

Data Path : Z:\HPCHEM1\BNA F\DATA\BF071416\
 Data File : BF088991.D
 Acq On : 14 Jul 2016 17:46
 Operator : UM/SJ
 Sample : PB92031BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB92031BL

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	1.656	4	6	15	rVB	315517	453229	14.48%	1.577%
2	1.781	15	17	27	rVV	649787	1043092	33.34%	3.628%
3	1.907	27	28	47	rVB	38279	111344	3.56%	0.387%
4	4.844	282	285	290	rVB	141242	207561	6.63%	0.722%
5	5.290	321	324	326	rBV	1335330	1703494	54.44%	5.926%
6	6.342	412	416	418	rBV	1457249	1824671	58.31%	6.347%
7	6.456	423	426	428	rBV	1763220	1660921	53.08%	5.777%
8	6.685	444	446	448	rBV	431132	361355	11.55%	1.257%
9	6.845	457	460	462	rBV	1355970	1403905	44.87%	4.883%
10	7.256	492	496	498	rBV	943760	1170205	37.40%	4.071%
11	7.965	556	558	561	rVB	401532	437353	13.98%	1.521%
12	9.051	649	653	655	rBV	2188459	2173947	69.48%	7.562%
13	9.656	703	706	709	rBV	53567	72232	2.31%	0.251%
14	9.713	709	711	714	rVB	578928	533479	17.05%	1.856%
15	9.931	727	730	736	rVB	46816	85858	2.74%	0.299%
16	10.502	777	780	782	rBV	953182	871048	27.84%	3.030%
17	10.765	802	803	808	rBV4	21999	51142	1.63%	0.178%
18	10.845	808	810	811	rVV2	23675	42430	1.36%	0.148%
19	10.868	811	812	818	rVB	54412	89884	2.87%	0.313%
20	11.188	838	840	843	rVB	642861	547194	17.49%	1.903%
21	11.599	873	876	880	rBV	27636	39639	1.27%	0.138%
22	11.748	882	889	897	rBV2	1280528	2688512	85.92%	9.352%
23	11.862	897	899	906	rVB3	38018	88862	2.84%	0.309%
24	11.954	906	907	910	rBV	28245	33579	1.07%	0.117%
25	12.388	942	945	947	rBV	79242	146501	4.68%	0.510%
26	12.434	947	949	953	rVV3	36076	85159	2.72%	0.296%
27	12.537	953	958	962	rVV	1793581	3129055	100.00%	10.884%
28	12.605	962	964	967	rVV	87703	200976	6.42%	0.699%
29	12.674	969	970	974	rVV2	63671	82206	2.63%	0.286%
30	12.777	976	979	982	rVB	1647810	1840871	58.83%	6.403%
31	13.005	997	999	1000	rBV	35811	36949	1.18%	0.129%
32	13.234	1017	1019	1023	rBV2	24500	45081	1.44%	0.157%
33	13.325	1023	1027	1033	rVB3	123486	263052	8.41%	0.915%
34	13.611	1049	1052	1054	rBV	234200	286889	9.17%	0.998%

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35	13.657	1054	1056	1057	rVV	258108	418911	13.39%	1.457%
36	13.680	1057	1058	1061	rVV	1044488	1448769	46.30%	5.039%
37	13.737	1061	1063	1065	rVB	58350	60010	1.92%	0.209%
38	13.817	1066	1070	1072	rVV2	410108	551385	17.62%	1.918%
39	13.988	1083	1085	1090	rVB2	64830	72056	2.30%	0.251%
40	14.297	1110	1112	1114	rBV	49225	68628	2.19%	0.239%
41	14.445	1123	1125	1127	rBV	184107	187723	6.00%	0.653%
42	14.560	1133	1135	1136	rBV	235709	235110	7.51%	0.818%
43	14.651	1141	1143	1146	rVV	427977	588050	18.79%	2.046%
44	14.880	1161	1163	1166	rBV	46107	62033	1.98%	0.216%
45	15.063	1175	1179	1181	rBV	32410	60598	1.94%	0.211%
46	15.177	1186	1189	1191	rVV	260217	293670	9.39%	1.022%
47	15.211	1191	1192	1196	rVB	31985	40103	1.28%	0.139%
48	15.485	1212	1216	1221	rVB	32069	83178	2.66%	0.289%
49	15.588	1222	1225	1228	rBV	23688	41212	1.32%	0.143%
50	15.703	1233	1235	1240	rVV4	24283	61310	1.96%	0.213%
51	15.806	1240	1244	1248	rVV5	13407	45063	1.44%	0.157%
52	15.874	1248	1250	1254	rVB3	16851	34893	1.12%	0.121%
53	15.988	1254	1260	1270	rBV2	31909	124370	3.97%	0.433%
54	16.606	1310	1314	1320	rVB	29933	88824	2.84%	0.309%
55	16.960	1339	1345	1352	rBV7	19792	85109	2.72%	0.296%
56	17.360	1375	1380	1390	rVB	31329	119776	3.83%	0.417%
57	17.531	1392	1395	1401	rVB5	9663	33552	1.07%	0.117%
58	18.286	1456	1461	1467	rVB2	22996	79513	2.54%	0.277%
59	19.429	1555	1561	1566	rBV5	13369	52809	1.69%	0.184%

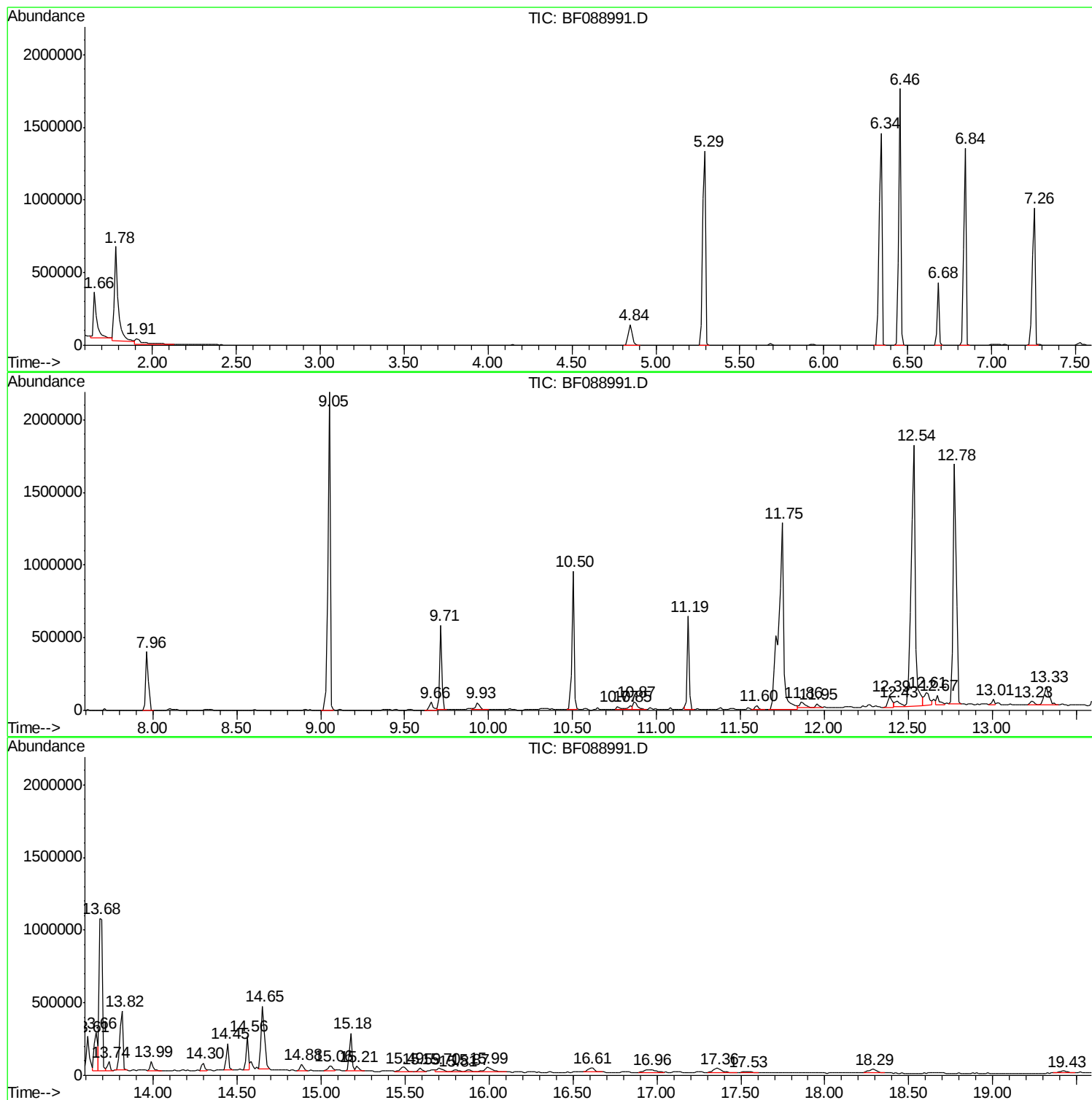
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF063016.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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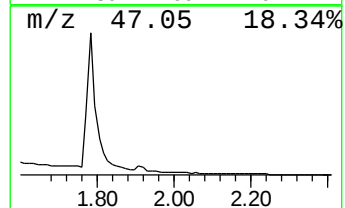
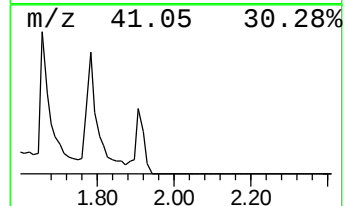
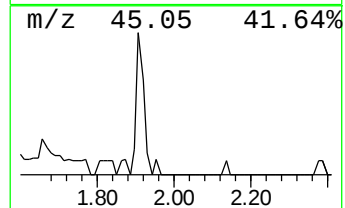
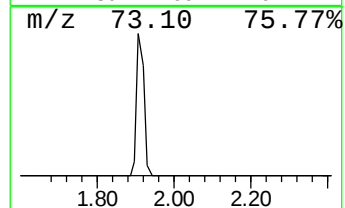
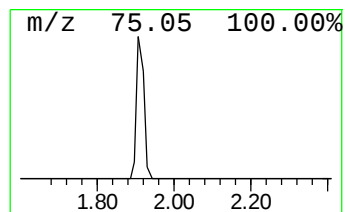
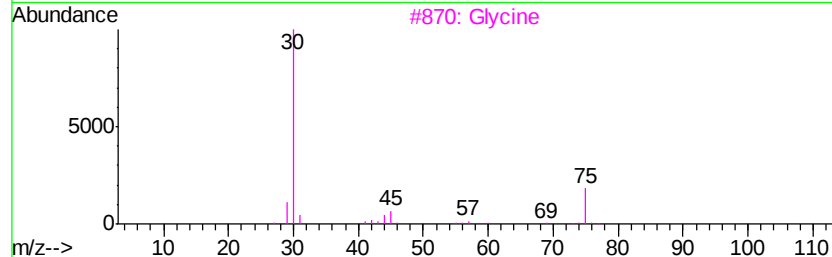
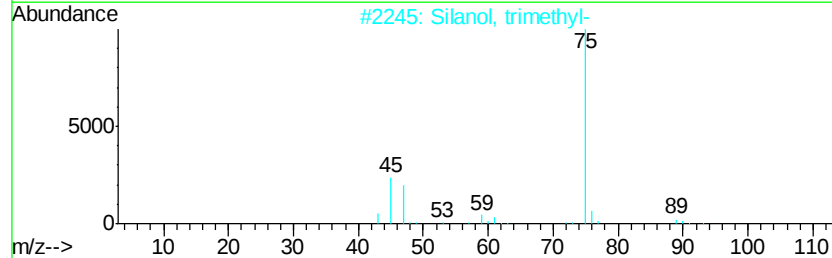
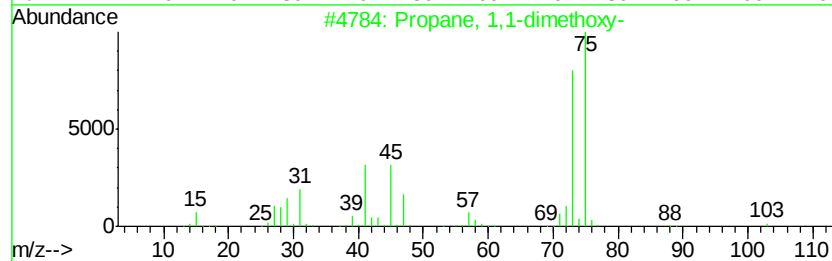
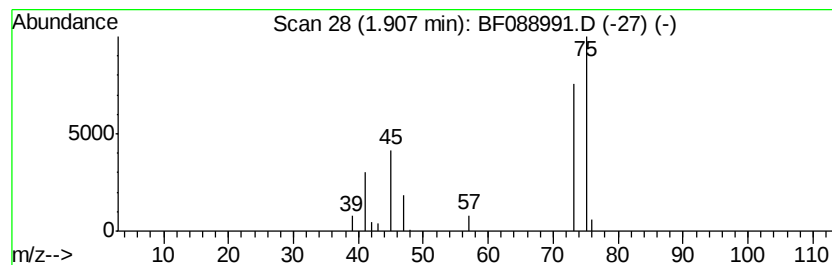
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF063016.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 Propane, 1,1-dimethoxy- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.91	6.16 ng	111344	1,4-Dichlorobenzene-d4	6.68

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propane, 1,1-dimethoxy-	104	C5H12O2	004744-10-9	83
2			Silanol, trimethyl-	90	C3H10OSi	001066-40-6	33
3			Glycine	75	C2H5NO2	000056-40-6	5
4			2-Bromomethyl-1,3-dioxolane	166	C4H7BrO2	004360-63-8	4
5			1,3-Dioxolane, 2-(chloromethyl)-	122	C4H7ClO2	002568-30-1	4



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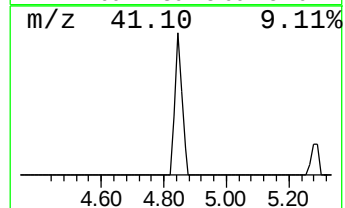
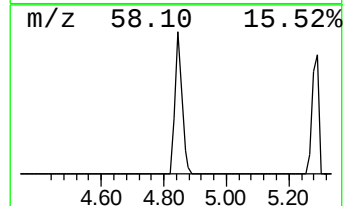
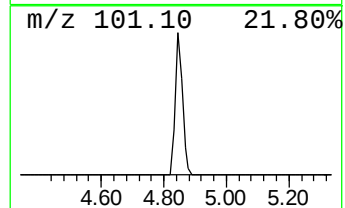
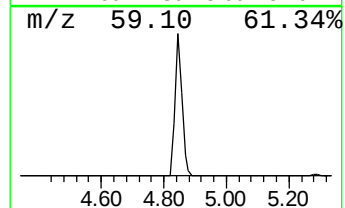
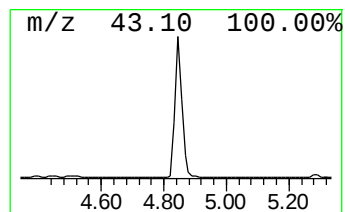
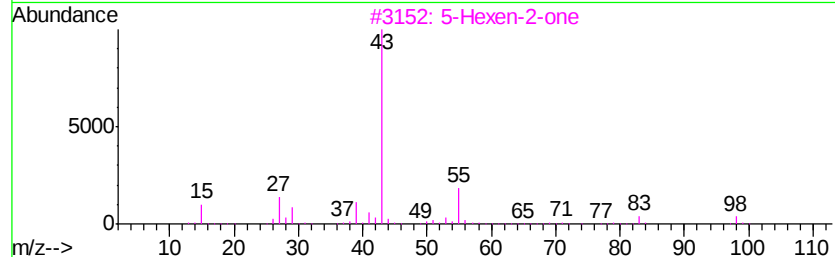
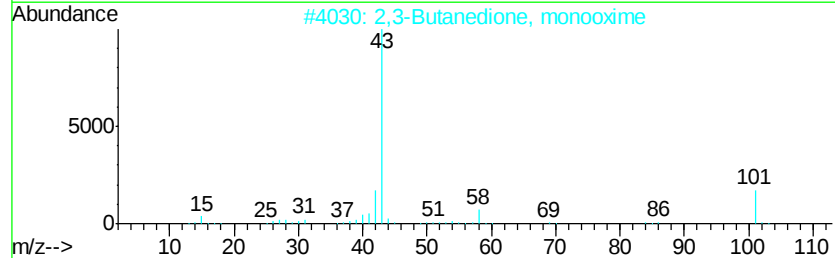
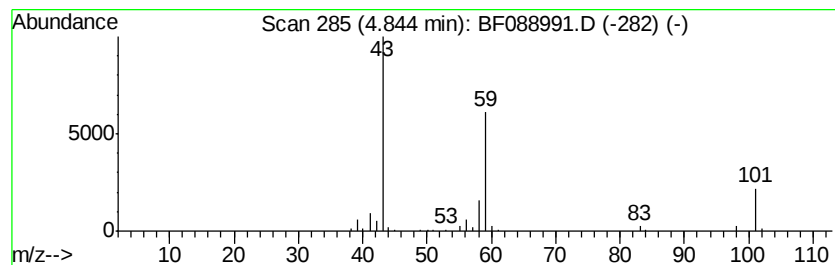
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.84	11.49 ng	207561	1,4-Dichlorobenzene-d4	6.68

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone,	4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56	
2	2,3-Butanedione,	monooxime	101	C4H7NO2	000057-71-6	25	
3	5-Hexen-2-one		98	C6H10O	000109-49-9	9	
4	Morpholine,	4-methyl-	101	C5H11NO	000109-02-4	9	
5	Butane,	1-ethoxy-	102	C6H14O	000628-81-9	9	



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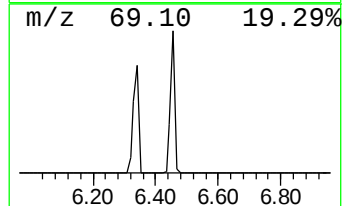
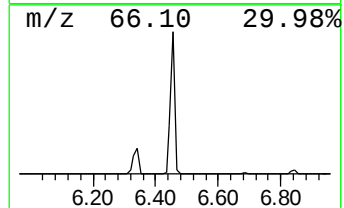
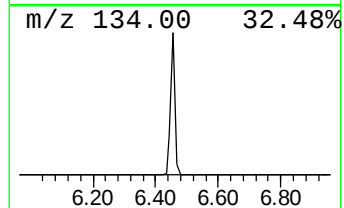
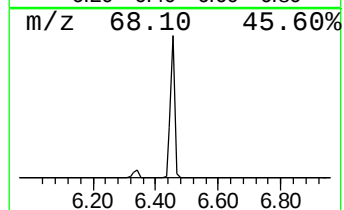
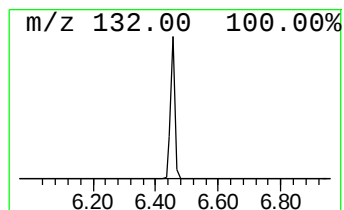
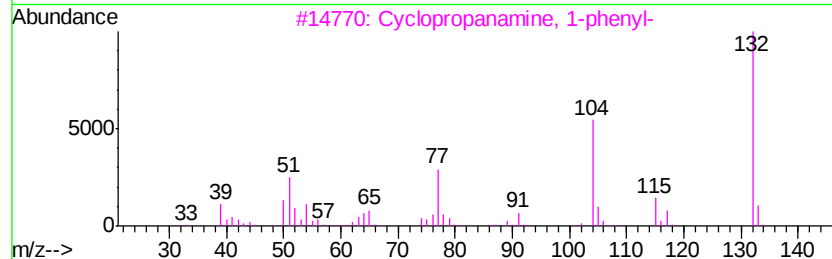
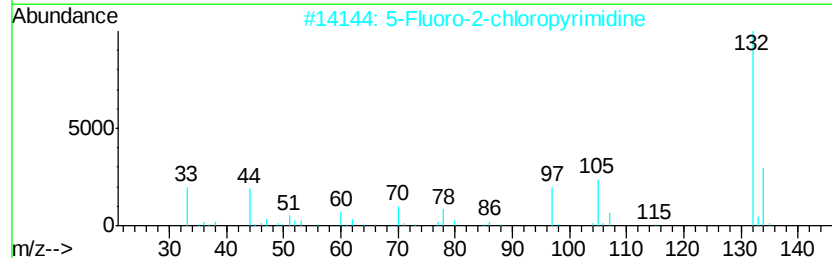
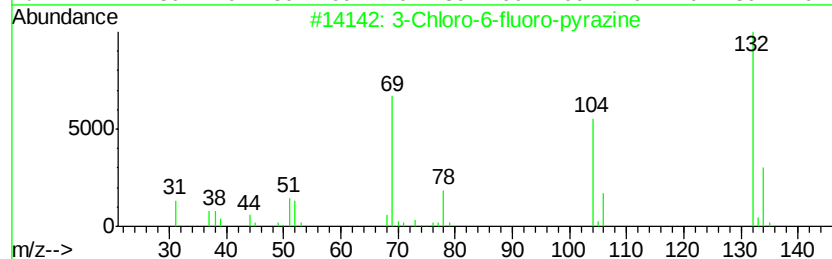
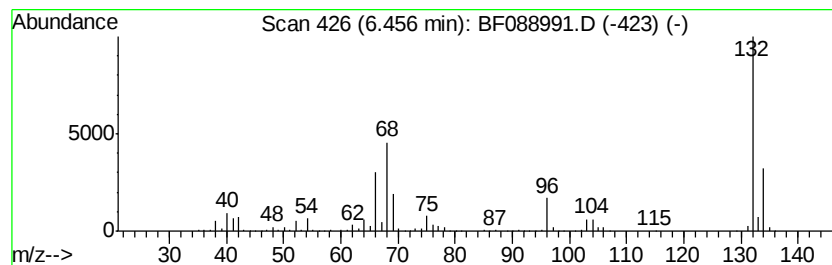
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TIC Library : C:\Database\NIST11.L
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Peak Number 5 unknown6.46 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.46	91.93 ng	1660920	1,4-Dichlorobenzene-d4	6.68

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
3		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	10
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		Xenon	132	Xe	007440-63-3	9



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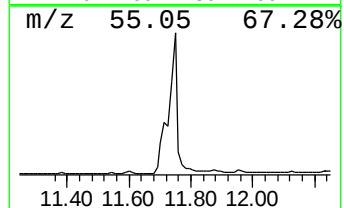
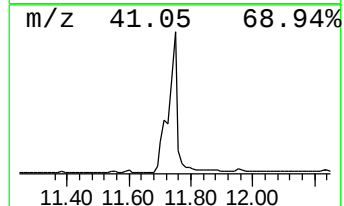
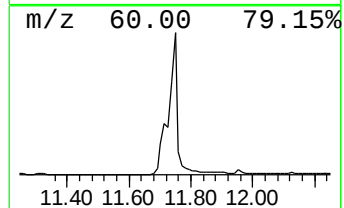
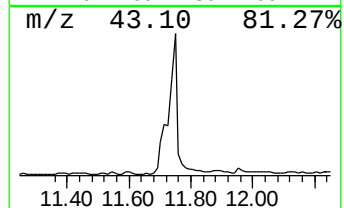
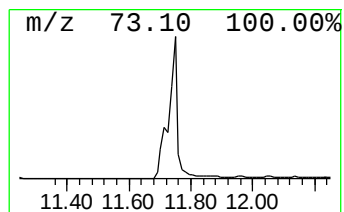
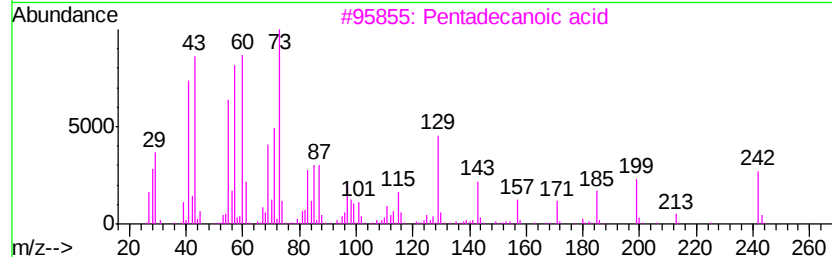
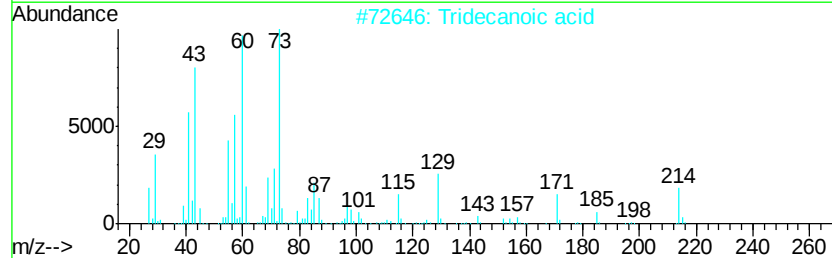
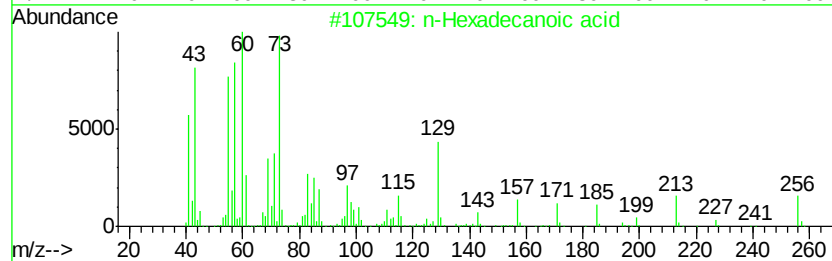
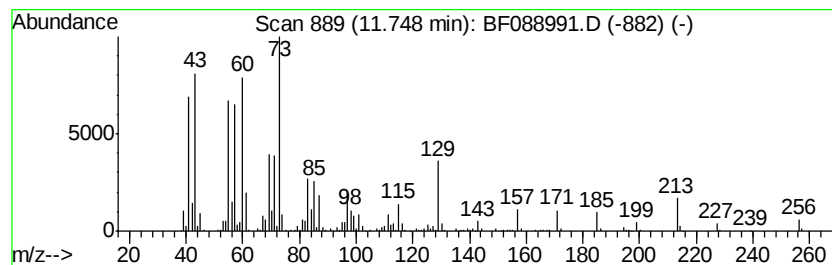
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Peak Number 7 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.75	98.27 ng	2688510	Phenanthrene-d10	11.19

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		Tridecanoic acid	214	C13H26O2	000638-53-9	91
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	91
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	80
5		Estra-1,3,5(10)-trien-17.beta.-ol	256	C18H24O	002529-64-8	30



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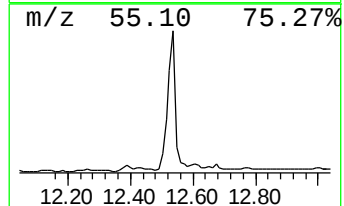
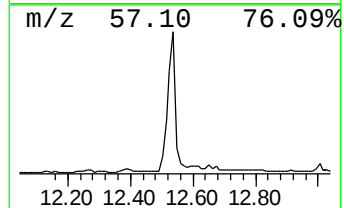
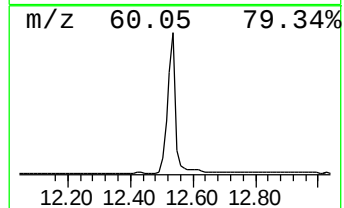
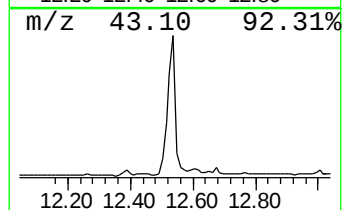
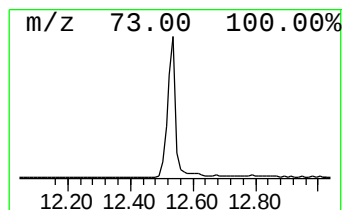
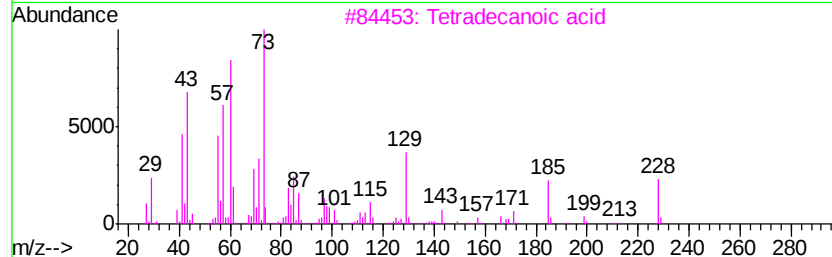
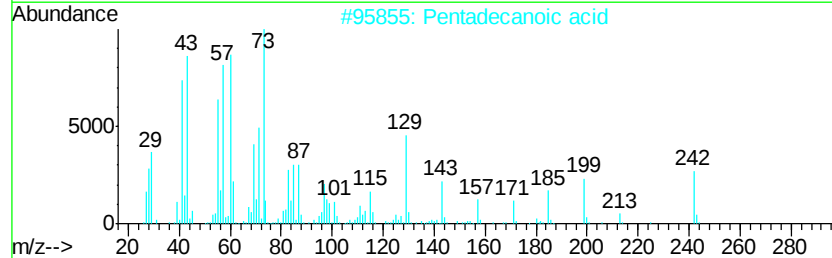
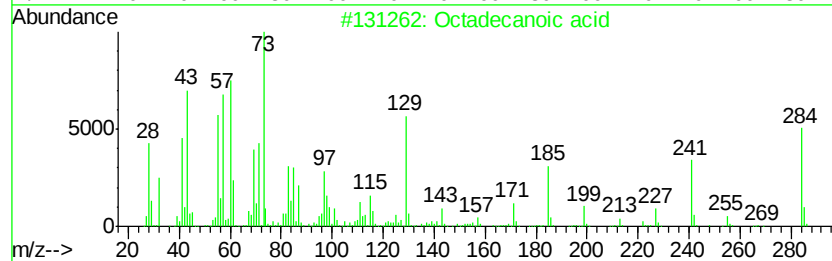
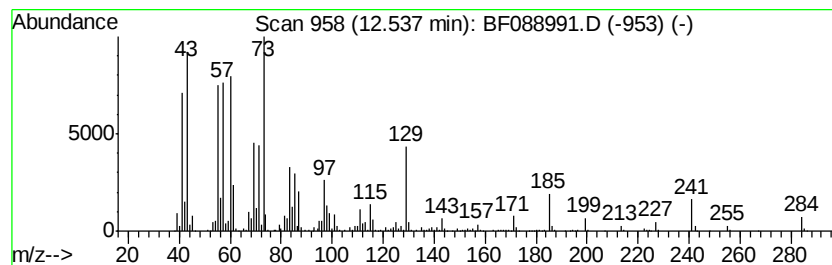
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Octadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.54	113.50 ng	3129060	Chrysene-d12	13.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	91
3		Tetradecanoic acid	228	C14H28O2	000544-63-8	83
4		Tridecanoic acid	214	C13H26O2	000638-53-9	76
5		n-Decanoic acid	172	C10H20O2	000334-48-5	68



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF071416\
Data File : BF088991.D
Acq On : 14 Jul 2016 17:46
Operator : UM/SJ
Sample : PB92031BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
PB92031BL

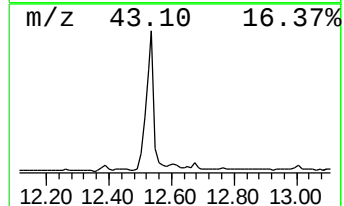
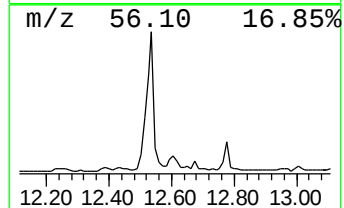
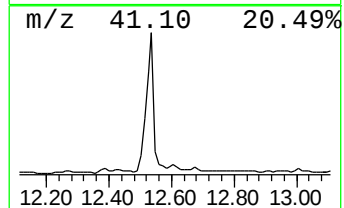
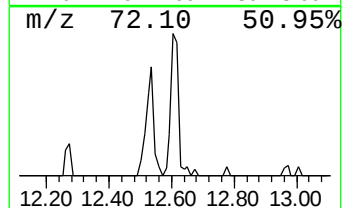
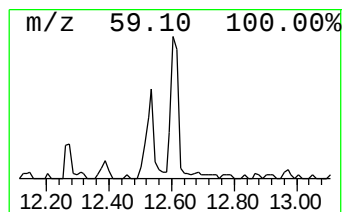
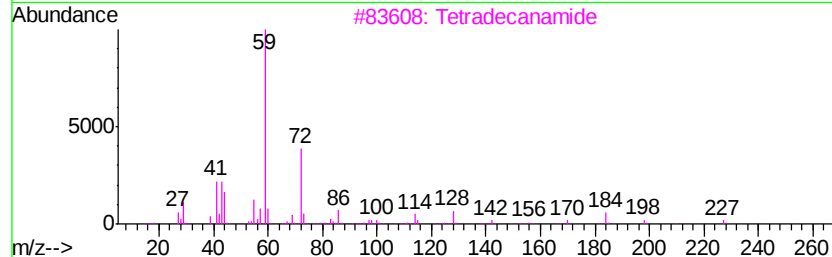
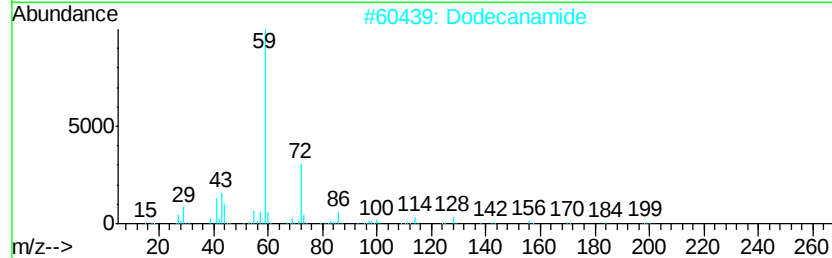
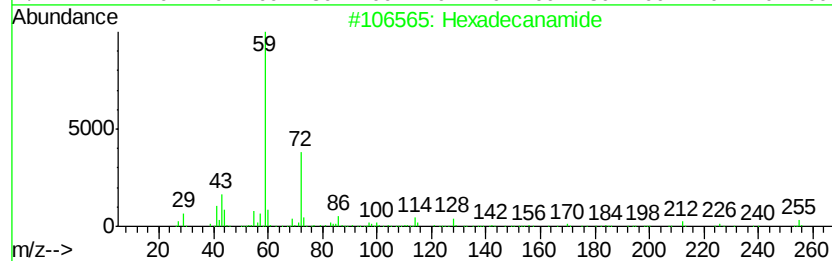
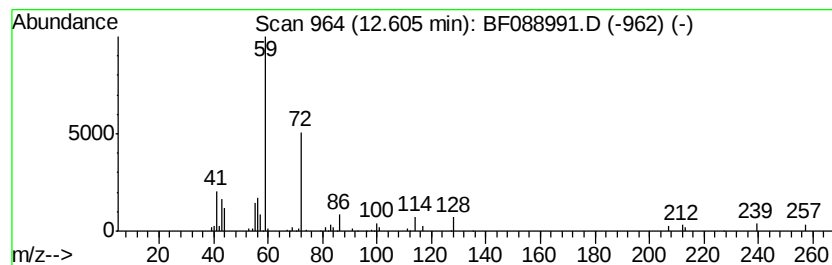
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Hexadecanamide Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.61	7.29 ng	200976	Chrysene-d12	13.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecanamide	255	C16H33NO	000629-54-9	72
2		Dodecanamide	199	C12H25NO	001120-16-7	72
3		Tetradecanamide	227	C14H29NO	000638-58-4	72
4		Decanamide-	171	C10H21NO	002319-29-1	72
5		9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	40



Data Path : Z:\HPCHEM1\BNA F\DATA\BF071416\
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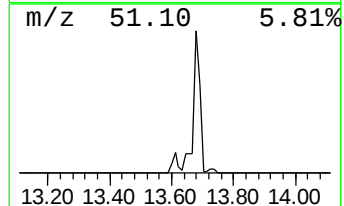
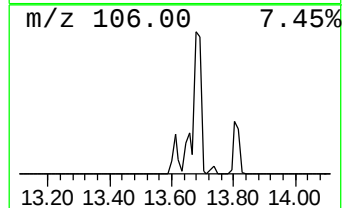
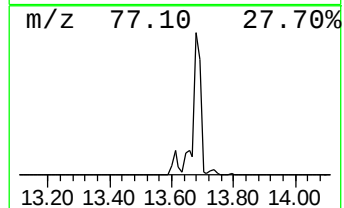
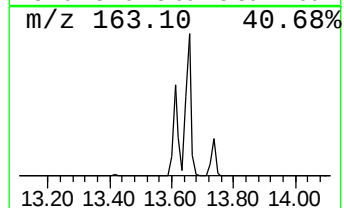
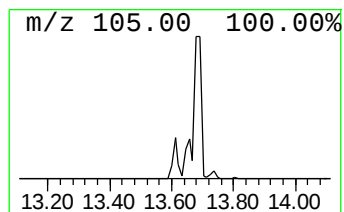
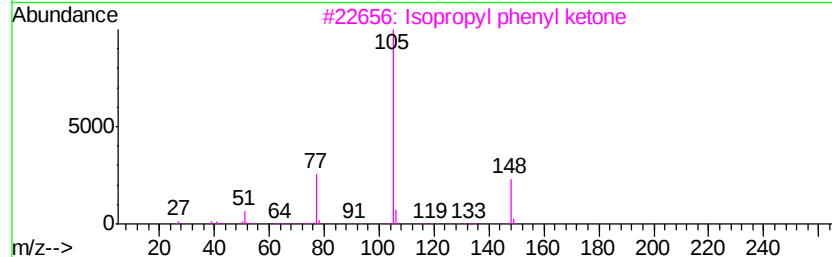
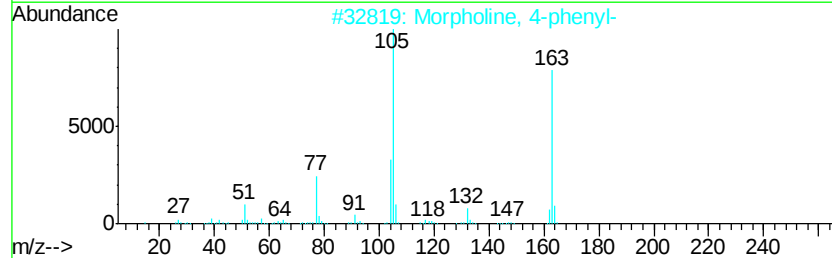
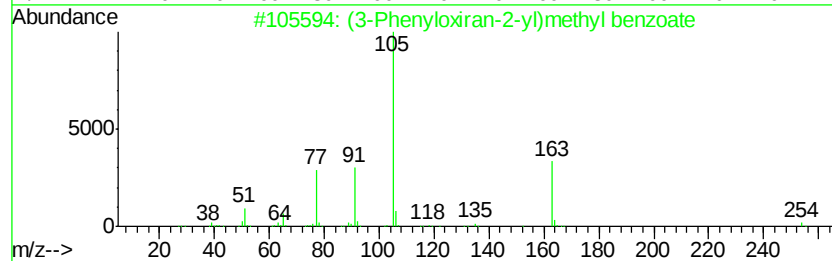
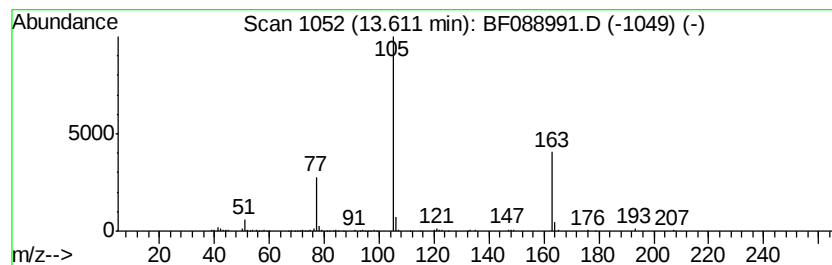
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 (3-Phenyloxiran-2-yl)methyl... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.61	10.41 ng	286889	Chrysene-d12	13.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(3-Phenyloxiran-2-yl)methyl benz...	254	C16H14O3	1000366-96-1	56
2		Morpholine, 4-phenyl-	163	C10H13NO	000092-53-5	53
3		Isopropyl phenyl ketone	148	C10H12O	000611-70-1	38
4		Benzoic acid, 2-propenyl ester	162	C10H10O2	000583-04-0	38
5		Benzo[b]thiophene-3-carboxylic a...	343	C18H17NO4S	096334-43-9	38



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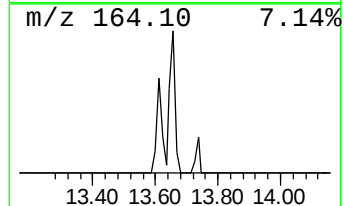
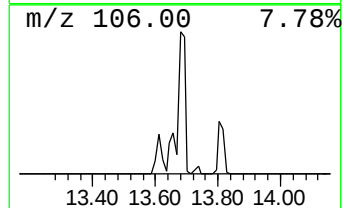
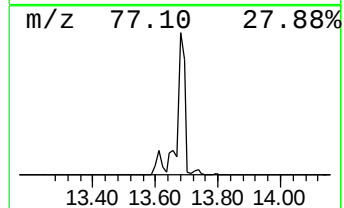
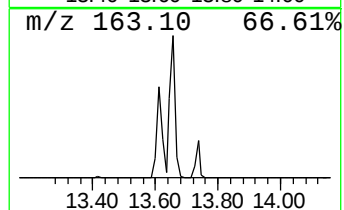
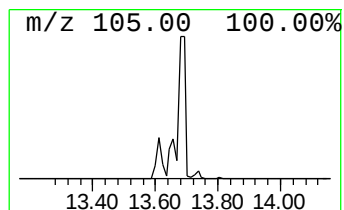
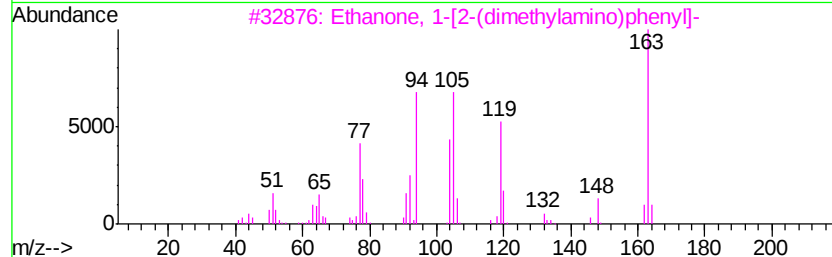
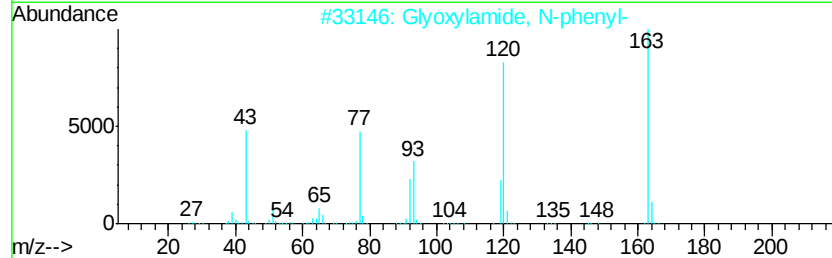
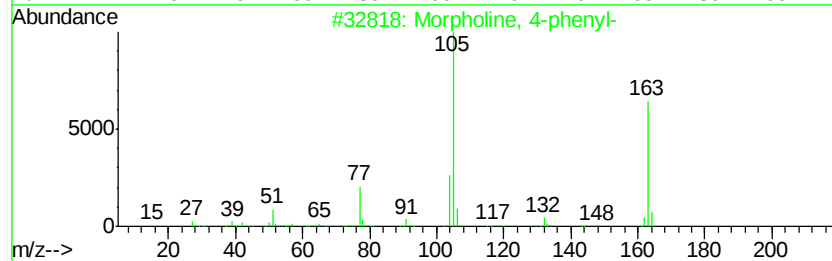
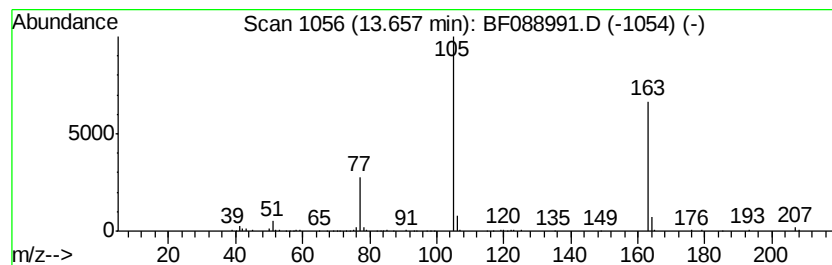
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Morpholine, 4-phenyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.66	15.19 ng	418911	Chrysene-d12	13.82
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Morpholine, 4-phenyl-	163 C10H13NO	000092-53-5 56
2		Glyoxylamide, N-phenyl-	163 C9H9NO2	032331-52-5 50
3		Ethanone, 1-[2-(dimethylamino)ph...	163 C10H13NO	010336-55-7 45
4		Benzoic acid, 2-benzoyloxy-3-brom...	348 C17H17BrO3	1000191-33-3 42
5		3-(Benzoylthio)-2-methylpropanoi...	224 C11H12O3S	067714-34-5 38



Data Path : Z:\HPCHEM1\BNA F\DATA\BF071416\
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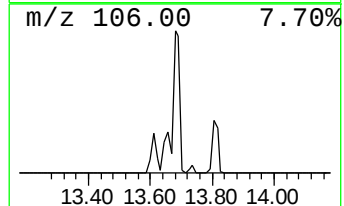
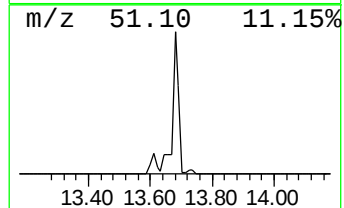
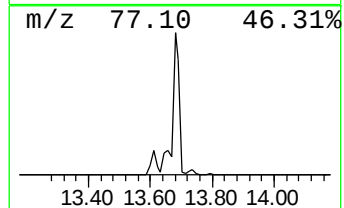
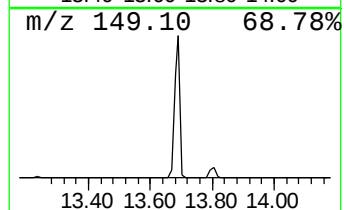
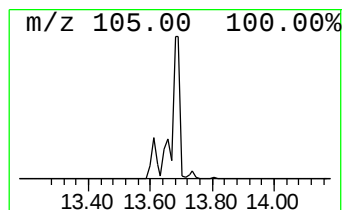
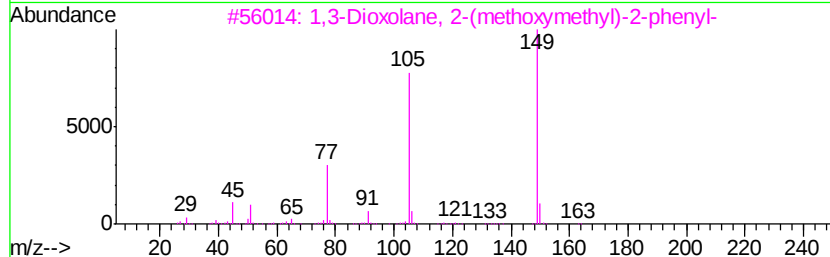
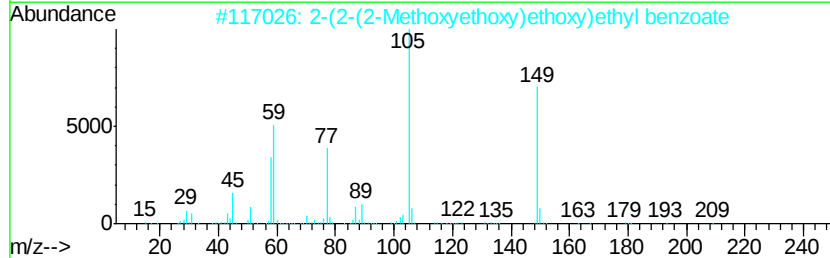
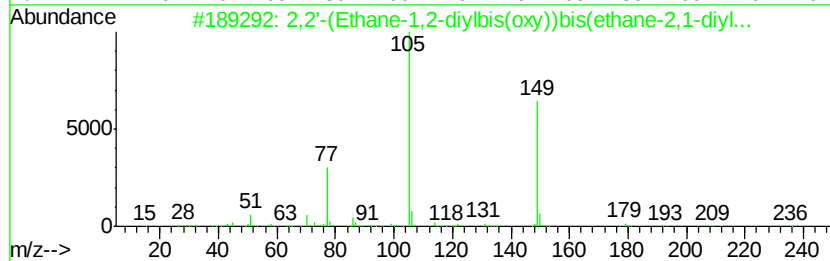
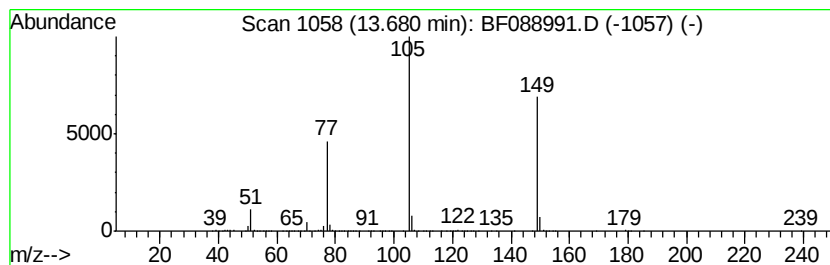
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 2,2'-(Ethane-1,2-diylbis(ox... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.68	52.55 ng	1448770	Chrysene-d12	13.82	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2'-(Ethane-1,2-diylbis(oxv))bi...	358	C20H22O6	1000366-99-4	74
2	2-(2-(2-Methoxvethoxy)ethoxy)eth...	268	C14H20O5	1000366-99-1	74
3	1,3-Dioxolane, 2-(methoxymethyl)...	194	C11H14O3	1000156-71-2	59
4	Diethylene glycol dibenzoate	314	C18H18O5	000120-55-8	59
5	Ethanol, 2-(4-phenoxyphenoxy)-, ...	334	C21H18O4	055191-59-8	56



Data Path : Z:\HPCHEM1\BNA F\DATA\BF071416\
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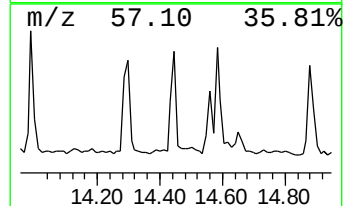
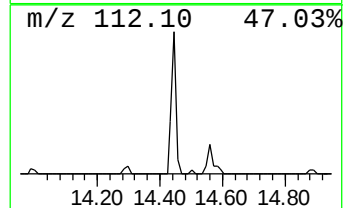
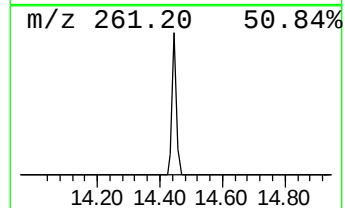
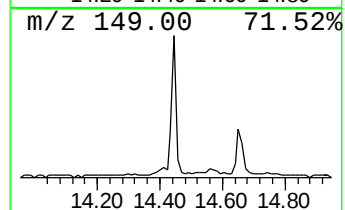
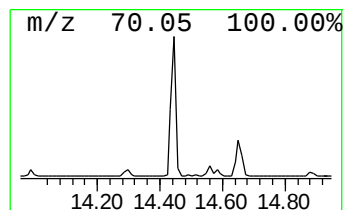
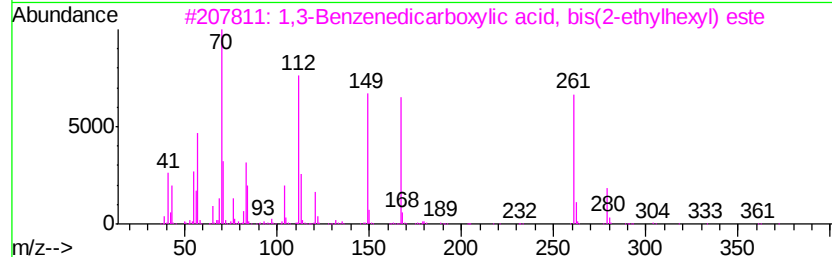
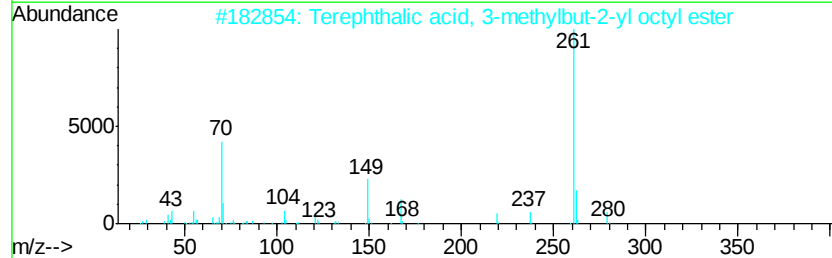
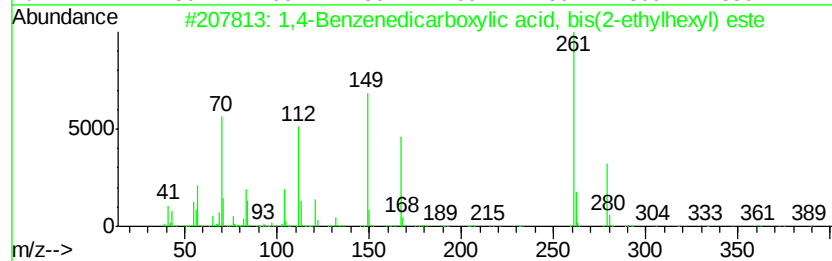
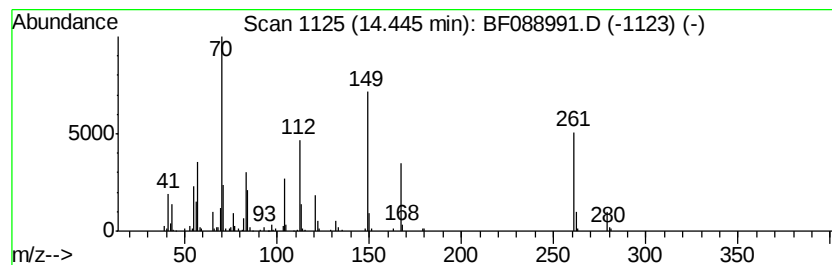
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 1,4-Benzenedicarboxylic aci... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.45	6.81 ng	187723	Chrysene-d12	13.82

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,4-Benzenedicarboxylic acid, bi...	390	C24H38O4	006422-86-2	72
2			Terephthalic acid, 3-methylbut-2...	348	C21H32O4	1000356-27-6	64
3			1,3-Benzenedicarboxylic acid, bi...	390	C24H38O4	000137-89-3	62
4			Terephthalic acid, monochloride,...	296	C16H21ClO3	1000324-22-9	43
5			Terephthalic acid, phenyl octyl ...	354	C22H26O4	1000324-06-6	30



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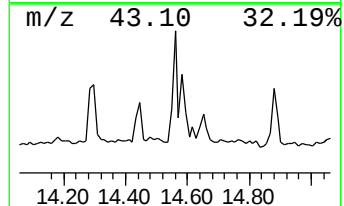
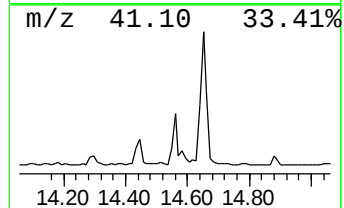
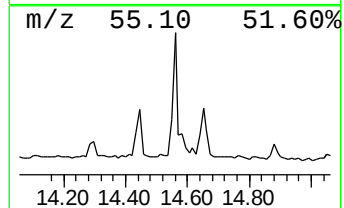
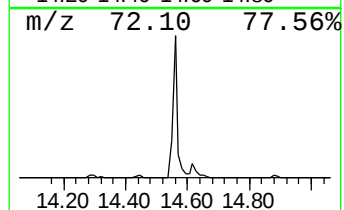
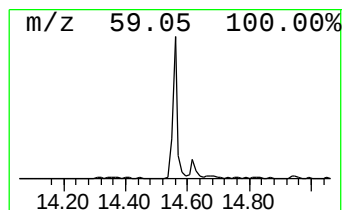
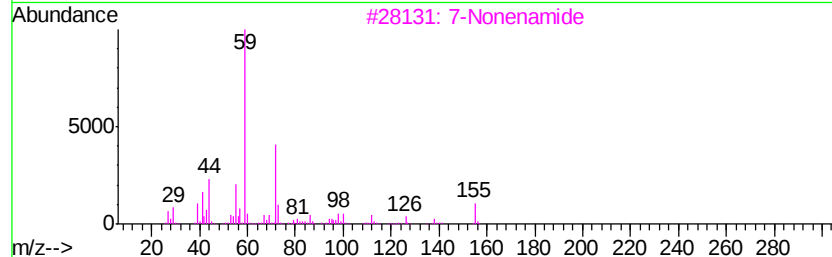
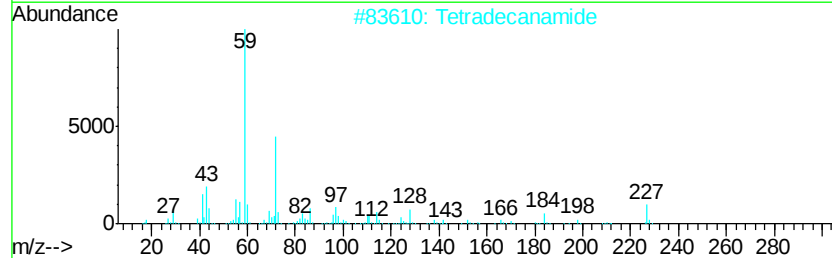
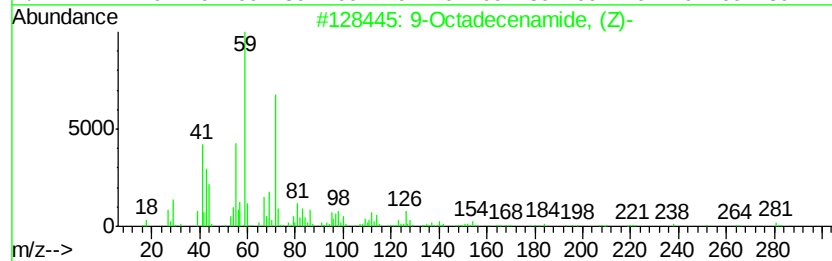
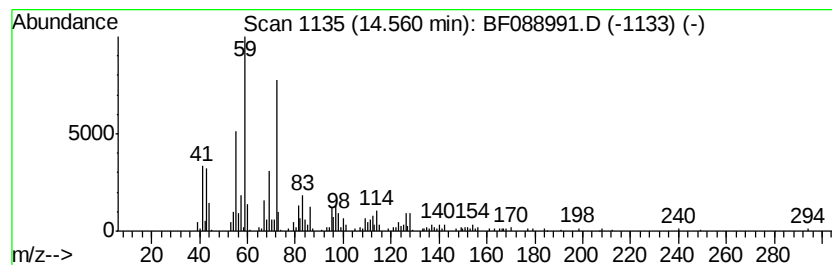
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 15 9-Octadecenamide, (Z)- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.56	16.01 ng	235110	Perylene-d12	15.18

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	91
2		Tetradecanamide	227	C14H29NO	000638-58-4	53
3		7-Nonenamide	155	C9H17NO	090949-53-4	47
4		1-Heptene, 2-methoxy-	128	C8H16O	061142-46-9	35
5		Acetic acid, cyanohydroxyimino-,...	128	C4H4N2O3	061295-92-9	35



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Data File : BF088991.D
Acq On : 14 Jul 2016 17:46
Operator : UM/SJ
Sample : PB92031BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

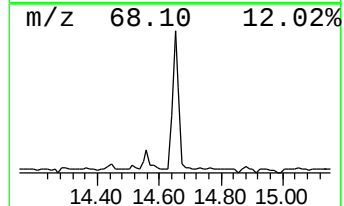
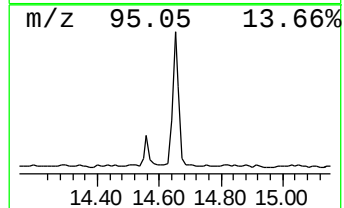
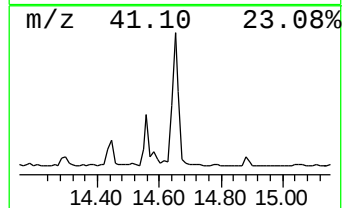
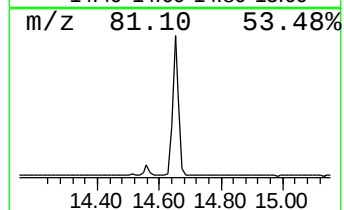
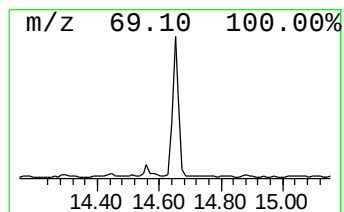
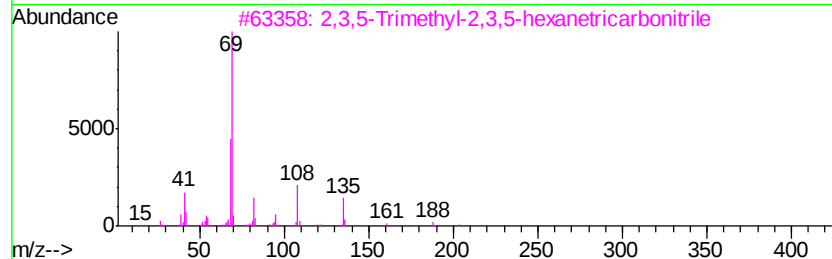
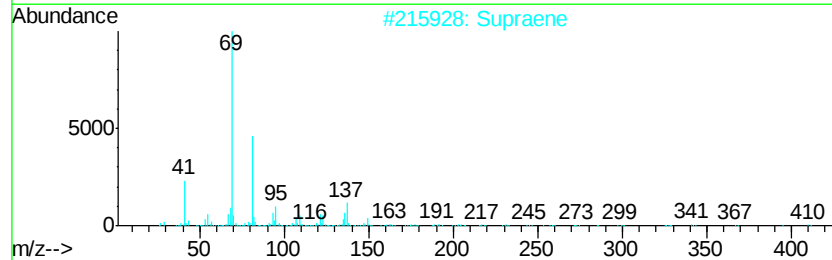
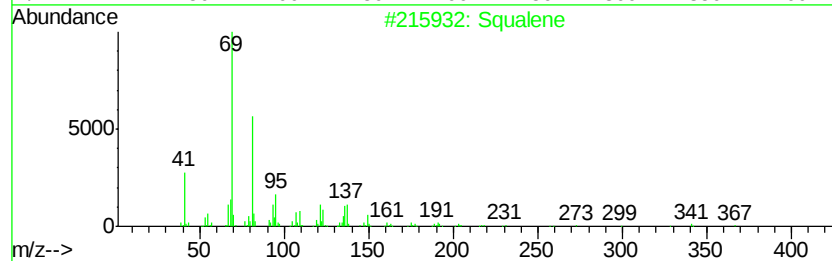
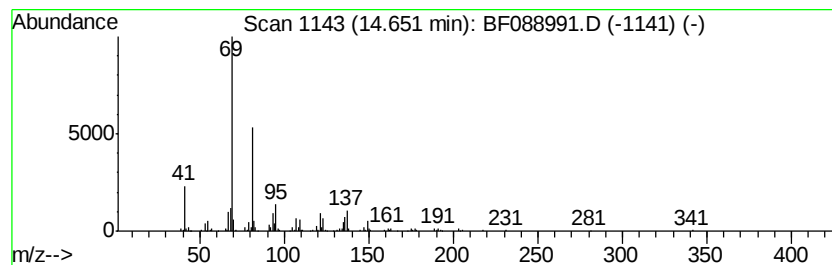
Instrument :
BNA_F
ClientSampleId :
PB92031BL

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

Peak Number 16 Squalene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.65	40.05 ng	588050	Perylene-d12	15.18
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Squalene	410 C30H50	000111-02-4	91
2	Supraene	410 C30H50	007683-64-9	87
3	2,3,5-Trimethyl-2,3,5-hexanetric...	203 C12H17N3	003974-79-6	50
4	Butanoic acid, 4-cyano-2-ethyl-2...	290 C15H18N2O4	203808-68-8	47
5	1,5-Heptadiene, 2,3,6-trimethyl-	138 C10H18	033501-88-1	47



Data Path : Z:\HPCHEM1\BNA F\DATA\BF071416\
Data File : BF088991.D
Acq On : 14 Jul 2016 17:46
Operator : UM/SJ
Sample : PB92031BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampled :
PB92031BL

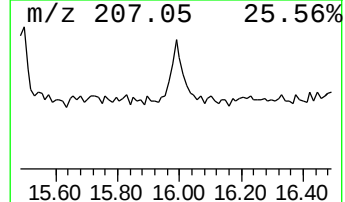
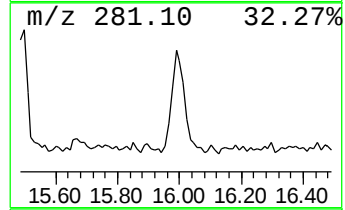
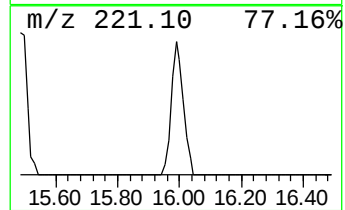
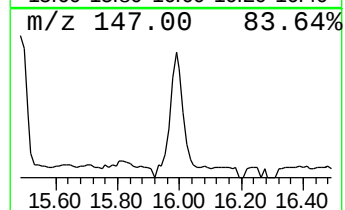
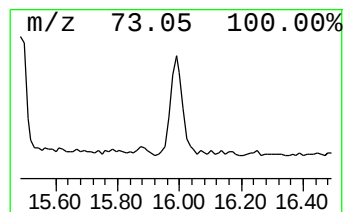
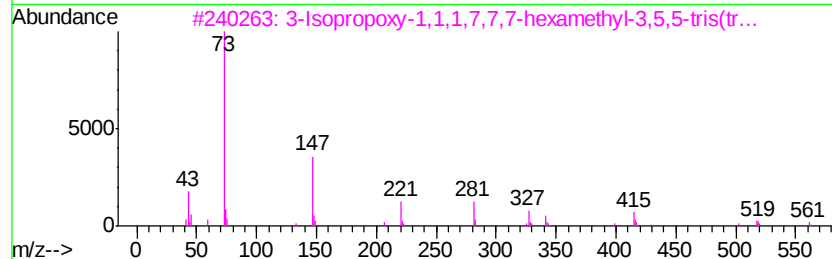
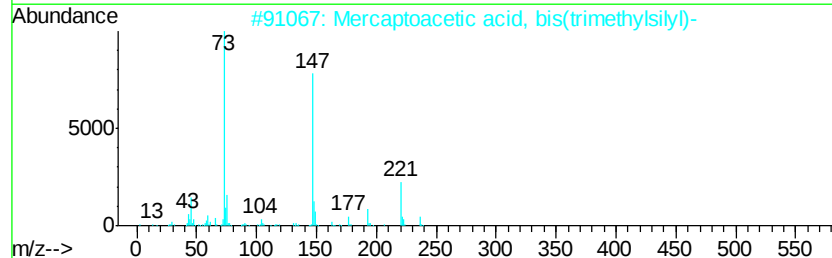
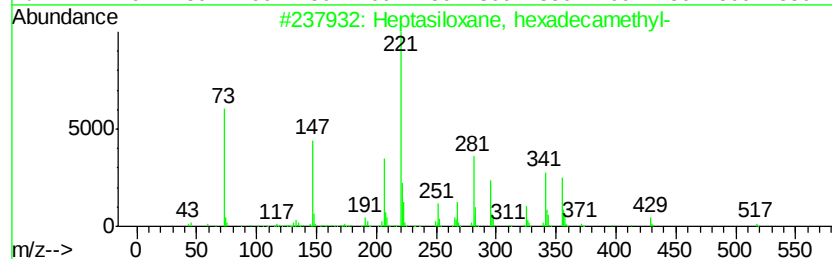
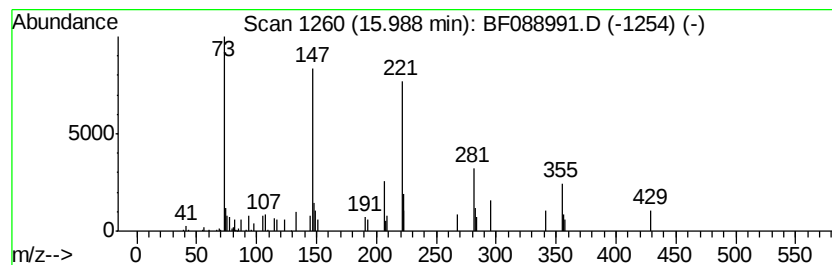
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF063016.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Heptasiloxane, hexadecamethyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.99	8.47 ng	124370	Perylene-d12	15.18

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptasiloxane, hexadecamethyl-	532	C16H48O6Si7	000541-01-5	58
2		Mercaptoacetic acid, bis(trimeth...	236	C8H20O2SSi2	006398-62-5	40
3		3-Isopropoxy-1,1,1,7,7,7-hexamet...	576	C18H52O7Si7	071579-69-6	32
4		N-Benzyl-N-ethyl-p-isopropylbenz...	281	C19H23NO	015089-22-2	27
5		Tris(trimethylsilyl)borate	278	C9H27B03Si3	004325-85-3	27



Data Path : Z:\HPCHEM1\BNA F\DATA\BF071416\
Data File : BF088991.D
Acq On : 14 Jul 2016 17:46
Operator : UM/SJ
Sample : PB92031BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampled :
PB92031BL

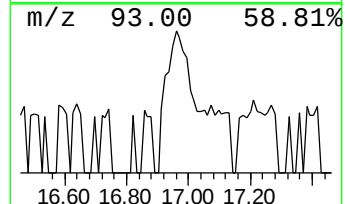
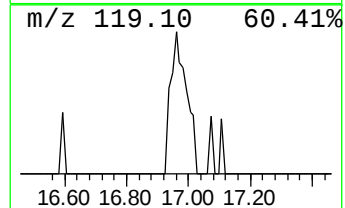
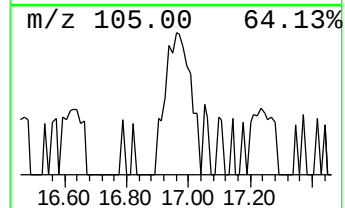
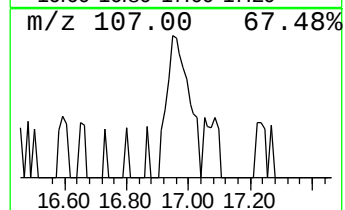
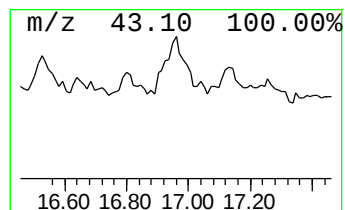
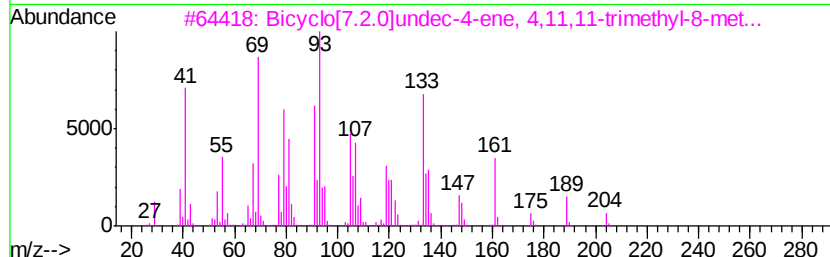
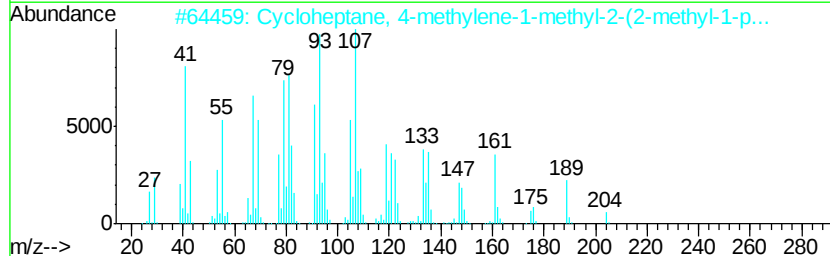
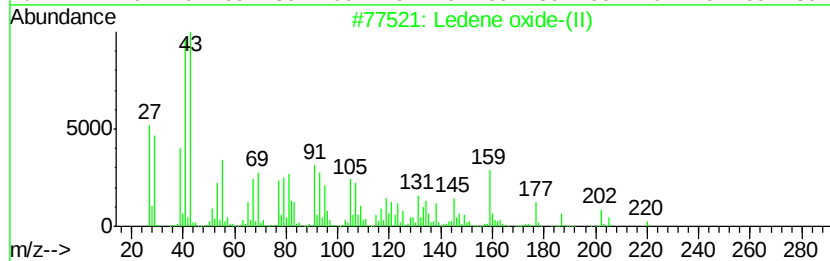
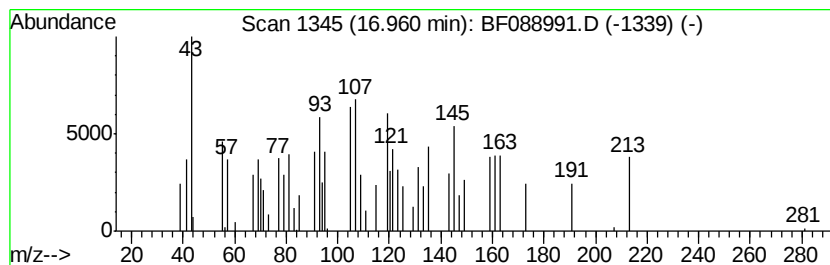
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF063016.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 unknown16.96 Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.96	5.80 ng	85109	Perylene-d12	15.18

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ledene oxide-(II)	220	C15H24O	1000159-36-7	27
2			Cycloheptane, 4-methylene-1-meth...	204	C15H24	1000159-38-5	16
3			Bicyclo[7.2.0]undec-4-ene, 4,11,...	204	C15H24	013877-93-5	16
4			(+)-2-Carene, 2-acetyl-	178	C12H18O	1000151-09-7	14
5			1-Cycloheptene, 1,4-dimethyl-3-(...	204	C15H24	1000159-38-6	14



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF071416\
 Data File : BF088991.D
 Acq On : 14 Jul 2016 17:46
 Operator : UM/SJ
 Sample : PB92031BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB92031BL

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
Propane, 1,1-dime...	1.91	6.2	ng	111344	1	6.68	361355	20.0
2-Pentanone, 4-hy...	4.84	11.5	ng	207561	1	6.68	361355	20.0
unknown6.46	6.46	91.9	ng	1660920	1	6.68	361355	20.0
n-Hexadecanoic acid	11.75	98.3	ng	2688510	4	11.19	547194	20.0
Octadecanoic acid	12.54	113.5	ng	3129060	5	13.82	551385	20.0
Hexadecanamide	12.61	7.3	ng	200976	5	13.82	551385	20.0
(3-Phenyloxiran-2...	13.61	10.4	ng	286889	5	13.82	551385	20.0
Morpholine, 4-phe...	13.66	15.2	ng	418911	5	13.82	551385	20.0
2,2'-(Ethane-1,2-...	13.68	52.5	ng	1448770	5	13.82	551385	20.0
1,4-Benzenedicarb...	14.45	6.8	ng	187723	5	13.82	551385	20.0
9-Octadecenamide,...	14.56	16.0	ng	235110	6	15.18	293670	20.0
Squalene	14.65	40.0	ng	588050	6	15.18	293670	20.0
Heptasiloxane, he...	15.99	8.5	ng	124370	6	15.18	293670	20.0
unknown16.96	16.96	5.8	ng	85109	6	15.18	293670	20.0