

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011521\
 Data File : BP004598.D
 Acq On : 16 Jan 2021 03:24
 Operator : CG/JU
 Sample : M1120-01 2X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 AU-04-011321

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-BP011221.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP004598.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.399	369	376	393	rBV	1129075	1794135	4.26%	1.313%
2	6.981	635	645	661	rBV	1112579	1890558	4.49%	1.384%
3	7.805	777	785	794	rBV	772821	1262883	3.00%	0.924%
4	8.963	976	982	990	rBV	697278	1172739	2.79%	0.858%
5	10.599	1249	1260	1275	rBV	1041114	1932793	4.59%	1.415%
6	12.257	1535	1542	1556	rBV	287737	515540	1.23%	0.377%
7	13.069	1673	1680	1691	rBV	2057652	3188097	7.58%	2.334%
8	14.457	1908	1916	1926	rBV	1740271	2590879	6.16%	1.897%
9	15.957	2164	2171	2178	rBV	1664214	2517487	5.98%	1.843%
10	17.222	2379	2386	2392	rBV	2043836	3078235	7.31%	2.253%
11	17.898	2494	2501	2505	rBV	12126668	17643824	41.93%	12.916%
12	18.116	2534	2538	2552	rBV	365627	657503	1.56%	0.481%
13	19.016	2682	2691	2693	rBV	18626021	42083734	100.00%	30.806%
14	19.045	2693	2696	2704	rVV3	17354762	29662905	70.49%	21.714%
15	19.110	2704	2707	2710	rVB2	417252	483384	1.15%	0.354%
16	19.157	2710	2715	2720	rVB	7384329	8905604	21.16%	6.519%
17	19.210	2720	2724	2726	rBV	450505	575800	1.37%	0.422%
18	19.510	2770	2775	2781	rVB2	397258	629915	1.50%	0.461%
19	19.786	2817	2822	2830	rVB	3834640	4845660	11.51%	3.547%
20	20.022	2859	2862	2864	rBV	430958	548046	1.30%	0.401%
21	20.110	2875	2877	2885	rVB	460734	524368	1.25%	0.384%
22	20.222	2892	2896	2901	rBV	781931	1057206	2.51%	0.774%
23	20.827	2993	2999	3003	rBV3	339534	696287	1.65%	0.510%
24	21.145	3050	3053	3059	rVB	866741	927615	2.20%	0.679%
25	21.316	3077	3082	3090	rVB	2706531	3586651	8.52%	2.626%
26	23.663	3476	3481	3496	rVB2	1854573	3835010	9.11%	2.807%

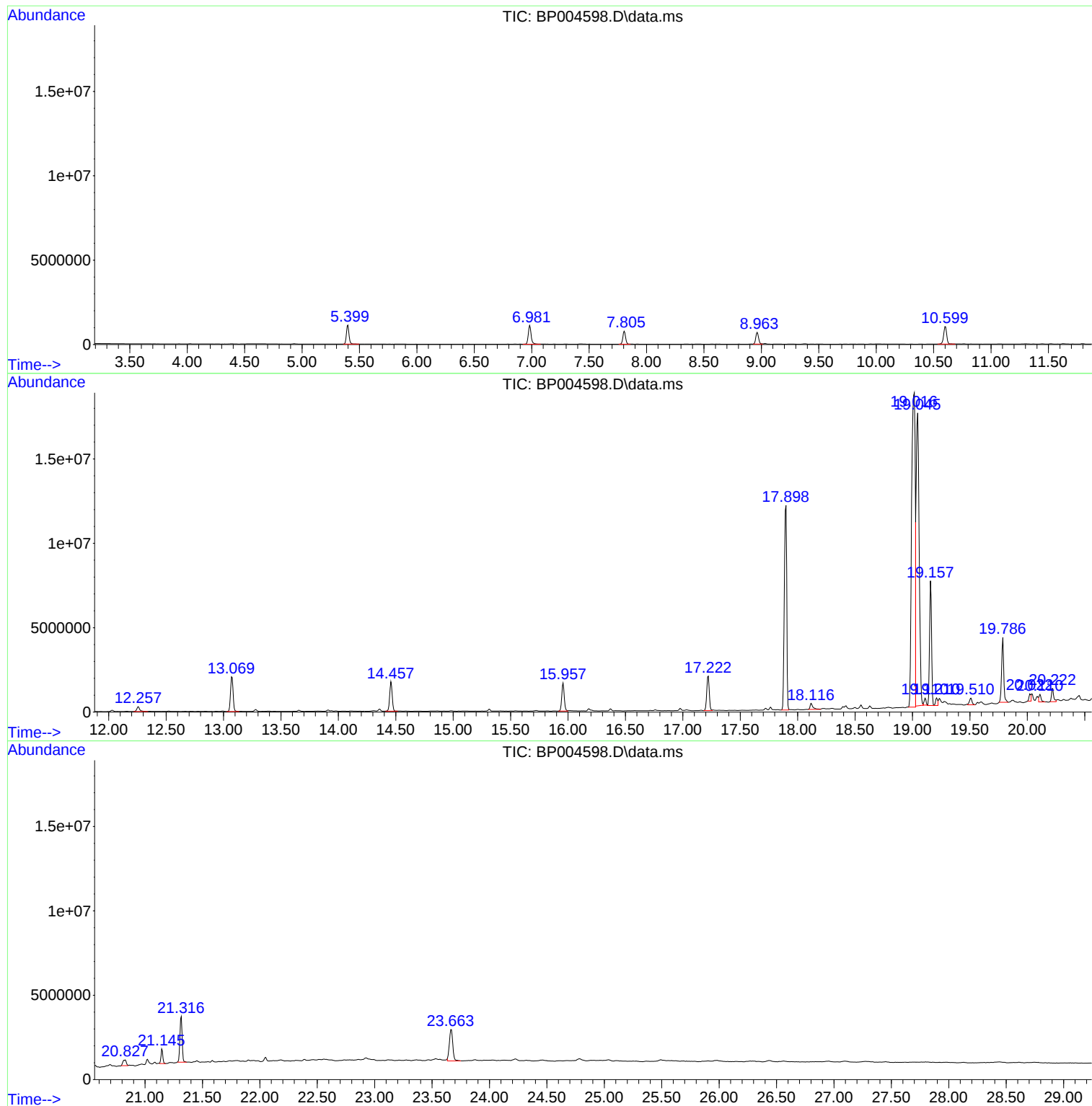
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270E-BP011221.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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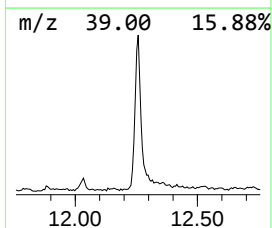
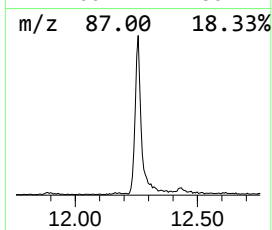
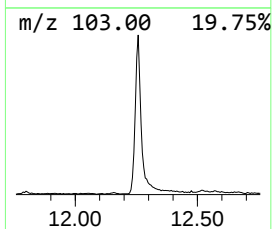
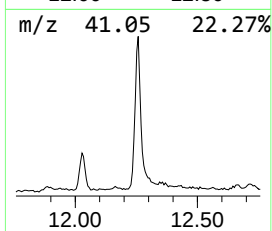
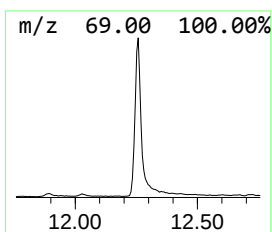
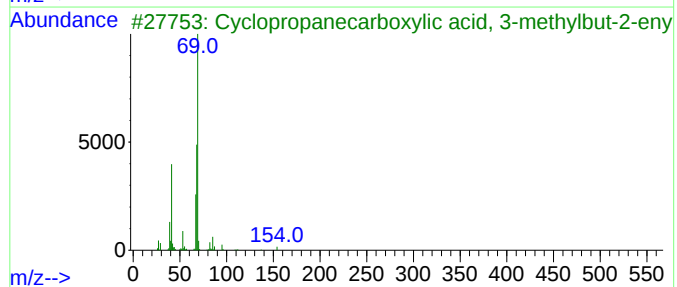
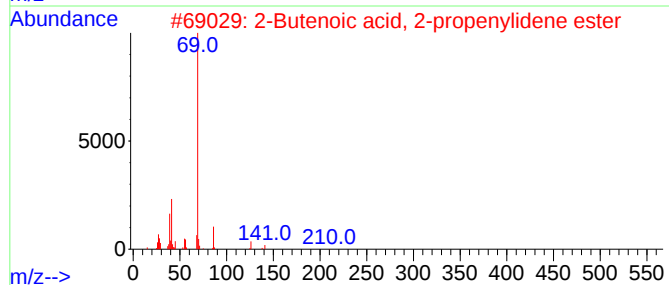
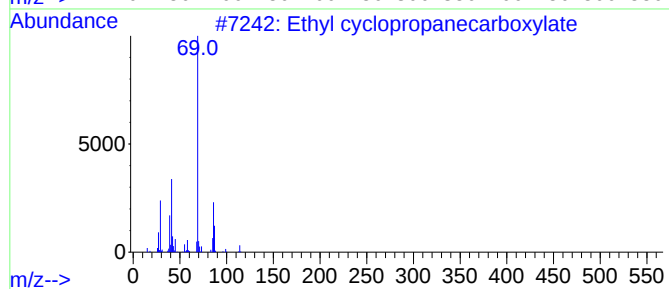
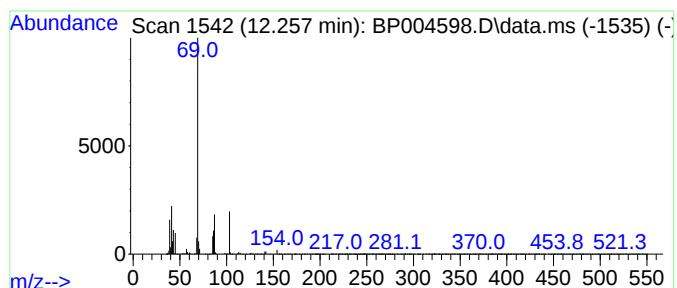
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown12.257 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.257	5.33 ng	515540	Naphthalene-d8	10.599

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethyl cyclopropanecarboxylate	114	C6H10O2	004606-07-9	38
2			2-Butenoic acid, 2-propenylidene...	210	C11H14O4	055030-70-1	33
3			Cyclopropanecarboxylic acid, 3-m...	154	C9H14O2	1000299-37-4	9
4			Pentanoic acid, 2-hydroxy-4-meth...	146	C7H14O3	040348-72-9	9
5			Cyclopropanecarboxylic acid, sil...	192	C4H5AgO2	075112-77-5	9



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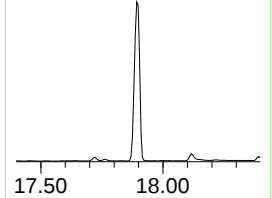
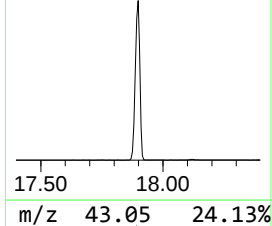
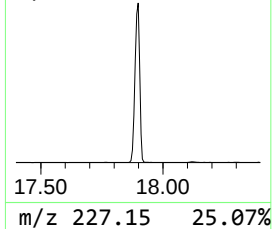
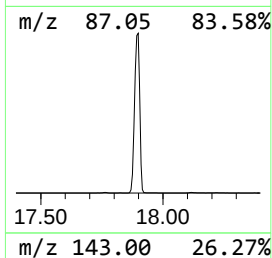
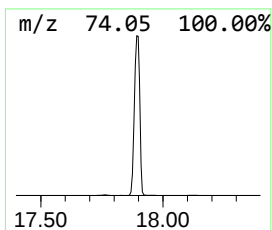
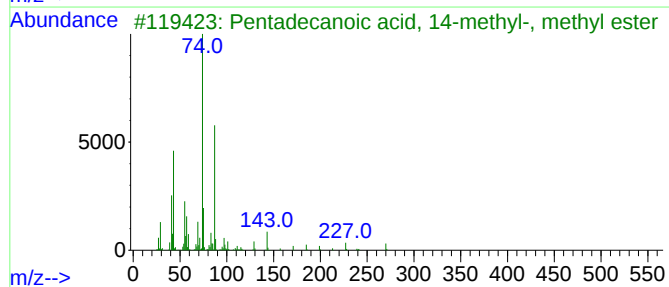
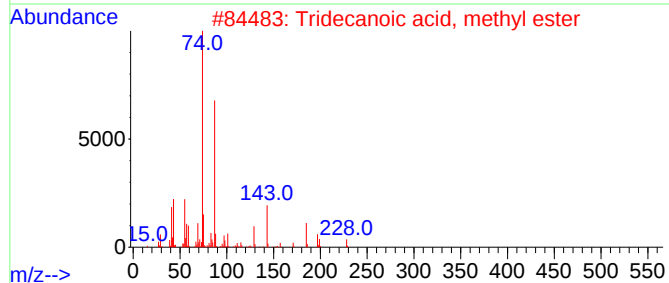
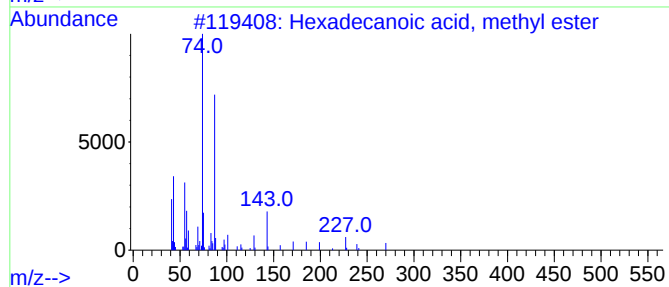
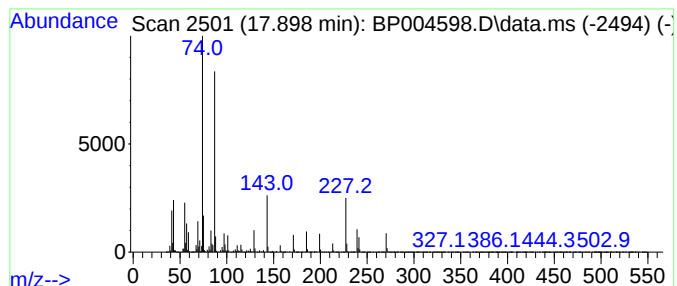
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Peak Number 2 Hexadecanoic acid, methyl e... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.898	114.64 ng	17643800	Phenanthrene-d10	17.222

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	97
2			Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	96
3			Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	95
4			Methyl tetradecanoate	242	C15H30O2	000124-10-7	89
5			Hexadecanoic acid, 2-methyl-	270	C17H34O2	027147-71-3	74



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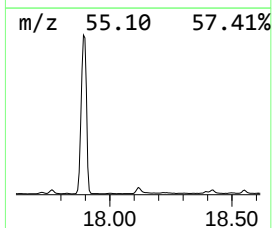
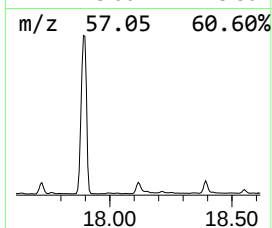
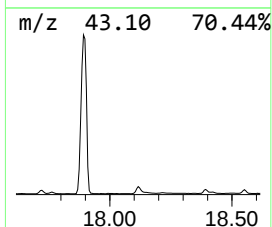
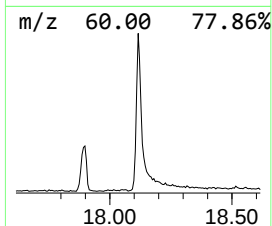
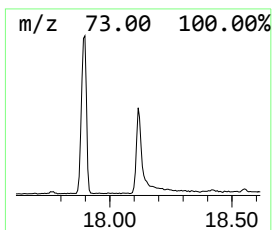
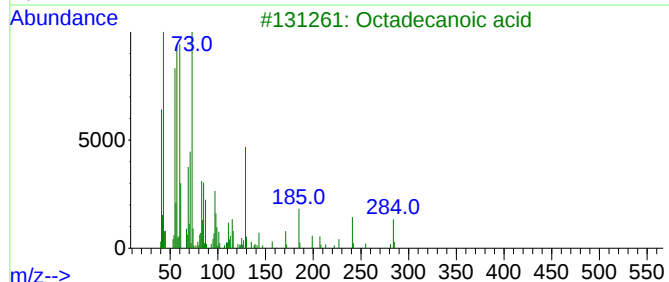
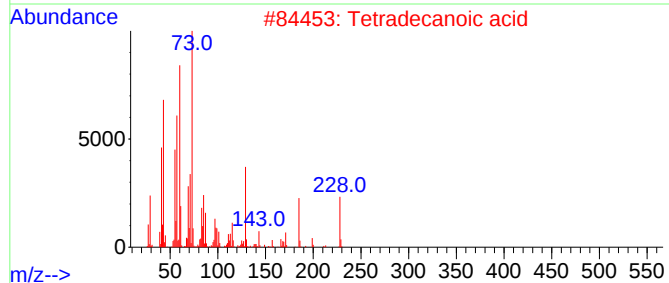
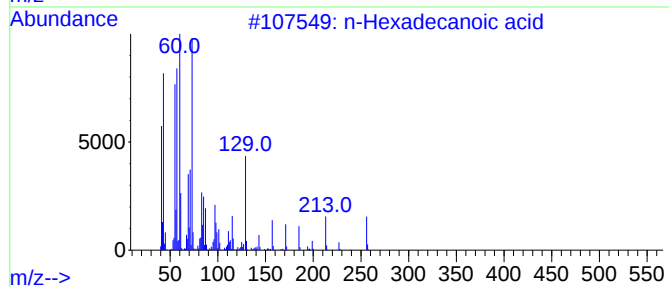
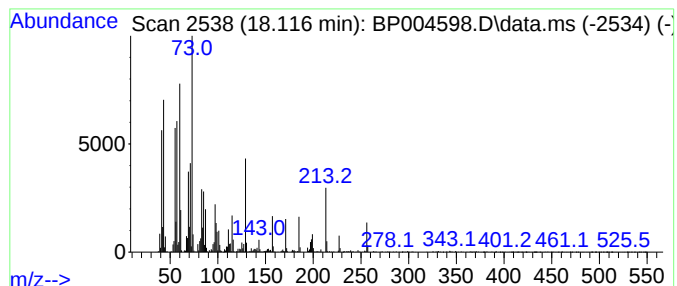
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 n-Hexadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.116	4.27 ng	657503	Phenanthrene-d10	17.222

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2			Tetradecanoic acid	228	C14H28O2	000544-63-8	91
3			Octadecanoic acid	284	C18H36O2	000057-11-4	87
4			Tridecanoic acid	214	C13H26O2	000638-53-9	86
5			Undecanoic acid	186	C11H22O2	000112-37-8	80



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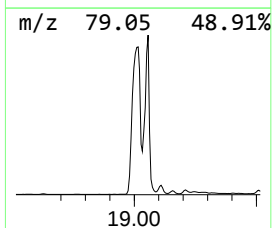
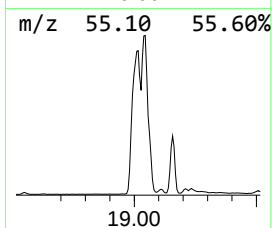
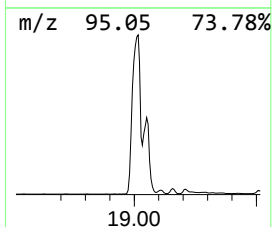
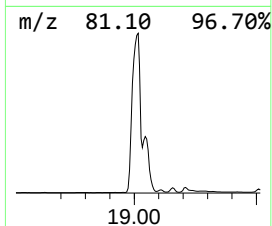
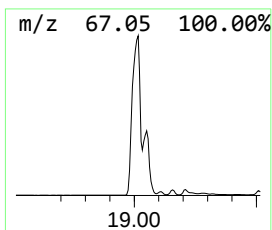
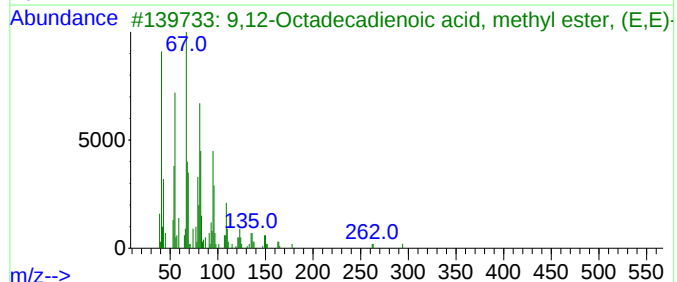
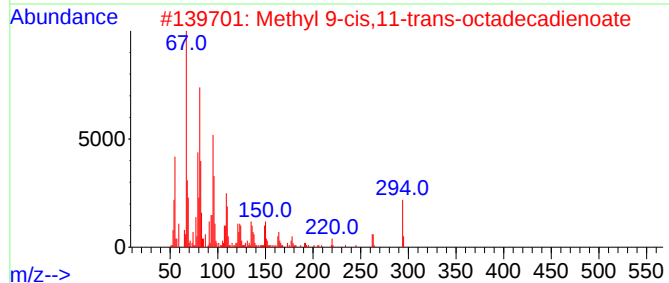
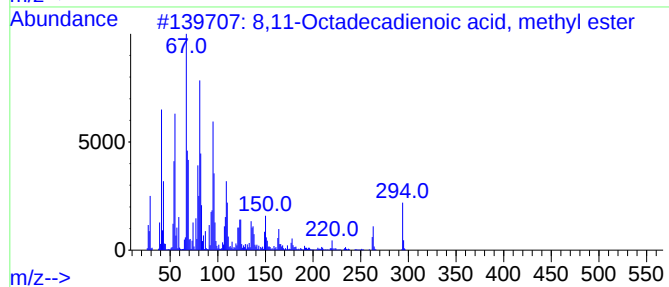
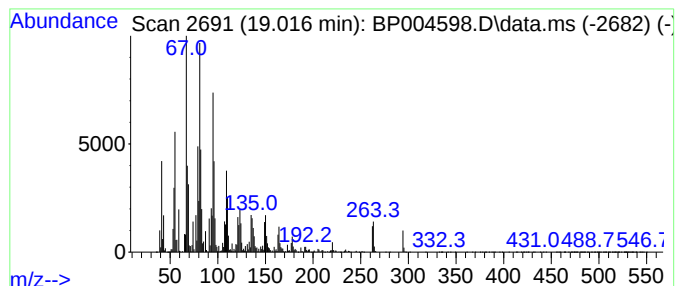
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Peak Number 4 8,11-Octadecadienoic acid, ... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.016	273.43 ng	42083700	Phenanthrene-d10	17.222

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			8,11-Octadecadienoic acid, methy...	294	C19H34O2	056599-58-7	99
2			Methyl 9-cis,11-trans-octadecadi...	294	C19H34O2	1000336-44-0	99
3			9,12-Octadecadienoic acid, methy...	294	C19H34O2	002566-97-4	99
4			9,12-Octadecadienoic acid (Z,Z)-...	294	C19H34O2	000112-63-0	99
5			10,13-Octadecadienoic acid, meth...	294	C19H34O2	056554-62-2	99



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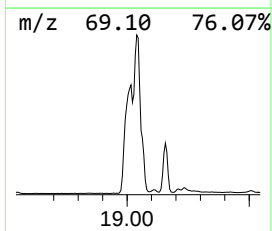
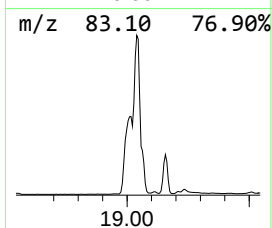
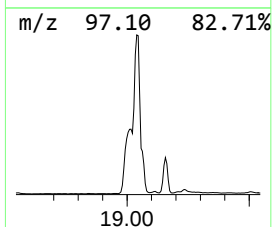
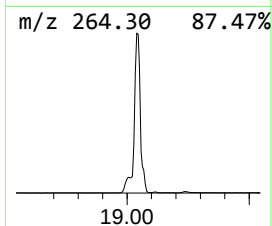
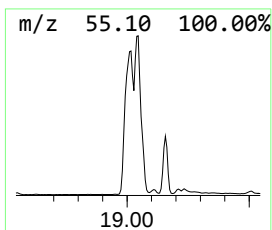
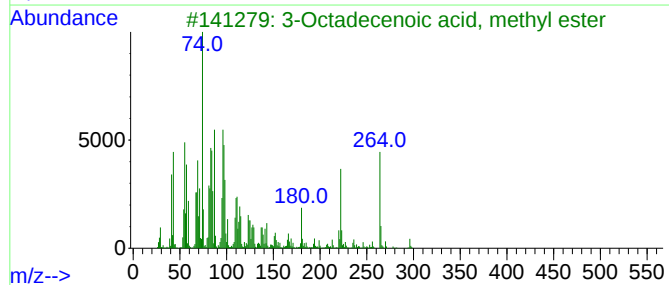
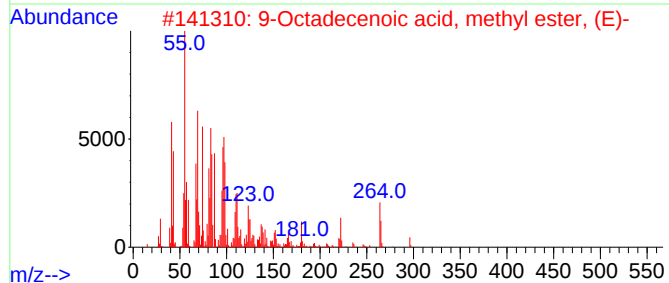
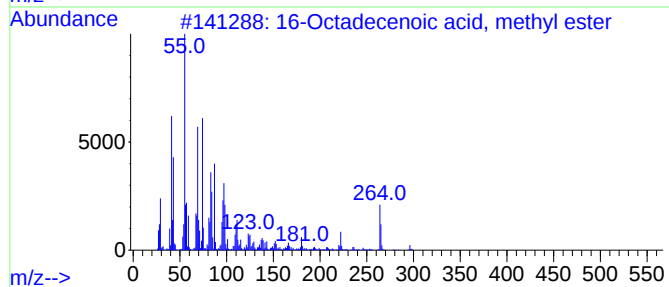
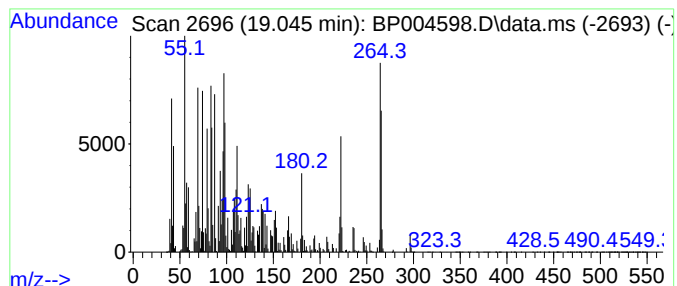
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Peak Number 5 16-Octadecenoic acid, methyl... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.045	192.73 ng	29662900	Phenanthrene-d10	17.222	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	16-Octadecenoic acid, methyl ester	296	C19H36O2	056554-49-5	97
2	9-Octadecenoic acid, methyl este...	296	C19H36O2	001937-62-8	92
3	3-Octadecenoic acid, methyl ester	296	C19H36O2	056599-33-8	89
4	cis-13-Octadecenoic acid, methyl...	296	C19H36O2	1000333-58-3	89
5	9-Octadecenoic acid (Z)-, methyl...	296	C19H36O2	000112-62-9	86



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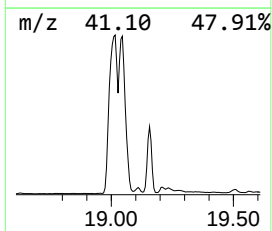
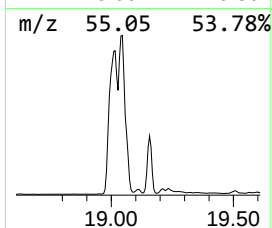
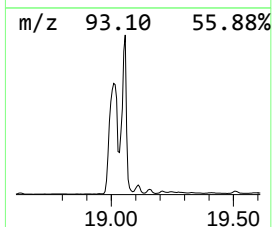
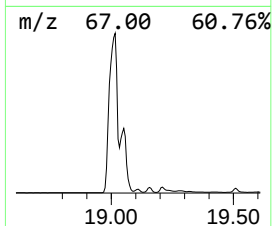
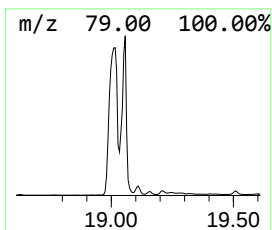
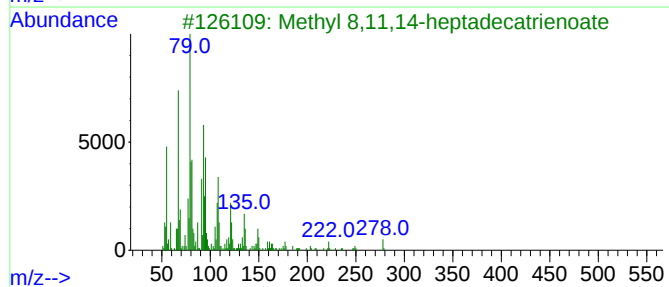
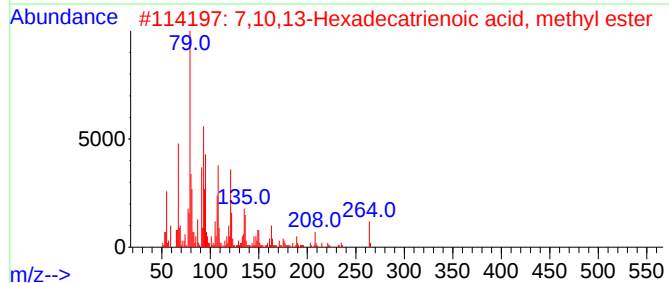
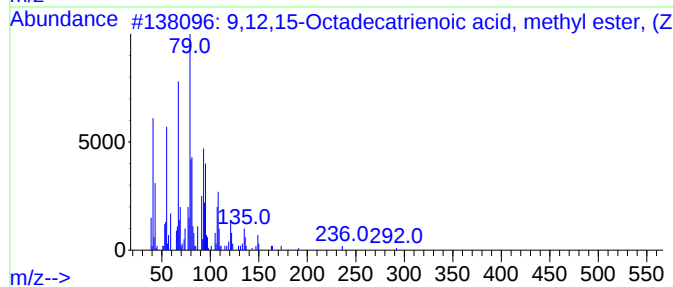
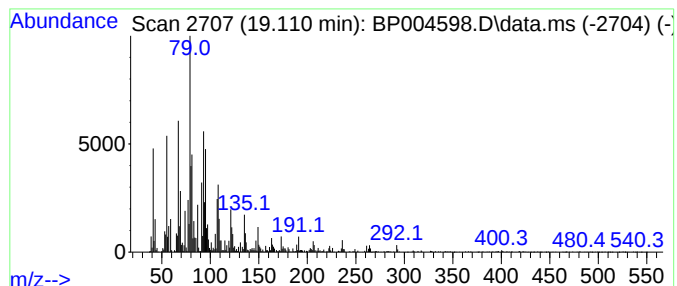
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 9,12,15-Octadecatrienoic ac... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.110	3.14 ng	483384	Phenanthrene-d10	17.222

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,12,15-Octadecatrienoic acid, m...	292	C19H32O2	000301-00-8	99		
2	7,10,13-Hexadecatrienoic acid, m...	264	C17H28O2	056554-30-4	94		
3	Methyl 8,11,14-heptadecatrienoate	278	C18H30O2	1000336-35-1	91		
4	9,12,15-Octadecatrienoic acid, (...)	278	C18H30O2	000463-40-1	90		
5	Ethyl 9,12,15-octadecatrienoate	306	C20H34O2	1000336-77-4	90		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011521\
Data File : BP004598.D
Acq On : 16 Jan 2021 03:24
Operator : CG/JU
Sample : M1120-01 2X
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
AU-04-011321

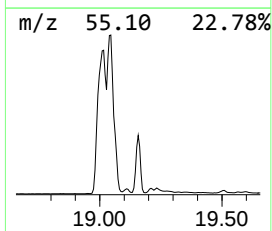
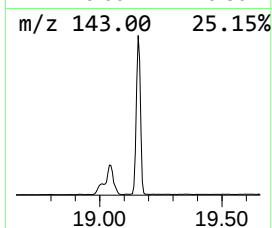
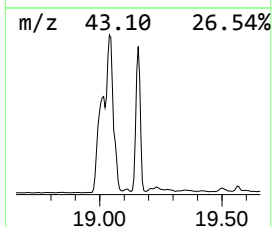
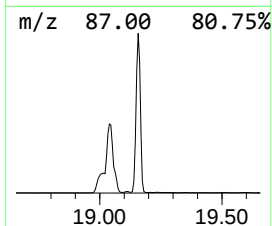
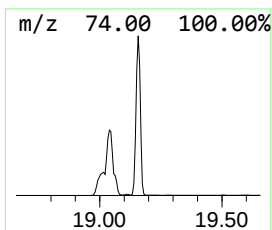
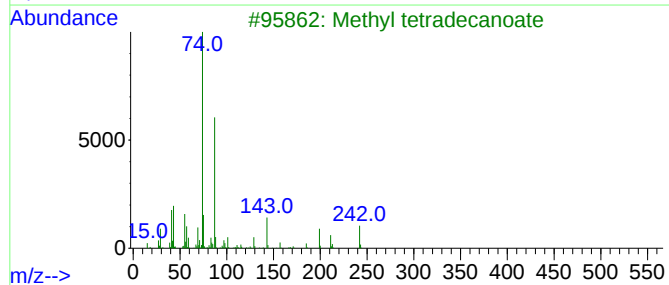
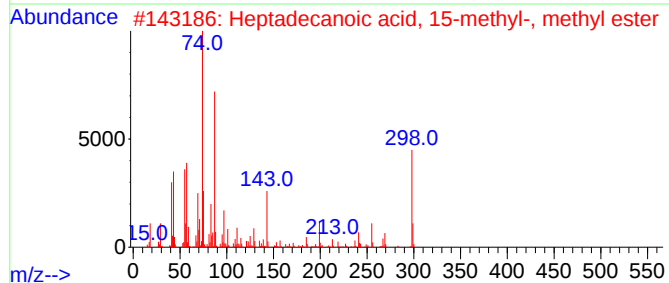
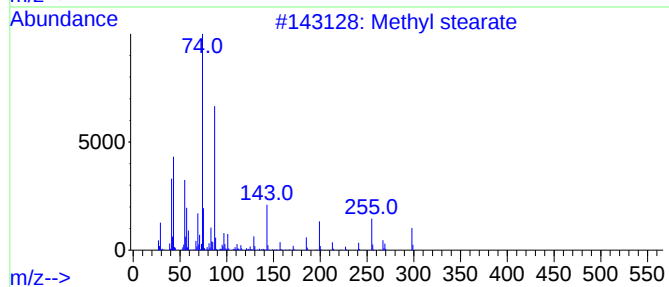
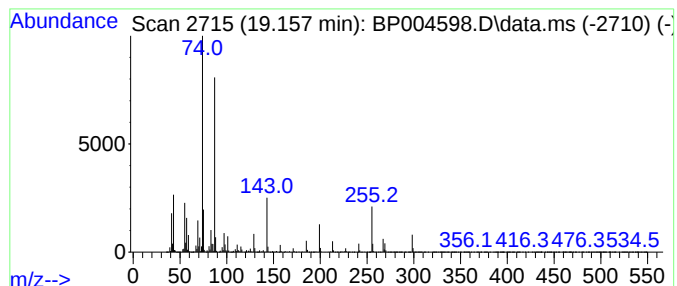
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Methyl stearate Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.157	57.86 ng	8905600	Phenanthrene-d10	17.222

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methyl stearate	298	C19H38O2	000112-61-8	98
2			Heptadecanoic acid, 15-methyl-, ...	298	C19H38O2	054833-55-5	93
3			Methyl tetradecanoate	242	C15H30O2	000124-10-7	91
4			Heptadecanoic acid, 16-methyl-, ...	298	C19H38O2	005129-61-3	90
5			Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011521\
Data File : BP004598.D
Acq On : 16 Jan 2021 03:24
Operator : CG/JU
Sample : M1120-01 2X
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
AU-04-011321

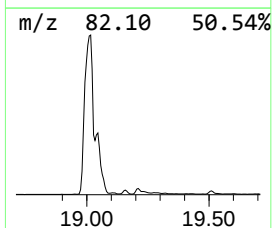
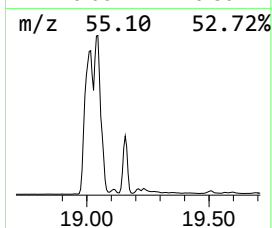
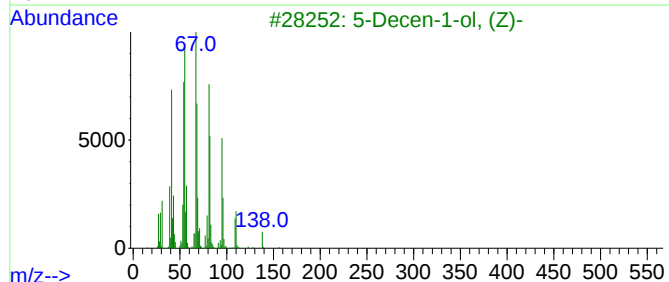
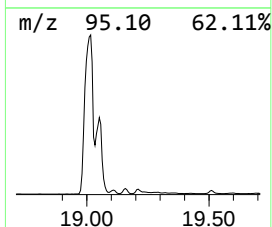
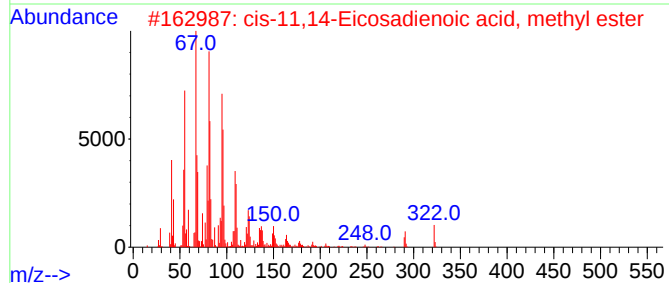
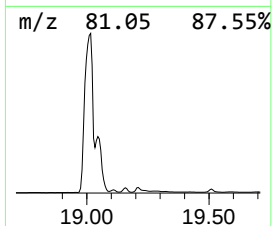
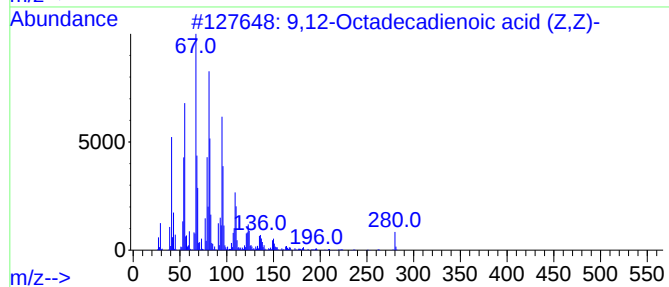
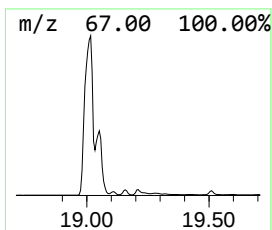
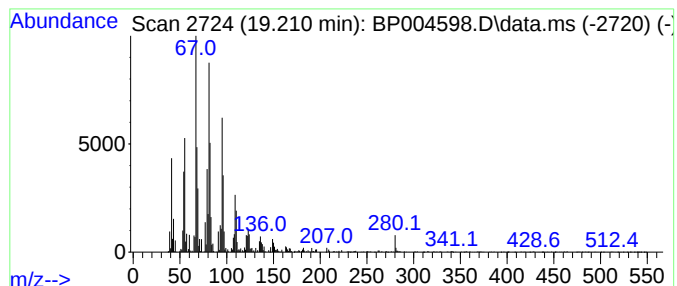
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 9,12-Octadecadienoic acid (... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.210	3.74 ng	575800	Phenanthrene-d10	17.222

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9,12-Octadecadienoic acid (Z,Z)-	280	C18H32O2	000060-33-3	99
2			cis-11,14-Eicosadienoic acid, me...	322	C21H38O2	1000333-61-8	96
3			5-Decen-1-ol, (Z)-	156	C10H20O	051652-47-2	91
4			3,4-Octadiene, 7-methyl-	124	C9H16	037050-05-8	87
5			5-Decen-1-ol, (E)-	156	C10H20O	056578-18-8	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011521\
Data File : BP004598.D
Acq On : 16 Jan 2021 03:24
Operator : CG/JU
Sample : M1120-01 2X
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
AU-04-011321

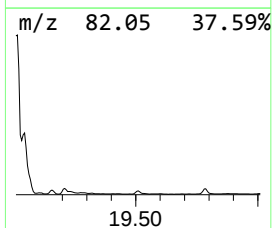
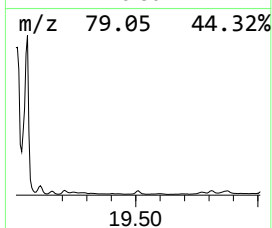
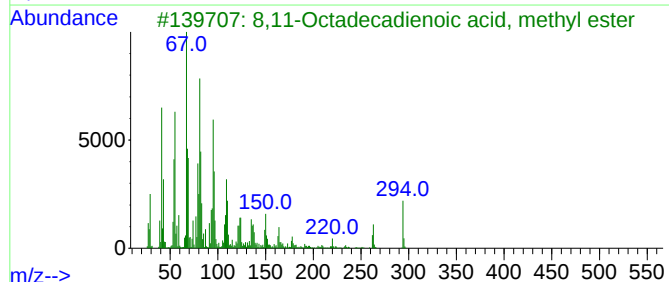
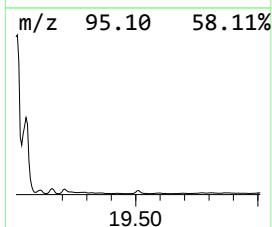
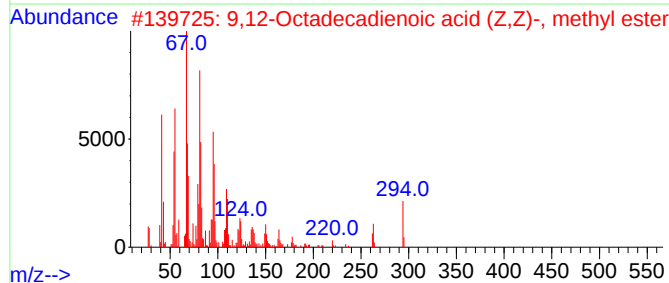
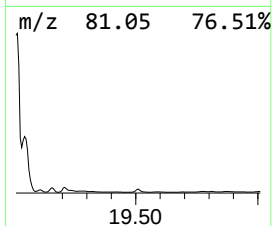
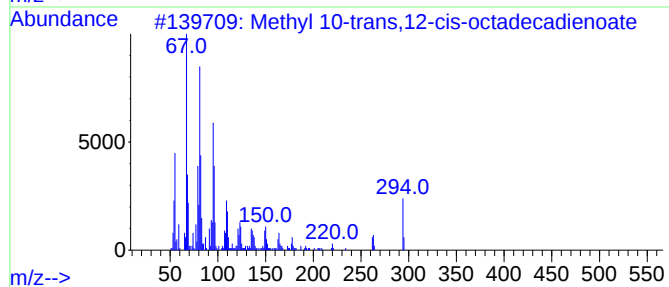
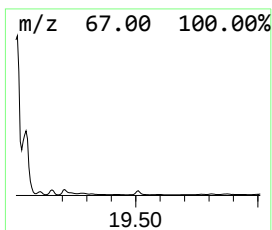
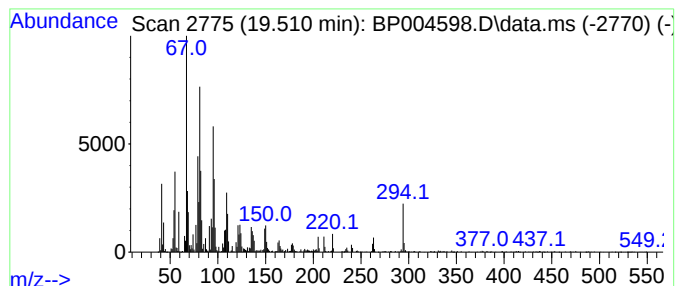
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Methyl 10-trans,12-cis-octa... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.510	3.51 ng	629915	Chrysene-d12	21.316

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methyl 10-trans,12-cis-octadecad...	294	C19H34O2	1000336-44-2	98
2			9,12-Octadecadienoic acid (Z,Z)-...	294	C19H34O2	000112-63-0	93
3			8,11-Octadecadienoic acid, methy...	294	C19H34O2	056599-58-7	91
4			Methyl 6,11-octadecadienoate	294	C19H34O2	1000336-43-2	91
5			Methyl 9-cis,11-trans-octadecadi...	294	C19H34O2	1000336-44-0	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011521\
Data File : BP004598.D
Acq On : 16 Jan 2021 03:24
Operator : CG/JU
Sample : M1120-01 2X
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
AU-04-011321

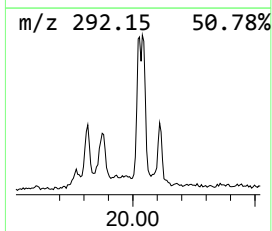
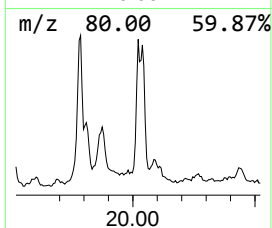
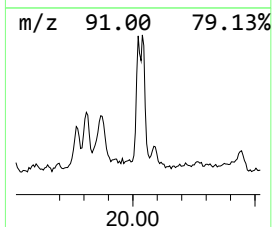
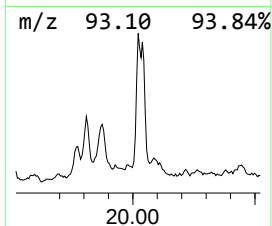
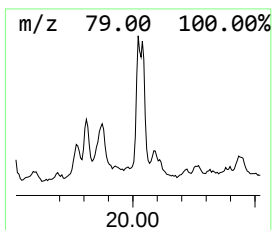
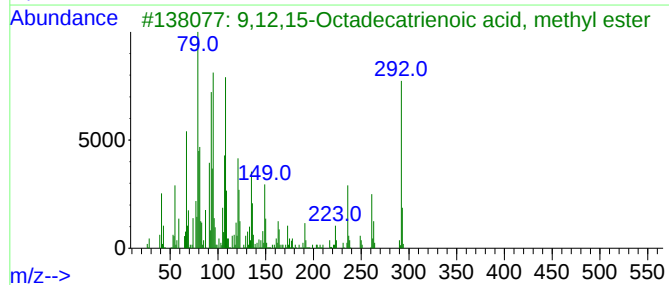
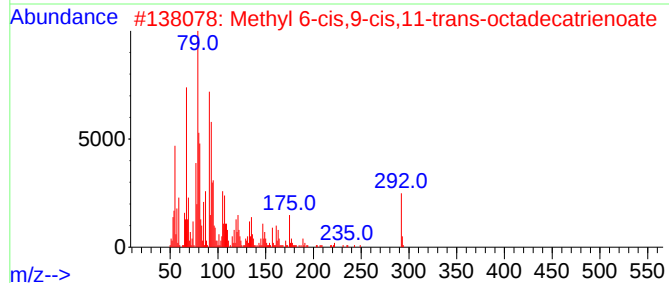
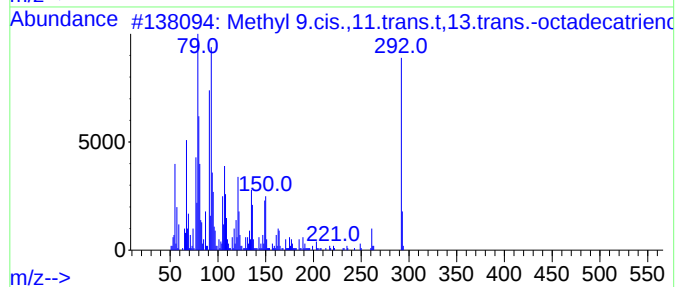
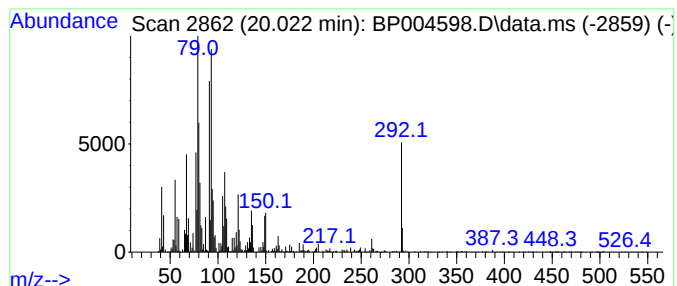
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Methyl 9.cis.,11.trans.t,13... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.022	3.06 ng	548046	Chrysene-d12	21.316

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methyl 9.cis.,11.trans.t,13.tran...	292	C19H32O2	1000336-42-6	99
2			Methyl 6-cis,9-cis,11-trans-octa...	292	C19H32O2	1000336-37-7	59
3			9,12,15-Octadecatrienoic acid, m...	292	C19H32O2	007361-80-0	58
4			Estr-5(10)-en-17-ol, 3-fluoro-6...	292	C19H29FO	024467-83-2	36
5			6,9,12-Octadecatrienoic acid, me...	292	C19H32O2	002676-41-7	32



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011521\
Data File : BP004598.D
Acq On : 16 Jan 2021 03:24
Operator : CG/JU
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Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
AU-04-011321

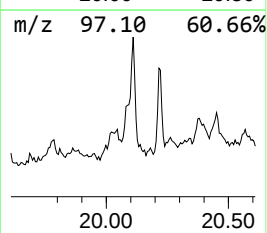
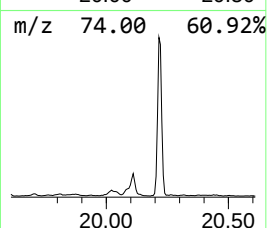
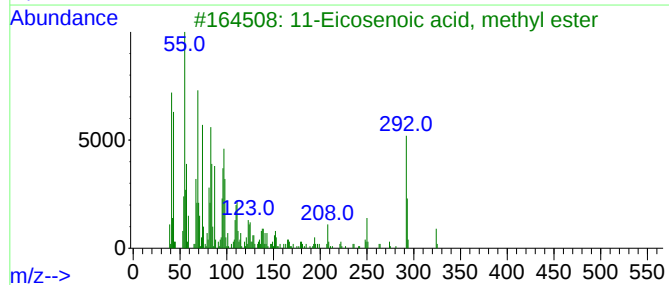
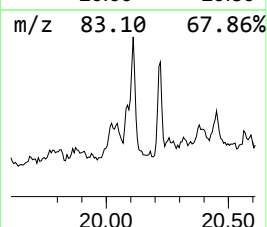
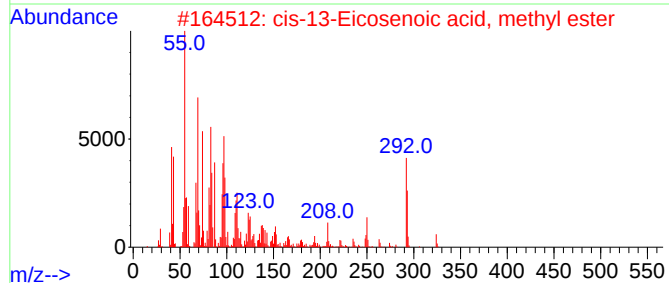
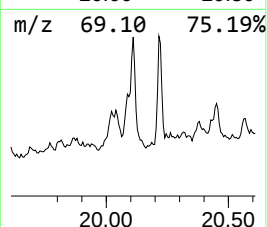
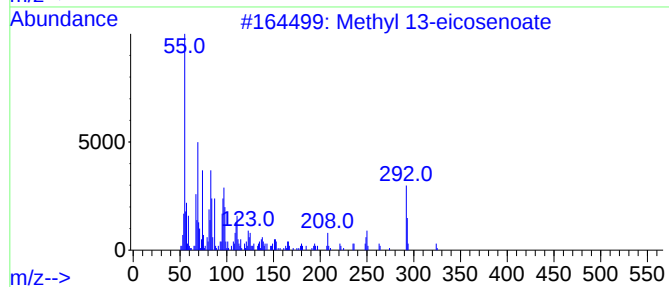
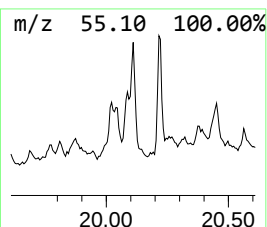
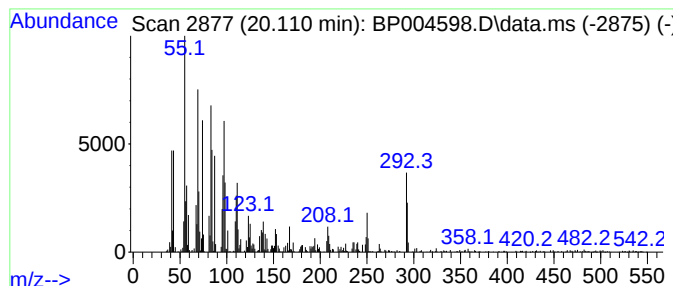
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Methyl 13-eicosenoate Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.110	2.92 ng	524368	Chrysene-d12	21.316

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methyl 13-eicosenoate	324	C21H40O2	1000336-48-4	86
2			cis-13-Eicosenoic acid, methyl e...	324	C21H40O2	1000333-52-1	83
3			11-Eicosenoic acid, methyl ester	324	C21H40O2	003946-08-5	62
4			cis-11-Eicosenoic acid, methyl e...	324	C21H40O2	1000333-63-8	58
5			Methyl 9-eicosenoate	324	C21H40O2	1000336-50-5	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011521\
Data File : BP004598.D
Acq On : 16 Jan 2021 03:24
Operator : CG/JU
Sample : M1120-01 2X
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
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ClientSampleId :
AU-04-011321

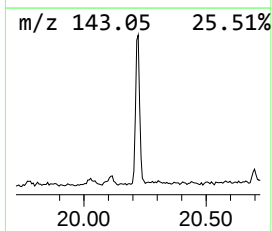
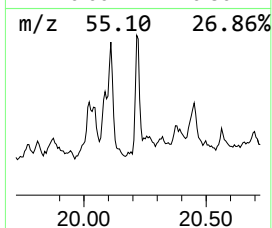
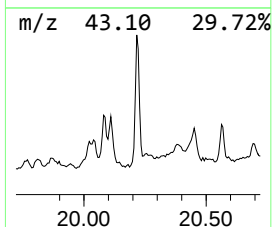
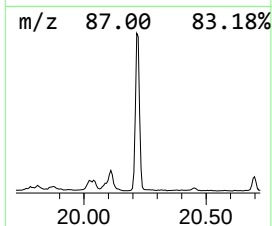
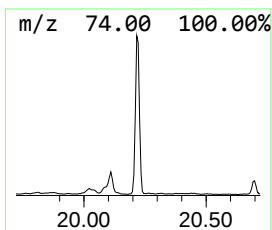
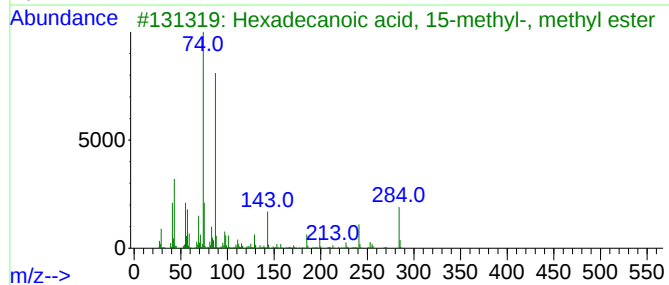
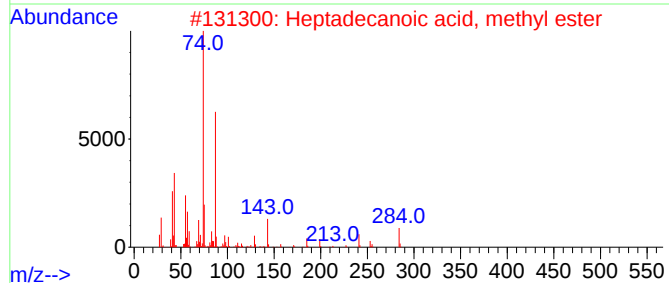
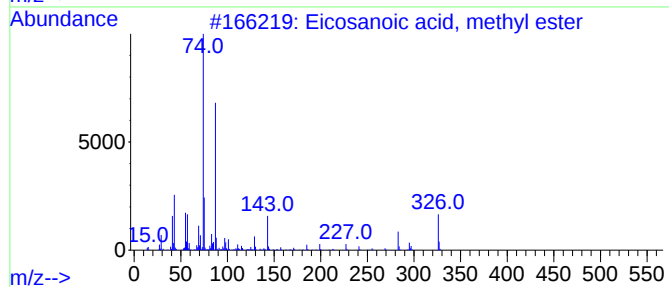
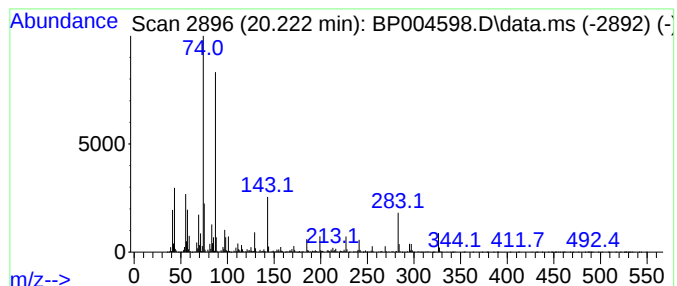
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Eicosanoic acid, methyl ester Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.221	5.90 ng	1057210	Chrysene-d12	21.316

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosanoic acid, methyl ester	326	C21H42O2	001120-28-1	98
2			Heptadecanoic acid, methyl ester	284	C18H36O2	001731-92-6	96
3			Hexadecanoic acid, 15-methyl-, m...	284	C18H36O2	006929-04-0	96
4			Hexadecanoic acid, 14-methyl-, m...	284	C18H36O2	002490-49-5	91
5			Nonadecanoic acid, methyl ester	312	C20H40O2	001731-94-8	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011521\
Data File : BP004598.D
Acq On : 16 Jan 2021 03:24
Operator : CG/JU
Sample : M1120-01 2X
Misc :
ALS Vial : 20 Sample Multiplier: 1

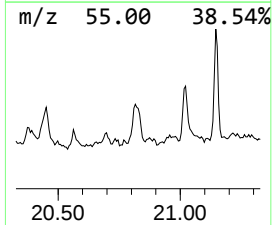
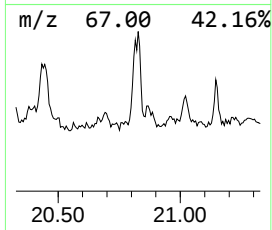
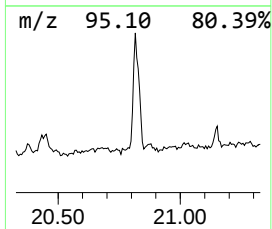
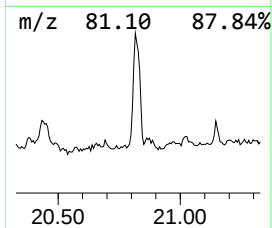
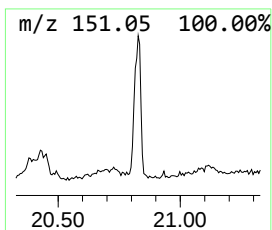
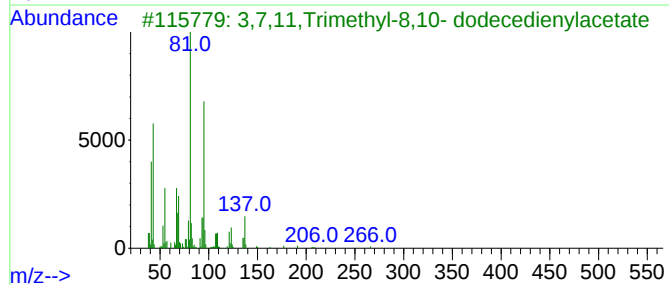
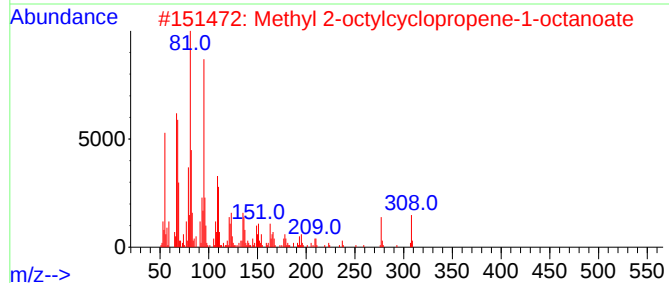
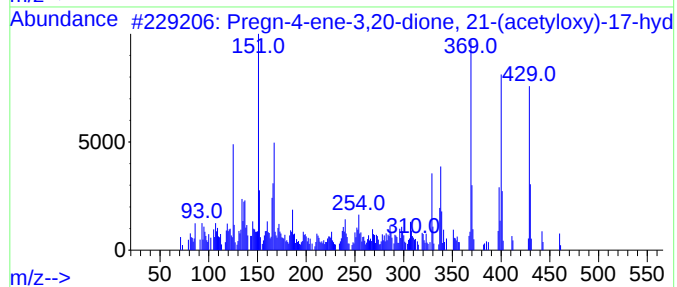
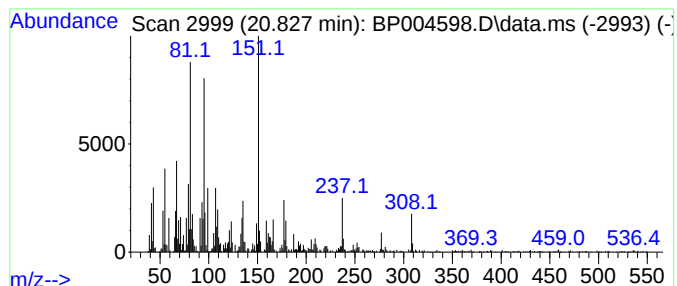
Instrument :
BNA_P
ClientSampleId :
AU-04-011321

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270E-BP011221.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Pregn-4-ene-3,20-dione, 21-... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD		R.T.	
20.827	3.88 ng	696287	Chrysene-d12		21.316	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Pregn-4-ene-3,20-dione, 21-(acet...		460	C26H40N2O5	074299-01-7	53
2	Methyl 2-octylcyclopropene-1-oct...		308	C20H36O2	003220-60-8	53
3	3,7,11,Trimethyl-8,10- dodecedie...		266	C17H30O2	1000374-10-3	46
4	Cedrol		222	C15H26O	000077-53-2	38
5	Phenol, 2,6-dimethyl-4-nitroso-		151	C8H9NO2	013331-93-6	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011521\
Data File : BP004598.D
Acq On : 16 Jan 2021 03:24
Operator : CG/JU
Sample : M1120-01 2X
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
AU-04-011321

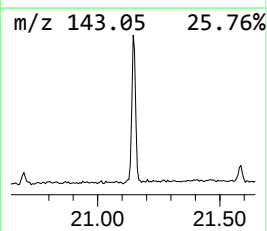
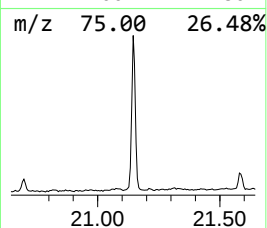
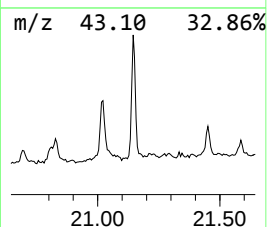
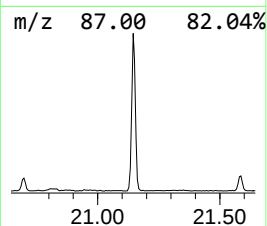
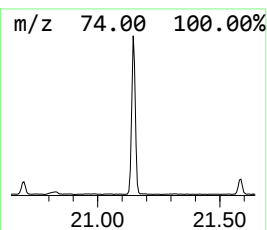
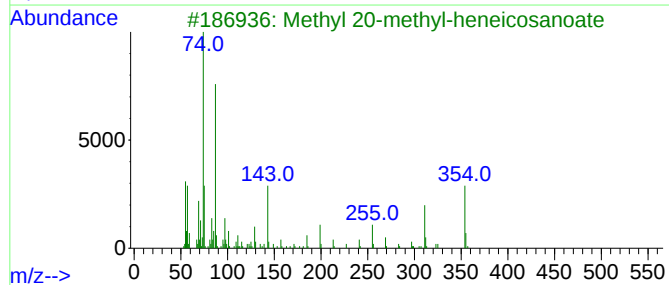
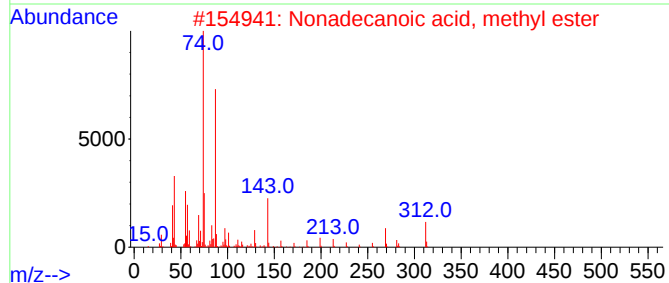
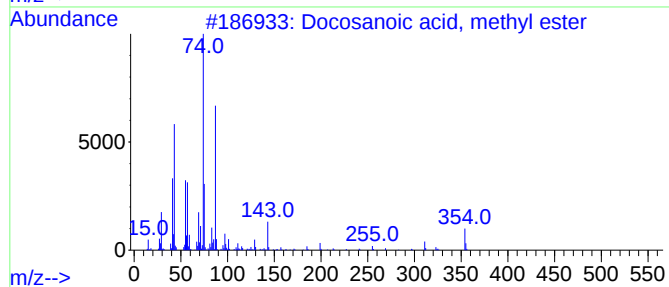
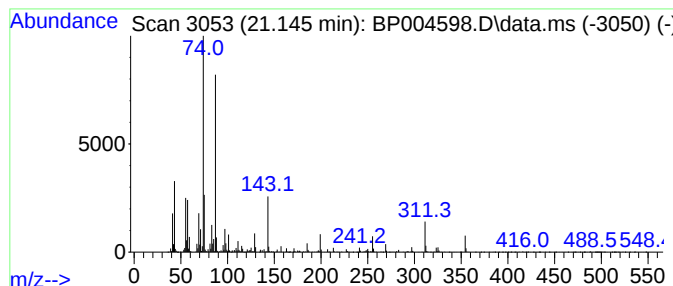
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270E-BP011221.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Docosanoic acid, methyl ester Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.145	5.17 ng	927615	Chrysene-d12	21.316

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Docosanoic acid, methyl ester	354	C23H46O2	000929-77-1	99
2			Nonadecanoic acid, methyl ester	312	C20H40O2	001731-94-8	93
3			Methyl 20-methyl-heneicosanoate	354	C23H46O2	1000336-47-4	91
4			Hexacosanoic acid, methyl ester	410	C27H54O2	005802-82-4	90
5			Heptadecanoic acid, 16-methyl-, ...	298	C19H38O2	005129-61-3	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011521\
 Data File : BP004598.D
 Acq On : 16 Jan 2021 03:24
 Operator : CG/JU
 Sample : M1120-01 2X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 AU-04-011321

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270E-BP011221.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
unknown12.257	12.257	5.3	ng	515540	2	10.599	1932790	20.0
Hexadecanoic ac...	17.898	114.6	ng	17643800	4	17.222	3078240	20.0
n-Hexadecanoic ...	18.116	4.3	ng	657503	4	17.222	3078240	20.0
8,11-Octadecadi...	19.016	273.4	ng	42083700	4	17.222	3078240	20.0
16-Octadecenoic...	19.045	192.7	ng	29662900	4	17.222	3078240	20.0
9,12,15-Octadec...	19.110	3.1	ng	483384	4	17.222	3078240	20.0
Methyl stearate	19.157	57.9	ng	8905600	4	17.222	3078240	20.0
9,12-Octadecadi...	19.210	3.7	ng	575800	4	17.222	3078240	20.0
Methyl 10-trans...	19.510	3.5	ng	629915	5	21.316	3586650	20.0
Methyl 9.cis.,1...	20.022	3.1	ng	548046	5	21.316	3586650	20.0
Methyl 13-eicos...	20.110	2.9	ng	524368	5	21.316	3586650	20.0
Eicosanoic acid...	20.221	5.9	ng	1057210	5	21.316	3586650	20.0
Pregn-4-ene-3,2...	20.827	3.9	ng	696287	5	21.316	3586650	20.0
Docosanoic acid...	21.145	5.2	ng	927615	5	21.316	3586650	20.0