

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF073018\
 Data File : BF107845.D
 Acq On : 31 Jul 2018 11:56
 Operator : JU/SJ
 Sample : J4221-06
 Misc :
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-A29

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:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF073018.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	5.190	482	485	500	rVB	216474	270272	4.74%	0.643%
2	5.601	551	555	580	rBV	1104008	1991350	34.92%	4.738%
3	6.607	722	726	744	rBV	512492	1354249	23.75%	3.222%
4	6.737	744	748	783	rVV	3254962	4559722	79.95%	10.849%
5	6.960	783	786	808	rVV	590453	1169723	20.51%	2.783%
6	7.119	808	813	841	rVV	3210037	4000620	70.15%	9.519%
7	7.531	879	883	904	rBV	1983860	3225076	56.55%	7.674%
8	7.678	905	908	921	rVB	132210	185623	3.25%	0.442%
9	8.242	1000	1004	1041	rBV	985750	1500210	26.31%	3.570%
10	9.319	1182	1187	1214	rBV	4218436	5703013	100.00%	13.570%
11	9.707	1249	1253	1264	rBV	150056	191363	3.36%	0.455%
12	10.001	1295	1303	1316	rBV	1364022	1535229	26.92%	3.653%
13	10.801	1434	1439	1451	rBV	2536202	3385741	59.37%	8.056%
14	11.501	1554	1558	1574	rBV	970092	1461855	25.63%	3.478%
15	12.001	1640	1643	1648	rBV	142183	176574	3.10%	0.420%
16	12.554	1734	1737	1743	rVB	84108	76355	1.34%	0.182%
17	12.930	1798	1801	1803	rBV	151601	125267	2.20%	0.298%
18	12.954	1803	1805	1811	rVB	156144	129521	2.27%	0.308%
19	13.083	1822	1827	1844	rBV	3807400	4569642	80.13%	10.873%
20	13.283	1858	1861	1865	rVB	210460	190555	3.34%	0.453%
21	13.542	1901	1905	1910	rBV5	53693	91309	1.60%	0.217%
22	13.630	1916	1920	1922	rBV	361146	330847	5.80%	0.787%
23	13.960	1972	1976	1984	rVB	476292	469978	8.24%	1.118%
24	14.154	2005	2009	2020	rBV	852301	1256682	22.04%	2.990%
25	14.277	2026	2030	2034	rBV	435627	411097	7.21%	0.978%
26	14.583	2078	2082	2088	rBV	470779	470398	8.25%	1.119%
27	14.901	2132	2136	2142	rBV	470707	476992	8.36%	1.135%
28	14.983	2147	2150	2154	rVB	70623	79449	1.39%	0.189%
29	15.254	2192	2196	2204	rBV	419878	497675	8.73%	1.184%
30	15.654	2260	2264	2267	rBV	348729	482506	8.46%	1.148%
31	15.689	2267	2270	2281	rVB	594384	983401	17.24%	2.340%
32	16.118	2338	2343	2350	rVB	247341	363934	6.38%	0.866%
33	16.654	2429	2434	2442	rVB	175494	311445	5.46%	0.741%

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Misc :
ALS Vial : 44 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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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

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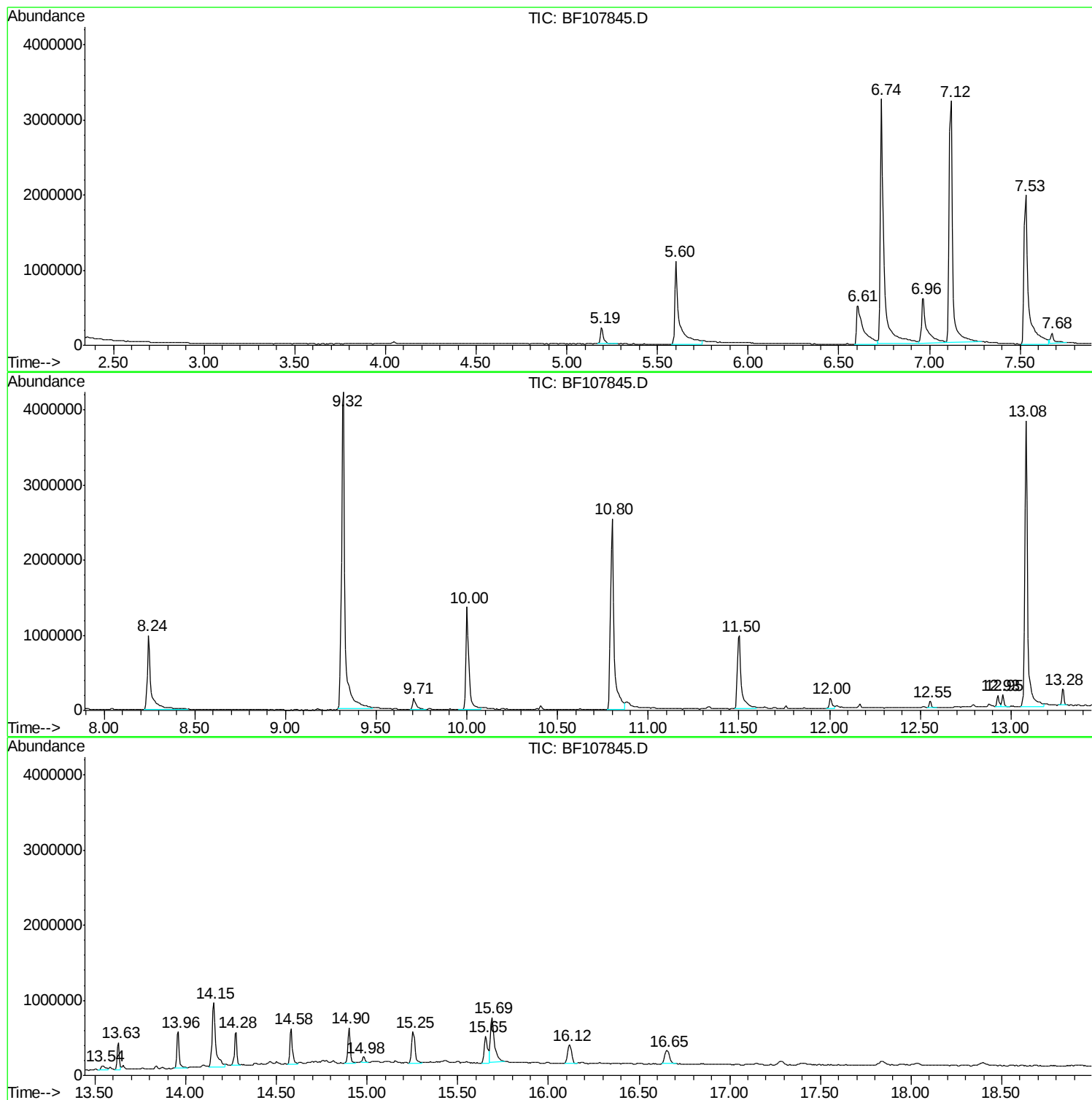
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073018.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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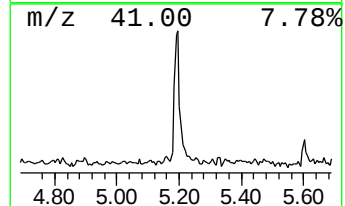
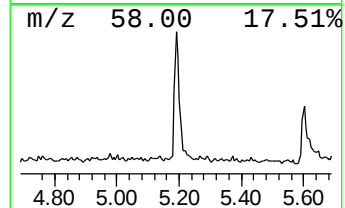
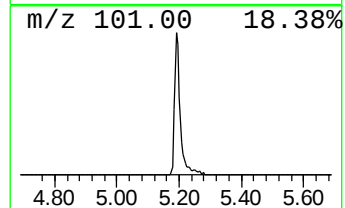
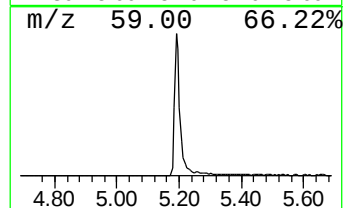
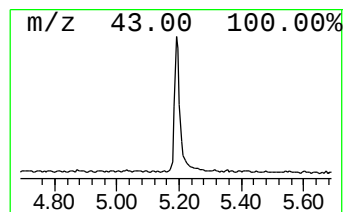
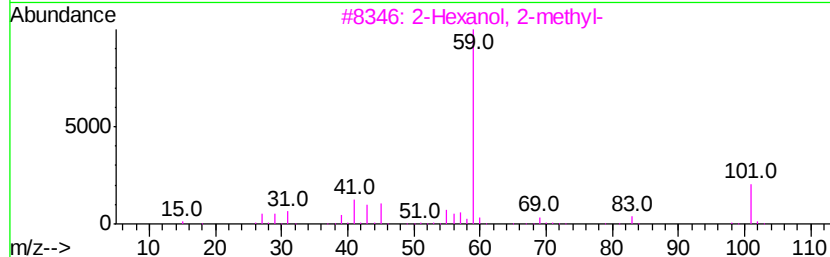
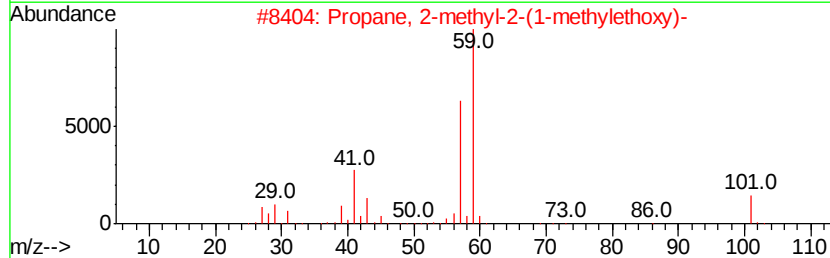
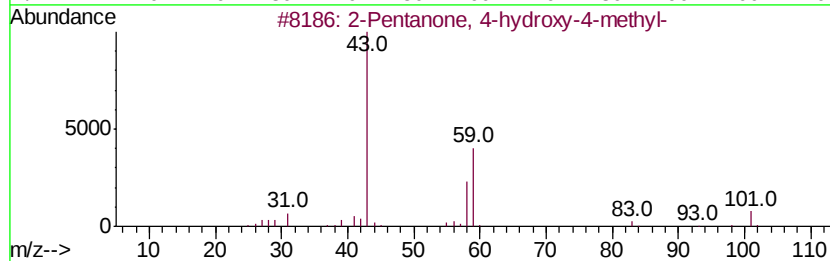
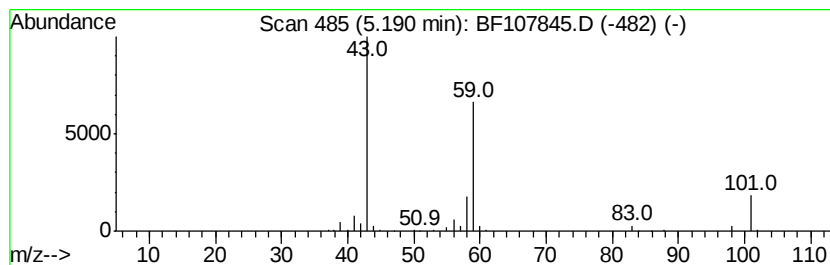
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.19	4.62 ng	270272	1,4-Dichlorobenzene-d4	6.97

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	53
3			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	38
4			3-Octanol	130	C8H18O	000589-98-0	33
5			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33



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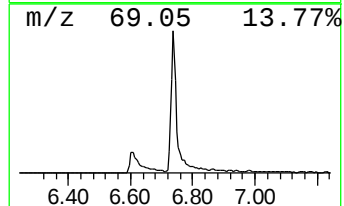
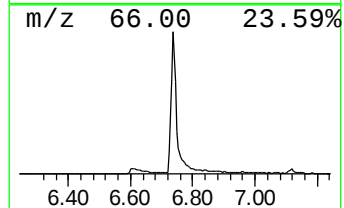
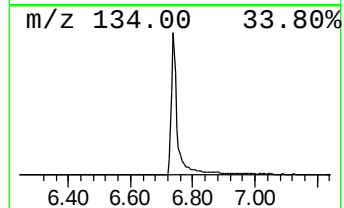
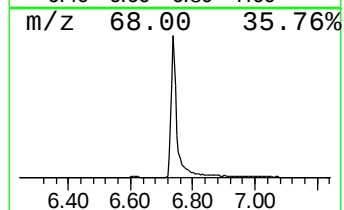
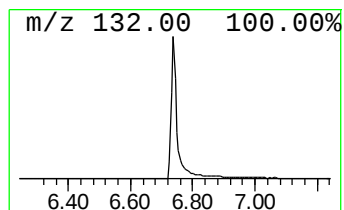
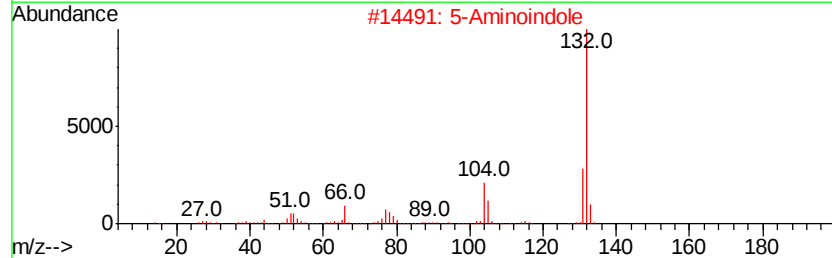
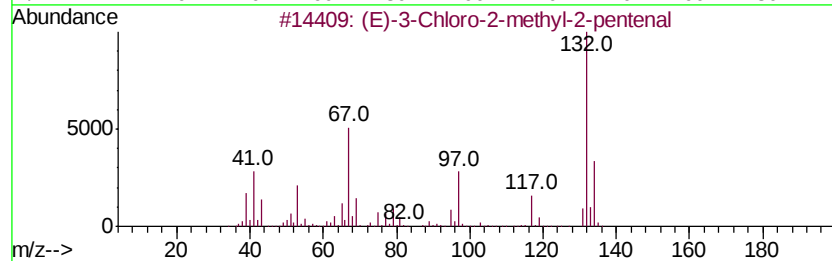
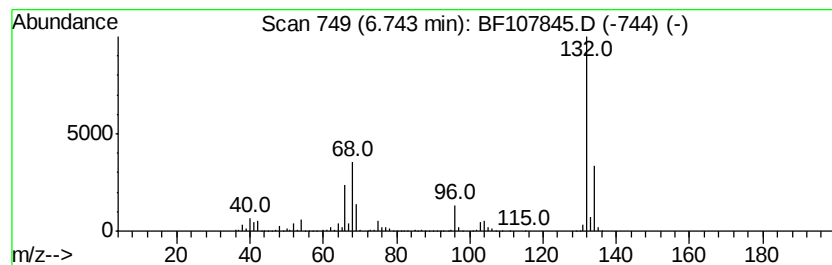
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown6.74 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.74	77.96 ng	4559720	1,4-Dichlorobenzene-d4	6.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	27
2		5-Aminoindole	132	C8H8N2	005192-03-0	12
3		Tranvlcyproline-propionyl	189	C12H15NO	1000123-86-3	10
4		2H-Cyclopentadipyridazine, 2-me...	132	C8H8N2	022291-85-6	9
5		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9



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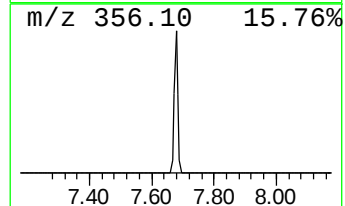
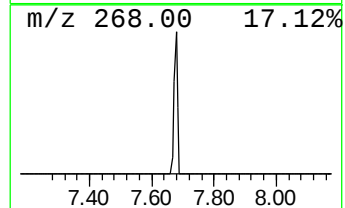
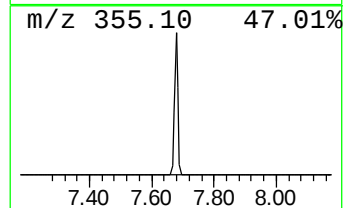
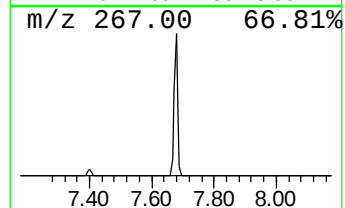
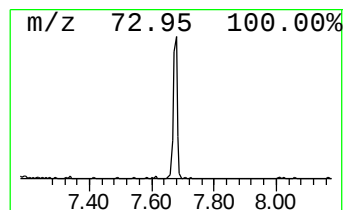
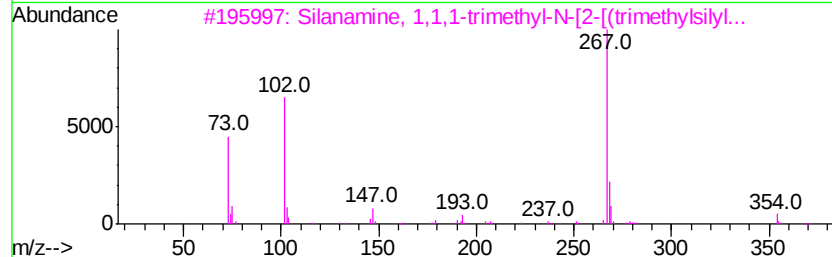
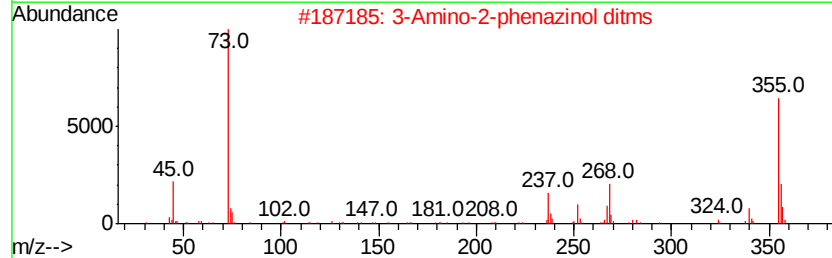
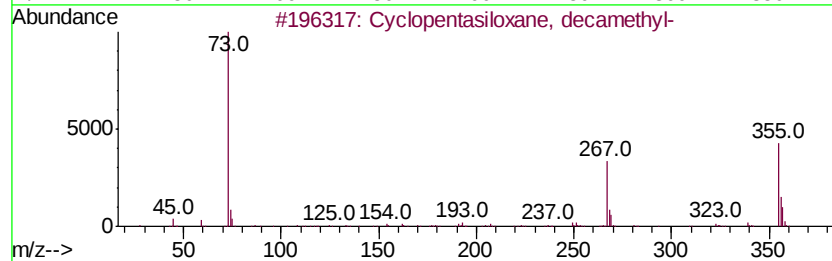
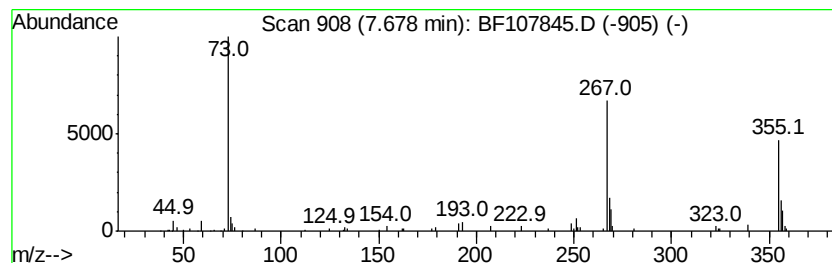
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Cyclopentasiloxane, decamet... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.68	2.47 ng	185623	Naphthalene-d8	8.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentasiloxane, decamethyl-	370	C10H30O5Si5	000541-02-6	90
2		3-Amino-2-phenazinol ditms	355	C18H25N3OSi2	1000332-15-5	49
3		Silanamine, 1,1,1-trimethyl-N-[2...	369	C17H35N02Si3	068595-60-8	43
4		Benzeneethanamine, N-butyl-.beta...	353	C18H35N02Si2	040629-66-1	38
5		Benzaldehyde, 2,5-bis[(trimethyl...	282	C13H22O3Si2	056114-69-3	38



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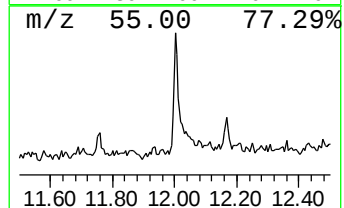
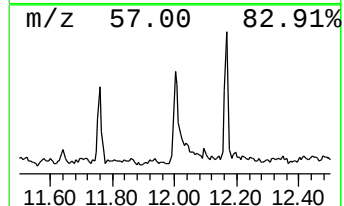
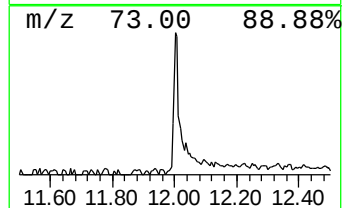
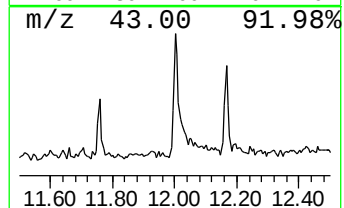
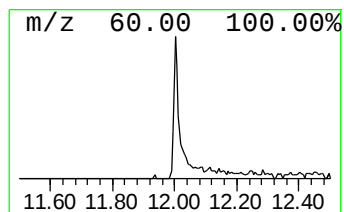
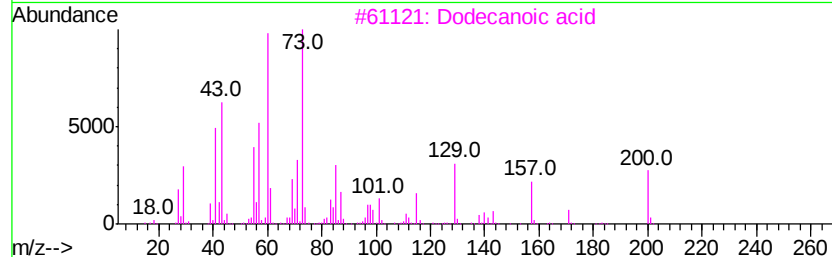
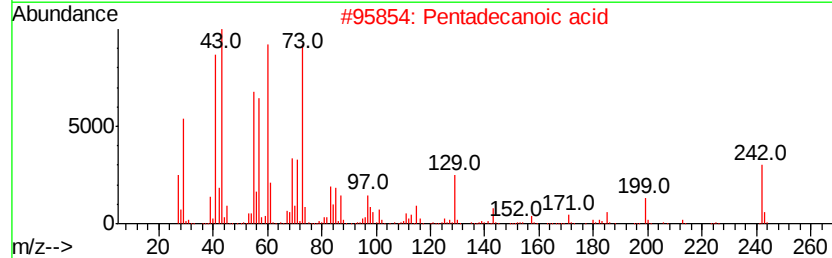
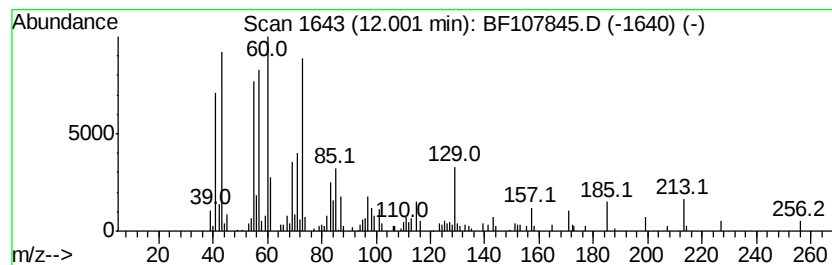
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 n-Hexadecanoic acid Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.00	2.42 ng	176574	Phenanthrene-d10	11.50

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	94
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	80
3		Dodecanoic acid	200	C12H24O2	000143-07-7	72
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	72
5		Tridecanoic acid	214	C13H26O2	000638-53-9	70



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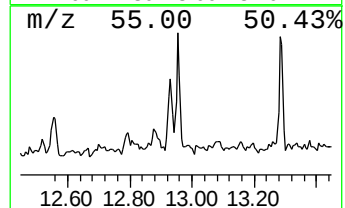
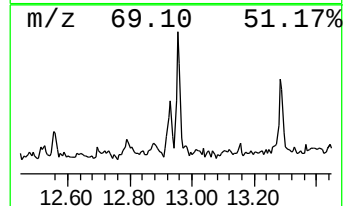
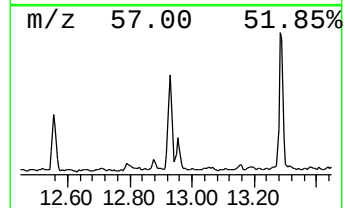
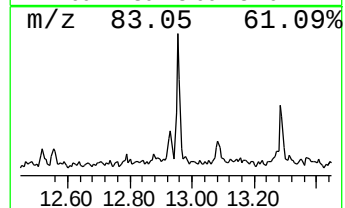
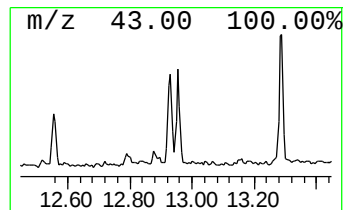
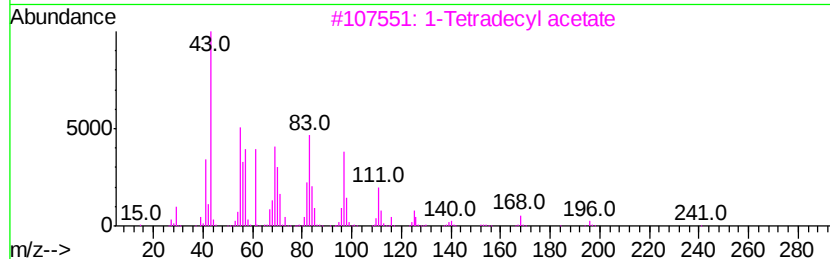
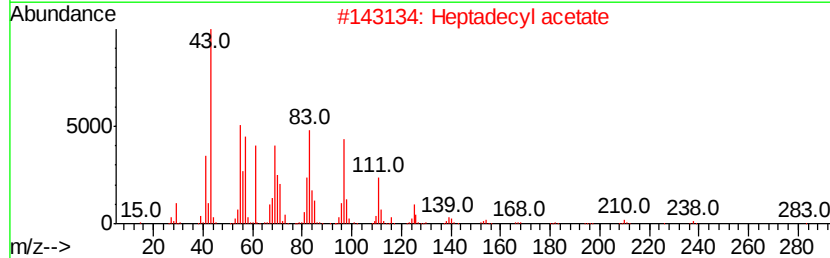
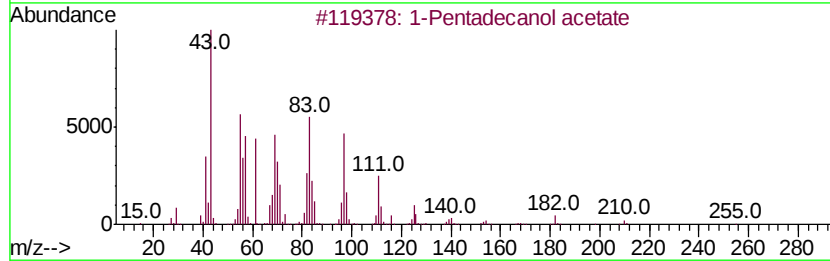
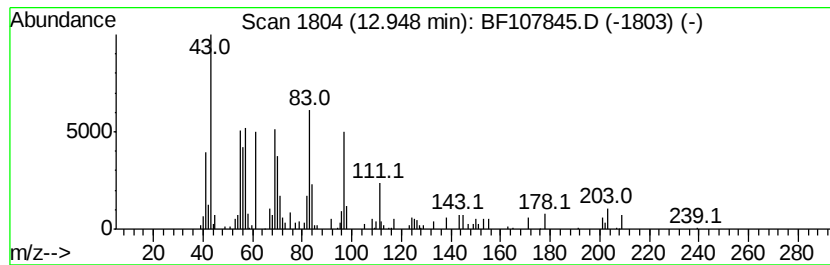
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 1-Pentadecanol acetate Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.95	2.06 ng	129521	Chrysene-d12	14.15

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Pentadecanol acetate	270	C17H34O2	000629-58-3	91
2		Heptadecyl acetate	298	C19H38O2	000822-20-8	90
3		1-Tetradecyl acetate	256	C16H32O2	000638-59-5	90
4		n-Heptadecanol-1	256	C17H36O	001454-85-9	76
5		1-Hexadecanol, acetate	284	C18H36O2	000629-70-9	74



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MW-A29

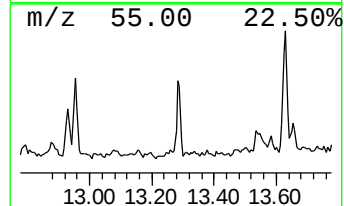
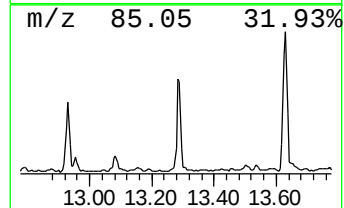
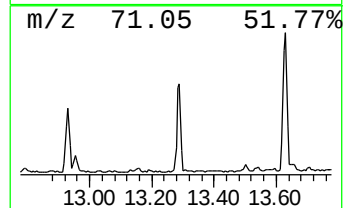
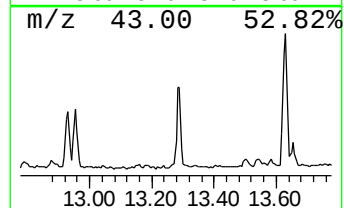
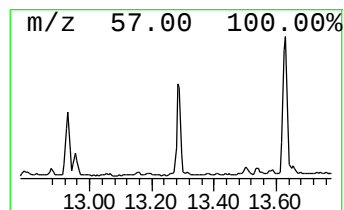
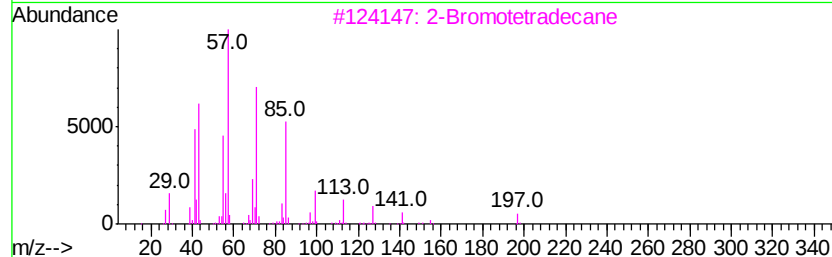
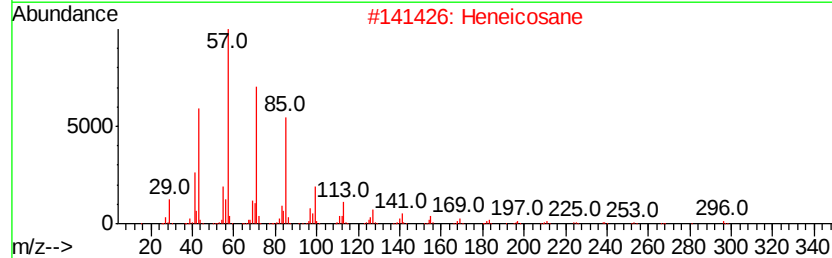
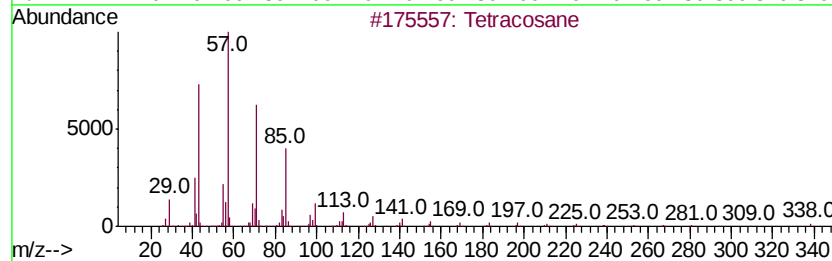
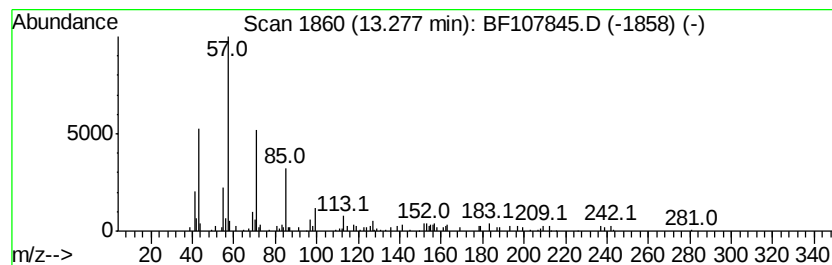
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Tetracosane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.28	3.03 ng	190555	Chrysene-d12	14.15

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetracosane	338	C24H50	000646-31-1	91
2		Heneicosane	296	C21H44	000629-94-7	90
3		2-Bromotetradecane	276	C14H29Br	074036-95-6	90
4		Eicosane	282	C20H42	000112-95-8	87
5		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF073018\
Data File : BF107845.D
Acq On : 31 Jul 2018 11:56
Operator : JU/SJ
Sample : J4221-06
Misc :
ALS Vial : 44 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-A29

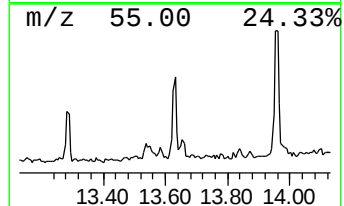
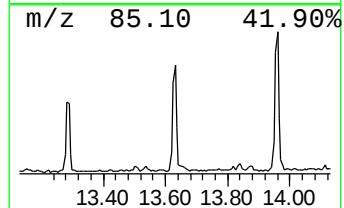
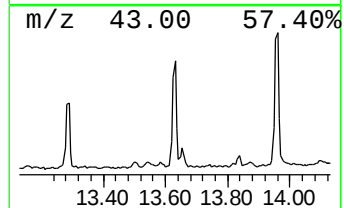
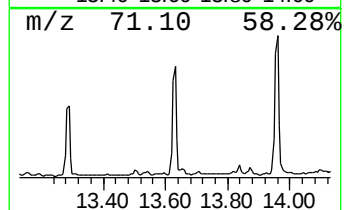
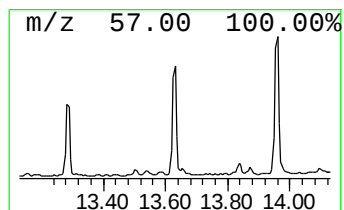
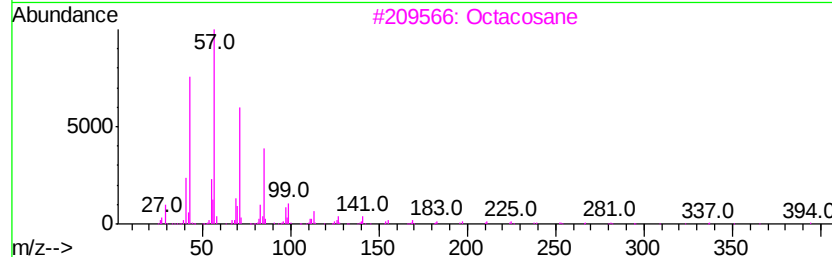
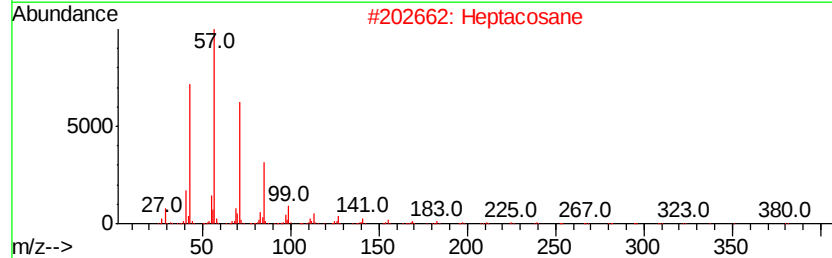
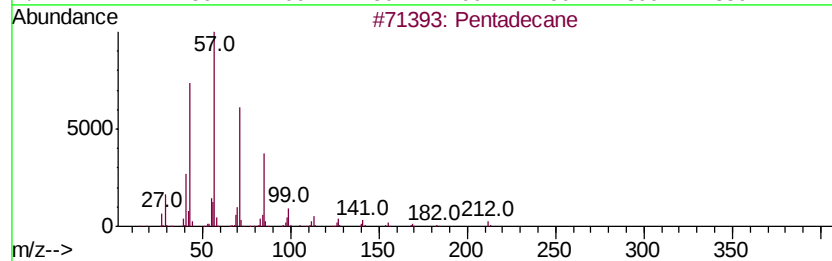
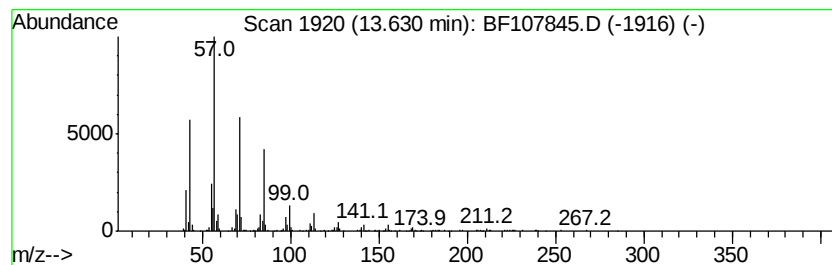
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Pentadecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.63	5.27 ng	330847	Chrysene-d12	14.15

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentadecane	212	C15H32	000629-62-9	93
2		Heptacosane	380	C27H56	000593-49-7	91
3		Octacosane	394	C28H58	000630-02-4	91
4		Tetracosane	338	C24H50	000646-31-1	91
5		Heneicosane	296	C21H44	000629-94-7	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF073018\
Data File : BF107845.D
Acq On : 31 Jul 2018 11:56
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Sample : J4221-06
Misc :
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Instrument :
BNA_F
ClientSampleId :
MW-A29

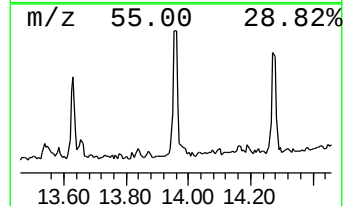
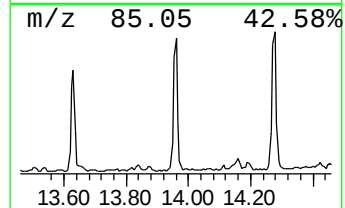
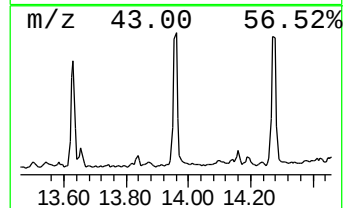
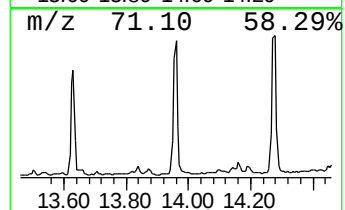
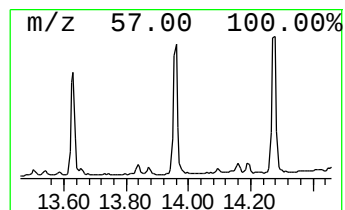
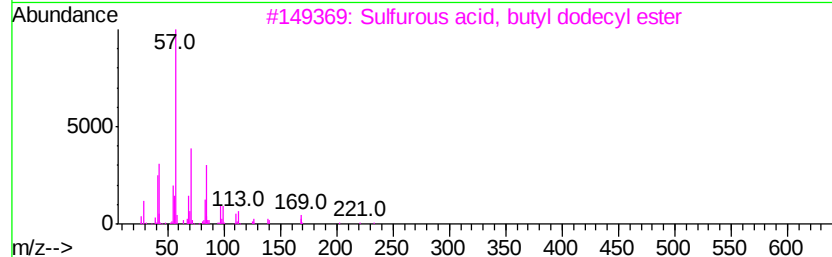
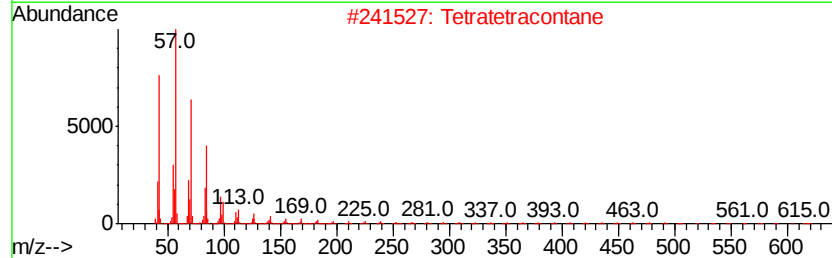
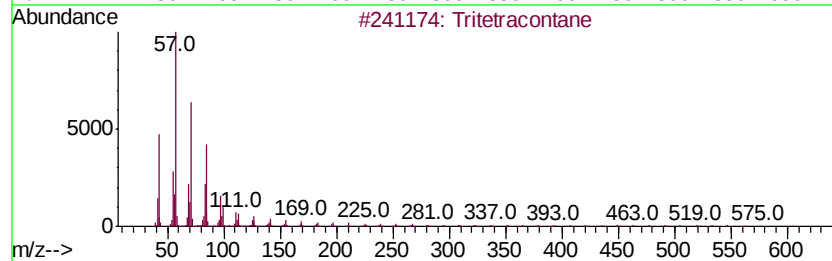
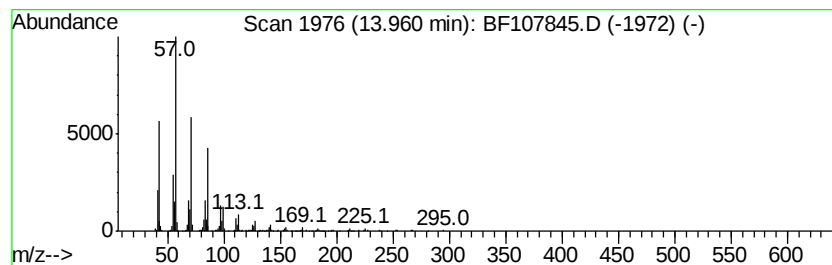
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Tritetracontane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.96	7.48 ng	469978	Chrysene-d12	14.15

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tritetracontane	605	C43H88	007098-21-7	91
2		Tetratetracontane	619	C44H90	007098-22-8	91
3		Sulfurous acid, butyl dodecyl ester	306	C16H34O3S	1000309-17-9	91
4		Sulfurous acid, butyl tetradecyl...	334	C18H38O3S	1000309-18-1	91
5		Sulfurous acid, butyl tridecyl e...	320	C17H36O3S	1000309-18-0	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF073018\
Data File : BF107845.D
Acq On : 31 Jul 2018 11:56
Operator : JU/SJ
Sample : J4221-06
Misc :
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Instrument :
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ClientSampleId :
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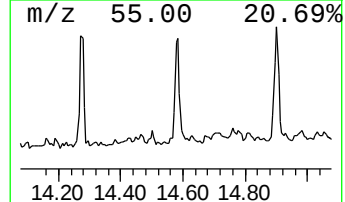
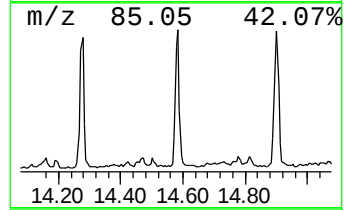
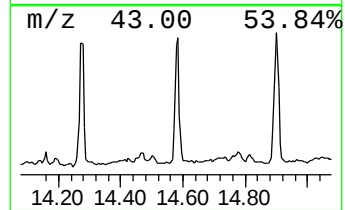
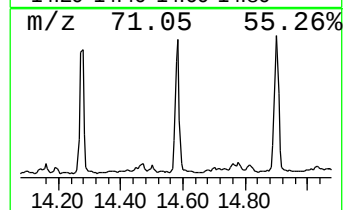
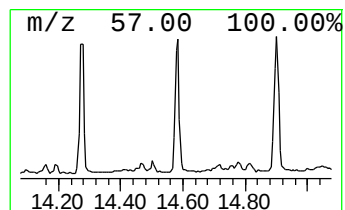
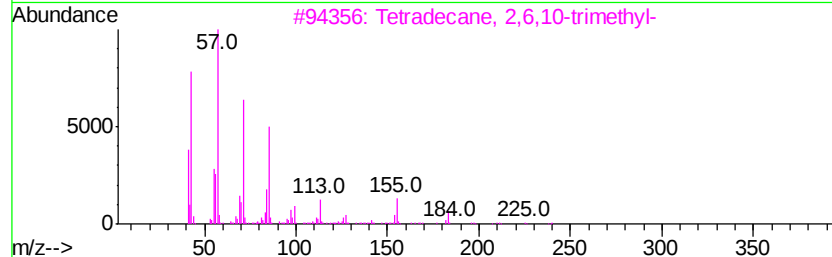
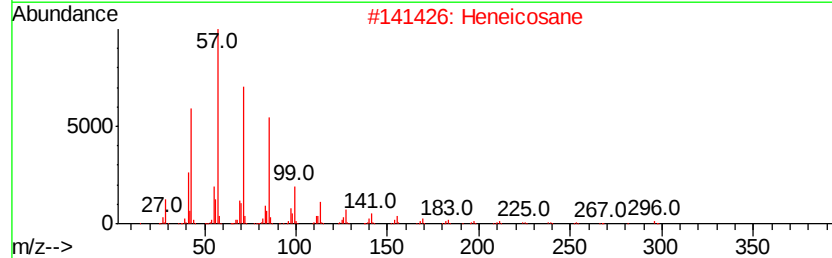
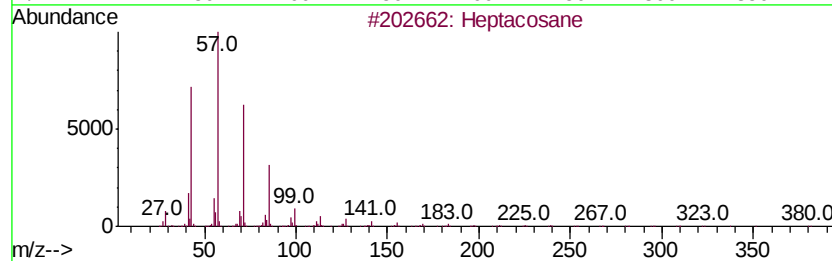
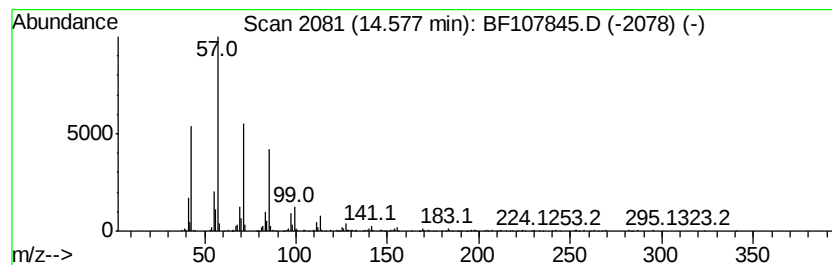
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Heptacosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.58	7.49 ng	470398	Chrysene-d12	14.15

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptacosane	380	C27H56	000593-49-7	91
2		Heneicosane	296	C21H44	000629-94-7	90
3		Tetradecane, 2,6,10-trimethyl-	240	C17H36	014905-56-7	90
4		Octacosane	394	C28H58	000630-02-4	90
5		Hexadecane	226	C16H34	000544-76-3	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF073018\
Data File : BF107845.D
Acq On : 31 Jul 2018 11:56
Operator : JU/SJ
Sample : J4221-06
Misc :
ALS Vial : 44 Sample Multiplier: 1

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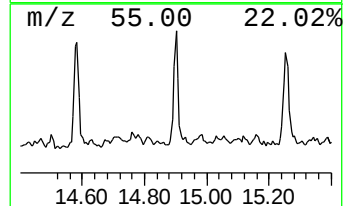
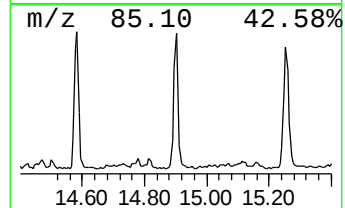
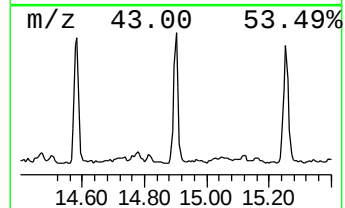
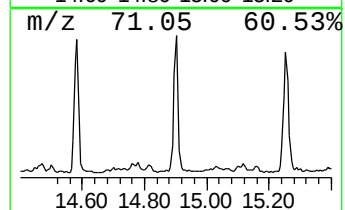
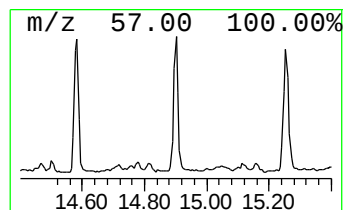
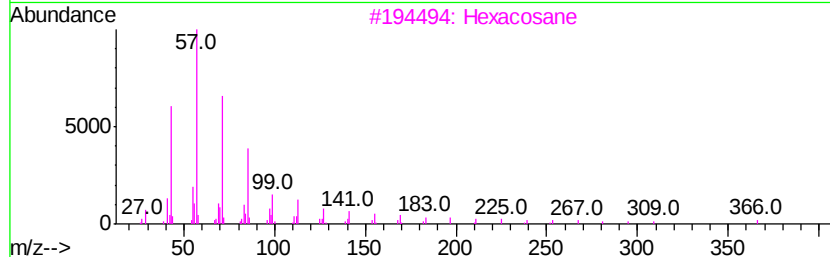
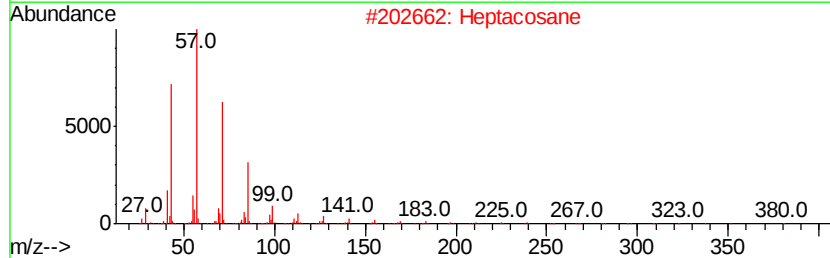
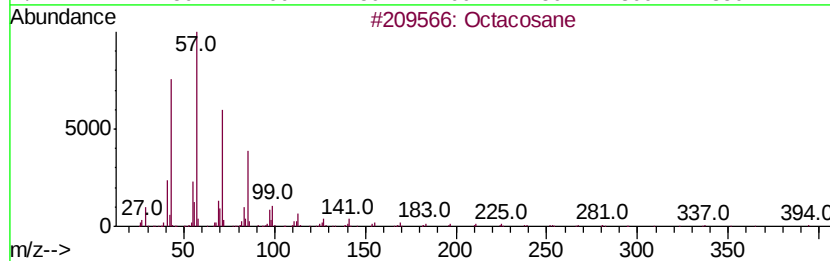
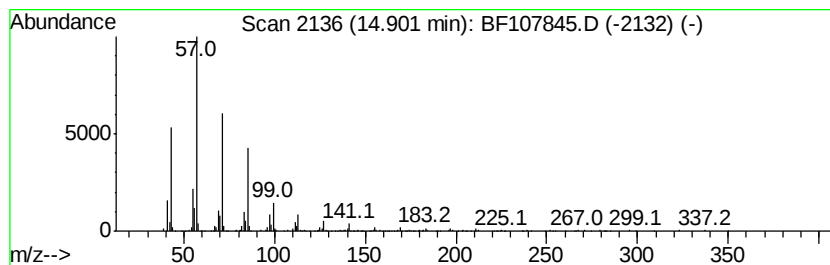
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Octacosane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.90	7.59 ng	476992	Chrysene-d12	14.15

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosane	394	C28H58	000630-02-4	91
2		Heptacosane	380	C27H56	000593-49-7	91
3		Hexacosane	366	C26H54	000630-01-3	90
4		Docosane	310	C22H46	000629-97-0	90
5		Tridecane, 6-propyl-	226	C16H34	055045-10-8	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF073018\
Data File : BF107845.D
Acq On : 31 Jul 2018 11:56
Operator : JU/SJ
Sample : J4221-06
Misc :
ALS Vial : 44 Sample Multiplier: 1

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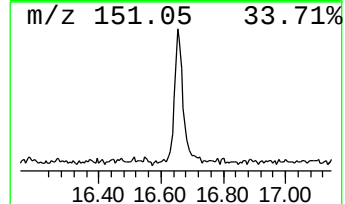
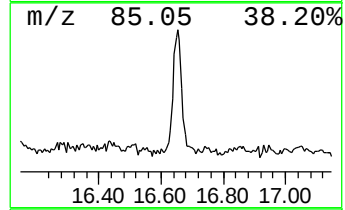
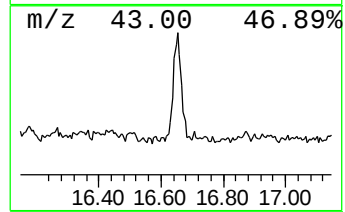
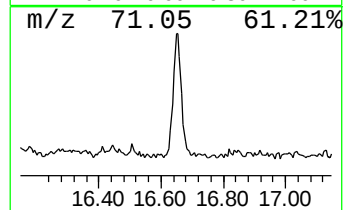
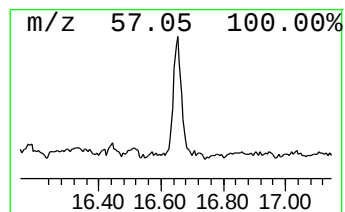
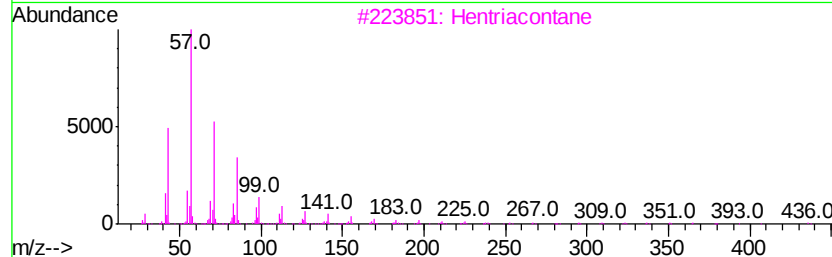
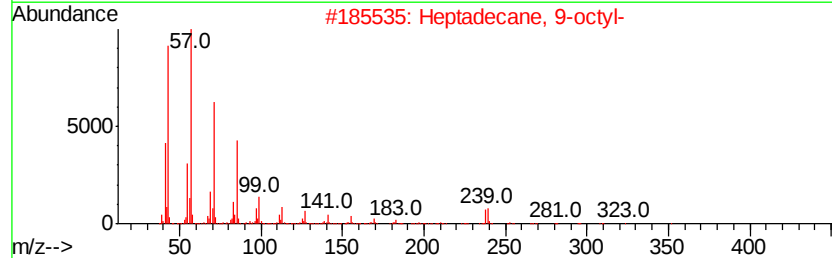
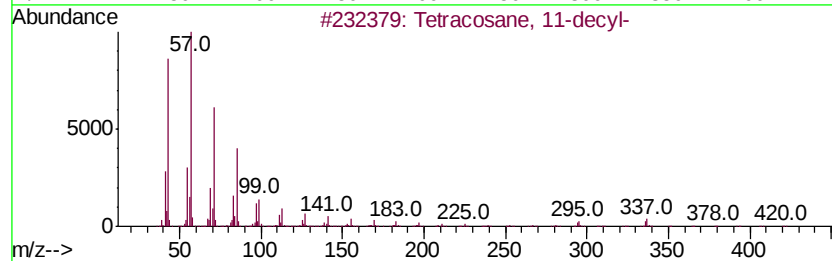
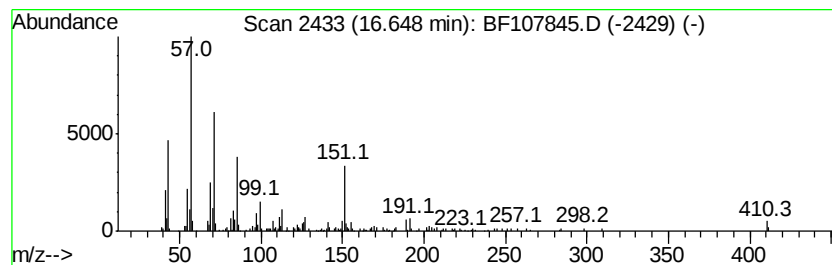
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Tetracosane, 11-decyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.65	6.33 ng	311445	Perylene-d12	15.69

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetracosane, 11-decyl-	479	C34H70	055429-84-0	62
2			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	62
3			Hentriacontane	437	C31H64	000630-04-6	62
4			2-methyloctacosane	408	C29H60	1000376-72-8	58
5			Tetratetracontane	619	C44H90	007098-22-8	58



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF073018\
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Acq On : 31 Jul 2018 11:56
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Sample : J4221-06
Misc :
ALS Vial : 44 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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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
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unknown6.74	6.74	78.0	ng	4559720	1	6.97	1169720	20.0
Cyclopentasiloxan...	7.68	2.5	ng	185623	2	8.24	1500210	20.0
n-Hexadecanoic acid	12.00	2.4	ng	176574	4	11.50	1461860	20.0
1-Pentadecanol ac...	12.95	2.1	ng	129521	5	14.15	1256680	20.0
Tetracosane	13.28	3.0	ng	190555	5	14.15	1256680	20.0
Pentadecane	13.63	5.3	ng	330847	5	14.15	1256680	20.0
Tritetracontane	13.96	7.5	ng	469978	5	14.15	1256680	20.0
Heptacosane	14.58	7.5	ng	470398	5	14.15	1256680	20.0
Octacosane	14.90	7.6	ng	476992	5	14.15	1256680	20.0
Tetracosane, 11-d...	16.65	6.3	ng	311445	6	15.69	983401	20.0