

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF110517\
 Data File : BF100219.D
 Acq On : 5 Nov 2017 13:10
 Operator : SJ/JU
 Sample : I6275-01
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 DRUM(1-2)

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-BF102317.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	3.675	266	270	281	rBV	56628	124758	5.08%	0.638%
2	4.987	490	493	524	rBV	166366	346684	14.11%	1.772%
3	5.393	559	562	597	rBV	1662112	2226802	90.65%	11.384%
4	6.422	733	737	754	rBV	1772599	2148691	87.47%	10.985%
5	6.539	754	757	771	rVB	2254018	2294066	93.39%	11.728%
6	6.763	792	795	804	rBV	536365	509186	20.73%	2.603%
7	6.828	804	806	816	rVV	86201	140135	5.70%	0.716%
8	6.916	818	821	836	rVB	1930766	1838678	74.85%	9.400%
9	7.339	889	893	910	rBV	1318460	1360229	55.38%	6.954%
10	8.045	1010	1013	1031	rVB	706984	684355	27.86%	3.499%
11	9.116	1191	1195	1198	rBV	2827903	2456383	100.00%	12.558%
12	9.463	1252	1254	1261	rBV	247029	415401	16.91%	2.124%
13	9.516	1261	1263	1274	rVB	129935	175983	7.16%	0.900%
14	9.798	1307	1311	1317	rBV	711845	651699	26.53%	3.332%
15	9.892	1323	1327	1332	rBV3	35897	41676	1.70%	0.213%
16	10.098	1359	1362	1364	rBV3	30045	37253	1.52%	0.190%
17	10.516	1430	1433	1436	rBV	46955	42550	1.73%	0.218%
18	10.592	1442	1446	1451	rBV	1201077	1157724	47.13%	5.919%
19	10.692	1459	1463	1466	rBV	85964	97028	3.95%	0.496%
20	11.122	1530	1536	1539	rBV2	39145	77271	3.15%	0.395%
21	11.157	1539	1542	1549	rVB	59985	81270	3.31%	0.415%
22	11.286	1561	1564	1572	rBV	546182	589708	24.01%	3.015%
23	11.657	1624	1627	1631	rVB3	87576	74718	3.04%	0.382%
24	11.722	1635	1638	1642	rVV4	53302	83224	3.39%	0.425%
25	11.798	1649	1651	1655	rBV4	47173	67632	2.75%	0.346%
26	11.969	1677	1680	1682	rBV	112368	115496	4.70%	0.590%
27	12.116	1703	1705	1707	rBV	98273	91042	3.71%	0.465%
28	12.245	1724	1727	1729	rBV	174860	190478	7.75%	0.974%
29	12.698	1802	1804	1807	rBV2	419505	426415	17.36%	2.180%
30	12.874	1831	1834	1837	rVB	1116617	1014421	41.30%	5.186%

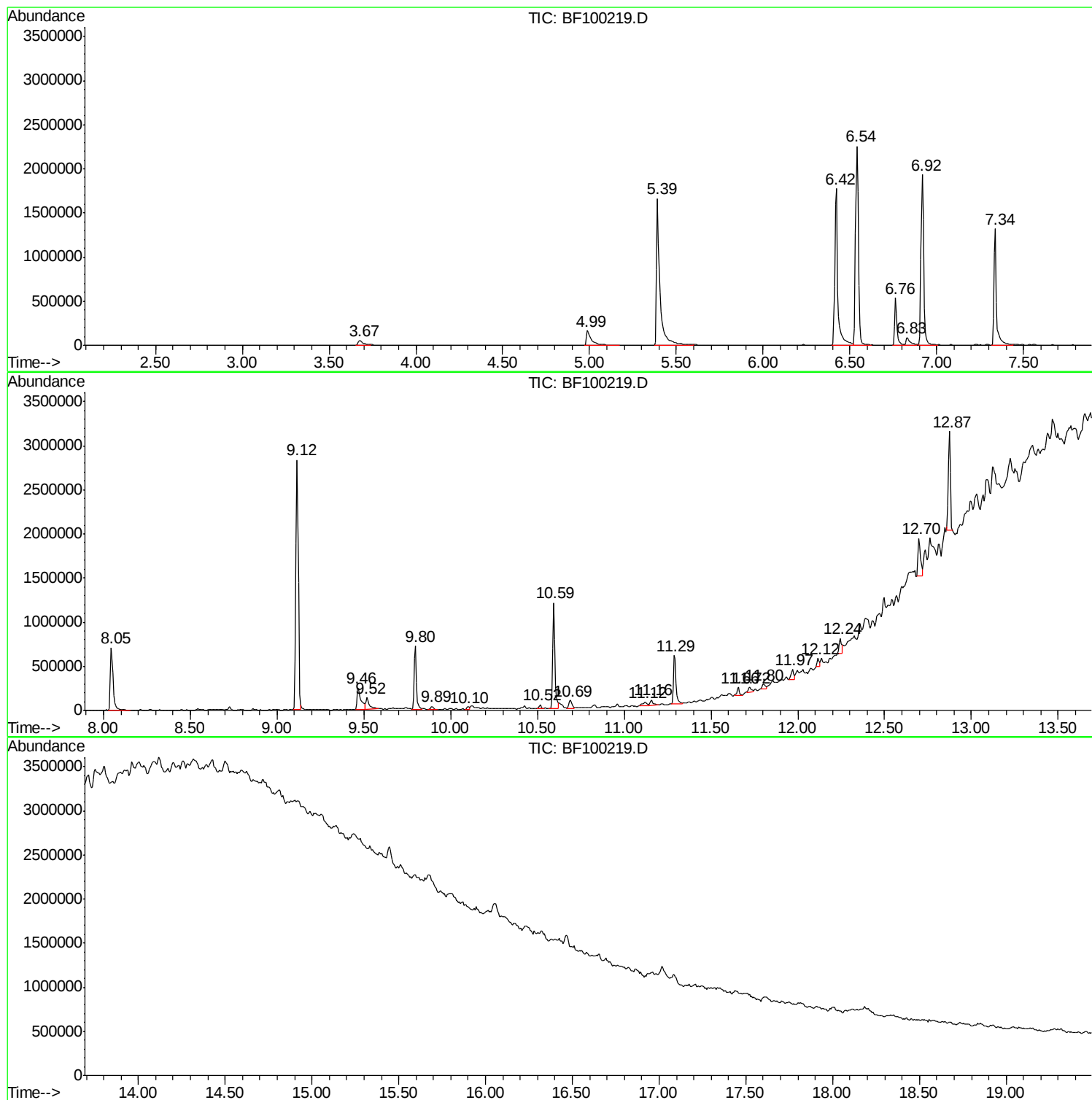
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Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
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ClientSampleId :
DRUM(1-2)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF102317.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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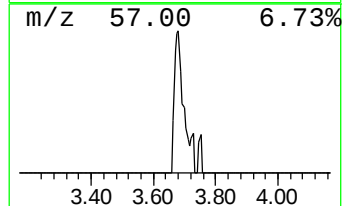
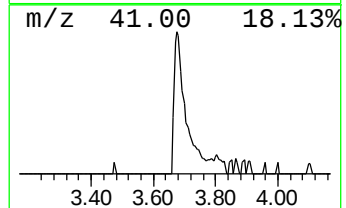
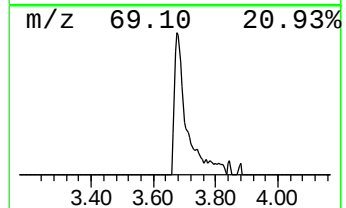
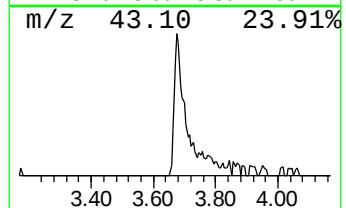
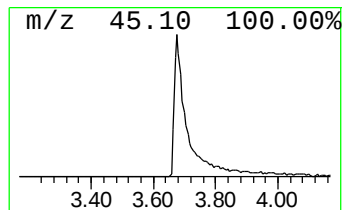
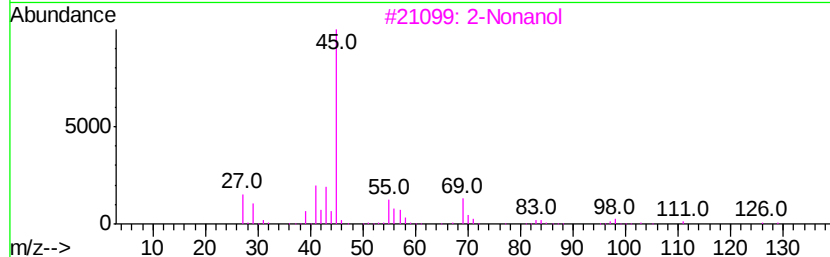
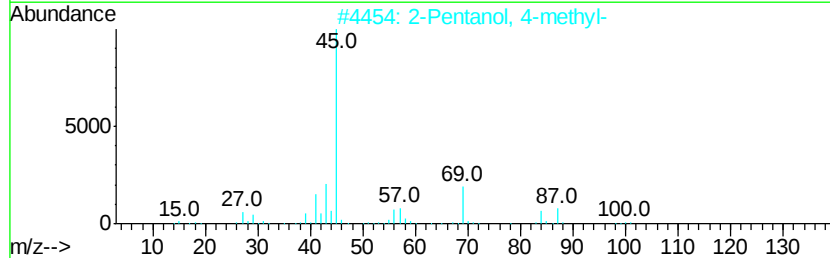
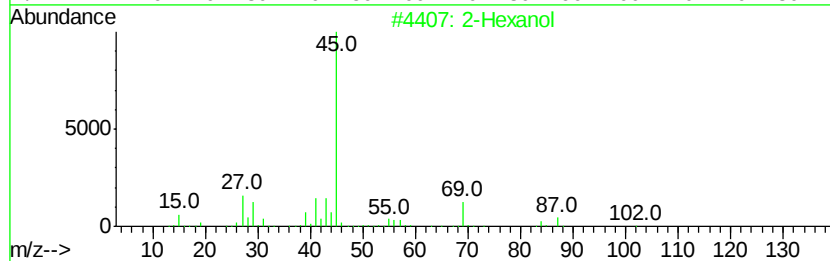
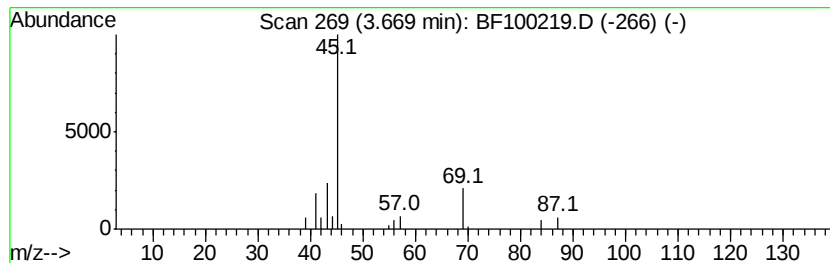
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF102317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Hexanol Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.67	4.90 ng	124758	1,4-Dichlorobenzene-d4	6.76

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Hexanol	102	C6H14O	000626-93-7	83
2			2-Pentanol, 4-methyl-	102	C6H14O	000108-11-2	83
3			2-Nonanol	144	C9H20O	000628-99-9	64
4			2-Pentanol, 3-methyl-	102	C6H14O	000565-60-6	40
5			2-Heptanol	116	C7H16O	000543-49-7	40



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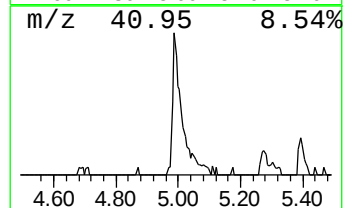
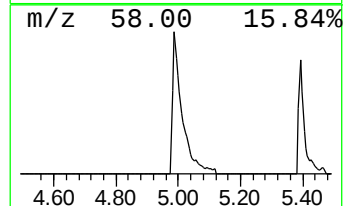
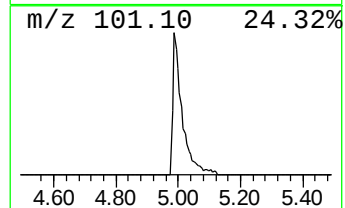
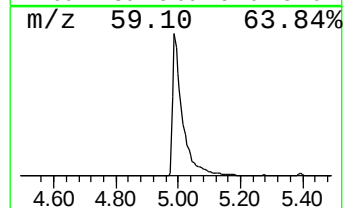
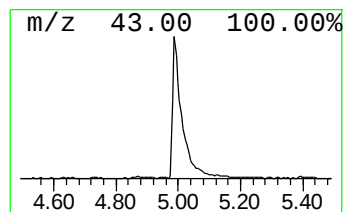
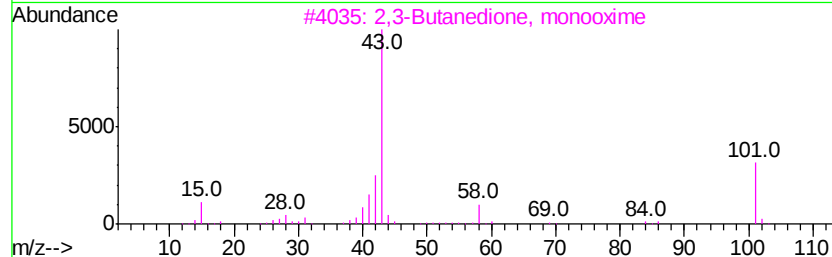
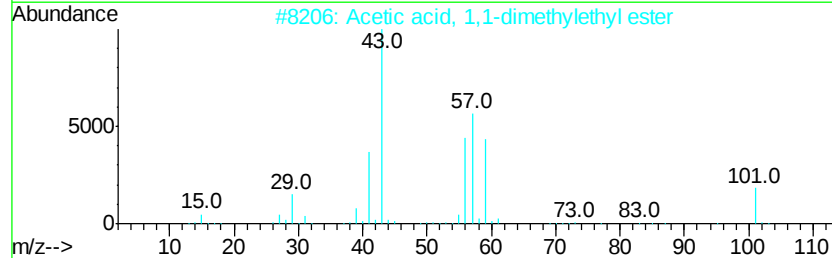
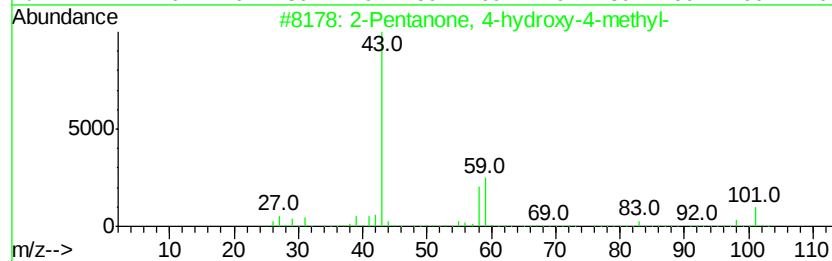
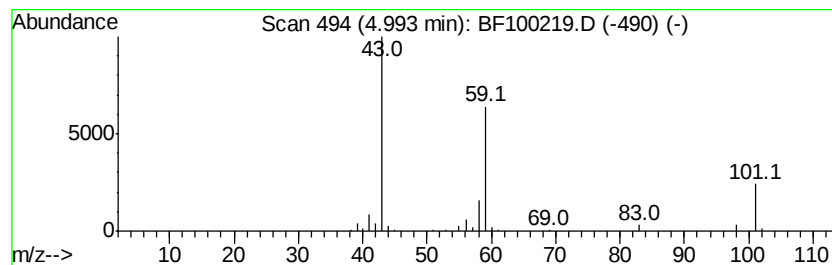
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.99	13.62 ng	346684	1,4-Dichlorobenzene-d4	6.76

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	17
4		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17
5		5-Hexen-2-one	98	C6H10O	000109-49-9	9



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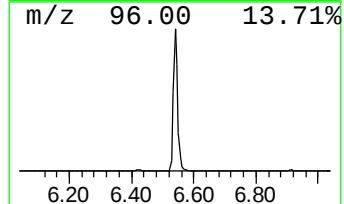
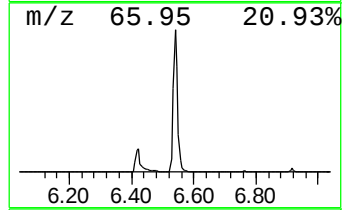
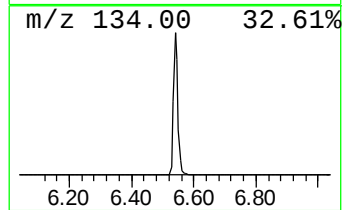
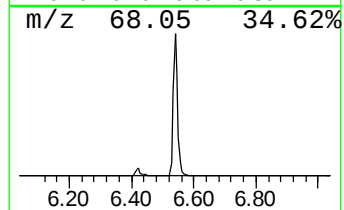
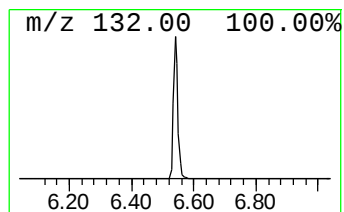
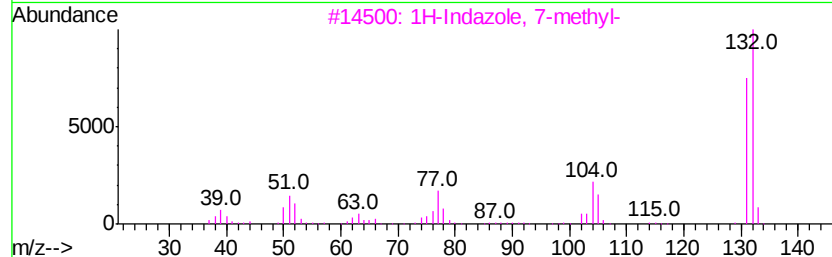
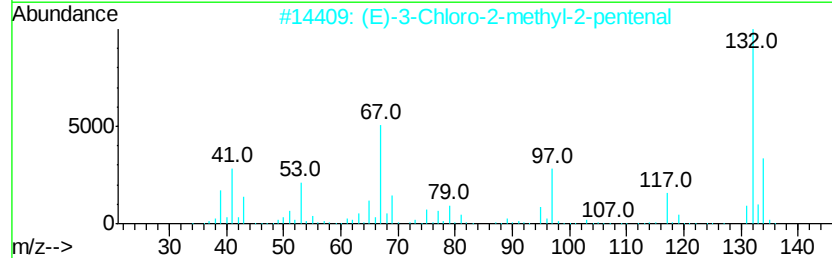
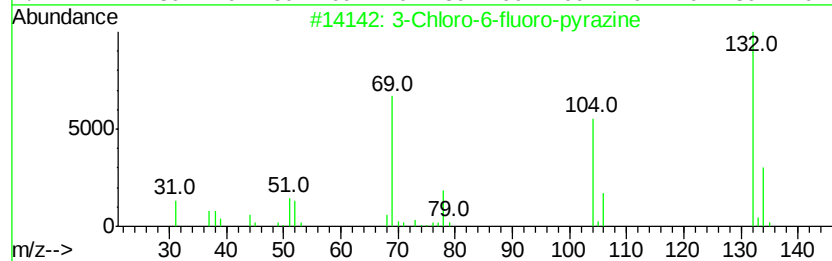
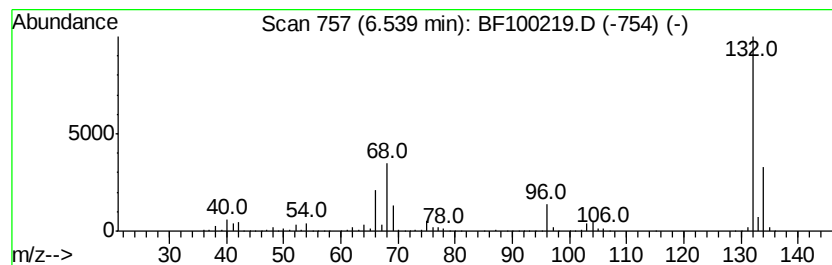
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown6.54 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.54	90.11 ng	2294070	1,4-Dichlorobenzene-d4	6.76

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	38
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	25
3		1H-Indazole, 7-methyl-	132	C8H8N2	1000316-00-7	10
4		Xenon	132	Xe	007440-63-3	9
5		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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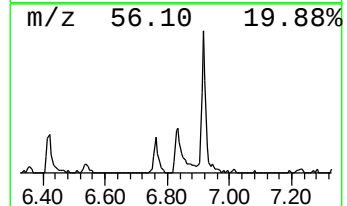
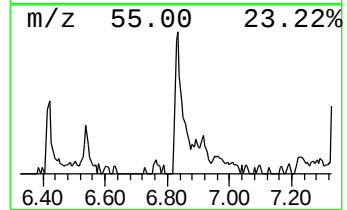
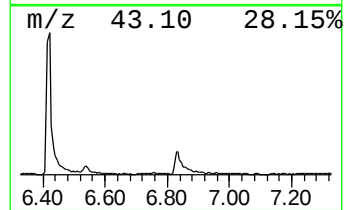
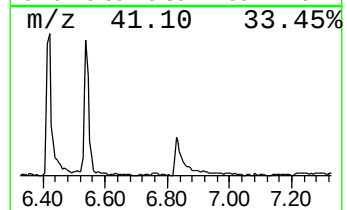
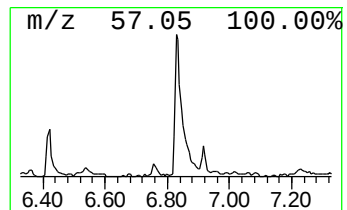
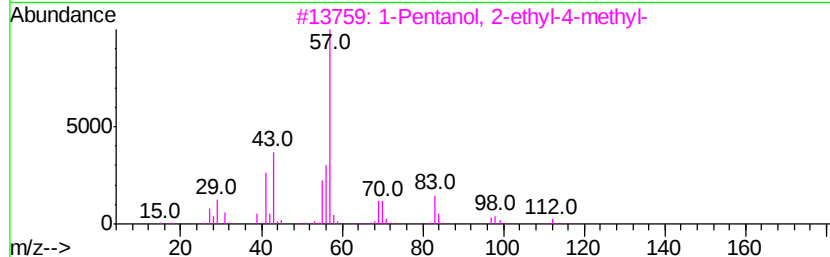
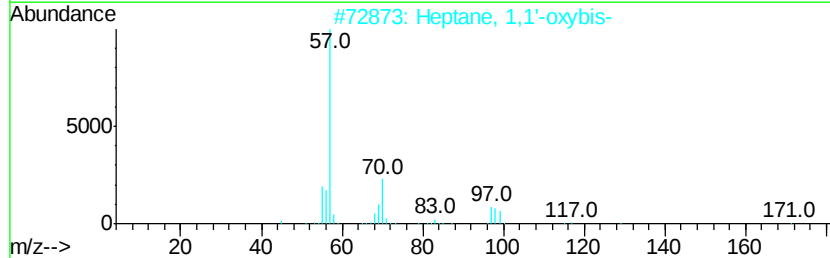
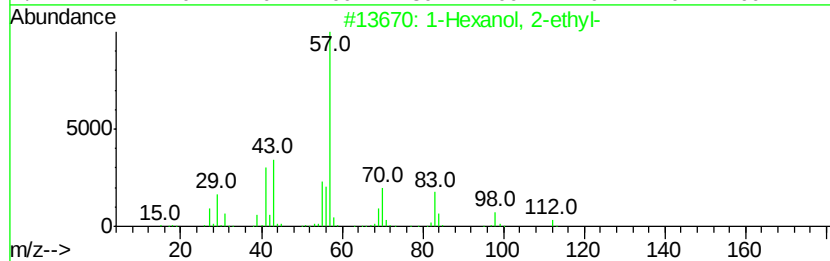
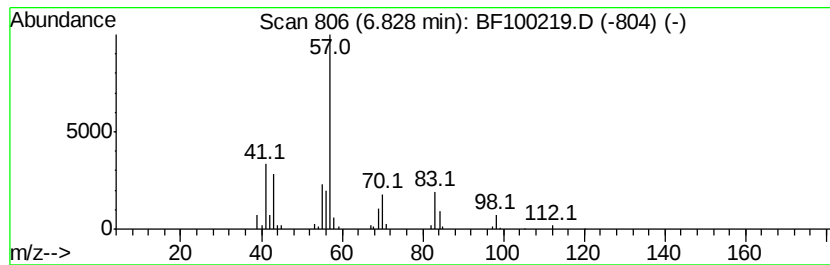
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 1-Hexanol, 2-ethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.83	5.50 ng	140135	1,4-Dichlorobenzene-d4	6.76

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	83
2			Heptane, 1,1'-oxybis-	214	C14H30O	000629-64-1	50
3			1-Pentanol, 2-ethyl-4-methyl-	130	C8H18O	000106-67-2	50
4			Decane, 2,5,6-trimethyl-	184	C13H28	062108-23-0	45
5			Oxalic acid, isobutyl heptyl ester	244	C13H24O4	1000309-37-2	43



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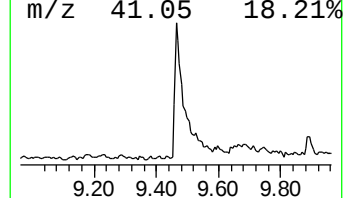
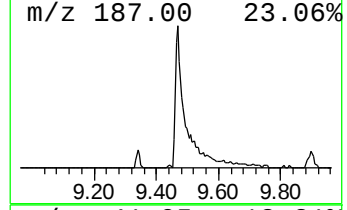
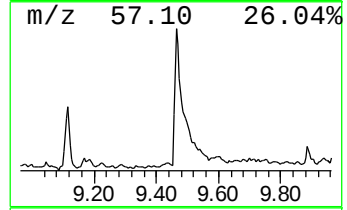
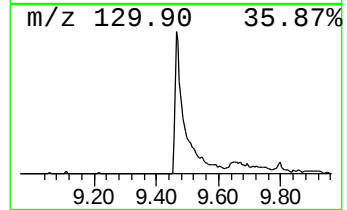
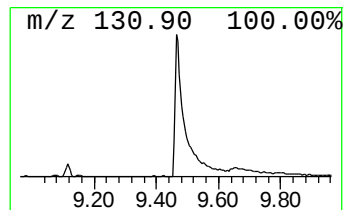
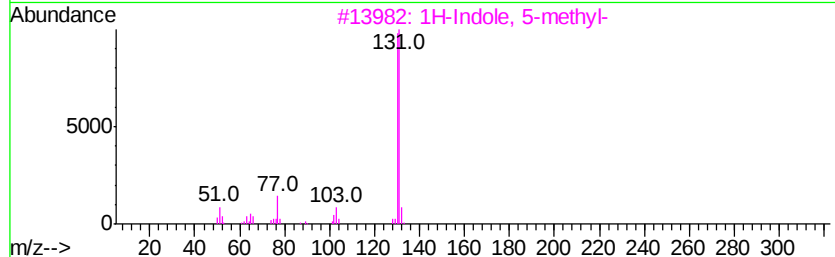
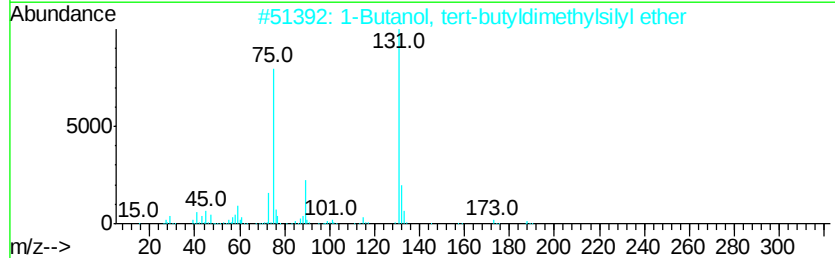
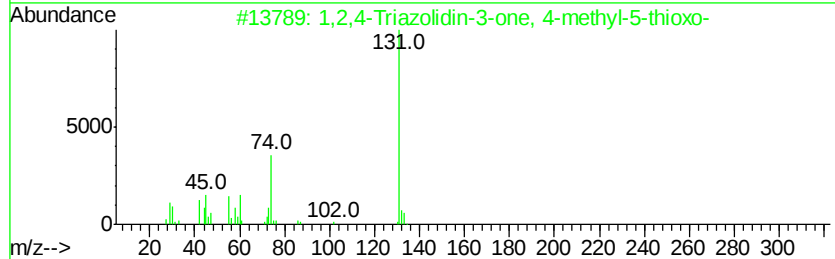
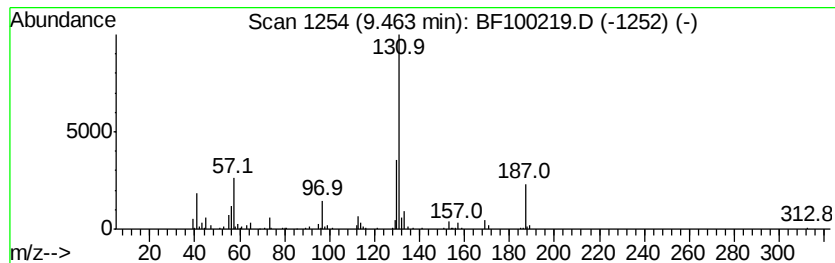
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 unknown9.46 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.46	12.75 ng	415401	Acenaphthene-d10	9.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,2,4-Triazolidin-3-one, 4-methy...	131	C3H5N3OS	022244-61-7	9
2		1-Butanol, tert-butyldimethylsil...	188	C10H24OSi	1000333-04-7	9
3		1H-Indole, 5-methyl-	131	C9H9N	000614-96-0	9
4		2-Oxazolidinethione, 4,4-dimethyl-	131	C5H9NOS	054013-55-7	9
5		2,4(3H,5H)-Thiazoledione, 5-methyl-	131	C4H5N02S	1000319-17-7	9



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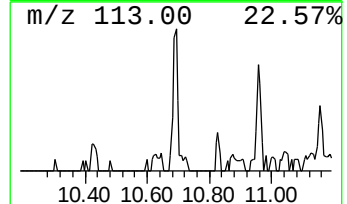
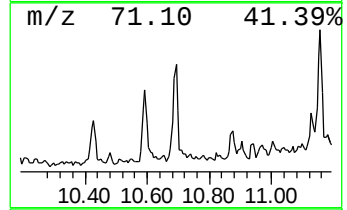
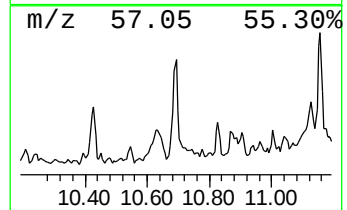
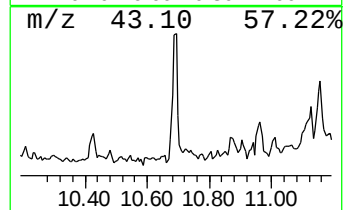
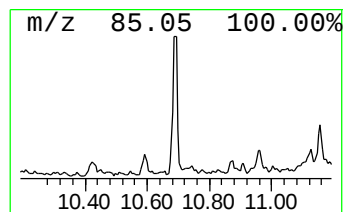
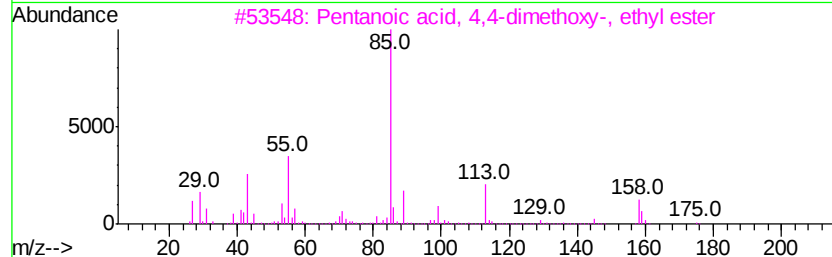
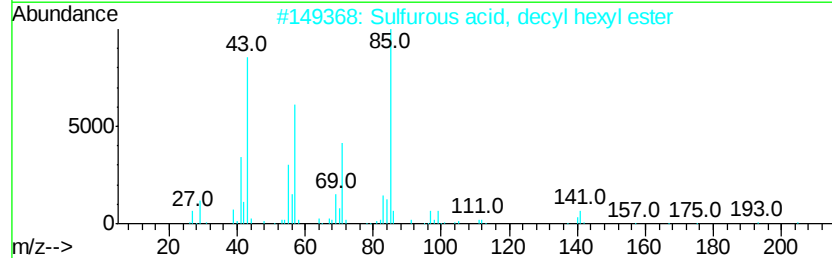
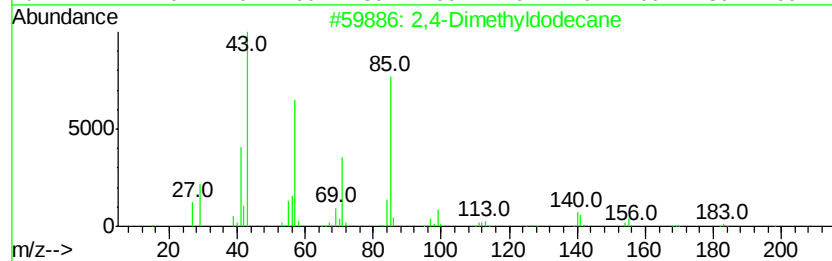
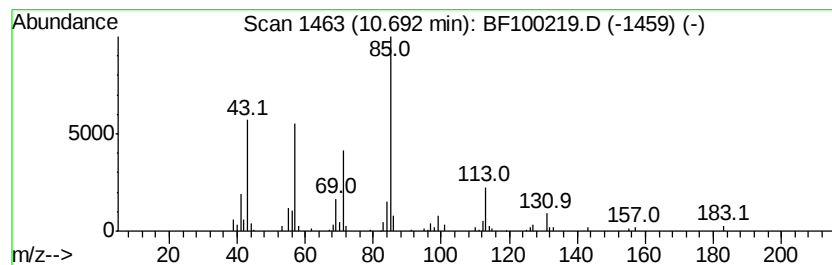
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BNA_F
ClientSampleId :
DRUM(1-2)

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF102317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 2,4-Dimethyldodecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.69	3.29 ng	97028	Phenanthrene-d10	11.29
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	2,4-Dimethyldodecane	198 C14H30	006117-99-3	59
2	Sulfurous acid, decyl hexyl ester	306 C16H34O3S	1000309-13-2	59
3	Pentanoic acid, 4,4-dimethoxy-, ...	190 C9H18O4	1000197-28-8	47
4	Heptane, 4,4-dimethyl-	128 C9H20	001068-19-5	43
5	2(3H)-Furanone, 5-heptyldihydro-	184 C11H20O2	000104-67-6	43



Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA F\DATA\BF110517\
Data File : BF100219.D
Acq On : 5 Nov 2017 13:10
Operator : SJ/JU
Sample : I6275-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
DRUM(1-2)

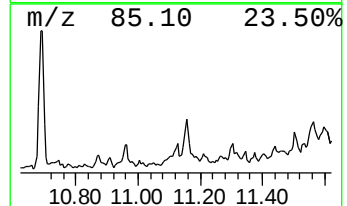
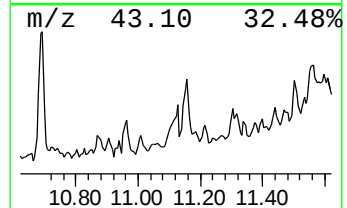
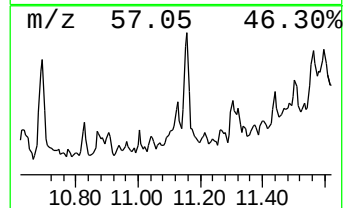
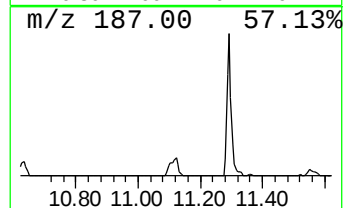
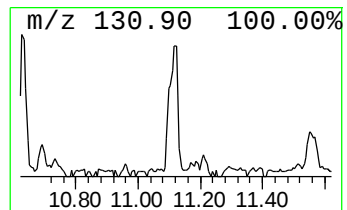
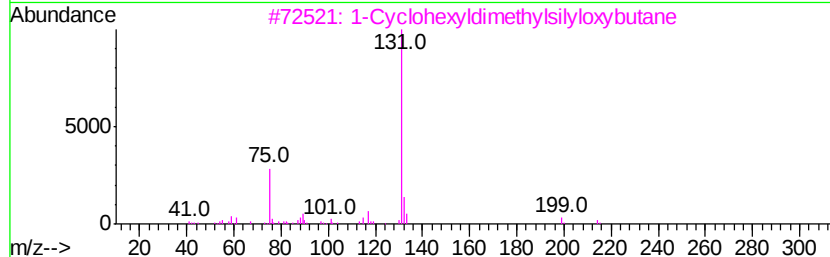
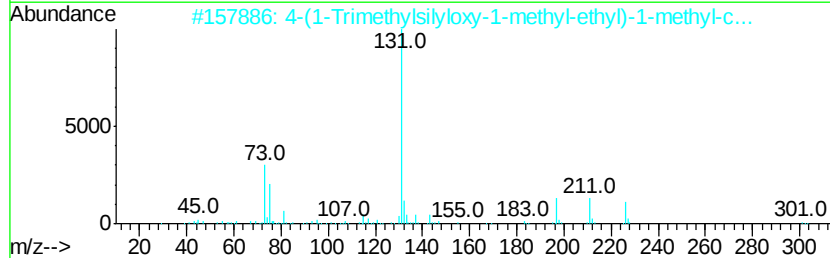
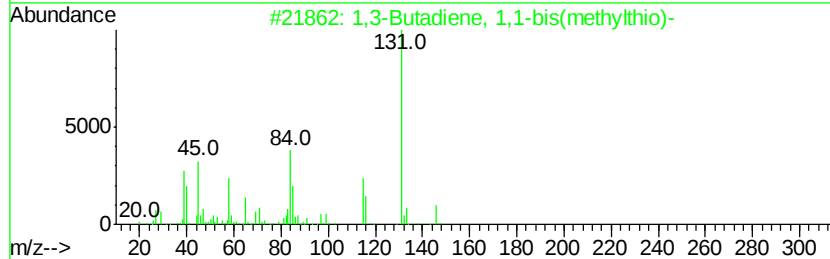
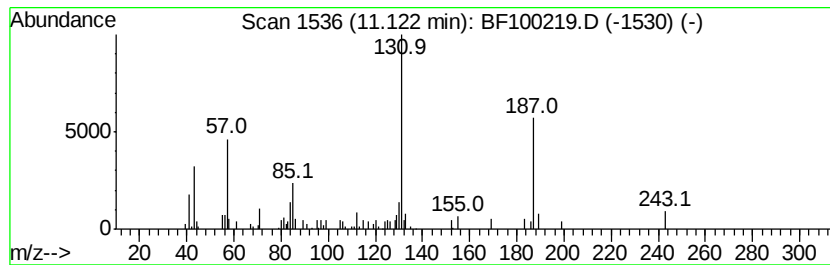
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 unknown11.12 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.12	2.62 ng	77271	Phenanthrene-d10	11.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,3-Butadiene, 1,1-bis(methylthio)-	146	C6H10S2	066612-70-2	25
2		4-(1-Trimethylsilyloxy-1-methyl-...	316	C16H36O2Si2	1000373-06-9	17
3		1-Cyclohexyldimethylsilyloxybutane	214	C12H26OSi	1000281-95-6	17
4		1H-Pyrrole-2,5-dione, 1-(4-methy...	187	C11H9NO2	001631-28-3	14
5		2-Propenoic acid, 3-(1H-indol-3-...	187	C11H9NO2	029953-71-7	14



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF110517\
Data File : BF100219.D
Acq On : 5 Nov 2017 13:10
Operator : SJ/JU
Sample : I6275-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

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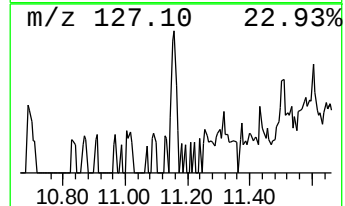
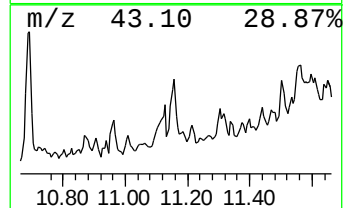
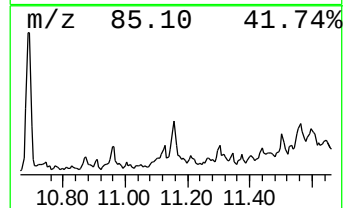
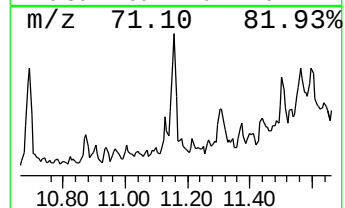
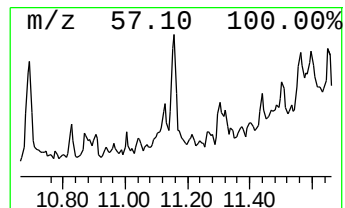
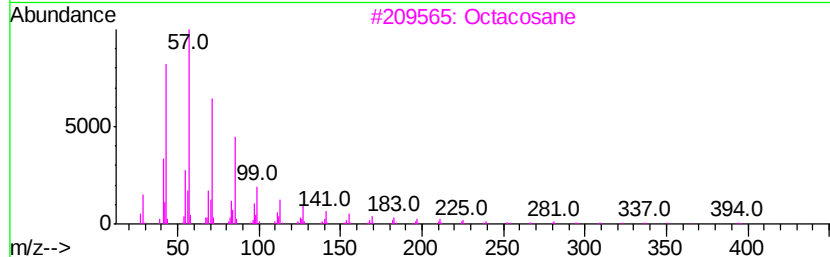
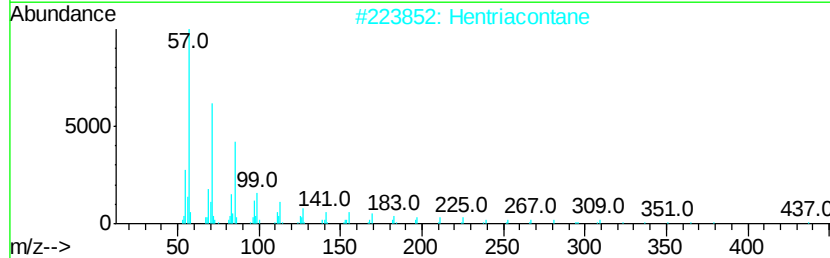
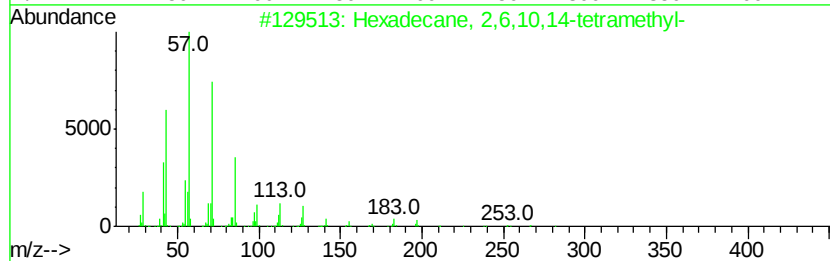
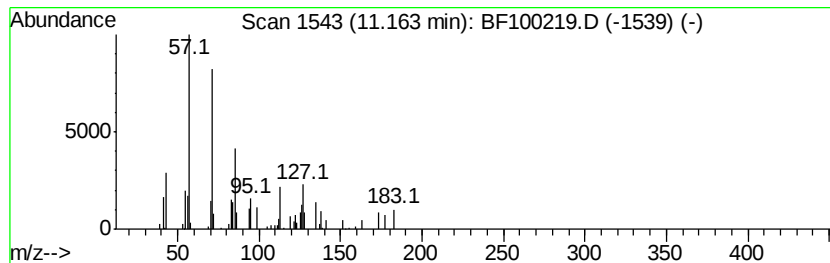
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Hexadecane, 2,6,10,14-tetra... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.16	2.76 ng	81270	Phenanthrene-d10	11.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	90
2		Hentriacontane	437	C31H64	000630-04-6	83
3		Octacosane	394	C28H58	000630-02-4	78
4		Hexacosane	366	C26H54	000630-01-3	78
5		Dodecane, 2-methyl-8-propyl-	226	C16H34	055045-07-3	78



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF110517\
Data File : BF100219.D
Acq On : 5 Nov 2017 13:10
Operator : SJ/JU
Sample : I6275-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
DRUM(1-2)

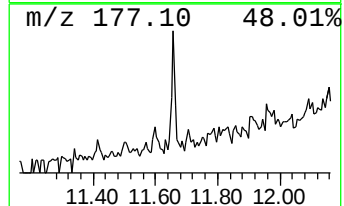
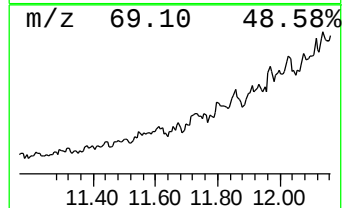
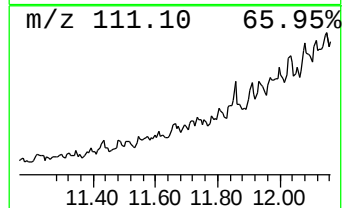
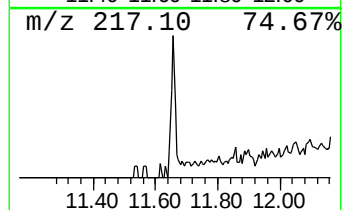
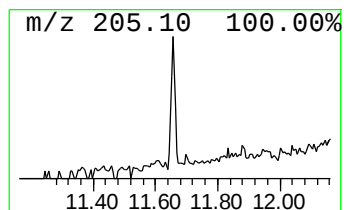
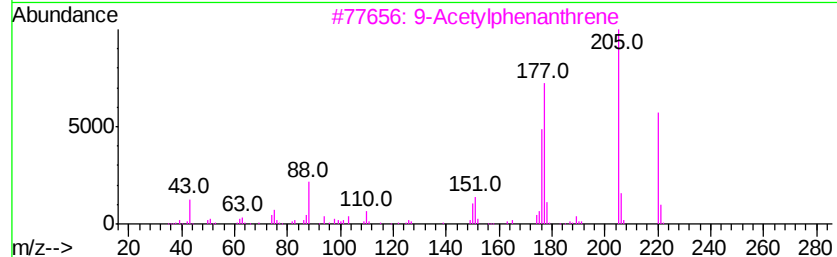
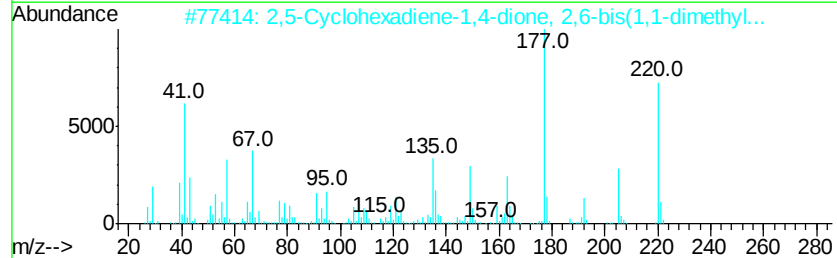
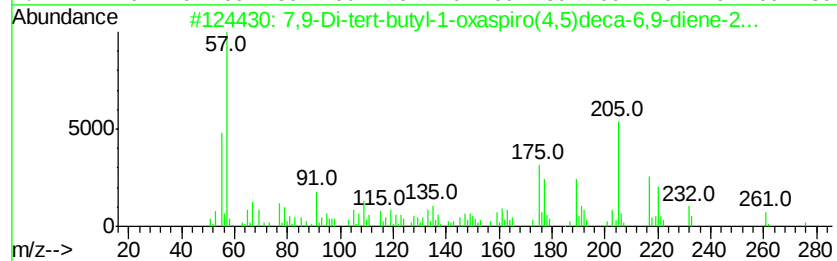
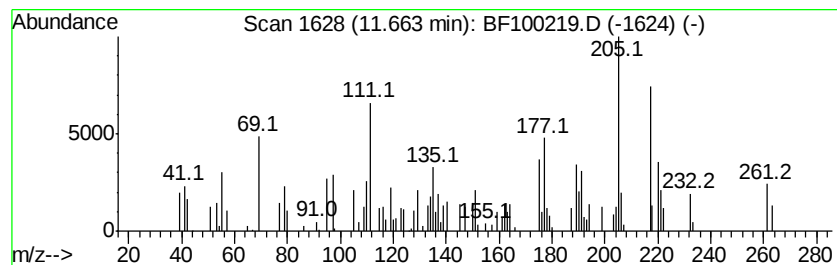
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 7,9-Di-tert-butyl-1-oxaspiro... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.66	2.53 ng	74718	Phenanthrene-d10	11.29

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			7,9-Di-tert-butyl-1-oxaspiro(4,5...	276	C17H24O3	082304-66-3	83
2			2,5-Cyclohexadiene-1,4-dione, 2,...	220	C14H20O2	000719-22-2	46
3			9-Acetylphenanthrene	220	C16H12O	002039-77-2	43
4			Ethanone, 1-(5,6,7,8-tetrahydro-...	220	C14H20O2	071596-88-8	38
5			Ethanone, 1-(9-anthracenyl)-	220	C16H12O	000784-04-3	38



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF110517\
Data File : BF100219.D
Acq On : 5 Nov 2017 13:10
Operator : SJ/JU
Sample : I6275-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

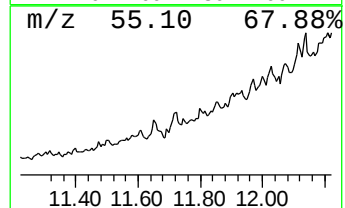
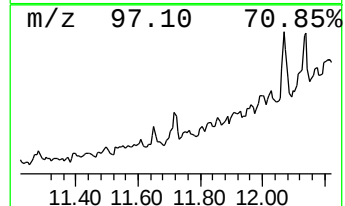
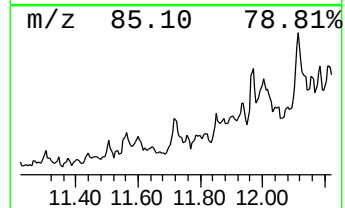
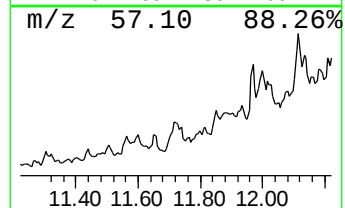
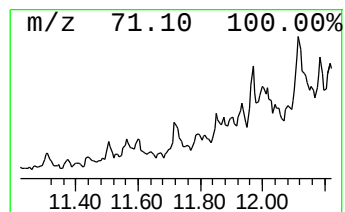
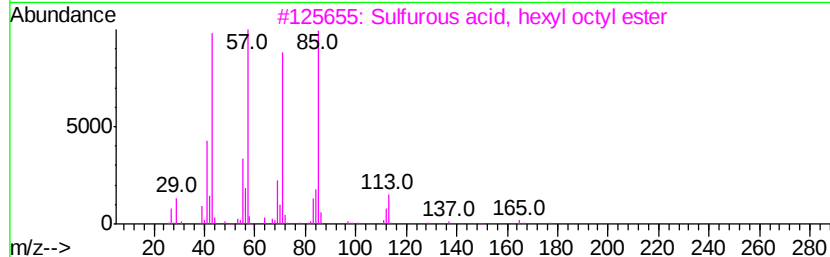
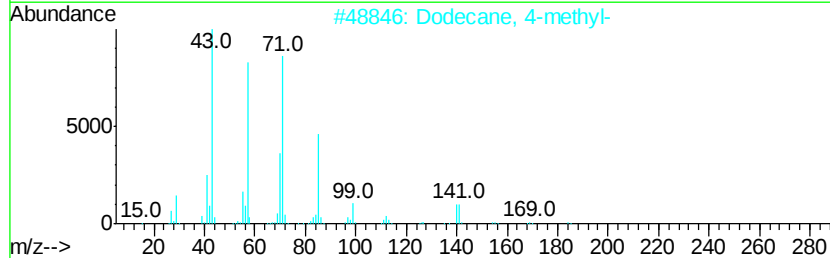
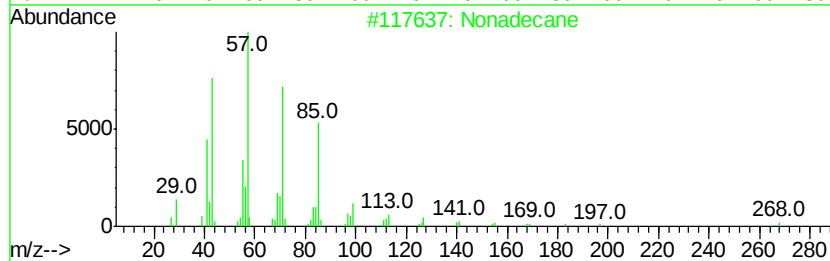
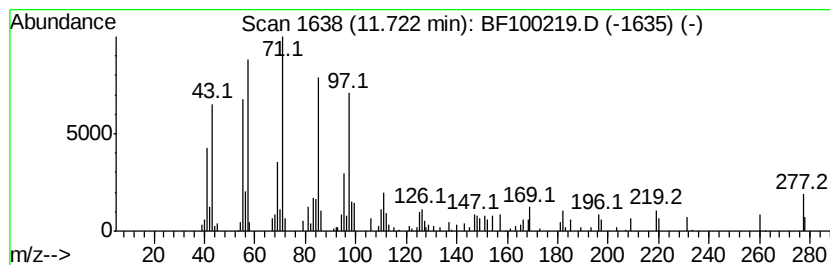
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF102317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 unknown11.72 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.72	2.82 ng	83224	Phenanthrene-d10	11.29	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonadecane	268	C19H40	000629-92-5	47
2	Dodecane, 4-methyl-	184	C13H28	006117-97-1	47
3	Sulfurous acid, hexyl octyl ester	278	C14H30O3S	1000309-13-0	43
4	Tetradecane	198	C14H30	000629-59-4	43
5	Oxalic acid, isohexyl neopentyl ...	244	C13H24O4	1000309-73-0	43



Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA F\DATA\BF110517\
Data File : BF100219.D
Acq On : 5 Nov 2017 13:10
Operator : SJ/JU
Sample : I6275-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
DRUM(1-2)

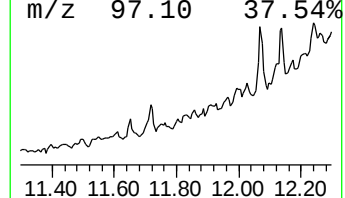
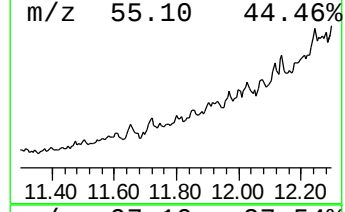
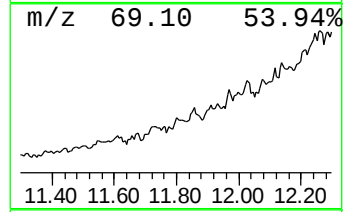
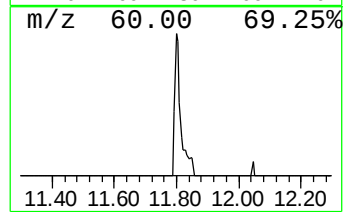
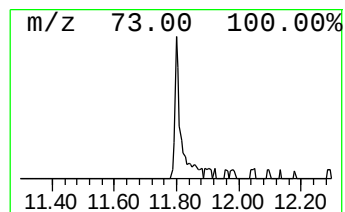
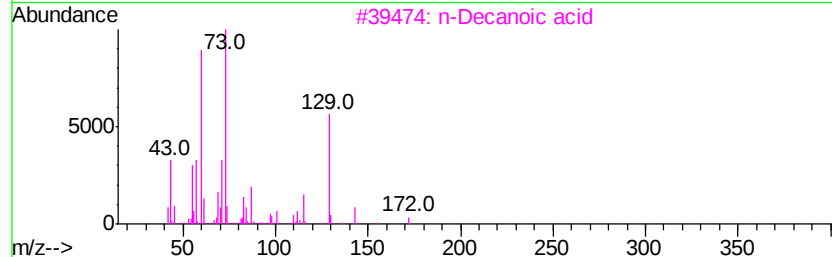
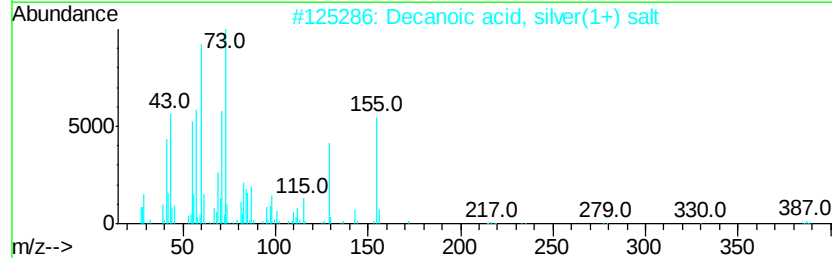
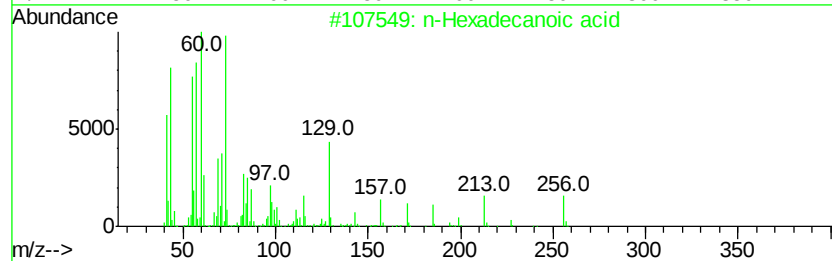
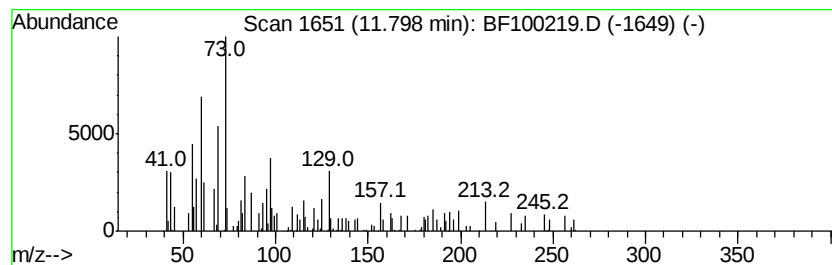
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 n-Hexadecanoic acid Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.80	2.29 ng	67632	Phenanthrene-d10	11.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	55
2		Decanoic acid, silver(1+) salt	278	C10H19AgO2	013126-67-5	47
3		n-Decanoic acid	172	C10H20O2	000334-48-5	47
4		Dodecanoic acid	200	C12H24O2	000143-07-7	43
5		.alpha.-D-Glucopyranose, 4-O-.be...	342	C12H22O11	014641-93-1	43



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF110517\
Data File : BF100219.D
Acq On : 5 Nov 2017 13:10
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ALS Vial : 5 Sample Multiplier: 1

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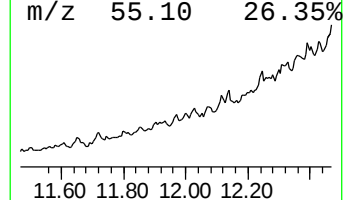
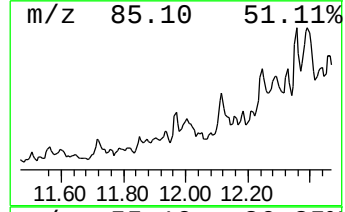
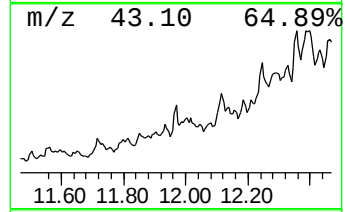
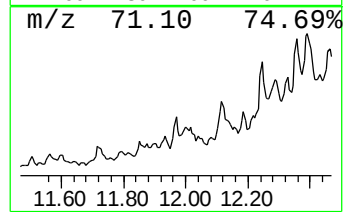
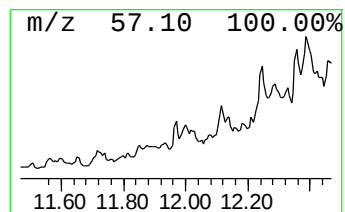
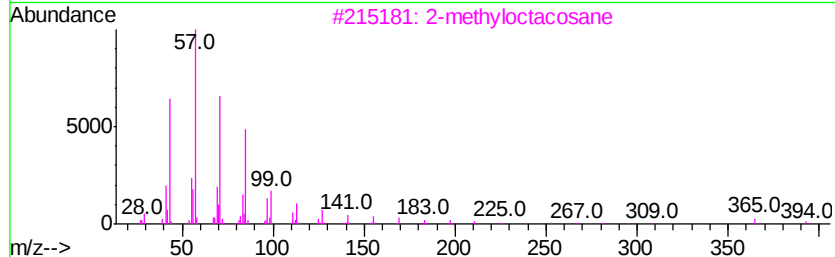
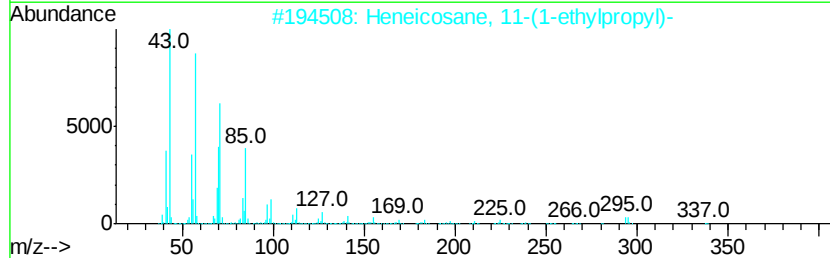
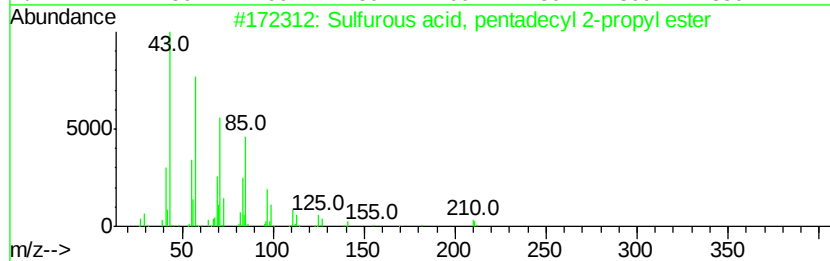
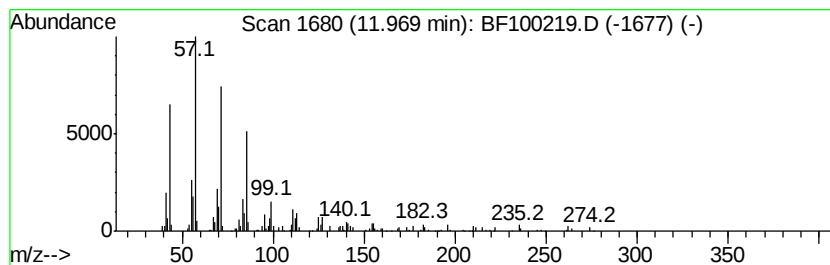
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Sulfurous acid, pentadecyl ... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.97	3.92 ng	115496	Phenanthrene-d10	11.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Sulfurous acid, pentadecyl 2-pro...	334	C18H38O3S	1000309-12-6	90
2		Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	90
3		2-methyloctacosane	408	C29H60	1000376-72-8	90
4		Hexacosane	366	C26H54	000630-01-3	87
5		Nonadecane	268	C19H40	000629-92-5	86



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Operator : SJ/JU
Sample : I6275-01
Misc :
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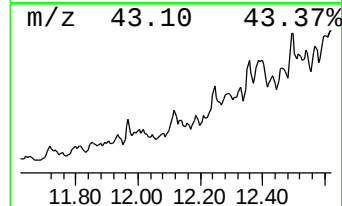
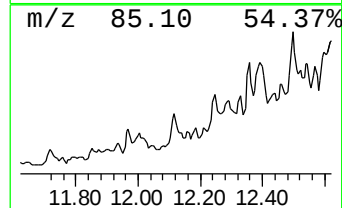
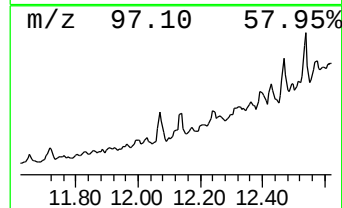
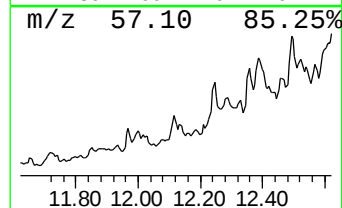
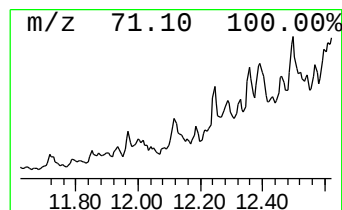
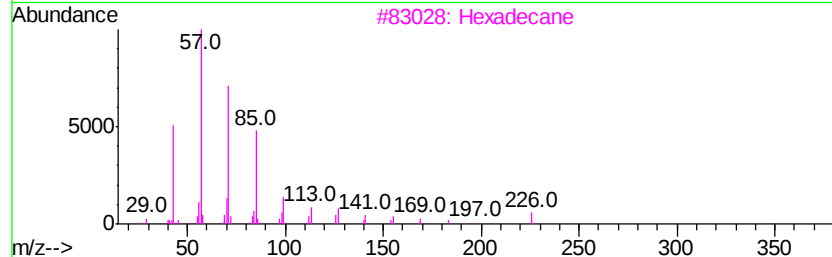
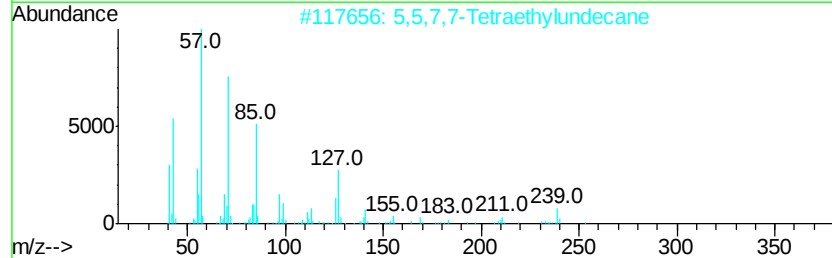
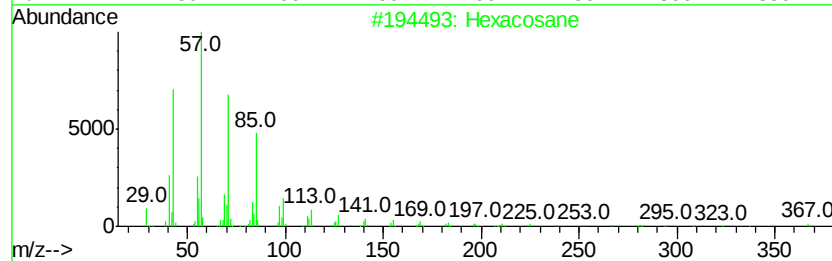
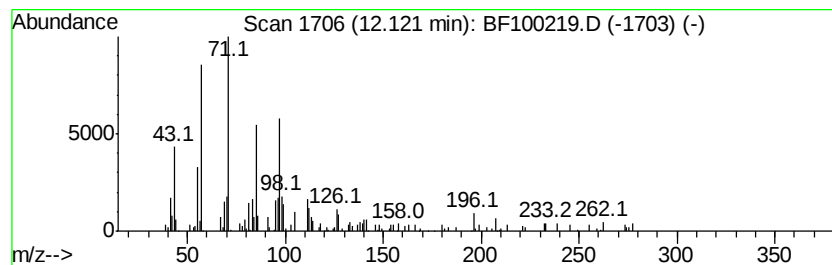
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ClientSampleId :
DRUM(1-2)

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF102317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Hexacosane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.12	3.09 ng	91042	Phenanthrene-d10	11.29
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Hexacosane		366 C26H54	000630-01-3 72
2	5,5,7,7-Tetraethylundecane		268 C19H40	1000360-42-0 64
3	Hexadecane		226 C16H34	000544-76-3 64
4	Heptadecane		240 C17H36	000629-78-7 64
5	Tridecane, 1-iodo-		310 C13H27I	035599-77-0 59



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF110517\
Data File : BF100219.D
Acq On : 5 Nov 2017 13:10
Operator : SJ/JU
Sample : I6275-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
DRUM(1-2)

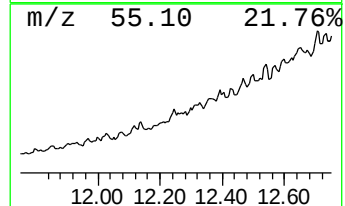
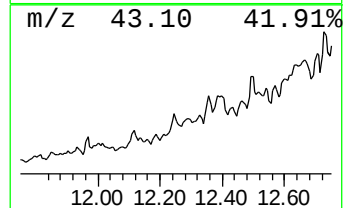
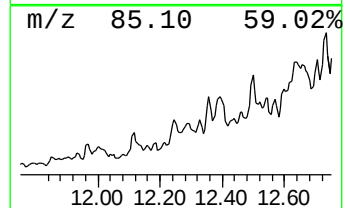
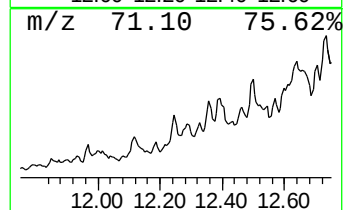
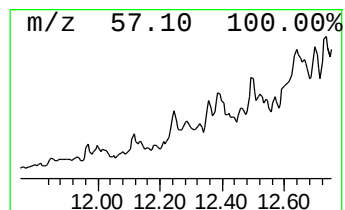
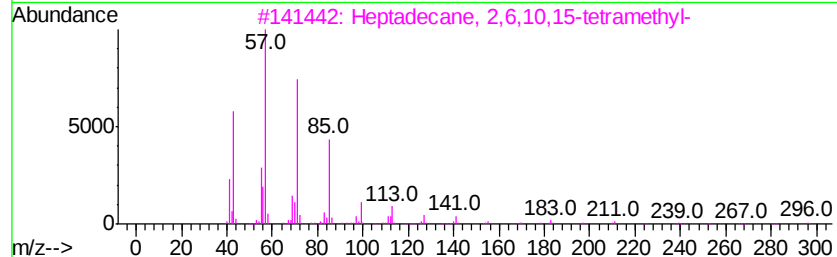
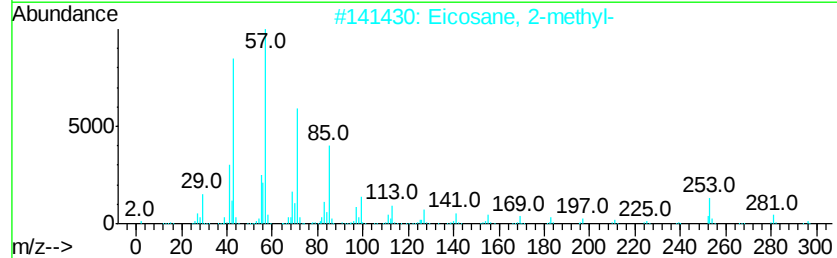
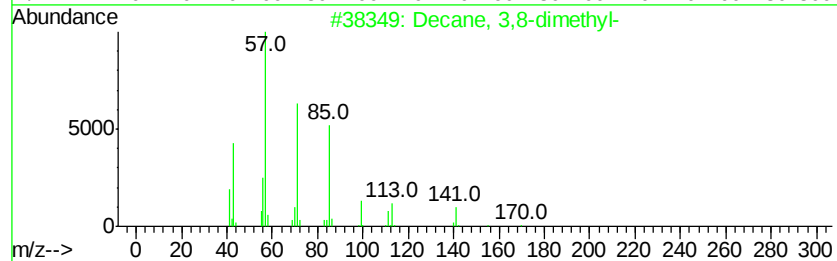
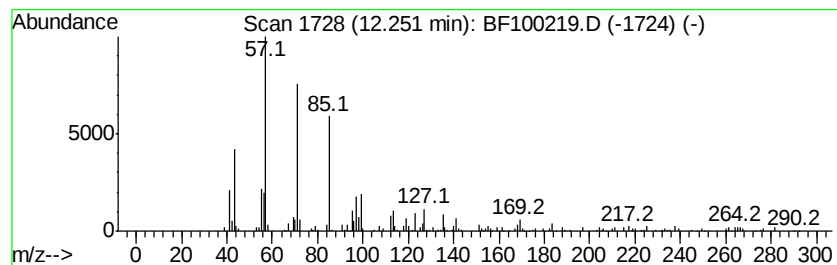
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Decane, 3,8-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.25	6.46 ng	190478	Phenanthrene-d10	11.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Decane, 3,8-dimethyl-	170	C12H26	017312-55-9	87
2		Eicosane, 2-methyl-	296	C21H44	001560-84-5	86
3		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	86
4		Hexacosane	366	C26H54	000630-01-3	83
5		Octadecane	254	C18H38	000593-45-3	80



Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA F\DATA\BF110517\
Data File : BF100219.D
Acq On : 5 Nov 2017 13:10
Operator : SJ/JU
Sample : I6275-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
DRUM(1-2)

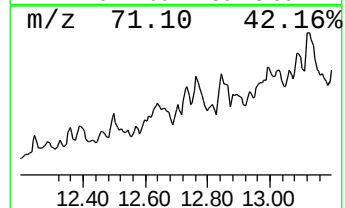
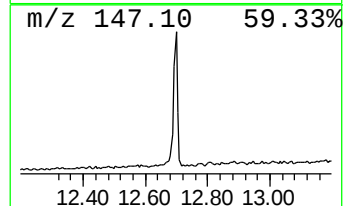
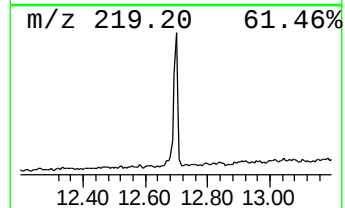
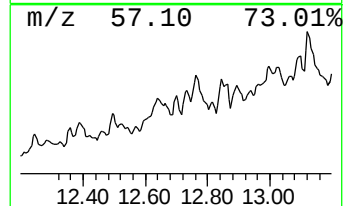
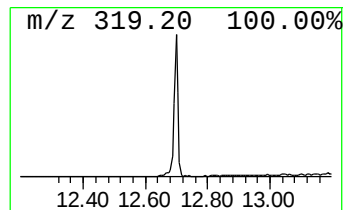
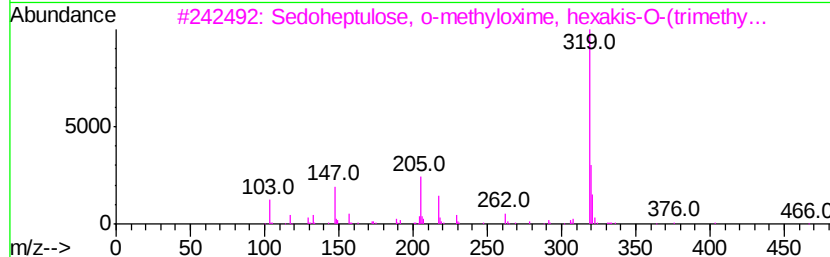
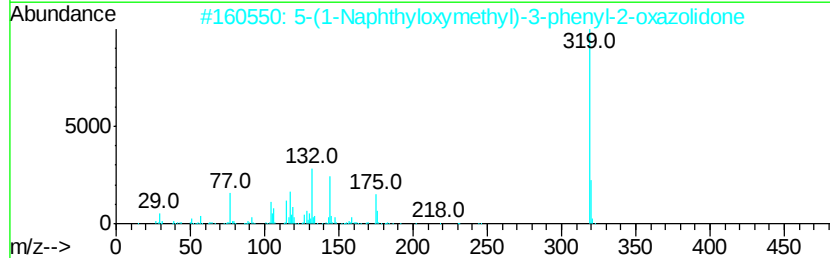
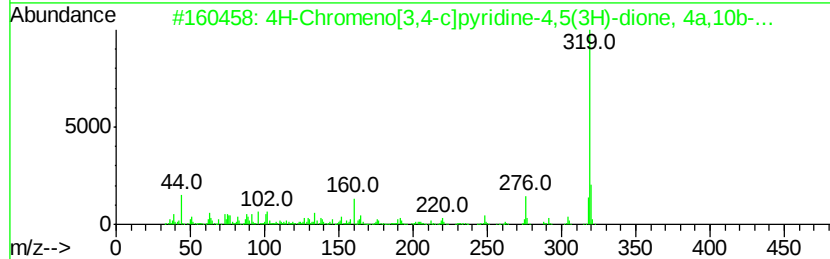
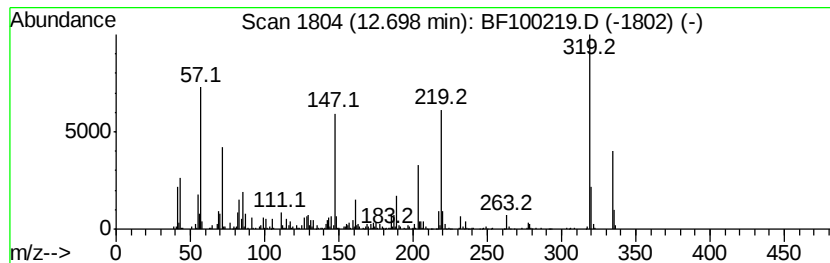
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF102317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 unknown12.70 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.70	8.41 ng	426415	Chrysene-d12	13.96

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4H-Chromeno[3,4-c]pyridine-4,5(3...	319	C19H13N04	029542-35-6	22
2		5-(1-Naphthyloxymethyl)-3-phenyl...	319	C20H17N03	005255-99-2	18
3		Sedoheptulose, o-methyloxime, he...	671	C26H65N07Si6	1000298-66-2	16
4		Methanone, (4-benzyl-1-piperidyl...	319	C22H25N0	1000271-12-8	14
5		1H,6H-5,11b-Ethano[1,3]dioxolo[4...	319	C17H21N05	003660-65-9	12



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF110517\
 Data File : BF100219.D
 Acq On : 5 Nov 2017 13:10
 Operator : SJ/JU
 Sample : I6275-01
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 DRUM(1-2)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF102317.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
2-Hexanol	3.67	4.9	ng	124758	1	6.76	509186	20.0
2-Pentanone, 4-hy...	4.99	13.6	ng	346684	1	6.76	509186	20.0
unknown6.54	6.54	90.1	ng	2294070	1	6.76	509186	20.0
1-Hexanol, 2-ethyl-	6.83	5.5	ng	140135	1	6.76	509186	20.0
unknown9.46	9.46	12.8	ng	415401	3	9.80	651699	20.0
2,4-Dimethyldodecane	10.69	3.3	ng	97028	4	11.29	589708	20.0
unknown11.12	11.12	2.6	ng	77271	4	11.29	589708	20.0
Hexadecane, 2,6,1...	11.16	2.8	ng	81270	4	11.29	589708	20.0
7,9-Di-tert-butyl...	11.66	2.5	ng	74718	4	11.29	589708	20.0
unknown11.72	11.72	2.8	ng	83224	4	11.29	589708	20.0
n-Hexadecanoic acid	11.80	2.3	ng	67632	4	11.29	589708	20.0
Sulfurous acid, p...	11.97	3.9	ng	115496	4	11.29	589708	20.0
Hexacosane	12.12	3.1	ng	91042	4	11.29	589708	20.0
Decane, 3,8-dimet...	12.25	6.5	ng	190478	4	11.29	589708	20.0
unknown12.70	12.70	8.4	ng	426415	5	13.96	1014420	20.0