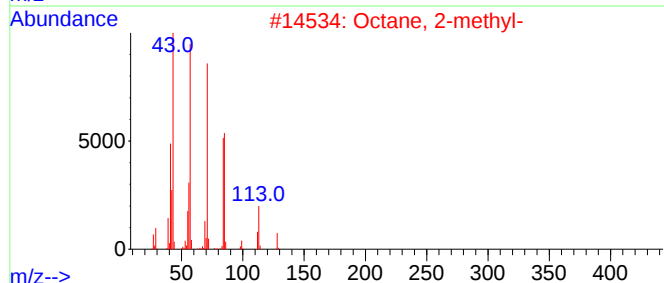
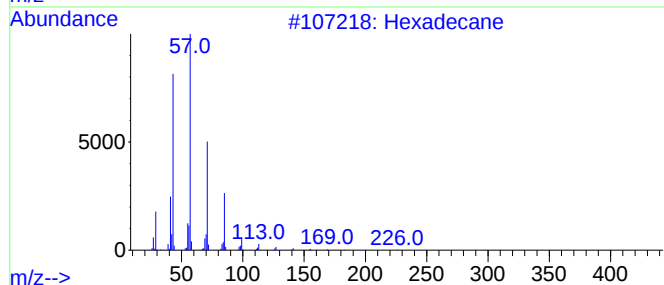
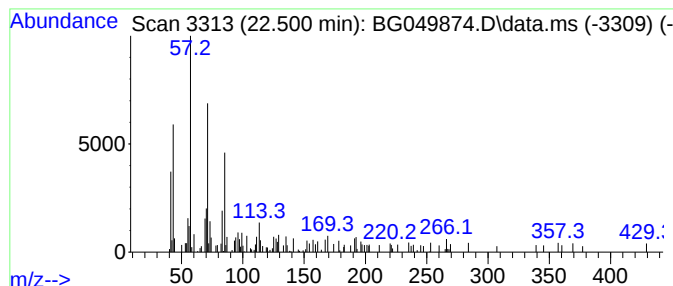
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Hexadecane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.500	2.69 ng	53157	Chrysene-d12	21.713

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane	226	C16H34	000544-76-3	58
2			Octane, 2-methyl-	128	C9H20	003221-61-2	52
3			Sulfurous acid, 2-ethylhexyl hex...	278	C14H30O3S	1000309-20-2	52
4			Octadecane, 1-iodo-	380	C18H37I	000629-93-6	50
5			Decane, 3,7-dimethyl-	170	C12H26	017312-54-8	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081721\
Data File : BG049874.D
Acq On : 17 Aug 2021 12:24
Operator : CG/JU
Sample : M3412-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
PG-2

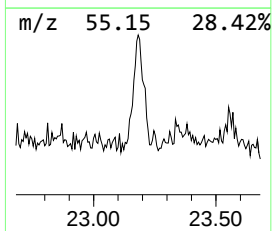
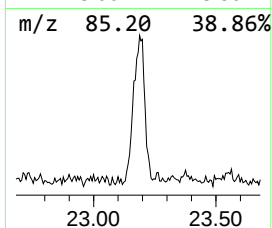
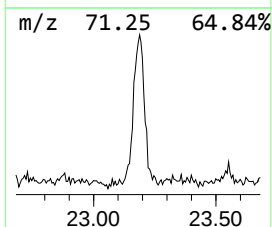
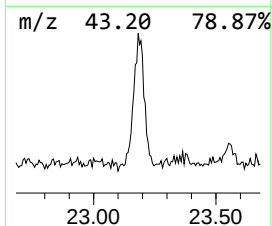
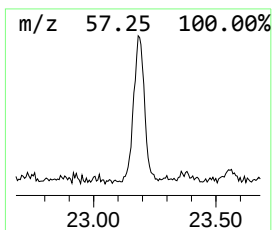
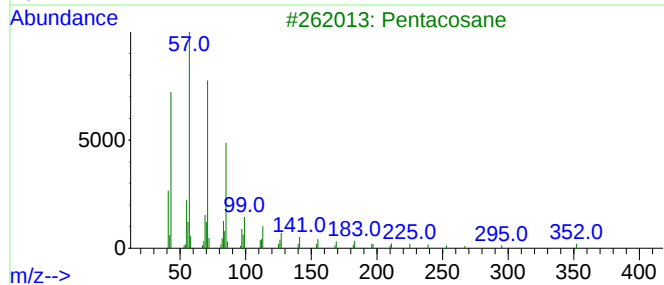
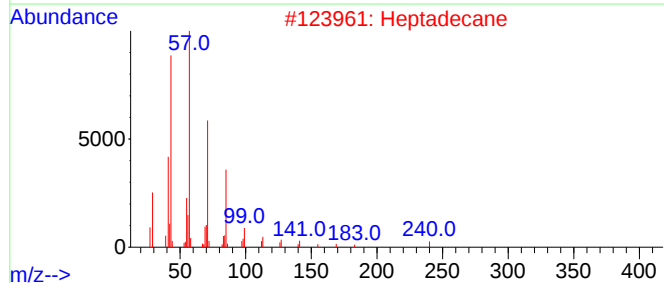
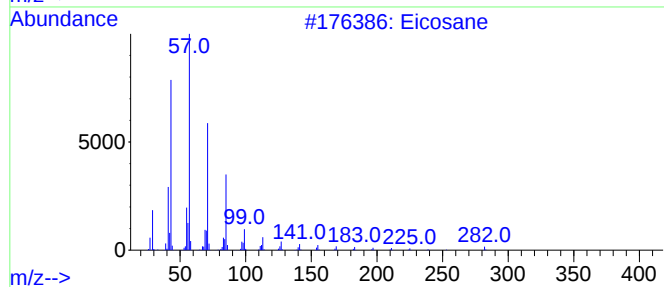
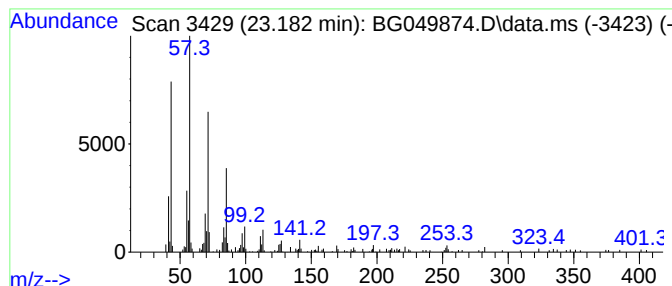
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Eicosane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.182	16.91 ng	333989	Chrysene-d12	21.713

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane	282	C20H42	000112-95-8	93
2			Heptadecane	240	C17H36	000629-78-7	91
3			Pentacosane	352	C25H52	000629-99-2	87
4			Octacosane	394	C28H58	000630-02-4	87
5			Hexacosane	366	C26H54	000630-01-3	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081721\
Data File : BG049874.D
Acq On : 17 Aug 2021 12:24
Operator : CG/JU
Sample : M3412-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
PG-2

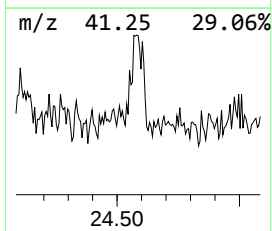
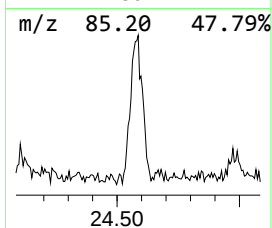
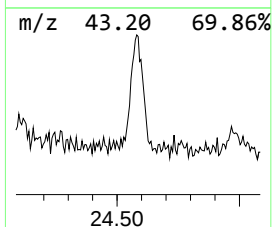
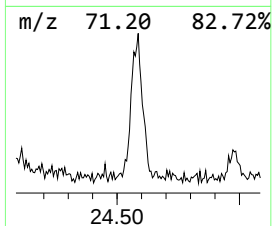
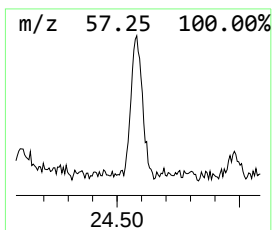
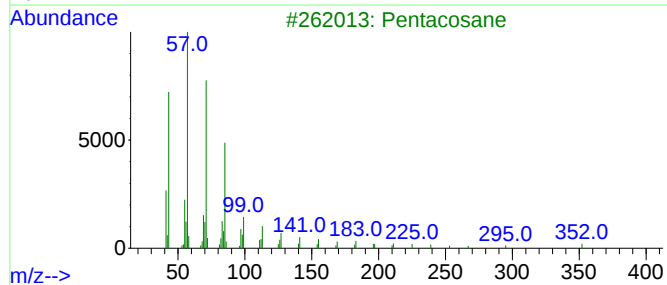
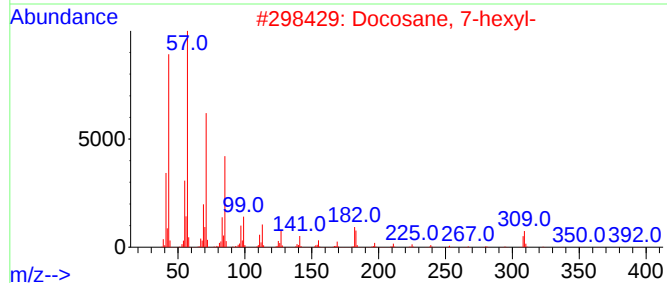
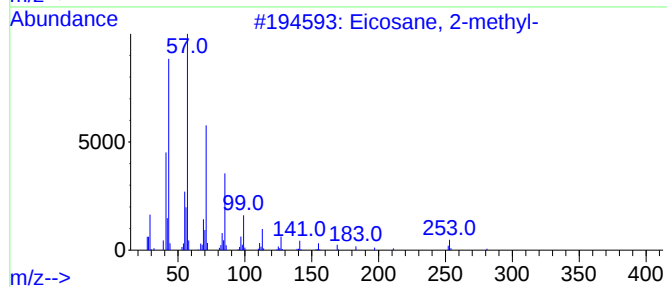
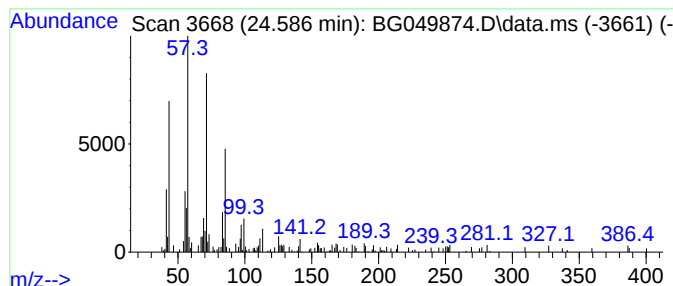
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Eicosane, 2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.586	10.40 ng	218157	Perylene-d12	24.985

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane, 2-methyl-	296	C21H44	001560-84-5	91
2			Docosane, 7-hexyl-	394	C28H58	055373-86-9	90
3			Pentacosane	352	C25H52	000629-99-2	87
4			Hexacosane	366	C26H54	000630-01-3	87
5			Tetratetracontane	619	C44H90	007098-22-8	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081721\
Data File : BG049874.D
Acq On : 17 Aug 2021 12:24
Operator : CG/JU
Sample : M3412-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
PG-2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG080321.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
Butane, 2-metho...	3.132	28.7	ng	265642	1	8.026	185272	20.0
2-Pentanone, 4-...	5.100	3.6	ng	33821	1	8.026	185272	20.0
Ethanol, 2-(2-b...	10.611	3.4	ng	43763	2	10.846	256198	20.0
Pentadecanoic acid	18.312	2.9	ng	54786	4	17.431	376249	20.0
Sulfurous acid,...	18.923	6.7	ng	126879	4	17.431	376249	20.0
Octacosane	20.121	15.3	ng	303107	5	21.713	395068	20.0
unknown20.345	20.345	4.8	ng	94262	5	21.713	395068	20.0
Dibenzo[b,e]thi...	20.585	2.5	ng	48934	5	21.713	395068	20.0
Tetracosane	21.038	18.7	ng	370194	5	21.713	395068	20.0
5-Octadecene, (E)-	21.343	5.2	ng	101660	5	21.713	395068	20.0
Diethylene glyc...	21.425	2.9	ng	58061	5	21.713	395068	20.0
Docosane	22.019	14.2	ng	280370	5	21.713	395068	20.0
Hexadecane	22.500	2.7	ng	53157	5	21.713	395068	20.0
Eicosane	23.182	16.9	ng	333989	5	21.713	395068	20.0
Eicosane, 2-met...	24.586	10.4	ng	218157	6	24.985	419450	20.0