

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052821\
 Data File : BP005640.D
 Acq On : 28 May 2021 15:48
 Operator : CG/JU
 Sample : M2532-01
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 MH-N-(MH-10.29)

Integration Parameters: rteint.p

Integrator: RTE

Smoother: 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_P\Methods\8270E-BP052721.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP005640.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.111	3	4	30	rVB	5068812	4959486	44.03%	7.366%
2	5.540	409	417	427	rBV	1478532	2138950	18.99%	3.177%
3	7.140	682	689	701	rBV	854837	1340334	11.90%	1.991%
4	7.987	824	833	839	rBV	654964	1038921	9.22%	1.543%
5	9.146	1021	1030	1043	rBV	2066655	3377770	29.99%	5.017%
6	10.793	1303	1310	1317	rBV	845588	1376059	12.22%	2.044%
7	13.240	1719	1726	1734	rBV	5416773	7534575	66.89%	11.191%
8	14.063	1862	1866	1870	rVB	96002	123945	1.10%	0.184%
9	14.457	1928	1933	1943	rVB2	54301	115624	1.03%	0.172%
10	14.604	1952	1958	1966	rVB	1149771	1731231	15.37%	2.571%
11	16.092	2204	2211	2219	rBV	3947512	5662024	50.27%	8.410%
12	16.733	2313	2320	2331	rBV5	150449	332416	2.95%	0.494%
13	17.351	2419	2425	2438	rVB	1468098	2093161	18.58%	3.109%
14	18.233	2569	2575	2583	rBV	1029902	1470001	13.05%	2.183%
15	18.492	2614	2619	2623	rBV	257020	402121	3.57%	0.597%
16	18.539	2623	2627	2635	rVV	544957	1119821	9.94%	1.663%
17	18.704	2637	2655	2681	rVB3	581304	3469248	30.80%	5.153%
18	19.457	2779	2783	2799	rBV	425983	625394	5.55%	0.929%
19	19.898	2852	2858	2863	rBV	9172305	11264311	100.00%	16.730%
20	20.469	2945	2955	2958	rBV2	708242	1956606	17.37%	2.906%
21	20.580	2968	2974	2980	rBV7	155801	476356	4.23%	0.708%
22	20.845	3013	3019	3029	rBV4	309764	703575	6.25%	1.045%
23	21.127	3063	3067	3073	rBV2	359310	517144	4.59%	0.768%
24	21.416	3112	3116	3122	rVB	2029189	2600944	23.09%	3.863%
25	21.557	3134	3140	3153	rVB	1477387	3060587	27.17%	4.546%
26	22.668	3322	3329	3340	rVB2	841221	2380310	21.13%	3.535%
27	23.545	3473	3478	3499	rVB3	575062	2484545	22.06%	3.690%
28	23.862	3525	3532	3539	rVB2	1464676	2973014	26.39%	4.416%

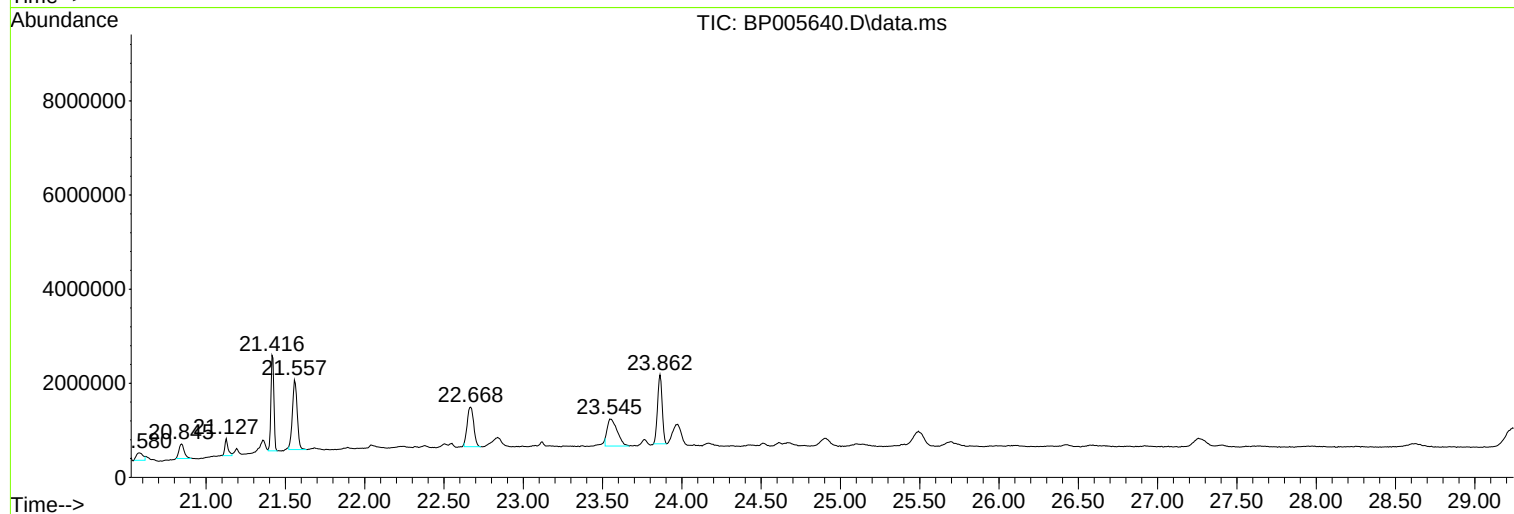
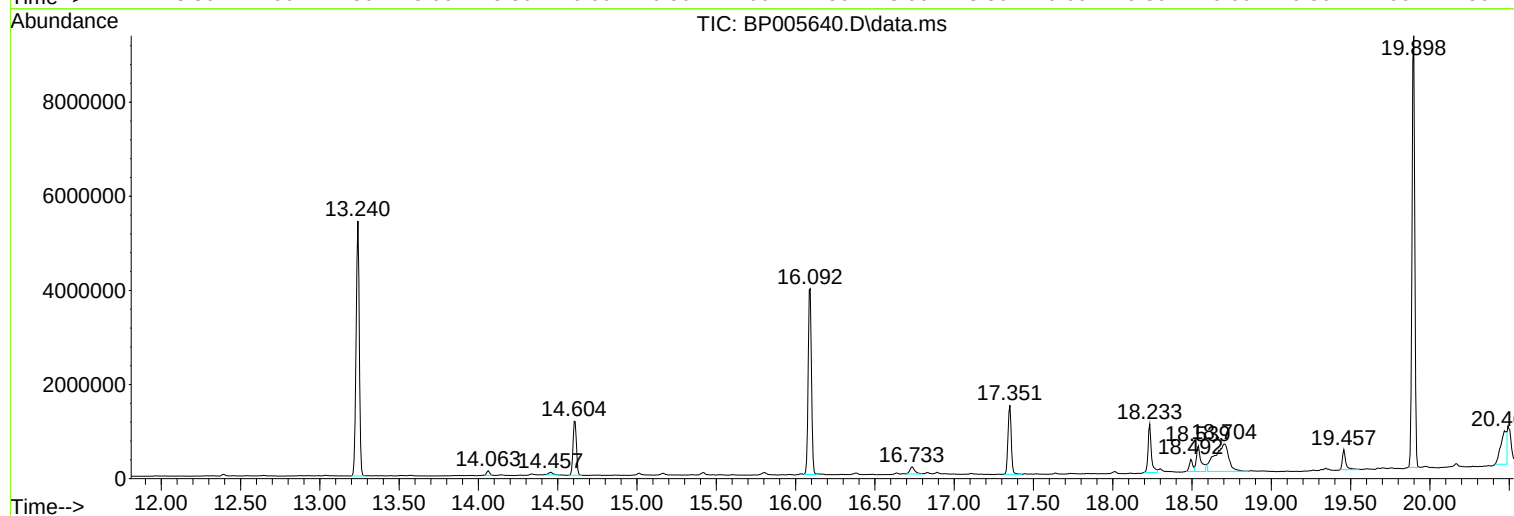
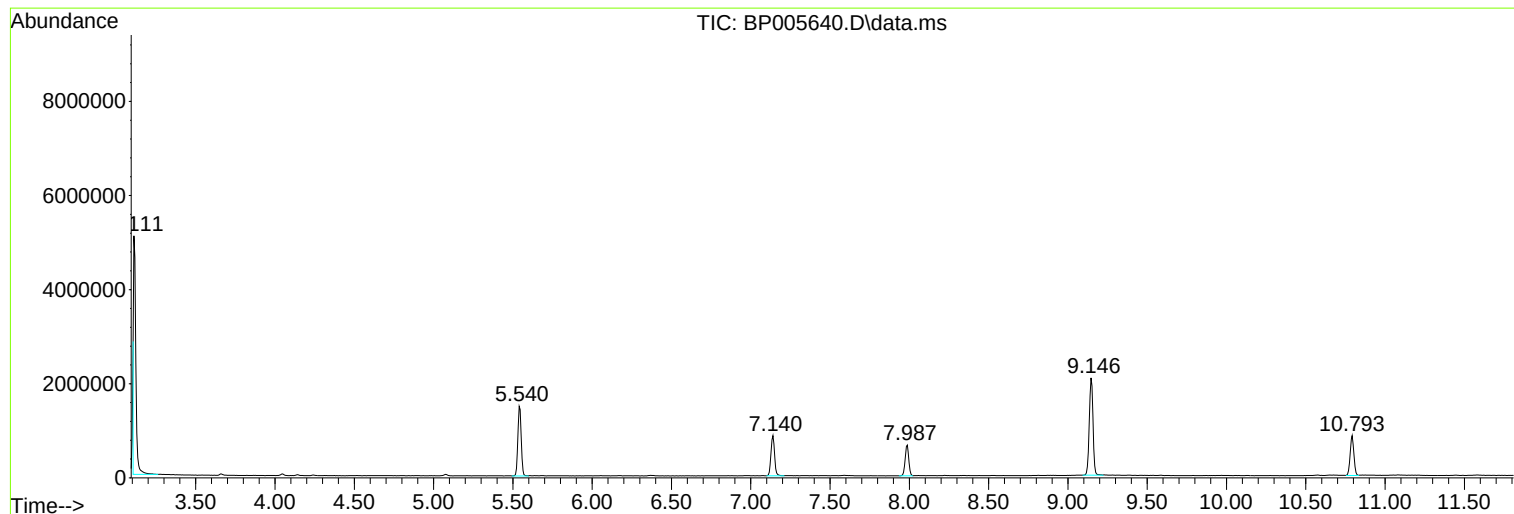
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP052721.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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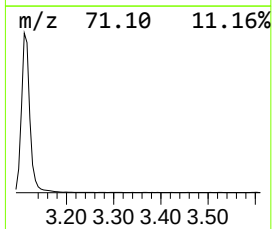
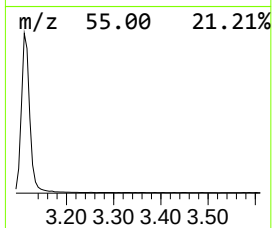
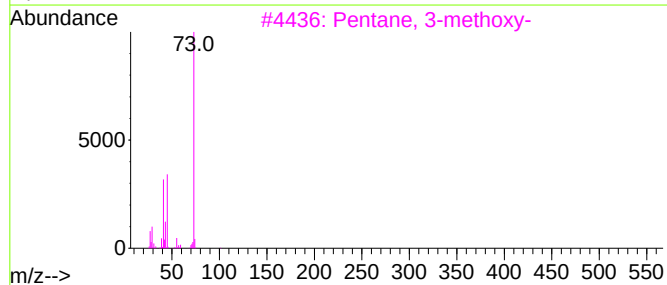
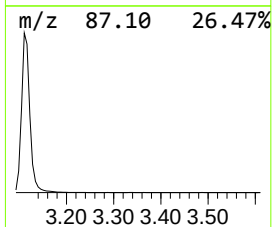
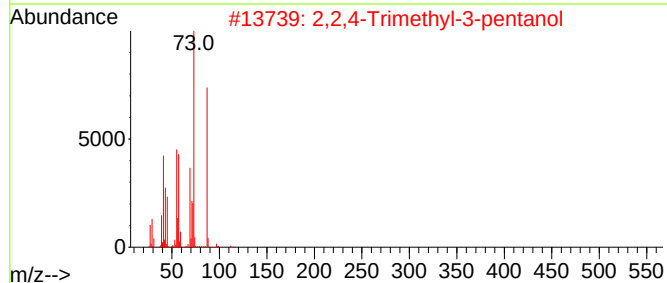
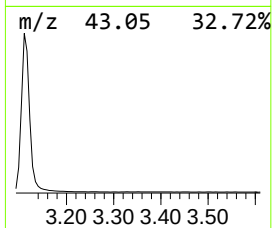
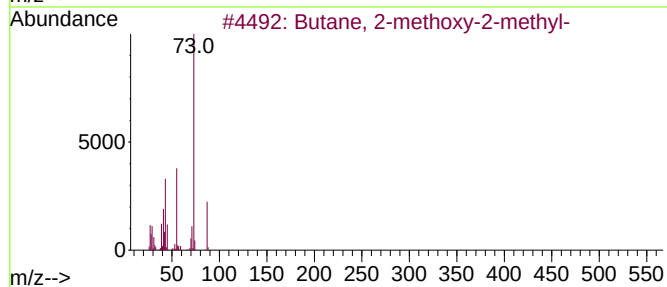
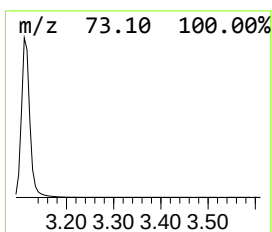
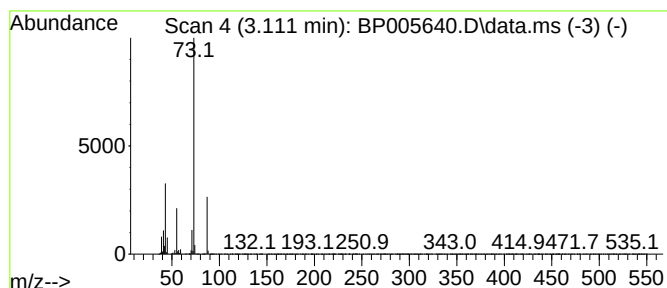
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.111	95.47 ng	4959490	1,4-Dichlorobenzene-d4	7.987

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
4			N-Ethylformamide	73	C3H7NO	000627-45-2	9
5			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9



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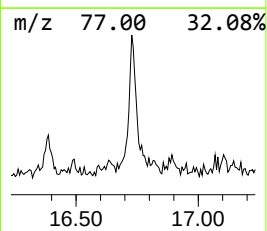
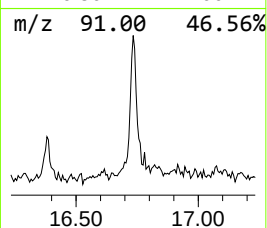
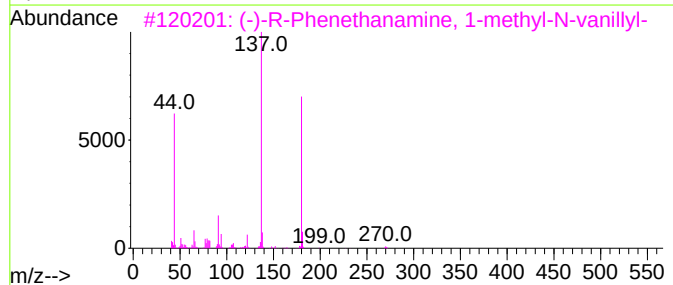
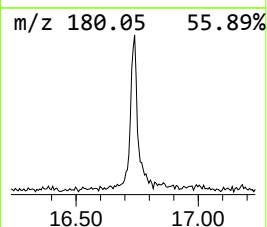
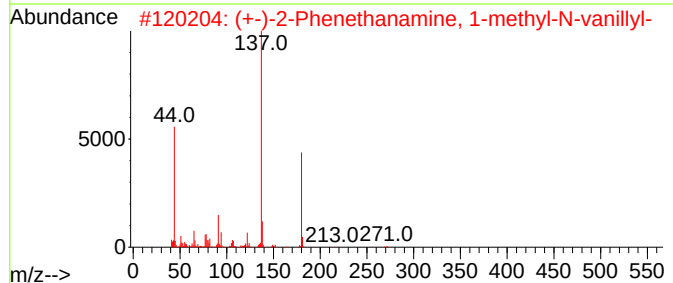
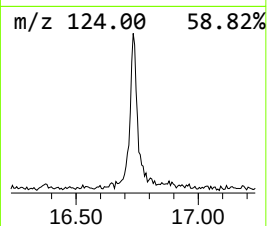
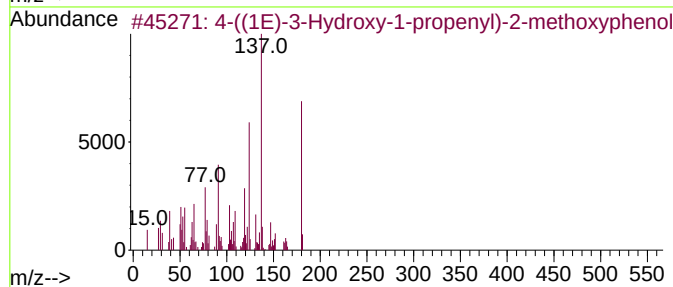
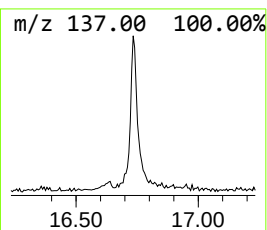
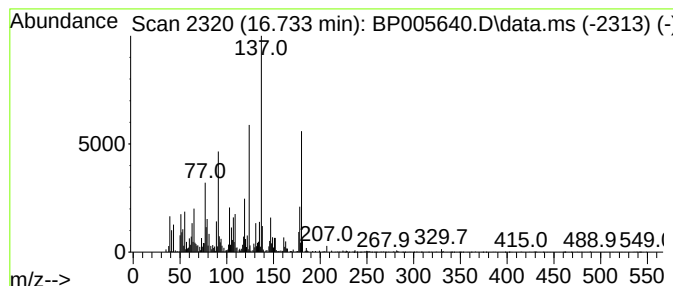
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 4-((1E)-3-Hydroxy-1-propenyl)-2-methoxyphenol Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.733	3.18 ng	332416	Phenanthrene-d10	17.351

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-((1E)-3-Hydroxy-1-propenyl)-2-methoxyphenol	180	C10H12O3	1000297-95-5	97
2			(+)-2-Phenethanamine, 1-methyl-N-vanillyl-	271	C17H21NO2	1000127-89-8	38
3			(-)-R-Phenethanamine, 1-methyl-N-vanillyl-	271	C17H21NO2	1000127-90-4	38
4			(+)-s-2-Phenethanamine, 1-methyl-N-vanillyl-	271	C17H21NO2	1000127-89-6	38
5			Pyrazine, 2-methoxy-3-(1-methylethyl)-	152	C8H12N2O	025773-40-4	35



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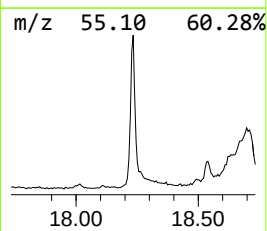
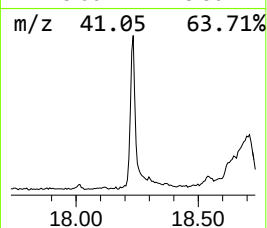
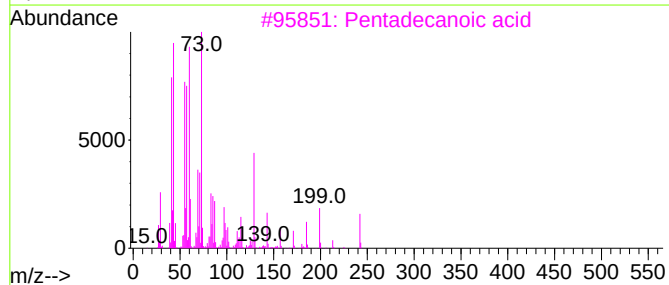
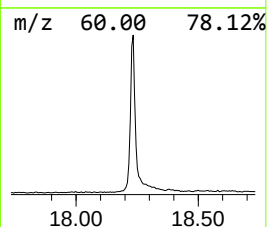
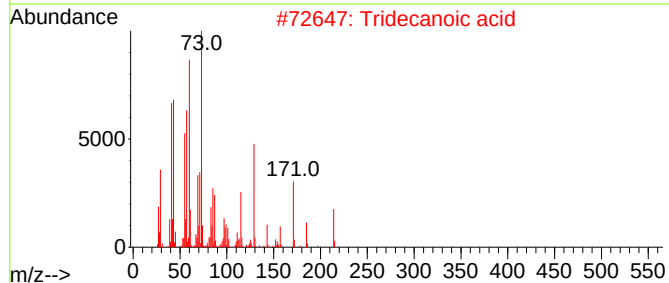
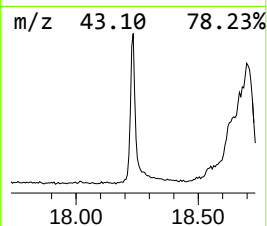
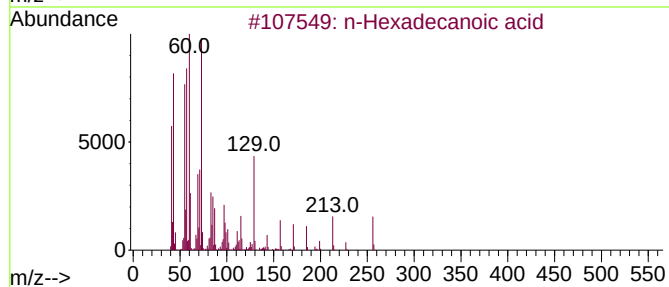
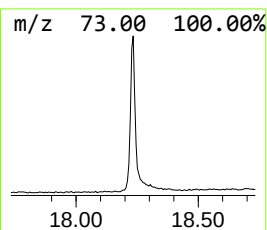
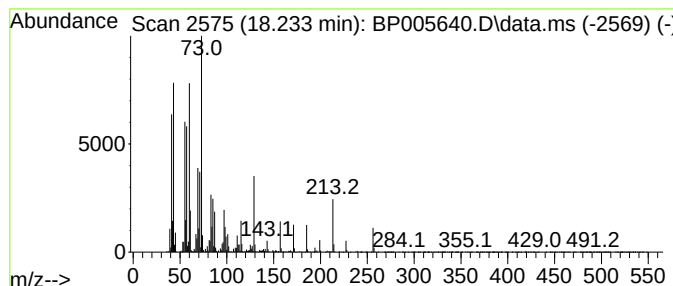
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 n-Hexadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.233	14.05 ng	1470000	Phenanthrene-d10	17.351

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2			Tridecanoic acid	214	C13H26O2	000638-53-9	92
3			Pentadecanoic acid	242	C15H30O2	001002-84-2	87
4			Octadecanoic acid	284	C18H36O2	000057-11-4	81
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	81



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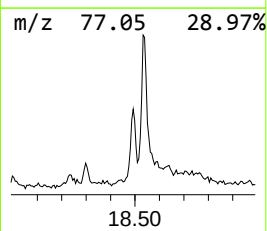
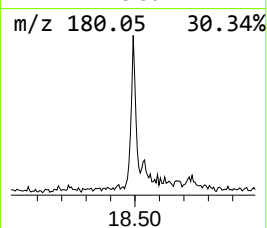
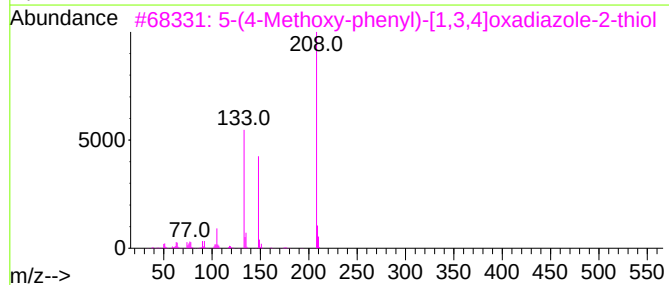
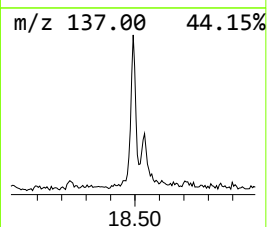
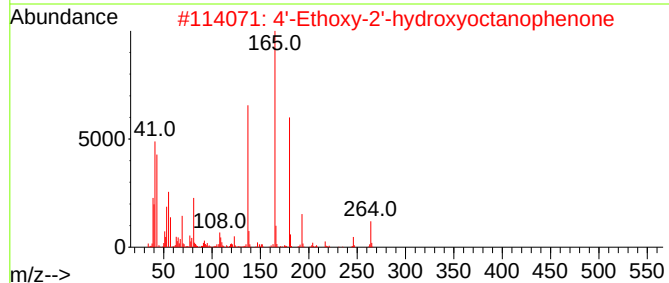
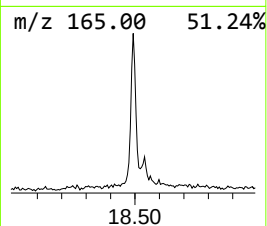
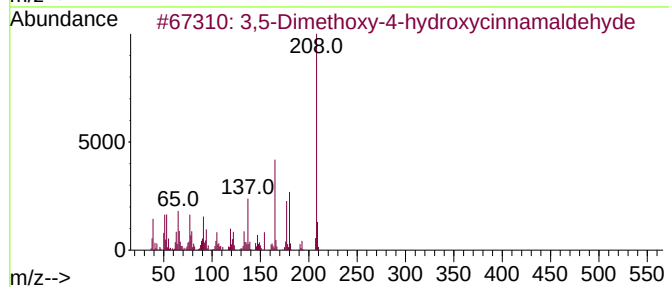
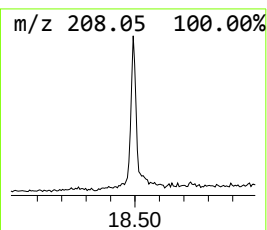
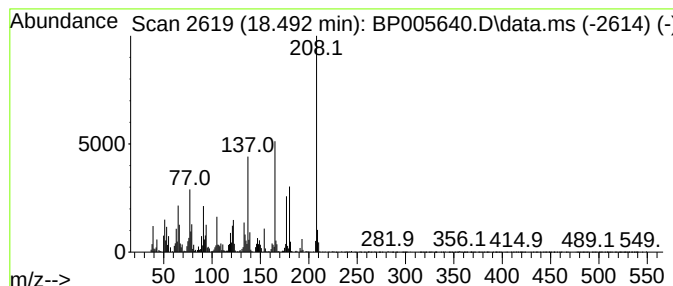
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 4 3,5-Dimethoxy-4-hydroxycinn... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.492	3.84 ng	402121	Phenanthrene-d10	17.351

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3,5-Dimethoxy-4-hydroxycinnamal... 208	C11H12O4	087345-53-7	96	
2			4'-Ethoxy-2'-hydroxyoctanophenone 264	C16H24O3	022198-47-6	42	
3			5-(4-Methoxy-phenyl)-[1,3,4]oxad... 208	C9H8N2O2S	023766-26-9	42	
4			2-Propenoic acid, 3-(2,3-dimetho... 208	C11H12O4	007345-82-6	41	
5			1,2-Dimethoxy-4-(3-methoxy-1-pro... 208	C12H16O3	1000313-35-0	41	



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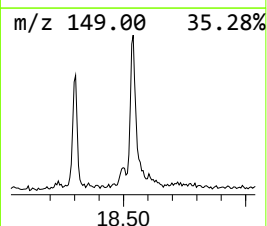
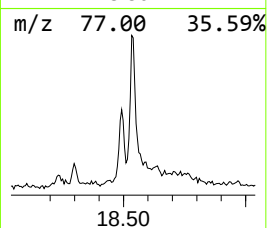
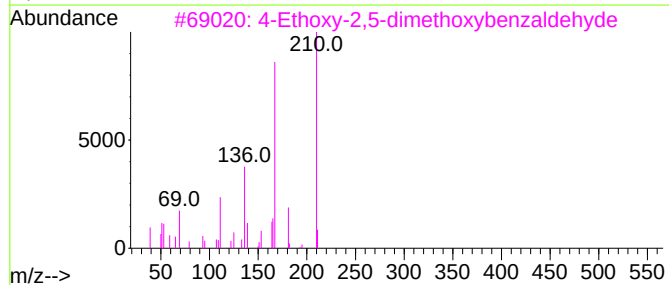
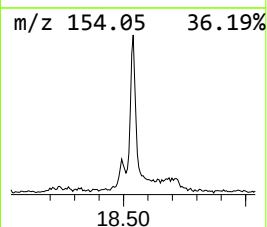
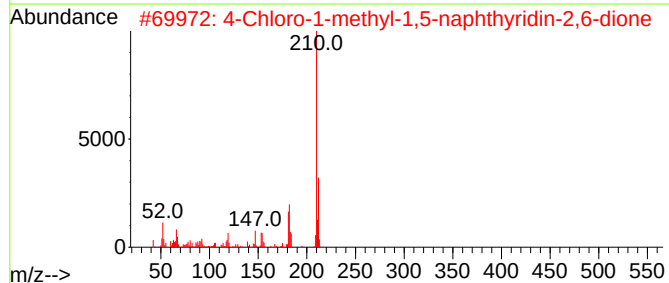
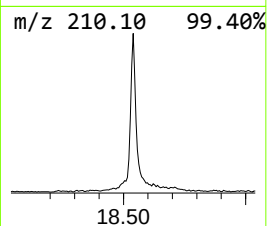
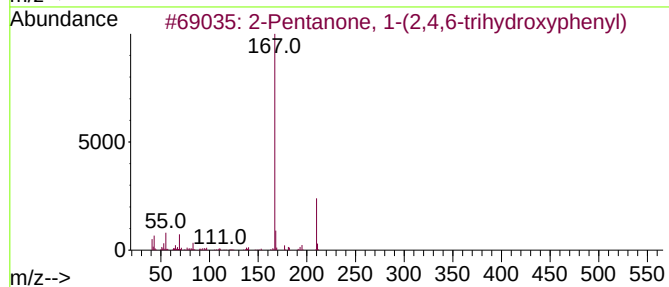
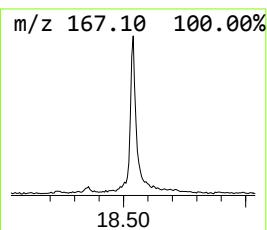
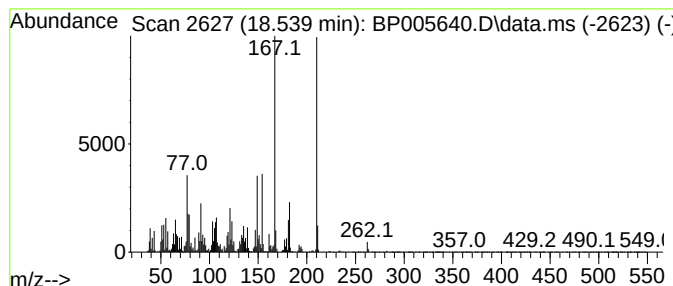
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 5 2-Pentanone, 1-(2,4,6-trihydroxy... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.539	10.70 ng	1119820	Phenanthrene-d10	17.351

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 1-(2,4,6-trihydroxy...	210	C11H14O4	1000116-22-3	58
2			4-Chloro-1-methyl-1,5-naphthyrid...	210	C9H7C1N2O2	1000213-59-5	45
3			4-Ethoxy-2,5-dimethoxybenzaldehyde	210	C11H14O4	1000121-30-2	43
4			Phenol, 2-(2-imidazo[1,2-a]pyrid...	210	C13H10N2O	057636-32-5	30
5			3,5-Pyridinedicarboxylic acid	167	C7H5NO4	000499-81-0	27



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MH-N-(MH-10.29)

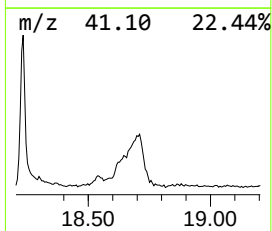
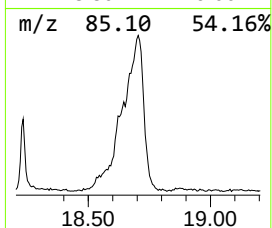
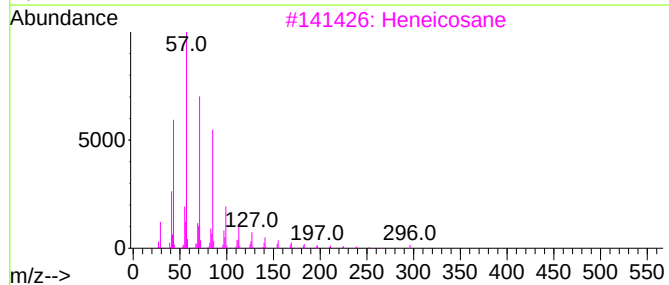
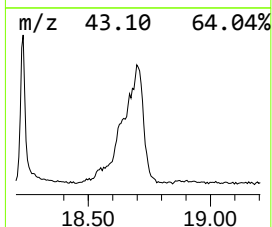
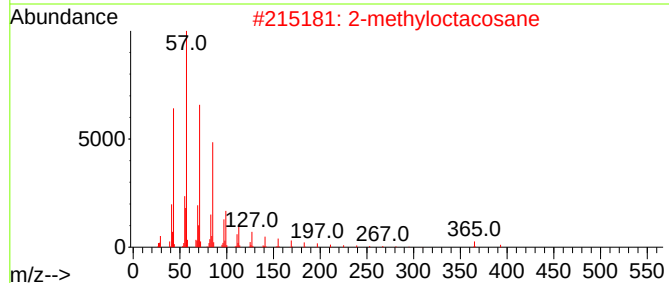
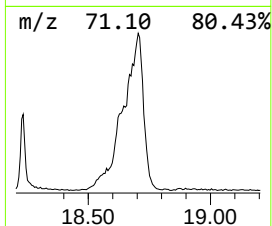
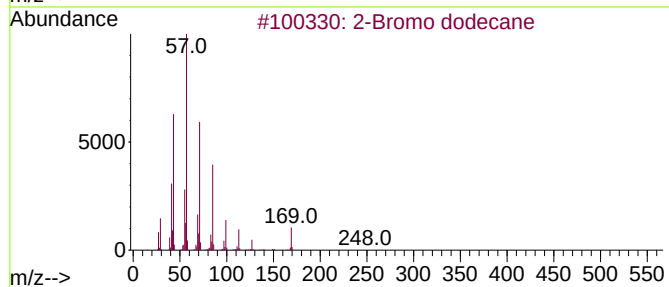
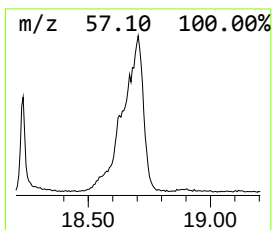
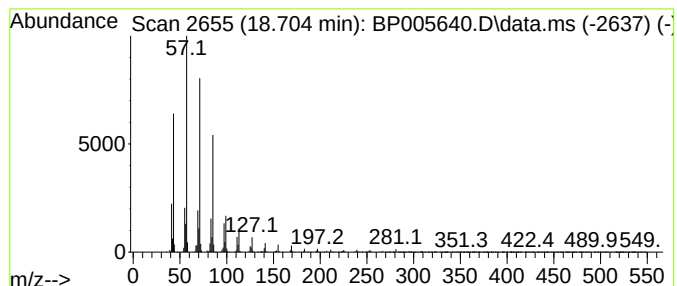
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 2-Bromo dodecane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.704	33.15 ng	3469250	Phenanthrene-d10	17.351

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Bromo dodecane	248	C12H25Br	013187-99-0	93
2		2-methyloctacosane	408	C29H60	1000376-72-8	91
3		Heneicosane	296	C21H44	000629-94-7	91
4		Heptadecane, 9-hexyl-	324	C23H48	055124-79-3	91
5		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052821\
Data File : BP005640.D
Acq On : 28 May 2021 15:48
Operator : CG/JU
Sample : M2532-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
MH-N-(MH-10.29)

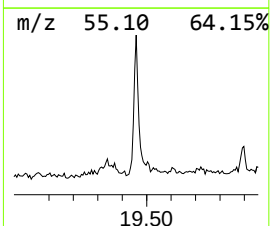
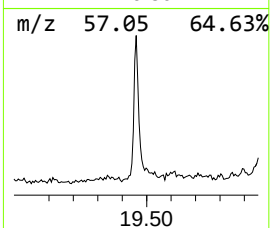
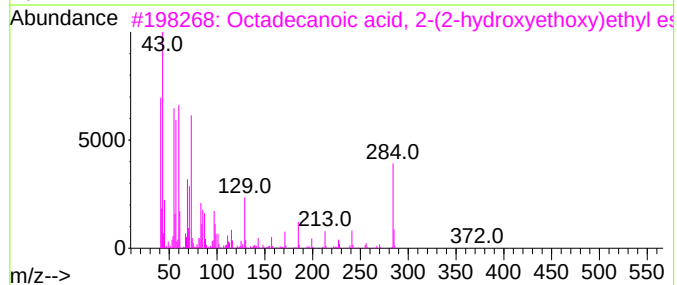
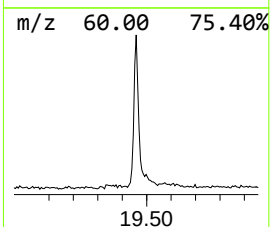
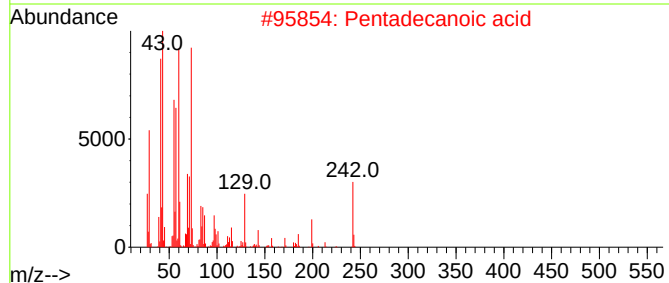
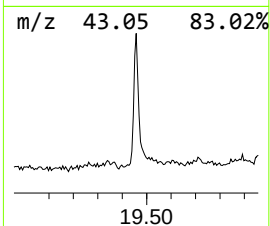
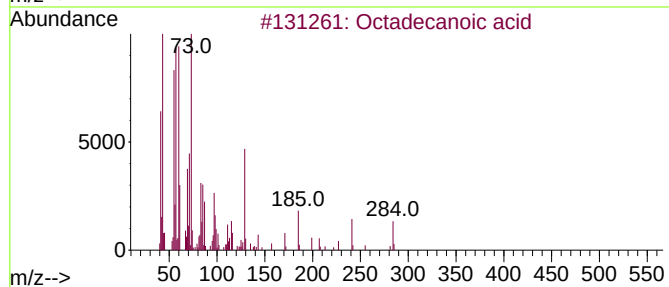
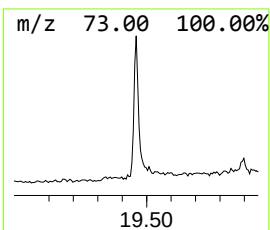
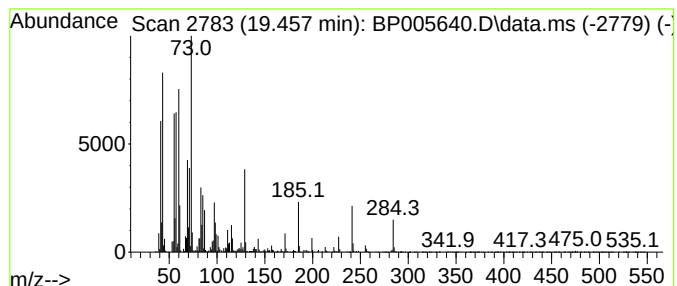
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Octadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.457	4.81 ng	625394	Chrysene-d12	21.416

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			Pentadecanoic acid	242	C15H30O2	001002-84-2	91
3			Octadecanoic acid, 2-(2-hydroxye...	372	C22H44O4	000106-11-6	87
4			Tetradecanoic acid	228	C14H28O2	000544-63-8	87
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052821\
Data File : BP005640.D
Acq On : 28 May 2021 15:48
Operator : CG/JU
Sample : M2532-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
MH-N-(MH-10.29)

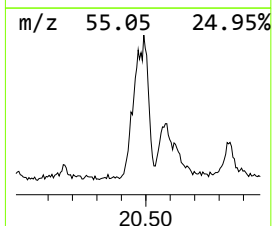
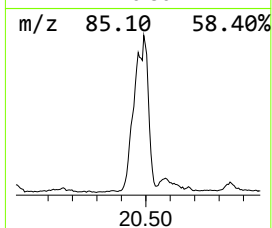
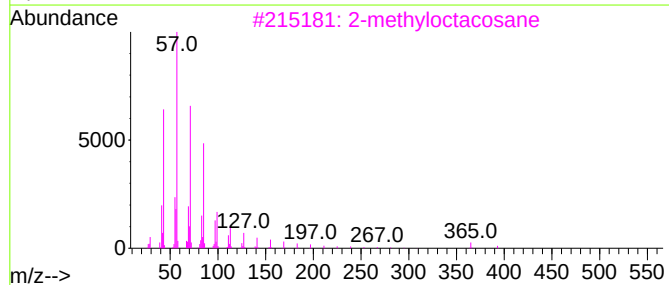
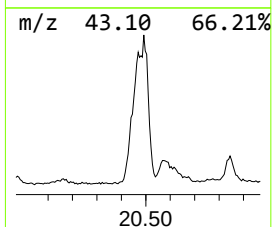
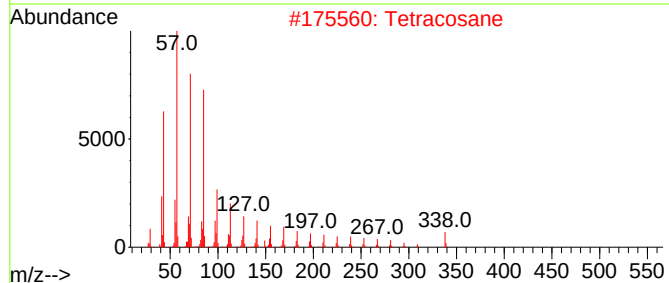
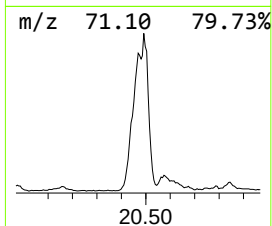
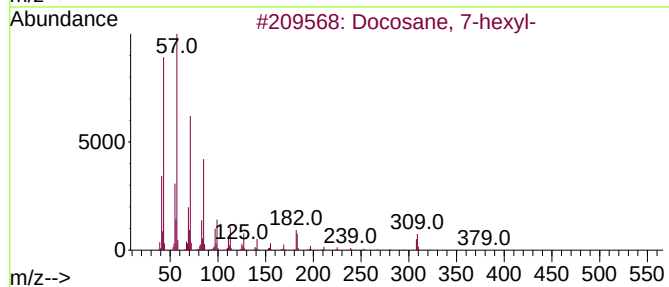
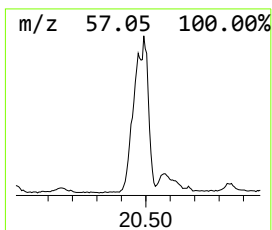
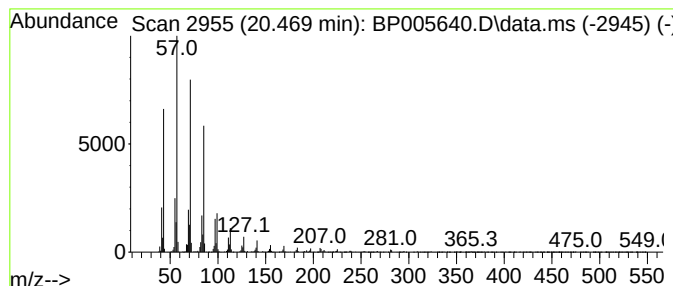
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Docosane, 7-hexyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.468	15.05 ng	1956610	Chrysene-d12	21.416

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Docosane, 7-hexyl-	394	C28H58	055373-86-9	94
2		Tetracosane	338	C24H50	000646-31-1	91
3		2-methyloctacosane	408	C29H60	1000376-72-8	91
4		Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	91
5		Octadecane	254	C18H38	000593-45-3	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052821\
Data File : BP005640.D
Acq On : 28 May 2021 15:48
Operator : CG/JU
Sample : M2532-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
MH-N-(MH-10.29)

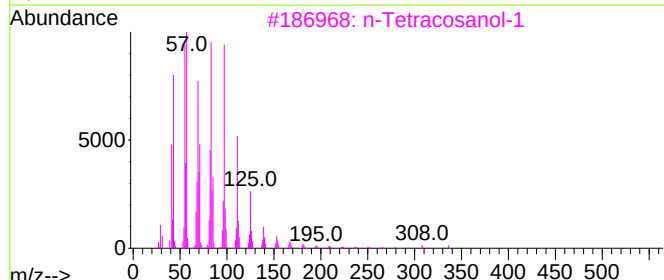
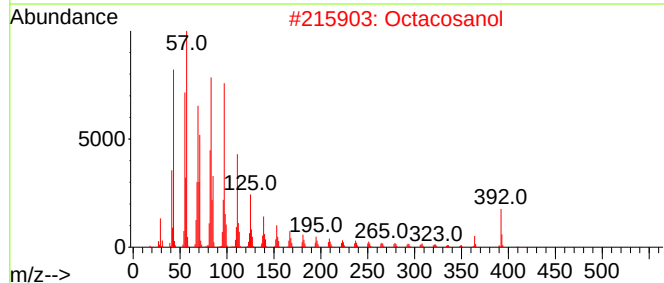
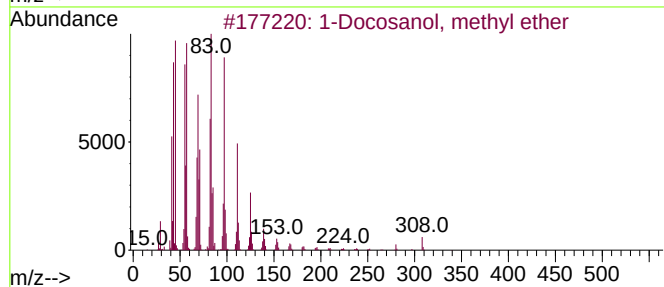
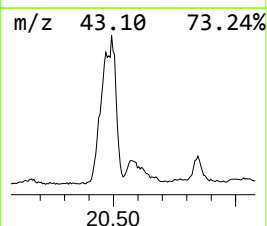
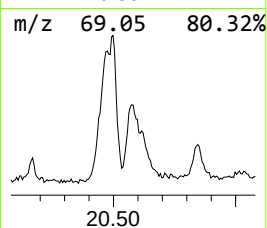
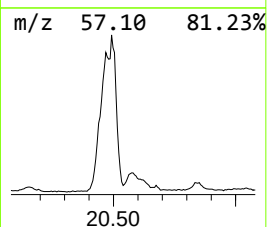
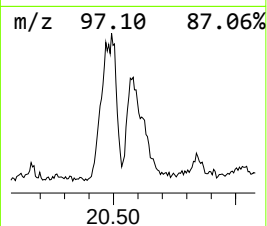
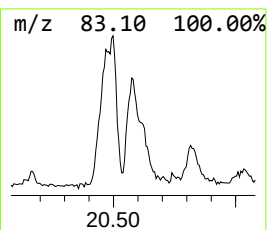
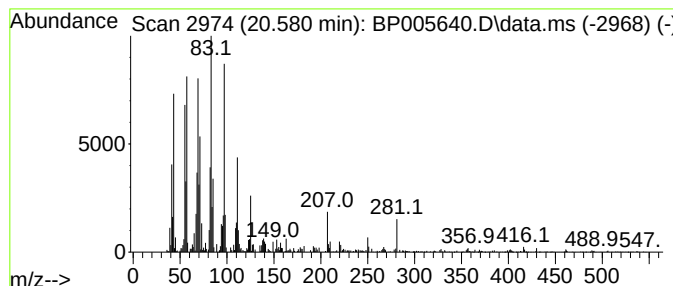
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 1-Docosanol, methyl ether Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.580	3.66 ng	476356	Chrysene-d12	21.416

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Docosanol, methyl ether			340	C23H48O	1000333-91-9	95
2	Octacosanol			410	C28H58O	000557-61-9	93
3	n-Tetracosanol-1			354	C24H50O	000506-51-4	93
4	Behenic alcohol			326	C22H46O	000661-19-8	87
5	1-Heptadecene			238	C17H34	006765-39-5	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052821\
Data File : BP005640.D
Acq On : 28 May 2021 15:48
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Misc :
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ClientSampleId :
MH-N-(MH-10.29)

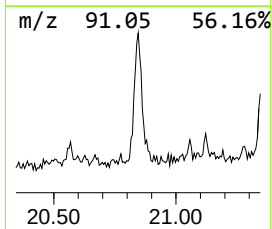
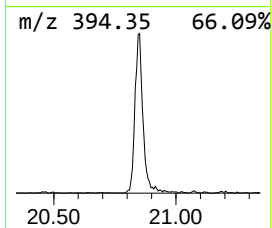
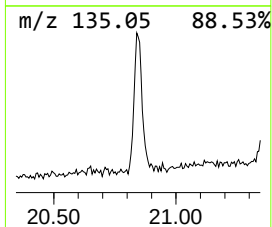
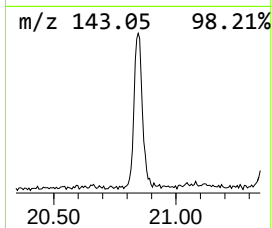
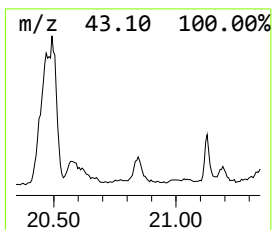
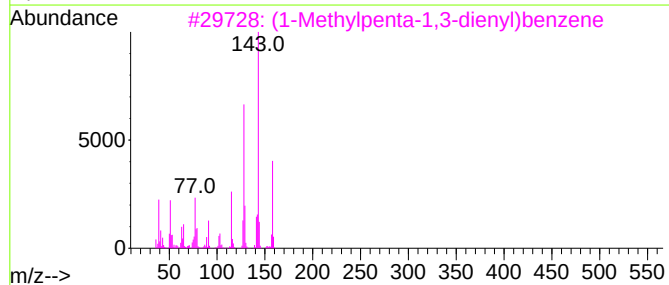
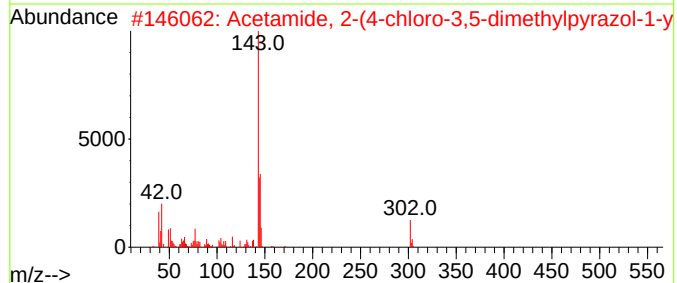
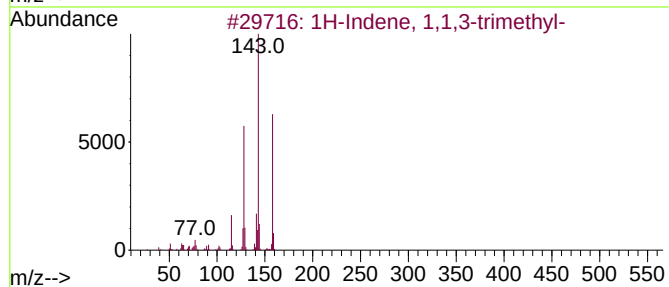
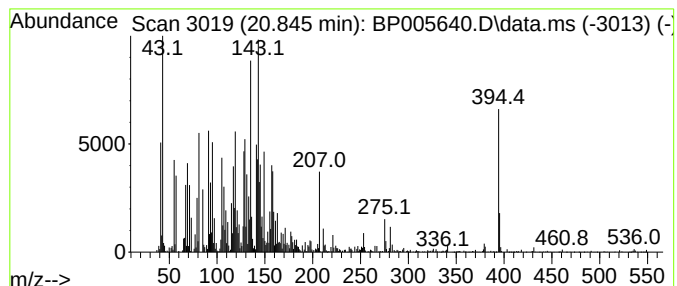
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 unknown20.84 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.845	5.41 ng	703575	Chrysene-d12	21.416

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 1,1,3-trimethyl-	158	C12H14	002177-45-9	18
2			Acetamide, 2-(4-chloro-3,5-dimet...	302	C15H15ClN4O	1000304-79-7	15
3			(1-Methylpenta-1,3-dienyl)benzene	158	C12H14	116669-49-9	14
4			(1-Methylenepent-2-enyl)benzene	158	C12H14	084313-24-6	14
5			Quinoline, 7-methyl-	143	C10H9N	000612-60-2	11



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052821\
Data File : BP005640.D
Acq On : 28 May 2021 15:48
Operator : CG/JU
Sample : M2532-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
MH-N-(MH-10.29)

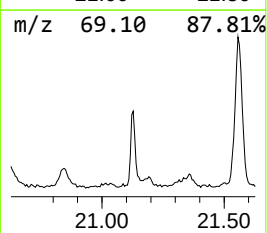
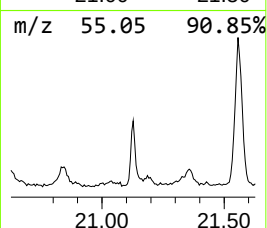
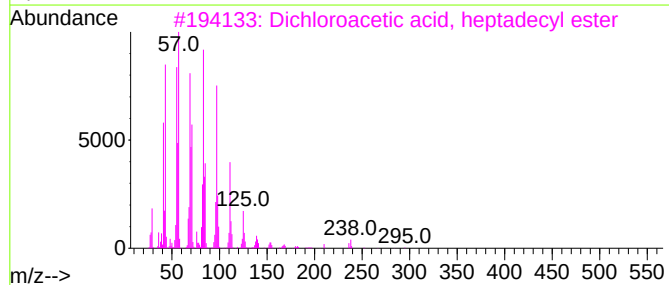
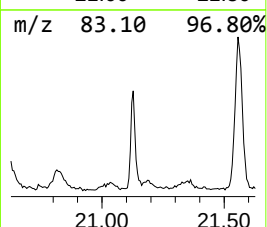
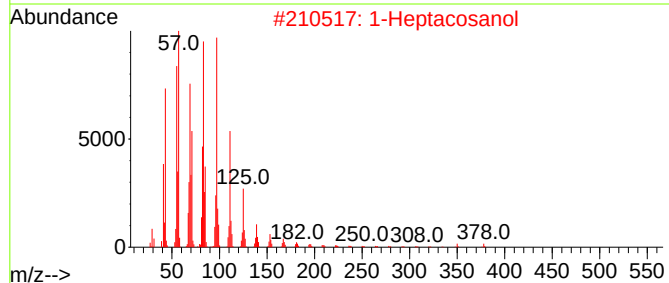
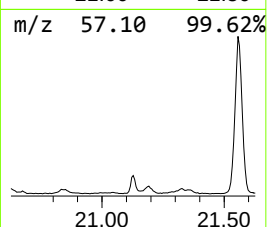
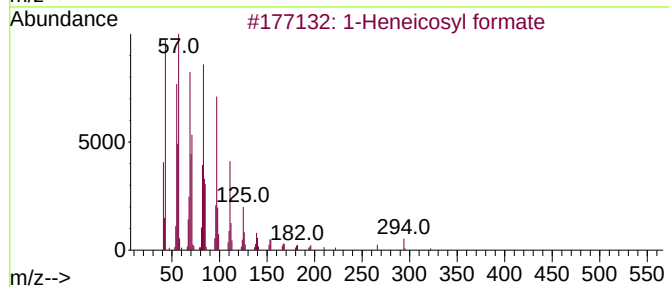
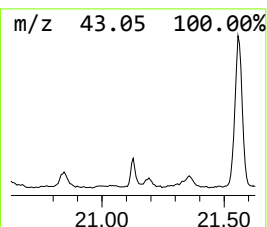
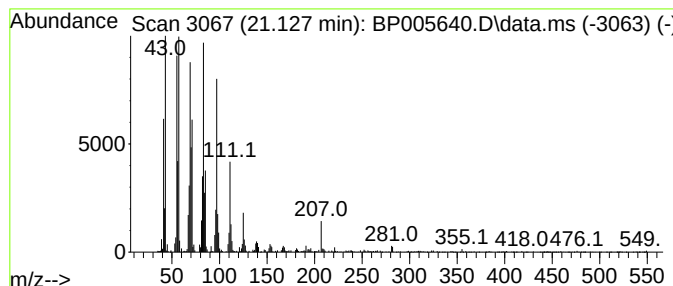
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 1-Heneicosyl formate Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.127	3.98 ng	517144	Chrysene-d12	21.416

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Heneicosyl formate	340	C22H44O2	077899-03-7	97
2			1-Heptacosanol	396	C27H56O	002004-39-9	94
3			Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	94
4			E-14-Hexadecenal	238	C16H30O	330207-53-9	93
5			Cyclopentadecane	210	C15H30	000295-48-7	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052821\
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Acq On : 28 May 2021 15:48
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Misc :
ALS Vial : 8 Sample Multiplier: 1

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ClientSampleId :
MH-N-(MH-10.29)

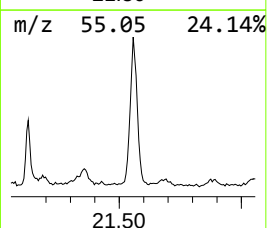
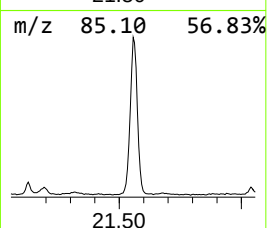
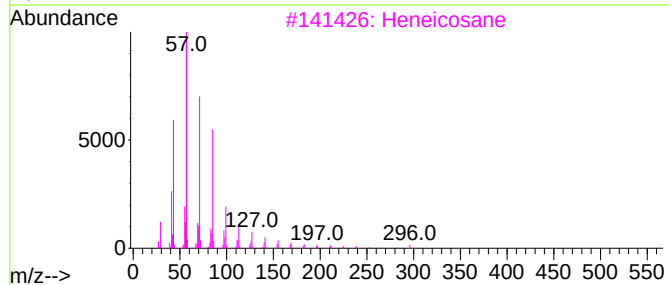
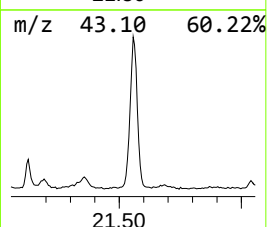
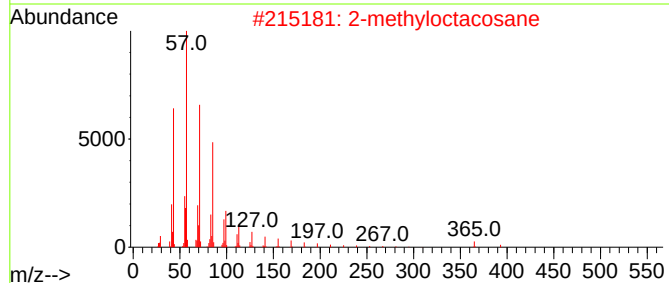
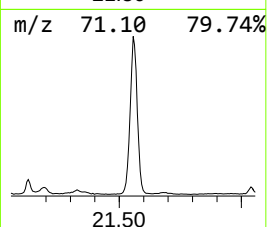
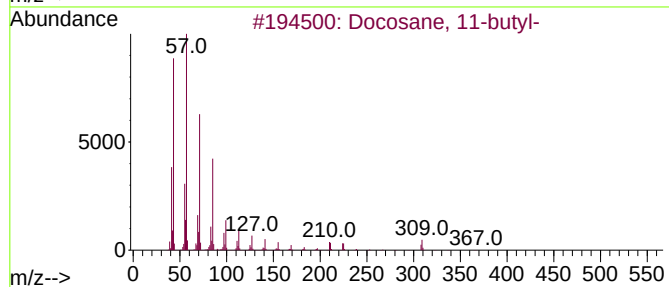
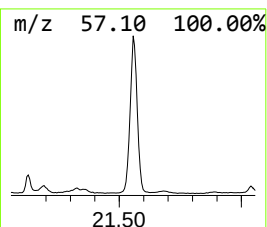
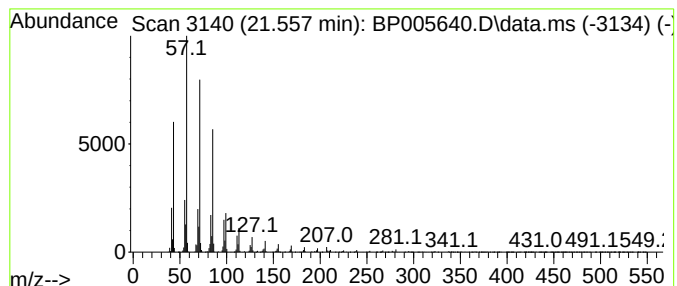
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Docosane, 11-butyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.557	23.53 ng	3060590	Chrysene-d12	21.416

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Docosane, 11-butyl-	366	C26H54	013475-76-8	95
2		2-methyloctacosane	408	C29H60	1000376-72-8	91
3		Heneicosane	296	C21H44	000629-94-7	91
4		Eicosane	282	C20H42	000112-95-8	91
5		Heptacosane	380	C27H56	000593-49-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052821\
Data File : BP005640.D
Acq On : 28 May 2021 15:48
Operator : CG/JU
Sample : M2532-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
MH-N-(MH-10.29)

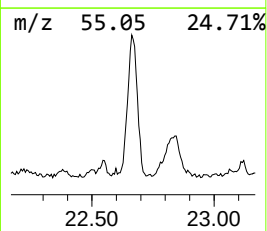
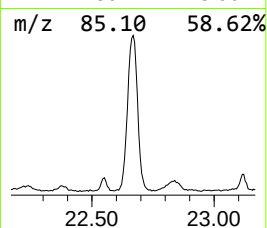
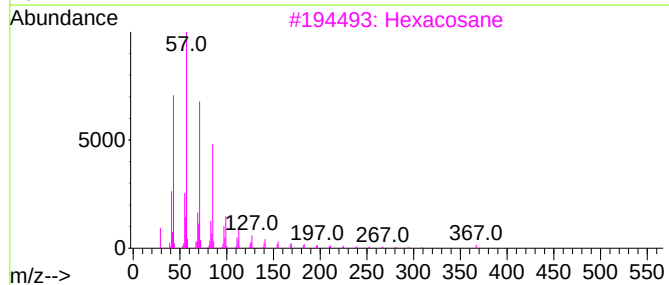
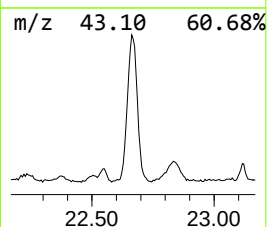
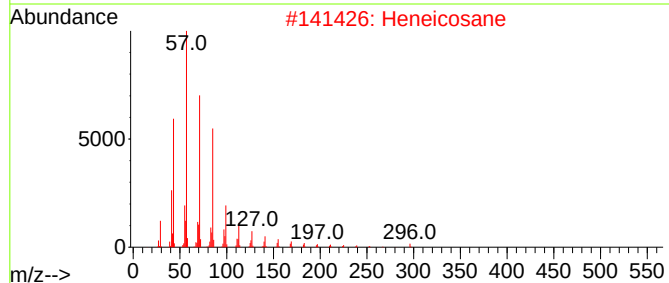
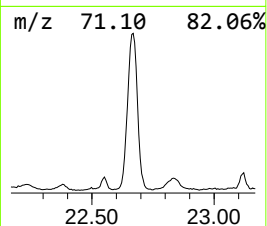
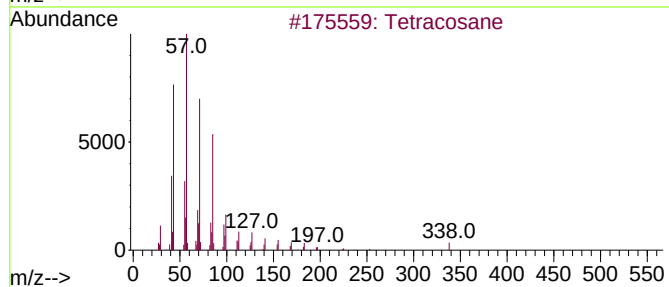
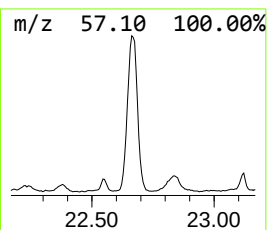
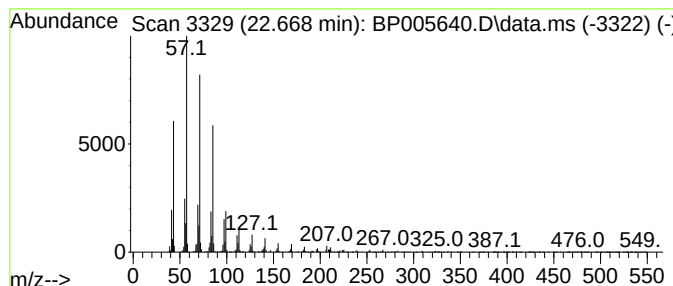
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Tetracosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.668	16.01 ng	2380310	Perylene-d12	23.863

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		Hexacosane	366	C26H54	000630-01-3	91
4		2-methyloctacosane	408	C29H60	1000376-72-8	91
5		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052821\
Data File : BP005640.D
Acq On : 28 May 2021 15:48
Operator : CG/JU
Sample : M2532-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
MH-N-(MH-10.29)

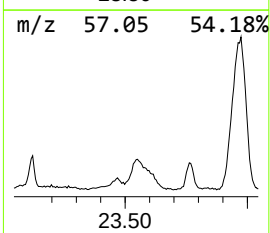
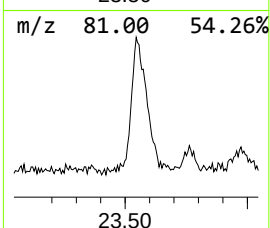
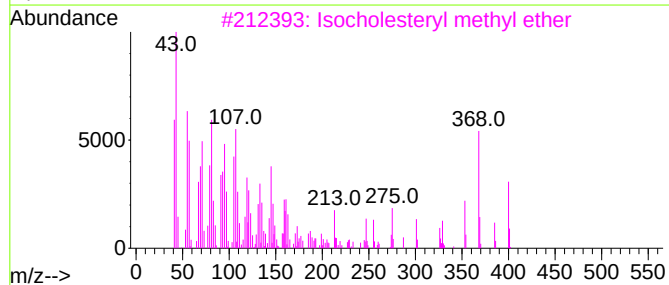
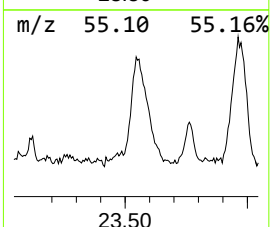
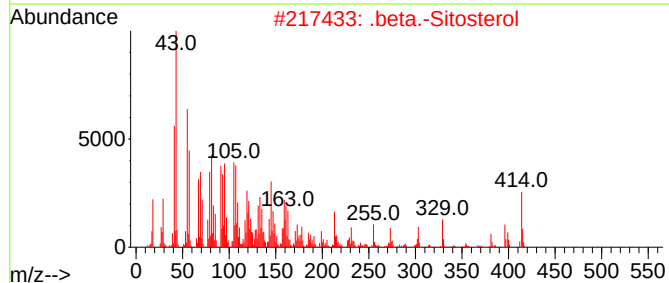
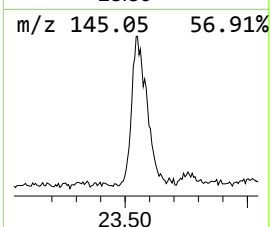
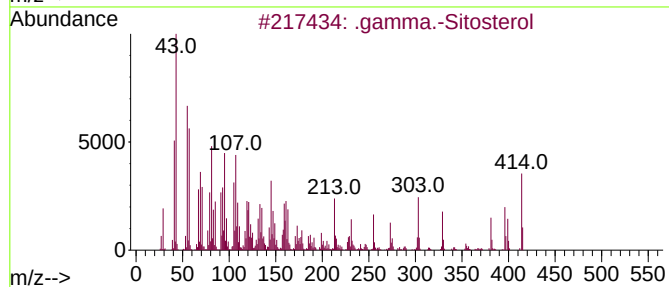
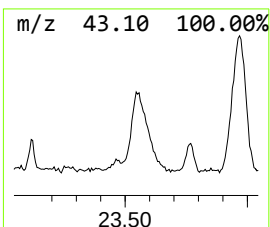
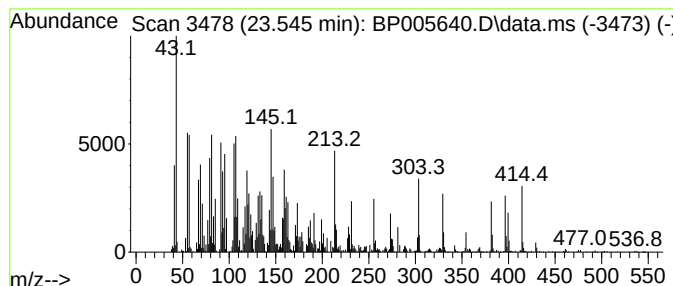
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 .gamma.-Sitosterol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.545	16.71 ng	2484550	Perylene-d12	23.863

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			.gamma.-Sitosterol	414	C29H50O	000083-47-6	93
2			.beta.-Sitosterol	414	C29H50O	000083-46-5	91
3			Isocholesteryl methyl ether	400	C28H48O	029944-53-4	30
4			17-(1,5-Dimethylhexyl)-10,13-dim...	414	C29H50O	1000210-86-9	20
5			Norflurazon	303	C12H9ClF3N3O	027314-13-2	15



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052821\
Data File : BP005640.D
Acq On : 28 May 2021 15:48
Operator : CG/JU
Sample : M2532-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
MH-N-(MH-10.29)

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP052721.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-metho...	3.111	95.5	ng	4959490	1	7.987	1038920	20.0
4-((1E)-3-Hydro...	16.733	3.2	ng	332416	4	17.351	2093160	20.0
n-Hexadecanoic ...	18.233	14.1	ng	1470000	4	17.351	2093160	20.0
3,5-Dimethoxy-4...	18.492	3.8	ng	402121	4	17.351	2093160	20.0
2-Pentanone, 1-...	18.539	10.7	ng	1119820	4	17.351	2093160	20.0
2-Bromo dodecane	18.704	33.1	ng	3469250	4	17.351	2093160	20.0
Octadecanoic acid	19.457	4.8	ng	625394	5	21.416	2600940	20.0
Docosane, 7-hexyl-	20.468	15.1	ng	1956610	5	21.416	2600940	20.0
1-Docosanol, me...	20.580	3.7	ng	476356	5	21.416	2600940	20.0
unknown20.84	20.845	5.4	ng	703575	5	21.416	2600940	20.0
1-Heneicosyl fo...	21.127	4.0	ng	517144	5	21.416	2600940	20.0
Docosane, 11-bu...	21.557	23.5	ng	3060590	5	21.416	2600940	20.0
Tetracosane	22.668	16.0	ng	2380310	6	23.863	2973010	20.0
.gamma.-Sitosterol	23.545	16.7	ng	2484550	6	23.863	2973010	20.0