

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040621\
 Data File : BP005195.D
 Acq On : 06 Apr 2021 14:20
 Operator : CG/JU
 Sample : M1894-03
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 0154-FIELD-BLANK

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-BP032321.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP005195.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.293	369	375	384	rVB	245332	347877	21.96%	4.006%
2	6.875	638	644	655	rVB	132746	217720	13.75%	2.507%
3	7.693	777	783	790	rBV	87198	133507	8.43%	1.538%
4	8.858	973	981	987	rBV	321021	512928	32.39%	5.907%
5	10.487	1250	1258	1265	rBV	110336	184846	11.67%	2.129%
6	12.969	1673	1680	1686	rBV	759383	1068524	67.47%	12.306%
7	13.810	1813	1823	1831	rBV2	14412	23030	1.45%	0.265%
8	14.351	1909	1915	1925	rVB	164736	243245	15.36%	2.801%
9	15.745	2144	2152	2160	rBV	24829	45251	2.86%	0.521%
10	15.851	2164	2170	2178	rVV	545681	790234	49.89%	9.101%
11	17.116	2378	2385	2392	rBV	197362	282218	17.82%	3.250%
12	18.016	2532	2538	2547	rBV	289752	411233	25.96%	4.736%
13	18.081	2547	2549	2559	rVB2	15080	25773	1.63%	0.297%
14	19.139	2725	2729	2744	rVB	239692	393744	24.86%	4.535%
15	19.257	2745	2749	2757	rVB	193864	232932	14.71%	2.683%
16	19.404	2771	2774	2778	rVB	34655	40679	2.57%	0.468%
17	19.692	2817	2823	2828	rBV	1352382	1583819	100.00%	18.241%
18	20.363	2928	2937	2951	rBV	253441	620694	39.19%	7.149%
19	20.463	2951	2954	2961	rVB	53817	70568	4.46%	0.813%
20	20.863	3019	3022	3029	rVB2	59250	95988	6.06%	1.105%
21	20.922	3029	3032	3041	rVB2	145773	208559	13.17%	2.402%
22	20.992	3041	3044	3049	rBV	36686	39374	2.49%	0.453%
23	21.122	3064	3066	3073	rVB3	39374	47967	3.03%	0.552%
24	21.210	3076	3081	3087	rBV	268588	329799	20.82%	3.798%
25	21.304	3094	3097	3100	rBV2	38897	47828	3.02%	0.551%
26	22.263	3256	3260	3273	rVB	239643	339947	21.46%	3.915%
27	23.492	3463	3469	3478	rVB	176305	344542	21.75%	3.968%

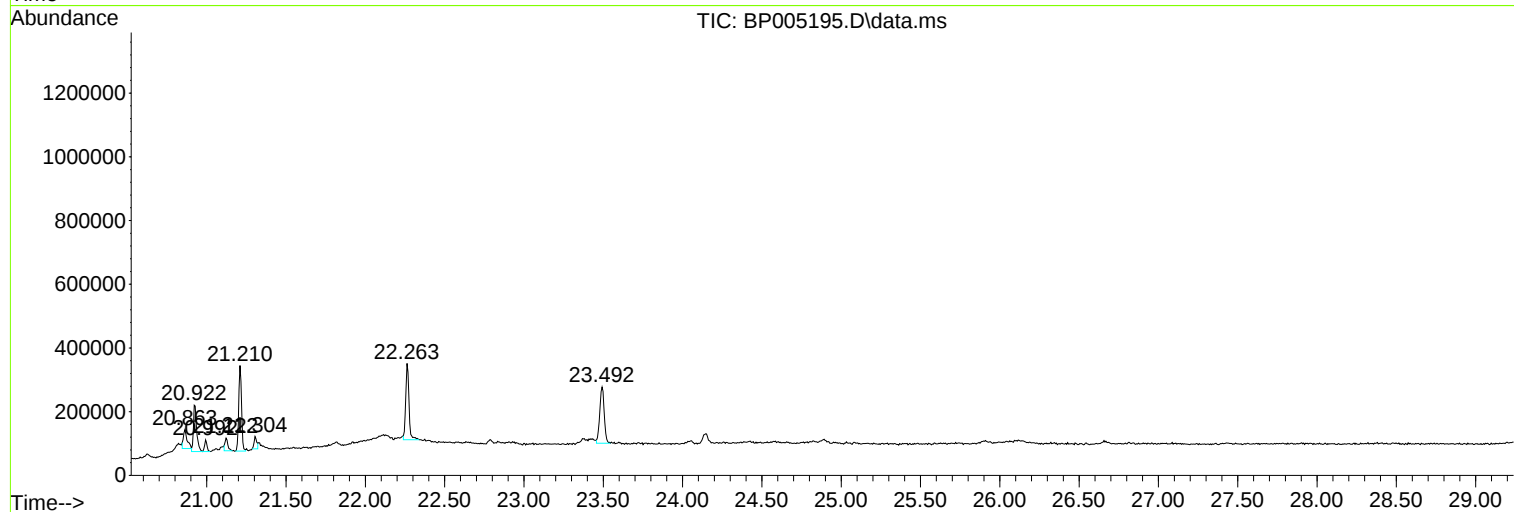
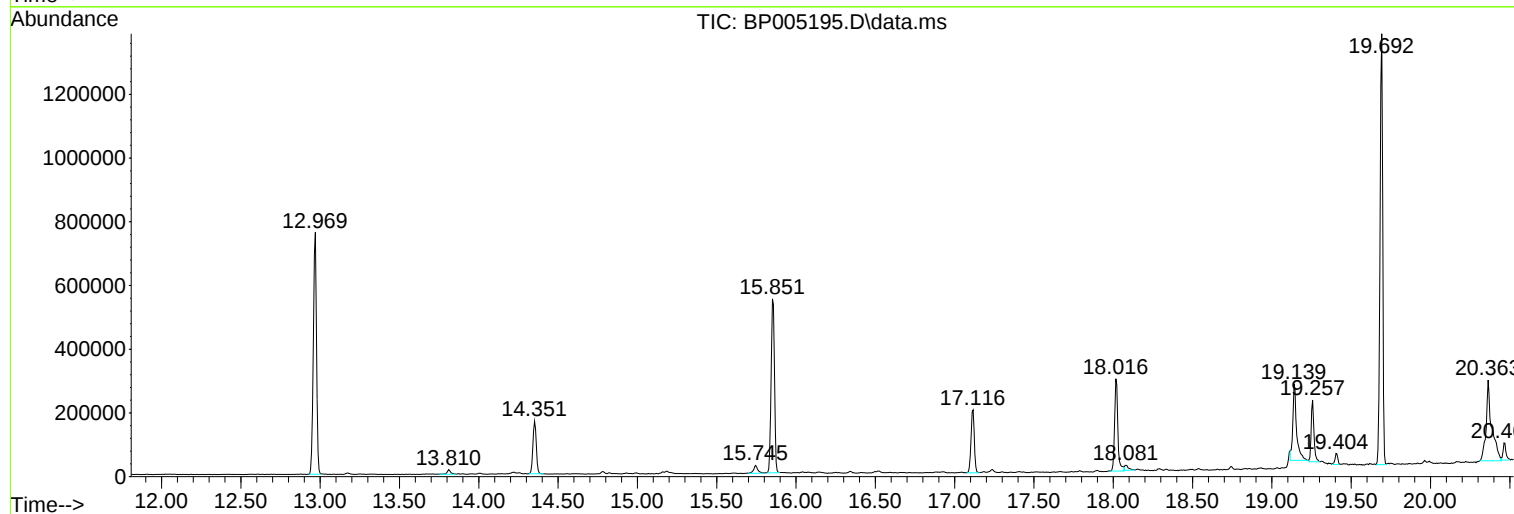
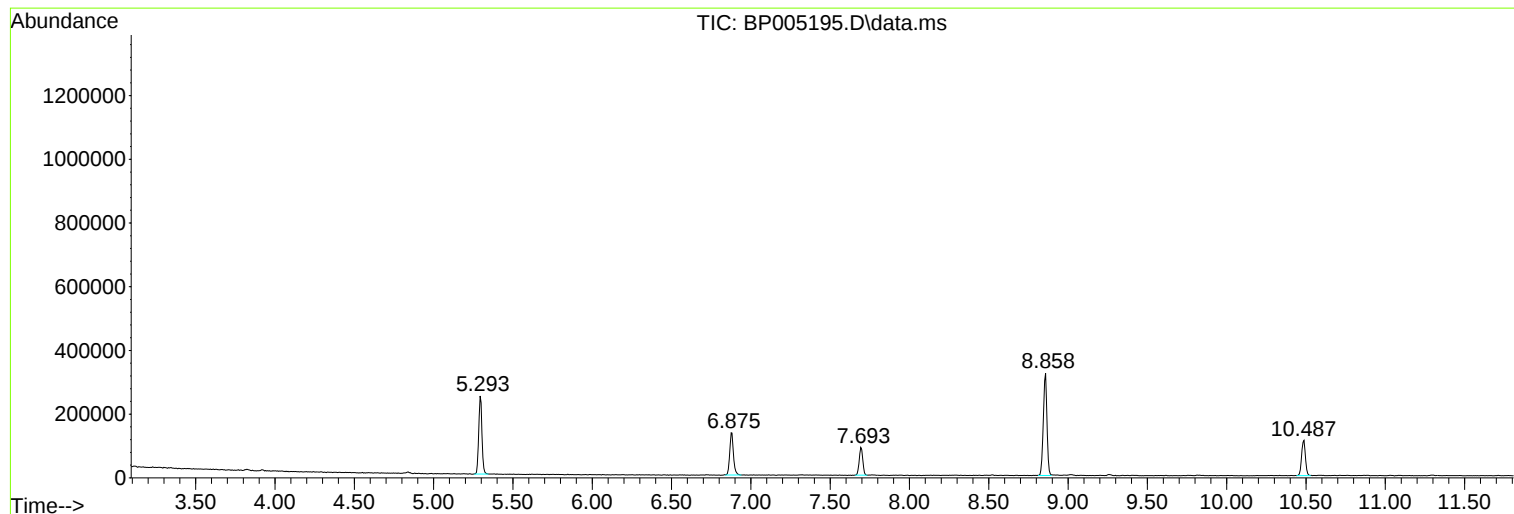
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Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP032321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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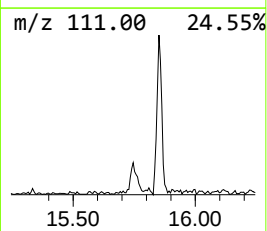
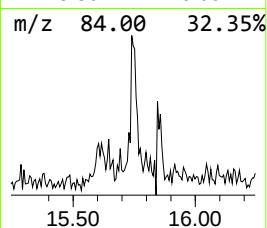
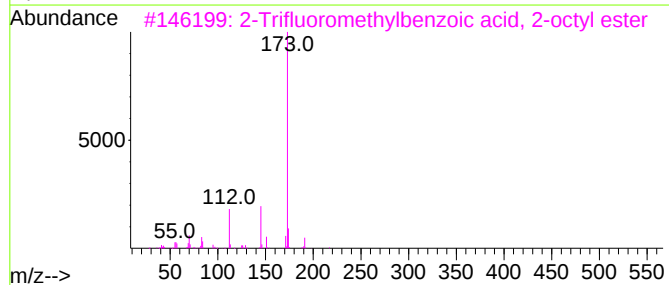
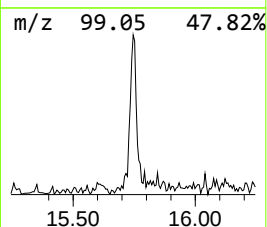
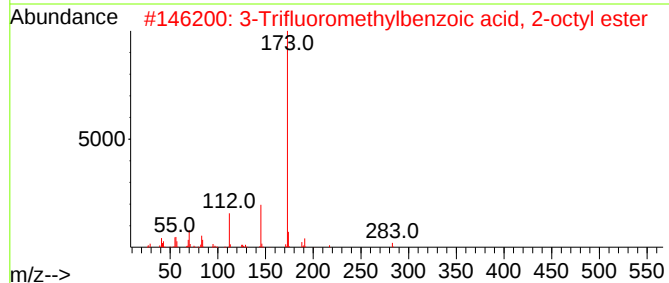
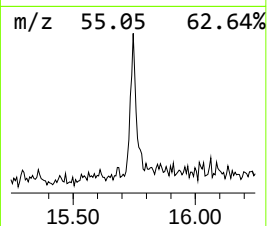
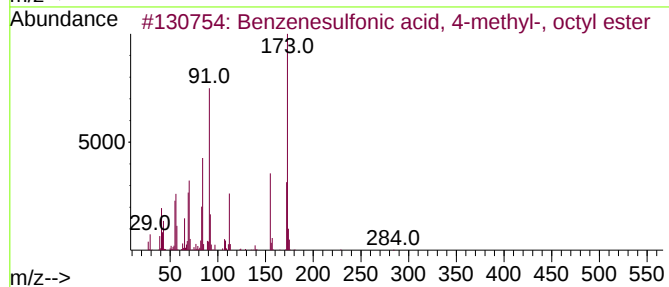
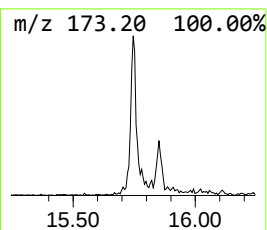
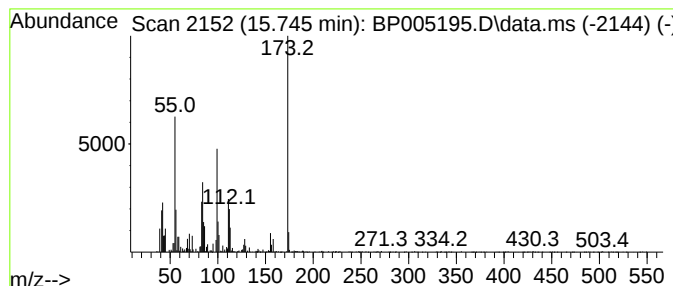
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown15.74 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.746	3.21 ng	45251	Phenanthrene-d10	17.116

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzenesulfonic acid, 4-methyl-,...	284	C15H24O3S	003386-35-4	47
2			3-Trifluoromethylbenzoic acid, 2...	302	C16H21F3O2	1000282-68-6	43
3			2-Trifluoromethylbenzoic acid, 2...	302	C16H21F3O2	1000282-64-8	37
4			2,6-Difluoro-3-methylbenzoic aci...	424	C26H42F2O2	1000338-85-5	35
5			2-Trifluoromethylbenzoic acid, 2...	400	C23H35F3O2	1000282-03-5	32



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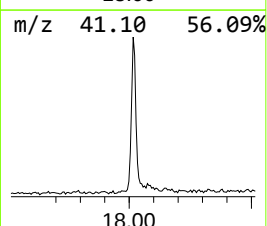
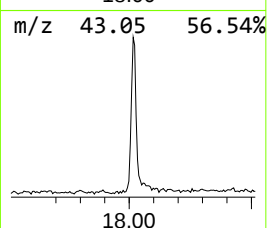
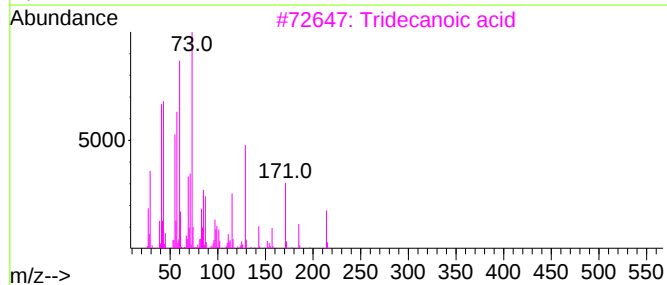
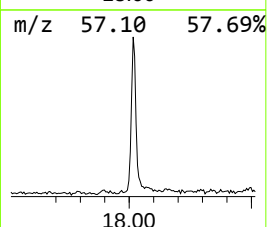
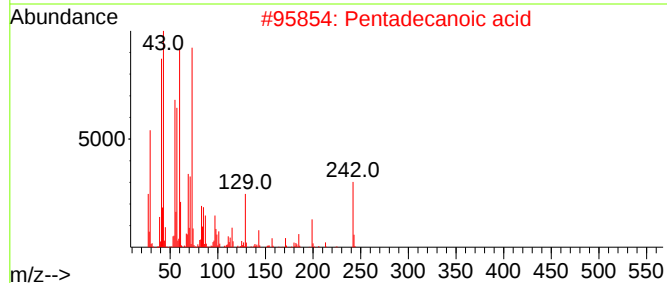
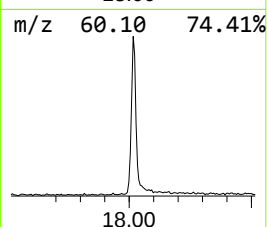
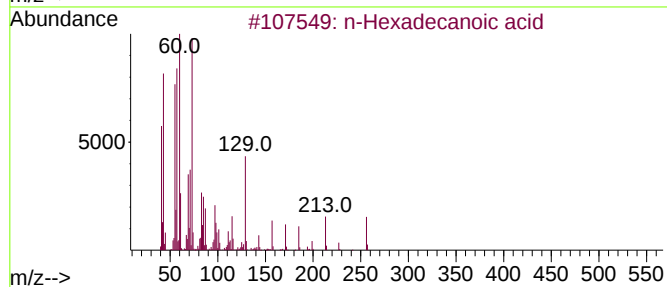
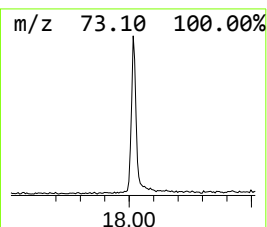
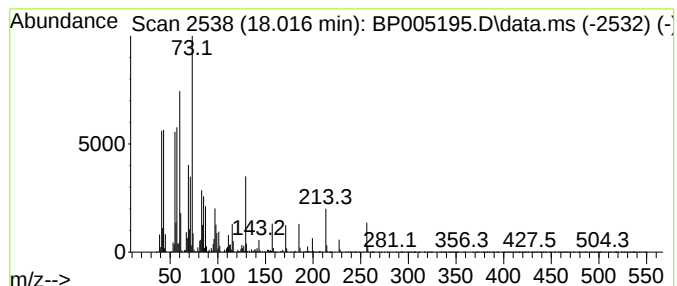
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP032321.M
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 2 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.016	29.14 ng	411233	Phenanthrene-d10	17.116

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2			Pentadecanoic acid	242	C15H30O2	001002-84-2	91
3			Tridecanoic acid	214	C13H26O2	000638-53-9	89
4			Tetradecanoic acid	228	C14H28O2	000544-63-8	59
5			Dodecanoic acid	200	C12H24O2	000143-07-7	58



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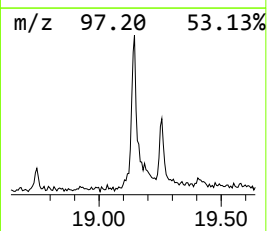
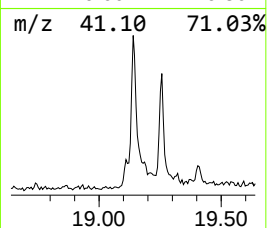
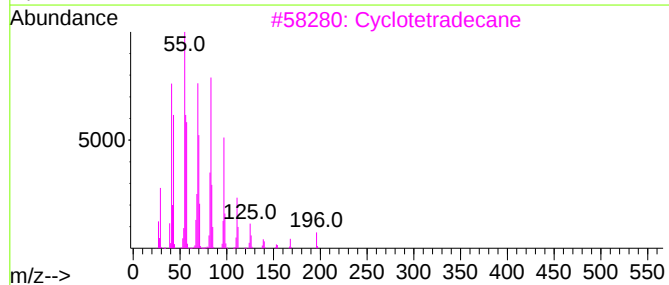
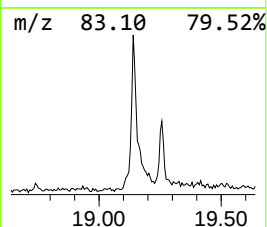
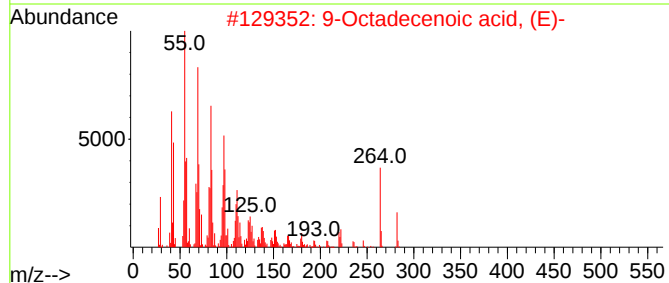
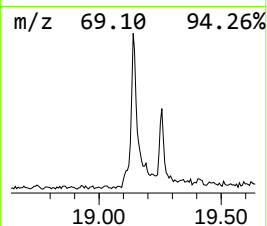
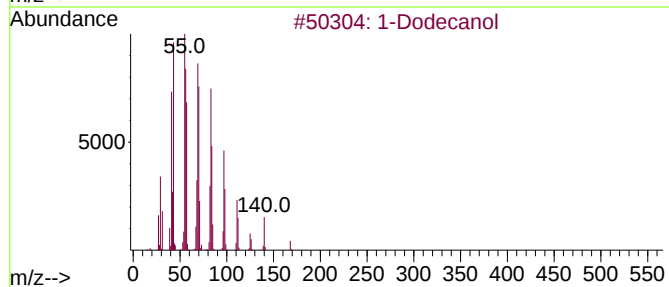
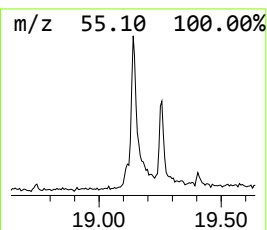
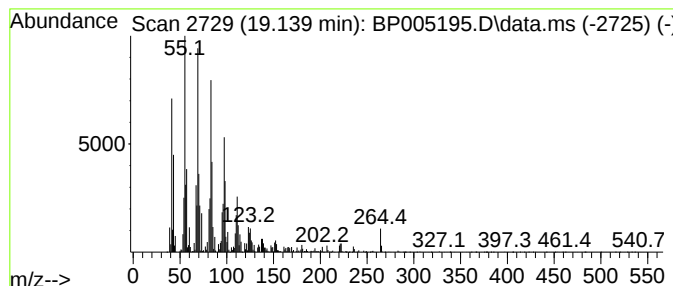
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 1-Dodecanol Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.139	27.90 ng	393744	Phenanthrene-d10	17.116

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Dodecanol	186	C12H26O	000112-53-8	83
2			9-Octadecenoic acid, (E)-	282	C18H34O2	000112-79-8	81
3			Cyclotetradecane	196	C14H28	000295-17-0	72
4			Erucic acid	338	C22H42O2	000112-86-7	62
5			Octadec-9-enoic acid	282	C18H34O2	1000190-13-7	60



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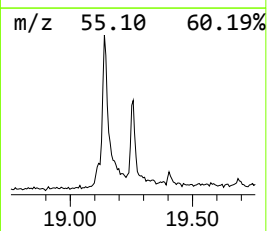
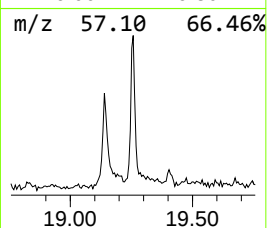
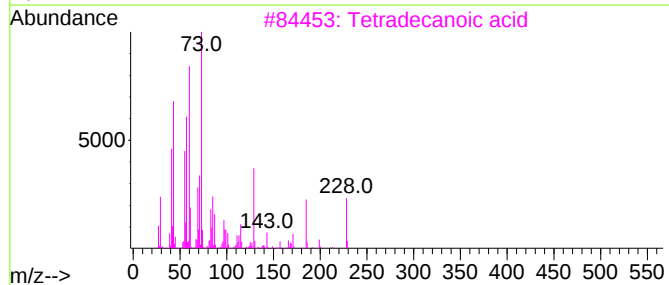
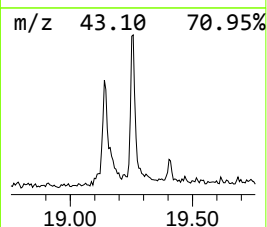
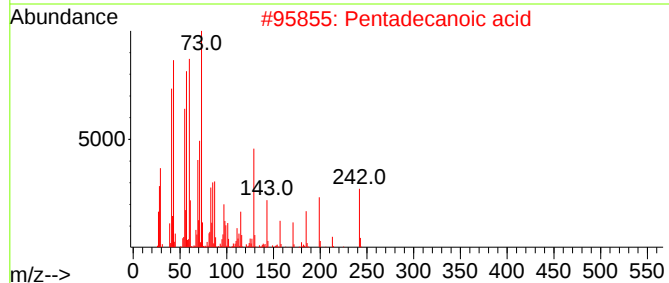
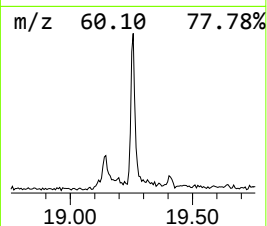
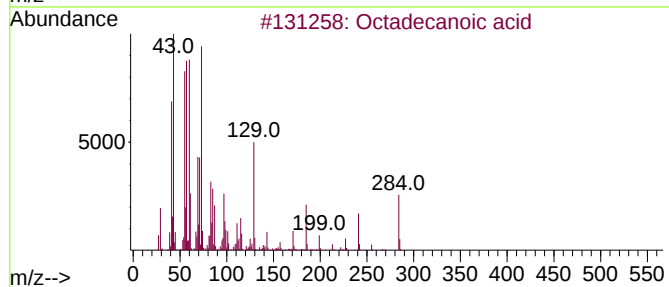
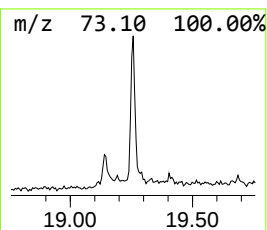
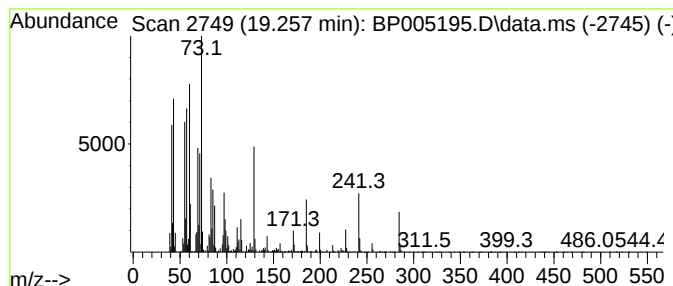
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 4 Octadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.257	14.13 ng	232932	Chrysene-d12	21.210

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	90
3			Tetradecanoic acid	228	C14H28O2	000544-63-8	83
4			n-Decanoic acid	172	C10H20O2	000334-48-5	70
5			Tridecanoic acid	214	C13H26O2	000638-53-9	68



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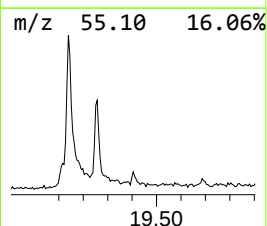
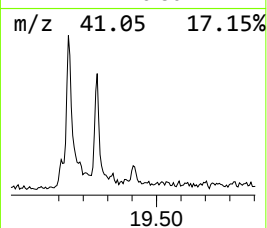
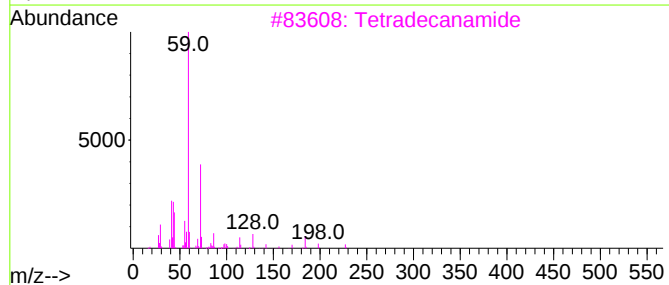
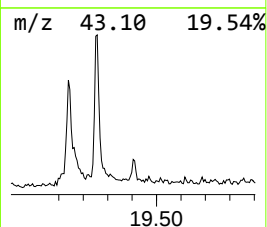
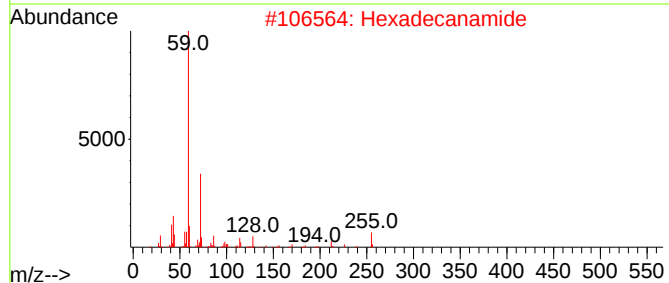
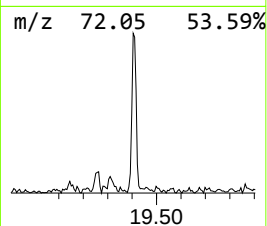
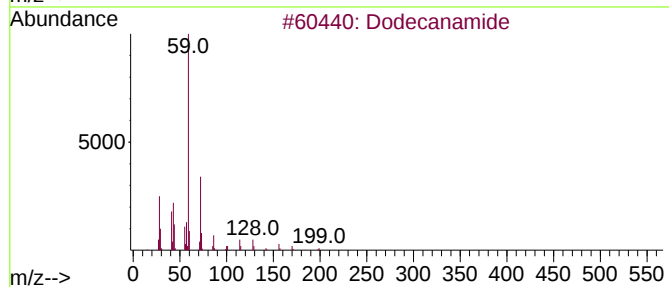
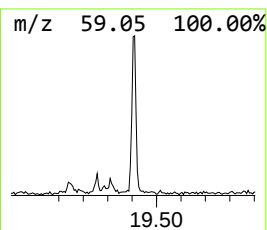
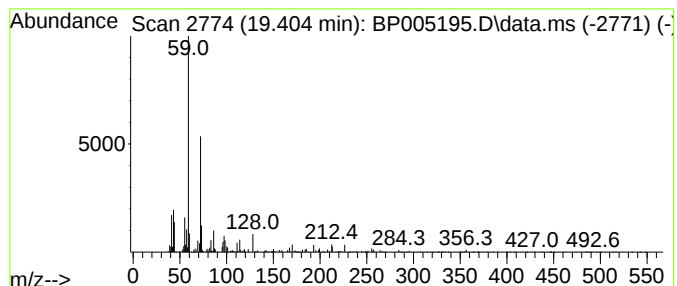
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Peak Number 5 Dodecanamide Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.404	2.47 ng	40679	Chrysene-d12	21.210

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecanamide	199	C12H25NO	001120-16-7	90
2			Hexadecanamide	255	C16H33NO	000629-54-9	87
3			Tetradecanamide	227	C14H29NO	000638-58-4	78
4			7-Nonenamide	155	C9H17NO	090949-53-4	64
5			Octadecanamide	283	C18H37NO	000124-26-5	64



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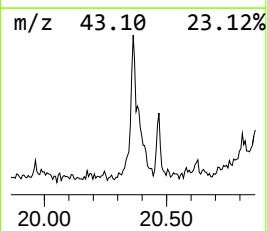
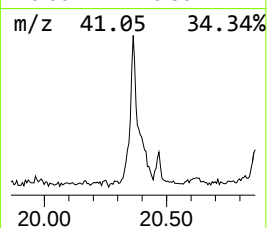
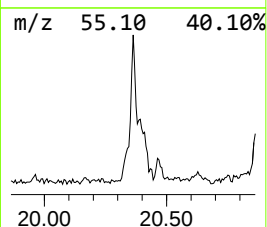
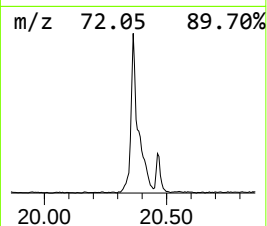
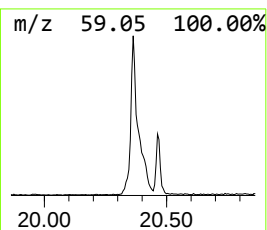
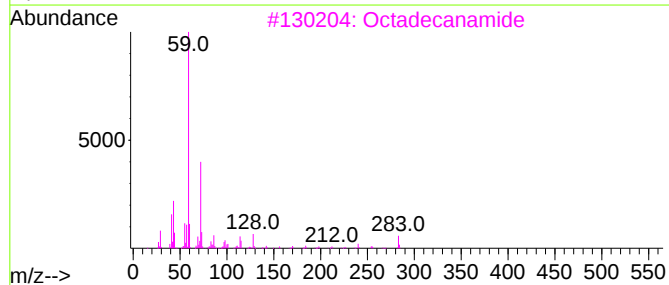
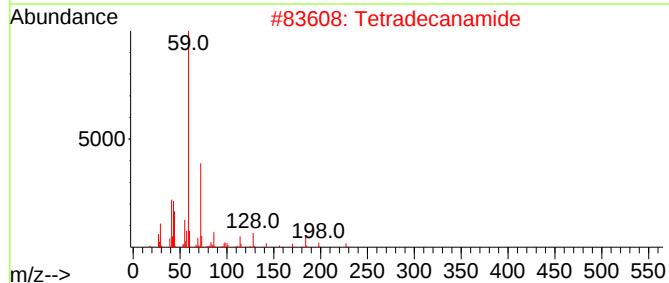
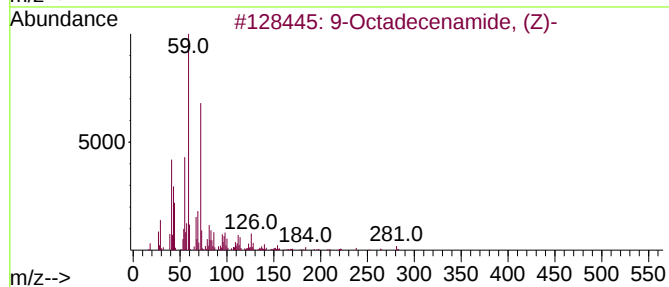
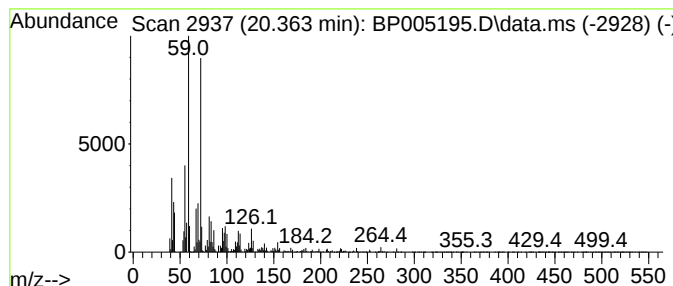
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 9-Octadecenamide, (Z)- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.363	37.64 ng	620694	Chrysene-d12	21.210

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	98
2			Tetradecanamide	227	C14H29NO	000638-58-4	53
3			Octadecanamide	283	C18H37NO	000124-26-5	38
4			Octanamide	143	C8H17NO	000629-01-6	38
5			Pentanamide, 4-methyl-	115	C6H13NO	001119-29-5	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040621\
Data File : BP005195.D
Acq On : 06 Apr 2021 14:20
Operator : CG/JU
Sample : M1894-03
Misc :
ALS Vial : 6 Sample Multiplier: 1

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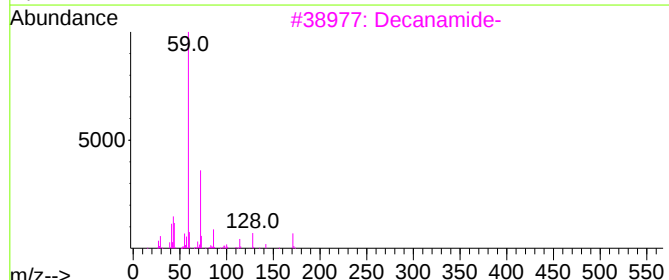
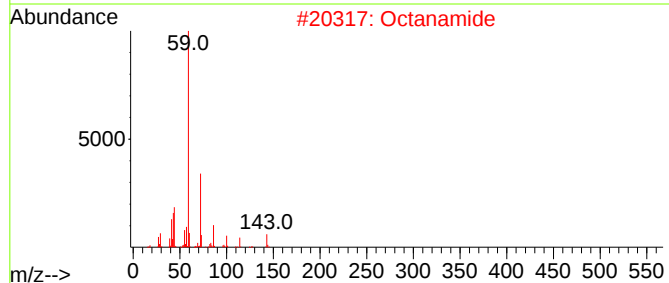
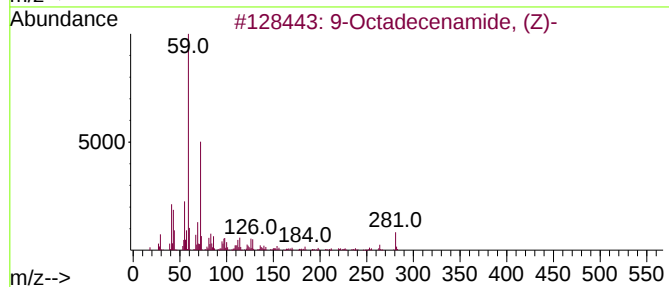
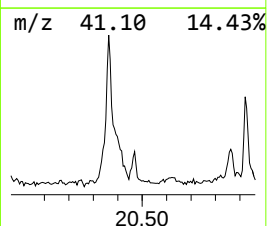
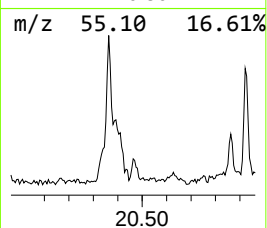
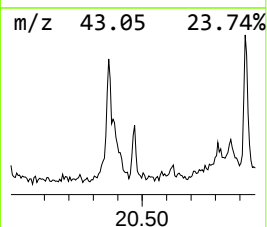
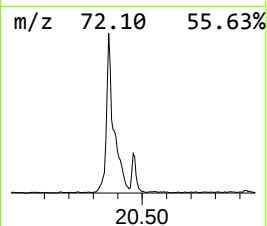
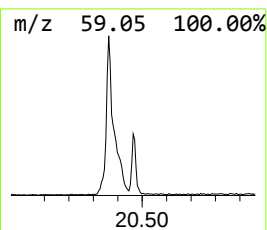
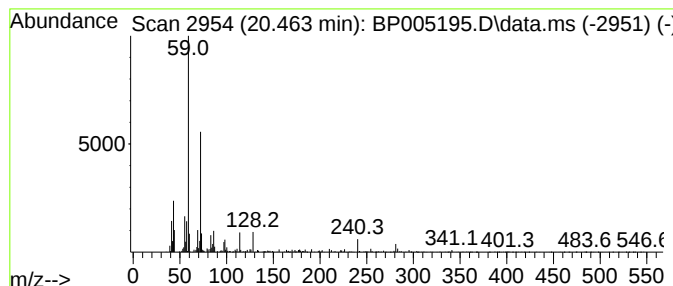
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Octanamide Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.463	4.28 ng	70568	Chrysene-d12	21.210

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Octadecenamide, (Z)-			281	C18H35NO	000301-02-0	90
2	Octanamide			143	C8H17NO	000629-01-6	59
3	Decanamide-			171	C10H21NO	002319-29-1	59
4	Dodecanamide			199	C12H25NO	001120-16-7	50
5	Undecanamide, 11-bromo-			263	C11H22BrNO	005875-26-3	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040621\
Data File : BP005195.D
Acq On : 06 Apr 2021 14:20
Operator : CG/JU
Sample : M1894-03
Misc :
ALS Vial : 6 Sample Multiplier: 1

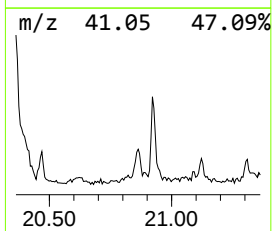
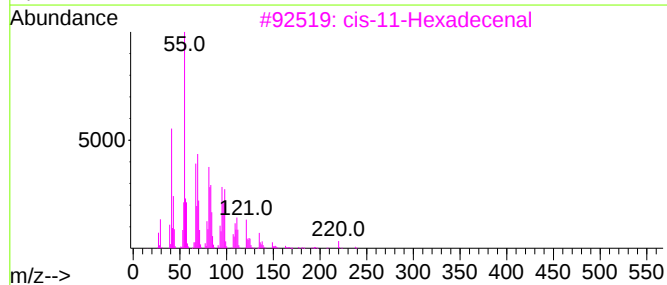
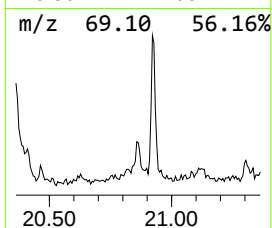
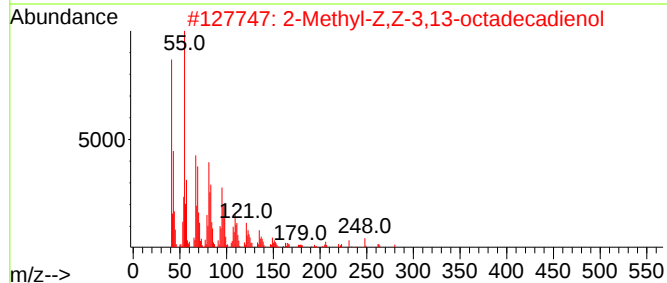
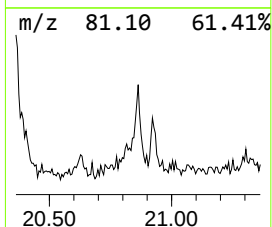
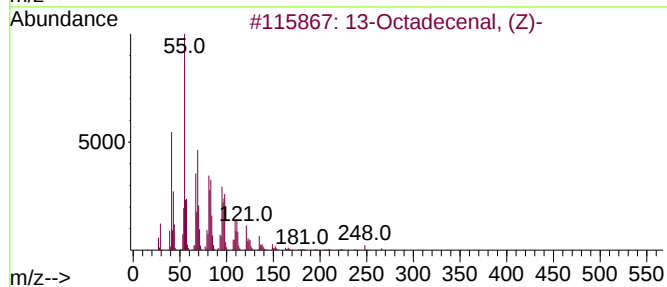
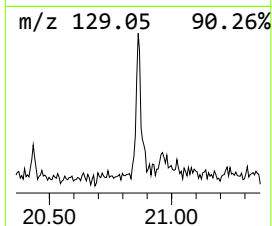
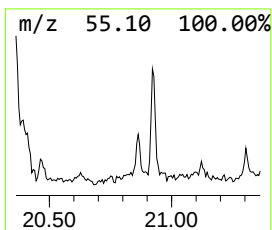
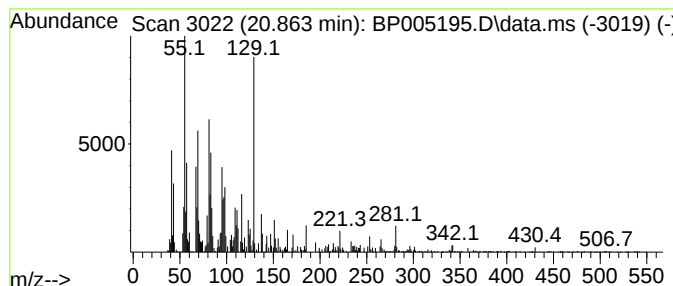
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP032321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 13-Octadecenal, (Z)- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD		R.T.	
20.863	5.82 ng	95988	Chrysene-d12		21.210	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	13-Octadecenal, (Z)-		266	C18H34O	058594-45-9	89
2	2-Methyl-Z,Z-3,13-octadecadienol		280	C19H36O	1000130-90-5	80
3	cis-11-Hexadecenal		238	C16H30O	053939-28-9	70
4	9-Octadecenal, (Z)-		266	C18H34O	002423-10-1	62
5	6-Octadecenoic acid		282	C18H34O2	1000336-66-8	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040621\
Data File : BP005195.D
Acq On : 06 Apr 2021 14:20
Operator : CG/JU
Sample : M1894-03
Misc :
ALS Vial : 6 Sample Multiplier: 1

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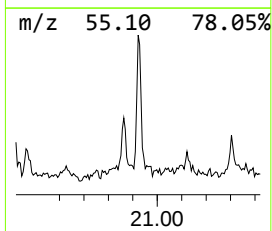
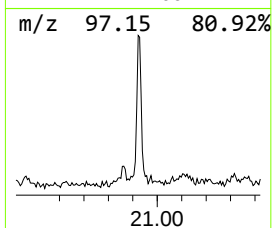
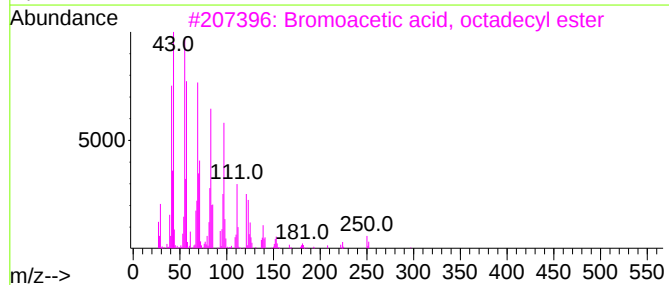
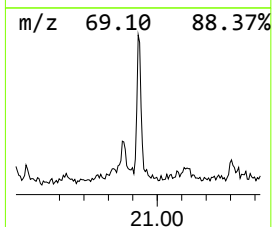
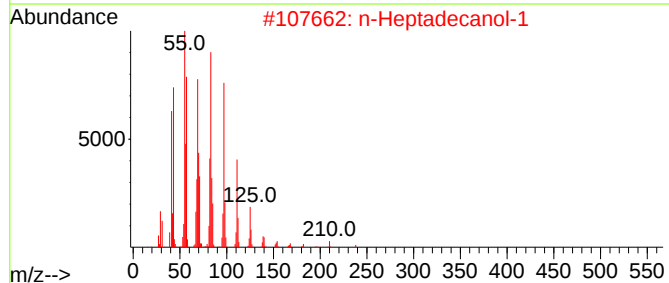
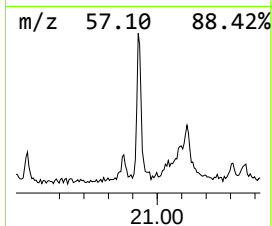
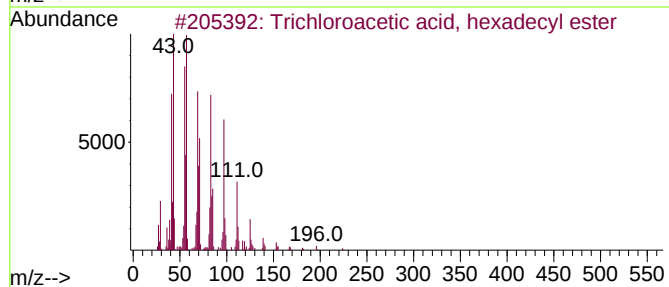
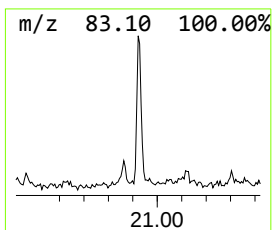
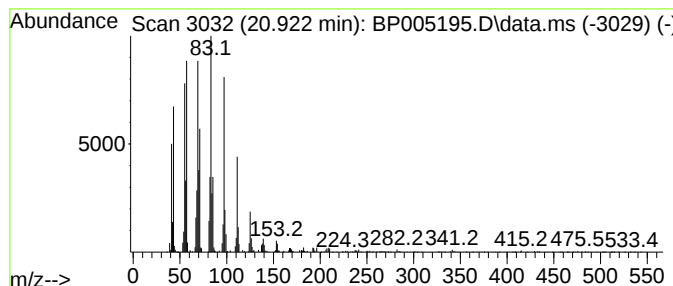
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Trichloroacetic acid, hexad... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.922	12.65 ng	208559	Chrysene-d12	21.210

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	91
2			n-Heptadecanol-1	256	C17H36O	001454-85-9	87
3			Bromoacetic acid, octadecyl ester	390	C20H39BrO2	018992-03-5	87
4			3-Eicosene, (E)-	280	C20H40	074685-33-9	87
5			Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040621\
Data File : BP005195.D
Acq On : 06 Apr 2021 14:20
Operator : CG/JU
Sample : M1894-03
Misc :
ALS Vial : 6 Sample Multiplier: 1

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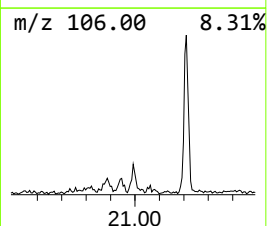
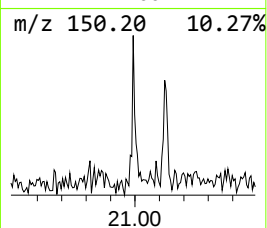
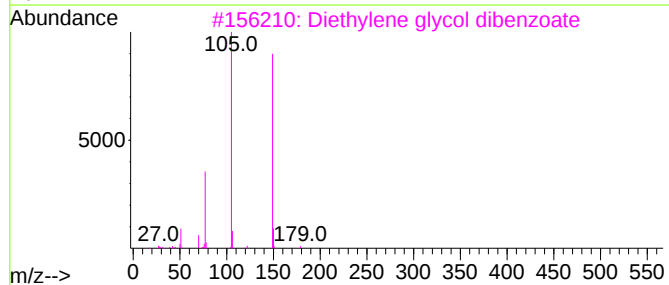
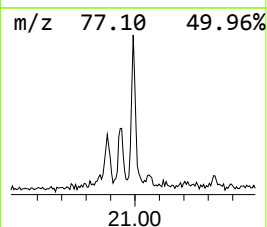
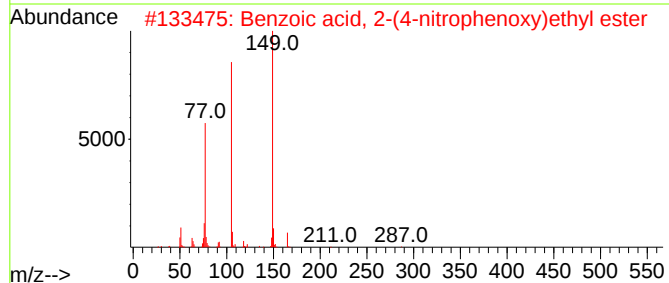
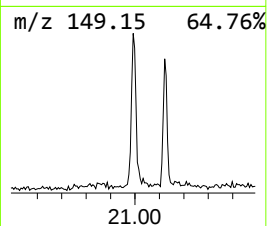
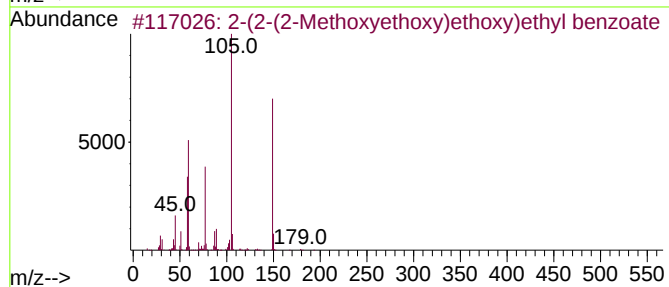
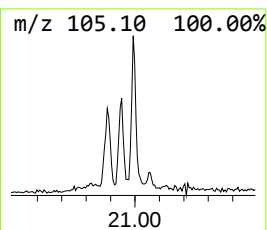
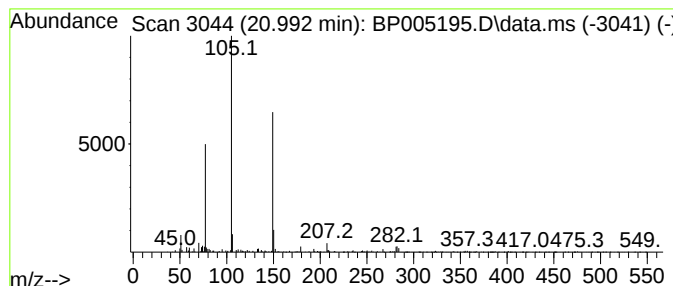
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 2-(2-(2-Methoxyethoxy)ethox... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.992	2.39 ng	39374	Chrysene-d12	21.210

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-(2-(2-Methoxyethoxy)ethoxy)eth...	268	C14H20O5	1000366-99-1	72		
2	Benzoic acid, 2-(4-nitrophenoxy)...	287	C15H13NO5	1000368-92-7	56		
3	Diethylene glycol dibenzoate	314	C18H18O5	000120-55-8	53		
4	Propiohydrazide, 3-phenyl-N2-(2-...	372	C23H20N2O3	1000295-37-0	43		
5	1,3-Dioxolane, 2-(methoxymethyl)...	194	C11H14O3	1000156-71-2	43		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040621\
Data File : BP005195.D
Acq On : 06 Apr 2021 14:20
Operator : CG/JU
Sample : M1894-03
Misc :
ALS Vial : 6 Sample Multiplier: 1

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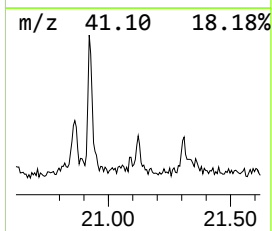
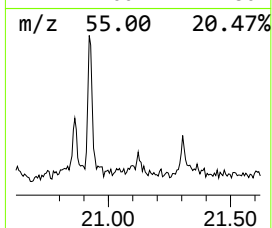
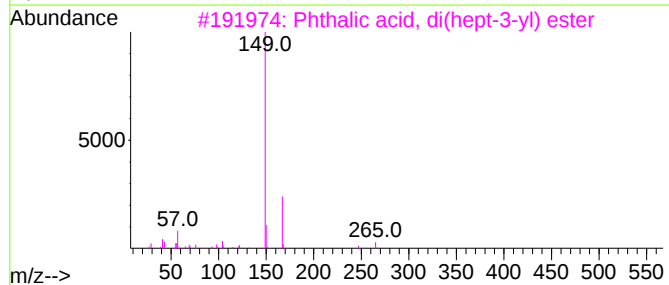
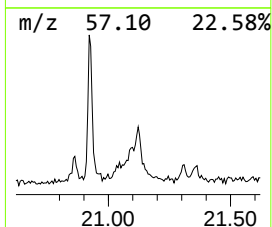
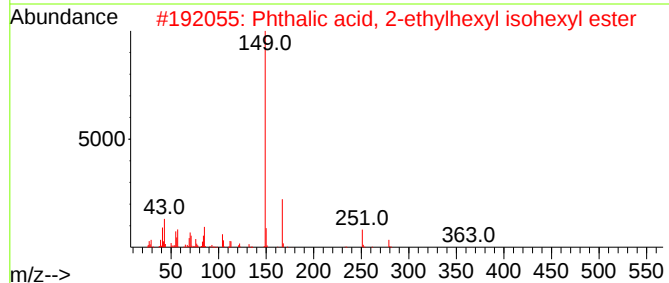
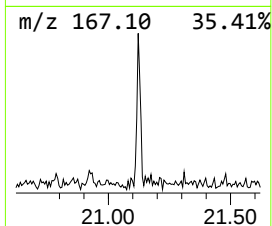
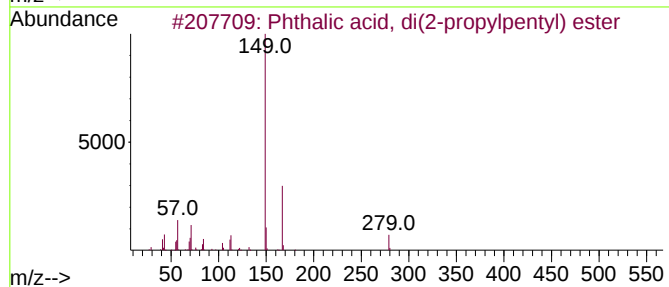
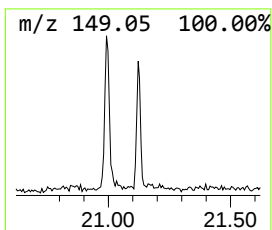
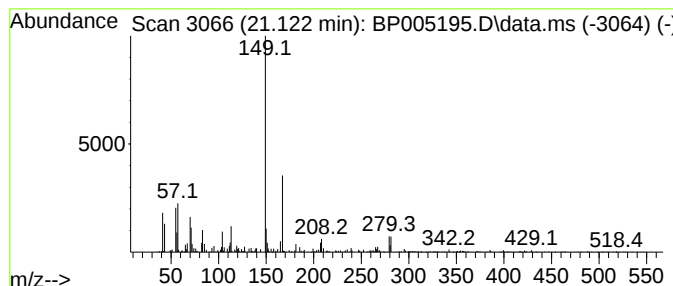
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Phthalic acid, di(2-propylp... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.122	2.91 ng	47967	Chrysene-d12	21.210

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phthalic acid, di(2-propylpentyl)...	390	C24H38O4	1000377-93-5	72
2			Phthalic acid, 2-ethylhexyl isoh...	362	C22H34O4	1000308-98-5	64
3			Phthalic acid, di(hept-3-yl) ester	362	C22H34O4	1000357-00-0	59
4			Phthalic acid, di(hept-4-yl) ester	362	C22H34O4	1000356-80-0	53
5			Phthalic acid, 5-methylhex-2-yl ...	446	C28H46O4	1000371-08-1	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040621\
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Sample : M1894-03
Misc :
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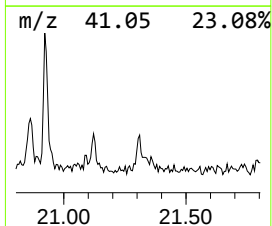
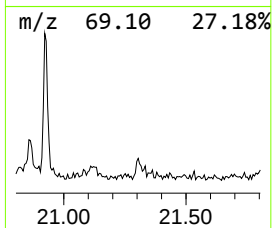
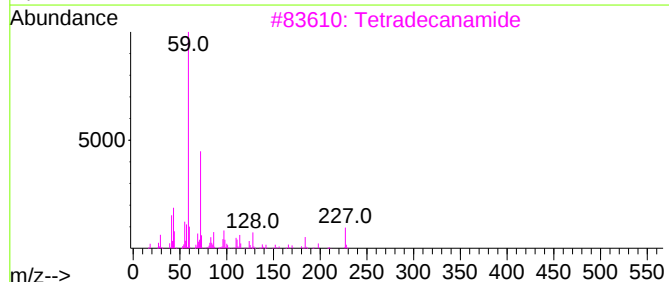
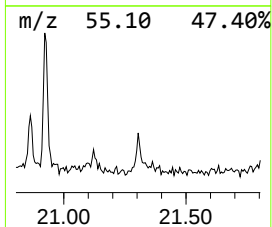
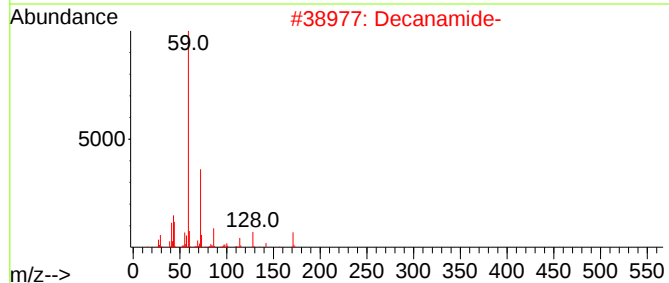
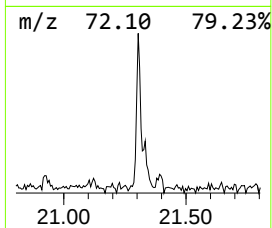
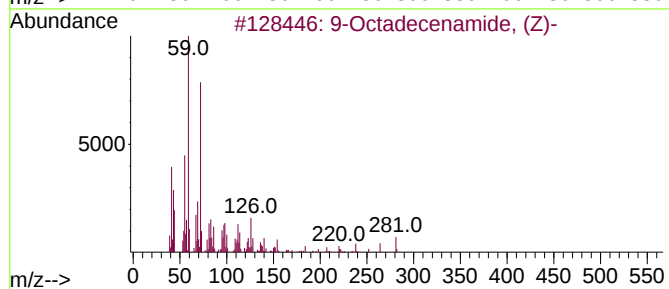
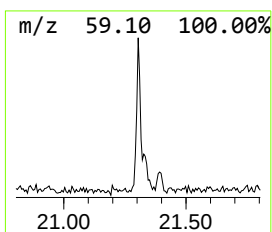
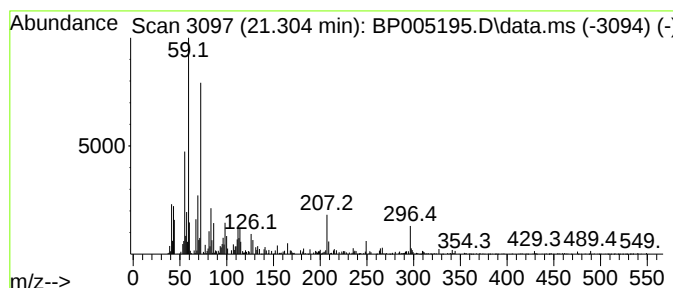
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP032321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Decanamide- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.304	2.90 ng	47828	Chrysene-d12	21.210
Hit#	of 5	Tentative ID	MW MolForm	CAS# Qual
1	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0 72
2	Decanamide-	171	C10H21NO	002319-29-1 72
3	Tetradecanamide	227	C14H29NO	000638-58-4 50
4	Pentadecanamide, 15-bromo-	319	C15H30BrNO	1000163-86-1 50
5	Dodecanamide	199	C12H25NO	001120-16-7 43



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040621\
Data File : BP005195.D
Acq On : 06 Apr 2021 14:20
Operator : CG/JU
Sample : M1894-03
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
0154-FIELD-BLANK

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP032321.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
unknown15.74	15.746	3.2	ng	45251	4	17.116	282218	20.0
n-Hexadecanoic ...	18.016	29.1	ng	411233	4	17.116	282218	20.0
1-Dodecanol	19.139	27.9	ng	393744	4	17.116	282218	20.0
Octadecanoic acid	19.257	14.1	ng	232932	5	21.210	329799	20.0
Dodecanamide	19.404	2.5	ng	40679	5	21.210	329799	20.0
9-Octadecenamid...	20.363	37.6	ng	620694	5	21.210	329799	20.0
Octanamide	20.463	4.3	ng	70568	5	21.210	329799	20.0
13-Octadecenal,...	20.863	5.8	ng	95988	5	21.210	329799	20.0
Trichloroacetic...	20.922	12.7	ng	208559	5	21.210	329799	20.0
2-(2-(2-Methoxy...	20.992	2.4	ng	39374	5	21.210	329799	20.0
Phthalic acid, ...	21.122	2.9	ng	47967	5	21.210	329799	20.0
Decanamide-	21.304	2.9	ng	47828	5	21.210	329799	20.0