

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP022224\
 Data File : BP019820.D
 Acq On : 22 Feb 2024 12:42
 Operator : MA/JU
 Sample : P1462-01
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 20240762-39N-N-FIELD-BLANK-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_P\Methods\8270E-BP013124.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP019820.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.464	398	404	413	rVB	246579	367042	2.63%	1.091%
2	7.063	669	676	687	rBV	199311	353194	2.53%	1.050%
3	7.887	810	816	824	rVB	189779	306656	2.20%	0.911%
4	9.057	1009	1015	1025	rVB	159168	282801	2.03%	0.840%
5	10.687	1284	1292	1300	rBV	227179	404840	2.90%	1.203%
6	13.151	1705	1711	1720	rBV	370907	607346	4.36%	1.805%
7	14.528	1938	1945	1955	rBV	385939	607458	4.36%	1.805%
8	16.028	2193	2200	2205	rBV	379560	612908	4.40%	1.822%
9	16.169	2218	2224	2244	rVB6	69091	274177	1.97%	0.815%
10	16.516	2266	2283	2309	rBV	1183526	3988929	28.61%	11.855%
11	17.286	2408	2414	2426	rBV	518330	787997	5.65%	2.342%
12	18.186	2561	2567	2582	rBV	647083	966699	6.93%	2.873%
13	19.004	2702	2706	2708	rBV	117512	152248	1.09%	0.452%
14	19.033	2708	2711	2721	rVB	973234	1178853	8.46%	3.504%
15	19.310	2754	2758	2767	rVB	198376	288134	2.07%	0.856%
16	19.428	2772	2778	2791	rBV	401872	616047	4.42%	1.831%
17	19.580	2799	2804	2810	rBV	839719	1005535	7.21%	2.988%
18	19.857	2845	2851	2858	rBV	973379	1204679	8.64%	3.580%
19	20.551	2959	2969	2980	rBV	9054947	13941645	100.00%	41.434%
20	20.645	2981	2985	2991	rVB	287839	351719	2.52%	1.045%
21	21.139	3065	3069	3076	rVB	493236	564193	4.05%	1.677%
22	21.386	3106	3111	3117	rBV	722326	977297	7.01%	2.904%
23	21.498	3126	3130	3142	rVB4	141536	273677	1.96%	0.813%
24	22.527	3299	3305	3323	rVB	1608450	2510561	18.01%	7.461%
25	23.804	3516	3522	3533	rVB	474077	1023173	7.34%	3.041%

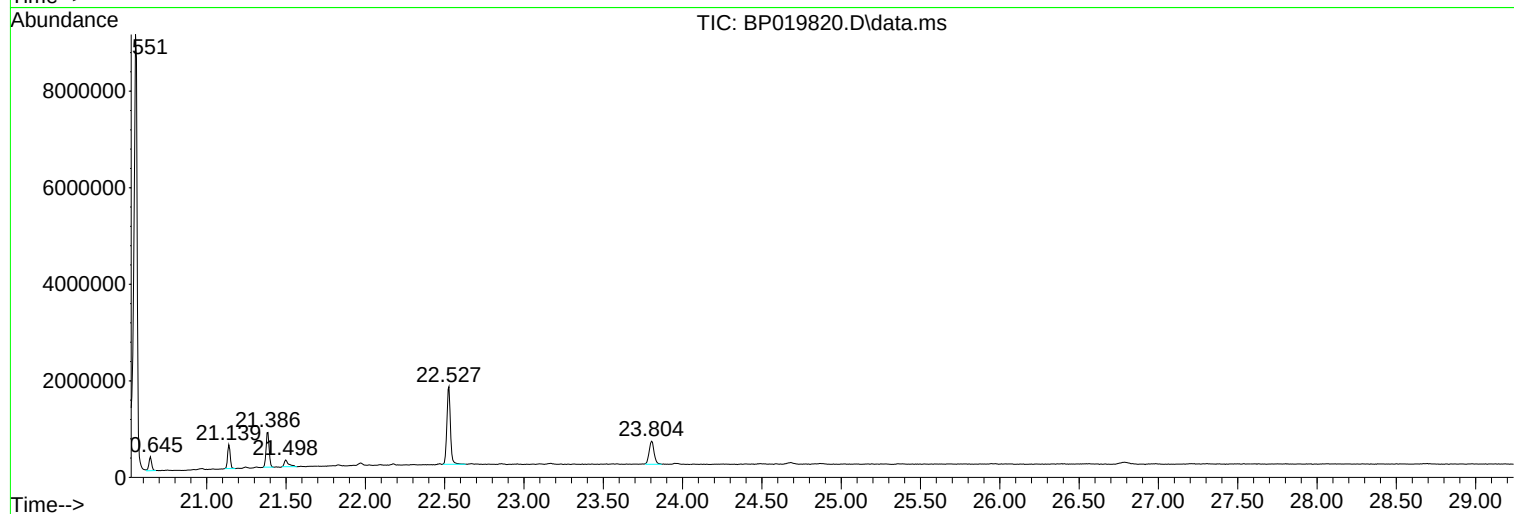
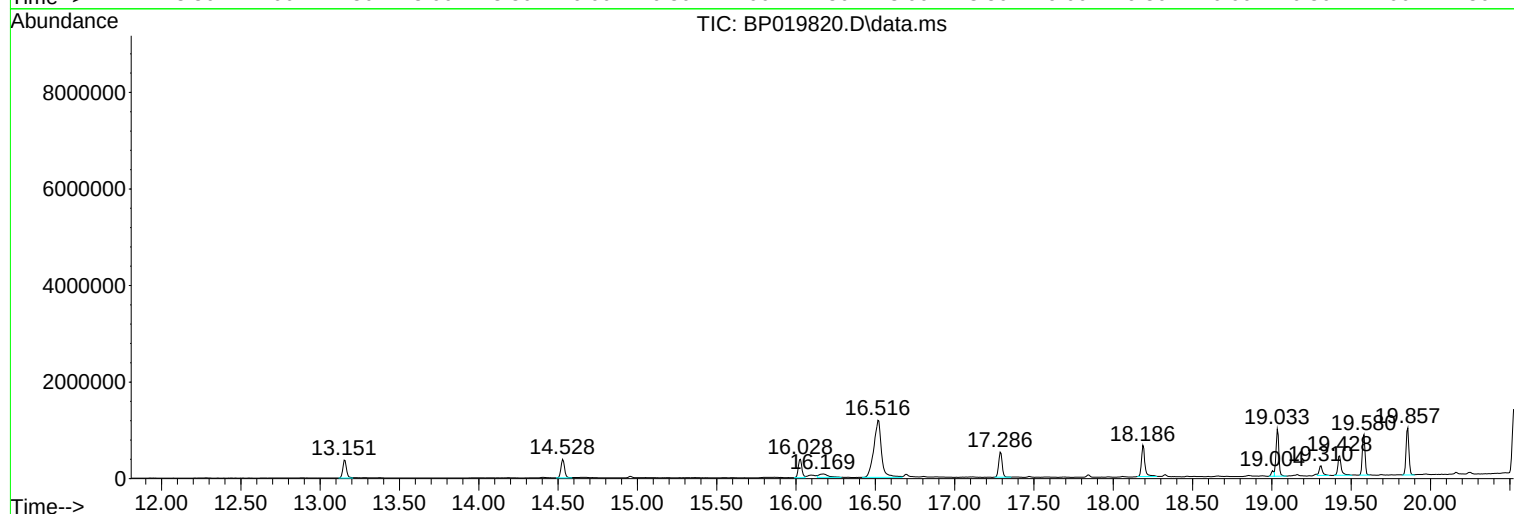
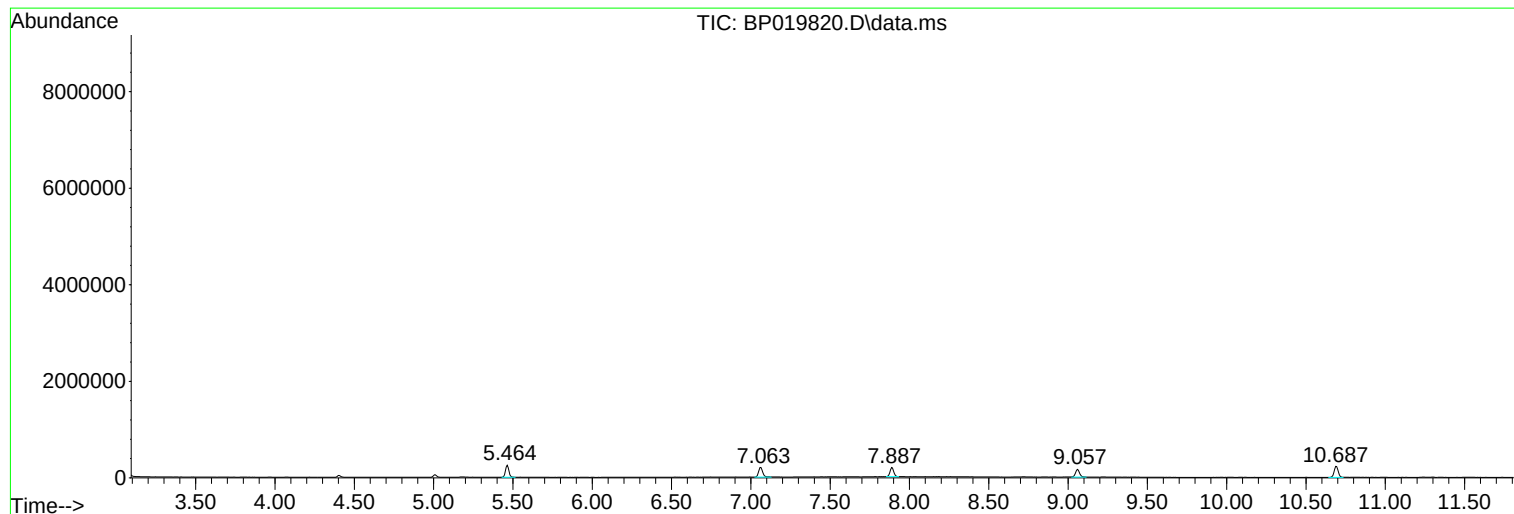
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Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP013124.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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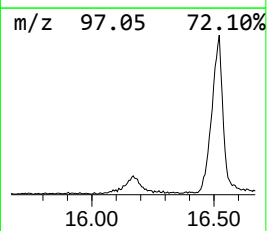
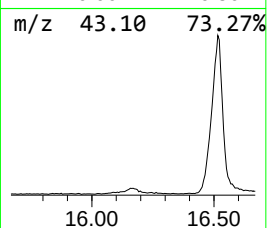
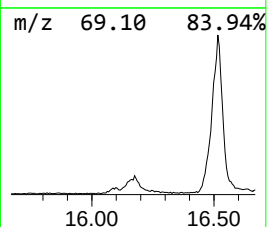
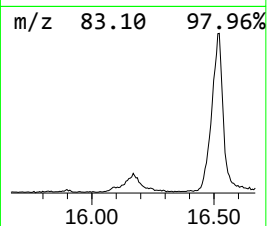
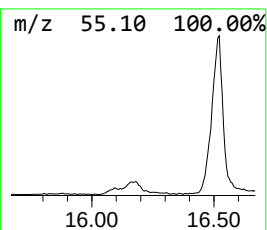
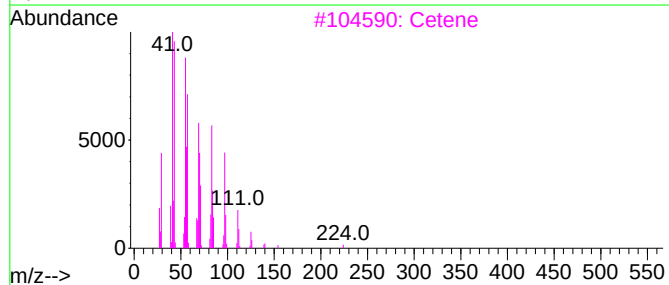
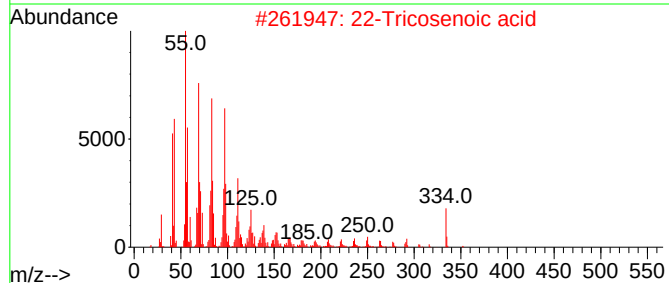
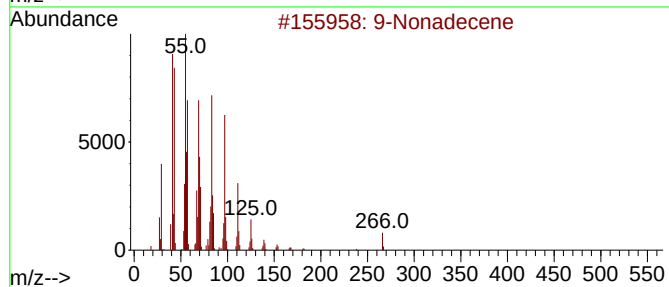
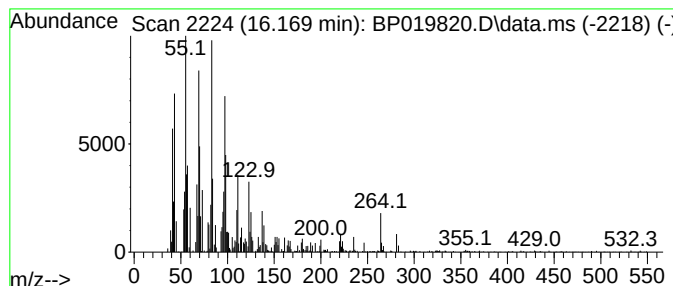
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 9-Nonadecene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.169	6.96 ng	274177	Phenanthrene-d10	17.287

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9-Nonadecene	266	C19H38	031035-07-1	86
2			22-Tricosenoic acid	352	C23H44O2	065119-95-1	83
3			Cetene	224	C16H32	000629-73-2	76
4			Bromoacetic acid, hexadecyl ester	362	C18H35BrO2	005454-48-8	72
5			7-Heptadecene, 1-chloro-	272	C17H33Cl	056554-78-0	64



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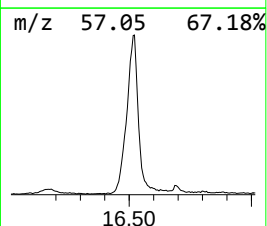
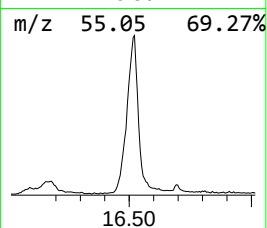
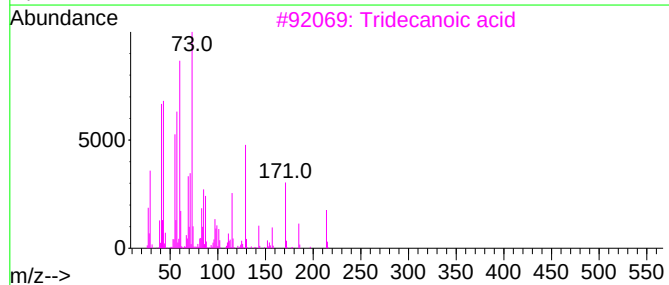
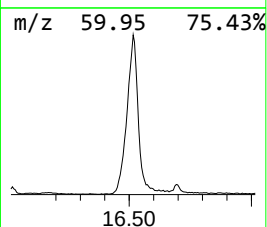
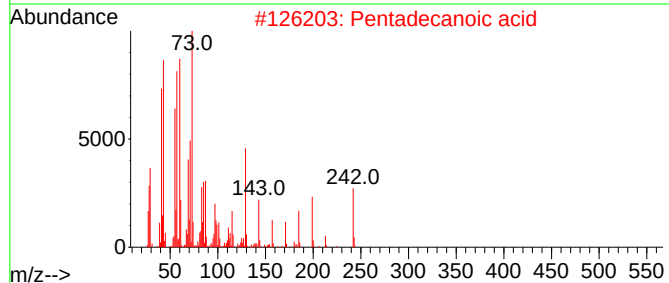
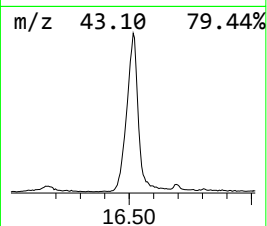
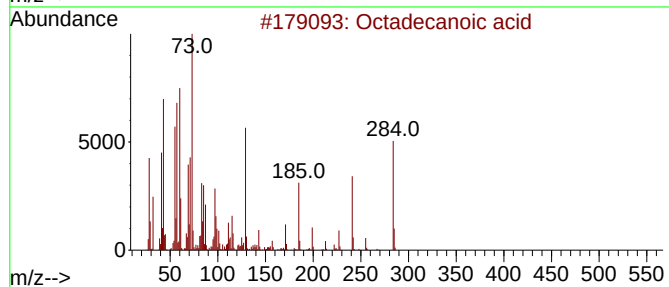
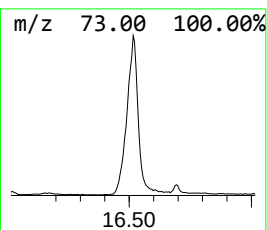
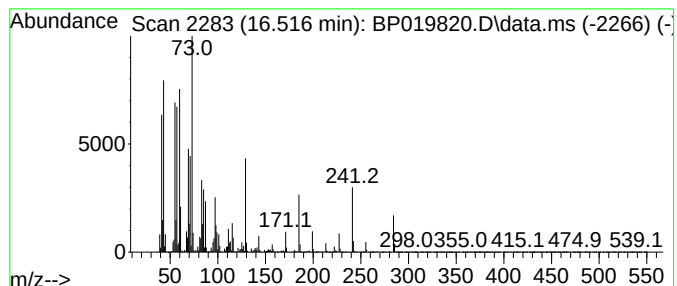
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP013124.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Octadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.516	101.24 ng	3988930	Phenanthrene-d10	17.287

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	94
3			Tridecanoic acid	214	C13H26O2	000638-53-9	90
4			Tetradecanoic acid	228	C14H28O2	000544-63-8	74
5			Octadecanoic acid, 2-(2-hydroxye...	372	C22H44O4	000106-11-6	72



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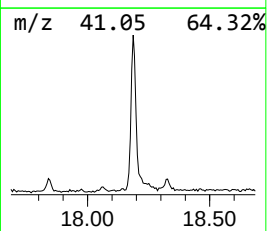
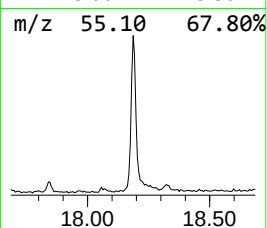
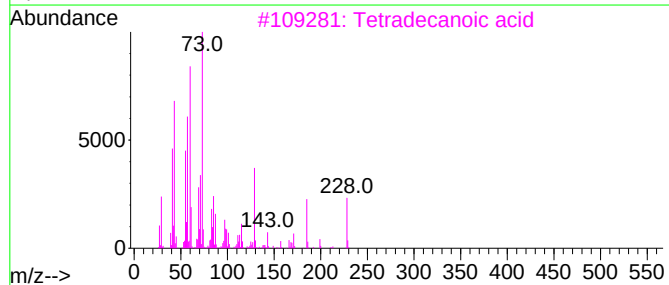
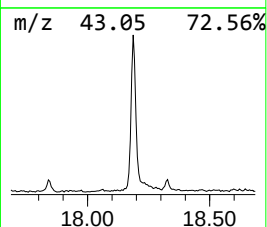
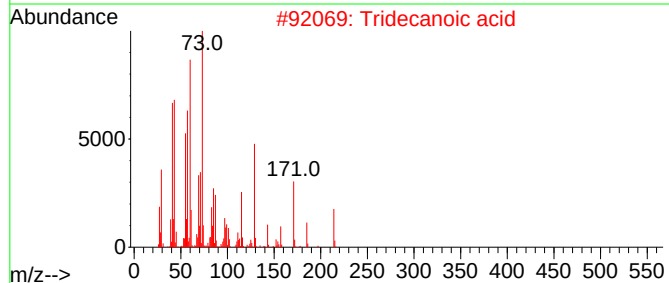
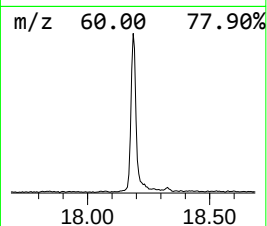
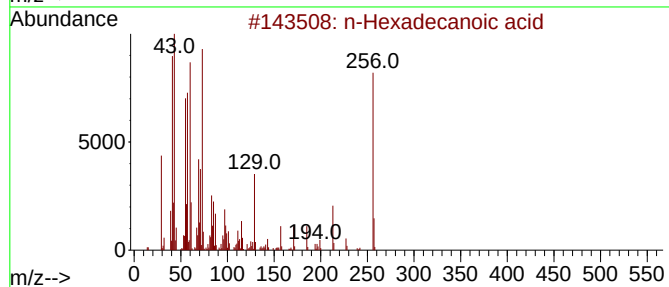
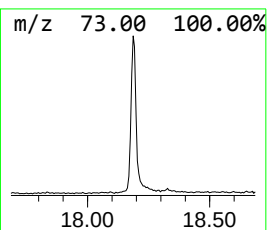
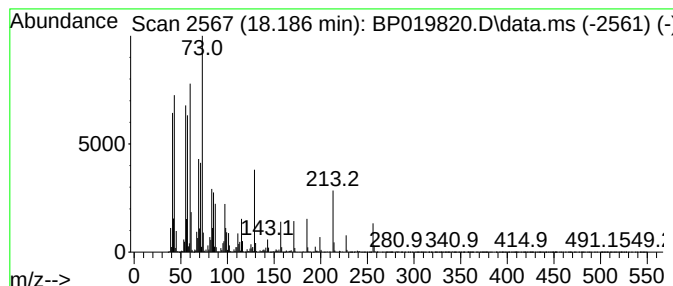
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.186	24.54 ng	966699	Phenanthrene-d10	17.287

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	91
3			Tetradecanoic acid	228	C14H28O2	000544-63-8	90
4			Pentadecanoic acid	242	C15H30O2	001002-84-2	87
5			n-Decanoic acid	172	C10H20O2	000334-48-5	76



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Misc :
ALS Vial : 5 Sample Multiplier: 1

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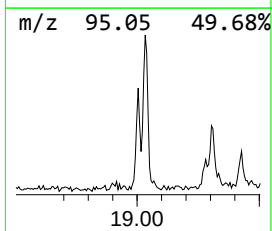
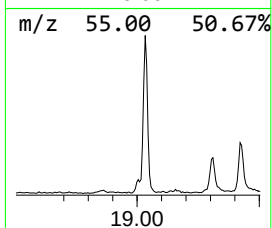
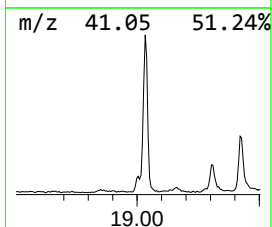
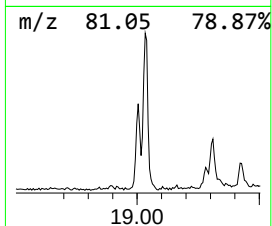
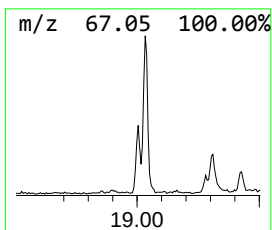
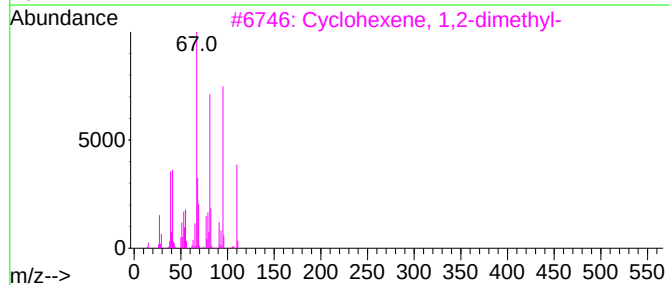
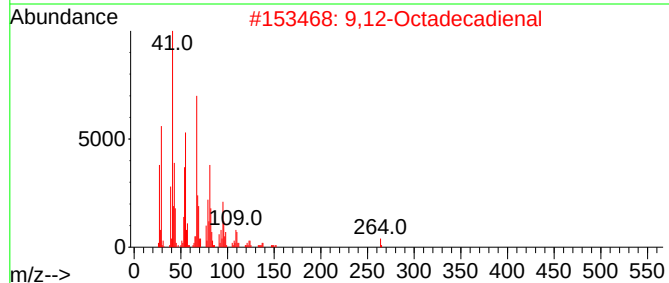
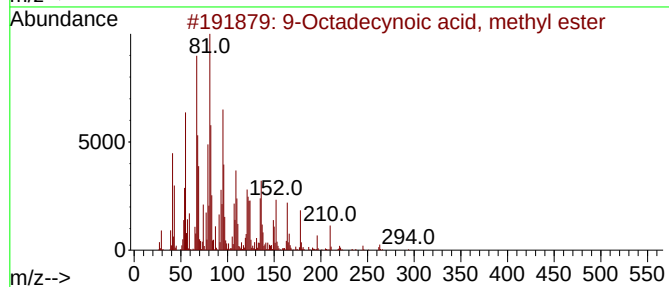
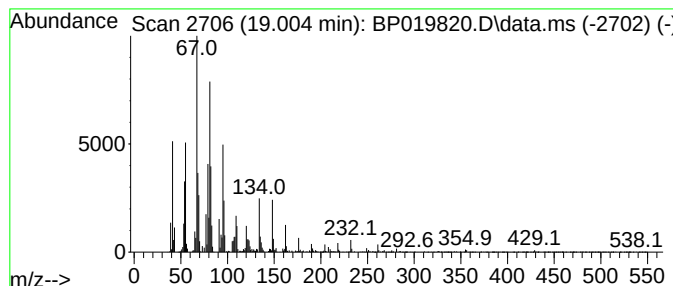
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 9-Octadecynoic acid, methyl... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.004	3.86 ng	152248	Phenanthrene-d10	17.287

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9-Octadecynoic acid, methyl ester	294	C19H34O2	001120-32-7	72
2			9,12-Octadecadienal	264	C18H32O	026537-70-2	64
3			Cyclohexene, 1,2-dimethyl-	110	C8H14	001674-10-8	52
4			Cyclooctane, ethenyl-	138	C10H18	061142-41-4	49
5			Spiro[2.5]octane	110	C8H14	000185-65-9	49



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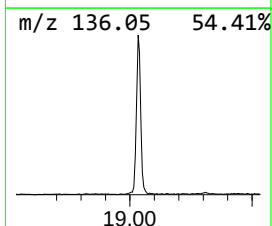
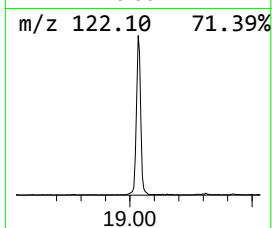
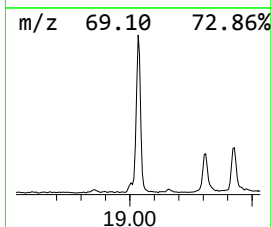
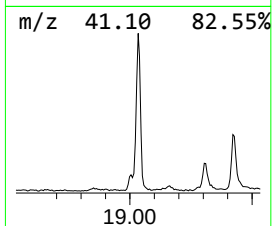
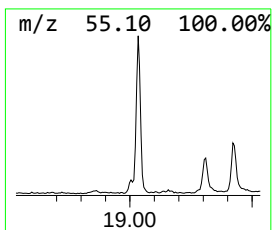
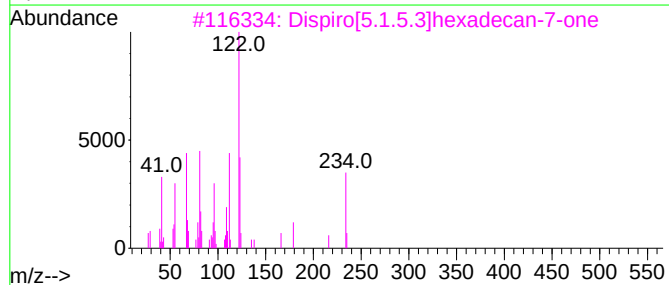
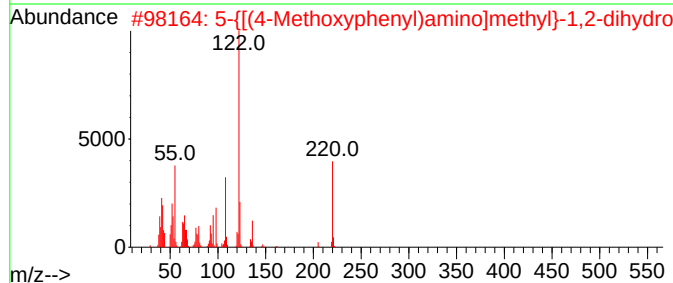
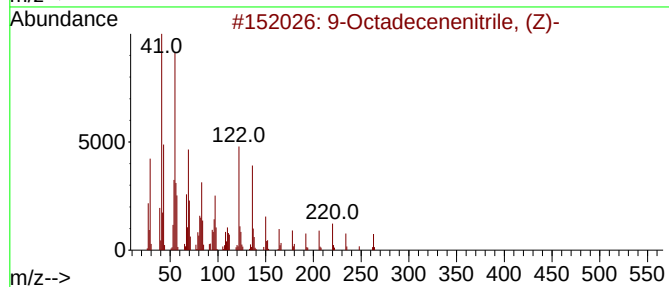
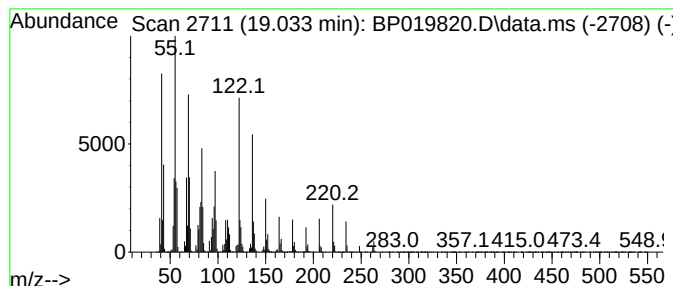
Instrument :
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ClientSampleId :
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 9-Octadecenitrile, (Z)- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD		R.T.	
19.033	29.92 ng	1178850	Phenanthrene-d10		17.287	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	9-Octadecenitrile, (Z)-		263	C18H33N	000112-91-4	94
2	5-[(4-Methoxyphenyl)amino]methy...		220	C10H12N4O2	1000434-75-9	20
3	Dispiro[5.1.5.3]hexadecan-7-one		234	C16H26O	001781-86-8	18
4	2,3-Dimethyl-5-n-propylpyrazine		150	C9H14N2	032262-98-9	18
5	Oleic Acid		282	C18H34O2	000112-80-1	18



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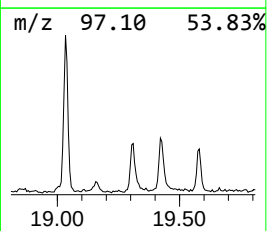
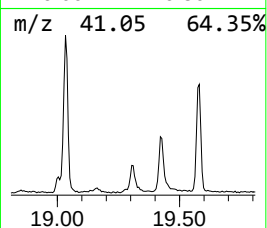
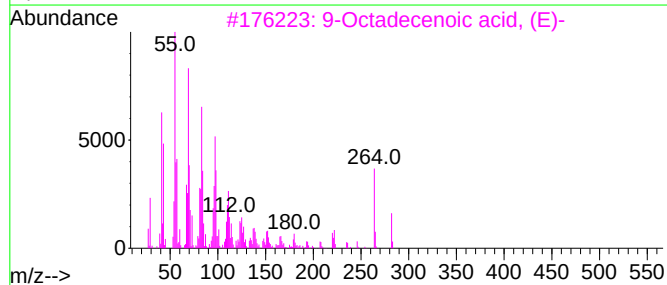
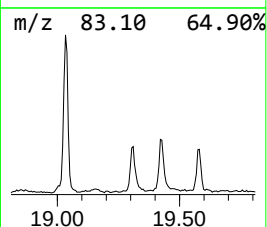
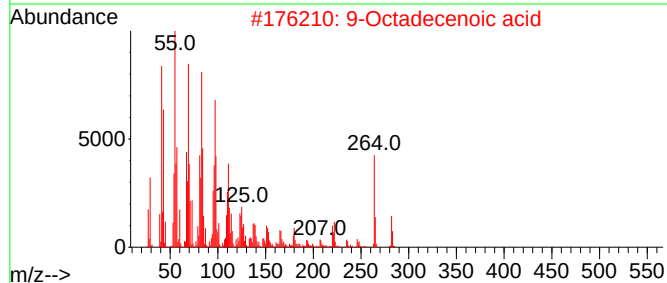
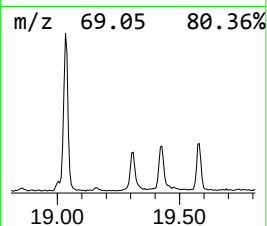
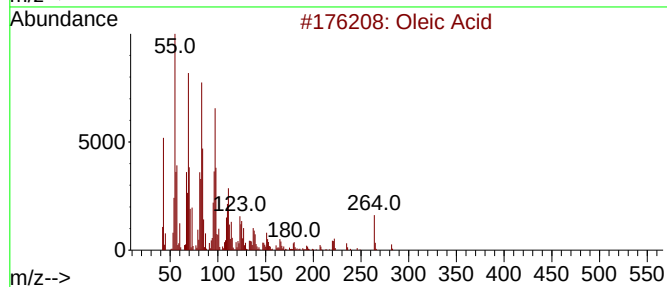
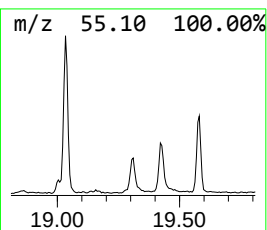
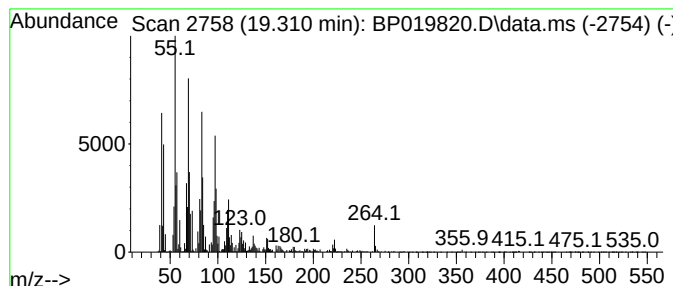
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Oleic Acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.310	7.31 ng	288134	Phenanthrene-d10	17.287

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Oleic Acid	282	C18H34O2	000112-80-1	91
2			9-Octadecenoic acid	282	C18H34O2	002027-47-6	91
3			9-Octadecenoic acid, (E)-	282	C18H34O2	000112-79-8	90
4			Octyl cis-vaccenate	394	C26H50O2	180252-12-4	87
5			2-Undecanol oleate	436	C29H56O2	1000465-66-9	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP022224\
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Acq On : 22 Feb 2024 12:42
Operator : MA/JU
Sample : P1462-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

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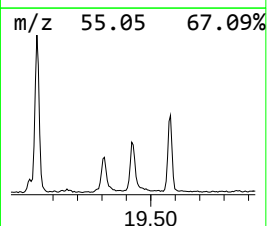
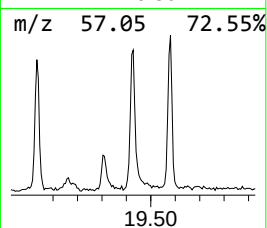
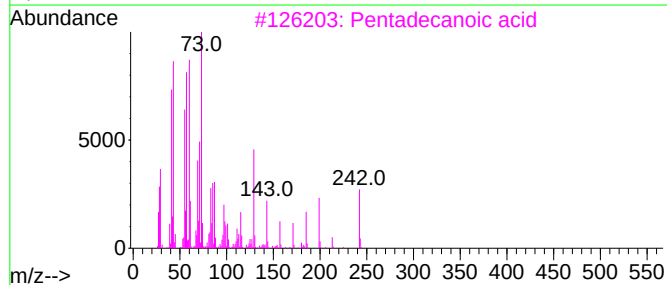
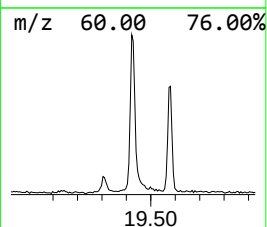
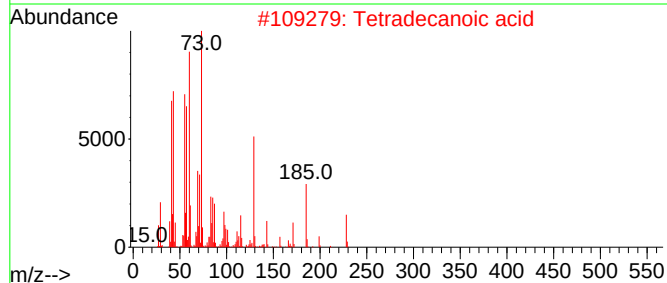
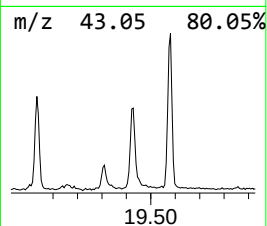
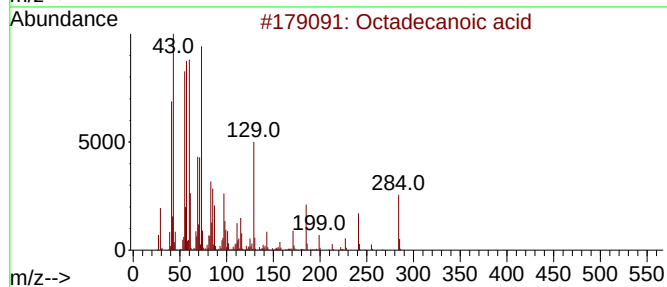
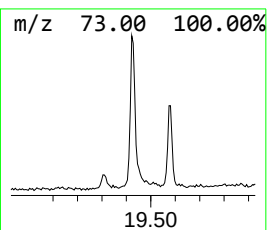
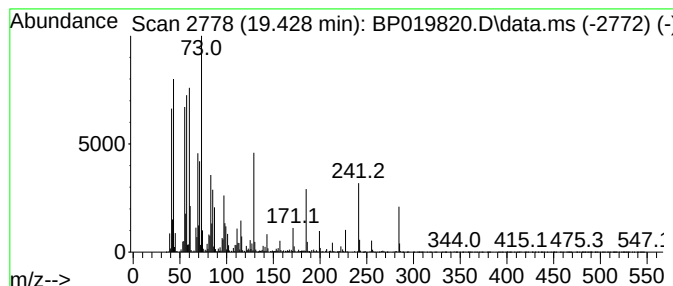
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Tetradecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.427	12.61 ng	616047	Chrysene-d12	21.380

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			Tetradecanoic acid	228	C14H28O2	000544-63-8	81
3			Pentadecanoic acid	242	C15H30O2	001002-84-2	76
4			Octadecanoic acid, 2-(2-hydroxye...	372	C22H44O4	000106-11-6	72
5			Tridecanoic acid	214	C13H26O2	000638-53-9	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP022224\
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Operator : MA/JU
Sample : P1462-01
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Instrument :
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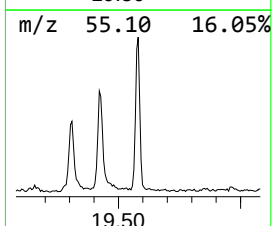
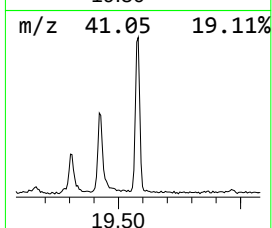
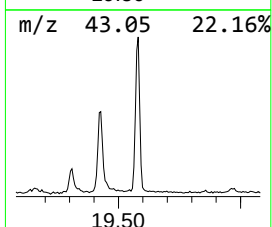
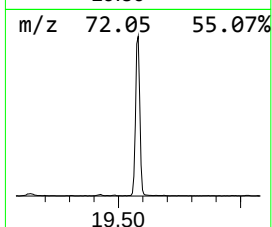
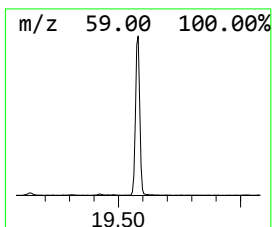
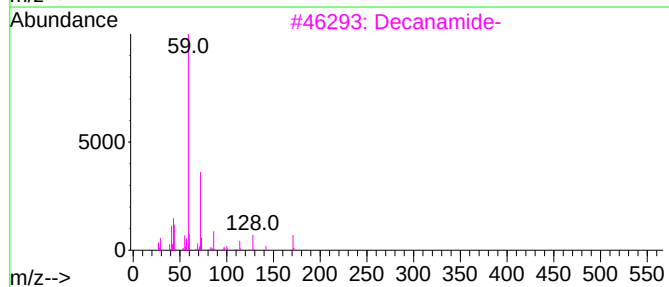
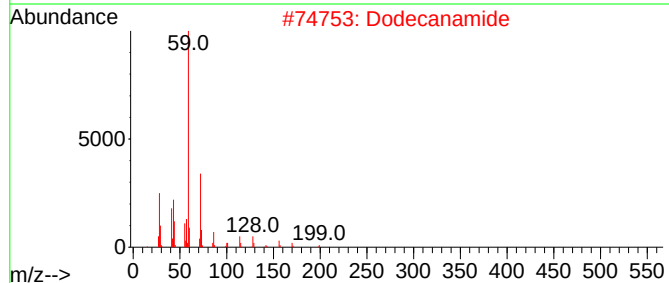
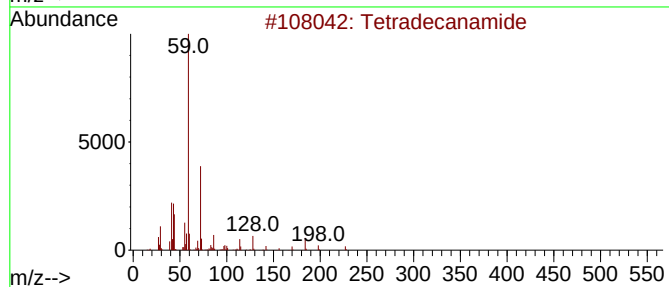
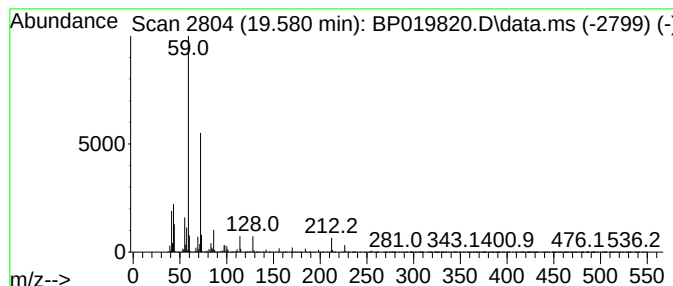
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Tetradecanamide Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.581	20.58 ng	1005540	Chrysene-d12	21.380

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecanamide	227	C14H29NO	000638-58-4	86
2			Dodecanamide	199	C12H25NO	001120-16-7	64
3			Decanamide-	171	C10H21NO	002319-29-1	59
4			Hexadecanamide	255	C16H33NO	000629-54-9	59
5			Heptanamide, 4-ethyl-5-methyl-	171	C10H21NO	054789-40-1	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP022224\
Data File : BP019820.D
Acq On : 22 Feb 2024 12:42
Operator : MA/JU
Sample : P1462-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

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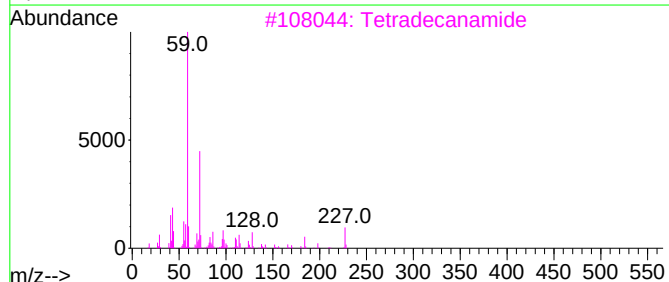
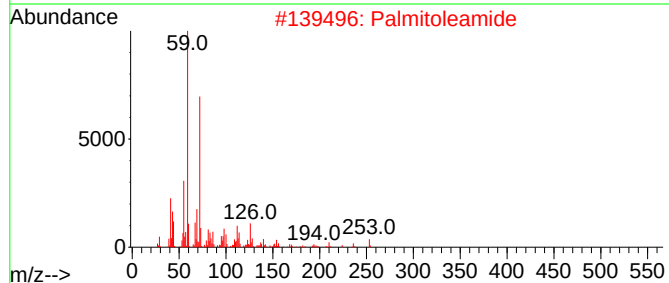
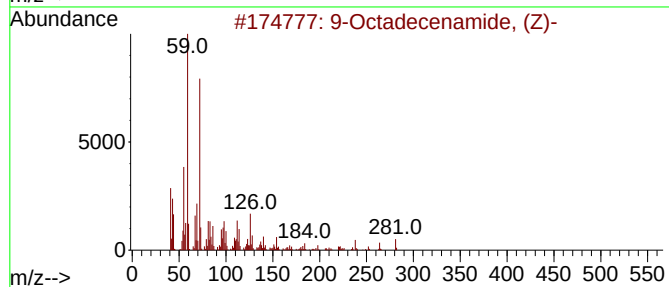
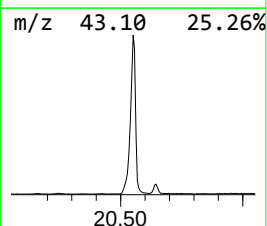
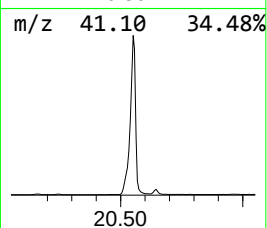
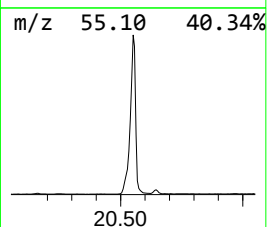
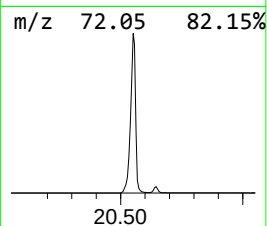
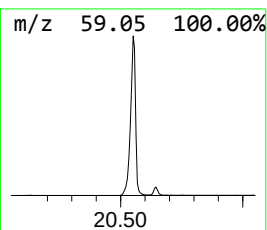
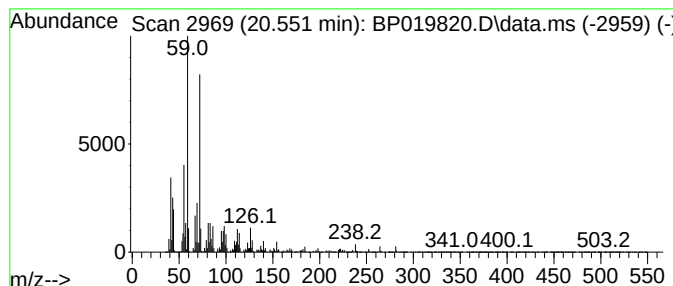
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 9-Octadecenamide, (Z)- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.551	285.31 ng	13941600	Chrysene-d12	21.380

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	99
2			Palmitoleamide	253	C16H31NO	106010-22-4	58
3			Tetradecanamide	227	C14H29NO	000638-58-4	43
4			Octadecanamide	283	C18H37NO	000124-26-5	38
5			Octanamide	143	C8H17NO	000629-01-6	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP022224\
Data File : BP019820.D
Acq On : 22 Feb 2024 12:42
Operator : MA/JU
Sample : P1462-01
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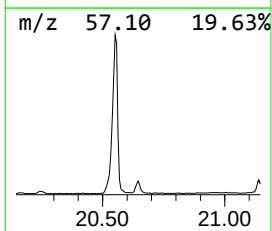
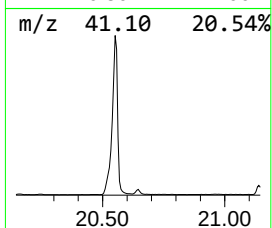
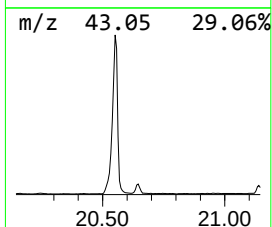
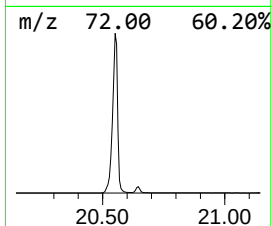
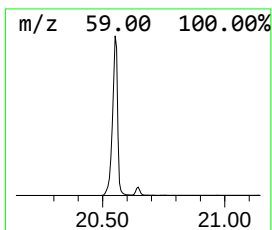
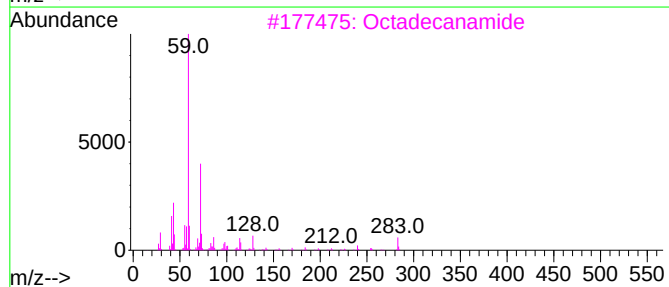
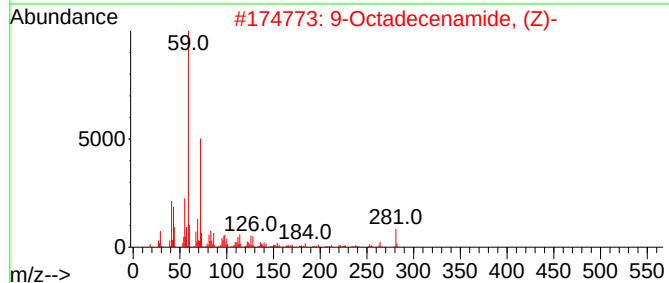
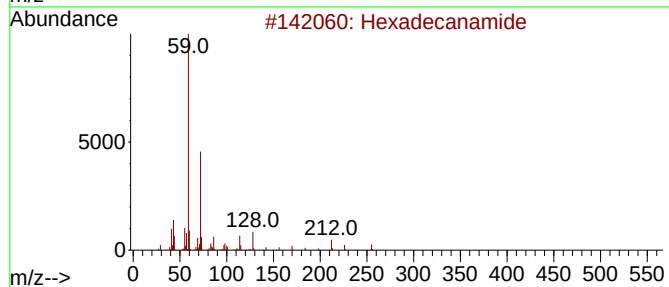
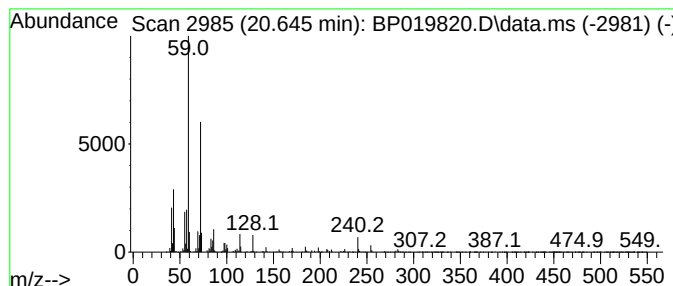
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Hexadecanamide Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.645	7.20 ng	351719	Chrysene-d12	21.380

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecanamide	255	C16H33NO	000629-54-9	86
2			9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	78
3			Octadecanamide	283	C18H37NO	000124-26-5	78
4			Tetradecanamide	227	C14H29NO	000638-58-4	64
5			Dodecanamide	199	C12H25NO	001120-16-7	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP022224\
Data File : BP019820.D
Acq On : 22 Feb 2024 12:42
Operator : MA/JU
Sample : P1462-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

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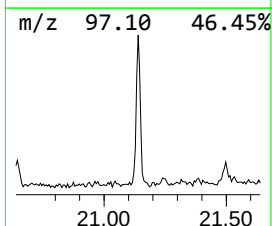
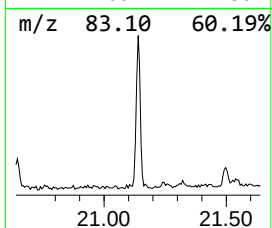
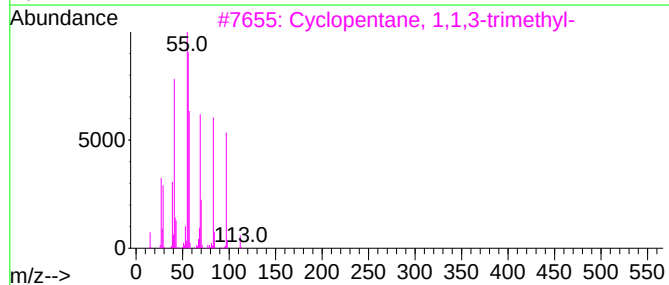
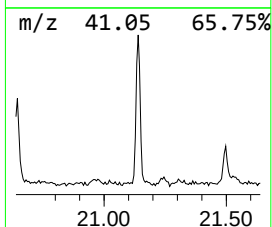
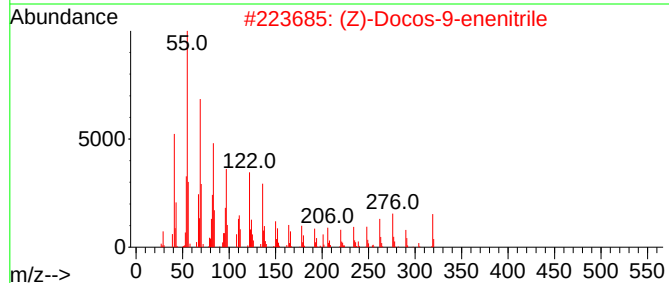
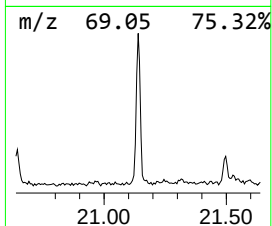
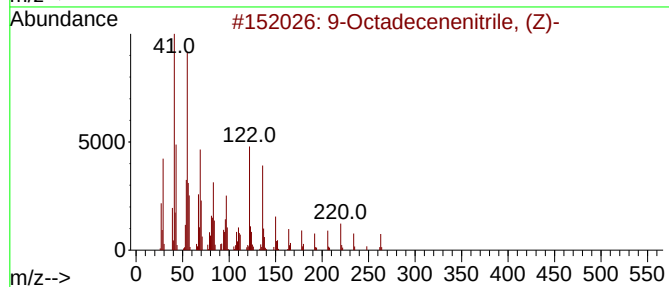
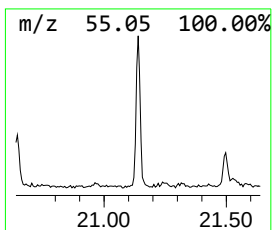
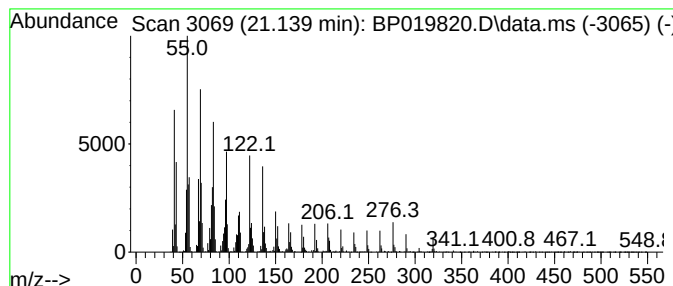
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 (Z)-Docos-9-enenitrile Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.139	11.55 ng	564193	Chrysene-d12	21.380

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9-Octadecenitrile, (Z)-	263	C18H33N	000112-91-4	90
2			(Z)-Docos-9-enenitrile	319	C22H41N	1000465-48-0	89
3			Cyclopentane, 1,1,3-trimethyl-	112	C8H16	004516-69-2	46
4			3-n-Hexylthiane, S,S-dioxide	218	C11H22O2S	061639-18-7	38
5			Palmitoleonitrile	235	C16H29N	056617-97-1	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP022224\
Data File : BP019820.D
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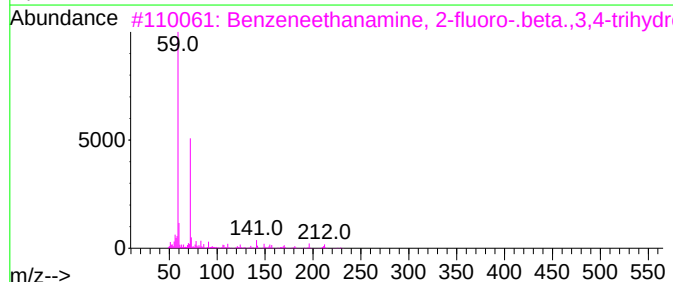
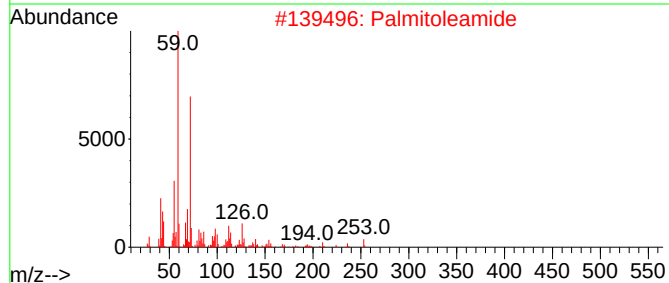
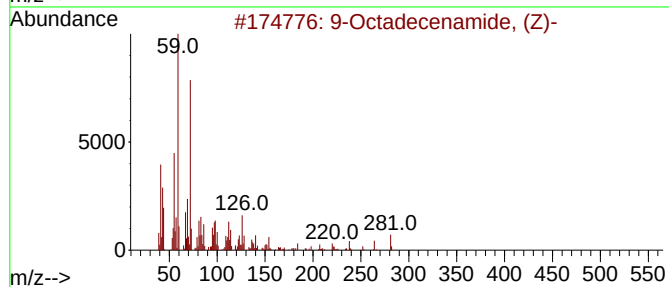
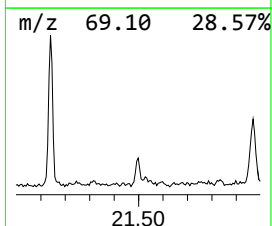
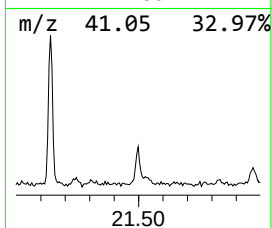
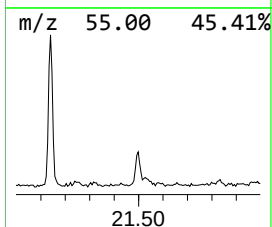
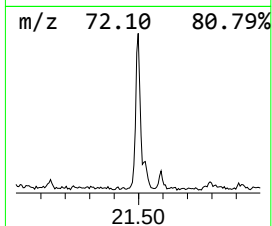
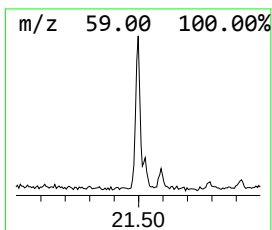
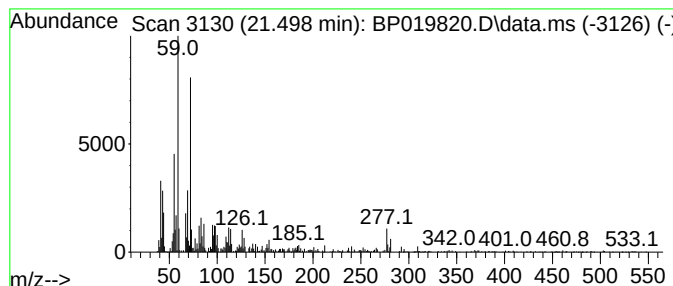
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Palmitoleamide Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.498	5.60 ng	273677	Chrysene-d12	21.380

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	93
2			Palmitoleamide	253	C16H31NO	106010-22-4	80
3			Benzeneethanamine, 2-fluoro-.bet...	229	C11H16FN03	061338-98-5	49
4			9-Octadecenamide	281	C18H35NO	003322-62-1	45
5			Dodecanamide	199	C12H25NO	001120-16-7	45



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP022224\
Data File : BP019820.D
Acq On : 22 Feb 2024 12:42
Operator : MA/JU
Sample : P1462-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

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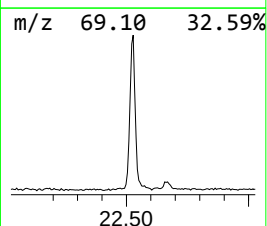
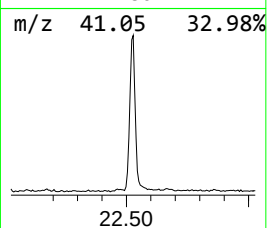
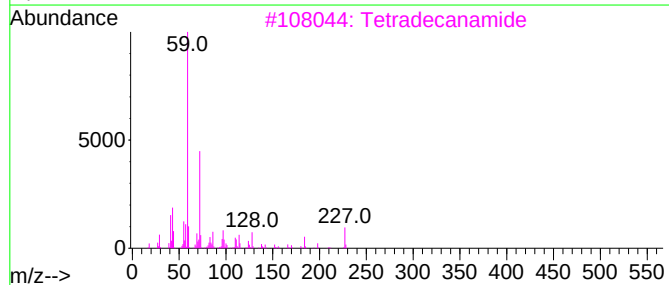
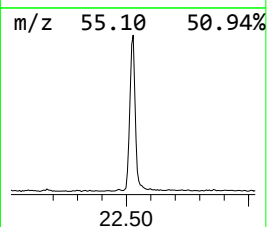
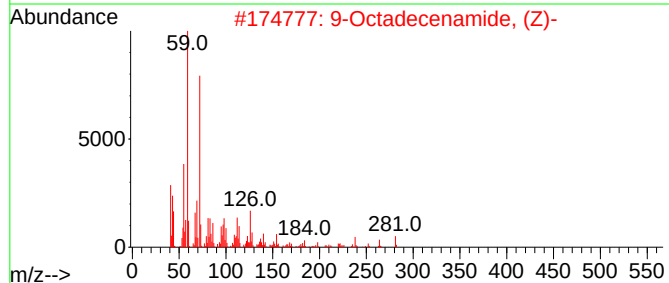
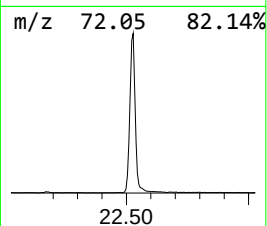
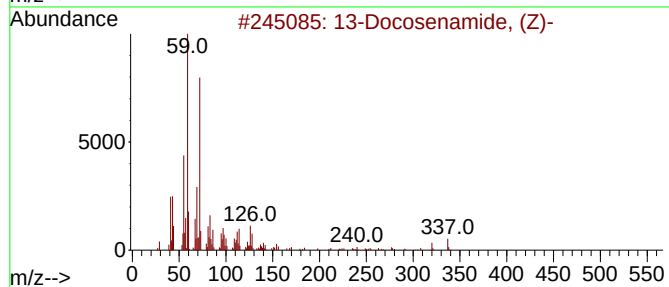
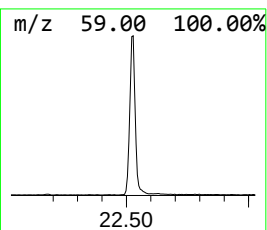
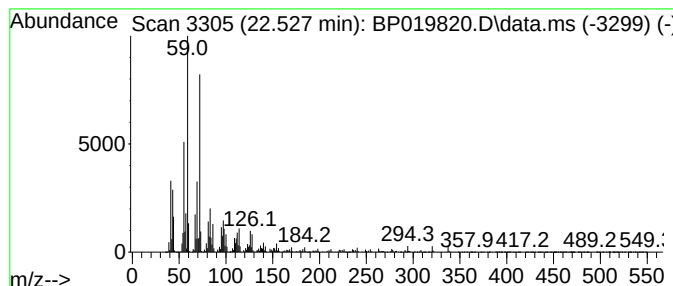
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 13-Docosenamide, (Z)- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.527	51.38 ng	2510560	Chrysene-d12	21.380

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	99
2			9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	94
3			Tetradecanamide	227	C14H29NO	000638-58-4	47
4			1-Propyne, 3-[(1-methylethyl)thio]-	114	C6H10S	014272-25-4	38
5			Pent-4-enoylamide, 2-methyl-N-pr...	211	C13H25NO	1000438-81-8	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP022224\
Data File : BP019820.D
Acq On : 22 Feb 2024 12:42
Operator : MA/JU
Sample : P1462-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
20240762-39N-N-FIELD-BLANK-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP013124.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
9-Nonadecene	16.169	7.0	ng	274177	4	17.287	787997	20.0
Octadecanoic acid	16.516	101.2	ng	3988930	4	17.287	787997	20.0
n-Hexadecanoic ...	18.186	24.5	ng	966699	4	17.287	787997	20.0
9-Octadecynoic ...	19.004	3.9	ng	152248	4	17.287	787997	20.0
9-Octadecenit...	19.033	29.9	ng	1178850	4	17.287	787997	20.0
Oleic Acid	19.310	7.3	ng	288134	4	17.287	787997	20.0
Tetradecanoic acid	19.427	12.6	ng	616047	5	21.380	977297	20.0
Tetradecanamide	19.581	20.6	ng	1005540	5	21.380	977297	20.0
9-Octadecenamid...	20.551	285.3	ng	13941600	5	21.380	977297	20.0
Hexadecanamide	20.645	7.2	ng	351719	5	21.380	977297	20.0
(Z)-Docos-9-ene...	21.139	11.6	ng	564193	5	21.380	977297	20.0
Palmitoleamide	21.498	5.6	ng	273677	5	21.380	977297	20.0
13-Docosenamide...	22.527	51.4	ng	2510560	5	21.380	977297	20.0