

Data Path : Z:\HPCHEM1\BNA F\DATA\BF071816\
 Data File : BF089085.D
 Acq On : 18 Jul 2016 16:40
 Operator : UM/SJ
 Sample : H4045-16 5X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 H10-B6(0-6)C

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-BF063016.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	1.747	13	14	54	rVB	1773775	4212815	100.00%	23.406%
2	5.222	315	318	320	rBV	171900	228266	5.42%	1.268%
3	6.285	408	411	413	rBV	305959	260014	6.17%	1.445%
4	6.399	419	421	424	rBV	234364	251593	5.97%	1.398%
5	6.639	440	442	444	rVB	305470	268841	6.38%	1.494%
6	6.787	453	455	458	rBV	181438	187946	4.46%	1.044%
7	6.970	469	471	474	rBV	145613	134354	3.19%	0.746%
8	7.199	489	491	493	rBV	201622	163747	3.89%	0.910%
9	7.850	546	548	552	rBV	90482	76976	1.83%	0.428%
10	7.919	552	554	557	rBV	297194	365230	8.67%	2.029%
11	8.502	604	605	607	rBV	36965	43657	1.04%	0.243%
12	8.833	629	634	638	rBV2	36961	66311	1.57%	0.368%
13	9.005	647	649	651	rVB	270576	304099	7.22%	1.690%
14	9.096	653	657	660	rBV	47872	59875	1.42%	0.333%
15	9.393	681	683	686	rBV2	47626	69586	1.65%	0.387%
16	9.668	705	707	710	rVV	404522	393528	9.34%	2.186%
17	9.896	724	727	730	rBV	128389	172750	4.10%	0.960%
18	10.456	774	776	778	rVB	115222	122242	2.90%	0.679%
19	10.834	804	809	812	rBV	143257	193724	4.60%	1.076%
20	11.039	824	827	832	rVB	45476	68995	1.64%	0.383%
21	11.142	834	836	840	rVB	418900	374932	8.90%	2.083%
22	11.714	881	886	888	rBV	1096634	1701828	40.40%	9.455%
23	12.342	939	941	943	rBV	44077	55018	1.31%	0.306%
24	12.411	943	947	950	rBV2	35220	100742	2.39%	0.560%
25	12.514	950	956	958	rBV	1116027	1986393	47.15%	11.036%
26	12.571	959	961	964	rVB2	40792	82979	1.97%	0.461%
27	12.731	973	975	977	rBV2	169120	207438	4.92%	1.153%
28	13.302	1020	1025	1031	rVB3	51809	107031	2.54%	0.595%
29	13.588	1046	1050	1051	rBV	315376	472282	11.21%	2.624%
30	13.668	1054	1057	1059	rVV	1766937	2519607	59.81%	13.999%
31	13.702	1059	1060	1063	rVB	84332	88400	2.10%	0.491%
32	13.771	1063	1066	1068	rBV	243804	280385	6.66%	1.558%
33	13.965	1080	1083	1085	rBV	47372	66643	1.58%	0.370%
34	14.262	1107	1109	1115	rVB	69426	81507	1.93%	0.453%

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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	14.422	1120	1123	1125	rBV	900435	1098850	26.08%	6.105%
36	14.537	1130	1133	1136	rBV2	265334	386560	9.18%	2.148%
37	14.617	1139	1140	1144	rVB	55038	50349	1.20%	0.280%
38	14.845	1158	1160	1162	rVV	56566	65612	1.56%	0.365%
39	15.120	1182	1184	1187	rVV	196288	238879	5.67%	1.327%
40	15.177	1187	1189	1192	rVB	39534	48878	1.16%	0.272%
41	15.543	1219	1221	1224	rBV	42957	68330	1.62%	0.380%
42	15.611	1224	1227	1229	rVV	19204	42390	1.01%	0.236%
43	15.668	1229	1232	1235	rVV2	33651	67970	1.61%	0.378%
44	15.748	1235	1239	1244	rVV7	18858	56063	1.33%	0.311%
45	16.880	1334	1338	1343	rVB2	43961	105307	2.50%	0.585%

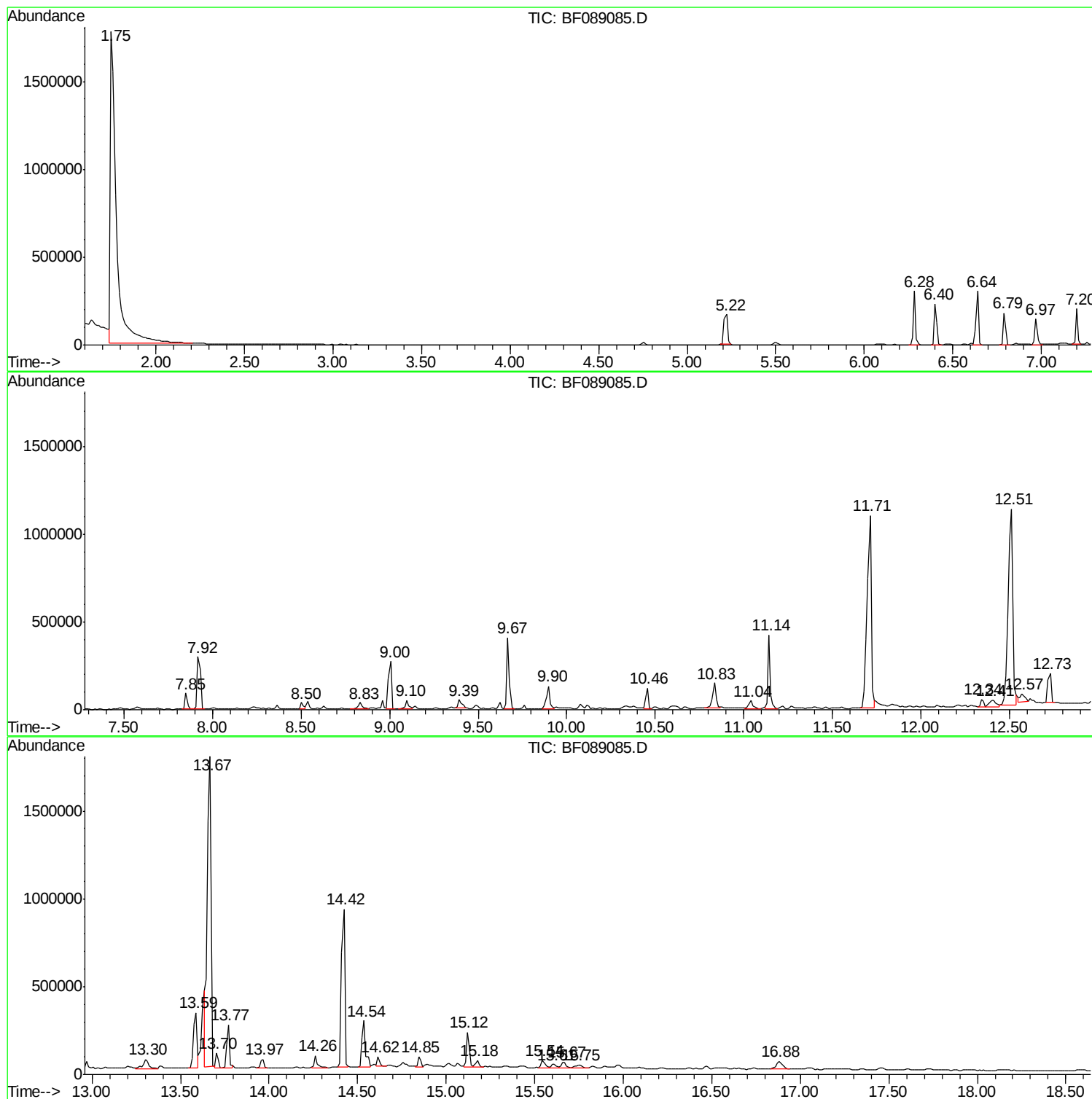
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF063016.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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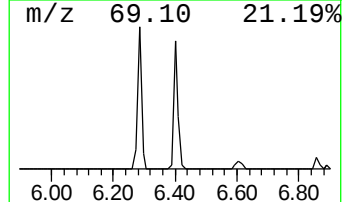
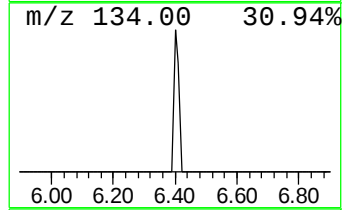
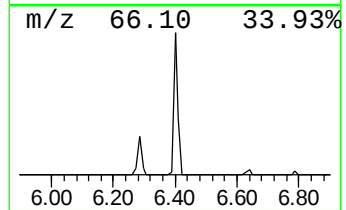
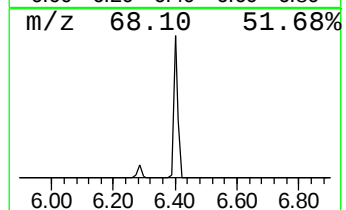
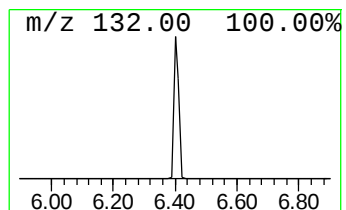
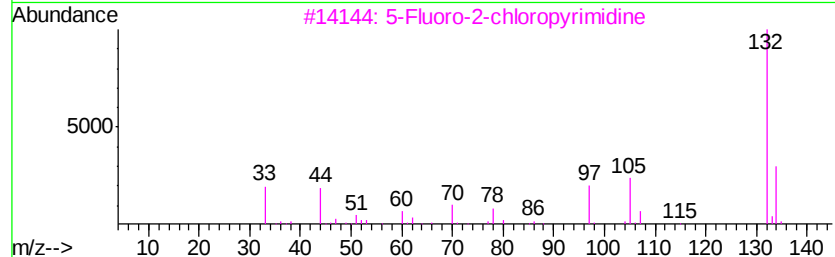
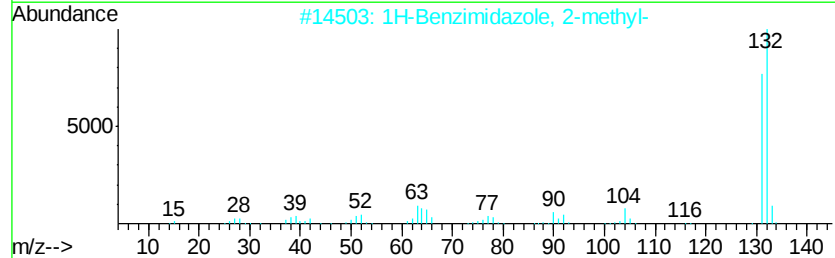
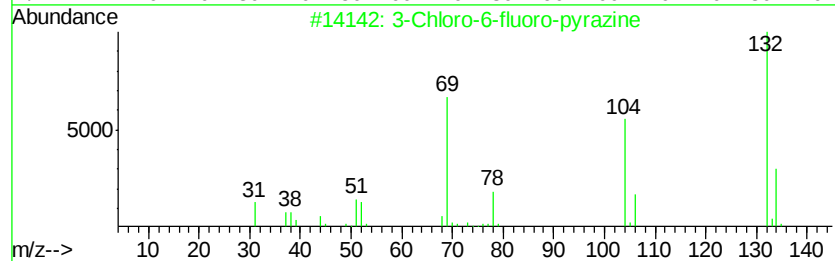
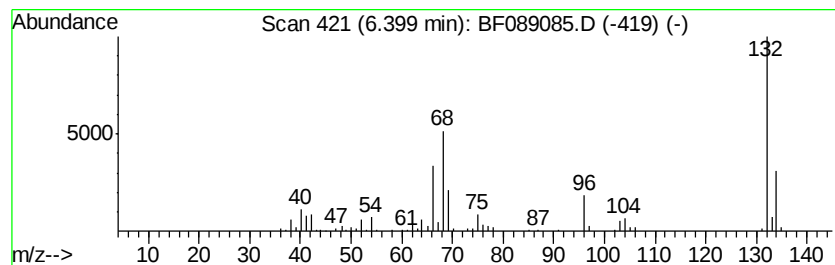
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 unknown6.40 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.40	18.72 ng	251593	1,4-Dichlorobenzene-d4	6.64

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2			1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
3			5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
4			(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	10
5			Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10



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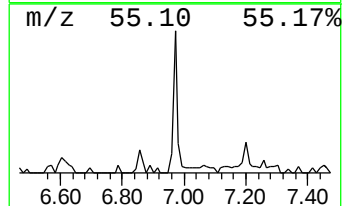
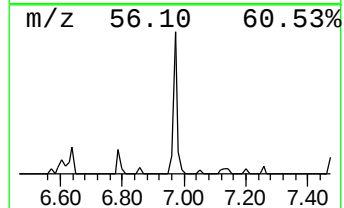
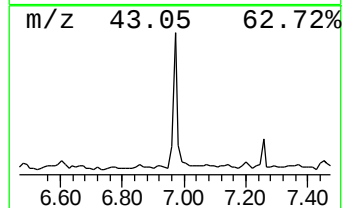
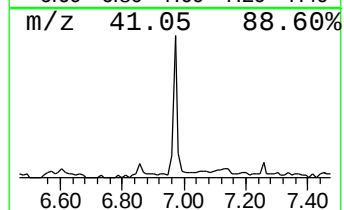
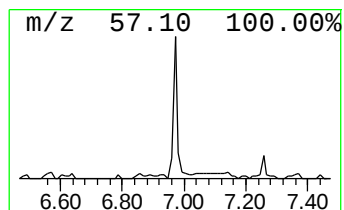
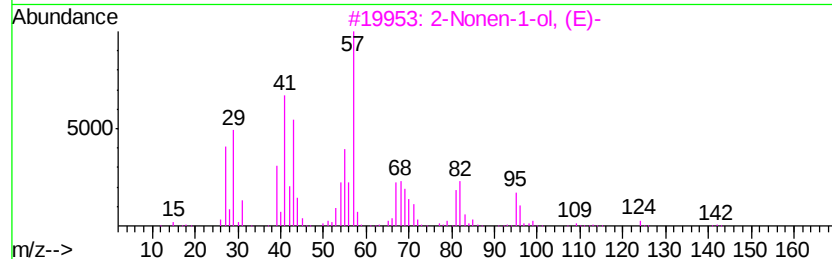
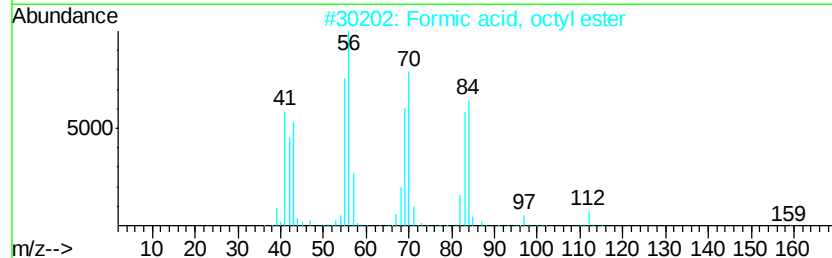
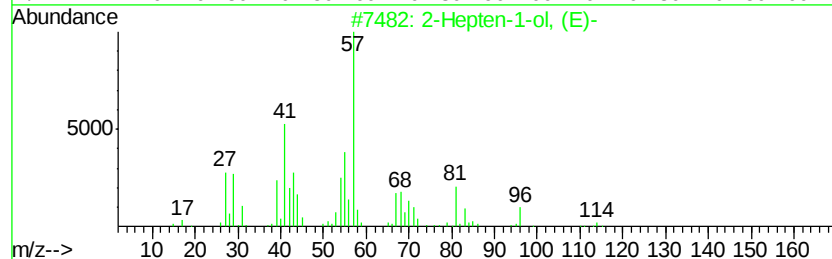
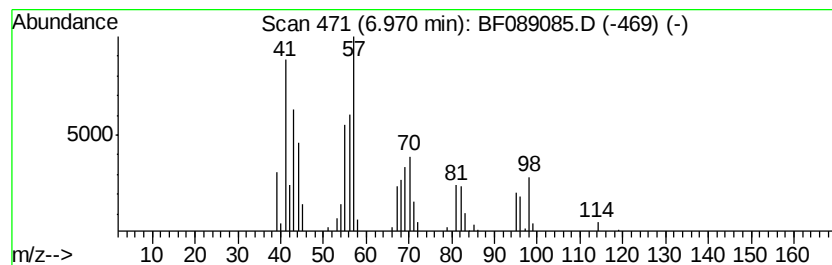
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Peak Number 4 unknown6.97 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.97	10.00 ng	134354	1,4-Dichlorobenzene-d4	6.64

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Hepten-1-ol, (E)-	114	C7H14O	033467-76-4	35
2		Formic acid, octyl ester	158	C9H18O2	000112-32-3	27
3		2-Nonen-1-ol, (E)-	142	C9H18O	031502-14-4	27
4		Heptane, 2-chloro-	134	C7H15Cl	001001-89-4	27
5		Cyclopentane, 1,2-dimethyl-	98	C7H14	002452-99-5	27



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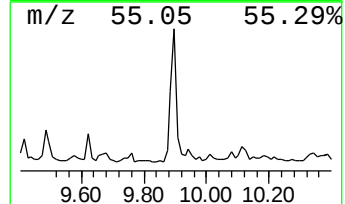
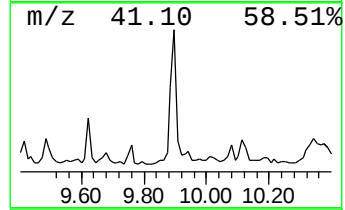
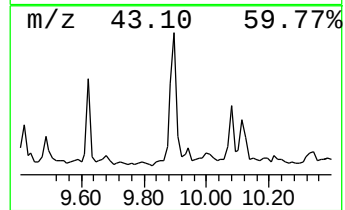
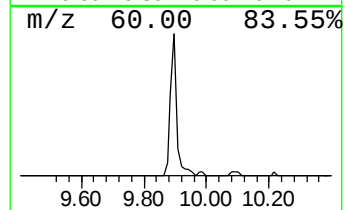
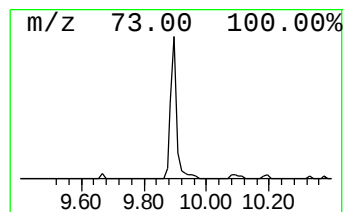
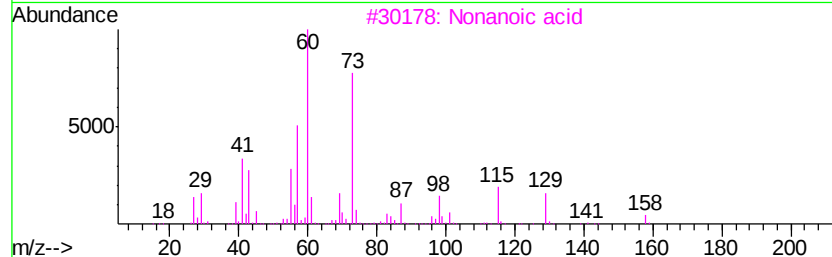
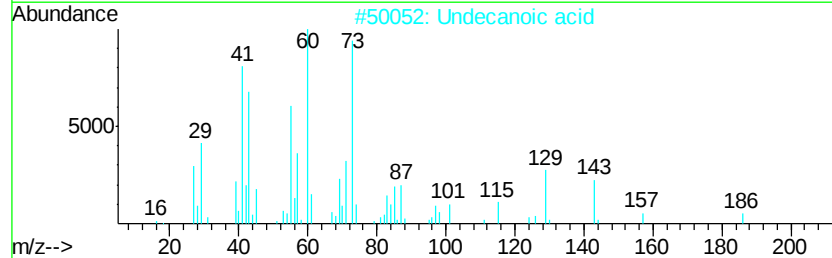
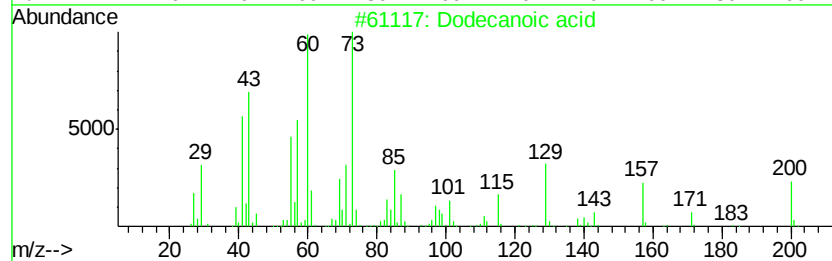
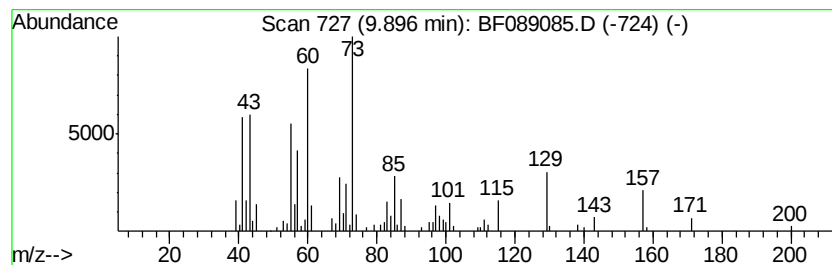
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Dodecanoic acid Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.90	8.78 ng	172750	Acenaphthene-d10	9.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dodecanoic acid	200	C12H24O2	000143-07-7	94
2		Undecanoic acid	186	C11H22O2	000112-37-8	80
3		Nonanoic acid	158	C9H18O2	000112-05-0	47
4		Ethanone, 1-(4,5-dihydro-2-thiaz...	129	C5H7NOS	029926-41-8	46
5		Hexanoic acid	116	C6H12O2	000142-62-1	43



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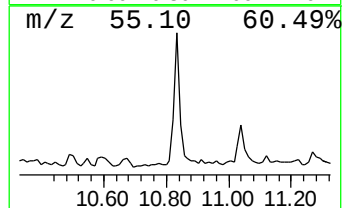
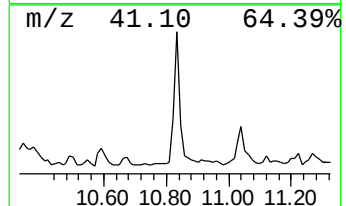
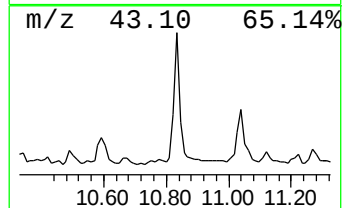
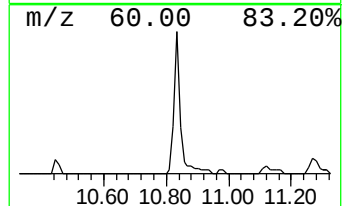
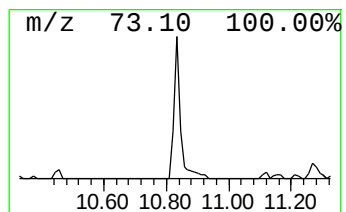
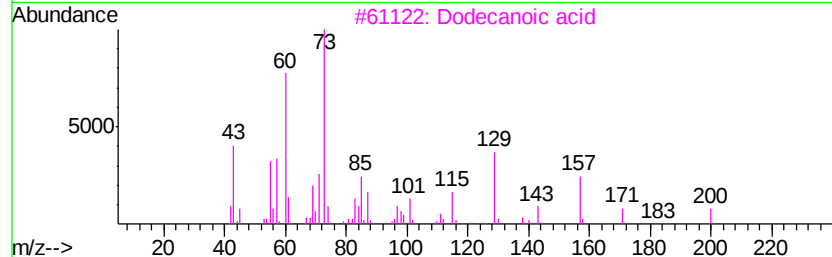
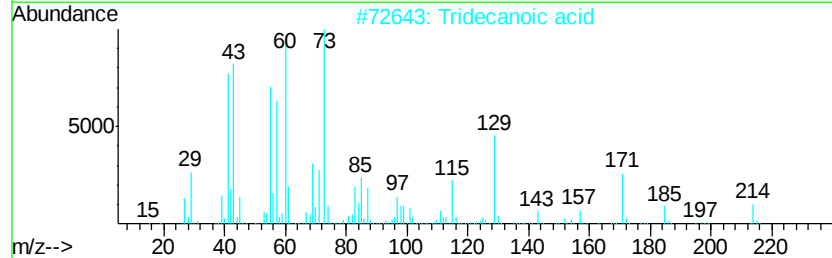
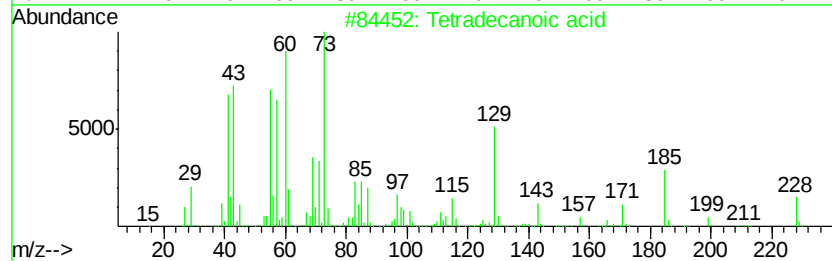
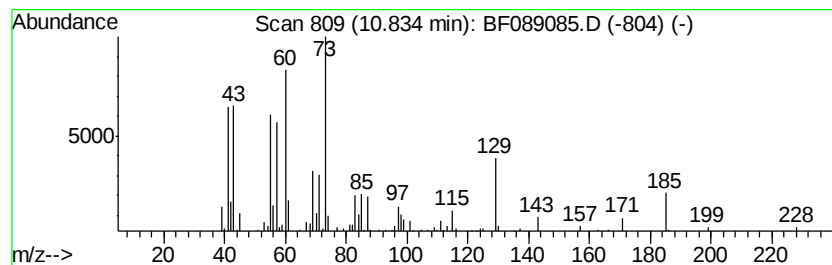
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TIC Library : C:\Database\NIST11.L
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Peak Number 6 Tetradecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD		R.T.	
10.83	10.33 ng	193724	Phenanthrene-d10		11.14	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetradecanoic acid	228	C14H28O2	000544-63-8	95
2		Tridecanoic acid	214	C13H26O2	000638-53-9	87
3		Dodecanoic acid	200	C12H24O2	000143-07-7	72
4		Decanoic acid, silver(1+) salt	278	C10H19AgO2	013126-67-5	72
5		n-PROPYL NONYL ETHER	186	C12H26O	1000130-69-3	25



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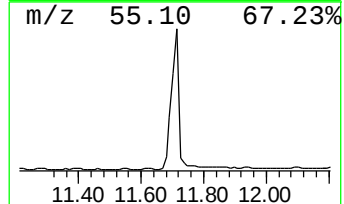
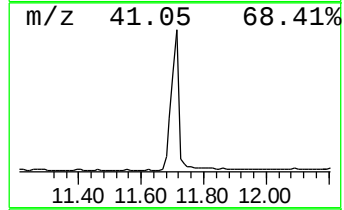
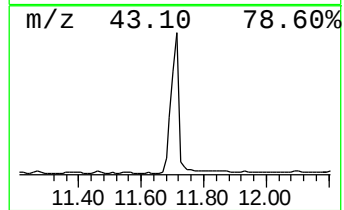
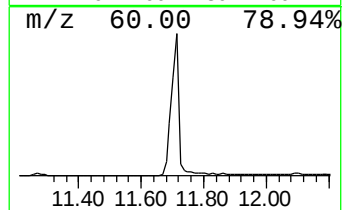
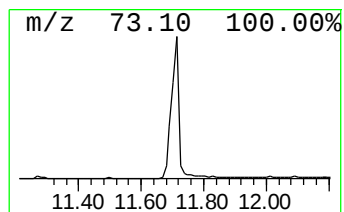
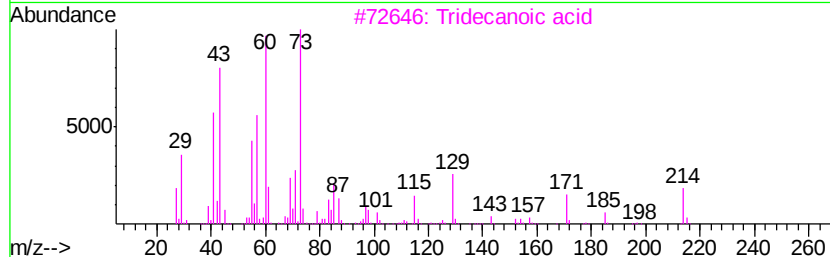
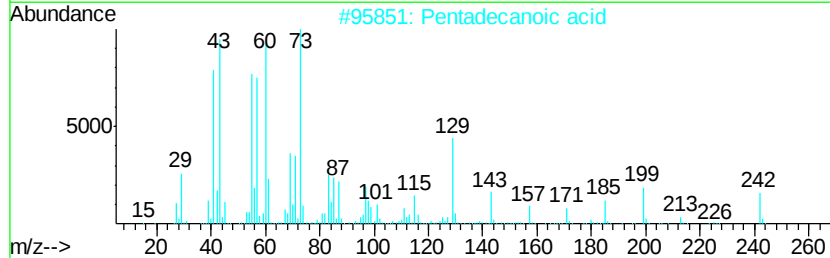
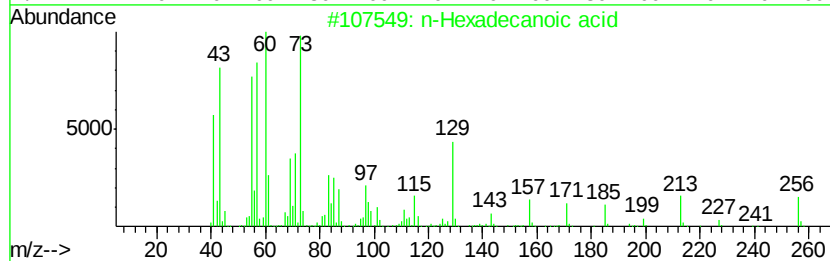
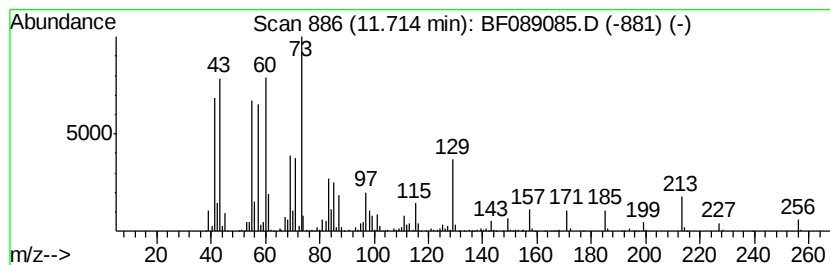
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.71	90.78 ng	1701830	Phenanthrene-d10	11.14

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	93
3		Tridecanoic acid	214	C13H26O2	000638-53-9	89
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	74
5		n-Decanoic acid	172	C10H20O2	000334-48-5	64



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Data File : BF089085.D
Acq On : 18 Jul 2016 16:40
Operator : UM/SJ
Sample : H4045-16 5X
Misc :
ALS Vial : 5 Sample Multiplier: 1

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ClientSampleId :
H10-B6(0-6)C

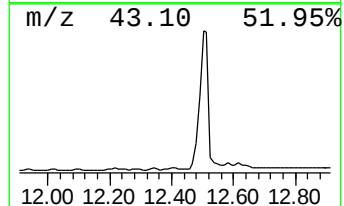
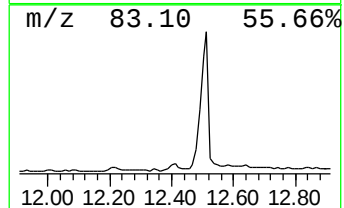
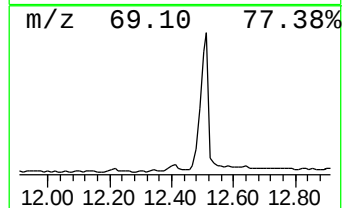
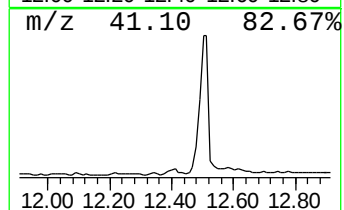
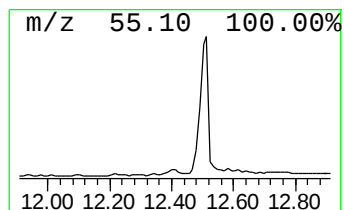
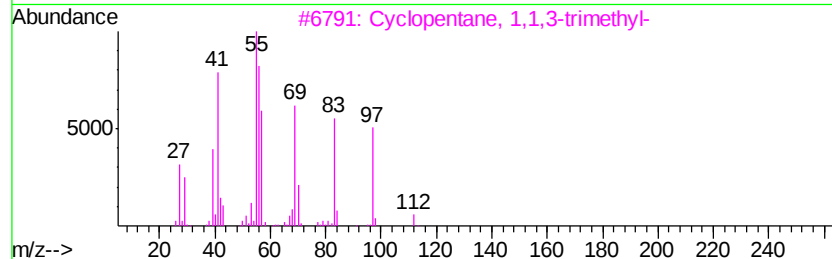
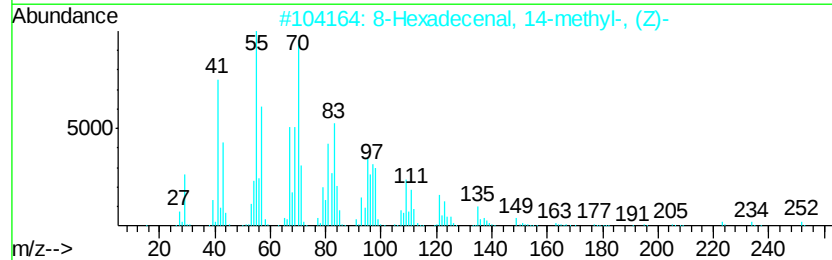
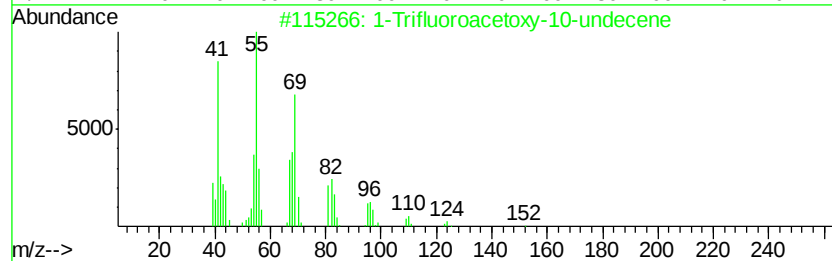
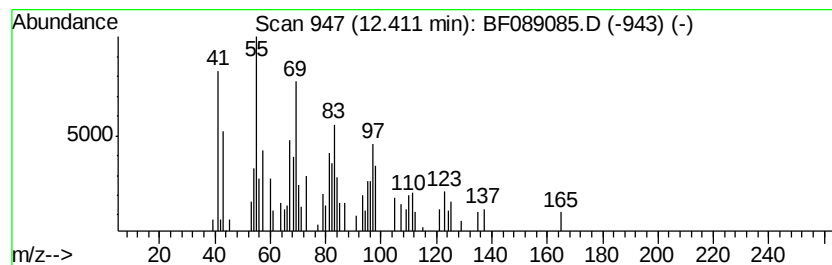
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 unknown12.41 Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.41	5.37 ng	100742	Phenanthrene-d10	11.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Trifluoroacetoxv-10-undecene	266	C13H21F3O2	001675-20-3	46
2		8-Hexadecenal, 14-methyl-, (Z)-	252	C17H32O	060609-53-2	46
3		Cyclopentane, 1,1,3-trimethyl-	112	C8H16	004516-69-2	46
4		Z-8-Methyl-9-tetradecenoic acid	240	C15H28O2	1000130-84-5	43
5		2-Hexenal, (E)-	98	C6H10O	006728-26-3	43



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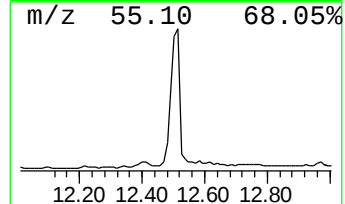
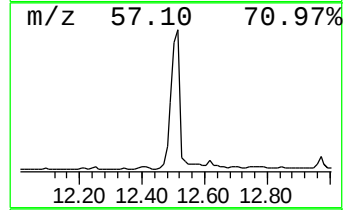
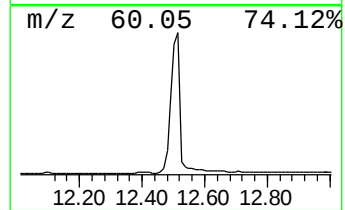
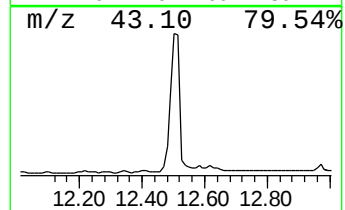
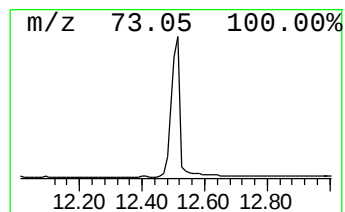
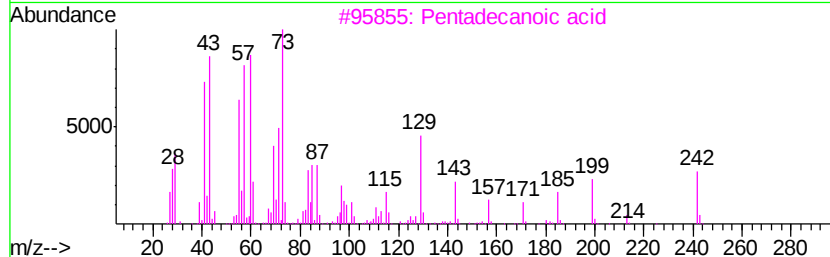
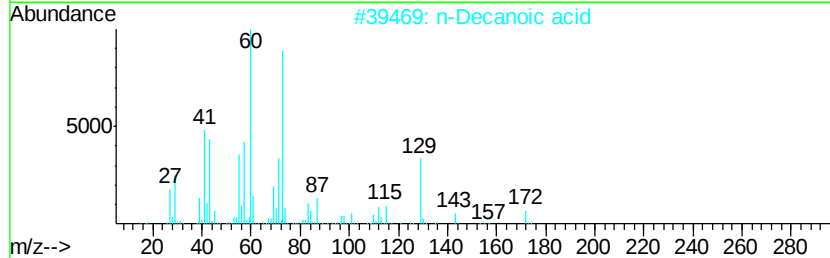
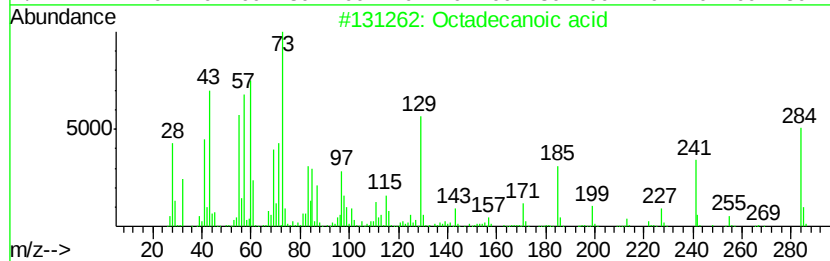
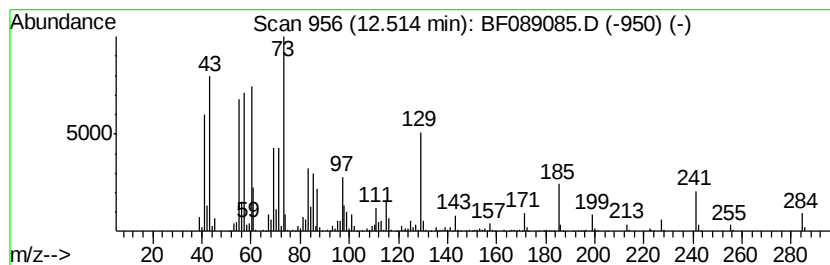
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TIC Library : C:\Database\NIST11.L
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Peak Number 9 Octadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.51	141.69 ng	1986390	Chrysene-d12	13.77

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2		n-Decanoic acid	172	C10H20O2	000334-48-5	70
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	58
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	41
5		Oxalic acid, isobutyl heptadecyl...	384	C23H44O4	1000309-38-2	18



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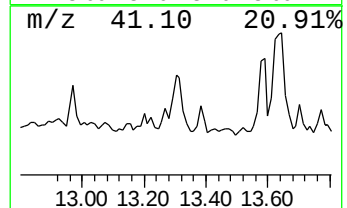
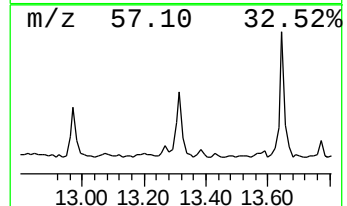
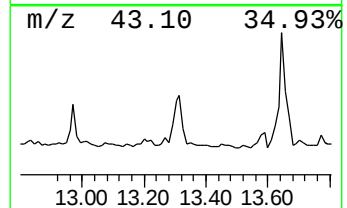
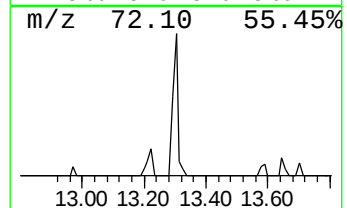
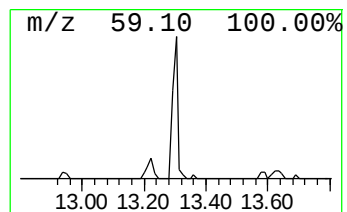
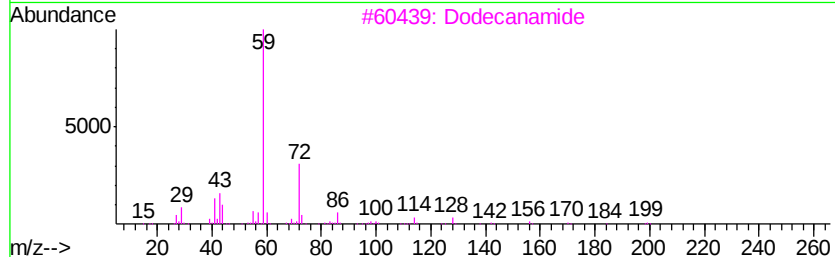
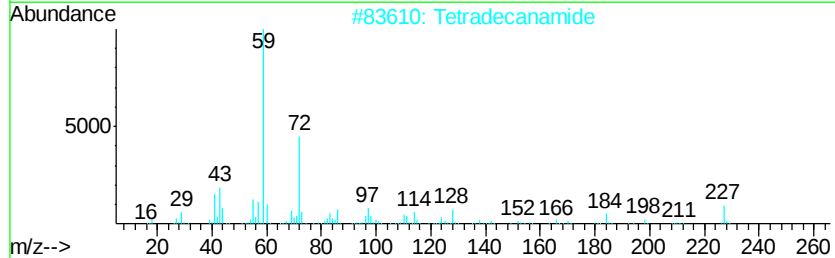
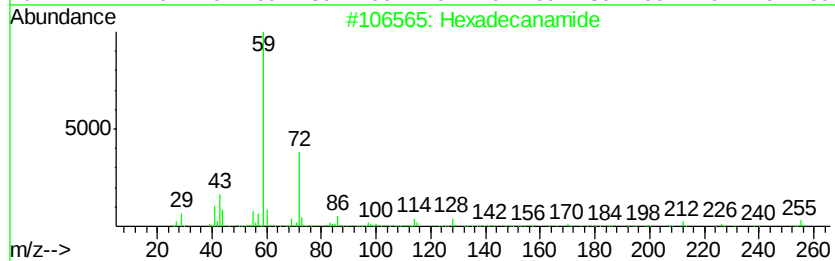
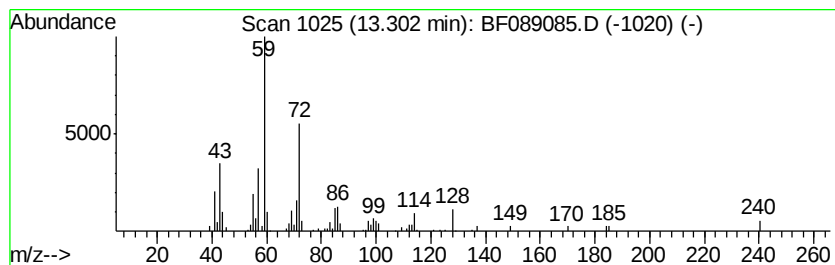
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TIC Library : C:\Database\NIST11.L
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Peak Number 11 Hexadecanamide Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.30	7.63 ng	107031	Chrysene-d12	13.77

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecanamide	255	C16H33NO	000629-54-9	78
2		Tetradecanamide	227	C14H29NO	000638-58-4	64
3		Dodecanamide	199	C12H25NO	001120-16-7	59
4		Decanamide-	171	C10H21NO	002319-29-1	56
5		Undecanamide, 11-bromo-	263	C11H22BrNO	005875-26-3	53



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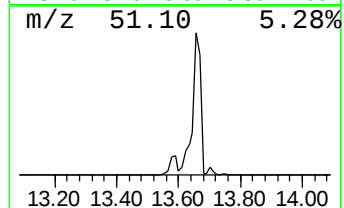
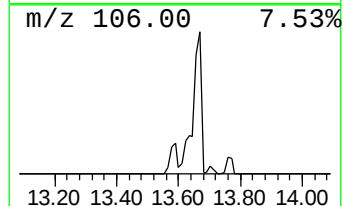
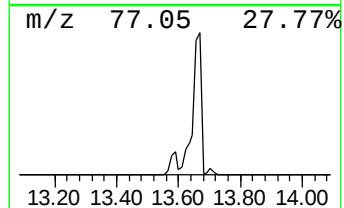
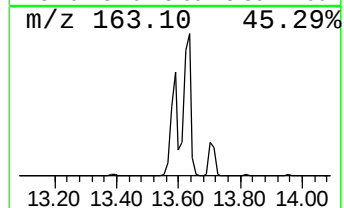
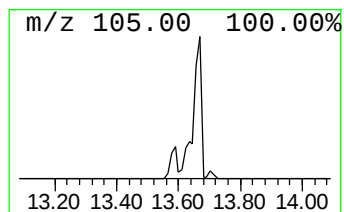
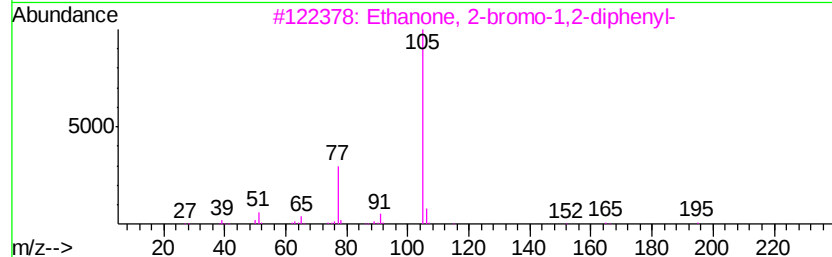
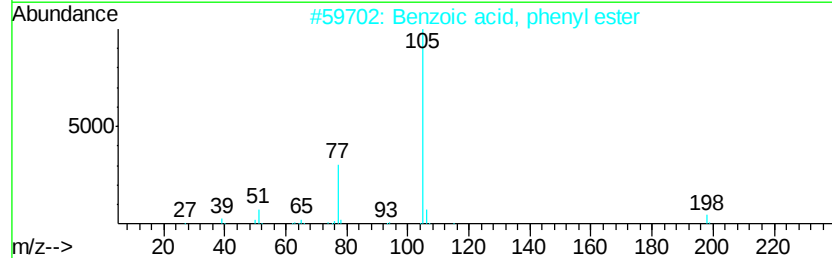
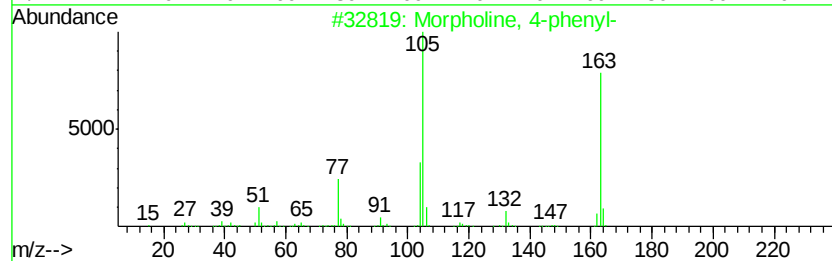
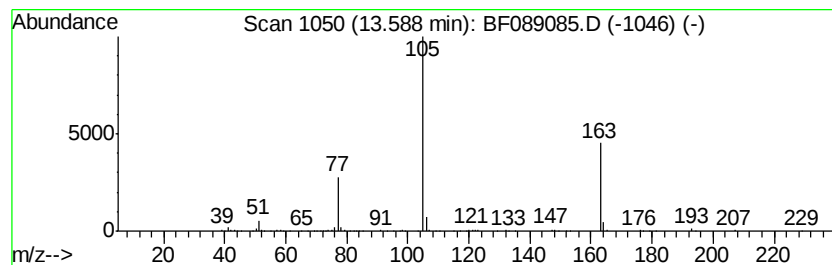
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Morpholine, 4-phenyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.59	33.69 ng	472282	Chrysene-d12	13.77

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Morpholine, 4-phenyl-	163	C10H13NO	000092-53-5	53
2		Benzoic acid, phenyl ester	198	C13H10O2	000093-99-2	38
3		Ethanone, 2-bromo-1,2-diphenyl-	274	C14H11BrO	001484-50-0	38
4		2,6-Dimethylphenyl benzoate	226	C15H14O2	001871-36-9	38
5		1,3-Benzenediol, monobenzoate	214	C13H10O3	000136-36-7	38



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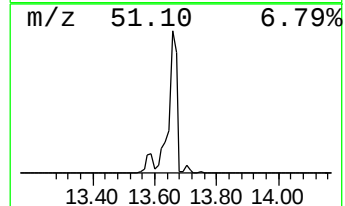
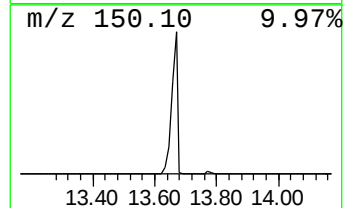
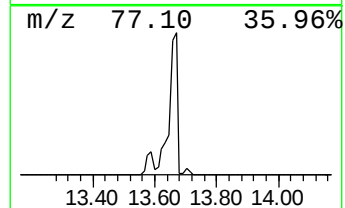
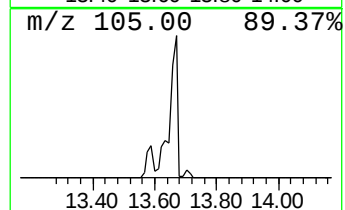
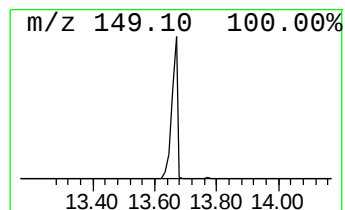
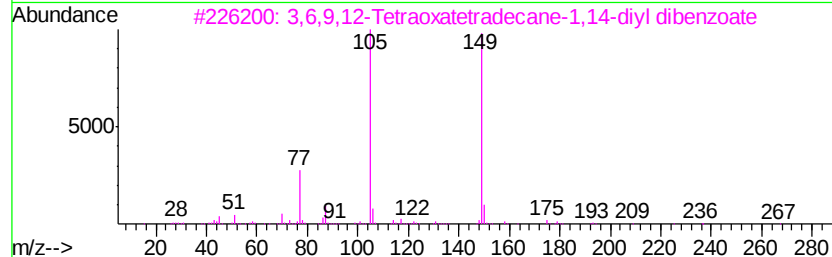
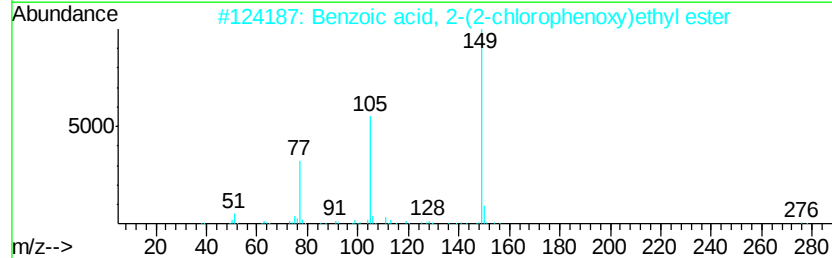
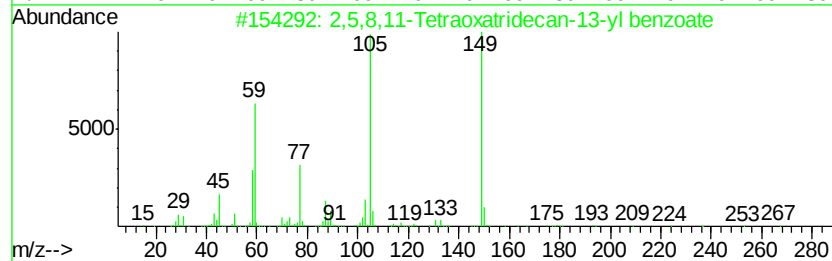
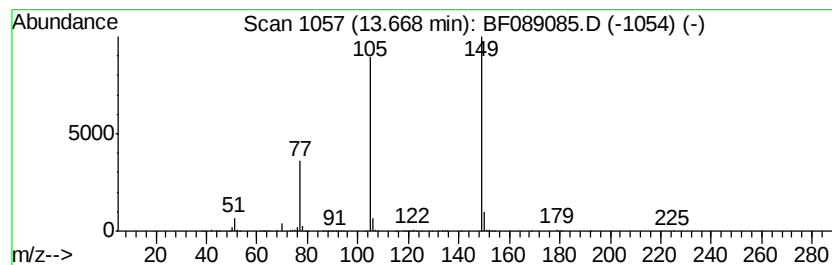
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 2,5,8,11-Tetraoxatridecan-1... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.67	179.73 ng	2519610	Chrysene-d12	13.77

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,5,8,11-Tetraoxatridecan-13-yl ...	312	C16H24O6	1000366-96-7	78
2		Benzoic acid, 2-(2-chlorophenoxy...	276	C15H13ClO3	1000368-25-9	72
3		3,6,9,12-Tetraoxatetradecane-1,1...	446	C24H30O8	1000366-97-0	64
4		1,3-Dioxolane, 2-(methoxymethyl)...	194	C11H14O3	1000156-71-2	64
5		2-(2-Carboxyvinyl)pyridine, trans	149	C8H7NO2	007340-22-9	58



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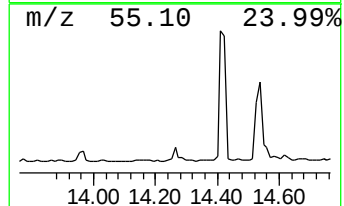
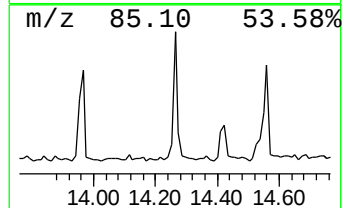
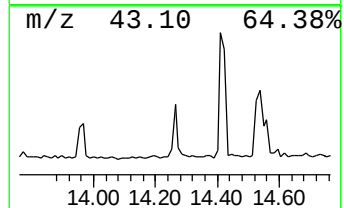
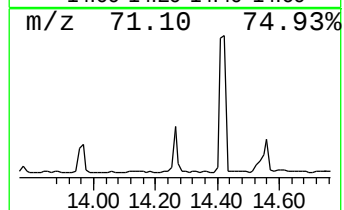
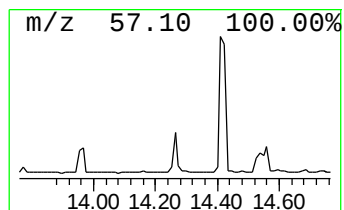
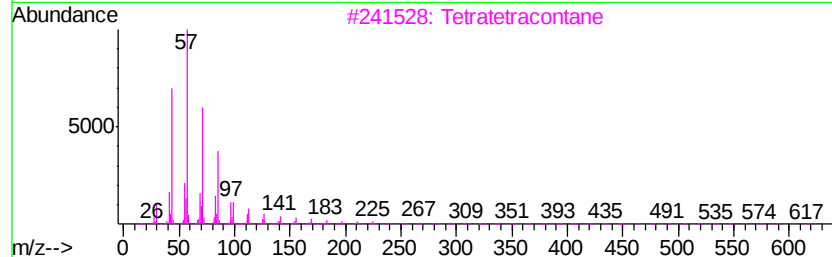
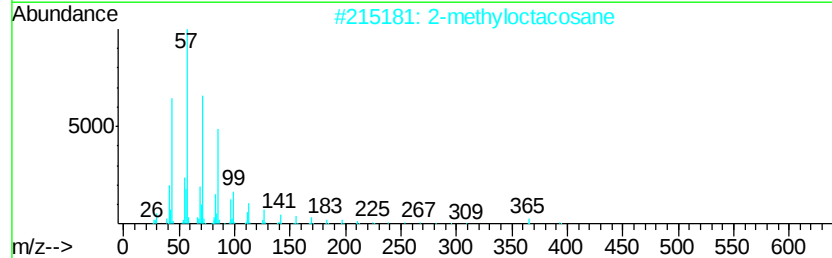
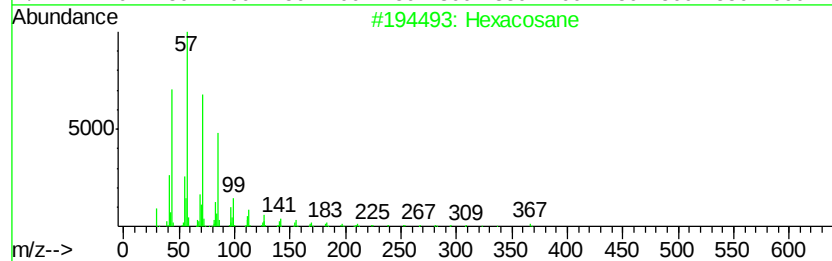
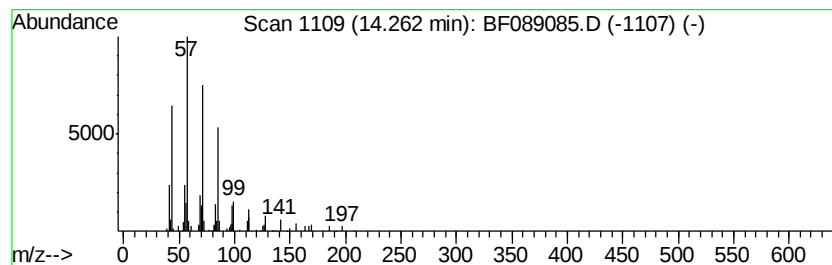
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Hexacosane Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.26	5.81 ng	81507	Chrysene-d12	13.77

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexacosane	366	C26H54	000630-01-3	91
2		2-methyloctacosane	408	C29H60	1000376-72-8	91
3		Tetratetracontane	619	C44H90	007098-22-8	91
4		triacontane	422	C30H62	000638-68-6	91
5		Heneicosane	296	C21H44	000629-94-7	91



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ALS Vial : 5 Sample Multiplier: 1

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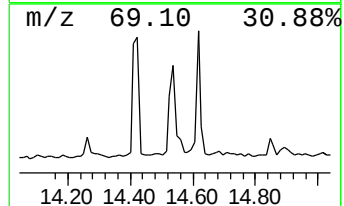
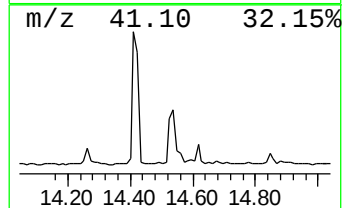
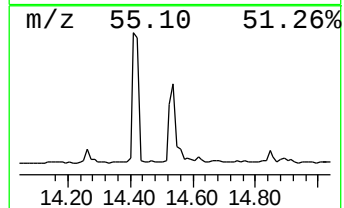
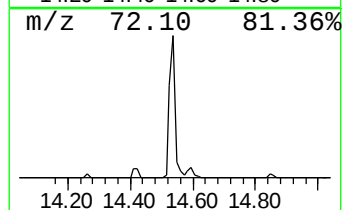
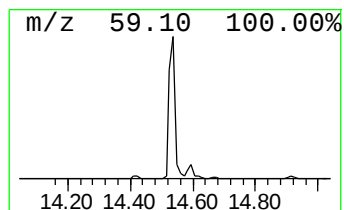
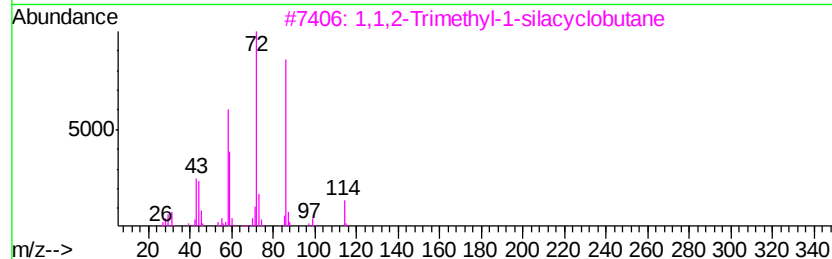
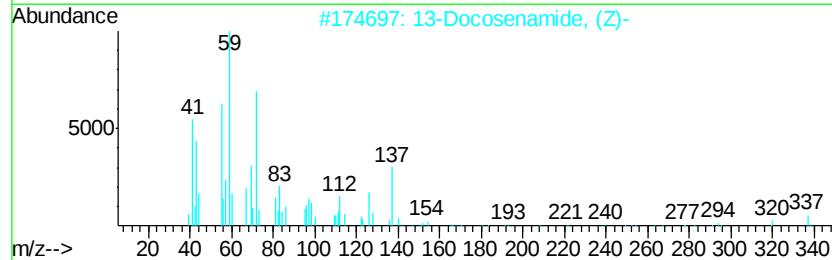
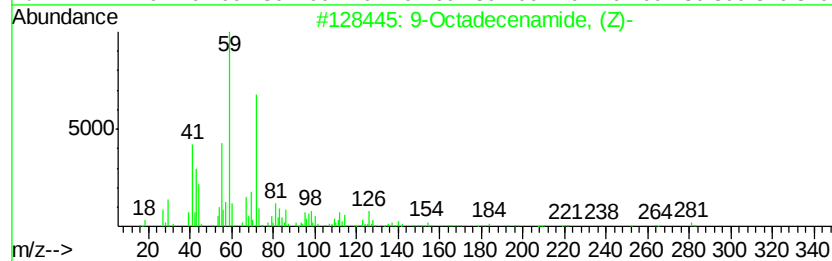
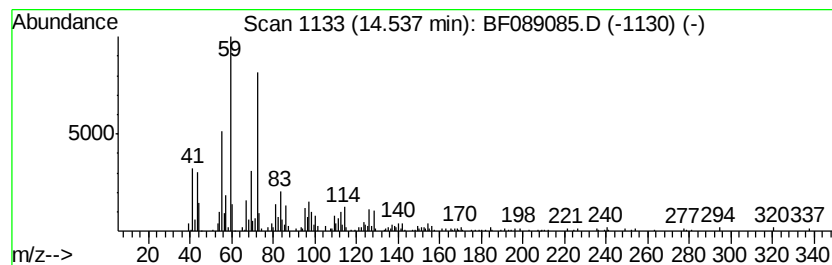
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 9-Octadecenamide, (Z)- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.54	32.36 ng	386560	Perylene-d12	15.12

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	99
2		13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	43
3		1,1,2-Trimethyl-1-silacyclobutane	114	C6H14Si	030681-90-4	38
4		Heptanamide, 4-ethyl-5-methyl-	171	C10H21NO	054789-40-1	38
5		Dodecanamide	199	C12H25NO	001120-16-7	35



Data Path : Z:\HPCHEM1\BNA F\DATA\BF071816\
Data File : BF089085.D
Acq On : 18 Jul 2016 16:40
Operator : UM/SJ
Sample : H4045-16 5X
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
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ClientSampleId :
H10-B6(0-6)C

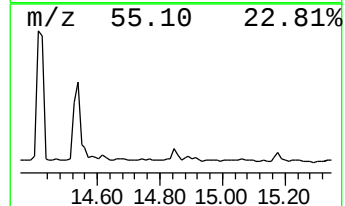
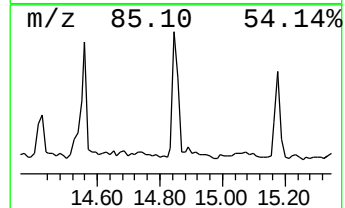
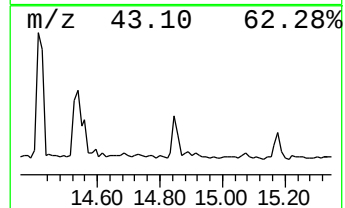
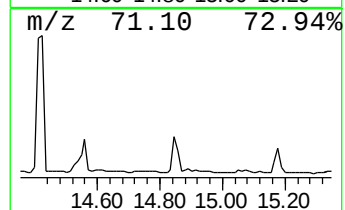
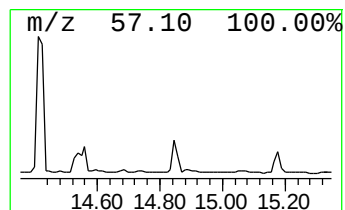
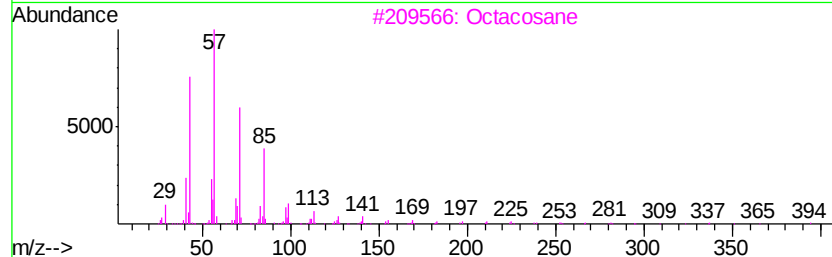
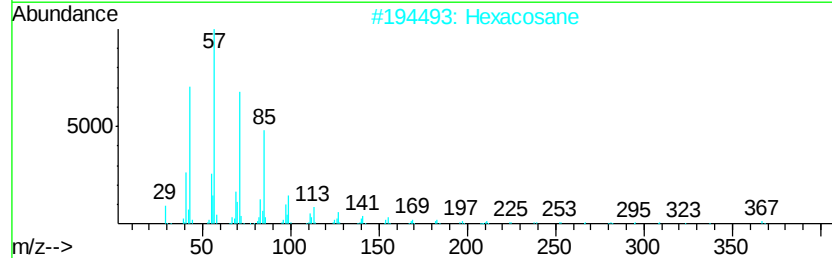
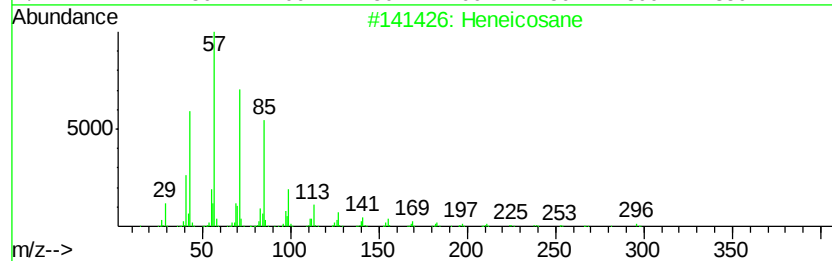
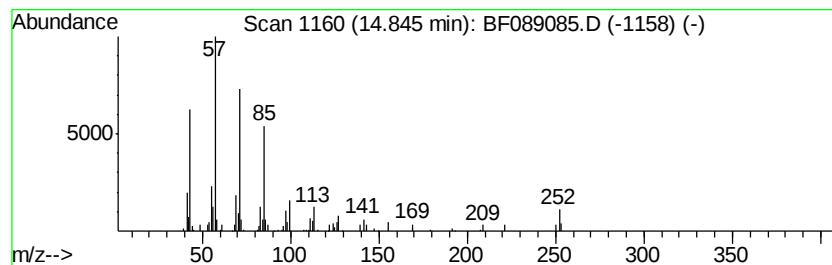
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Heneicosane Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.85	5.49 ng	65612	Perylene-d12	15.12

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	87
2			Hexacosane	366	C26H54	000630-01-3	87
3			Octacosane	394	C28H58	000630-02-4	83
4			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	83
5			Triacontane	422	C30H62	000638-68-6	83



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF071816\
Data File : BF089085.D
Acq On : 18 Jul 2016 16:40
Operator : UM/SJ
Sample : H4045-16 5X
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
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ClientSampleId :
H10-B6(0-6)C

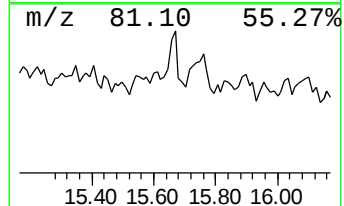
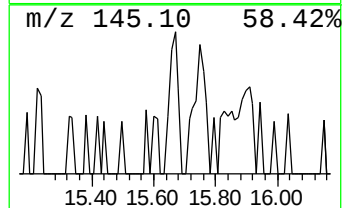
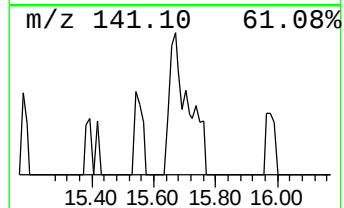
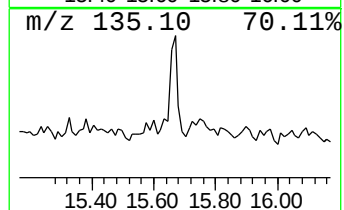
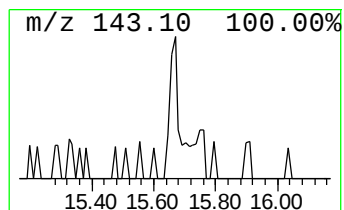
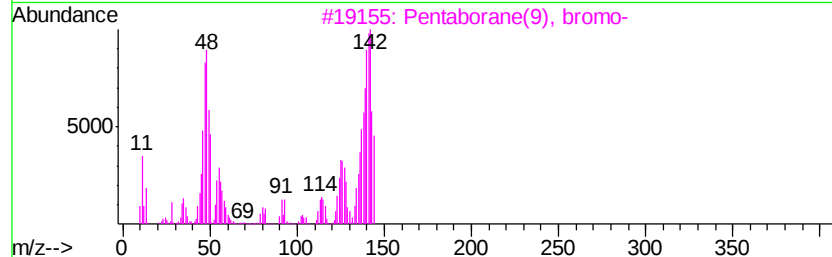
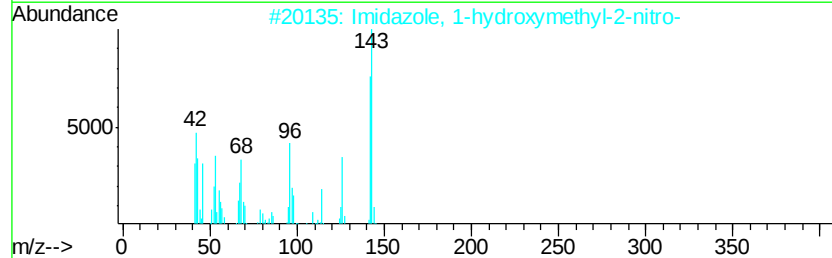
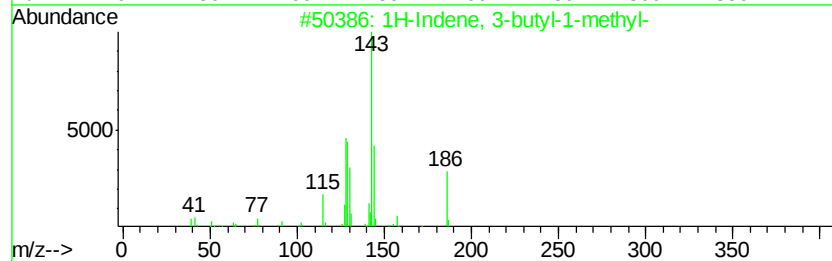
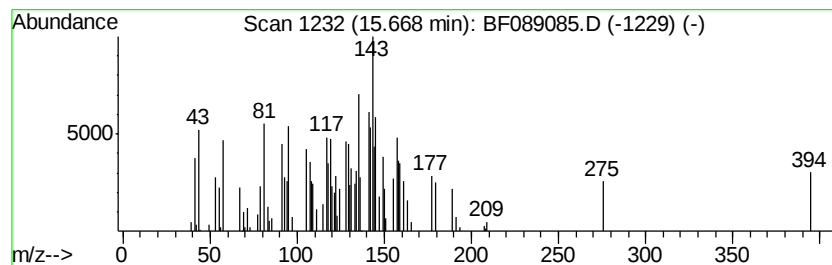
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 unknown15.67 Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.67	5.69 ng	67970	Perylene-d12	15.12

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 3-butyl-1-methyl-	186	C14H18	111400-84-1	22
2		Imidazole, 1-hydroxymethyl-2-nitro-	143	C4H5N3O3	097987-87-6	18
3		Pentaborane(9), bromo-	142	B5BrH8	066794-48-7	18
4		Propanedioic acid, (hydroxvimino...	189	C7H11NO5	006829-41-0	14
5		1,4-Bis(2-hydroxyethyl)piperazine	174	C8H18N2O2	000122-96-3	11



Data Path : Z:\HPCHEM1\BNA F\DATA\BF071816\
Data File : BF089085.D
Acq On : 18 Jul 2016 16:40
Operator : UM/SJ
Sample : H4045-16 5X
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
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ClientSampleId :
H10-B6(0-6)C

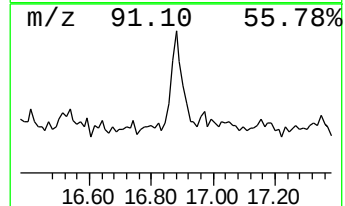
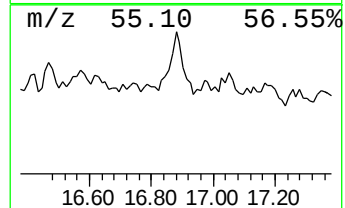
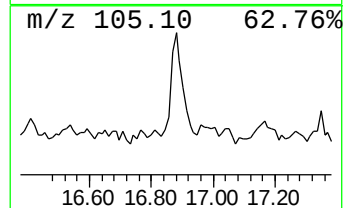
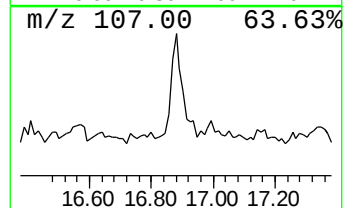
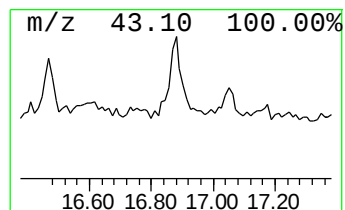
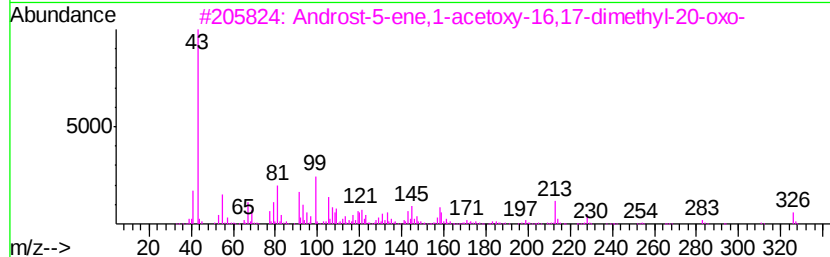
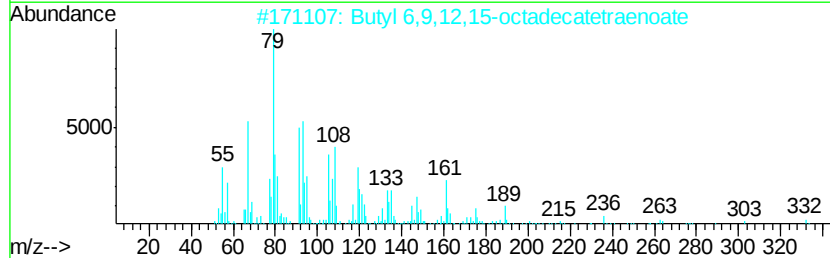
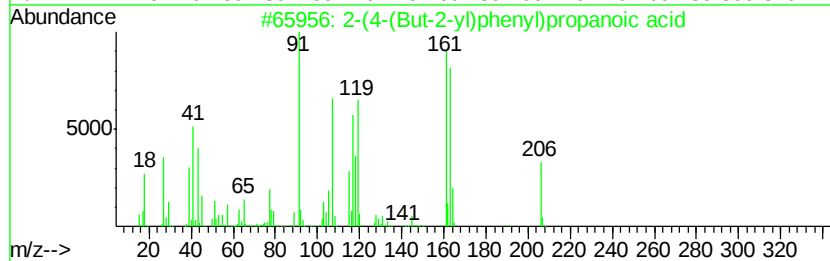
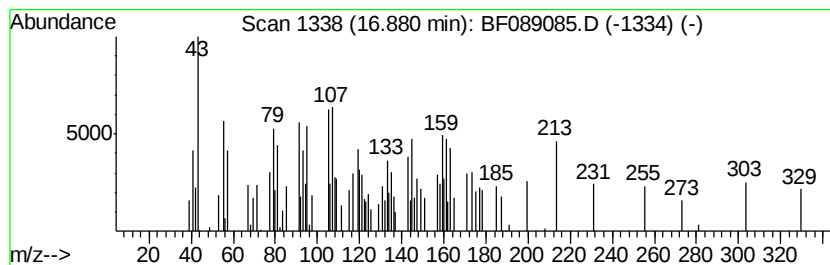
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 unknown16.88 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.88	8.82 ng	105307	Perylene-d12	15.12

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-(4-(But-2-yl)phenyl)propanoic ...	206	C13H18O2	064451-76-9	27		
2	Butyl 6,9,12,15-octadecatetraenoate	332	C22H36O2	1000336-71-3	14		
3	Androst-5-ene,1-acetoxv-16,17-di...	386	C25H38O3	294866-91-4	14		
4	Benzeneacetamide, .alpha.-ethyl-	163	C10H13NO	000090-26-6	10		
5	1H-Cyclopenta[a]phenanthrene-16-...	398	C24H34N2O3	1000337-25-3	10		



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF071816\
 Data File : BF089085.D
 Acq On : 18 Jul 2016 16:40
 Operator : UM/SJ
 Sample : H4045-16 5X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF063016.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
unknown6.40	6.40	18.7	ng	251593	1	6.64	268841	20.0
unknown6.97	6.97	10.0	ng	134354	1	6.64	268841	20.0
Dodecanoic acid	9.90	8.8	ng	172750	3	9.67	393528	20.0
Tetradecanoic acid	10.83	10.3	ng	193724	4	11.14	374932	20.0
n-Hexadecanoic acid	11.71	90.8	ng	1701830	4	11.14	374932	20.0
unknown12.41	12.41	5.4	ng	100742	4	11.14	374932	20.0
Octadecanoic acid	12.51	141.7	ng	1986390	5	13.77	280385	20.0
Hexadecanamide	13.30	7.6	ng	107031	5	13.77	280385	20.0
Morpholine, 4-phe...	13.59	33.7	ng	472282	5	13.77	280385	20.0
2,5,8,11-Tetraoxa...	13.67	179.7	ng	2519610	5	13.77	280385	20.0
Hexacosane	14.26	5.8	ng	81507	5	13.77	280385	20.0
9-Octadecenamide, ...	14.54	32.4	ng	386560	6	15.12	238879	20.0
Heneicosane	14.85	5.5	ng	65612	6	15.12	238879	20.0
unknown15.67	15.67	5.7	ng	67970	6	15.12	238879	20.0
unknown16.88	16.88	8.8	ng	105307	6	15.12	238879	20.0