

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052816\
 Data File : BF087512.D
 Acq On : 29 May 2016 10:00
 Operator : UM/SJ
 Sample : H3222-01
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-10

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-BF052316.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	4.730	274	275	283	rBV	22270	87339	2.15%	0.241%
2	5.370	329	331	352	rBV	777506	1786961	43.96%	4.923%
3	6.159	397	400	406	rBV	30717	69123	1.70%	0.190%
4	6.867	459	462	464	rBV	37455	72940	1.79%	0.201%
5	6.982	470	472	475	rBV	32947	51883	1.28%	0.143%
6	7.027	475	476	482	rVV	513345	1029991	25.34%	2.838%
7	7.119	482	484	492	rVB	2097295	3269154	80.43%	9.006%
8	7.404	505	509	512	rVB3	32781	78889	1.94%	0.217%
9	7.473	512	515	527	rBV	416084	867807	21.35%	2.391%
10	7.702	532	535	542	rBV	2080943	3025560	74.44%	8.335%
11	7.804	542	544	549	rVB	38278	50434	1.24%	0.139%
12	7.885	549	551	559	rVB2	112279	225066	5.54%	0.620%
13	8.022	559	563	573	rBV4	93065	295452	7.27%	0.814%
14	8.170	573	576	579	rBV2	47151	85759	2.11%	0.236%
15	8.239	579	582	585	rVB3	22227	43460	1.07%	0.120%
16	8.319	585	589	590	rBV3	47408	128553	3.16%	0.354%
17	8.387	590	595	602	rVV2	1809127	2912973	71.67%	8.025%
18	8.559	609	610	613	rVB	60094	74333	1.83%	0.205%
19	8.719	622	624	629	rVB2	91600	144896	3.56%	0.399%
20	8.822	629	633	635	rBV2	42234	57584	1.42%	0.159%
21	8.867	635	637	639	rVB2	33714	44788	1.10%	0.123%
22	8.970	643	646	647	rBV	55595	95209	2.34%	0.262%
23	8.993	647	648	653	rVB2	110301	217552	5.35%	0.599%
24	9.096	654	657	661	rVB2	56283	113174	2.78%	0.312%
25	9.187	663	665	669	rBV2	30019	74022	1.82%	0.204%
26	9.279	671	673	678	rVB5	31930	73458	1.81%	0.202%
27	9.382	678	682	683	rBV3	25705	57678	1.42%	0.159%
28	9.416	683	685	687	rVB	41838	52254	1.29%	0.144%
29	9.496	690	692	699	rVV	789241	1349609	33.20%	3.718%
30	9.599	699	701	709	rVB3	98476	218778	5.38%	0.603%
31	9.908	726	728	731	rVB3	33910	52217	1.28%	0.144%
32	9.976	731	734	737	rBV4	39102	80833	1.99%	0.223%
33	10.216	752	755	757	rBV3	33896	86342	2.12%	0.238%
34	10.273	757	760	762	rVV2	122796	229601	5.65%	0.633%

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35	10.342	763	766	770	rVV	139252	264514	6.51%	0.729%
36	10.445	770	775	778	rVV2	52363	160656	3.95%	0.443%
37	10.490	778	779	782	rVB	41806	47682	1.17%	0.131%
38	10.548	782	784	785	rBV2	28925	45642	1.12%	0.126%
39	10.582	785	787	791	rVB2	56859	89237	2.20%	0.246%
40	10.650	791	793	796	rBV3	27560	65304	1.61%	0.180%
41	10.776	799	804	805	rBV5	33165	92051	2.26%	0.254%
42	10.810	805	807	810	rVV	306108	416802	10.25%	1.148%
43	10.868	810	812	816	rVB4	30952	51857	1.28%	0.143%
44	10.982	820	822	826	rVB3	34697	58934	1.45%	0.162%
45	11.085	828	831	832	rBV	52387	74388	1.83%	0.205%
46	11.153	835	837	840	rVB	46821	73848	1.82%	0.203%
47	11.211	840	842	845	rBV3	29974	46706	1.15%	0.129%
48	11.279	845	848	851	rBV	3188511	4064576	100.00%	11.198%
49	11.485	863	866	868	rBV3	79650	158041	3.89%	0.435%
50	11.531	868	870	873	rVV3	74578	158415	3.90%	0.436%
51	11.576	873	874	878	rVB3	58429	96638	2.38%	0.266%
52	11.748	886	889	890	rBV3	37224	69303	1.71%	0.191%
53	11.782	890	892	893	rVV	49585	81939	2.02%	0.226%
54	11.805	893	894	896	rVV2	75090	105438	2.59%	0.290%
55	11.851	896	898	901	rVB5	49684	89858	2.21%	0.248%
56	11.976	906	909	916	rVV3	239585	474693	11.68%	1.308%
57	12.113	919	921	926	rVB3	92832	181439	4.46%	0.500%
58	12.216	929	930	932	rBV2	26154	42186	1.04%	0.116%
59	12.319	937	939	943	rBV	718338	1144025	28.15%	3.152%
60	12.376	943	944	947	rVB2	39678	62976	1.55%	0.173%
61	12.594	961	963	967	rBV4	29242	77262	1.90%	0.213%
62	12.834	980	984	985	rBV4	61476	131129	3.23%	0.361%
63	12.868	985	987	989	rVB3	56986	98500	2.42%	0.271%
64	13.096	1005	1007	1010	rBV4	38390	77303	1.90%	0.213%
65	13.154	1010	1012	1017	rVV4	47436	101372	2.49%	0.279%
66	13.234	1017	1019	1022	rVB2	46619	74726	1.84%	0.206%
67	13.508	1040	1043	1046	rBV	161046	228670	5.63%	0.630%
68	13.622	1050	1053	1062	rBV	1710965	2975810	73.21%	8.198%
69	13.931	1076	1080	1085	rVB7	66110	215223	5.30%	0.593%
70	14.399	1119	1121	1124	rBV3	26220	52490	1.29%	0.145%
71	14.514	1129	1131	1134	rVB	74272	103867	2.56%	0.286%

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72	14.731	1147	1150	1156	rBV	764014	1001881	24.65%	2.760%
73	15.725	1235	1237	1244	rVB	144574	269496	6.63%	0.742%
74	16.823	1331	1333	1337	rBV2	68416	146088	3.59%	0.402%
75	16.925	1340	1342	1345	rVB	94249	118544	2.92%	0.327%
76	17.085	1353	1356	1359	rBV	3217624	3676211	90.45%	10.128%
77	17.851	1421	1423	1426	rBV	46051	66493	1.64%	0.183%
78	17.920	1427	1429	1434	rVB2	47222	67873	1.67%	0.187%
79	18.354	1464	1467	1472	rBV4	29683	70049	1.72%	0.193%
80	18.434	1472	1474	1484	rVB	582256	829826	20.42%	2.286%
81	20.114	1619	1621	1626	rBV	357263	594915	14.64%	1.639%

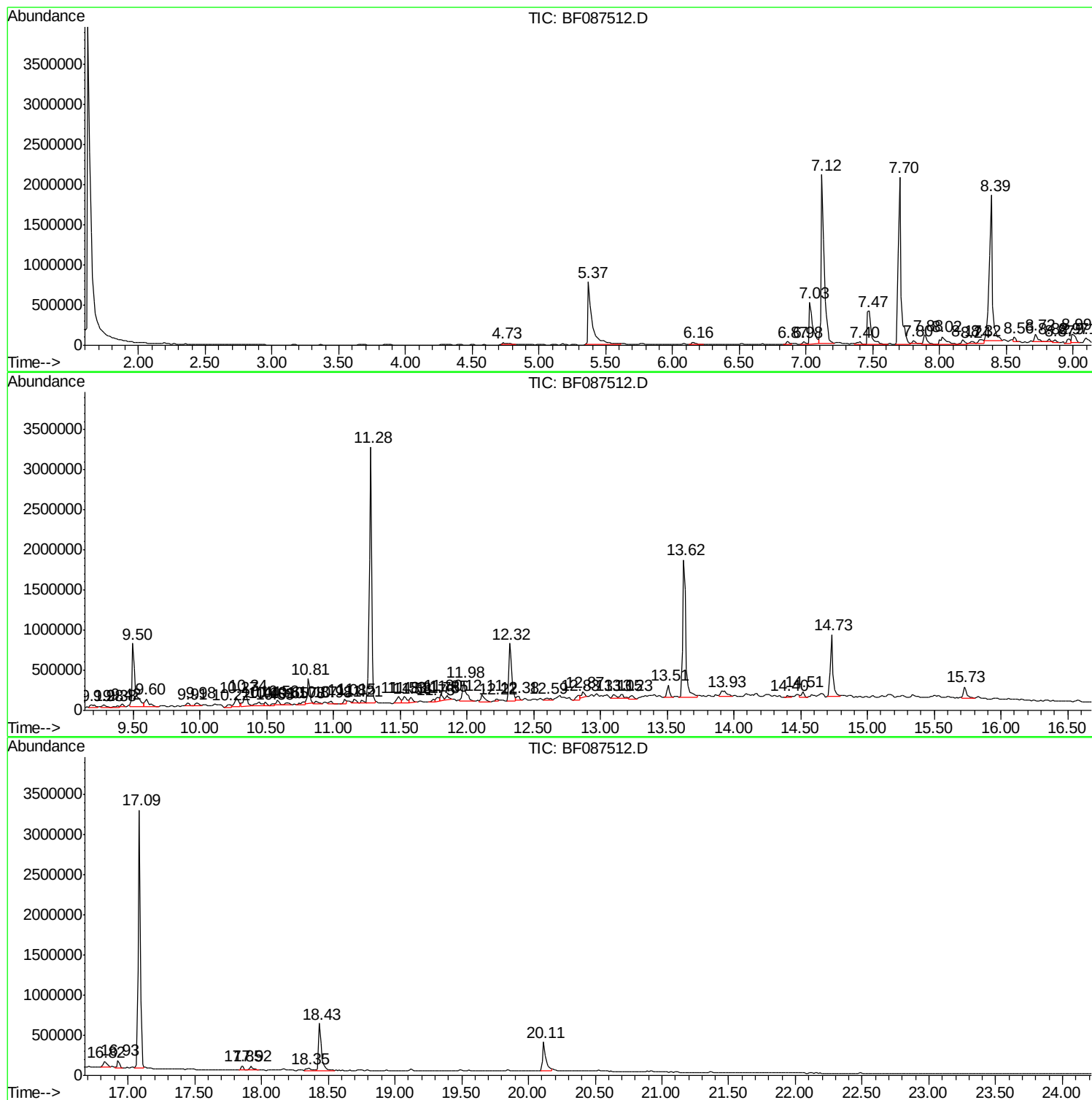
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF052316.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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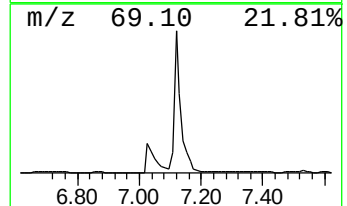
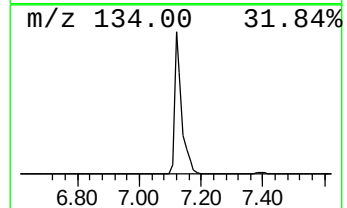
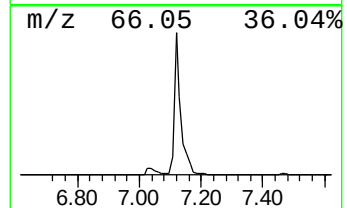
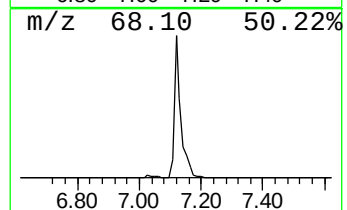
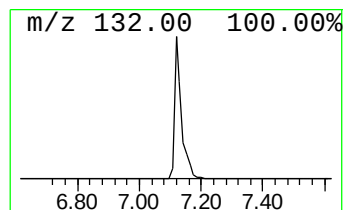
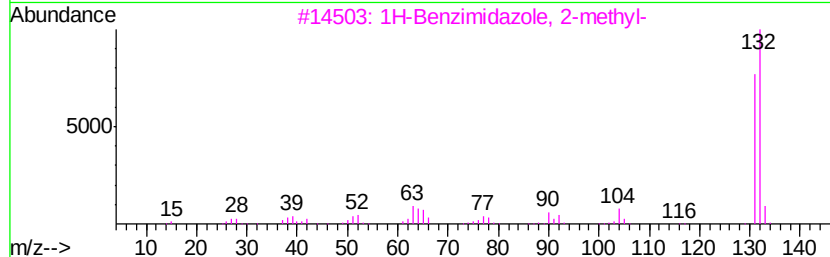
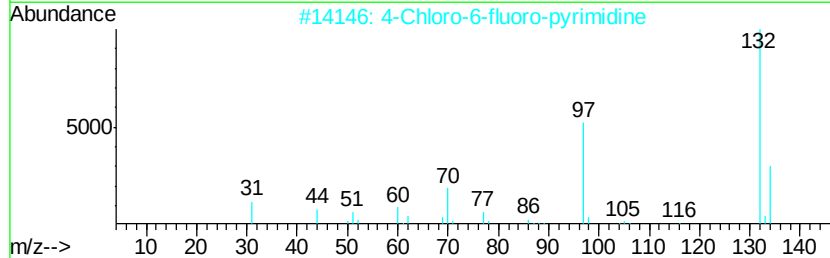
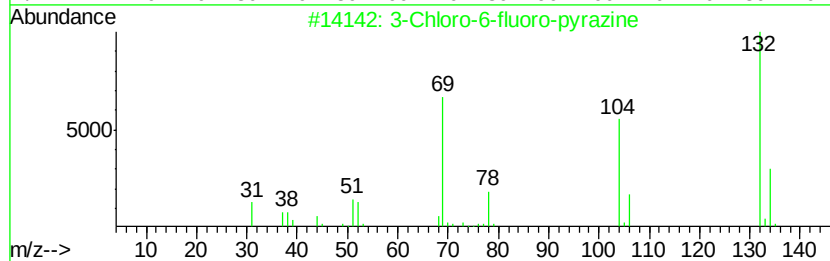
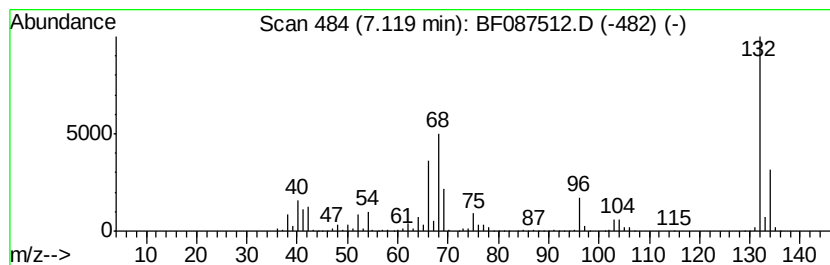
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown7.12 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.12	75.34 ng	3269150	1,4-Dichlorobenzene-d4	7.47

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
4		Pyrazolo(2,3-a)pyridine, 7-methyl-	132	C8H8N2	034760-58-2	22
5		Tranlylcypromine-propionyl	189	C12H15NO	1000123-86-3	22



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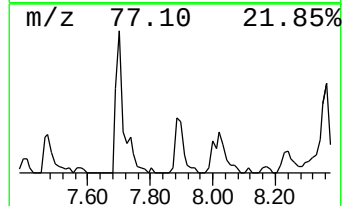
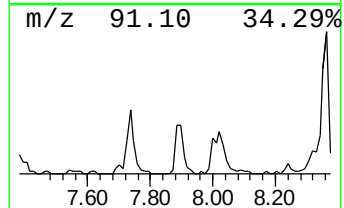
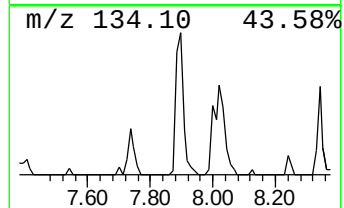
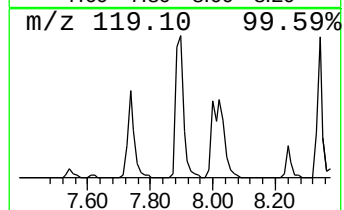
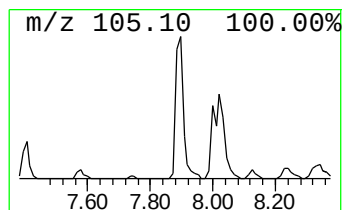
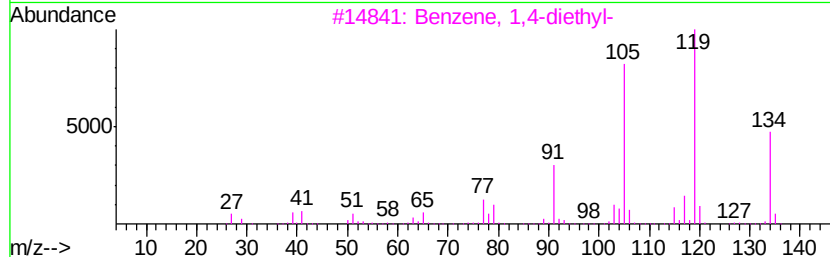
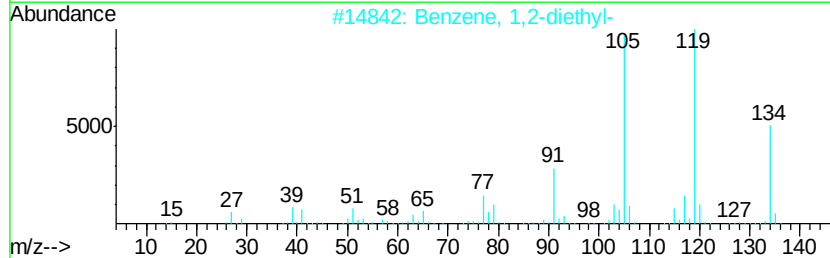
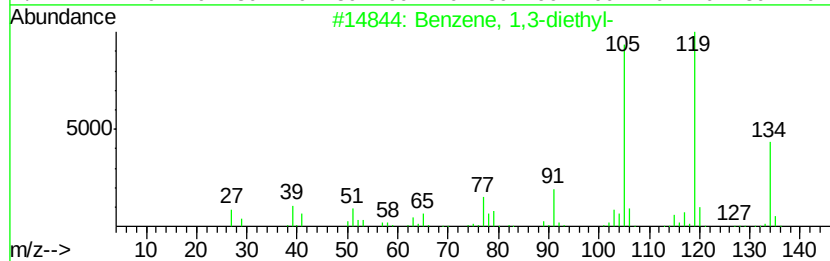
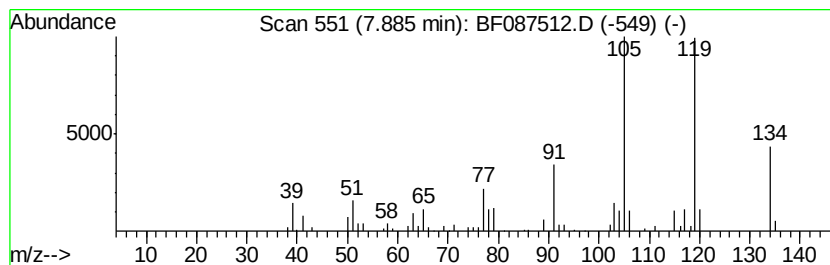
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TIC Library : C:\Database\NIST11.L
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Peak Number 3 Benzene, 1,3-diethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.88	5.19 ng	225066	1,4-Dichlorobenzene-d4	7.47

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	97
2		Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	97
3		Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	93
4		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	87
5		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	87



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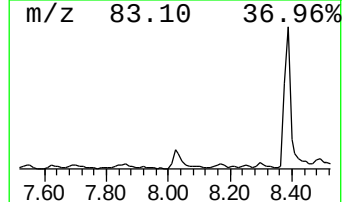
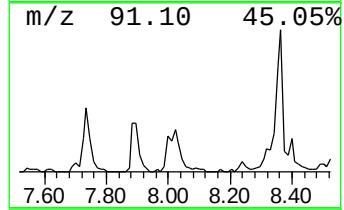
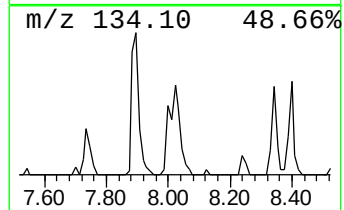
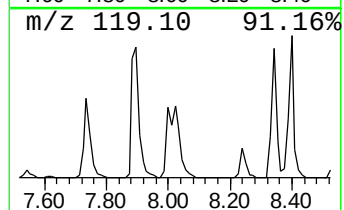
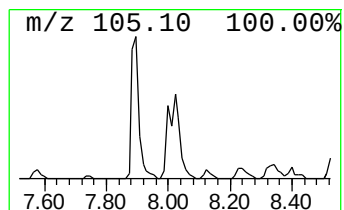
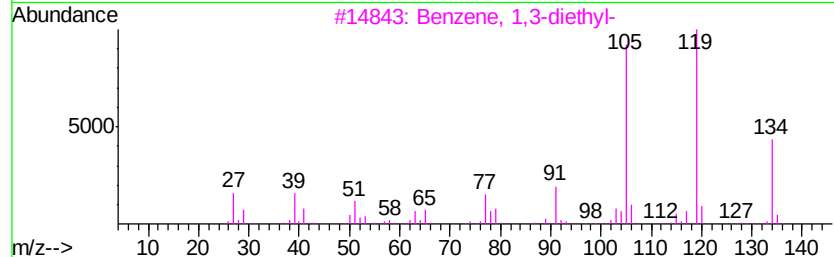
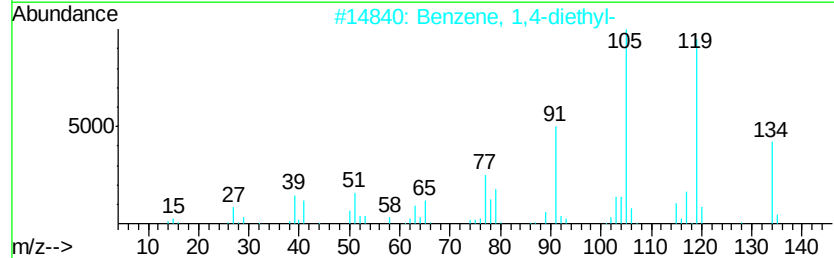
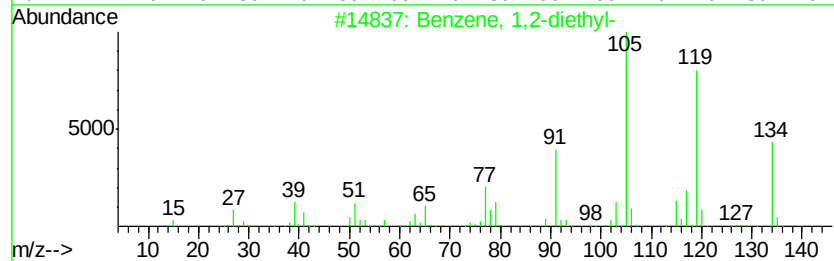
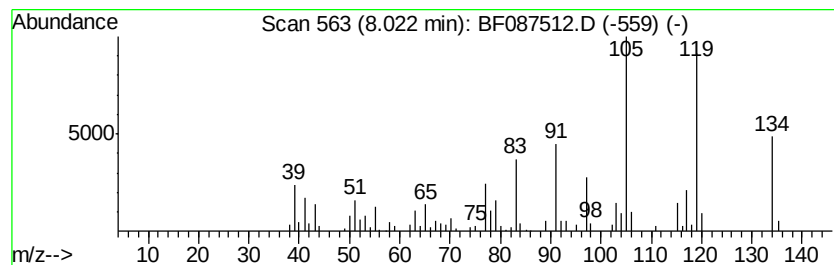
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TIC Library : C:\Database\NIST11.L
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Peak Number 4 Benzene, 1,2-diethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.02	6.81 ng	295452	1,4-Dichlorobenzene-d4	7.47

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	96
2			Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	90
3			Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	76
4			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	76
5			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	70



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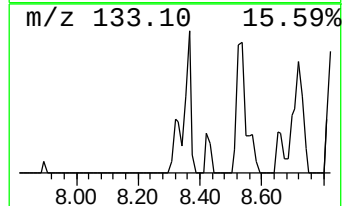
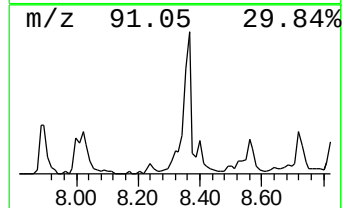
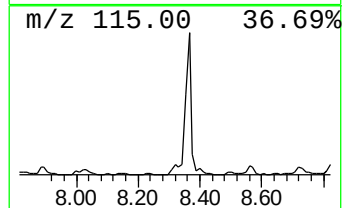
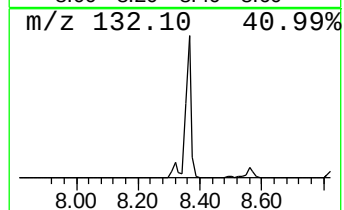
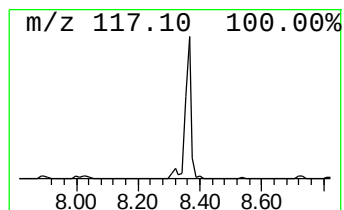
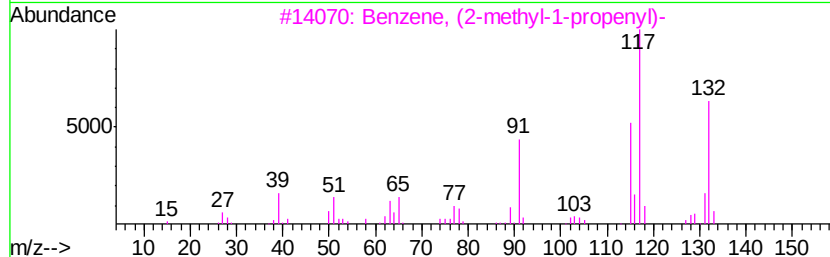
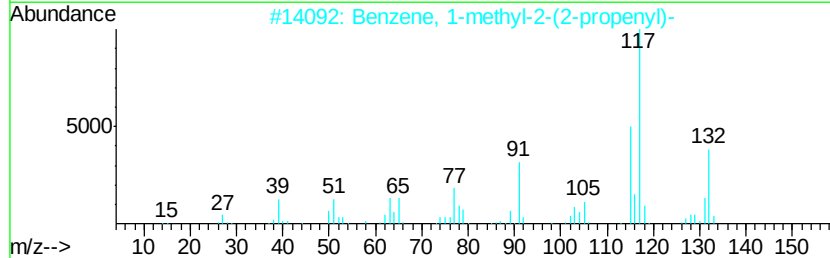
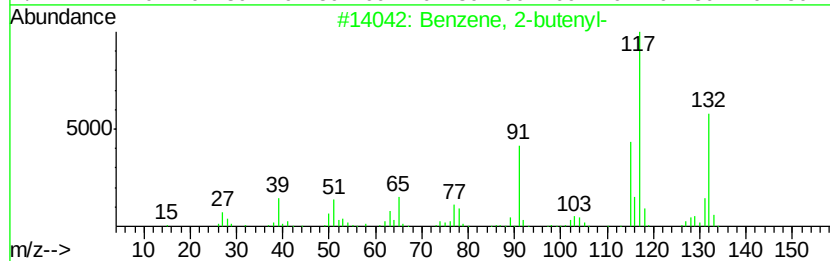
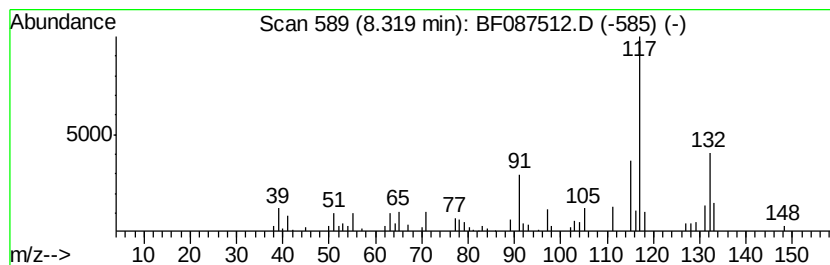
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Benzene, 2-butenyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.32	2.96 ng	128553	1,4-Dichlorobenzene-d4	7.47

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-butenyl-	132	C10H12	001560-06-1	96
2		Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	93
3		Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	93
4		1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	93
5		(E)-1-Phenyl-1-butene	132	C10H12	001005-64-7	91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052816\
Data File : BF087512.D
Acq On : 29 May 2016 10:00
Operator : UM/SJ
Sample : H3222-01
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-10

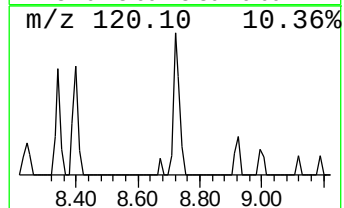
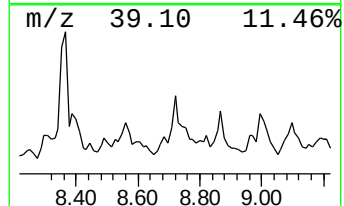
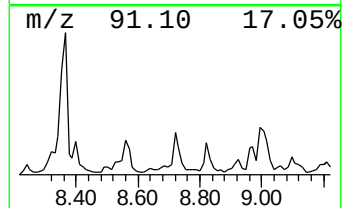
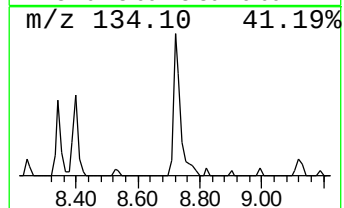
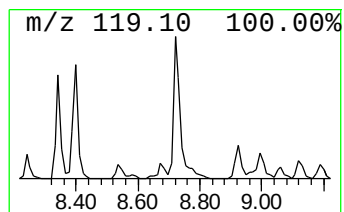
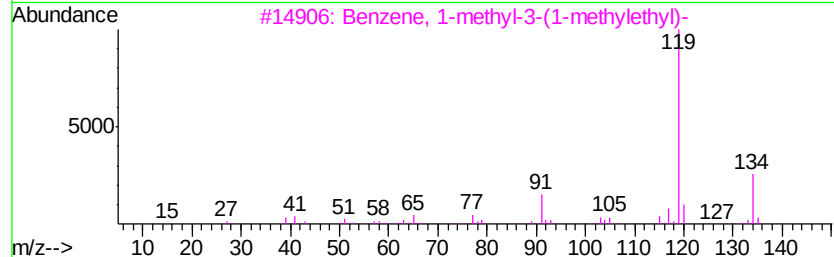
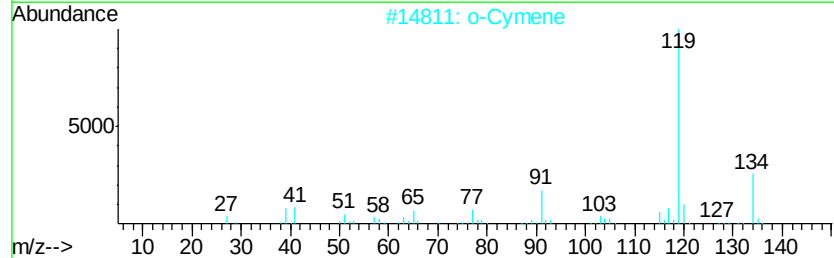
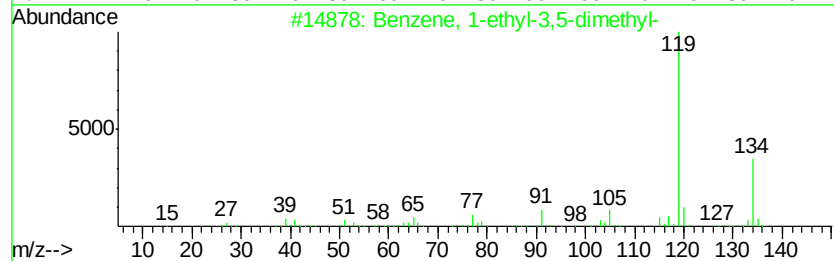
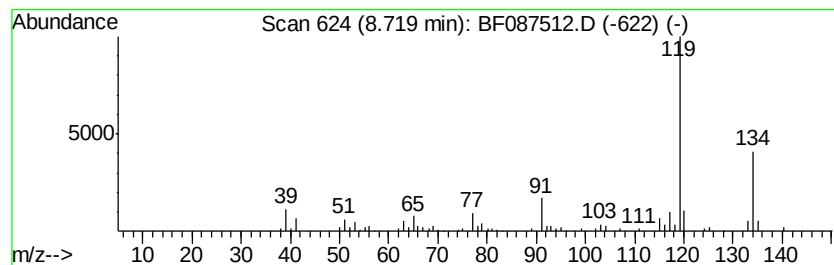
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Benzene, 1-ethyl-3,5-dimethyl- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.72	2.15 ng	144896	Napthalene-d8	9.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	95
2		o-Cymene	134	C10H14	000527-84-4	95
3		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94
4		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	94
5		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	94



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052816\
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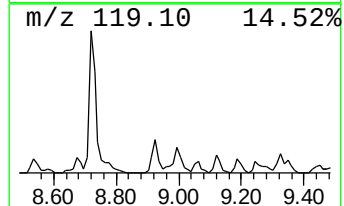
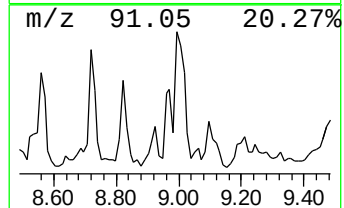
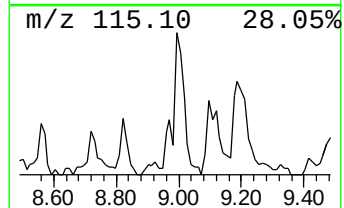
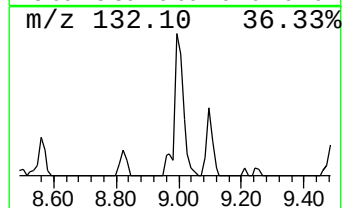
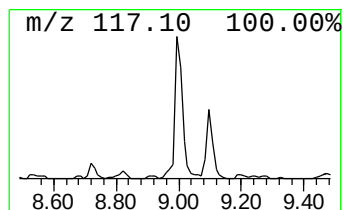
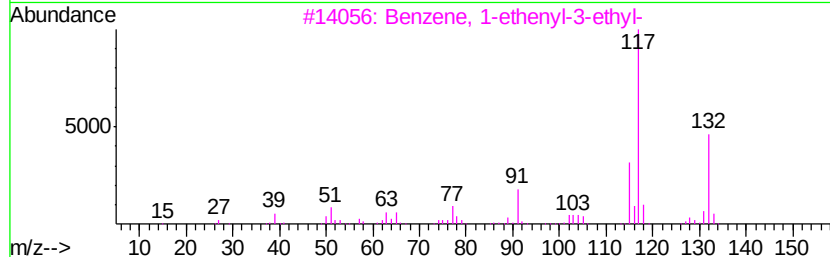
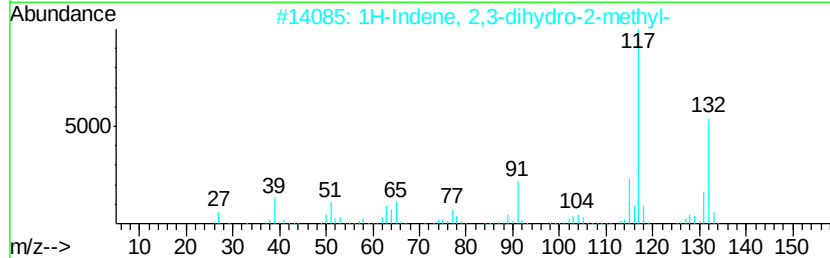
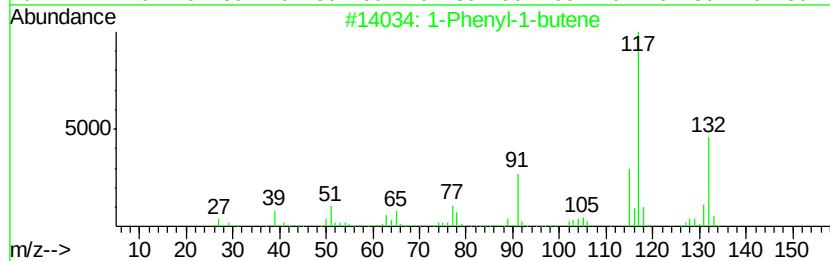
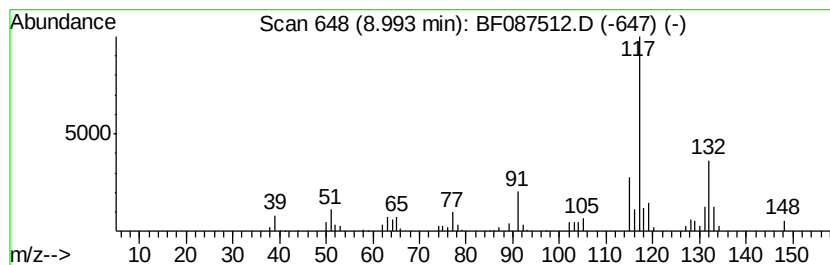
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 1-Phenyl-1-butene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.99	3.22 ng	217552	Napthalene-d8	9.50
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		1-Phenyl-1-butene	132 C10H12	000824-90-8 95
2		1H-Indene, 2,3-dihydro-2-methyl-	132 C10H12	000824-63-5 93
3		Benzene, 1-ethenyl-3-ethyl-	132 C10H12	007525-62-4 93
4		Benzene, 1-methyl-2-(2-propenyl)-	132 C10H12	001587-04-8 91
5		Benzene, 2-ethenyl-1,3-dimethyl-	132 C10H12	002039-90-9 87



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052816\
Data File : BF087512.D
Acq On : 29 May 2016 10:00
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Misc :
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Instrument :
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ClientSampleId :
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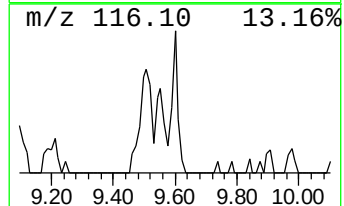
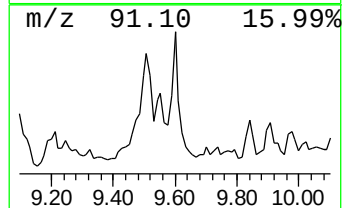
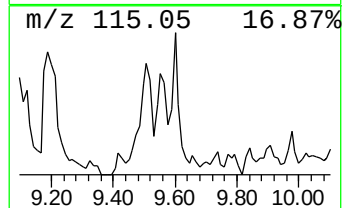
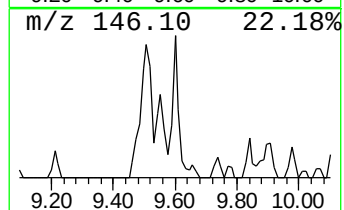
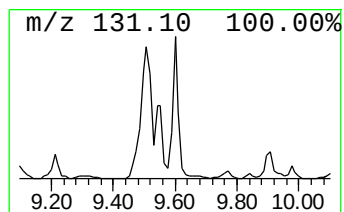
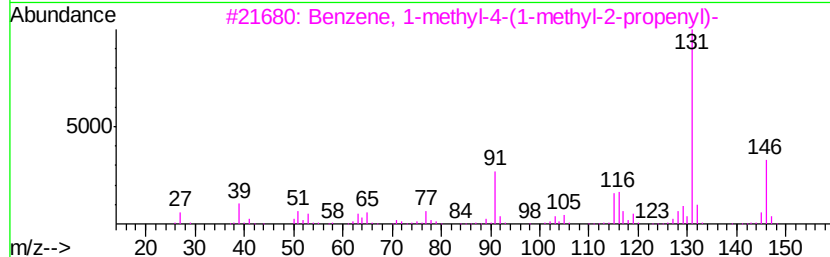
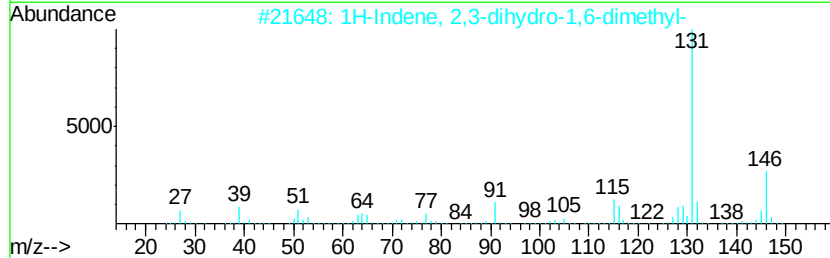
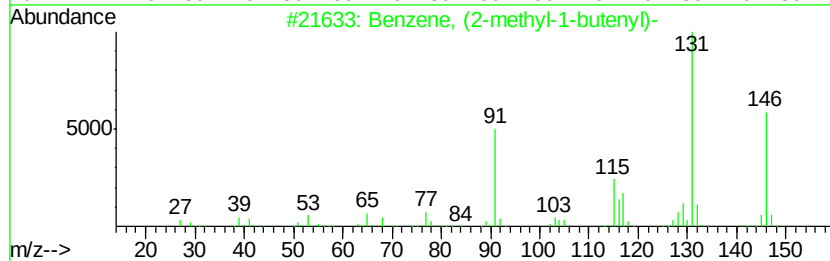
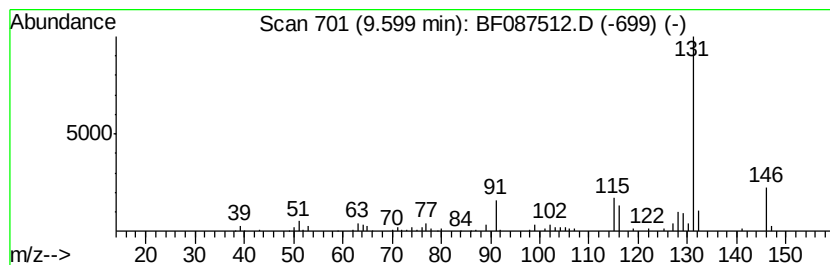
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Benzene, (2-methyl-1-butenyl)- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.60	3.24 ng	218778	Napthalene-d8	9.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	91
2		1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	91
3		Benzene, 1-methyl-4-(1-methyl-2-...	146	C11H14	097664-18-1	91
4		1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	91
5		Benzene, 1-methyl-3-(1-methyl-2-...	146	C11H14	052161-57-6	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052816\
Data File : BF087512.D
Acq On : 29 May 2016 10:00
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Sample : H3222-01
Misc :
ALS Vial : 36 Sample Multiplier: 1

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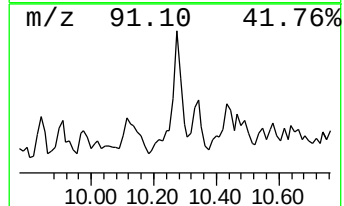
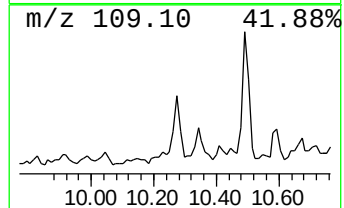
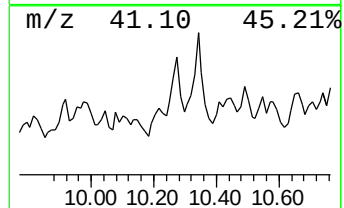
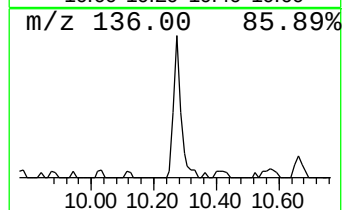
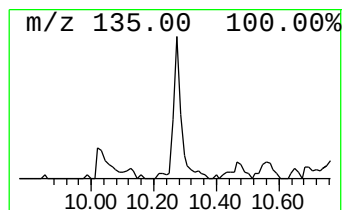
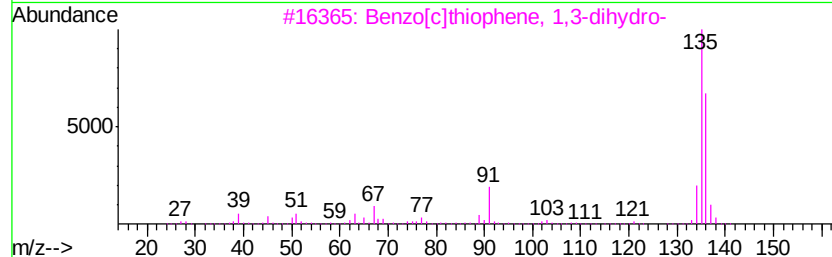
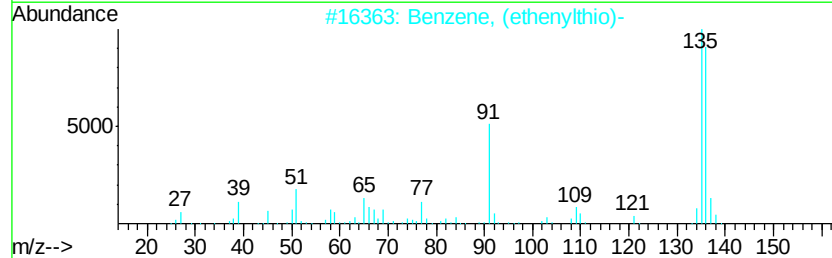
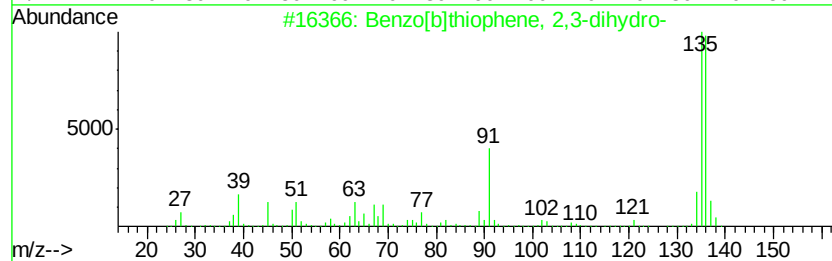
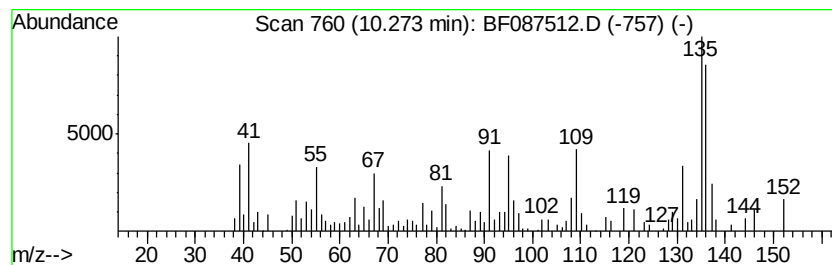
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Benzo[b]thiophene, 2,3-dihy... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.27	3.40 ng	229601	Napthalene-d8	9.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[b]thiophene, 2,3-dihydro-	136	C8H8S	004565-32-6	55
2		Benzene, (ethenylthio)-	136	C8H8S	001822-73-7	55
3		Benzo[c]thiophene, 1,3-dihydro-	136	C8H8S	002471-92-3	45
4		Pyrazine, 2,6-diethyl-	136	C8H12N2	013067-27-1	45
5		4-Hydroxy-2-methylbenzaldehyde	136	C8H8O2	041438-18-0	45



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052816\
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Sample : H3222-01
Misc :
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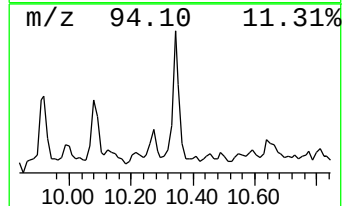
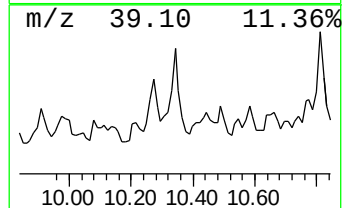
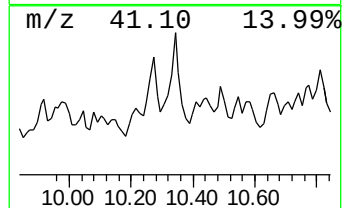
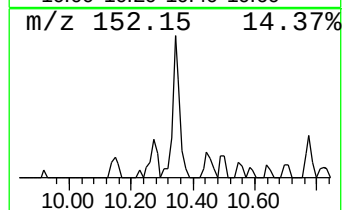
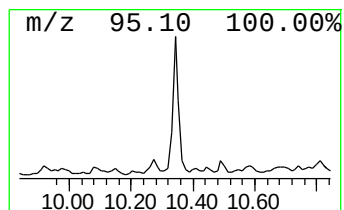
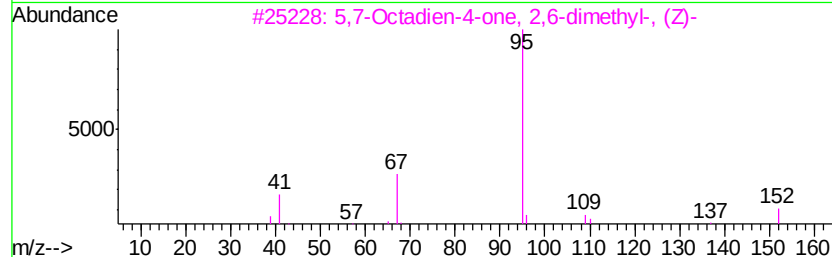
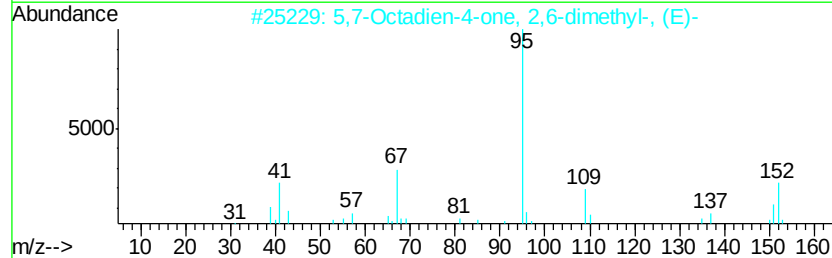
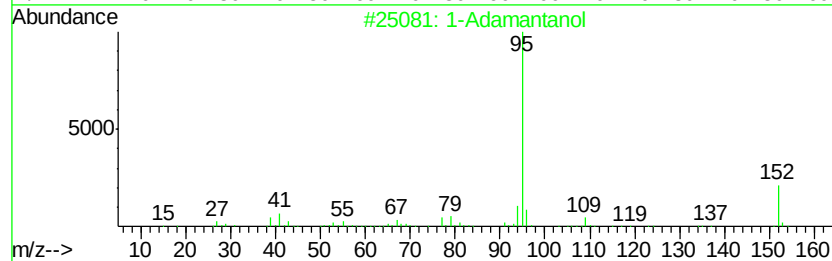
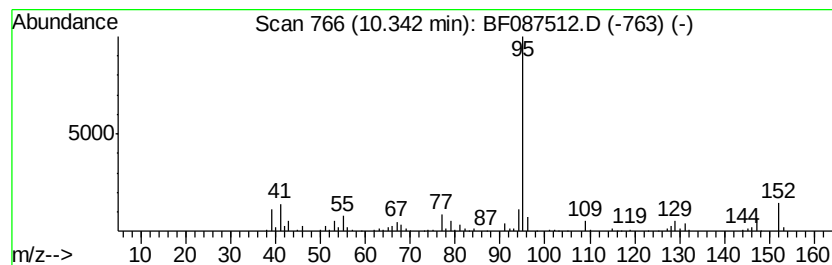
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 1-Adamantanol Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.34	3.92 ng	264514	Napthalene-d8	9.50	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Adamantanol	152	C10H16O	000768-95-6	87
2	5,7-Octadien-4-one, 2,6-dimethyl...	152	C10H16O	006752-80-3	64
3	5,7-Octadien-4-one, 2,6-dimethyl...	152	C10H16O	003588-18-9	53
4	4-Pyridinol	95	C5H5NO	000626-64-2	50
5	Furan, 2-ethyl-5-methyl-	110	C7H10O	001703-52-2	33



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052816\
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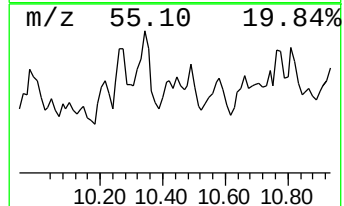
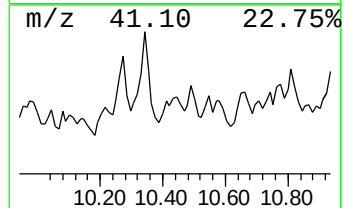
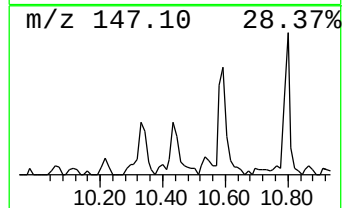
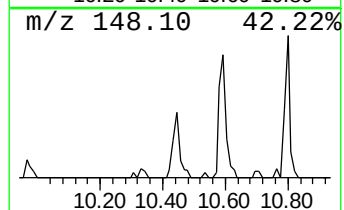
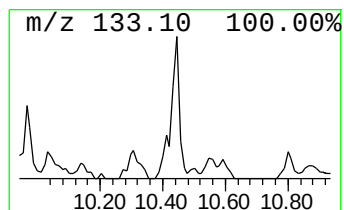
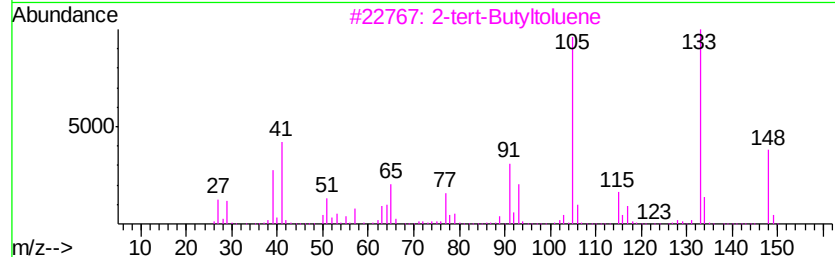
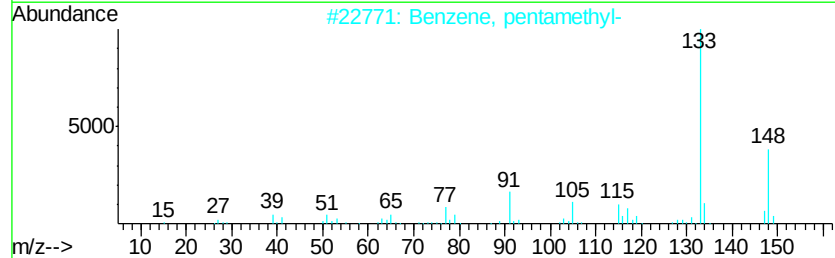
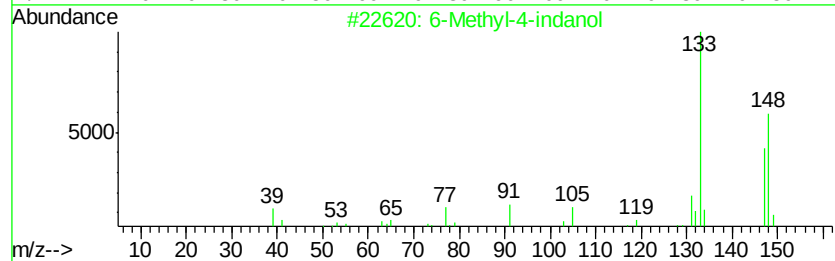
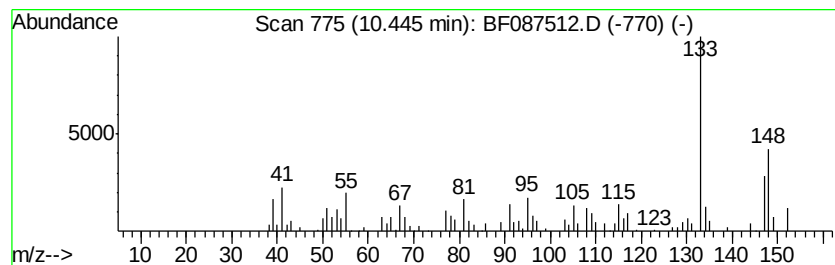
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 6-Methyl-4-indanol Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.44	2.38 ng	160656	Napthalene-d8	9.50	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	6-Methyl-4-indanol	148	C10H12O	020294-32-0	76
2	Benzene, pentamethyl-	148	C11H16	000700-12-9	70
3	2-tert-Butyltoluene	148	C11H16	001074-92-6	64
4	Benzene, 1-ethyl-3-(1-methylethyl)-	148	C11H16	004920-99-4	58
5	Benzene, 1-ethyl-2,4,5-trimethyl-	148	C11H16	017851-27-3	52



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052816\
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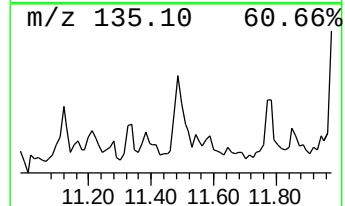
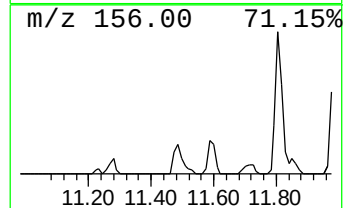
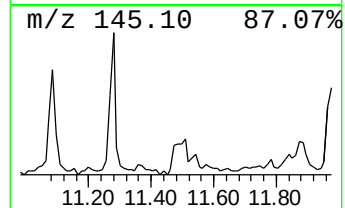
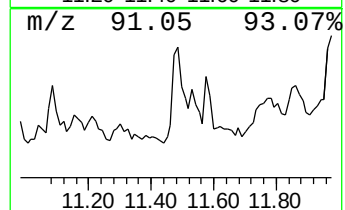
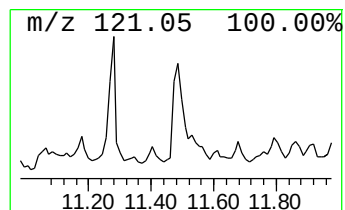
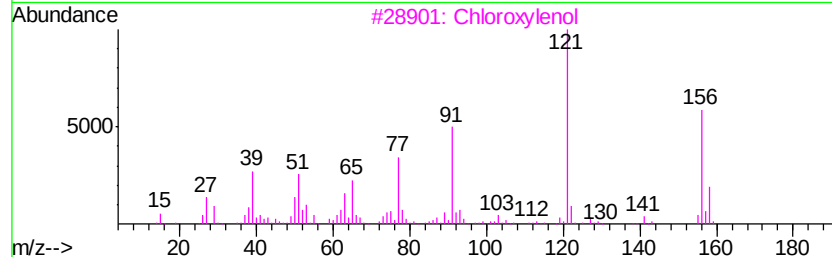
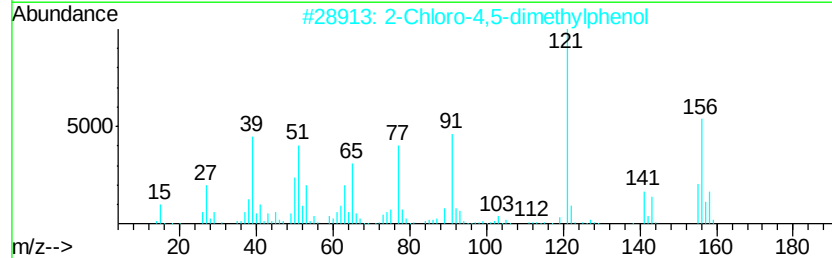
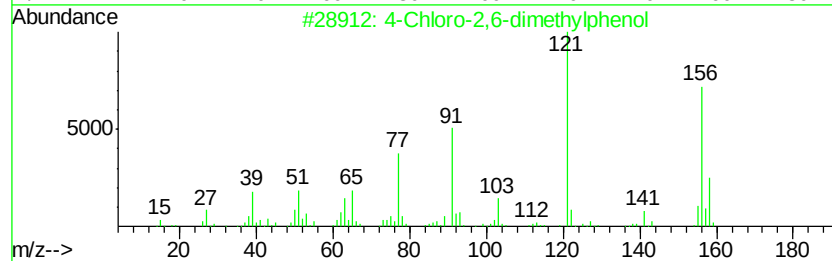
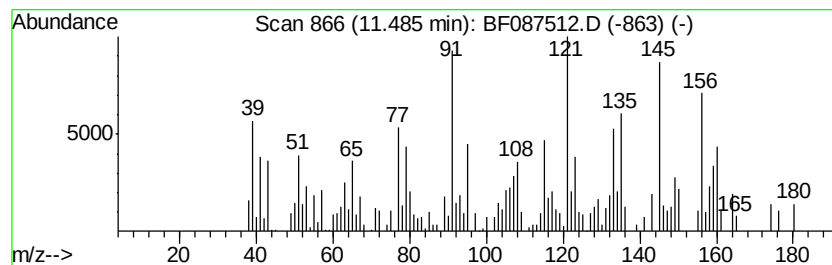
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF052316.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 unknown11.48 Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.48	2.76 ng	158041	Acenaphthene-d10	12.32
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		4-Chloro-2,6-dimethylphenol	156 C8H9ClO	001123-63-3 40
2		2-Chloro-4,5-dimethylphenol	156 C8H9ClO	001124-04-5 38
3		Chloroxlenol	156 C8H9ClO	000088-04-0 25
4		Benzene, 1-(2-butenyl)-2,3-dimet...	160 C12H16	054340-85-1 25
5		1H-Benzimidazole, 2-(1-methyleth...	160 C10H12N2	005851-43-4 25



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052816\
Data File : BF087512.D
Acq On : 29 May 2016 10:00
Operator : UM/SJ
Sample : H3222-01
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-10

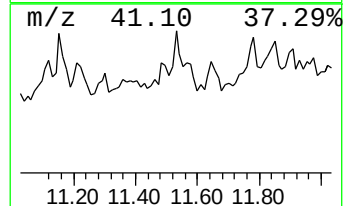
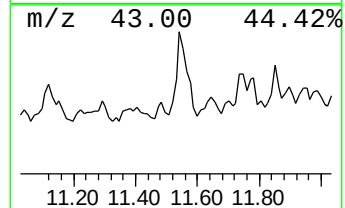
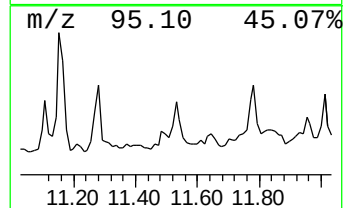
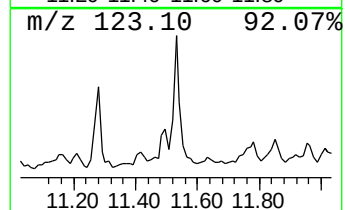
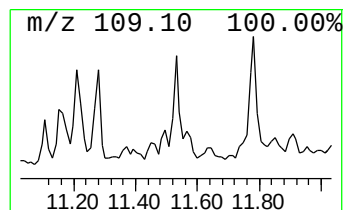
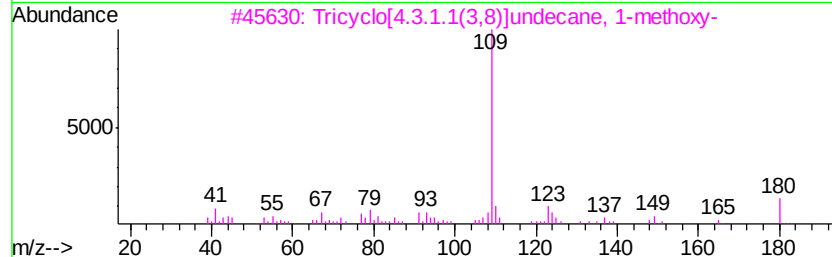
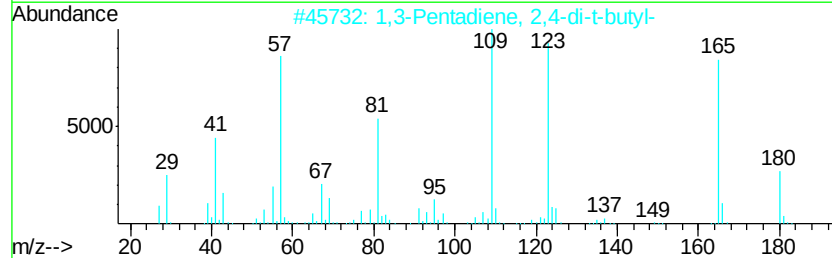
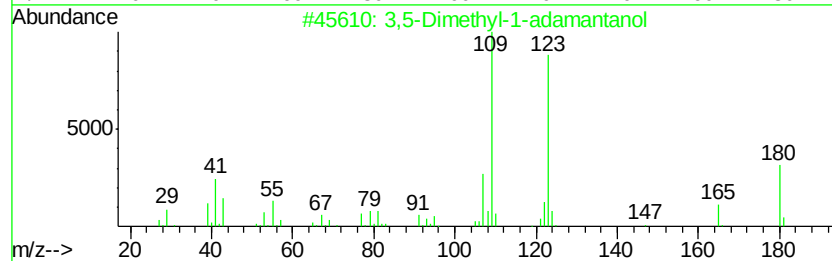
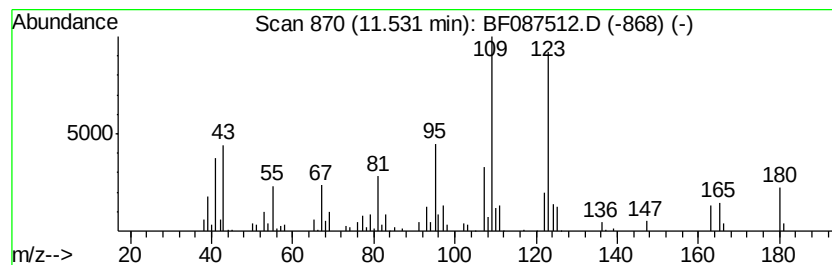
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 3,5-Dimethyl-1-adamantanol Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.53	2.77 ng	158415	Acenaphthene-d10	12.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3,5-Dimethyl-1-adamantanol	180	C12H20O	000707-37-9	68
2		1,3-Pentadiene, 2,4-di-t-butyl-	180	C13H24	132850-21-6	62
3		Tricyclo[4.3.1.1(3,8)]undecane, ...	180	C12H20O	021898-95-3	38
4		Methyl 4-fluorobenzoate	154	C8H7FO2	000403-33-8	35
5		4-Fluorobenzoic acid, 2-ethylcyc...	250	C15H19FO2	1000279-06-0	35



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052816\
Data File : BF087512.D
Acq On : 29 May 2016 10:00
Operator : UM/SJ
Sample : H3222-01
Misc :
ALS Vial : 36 Sample Multiplier: 1

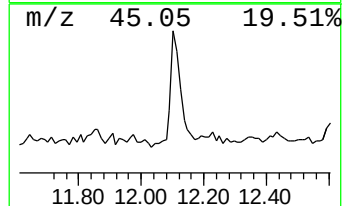
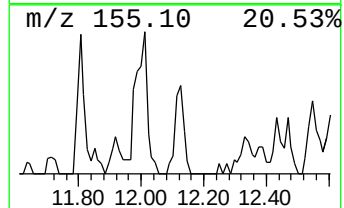
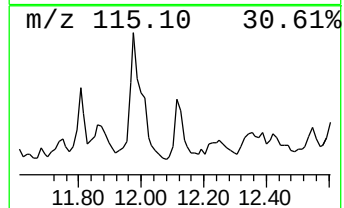
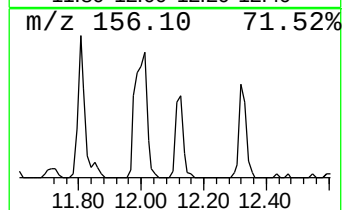
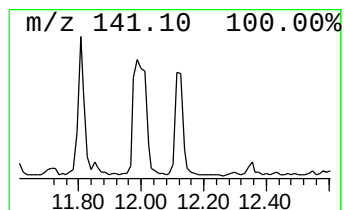
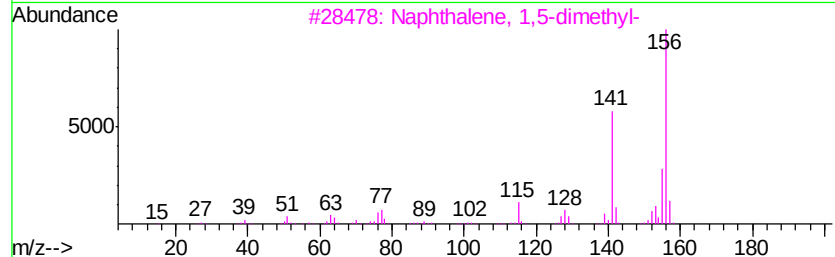
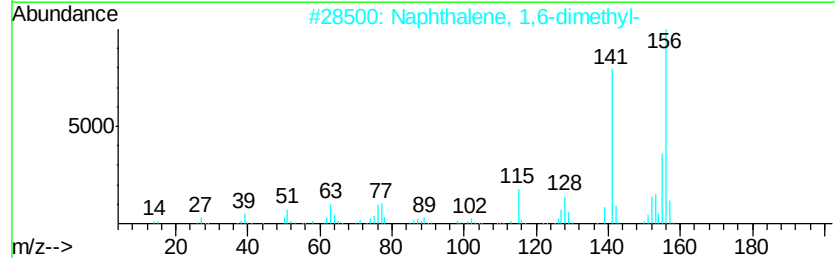
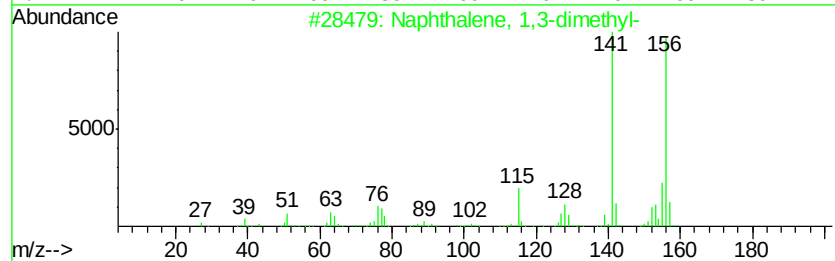
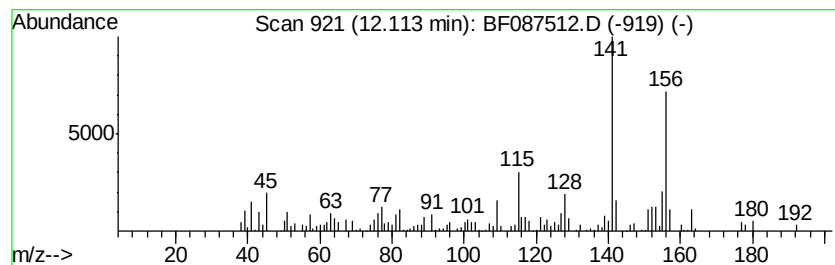
Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF052316.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Naphthalene, 1,3-dimethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD		R.T.
12.11	3.17 ng	181439	Acenaphthene-d10		12.32
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	96
2	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	95
3	Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	95
4	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	94
5	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	94



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052816\
Data File : BF087512.D
Acq On : 29 May 2016 10:00
Operator : UM/SJ
Sample : H3222-01
Misc :
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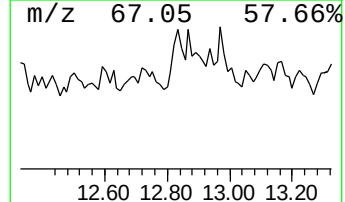
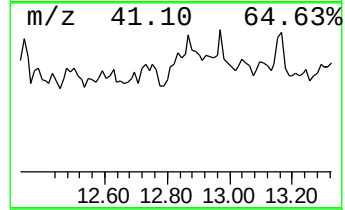
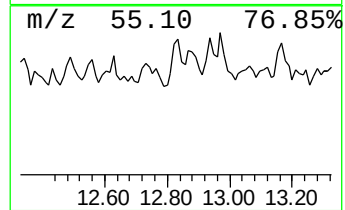
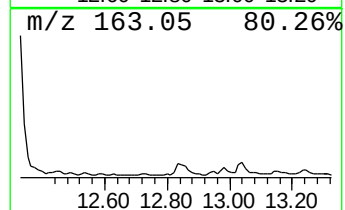
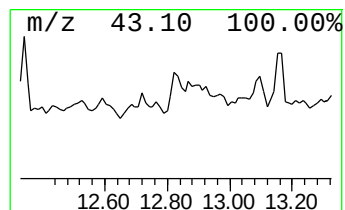
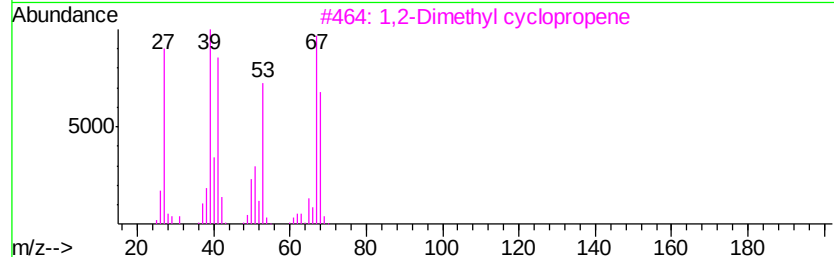
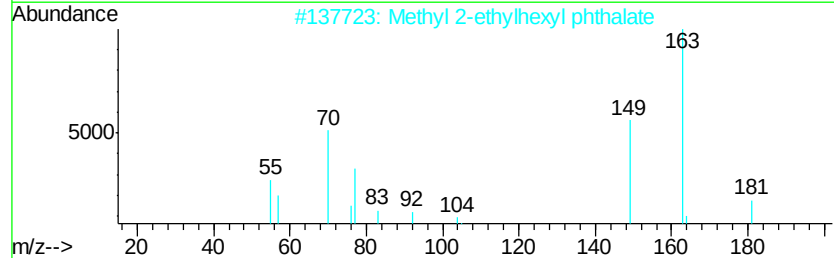
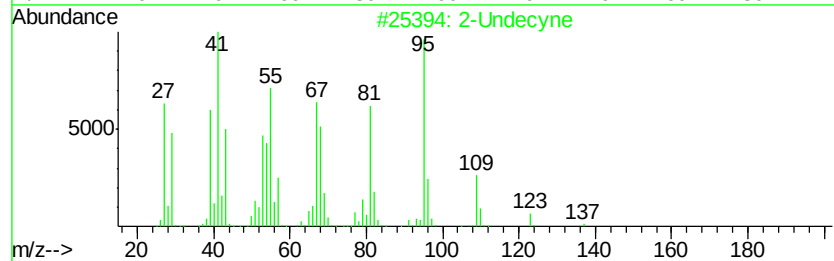
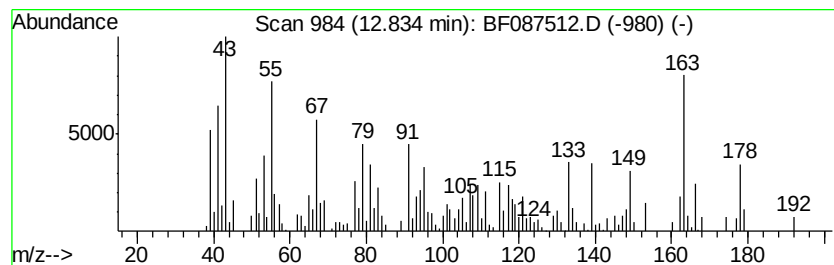
Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF052316.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 unknown12.83 Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.83	2.29 ng	131129	Acenaphthene-d10	12.32	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Undecyne	152	C11H20	060212-29-5	35
2	Methyl 2-ethylhexyl phthalate	292	C17H24O4	056166-83-7	27
3	1,2-Dimethyl cyclopropene	68	C5H8	014309-32-1	27
4	3-Cyclopropenoic acid,1-butyl, m...	154	C9H14O2	067500-40-7	27
5	Pyridine, 2,6-epidioxy-5-ethyl-3...	178	C9H10N2O2	1000263-31-1	25



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052816\
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Sample : H3222-01
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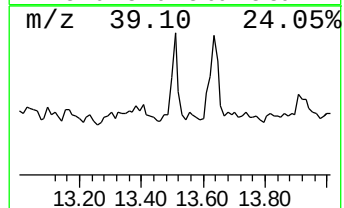
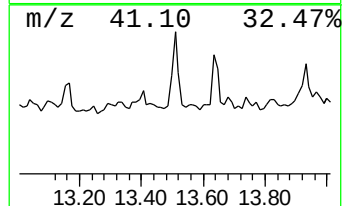
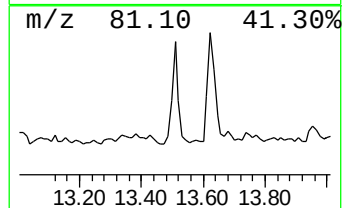
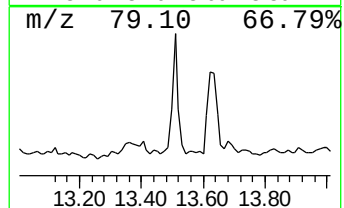
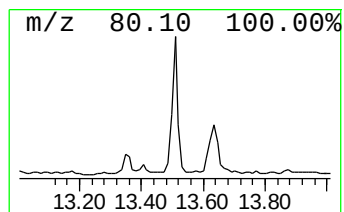
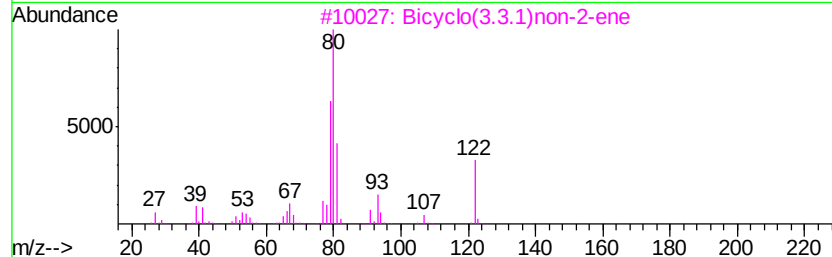
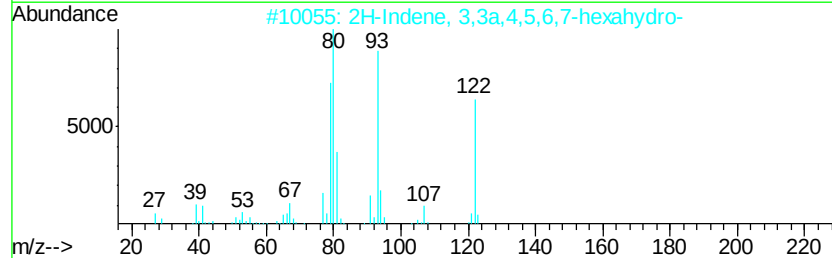
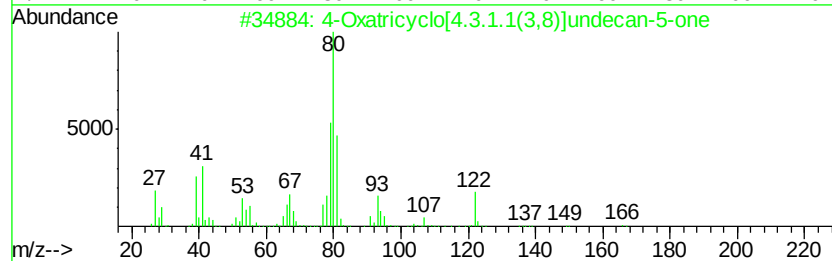
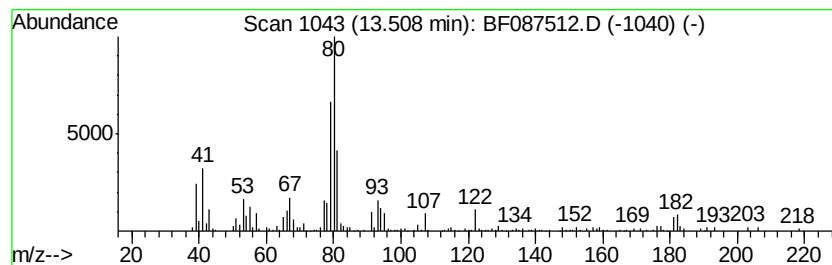
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF052316.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 4-Oxatricyclo[4.3.1.1(3,8)]... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.51	4.00 ng	228670	Acenaphthene-d10	12.32	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Oxatricyclo[4.3.1.1(3,8)]undec...	166	C10H14O2	021898-84-0	91
2	2H-Indene, 3,3a,4,5,6,7-hexahydro-	122	C9H14	016189-41-6	90
3	Bicyclo(3.3.1)non-2-ene	122	C9H14	006671-66-5	90
4	Bicyclo[2.2.2]oct-5-en-2-one	122	C8H10O	002220-40-8	86
5	Bicyclo[3.3.1]nonan-3-ol, endo-	140	C9H16O	010036-10-9	86



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052816\
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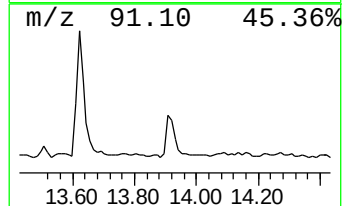
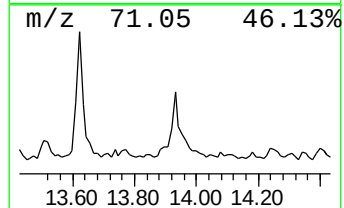
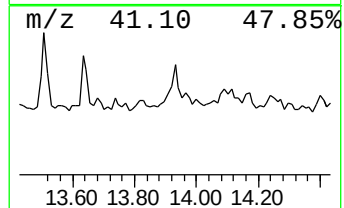
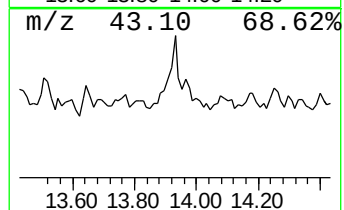
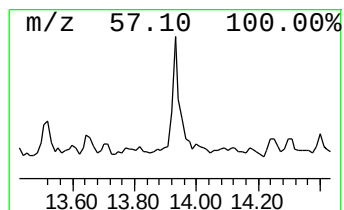
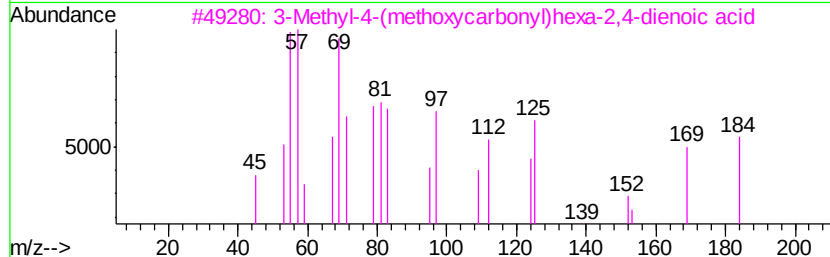
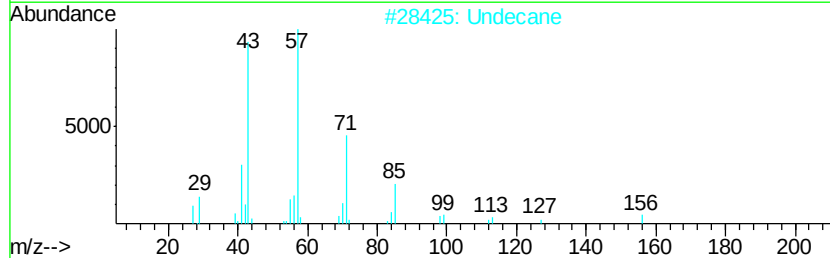
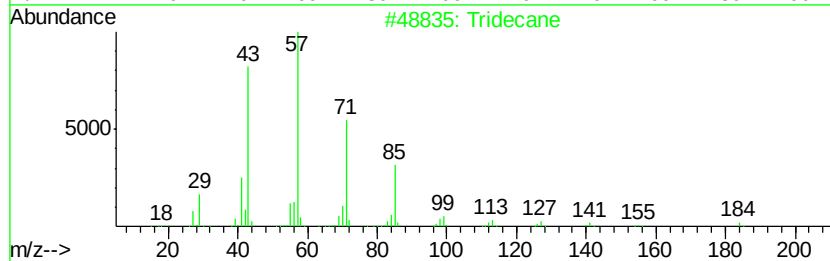
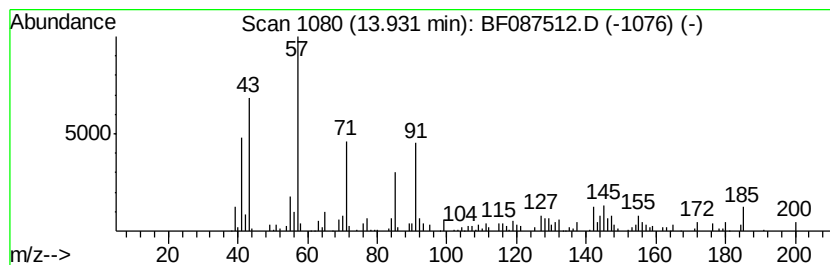
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Tridecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.93	4.30 ng	215223	Phenanthrene-d10	14.73	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecane	184	C13H28	000629-50-5	64
2	Undecane	156	C11H24	001120-21-4	55
3	3-Methyl-4-(methoxycarbonyl)hexa...	184	C9H12O4	1000104-10-8	50
4	Hexadecane	226	C16H34	000544-76-3	49
5	Octane, 2-methyl-	128	C9H20	003221-61-2	49



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052816\
Data File : BF087512.D
Acq On : 29 May 2016 10:00
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Misc :
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Instrument :
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ClientSampleId :
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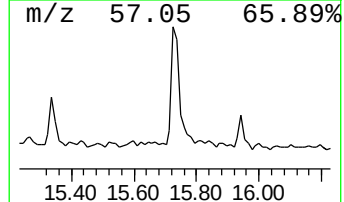
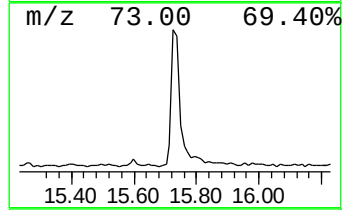
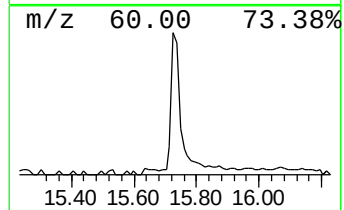
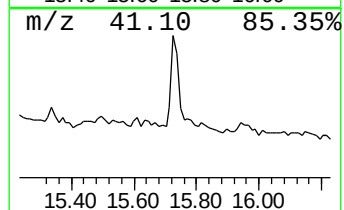
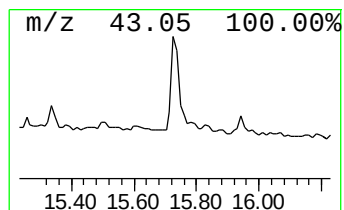
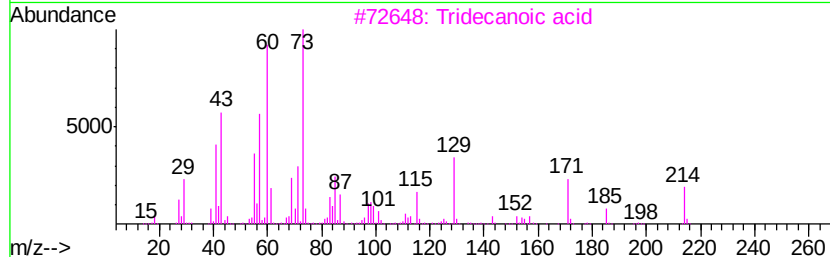
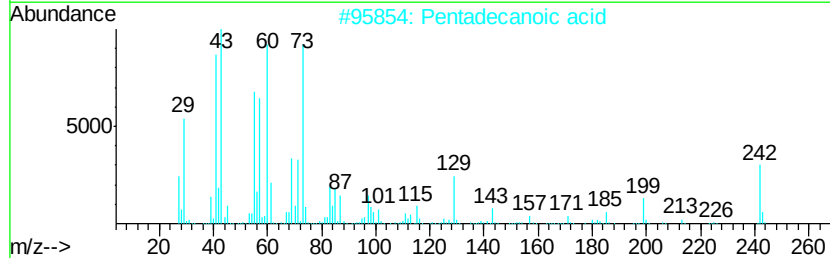
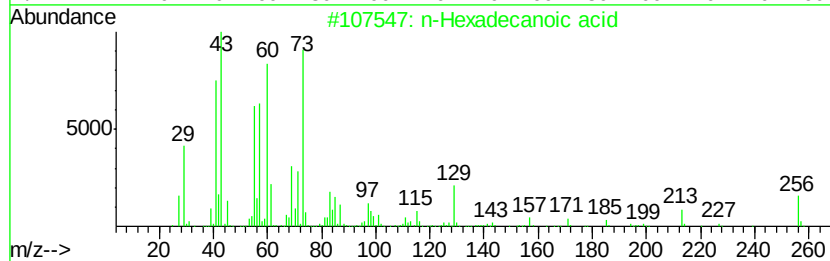
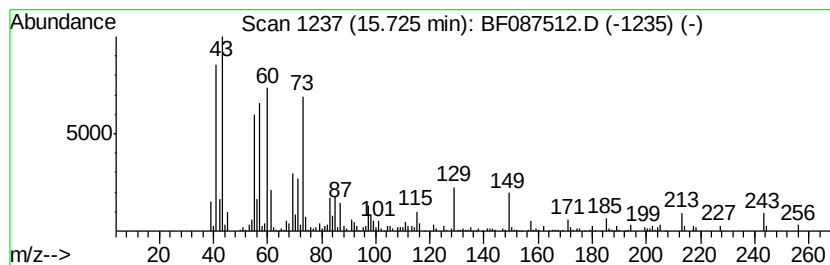
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.73	5.38 ng	269496	Phenanthrene-d10	14.73

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	90
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	83
3	Tridecanoic acid	214	C13H26O2	000638-53-9	76
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	72
5	Dodecanoic acid	200	C12H24O2	000143-07-7	64



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052816\
Data File : BF087512.D
Acq On : 29 May 2016 10:00
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Sample : H3222-01
Misc :
ALS Vial : 36 Sample Multiplier: 1

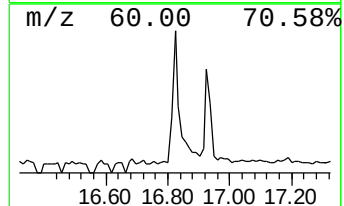
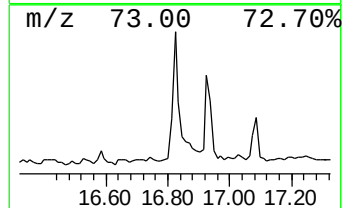
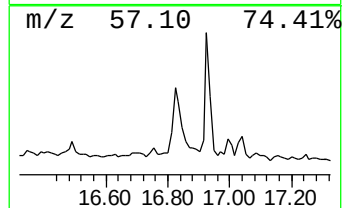
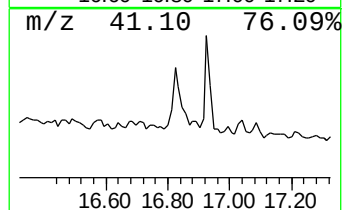
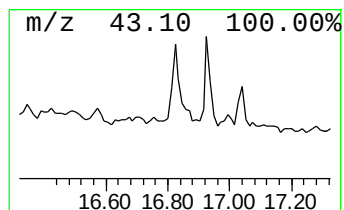
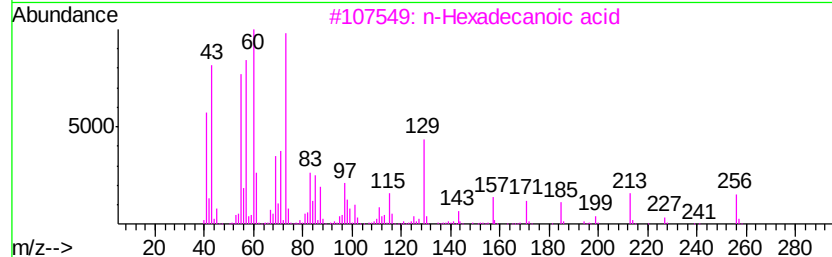
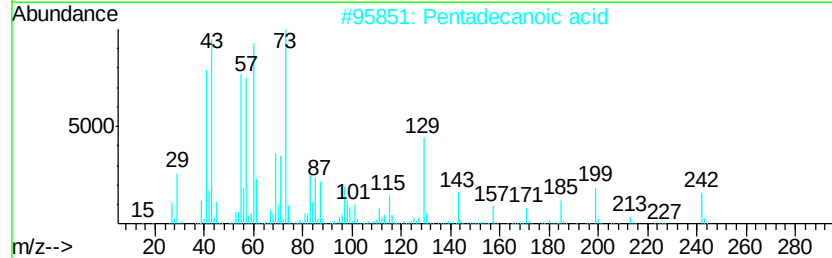
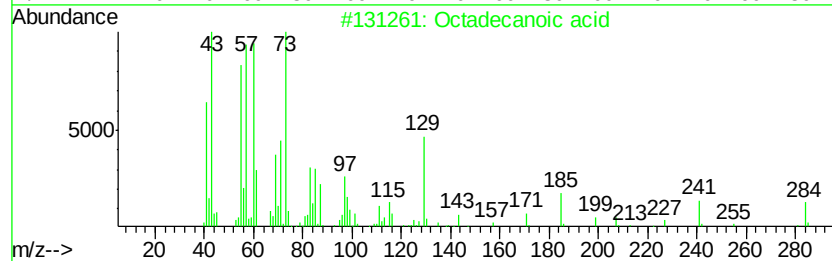
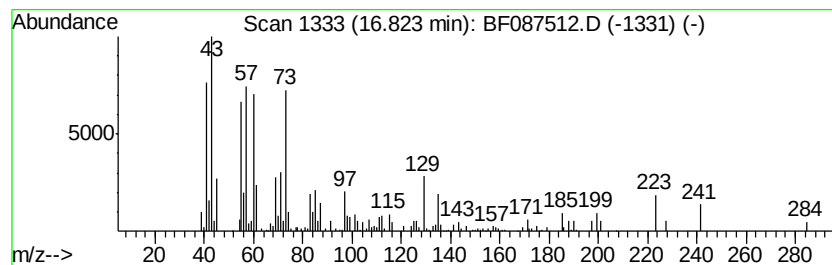
Instrument :
BNA_F
ClientSampleId :
MW-10

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF052316.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 Octadecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD		R.T.
16.82	3.52 ng	146088	Chrysene-d12		18.43
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	68
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	58
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	53
4	Dodecanoic acid	200	C12H24O2	000143-07-7	46
5	Tetradecanoic acid	228	C14H28O2	000544-63-8	43



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052816\
Data File : BF087512.D
Acq On : 29 May 2016 10:00
Operator : UM/SJ
Sample : H3222-01
Misc :
ALS Vial : 36 Sample Multiplier: 1

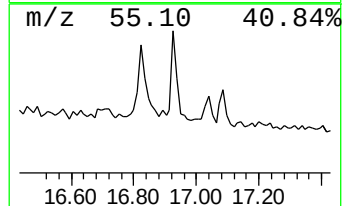
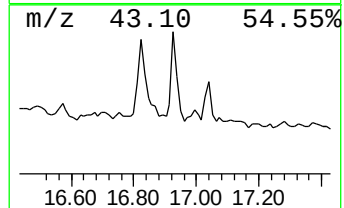
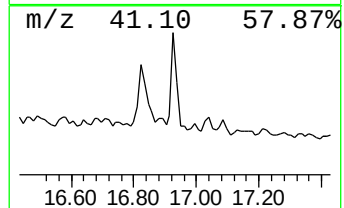
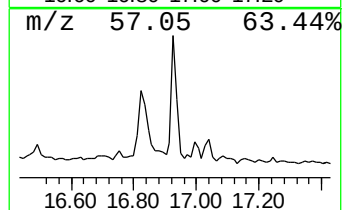
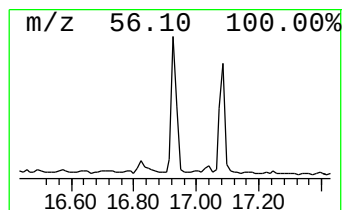
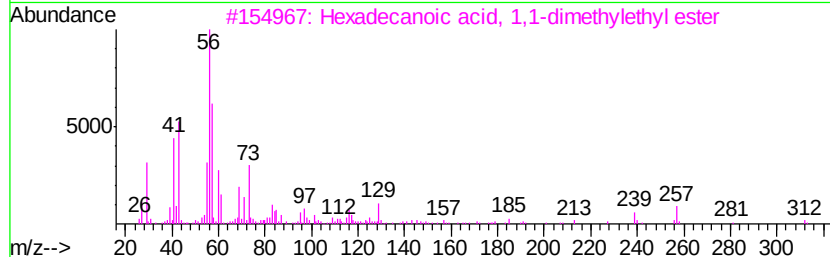
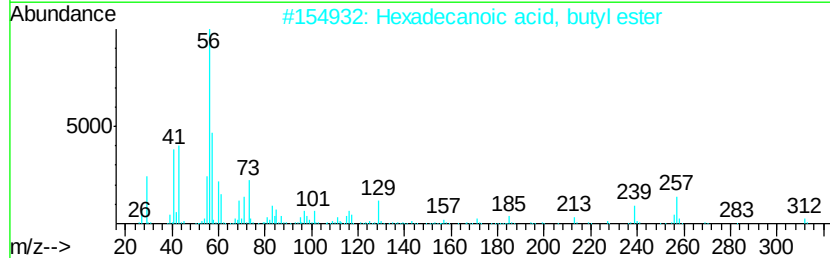
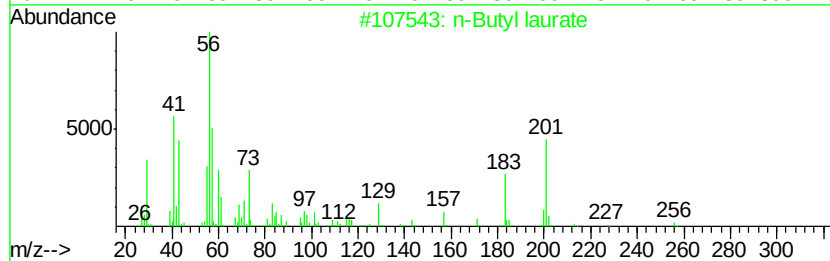
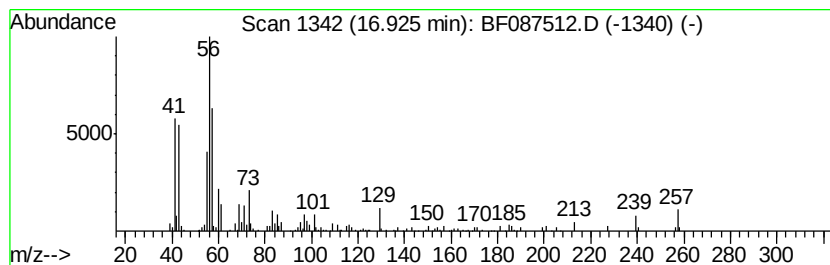
Instrument :
BNA_F
ClientSampleId :
MW-10

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF052316.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 n-Butyl laurate Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.	
16.93	2.86 ng	118544	Chrysene-d12	18.43	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Butyl laurate	256	C16H32O2	000106-18-3	91
2	Hexadecanoic acid, butyl ester	312	C20H40O2	000111-06-8	91
3	Hexadecanoic acid, 1,1-dimethyle...	312	C20H40O2	031158-91-5	58
4	Cyclopentanecarboxylic acid, 2-a...	129	C6H11NO2	040482-05-1	43
5	1-Hexene, 3,4-dimethyl-	112	C8H16	016745-94-1	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052816\
 Data File : BF087512.D
 Acq On : 29 May 2016 10:00
 Operator : UM/SJ
 Sample : H3222-01
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-10

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF052316.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
unknown7.12	7.12	75.3	ng	3269150	1	7.47	867807	20.0
Benzene, 1,3-diet...	7.88	5.2	ng	225066	1	7.47	867807	20.0
Benzene, 1,2-diet...	8.02	6.8	ng	295452	1	7.47	867807	20.0
Benzene, 2-butenyl-	8.32	3.0	ng	128553	1	7.47	867807	20.0
Benzene, 1-ethyl-...	8.72	2.1	ng	144896	2	9.50	1349610	20.0
1-Phenyl-1-butene	8.99	3.2	ng	217552	2	9.50	1349610	20.0
Benzene, (2-methy...	9.60	3.2	ng	218778	2	9.50	1349610	20.0
Benzo[b]thiophene...	10.27	3.4	ng	229601	2	9.50	1349610	20.0
1-Adamantanol	10.34	3.9	ng	264514	2	9.50	1349610	20.0
6-Methyl-4-indanol	10.44	2.4	ng	160656	2	9.50	1349610	20.0
unknown11.48	11.48	2.8	ng	158041	3	12.32	1144030	20.0
3,5-Dimethyl-1-ad...	11.53	2.8	ng	158415	3	12.32	1144030	20.0
Naphthalene, 1,3-...	12.11	3.2	ng	181439	3	12.32	1144030	20.0
unknown12.83	12.83	2.3	ng	131129	3	12.32	1144030	20.0
4-Oxatricyclo[4.3...	13.51	4.0	ng	228670	3	12.32	1144030	20.0
Tridecane	13.93	4.3	ng	215223	4	14.73	1001880	20.0
n-Hexadecanoic acid	15.73	5.4	ng	269496	4	14.73	1001880	20.0
Octadecanoic acid	16.82	3.5	ng	146088	5	18.43	829826	20.0
n-Butyl laurate	16.93	2.9	ng	118544	5	18.43	829826	20.0