

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF072417\
 Data File : BF097010.D
 Acq On : 25 Jul 2017 2:49
 Operator : SJ/JU
 Sample : I4361-09
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RAMP-D-5-(0-2)

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:\HPCHEM1\BNA_F\METHODS\8270-BF071317.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	2.351	73	79	94	rVB	154389	252713	7.70%	0.924%
2	2.751	144	147	158	rBV2	15955	49344	1.50%	0.180%
3	4.675	470	474	489	rVB	230807	385898	11.76%	1.411%
4	5.122	546	550	562	rBV	2861830	3281887	100.00%	12.000%
5	6.181	725	730	733	rBV	2842644	3132518	95.45%	11.454%
6	6.287	744	748	752	rBV	2769081	2912907	88.76%	10.651%
7	6.510	783	786	792	rBV	744988	672330	20.49%	2.458%
8	6.669	809	813	817	rBV	2346284	2206400	67.23%	8.067%
9	7.081	878	883	886	rBV	1940122	1968266	59.97%	7.197%
10	7.792	1000	1004	1008	rBV	1011239	956542	29.15%	3.497%
11	8.875	1183	1188	1191	rBV	2953581	2941142	89.62%	10.754%
12	9.269	1251	1255	1259	rBV	376508	315538	9.61%	1.154%
13	9.539	1297	1301	1305	rBV	1144505	998749	30.43%	3.652%
14	9.992	1376	1378	1381	rVB	50738	43379	1.32%	0.159%
15	10.327	1431	1435	1446	rVB	1713814	1697029	51.71%	6.205%
16	10.463	1455	1458	1464	rBV	84450	89870	2.74%	0.329%
17	10.910	1529	1534	1537	rBV2	44793	40711	1.24%	0.149%
18	11.016	1547	1552	1555	rBV	1035565	920436	28.05%	3.365%
19	11.569	1643	1646	1653	rBV2	80458	84973	2.59%	0.311%
20	12.598	1816	1821	1825	rBV	1983085	2115708	64.47%	7.736%
21	13.510	1971	1976	1985	rBV	102024	150491	4.59%	0.550%
22	13.639	1993	1998	2005	rBV	674787	684020	20.84%	2.501%
23	13.951	2048	2051	2057	rVB	49676	47061	1.43%	0.172%
24	14.551	2150	2153	2157	rBV2	53244	52818	1.61%	0.193%
25	14.721	2178	2182	2185	rBV	45272	45668	1.39%	0.167%
26	14.986	2223	2227	2233	rBV	476009	563889	17.18%	2.062%
27	15.039	2233	2236	2241	rVB	47651	59408	1.81%	0.217%
28	15.209	2261	2265	2271	rVB2	52384	72911	2.22%	0.267%
29	15.404	2294	2298	2303	rBV	73989	107608	3.28%	0.393%
30	15.821	2364	2369	2376	rBV	66913	120773	3.68%	0.442%
31	16.056	2405	2409	2416	rBV7	29024	51181	1.56%	0.187%
32	16.309	2446	2452	2461	rBV2	63018	122756	3.74%	0.449%
33	16.880	2544	2549	2556	rVB2	46624	98532	3.00%	0.360%
34	17.551	2657	2663	2664	rBV3	28993	45259	1.38%	0.165%

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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

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35	18.362	2789	2801	2803	rBV6	23756	60817	1.85%	0.222%
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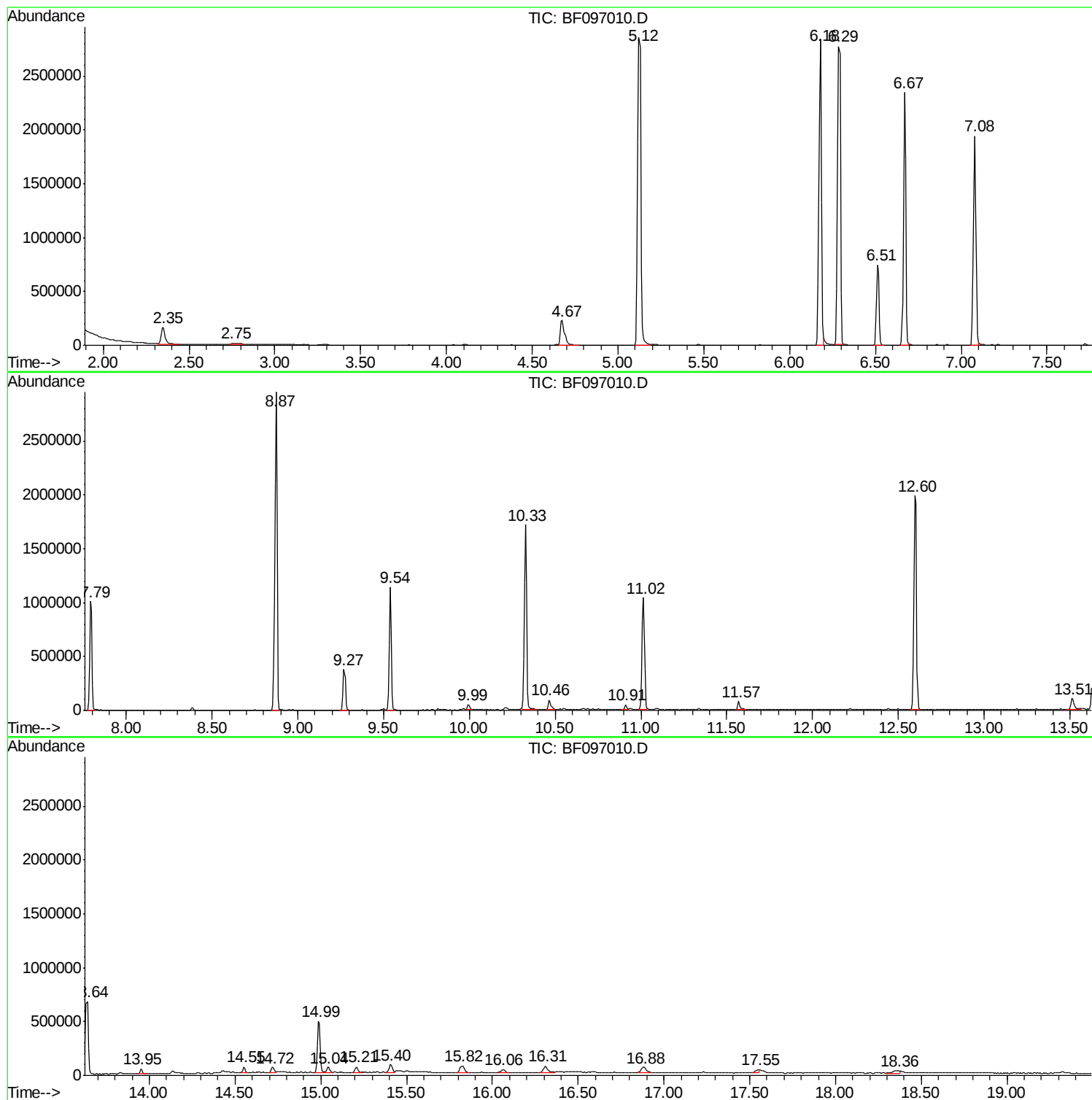
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF071317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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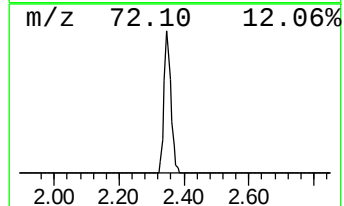
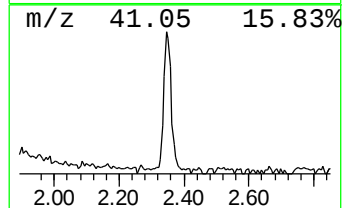
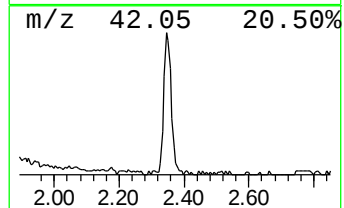
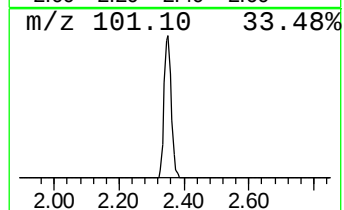
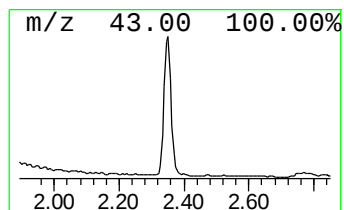
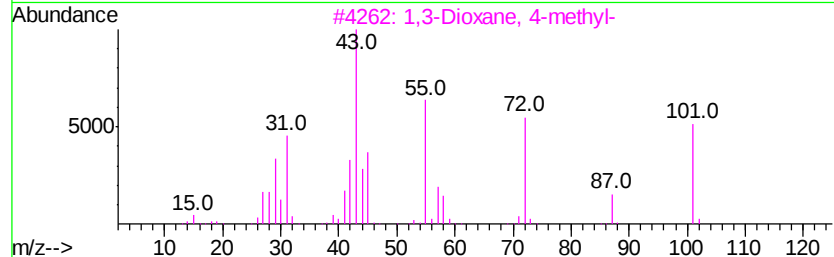
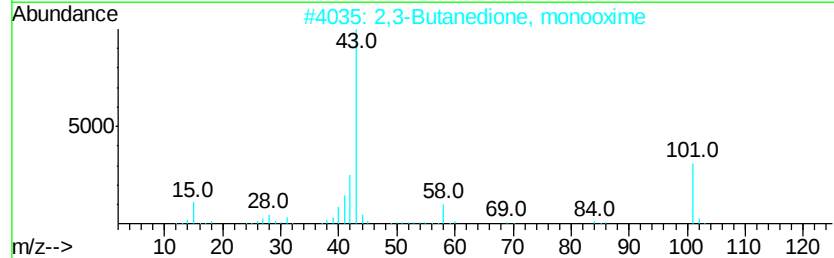
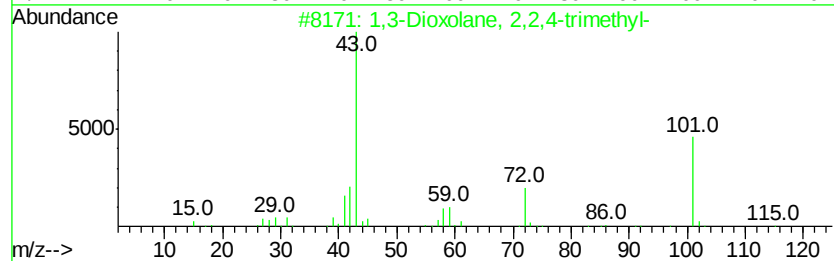
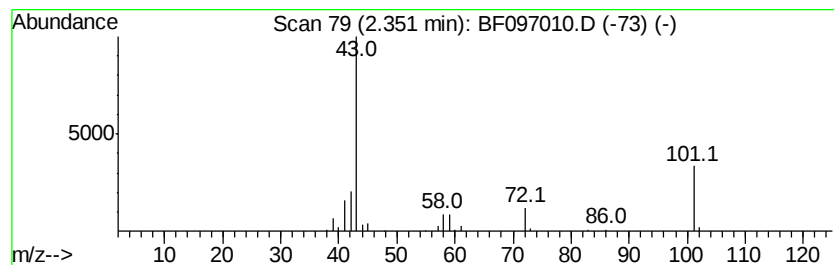
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 1,3-Dioxolane, 2,2,4-trimet... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.35	7.52 ng	252713	1,4-Dichlorobenzene-d4	6.52

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,3-Dioxolane, 2,2,4-trimethyl-	116	C6H12O2	001193-11-9	59
2			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	36
3			1,3-Dioxane, 4-methyl-	102	C5H10O2	001120-97-4	33
4			1,3-Dioxolane-2-methanol, 2,4-di...	132	C6H12O3	053951-43-2	28
5			1-Propanol, 2-methyl-	74	C4H10O	000078-83-1	10



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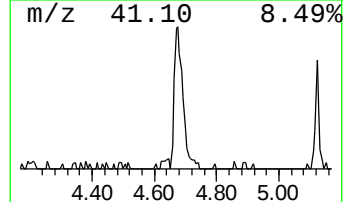
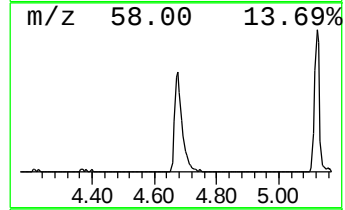
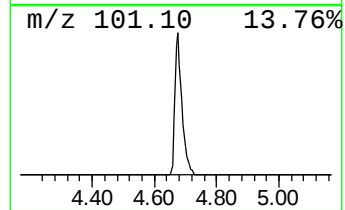
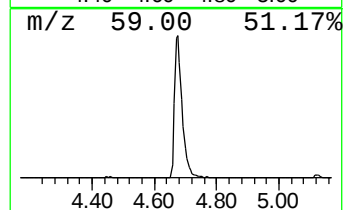
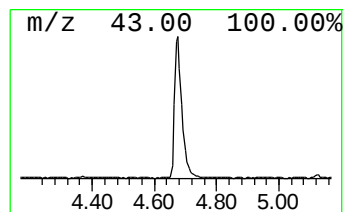
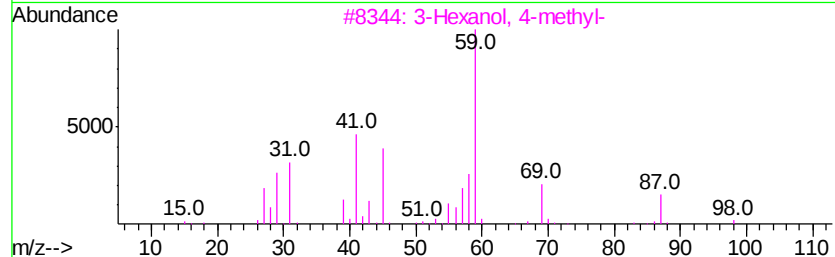
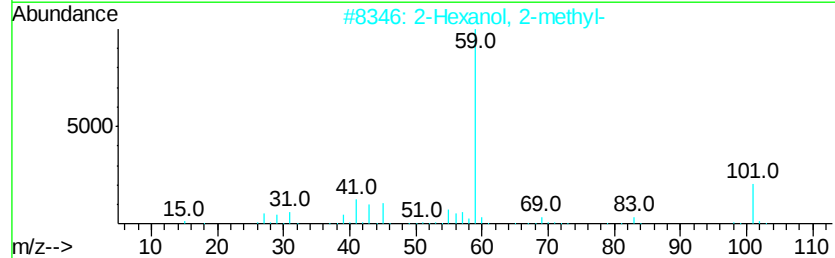
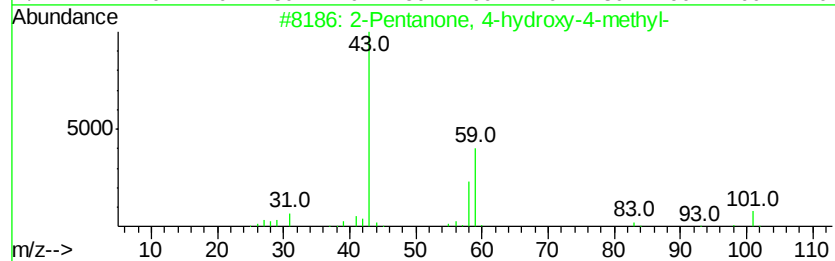
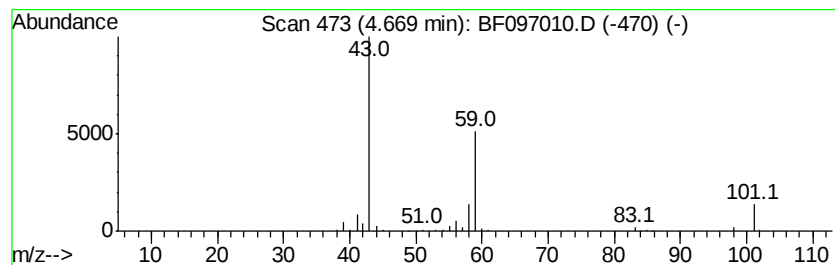
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.67	11.48 ng	385898	1,4-Dichlorobenzene-d4	6.52

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	33
3		3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
4		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	25
5		2-Propanone, 1-(1-methylethoxy)-	116	C6H12O2	042781-12-4	9



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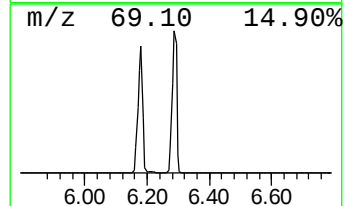
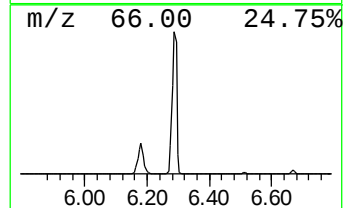
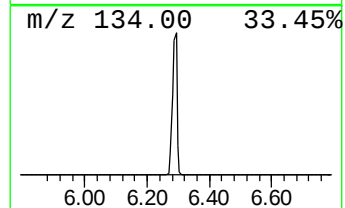
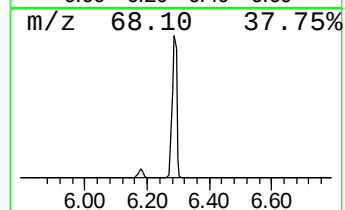
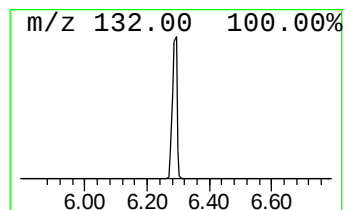
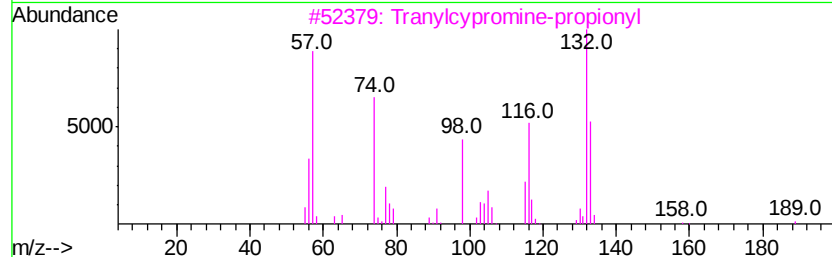
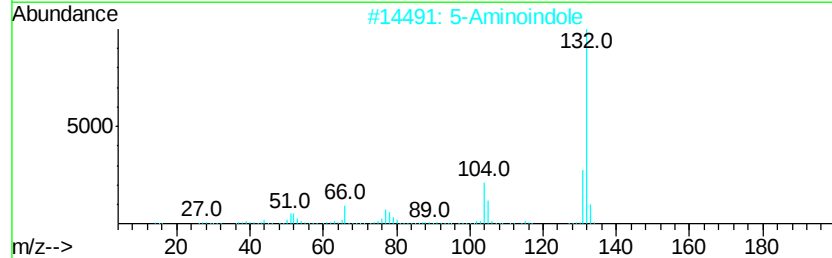
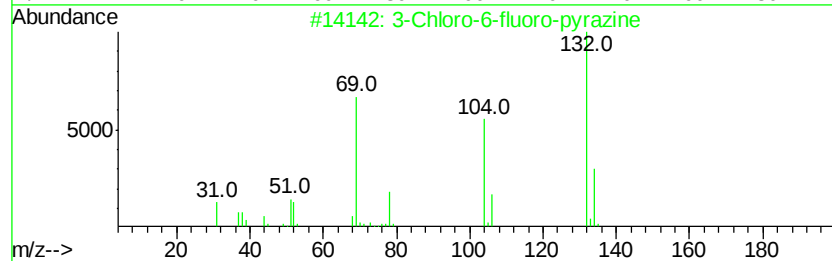
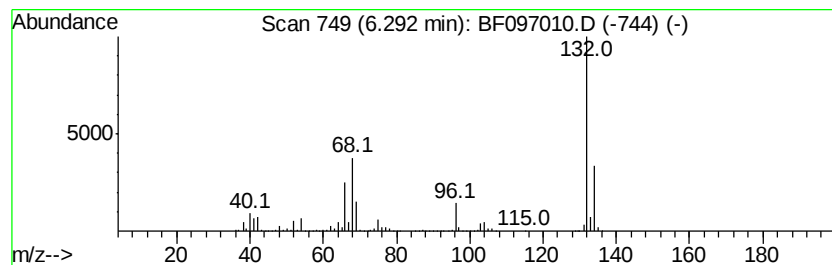
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown6.29 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.29	86.65 ng	2912910	1,4-Dichlorobenzene-d4	6.52

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2			5-Aminoindole	132	C8H8N2	005192-03-0	12
3			Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	10
4			1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9
5			Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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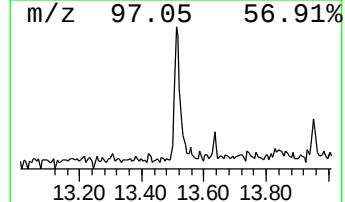
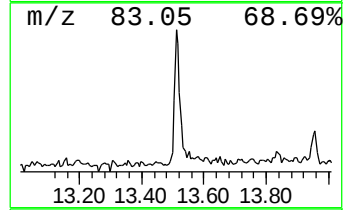
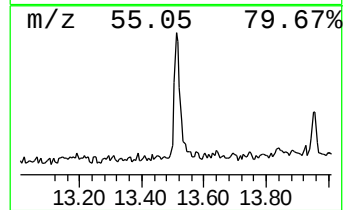
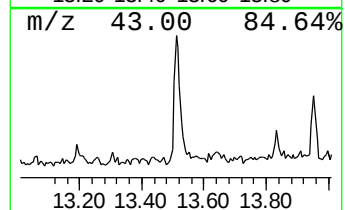
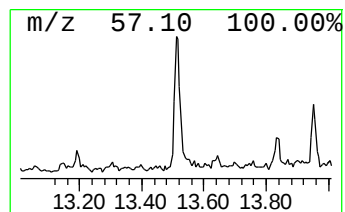
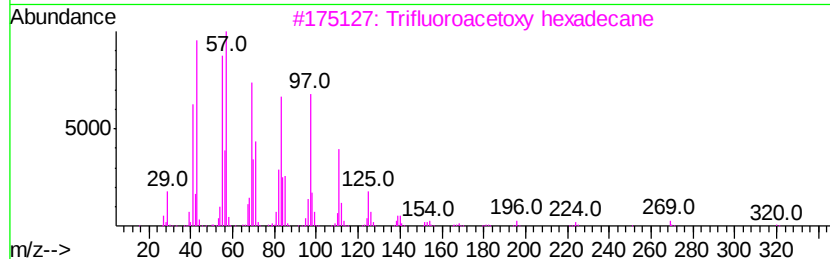
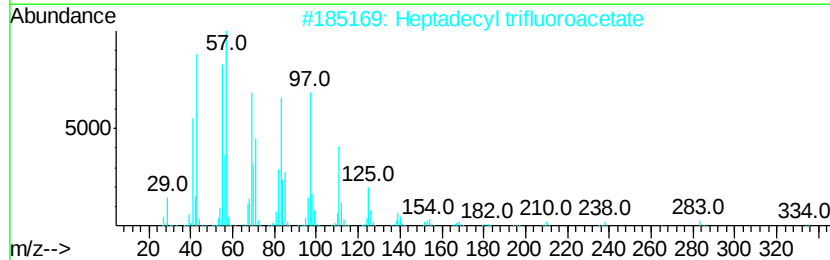
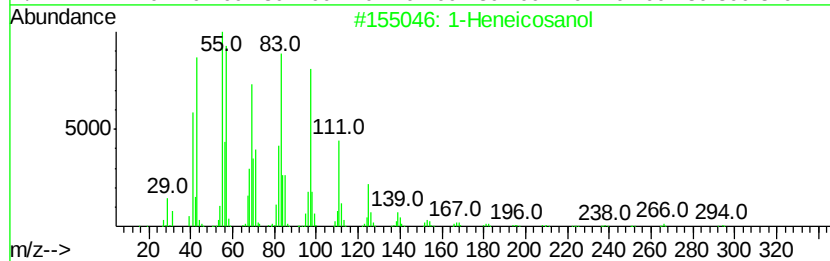
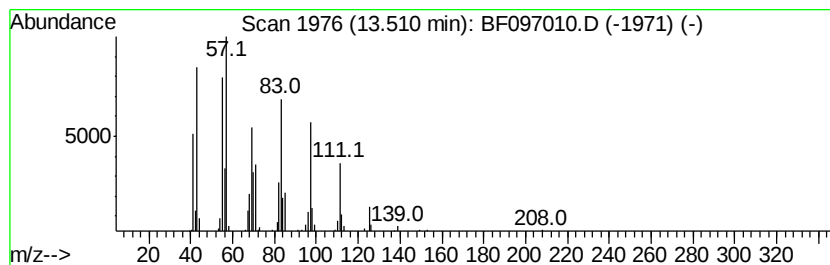
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 1-Heneicosanol Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.51	4.40 ng	150491	Chrysene-d12	13.64	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heneicosanol	312	C21H44O	015594-90-8	95
2	Heptadecyl trifluoroacetate	352	C19H35F3O2	1000351-87-0	95
3	Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	95
4	Behenic alcohol	326	C22H46O	000661-19-8	91
5	1-Octadecanol	270	C18H38O	000112-92-5	91



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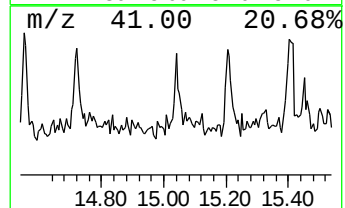
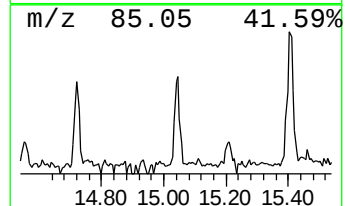
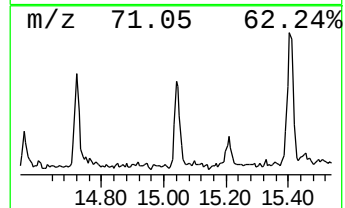
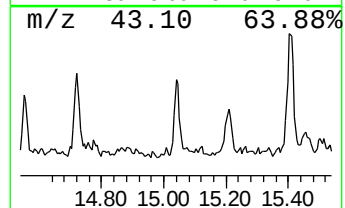
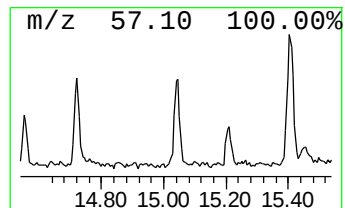
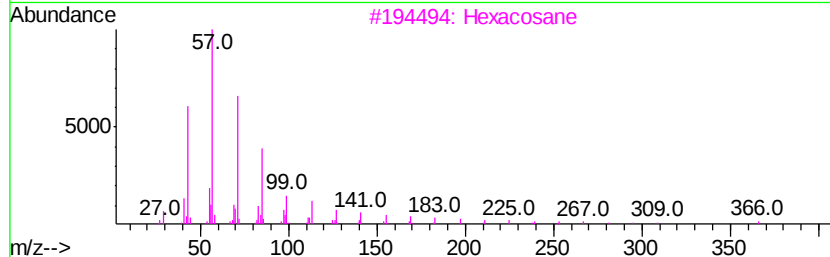
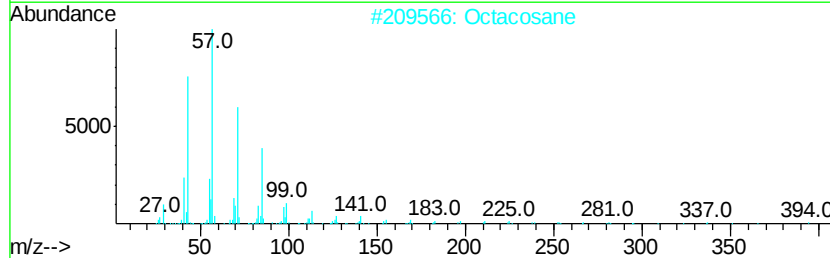
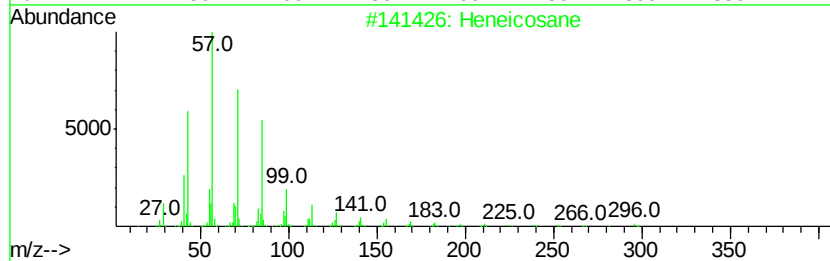
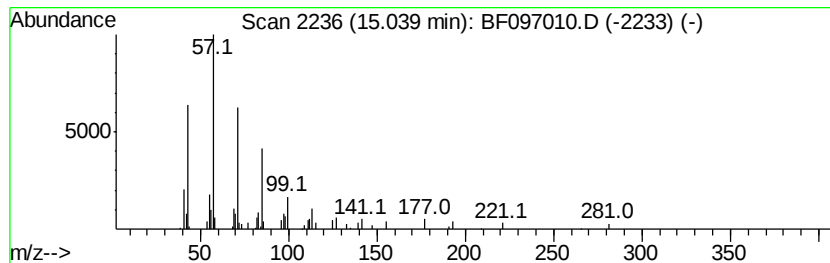
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Peak Number 6 Heneicosane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.04	2.11 ng	59408	Perylene-d12	14.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	90
2		Octacosane	394	C28H58	000630-02-4	87
3		Hexacosane	366	C26H54	000630-01-3	87
4		Heptacosane	380	C27H56	000593-49-7	86
5		Heptadecane	240	C17H36	000629-78-7	86



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Sample : I4361-09
Misc :
ALS Vial : 20 Sample Multiplier: 1

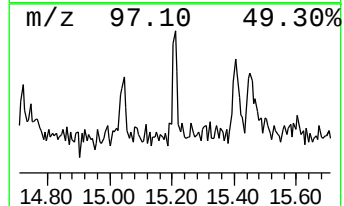
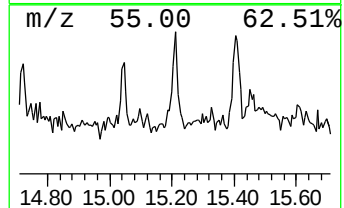
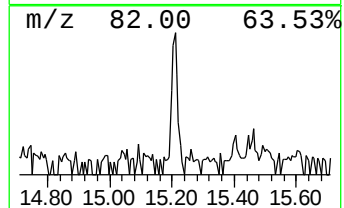
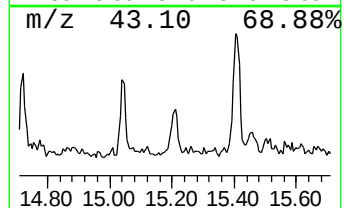
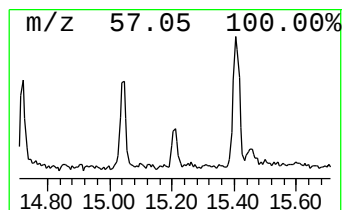
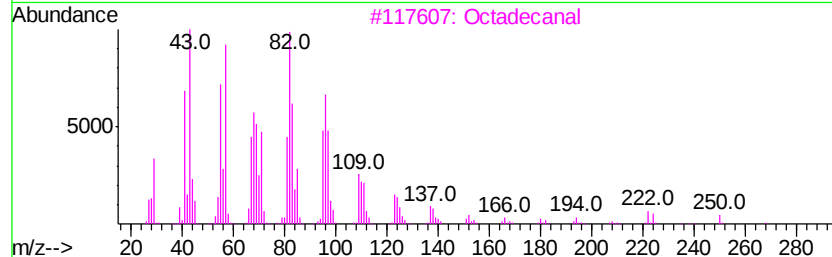
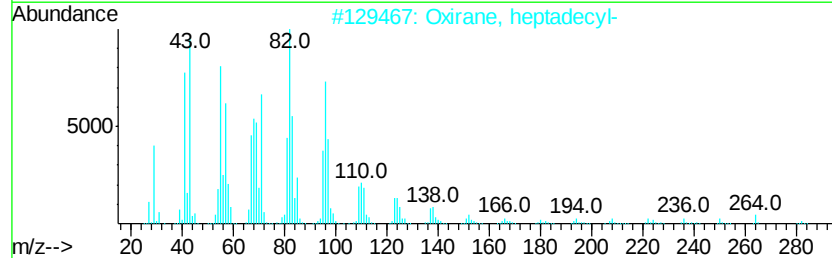
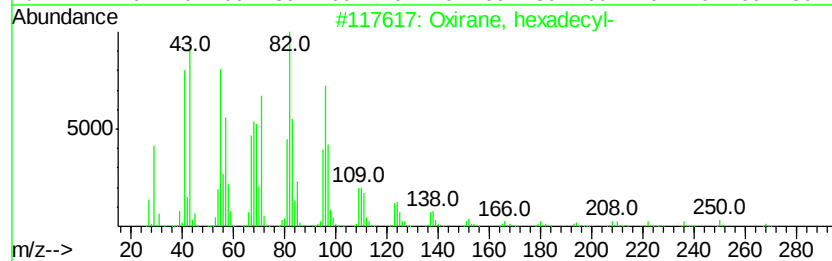
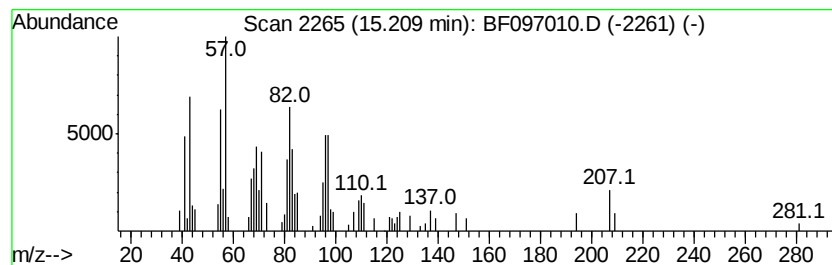
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ClientSampleId :
RAMP-D-5-(0-2)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF071317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Oxirane, hexadecyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD		R.T.
15.21	2.59 ng	72911	Perylene-d12		14.99
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Oxirane, hexadecyl-	268	C18H36O	007390-81-0	72
2	Oxirane, heptadecyl-	282	C19H38O	067860-04-2	64
3	Octadecanal	268	C18H36O	000638-66-4	62
4	15-Octadecenal	266	C18H34O	056554-93-9	62
5	Hexadecanal	240	C16H32O	000629-80-1	58



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF072417\
Data File : BF097010.D
Acq On : 25 Jul 2017 2:49
Operator : SJ/JU
Sample : I4361-09
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
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ClientSampleId :
RAMP-D-5-(0-2)

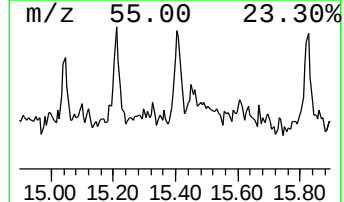
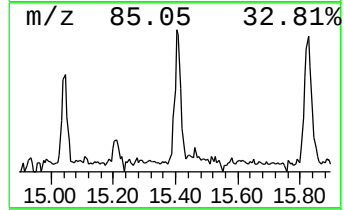
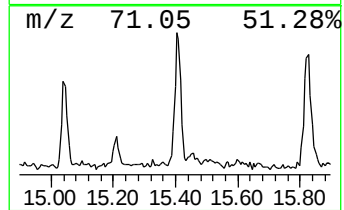
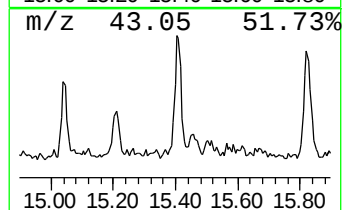
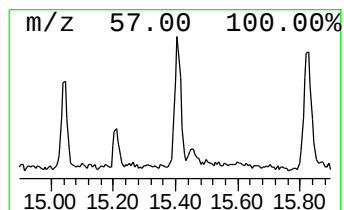
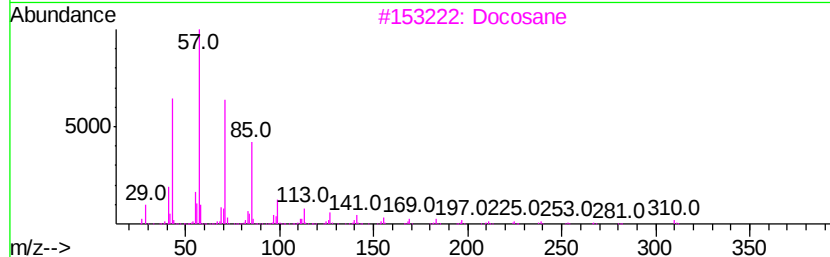
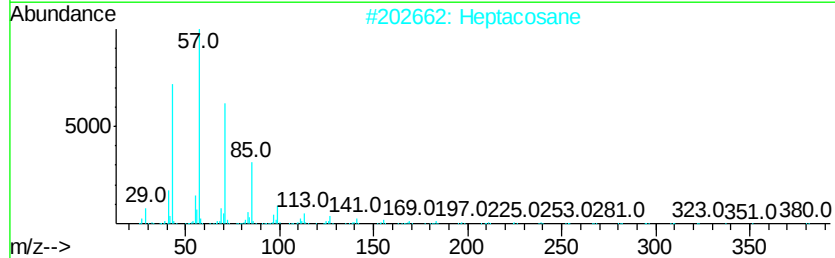
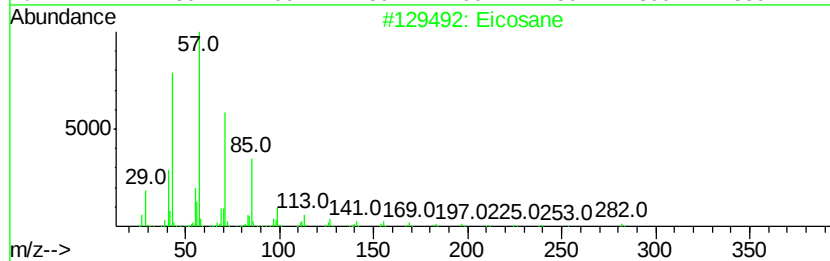
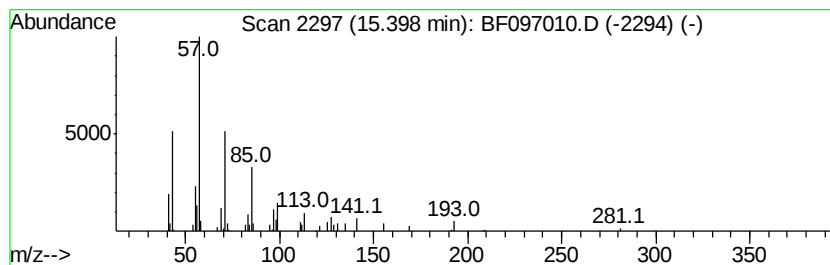
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Eicosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.40	3.82 ng	107608	Perylene-d12	14.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	91
2		Heptacosane	380	C27H56	000593-49-7	86
3		Docosane	310	C22H46	000629-97-0	86
4		Undecane, 2,9-dimethyl-	184	C13H28	017301-26-7	72
5		Heneicosane	296	C21H44	000629-94-7	64



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF072417\
Data File : BF097010.D
Acq On : 25 Jul 2017 2:49
Operator : SJ/JU
Sample : I4361-09
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
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ClientSampleId :
RAMP-D-5-(0-2)

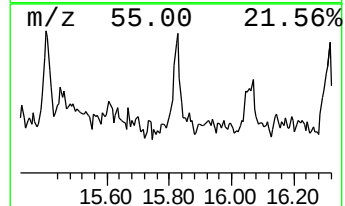
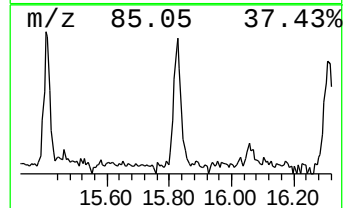
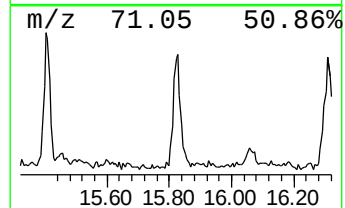
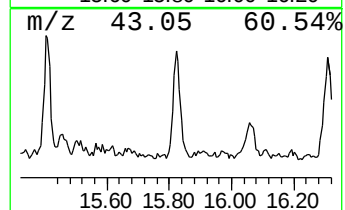
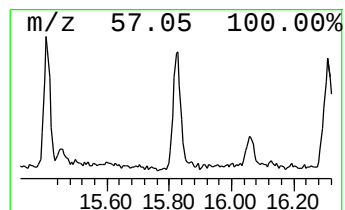
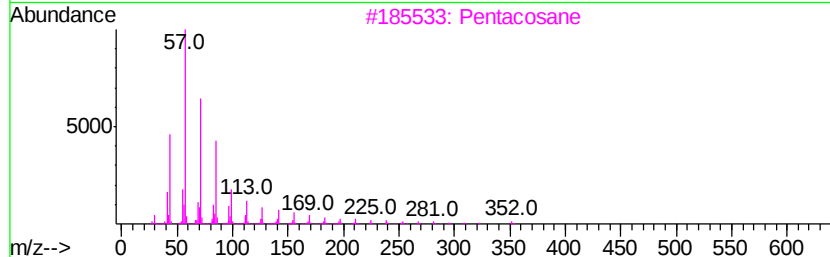
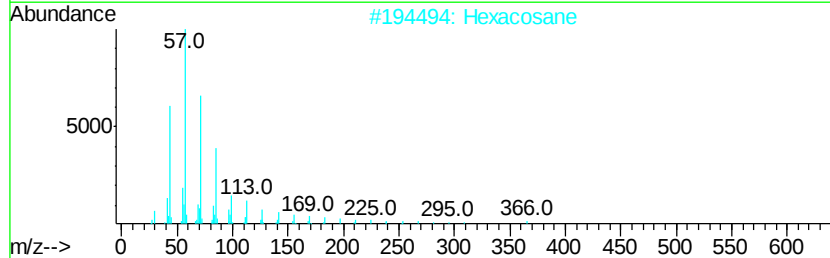
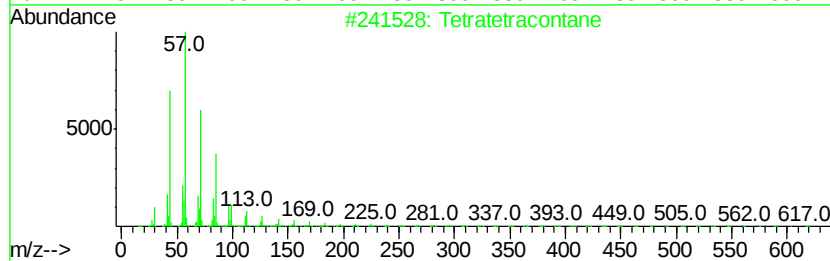
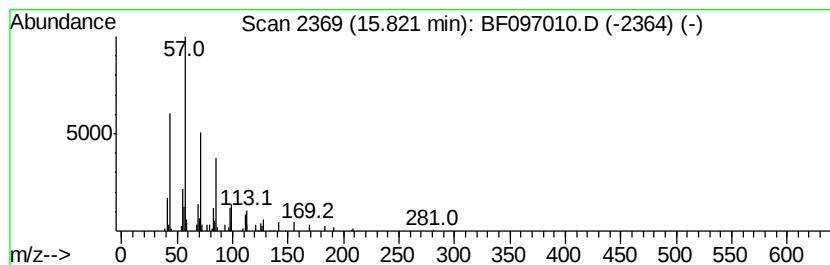
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Tetratetracontane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.82	4.28 ng	120773	Perylene-d12	14.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetratetracontane	619	C44H90	007098-22-8	91
2		Hexacosane	366	C26H54	000630-01-3	91
3		Pentacosane	352	C25H52	000629-99-2	91
4		Hentriacontane	437	C31H64	000630-04-6	91
5		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF072417\
Data File : BF097010.D
Acq On : 25 Jul 2017 2:49
Operator : SJ/JU
Sample : I4361-09
Misc :
ALS Vial : 20 Sample Multiplier: 1

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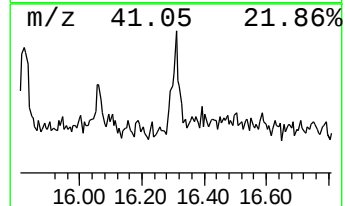
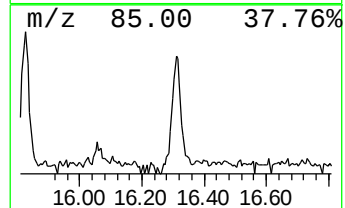
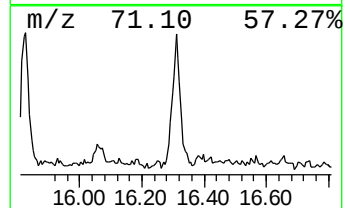
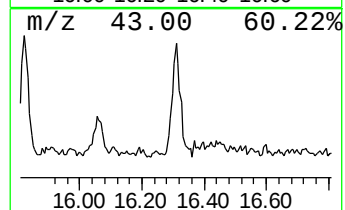
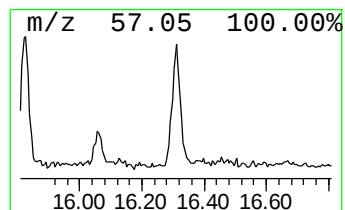
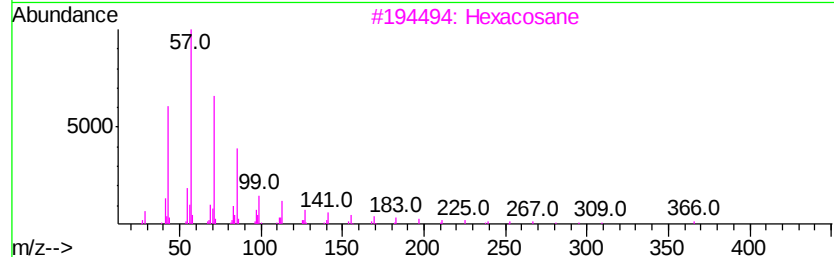
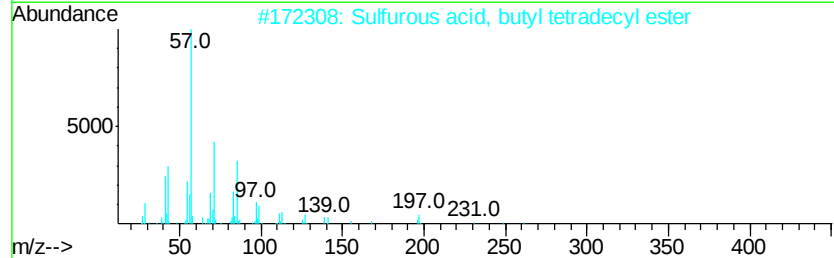
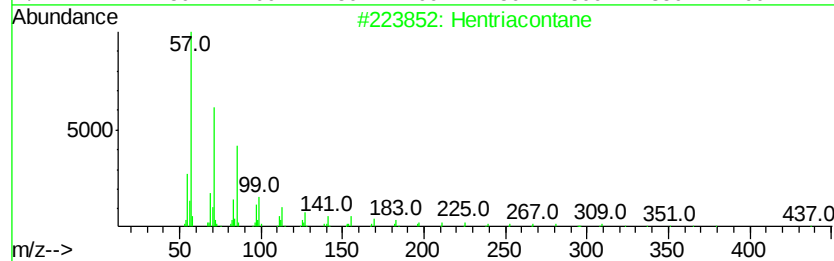
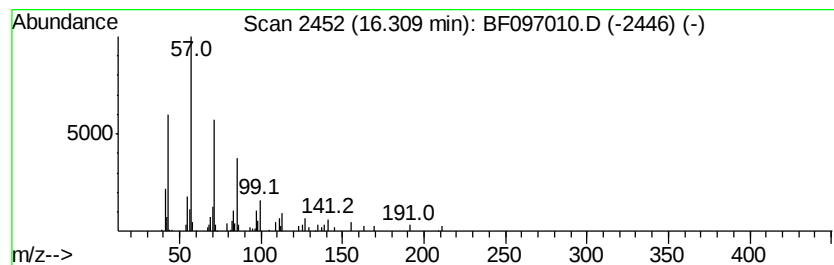
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Hentriacontane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.31	4.35 ng	122756	Perylene-d12	14.99

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hentriacontane	437	C31H64	000630-04-6	90
2			Sulfurous acid, butyl tetradecyl...	334	C18H38O3S	1000309-18-1	90
3			Hexacosane	366	C26H54	000630-01-3	87
4			Tetratetracontane	619	C44H90	007098-22-8	87
5			Triacontane	422	C30H62	000638-68-6	87



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF072417\
Data File : BF097010.D
Acq On : 25 Jul 2017 2:49
Operator : SJ/JU
Sample : I4361-09
Misc :
ALS Vial : 20 Sample Multiplier: 1

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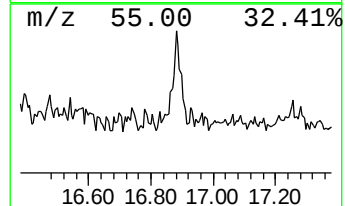
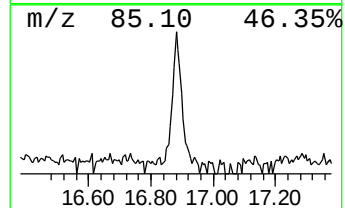
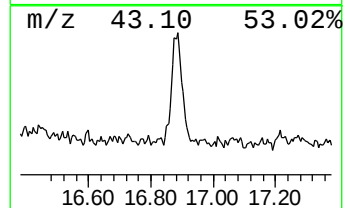
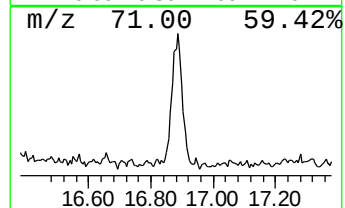
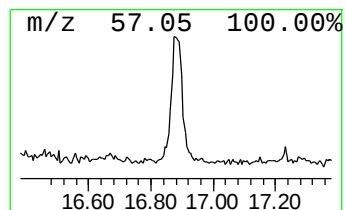
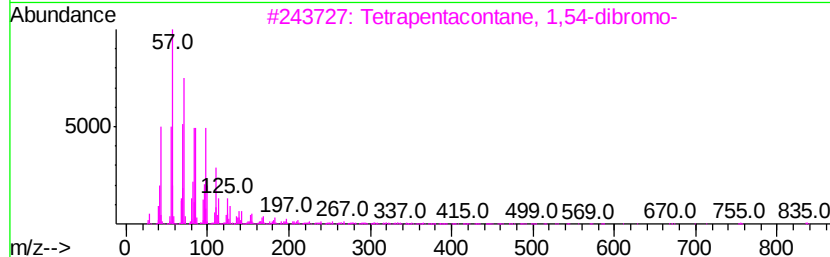
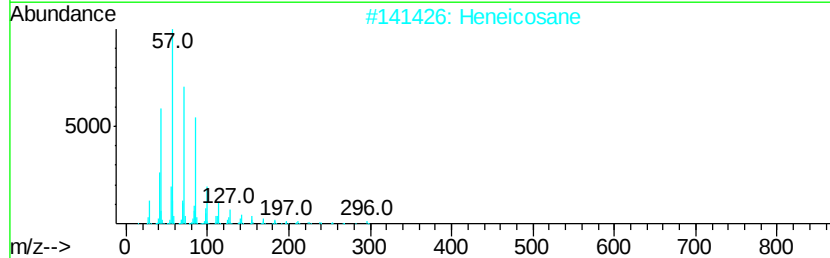
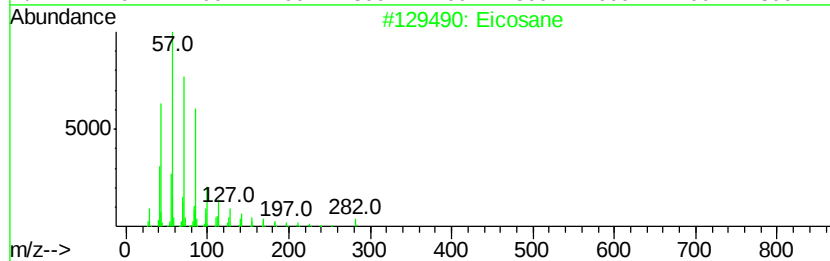
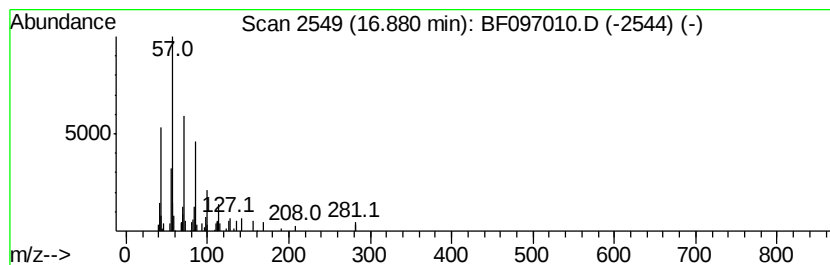
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Tetrapentacontane, 1,54-dib... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.88	3.49 ng	98532	Perylene-d12	14.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	96
2		Heneicosane	296	C21H44	000629-94-7	87
3		Tetrapentacontane, 1,54-dibromo-	915	C54H108Br2	1000156-09-4	86
4		Hexacosane	366	C26H54	000630-01-3	83
5		Octacosane	394	C28H58	000630-02-4	83



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF072417\
Data File : BF097010.D
Acq On : 25 Jul 2017 2:49
Operator : SJ/JU
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Misc :
ALS Vial : 20 Sample Multiplier: 1

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ClientSampleId :
RAMP-D-5-(0-2)

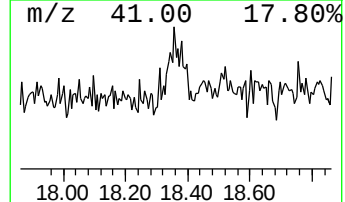
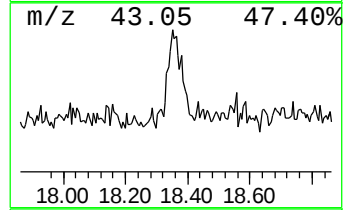
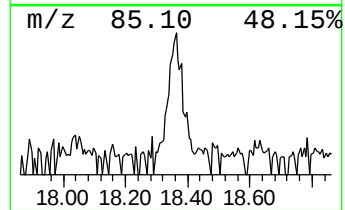
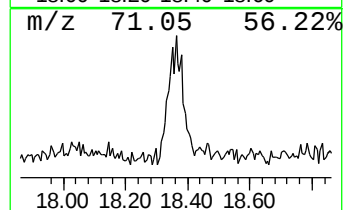
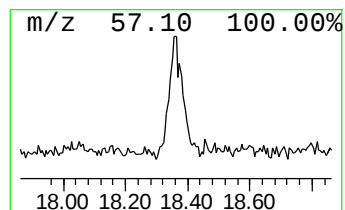
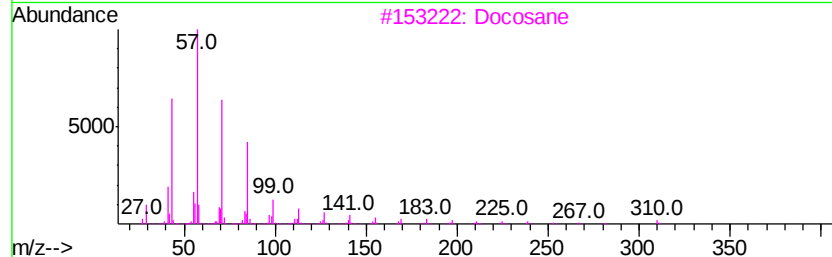
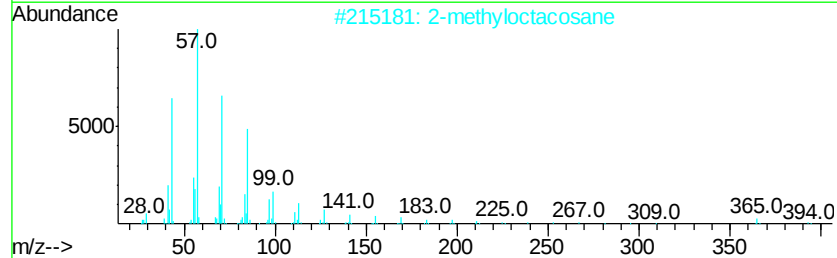
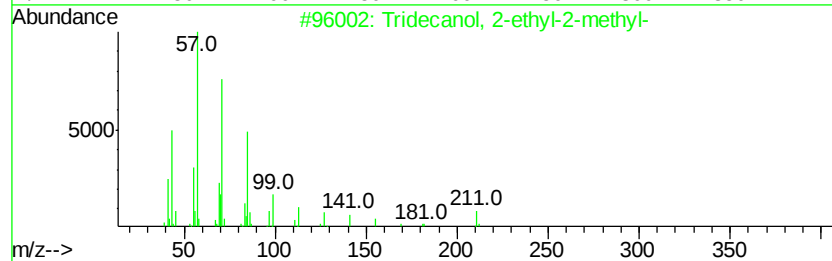
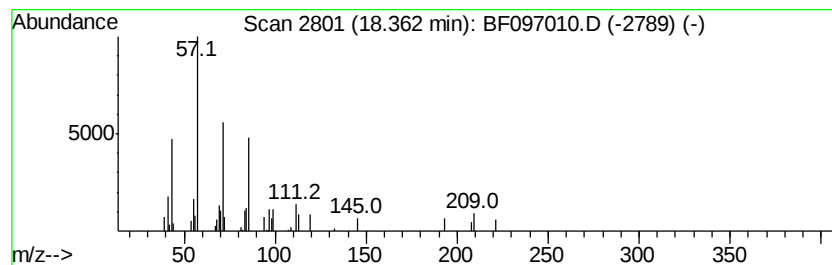
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Tridecanol, 2-ethyl-2-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.36	2.16 ng	60817	Perylene-d12	14.99

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecanol, 2-ethyl-2-methyl-	242	C16H34O	1000115-66-1	64
2			2-methyloctacosane	408	C29H60	1000376-72-8	64
3			Docosane	310	C22H46	000629-97-0	59
4			Octacosane	394	C28H58	000630-02-4	59
5			Methoxyacetic acid, 2-tetradecyl...	286	C17H34O3	1000282-04-8	53



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF072417\
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Acq On : 25 Jul 2017 2:49
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Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
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ClientSampleId :
RAMP-D-5-(0-2)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF071317.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
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2-Pentanone, 4-hy...	4.67	11.5	ng	385898	1	6.52	672330	20.0
unknown6.29	6.29	86.7	ng	2912910	1	6.52	672330	20.0
1-Heneicosanol	13.51	4.4	ng	150491	5	13.64	684020	20.0
Heneicosane	15.04	2.1	ng	59408	6	14.99	563889	20.0
Oxirane, hexadecyl-	15.21	2.6	ng	72911	6	14.99	563889	20.0
Eicosane	15.40	3.8	ng	107608	6	14.99	563889	20.0
Tetratetracontane	15.82	4.3	ng	120773	6	14.99	563889	20.0
Hentriacontane	16.31	4.3	ng	122756	6	14.99	563889	20.0
Tetrapentacontane...	16.88	3.5	ng	98532	6	14.99	563889	20.0
Tridecanol, 2-eth...	18.36	2.2	ng	60817	6	14.99	563889	20.0