

Data Path : Z:\HPCHEM1\BNA F\DATA\BF042517\
 Data File : BF094614.D
 Acq On : 25 Apr 2017 19:20
 Operator : SJ/MA
 Sample : I2824-02
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 E2-CONCRETE-BATCH7

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-BF041717.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	2.702	126	130	139	rVB	101635	124771	1.68%	0.394%
2	3.366	236	243	246	rBV	3790182	5663481	76.38%	17.866%
3	3.937	328	340	342	rBV	3347515	7414545	100.00%	23.390%
4	5.372	580	584	603	rBV	848035	856798	11.56%	2.703%
5	5.501	603	606	610	rBV	107030	102150	1.38%	0.322%
6	5.748	645	648	653	rBV	736966	683994	9.23%	2.158%
7	5.931	674	679	682	rBV	2375336	2180854	29.41%	6.880%
8	6.072	700	703	710	rBV	258964	229308	3.09%	0.723%
9	6.260	731	735	747	rVB	280536	345323	4.66%	1.089%
10	6.401	754	759	762	rBV	1715100	1681595	22.68%	5.305%
11	6.890	837	842	850	rVB	137920	166342	2.24%	0.525%
12	7.019	860	864	866	rBV	161346	204808	2.76%	0.646%
13	7.037	866	867	875	rVV	132257	102355	1.38%	0.323%
14	7.242	898	902	906	rBV	1015321	981942	13.24%	3.098%
15	8.572	1123	1128	1131	rVB	2972088	3074967	41.47%	9.700%
16	8.707	1148	1151	1156	rVB	91629	83192	1.12%	0.262%
17	9.066	1209	1212	1216	rBV	209976	204030	2.75%	0.644%
18	9.378	1259	1265	1274	rBV3	978541	1141815	15.40%	3.602%
19	9.978	1364	1367	1372	rBV	264749	239571	3.23%	0.756%
20	10.254	1409	1414	1417	rBV	100261	121546	1.64%	0.383%
21	10.301	1419	1422	1427	rVB2	63547	83796	1.13%	0.264%
22	10.560	1463	1466	1468	rBV	335321	346503	4.67%	1.093%
23	10.578	1468	1469	1473	rVB	188891	166075	2.24%	0.524%
24	10.801	1504	1507	1511	rBV3	60031	80119	1.08%	0.253%
25	11.113	1557	1560	1564	rBV	290849	291891	3.94%	0.921%
26	11.154	1564	1567	1571	rVB	108806	118313	1.60%	0.373%
27	11.201	1571	1575	1578	rBV	852913	827570	11.16%	2.611%
28	11.448	1614	1617	1621	rVB4	61635	81519	1.10%	0.257%
29	11.642	1647	1650	1653	rVB	231808	211092	2.85%	0.666%
30	12.148	1733	1736	1739	rVB	165854	148883	2.01%	0.470%
31	12.630	1814	1818	1821	rVB	93220	96519	1.30%	0.304%
32	13.177	1907	1911	1915	rVB	1935947	2158126	29.11%	6.808%
33	14.348	2105	2110	2122	rBV3	77042	150164	2.03%	0.474%
34	14.448	2123	2127	2135	rBV	609408	669234	9.03%	2.111%

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Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
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ClientSampleId :
E2-CONCRETE-BATCH7

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

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Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	16.077	2399	2404	2416	rBV	584512	666782	8.99%	2.103%
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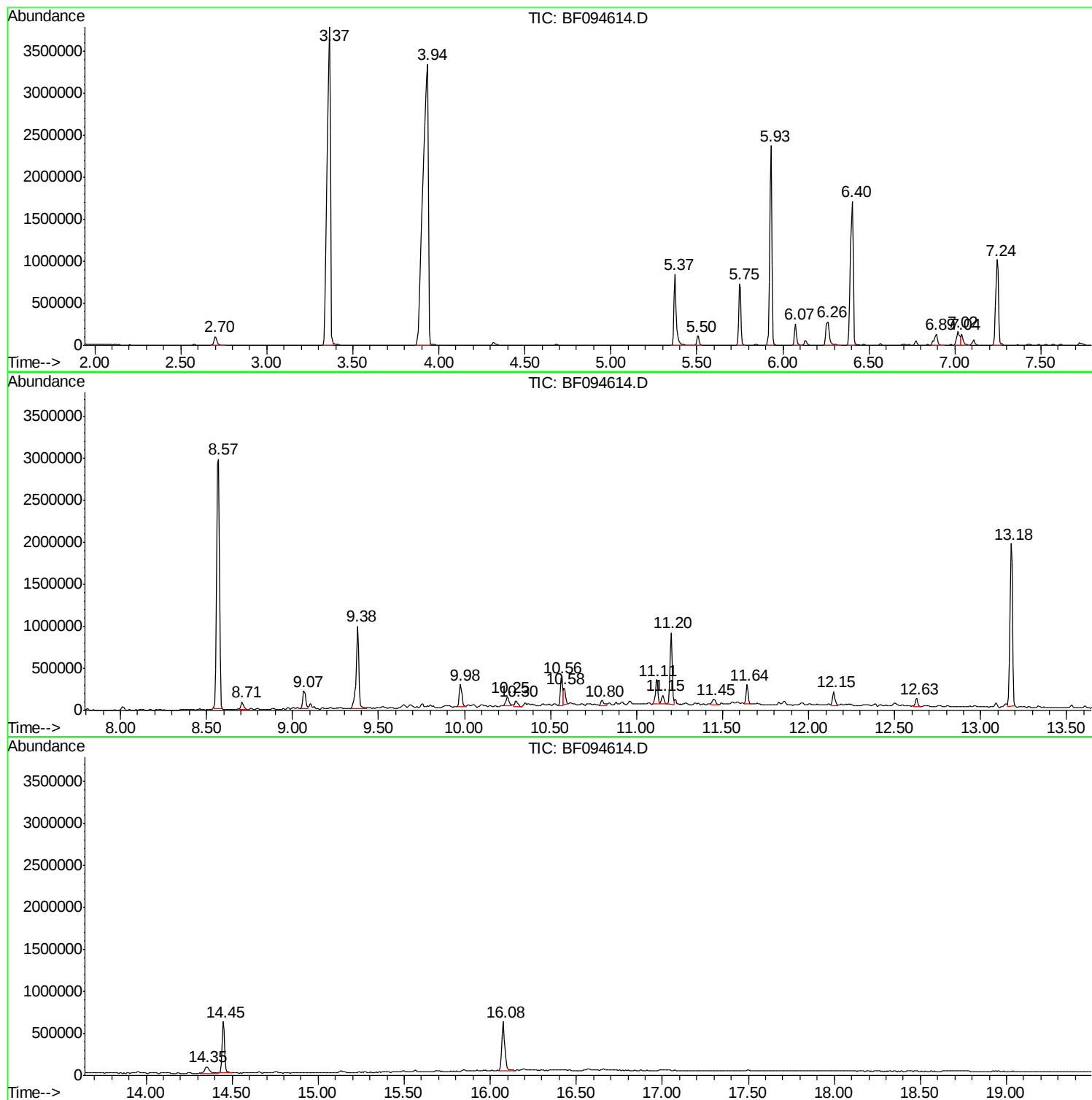
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF041717.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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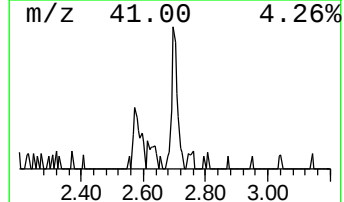
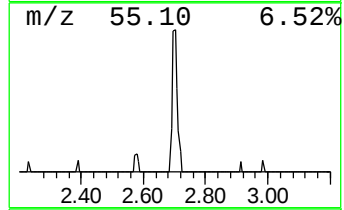
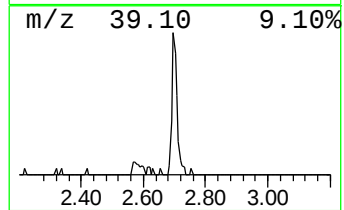
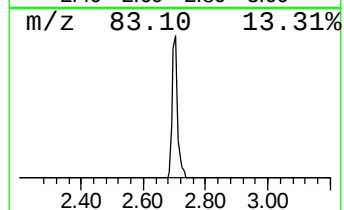
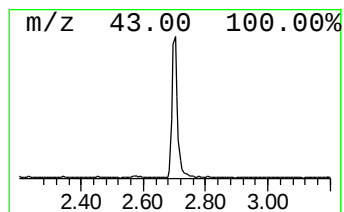
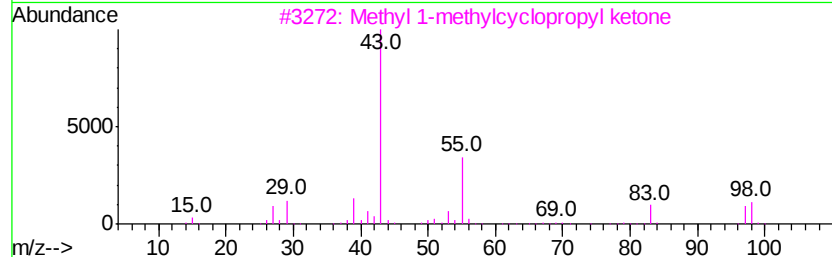
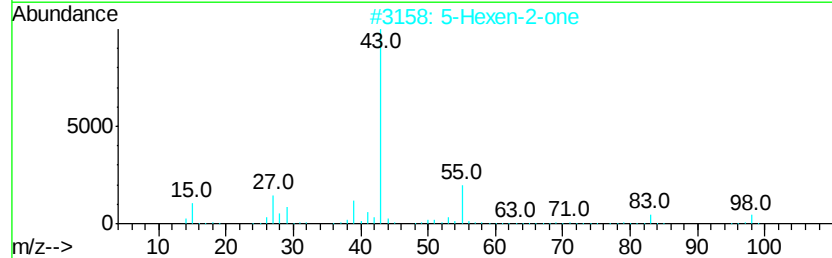
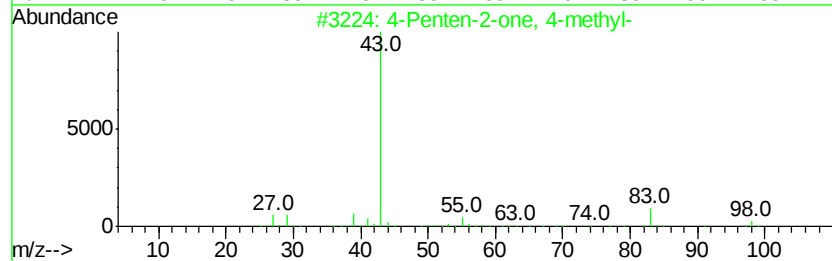
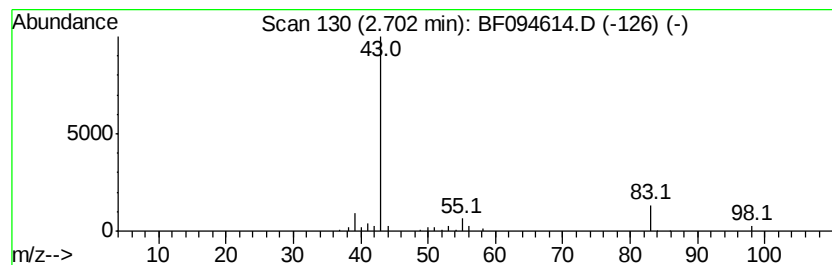
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 4-Penten-2-one, 4-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.70	3.65 ng	124771	1,4-Dichlorobenzene-d4	5.75

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Penten-2-one, 4-methyl-	98	C6H10O	003744-02-3	80
2			5-Hexen-2-one	98	C6H10O	000109-49-9	58
3			Methyl 1-methylcyclopropyl ketone	98	C6H10O	001567-75-5	33
4			2-Vinylethyl acetate	114	C6H10O2	001576-84-7	23
5			2-Pentanone, 3-methylene-	98	C6H10O	004359-77-7	17



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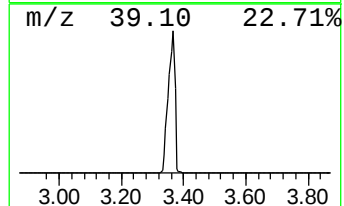
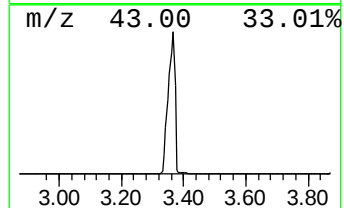
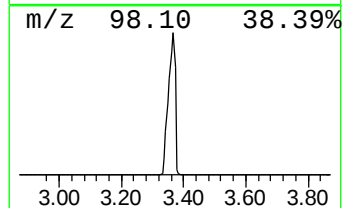
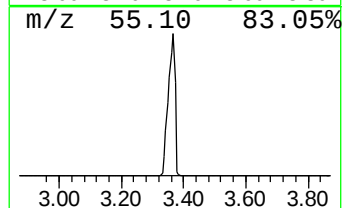
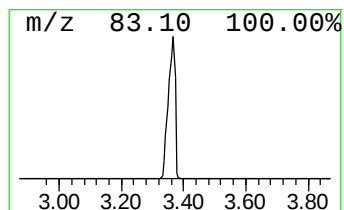
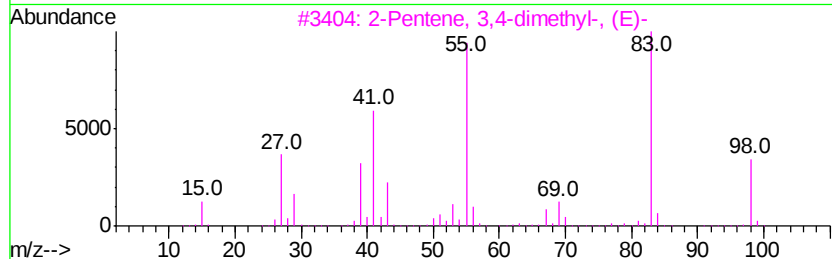
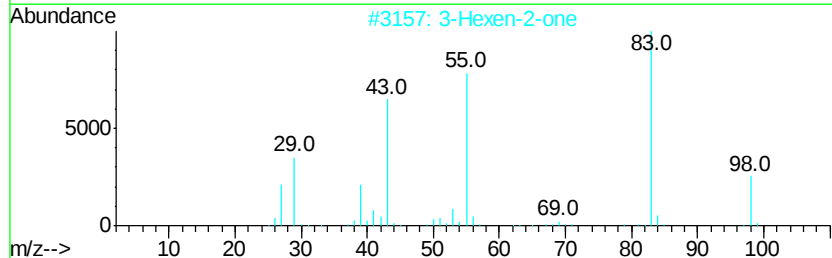
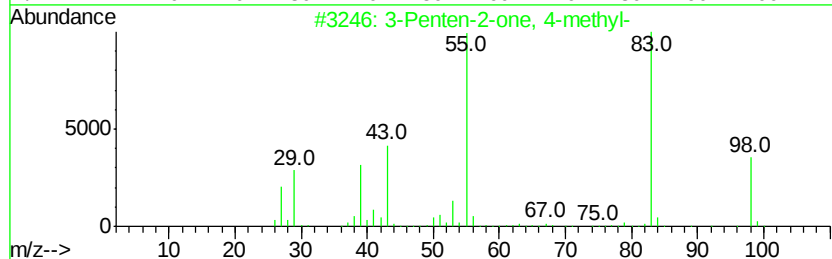
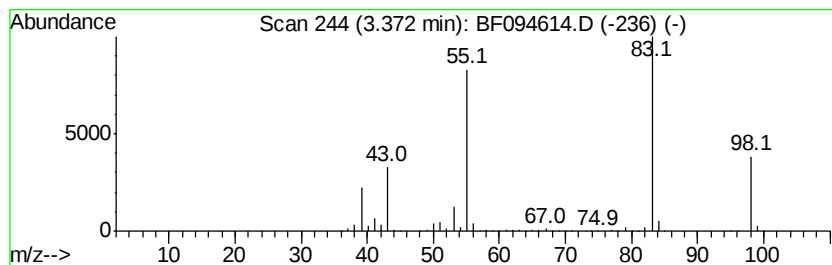
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF041717.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 3-Penten-2-one, 4-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.37	165.60 ng	5663480	1,4-Dichlorobenzene-d4	5.75

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	95
2		3-Hexen-2-one	98	C6H10O	000763-93-9	91
3		2-Pentene, 3,4-dimethyl-, (E)-	98	C7H14	004914-92-5	90
4		2-Pentene, 3,4-dimethyl-, (Z)-	98	C7H14	004914-91-4	86
5		2-Pentene, 2,4-dimethyl-	98	C7H14	000625-65-0	80



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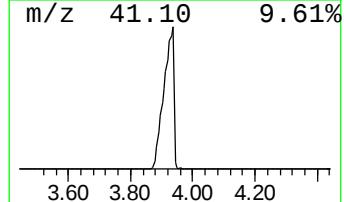
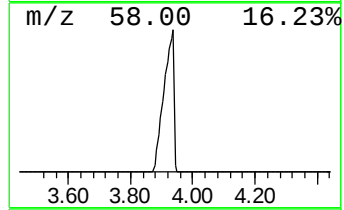
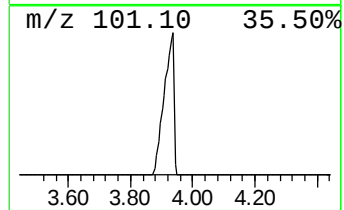
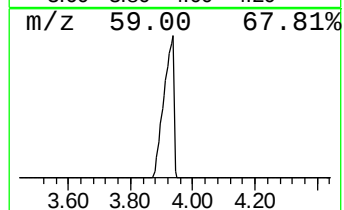
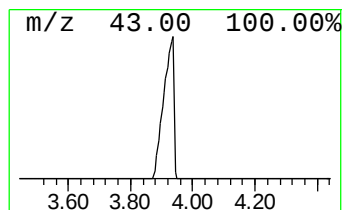
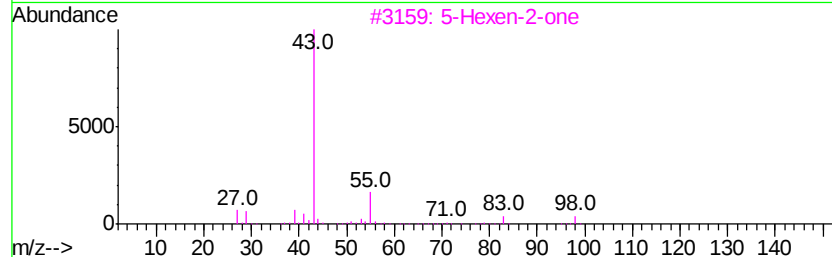
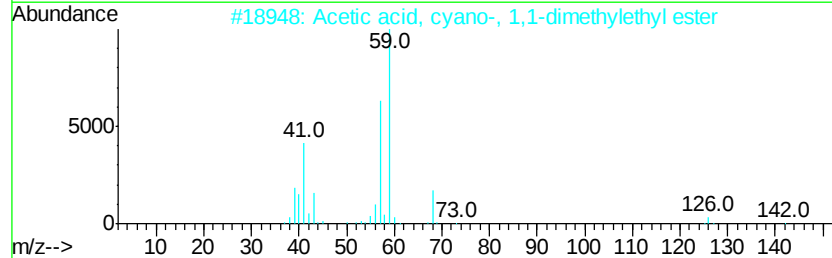
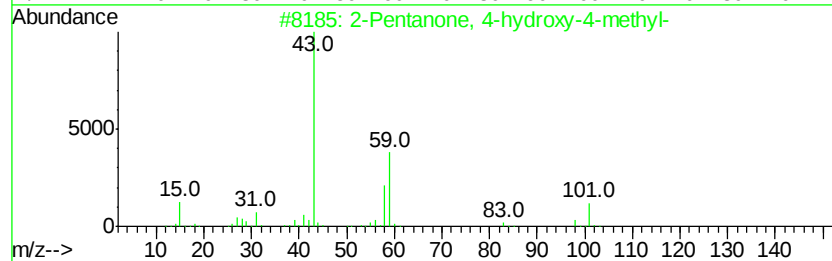
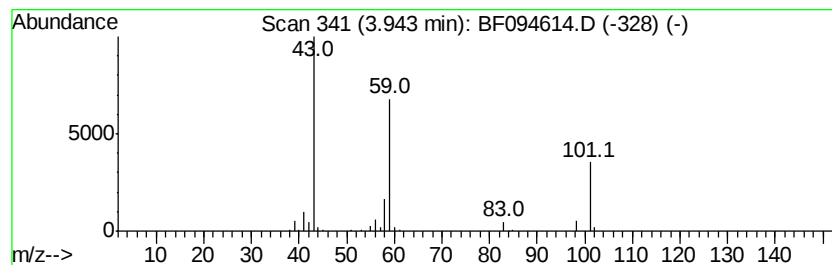
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TIC Library : C:\DATABASE\NIST11.L
 TIC Integration Parameters: LSCINT.P

 Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.94	216.80 ng	7414550	1,4-Dichlorobenzene-d4	5.75

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72
2		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	25
3		5-Hexen-2-one	98	C6H10O	000109-49-9	9
4		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5		Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	9



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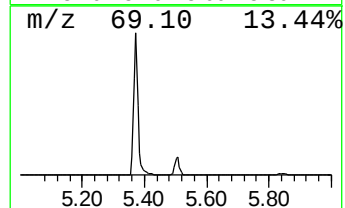
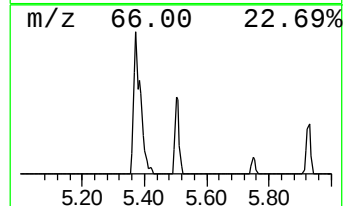
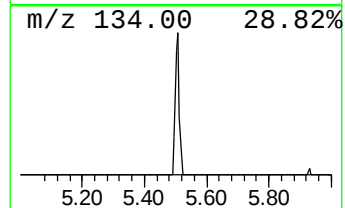
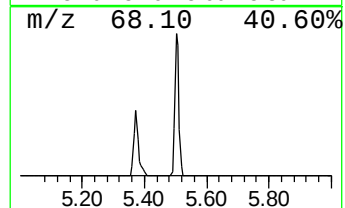
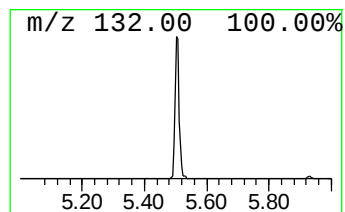
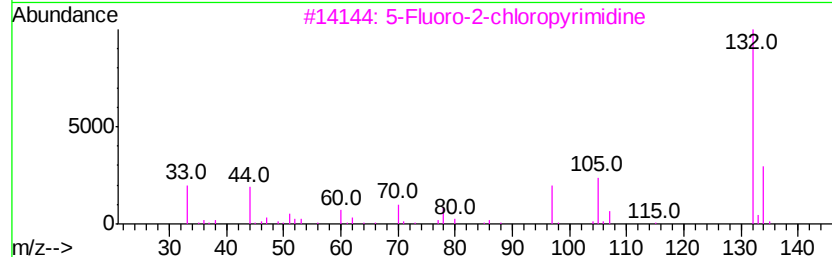
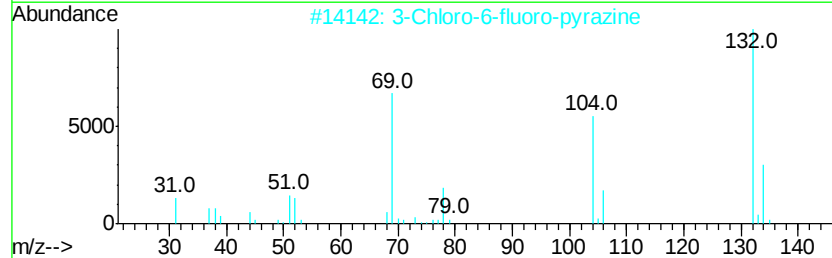
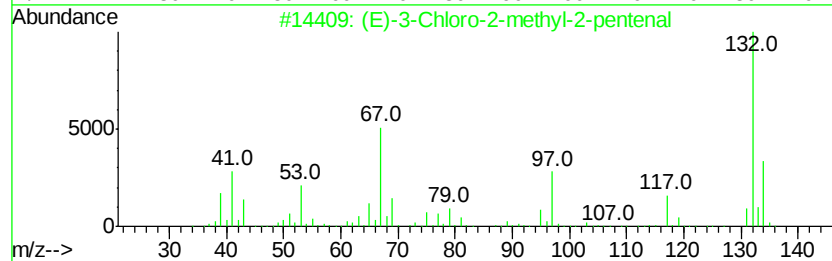
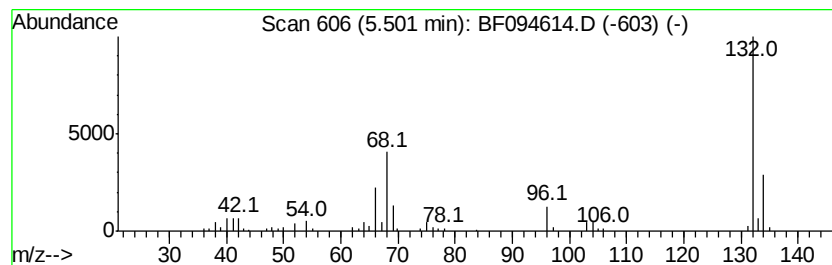
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 4 unknown5.50 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.50	2.99 ng	102150	1,4-Dichlorobenzene-d4	5.75

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	25	
2	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16	
3	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12	
4	1H-Indazole, 7-methyl-	132	C8H8N2	1000316-00-7	10	
5	5-Aminoindole	132	C8H8N2	005192-03-0	10	



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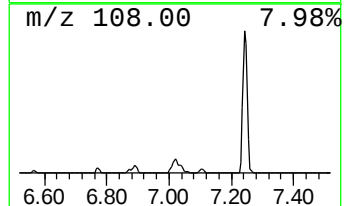
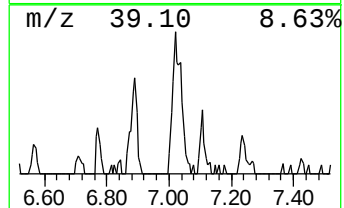
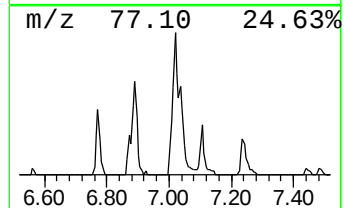
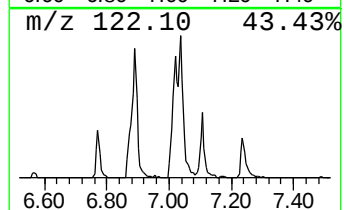
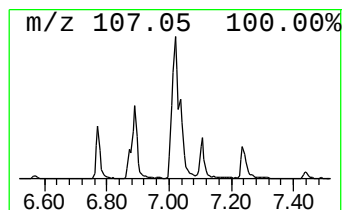
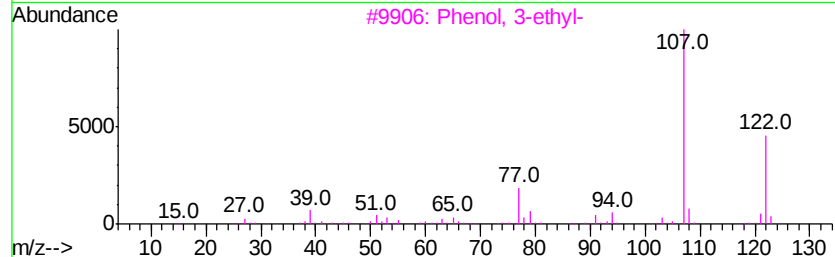
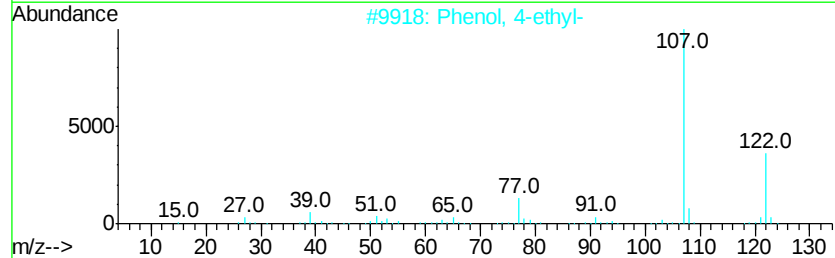
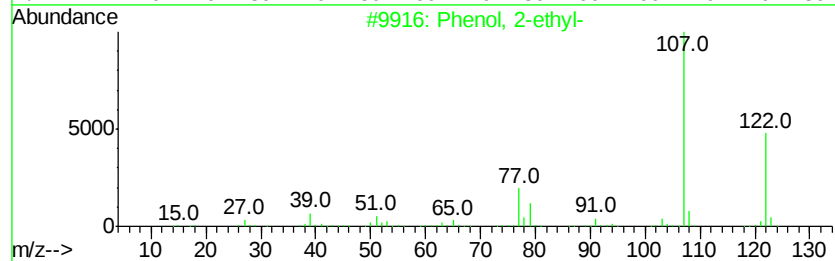
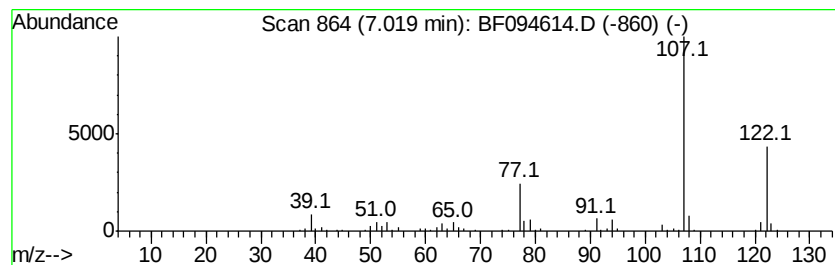
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TIC Integration Parameters: LSCINT.P

Peak Number 6 Phenol, 2-ethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.02	4.17 ng	204808	Napthalene-d8	7.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenol, 2-ethyl-	122	C8H10O	000090-00-6	91
2		Phenol, 4-ethyl-	122	C8H10O	000123-07-9	90
3		Phenol, 3-ethyl-	122	C8H10O	000620-17-7	90
4		Phenol, 3,4-dimethyl-	122	C8H10O	000095-65-8	78
5		Phenol, 2,3-dimethyl-	122	C8H10O	000526-75-0	78



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042517\
Data File : BF094614.D
Acq On : 25 Apr 2017 19:20
Operator : SJ/MA
Sample : I2824-02
Misc :
ALS Vial : 19 Sample Multiplier: 1

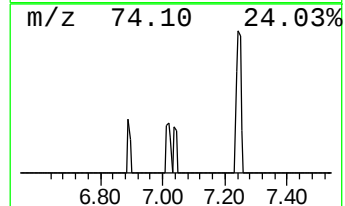
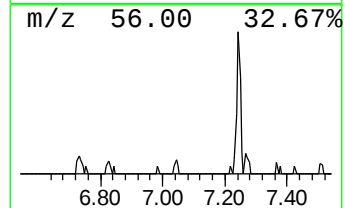
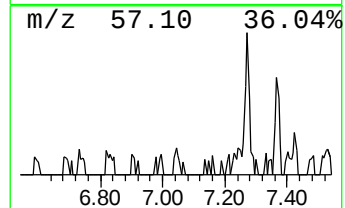
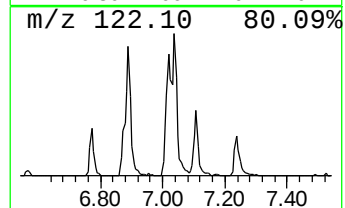
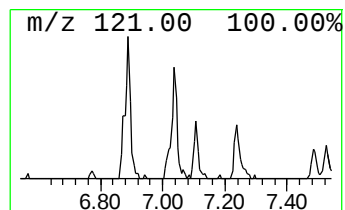
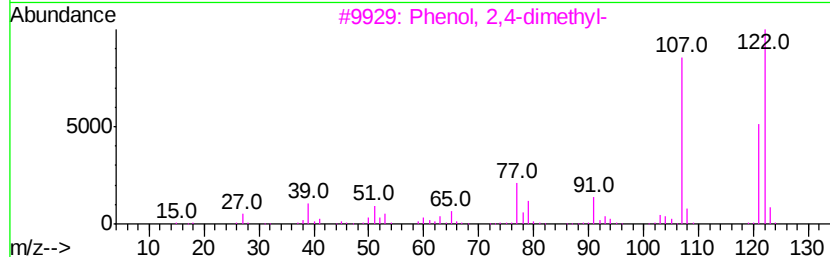
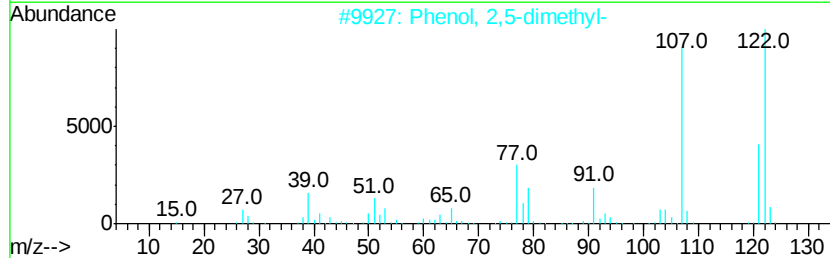
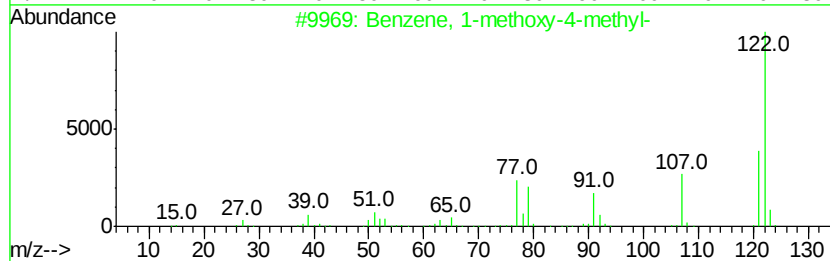
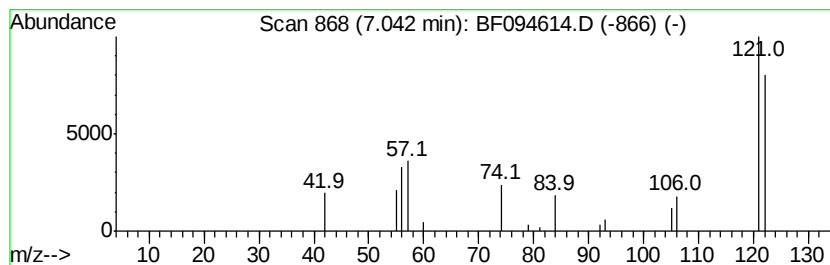
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ClientSampleId :
E2-CONCRETE-BATCH7

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF041717.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Benzene, 1-methoxy-4-methyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.04	2.08 ng	102355	Napthalene-d8	7.24
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Benzene, 1-methoxy-4-methyl-	122 C8H10O	000104-93-8	90
2	Phenol, 2,5-dimethyl-	122 C8H10O	000095-87-4	78
3	Phenol, 2,4-dimethyl-	122 C8H10O	000105-67-9	78
4	Phenol, 3,5-dimethyl-	122 C8H10O	000108-68-9	72
5	Phenol, 2,6-dimethyl-	122 C8H10O	000576-26-1	72



Data Path : Z:\HPCHEM1\BNA F\DATA\BF042517\
Data File : BF094614.D
Acq On : 25 Apr 2017 19:20
Operator : SJ/MA
Sample : I2824-02
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
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ClientSampleId :
E2-CONCRETE-BATCH7

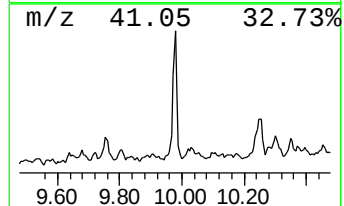
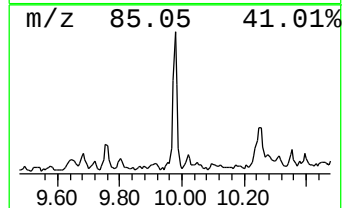
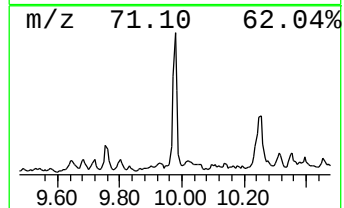
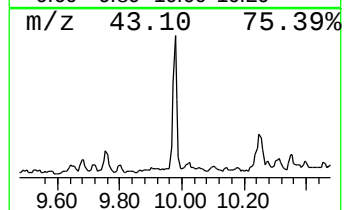
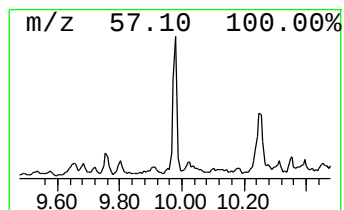
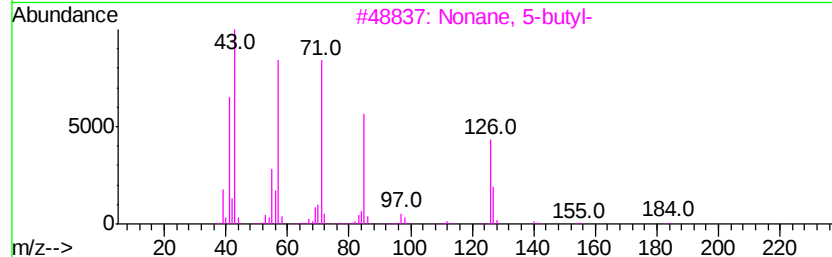
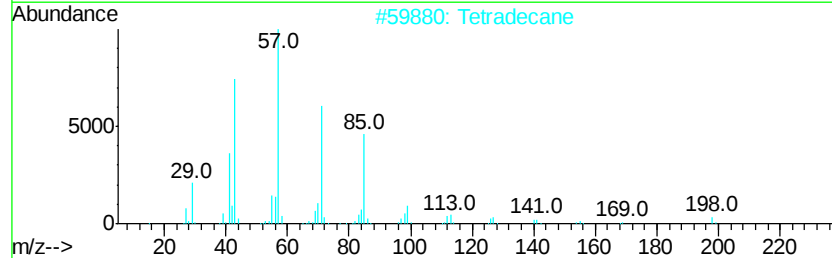
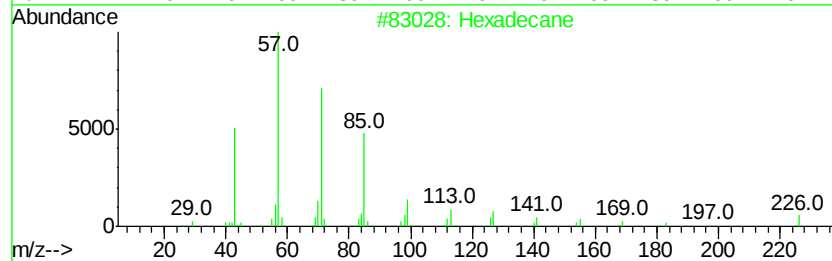
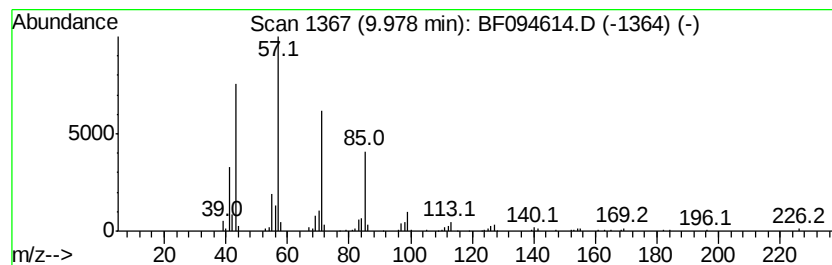
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Hexadecane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.98	4.20 ng	239571	Acenaphthene-d10	9.38

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane	226	C16H34	000544-76-3	95
2		Tetradecane	198	C14H30	000629-59-4	91
3		Nonane, 5-butyl-	184	C13H28	017312-63-9	90
4		Heneicosane	296	C21H44	000629-94-7	90
5		Pentadecane	212	C15H32	000629-62-9	90



Data Path : Z:\HPCHEM1\BNA F\DATA\BF042517\
Data File : BF094614.D
Acq On : 25 Apr 2017 19:20
Operator : SJ/MA
Sample : I2824-02
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
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ClientSampleId :
E2-CONCRETE-BATCH7

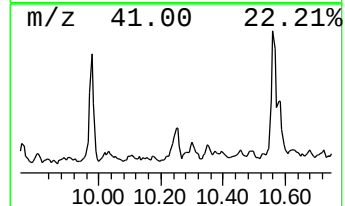
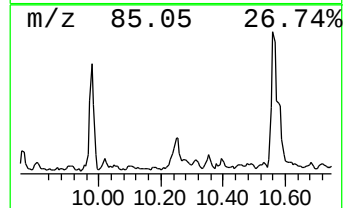
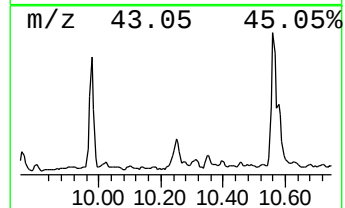
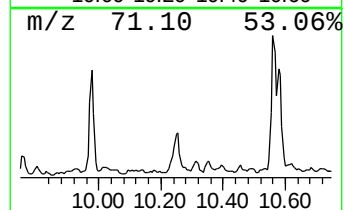
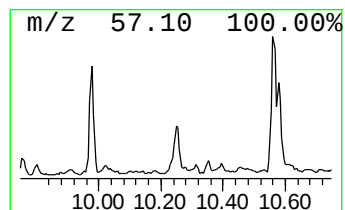
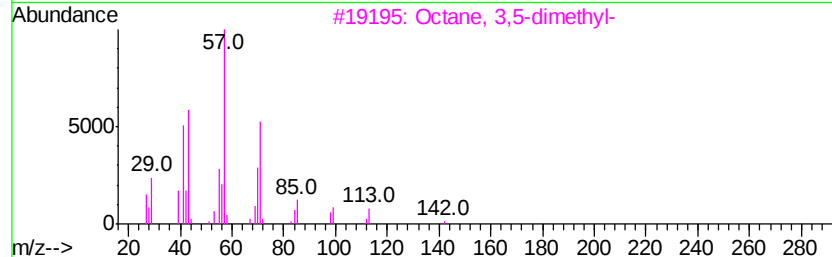
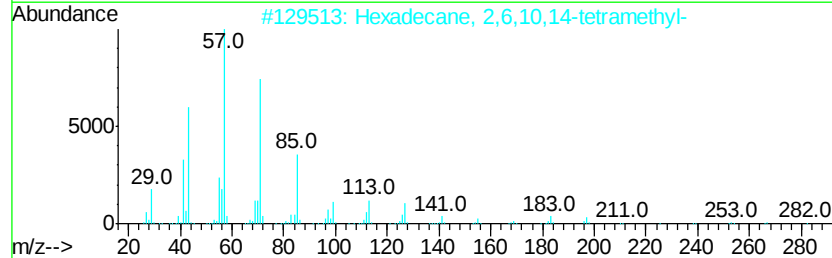
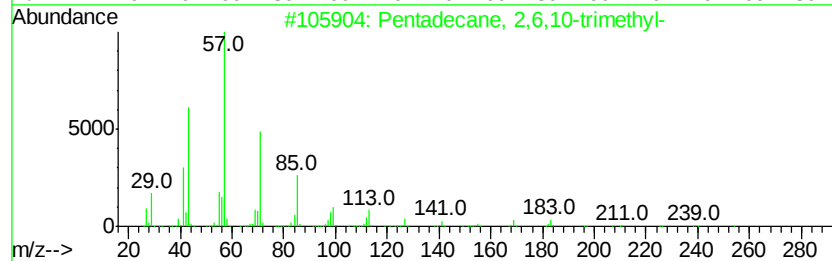
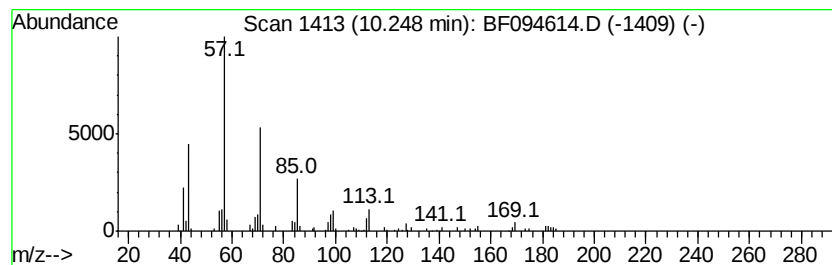
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Pentadecane, 2,6,10-trimethyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.25	2.13 ng	121546	Acenaphthene-d10	9.38

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentadecane, 2,6,10-trimethyl-	254	C18H38	003892-00-0	86
2		Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	74
3		Octane, 3,5-dimethyl-	142	C10H22	015869-93-9	72
4		Heptadecane, 7-methyl-	254	C18H38	020959-33-5	72
5		Undecane, 6-ethyl-	184	C13H28	017312-60-6	64



Data Path : Z:\HPCHEM1\BNA F\DATA\BF042517\
Data File : BF094614.D
Acq On : 25 Apr 2017 19:20
Operator : SJ/MA
Sample : I2824-02
Misc :
ALS Vial : 19 Sample Multiplier: 1

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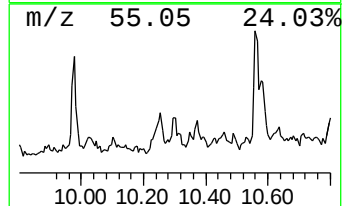
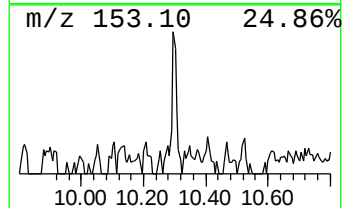
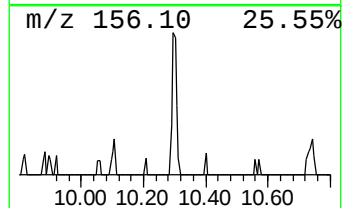
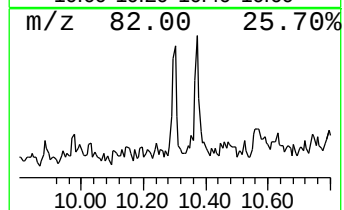
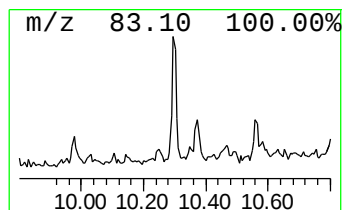
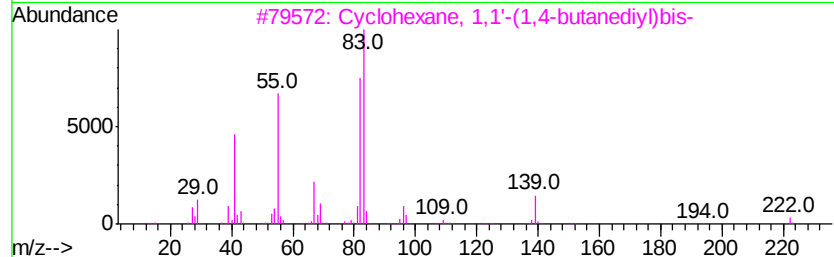
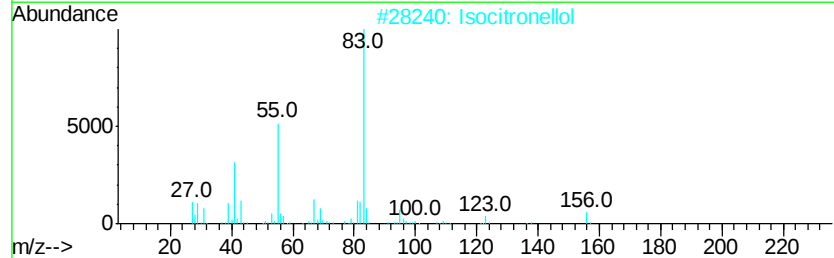
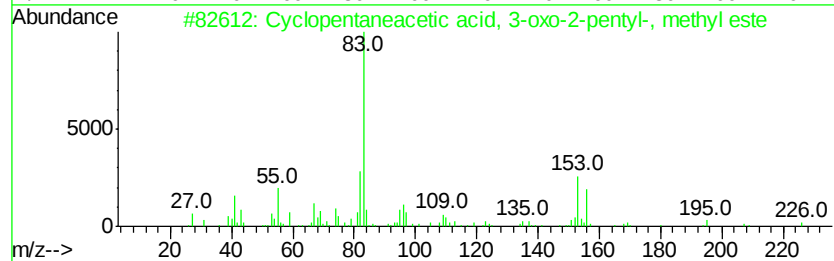
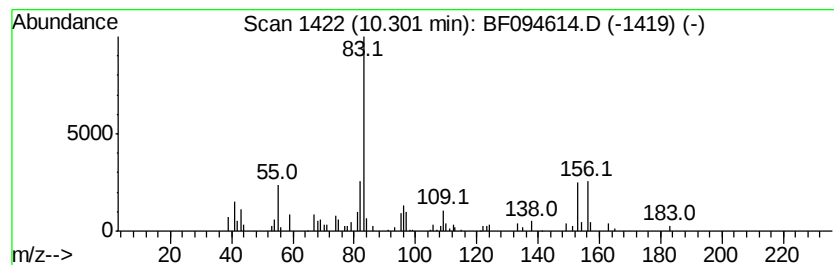
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Cyclopentaneacetic acid, 3-... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.30	2.03 ng	83796	Phenanthrene-d10	11.20

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentaneacetic acid, 3-oxo-2...	226	C13H22O3	024851-98-7	80
2		Isocitronellol	156	C10H20O	018479-52-2	50
3		Cyclohexane, 1,1'-(1,4-butanediol...	222	C16H30	006165-44-2	43
4		1,1-Dicyclohexylbutane	222	C16H30	054890-00-5	40
5		Oxalic acid, cyclohexyl pentyl e...	242	C13H22O4	1000309-30-6	35



Data Path : Z:\HPCHEM1\BNA F\DATA\BF042517\
Data File : BF094614.D
Acq On : 25 Apr 2017 19:20
Operator : SJ/MA
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Instrument :
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ClientSampleId :
E2-CONCRETE-BATCH7

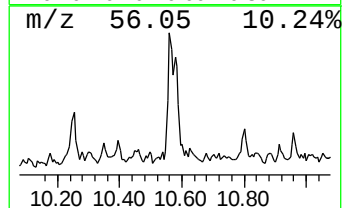
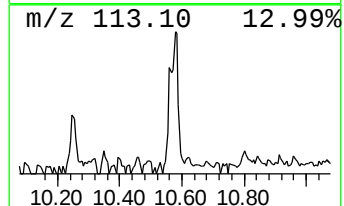
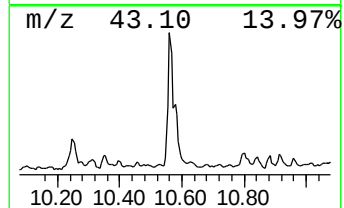
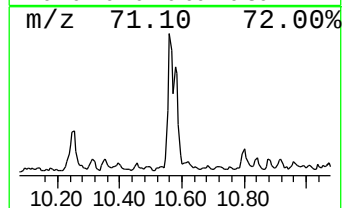
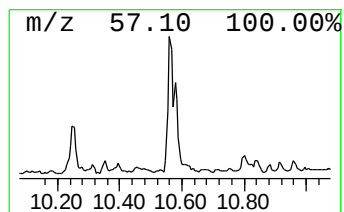
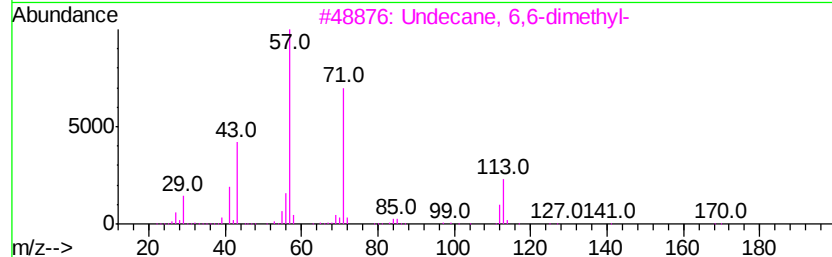
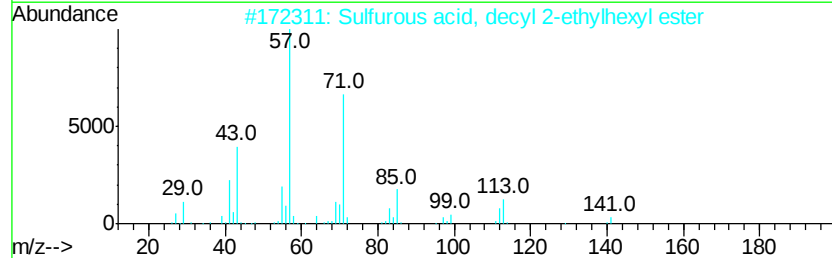
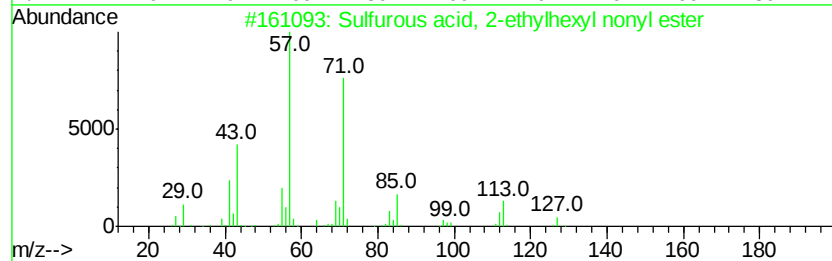
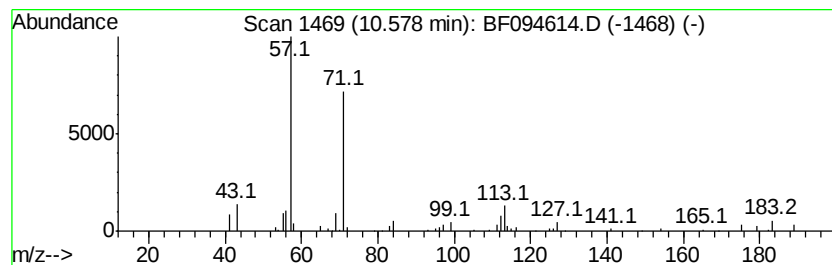
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Sulfurous acid, 2-ethylhexy... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.58	4.01 ng	166075	Phenanthrene-d10	11.20

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Sulfurous acid, 2-ethylhexyl non...	320	C17H36O3S	1000309-19-2	64
2		Sulfurous acid, decyl 2-ethylhex...	334	C18H38O3S	1000309-19-3	59
3		Undecane, 6,6-dimethyl-	184	C13H28	017312-76-4	53
4		Undecane, 4,6-dimethyl-	184	C13H28	017312-82-2	53
5		Sulfurous acid, 2-ethylhexyl tri...	376	C21H44O3S	1000309-19-6	50



Data Path : Z:\HPCHEM1\BNA F\DATA\BF042517\
Data File : BF094614.D
Acq On : 25 Apr 2017 19:20
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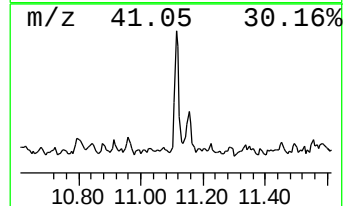
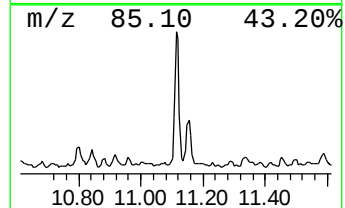
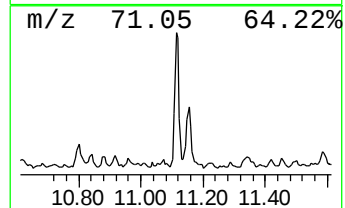
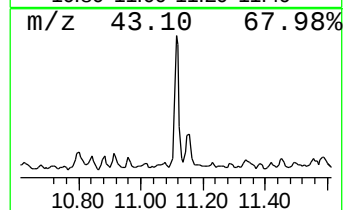
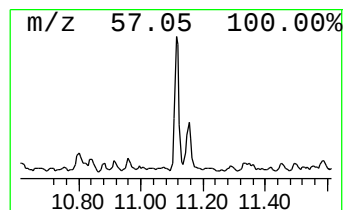
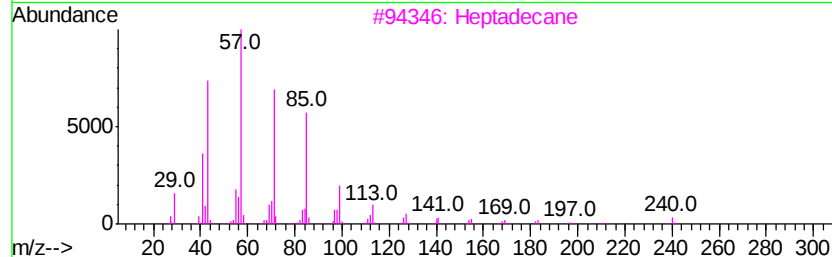
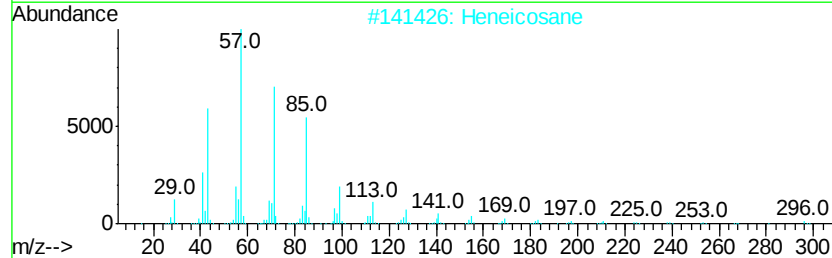
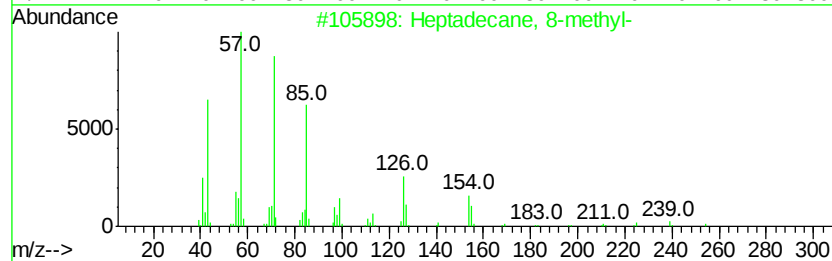
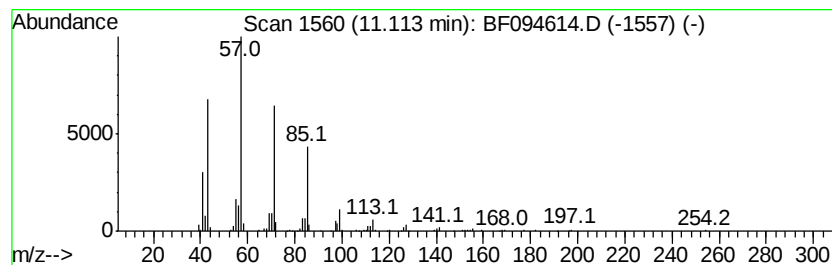
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Heptadecane, 8-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.11	7.05 ng	291891	Phenanthrene-d10	11.20

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane, 8-methyl-	254	C18H38	013287-23-5	90
2		Heneicosane	296	C21H44	000629-94-7	90
3		Heptadecane	240	C17H36	000629-78-7	90
4		Heptacosane	380	C27H56	000593-49-7	90
5		Tetracosane	338	C24H50	000646-31-1	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042517\
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Acq On : 25 Apr 2017 19:20
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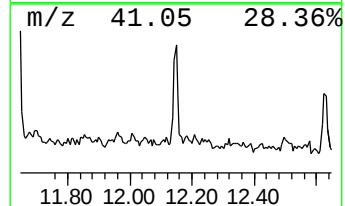
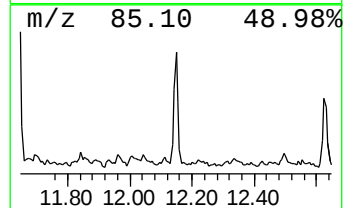
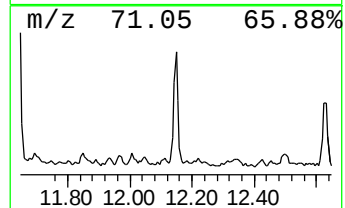
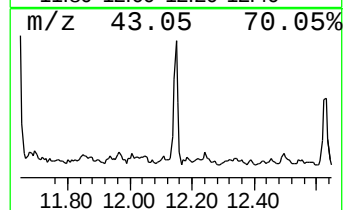
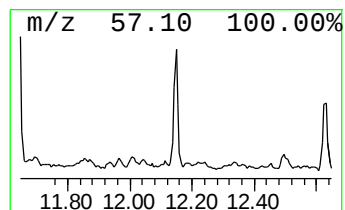
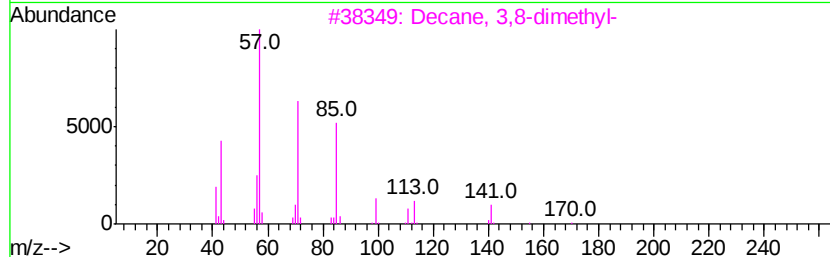
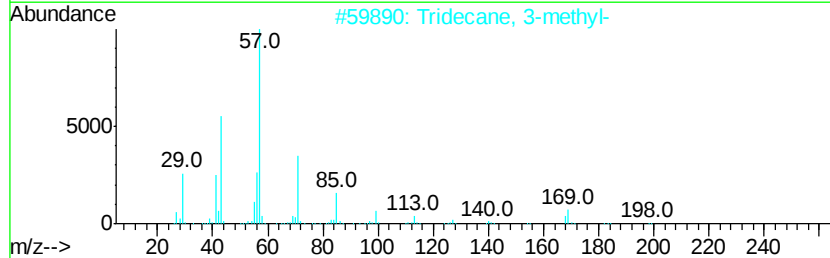
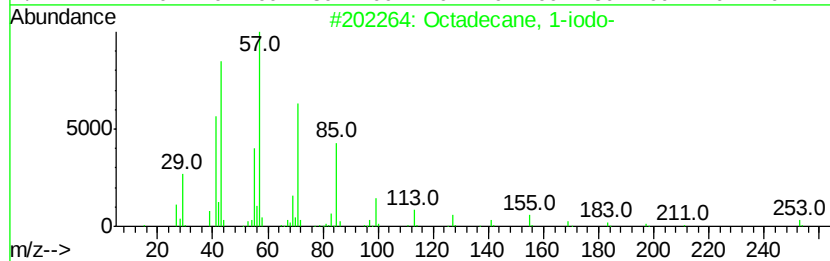
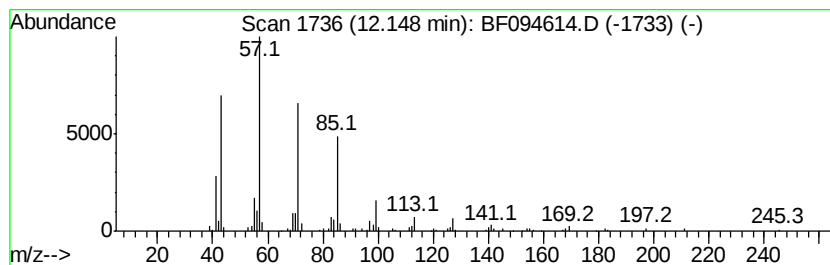
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Octadecane, 1-iodo- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.15	3.60 ng	148883	Phenanthrene-d10	11.20

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecane, 1-iodo-	380	C18H37I	000629-93-6	90
2			Tridecane, 3-methyl-	198	C14H30	006418-41-3	90
3			Decane, 3,8-dimethyl-	170	C12H26	017312-55-9	78
4			Heneicosane	296	C21H44	000629-94-7	74
5			Undecane, 2,7-dimethyl-	184	C13H28	017301-24-5	72



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042517\
Data File : BF094614.D
Acq On : 25 Apr 2017 19:20
Operator : SJ/MA
Sample : I2824-02
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
E2-CONCRETE-BATCH7

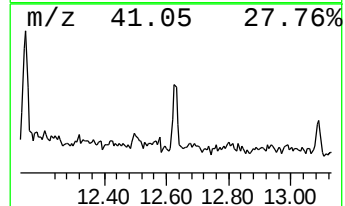
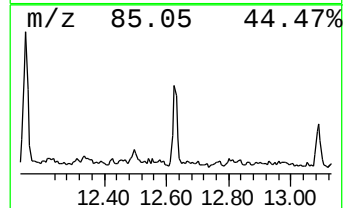
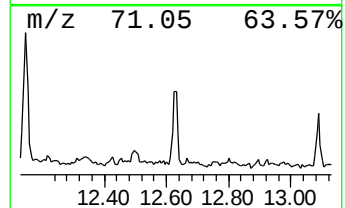
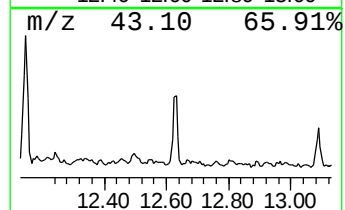
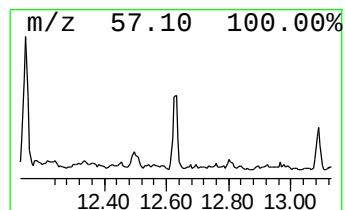
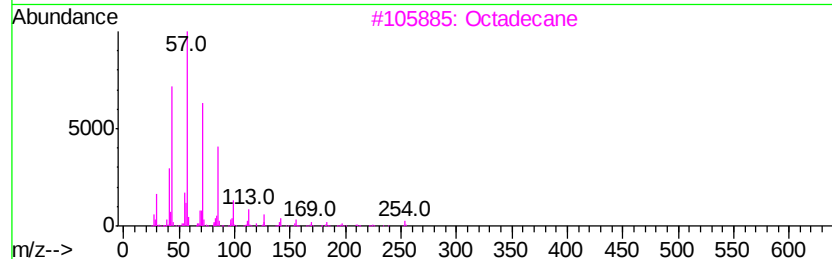
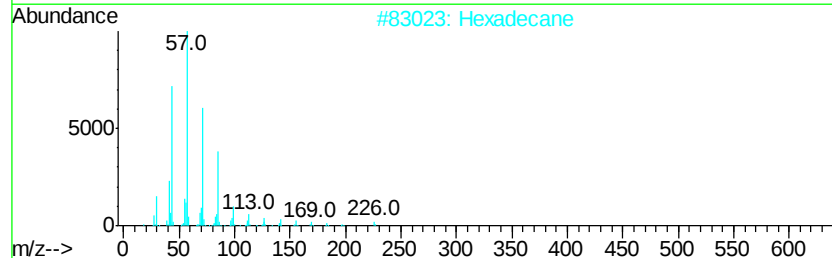
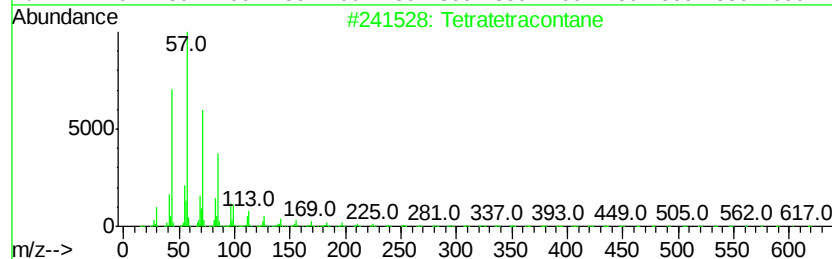
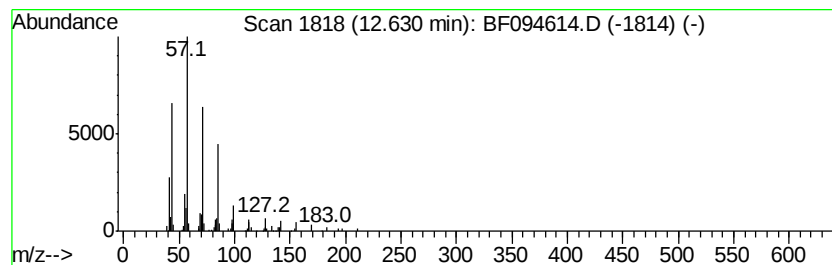
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Tetratetracontane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.63	2.33 ng	96519	Phenanthrene-d10	11.20

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetratetracontane	619	C44H90	007098-22-8	91
2			Hexadecane	226	C16H34	000544-76-3	91
3			Octadecane	254	C18H38	000593-45-3	91
4			Docosane	310	C22H46	000629-97-0	91
5			Octacosane	394	C28H58	000630-02-4	90



Data Path : Z:\HPCHEM1\BNA F\DATA\BF042517\
Data File : BF094614.D
Acq On : 25 Apr 2017 19:20
Operator : SJ/MA
Sample : I2824-02
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
E2-CONCRETE-BATCH7

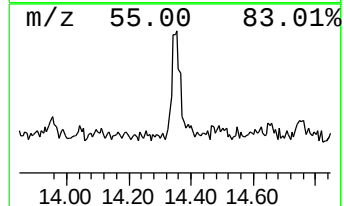
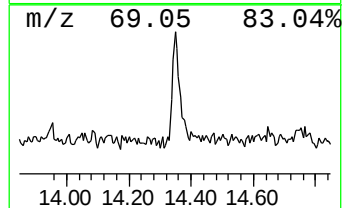
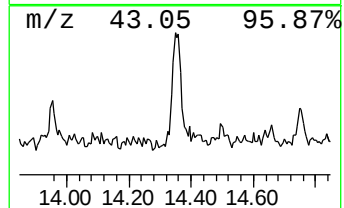
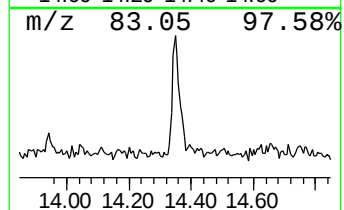
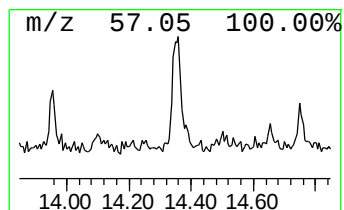
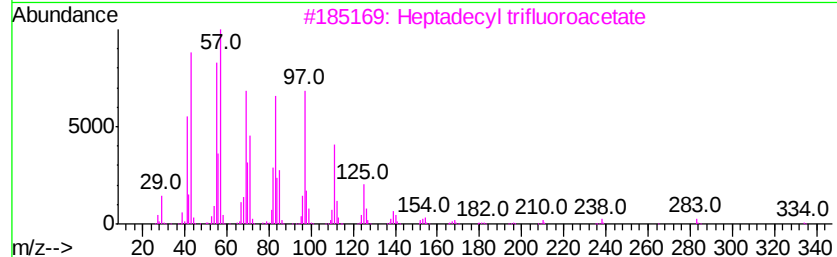
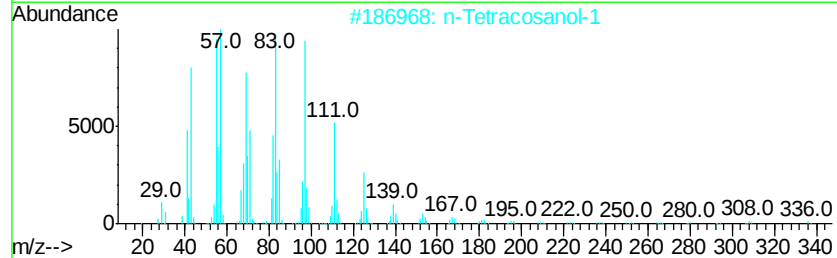
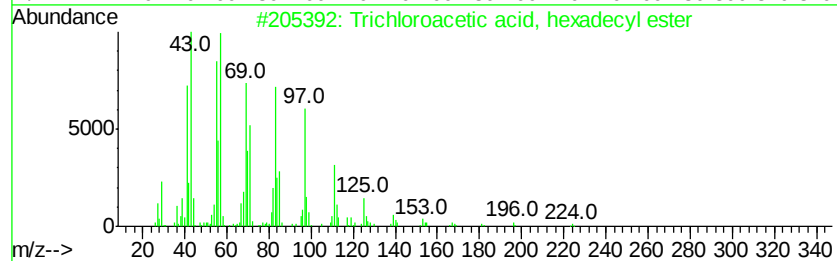
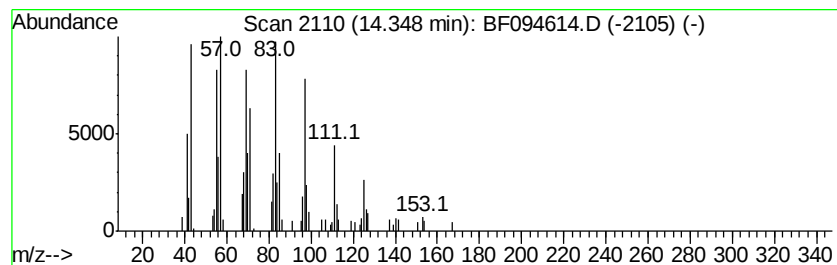
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Trichloroacetic acid, hexadecyl... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.35	4.49 ng	150164	Chrysene-d12	14.45

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	91
2			n-Tetracosanol-1	354	C24H50O	000506-51-4	91
3			Heptadecyl trifluoroacetate	352	C19H35F3O2	1000351-87-0	90
4			Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	87
5			Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	86



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042517\
 Data File : BF094614.D
 Acq On : 25 Apr 2017 19:20
 Operator : SJ/MA
 Sample : I2824-02
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 E2-CONCRETE-BATCH7

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF041717.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
4-Penten-2-one, 4...	2.70	3.6	ng	124771	1	5.75	683994	20.0
3-Penten-2-one, 4...	3.37	165.6	ng	5663480	1	5.75	683994	20.0
2-Pentanone, 4-hy...	3.94	216.8	ng	7414550	1	5.75	683994	20.0
unknown5.50	5.50	3.0	ng	102150	1	5.75	683994	20.0
Phenol, 2-ethyl-	7.02	4.2	ng	204808	2	7.24	981942	20.0
Benzene, 1-methox...	7.04	2.1	ng	102355	2	7.24	981942	20.0
Hexadecane	9.98	4.2	ng	239571	3	9.38	1141820	20.0
Pentadecane, 2,6,...	10.25	2.1	ng	121546	3	9.38	1141820	20.0
Cyclopentaneaceti...	10.30	2.0	ng	83796	4	11.20	827570	20.0
Sulfurous acid, 2...	10.58	4.0	ng	166075	4	11.20	827570	20.0
Heptadecane, 8-me...	11.11	7.0	ng	291891	4	11.20	827570	20.0
Octadecane, 1-iodo-	12.15	3.6	ng	148883	4	11.20	827570	20.0
Tetratetracontane	12.63	2.3	ng	96519	4	11.20	827570	20.0
Trichloroacetic a...	14.35	4.5	ng	150164	5	14.45	669234	20.0