

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP011520\
 Data File : BP001619.D
 Acq On : 15 Jan 2020 20:47
 Operator : JU
 Sample : L1126-07
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 FB-01142020

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-BP010820.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	2.875	10	15	24	rBV	62271	116250	4.54%	0.790%
2	2.952	24	28	44	rVB	687661	853831	33.37%	5.804%
3	5.257	415	420	431	rVB	247637	362733	14.18%	2.466%
4	6.822	680	686	702	rVB	169336	277852	10.86%	1.889%
5	7.175	738	746	761	rBV	429541	685813	26.80%	4.662%
6	7.634	817	824	830	rBV	147659	228505	8.93%	1.553%
7	7.951	871	878	888	rVB	560321	865623	33.83%	5.884%
8	8.781	1011	1019	1036	rBV	487559	821483	32.11%	5.584%
9	10.404	1288	1295	1306	rBV	157662	279290	10.92%	1.898%
10	12.898	1712	1719	1732	rBV	1079141	1571207	61.41%	10.680%
11	13.745	1856	1863	1875	rBV	54091	79253	3.10%	0.539%
12	14.280	1947	1954	1961	rBV	253204	378466	14.79%	2.573%
13	15.780	2203	2209	2220	rBV	623918	869669	33.99%	5.912%
14	17.039	2417	2423	2428	rBV	312660	450575	17.61%	3.063%
15	17.974	2575	2582	2588	rBV2	53981	82765	3.23%	0.563%
16	19.215	2789	2793	2799	rBV5	16105	26314	1.03%	0.179%
17	19.633	2858	2864	2879	rBV	2049794	2558615	100.00%	17.392%
18	20.321	2978	2981	2987	rVB3	18795	26146	1.02%	0.178%
19	20.451	2997	3003	3006	rBV2	31559	39639	1.55%	0.269%
20	20.904	3074	3080	3086	rBV	126541	185853	7.26%	1.263%
21	21.145	3116	3121	3128	rBV	375928	512190	20.02%	3.482%
22	21.333	3149	3153	3157	rBV	132383	150919	5.90%	1.026%
23	21.768	3223	3227	3235	rBV	203187	237157	9.27%	1.612%
24	22.239	3302	3307	3313	rVB	236654	324160	12.67%	2.204%
25	22.751	3389	3394	3400	rVB	248181	350166	13.69%	2.380%
26	22.921	3418	3423	3430	rVB	173874	261397	10.22%	1.777%
27	23.333	3487	3493	3496	rBV	202477	350174	13.69%	2.380%
28	23.374	3496	3500	3509	rVB2	236177	427941	16.73%	2.909%
29	23.992	3598	3605	3612	rBV	156548	320119	12.51%	2.176%
30	24.133	3622	3629	3638	rVB	246601	527387	20.61%	3.585%
31	24.750	3728	3734	3743	rVB2	77510	171308	6.70%	1.164%
32	25.721	3891	3899	3909	rVB2	113902	318337	12.44%	2.164%

LSC Area Percent Report

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Misc :
ALS Vial : 15 Sample Multiplier: 1

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ClientSampleId :
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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\8270-BP010820.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

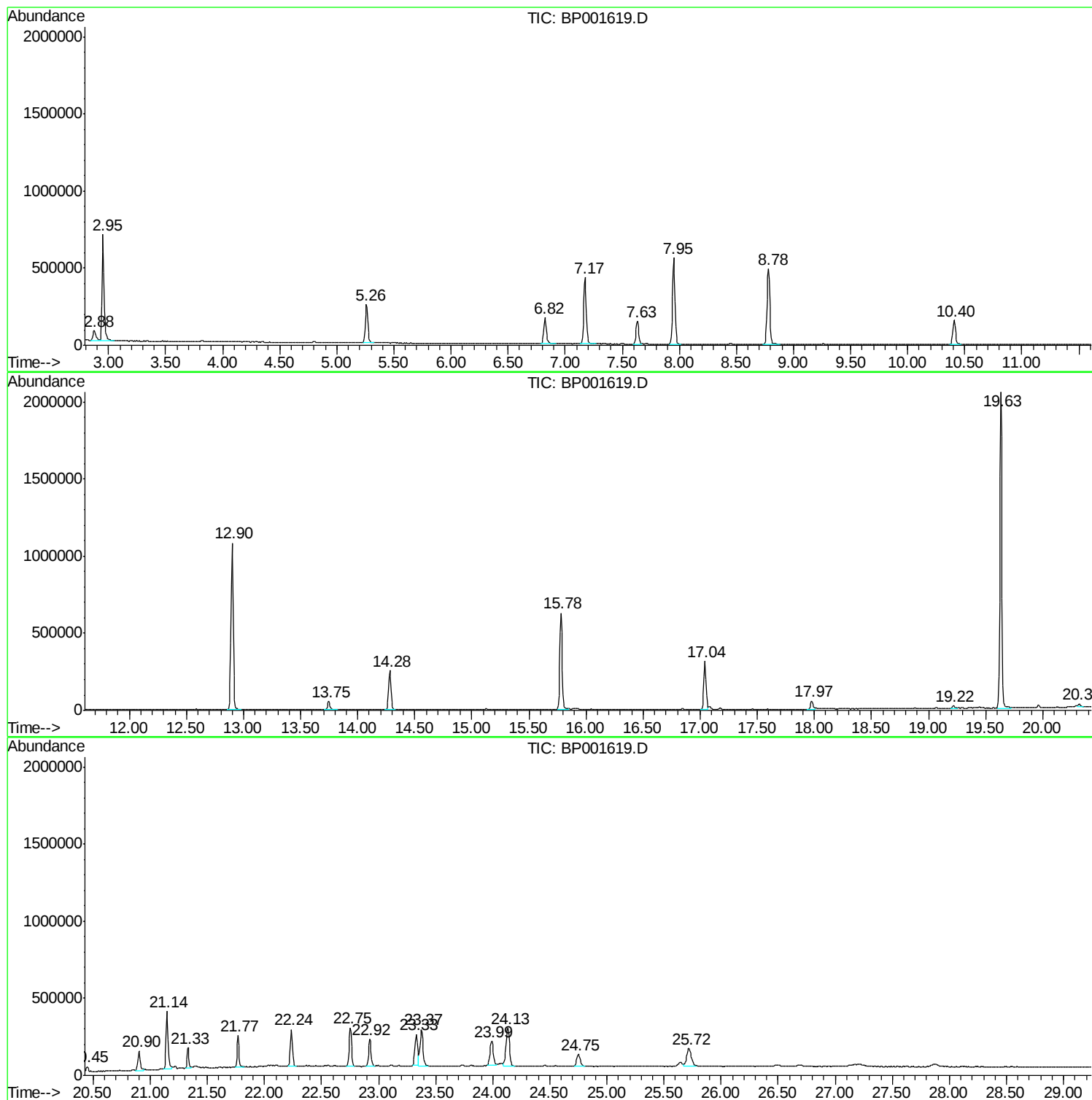
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP010820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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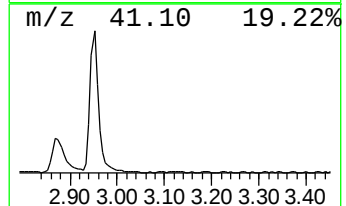
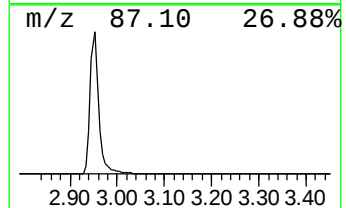
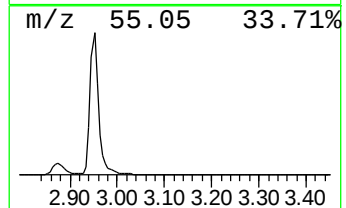
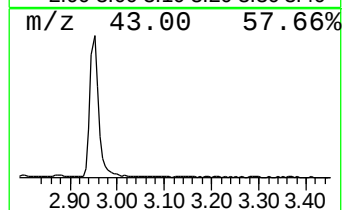
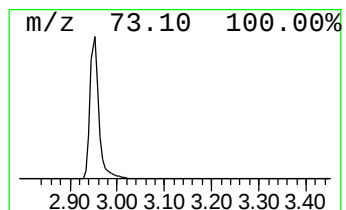
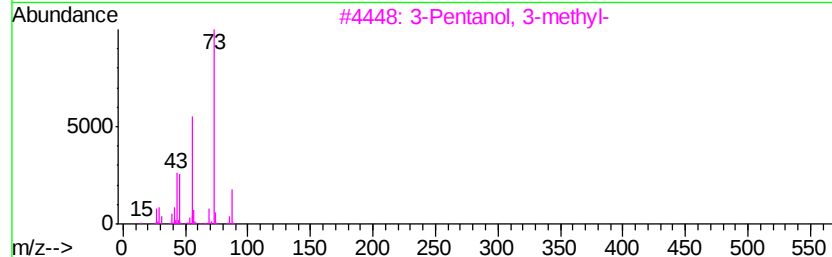
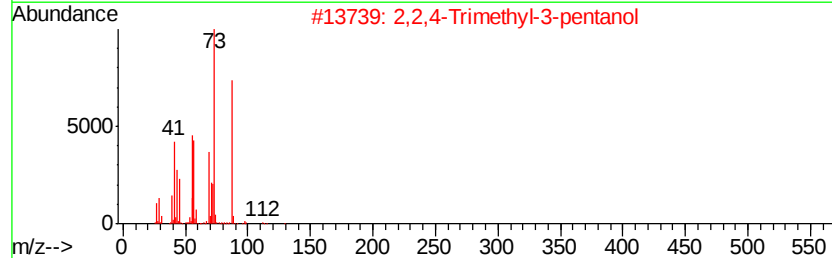
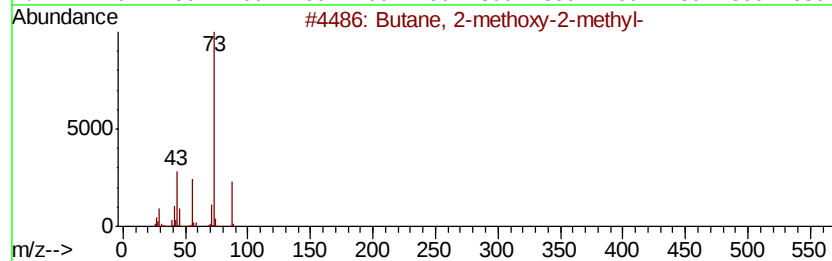
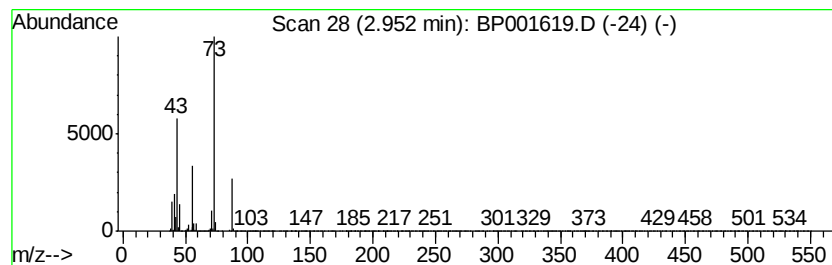
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.95	74.73 ng	853831	1,4-Dichlorobenzene-d4	7.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
3		3-Pentanol, 3-methyl-	102	C6H14O	000077-74-7	25
4		Silane, tetramethyl-	88	C4H12Si	000075-76-3	23
5		2-Methyl-5-hexen-3-ol	114	C7H14O	032815-70-6	23



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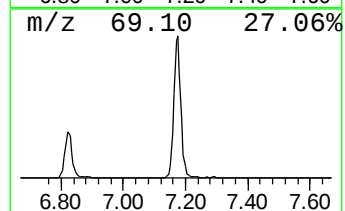
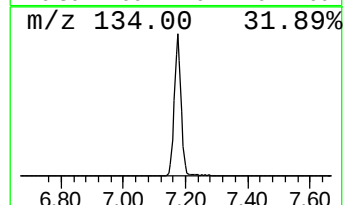
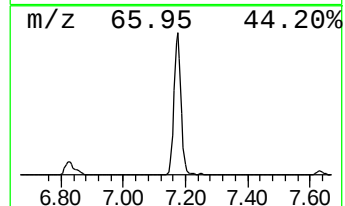
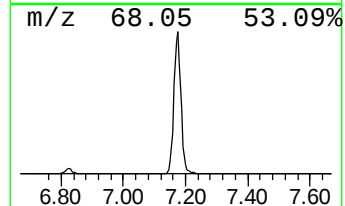
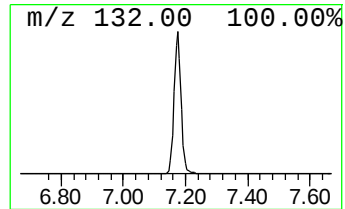
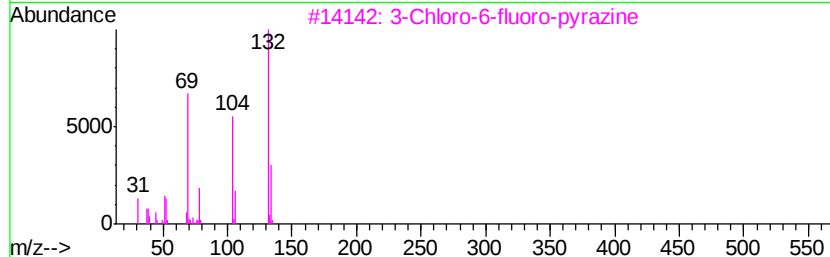
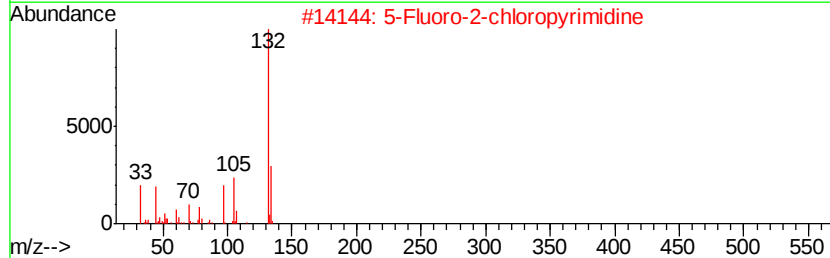
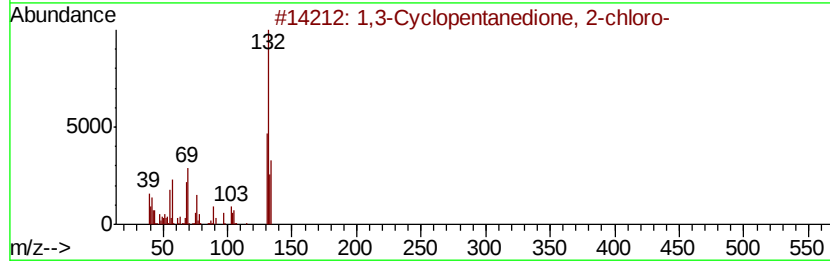
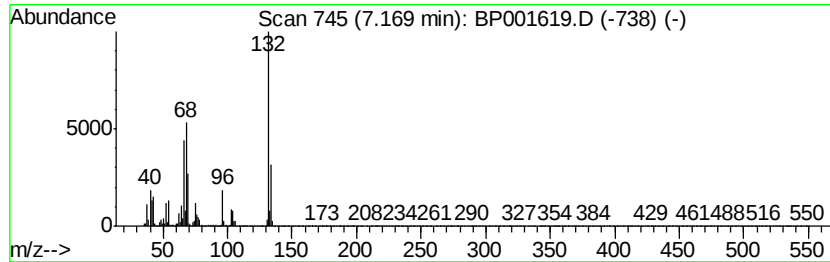
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown7.17 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.17	60.03 ng	685813	1,4-Dichlorobenzene-d4	7.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	22
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	14
3		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	14
4		1,3,4-Thiadiazole-2(3H)-thione, ...	132	C3H4N2S2	029490-19-5	14
5		Pyrazolo(2,3-a)pyridine, 7-methyl-	132	C8H8N2	034760-58-2	14



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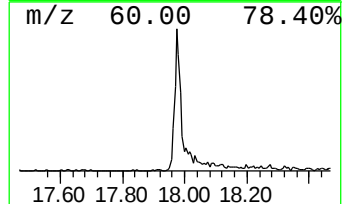
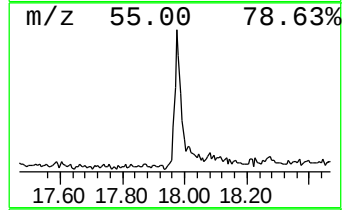
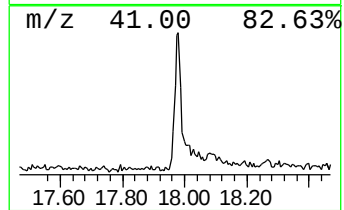
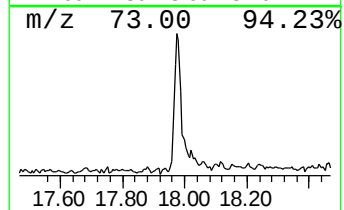
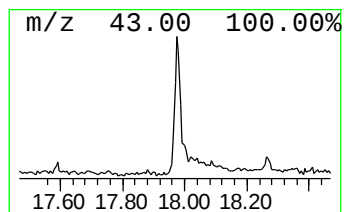
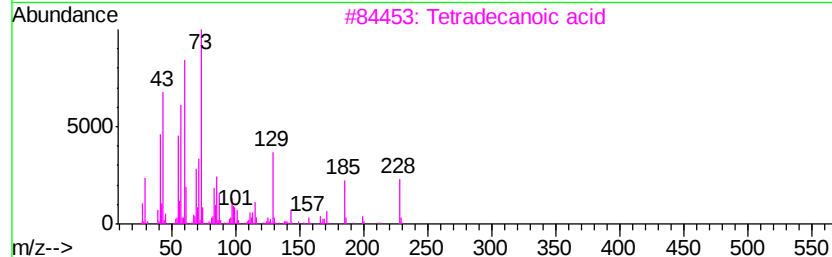
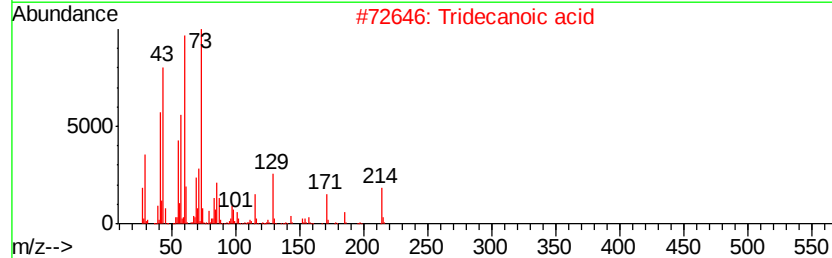
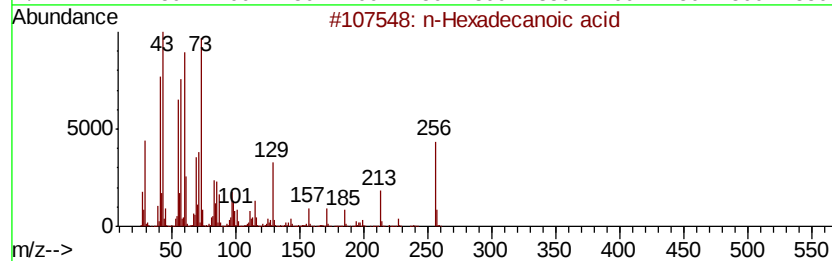
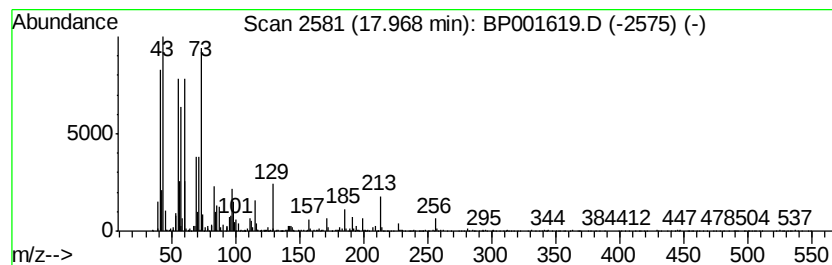
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TIC Library : C:\DATABASE\NIST11.L
 TIC Integration Parameters: LSCINT.P

 Peak Number 5 n-Hexadecanoic acid Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.97	3.67 ng	82765	Phenanthrene-d10	17.04

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	83
2		Tridecanoic acid	214	C13H26O2	000638-53-9	74
3		Tetradecanoic acid	228	C14H28O2	000544-63-8	72
4		Undecanoic acid	186	C11H22O2	000112-37-8	64
5		Octadecanoic acid	284	C18H36O2	000057-11-4	60



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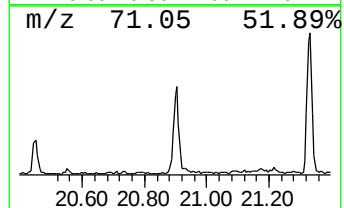
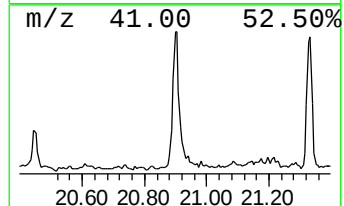
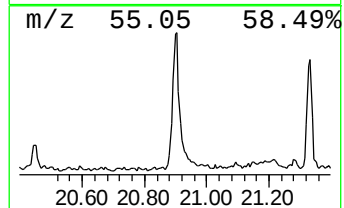
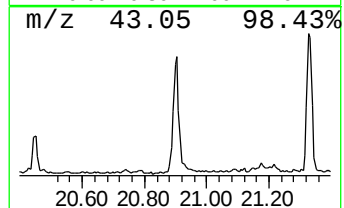
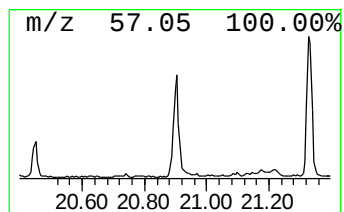
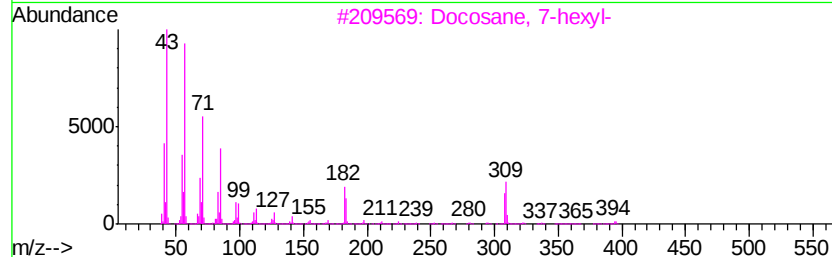
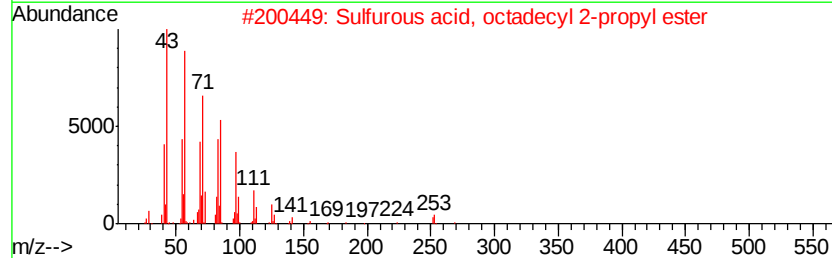
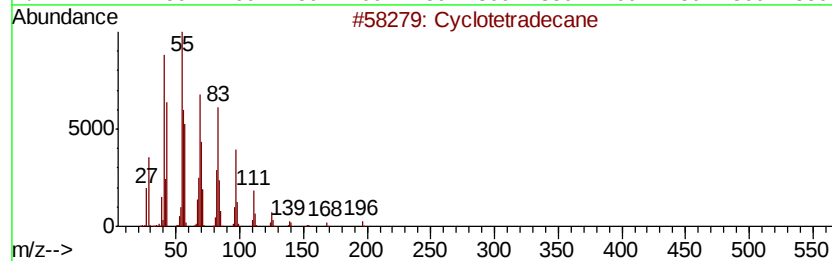
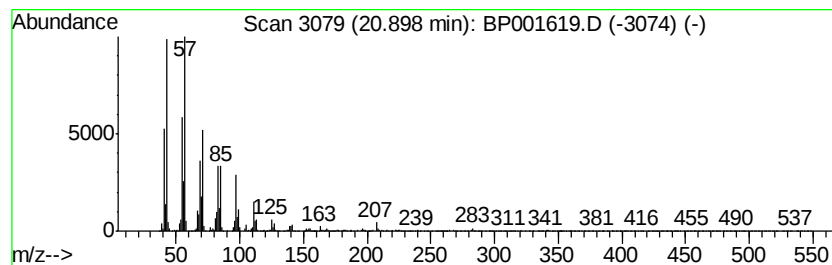
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Cyclotetradecane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.90	7.26 ng	185853	Chrysene-d12	21.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclotetradecane	196	C14H28	000295-17-0	83
2		Sulfurous acid, octadecyl 2-prop...	376	C21H44O3S	1000309-12-7	83
3		Docosane, 7-hexyl-	394	C28H58	055373-86-9	76
4		1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	70
5		8-Heptadecene	238	C17H34	002579-04-6	64



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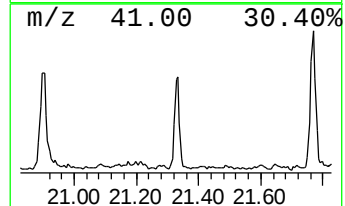
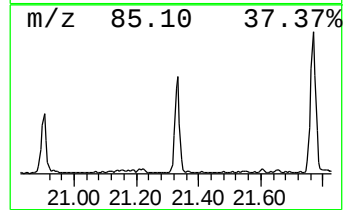
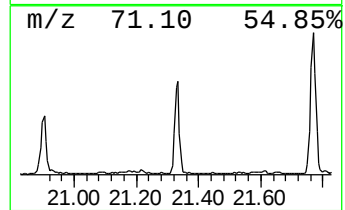
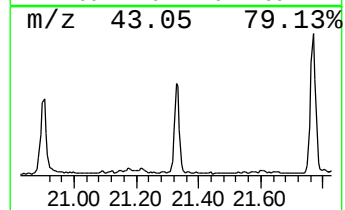
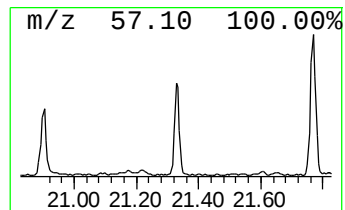
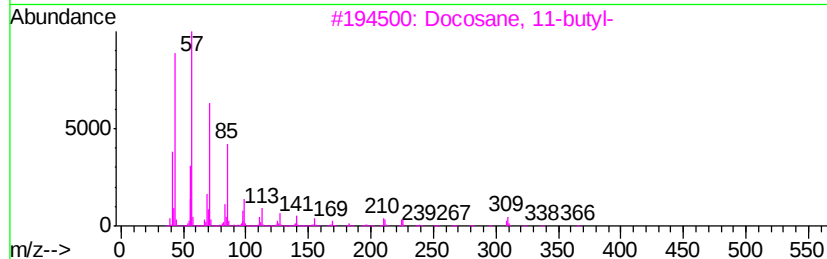
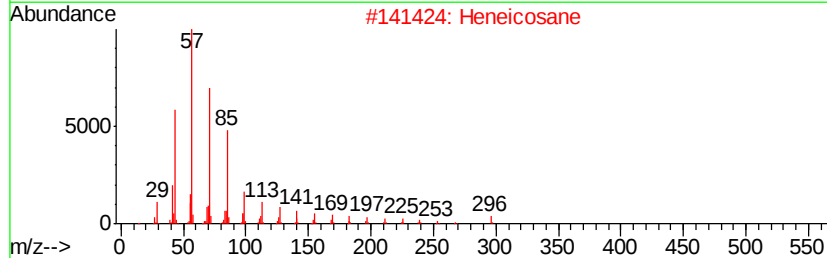
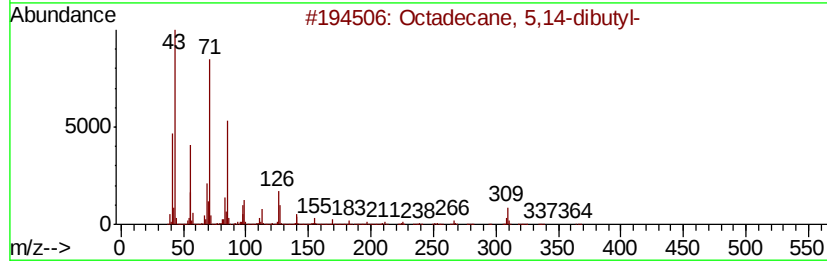
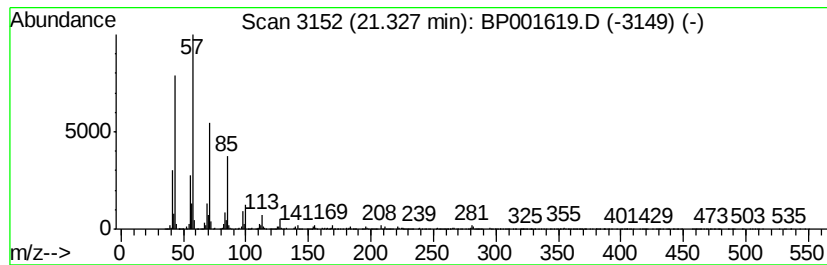
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TIC Library : C:\DATABASE\NIST11.L
 TIC Integration Parameters: LSCINT.P

 Peak Number 7 Octadecane, 5,14-dibutyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.33	5.89 ng	150919	Chrysene-d12	21.14

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecane, 5,14-dibutyl-	366	C26H54	055282-13-8	93
2			Heneicosane	296	C21H44	000629-94-7	93
3			Docosane, 11-butyl-	366	C26H54	013475-76-8	91
4			2-Bromo dodecane	248	C12H25Br	013187-99-0	91
5			Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	91



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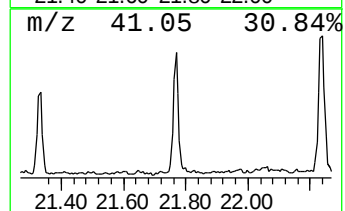
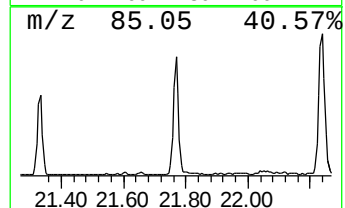
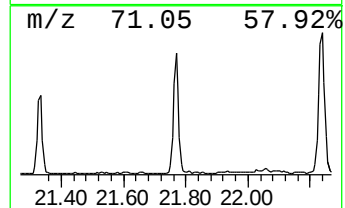
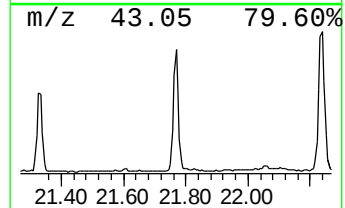
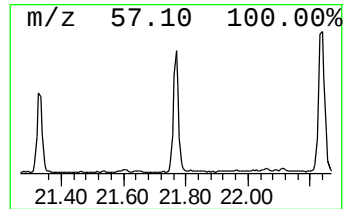
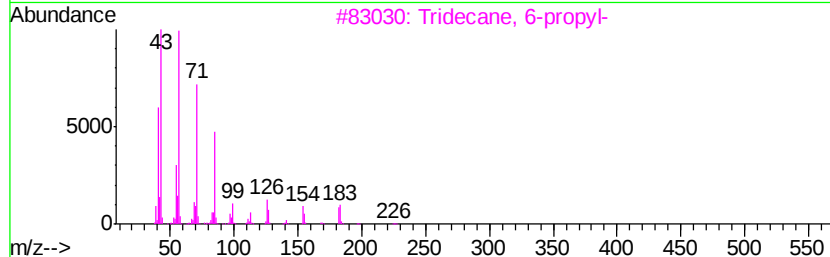
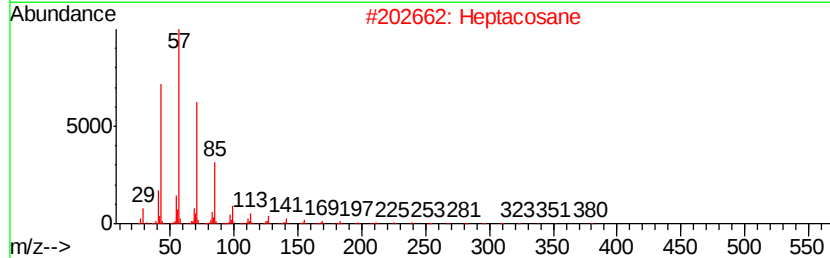
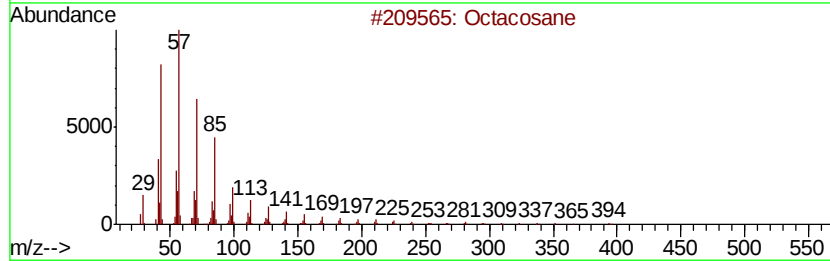
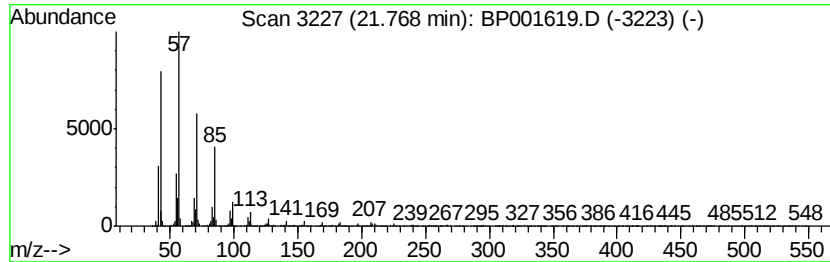
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Octacosane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.77	9.26 ng	237157	Chrysene-d12	21.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosane	394	C28H58	000630-02-4	96
2		Heptacosane	380	C27H56	000593-49-7	91
3		Tridecane, 6-propyl-	226	C16H34	055045-10-8	91
4		Heneicosane	296	C21H44	000629-94-7	90
5		Heneicosane, 11-decyl-	437	C31H64	055320-06-4	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP011520\
Data File : BP001619.D
Acq On : 15 Jan 2020 20:47
Operator : JU
Sample : L1126-07
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
FB-01142020

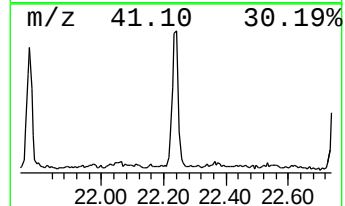
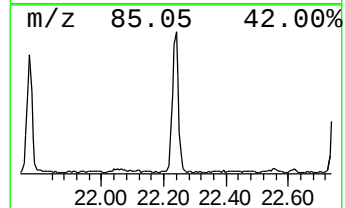
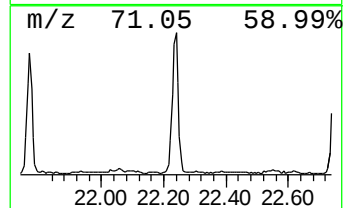
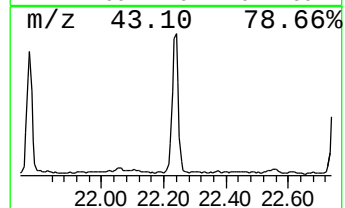
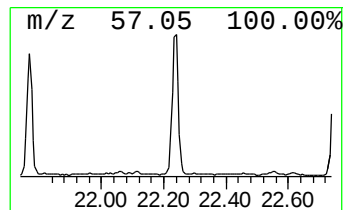
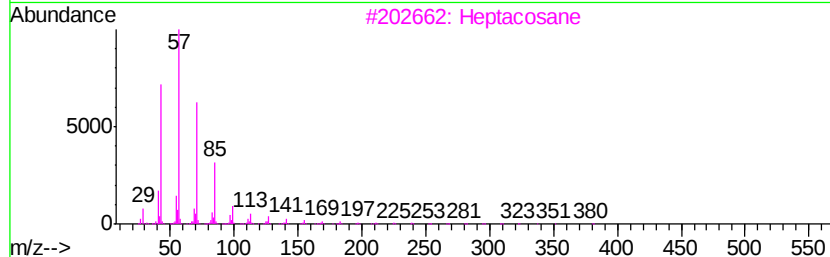
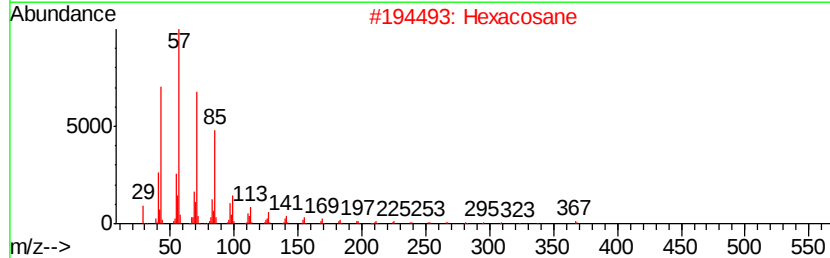
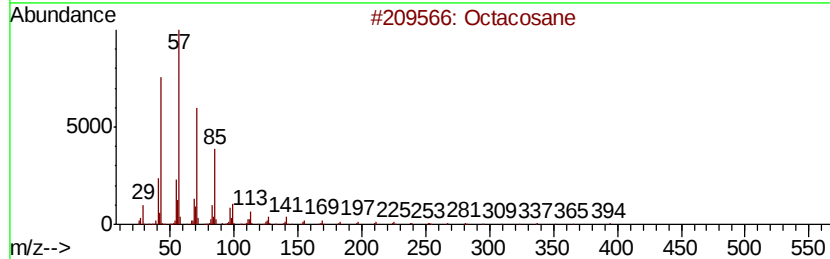
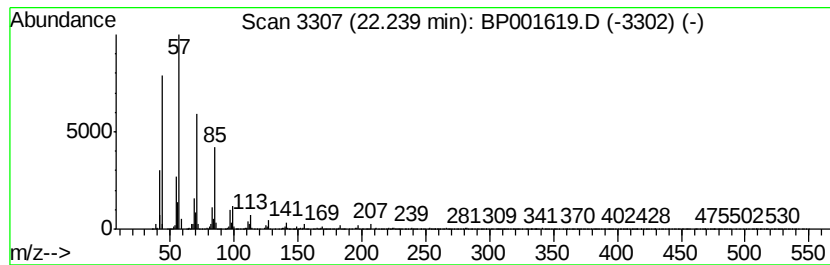
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Heptacosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.24	12.66 ng	324160	Chrysene-d12	21.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosane	394	C28H58	000630-02-4	91
2		Hexacosane	366	C26H54	000630-01-3	91
3		Heptacosane	380	C27H56	000593-49-7	91
4		Eicosane, 7-hexyl-	366	C26H54	055333-99-8	91
5		Tetracosane	338	C24H50	000646-31-1	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP011520\
Data File : BP001619.D
Acq On : 15 Jan 2020 20:47
Operator : JU
Sample : L1126-07
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
FB-01142020

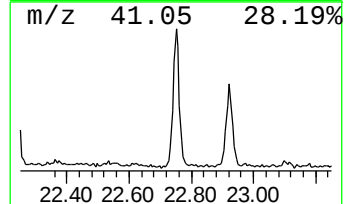
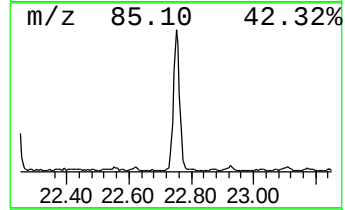
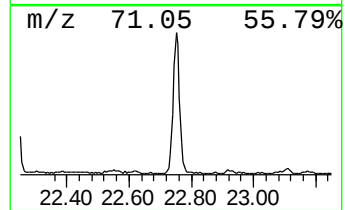
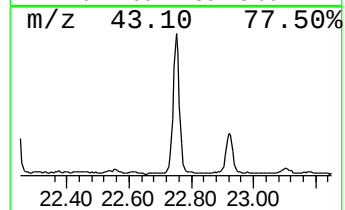
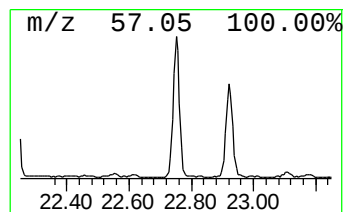
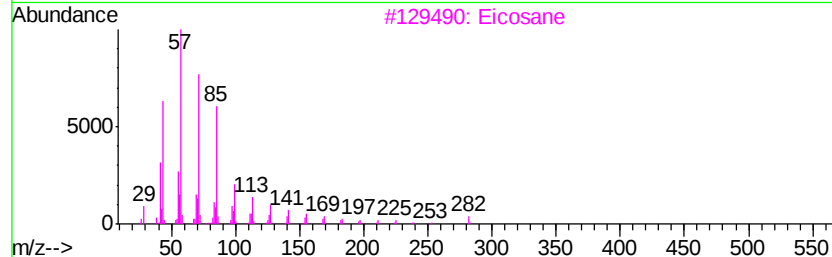
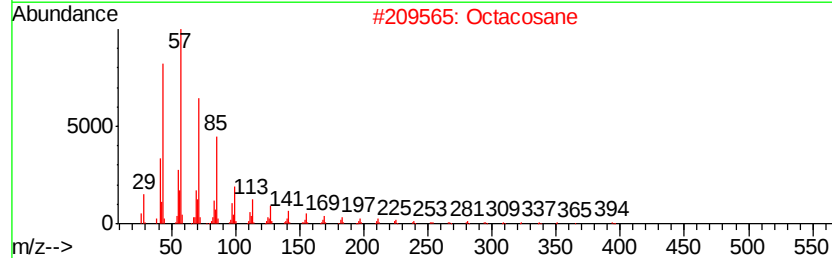
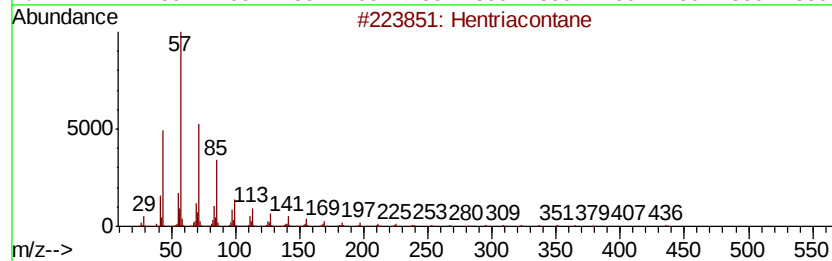
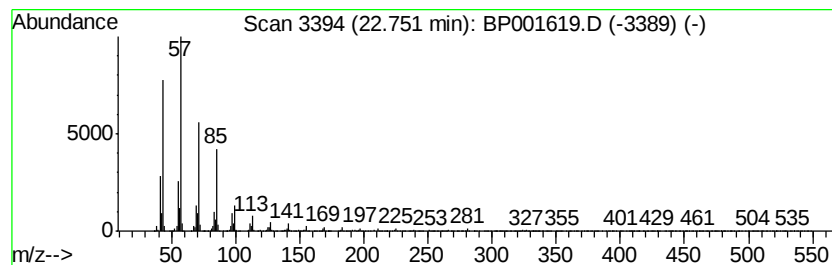
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Hentriacontane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.75	16.37 ng	350166	Perylene-d12	23.37

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hentriacontane		437	C31H64	000630-04-6	98
2	Octacosane		394	C28H58	000630-02-4	97
3	Eicosane		282	C20H42	000112-95-8	94
4	Hexacosane		366	C26H54	000630-01-3	91
5	Tetracosane		338	C24H50	000646-31-1	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP011520\
Data File : BP001619.D
Acq On : 15 Jan 2020 20:47
Operator : JU
Sample : L1126-07
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
FB-01142020

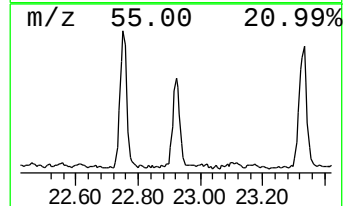
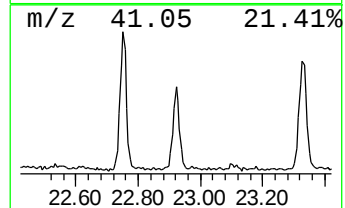
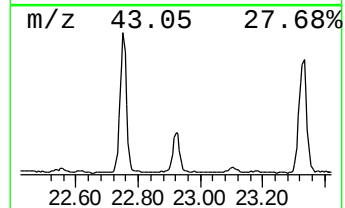
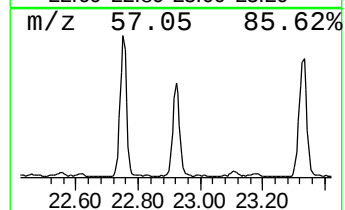
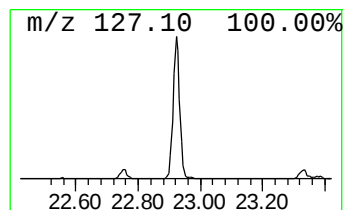
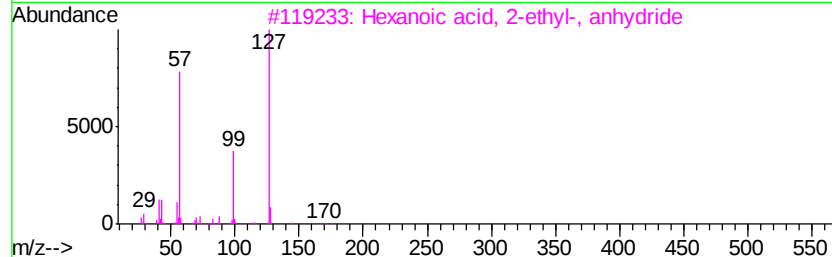
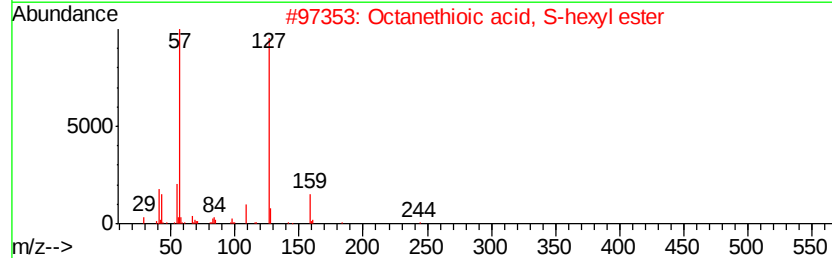
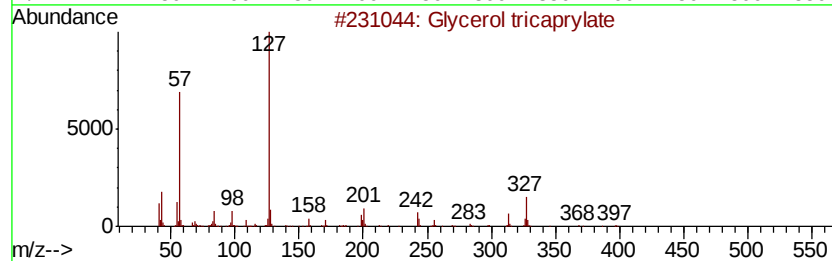
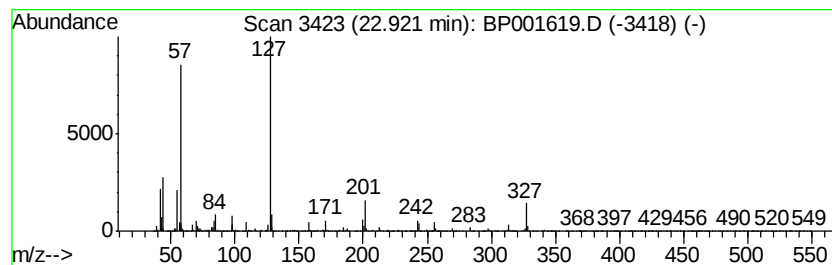
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Glycerol tricaprlylate Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.92	12.22 ng	261397	Perylene-d12	23.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Glycerol tricaprlylate	470	C27H50O6	000538-23-8	80
2		Octanethioic acid, S-hexyl ester	244	C14H28OS	055590-85-7	58
3		Hexanoic acid, 2-ethyl-, anhydride	270	C16H30O3	036765-89-6	53
4		Caprylic anhydride	270	C16H30O3	000623-66-5	53
5		Octanoic acid, 2-formyl-4,6-dich...	316	C15H18Cl2O3	1000330-95-4	53



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP011520\
Data File : BP001619.D
Acq On : 15 Jan 2020 20:47
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Misc :
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Instrument :
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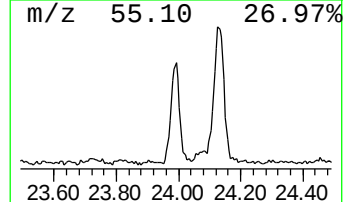
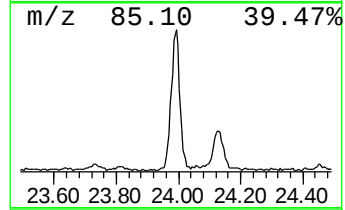
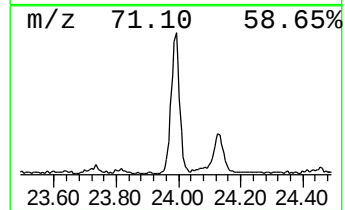
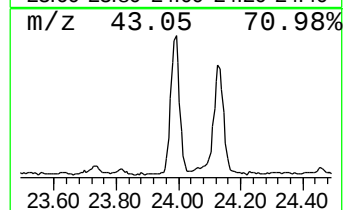
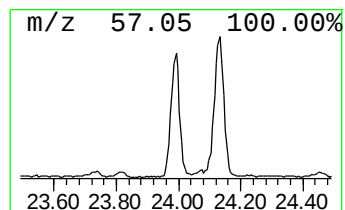
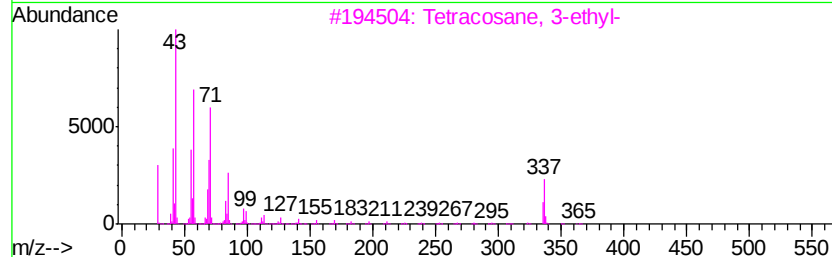
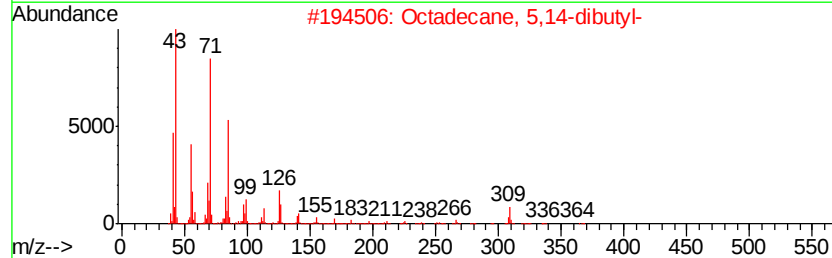
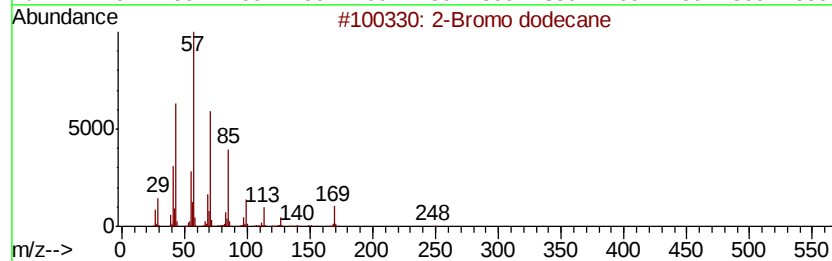
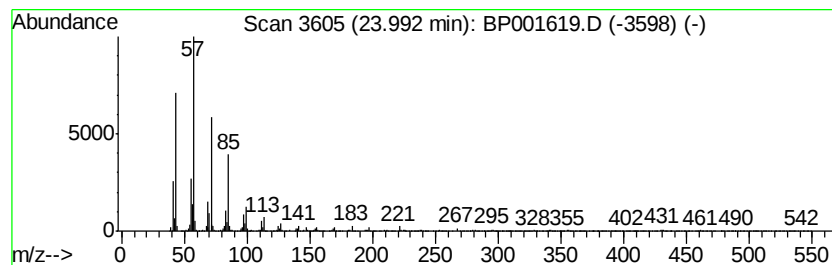
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 2-Bromo dodecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.99	14.96 ng	320119	Perylene-d12	23.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Bromo dodecane	248	C12H25Br	013187-99-0	93
2		Octadecane, 5,14-dibutyl-	366	C26H54	055282-13-8	92
3		Tetracosane, 3-ethyl-	366	C26H54	055282-17-2	91
4		Heptacosane	380	C27H56	000593-49-7	90
5		Heneicosane	296	C21H44	000629-94-7	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP011520\
 Data File : BP001619.D
 Acq On : 15 Jan 2020 20:47
 Operator : JU
 Sample : L1126-07
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

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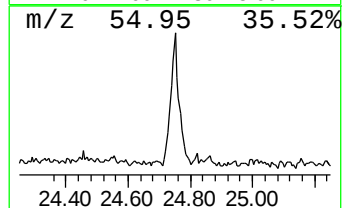
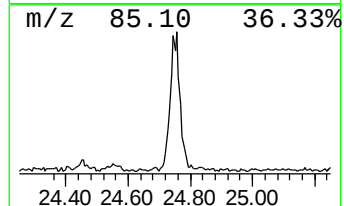
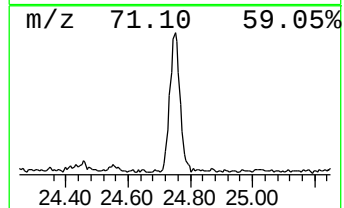
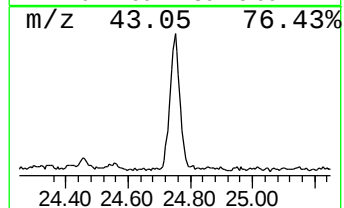
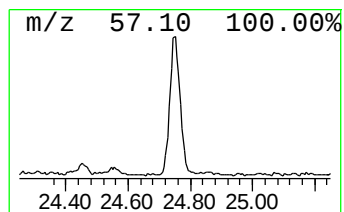
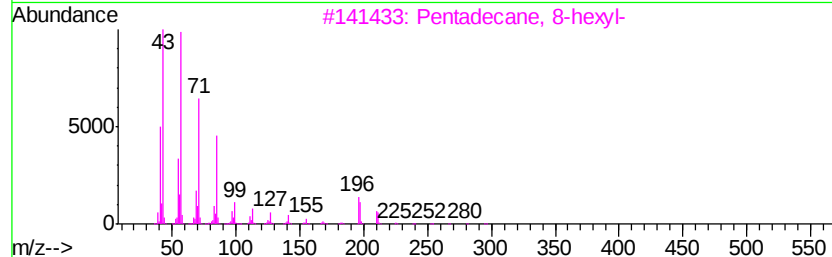
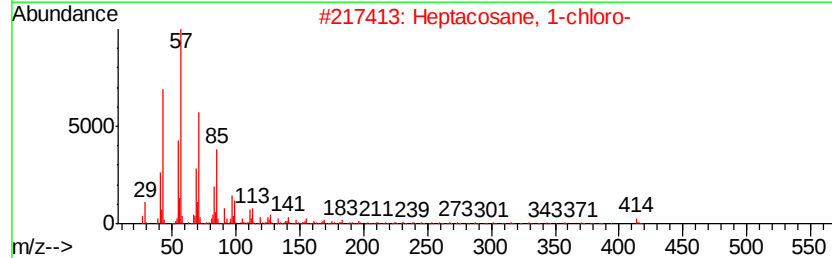
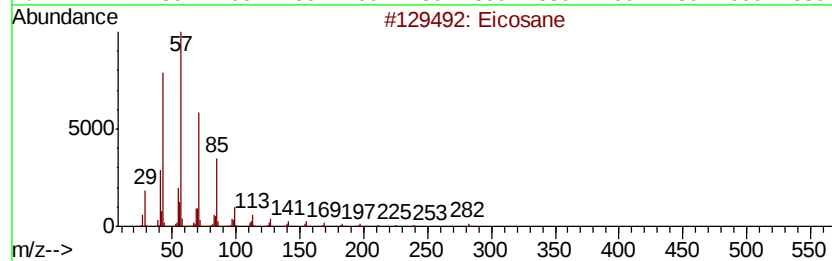
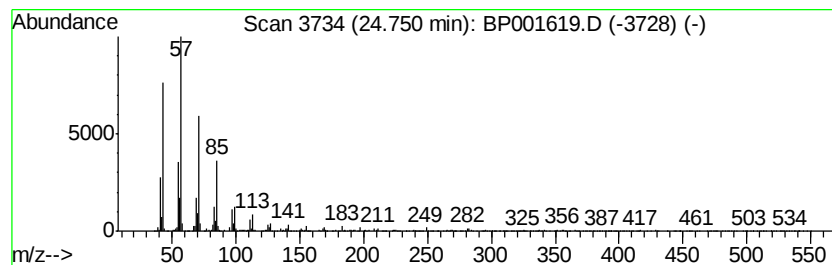
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 14 Eicosane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.75	8.01 ng	171308	Perylene-d12	23.37

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane	282	C20H42	000112-95-8	96
2			Heptacosane, 1-chloro-	414	C27H55Cl	062016-79-9	95
3			Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	91
4			Docosane, 7-hexyl-	394	C28H58	055373-86-9	91
5			Tetracosane, 3-ethyl-	366	C26H54	055282-17-2	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP011520\
Data File : BP001619.D
Acq On : 15 Jan 2020 20:47
Operator : JU
Sample : L1126-07
Misc :
ALS Vial : 15 Sample Multiplier: 1

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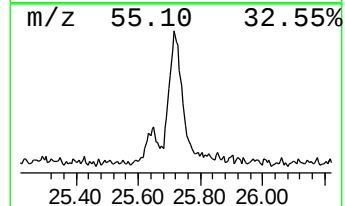
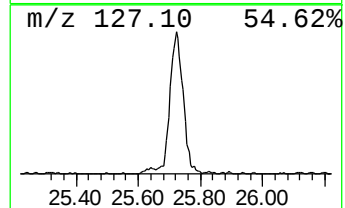
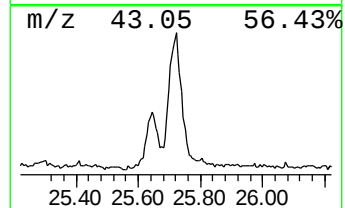
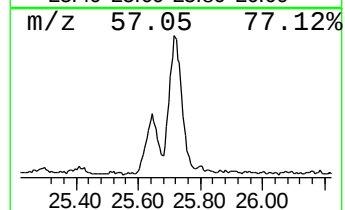
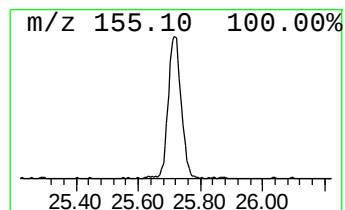
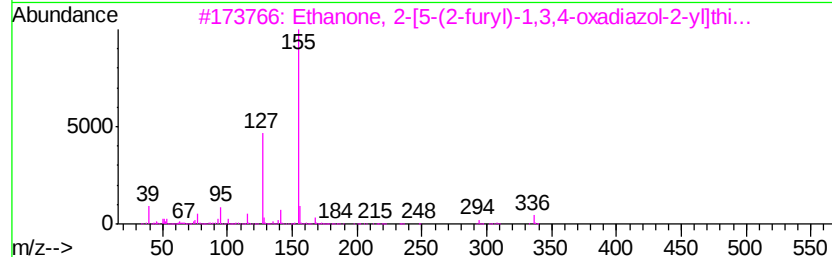
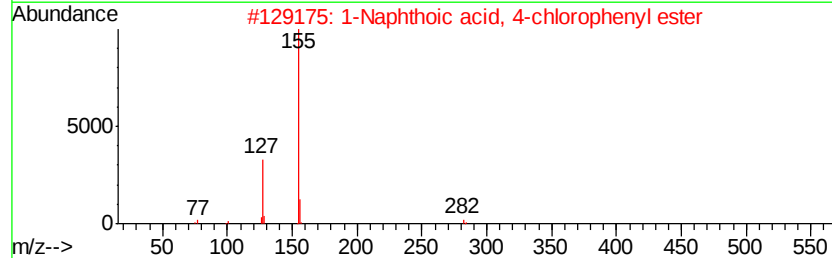
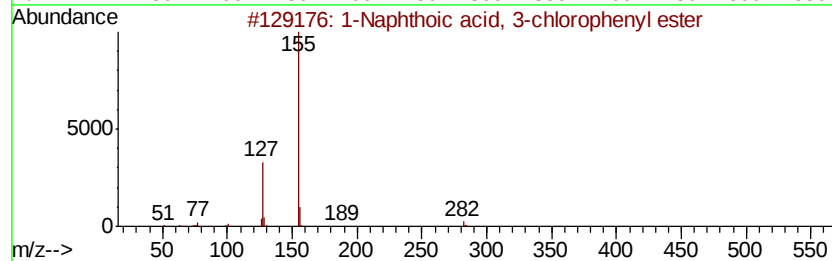
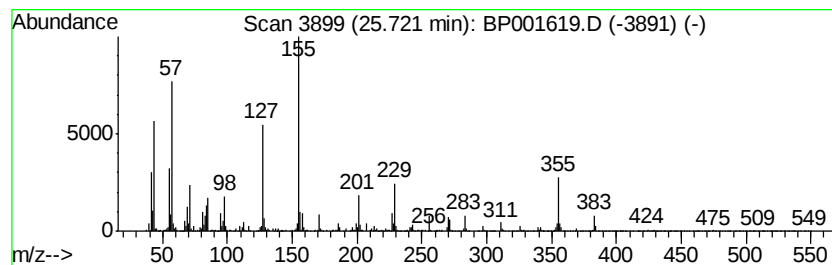
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 1-Naphthoic acid, 3-chlorop... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.72	14.88 ng	318337	Perylene-d12	23.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Naphthoic acid, 3-chlorophenyl...	282	C17H11ClO2	1000325-54-1	50
2		1-Naphthoic acid, 4-chlorophenyl...	282	C17H11ClO2	1000307-82-4	50
3		Ethanone, 2-[5-(2-furyl)-1,3,4-o...	336	C18H12N2O3S	296243-46-4	50
4		2H-Pyran-2-carboxylic acid, 5-et...	200	C10H16O4	019776-82-0	50
5		1-Naphthoic acid, 2,4,6-trichlor...	350	C17H9Cl3O2	1000325-54-3	46



Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP011520\
 Data File : BP001619.D
 Acq On : 15 Jan 2020 20:47
 Operator : JU
 Sample : L1126-07
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 FB-01142020

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP010820.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	74.7	ng	853831	1	7.63	228505	20.0
unknown7.17	7.17	60.0	ng	685813	1	7.63	228505	20.0
n-Hexadecanoic acid	17.97	3.7	ng	82765	4	17.04	450575	20.0
Cyclotetradecane	20.90	7.3	ng	185853	5	21.14	512190	20.0
Octadecane, 5,14-...	21.33	5.9	ng	150919	5	21.14	512190	20.0
Octacosane	21.77	9.3	ng	237157	5	21.14	512190	20.0
Heptacosane	22.24	12.7	ng	324160	5	21.14	512190	20.0
Hentriacontane	22.75	16.4	ng	350166	6	23.37	427941	20.0
Glycerol tricapry...	22.92	12.2	ng	261397	6	23.37	427941	20.0
2-Bromo dodecane	23.99	15.0	ng	320119	6	23.37	427941	20.0
Eicosane	24.75	8.0	ng	171308	6	23.37	427941	20.0
1-Naphthoic acid,...	25.72	14.9	ng	318337	6	23.37	427941	20.0