

Data Path : Z:\HPCHEM1\BNA F\DATA\BF010417\
 Data File : BF092082.D
 Acq On : 4 Jan 2017 21:50
 Operator : UM/SJ
 Sample : I1004-05
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-36

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-BF122116.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	2.013	8	11	14	rBV2	33775	121092	2.44%	0.118%
2	2.618	61	64	66	rBV	60554	99653	2.01%	0.097%
3	4.801	252	255	258	rVV	517891	851340	17.13%	0.832%
4	4.859	258	260	264	rVB2	76739	142470	2.87%	0.139%
5	4.927	264	266	270	rVB	113055	129193	2.60%	0.126%
6	5.259	293	295	298	rBV	1963006	2554095	51.39%	2.496%
7	5.350	302	303	307	rBV3	112101	243908	4.91%	0.238%
8	5.407	307	308	312	rVV3	156282	243257	4.89%	0.238%
9	5.476	312	314	316	rVV2	203332	239980	4.83%	0.235%
10	5.521	316	318	319	rVV2	114477	119803	2.41%	0.117%
11	5.544	319	320	323	rVB	218257	221885	4.46%	0.217%
12	5.647	326	329	331	rBV2	62417	120077	2.42%	0.117%
13	5.681	331	332	335	rBV	281476	380487	7.66%	0.372%
14	5.807	339	343	345	rBV	336365	416878	8.39%	0.407%
15	5.853	345	347	348	rBV	108395	158394	3.19%	0.155%
16	5.899	348	351	354	rVV3	189483	465728	9.37%	0.455%
17	5.979	354	358	362	rVB4	197300	378689	7.62%	0.370%
18	6.059	362	365	368	rBV4	162913	394946	7.95%	0.386%
19	6.104	368	369	372	rVB	406240	524300	10.55%	0.512%
20	6.173	372	375	378	rBV3	340347	632207	12.72%	0.618%
21	6.264	381	383	384	rVV	409384	535204	10.77%	0.523%
22	6.299	384	386	387	rVV	506137	744705	14.98%	0.728%
23	6.333	387	389	393	rVV2	1286437	2476587	49.83%	2.421%
24	6.436	395	398	401	rVV	4308809	4970080	100.00%	4.858%
25	6.504	403	404	406	rVV2	261078	381854	7.68%	0.373%
26	6.539	406	407	410	rVV2	241353	515325	10.37%	0.504%
27	6.619	410	414	416	rVV3	473978	1310490	26.37%	1.281%
28	6.664	416	418	422	rVV3	1079481	1788212	35.98%	1.748%
29	6.824	429	432	433	rBV	1797209	2626863	52.85%	2.567%
30	6.916	438	440	442	rVB	336864	312856	6.29%	0.306%
31	7.007	445	448	451	rBV3	285614	546019	10.99%	0.534%
32	7.053	451	452	455	rBV3	132658	246027	4.95%	0.240%
33	7.110	455	457	460	rVB	400311	439142	8.84%	0.429%
34	7.202	460	465	466	rBV2	630593	1488291	29.95%	1.455%

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35	7.236	466	468	471	rVB	1927539	2055403	41.36%	2.009%
36	7.316	473	475	477	rBV3	153693	210978	4.24%	0.206%
37	7.396	480	482	483	rVB	153052	206013	4.15%	0.201%
38	7.442	483	486	488	rBV	724225	875230	17.61%	0.855%
39	7.533	491	494	496	rBV3	150607	291653	5.87%	0.285%
40	7.567	496	497	499	rBV2	107320	184961	3.72%	0.181%
41	7.625	499	502	504	rVB4	316637	680910	13.70%	0.665%
42	7.693	504	508	510	rBV2	979814	1510597	30.39%	1.476%
43	7.727	510	511	514	rVB3	77093	97459	1.96%	0.095%
44	7.785	514	516	518	rBV2	320728	455001	9.15%	0.445%
45	7.842	519	521	522	rVB2	127369	144975	2.92%	0.142%
46	7.899	524	526	528	rBV3	157235	186595	3.75%	0.182%
47	7.956	528	531	534	rBV3	1126117	2467049	49.64%	2.411%
48	8.070	538	541	543	rVB3	245437	457623	9.21%	0.447%
49	8.116	543	545	546	rBV	222855	290921	5.85%	0.284%
50	8.150	546	548	549	rBV2	297482	424473	8.54%	0.415%
51	8.185	549	551	554	rVB2	296601	504487	10.15%	0.493%
52	8.265	554	558	559	rBV4	150877	343209	6.91%	0.335%
53	8.310	559	562	563	rBV2	521256	759749	15.29%	0.743%
54	8.390	566	569	570	rBV	491405	777524	15.64%	0.760%
55	8.470	574	576	577	rVV2	625925	906183	18.23%	0.886%
56	8.493	577	578	580	rVB	1130897	793154	15.96%	0.775%
57	8.562	582	584	585	rVB2	230951	315075	6.34%	0.308%
58	8.607	585	588	590	rBV3	554769	1314794	26.45%	1.285%
59	8.653	590	592	594	rVB2	1481526	1614410	32.48%	1.578%
60	8.767	594	602	603	rBV4	1683959	3872839	77.92%	3.785%
61	8.802	603	605	606	rVB2	413504	499798	10.06%	0.488%
62	8.836	606	608	609	rBV2	311502	547412	11.01%	0.535%
63	8.905	613	614	616	rVV2	420432	493725	9.93%	0.483%
64	8.985	616	621	623	rVV3	1093050	2857968	57.50%	2.793%
65	9.030	623	625	631	rVB	2625539	4744776	95.47%	4.637%
66	9.202	638	640	641	rBV2	379142	536717	10.80%	0.525%
67	9.362	653	654	656	rBV2	511776	705359	14.19%	0.689%
68	9.442	659	661	662	rVV2	436276	670384	13.49%	0.655%
69	9.476	662	664	666	rVV	969504	1837575	36.97%	1.796%
70	9.568	670	672	673	rBV2	227769	248689	5.00%	0.243%
71	9.705	680	684	685	rVB2	959427	1808466	36.39%	1.768%

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72	9.750	685	688	694	rBV2	1735788	4261950	85.75%	4.165%
73	9.830	694	695	696	rVB	359714	246764	4.96%	0.241%
74	9.888	696	700	701	rBV2	536589	878740	17.68%	0.859%
75	9.933	702	704	705	rVB	351867	391840	7.88%	0.383%
76	10.025	710	712	715	rVV2	498057	958355	19.28%	0.937%
77	10.105	717	719	721	rVV2	772101	1718332	34.57%	1.679%
78	10.173	721	725	728	rVB4	1658298	4088600	82.26%	3.996%
79	10.242	728	731	735	rBV4	1201389	2300965	46.30%	2.249%
80	10.310	735	737	738	rBV2	253966	281783	5.67%	0.275%
81	10.368	739	742	746	rBV5	446253	1409279	28.36%	1.377%
82	10.505	751	754	756	rVB2	1664662	1975299	39.74%	1.931%
83	10.539	756	757	760	rBV3	494628	682257	13.73%	0.667%
84	10.733	772	774	775	rBV2	454032	658559	13.25%	0.644%
85	10.779	776	778	780	rVV2	510371	836974	16.84%	0.818%
86	10.825	780	782	784	rVV3	271622	449827	9.05%	0.440%
87	10.871	784	786	787	rVV	347266	478716	9.63%	0.468%
88	10.951	790	793	795	rBV3	387165	886751	17.84%	0.867%
89	11.076	802	804	807	rBV4	1234908	3114027	62.66%	3.044%
90	11.191	811	814	818	rVB3	1182653	2733686	55.00%	2.672%
91	11.465	836	838	839	rBV2	469069	532964	10.72%	0.521%
92	11.545	843	845	847	rVB	687885	944362	19.00%	0.923%
93	11.751	862	863	865	rBV2	364781	508106	10.22%	0.497%
94	11.785	865	866	872	rVB2	607202	851929	17.14%	0.833%
95	12.276	908	909	914	rBV4	195063	700226	14.09%	0.684%
96	12.516	928	930	943	rVB	758993	1332396	26.81%	1.302%
97	12.768	949	952	954	rVB	4626674	4515528	90.85%	4.413%
98	13.659	1028	1030	1035	rVB	84233	135012	2.72%	0.132%
99	13.808	1040	1043	1045	rBV	482078	525367	10.57%	0.513%
100	15.180	1160	1163	1166	rVB	337570	404873	8.15%	0.396%

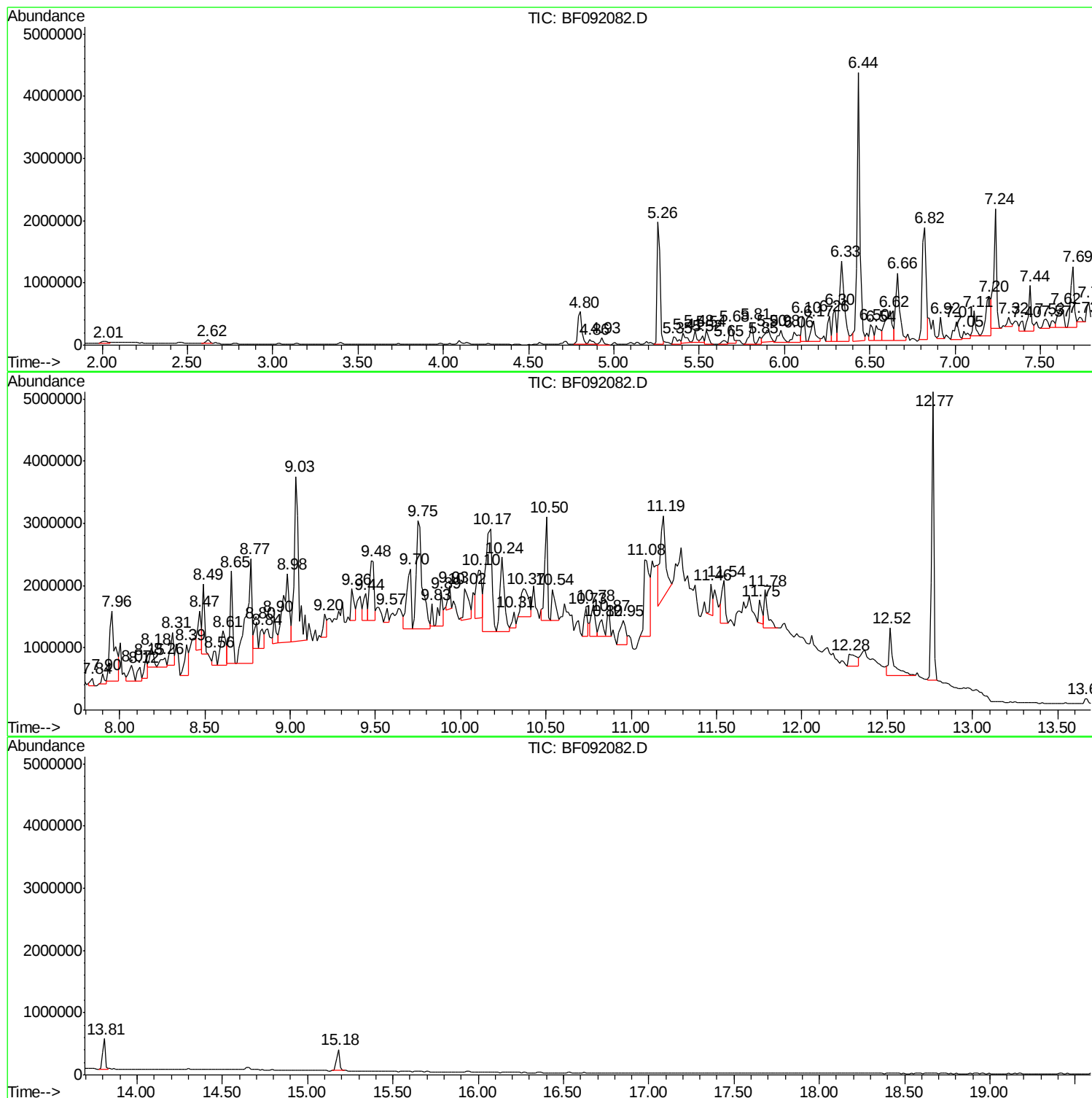
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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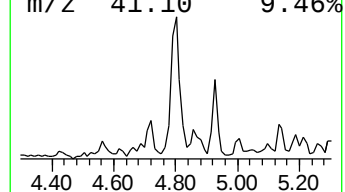
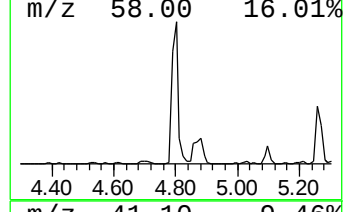
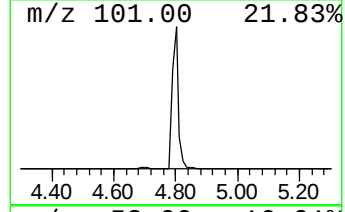
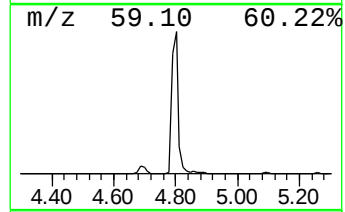
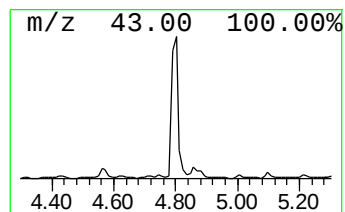
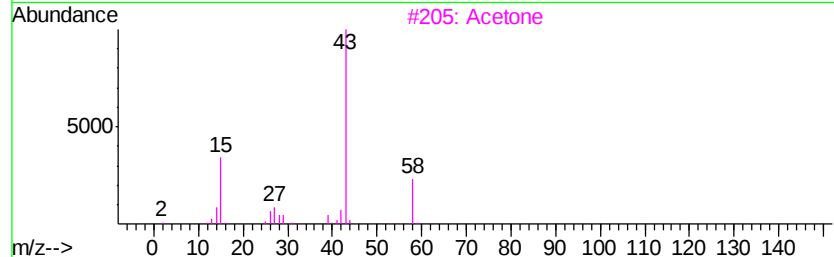
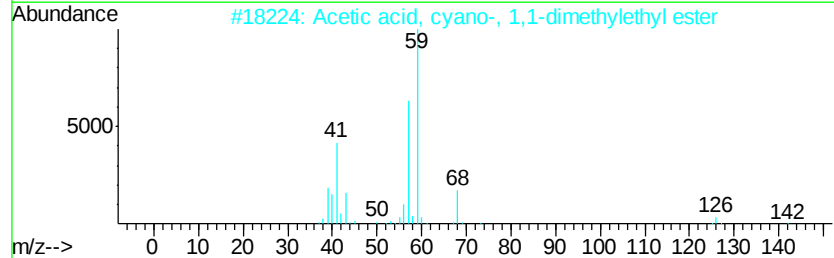
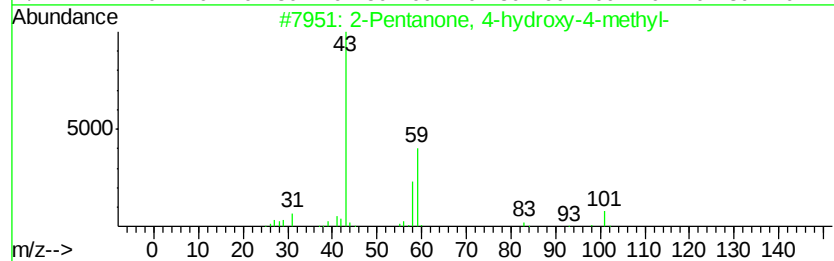
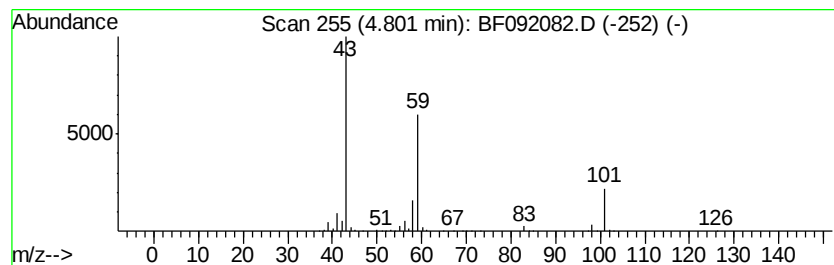
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF122116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.80	9.52 ng	851340	1,4-Dichlorobenzene-d4	6.66

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2	Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17
3	Acetone	58	C3H6O	000067-64-1	10
4	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5	5-Oxohexanenitrile	111	C6H9NO	010412-98-3	9



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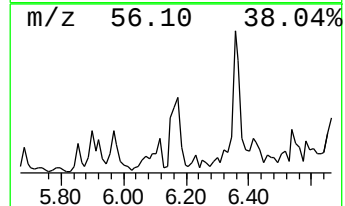
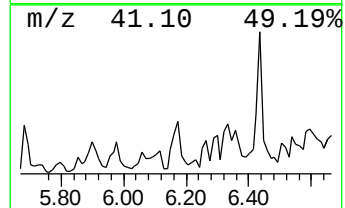
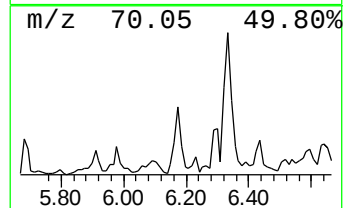
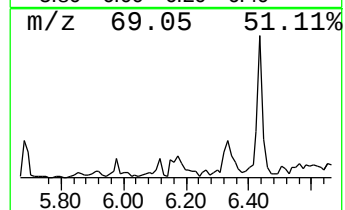
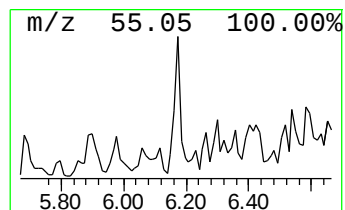
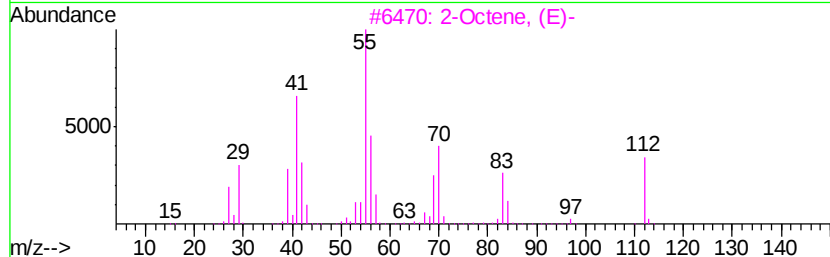
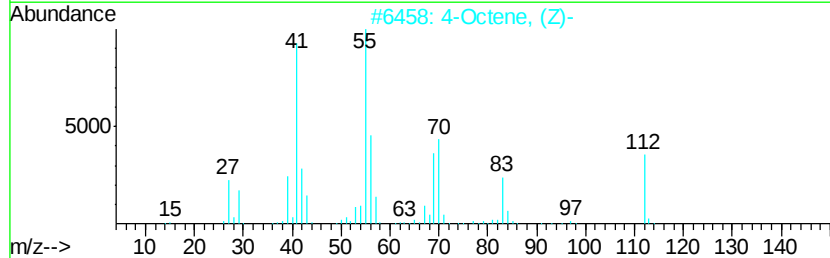
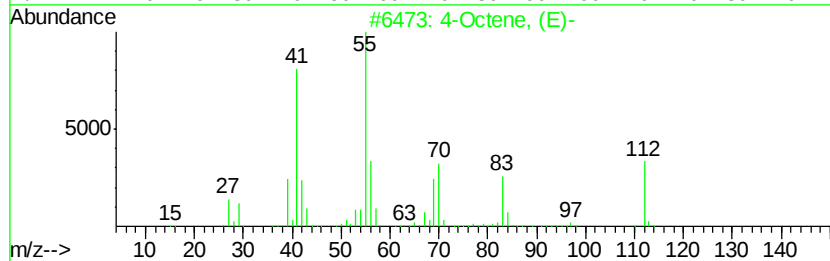
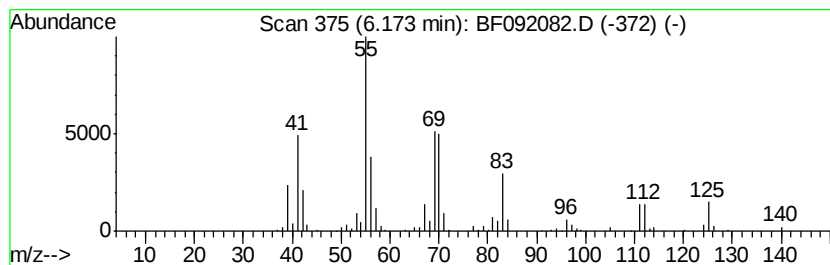
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 4-Octene, (E)- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.17	7.07 ng	632207	1,4-Dichlorobenzene-d4	6.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Octene, (E)-	112	C8H16	014850-23-8	81
2		4-Octene, (Z)-	112	C8H16	007642-15-1	74
3		2-Octene, (E)-	112	C8H16	013389-42-9	74
4		3-Octene, (Z)-	112	C8H16	014850-22-7	62
5		2-Octene, (Z)-	112	C8H16	007642-04-8	58



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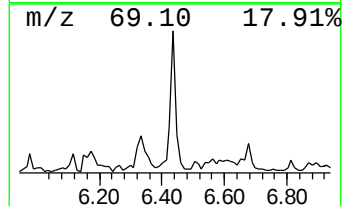
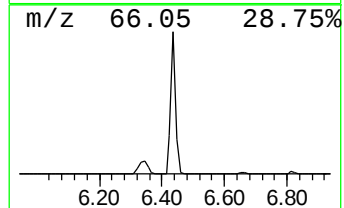
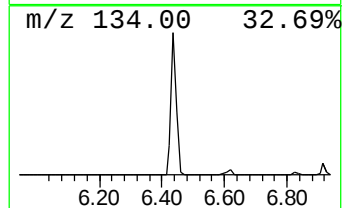
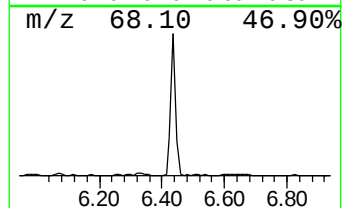
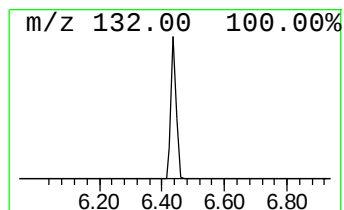
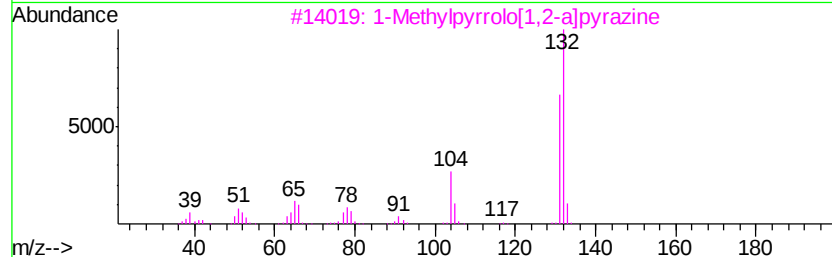
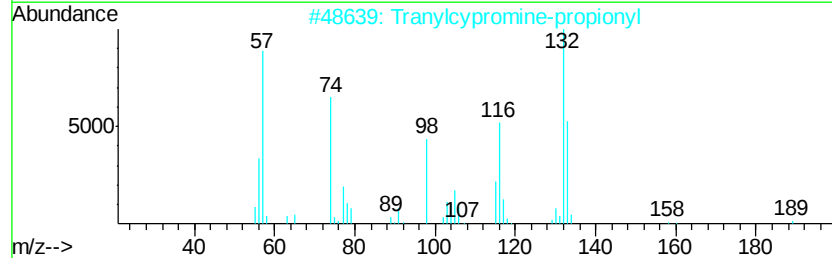
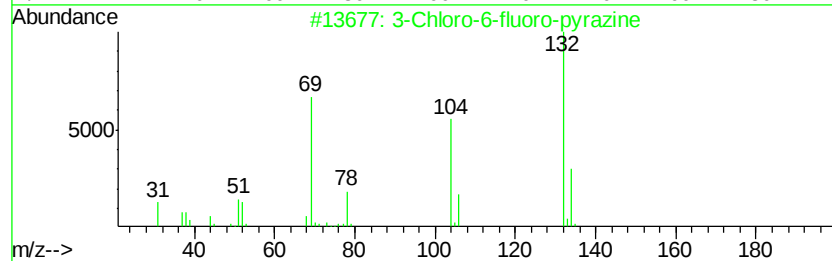
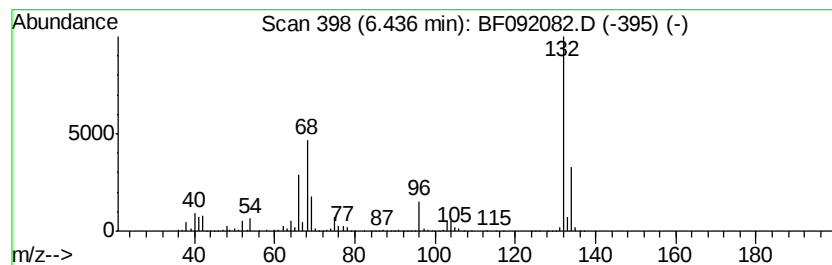
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Peak Number 3 unknown6.44 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.44	55.59 ng	4970080	1,4-Dichlorobenzene-d4	6.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9
4		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF010417\
Data File : BF092082.D
Acq On : 4 Jan 2017 21:50
Operator : UM/SJ
Sample : I1004-05
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
MW-36

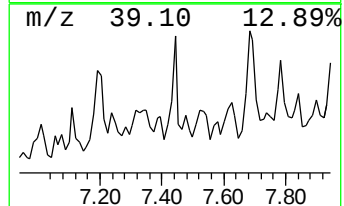
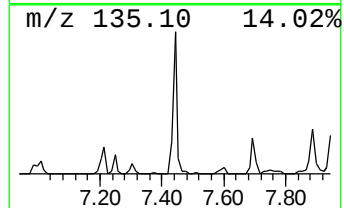
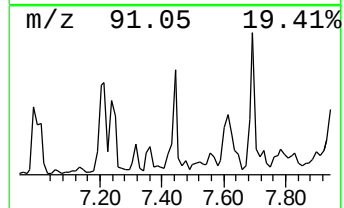
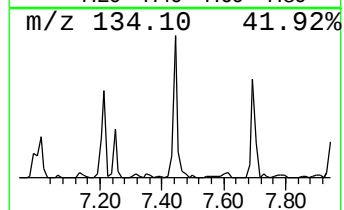
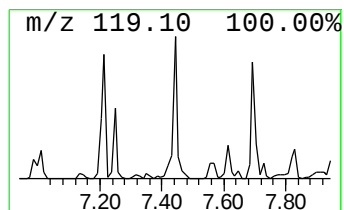
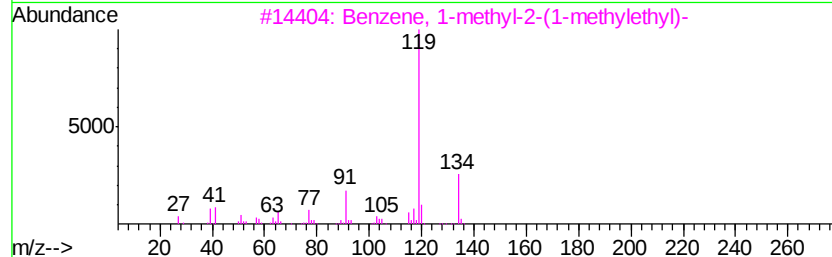
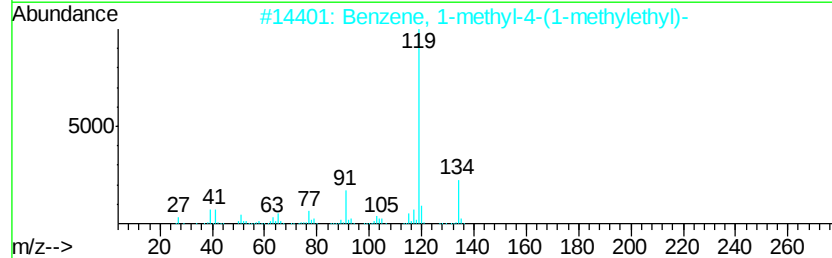
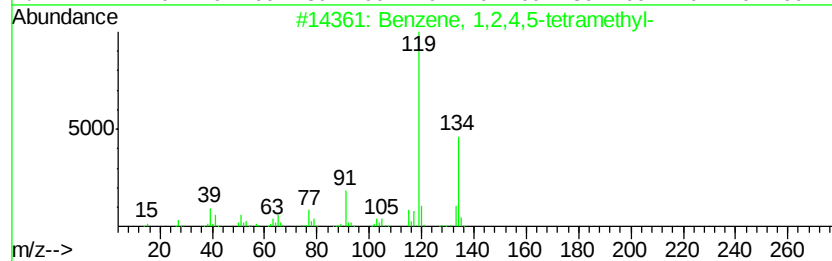
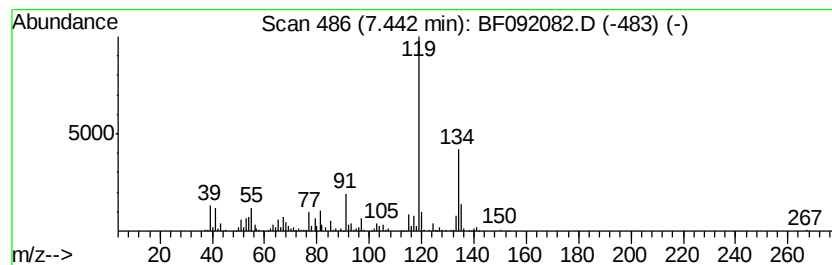
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.44	7.10 ng	875230	Naphthalene-d8	7.96

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	93
2			Benzene, 1-methyl-4-(1-methylethyl)-	134	C10H14	000099-87-6	87
3			Benzene, 1-methyl-2-(1-methylethyl)-	134	C10H14	000527-84-4	87
4			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	87
5			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	87



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF010417\
Data File : BF092082.D
Acq On : 4 Jan 2017 21:50
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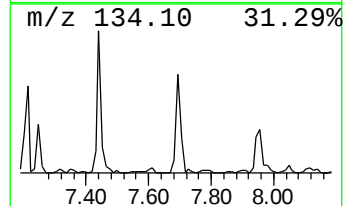
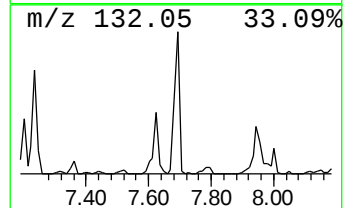
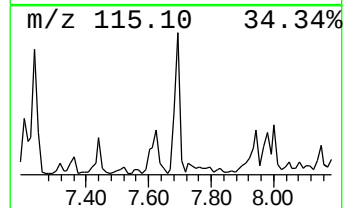
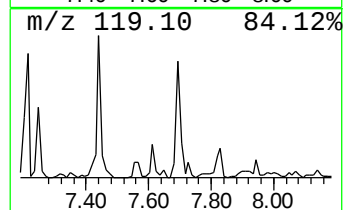
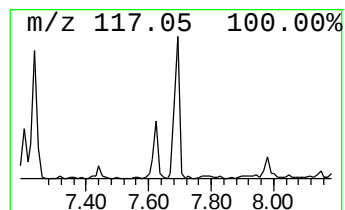
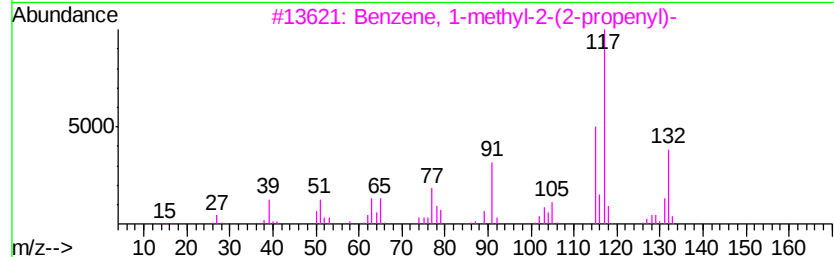
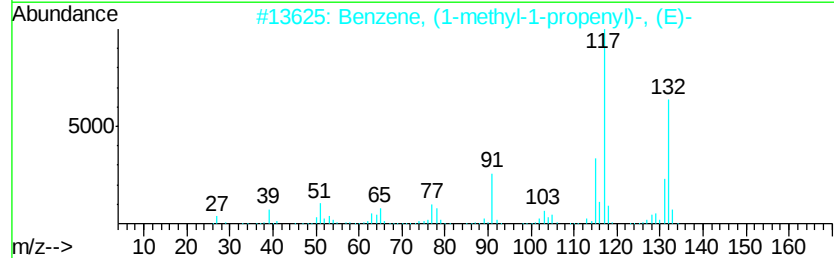
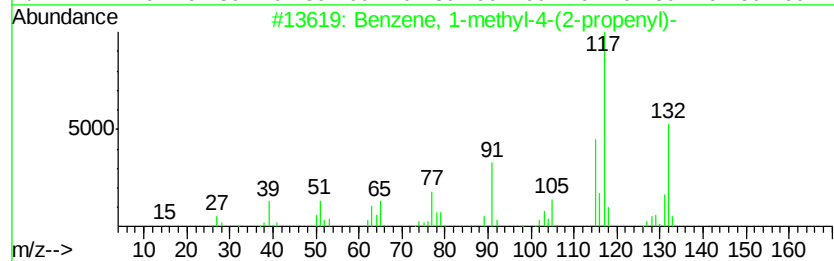
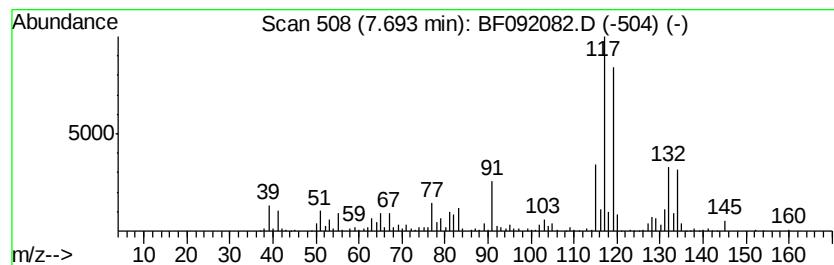
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Benzene, 1-methyl-4-(2-prop... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.69	12.25 ng	1510600	Napthalene-d8	7.96

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	55
2		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	55
3		Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	55
4		1-Phenyl-1-butene	132	C10H12	000824-90-8	55
5		Benzene, 1-methyl-2-(1-methyleth...	134	C10H14	000527-84-4	46



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF010417\
Data File : BF092082.D
Acq On : 4 Jan 2017 21:50
Operator : UM/SJ
Sample : I1004-05
Misc :
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Instrument :
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ClientSampleId :
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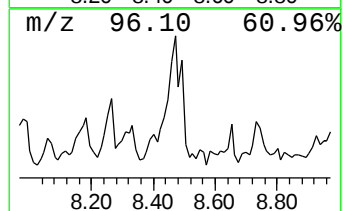
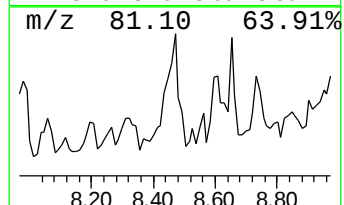
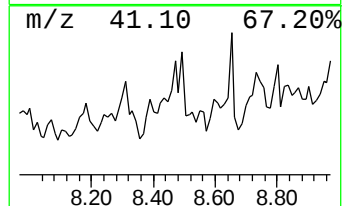
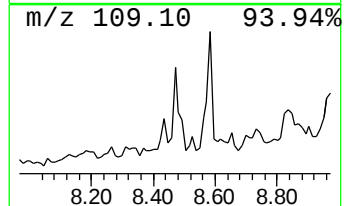
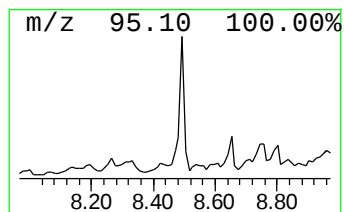
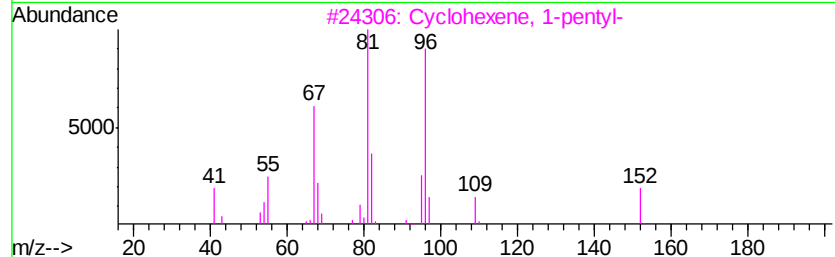
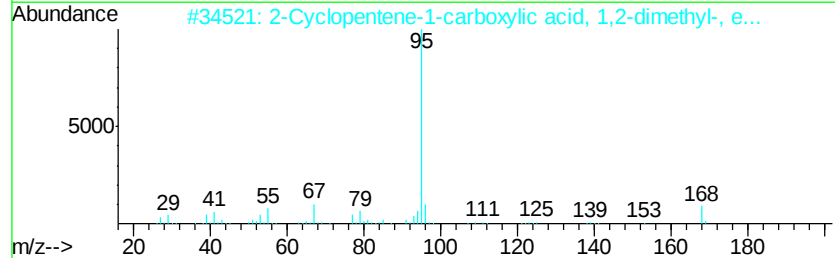
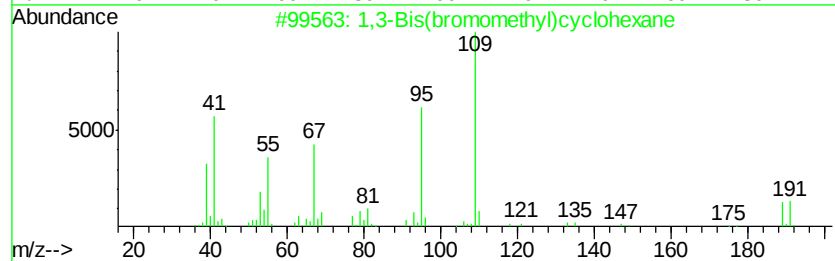
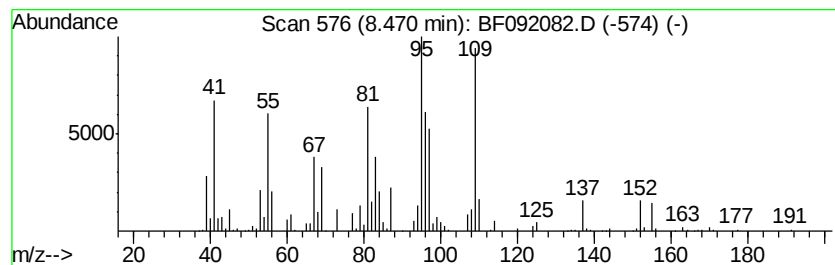
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 unknown8.47 Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.47	7.35 ng	906183	Naphthalene-d8	7.96

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,3-Bis(bromomethyl)cyclohexane	268	C8H14Br2	1000216-89-9	35
2			2-Cyclopentene-1-carboxylic acid...	168	C10H16O2	005809-03-0	30
3			Cyclohexene, 1-pentyl-	152	C11H20	015232-85-6	30
4			1,3-Dimethyl-1-cyclohexene	110	C8H14	002808-76-6	25
5			5-Octen-2-yn-4-ol	124	C8H12O	1000196-87-1	25



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF010417\
Data File : BF092082.D
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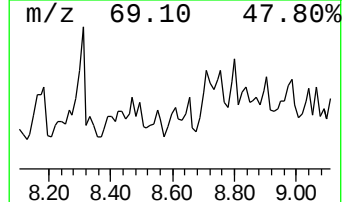
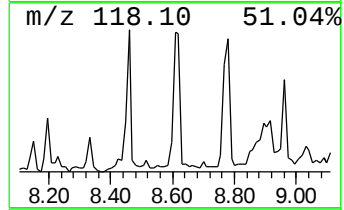
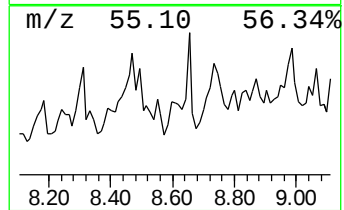
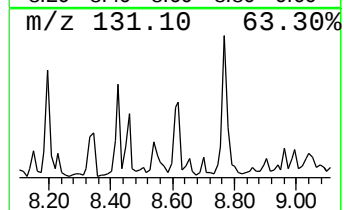
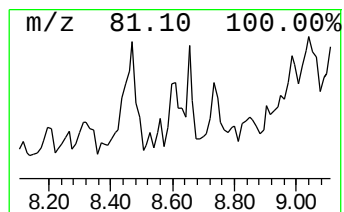
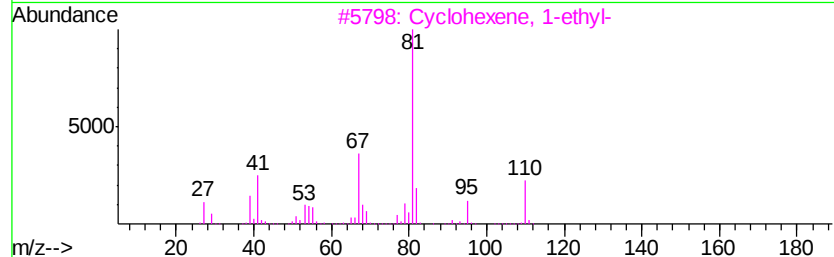
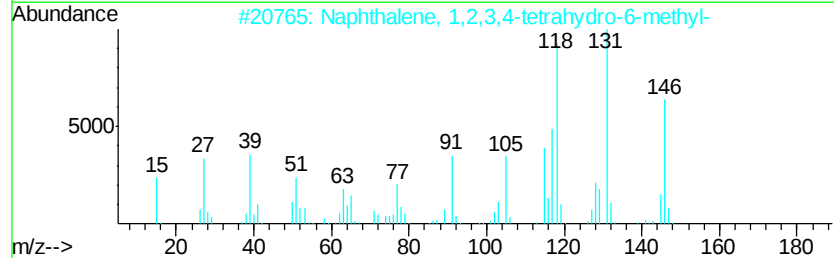
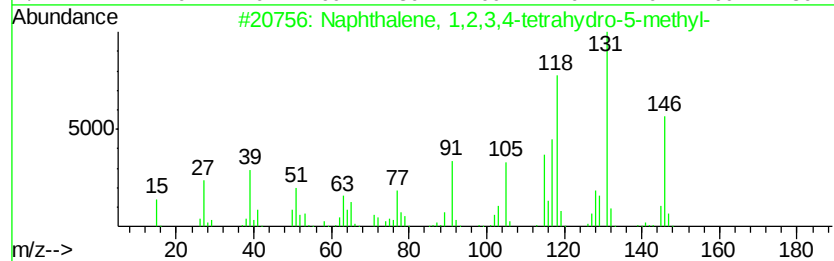
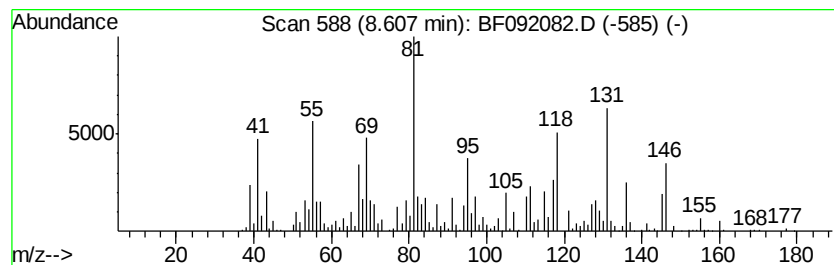
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Naphthalene, 1,2,3,4-tetra... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.61	10.66 ng	1314790	Naphthalene-d8	7.96

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	89
2		Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	60
3		Cyclohexene, 1-ethyl-	110	C8H14	001453-24-3	35
4		Cyclohexene, 4-ethyl-	110	C8H14	003742-42-5	35
5		Cyclohexene, 3-ethyl-	110	C8H14	002808-71-1	20



Data Path : Z:\HPCHEM1\BNA F\DATA\BF010417\
Data File : BF092082.D
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Instrument :
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ClientSampleId :
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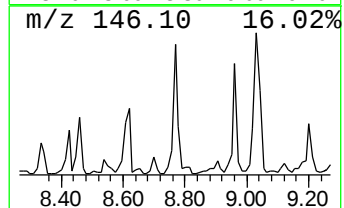
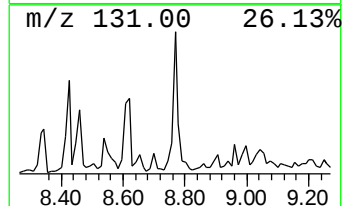
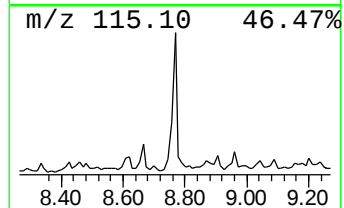
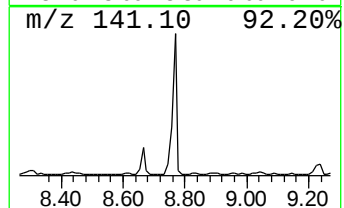
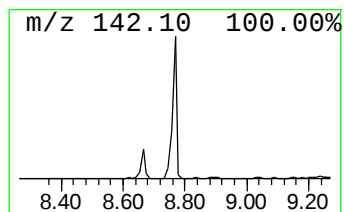
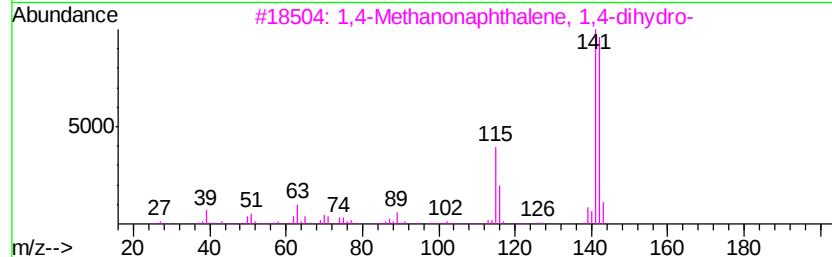
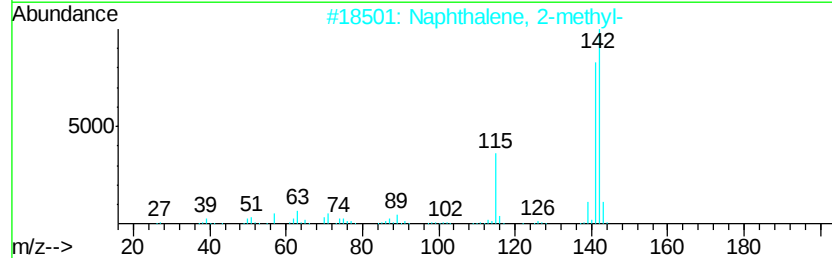
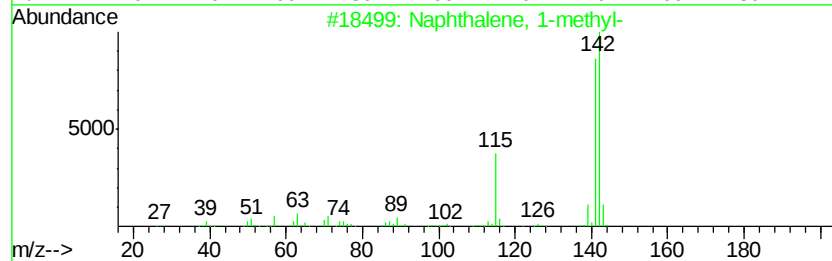
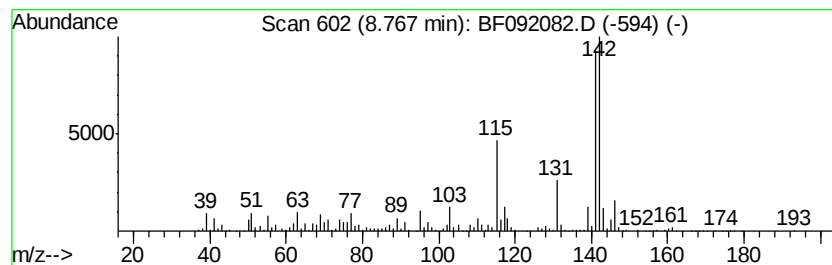
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Naphthalene, 1-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.77	31.40 ng	3872840	Naphthalene-d8	7.96

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	95
2		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	95
3		1,4-Methanonaphthalene, 1,4-dihv...	142	C11H10	004453-90-1	89
4		Benzocycloheptatriene	142	C11H10	000264-09-5	76
5		1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	76



Data Path : Z:\HPCHEM1\BNA F\DATA\BF010417\
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ClientSampleId :
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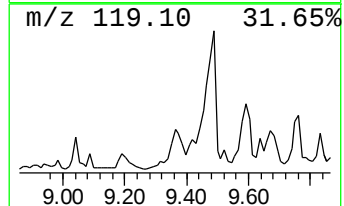
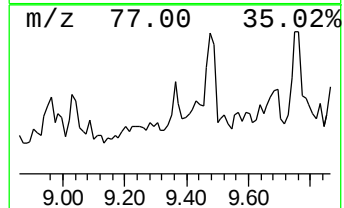
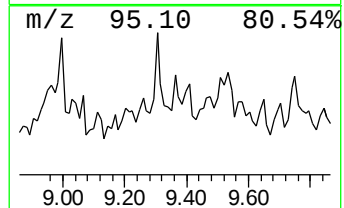
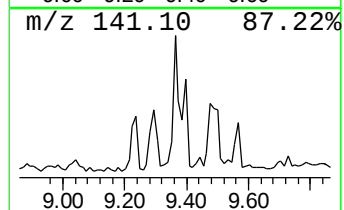
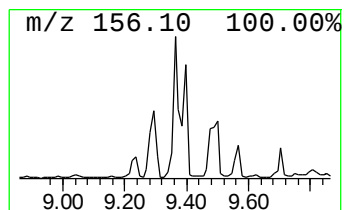
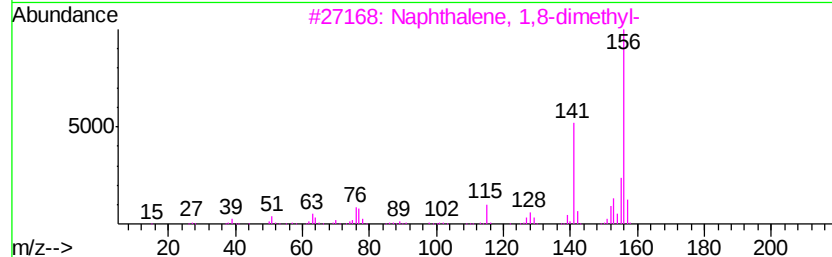
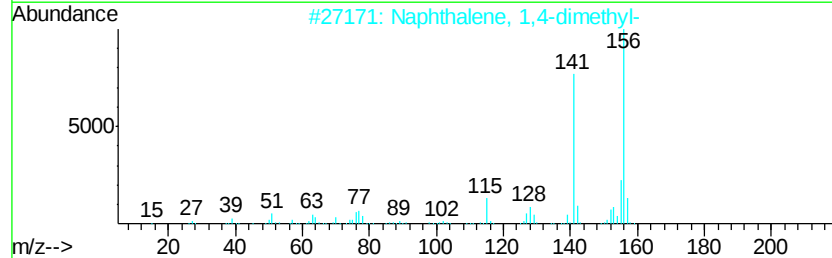
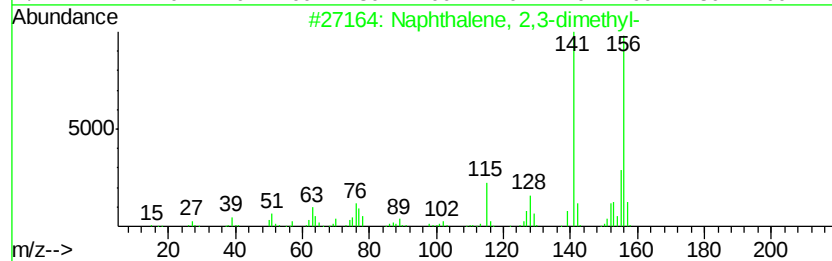
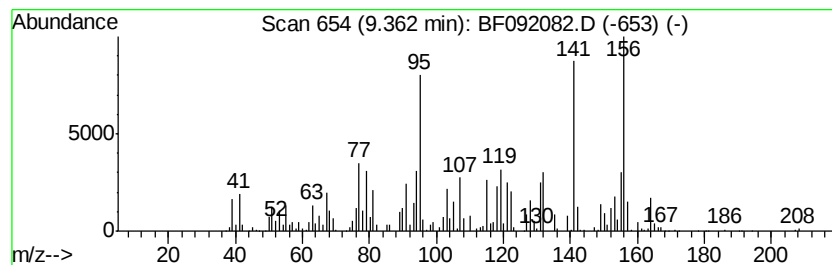
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Naphthalene, 2,3-dimethyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.36	7.80 ng	705359	Acenaphthene-d10	9.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	94
2		Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	89
3		Naphthalene, 1,8-dimethyl-	156	C12H12	000569-41-5	87
4		Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	87
5		Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	86



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF010417\
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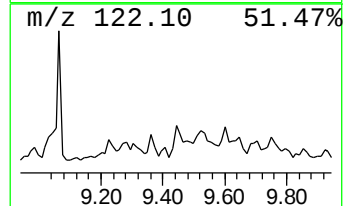
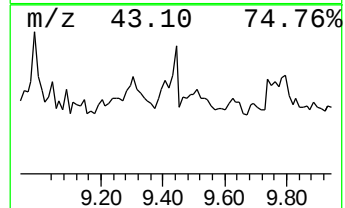
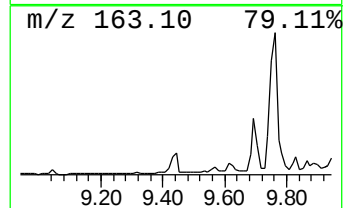
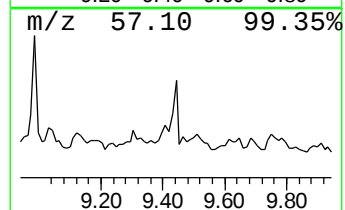
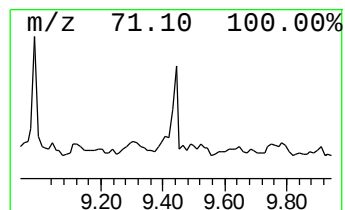
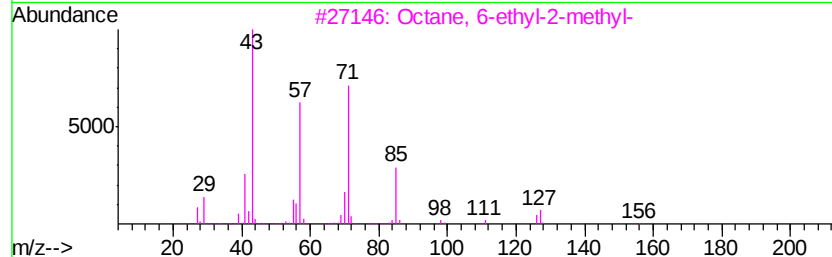
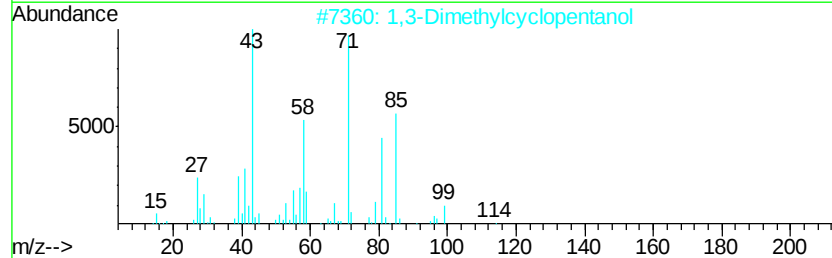
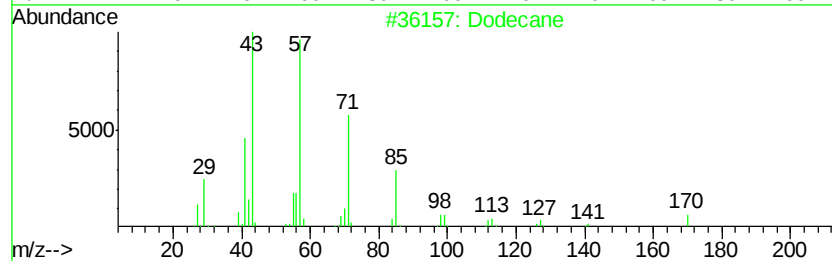
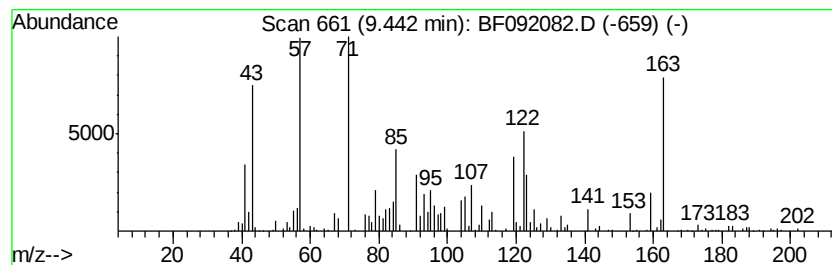
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 unknown9.44 Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.44	7.41 ng	670384	Acenaphthene-d10	9.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dodecane	170	C12H26	000112-40-3	25
2		1,3-Dimethylcyclopentanol	114	C7H14O	019550-46-0	18
3		Octane, 6-ethyl-2-methyl-	156	C11H24	062016-19-7	14
4		Dodecane, 1-iodo-	296	C12H25I	004292-19-7	14
5		Undecane, 4,8-dimethyl-	184	C13H28	017301-33-6	14



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF010417\
Data File : BF092082.D
Acq On : 4 Jan 2017 21:50
Operator : UM/SJ
Sample : I1004-05
Misc :
ALS Vial : 11 Sample Multiplier: 1

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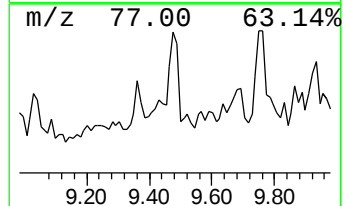
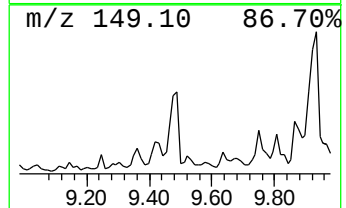
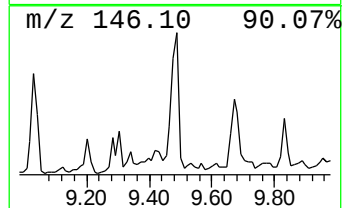
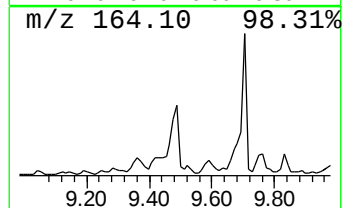
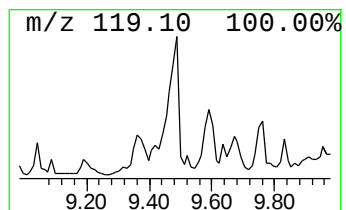
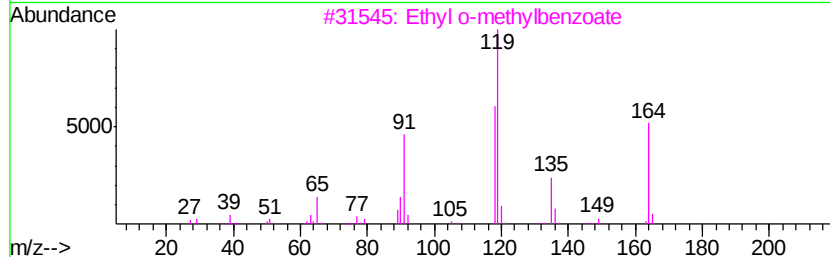
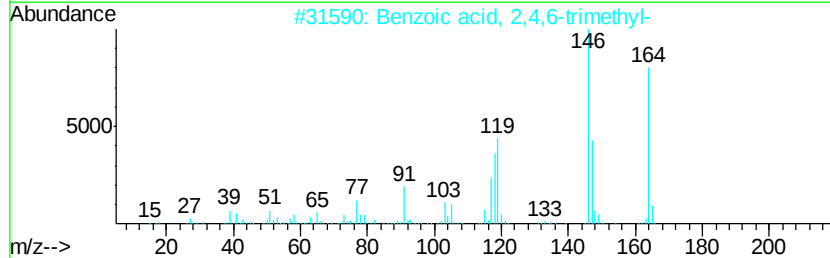
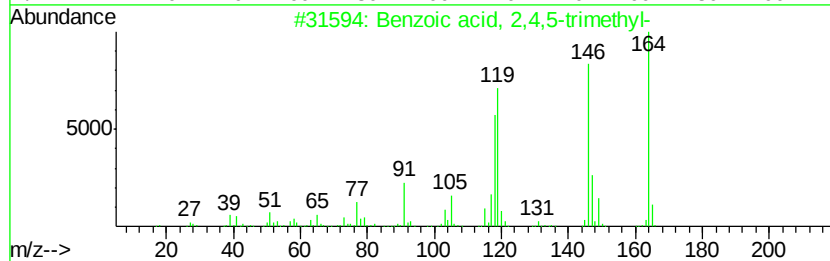
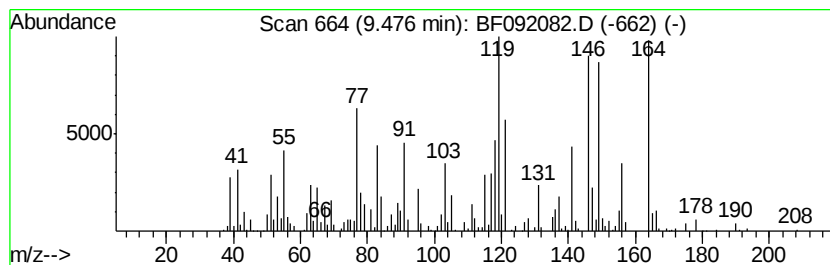
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Benzoic acid, 2,4,5-trimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.48	20.32 ng	1837580	Acenaphthene-d10	9.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzoic acid, 2,4,5-trimethyl-	164	C10H12O2	000528-90-5	50
2		Benzoic acid, 2,4,6-trimethyl-	164	C10H12O2	000480-63-7	46
3		Ethyl o-methylbenzoate	164	C10H12O2	000087-24-1	42
4		3,6-Dimethyl-4H-furo[3,2-c]pyran...	164	C9H8O3	036745-38-7	30
5		Benzofuroxan-6-carboxaldehyde	164	C7H4N2O3	057948-15-9	25



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF010417\
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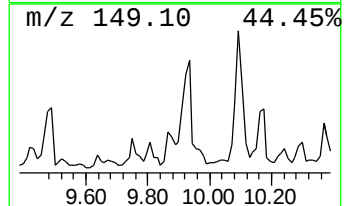
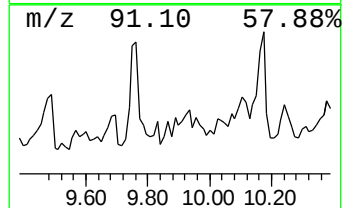
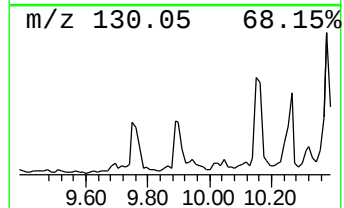
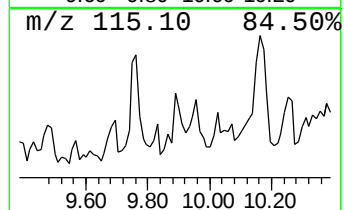
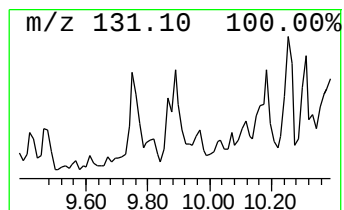
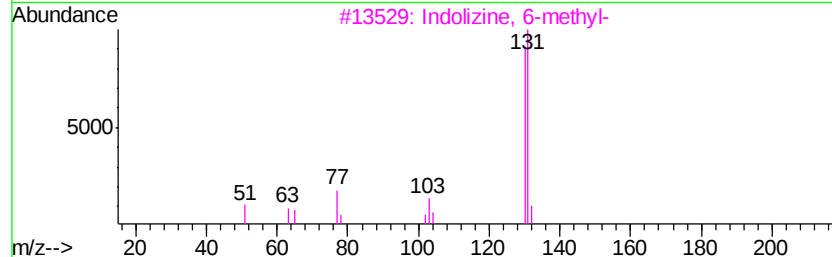
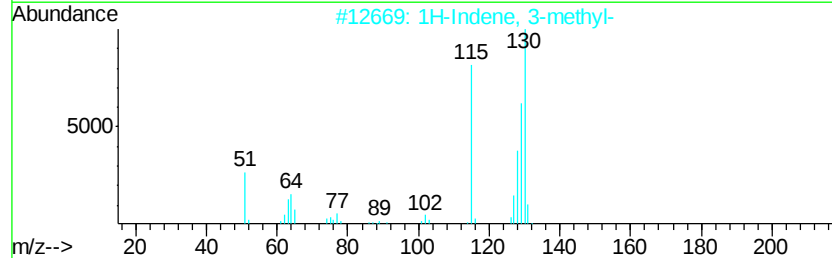
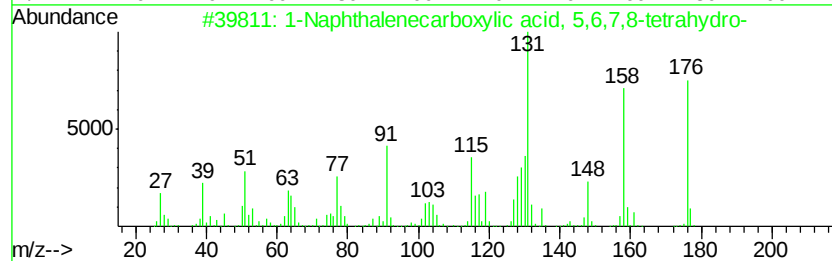
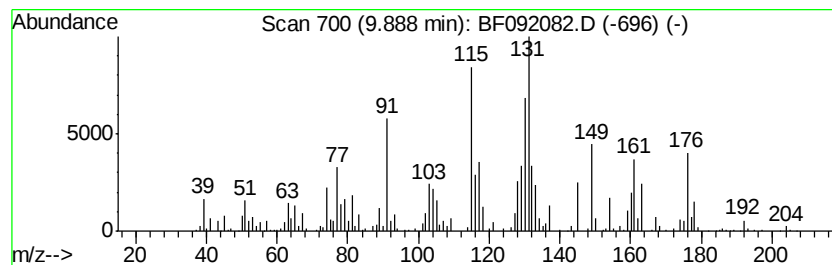
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 unknown9.89 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.89	9.72 ng	878740	Acenaphthene-d10	9.70

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Naphthalenecarboxylic acid, 5,...	176	C11H12O2	004242-18-6	46
2			1H-Indene, 3-methyl-	130	C10H10	000767-60-2	35
3			Indolizine, 6-methyl-	131	C9H9N	001761-11-1	20
4			2-Propenoic acid, 3-phenyl-, eth...	176	C11H12O2	000103-36-6	18
5			Isoquinoline, 3,4-dihydro-	131	C9H9N	003230-65-7	18



Data Path : Z:\HPCHEM1\BNA F\DATA\BF010417\
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Misc :
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Instrument :
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ClientSampleId :
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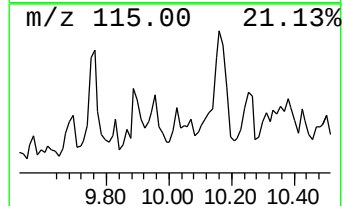
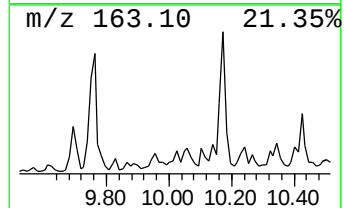
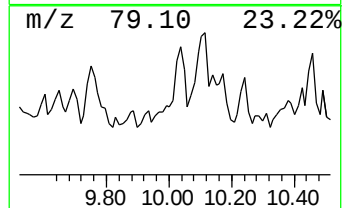
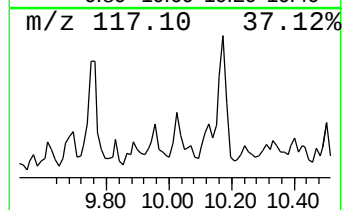
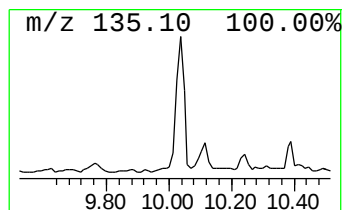
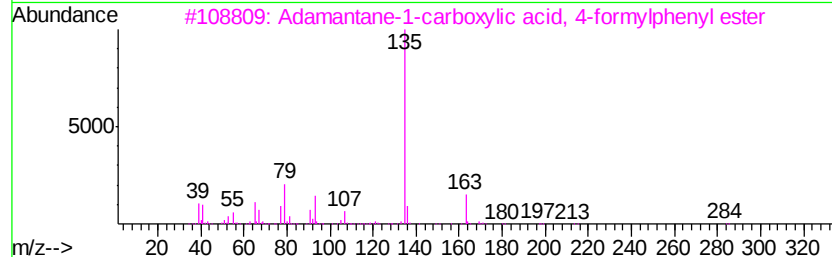
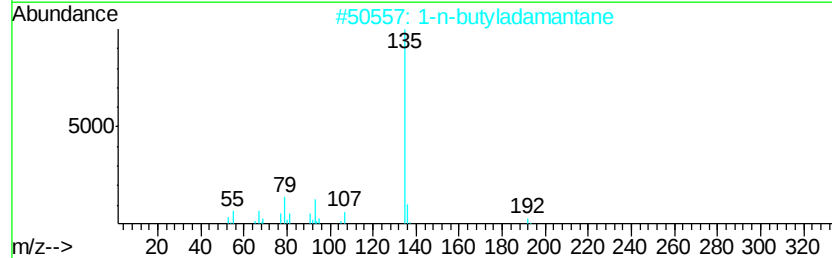
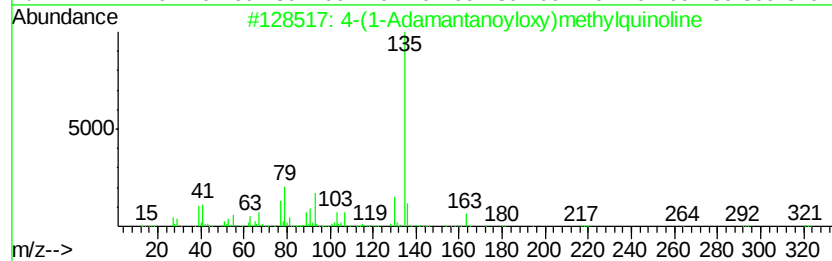
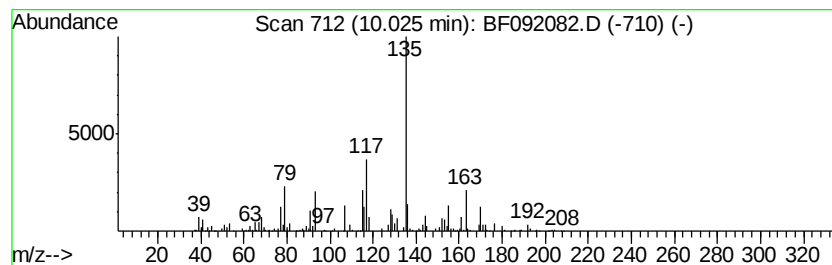
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 unknown10.02 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.02	10.60 ng	958355	Acenaphthene-d10	9.70

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	4		(1-Adamantanovloxy)methylquino...	321	C21H23NO2	1000287-18-6	43
2	1		n-butyladamantane	192	C14H24	014449-41-3	43
3			Adamantane-1-carboxylic acid, 4-...	284	C18H20O3	1000276-61-1	43
4			(2-Oxazolidinylidene)malononitrile	135	C6H5N3O	002733-51-9	38
5			Benzoic acid, 4-methoxy-, diethy...	220	C12H17BO3	146681-61-0	38



Data Path : Z:\HPCHEM1\BNA F\DATA\BF010417\
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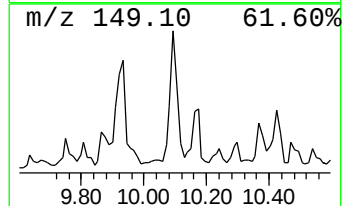
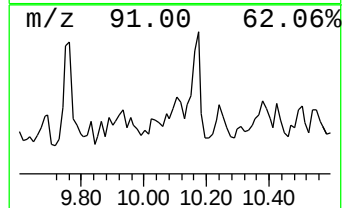
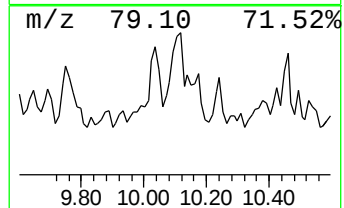
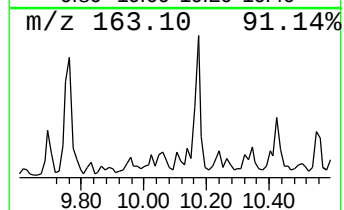
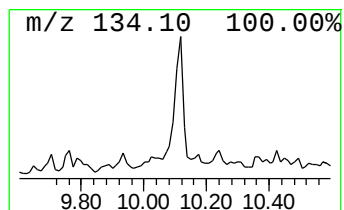
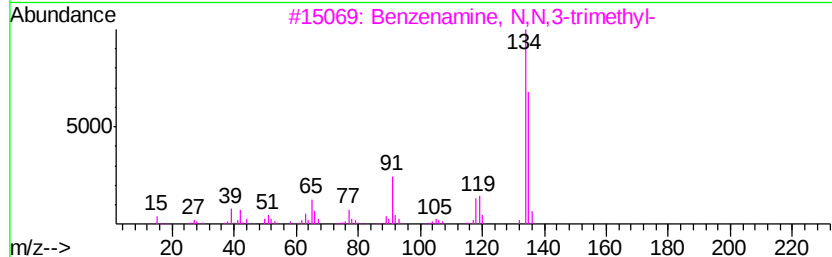
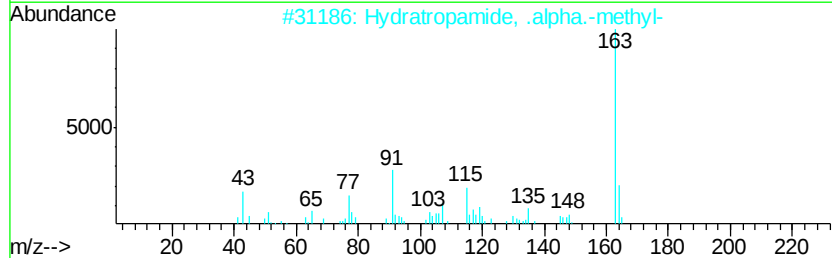
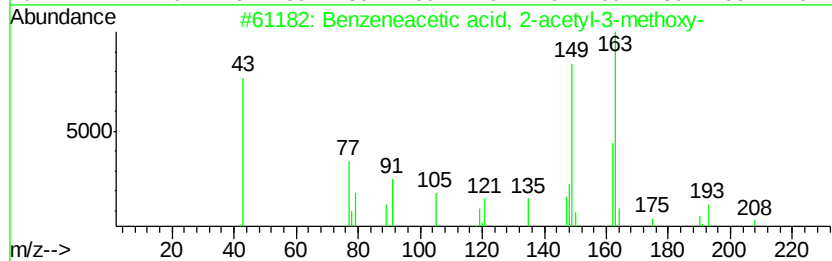
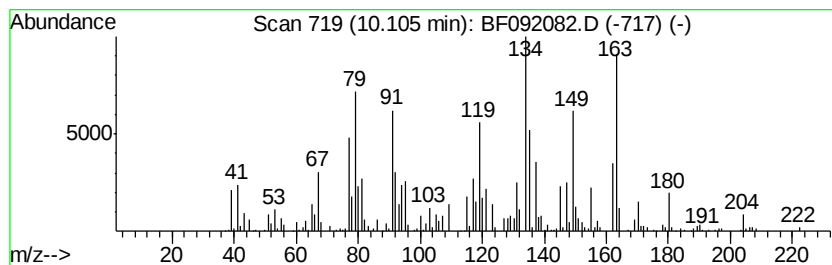
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 unknown10.10 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.10	19.00 ng	1718330	Acenaphthene-d10	9.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzeneacetic acid, 2-acetyl-3-m...	208	C11H12O4	120714-35-4	38
2		Hydratropamide, .alpha.-methyl-	163	C10H13NO	000826-54-0	30
3		Benzenamine, N,N,3-trimethyl-	135	C9H13N	000121-72-2	30
4		o-Toluidine, 5-isopropyl-	149	C10H15N	002051-53-8	27
5		Benzenamine, N,N,4-trimethyl-	135	C9H13N	000099-97-8	25



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF010417\
Data File : BF092082.D
Acq On : 4 Jan 2017 21:50
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Sample : I1004-05
Misc :
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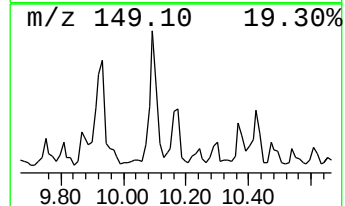
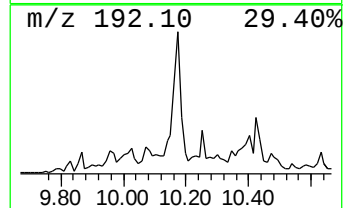
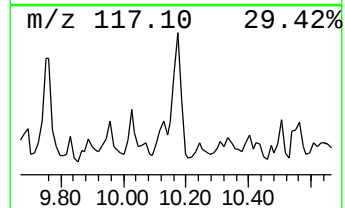
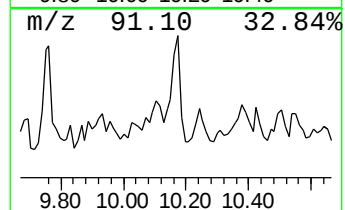
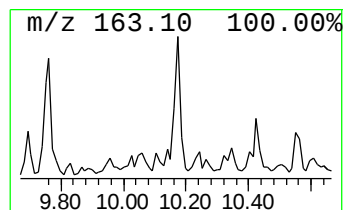
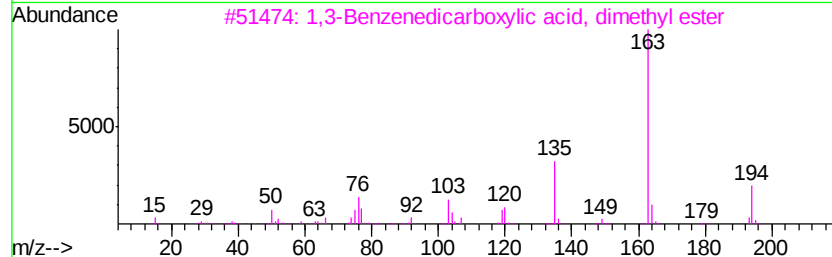
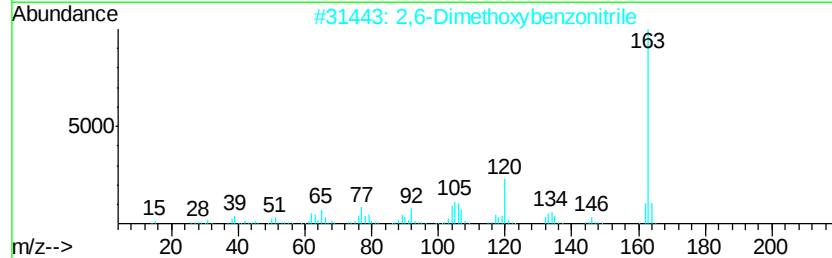
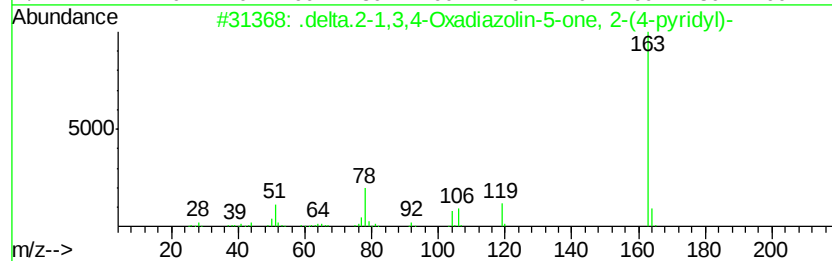
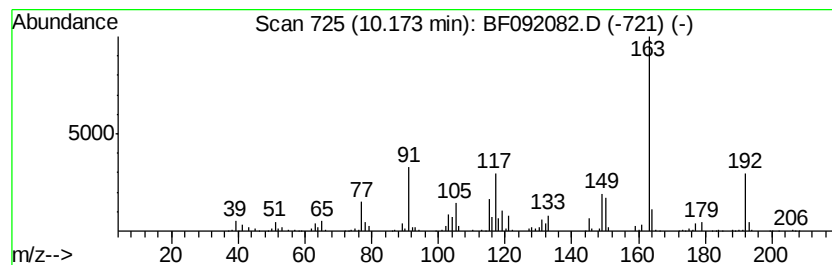
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF122116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 unknown10.17 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.17	45.22 ng	4088600	Acenaphthene-d10	9.70
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		.delta.2-1,3,4-Oxadiazolin-5-one...	163 C7H5N3O2	002845-82-1 43
2		2,6-Dimethoxybenzonitrile	163 C9H9NO2	016932-49-3 43
3		1,3-Benzenedicarboxylic acid, di...	194 C10H10O4	001459-93-4 43
4		Propofol	178 C12H18O	002078-54-8 38
5		1H-Benzimidazole, 5-nitro-	163 C7H5N3O2	000094-52-0 38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF010417\
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Misc :
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ClientSampleId :
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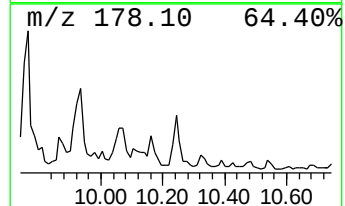
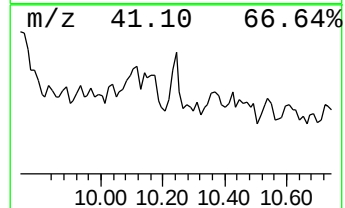
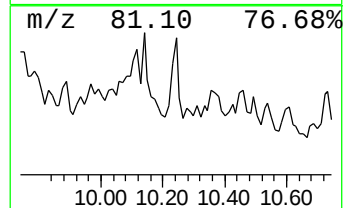
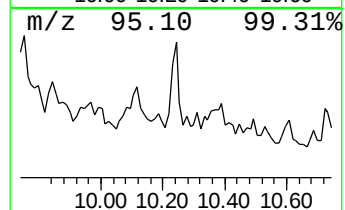
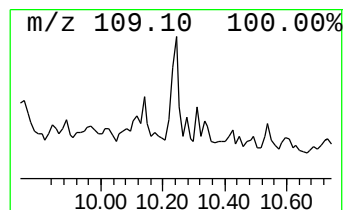
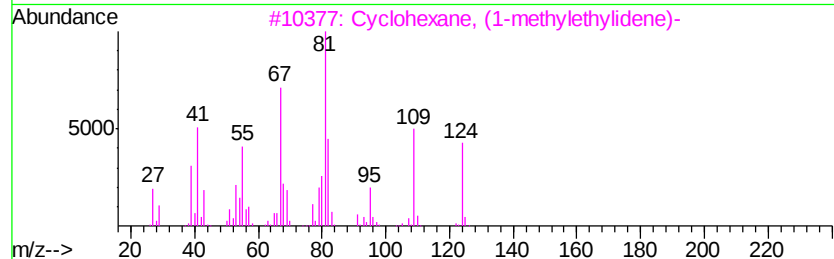
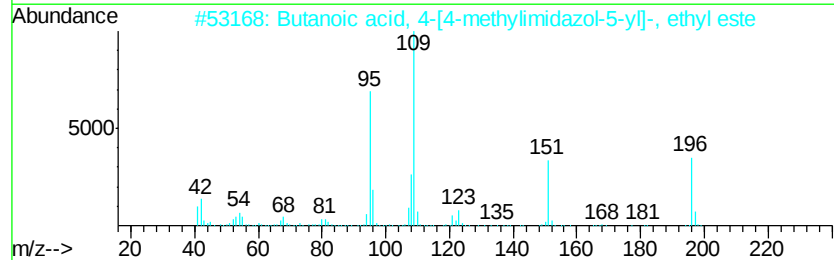
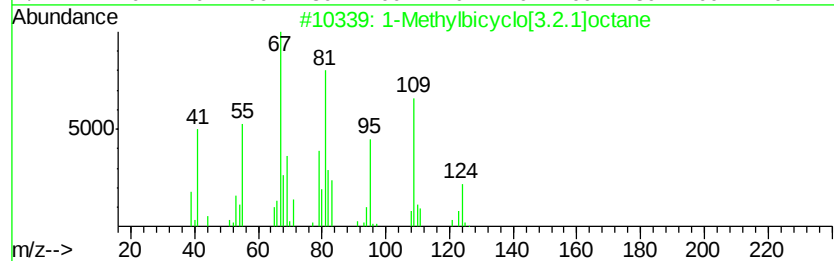
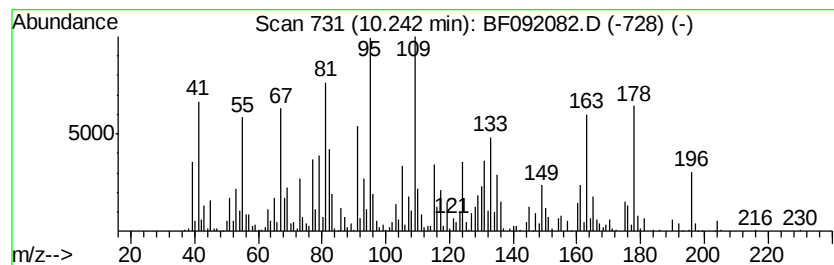
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 unknown10.24 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.24	25.45 ng	2300970	Acenaphthene-d10	9.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Methylbicyclo[3.2.1]octane	124	C9H16	119972-41-7	41
2		Butanoic acid, 4-[4-methylimidaz...	196	C10H16N2O2	1000116-74-5	38
3		Cyclohexane, (1-methylethylidene)-	124	C9H16	005749-72-4	18
4		Ethanone, 1-(2-methyl-1-cyclopent...	124	C8H12O	003168-90-9	15
5		Bicyclo[4.3.0]nonan-2-one, 8-iso...	178	C12H18O	1000159-42-7	15



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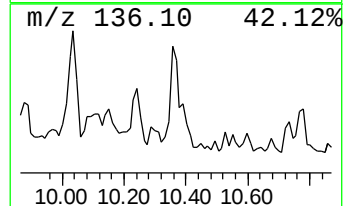
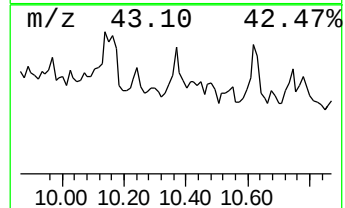
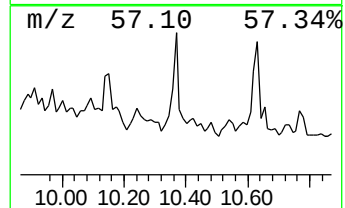
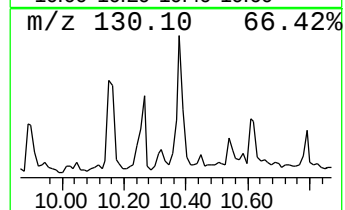
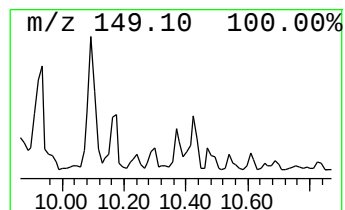
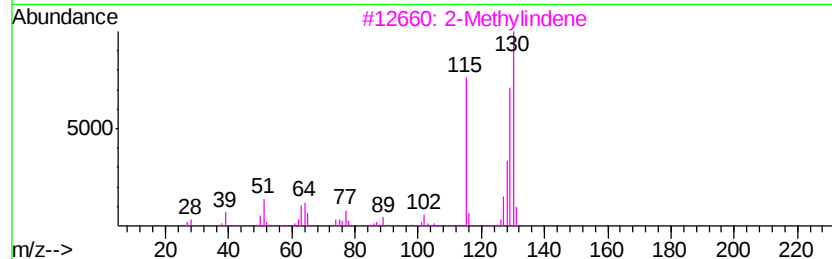
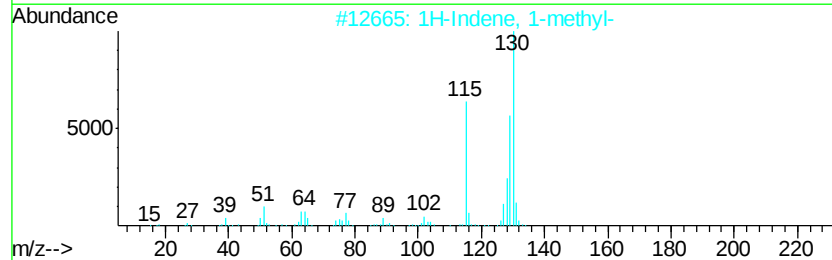
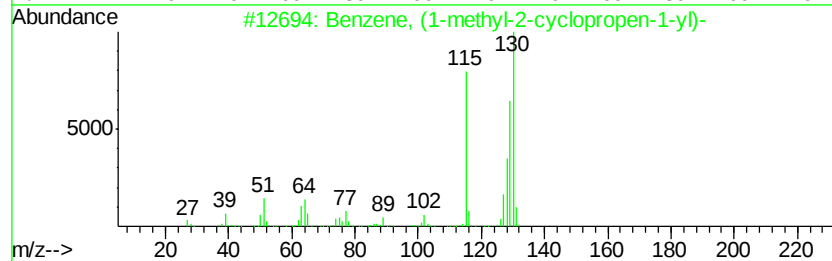
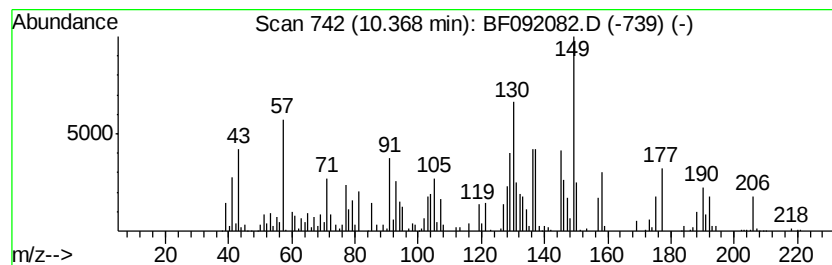
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 unknown10.37 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.37	15.59 ng	1409280	Acenaphthene-d10	9.70

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (1-methyl-2-cyclopropen...	130	C10H10	065051-83-4	15
2			1H-Indene, 1-methyl-	130	C10H10	000767-59-9	15
3			2-Methylindene	130	C10H10	002177-47-1	15
4			Butanoic acid, 2-(aminocarbonyl)...	187	C9H17NO3	007499-15-2	11
5			Benzonitrile, 2,4,6-trimethyl-	145	C10H11N	002571-52-0	11



Data Path : Z:\HPCHEM1\BNA F\DATA\BF010417\
Data File : BF092082.D
Acq On : 4 Jan 2017 21:50
Operator : UM/SJ
Sample : I1004-05
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-36

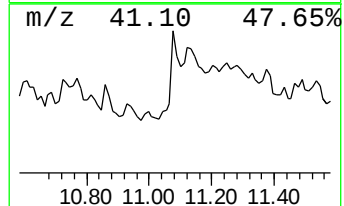
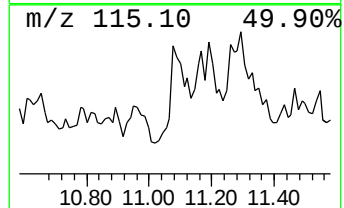
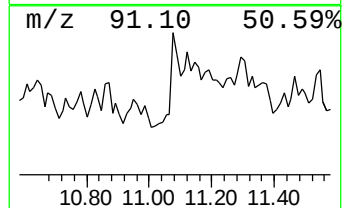
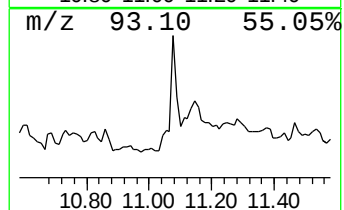
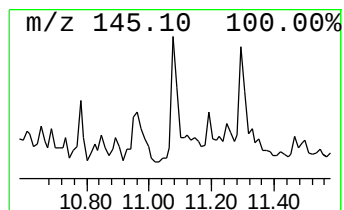
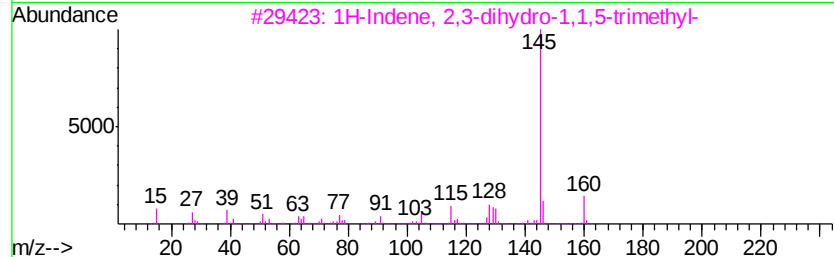
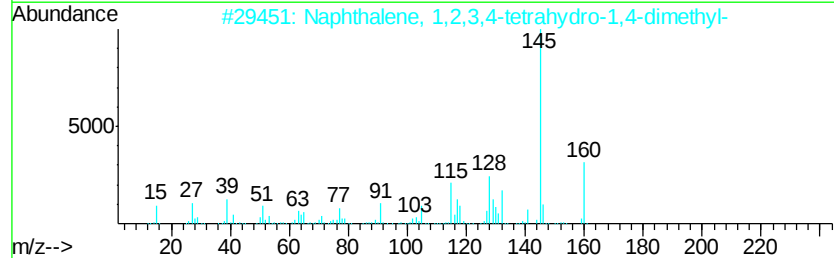
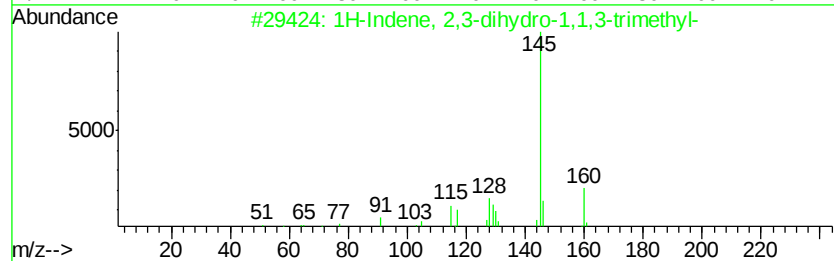
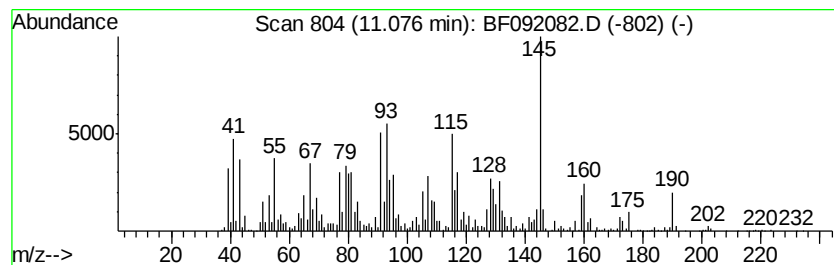
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 1H-Indene, 2,3-dihydro-1,1,... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.08	22.78 ng	3114030	Phenanthrene-d10	11.19

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-1,1,3-tri...	160	C12H16	002613-76-5	93
2		Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	004175-54-6	70
3		1H-Indene, 2,3-dihydro-1,1,5-tri...	160	C12H16	040650-41-7	70
4		Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	025419-33-4	62
5		Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	021564-91-0	60



Data Path : Z:\HPCHEM1\BNA F\DATA\BF010417\
Data File : BF092082.D
Acq On : 4 Jan 2017 21:50
Operator : UM/SJ
Sample : I1004-05
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-36

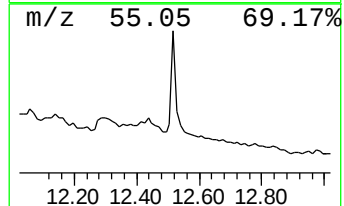
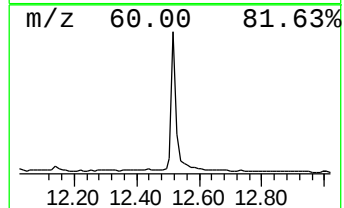
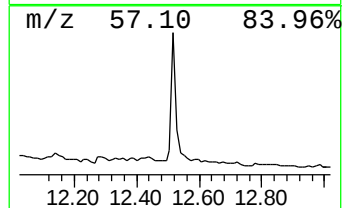
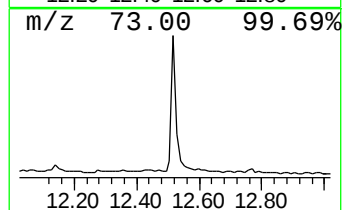
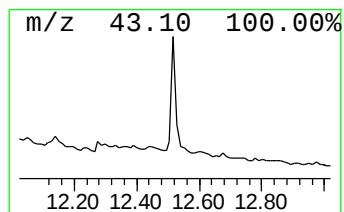
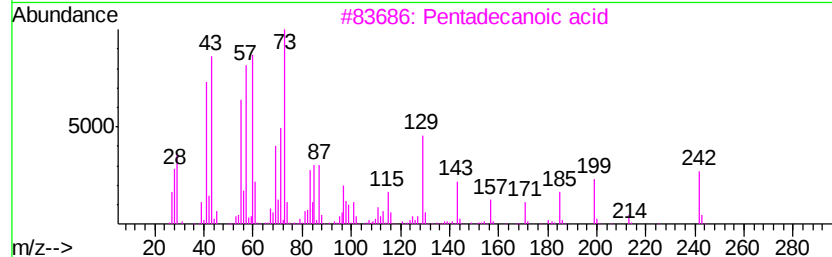
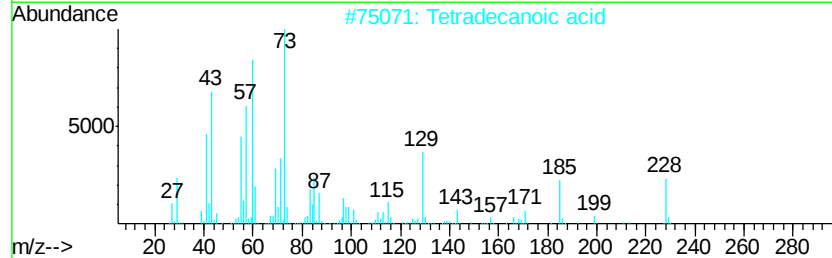
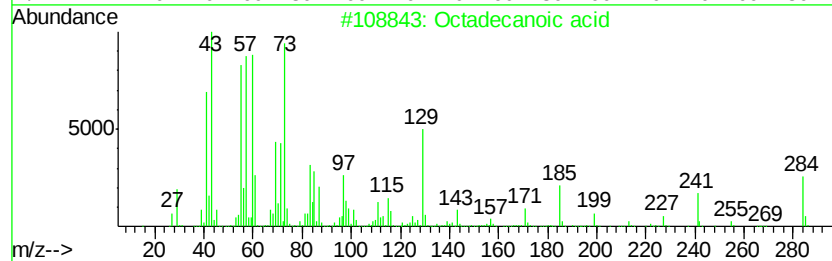
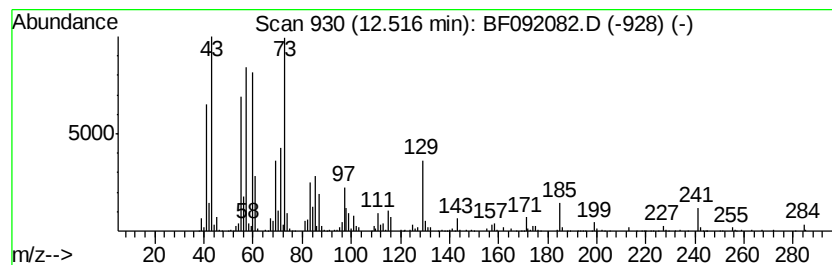
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF122116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 Octadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.52	50.72 ng	1332400	Chrysene-d12	13.81

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2		Tetradecanoic acid	228	C14H28O2	000544-63-8	94
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	91
4		Octadecanoic acid, 2-(2-hydroxyve...	372	C22H44O4	000106-11-6	72
5		Tridecanoic acid	214	C13H26O2	000638-53-9	43



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF010417\
 Data File : BF092082.D
 Acq On : 4 Jan 2017 21:50
 Operator : UM/SJ
 Sample : I1004-05
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-36

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF122116.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
2-Pentanone, 4-hy...	4.80	9.5 ng		851340	1	6.66	1788210	20.0
4-Octene, (E)-	6.17	7.1 ng		632207	1	6.66	1788210	20.0
unknown6.44	6.44	55.6 ng		4970080	1	6.66	1788210	20.0
Benzene, 1,2,4,5-...	7.44	7.1 ng		875230	2	7.96	2467050	20.0
Benzene, 1-methyl...	7.69	12.3 ng		1510600	2	7.96	2467050	20.0
unknown8.47	8.47	7.3 ng		906183	2	7.96	2467050	20.0
Naphthalene, 1,2,...	8.61	10.7 ng		1314790	2	7.96	2467050	20.0
Naphthalene, 1-me...	8.77	31.4 ng		3872840	2	7.96	2467050	20.0
Naphthalene, 2,3-...	9.36	7.8 ng		705359	3	9.70	1808470	20.0
unknown9.44	9.44	7.4 ng		670384	3	9.70	1808470	20.0
Benzoic acid, 2,4...	9.48	20.3 ng		1837580	3	9.70	1808470	20.0
unknown9.89	9.89	9.7 ng		878740	3	9.70	1808470	20.0
unknown10.02	10.02	10.6 ng		958355	3	9.70	1808470	20.0
unknown10.10	10.10	19.0 ng		1718330	3	9.70	1808470	20.0
unknown10.17	10.17	45.2 ng		4088600	3	9.70	1808470	20.0
unknown10.24	10.24	25.4 ng		2300970	3	9.70	1808470	20.0
unknown10.37	10.37	15.6 ng		1409280	3	9.70	1808470	20.0
1H-Indene, 2,3-di...	11.08	22.8 ng		3114030	4	11.19	2733690	20.0
Octadecanoic acid	12.52	50.7 ng		1332400	5	13.81	525367	20.0