

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA M\DATA\BM021318\
 Data File : BM014225.D
 Acq On : 13 Feb 2018 17:33
 Operator : SJ/JU
 Sample : J1518-08
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MW-5A

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_M\METHODS\8270-BM020718.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	3.810	134	140	152	rBV	256743	410427	3.38%	0.781%
2	5.157	361	369	375	rBV	899995	1410818	11.61%	2.683%
3	5.210	375	378	395	rVB	398870	918045	7.56%	1.746%
4	5.575	434	440	446	rBV	298513	440257	3.62%	0.837%
5	6.728	632	636	641	rBV	497195	794613	6.54%	1.511%
6	6.928	664	670	683	rVB	117101	205290	1.69%	0.390%
7	7.075	688	695	703	rBV	2200834	3330247	27.41%	6.334%
8	7.181	708	713	721	rVB	234105	366194	3.01%	0.696%
9	7.539	766	774	782	rBV	430465	741656	6.10%	1.411%
10	7.645	787	792	800	rVB	197147	336036	2.77%	0.639%
11	7.851	820	827	834	rBV	2038088	3172411	26.11%	6.034%
12	8.169	875	881	884	rBV	75883	128901	1.06%	0.245%
13	8.698	963	971	984	rBV	1709639	2859581	23.53%	5.439%
14	8.875	988	1001	1010	rBV3	211950	534031	4.39%	1.016%
15	9.716	1137	1144	1151	rVB7	69724	145320	1.20%	0.276%
16	10.310	1233	1245	1250	rBV	621858	1101575	9.07%	2.095%
17	10.357	1250	1253	1259	rVV2	121813	219842	1.81%	0.418%
18	11.563	1453	1458	1464	rBV4	81827	195648	1.61%	0.372%
19	12.051	1536	1541	1550	rVB7	62612	133262	1.10%	0.253%
20	12.804	1662	1669	1676	rBV	4670601	6699040	55.13%	12.741%
21	13.574	1794	1800	1807	rVB3	69340	128525	1.06%	0.244%
22	13.668	1810	1816	1821	rBV2	100670	180332	1.48%	0.343%
23	14.098	1884	1889	1893	rBV2	102663	137730	1.13%	0.262%
24	14.192	1898	1905	1914	rBV2	1005762	1596534	13.14%	3.036%
25	14.692	1986	1990	2004	rVB5	69014	157163	1.29%	0.299%
26	15.057	2047	2052	2056	rBV	111092	148357	1.22%	0.282%
27	15.698	2154	2161	2175	rBV	4459856	6315115	51.97%	12.011%
28	16.951	2368	2374	2387	rBV	1419009	2063853	16.98%	3.925%
29	17.862	2523	2529	2538	rBV	423272	649699	5.35%	1.236%
30	19.150	2743	2748	2758	rBV	217062	354562	2.92%	0.674%
31	19.598	2818	2824	2835	rBV	9959765	12151419	100.00%	23.110%
32	20.880	3037	3042	3050	rBV	519286	696401	5.73%	1.324%
33	21.156	3084	3089	3096	rBV	1575870	2226978	18.33%	4.235%
34	23.333	3453	3459	3469	rVB	856260	1630069	13.41%	3.100%

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Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\HPCHEM1\BNA_M\METHODS\8270-BM020718.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

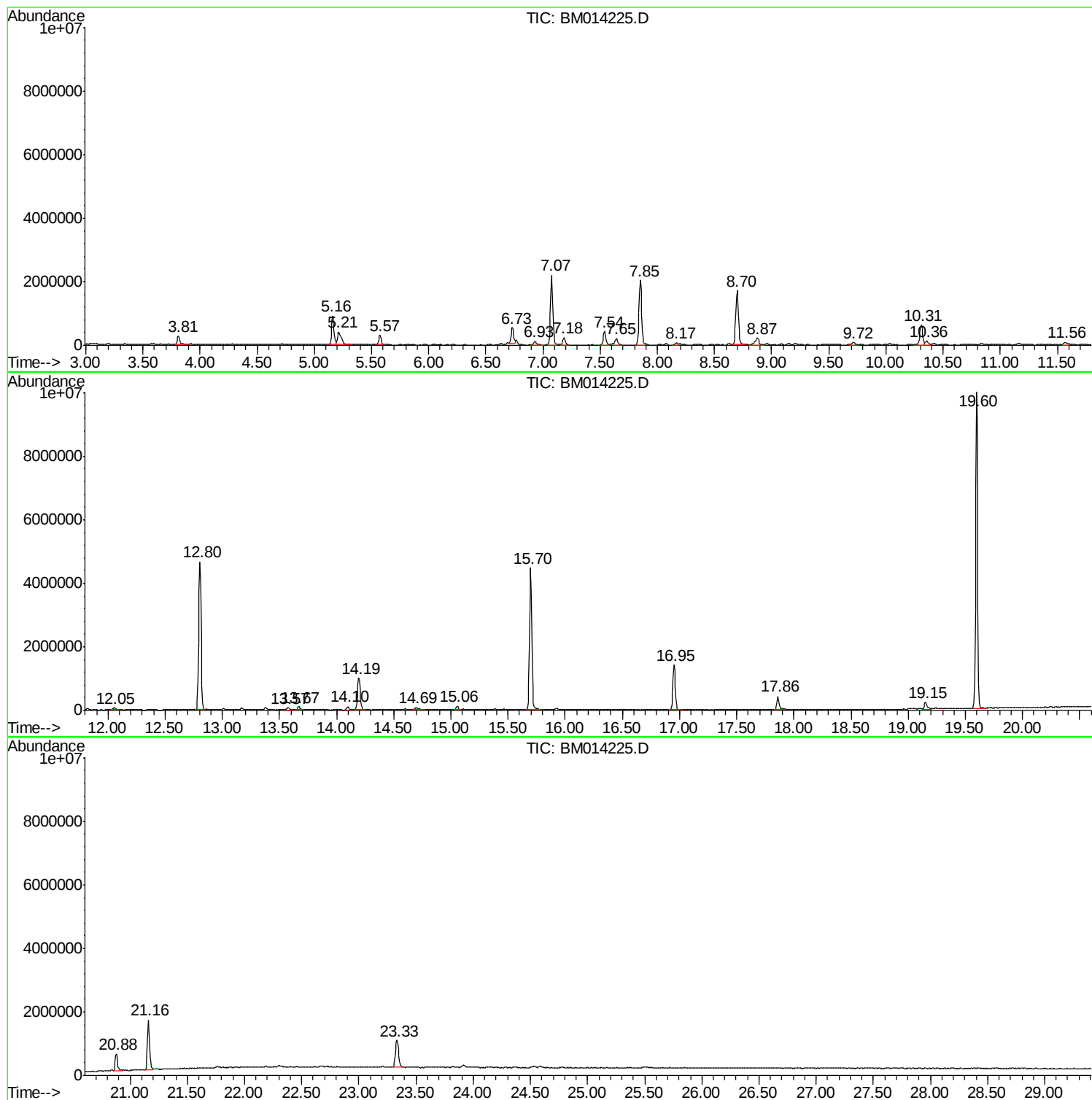
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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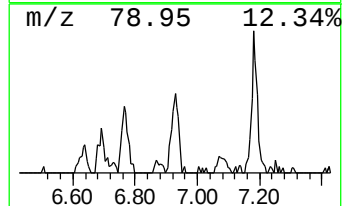
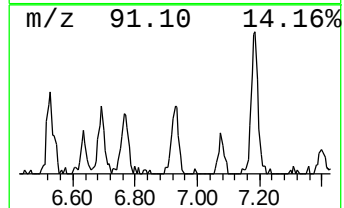
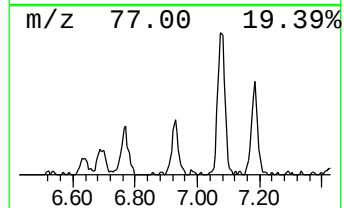
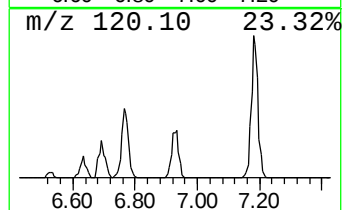
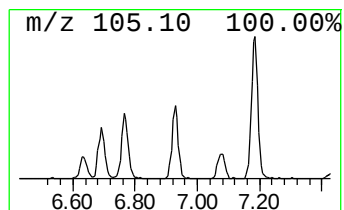
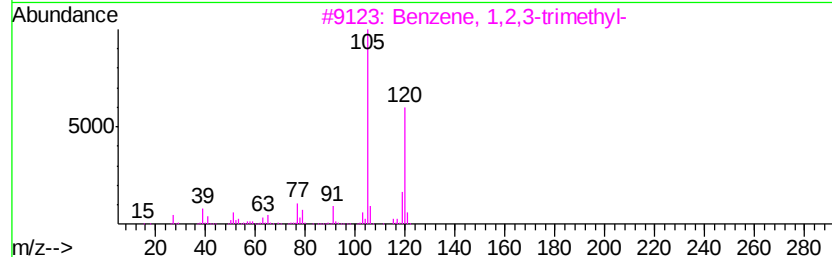
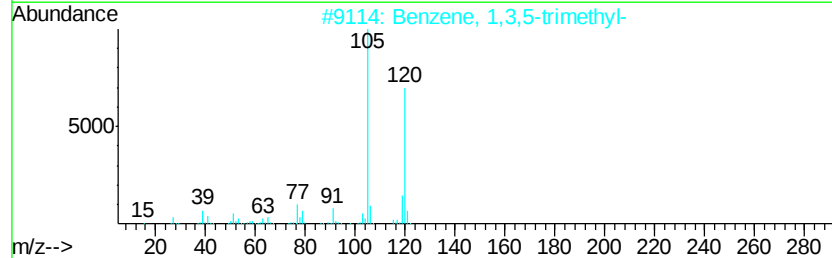
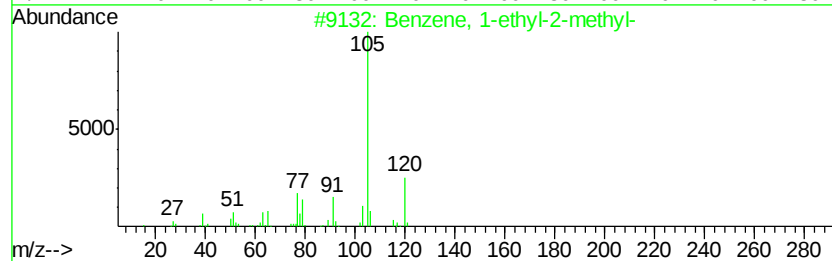
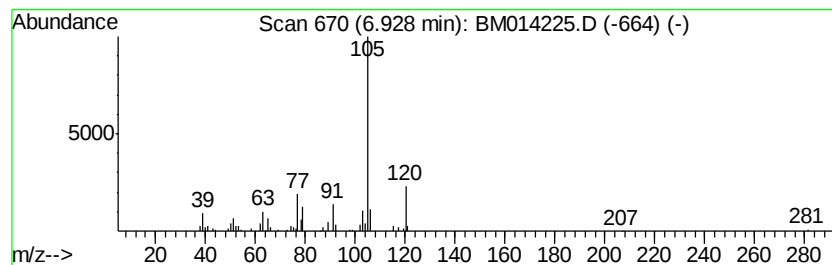
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Benzene, 1-ethyl-2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.93	5.54 ng	205290	1,4-Dichlorobenzene-d4	7.54

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	87
2		Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	87
3		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	87
4		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	86
5		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	83



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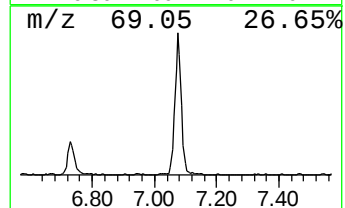
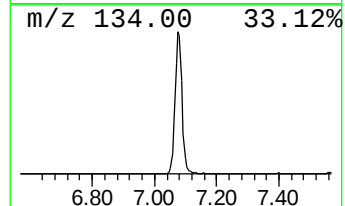
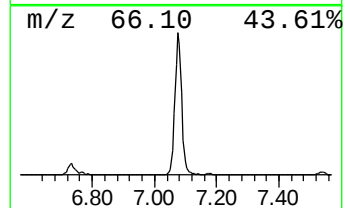
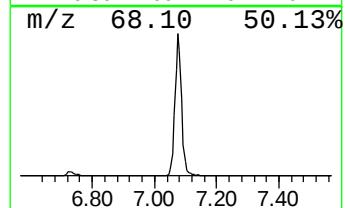
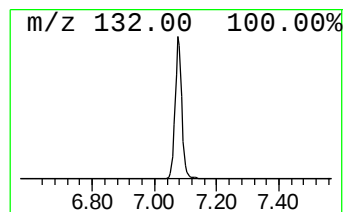
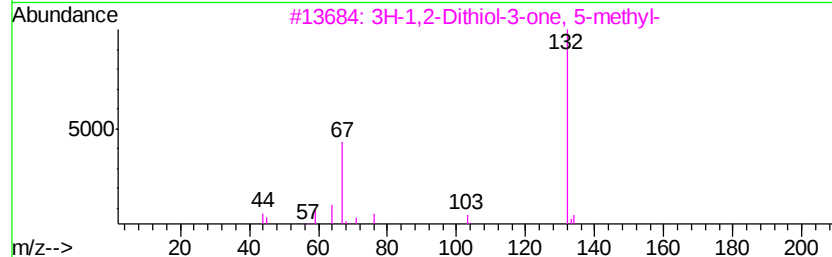
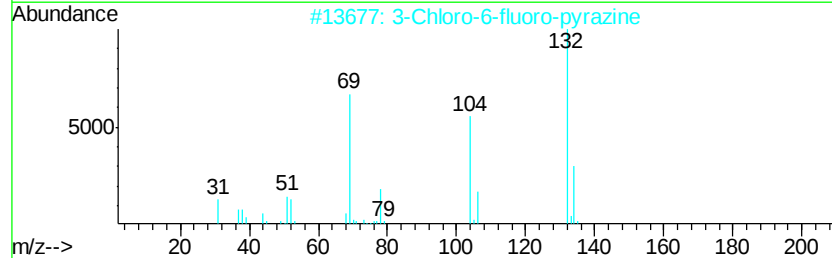
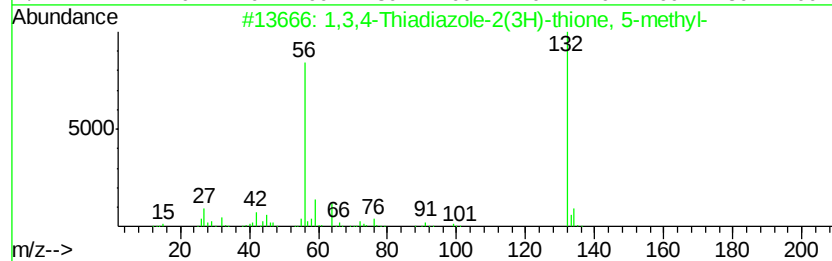
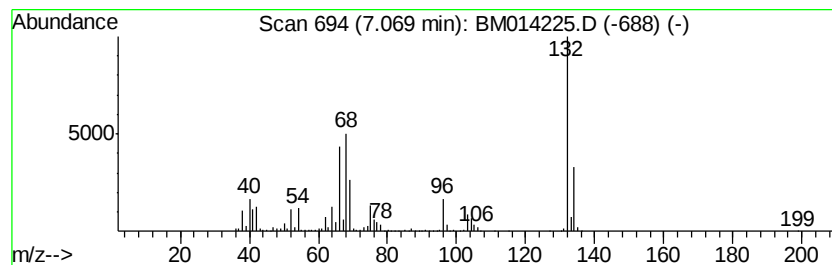
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 unknown7.07 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.07	89.81 ng	3330250	1,4-Dichlorobenzene-d4	7.54

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,3,4-Thiadiazole-2(3H)-thione, ...	132	C3H4N2S2	029490-19-5	14
2			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	14
3			3H-1,2-Dithiol-3-one, 5-methyl-	132	C4H4OS2	003620-08-4	14
4			1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	12
5			1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	9



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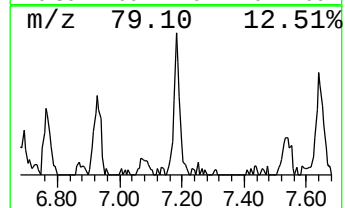
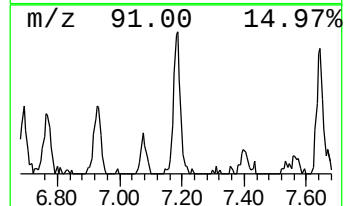
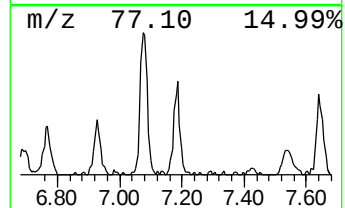
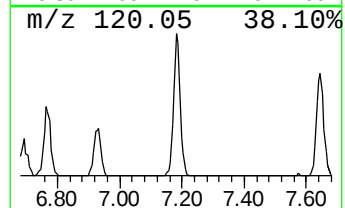
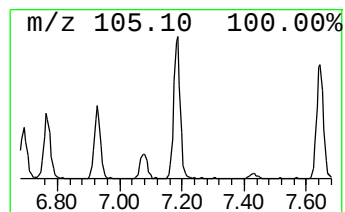
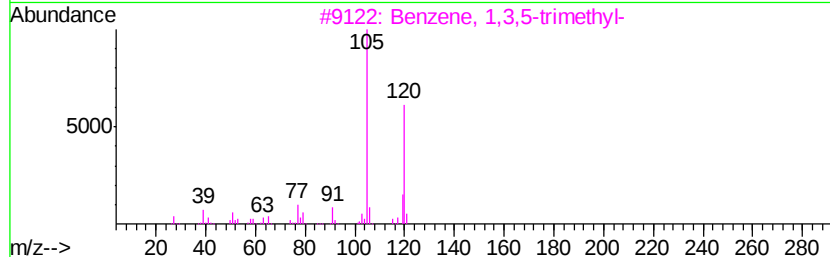
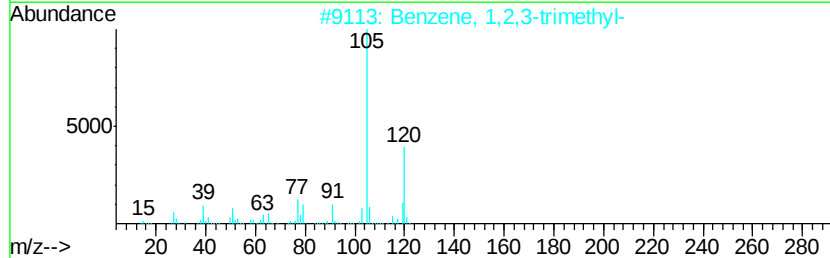
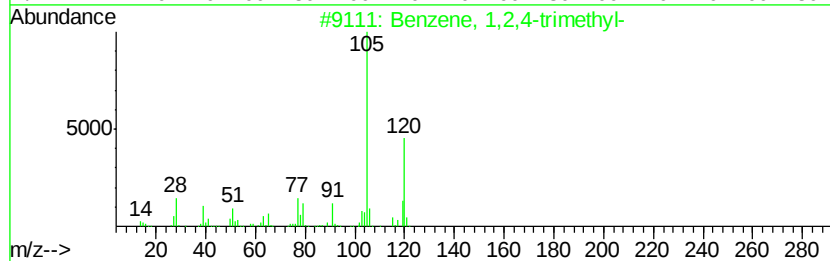
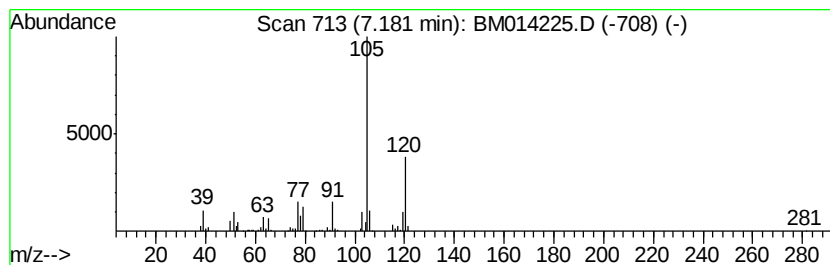
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Benzene, 1,2,4-trimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.18	9.88 ng	366194	1,4-Dichlorobenzene-d4	7.54

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	96
2		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
3		Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	93
4		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	87



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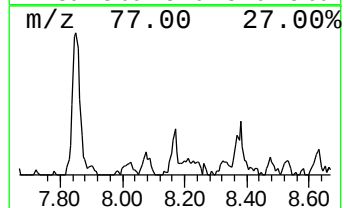
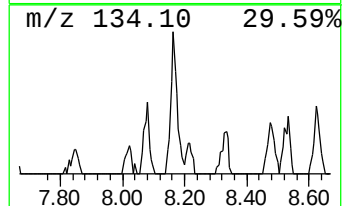
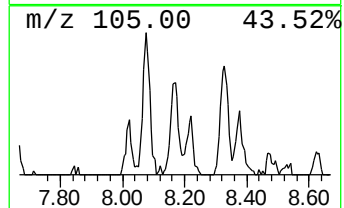
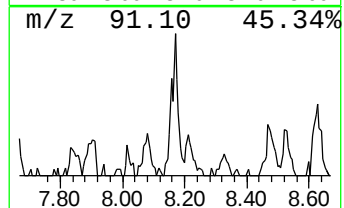
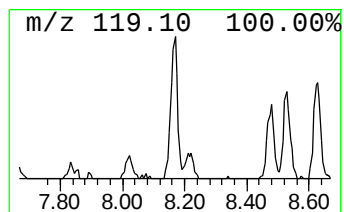
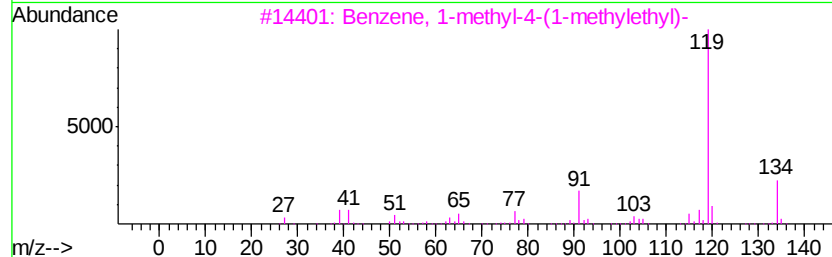
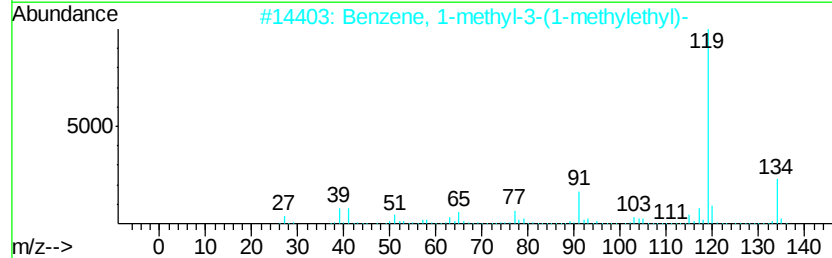
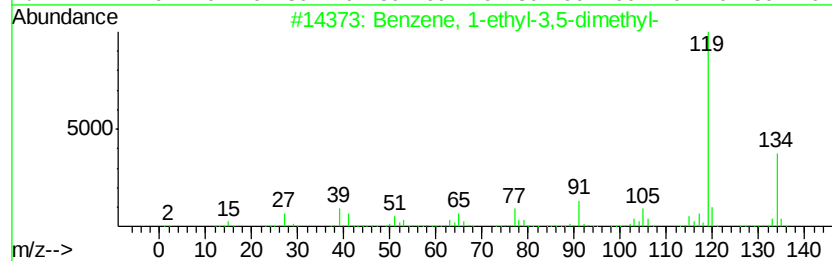
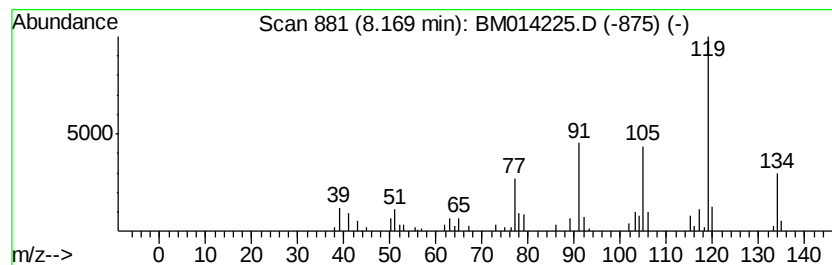
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Benzene, 1-ethyl-3,5-dimethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.17	3.48 ng	128901	1,4-Dichlorobenzene-d4	7.54

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	94
2		Benzene, 1-methyl-3-(1-methylethyl-)	134	C10H14	000535-77-3	94
3		Benzene, 1-methyl-4-(1-methylethyl-)	134	C10H14	000099-87-6	93
4		Benzene, 1-methyl-2-(1-methylethyl-)	134	C10H14	000527-84-4	93
5		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	93



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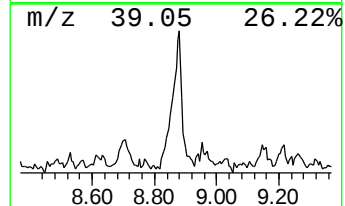
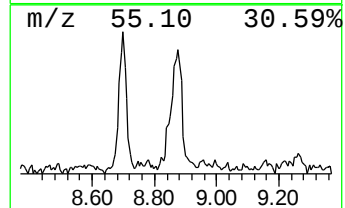
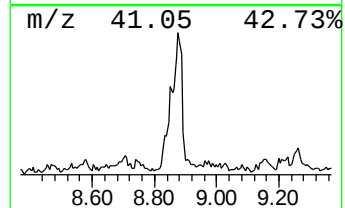
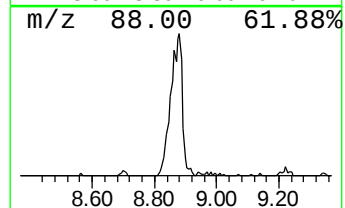
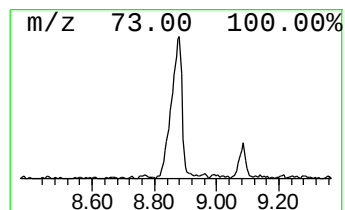
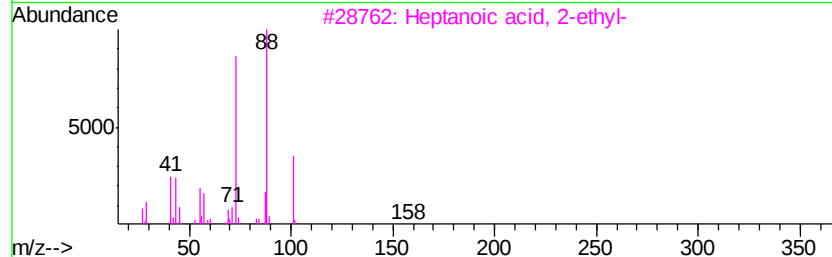
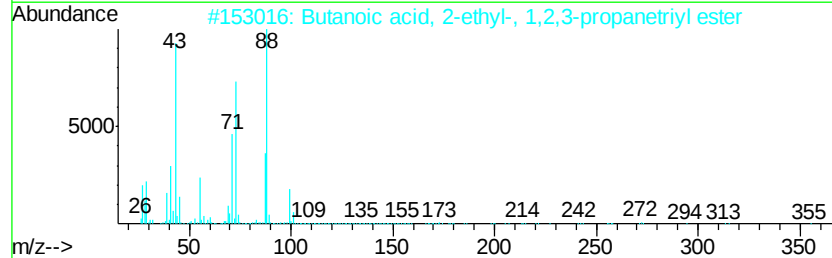
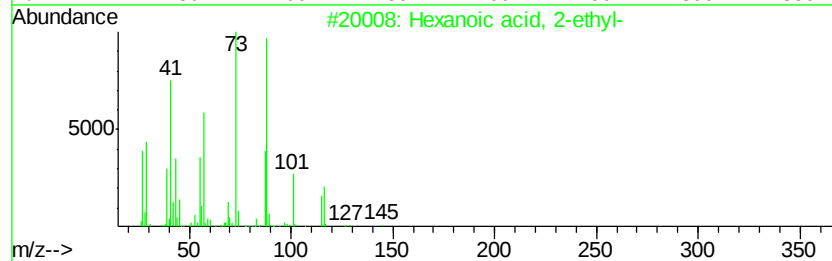
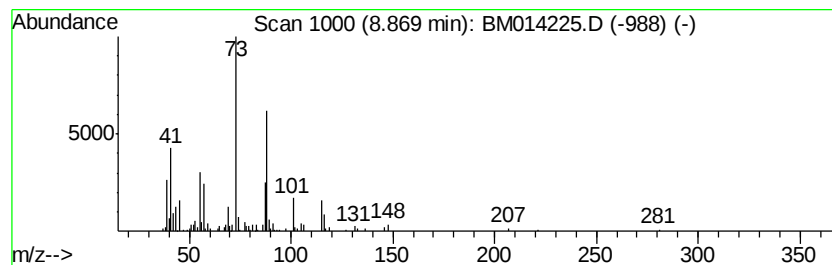
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Peak Number 9 Hexanoic acid, 2-ethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.87	14.40 ng	534031	1,4-Dichlorobenzene-d4	7.54

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	80
2		Butanoic acid, 2-ethyl-, 1,2,3-p...	386	C21H38O6	056554-54-2	50
3		Heptanoic acid, 2-ethyl-	158	C9H18O2	003274-29-1	47
4		3-Aza-2-oxabicyclo[2.2.2]oct-5-e...	267	C10H12F3N04	102698-35-1	38
5		Pentanoic acid, 2,4-dimethyl-, m...	144	C8H16O2	071672-33-8	37



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Acq On : 13 Feb 2018 17:33
Operator : SJ/JU
Sample : J1518-08
Misc :
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Instrument :
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ClientSampleId :
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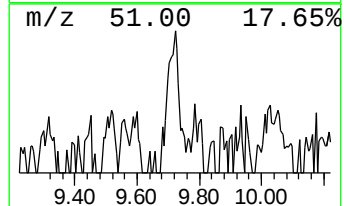
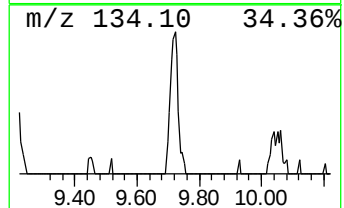
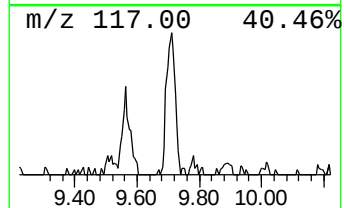
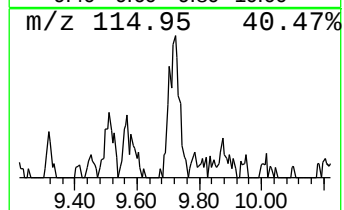
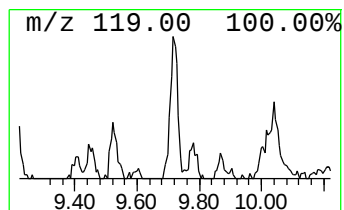
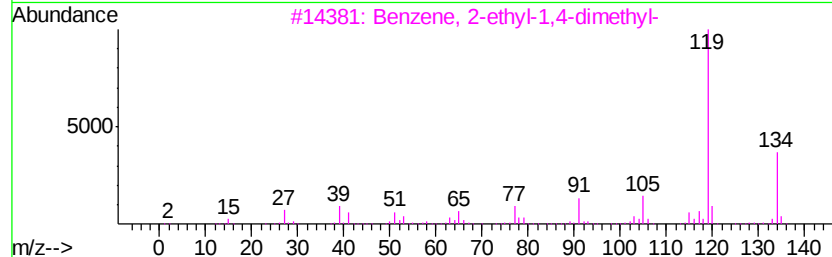
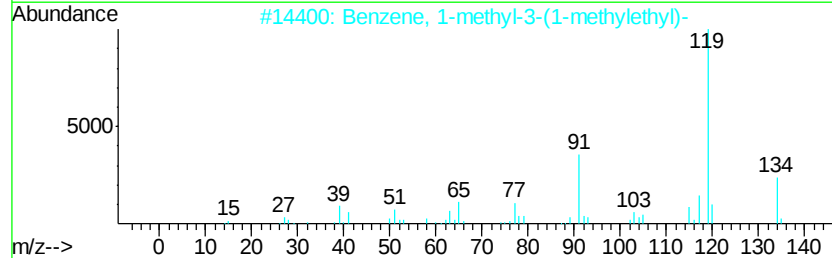
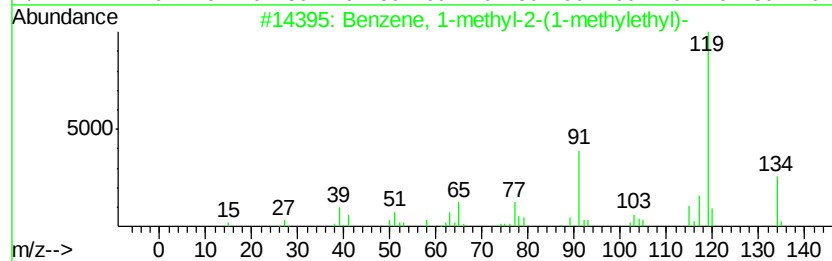
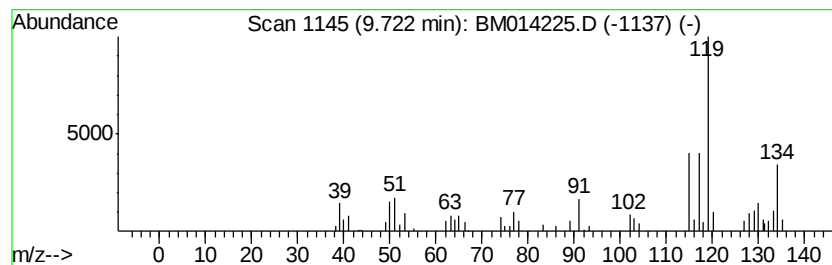
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 unknown9.72 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.72	2.64 ng	145320	Naphthalene-d8	10.31

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-2-(1-methylethyl)-	134	C10H14	000527-84-4	49
2		Benzene, 1-methyl-3-(1-methylethyl)-	134	C10H14	000535-77-3	46
3		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	43
4		Benzene, 1-methyl-4-(1-methylethyl)-	134	C10H14	000099-87-6	43
5		Ethanone, 1-(4-methylphenyl)-	134	C9H10O	000122-00-9	35



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA M\DATA\BM021318\
Data File : BM014225.D
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Misc :
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Instrument :
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ClientSampleId :
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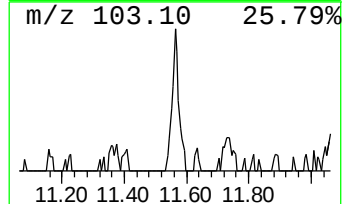
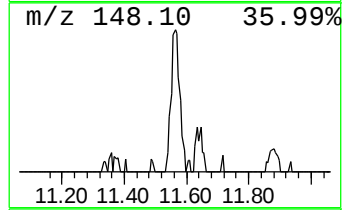
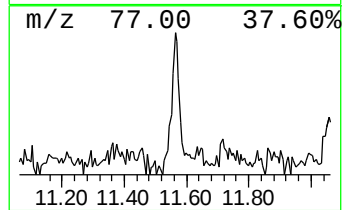
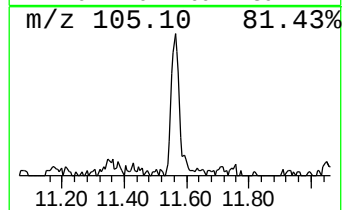
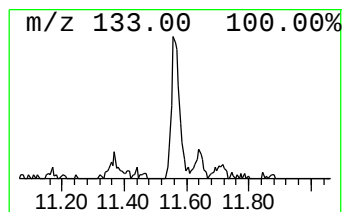
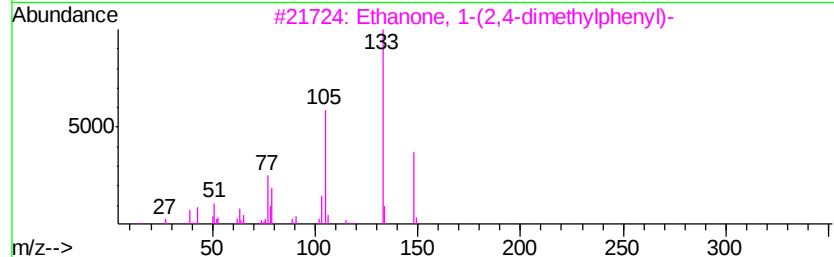
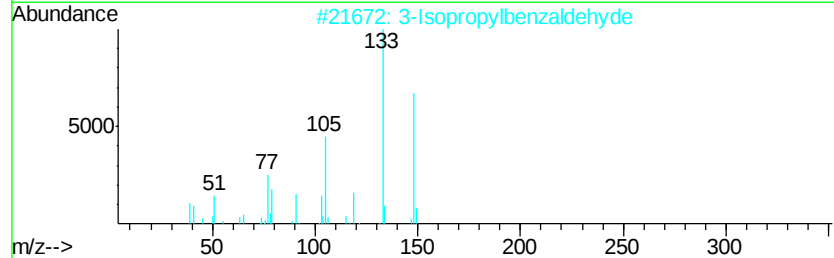
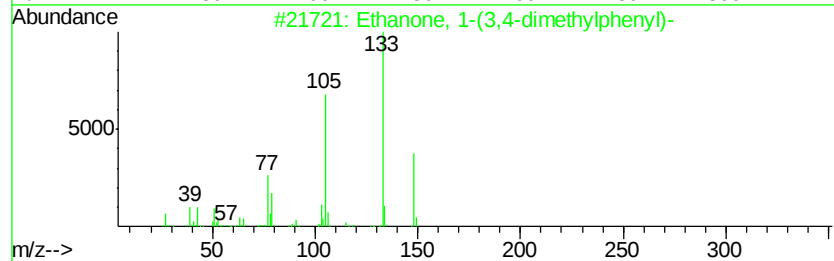
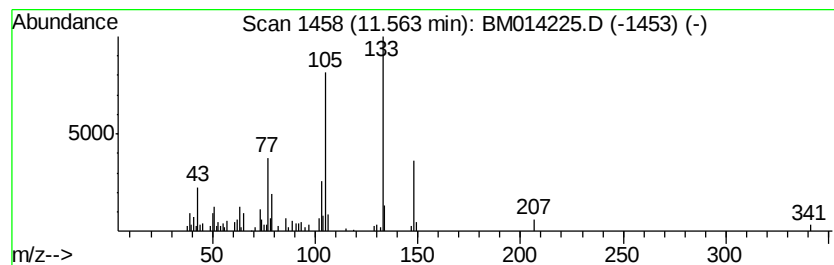
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Ethanone, 1-(3,4-dimethylph... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.56	3.55 ng	195648	Napthalene-d8	10.31

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethanone, 1-(3,4-dimethylphenyl)-	148	C10H12O	003637-01-2	93
2		3-Isopropylbenzaldehyde	148	C10H12O	034246-57-6	90
3		Ethanone, 1-(2,4-dimethylphenyl)-	148	C10H12O	000089-74-7	90
4		Benzene, 1-ethyl-4-(1-methylethyl)-	148	C11H16	004218-48-8	87
5		1-(2,3-Dimethylphenyl)ethanone	148	C10H12O	002142-71-4	80



Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA M\DATA\BM021318\
Data File : BM014225.D
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Operator : SJ/JU
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Misc :
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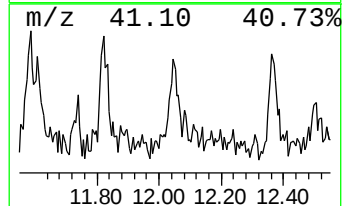
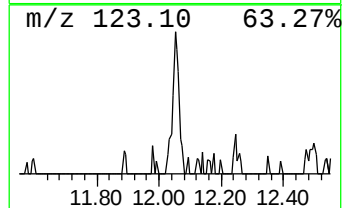
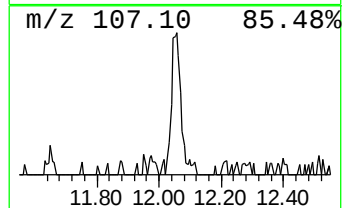
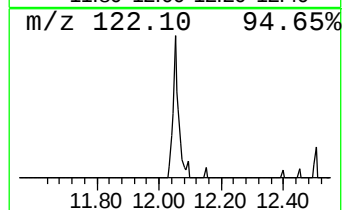
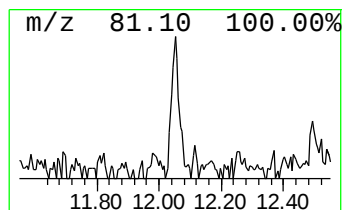
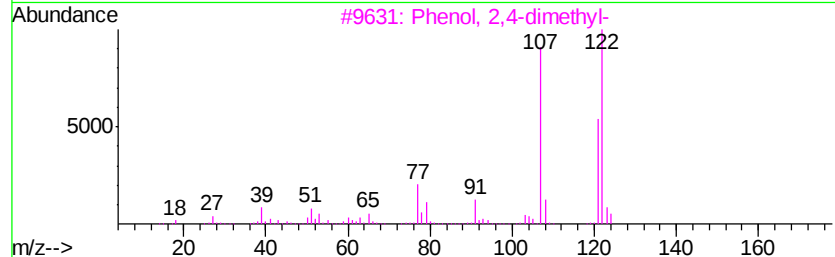
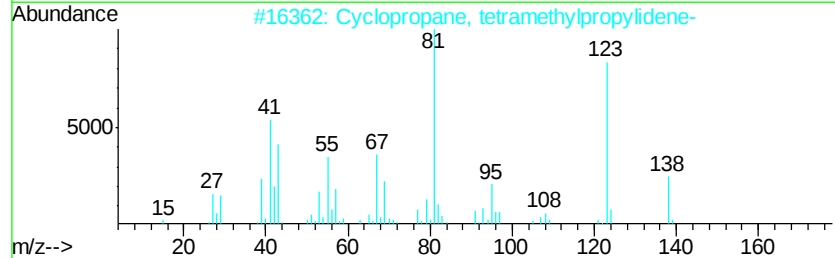
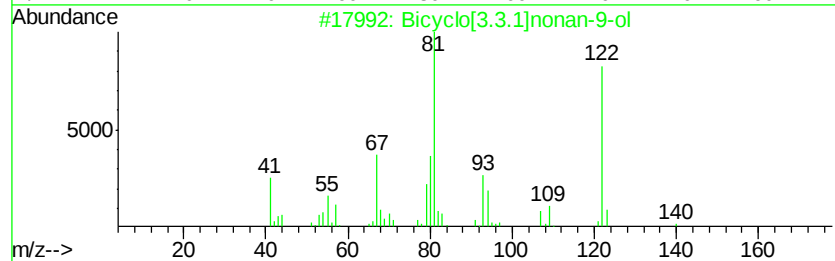
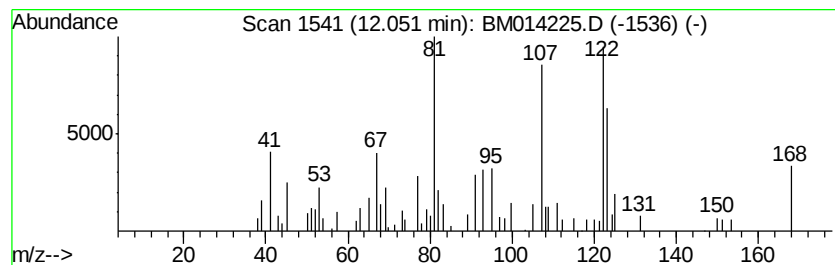
Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-BM020718.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Bicyclo[3.3.1]nonan-9-ol Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.05	2.42 ng	133262	Naphthalene-d8	10.31
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Bicyclo[3.3.1]nonan-9-ol	140 C9H16O	015598-80-8 43
2		Cyclopropane, tetramethylpropyli...	138 C10H18	024519-04-8 35
3		Phenol, 2,4-dimethyl-	122 C8H10O	000105-67-9 30
4		Phenol, 3,4-dimethyl-	122 C8H10O	000095-65-8 30
5		Phenol, 3-ethyl-	122 C8H10O	000620-17-7 30



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA M\DATA\BM021318\
Data File : BM014225.D
Acq On : 13 Feb 2018 17:33
Operator : SJ/JU
Sample : J1518-08
Misc :
ALS Vial : 9 Sample Multiplier: 1

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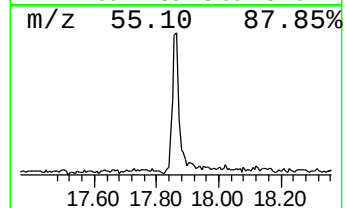
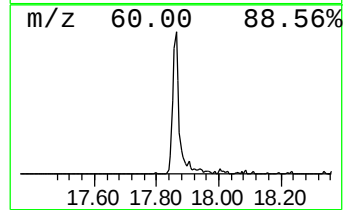
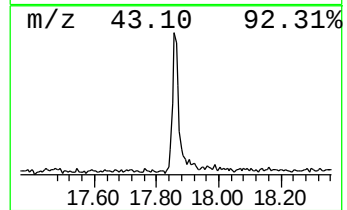
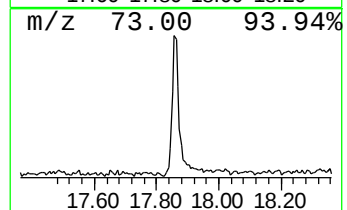
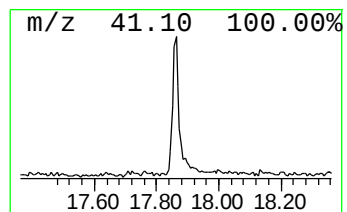
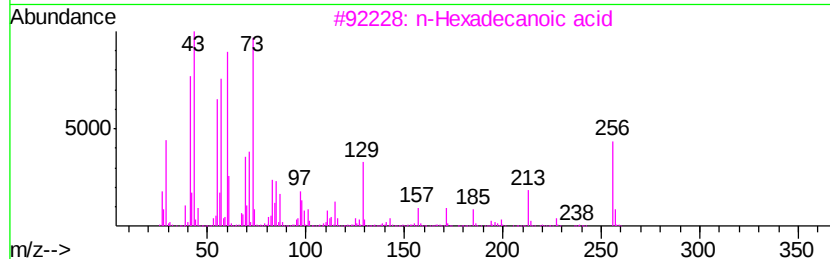
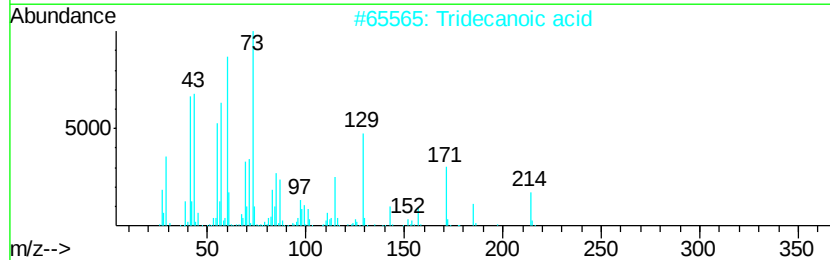
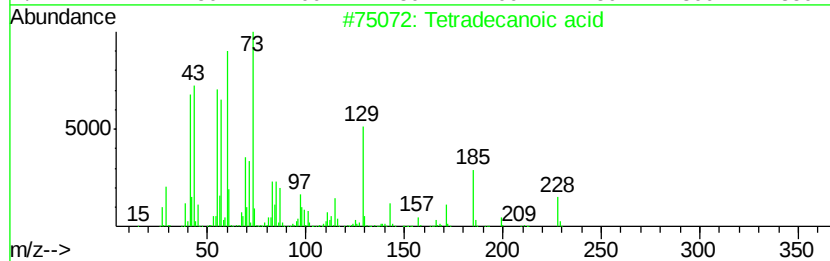
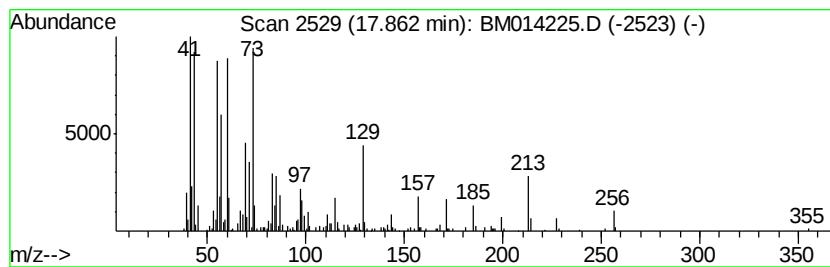
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Tetradecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.86	6.30 ng	649699	Phenanthrene-d10	16.95

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetradecanoic acid	228	C14H28O2	000544-63-8	95
2		Tridecanoic acid	214	C13H26O2	000638-53-9	90
3		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	80
4		Dodecanoic acid	200	C12H24O2	000143-07-7	58
5		n-Decanoic acid	172	C10H20O2	000334-48-5	47



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA M\DATA\BM021318\
Data File : BM014225.D
Acq On : 13 Feb 2018 17:33
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Sample : J1518-08
Misc :
ALS Vial : 9 Sample Multiplier: 1

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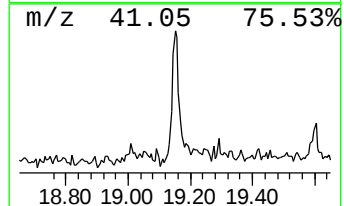
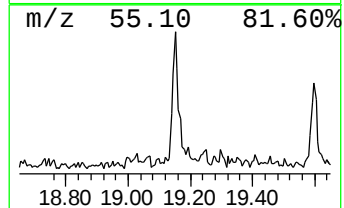
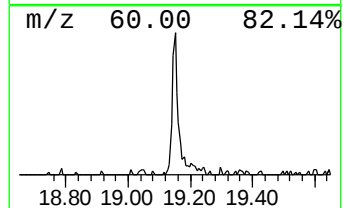
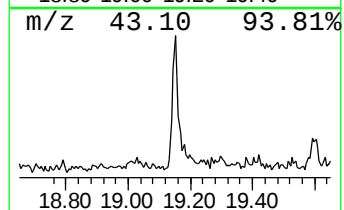
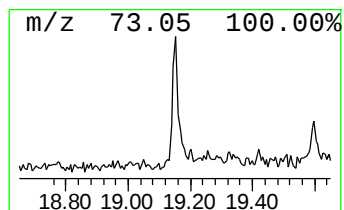
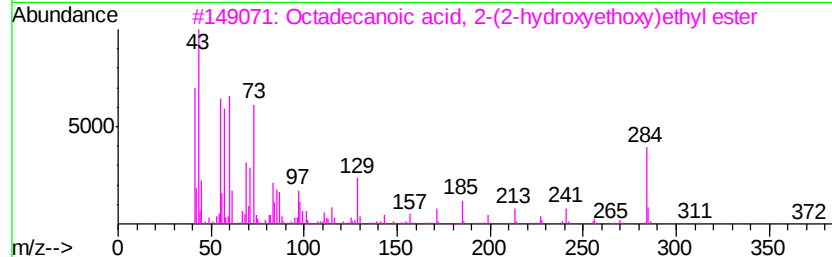
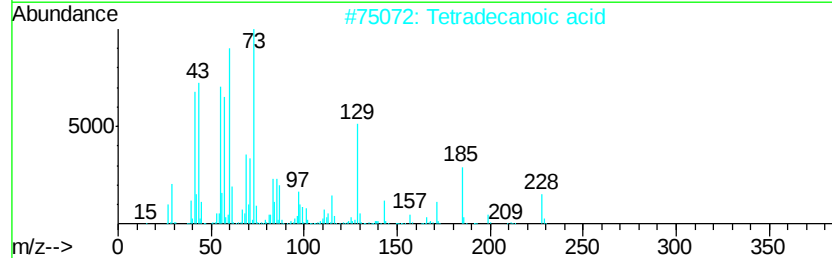
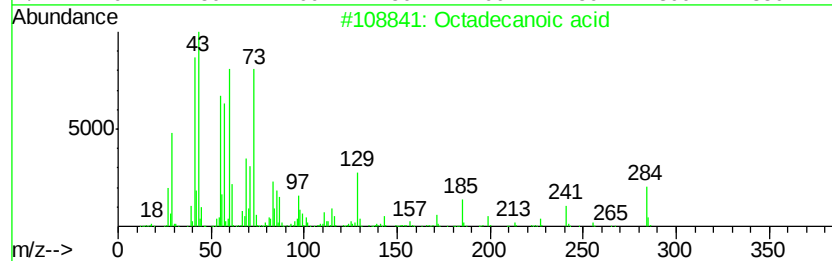
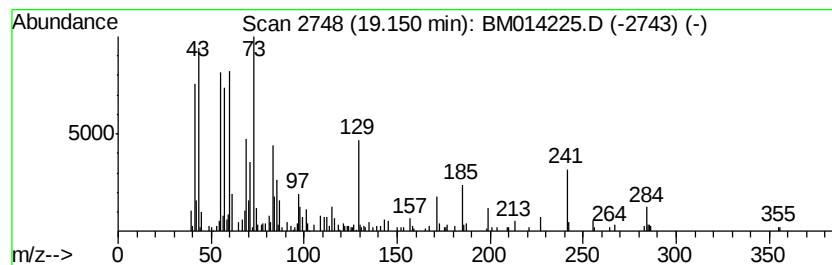
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Octadecanoic acid Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.15	3.18 ng	354562	Chrysene-d12	21.16

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	95
2			Tetradecanoic acid	228	C14H28O2	000544-63-8	74
3			Octadecanoic acid, 2-(2-hydroxyve...	372	C22H44O4	000106-11-6	59
4			12-Bromododecanoic acid	278	C12H23BrO2	073367-80-3	53
5			11-Bromoundecanoic acid	264	C11H21BrO2	002834-05-1	50



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Data File : BM014225.D
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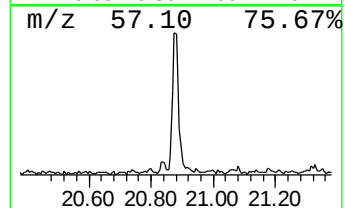
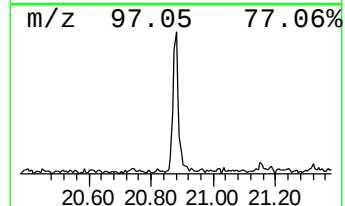
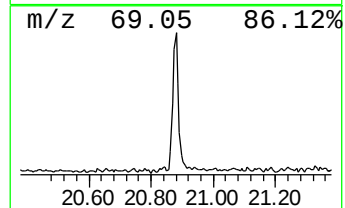
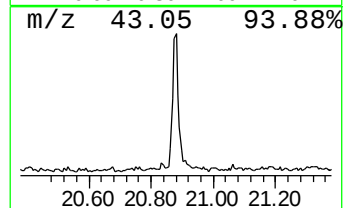
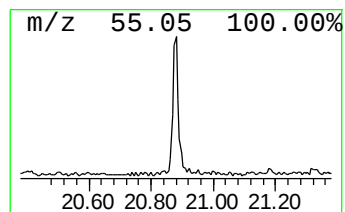
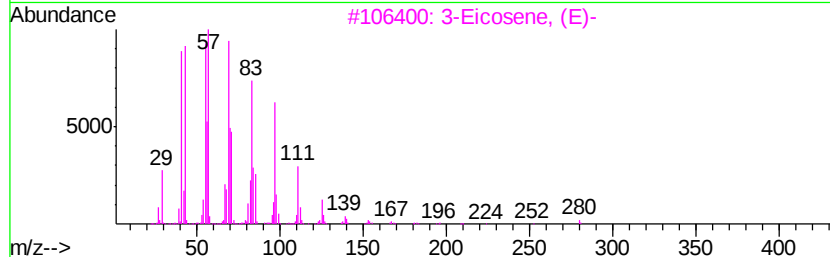
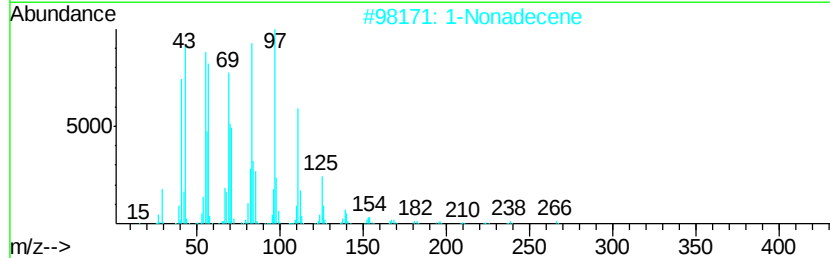
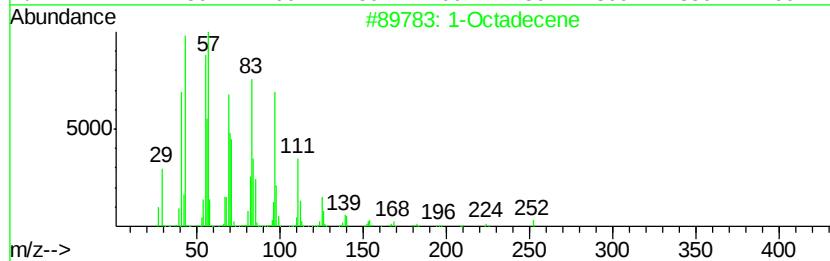
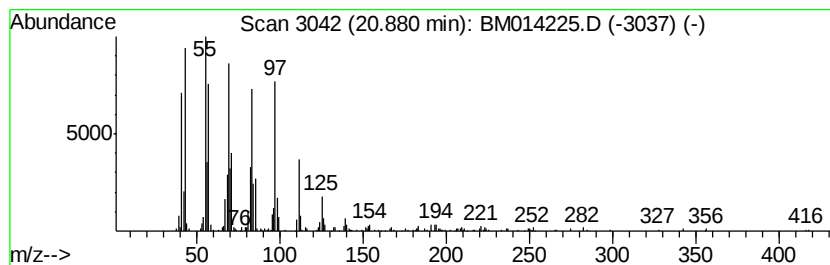
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 1-Octadecene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.88	6.25 ng	696401	Chrysene-d12	21.16

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Octadecene	252	C18H36	000112-88-9	95
2			1-Nonadecene	266	C19H38	018435-45-5	92
3			3-Eicosene, (E)-	280	C20H40	074685-33-9	83
4			5-Eicosene, (E)-	280	C20H40	074685-30-6	70
5			1-Hexadecanol	242	C16H34O	036653-82-4	68



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_M\DATA\BM021318\
Data File : BM014225.D
Acq On : 13 Feb 2018 17:33
Operator : SJ/JU
Sample : J1518-08
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-BM020718.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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unknown7.07	7.07	89.8	ng	3330250	1	7.54	741656	20.0
Benzene, 1,2,4-tr...	7.18	9.9	ng	366194	1	7.54	741656	20.0
Benzene, 1-ethyl-...	8.17	3.5	ng	128901	1	7.54	741656	20.0
Hexanoic acid, 2-...	8.87	14.4	ng	534031	1	7.54	741656	20.0
unknown9.72	9.72	2.6	ng	145320	2	10.31	1101580	20.0
Ethanone, 1-(3,4-...	11.56	3.5	ng	195648	2	10.31	1101580	20.0
Bicyclo[3.3.1]non...	12.05	2.4	ng	133262	2	10.31	1101580	20.0
Tetradecanoic acid	17.86	6.3	ng	649699	4	16.95	2063850	20.0
Octadecanoic acid	19.15	3.2	ng	354562	5	21.16	2226980	20.0
1-Octadecene	20.88	6.3	ng	696401	5	21.16	2226980	20.0