

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM122018\
 Data File : BM018342.D
 Acq On : 20 Dec 2018 22:23
 Operator : JU/SJ
 Sample : J6386-03
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 P001-SS016-0006-02

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_M\METHODS\8270-BM121518.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	3.111	34	38	49	rBV	409608	506412	1.65%	0.577%
2	5.134	375	382	395	rBV	2898514	4211568	13.76%	4.802%
3	6.705	638	649	662	rBV	2746520	4471196	14.61%	5.098%
4	7.034	697	705	715	rBV	3041281	4566868	14.92%	5.207%
5	7.487	776	782	790	rVB	554443	855035	2.79%	0.975%
6	7.799	828	835	847	rVB	2229914	3460911	11.31%	3.946%
7	8.646	972	979	989	rBV	1615314	2653730	8.67%	3.026%
8	10.251	1244	1252	1263	rVB	611857	1026537	3.35%	1.170%
9	12.751	1670	1677	1685	rBV	2969526	4350344	14.22%	4.960%
10	13.616	1815	1824	1834	rBV2	225127	426235	1.39%	0.486%
11	14.145	1908	1914	1922	rVB3	718179	1140357	3.73%	1.300%
12	15.657	2165	2171	2178	rBV	2248704	3308282	10.81%	3.772%
13	16.916	2379	2385	2390	rBV	788267	1184396	3.87%	1.350%
14	17.769	2525	2530	2535	rBV	225646	359002	1.17%	0.409%
15	17.827	2536	2540	2549	rVV	692390	991762	3.24%	1.131%
16	19.569	2831	2836	2843	rVB	3169224	3846207	12.57%	4.385%
17	19.857	2881	2885	2887	rBV2	310741	427205	1.40%	0.487%
18	20.851	3045	3054	3074	rVB	7373298	30600853	100.00%	34.889%
19	21.139	3098	3103	3111	rBV	1114463	1896362	6.20%	2.162%
20	21.474	3157	3160	3164	rVB	291065	324039	1.06%	0.369%
21	21.721	3198	3202	3203	rBV	1417746	1746457	5.71%	1.991%
22	21.739	3203	3205	3214	rVB	1964897	2190518	7.16%	2.497%
23	22.386	3311	3315	3319	rVB	527467	712346	2.33%	0.812%
24	22.645	3355	3359	3363	rVV	1199170	1590481	5.20%	1.813%
25	22.692	3363	3367	3375	rVB	832501	1264789	4.13%	1.442%
26	23.298	3466	3470	3480	rVB	812968	1549317	5.06%	1.766%
27	23.480	3497	3501	3509	rVB	920745	1469145	4.80%	1.675%
28	23.786	3549	3553	3560	rVB	734625	1260913	4.12%	1.438%
29	29.303	4485	4491	4505	rVB6	1478227	5318292	17.38%	6.064%

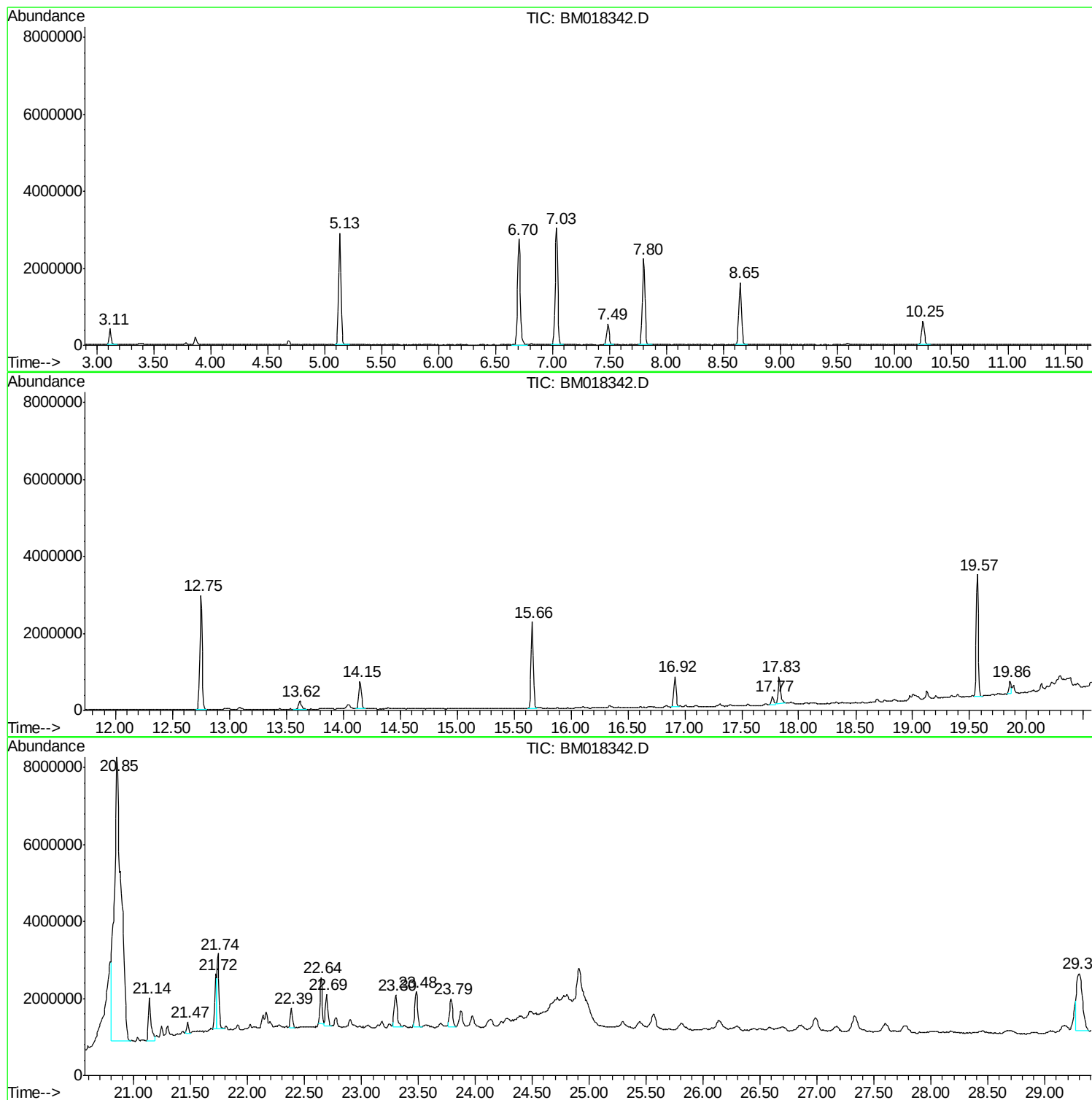
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM121518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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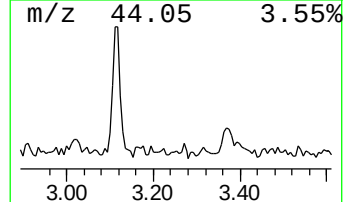
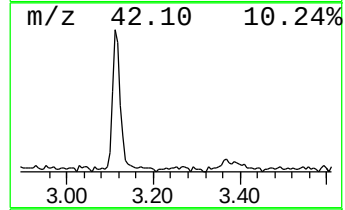
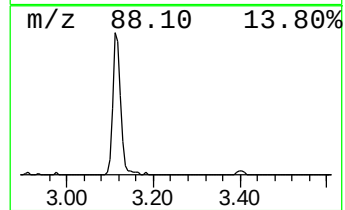
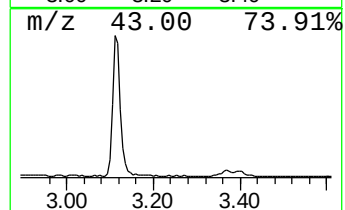
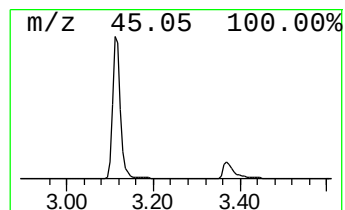
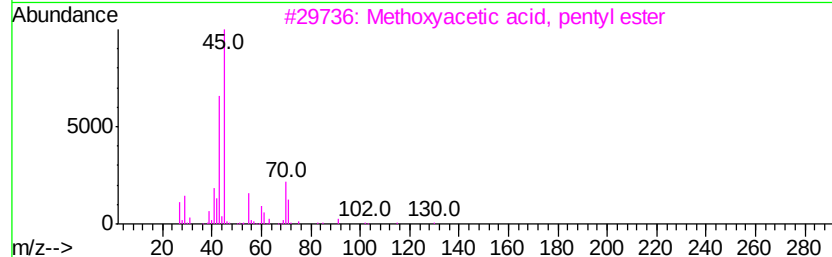
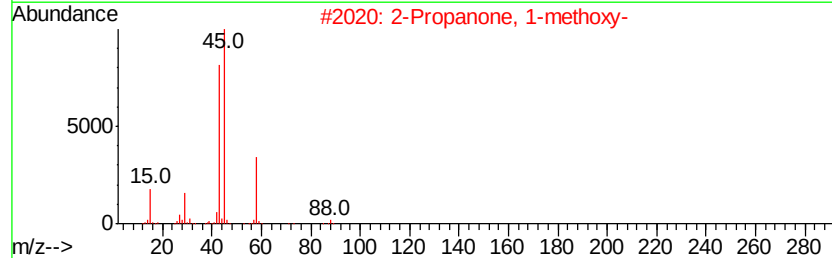
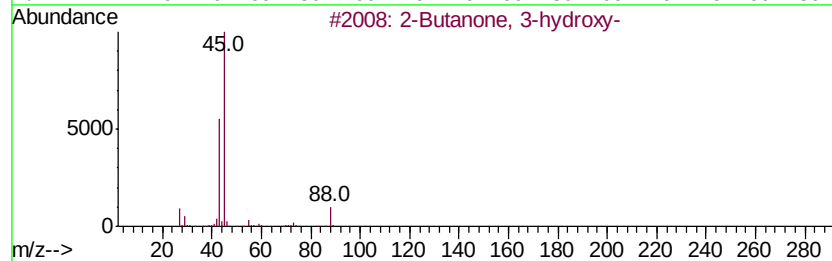
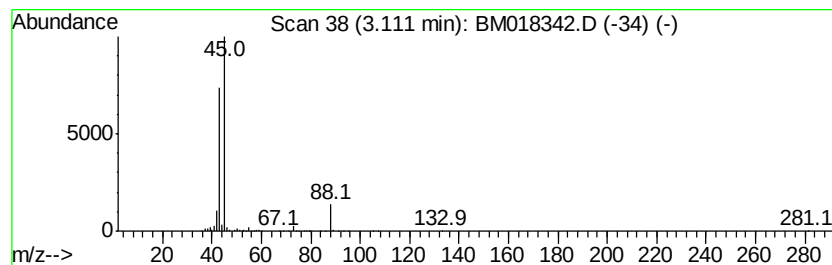
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
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Peak Number 1 2-Butanone, 3-hydroxy- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.11	11.85 ng	506412	1,4-Dichlorobenzene-d4	7.49

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Butanone, 3-hydroxy-	88	C4H8O2	000513-86-0	78
2			2-Propanone, 1-methoxy-	88	C4H8O2	005878-19-3	9
3			Methoxyacetic acid, pentyl ester	160	C8H16O3	168920-35-2	9
4			Methoxyacetic acid, hexyl ester	174	C9H18O3	145747-16-6	9
5			4-Penten-2-ol	86	C5H10O	000625-31-0	9



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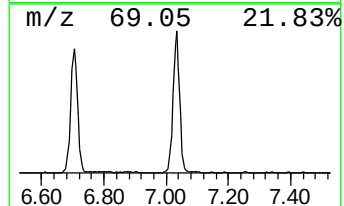
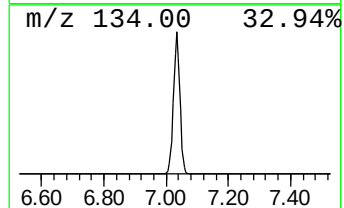
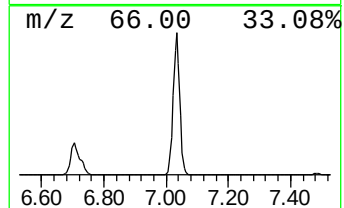
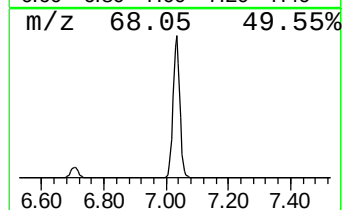
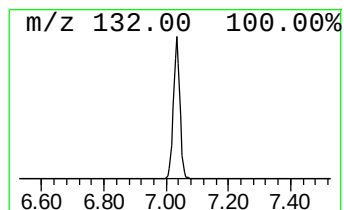
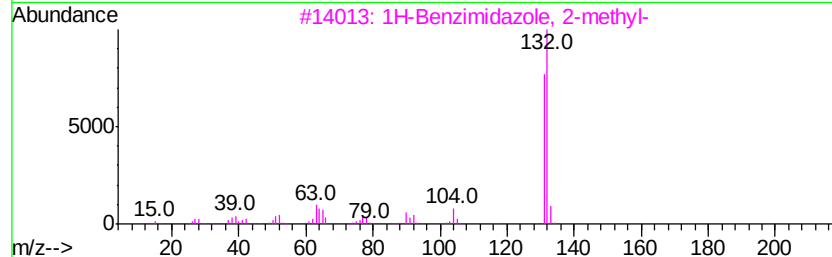
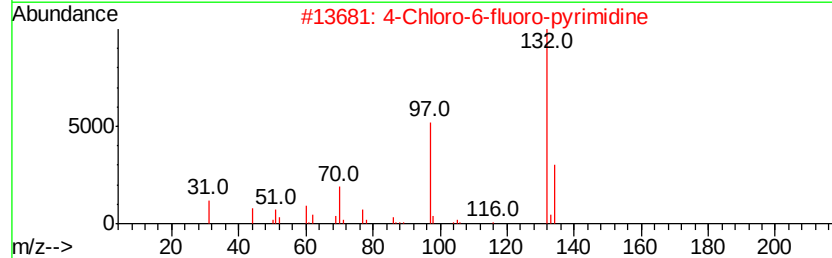
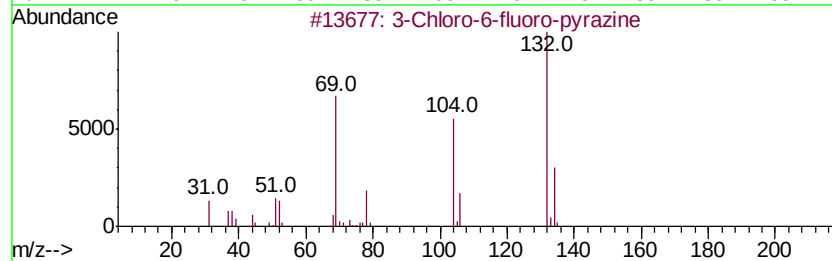
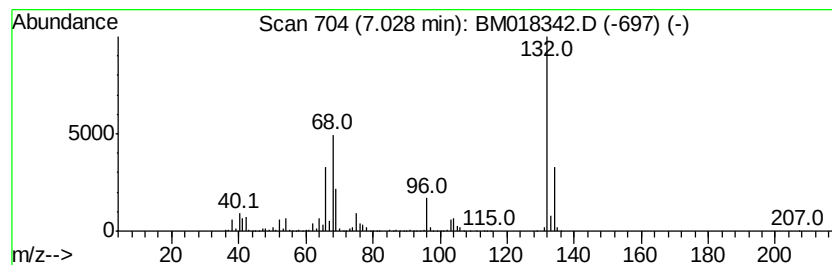
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Peak Number 2 unknown7.03 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.03	106.82 ng	4566870	1,4-Dichlorobenzene-d4	7.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
5		(5-METHYL-2-PYRIDYL)ACETONITRILE	132	C8H8N2	1000241-93-9	10



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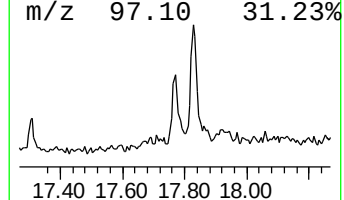
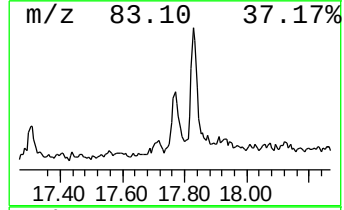
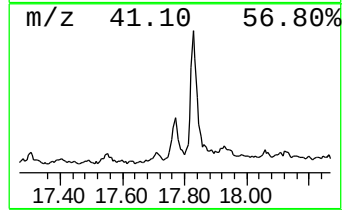
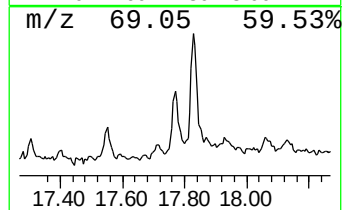
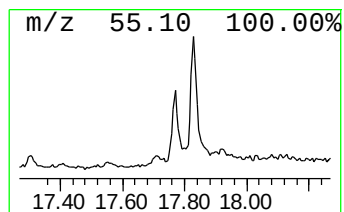
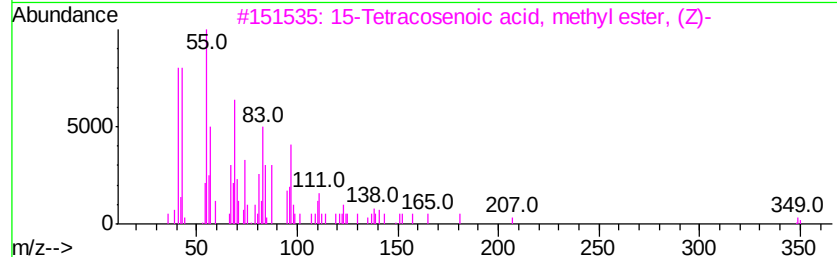
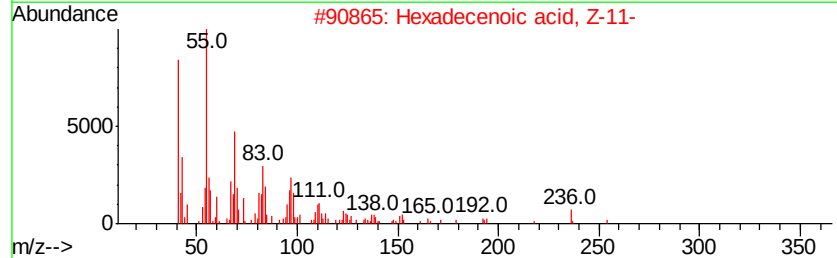
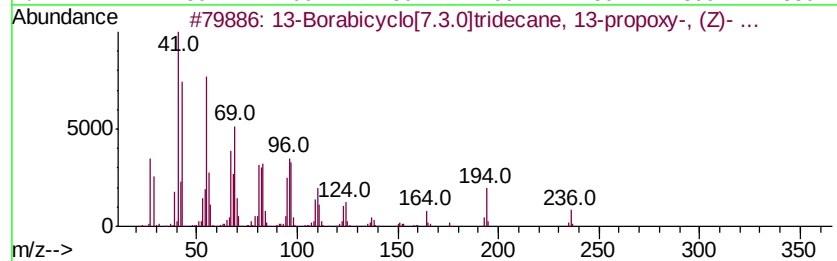
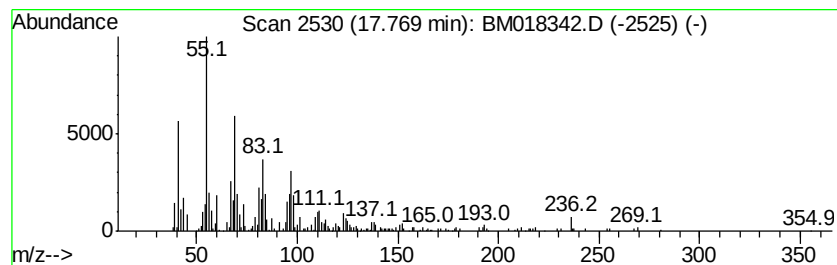
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Peak Number 4 13-Borabicyclo[7.3.0]tridec... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.77	6.06 ng	359002	Phenanthrene-d10	16.92	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	13-Borabicyclo[7.3.0]tridecane, ...	236	C15H29BO	1000156-41-7	96
2	Hexadecenoic acid, Z-11-	254	C16H30O2	002416-20-8	93
3	15-Tetracosenoic acid, methyl es...	380	C25H48O2	002733-88-2	74
4	Cyclopentadecanone, 2-hydroxy-	240	C15H28O2	004727-18-8	72
5	Z-11-Tetradecenoic acid	226	C14H26O2	1000130-83-3	64



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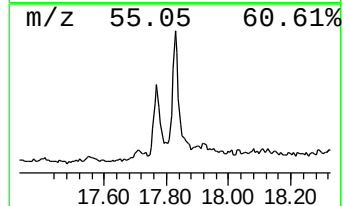
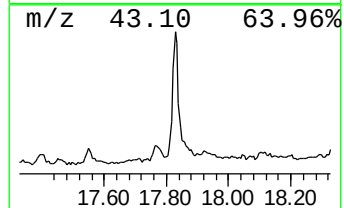
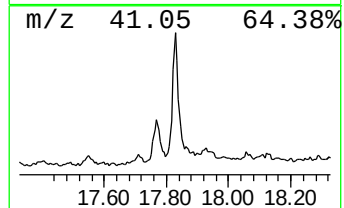
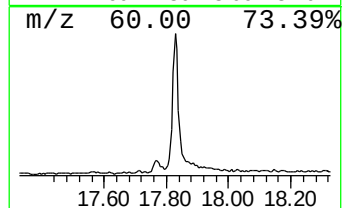
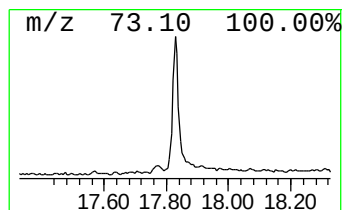
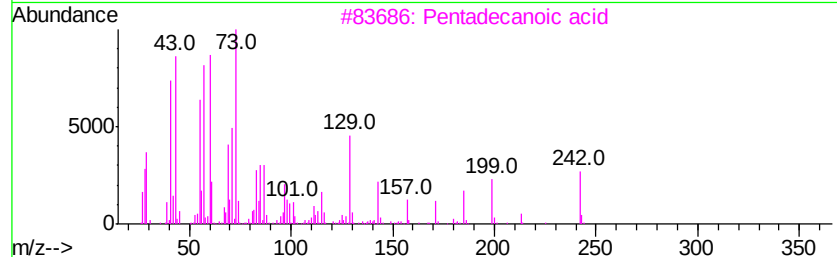
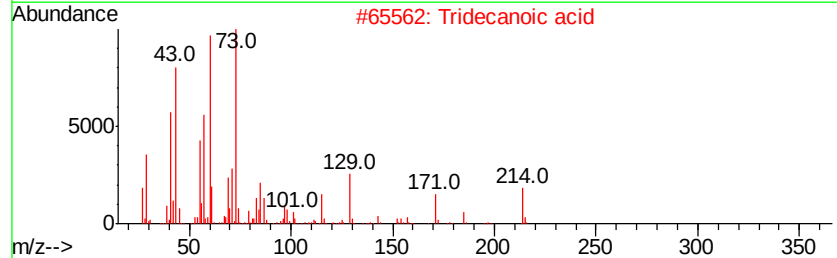
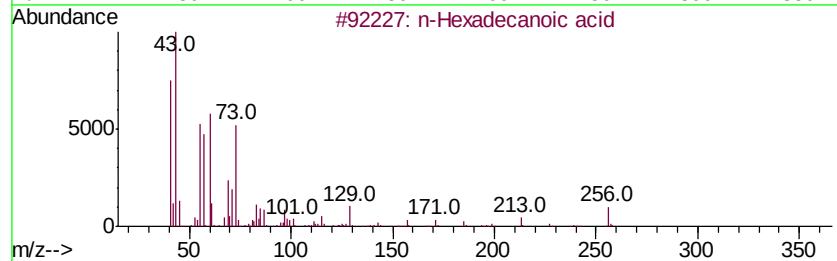
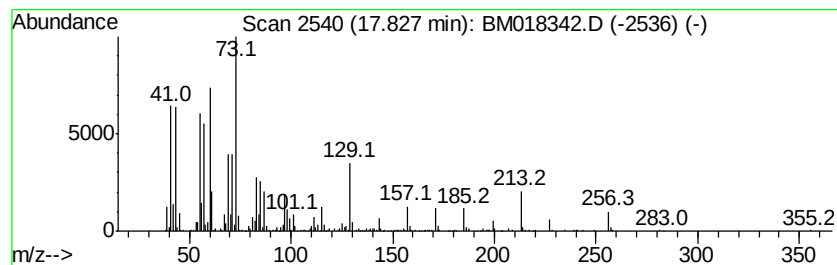
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Peak Number 5 n-Hexadecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.83	16.75 ng	991762	Phenanthrene-d10	16.92

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2		Tridecanoic acid	214	C13H26O2	000638-53-9	91
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	83
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	72
5		l-Tyrosinol	167	C9H13NO2	1000131-27-0	18



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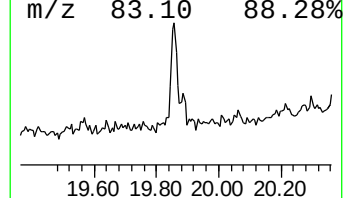
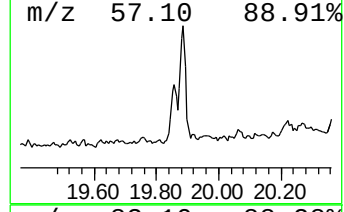
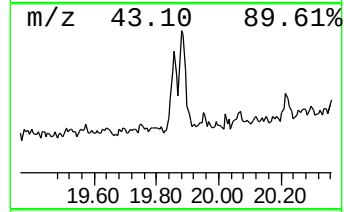
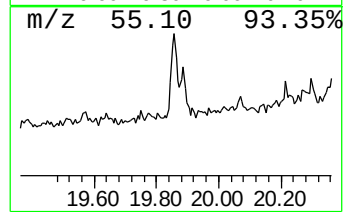
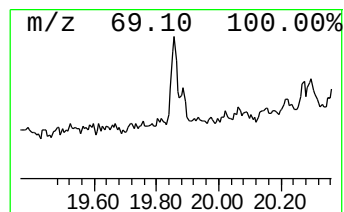
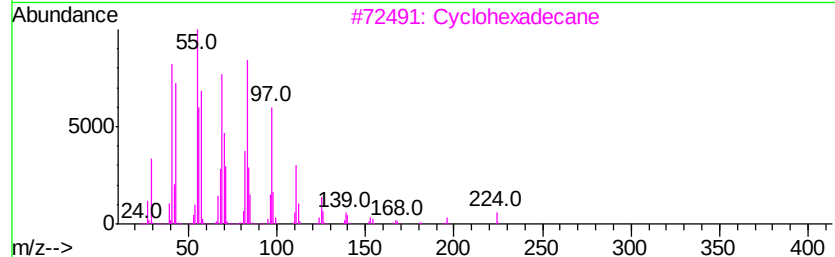
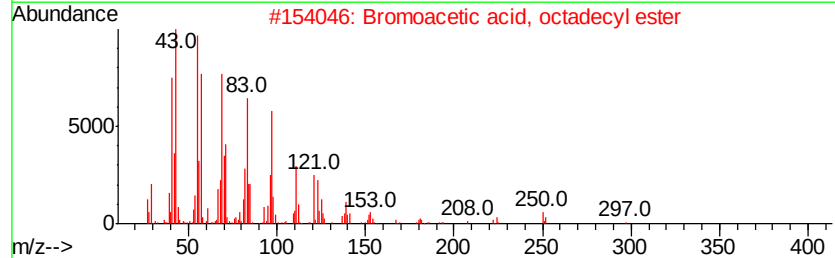
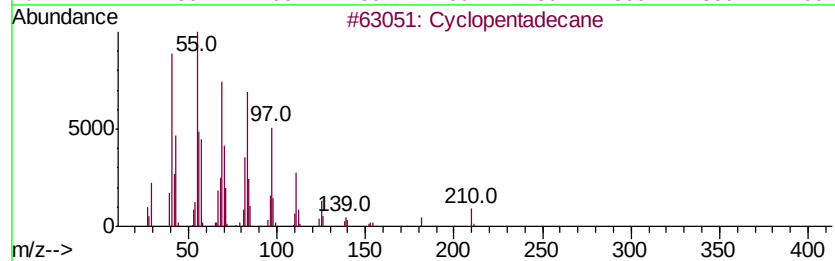
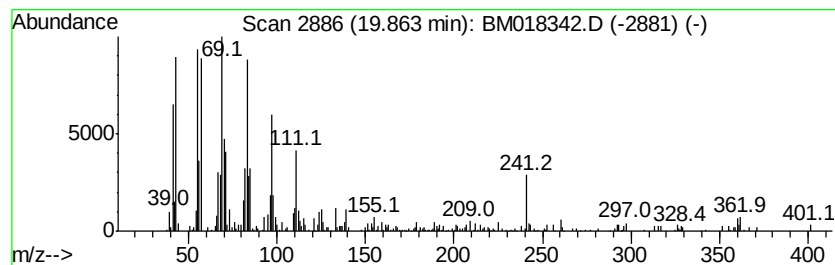
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Peak Number 6 Cyclopentadecane Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.86	4.51 ng	427205	Chrysene-d12	21.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentadecane	210	C15H30	000295-48-7	98
2		Bromoacetic acid, octadecyl ester	390	C20H39BrO2	018992-03-5	81
3		Cyclohexadecane	224	C16H32	000295-65-8	78
4		Bromoacetic acid, hexadecyl ester	362	C18H35BrO2	005454-48-8	76
5		9-Eicosene, (E)-	280	C20H40	074685-29-3	74



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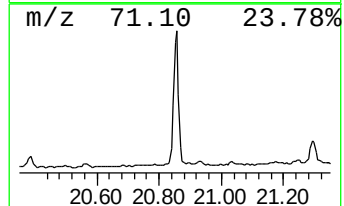
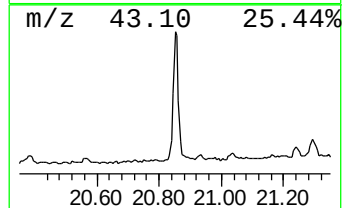
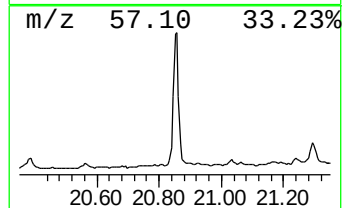
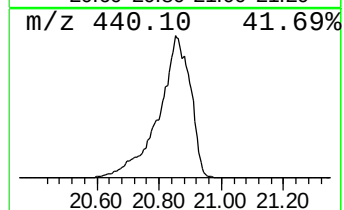
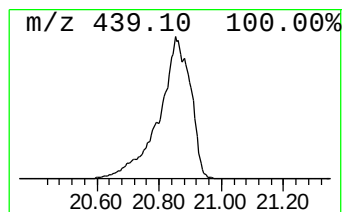
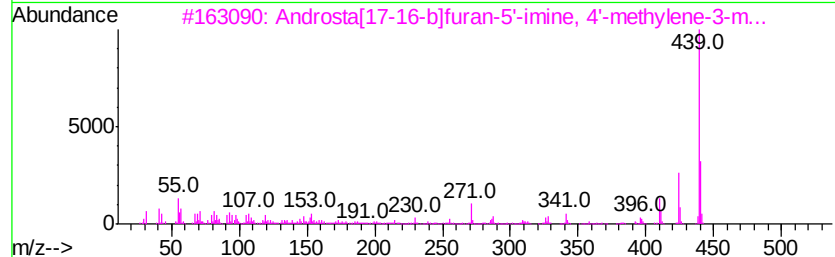
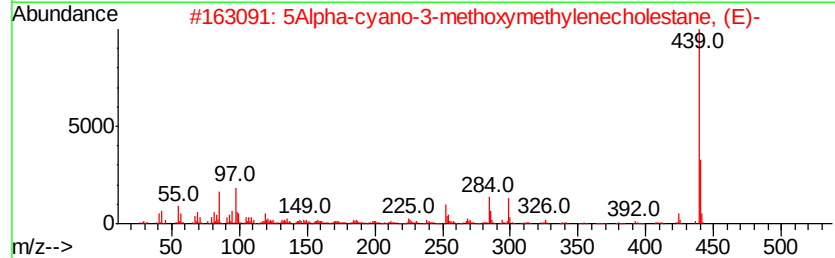
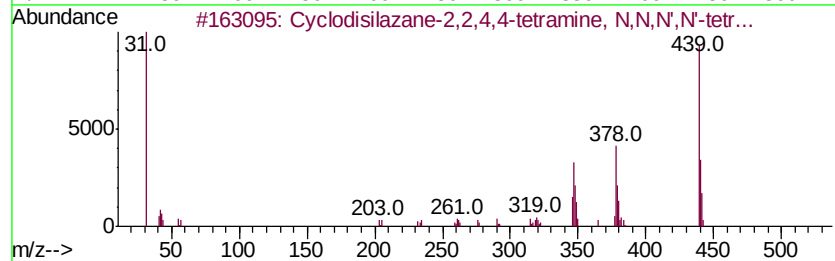
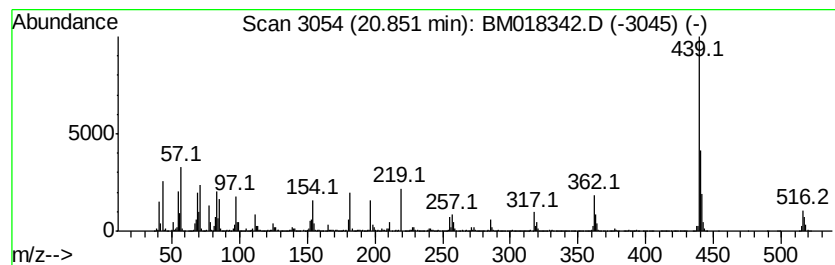
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Cyclodisilazane-2,2,4,4-tet... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.		
20.85	322.73 ng	30600900	Chrysene-d12	21.14		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cvcclodisilazane-2,2,4,4-tetramin...	440	C10H40N12Si4	034665-55-9	59
2		5Alpha-cyano-3-methoxymethylenec...	439	C30H49NO	021157-23-3	46
3		Androsta[17-16-blfuran-5'-imine...	439	C29H45N02	1000196-04-8	7
4		2,3-Pyridinedicarboxylic acid, 1...	439	C27H21N05	088078-86-8	5
5		Acetamide, 2,2,2-trifluoro-N-(5,...	439	C21H20F3N06	086436-46-6	3



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM122018\
Data File : BM018342.D
Acq On : 20 Dec 2018 22:23
Operator : JU/SJ
Sample : J6386-03
Misc :
ALS Vial : 13 Sample Multiplier: 1

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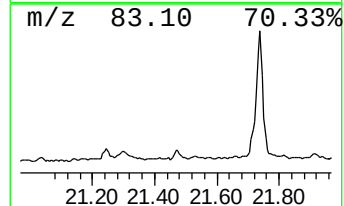
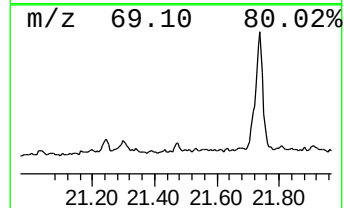
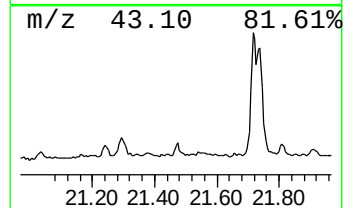
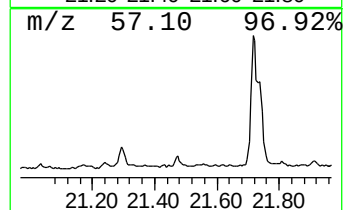
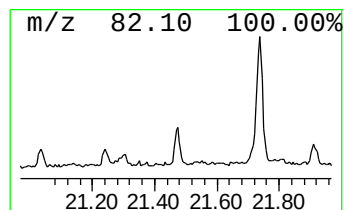
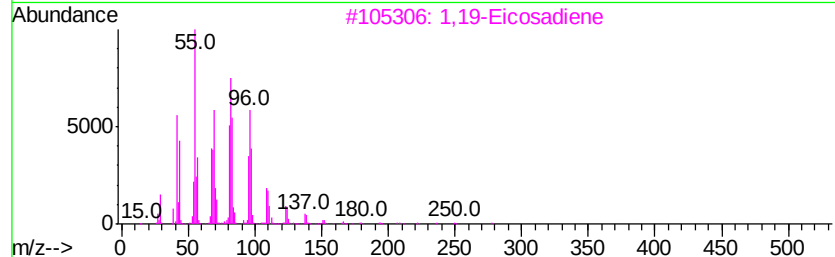
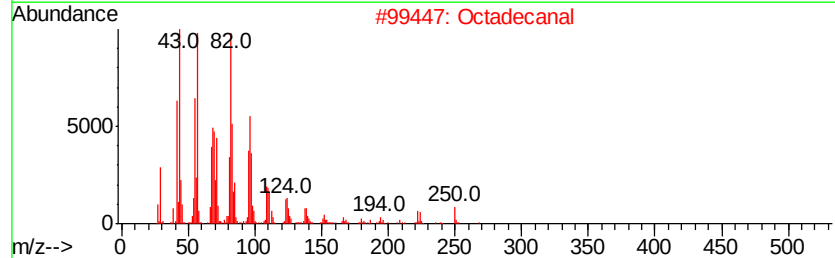
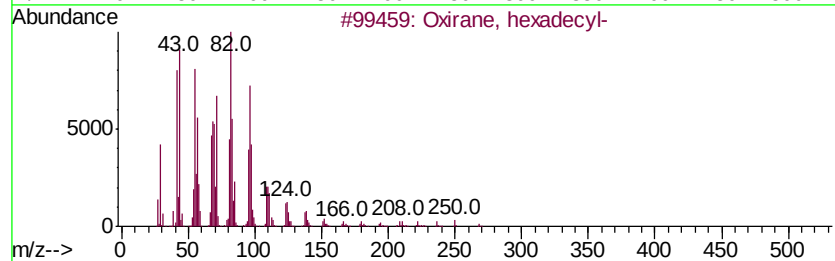
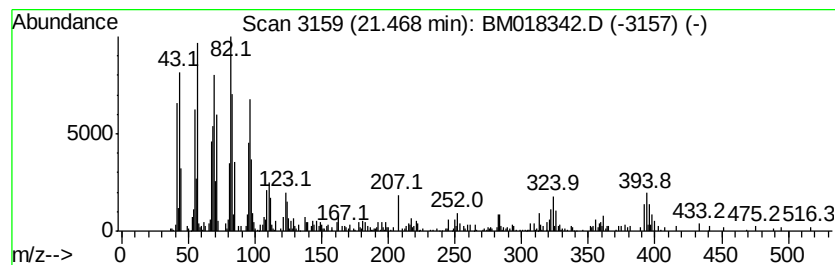
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Oxirane, hexadecyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.47	3.42 ng	324039	Chrysene-d12	21.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Oxirane, hexadecyl-	268	C18H36O	007390-81-0	94
2		Octadecanal	268	C18H36O	000638-66-4	93
3		1,19-Eicosadiene	278	C20H38	014811-95-1	93
4		16-Heptadecenal	252	C17H32O	1000144-57-9	83
5		Z-2-Octadecen-1-ol	268	C18H36O	1000131-11-0	70



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM122018\
Data File : BM018342.D
Acq On : 20 Dec 2018 22:23
Operator : JU/SJ
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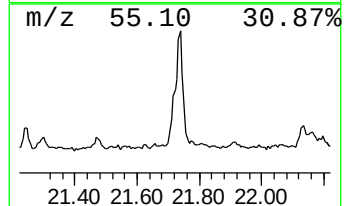
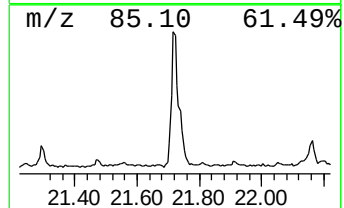
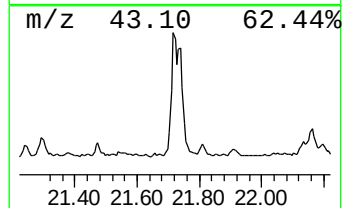
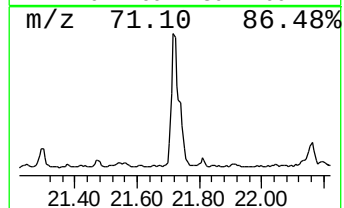
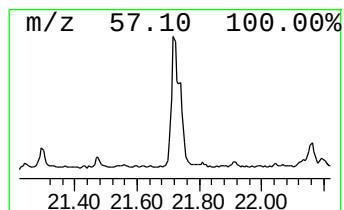
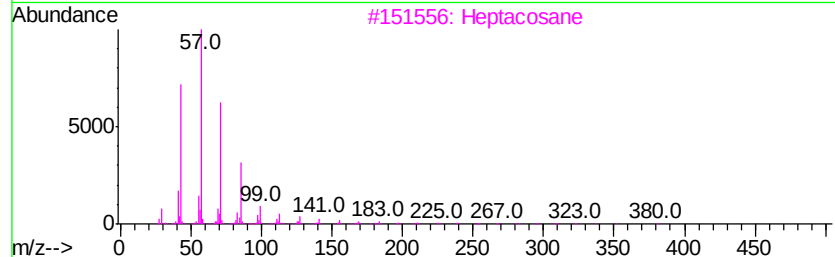
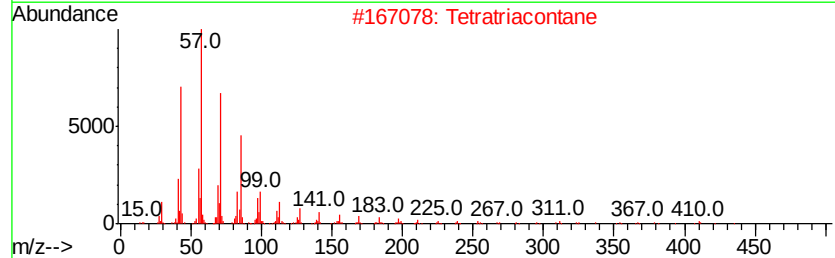
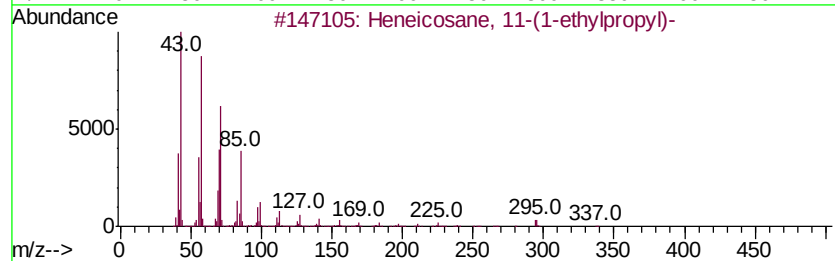
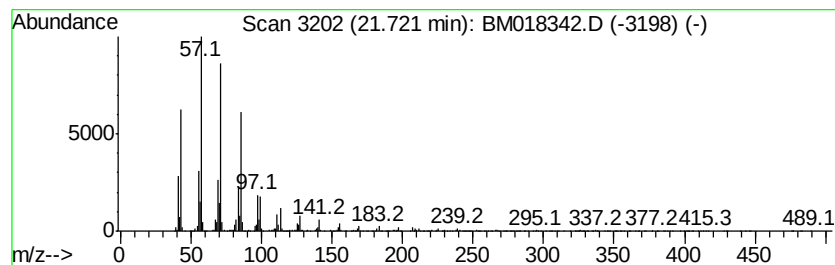
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Heneicosane, 11-(1-ethylpro... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.72	18.42 ng	1746460	Chrysene-d12	21.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	91
2		Tetratriacontane	479	C34H70	014167-59-0	91
3		Heptacosane	380	C27H56	000593-49-7	89
4		Tritetracontane	605	C43H88	007098-21-7	83
5		Eicosane	282	C20H42	000112-95-8	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM122018\
Data File : BM018342.D
Acq On : 20 Dec 2018 22:23
Operator : JU/SJ
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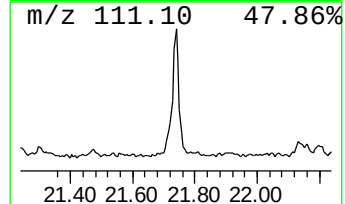
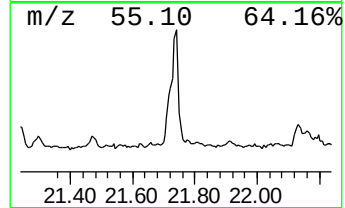
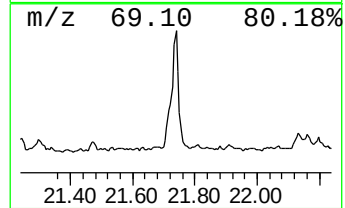
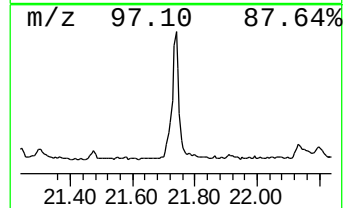
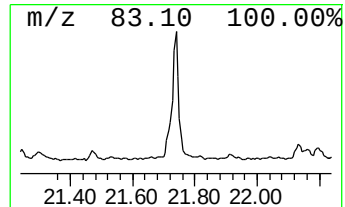
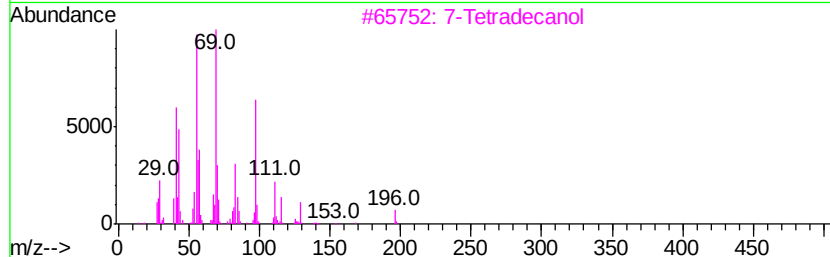
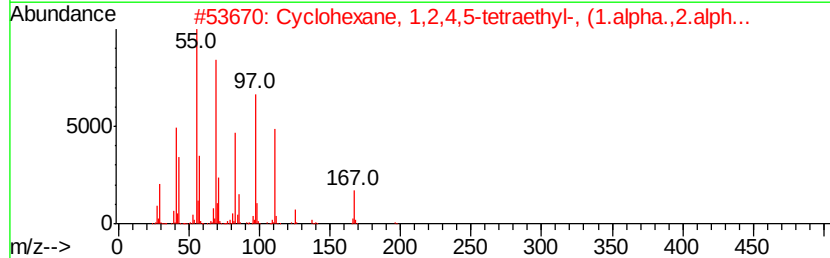
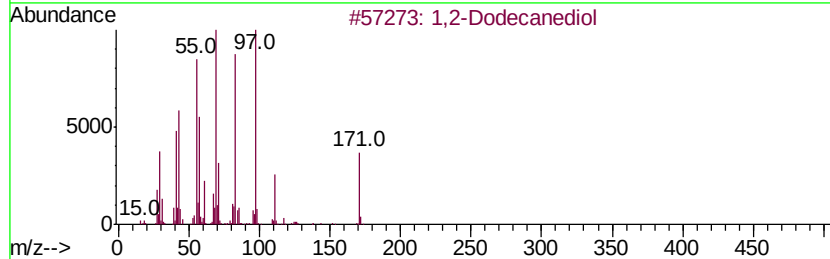
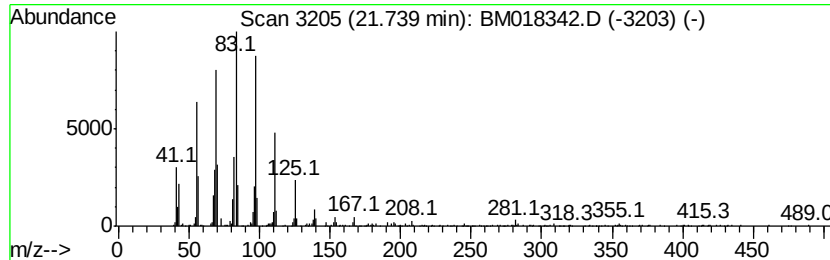
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 1,2-Dodecanediol Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.74	23.10 ng	2190520	Chrysene-d12	21.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,2-Dodecanediol	202	C12H26O2	001119-87-5	59
2		Cyclohexane, 1,2,4,5-tetraethyl-...	196	C14H28	061142-24-3	50
3		7-Tetradecanol	214	C14H30O	003981-79-1	47
4		Cyclopentane, 1,1,3,4-tetramethy...	126	C9H18	020309-77-7	47
5		Cyclopentane, (2-methylbutyl)-	140	C10H20	053366-38-4	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM122018\
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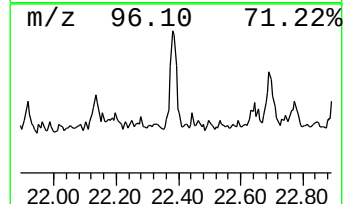
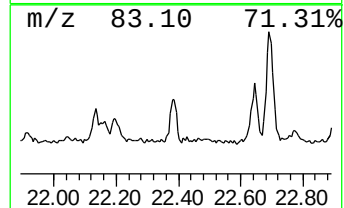
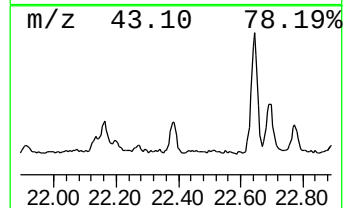
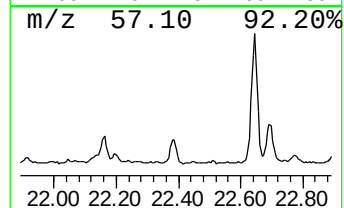
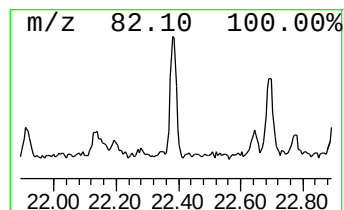
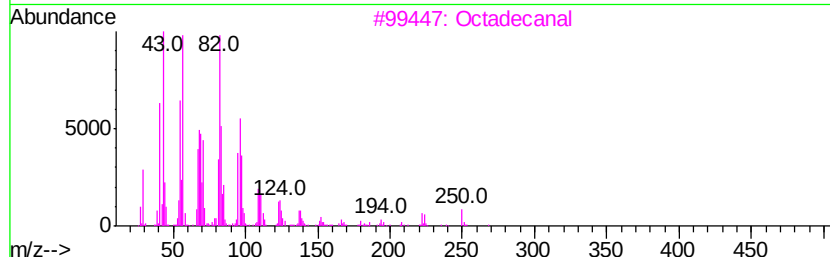
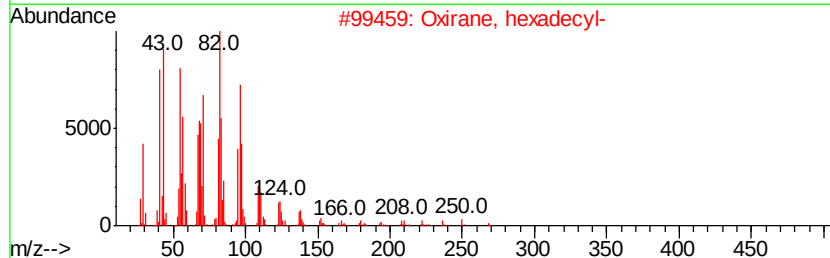
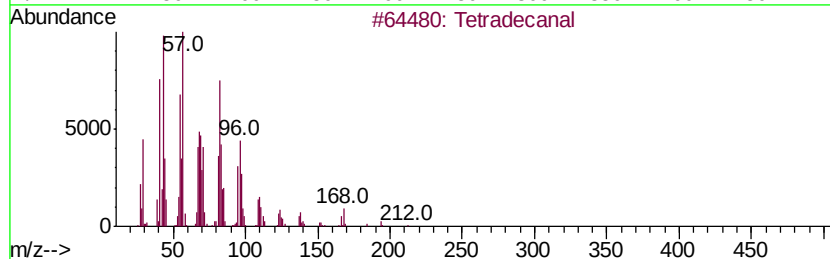
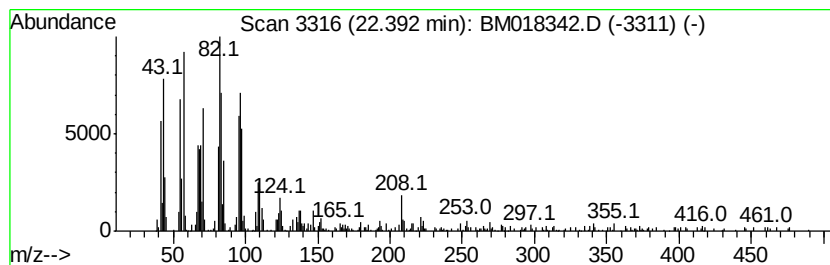
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Tetradeanal Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.39	9.20 ng	712346	Perylene-d12	23.30
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Tetradeanal	212 C14H28O	000124-25-4 91
2		Oxirane, hexadecyl-	268 C18H36O	007390-81-0 91
3		Octadecanal	268 C18H36O	000638-66-4 91
4		Oxirane, heptadecyl-	282 C19H38O	067860-04-2 90
5		Pentadecanal-	226 C15H30O	002765-11-9 87



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM122018\
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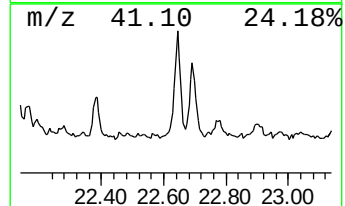
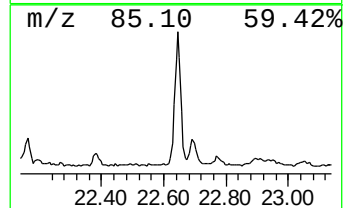
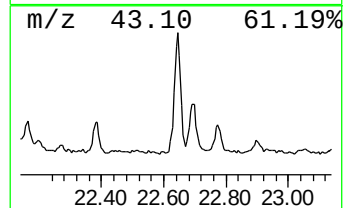
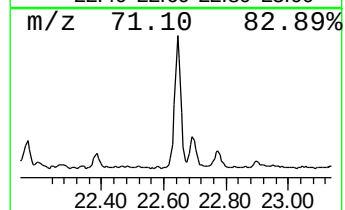
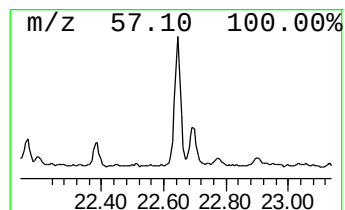
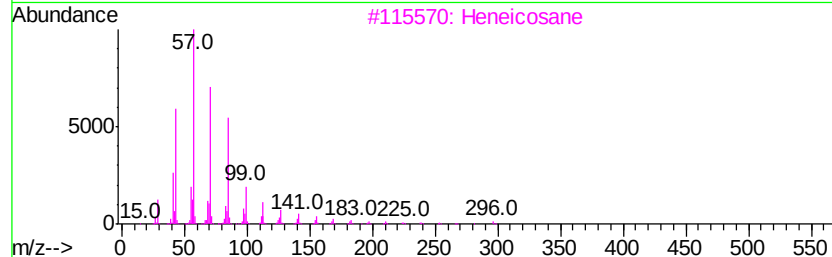
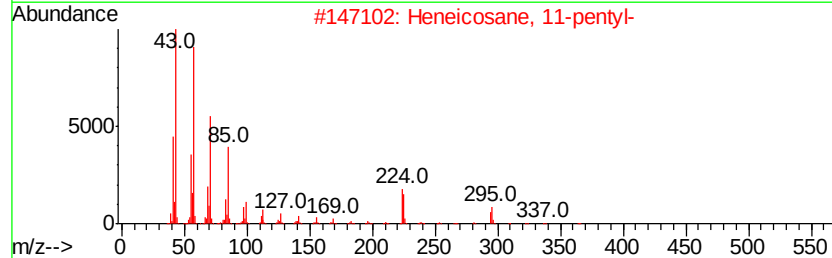
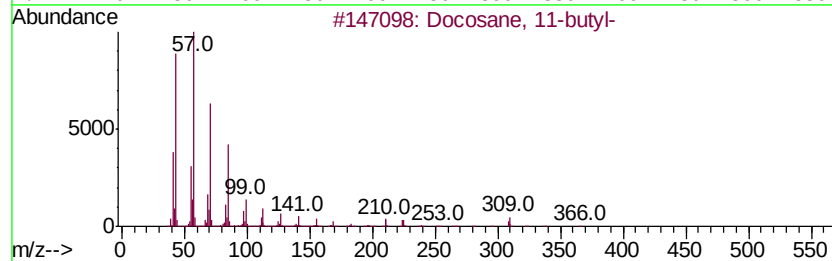
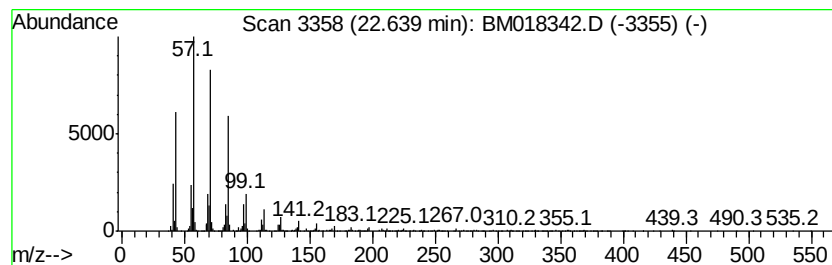
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Docosane, 11-butyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.64	20.53 ng	1590480	Perylene-d12	23.30
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Docosane, 11-butyl-	366 C26H54	013475-76-8 94
2		Heneicosane, 11-pentyl-	366 C26H54	014739-72-1 91
3		Heneicosane	296 C21H44	000629-94-7 91
4		Heneicosane, 11-(1-ethylpropyl)-	366 C26H54	055282-11-6 91
5		Heptacosane	380 C27H56	000593-49-7 91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM122018\
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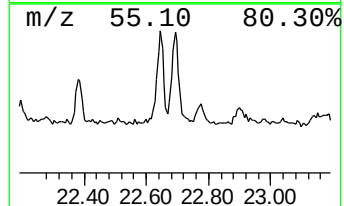
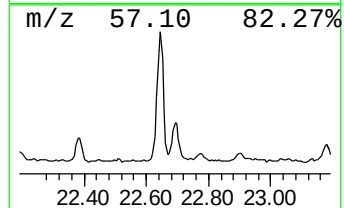
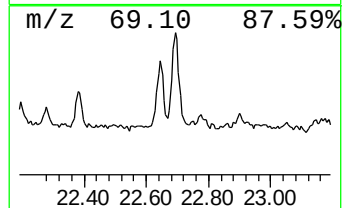
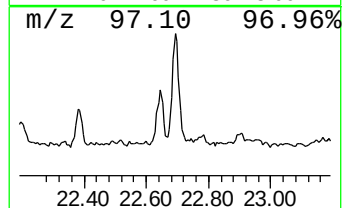
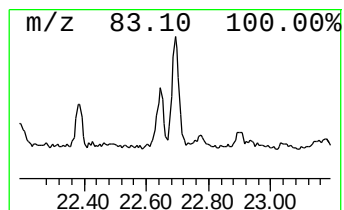
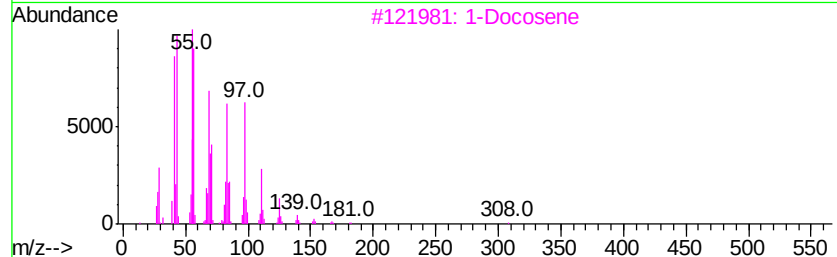
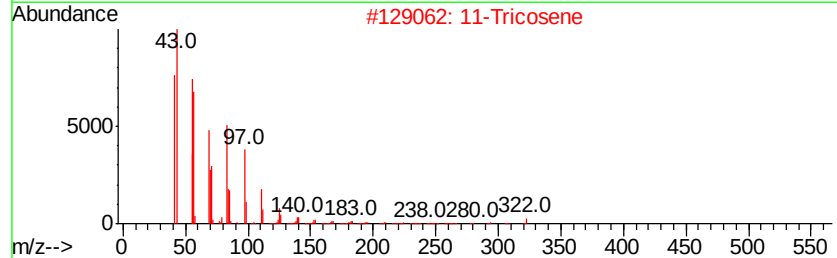
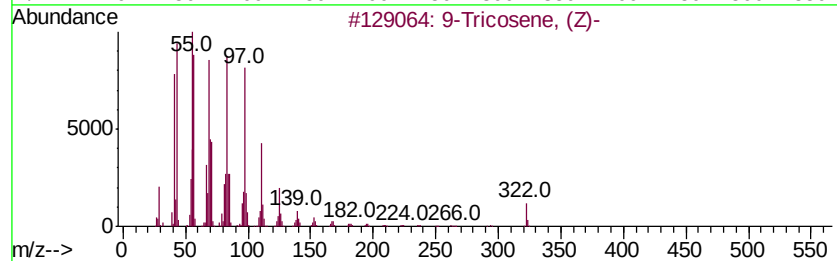
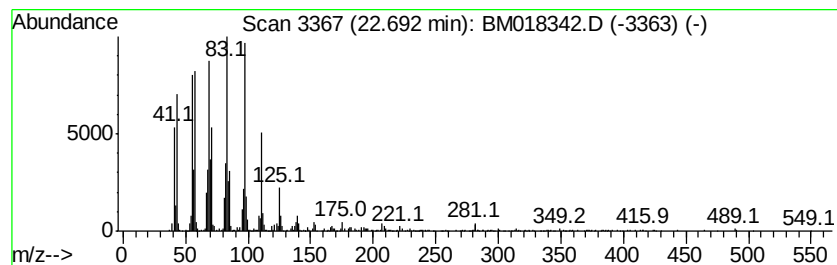
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 9-Tricosene, (Z)- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.69	16.33 ng	1264790	Perylene-d12	23.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9-Tricosene, (Z)-	322	C23H46	027519-02-4	87
2		11-Tricosene	322	C23H46	052078-56-5	87
3		1-Docosene	308	C22H44	001599-67-3	74
4		1-Tetracosanol	354	C24H50O	000506-51-4	58
5		Cyclopentane, 1,1'-[3-(2-cyclope...	304	C22H40	055255-85-1	52



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM122018\
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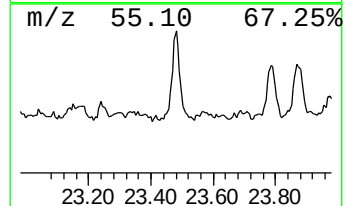
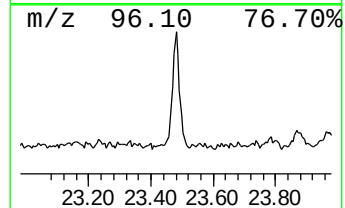
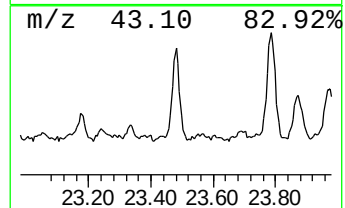
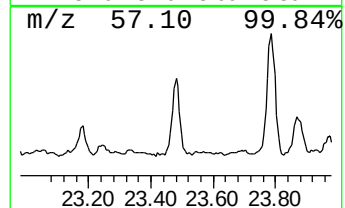
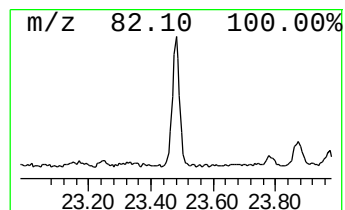
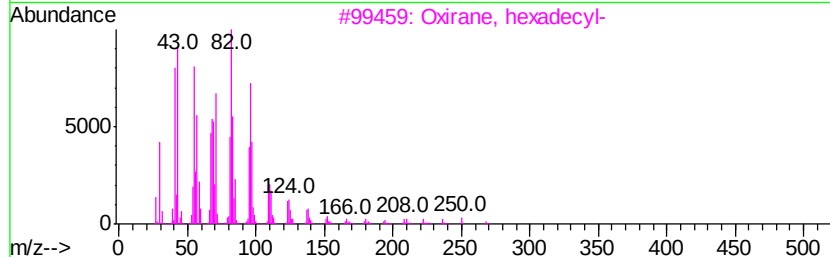
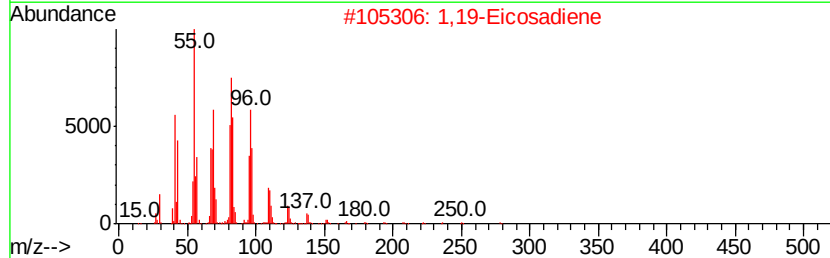
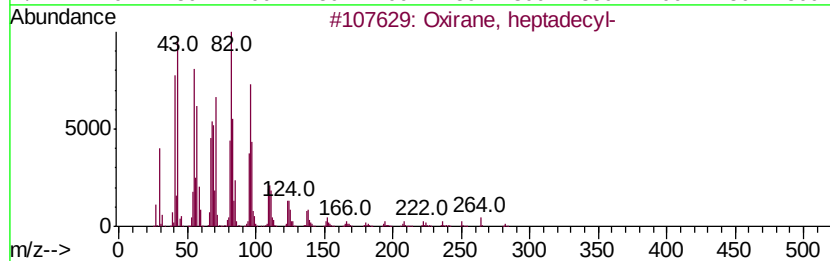
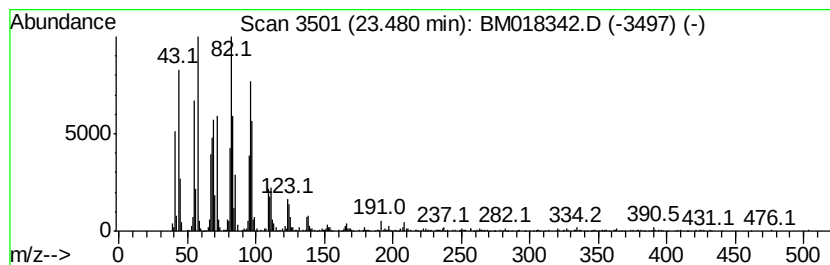
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ClientSampleId :
P001-SS016-0006-02

Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM121518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Oxirane, heptadecyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.48	18.97 ng	1469150	Perylene-d12	23.30
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Oxirane, heptadecyl-	282 C19H380	067860-04-2	96
2	1,19-Eicosadiene	278 C20H38	014811-95-1	94
3	Oxirane, hexadecyl-	268 C18H360	007390-81-0	87
4	Pentadecanal-	226 C15H300	002765-11-9	81
5	Tetradecanal	212 C14H280	000124-25-4	72



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM122018\
Data File : BM018342.D
Acq On : 20 Dec 2018 22:23
Operator : JU/SJ
Sample : J6386-03
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
P001-SS016-0006-02

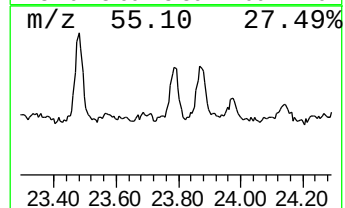
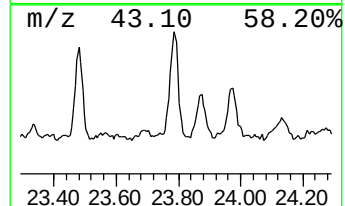
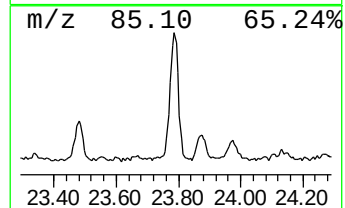
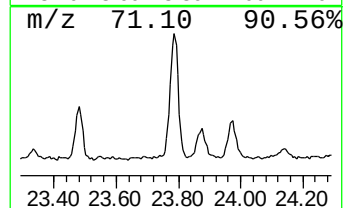
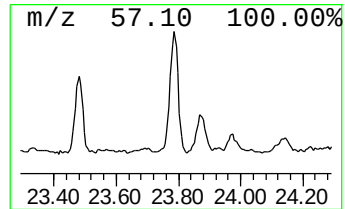
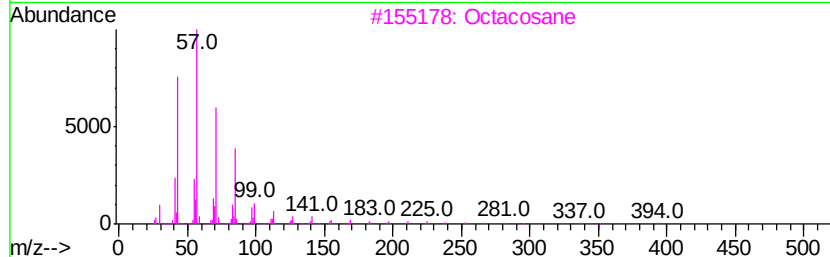
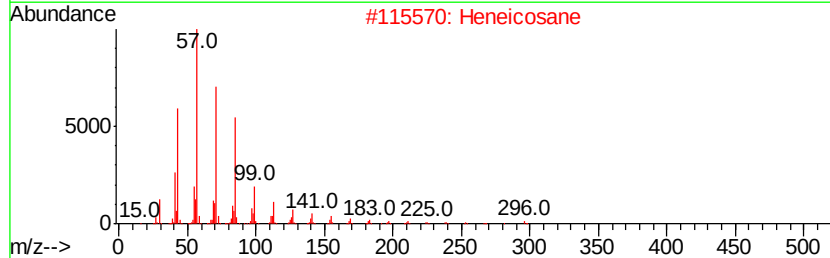
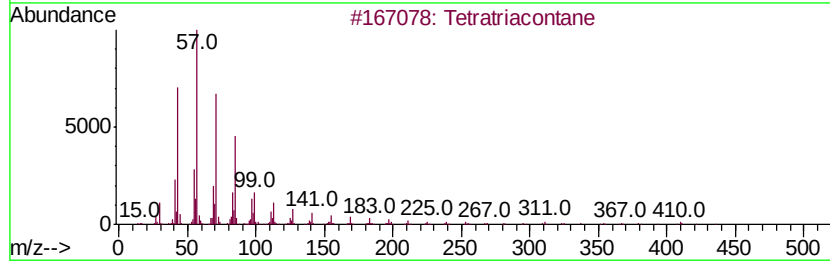
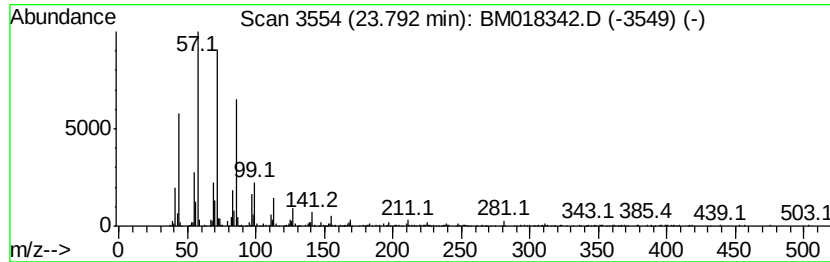
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Tetratriacontane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.79	16.28 ng	1260910	Perylene-d12	23.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetratriacontane	479	C34H70	014167-59-0	91
2		Heneicosane	296	C21H44	000629-94-7	90
3		Octacosane	394	C28H58	000630-02-4	86
4		Docosane, 11-decyl-	451	C32H66	055401-55-3	86
5		Triacontane	422	C30H62	000638-68-6	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM122018\
Data File : BM018342.D
Acq On : 20 Dec 2018 22:23
Operator : JU/SJ
Sample : J6386-03
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
P001-SS016-0006-02

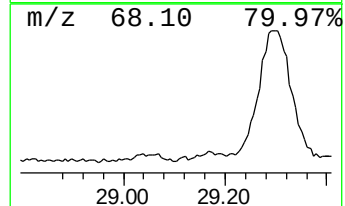
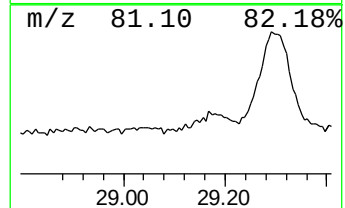
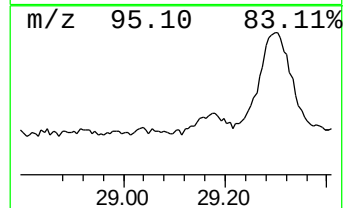
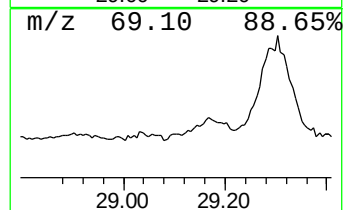
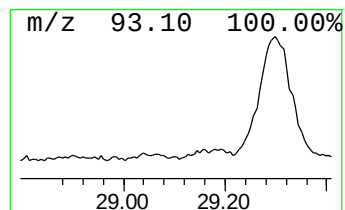
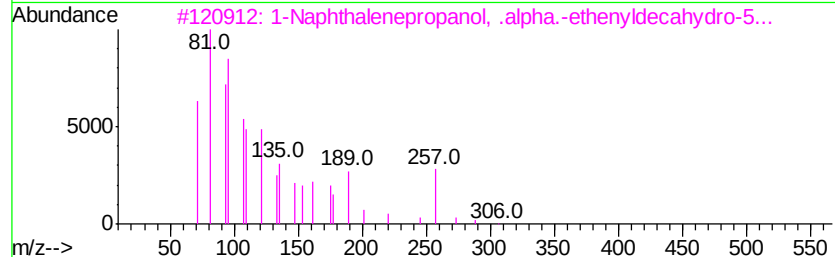
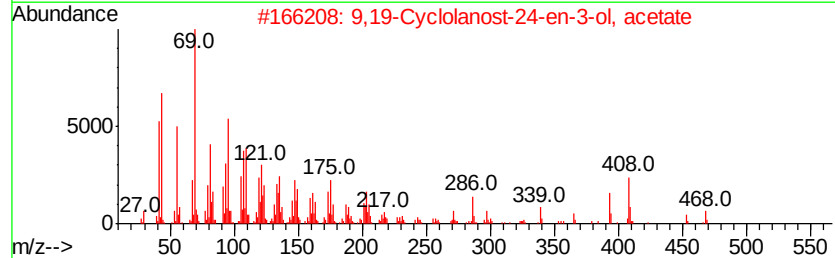
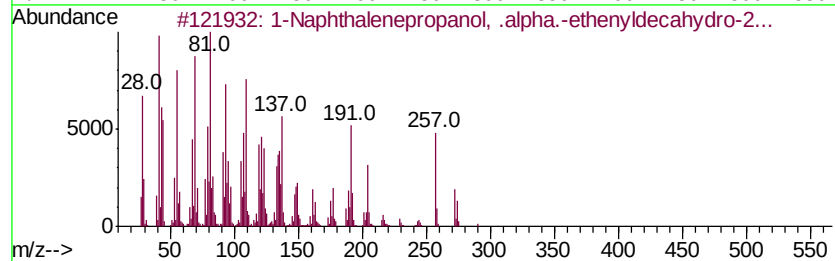
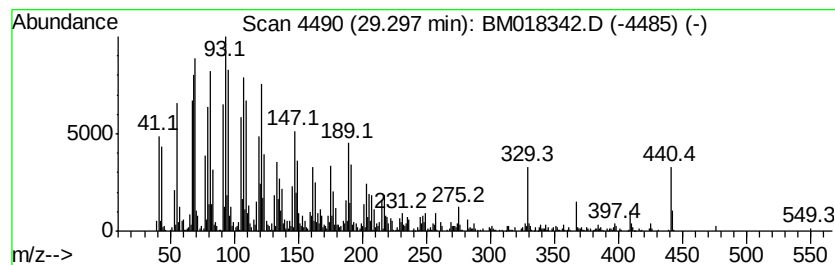
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 unknown29.30 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
29.30	68.65 ng	5318290	Perylene-d12	23.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Naphthalenepropanol, .alpha.-e...	308	C20H36O2	000515-03-7	38
2		9,19-Cyclolanost-24-en-3-ol, ace...	468	C32H52O2	124094-10-6	37
3		1-Naphthalenepropanol, .alpha.-e...	306	C20H34O2	003650-30-4	37
4		2,4,4-Trimethyl-3-hydroxymethyl-...	222	C15H26O	1000144-10-5	35
5		C(14a)-Homo-27-nor-14.beta.-gamm...	428	C30H52O	024739-08-0	35



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM122018\
 Data File : BM018342.D
 Acq On : 20 Dec 2018 22:23
 Operator : JU/SJ
 Sample : J6386-03
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 P001-SS016-0006-02

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\8270-BM121518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Butanone, 3-hyd...	3.11	11.9	ng	506412	1	7.49	855035	20.0
unknown7.03	7.03	106.8	ng	4566870	1	7.49	855035	20.0
13-Borabicyclo[7....	17.77	6.1	ng	359002	4	16.92	1184400	20.0
n-Hexadecanoic acid	17.83	16.8	ng	991762	4	16.92	1184400	20.0
Cyclopentadecane	19.86	4.5	ng	427205	5	21.14	1896360	20.0
Cyclodisilazane-2...	20.85	322.7	ng	30600900	5	21.14	1896360	20.0
Oxirane, hexadecyl-	21.47	3.4	ng	324039	5	21.14	1896360	20.0
Heneicosane, 11-(...	21.72	18.4	ng	1746460	5	21.14	1896360	20.0
1,2-Dodecanediol	21.74	23.1	ng	2190520	5	21.14	1896360	20.0
Tetradecanal	22.39	9.2	ng	712346	6	23.30	1549320	20.0
Docosane, 11-butyl-	22.64	20.5	ng	1590480	6	23.30	1549320	20.0
9-Tricosene, (Z)-	22.69	16.3	ng	1264790	6	23.30	1549320	20.0
Oxirane, heptadecyl-	23.48	19.0	ng	1469150	6	23.30	1549320	20.0
Tetratriacontane	23.79	16.3	ng	1260910	6	23.30	1549320	20.0
unknown29.30	29.30	68.7	ng	5318290	6	23.30	1549320	20.0