

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF033016\
 Data File : BF085879.D
 Acq On : 30 Mar 2016 11:28
 Operator : UM/SJ
 Sample : H1937-02
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FIELDBLANK-3-23-16

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-BF032416.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.344	4	5	6	rBV	51576	39530	1.07%	0.149%
2	2.378	6	8	9	rVV2	82806	103512	2.81%	0.391%
3	2.413	9	11	15	rVB4	46255	76916	2.09%	0.291%
4	2.516	18	20	21	rBV2	66374	89753	2.44%	0.339%
5	2.538	21	22	25	rVV3	47663	77138	2.10%	0.291%
6	2.630	29	30	32	rVB	57567	70198	1.91%	0.265%
7	2.744	37	40	42	rBV3	149843	202578	5.51%	0.765%
8	2.813	44	46	48	rVB3	160201	213148	5.80%	0.805%
9	2.893	52	53	60	rVB5	31429	60720	1.65%	0.229%
10	3.144	70	75	77	rVB4	69597	151070	4.11%	0.571%
11	3.327	88	91	92	rBV3	20168	40053	1.09%	0.151%
12	3.361	92	94	95	rVV	50284	55800	1.52%	0.211%
13	3.384	95	96	101	rVB2	133458	137370	3.74%	0.519%
14	3.544	104	110	111	rBV6	23330	51983	1.41%	0.196%
15	3.590	111	114	117	rBV3	28386	46982	1.28%	0.178%
16	3.704	122	124	128	rVB3	91360	131225	3.57%	0.496%
17	3.773	128	130	131	rBV	94460	82419	2.24%	0.311%
18	3.887	135	140	142	rBV3	15973	47279	1.29%	0.179%
19	3.967	146	147	150	rVV2	23801	46877	1.27%	0.177%
20	4.059	153	155	157	rBV2	35513	49447	1.34%	0.187%
21	4.116	157	160	163	rVB4	35250	53015	1.44%	0.200%
22	4.310	172	177	180	rBV	32645	101361	2.76%	0.383%
23	4.379	180	183	186	rVV4	107036	171886	4.67%	0.650%
24	4.424	186	187	189	rVV2	91270	67275	1.83%	0.254%
25	4.516	192	195	197	rVV4	38368	80424	2.19%	0.304%
26	4.550	197	198	201	rVB	103565	134908	3.67%	0.510%
27	4.824	220	222	223	rBV2	72376	60065	1.63%	0.227%
28	4.962	232	234	238	rBV	354807	444526	12.09%	1.680%
29	5.396	270	272	274	rBV	1054409	1272901	34.61%	4.810%
30	5.453	276	277	280	rVV2	44761	55460	1.51%	0.210%
31	5.590	284	289	292	rVB5	18282	38100	1.04%	0.144%
32	5.670	292	296	300	rBV5	46190	90280	2.45%	0.341%
33	5.796	305	307	312	rVB	149398	136914	3.72%	0.517%
34	5.922	314	318	319	rBV3	42755	49644	1.35%	0.188%

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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

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Peak separation: 5

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35	6.436	360	363	365 rBV	757667	800298	21.76%	3.024%
36	6.516	368	370	372 rVV3	107030	222480	6.05%	0.841%
37	6.562	372	374	377 rVB	2209428	2400659	65.28%	9.071%
38	6.802	393	395	397 rBV	470616	619118	16.84%	2.339%
39	6.893	401	403	406 rVB3	77361	124447	3.38%	0.470%
40	6.950	406	408	411 rBV	2147013	2236514	60.82%	8.451%
41	7.247	432	434	438 rVB3	40627	56428	1.53%	0.213%
42	7.362	441	444	447 rBV	1534967	1853651	50.41%	7.004%
43	7.853	482	487	490 rBV	64045	98125	2.67%	0.371%
44	8.082	505	507	510 rVB	957925	837720	22.78%	3.165%
45	8.333	526	529	530 rBV3	28570	43095	1.17%	0.163%
46	8.368	530	532	535 rVV2	49946	85421	2.32%	0.323%
47	8.425	535	537	540 rVB3	56193	50422	1.37%	0.191%
48	9.168	599	602	604 rBV	3054003	3677498	100.00%	13.896%
49	9.842	658	661	663 rBV	737254	885596	24.08%	3.346%
50	10.631	727	730	732 rBV	2142005	2111209	57.41%	7.978%
51	11.316	788	790	793 rBV	920826	901122	24.50%	3.405%
52	12.905	926	929	932 rBV	2796411	3453215	93.90%	13.049%
53	13.957	1018	1021	1023 rBV	932535	843581	22.94%	3.188%
54	15.374	1142	1145	1151 rBV	495195	632885	17.21%	2.391%

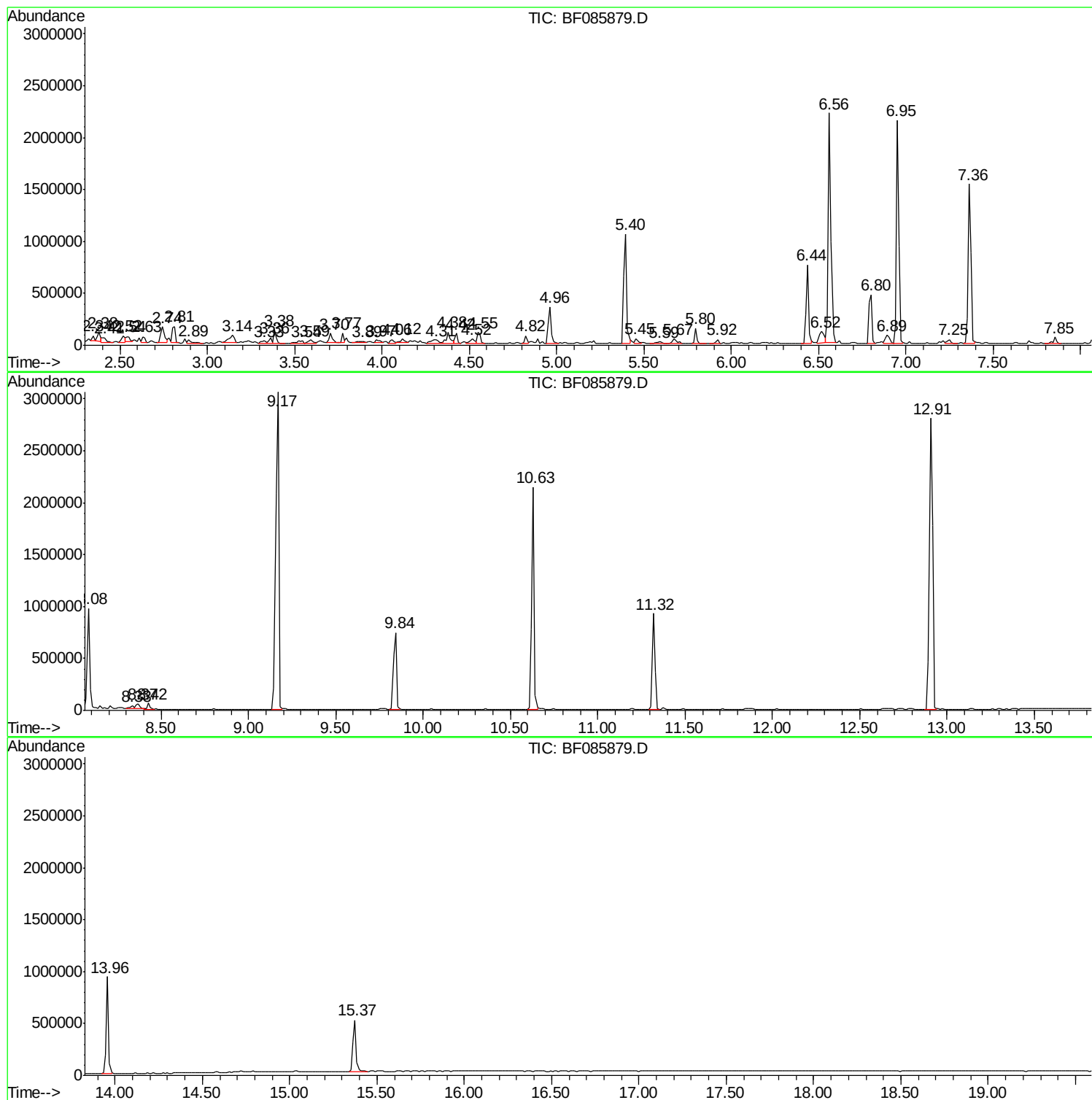
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032416.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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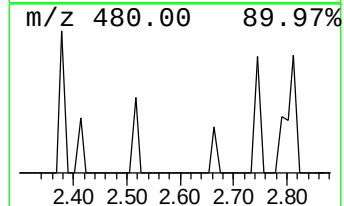
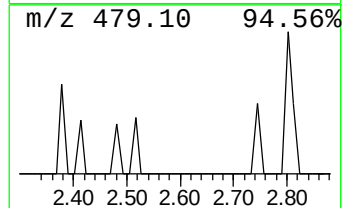
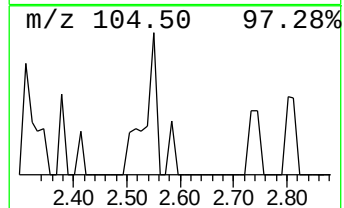
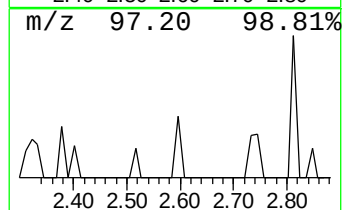
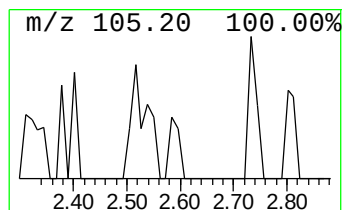
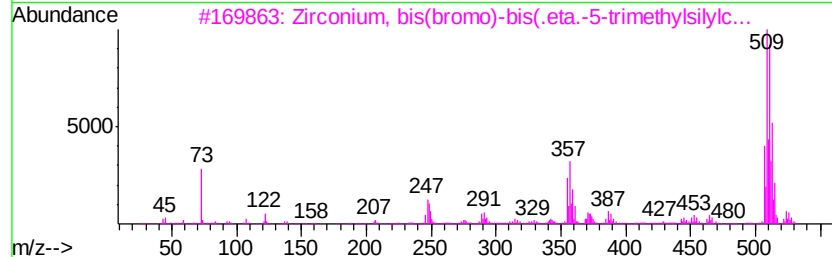
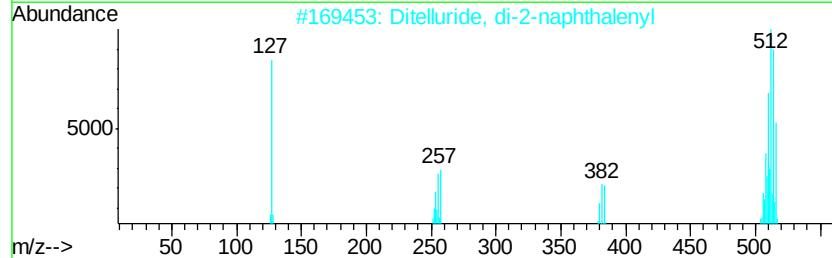
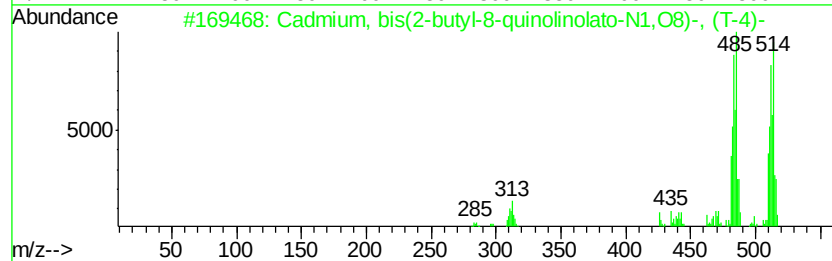
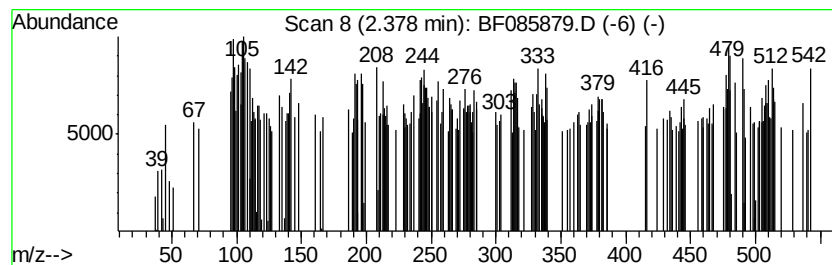
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032416.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown2.38 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.38	3.34 ng	103512	1,4-Dichlorobenzene-d4	6.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cadmium, bis(2-butyl-8-quinolino...	514	C26H28CdN2O2	015685-80-0	12
2		Ditelluride, di-2-naphthalenyl	514	C20H14Te2	001666-12-2	9
3		Zirconium, bis(bromo)-bis(.eta.-...	522	C16H26Br2Si2Zr	1000163-27-2	9
4		Cholestan-3-one, 2-bromo-, cycli...	508	C29H49BrO2	056052-89-2	7
5		5,10,15,20(22H,24H)-Porphinetetr...	540	C32H36N4O4	027800-02-8	4



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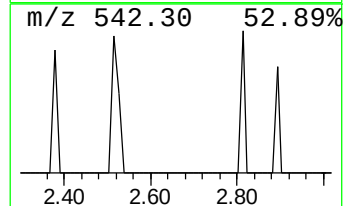
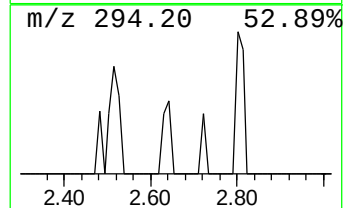
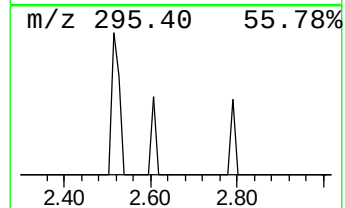
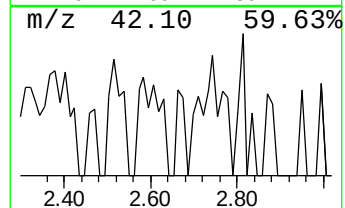
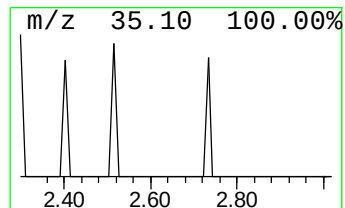
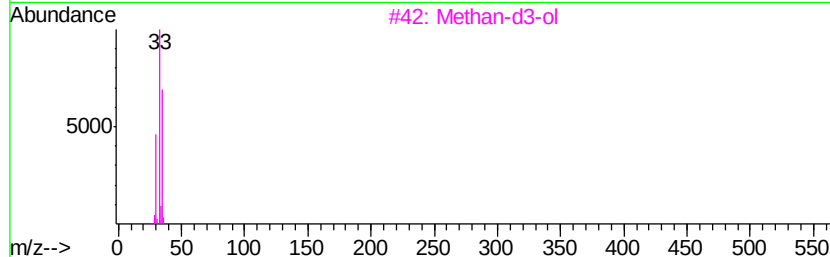
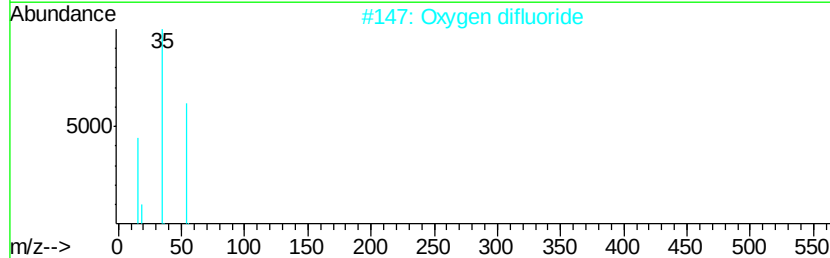
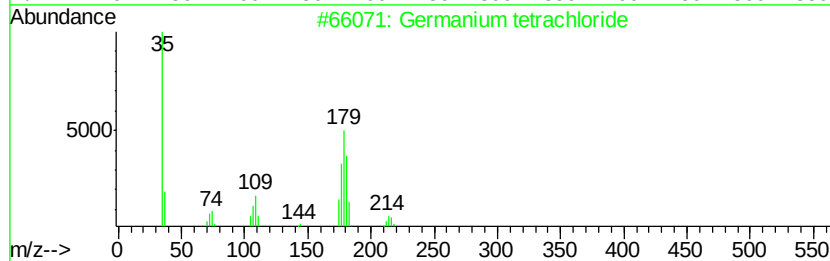
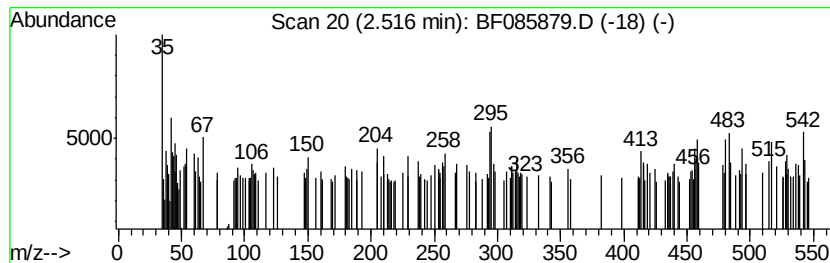
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown2.52 Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.52	2.90 ng	89753	1,4-Dichlorobenzene-d4	6.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Germanium tetrachloride	214	Cl4Ge	010038-98-9	16
2		Oxygen difluoride	54	F2O	007783-41-7	7
3		Methan-d3-ol	35	CHD3O	001849-29-2	3
4		Carbonochloridothioic acid, S-et...	124	C3H5ClOS	002941-64-2	2
5		2-Chloroethanol	80	C2H5ClO	000107-07-3	2



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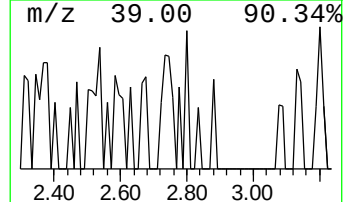
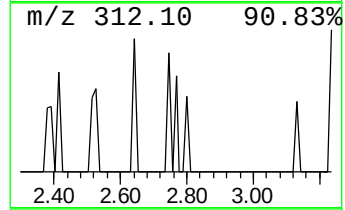
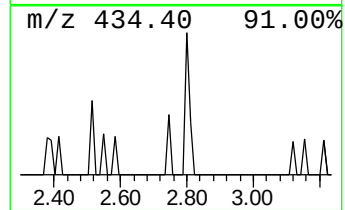
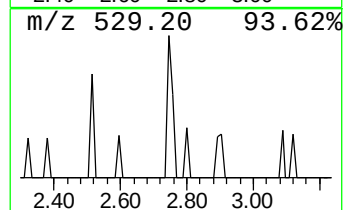
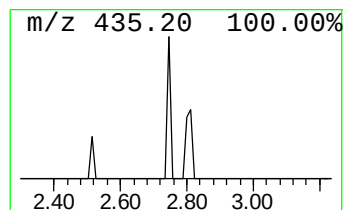
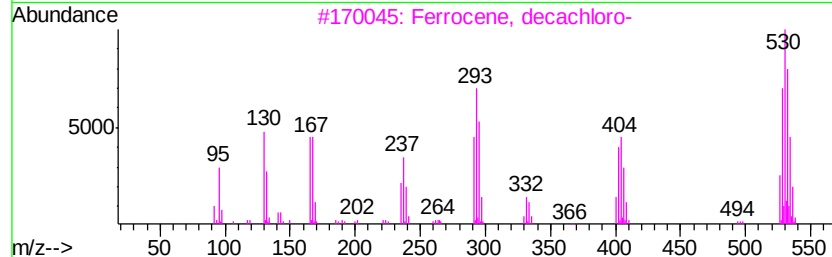
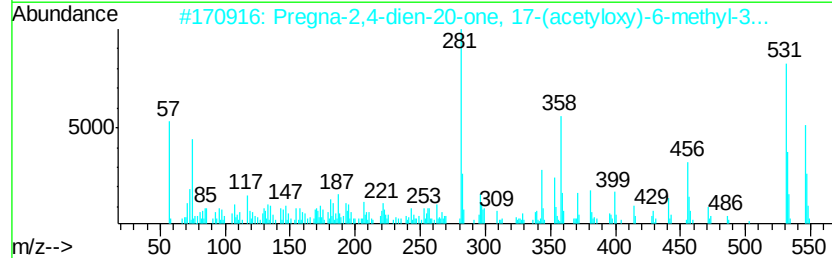
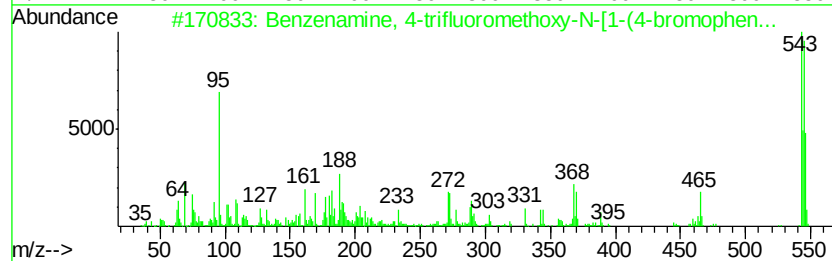
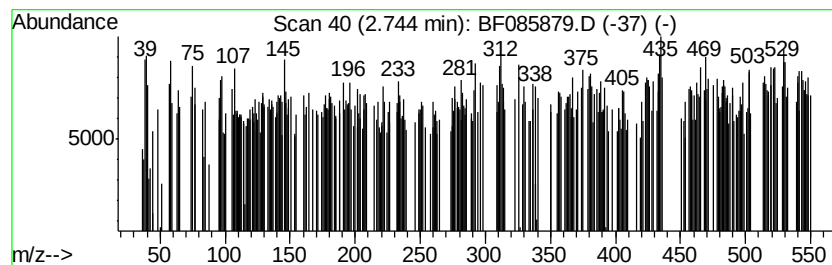
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 unknown2.74 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.74	6.54 ng	202578	1,4-Dichlorobenzene-d4	6.80

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzenamine, 4-trifluoromethoxy-...	544	C23H15BrF6N2O2	284495-96-1	20
2			Pregna-2,4-dien-20-one, 17-(acet...	546	C30H50O5Si2	069833-57-4	18
3			Ferrocene, decachloro-	526	C10Cl10Fe	011121-63-4	15
4			Distannoxane, hexabutyl-	598	C24H54O2Sn2	000056-35-9	10
5			Mercury iodo, (o-iodophenyl)-	532	C6H4HgI2	089487-89-8	10



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ClientSampled :
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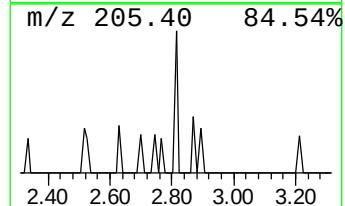
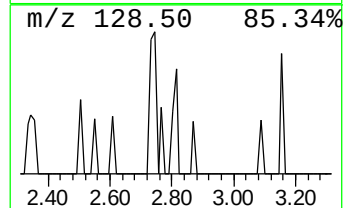
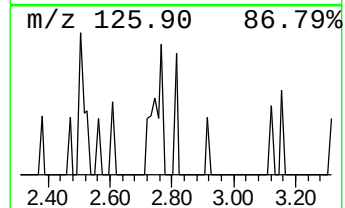
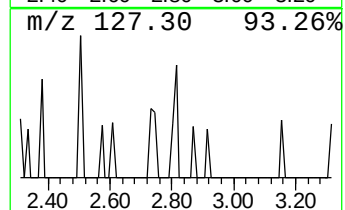
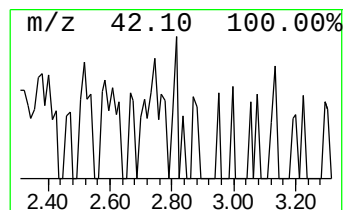
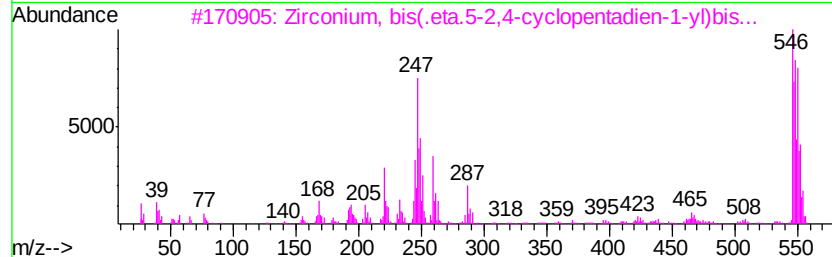
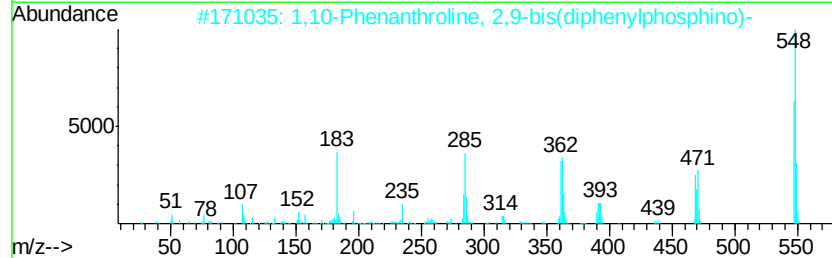
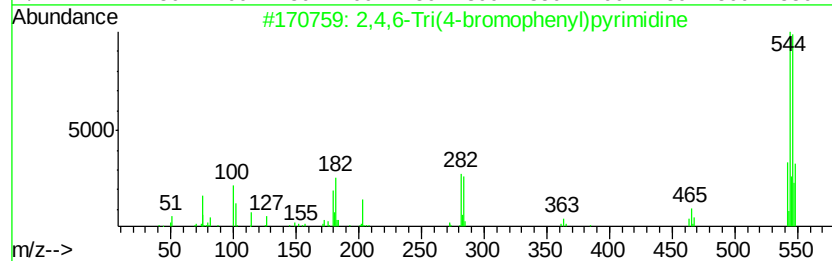
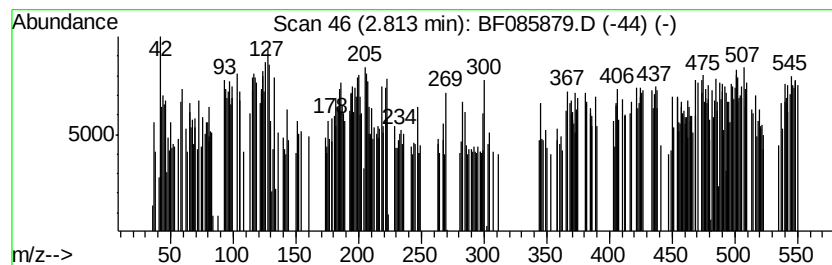
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 unknown2.81 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.81	6.89 ng	213148	1,4-Dichlorobenzene-d4	6.80

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,6-Tri(4-bromophenyl)pyrimidine	542	C22H13Br3N2	1000070-62-6	12		
2	1,10-Phenanthroline, 2,9-bis(dip...	548	C36H26N2P2	122768-46-1	11		
3	Zirconium, bis(.eta.5-2,4-cyclop...	546	C28H30Zr2	119144-91-1	10		
4	Iron, (.eta.5-2,4-cyclopentadien...	520	C31H30FeP2	051509-20-7	9		
5	Copper, [2,8,12,18-tetraethyl-3,...	539	C32H36CuN4	014055-18-6	4		



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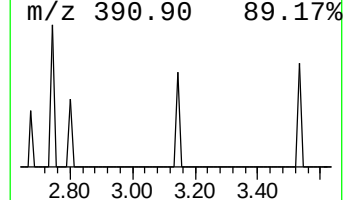
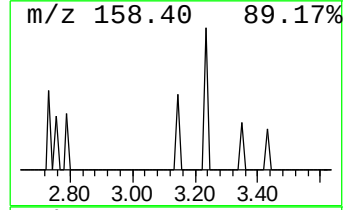
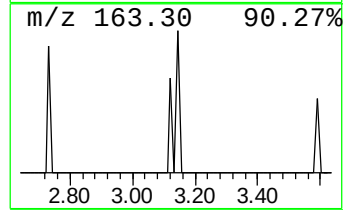
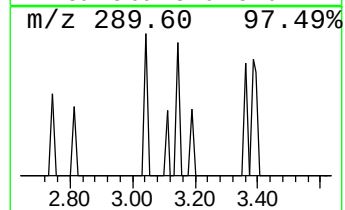
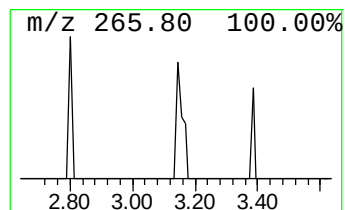
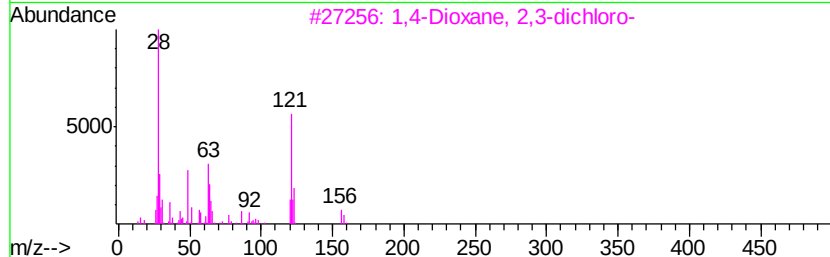
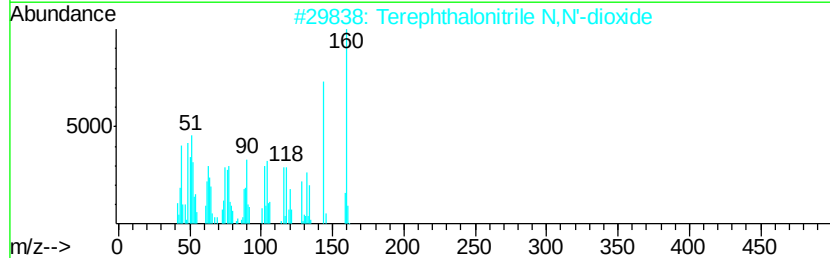
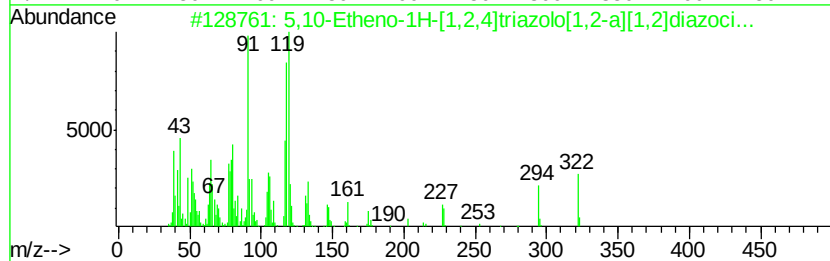
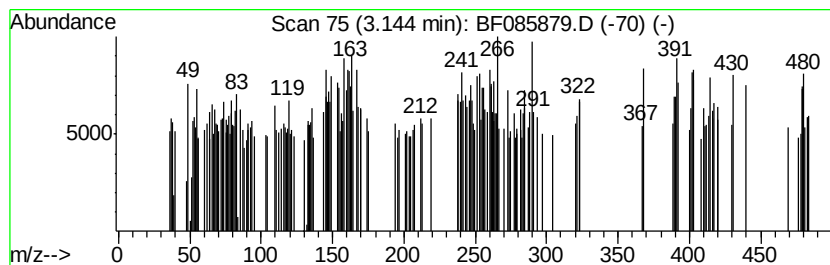
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 unknown3.14 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.14	4.88 ng	151070	1,4-Dichlorobenzene-d4	6.80

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	5,10-Etheno-1H-[1,2,4]triazolo[1...	322	C16H14N6O2	099140-99-5	20		
2	Terephthalonitrile N,N'-dioxide	160	C8H4N2O2	003729-34-8	20		
3	1,4-Dioxane, 2,3-dichloro-	156	C4H6Cl2O2	000095-59-0	14		
4	4,6-Dibromo-2-benzoxazolinone	291	C7H3Br2N02	1000260-92-6	14		
5	4-Bromo-6-chloro-2-benzoxazolinone	247	C7H3BrClN02	313503-21-8	14		



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF033016\
Data File : BF085879.D
Acq On : 30 Mar 2016 11:28
Operator : UM/SJ
Sample : H1937-02
Misc :
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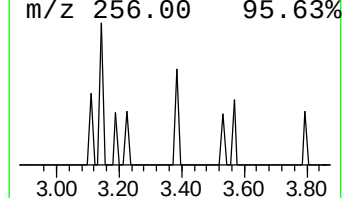
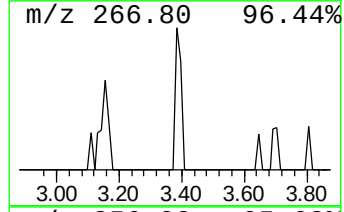
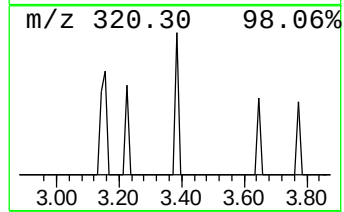
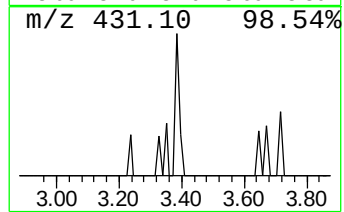
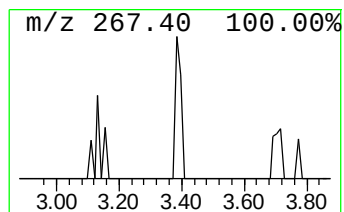
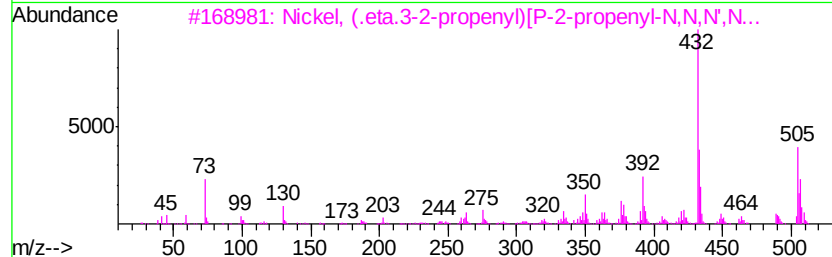
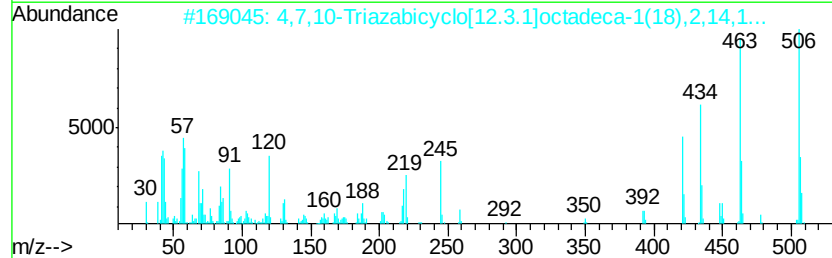
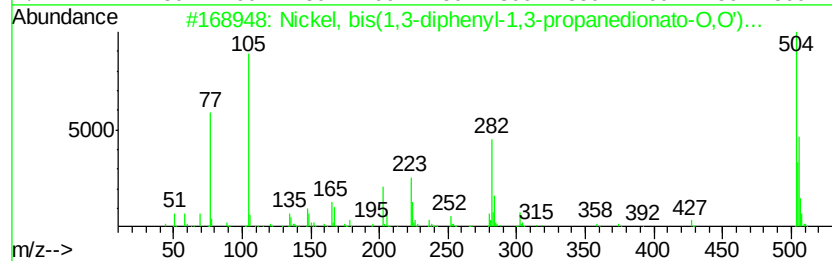
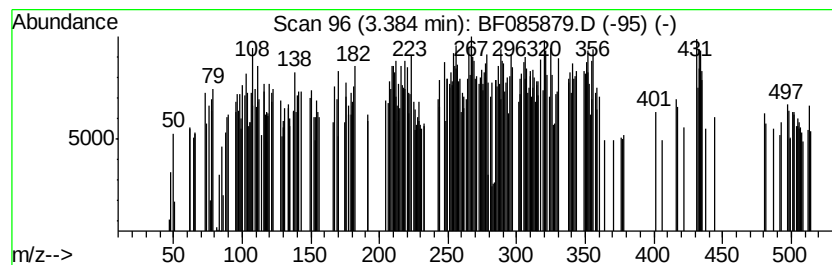
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 unknown3.38 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.38	4.44 ng	137370	1,4-Dichlorobenzene-d4	6.80

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nickel, bis(1,3-diphenyl-1,3-pro...	504	C30H22NiO4	014405-47-1	10		
2	4,7,10-Triazabicyclo[12.3.1]octa...	506	C29H38N4O4	038840-25-4	10		
3	Nickel, (.eta.3-2-propenyl)[P-2-...	505	C18H46N3NiPSi4	075949-58-5	10		
4	4,4'-Di(4-chlorophenylsulfone)bi...	502	C24H16Cl2O4S2	022287-56-5	9		
5	Cobalt, bis(1,3-diphenyl-1,3-pro...	505	C30H22CoO4	014405-50-6	9		



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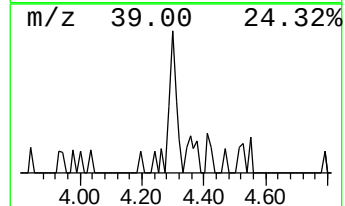
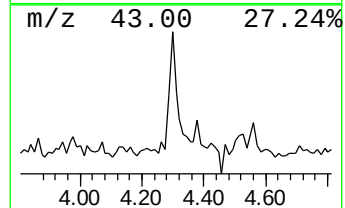
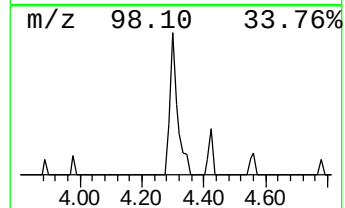
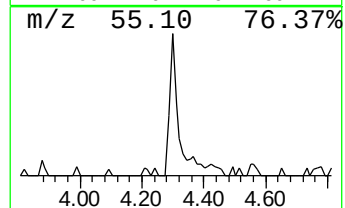
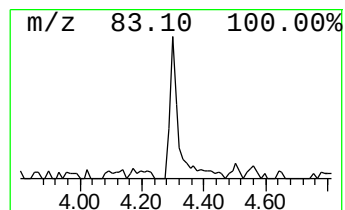
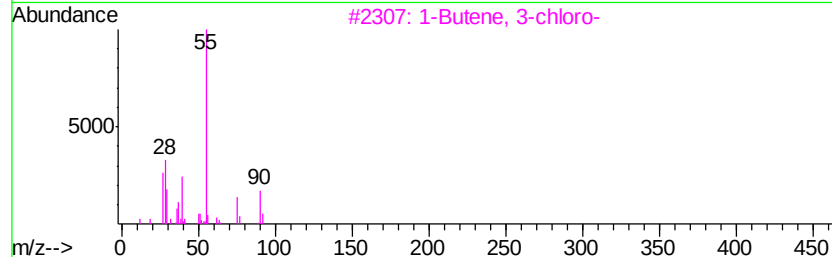
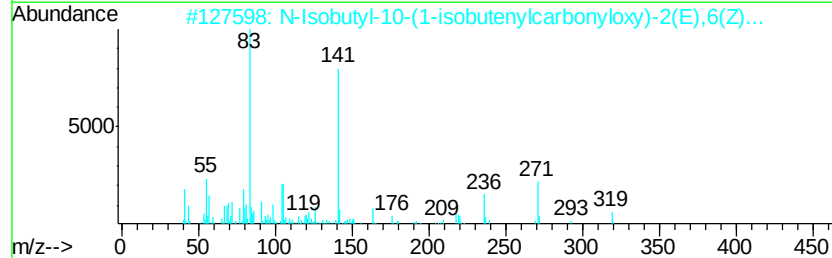
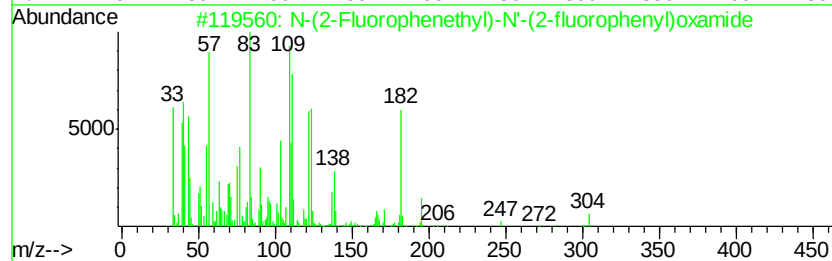
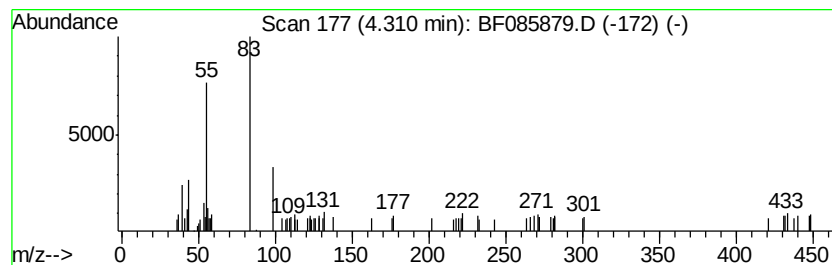
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 unknown4.31 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.31	3.27 ng	101361	1,4-Dichlorobenzene-d4	6.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	N-(2-Fluorophenethyl)-N'-(2-fluo...	304	C16H14F2N2O2	339006-39-2	35
2		N-Isobutyl-10-(1-isobutenylcarbo...	319	C19H29NO3	1000106-15-9	25
3		1-Butene, 3-chloro-	90	C4H7Cl	000563-52-0	9
4		Hexanal, (2,4-dinitrophenyl)hydr...	280	C12H16N4O4	001527-97-5	7
5		2,2'-Bioxirane, (R*,R*)-(./+/-.)-	86	C4H6O2	000298-18-0	2



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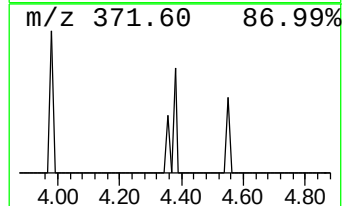
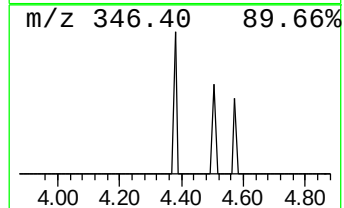
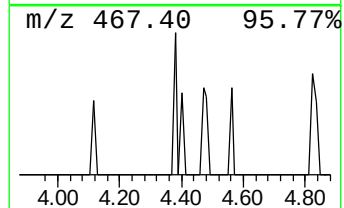
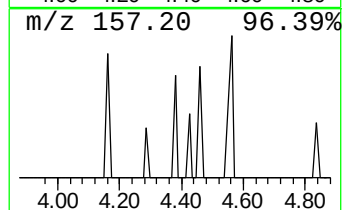
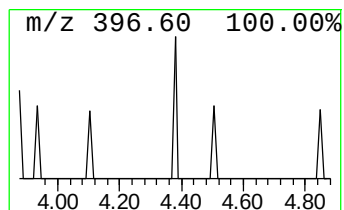
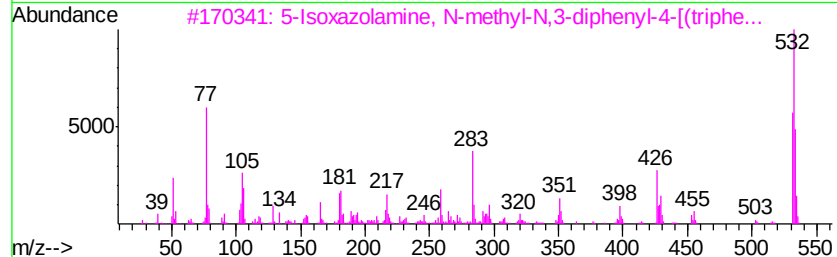
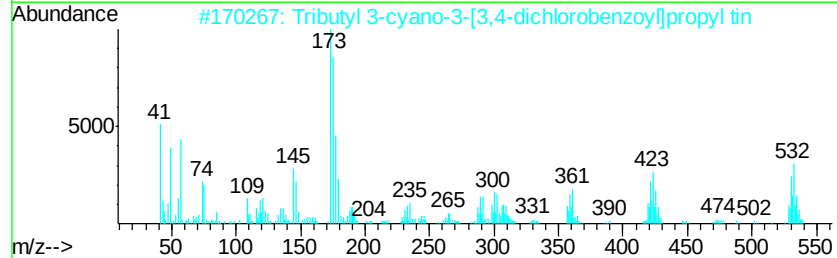
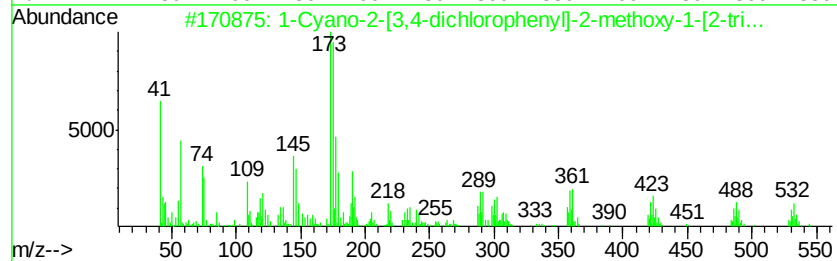
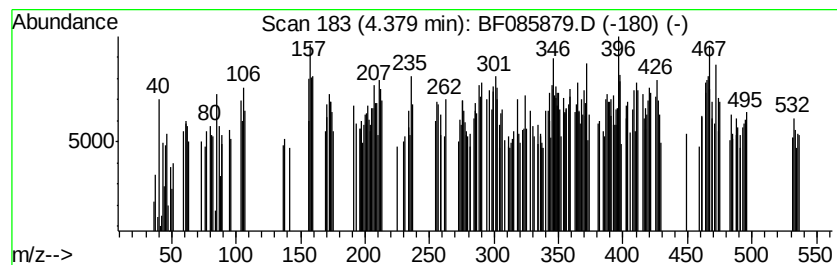
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 unknown4.38 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.38	5.55 ng	171886	1,4-Dichlorobenzene-d4	6.80

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Cyano-2-[3,4-dichlorophenyl]-2...	545	C24H37Cl2N0Sn	1000253-30-5	14
2			Tributyl 3-cyano-3-[3,4-dichloro...	531	C23H35Cl2N0Sn	1000253-30-2	10
3			5-Isoxazamine, N-methyl-N,3-di...	532	C36H28N20Si	135703-02-5	10
4			5-[2-Chlorophenylthio]-6-methoxy...	531	C29H26ClN303S	064895-64-3	9
5			Mercury iodo, (o-iodophenyl)-	532	C6H4HgI2	089487-89-8	9



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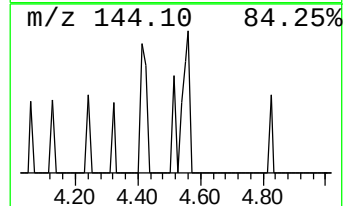
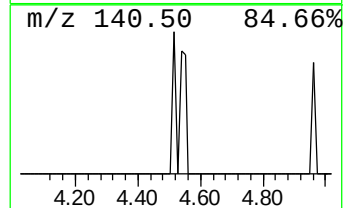
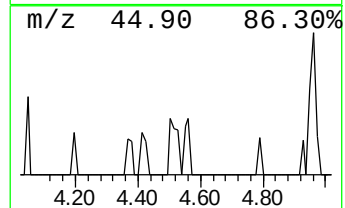
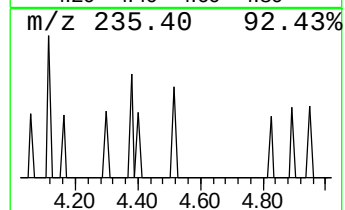
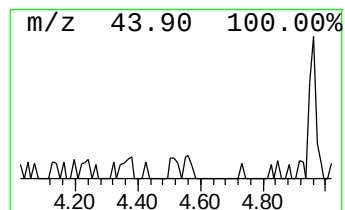
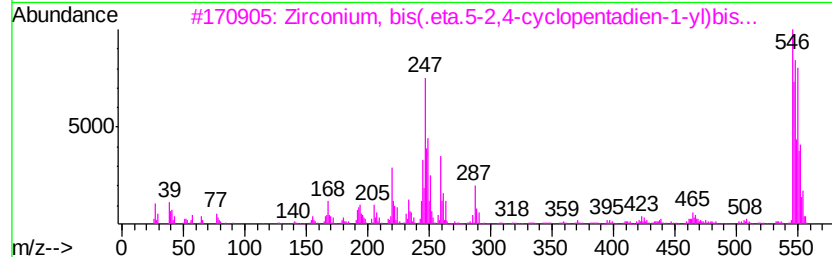
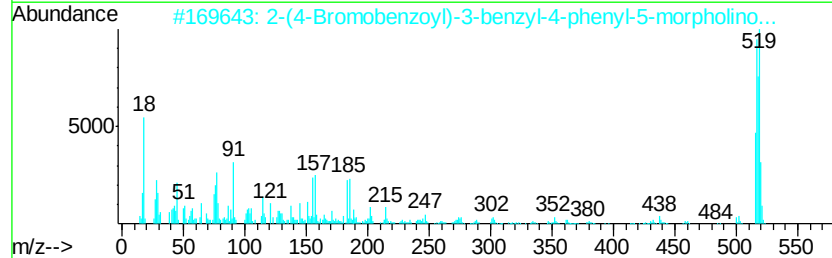
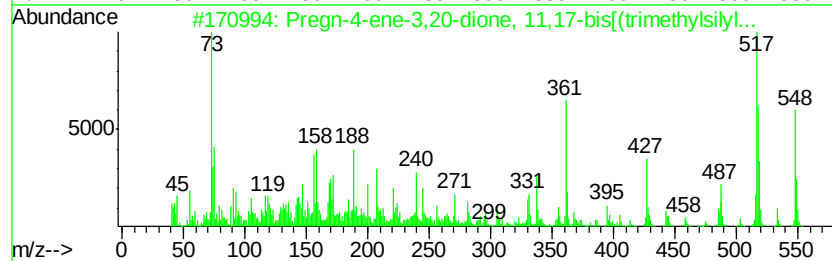
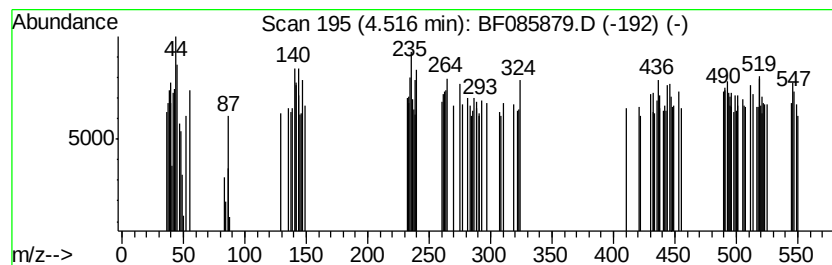
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 unknown4.52 Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.52	2.60 ng	80424	1,4-Dichlorobenzene-d4	6.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pregn-4-ene-3,20-dione, 11,17-bi...	548	C29H52N2O4Si2	057305-37-0	14
2		2-(4-Bromobenzoyl)-3-benzyl-4-ph...	517	C28H24BrN02S	1000286-90-3	14
3		Zirconium, bis(.eta.5-2,4-cyclop...	546	C28H30Zr2	119144-91-1	10
4		5-[(2-Pyridylmethyleme)amino]-2'...	518	C32H22N8	294651-18-6	9
5		BIS(2-(P-ETHOXYBENZYLIMINOMETHYL...	517	C28H30CuN4O2	1000243-53-5	5



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF033016\
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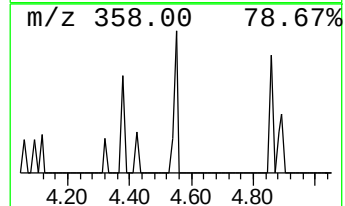
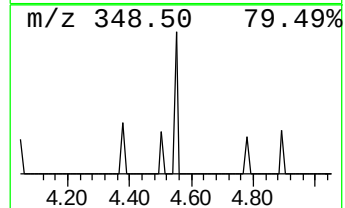
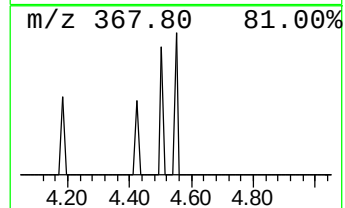
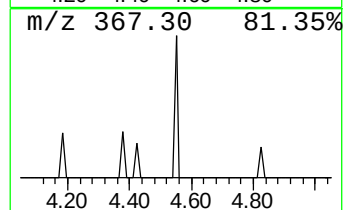
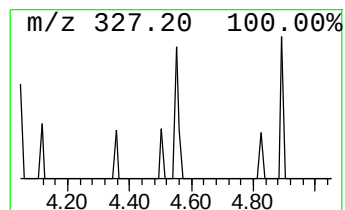
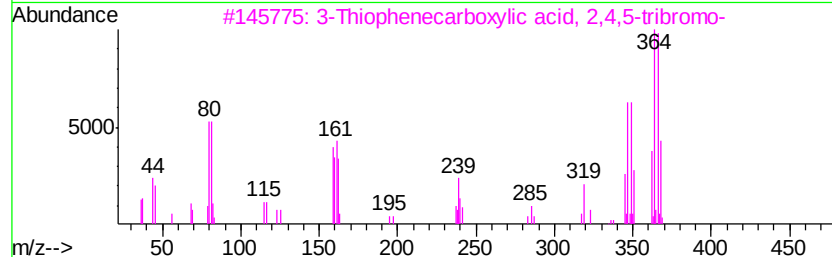
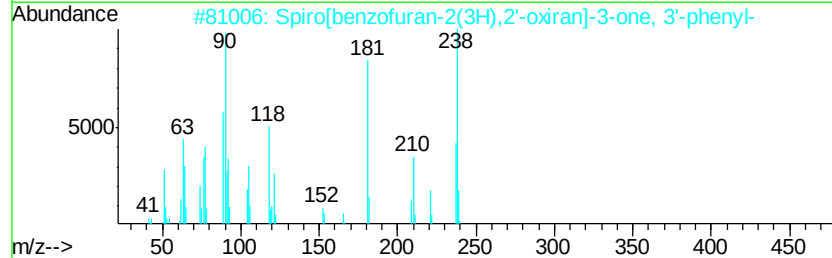
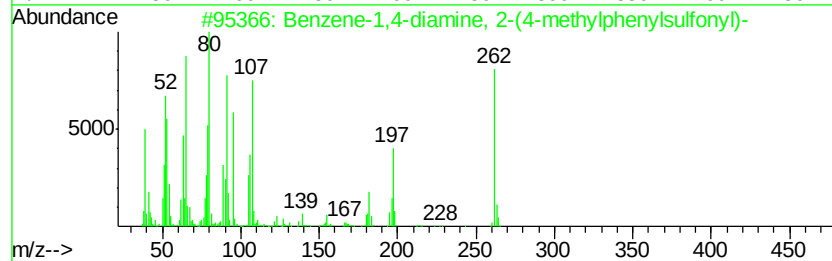
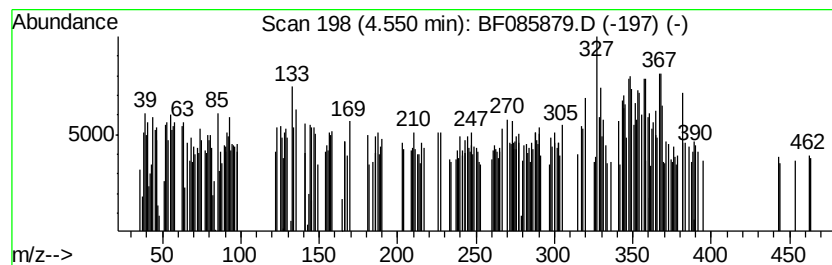
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 unknown4.55 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.55	4.36 ng	134908	1,4-Dichlorobenzene-d4	6.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene-1,4-diamine, 2-(4-methyl...	262	C13H14N2O2S	1000294-50-0	35
2		Spiro[benzofuran-2(3H),2'-oxiran...	238	C15H10O3	035405-25-5	25
3		3-Thiophenecarboxylic acid, 2,4,...	362	C5HBr3O2S	016694-22-7	18
4		6-[(3-Allylsalicylidene)amino]be...	277	C17H15N3O	294648-34-3	15
5		Pyridinium, 1-(m-chloroanilino)-...	204	C11H9ClN2	031378-94-6	14



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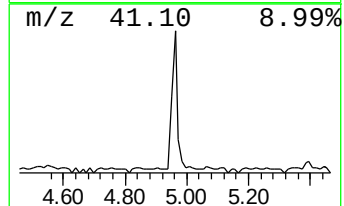
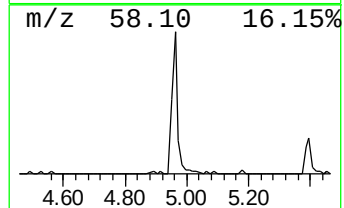
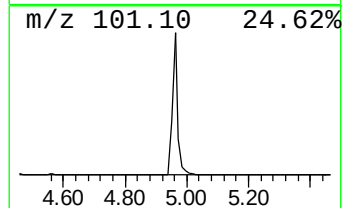
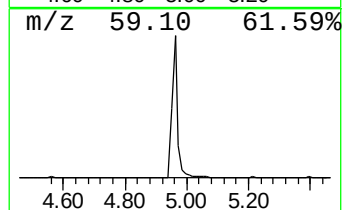
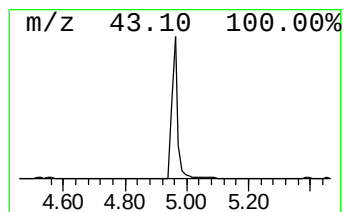
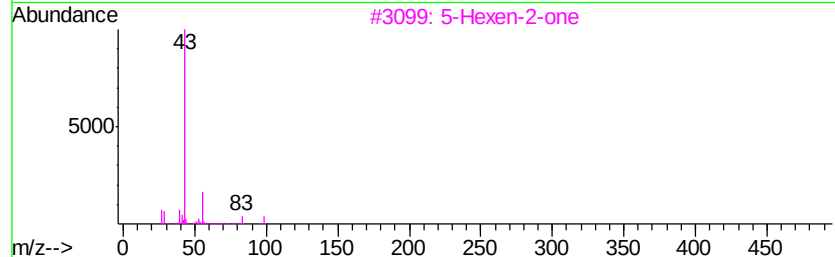
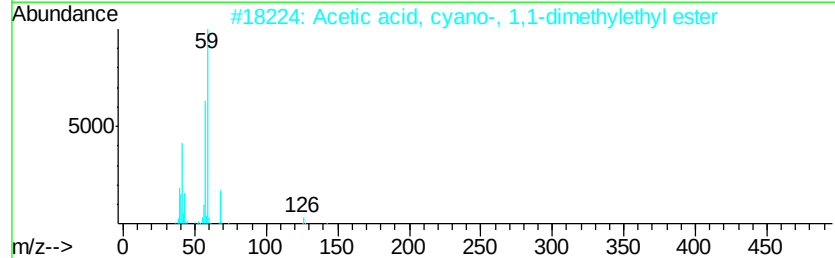
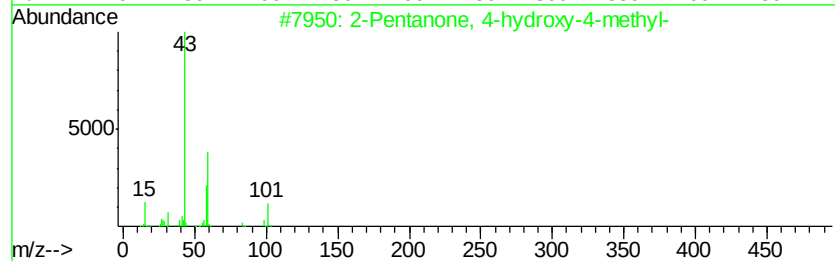
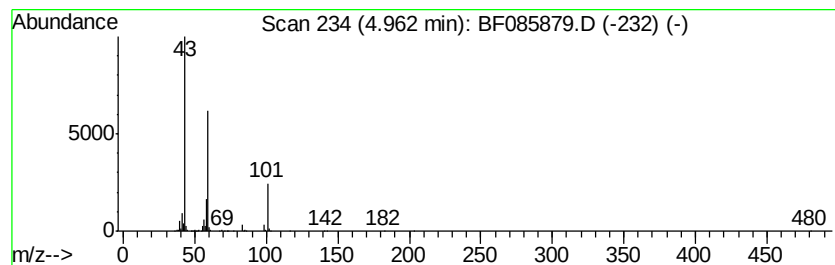
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.96	14.36 ng	444526	1,4-Dichlorobenzene-d4	6.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17
3		5-Hexen-2-one	98	C6H10O	000109-49-9	9
4		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5		1,3-Dioxane, 4,4-dimethyl-	116	C6H12O2	000766-15-4	9



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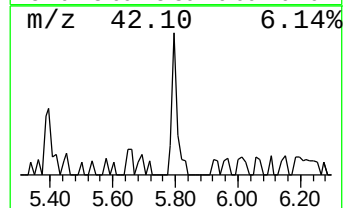
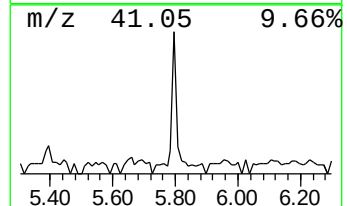
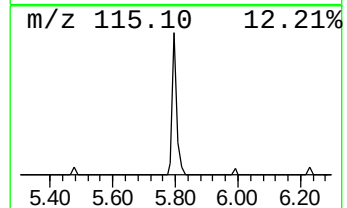
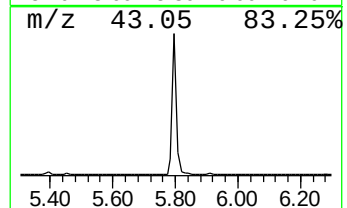
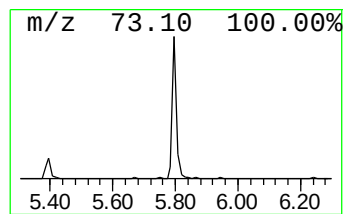
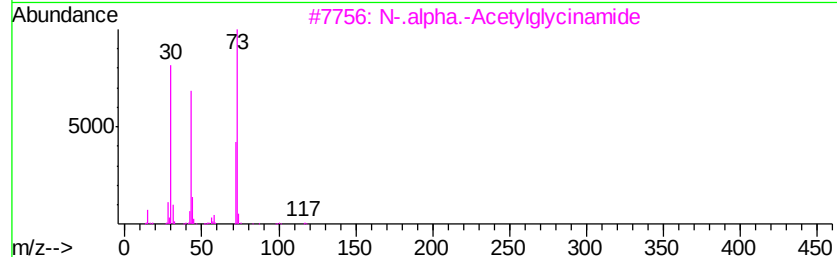
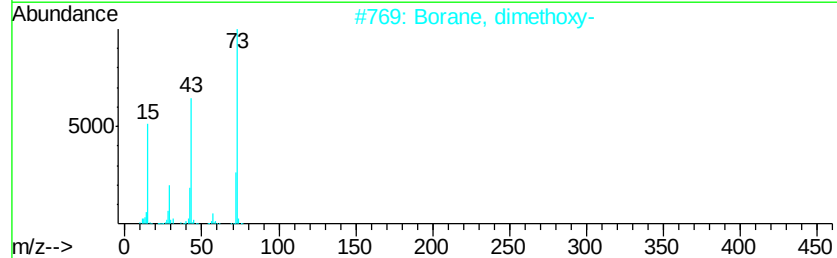
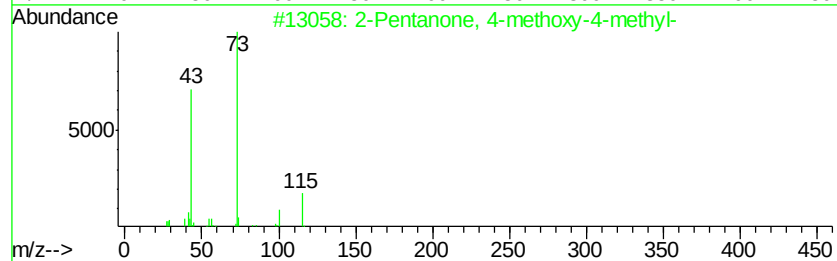
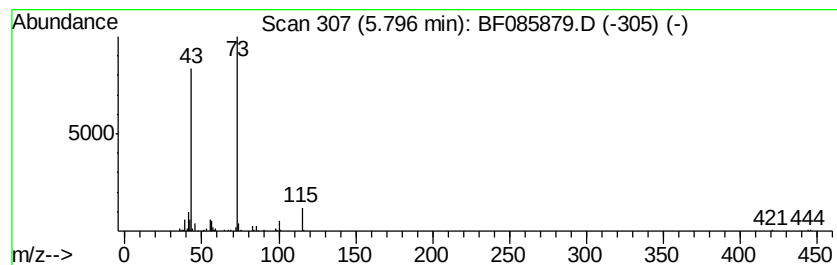
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 16 2-Pentanone, 4-methoxy-4-me... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.80	4.42 ng	136914	1,4-Dichlorobenzene-d4	6.80

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-methoxy-4-methyl-	130	C7H14O2	000107-70-0	72
2			Borane, dimethoxy-	74	C2H7BO2	004542-61-4	25
3			N-.alpha.-Acetylglycinamide	116	C4H8N2O2	002620-63-5	25
4			2-Pentanone, 3-methyl-	100	C6H12O	000565-61-7	22
5			4-Acetoxytridecane	242	C15H30O2	060826-30-4	17



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF033016\
Data File : BF085879.D
Acq On : 30 Mar 2016 11:28
Operator : UM/SJ
Sample : H1937-02
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
FIELDBLANK-3-23-16

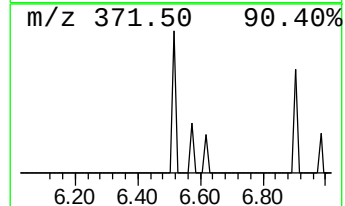
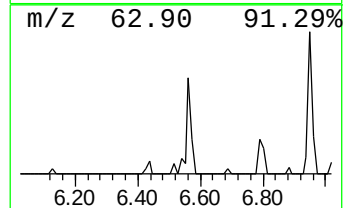
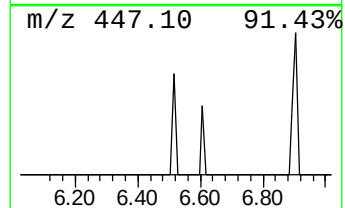
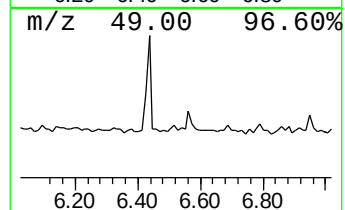
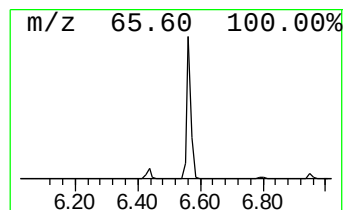
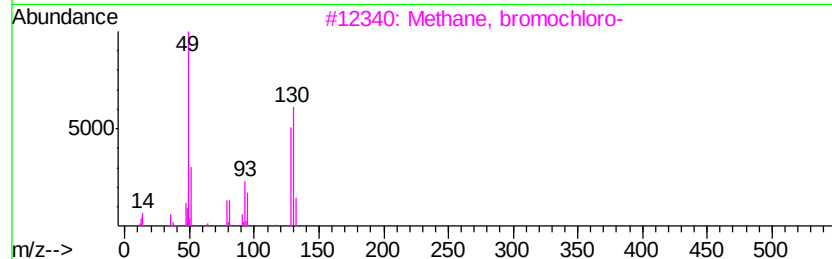
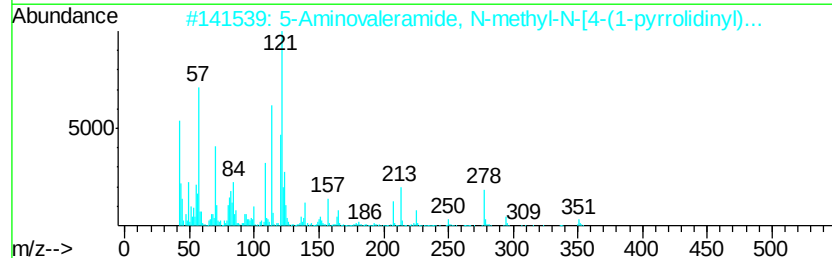
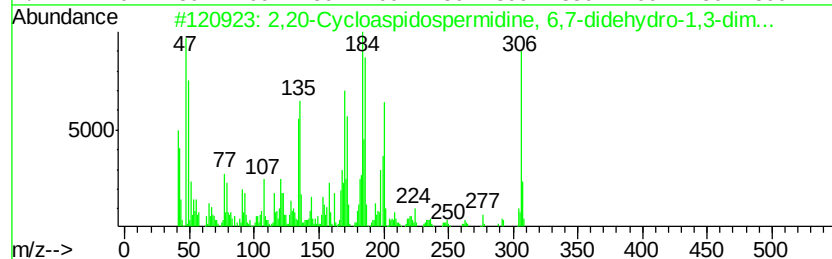
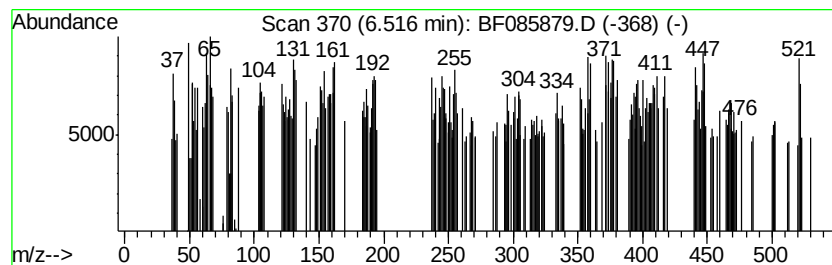
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 17 unknown6.52 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.52	7.19 ng	222480	1,4-Dichlorobenzene-d4	6.80

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2	20-Cycloaspidospermidine, 6,7-...	306	C21H26N2	056053-37-3	42	
2	5	Aminovaleramide, N-methyl-N-[4...	351	C19H33N3O3	124045-40-5	27	
3		Methane, bromochloro-	128	CH2BrCl	000074-97-5	27	
4		N-[4-Chloro-2-chloroacetamidophe...	286	C13H16Cl2N2O	1000254-96-6	27	
5		Benzamide, 2,2'-sulfinylbis[N-(1...	372	C20H24N2O3S	065838-71-3	27	



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF033016\
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Acq On : 30 Mar 2016 11:28
Operator : UM/SJ
Sample : H1937-02
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
FIELDBLANK-3-23-16

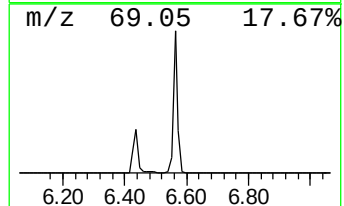
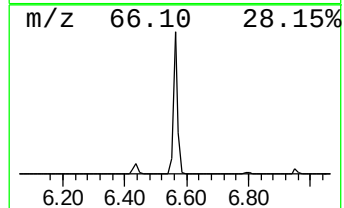
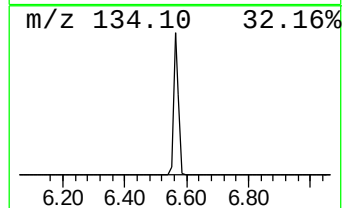
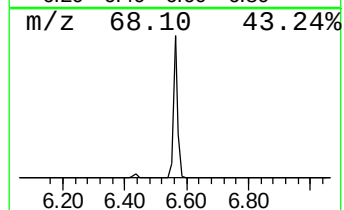
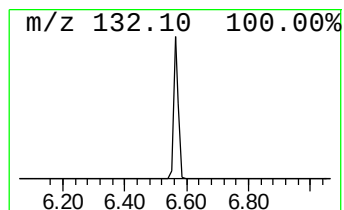
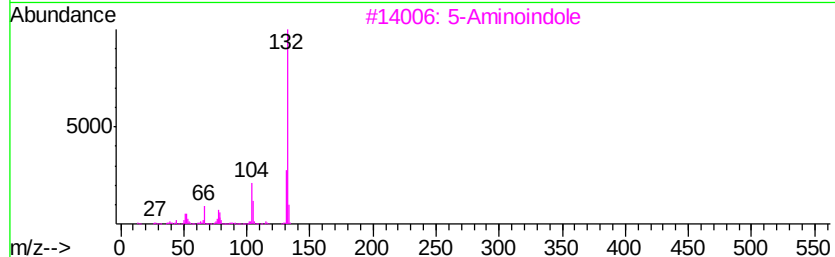
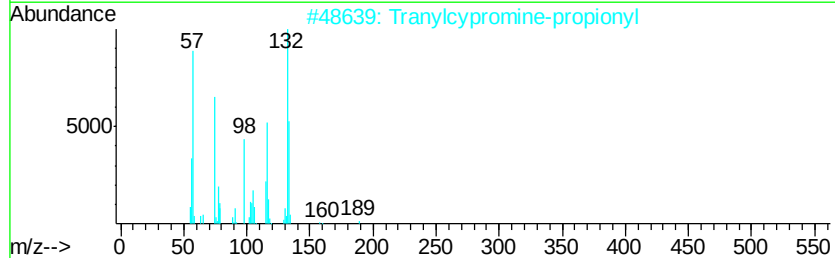
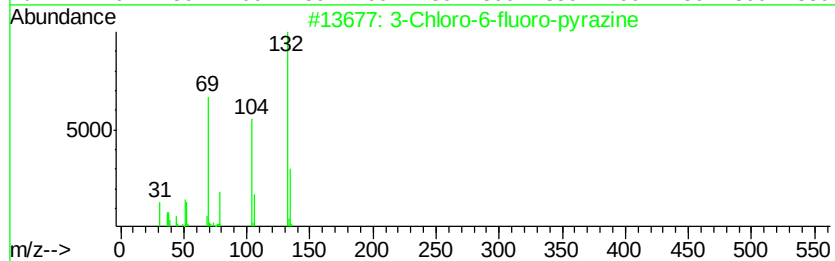
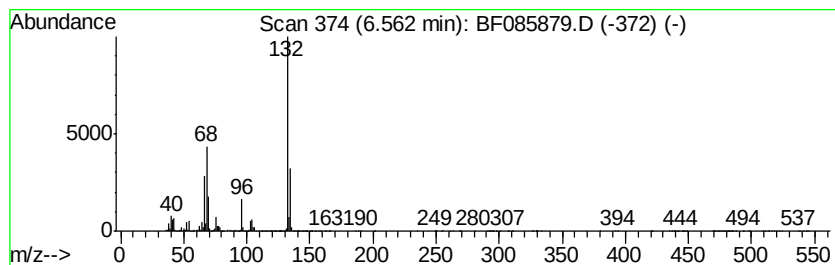
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032416.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 18 unknown6.56 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.56	77.55 ng	2400660	1,4-Dichlorobenzene-d4	6.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		5-Aminoindole	132	C8H8N2	005192-03-0	9
4		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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 Sample : H1937-02
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FIELDBLANK-3-23-16

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032416.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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unknown2.52	2.52	2.9	ng	89753	1	6.80	619118	20.0
unknown2.74	2.74	6.5	ng	202578	1	6.80	619118	20.0
unknown2.81	2.81	6.9	ng	213148	1	6.80	619118	20.0
unknown3.14	3.14	4.9	ng	151070	1	6.80	619118	20.0
unknown3.38	3.38	4.4	ng	137370	1	6.80	619118	20.0
unknown4.31	4.31	3.3	ng	101361	1	6.80	619118	20.0
unknown4.38	4.38	5.5	ng	171886	1	6.80	619118	20.0
unknown4.52	4.52	2.6	ng	80424	1	6.80	619118	20.0
unknown4.55	4.55	4.4	ng	134908	1	6.80	619118	20.0
2-Pentanone, 4-hy...	4.96	14.4	ng	444526	1	6.80	619118	20.0
2-Pentanone, 4-me...	5.80	4.4	ng	136914	1	6.80	619118	20.0
unknown6.52	6.52	7.2	ng	222480	1	6.80	619118	20.0
unknown6.56	6.56	77.5	ng	2400660	1	6.80	619118	20.0