

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF071416\
 Data File : BF088997.D
 Acq On : 14 Jul 2016 20:50
 Operator : UM/SJ
 Sample : H3996-09
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 U2-B2(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	1.655	5	6	15	rVB	147599	287445	16.14%	1.676%
2	1.781	15	17	23	rVV	1048271	1368918	76.84%	7.984%
3	1.861	23	24	67	rVB2	78766	371151	20.83%	2.165%
4	2.821	106	108	120	rBV	17184	34384	1.93%	0.201%
5	4.833	282	284	291	rBB	124562	178841	10.04%	1.043%
6	5.279	321	323	326	rBV	1348907	1394170	78.26%	8.131%
7	6.342	412	416	418	rBV	1432025	1716030	96.33%	10.008%
8	6.456	423	426	428	rBV	1714948	1545579	86.76%	9.014%
9	6.685	444	446	448	rBV	447695	353294	19.83%	2.060%
10	6.845	457	460	462	rBB	1449780	1316118	73.88%	7.676%
11	7.256	493	496	498	rBV	1014137	1066422	59.86%	6.220%
12	7.553	520	522	524	rVB	41864	55637	3.12%	0.324%
13	7.965	556	558	561	rBV	325471	413491	23.21%	2.412%
14	8.193	576	578	580	rBV	31374	25981	1.46%	0.152%
15	8.250	580	583	585	rVB	42539	35717	2.00%	0.208%
16	9.050	650	653	655	rBV	1958073	1781486	100.00%	10.390%
17	9.165	662	663	665	rBV	28844	28356	1.59%	0.165%
18	9.439	685	687	690	rBV	232278	206516	11.59%	1.204%
19	9.713	709	711	713	rVB	441599	400160	22.46%	2.334%
20	10.502	777	780	782	rBV	817484	757048	42.50%	4.415%
21	10.628	790	791	796	rBV	18654	32826	1.84%	0.191%
22	11.085	829	831	832	rBV	26826	38480	2.16%	0.224%
23	11.188	838	840	844	rVB	401404	377383	21.18%	2.201%
24	11.268	844	847	848	rBV2	13494	19544	1.10%	0.114%
25	11.508	865	868	870	rBV	17652	22544	1.27%	0.131%
26	11.805	892	894	901	rVB4	10104	22941	1.29%	0.134%
27	12.388	943	945	948	rBV	57935	73913	4.15%	0.431%
28	12.617	962	965	967	rBV	106770	111264	6.25%	0.649%
29	12.777	976	979	981	rBV	1326773	1213176	68.10%	7.075%
30	12.948	991	994	996	rBV	16286	27664	1.55%	0.161%
31	13.017	996	1000	1001	rVV2	24292	45226	2.54%	0.264%
32	13.051	1001	1003	1006	rVV	33667	42941	2.41%	0.250%
33	13.142	1009	1011	1013	rBV2	13861	24969	1.40%	0.146%
34	13.314	1024	1026	1027	rBV	30380	30486	1.71%	0.178%

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Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

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35	13.417	1034	1035	1036	rBV	98278	67567	3.79%	0.394%
36	13.451	1036	1038	1039	rBV	20122	19531	1.10%	0.114%
37	13.554	1046	1047	1049	rBV2	19358	20623	1.16%	0.120%
38	13.680	1057	1058	1062	rVB	87743	107124	6.01%	0.625%
39	13.817	1067	1070	1072	rBV	467349	591357	33.19%	3.449%
40	14.000	1084	1086	1091	rBV4	13690	25556	1.43%	0.149%
41	14.320	1111	1114	1119	rBV2	55099	115880	6.50%	0.676%
42	14.457	1124	1126	1128	rVB	98509	88982	4.99%	0.519%
43	14.731	1147	1150	1151	rVB	33521	47503	2.67%	0.277%
44	14.800	1154	1156	1161	rVB	43920	96580	5.42%	0.563%
45	14.902	1163	1165	1166	rBV	28017	45191	2.54%	0.264%
46	15.063	1177	1179	1182	rVB	29159	43439	2.44%	0.253%
47	15.120	1182	1184	1186	rVB	31930	33431	1.88%	0.195%
48	15.177	1187	1189	1192	rBV	205126	258490	14.51%	1.508%
49	15.611	1225	1227	1229	rBV	42249	56799	3.19%	0.331%
50	16.457	1298	1301	1302	rBV2	14355	28994	1.63%	0.169%
51	16.834	1331	1334	1339	rBV	14202	36149	2.03%	0.211%
52	17.668	1403	1407	1409	rVB4	23593	42864	2.41%	0.250%

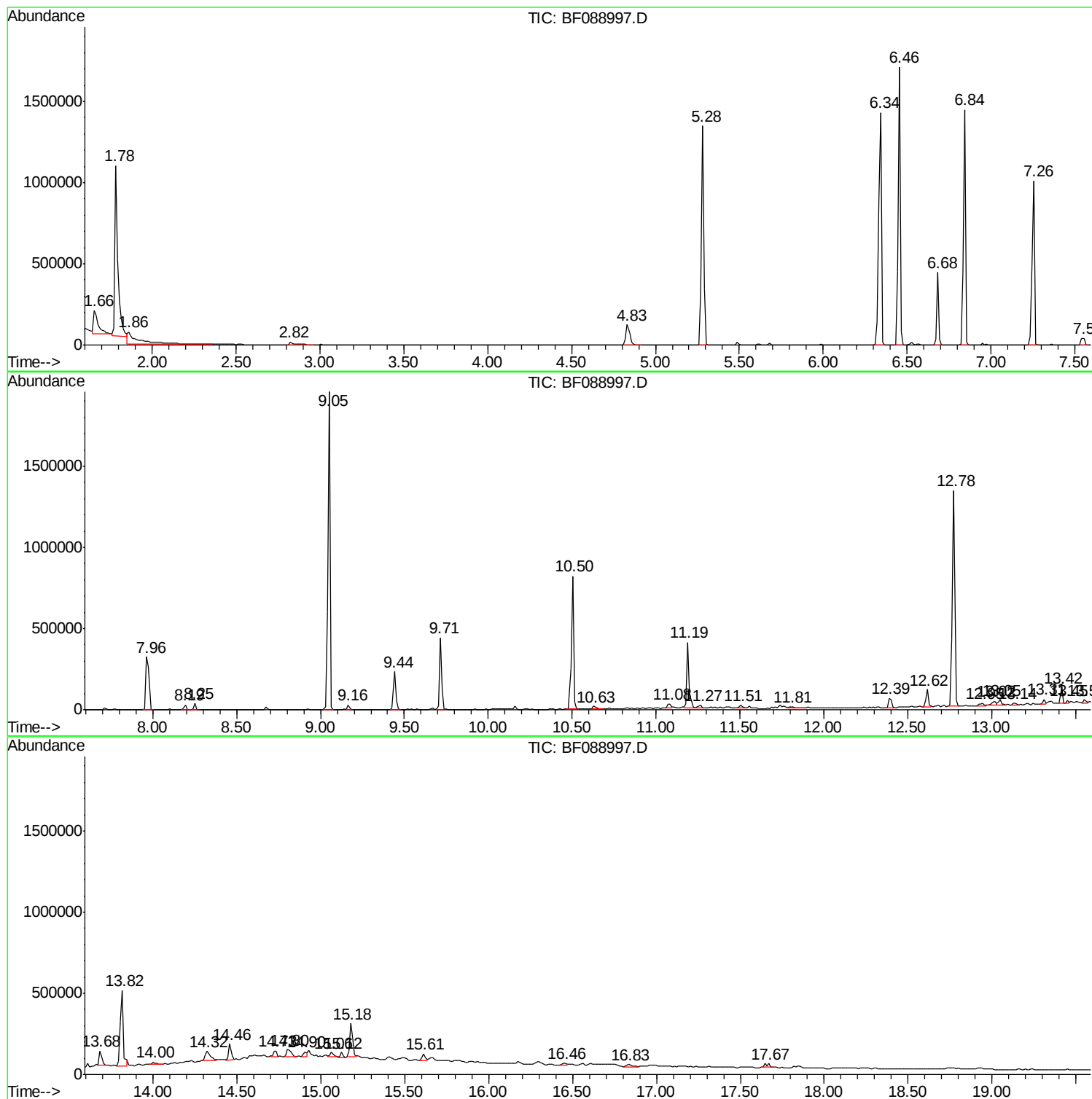
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF063016.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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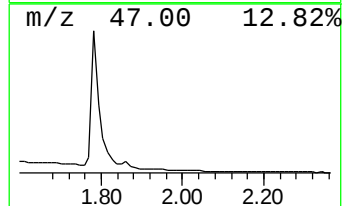
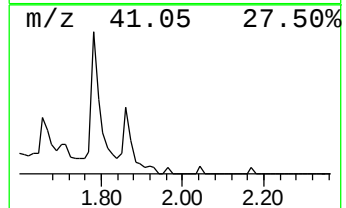
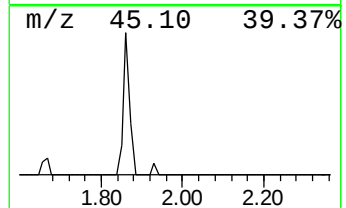
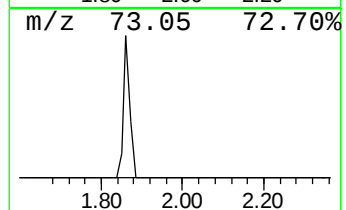
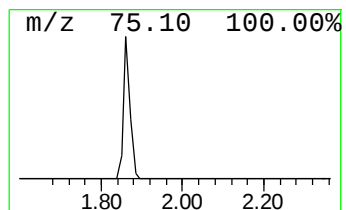
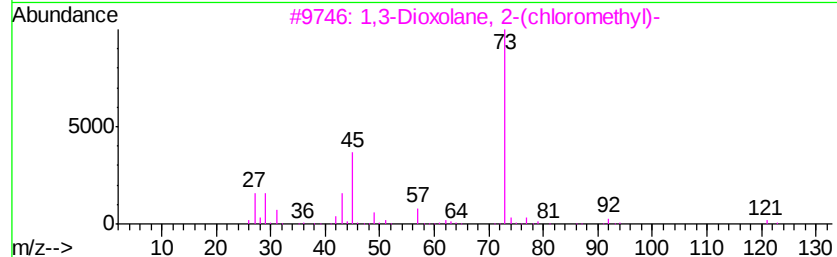
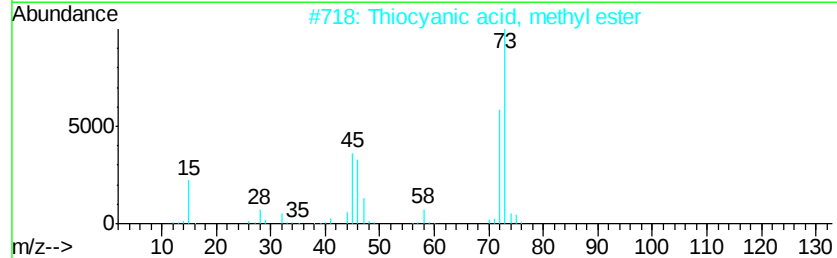
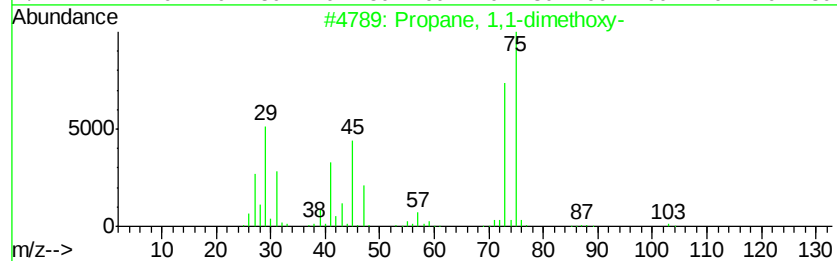
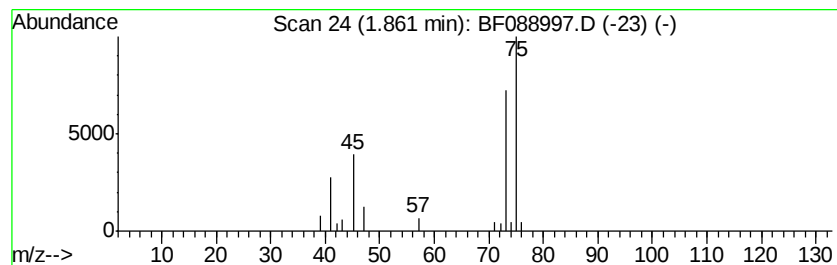
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 Propane, 1,1-dimethoxy- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.86	21.01 ng	371151	1,4-Dichlorobenzene-d4	6.68

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propane, 1,1-dimethoxy-	104	C5H12O2	004744-10-9	72
2		Thiocyanic acid, methyl ester	73	C2H3NS	000556-64-9	25
3		1,3-Dioxolane, 2-(chloromethyl)-	122	C4H7ClO2	002568-30-1	9
4		Glycine	75	C2H5NO2	000056-40-6	7
5		Methane, isothiocyanato-	73	C2H3NS	000556-61-6	4



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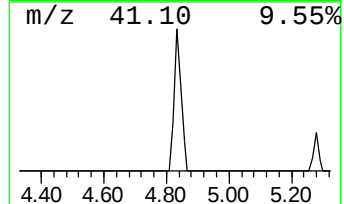
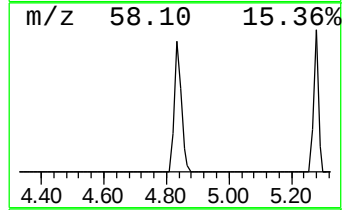
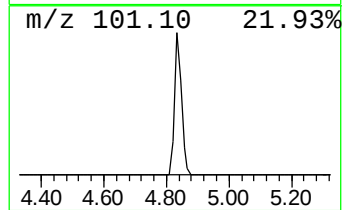
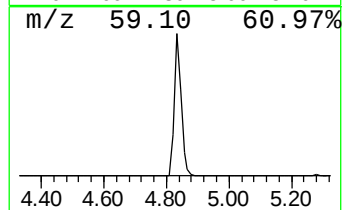
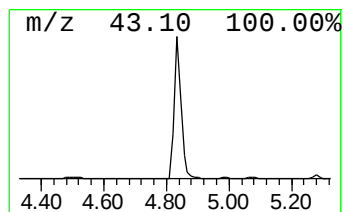
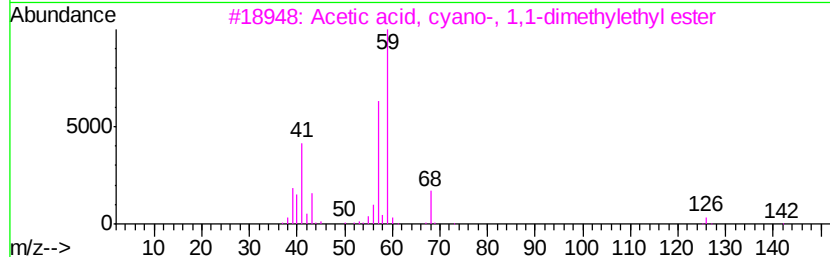
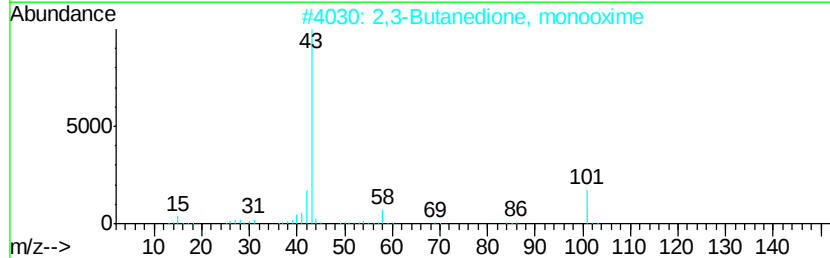
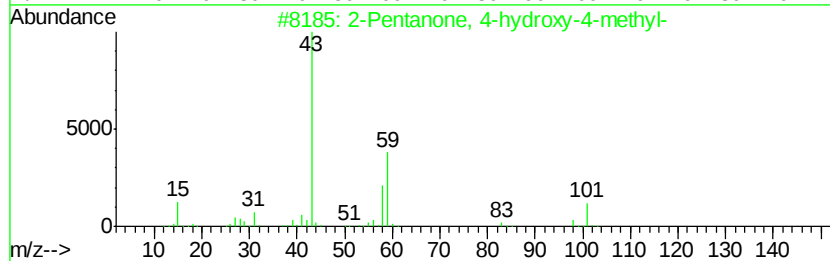
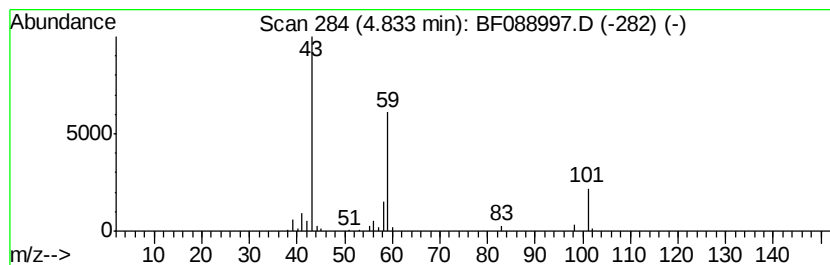
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TIC Integration Parameters: LSCINT.P

Peak Number 4 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.83	10.12 ng	178841	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	35
3			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17
4			(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	9
5			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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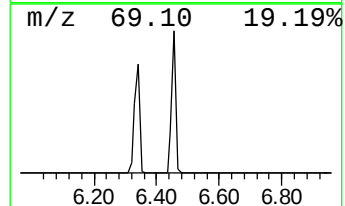
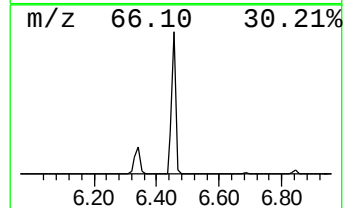
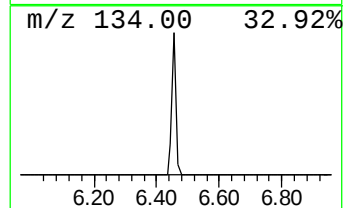
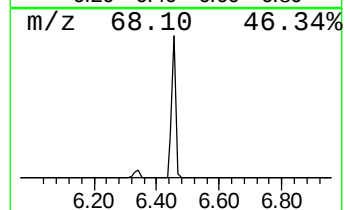
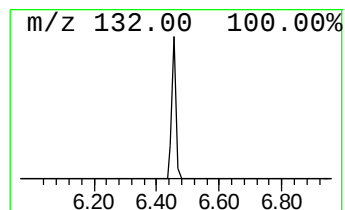
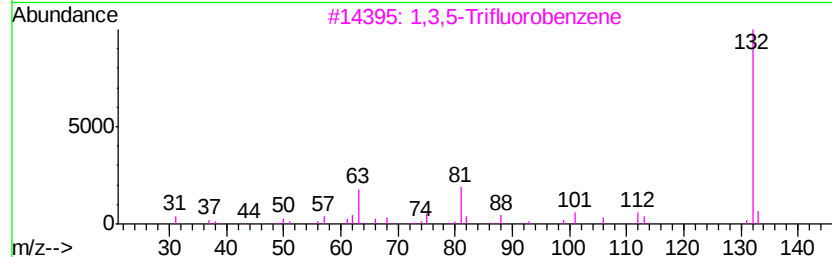
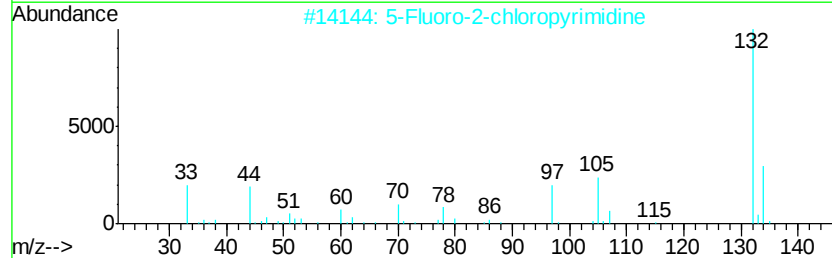
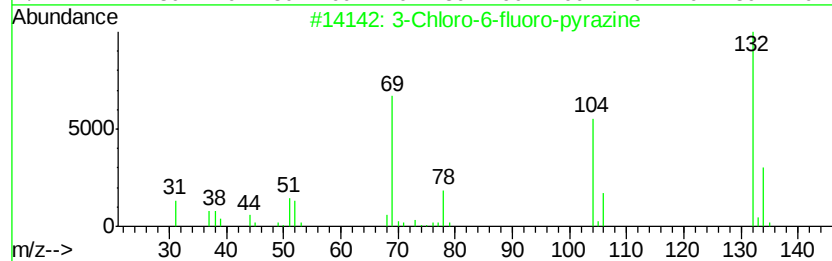
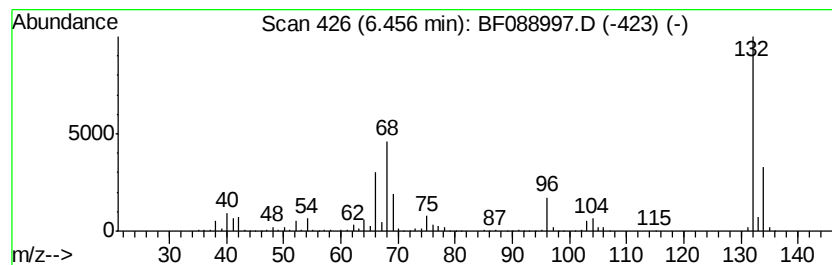
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TIC Library : C:\Database\NIST11.L
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Peak Number 5 unknown6.46 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.46	87.50 ng	1545580	1,4-Dichlorobenzene-d4	6.68

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27	
2	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16	
3	1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	14	
4	Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10	
5	Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	10	



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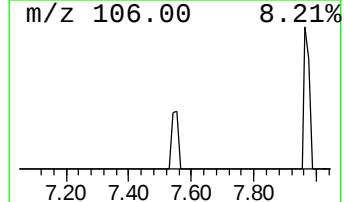
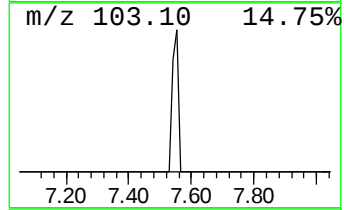
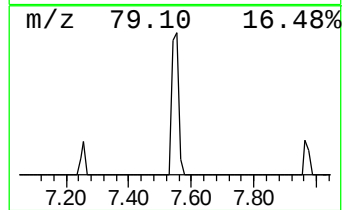
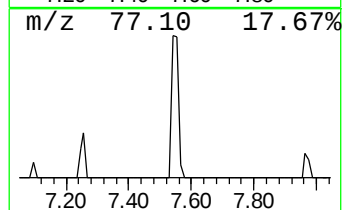
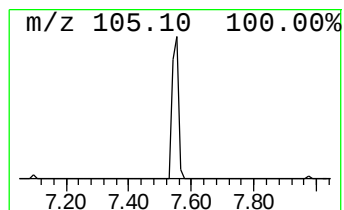
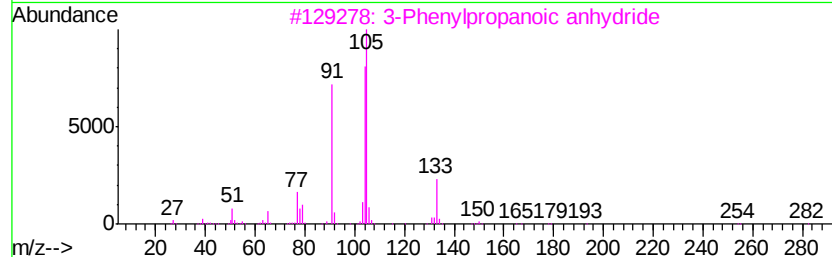
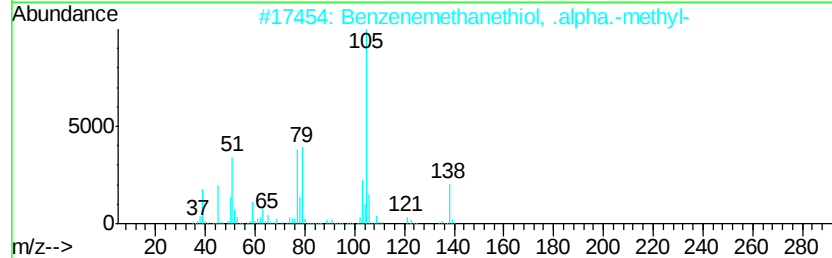
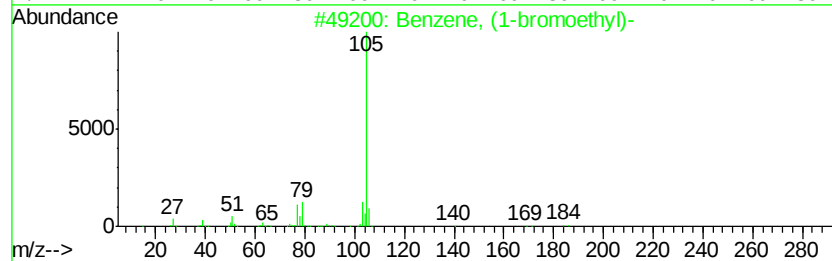
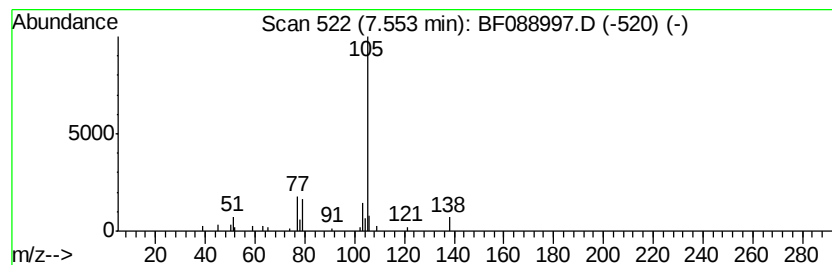
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Peak Number 7 Benzene, (1-bromoethyl)- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.55	2.69 ng	55637	Napthalene-d8	7.96

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, (1-bromoethyl)-	184	C8H9Br	000585-71-7	72
2		Benzenemethanethiol, .alpha.-met...	138	C8H10S	006263-65-6	64
3		3-Phenylpropanoic anhydride	282	C18H18O3	1000333-89-6	64
4		Benzene, 1-(bromomethyl)-4-methyl-	184	C8H9Br	000104-81-4	64
5		Benzene, (1,3-dimethyl-3-butenyl)-	160	C12H16	056851-51-5	64



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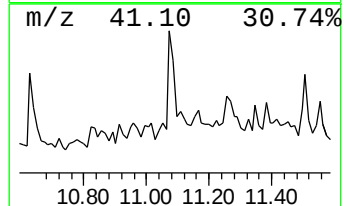
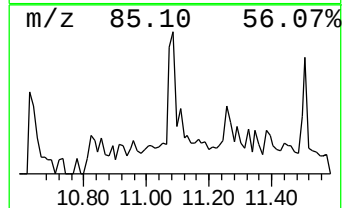
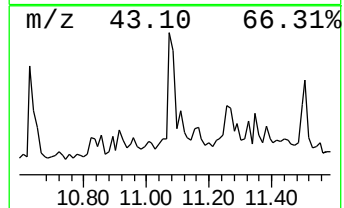
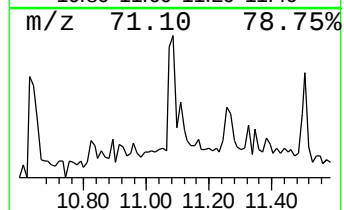
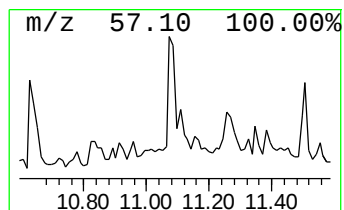
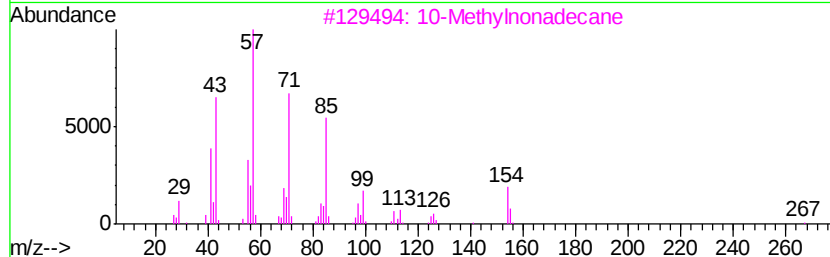
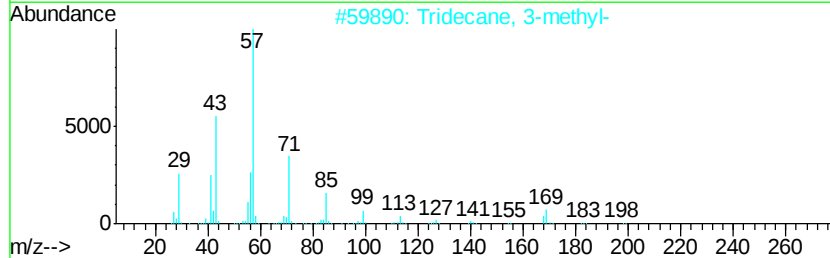
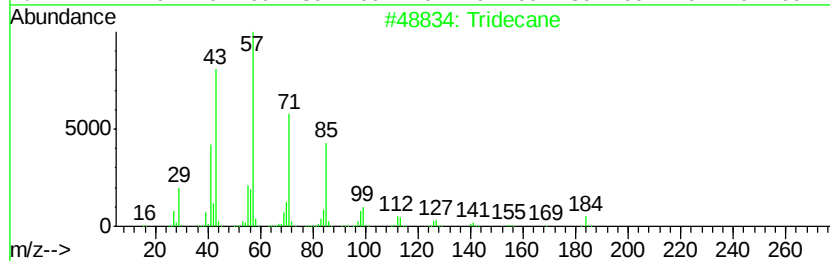
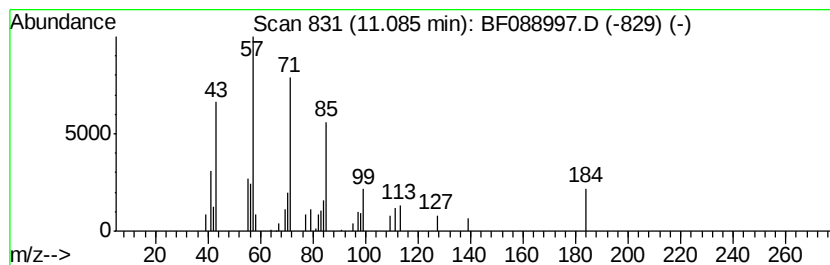
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Tridecane Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.09	2.04 ng	38480	Phenanthrene-d10	11.19	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecane	184	C13H28	000629-50-5	93
2	Tridecane, 3-methyl-	198	C14H30	006418-41-3	72
3	10-Methylnonadecane	282	C20H42	056862-62-5	72
4	Undecane, 2,3-dimethyl-	184	C13H28	017312-77-5	64
5	Heneicosane	296	C21H44	000629-94-7	64



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF071416\
Data File : BF088997.D
Acq On : 14 Jul 2016 20:50
Operator : UM/SJ
Sample : H3996-09
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
U2-B2(0-6)C

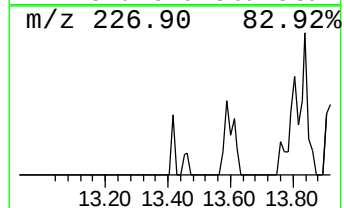
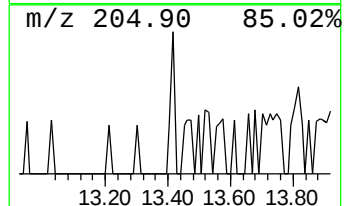
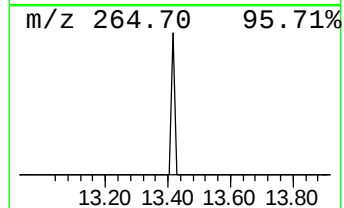
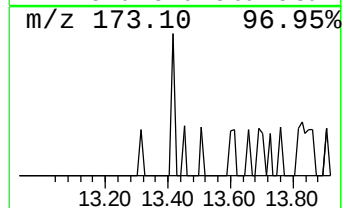
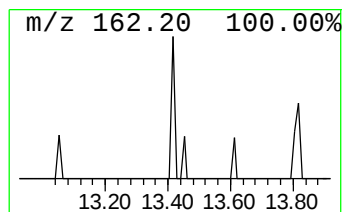
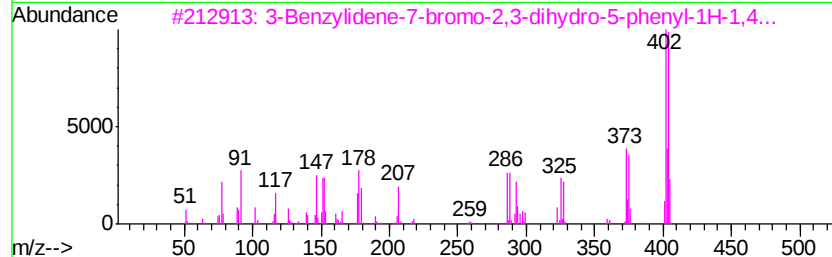
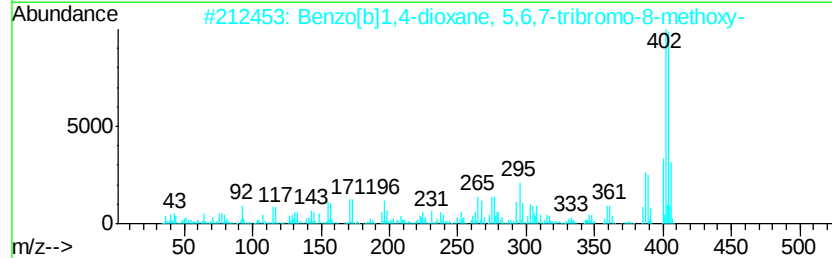
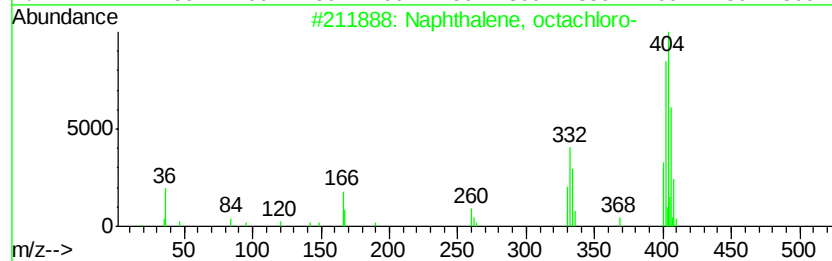
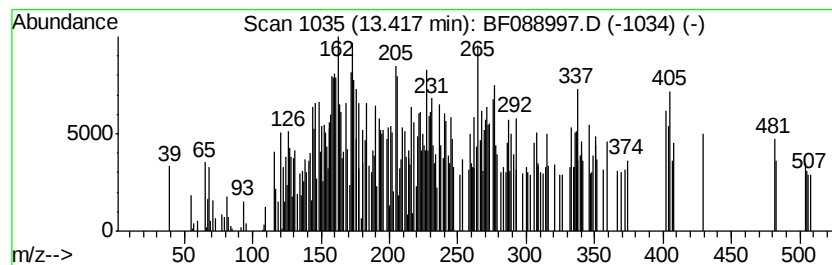
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 unknown13.42 Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.42	2.29 ng	67567	Chrysene-d12	13.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, octachloro-	400	C10Cl8	002234-13-1	10
2		Benzo[b]1,4-dioxane, 5,6,7-tribr...	400	C9H7Br3O3	273217-45-1	9
3		3-Benzylidene-7-bromo-2,3-dihvdr...	402	C22H15BrN2O	062642-82-4	9
4		2-Phenyl-4-(4-bromophenyl)-6-(2-...	402	C22H15BrN2O	313498-93-0	9
5		N-(3-Hydroxyphenyl)-2-(2,4,5-tri...	459	C20H24Cl3N03Si	1000332-17-7	9



Data Path : Z:\HPCHEM1\BNA F\DATA\BF071416\
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Operator : UM/SJ
Sample : H3996-09
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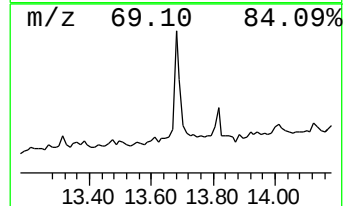
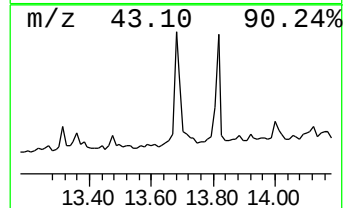
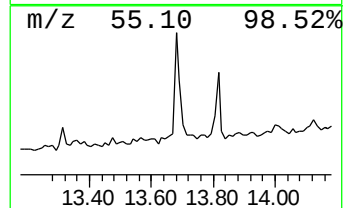
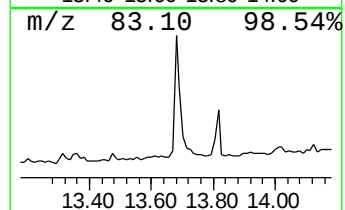
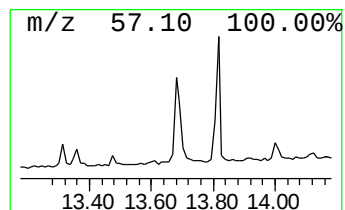
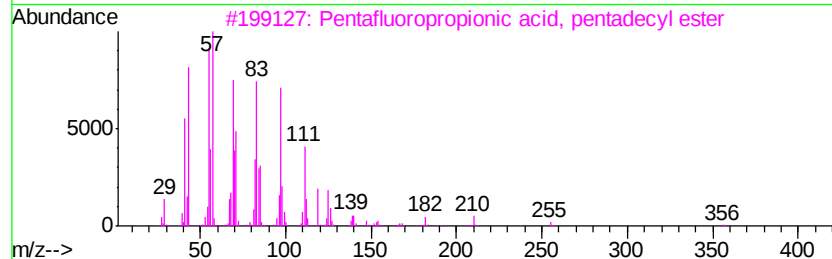
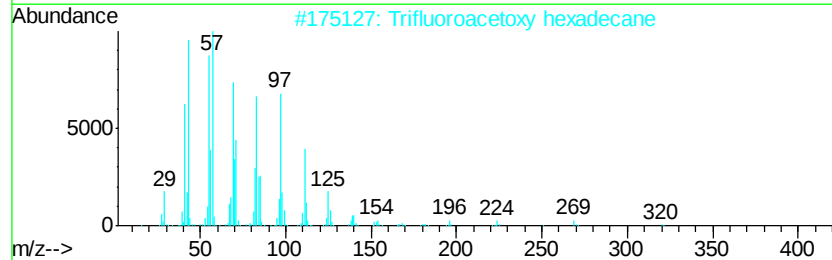
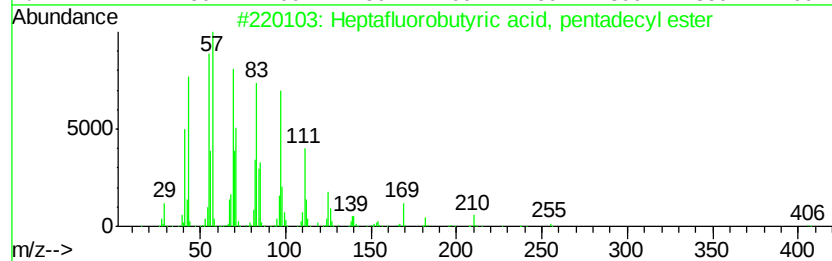
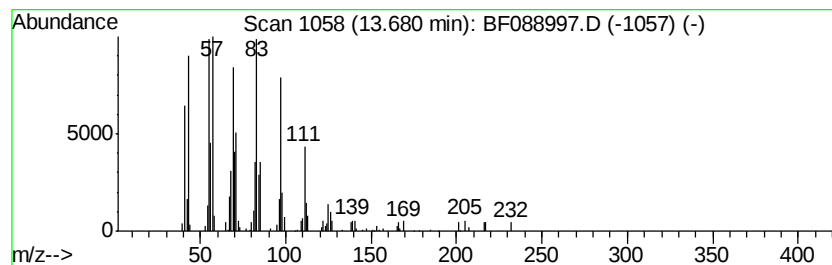
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Heptafluorobutyric acid, pe... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.68	3.62 ng	107124	Chrysene-d12	13.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	94
2		Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	94
3		Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	94
4		Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	94
5		Trifluoroacetic acid, pentadecyl...	324	C17H31F3O2	959010-23-2	93



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF071416\
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Acq On : 14 Jul 2016 20:50
Operator : UM/SJ
Sample : H3996-09
Misc :
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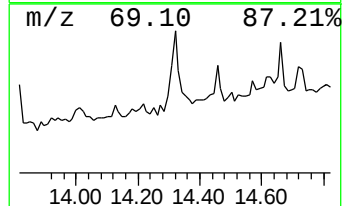
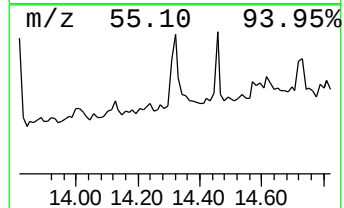
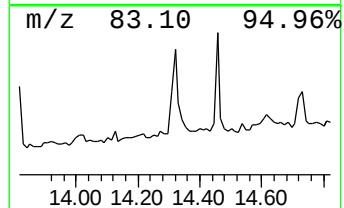
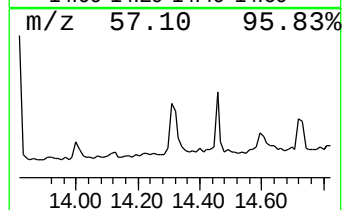
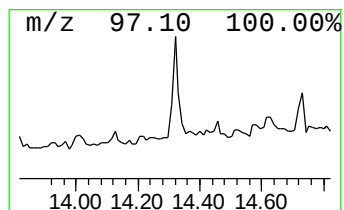
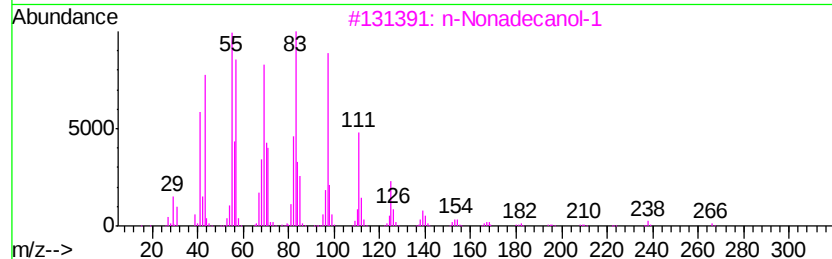
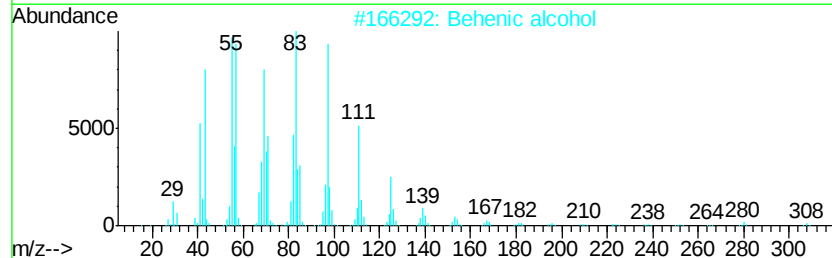
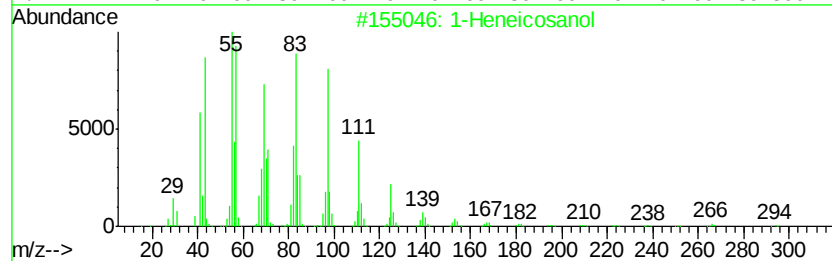
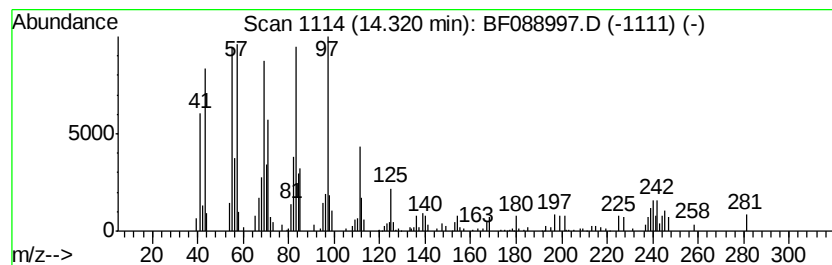
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 1-Heneicosanol Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.		
14.32	3.92 ng	115880	Chrysene-d12	13.82		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Heneicosanol	312	C21H44O	015594-90-8	93
2		Behenic alcohol	326	C22H46O	000661-19-8	93
3		n-Nonadecanol-1	284	C19H40O	001454-84-8	93
4		Trifluoroacetoxv hexadecane	338	C18H33F3O2	006222-03-3	91
5		Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF071416\
Data File : BF088997.D
Acq On : 14 Jul 2016 20:50
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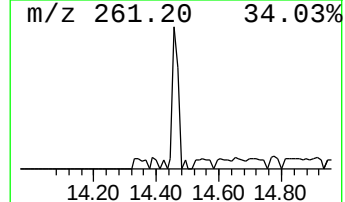
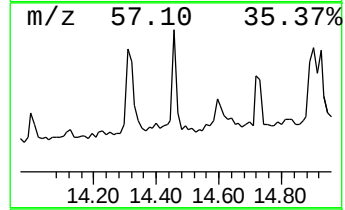
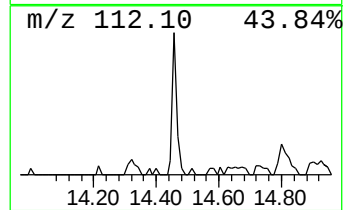
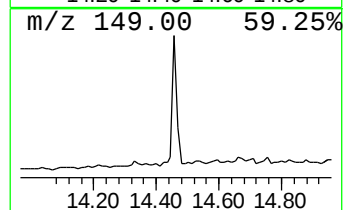
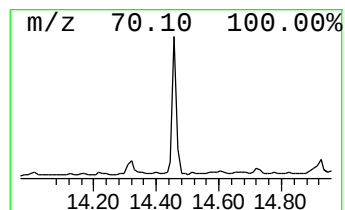
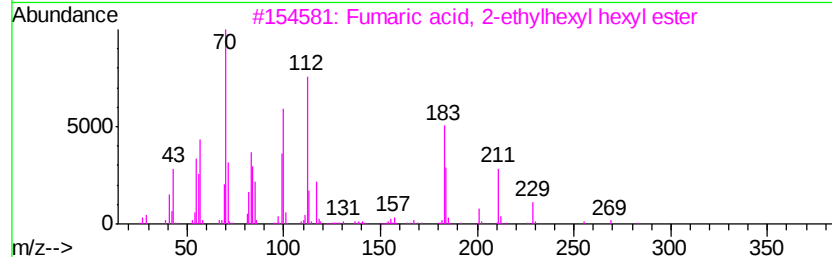
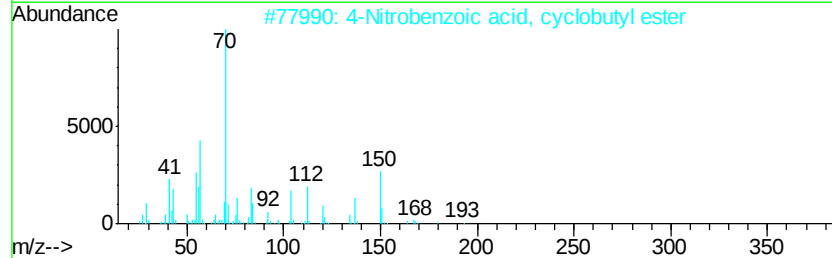
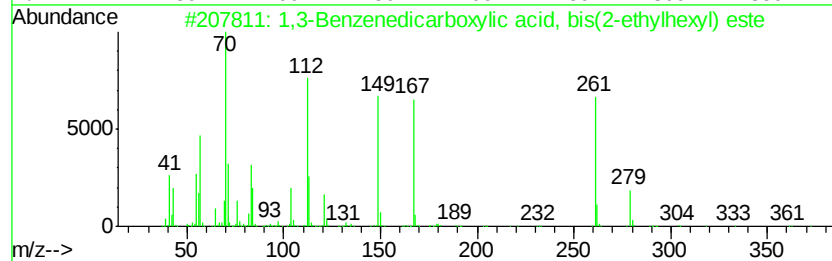
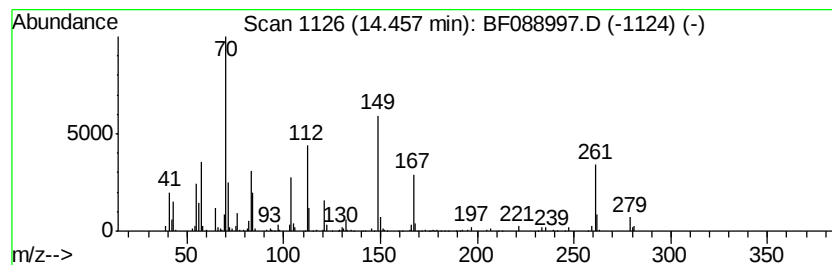
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 unknown14.46 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.46	3.01 ng	88982	Chrysene-d12	13.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,3-Benzenedicarboxylic acid, bi...	390	C24H38O4	000137-89-3	35
2		4-Nitrobenzoic acid, cyclobutyl ...	221	C11H11NO4	070335-00-1	32
3		Fumaric acid, 2-ethylhexyl hexyl...	312	C18H32O4	1000339-14-1	22
4		Diisooctyl phthalate	390	C24H38O4	000131-20-4	22
5		Carbonic acid, propargyl 2-ethyl...	212	C12H20O3	1000357-90-9	16



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF071416\
Data File : BF088997.D
Acq On : 14 Jul 2016 20:50
Operator : UM/SJ
Sample : H3996-09
Misc :
ALS Vial : 9 Sample Multiplier: 1

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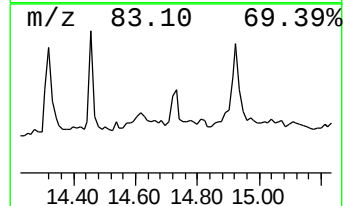
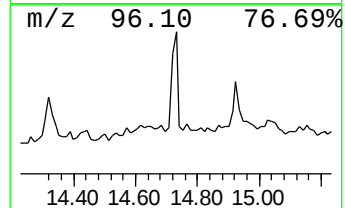
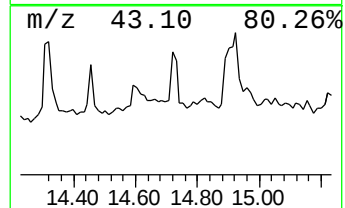
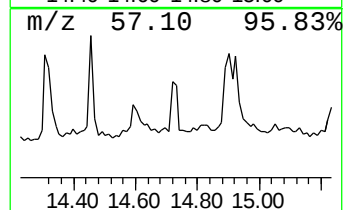
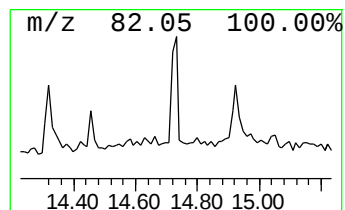
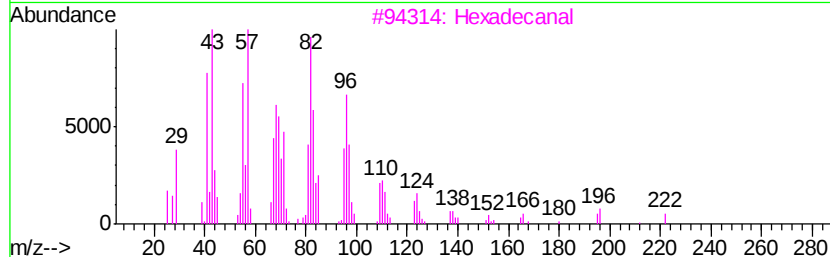
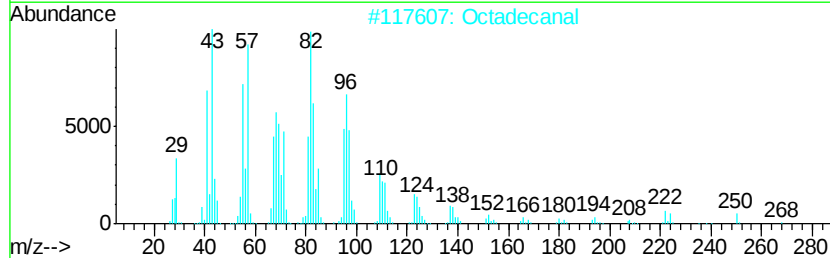
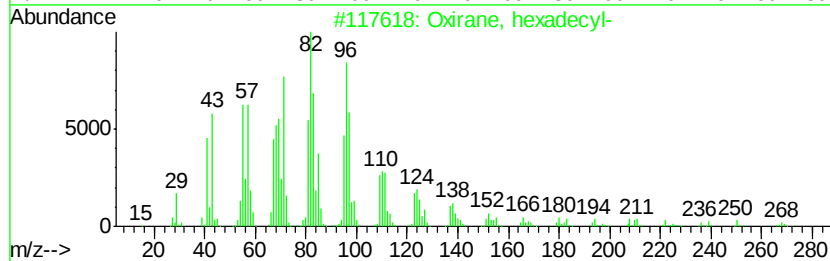
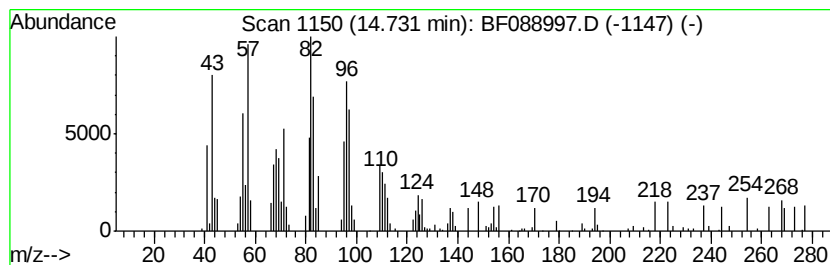
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Oxirane, hexadecyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.73	3.68 ng	47503	Perylene-d12	15.18

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Oxirane, hexadecyl-	268	C18H36O	007390-81-0	74
2			Octadecanal	268	C18H36O	000638-66-4	72
3			Hexadecanal	240	C16H32O	000629-80-1	64
4			1,16-Hexadecanediol	258	C16H34O2	007735-42-4	58
5			Cyclododecanol	184	C12H24O	001724-39-6	58



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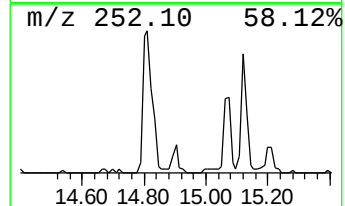
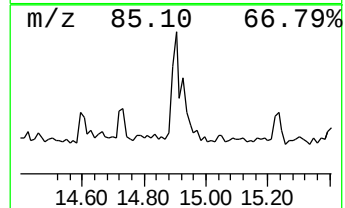
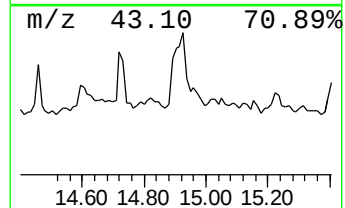
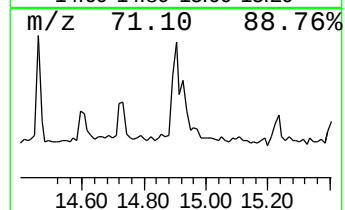
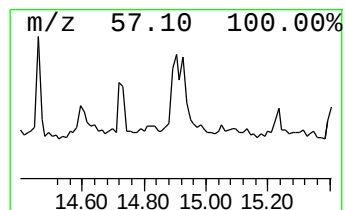
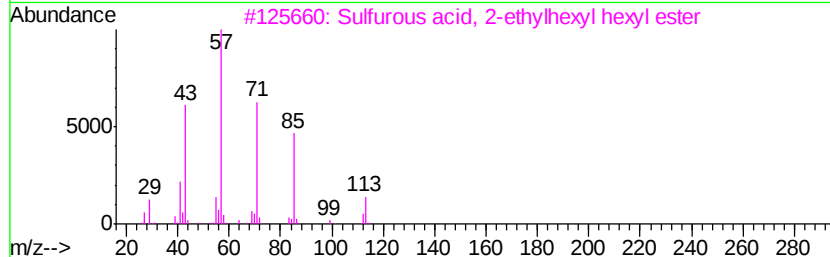
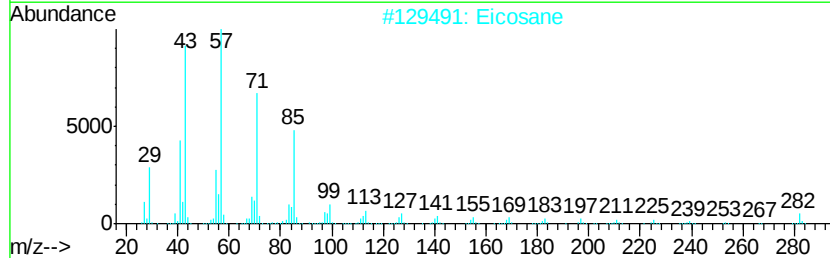
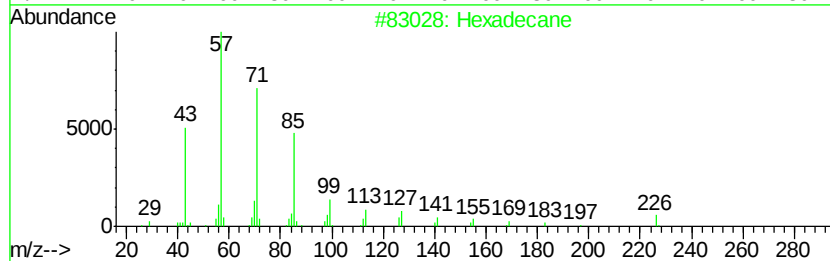
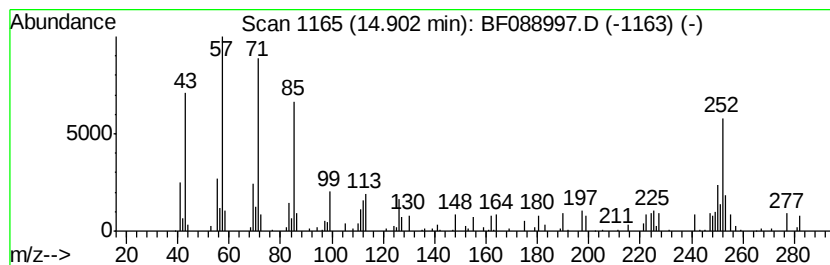
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Hexadecane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.90	3.50 ng	45191	Perylene-d12	15.18	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	53
2	Eicosane	282	C20H42	000112-95-8	49
3	Sulfurous acid, 2-ethylhexyl hex...	278	C14H30O3S	1000309-20-2	43
4	Sulfurous acid, nonyl 2-propyl e...	250	C12H26O3S	1000309-12-0	35
5	Sulfurous acid, 2-ethylhexyl tri...	376	C21H44O3S	1000309-19-6	35



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF071416\
Data File : BF088997.D
Acq On : 14 Jul 2016 20:50
Operator : UM/SJ
Sample : H3996-09
Misc :
ALS Vial : 9 Sample Multiplier: 1

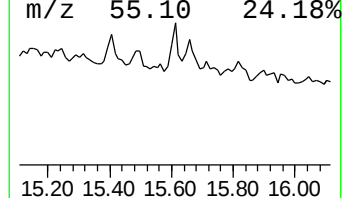
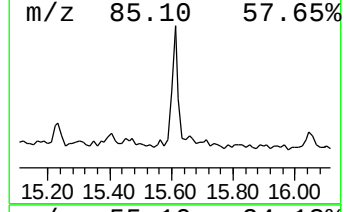
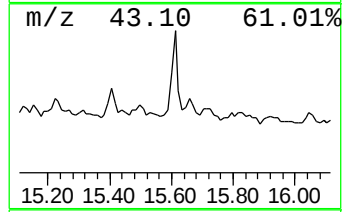
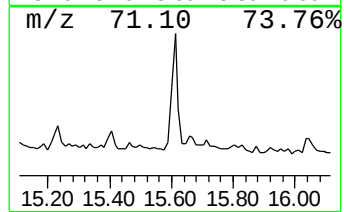
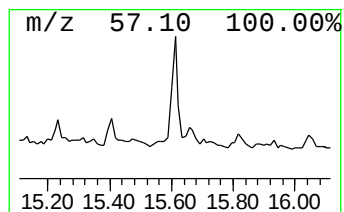
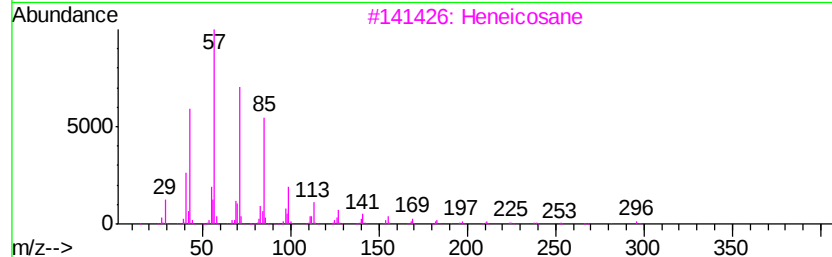
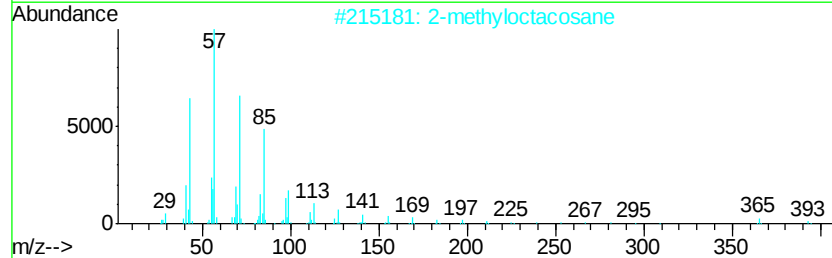
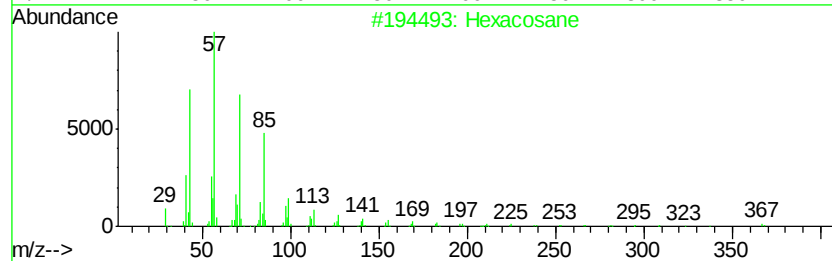
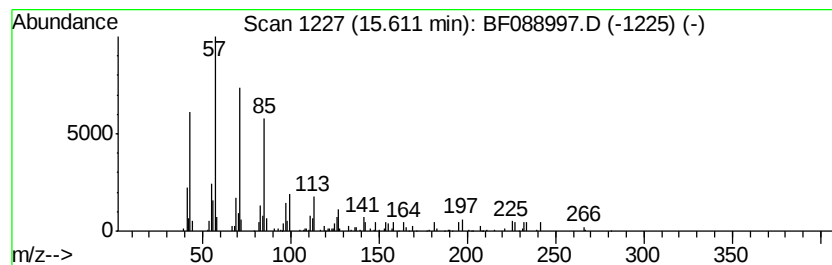
Instrument :
BNA_F
ClientSampleId :
U2-B2(0-6)C

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 Hexacosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD		R.T.
15.61	4.39 ng	56799	Perylene-d12		15.18
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexacosane	366	C26H54	000630-01-3	87
2	2-methyloctacosane	408	C29H60	1000376-72-8	87
3	Heneicosane	296	C21H44	000629-94-7	87
4	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	87
5	2-Bromotetradecane	276	C14H29Br	074036-95-6	86



Data Path : Z:\HPCHEM1\BNA F\DATA\BF071416\
Data File : BF088997.D
Acq On : 14 Jul 2016 20:50
Operator : UM/SJ
Sample : H3996-09
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
U2-B2(0-6)C

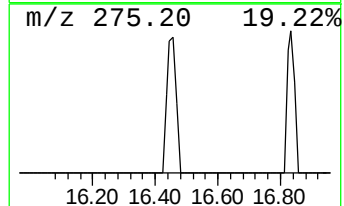
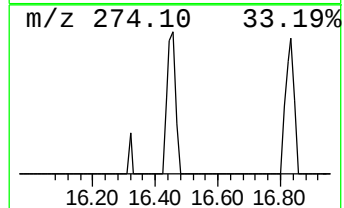
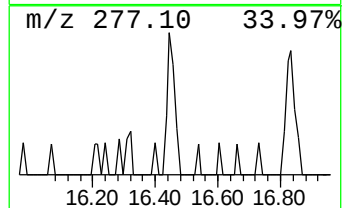
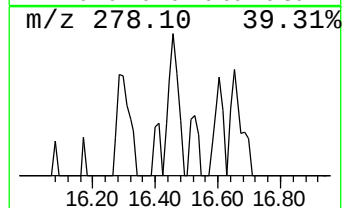
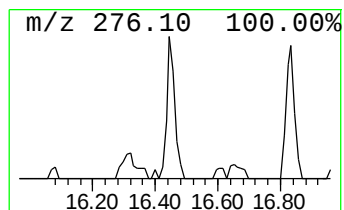
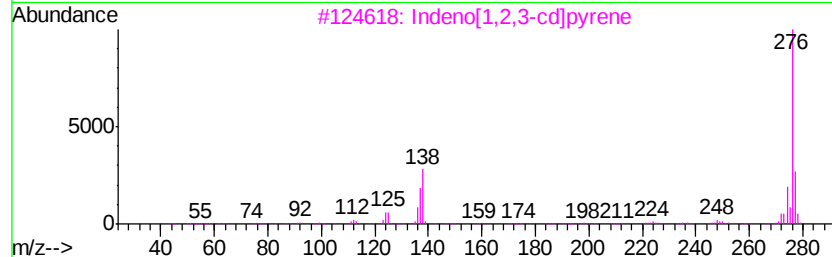
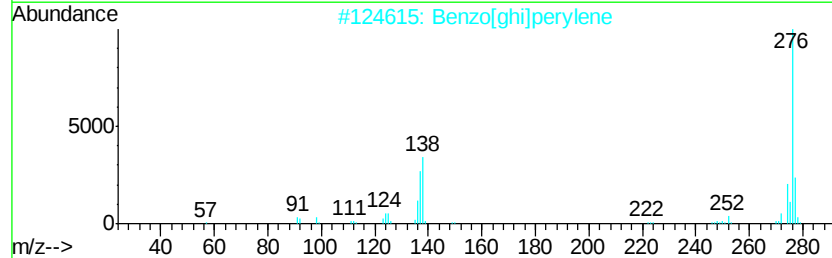
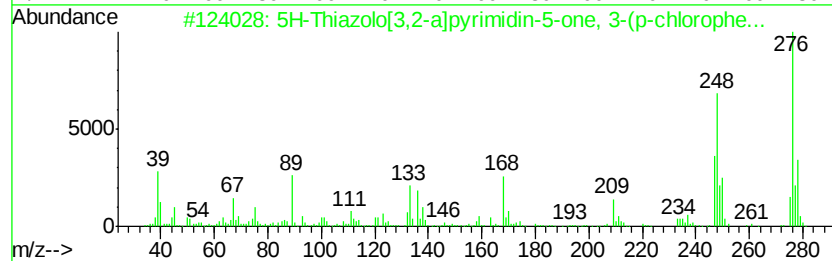
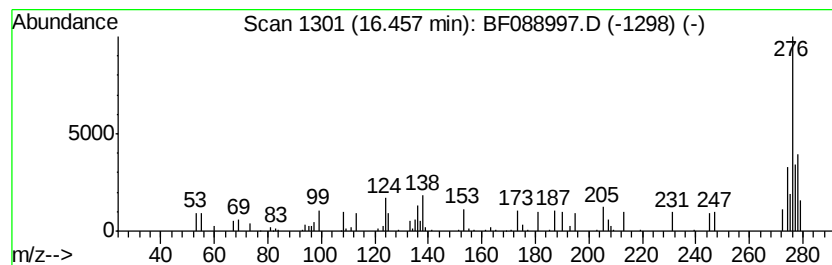
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 unknown16.46 Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.46	2.24 ng	28994	Perylene-d12	15.18

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	5H-Thiazolo[3,2-a]pvrimidin-5-on...	276	C13H9ClN2OS	023429-91-6	45		
2	Benzo[ghi]perylene	276	C22H12	000191-24-2	43		
3	Indeno[1,2,3-cd]pyrene	276	C22H12	000193-39-5	43		
4	Dibenzo[def,mno]chrysene	276	C22H12	000191-26-4	43		
5	5-(3,4-Dimethoxy-benzylidene)-py...	276	C13H12N2O5	1000318-16-4	43		



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF071416\
Data File : BF088997.D
Acq On : 14 Jul 2016 20:50
Operator : UM/SJ
Sample : H3996-09
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
U2-B2(0-6)C

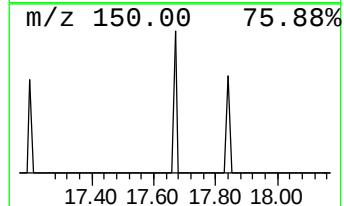
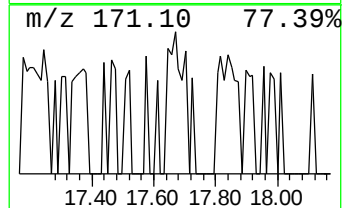
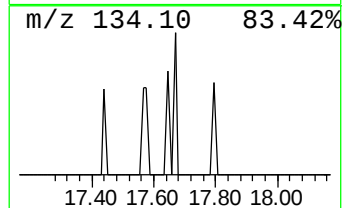
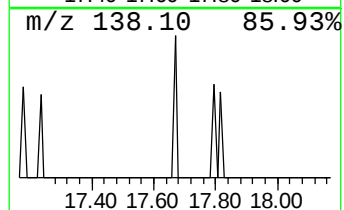
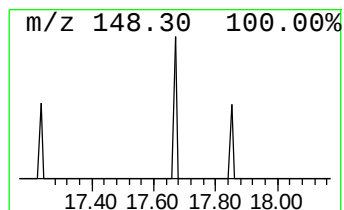
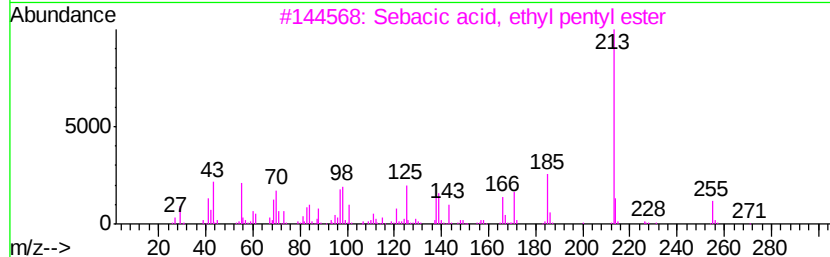
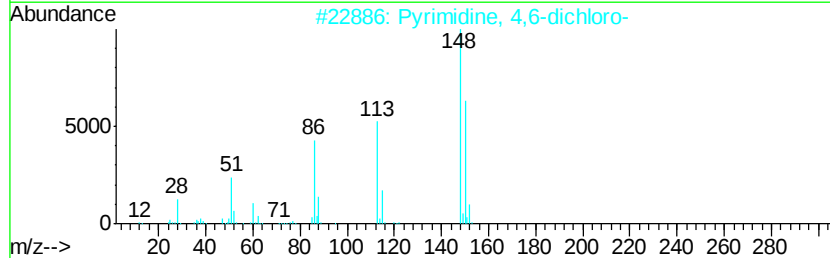
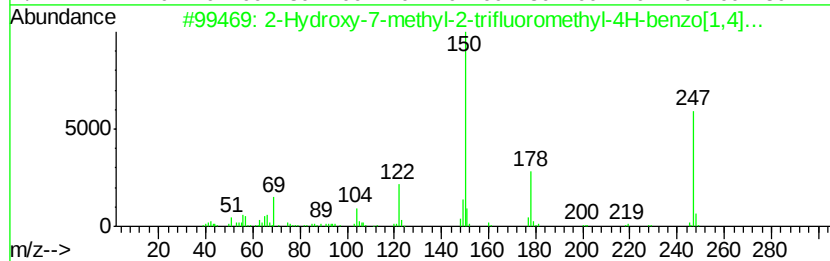
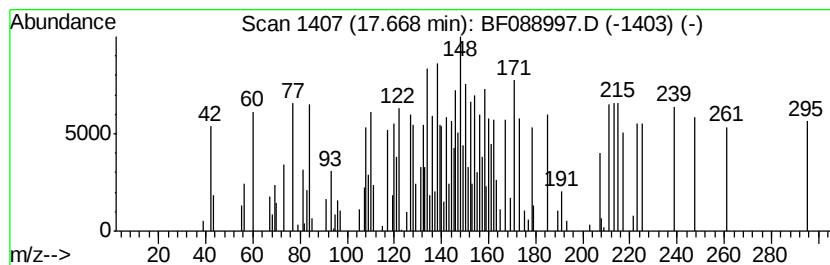
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF063016.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 unknown17.67 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.67	3.32 ng	42864	Perylene-d12	15.18

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Hydroxy-7-methyl-2-trifluorome...	247	C10H8F3NO3	1000306-79-2	9
2		Pyrimidine, 4,6-dichloro-	148	C4H2Cl2N2	001193-21-1	9
3		Sebacic acid, ethyl pentyl ester	300	C17H32O4	1000355-90-4	9
4		2-[3-Hydroxypropylamino]quinoline	202	C12H14N2O	092959-28-9	9
5		Methanamine, (3-chloro-4-ethoxyp...	185	C9H12ClNO	007569-87-1	9



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF071416\
Data File : BF088997.D
Acq On : 14 Jul 2016 20:50
Operator : UM/SJ
Sample : H3996-09
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
U2-B2(0-6)C

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...	1.86	21.0	ng	371151	1	6.68	353294	20.0
2-Pentanone, 4-hy...	4.83	10.1	ng	178841	1	6.68	353294	20.0
unknown6.46	6.46	87.5	ng	1545580	1	6.68	353294	20.0
Benzene, (1-bromo...	7.55	2.7	ng	55637	2	7.96	413491	20.0
Tridecane	11.09	2.0	ng	38480	4	11.19	377383	20.0
unknown13.42	13.42	2.3	ng	67567	5	13.82	591357	20.0
Heptafluorobutyri...	13.68	3.6	ng	107124	5	13.82	591357	20.0
1-Heneicosanol	14.32	3.9	ng	115880	5	13.82	591357	20.0
unknown14.46	14.46	3.0	ng	88982	5	13.82	591357	20.0
Oxirane, hexadecyl-	14.73	3.7	ng	47503	6	15.18	258490	20.0
Hexadecane	14.90	3.5	ng	45191	6	15.18	258490	20.0
Hexacosane	15.61	4.4	ng	56799	6	15.18	258490	20.0
unknown16.46	16.46	2.2	ng	28994	6	15.18	258490	20.0
unknown17.67	17.67	3.3	ng	42864	6	15.18	258490	20.0