

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP081820\
 Data File : BP003014.D
 Acq On : 18 Aug 2020 12:37
 Operator : CG/JU
 Sample : L3659-07
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SP1-3

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0
 Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP073020.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.952	7	11	29	rVB	881987	1211226	11.12%	1.839%
2	5.293	402	409	424	rBV	3044748	4387300	40.29%	6.662%
3	6.863	667	676	690	rBV	2801491	4493617	41.26%	6.823%
4	7.687	808	816	825	rBV	926870	1479790	13.59%	2.247%
5	8.828	1002	1010	1024	rBV	1662463	2710326	24.89%	4.115%
6	10.463	1279	1288	1302	rBV	1110629	1843750	16.93%	2.799%
7	12.945	1703	1710	1730	rBV	5578178	8279383	76.03%	12.571%
8	13.781	1840	1852	1861	rBV	233064	358124	3.29%	0.544%
9	14.322	1937	1944	1953	rBV	1738071	2511109	23.06%	3.813%
10	15.816	2191	2198	2210	rBV	3378772	4862163	44.65%	7.383%
11	17.075	2406	2412	2423	rBV	2094424	2897653	26.61%	4.400%
12	17.998	2563	2569	2577	rBV	629650	994433	9.13%	1.510%
13	19.233	2775	2779	2797	rBV	199120	385484	3.54%	0.585%
14	19.563	2831	2835	2843	rVB	205945	252765	2.32%	0.384%
15	19.651	2844	2850	2861	rBV	9061622	10890094	100.00%	16.535%
16	20.486	2989	2992	2996	rVB	142919	156713	1.44%	0.238%
17	20.733	3030	3034	3041	rVB	487235	523801	4.81%	0.795%
18	20.857	3050	3055	3059	rBV	202301	341434	3.14%	0.518%
19	20.904	3059	3063	3069	rVV2	1308719	1931665	17.74%	2.933%
20	20.957	3069	3072	3081	rVB	660894	850215	7.81%	1.291%
21	21.168	3102	3108	3112	rBV	2832757	3644737	33.47%	5.534%
22	21.345	3134	3138	3144	rBV	134846	168449	1.55%	0.256%
23	21.786	3208	3213	3224	rBV3	268614	609777	5.60%	0.926%
24	22.251	3288	3292	3296	rVB	251817	327018	3.00%	0.497%
25	22.380	3309	3314	3326	rVB	2107963	3046554	27.98%	4.626%
26	22.768	3375	3380	3384	rBV	499393	730457	6.71%	1.109%
27	23.351	3474	3479	3483	rBV	305336	498777	4.58%	0.757%
28	23.415	3483	3490	3499	rVB2	2195499	4139759	38.01%	6.286%
29	24.015	3587	3592	3600	rVB	421596	823116	7.56%	1.250%
30	24.098	3602	3606	3619	rVB	219460	510848	4.69%	0.776%

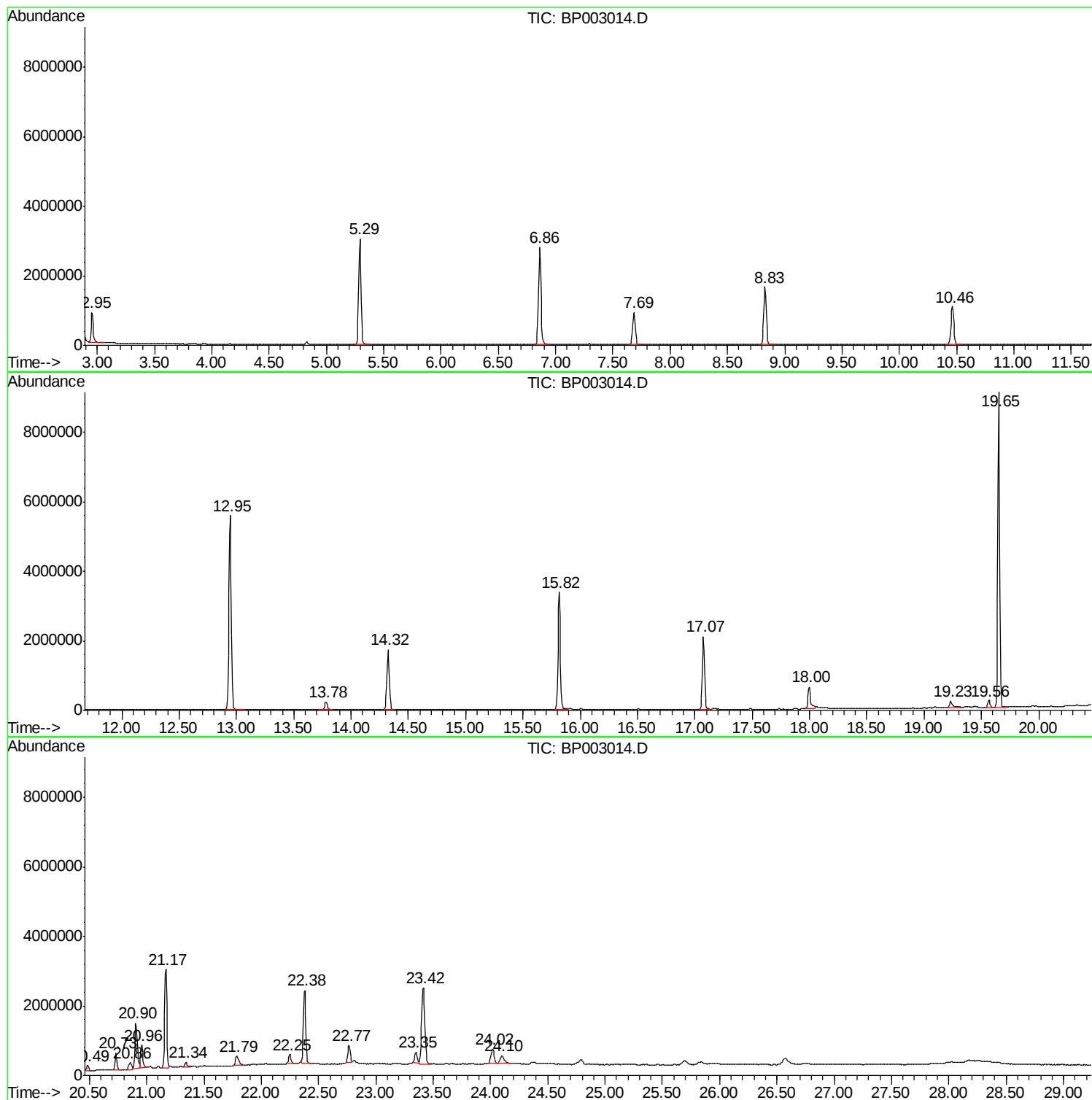
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP073020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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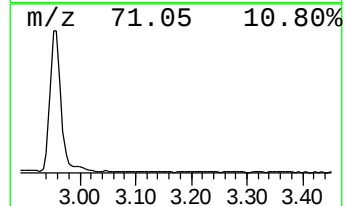
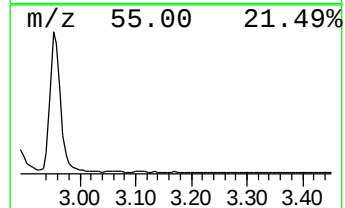
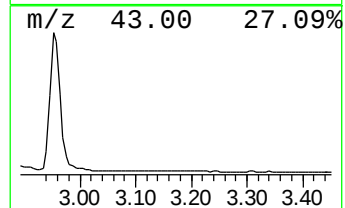
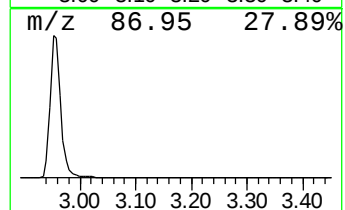
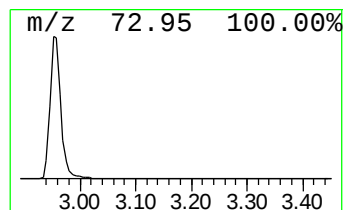
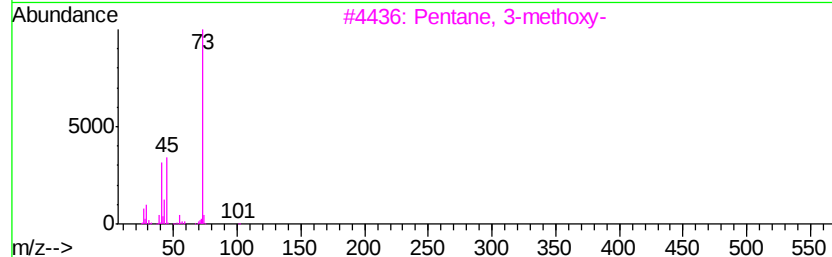
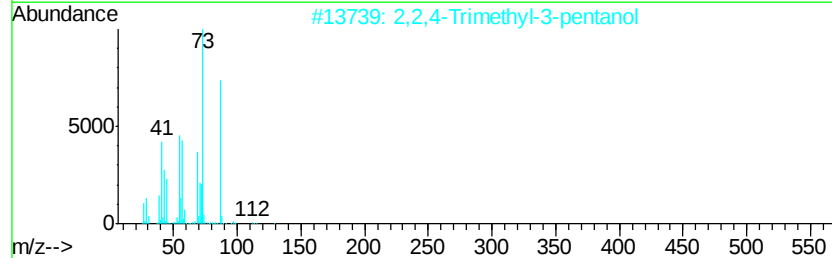
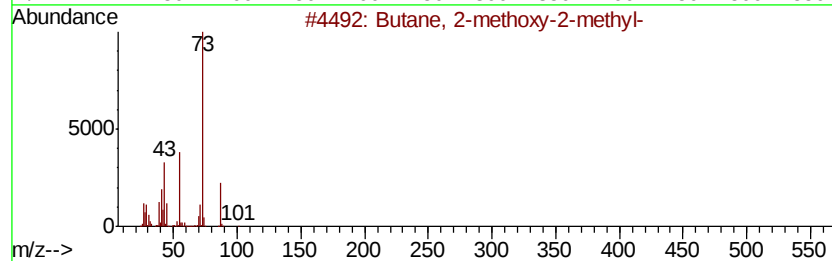
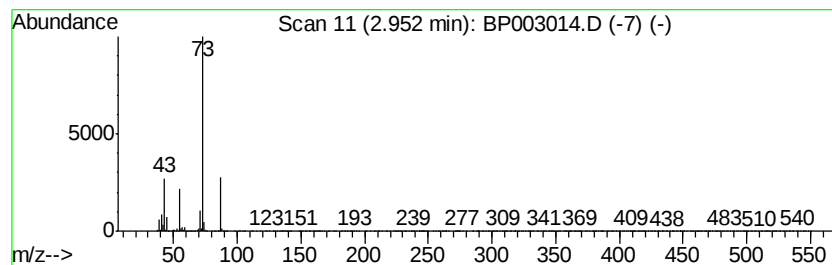
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.95	16.37 ng	1211230	1,4-Dichlorobenzene-d4	7.69

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	43
3		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
4		N-Ethylformamide	73	C3H7NO	000627-45-2	9
5		Silane, tetramethyl-	88	C4H12Si	000075-76-3	9



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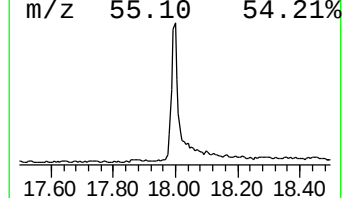
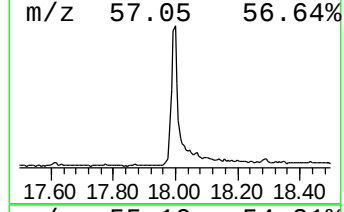
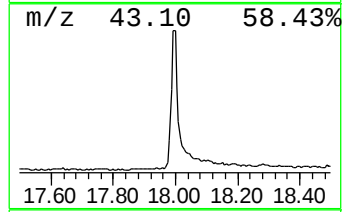
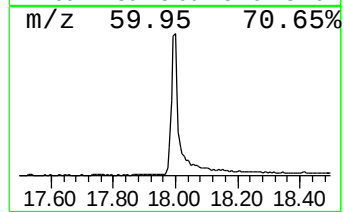
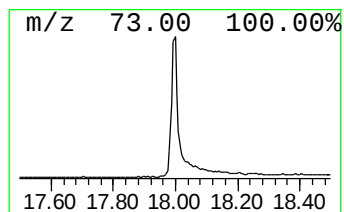
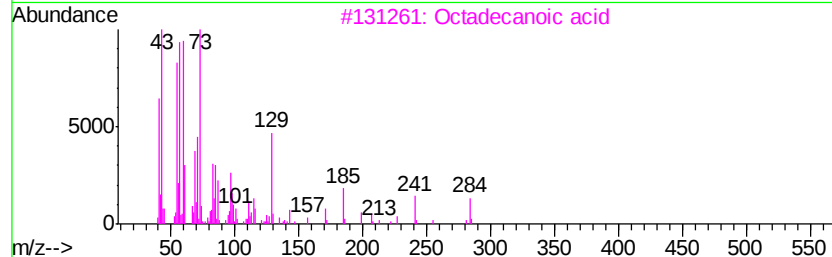
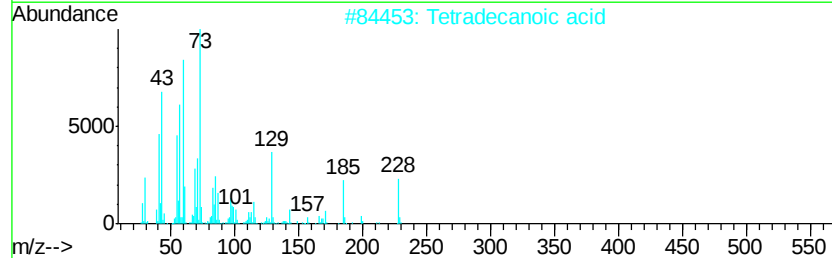
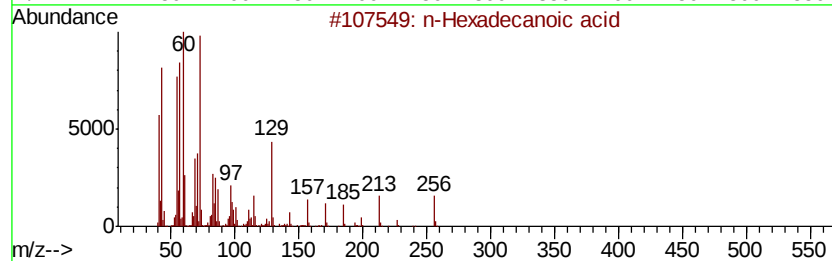
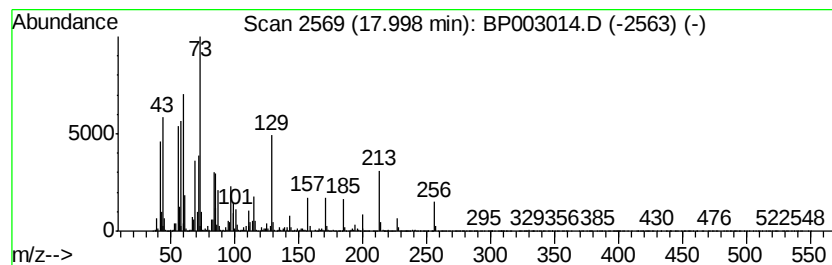
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 Peak Number 2 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.00	6.86 ng	994433	Phenanthrene-d10	17.07

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		Tetradecanoic acid	228	C14H28O2	000544-63-8	90
3		Octadecanoic acid	284	C18H36O2	000057-11-4	87
4		Pentadecanoic acid	242	C15H30O2	001002-84-2	80
5		Tridecanoic acid	214	C13H26O2	000638-53-9	74



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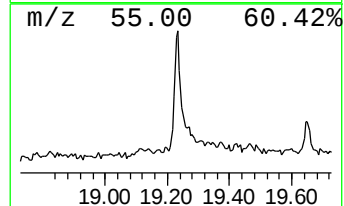
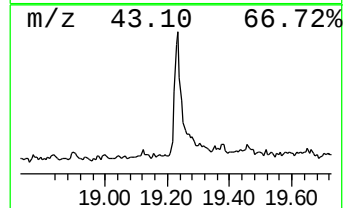
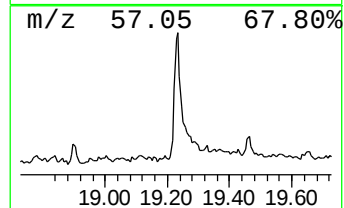
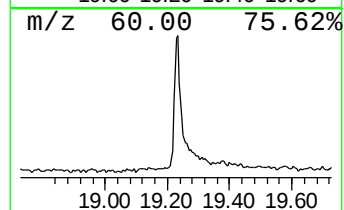
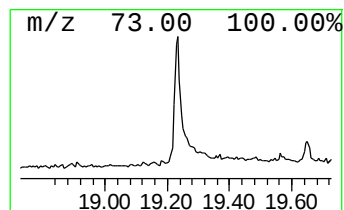
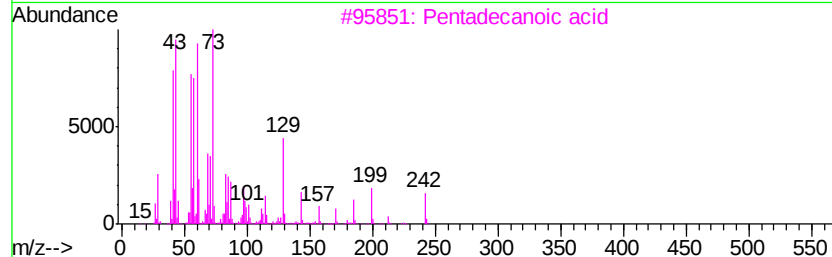
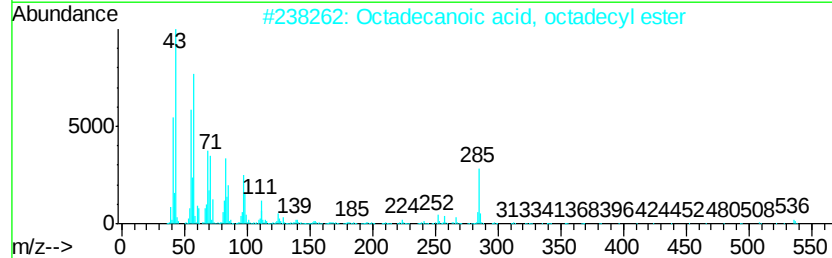
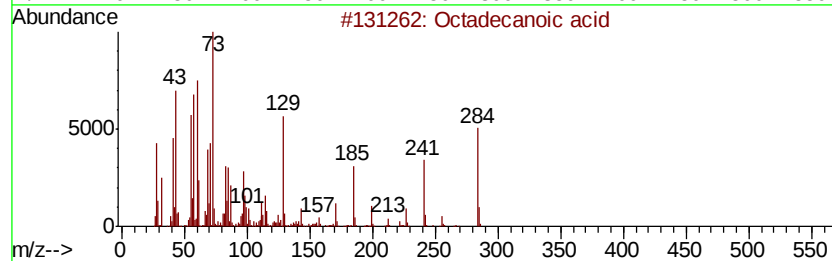
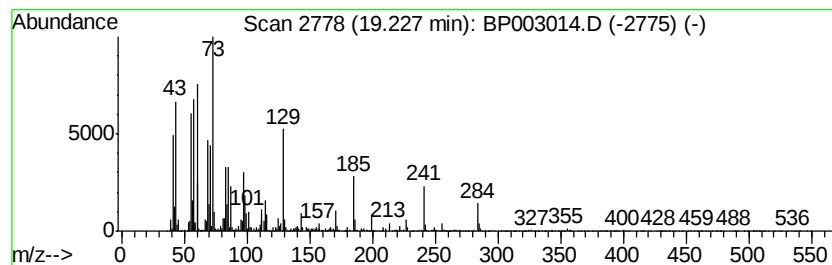
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Peak Number 3 Octadecanoic acid Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.23	2.12 ng	385484	Chrysene-d12	21.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2		Octadecanoic acid, octadecyl ester	537	C36H72O2	002778-96-3	81
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	53
4		Ethanol, 2-(octadecyloxy)-	314	C20H42O2	002136-72-3	50
5		Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	35



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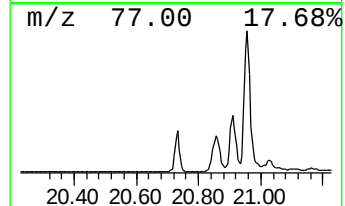
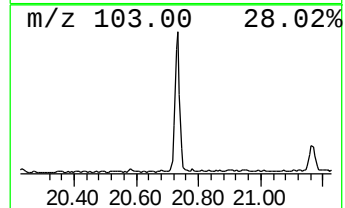
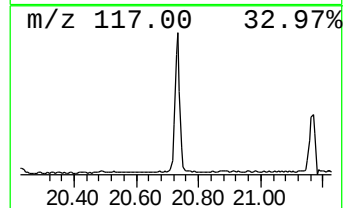
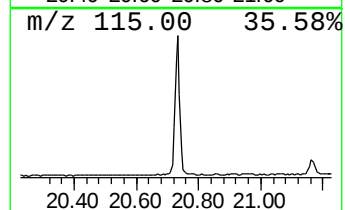
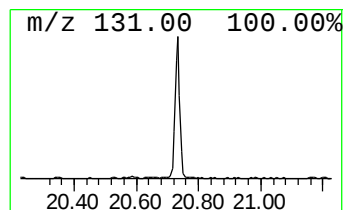
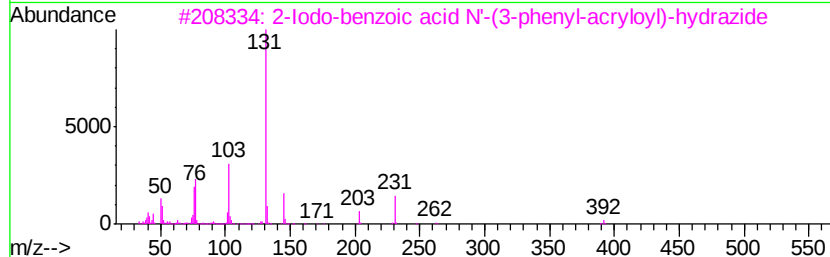
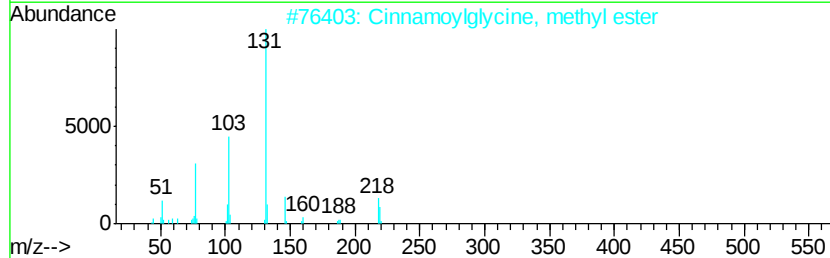
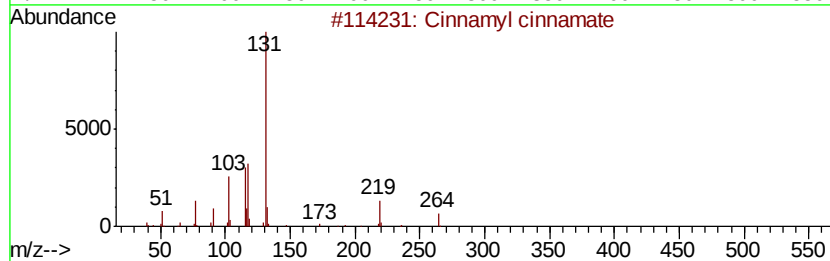
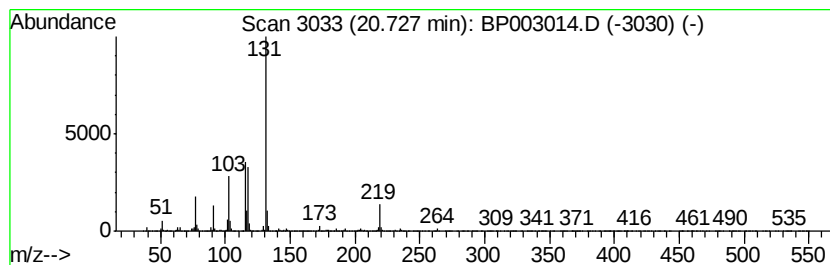
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Peak Number 4 Cinnamyl cinnamate Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.73	2.87 ng	523801	Chrysene-d12	21.17

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cinnamyl cinnamate	264	C18H16O2	000122-69-0	91	
2	Cinnamoylglycine, methyl ester	219	C12H13NO3	040778-04-9	50	
3	2-Iodo-benzoic acid N'-(3-phenyl...	392	C16H13IN2O2	315672-62-9	47	
4	2-Propenoic acid, 3-phenyl-, phe...	224	C15H12O2	025695-77-6	47	
5	2,2-Dimethyl-1-(2-vinylphenyl)pr...	188	C13H16O	1000210-99-9	47	



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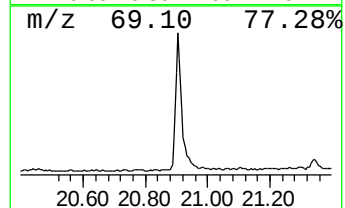
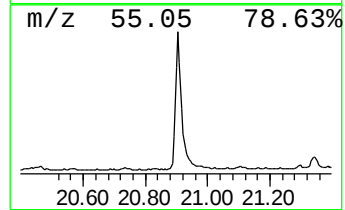
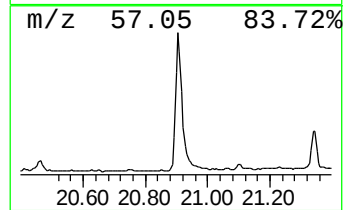
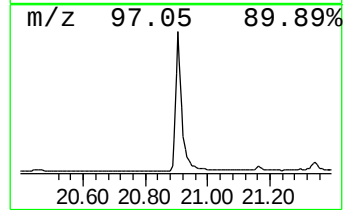
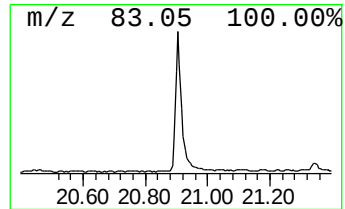
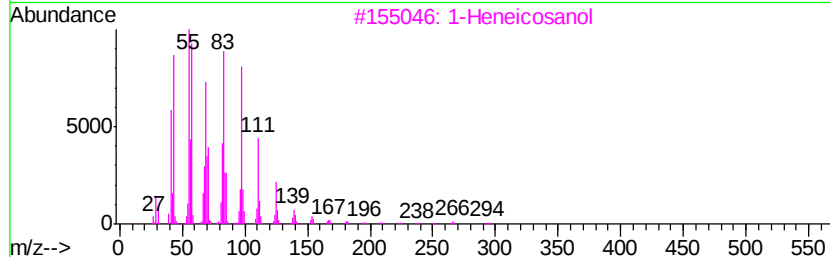
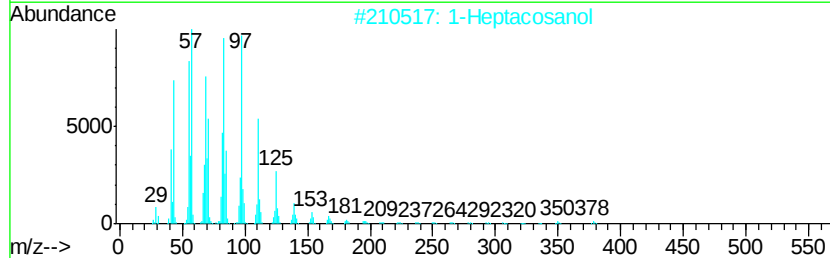
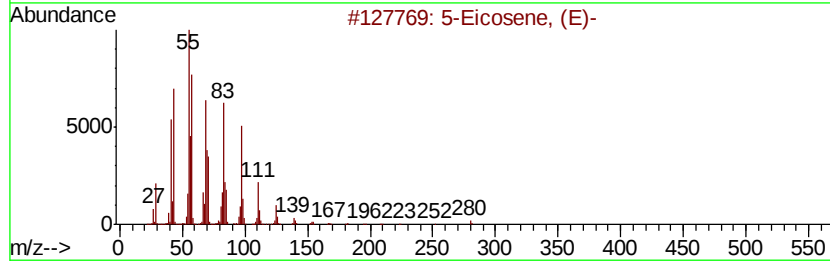
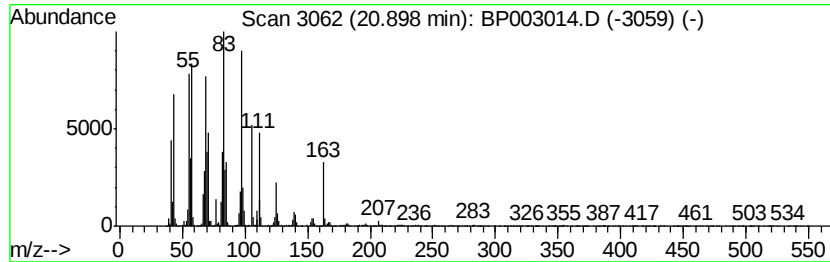
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 5 5-Eicosene, (E)- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.90	10.60 ng	1931670	Chrysene-d12	21.17

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Eicosene, (E)-		280	C20H40	074685-30-6	99
2	1-Heptacosanol		396	C27H56O	002004-39-9	93
3	1-Heneicosanol		312	C21H44O	015594-90-8	93
4	Carbonic acid, octadecyl 2,2,2-t...		444	C21H39Cl3O3	1000314-56-3	93
5	n-Nonadecanol-1		284	C19H40O	001454-84-8	93



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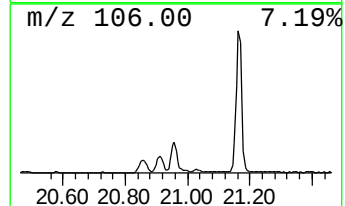
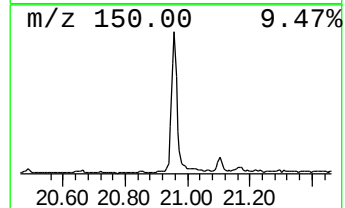
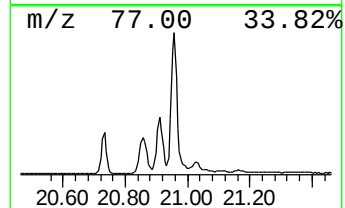
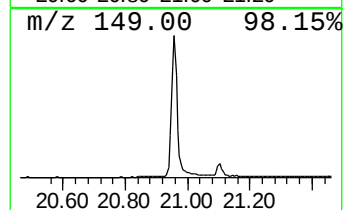
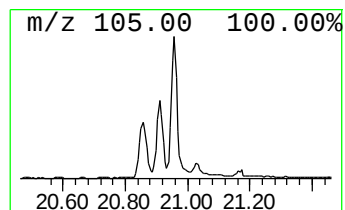
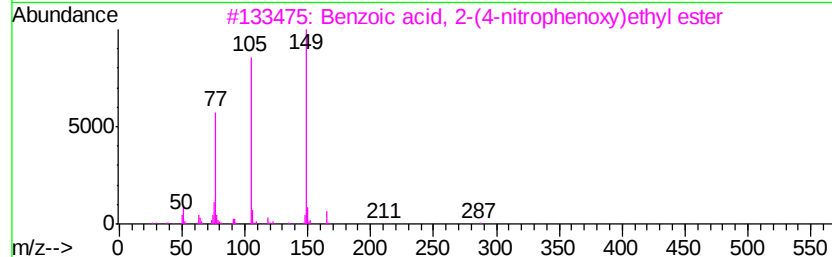
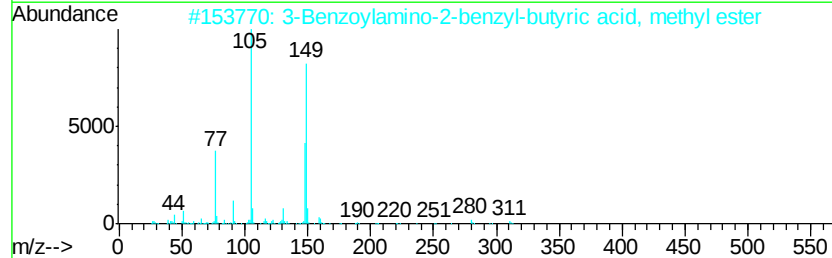
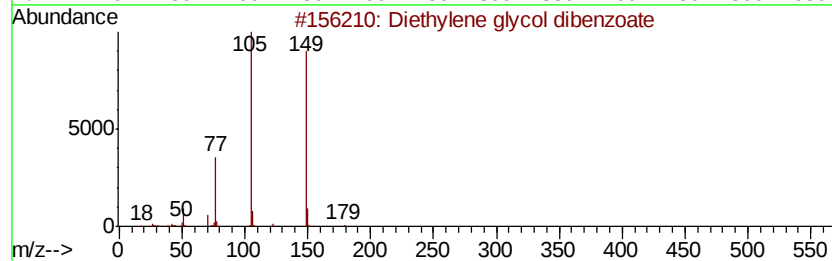
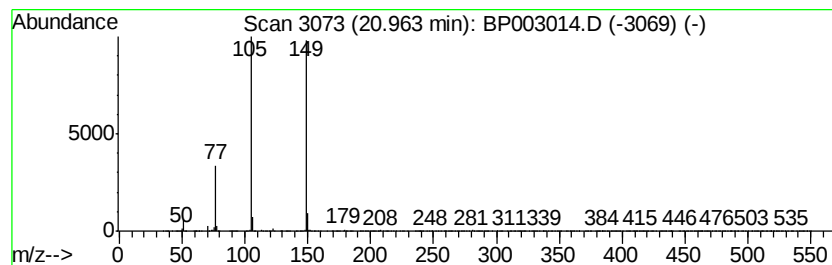
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Diethylene glycol dibenzoate Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.96	4.67 ng	850215	Chrysene-d12	21.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Diethylene glycol dibenzoate	314	C18H18O5	000120-55-8	83
2		3-Benzoylamino-2-benzyl-butyrlic ...	311	C19H21NO3	1000193-12-7	78
3		Benzoic acid, 2-(4-nitrophenoxy)...	287	C15H13NO5	1000368-92-7	78
4		2,5,8,11-Tetraoxatridecan-13-yl ...	312	C16H24O6	1000366-96-7	78
5		2-(2-(2-Butoxyethoxy)ethoxy)ethy...	310	C17H26O5	1000366-97-2	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP081820\
Data File : BP003014.D
Acq On : 18 Aug 2020 12:37
Operator : CG/JU
Sample : L3659-07
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SP1-3

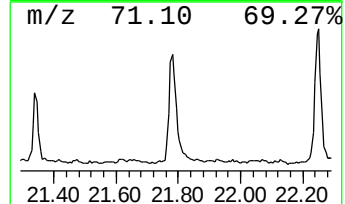
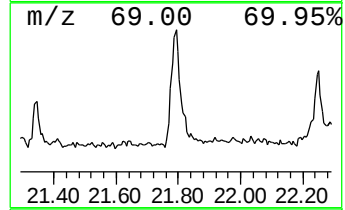
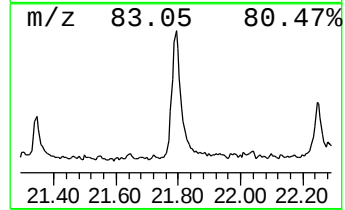
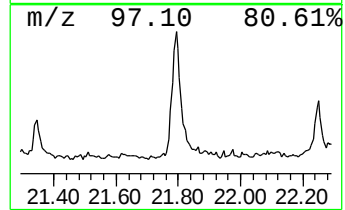
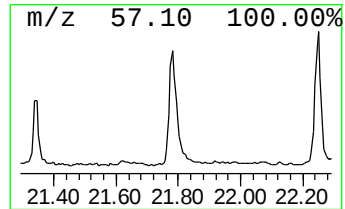
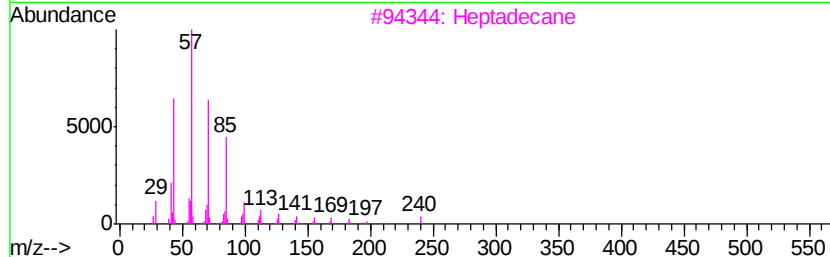
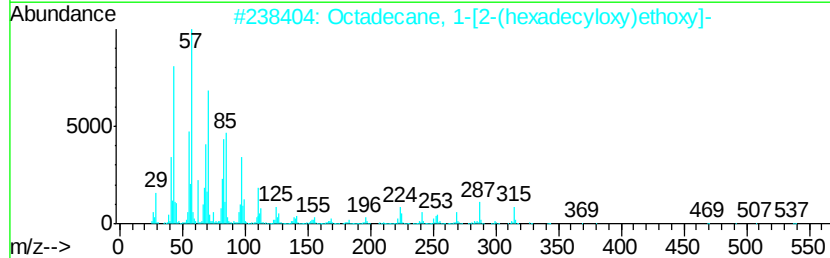
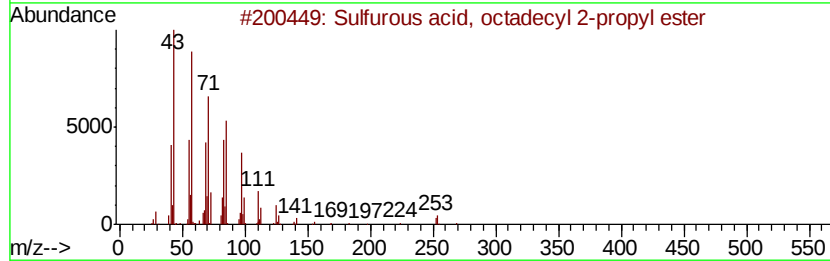
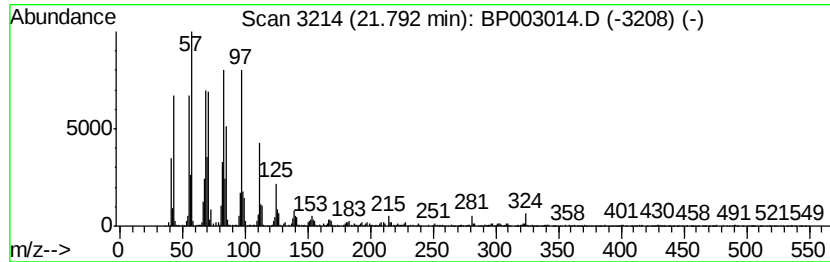
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Sulfurous acid, octadecyl 2... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.79	3.35 ng	609777	Chrysene-d12	21.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Sulfurous acid, octadecyl 2-prop...	376	C21H44O3S	1000309-12-7	91
2		Octadecane, 1-[2-(hexadecyloxy)e...	539	C36H74O2	017367-10-1	91
3		Heptadecane	240	C17H36	000629-78-7	87
4		Eicosane	282	C20H42	000112-95-8	87
5		Octacosyl acetate	452	C30H60O2	018206-97-8	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP081820\
Data File : BP003014.D
Acq On : 18 Aug 2020 12:37
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Misc :
ALS Vial : 4 Sample Multiplier: 1

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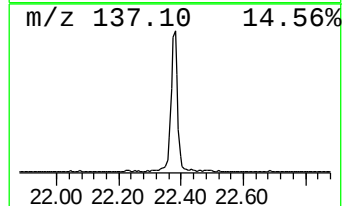
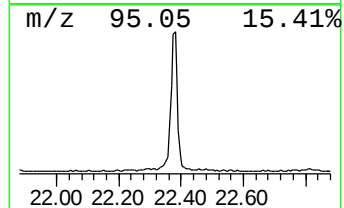
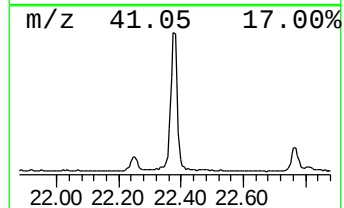
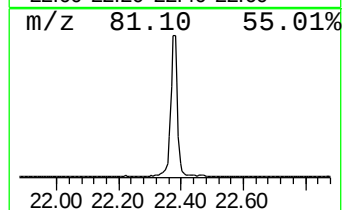
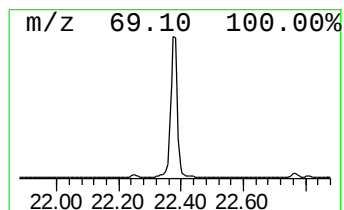
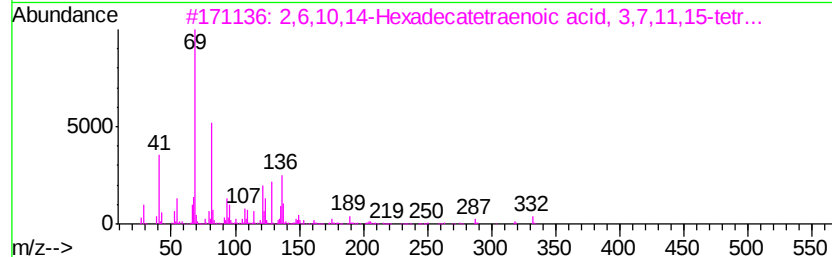
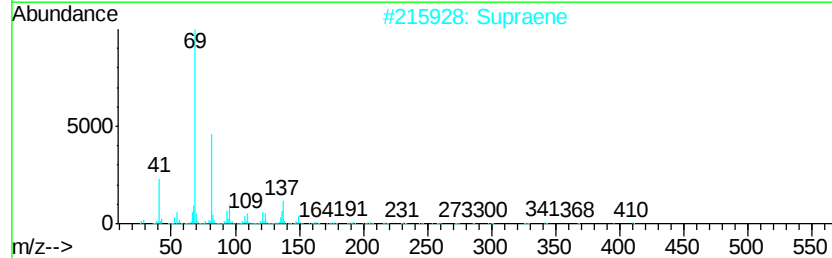
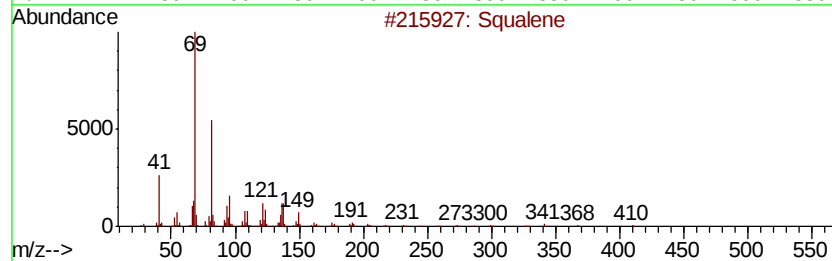
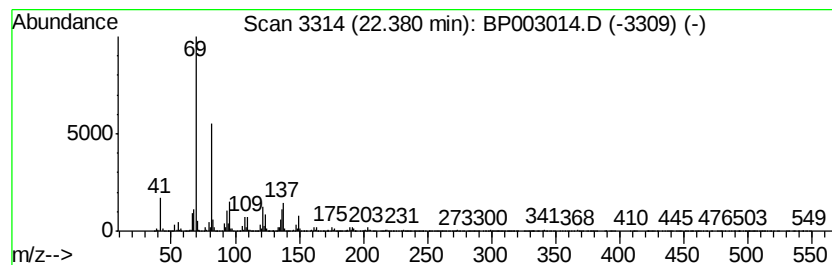
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Squalene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.38	14.72 ng	3046550	Perylene-d12	23.42

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Squalene	410	C30H50	000111-02-4	94
2		Supraene	410	C30H50	007683-64-9	91
3		2,6,10,14-Hexadecatetraenoic aci...	332	C22H36O2	024035-35-6	90
4		1,5,9-Undecatriene, 2,6,10-trime...	192	C14H24	062951-96-6	74
5		7,11-Dimethyldodeca-2,6,10-trien...	208	C14H24O	125529-10-4	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP081820\
Data File : BP003014.D
Acq On : 18 Aug 2020 12:37
Operator : CG/JU
Sample : L3659-07
Misc :
ALS Vial : 4 Sample Multiplier: 1

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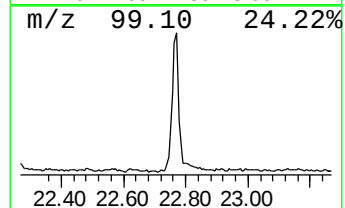
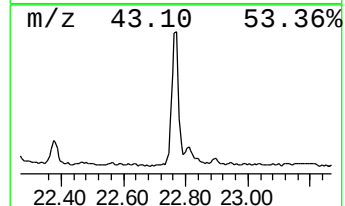
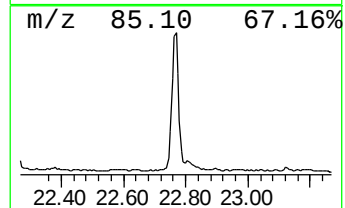
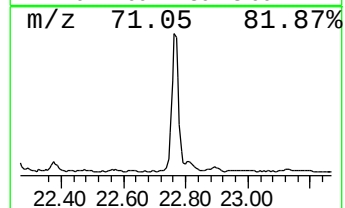
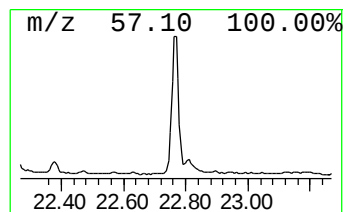
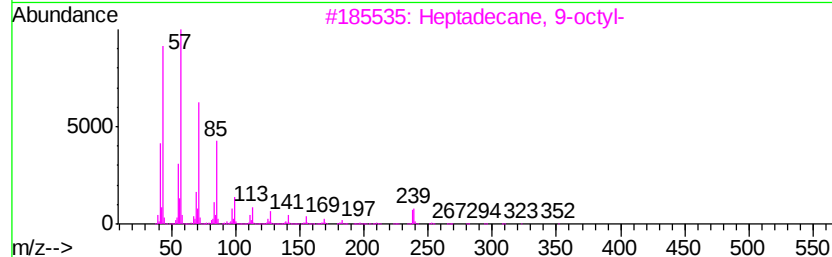
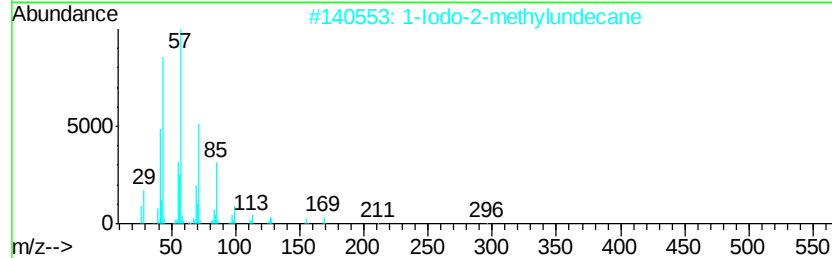
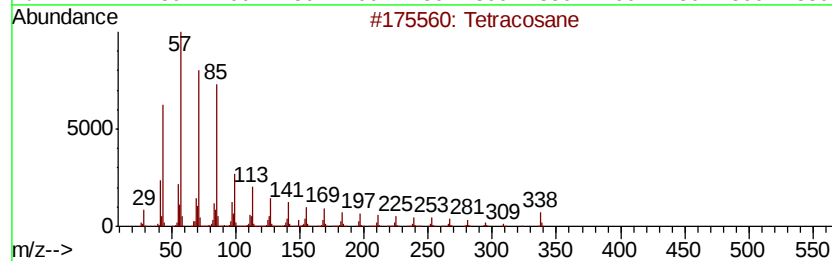
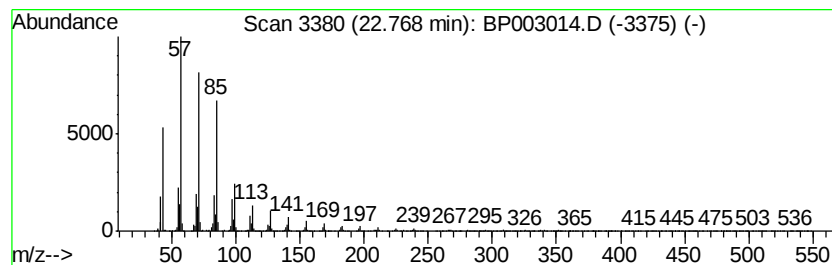
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Tetracosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.77	3.53 ng	730457	Perylene-d12	23.42

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetracosane	338	C24H50	000646-31-1	94
2		1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	93
3		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	93
4		Heneicosane	296	C21H44	000629-94-7	91
5		2-methyloctacosane	408	C29H60	1000376-72-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP081820\
Data File : BP003014.D
Acq On : 18 Aug 2020 12:37
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Sample : L3659-07
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SP1-3

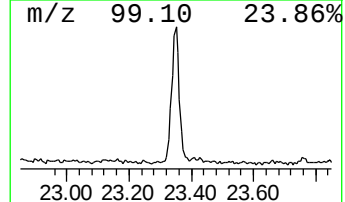
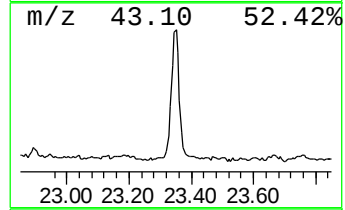
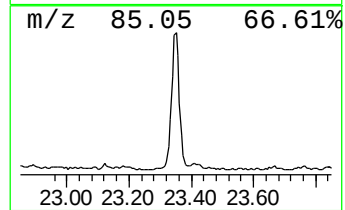
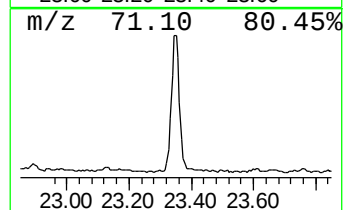
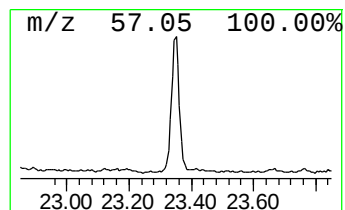
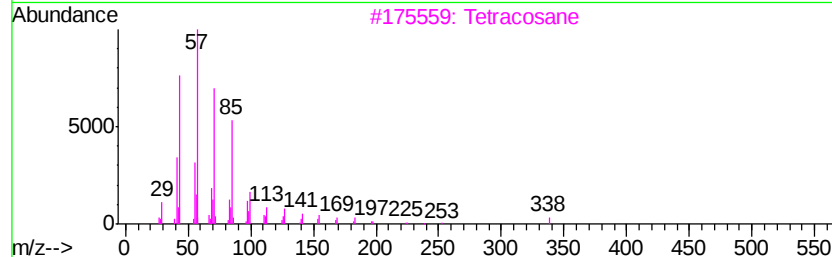
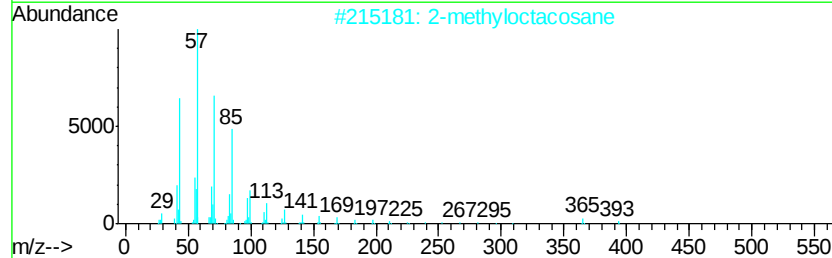
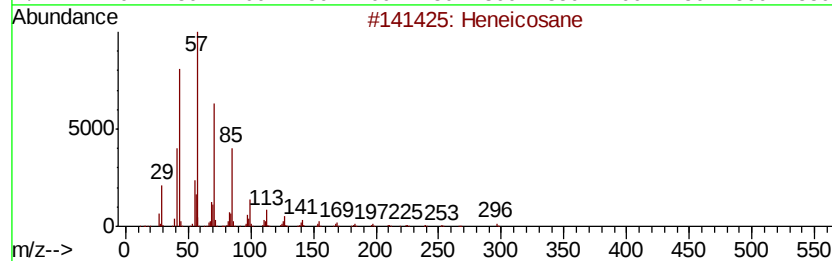
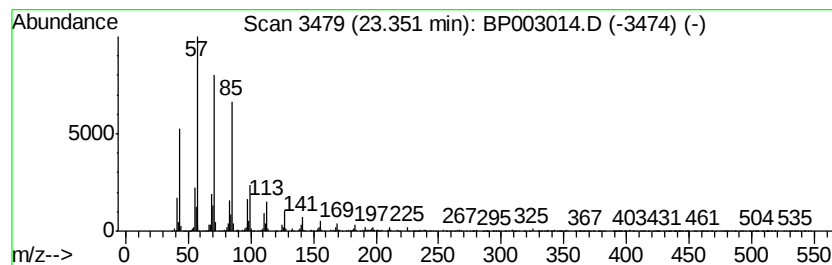
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Heneicosane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.35	2.41 ng	498777	Perylene-d12	23.42

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	97
2			2-methyloctacosane	408	C29H60	1000376-72-8	91
3			Tetracosane	338	C24H50	000646-31-1	91
4			Docosane, 11-butyl-	366	C26H54	013475-76-8	91
5			1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP081820\
Data File : BP003014.D
Acq On : 18 Aug 2020 12:37
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Misc :
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Instrument :
BNA_P
ClientSampleId :
SP1-3

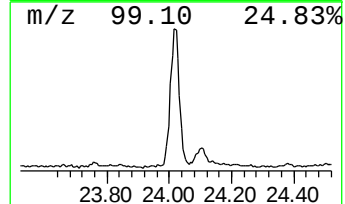
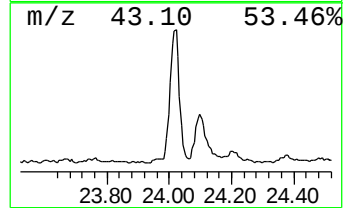
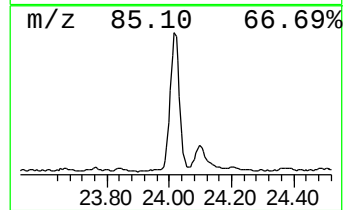
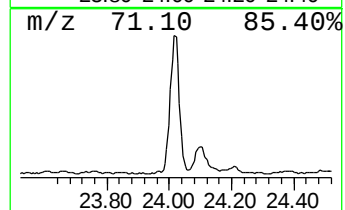
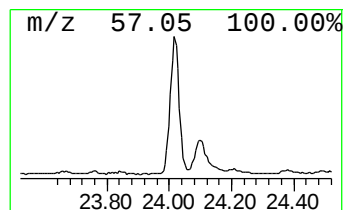
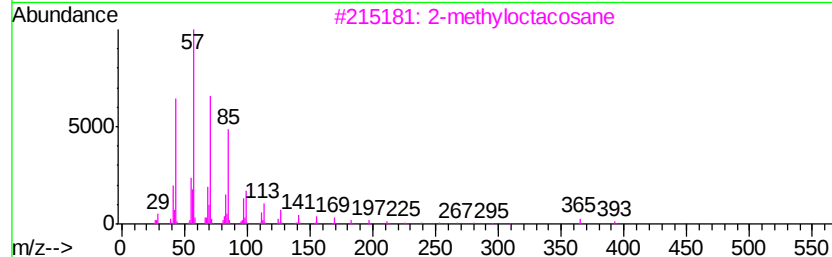
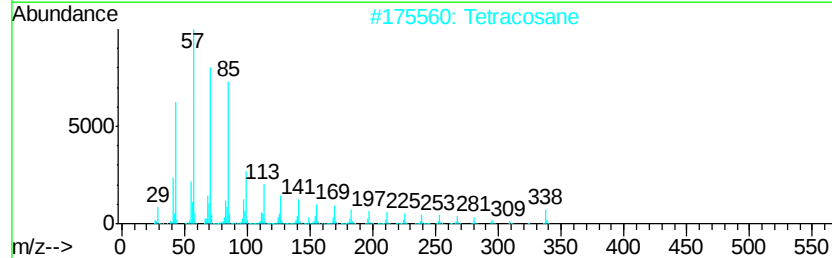
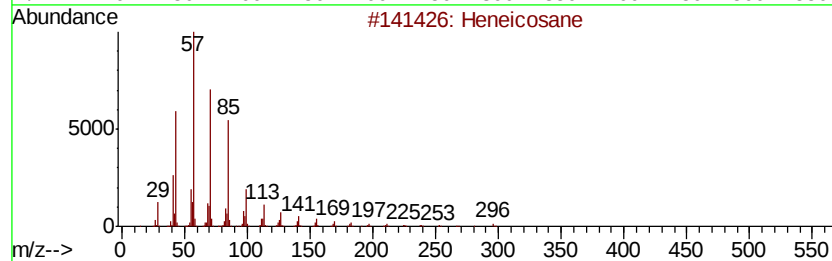
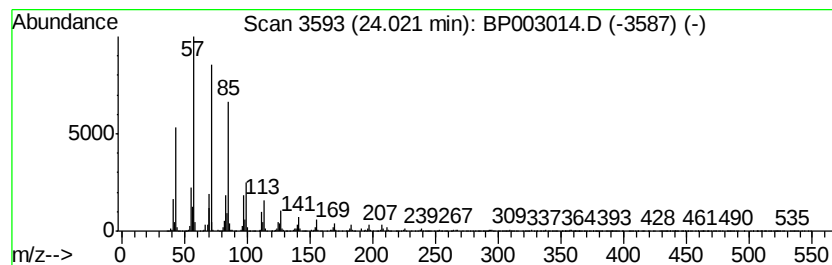
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 2-methyloctacosane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.02	3.98 ng	823116	Perylene-d12	23.42

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	97
2		Tetracosane	338	C24H50	000646-31-1	95
3		2-methyloctacosane	408	C29H60	1000376-72-8	91
4		Heneicosane, 11-decyl-	437	C31H64	055320-06-4	91
5		Octacosane	394	C28H58	000630-02-4	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP081820\
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Acq On : 18 Aug 2020 12:37
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Sample : L3659-07
Misc :
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ClientSampleId :
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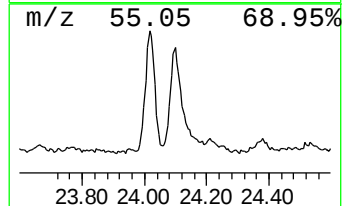
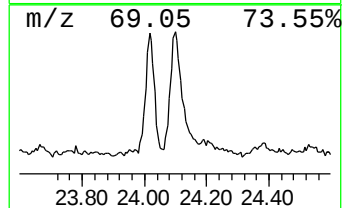
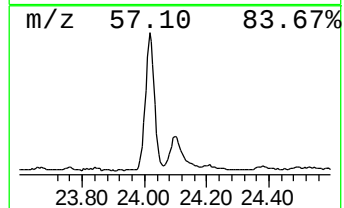
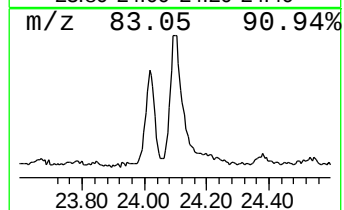
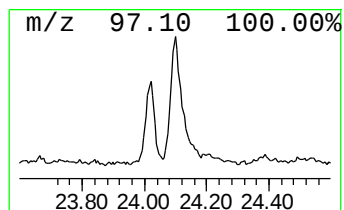
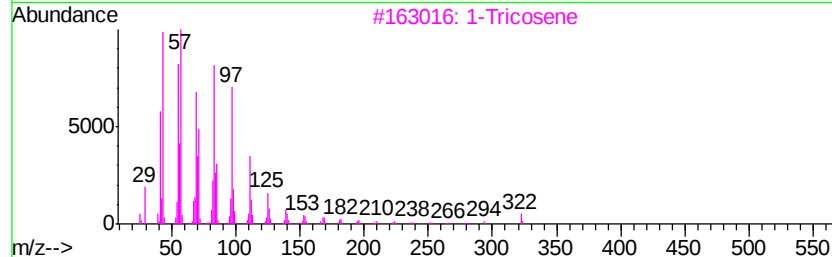
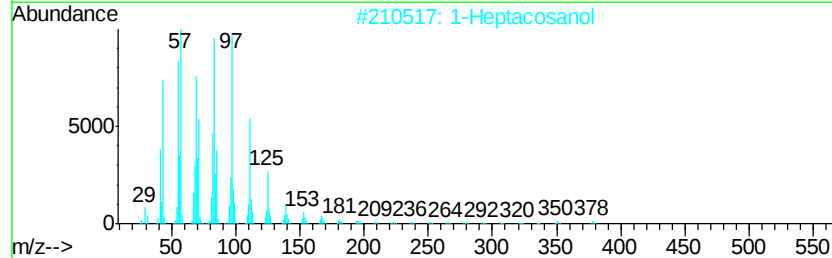
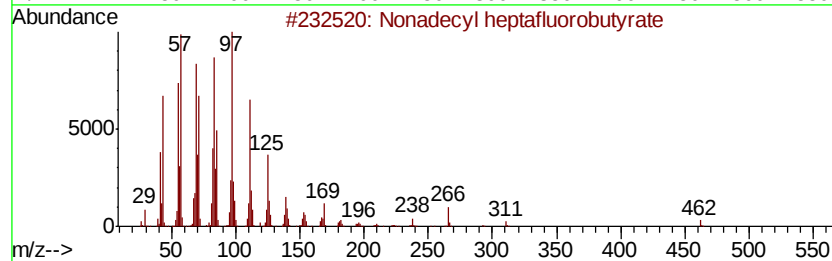
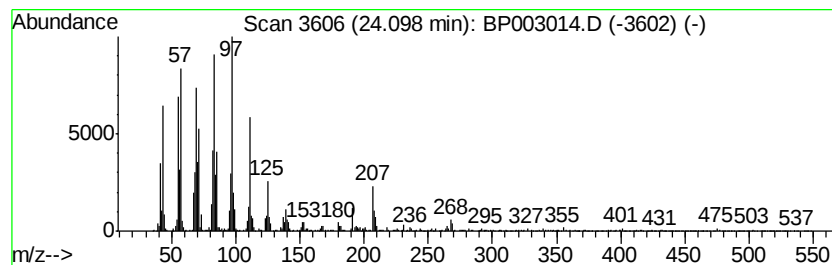
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Nonadecyl heptafluorobutyrate Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.10	2.47 ng	510848	Perylene-d12	23.42

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonadecyl heptafluorobutyrate	480	C23H39F7O2	1000351-83-9	93
2		1-Heptacosanol	396	C27H56O	002004-39-9	93
3		1-Tricosene	322	C23H46	018835-32-0	92
4		1-Nonadecene	266	C19H38	018435-45-5	91
5		9-Tricosene, (Z)-	322	C23H46	027519-02-4	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP081820\
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 Acq On : 18 Aug 2020 12:37
 Operator : CG/JU
 Sample : L3659-07
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SP1-3

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP073020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	2.95	16.4	ng	1211230	1	7.69	1479790	20.0
n-Hexadecanoic acid	18.00	6.9	ng	994433	4	17.07	2897650	20.0
Octadecanoic acid	19.23	2.1	ng	385484	5	21.17	3644740	20.0
Cinnamyl cinnamate	20.73	2.9	ng	523801	5	21.17	3644740	20.0
5-Eicosene, (E)-	20.90	10.6	ng	1931670	5	21.17	3644740	20.0
Diethylene glycol...	20.96	4.7	ng	850215	5	21.17	3644740	20.0
Sulfurous acid, o...	21.79	3.4	ng	609777	5	21.17	3644740	20.0
Squalene	22.38	14.7	ng	3046550	6	23.42	4139760	20.0
Tetracosane	22.77	3.5	ng	730457	6	23.42	4139760	20.0
Heneicosane	23.35	2.4	ng	498777	6	23.42	4139760	20.0
2-methyloctacosane	24.02	4.0	ng	823116	6	23.42	4139760	20.0
Nonadecyl heptafl...	24.10	2.5	ng	510848	6	23.42	4139760	20.0