

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040323\
 Data File : BM039279.D
 Acq On : 03 Apr 2023 15:06
 Operator : CG/JU
 Sample : 02161-04
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

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_M\Methods\8270-BM032923.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BM039279.D\data.ms

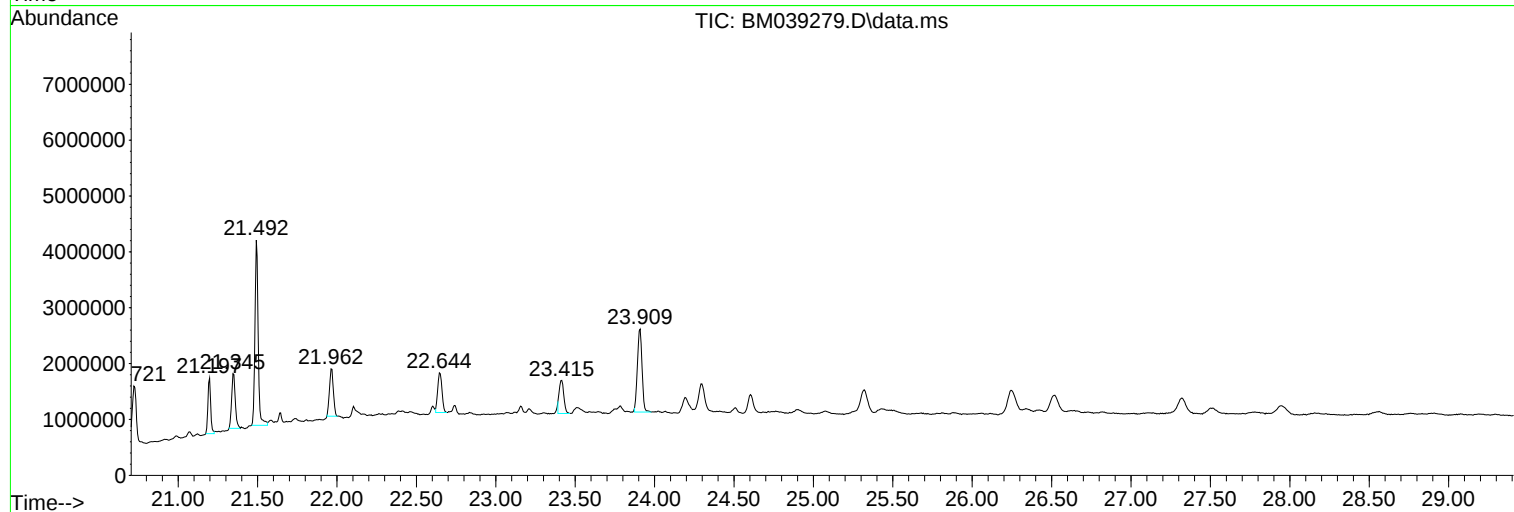
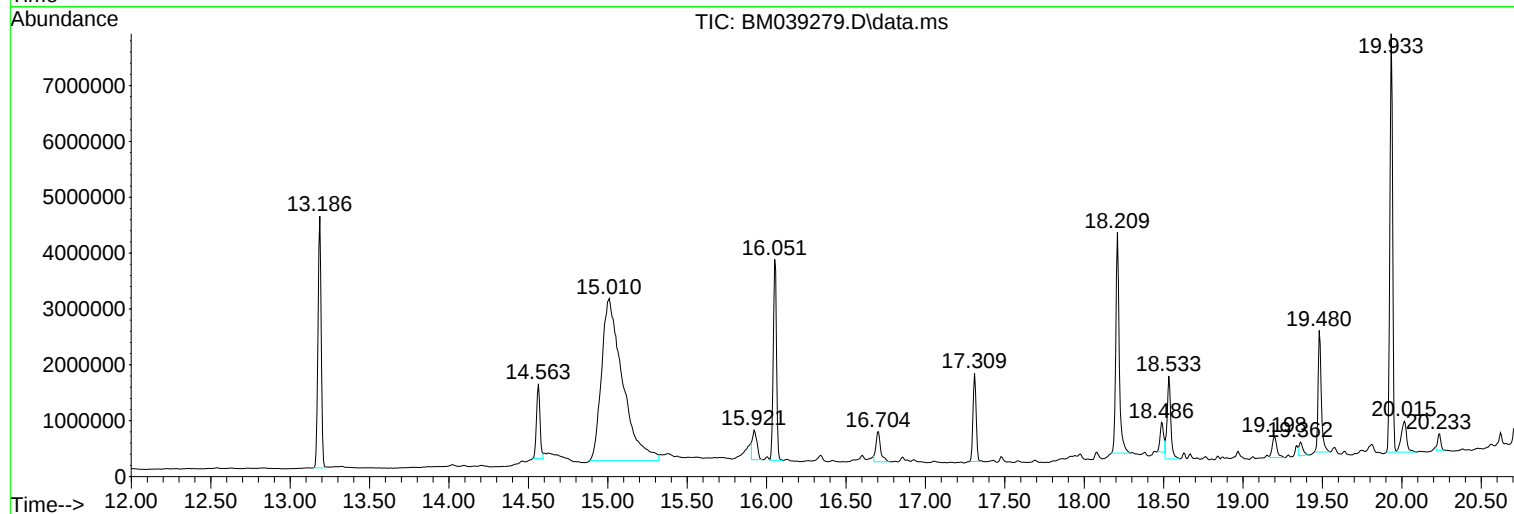
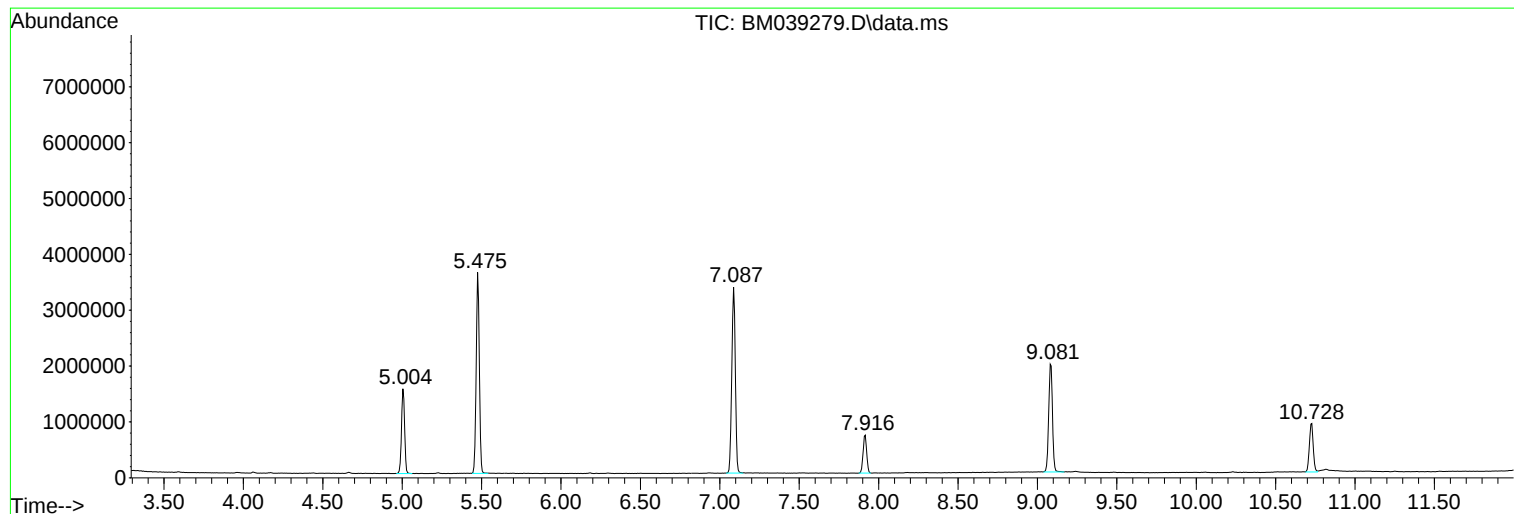
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.004	285	292	302	rBV	1516809	2057433	7.71%	2.008%
2	5.475	366	372	383	rBV	3597247	4991058	18.70%	4.871%
3	7.087	638	646	656	rBV	3328560	5067900	18.99%	4.946%
4	7.916	780	787	793	rBV	679984	1061858	3.98%	1.036%
5	9.081	978	985	1000	rVB	1930398	3181556	11.92%	3.105%
6	10.728	1258	1265	1271	rBV	867879	1469385	5.51%	1.434%
7	13.186	1676	1683	1689	rBV	4506349	6387084	23.93%	6.233%
8	14.563	1911	1917	1922	rBV	1335926	2009761	7.53%	1.961%
9	15.010	1972	1993	2046	rBV3	2900853	26688201	100.00%	26.044%
10	15.921	2145	2148	2157	rVB2	527160	1032549	3.87%	1.008%
11	16.051	2165	2170	2179	rVB	3596584	5116945	19.17%	4.993%
12	16.704	2276	2281	2293	rVB2	544611	1098827	4.12%	1.072%
13	17.309	2379	2384	2390	rBV	1579266	2130183	7.98%	2.079%
14	18.209	2531	2537	2551	rVB	3952099	5933097	22.23%	5.790%
15	18.486	2580	2584	2588	rBV	516389	866172	3.25%	0.845%
16	18.533	2588	2592	2604	rVB	1482295	2326206	8.72%	2.270%
17	19.198	2700	2705	2716	rVB	376246	773615	2.90%	0.755%
18	19.362	2731	2733	2742	rVB3	239419	379359	1.42%	0.370%
19	19.480	2749	2753	2765	rVB	2176969	3019616	11.31%	2.947%
20	19.933	2825	2830	2835	rBV	7490500	8972294	33.62%	8.756%
21	20.015	2837	2844	2857	rVB2	553733	1311973	4.92%	1.280%
22	20.233	2878	2881	2887	rVB	295351	413250	1.55%	0.403%
23	20.721	2959	2964	2971	rVB2	997812	1737593	6.51%	1.696%
24	21.197	3041	3045	3050	rBV	994120	1203782	4.51%	1.175%
25	21.345	3066	3070	3078	rVB	976790	1556265	5.83%	1.519%
26	21.492	3090	3095	3107	rBV2	3315301	4737748	17.75%	4.623%
27	21.962	3171	3175	3182	rVB	846886	1439926	5.40%	1.405%
28	22.644	3287	3291	3299	rVB	711388	1314550	4.93%	1.283%
29	23.415	3418	3422	3431	rVB	594031	1177955	4.41%	1.150%
30	23.909	3500	3506	3518	rVB2	1481121	3016899	11.30%	2.944%

Sum of corrected areas: 102473040

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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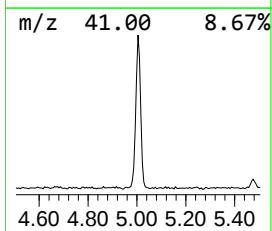
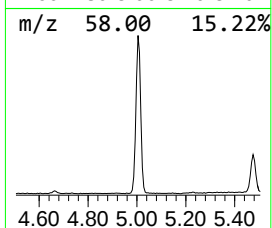
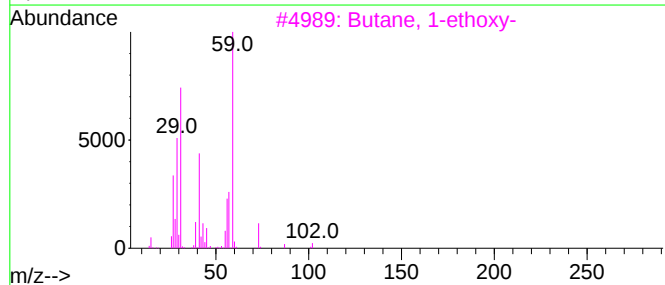
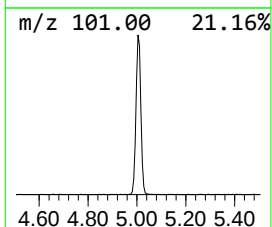
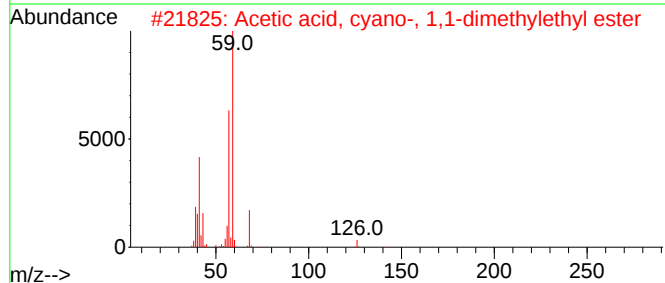
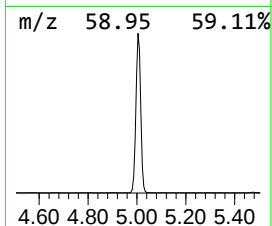
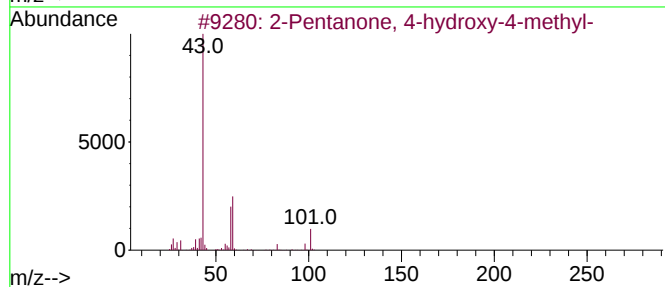
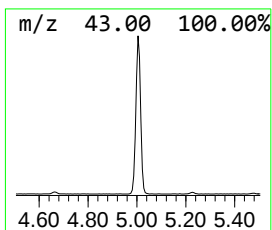
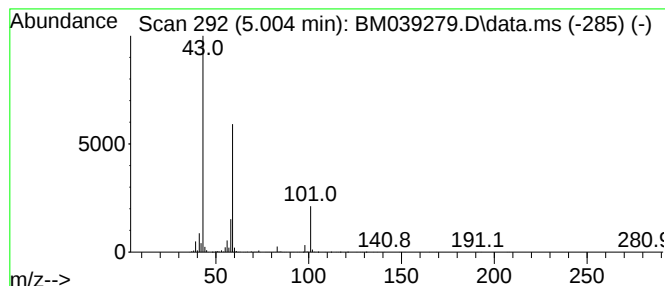
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TIC Library : C:\Database\NIST20.L
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.004	38.75 ng	2057430	1,4-Dichlorobenzene-d4	7.916

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
3		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9
4		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5		Formamide, N,N-diethyl-	101	C5H11NO	000617-84-5	9



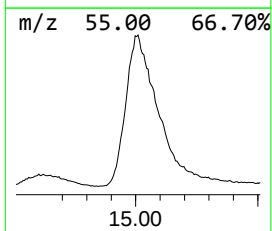
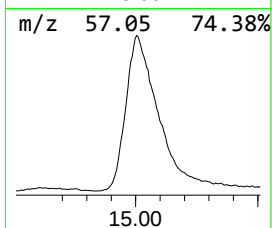
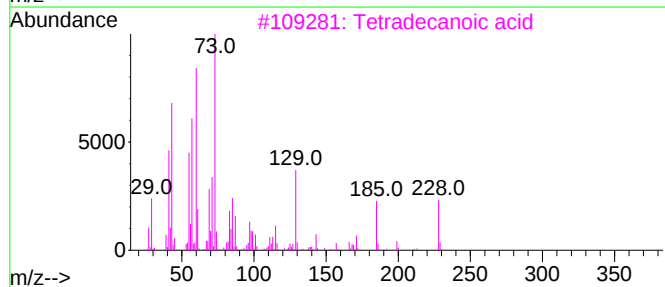
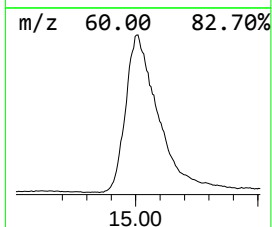
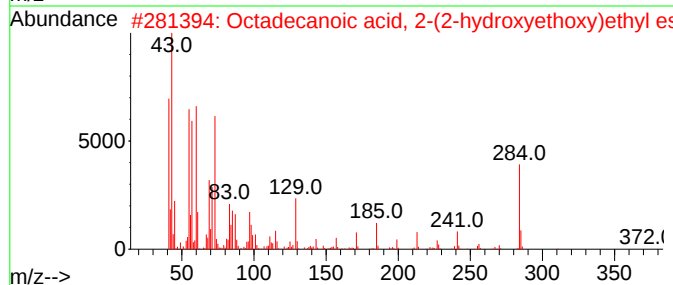
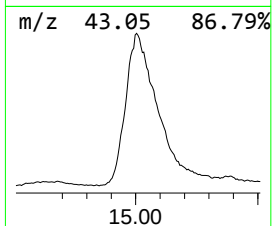
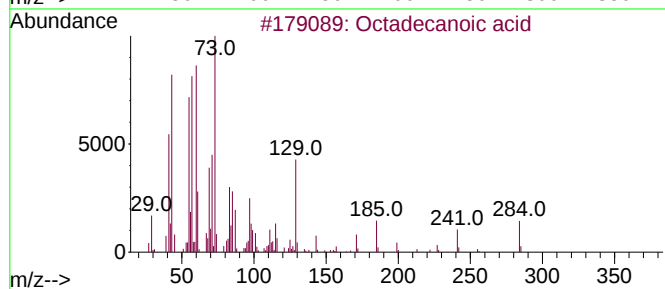
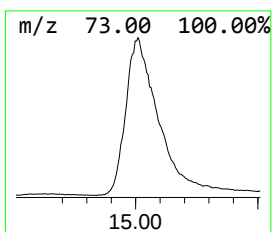
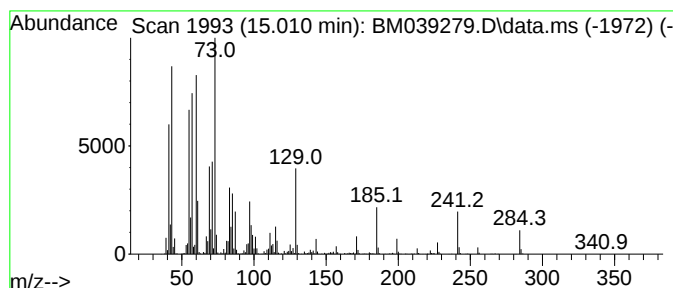
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Peak Number 2 Octadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.010	265.59 ng	26688200	Acenaphthene-d10	14.563	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2	Octadecanoic acid, 2-(2-hydroxye...	372	C22H44O4	000106-11-6	87
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	87
4	Pentadecanoic acid	242	C15H30O2	001002-84-2	86
5	n-Decanoic acid	172	C10H20O2	000334-48-5	62



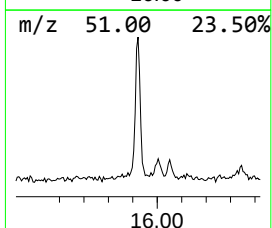
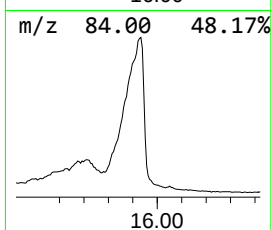
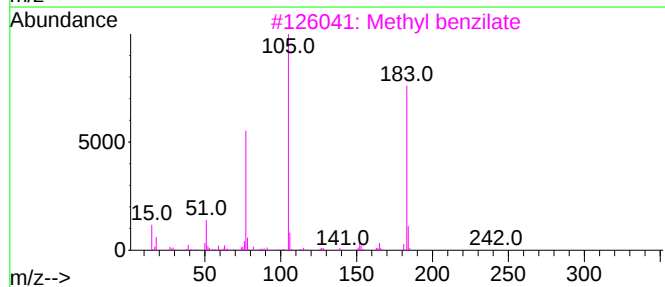
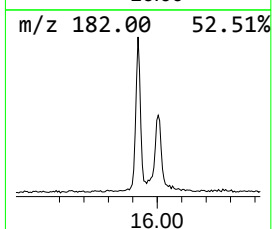
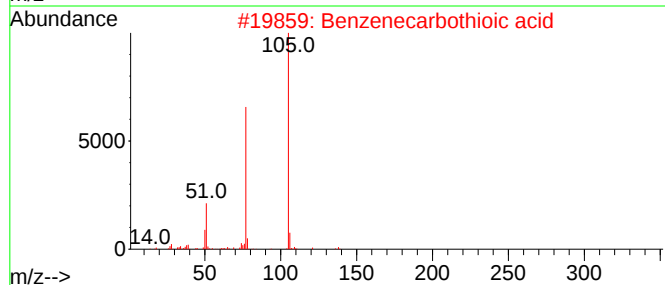
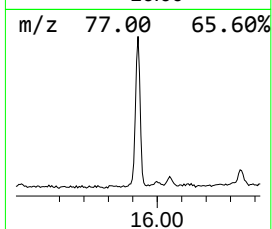
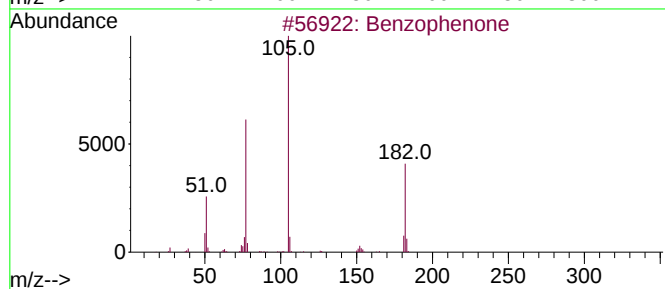
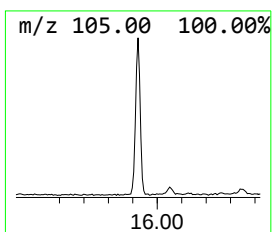
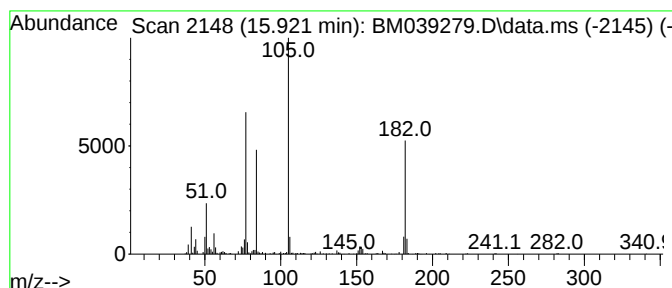
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
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Peak Number 3 Benzophenone Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.921	10.28 ng	1032550	Acenaphthene-d10	14.563	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzophenone	182	C13H10O	000119-61-9	96
2	Benzenecarbothioic acid	138	C7H6OS	000098-91-9	38
3	Methyl benzilate	242	C15H14O3	000076-89-1	38
4	Benzenecarbothioic acid, S-methy...	152	C8H8OS	005925-68-8	38
5	1,2-Propanedione, 1-phenyl-	148	C9H8O2	000579-07-7	35



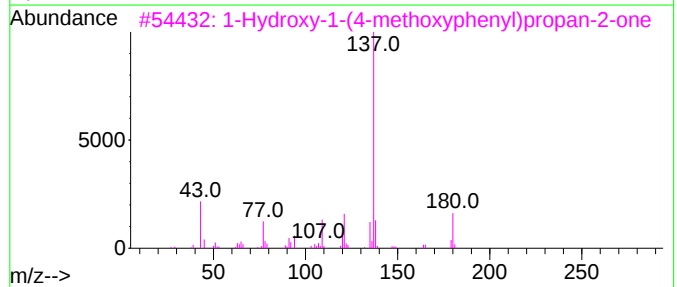
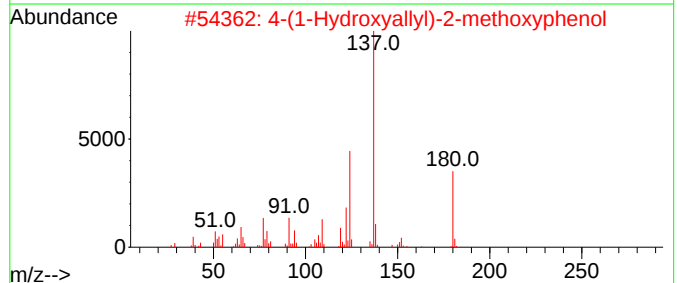
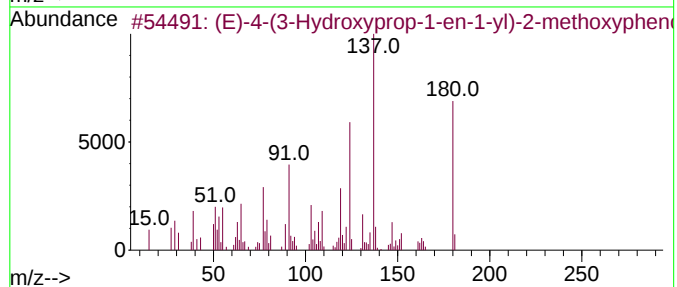
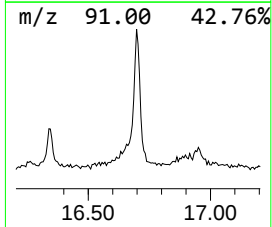
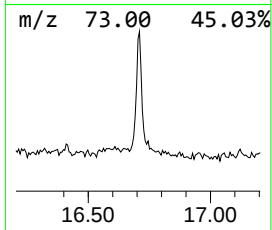
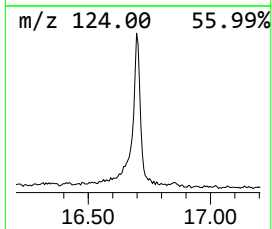
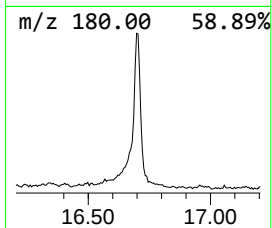
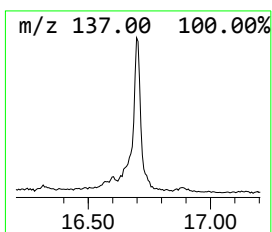
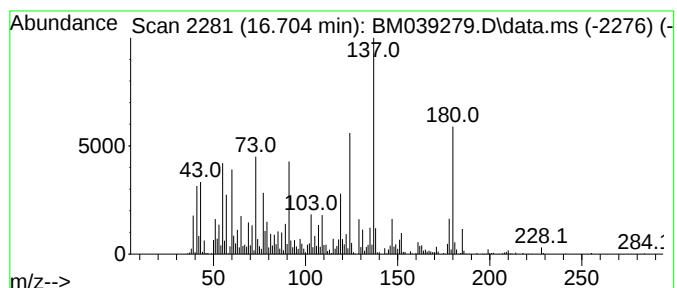
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Peak Number 4 (E)-4-(3-Hydroxyprop-1-en-1-yl)-2-methoxyphenol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
16.704	10.32 ng	1098830	Phenanthrene-d10	17.310	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(E)-4-(3-Hydroxyprop-1-en-1-yl)-...	180	C10H12O3	032811-40-8	99
2	4-(1-Hydroxyallyl)-2-methoxyphenol	180	C10H12O3	112465-50-6	58
3	1-Hydroxy-1-(4-methoxyphenyl)pro...	180	C10H12O3	015482-29-8	38
4	Methyl 3-(4-hydroxy-3-methoxyphe...	210	C11H14O4	056024-44-3	38
5	(-)-R-Phenethanamine, 1-methyl-N...	271	C17H21NO2	1000127-90-4	35



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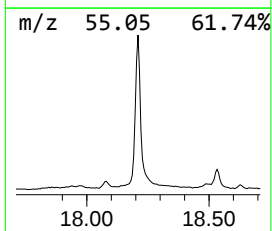
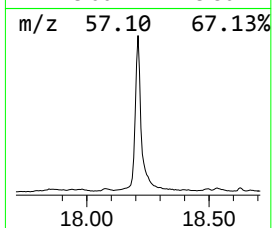
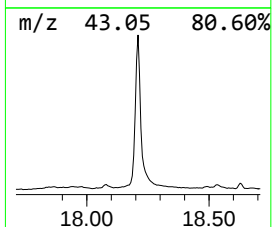
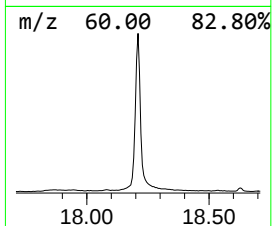
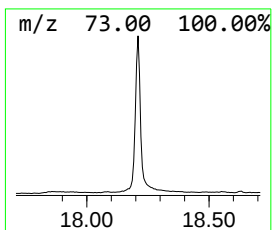
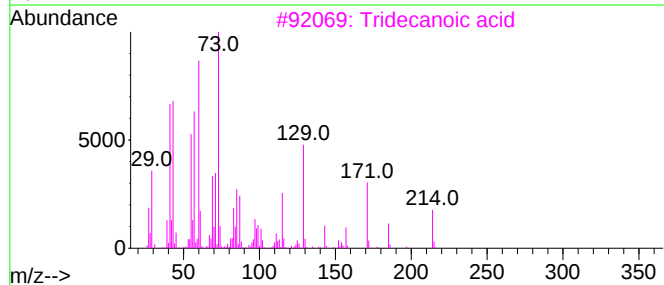
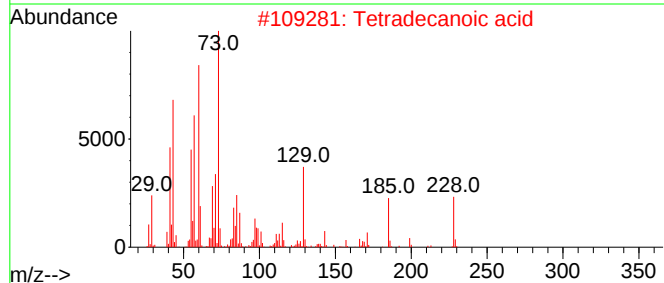
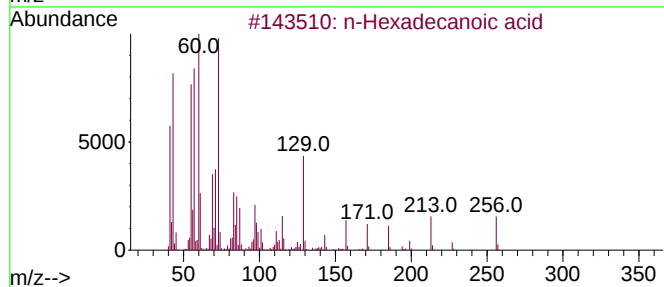
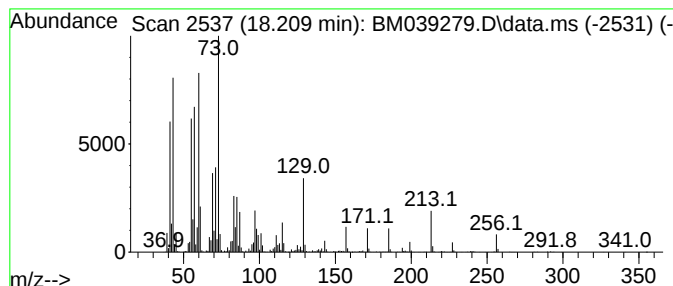
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 5 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.209	55.70 ng	5933100	Phenanthrene-d10	17.310

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2			Tetradecanoic acid	228	C14H28O2	000544-63-8	93
3			Tridecanoic acid	214	C13H26O2	000638-53-9	92
4			Pentadecanoic acid	242	C15H30O2	001002-84-2	87
5			Octadecanoic acid	284	C18H36O2	000057-11-4	83



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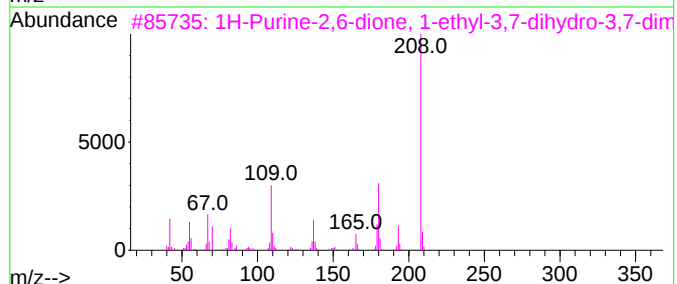
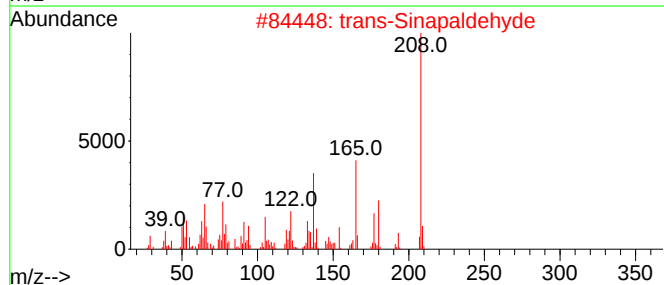
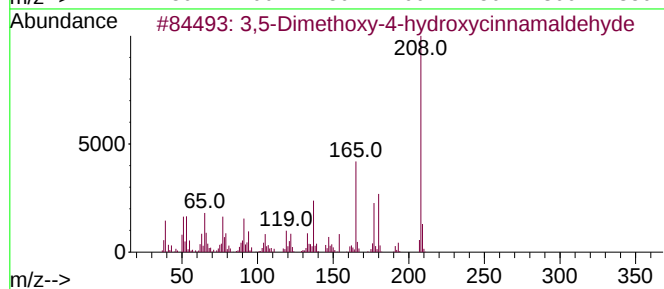
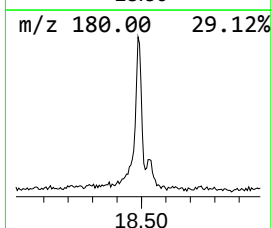
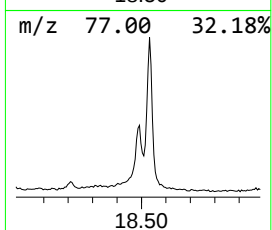
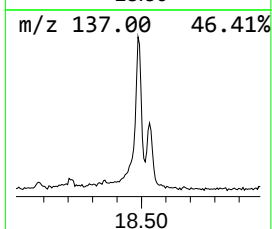
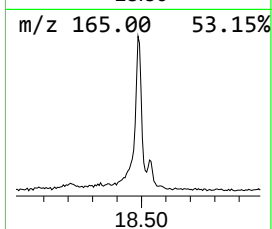
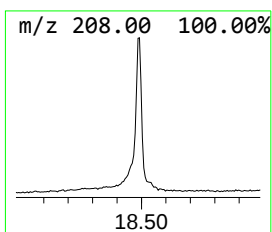
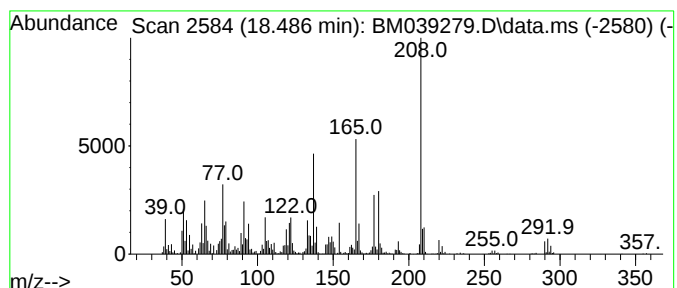
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 3,5-Dimethoxy-4-hydroxycinn... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.486	8.13 ng	866172	Phenanthrene-d10	17.310

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		3,5-Dimethoxy-4-hydroxycinnamal... 208	C11H12O4		087345-53-7	96
2		trans-Sinapaldehyde 208	C11H12O4		004206-58-0	94
3		1H-Purine-2,6-dione, 1-ethyl-3,7... 208	C9H12N4O2		039832-36-5	43
4		3-tert-Butyl-4-hydroxyanisole 180	C11H16O2		000121-00-6	42
5		1,2-Dimethoxy-4-(3-methoxy-1-pro... 208	C12H16O3		1000313-35-0	41



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Acq On : 03 Apr 2023 15:06
Operator : CG/JU
Sample : 02161-04
Misc :
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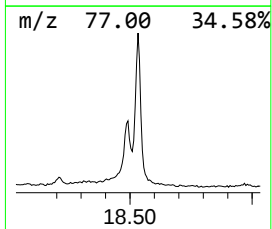
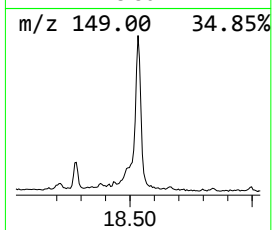
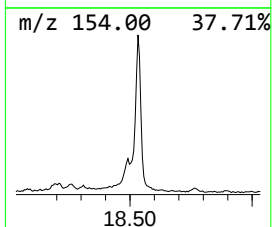
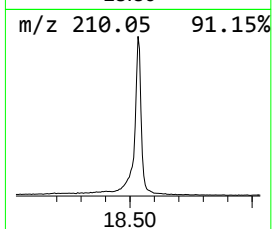
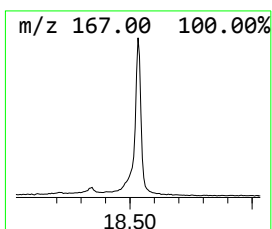
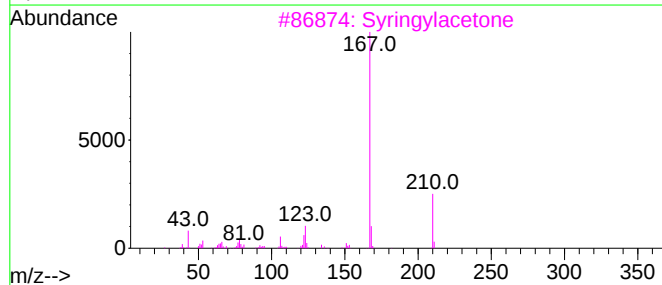
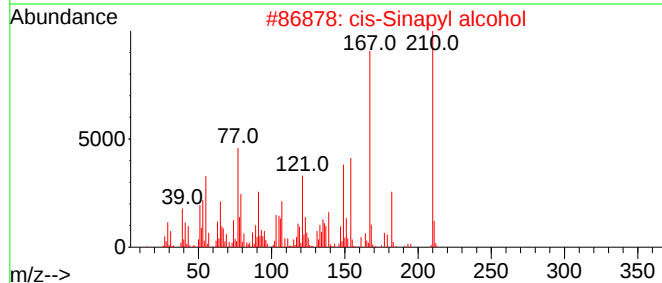
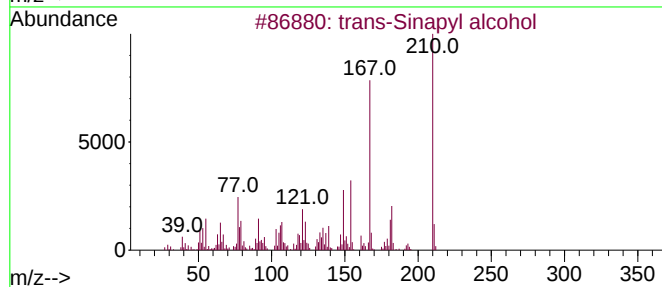
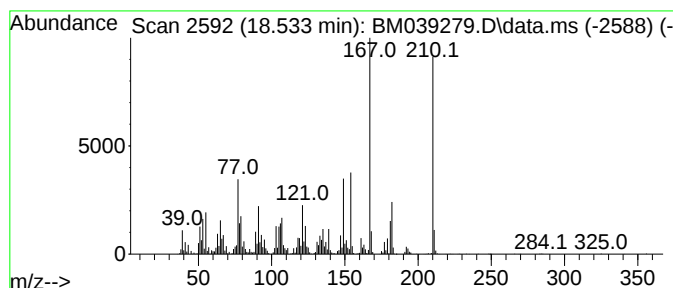
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 trans-Sinapyl alcohol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.533	21.84 ng	2326210	Phenanthrene-d10	17.310

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		trans-Sinapyl alcohol	210	C11H14O4	020675-96-1	96
2		cis-Sinapyl alcohol	210	C11H14O4	104330-63-4	93
3		Syringylacetone	210	C11H14O4	019037-58-2	64
4		4-Ethyl-2,6-dimethoxyphenol	182	C10H14O3	014059-92-8	62
5		1-Butanone, 1-(2,4,6-trihydroxy-...	210	C11H14O4	001509-06-4	52



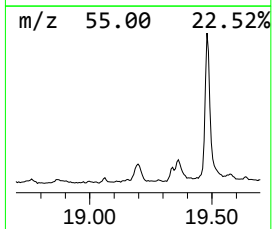
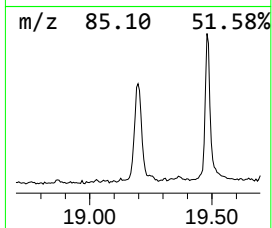
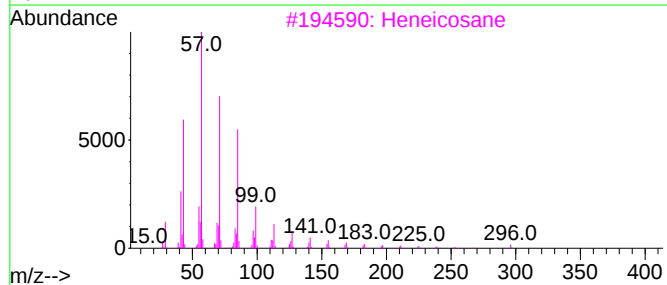
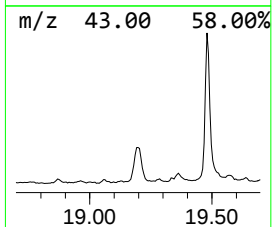
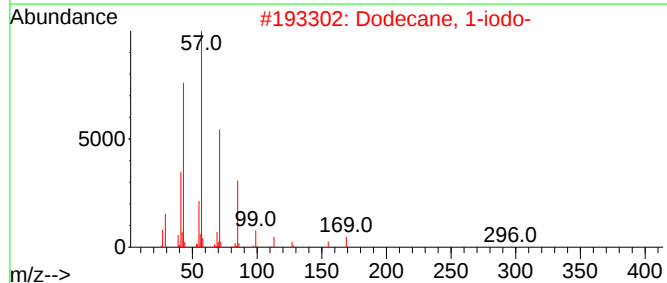
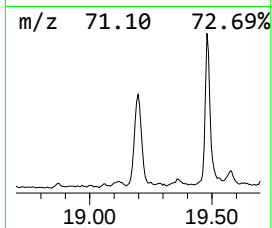
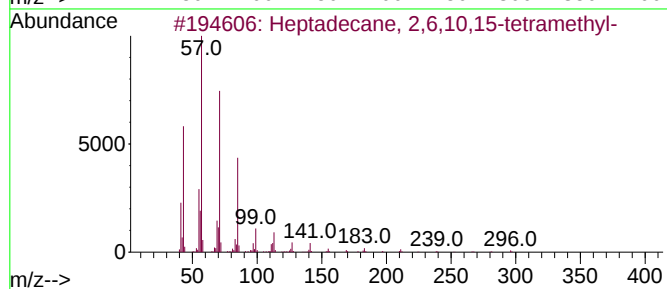
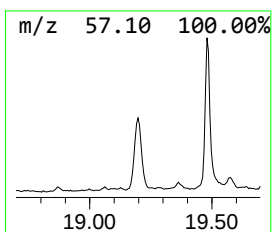
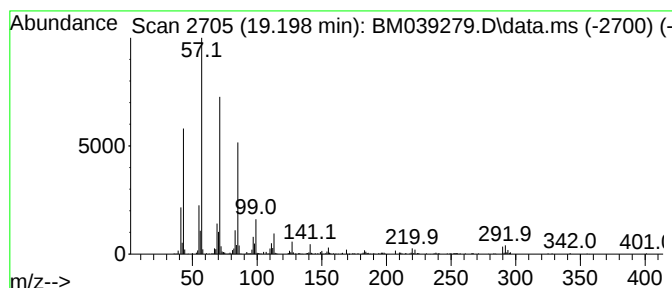
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Acq On : 03 Apr 2023 15:06
Operator : CG/JU
Sample : 02161-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

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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 Heptadecane, 2,6,10,15-tetr... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD		R.T.	
19.198	7.26 ng	773615	Phenanthrene-d10		17.310	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Heptadecane, 2,6,10,15-tetramethyl-		296	C21H44	054833-48-6	91
2	Dodecane, 1-iodo-		296	C12H25I	004292-19-7	87
3	Heneicosane		296	C21H44	000629-94-7	83
4	Dodecane, 2-methyl-		184	C13H28	001560-97-0	83
5	Octadecane, 1-iodo-		380	C18H37I	000629-93-6	80



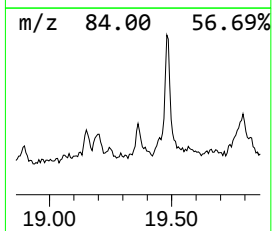
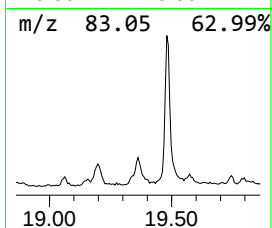
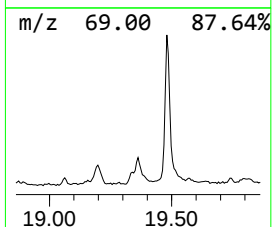
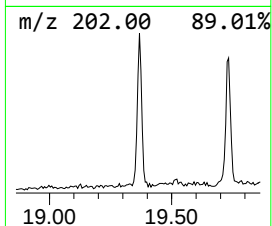
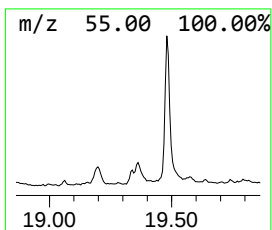
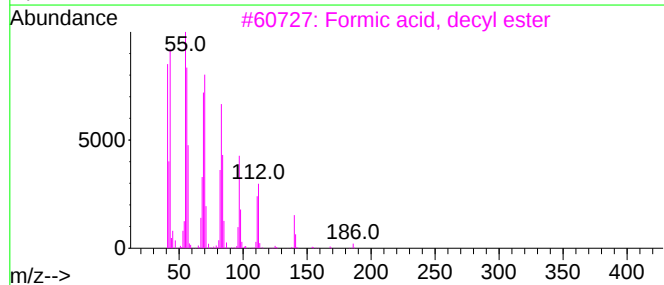
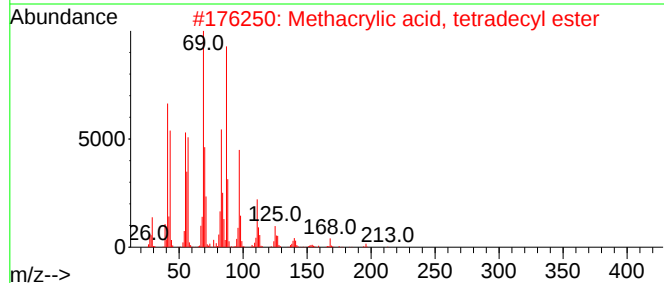
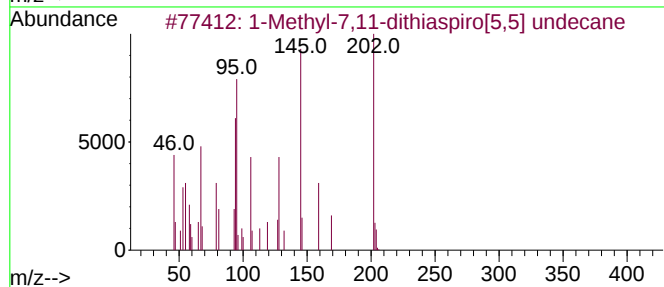
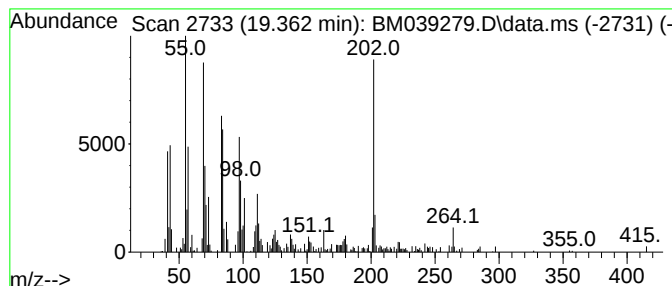
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Misc :
ALS Vial : 8 Sample Multiplier: 1

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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 1-Methyl-7,11-dithiaspiro[5... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.362	3.56 ng	379359	Phenanthrene-d10	17.310	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Methyl-7,11-dithiaspiro[5,5] u...	202	C10H18S2	107678-22-8	53
2	Methacrylic acid, tetradecyl ester	282	C18H34O2	1000340-29-0	46
3	Formic acid, decyl ester	186	C11H22O2	005451-52-5	25
4	1-Hexadecanol	242	C16H34O	036653-82-4	25
5	4-Octenoic acid, 2,3,7-trimethyl...	198	C12H22O2	1000196-85-3	20



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040323\
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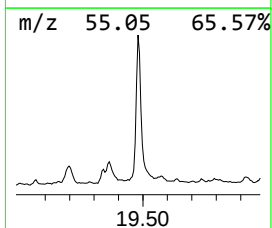
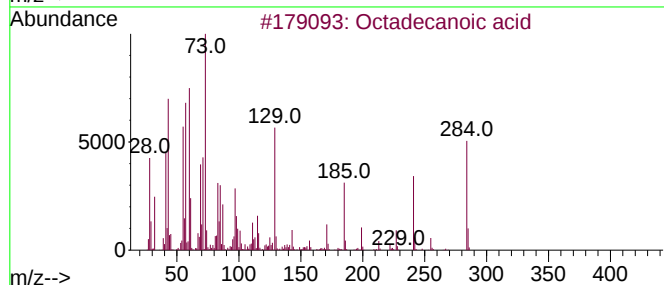
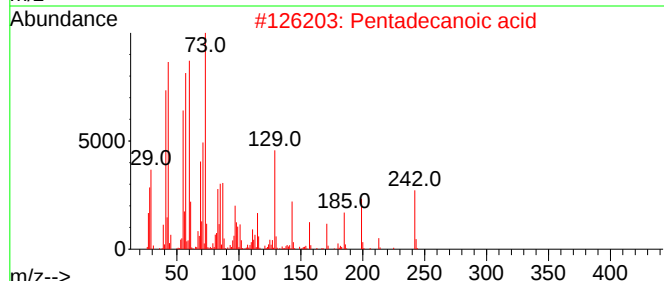
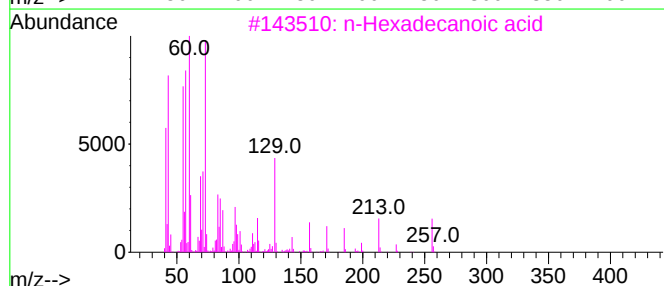
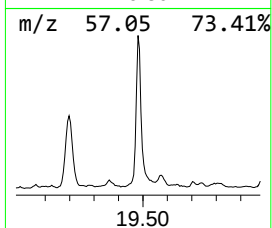
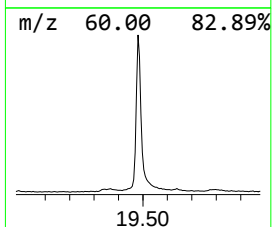
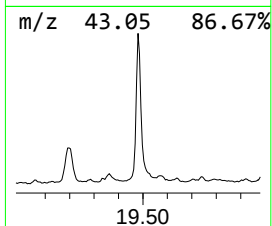
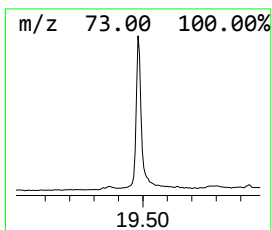
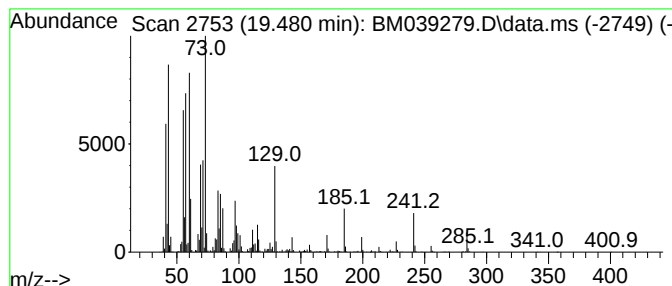
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Pentadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.480	12.75 ng	3019620	Chrysene-d12	21.492

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	89
3		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	87
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	87
5		n-Decanoic acid	172	C10H20O2	000334-48-5	68



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040323\
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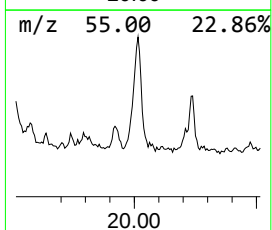
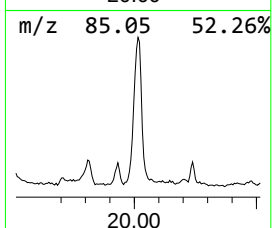
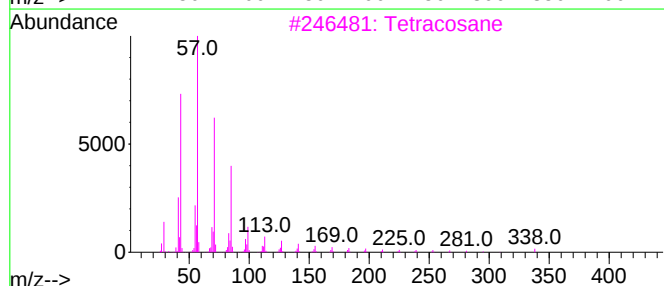
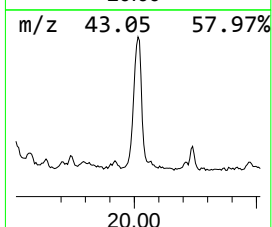
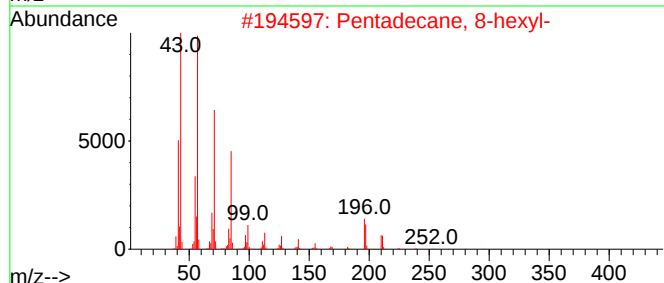
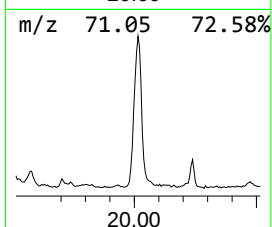
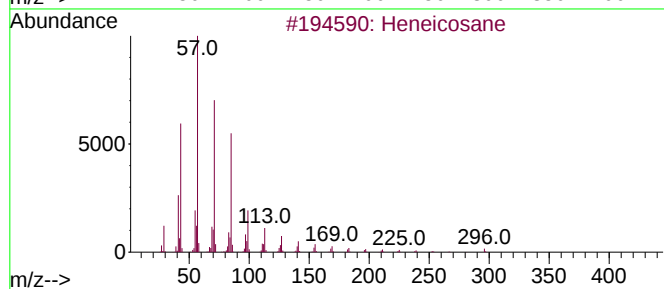
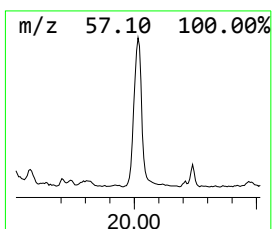
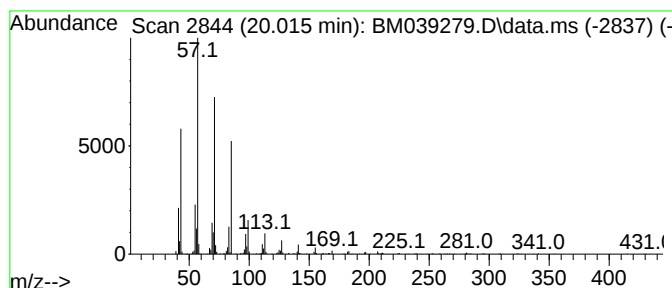
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Heneicosane Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.015	5.54 ng	1311970	Chrysene-d12	21.492

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	91
2			Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	90
3			Tetracosane	338	C24H50	000646-31-1	90
4			Tricosane	324	C23H48	000638-67-5	87
5			Octadecane, 1-iodo-	380	C18H37I	000629-93-6	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040323\
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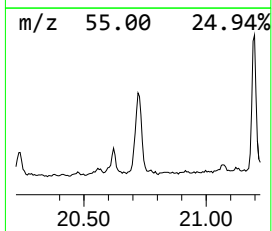
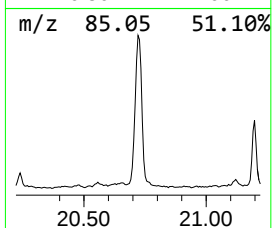
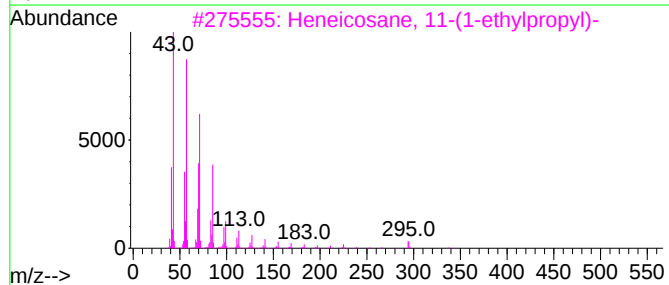
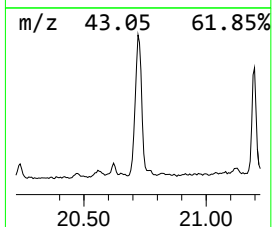
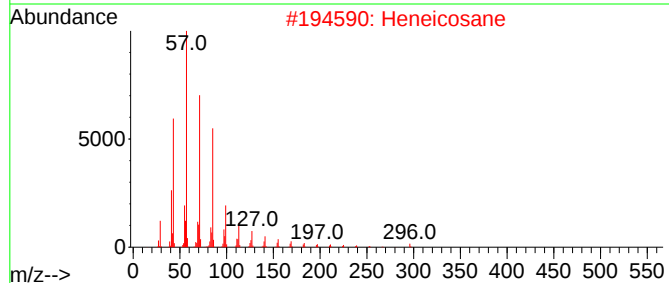
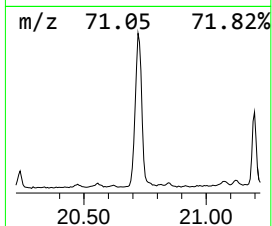
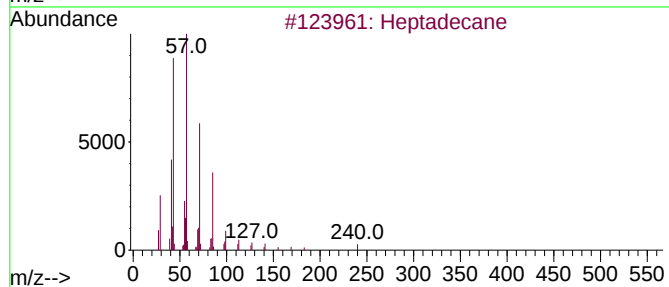
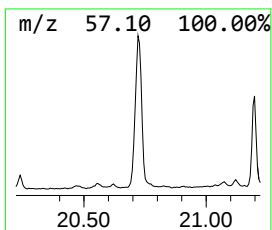
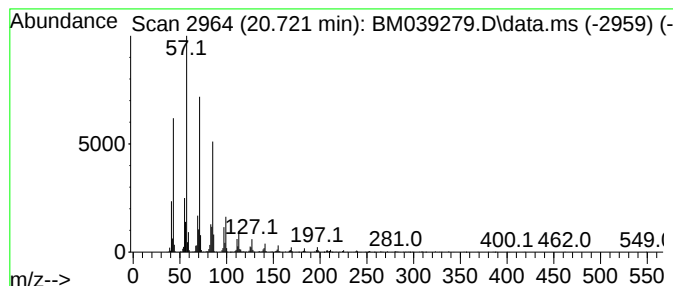
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 Heptadecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.721	7.34 ng	1737590	Chrysene-d12	21.492

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane	240	C17H36	000629-78-7	94
2		Heneicosane	296	C21H44	000629-94-7	90
3		Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	90
4		Hexacosane	366	C26H54	000630-01-3	90
5		Octacosane	394	C28H58	000630-02-4	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040323\
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ALS Vial : 8 Sample Multiplier: 1

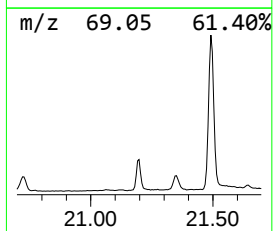
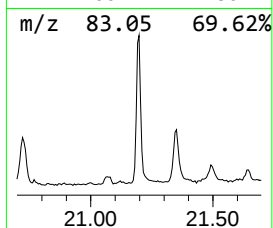
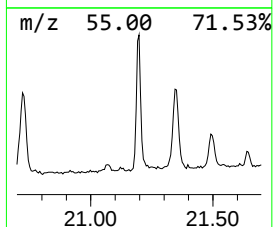
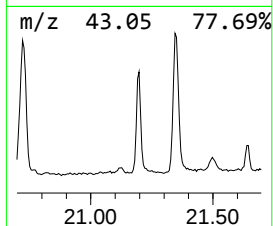
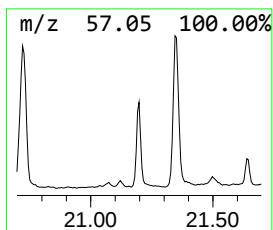
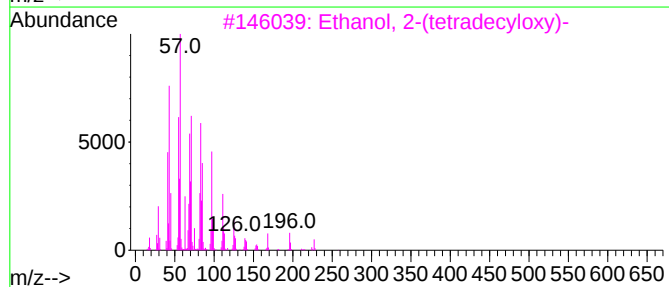
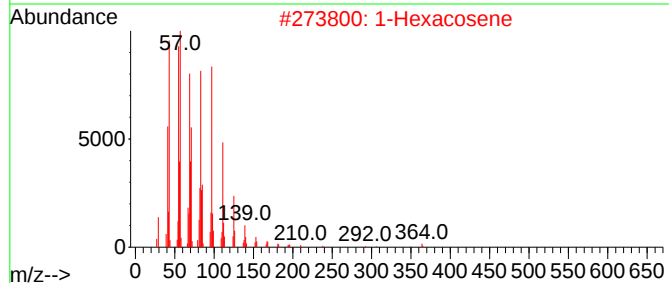
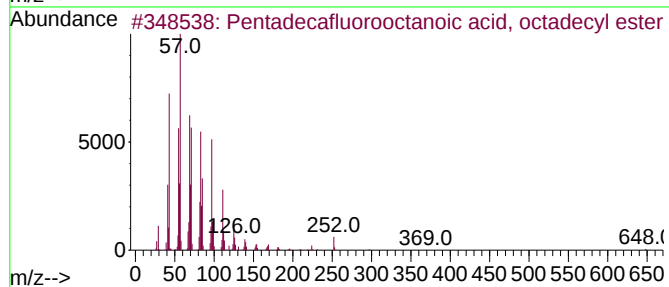
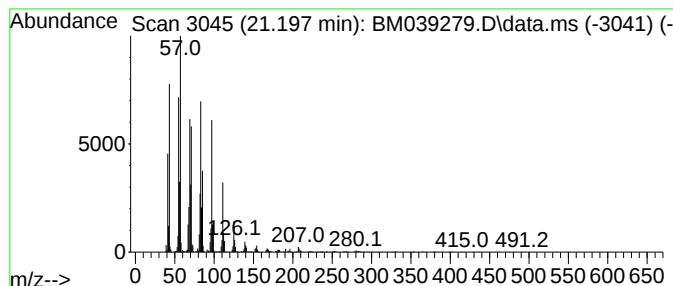
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Pentadecafluorooctanoic aci... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.198	5.08 ng	1203780	Chrysene-d12	21.492

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	94
2			1-Hexacosene	364	C26H52	018835-33-1	94
3			Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	93
4			Nonadecyl pentafluoropropionate	430	C22H39F5O2	1000351-88-8	91
5			Heptafluorobutyric acid, n-octad...	466	C22H37F7O2	000400-57-7	91



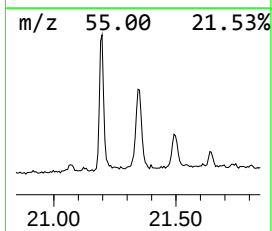
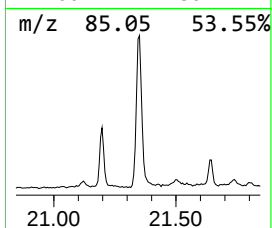
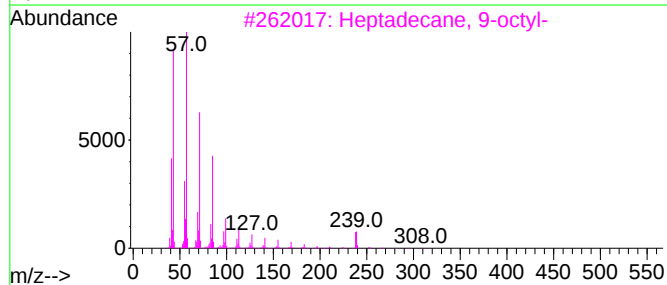
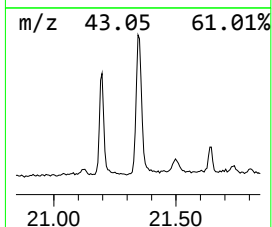
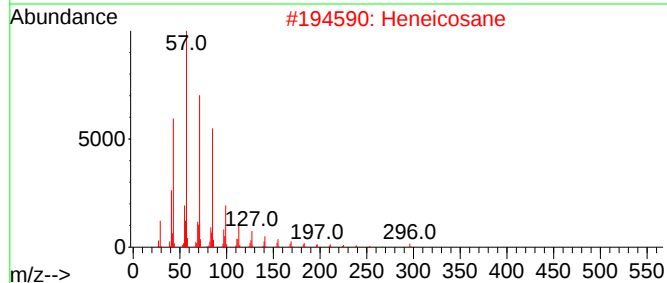
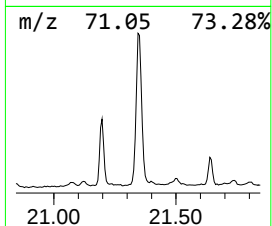
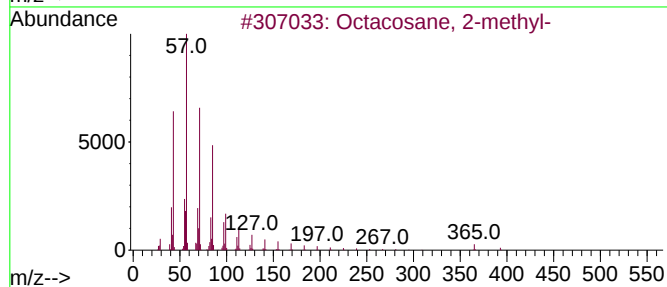
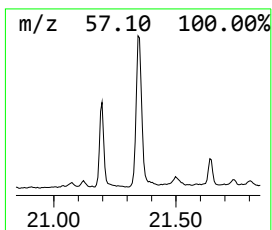
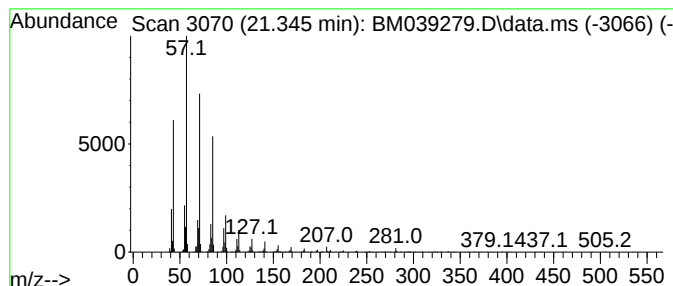
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Acq On : 03 Apr 2023 15:06
Operator : CG/JU
Sample : 02161-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM032923.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Octacosane, 2-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD		R.T.	
21.345	6.57 ng	1556270	Chrysene-d12		21.492	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Octacosane, 2-methyl-		408	C29H60	001560-98-1	91
2	Heneicosane		296	C21H44	000629-94-7	91
3	Heptadecane, 9-octyl-		352	C25H52	007225-64-1	91
4	Docosane, 1-iodo-		436	C22H45I	1000406-31-9	90
5	Octadecane, 1-iodo-		380	C18H37I	000629-93-6	87



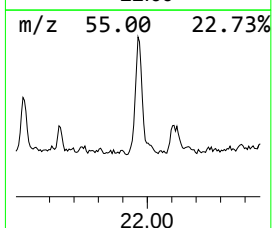
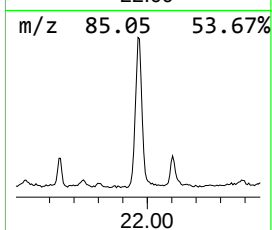
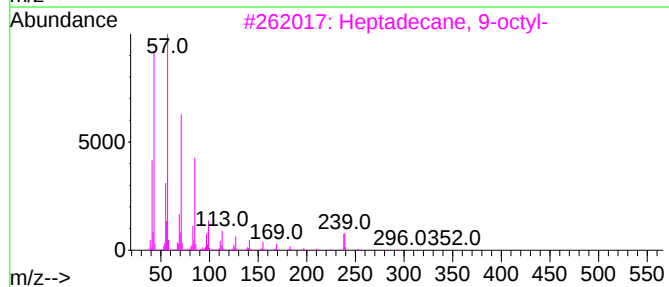
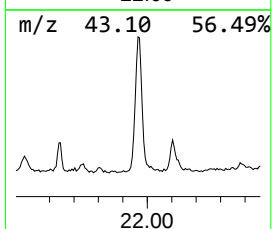
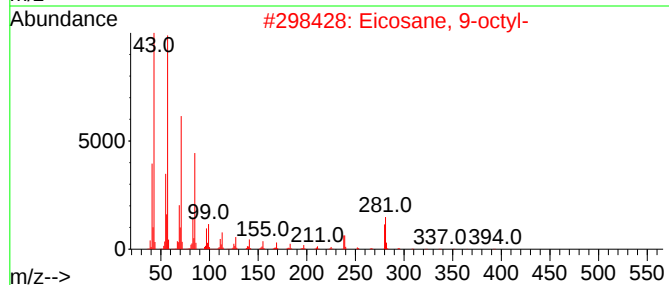
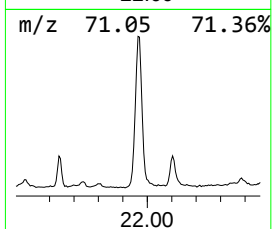
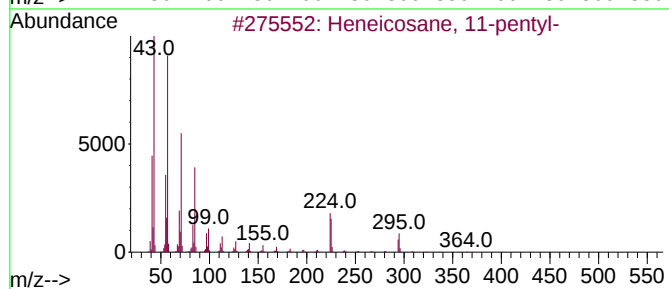
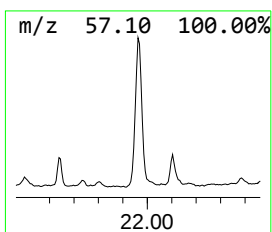
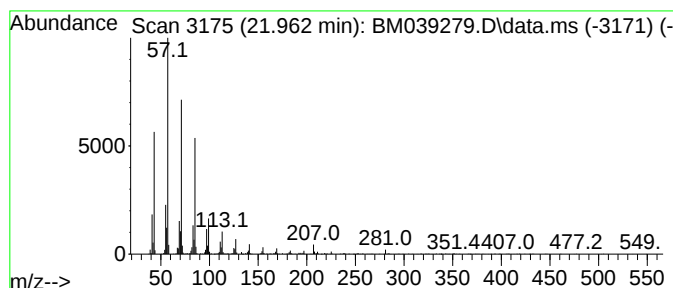
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Operator : CG/JU
Sample : 02161-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM032923.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Heneicosane, 11-pentyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.962	6.08 ng	1439930	Chrysene-d12	21.492	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane, 11-pentyl-	366	C26H54	014739-72-1	93
2	Eicosane, 9-octyl-	394	C28H58	013475-77-9	93
3	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
4	Octacosane	394	C28H58	000630-02-4	90
5	Heneicosane	296	C21H44	000629-94-7	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040323\
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Acq On : 03 Apr 2023 15:06
Operator : CG/JU
Sample : 02161-04
Misc :
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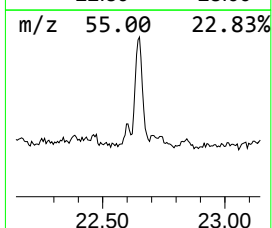
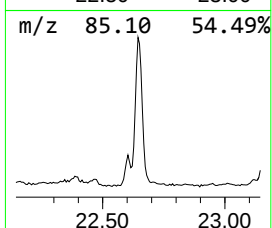
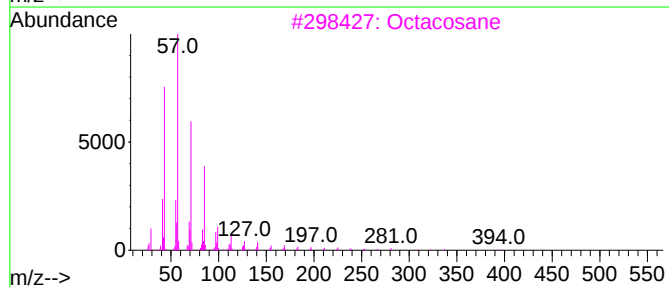
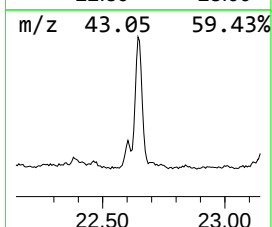
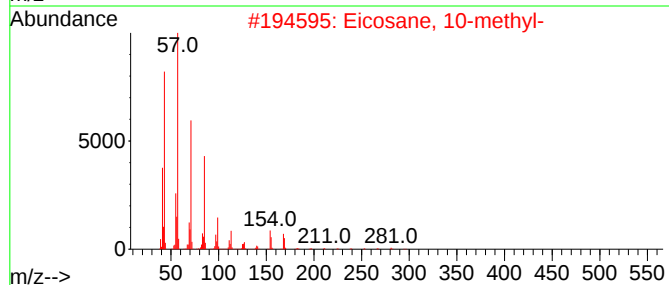
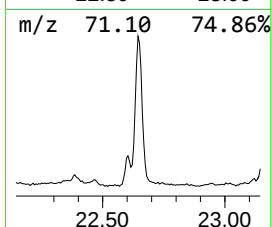
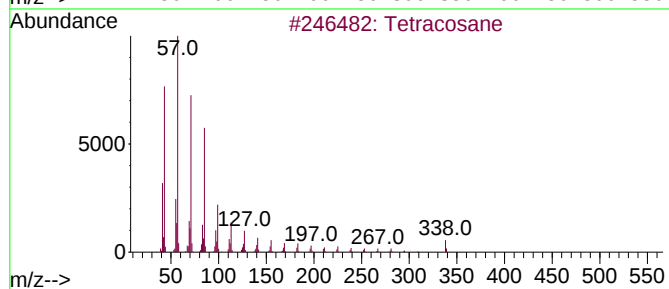
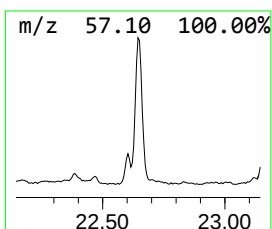
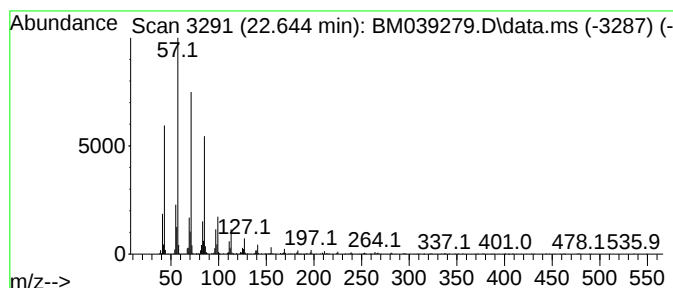
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Tetracosane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.644	5.55 ng	1314550	Chrysene-d12	21.492

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetracosane	338	C24H50	000646-31-1	97
2		Eicosane, 10-methyl-	296	C21H44	054833-23-7	94
3		Octacosane	394	C28H58	000630-02-4	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_M\Data\BM040323\
Data File : BM039279.D
Acq On : 03 Apr 2023 15:06
Operator : CG/JU
Sample : 02161-04
Misc :
ALS Vial : 8 Sample Multiplier: 1

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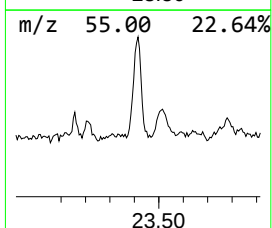
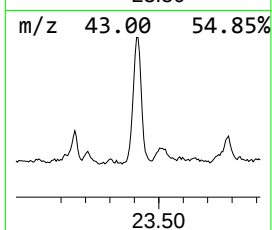
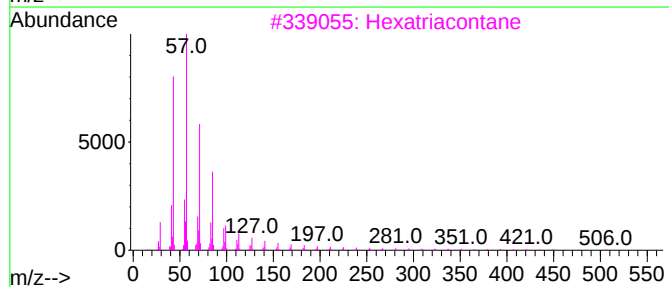
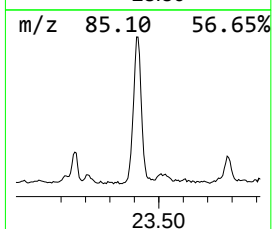
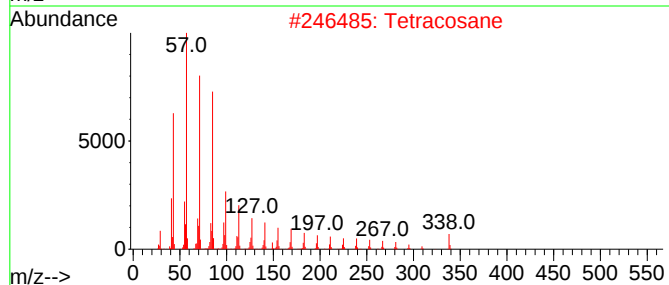
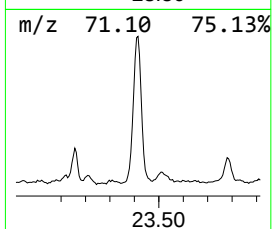
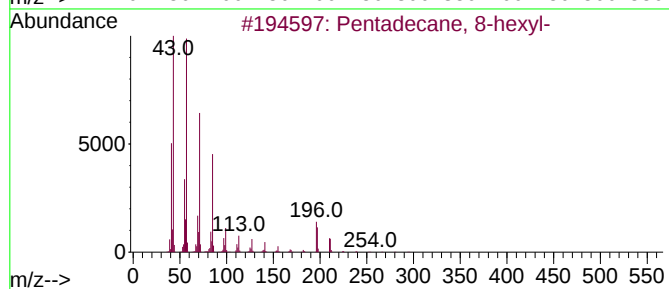
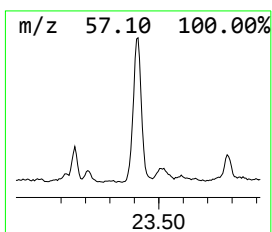
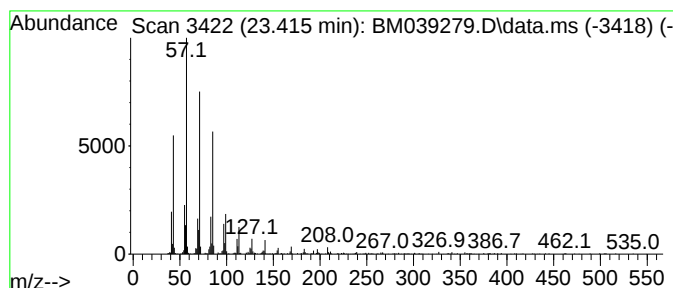
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 17 Pentadecane, 8-hexyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.415	7.81 ng	1177960	Perylene-d12	23.909

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	93
2		Tetracosane	338	C24H50	000646-31-1	91
3		Hexatriacontane	507	C36H74	000630-06-8	91
4		Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
5		Tetratetracontane	619	C44H90	007098-22-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040323\
 Data File : BM039279.D
 Acq On : 03 Apr 2023 15:06
 Operator : CG/JU
 Sample : 02161-04
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM032923.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
2-Pentanone, 4-...	5.004	38.8	ng	2057430	1	7.916	1061860	20.0
Octadecanoic acid	15.010	265.6	ng	26688200	3	14.563	2009760	20.0
Benzophenone	15.921	10.3	ng	1032550	3	14.563	2009760	20.0
(E)-4-(3-Hydrox...	16.704	10.3	ng	1098830	4	17.310	2130180	20.0
n-Hexadecanoic ...	18.209	55.7	ng	5933100	4	17.310	2130180	20.0
3,5-Dimethoxy-4...	18.486	8.1	ng	866172	4	17.310	2130180	20.0
trans-Sinapyl a...	18.533	21.8	ng	2326210	4	17.310	2130180	20.0
Heptadecane, 2,...	19.198	7.3	ng	773615	4	17.310	2130180	20.0
1-Methyl-7,11-d...	19.362	3.6	ng	379359	4	17.310	2130180	20.0
Pentadecanoic acid	19.480	12.8	ng	3019620	5	21.492	4737750	20.0
Heneicosane	20.015	5.5	ng	1311970	5	21.492	4737750	20.0
Heptadecane	20.721	7.3	ng	1737590	5	21.492	4737750	20.0
Pentadecafluoro...	21.198	5.1	ng	1203780	5	21.492	4737750	20.0
Octacosane, 2-m...	21.345	6.6	ng	1556270	5	21.492	4737750	20.0
Heneicosane, 11...	21.962	6.1	ng	1439930	5	21.492	4737750	20.0
Tetracosane	22.644	5.5	ng	1314550	5	21.492	4737750	20.0
Pentadecane, 8-...	23.415	7.8	ng	1177960	6	23.909	3016900	20.0