

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121818\
 Data File : BM018285.D
 Acq On : 18 Dec 2018 23:36
 Operator : JU/SJ
 Sample : J6389-04
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 P001-SS012-1824-01

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	3.357	77	80	88	rBV	55181	75600	1.53%	0.157%
2	4.681	299	305	312	rBV	96881	133221	2.69%	0.276%
3	5.122	373	380	392	rBV	2790281	3849309	77.73%	7.981%
4	6.692	639	647	661	rBV	2568792	4272062	86.26%	8.858%
5	7.028	696	704	718	rBV	2840144	4329897	87.43%	8.978%
6	7.486	775	782	789	rBV	626800	977657	19.74%	2.027%
7	7.798	827	835	847	rBV	2061124	3182304	64.26%	6.598%
8	8.645	972	979	998	rBV	1519725	2545337	51.40%	5.278%
9	10.251	1245	1252	1270	rBV	724538	1277835	25.80%	2.650%
10	12.751	1670	1677	1689	rBV	3450123	4952449	100.00%	10.269%
11	13.615	1818	1824	1839	rBV	326290	512963	10.36%	1.064%
12	14.145	1906	1914	1923	rBV2	990703	1495330	30.19%	3.101%
13	15.651	2164	2170	2180	rBV	2332989	3366054	67.97%	6.979%
14	15.880	2204	2209	2219	rBV5	28787	69051	1.39%	0.143%
15	16.909	2377	2384	2393	rBV	965378	1479274	29.87%	3.067%
16	17.762	2525	2529	2534	rBV5	44218	59088	1.19%	0.123%
17	17.827	2534	2540	2550	rBV	747972	1123644	22.69%	2.330%
18	18.974	2731	2735	2736	rBV3	46029	63460	1.28%	0.132%
19	19.121	2755	2760	2767	rBV	339495	509582	10.29%	1.057%
20	19.345	2795	2798	2801	rVB4	45629	53405	1.08%	0.111%
21	19.386	2801	2805	2810	rBV4	62059	103941	2.10%	0.216%
22	19.562	2830	2835	2841	rVB	3632893	4342195	87.68%	9.003%
23	19.844	2879	2883	2887	rBV4	137625	222699	4.50%	0.462%
24	20.127	2928	2931	2935	rVB	143900	169002	3.41%	0.350%
25	20.380	2971	2974	2978	rBV	120050	147660	2.98%	0.306%
26	20.591	3007	3010	3013	rBV4	88650	126344	2.55%	0.262%
27	20.850	3050	3054	3061	rBV	928650	1193019	24.09%	2.474%
28	21.127	3097	3101	3112	rBV2	1087529	1689557	34.12%	3.503%
29	21.291	3126	3129	3137	rVB2	246401	266808	5.39%	0.553%
30	21.715	3197	3201	3203	rBV	511521	606453	12.25%	1.257%
31	21.738	3203	3205	3214	rVB2	441576	559111	11.29%	1.159%
32	22.156	3273	3276	3281	rVB	453394	542989	10.96%	1.126%
33	22.638	3354	3358	3363	rBV	564337	800704	16.17%	1.660%
34	22.691	3364	3367	3374	rVB3	219585	352923	7.13%	0.732%

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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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35	23.180	3445	3450	3456	rVB	392560	619839	12.52%	1.285%
36	23.285	3463	3468	3475	rVB	815594	1462536	29.53%	3.033%
37	23.780	3548	3552	3558	rVB	406630	695024	14.03%	1.441%

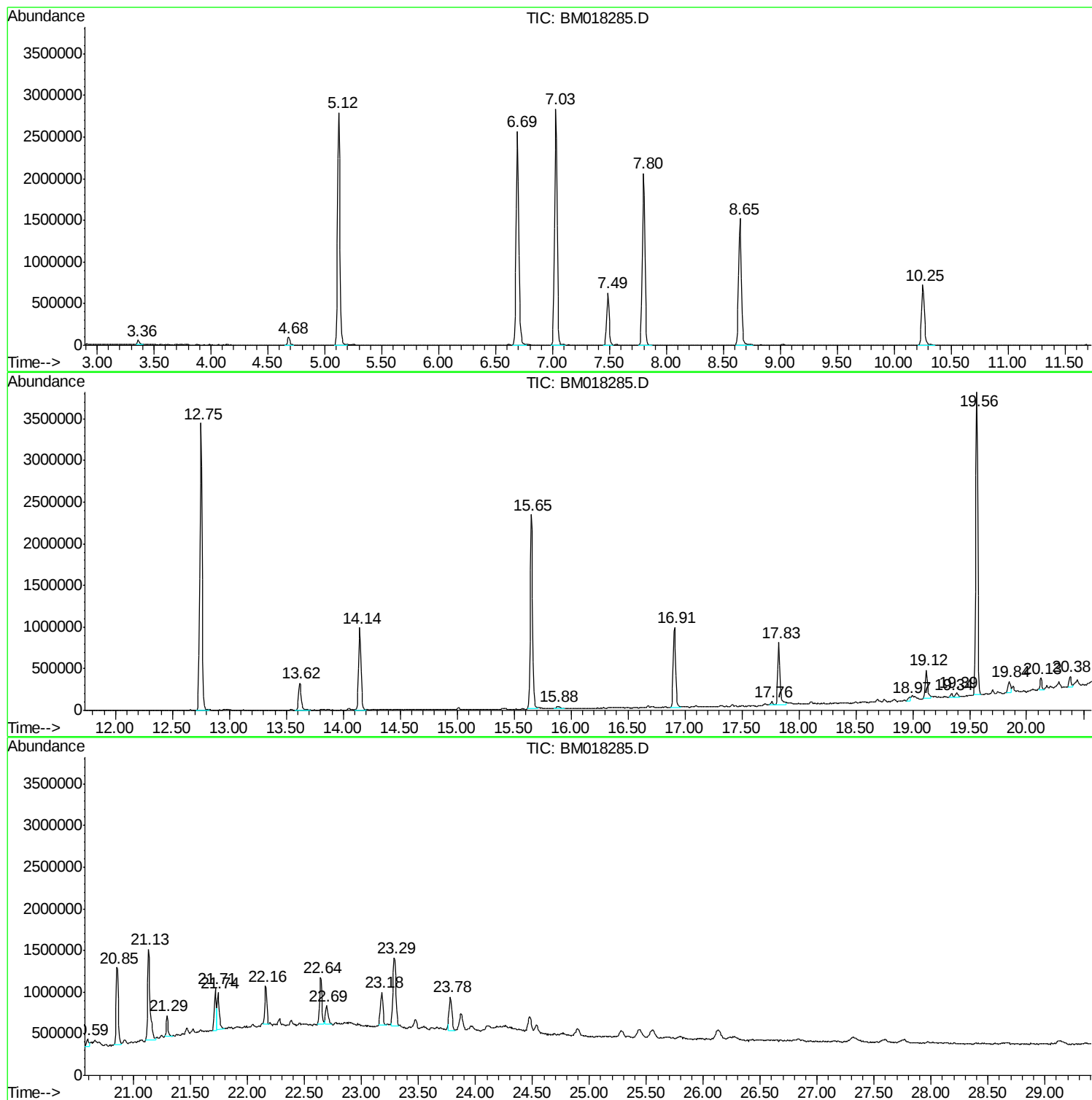
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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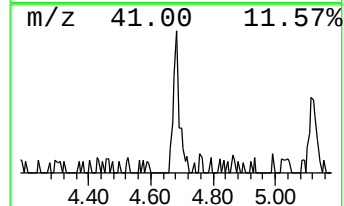
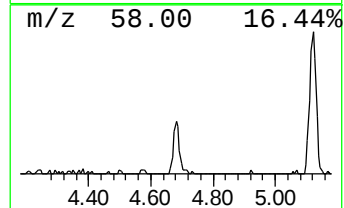
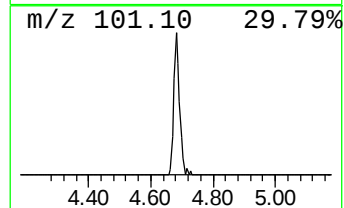
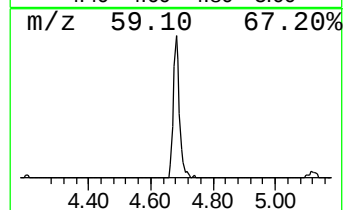
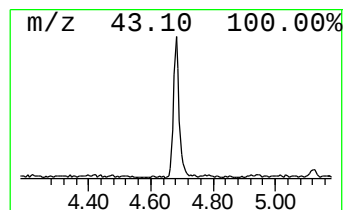
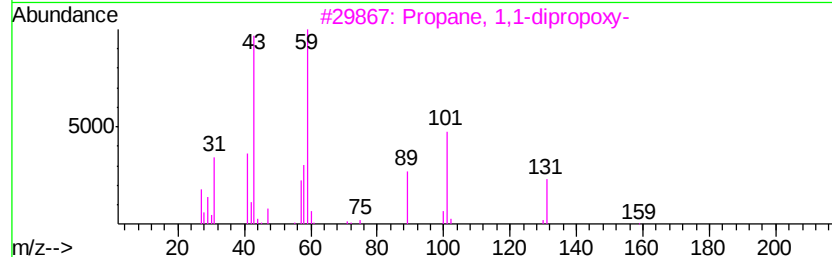
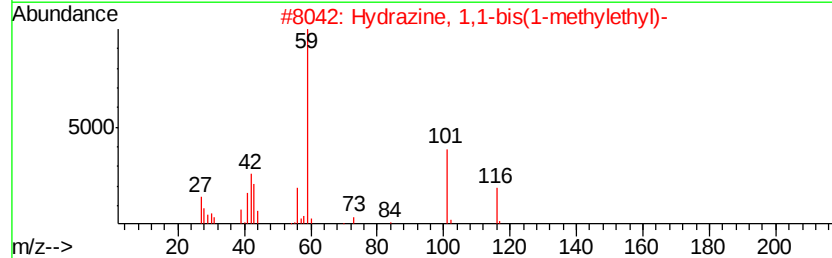
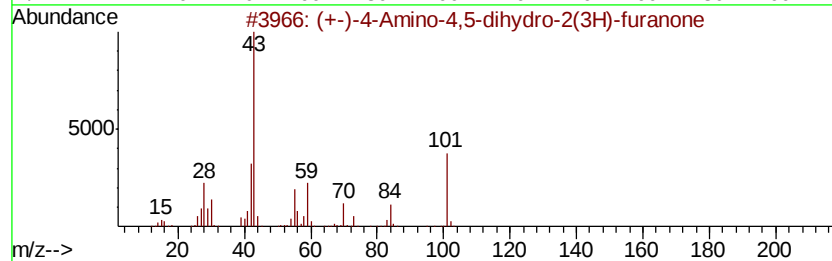
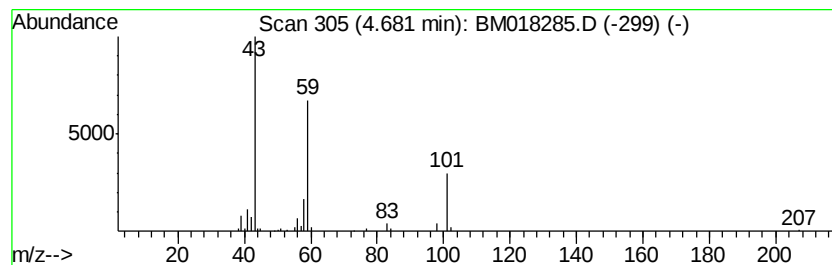
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown4.68 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.68	2.73 ng	133221	1,4-Dichlorobenzene-d4	7.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	(+)-	4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	47
2		Hydrazine, 1,1-bis(1-methylethyl)-	116	C6H16N2	000921-14-2	47
3		Propane, 1,1-dipropoxy-	160	C9H20O2	004744-11-0	36
4		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	28
5		3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	27



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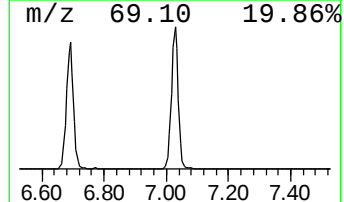
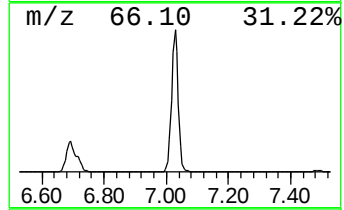
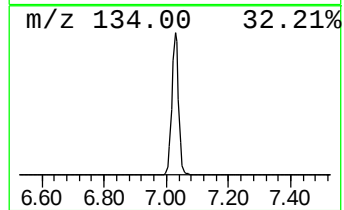
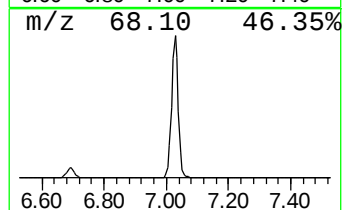
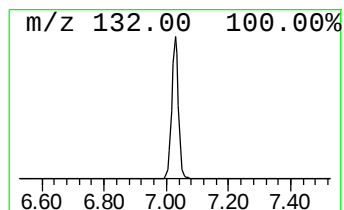
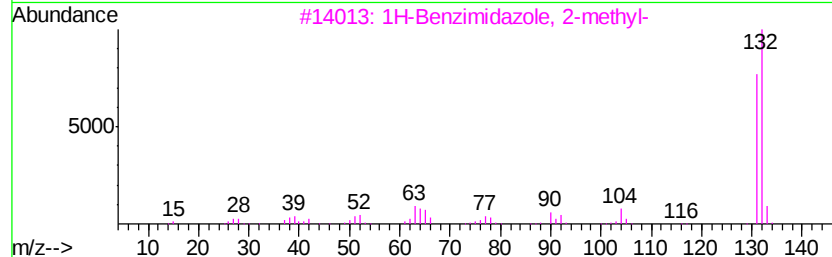
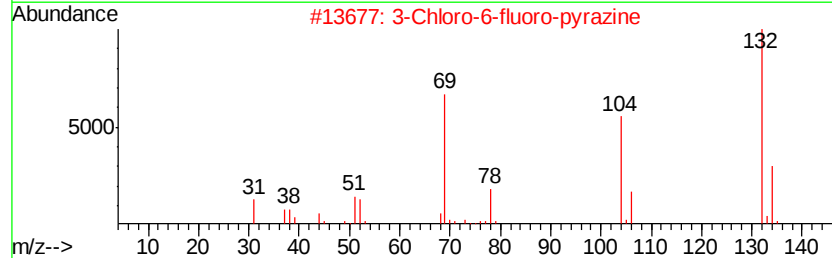
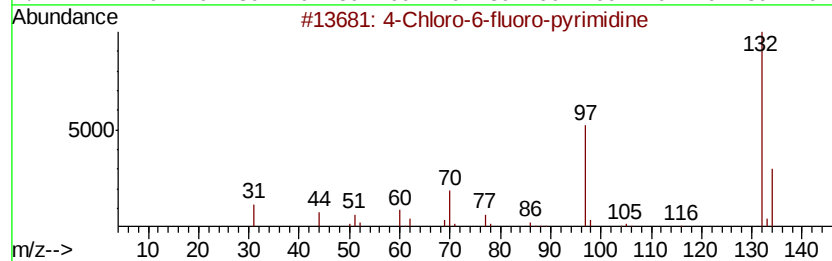
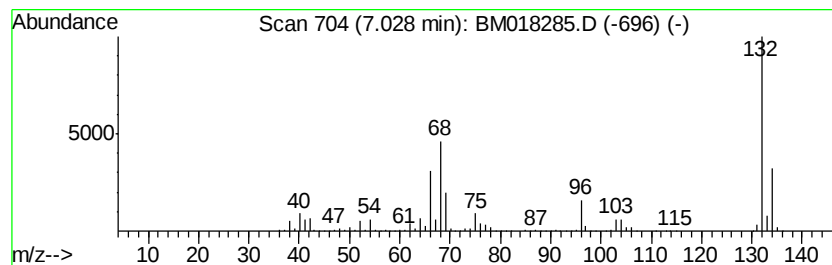
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 2 unknown7.03 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.03	88.58 ng	4329900	1,4-Dichlorobenzene-d4	7.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
5		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	10



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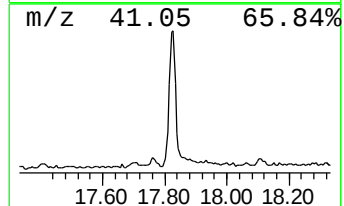
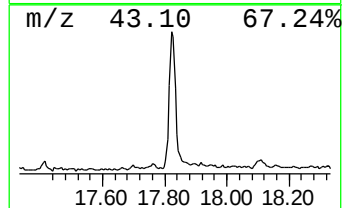
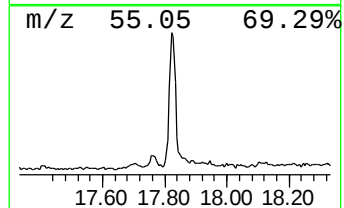
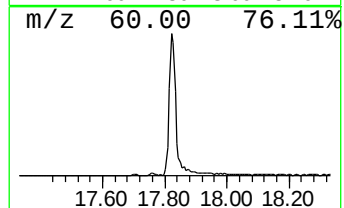
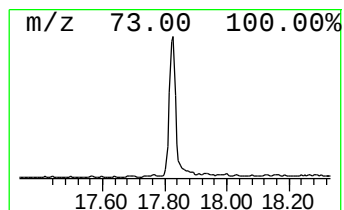
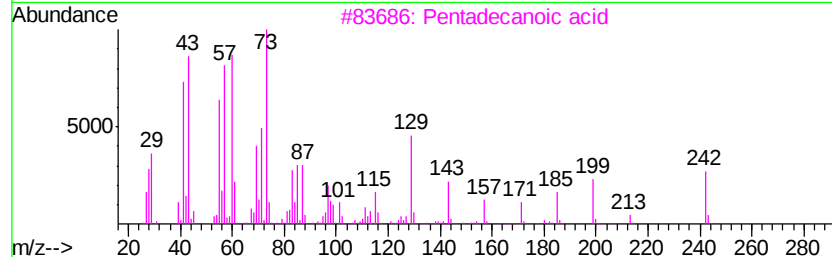
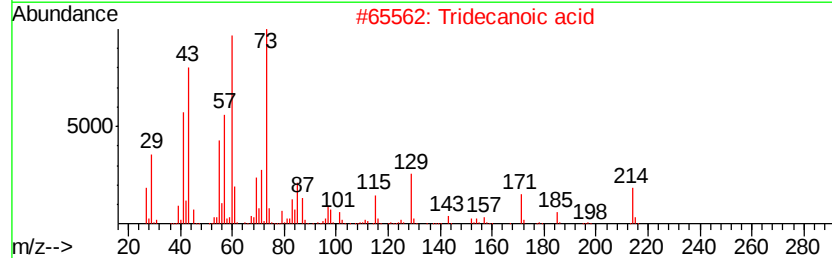
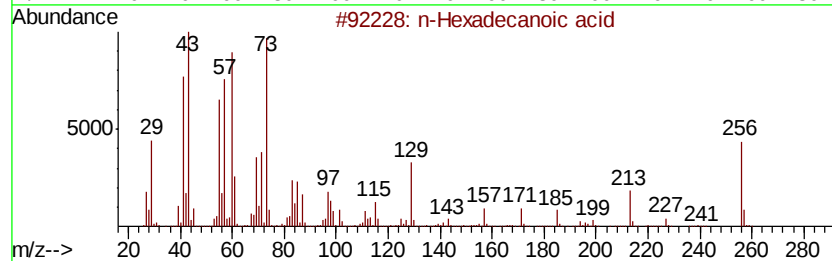
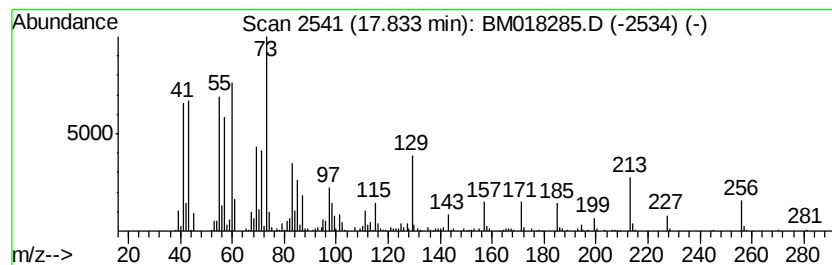
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD		R.T.
17.83	15.19 ng	1123640	Phenanthrene-d10		16.91
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	68
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	68
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	53
5	Dodecanoic acid	200	C12H24O2	000143-07-7	43



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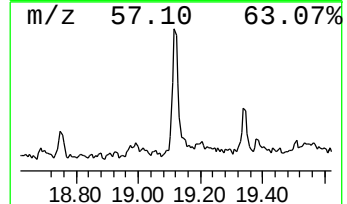
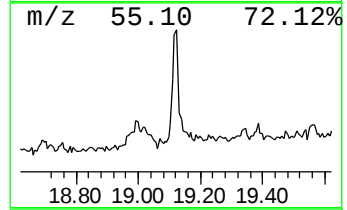
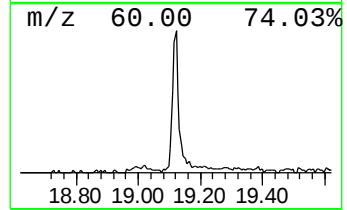
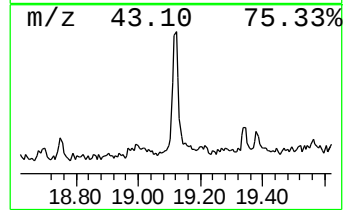
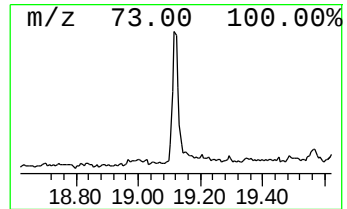
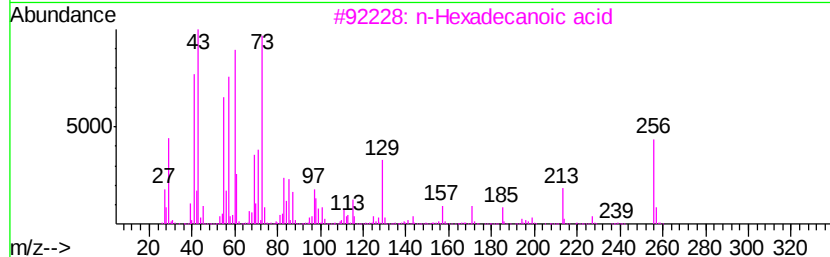
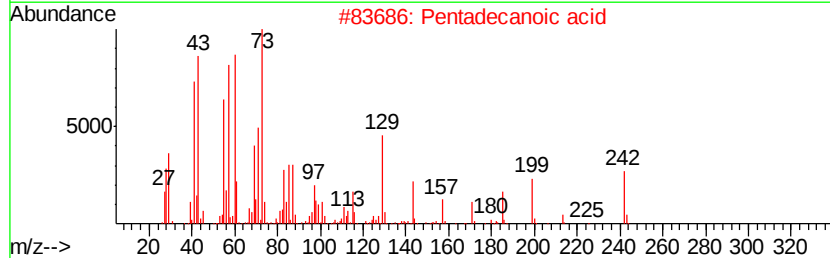
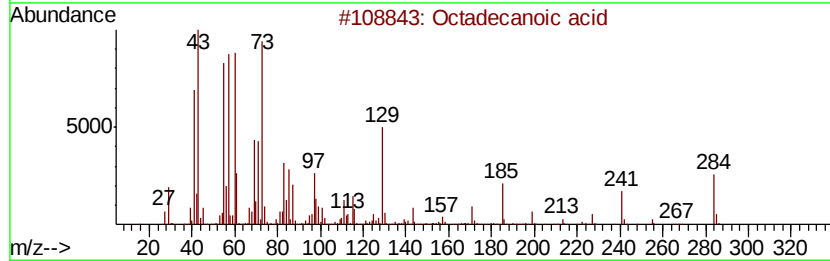
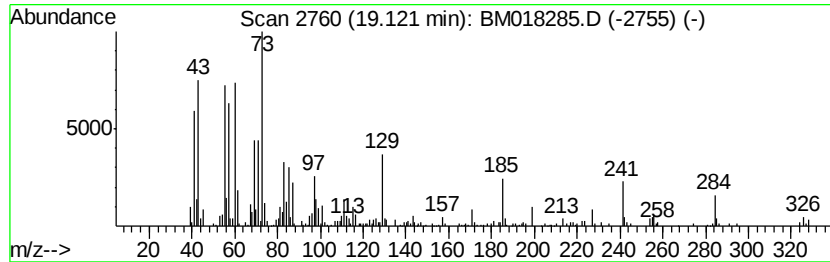
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Octadecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD		R.T.		
19.12	6.03 ng	509582	Chrysene-d12		21.13		
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	96
2			Pentadecanoic acid	242	C15H30O2	001002-84-2	90
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	89
4			Tetradecanoic acid	228	C14H28O2	000544-63-8	72
5			Octadecanoic acid, 2-(2-hydroxy...	372	C22H44O4	000106-11-6	72



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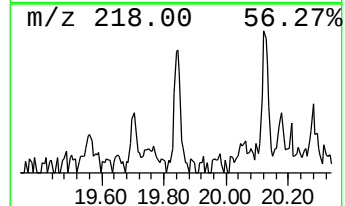
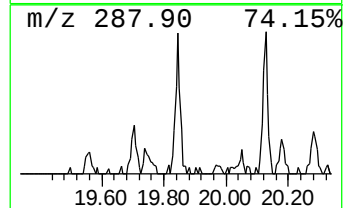
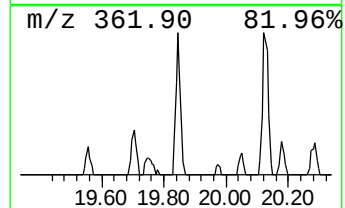
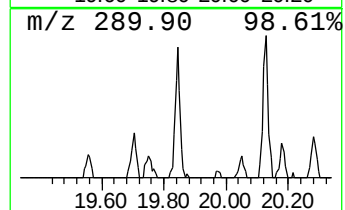
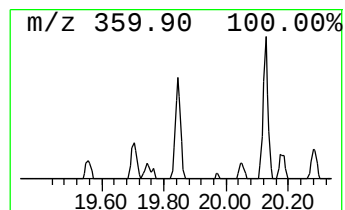
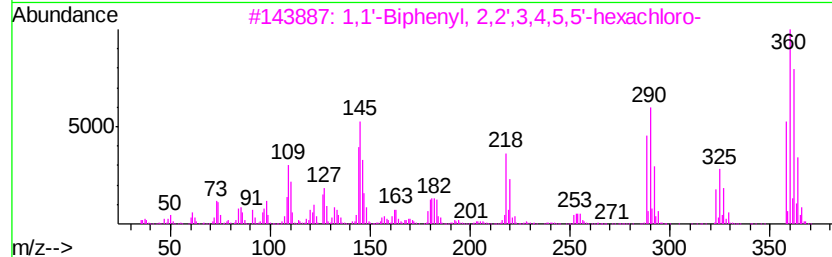
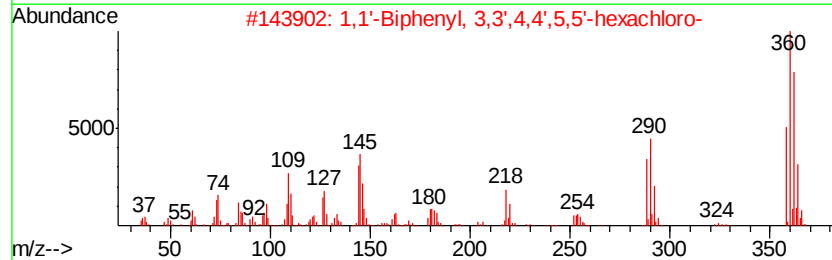
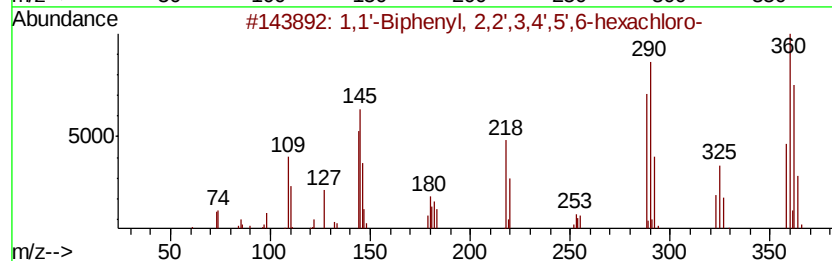
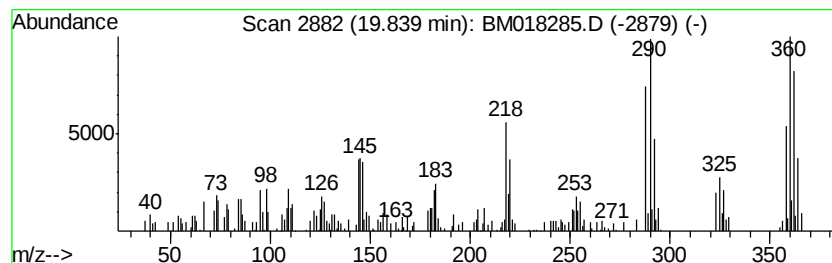
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 1,1'-Biphenyl, 2,2',3,4',5'... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.84	2.64 ng	222699	Chrysene-d12	21.13

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1'-Biphenyl, 2,2',3,4',5',6-he...	358	C12H4Cl6	038380-04-0	99
2		1,1'-Biphenyl, 3,3',4,4',5,5'-he...	358	C12H4Cl6	032774-16-6	99
3		1,1'-Biphenyl, 2,2',3,4,5,5'-hex...	358	C12H4Cl6	052712-04-6	99
4		1,1'-Biphenyl, 2,3,3',4,4',5-hex...	358	C12H4Cl6	038380-08-4	99
5		1,1'-Biphenyl, 2,2',3,3',4,4'-he...	358	C12H4Cl6	038380-07-3	98



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121818\
Data File : BM018285.D
Acq On : 18 Dec 2018 23:36
Operator : JU/SJ
Sample : J6389-04
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
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ClientSampleId :
P001-SS012-1824-01

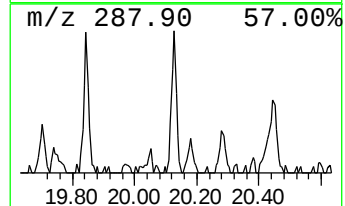
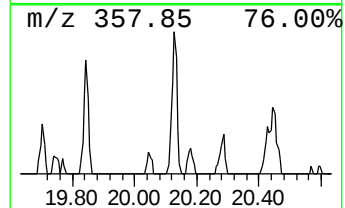
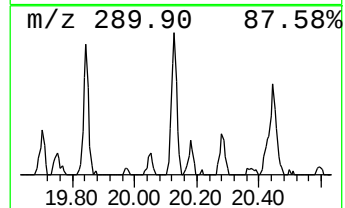
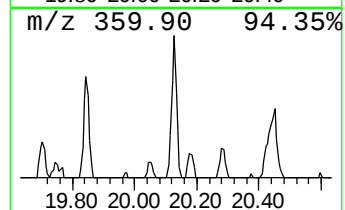
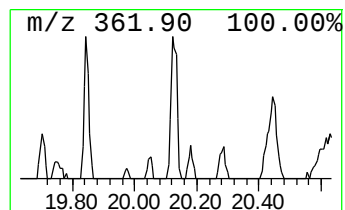
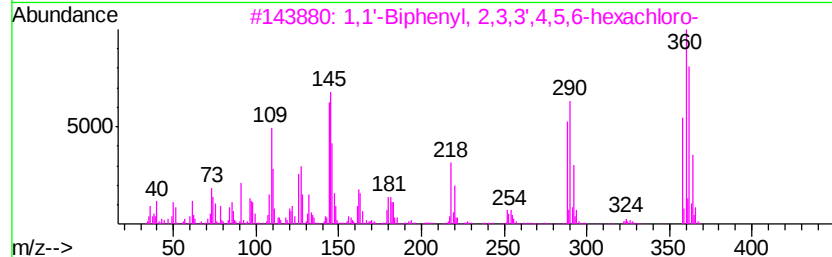
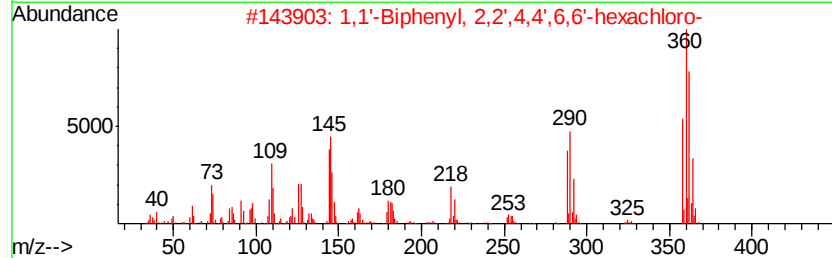
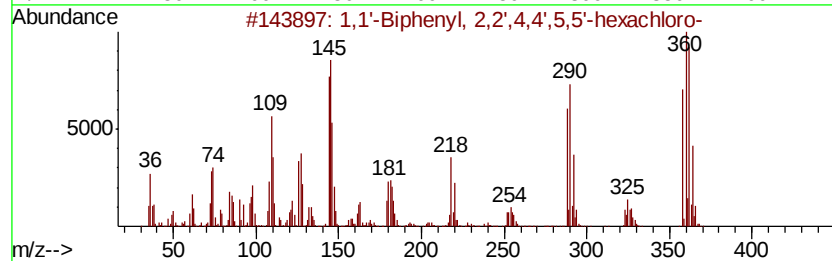
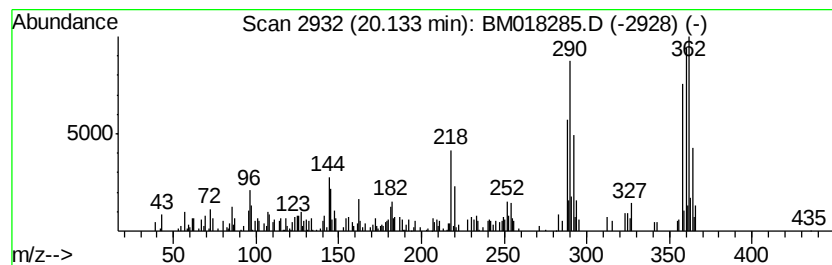
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 1,1'-Biphenyl, 2,2',4,4',5,... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.13	2.00 ng	169002	Chrysene-d12	21.13

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1'-Biphenyl, 2,2',4,4',5,5'-he...	358	C12H4Cl6	035065-27-1	99
2		1,1'-Biphenyl, 2,2',4,4',6,6'-he...	358	C12H4Cl6	033979-03-2	99
3		1,1'-Biphenyl, 2,3,3',4,5,6-hexa...	358	C12H4Cl6	041411-62-5	99
4		1,1'-Biphenyl, 3,3',4,4',5,5'-he...	358	C12H4Cl6	032774-16-6	99
5		1,1'-Biphenyl, 2,2',3,3',4,4'-he...	358	C12H4Cl6	038380-07-3	96



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121818\
Data File : BM018285.D
Acq On : 18 Dec 2018 23:36
Operator : JU/SJ
Sample : J6389-04
Misc :
ALS Vial : 16 Sample Multiplier: 1

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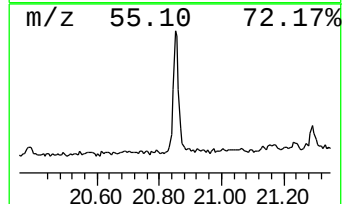
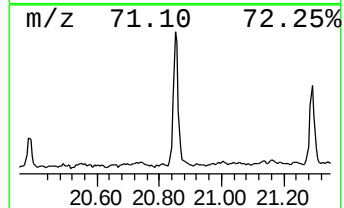
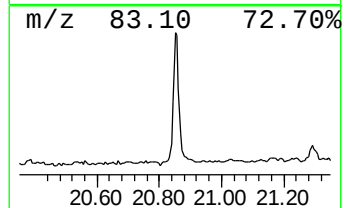
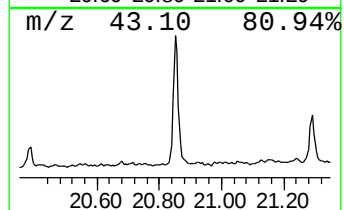
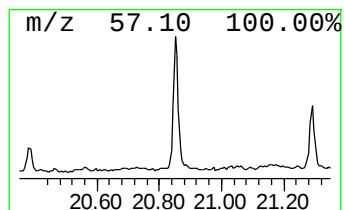
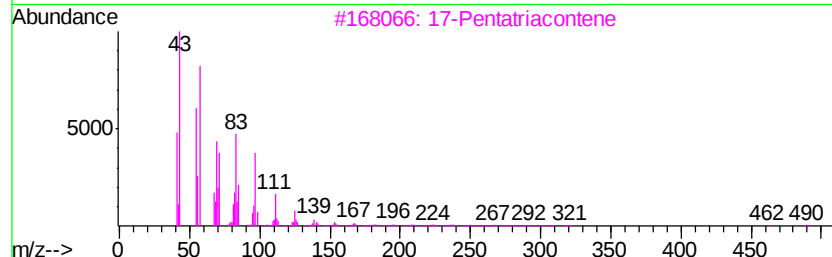
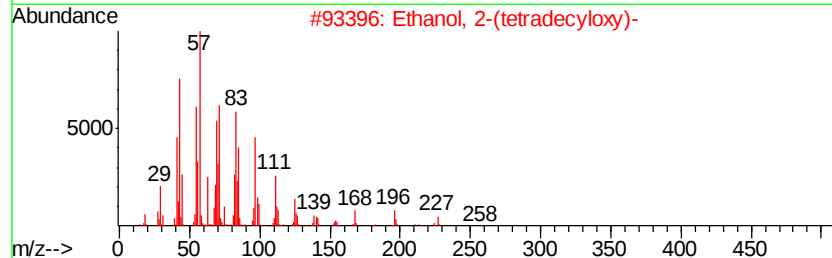
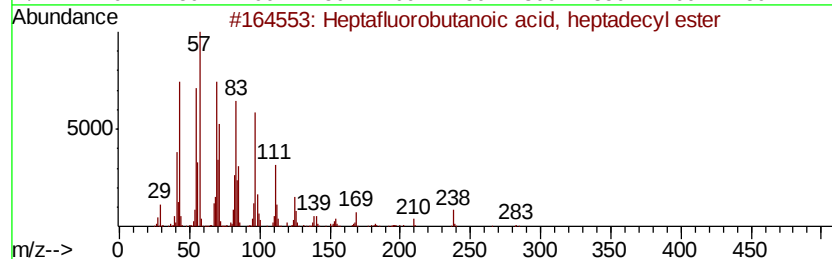
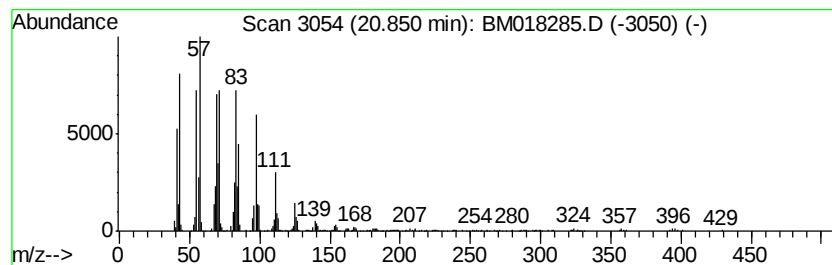
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Heptafluorobutanoic acid, h... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.85	14.12 ng	1193020	Chrysene-d12	21.13

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptafluorobutanoic acid, heptad...	452	C21H35F7O2	1000282-97-3	91
2		Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	91
3		17-Pentatriacontene	491	C35H70	006971-40-0	91
4		1-Hexadecene	224	C16H32	000629-73-2	86
5		1-Decanol, 2-hexyl-	242	C16H34O	002425-77-6	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121818\
Data File : BM018285.D
Acq On : 18 Dec 2018 23:36
Operator : JU/SJ
Sample : J6389-04
Misc :
ALS Vial : 16 Sample Multiplier: 1

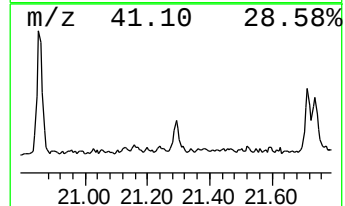
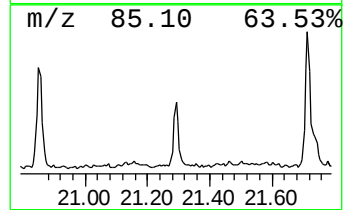
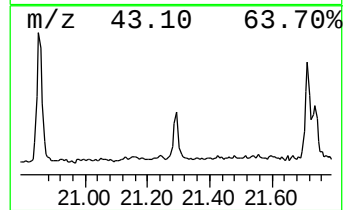
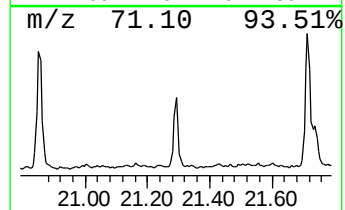
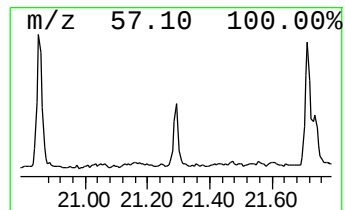
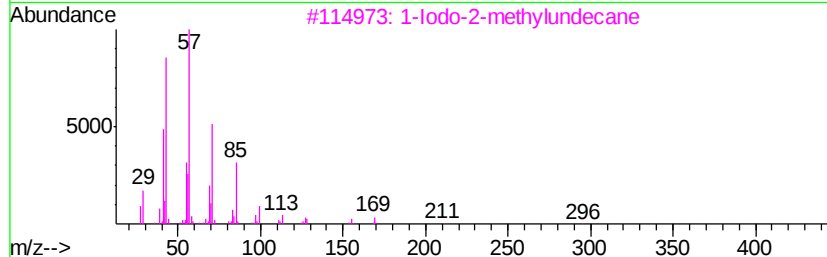
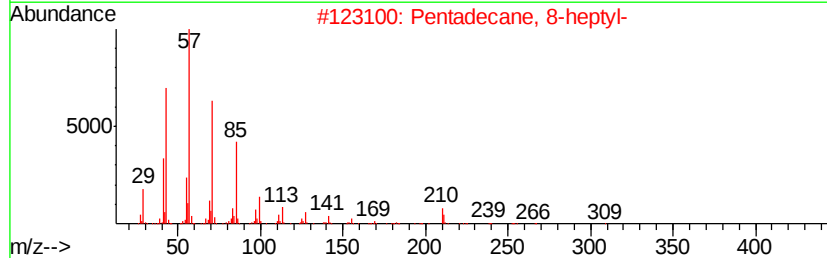
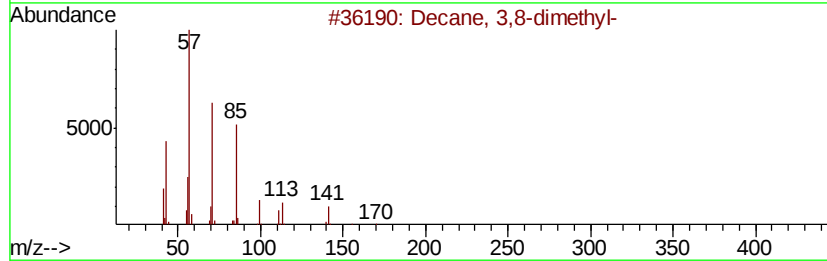
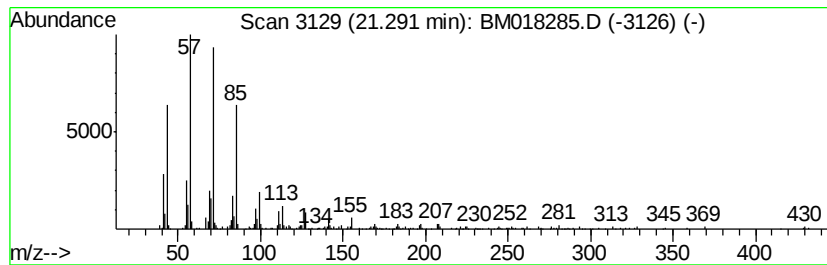
Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM121518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Decane, 3,8-dimethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.		
21.29	3.16 ng	266808	Chrysene-d12	21.13		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decane, 3,8-dimethyl-		170	C12H26	017312-55-9	86
2	Pentadecane, 8-heptyl-		310	C22H46	071005-15-7	83
3	1-Iodo-2-methylundecane		296	C12H25I	073105-67-6	81
4	Heptadecane, 2-methyl-		254	C18H38	001560-89-0	81
5	Octadecane		254	C18H38	000593-45-3	80



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121818\
Data File : BM018285.D
Acq On : 18 Dec 2018 23:36
Operator : JU/SJ
Sample : J6389-04
Misc :
ALS Vial : 16 Sample Multiplier: 1

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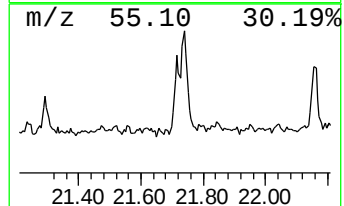
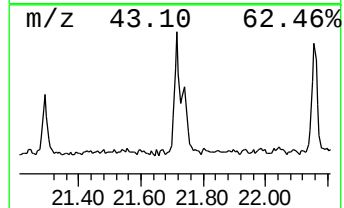
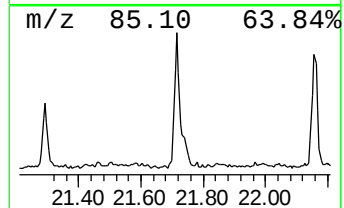
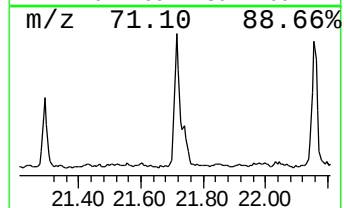
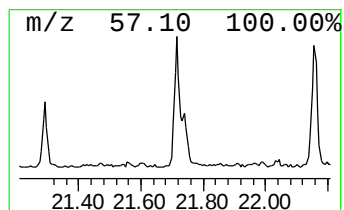
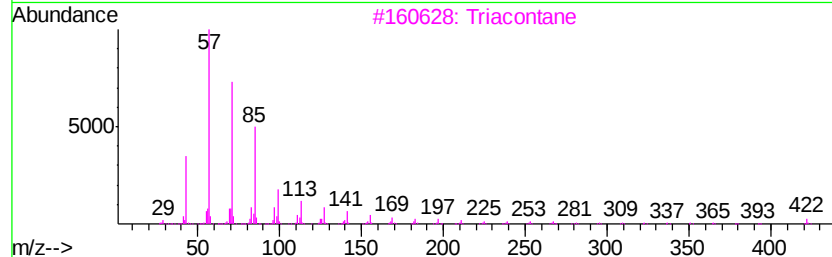
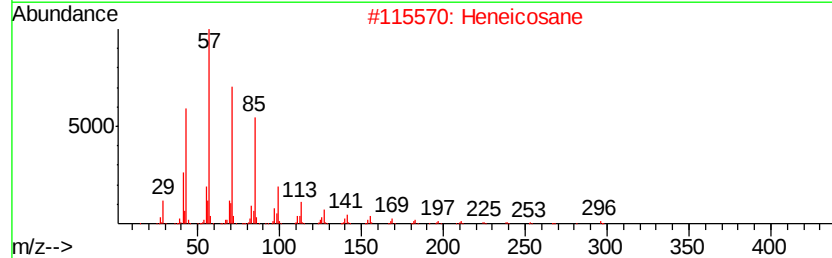
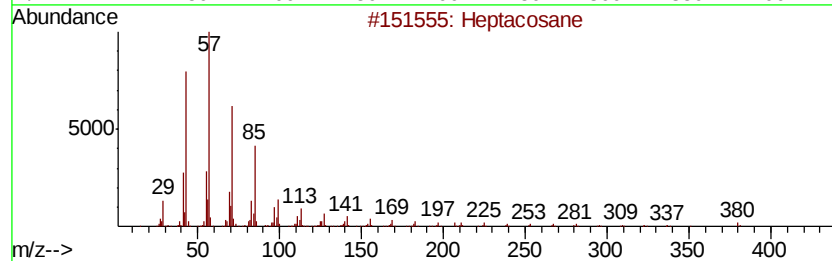
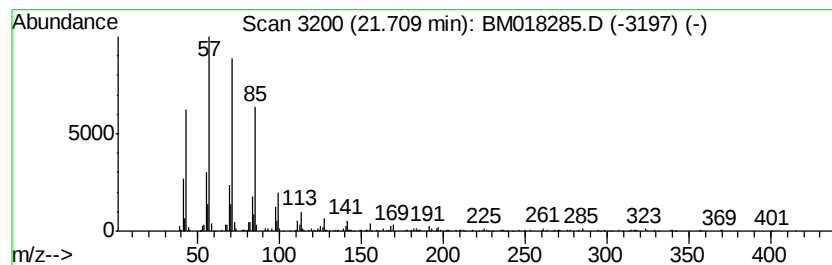
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Heptacosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.71	7.18 ng	606453	Chrysene-d12	21.13

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptacosane	380	C27H56	000593-49-7	91
2		Heneicosane	296	C21H44	000629-94-7	91
3		Triacontane	422	C30H62	000638-68-6	91
4		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
5		Eicosane	282	C20H42	000112-95-8	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121818\
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Misc :
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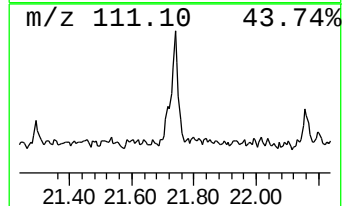
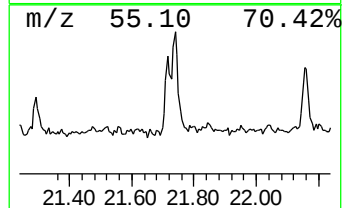
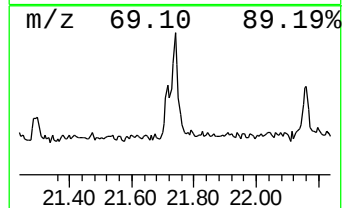
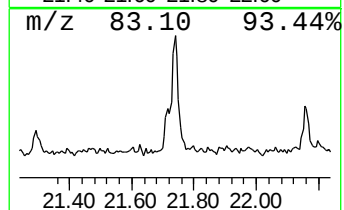
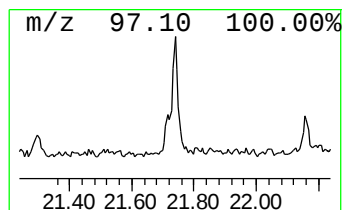
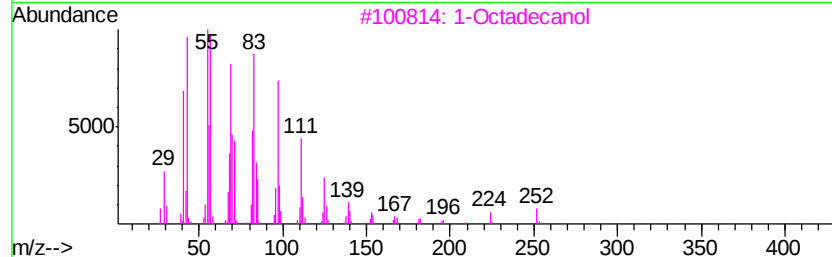
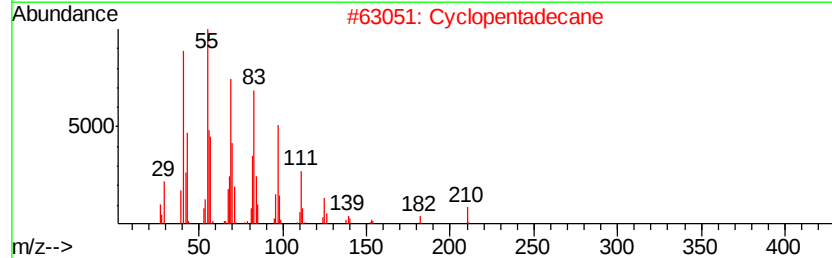
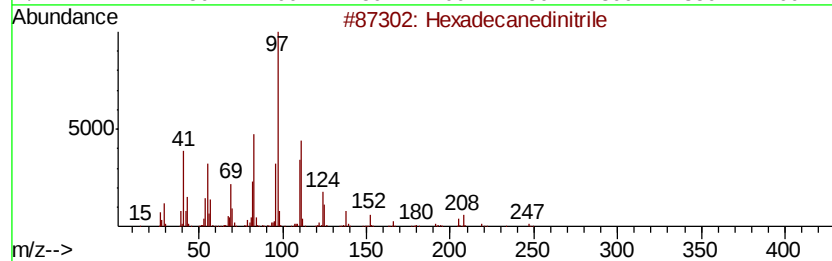
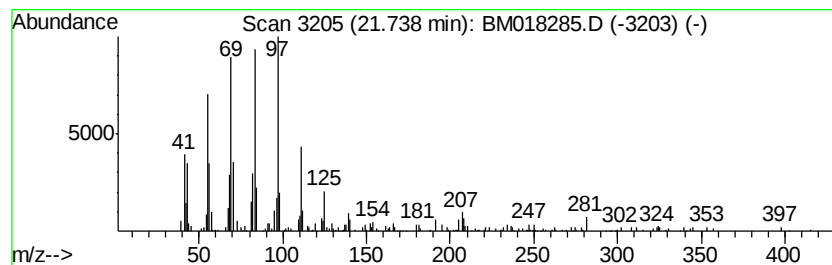
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Hexadecanedinitrile Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD		R.T.
21.74	6.62 ng	559111	Chrysene-d12		21.13
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecanedinitrile	248	C16H28N2	006812-44-8	43
2	Cyclopentadecane	210	C15H30	000295-48-7	41
3	1-Octadecanol	270	C18H38O	000112-92-5	38
4	Cycloheptane, methyl-	112	C8H16	004126-78-7	30
5	Cyclopentane, 1-pentyl-2-propyl-	182	C13H26	062199-51-3	25



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121818\
Data File : BM018285.D
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Operator : JU/SJ
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ALS Vial : 16 Sample Multiplier: 1

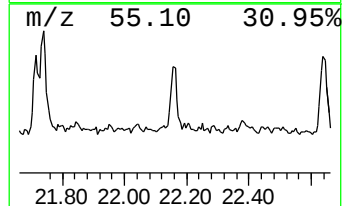
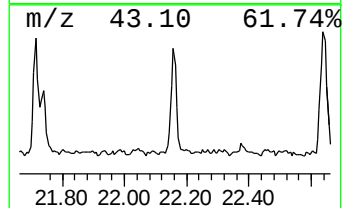
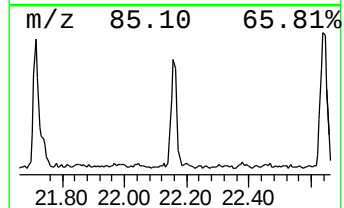
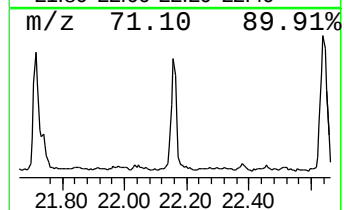
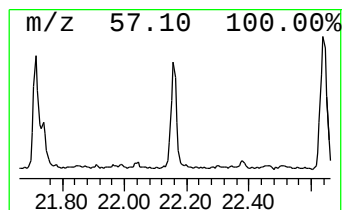
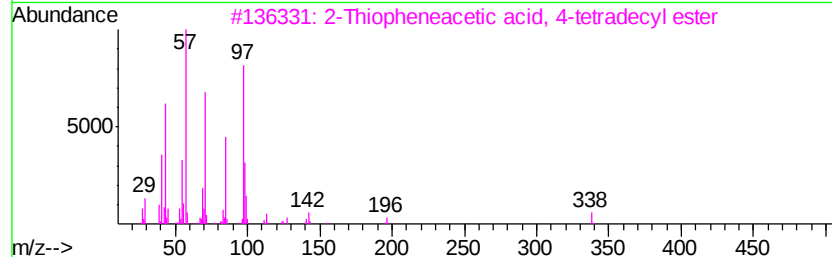
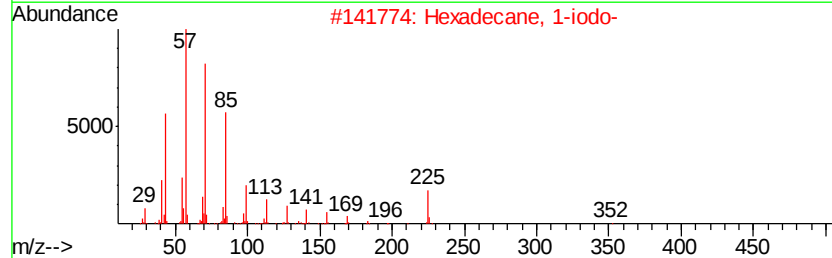
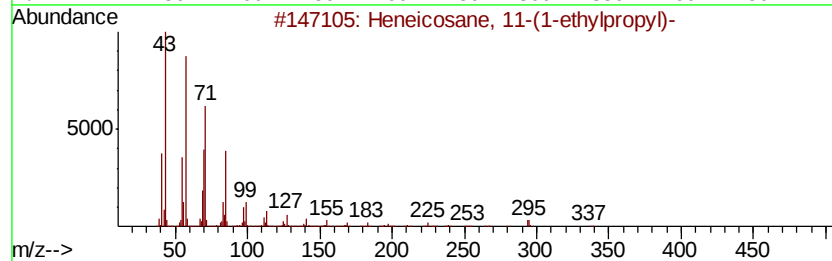
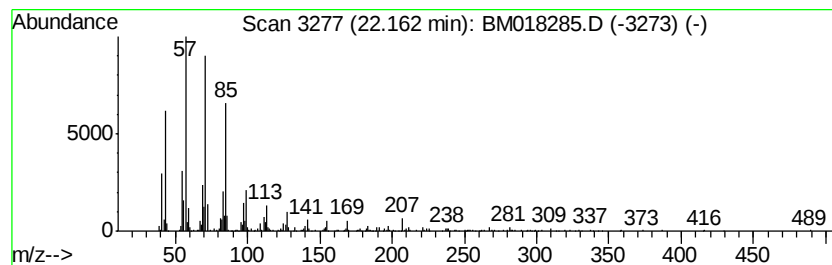
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM121518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Heneicosane, 11-(1-ethylpro... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
22.16	6.43 ng	542989	Chrysene-d12	21.13	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	83
2	Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	83
3	2-Thiopheneacetic acid, 4-tetrad...	338	C20H34O2S	1000280-67-0	80
4	Octadecane, 1-iodo-	380	C18H37I	000629-93-6	80
5	Heptacosane	380	C27H56	000593-49-7	74



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121818\
Data File : BM018285.D
Acq On : 18 Dec 2018 23:36
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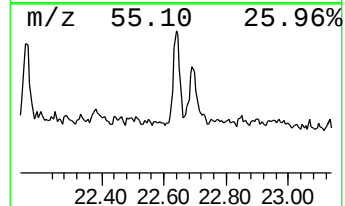
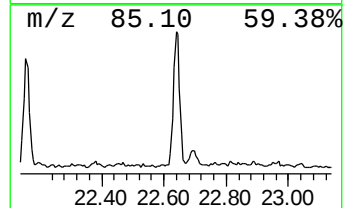
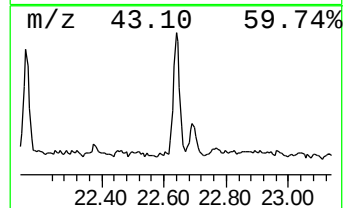
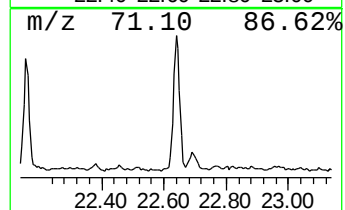
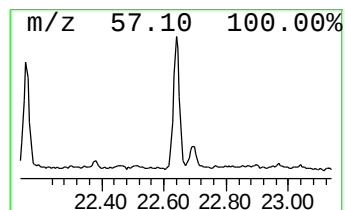
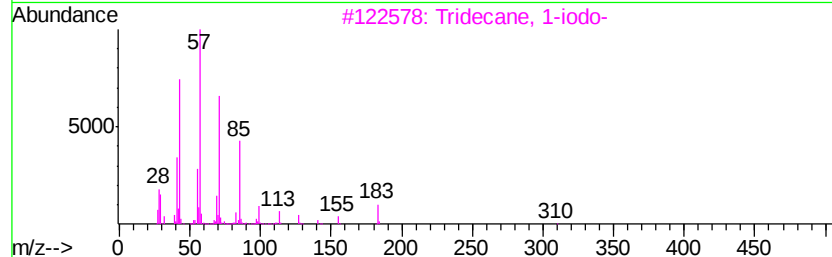
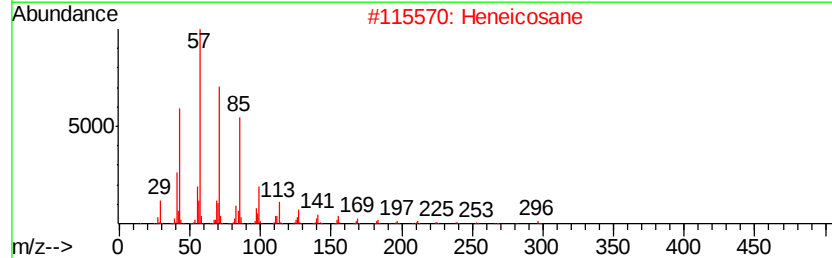
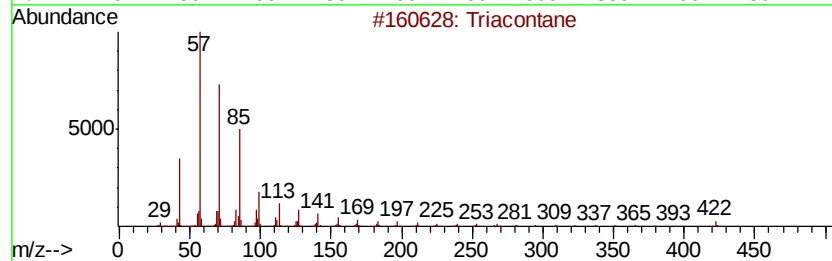
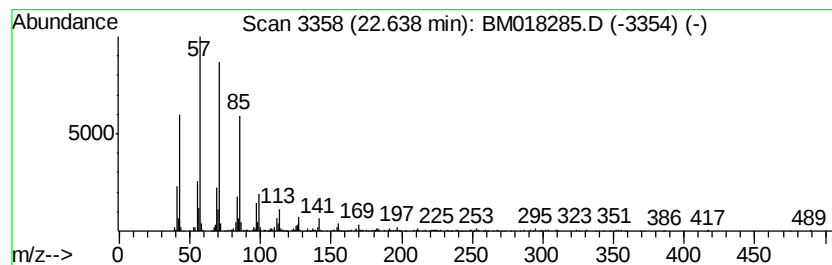
Instrument :
BNA_M
ClientSampleId :
P001-SS012-1824-01

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 13 Triacontane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
22.64	10.95 ng	800704	Perylene-d12	23.29	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Triacontane	422	C30H62	000638-68-6	91
2	Heneicosane	296	C21H44	000629-94-7	90
3	Tridecane, 1-iodo-	310	C13H27I	035599-77-0	87
4	Nonacosane	408	C29H60	000630-03-5	86
5	Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121818\
Data File : BM018285.D
Acq On : 18 Dec 2018 23:36
Operator : JU/SJ
Sample : J6389-04
Misc :
ALS Vial : 16 Sample Multiplier: 1

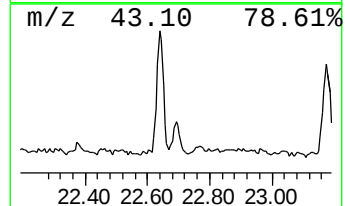
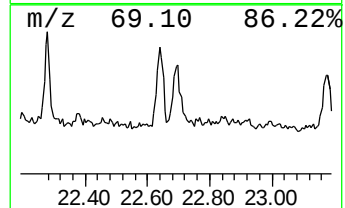
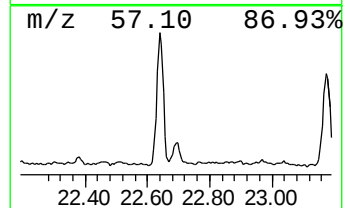
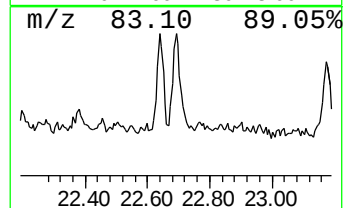
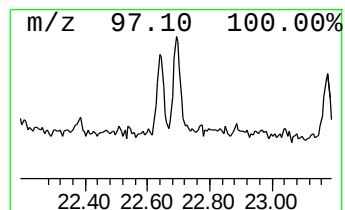
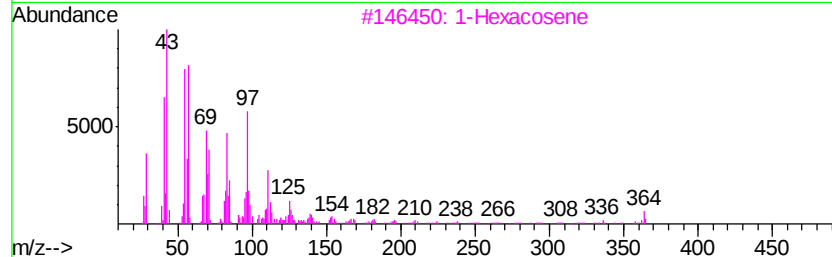
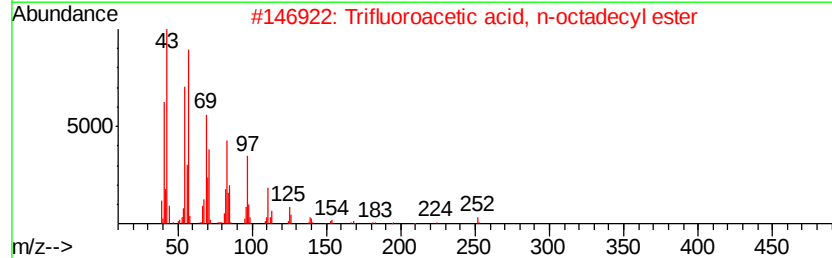
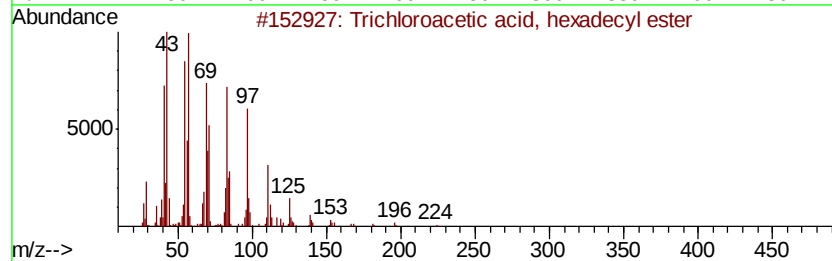
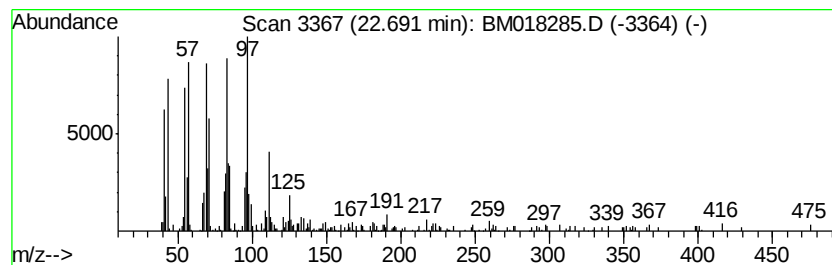
Instrument :
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ClientSampleId :
P001-SS012-1824-01

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 Trichloroacetic acid, hexad... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD		R.T.		
22.69	4.83 ng	352923	Perylene-d12		23.29		
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	80
2			Trifluoroacetic acid, n-octadecyl...	366	C20H37F3O2	079392-43-1	76
3			1-Hexacosene	364	C26H52	018835-33-1	58
4			1-Docosene	308	C22H44	001599-67-3	58
5			Cyclopentane, 1,1,3-trimethyl-	112	C8H16	004516-69-2	52



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121818\
Data File : BM018285.D
Acq On : 18 Dec 2018 23:36
Operator : JU/SJ
Sample : J6389-04
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
P001-SS012-1824-01

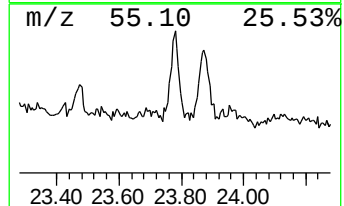
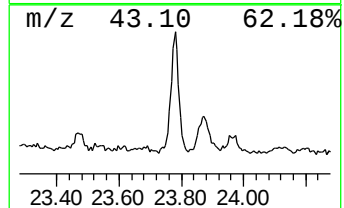
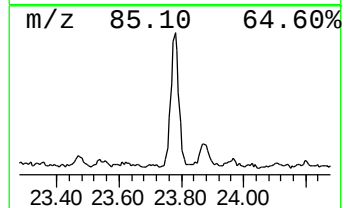
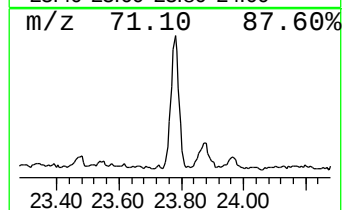
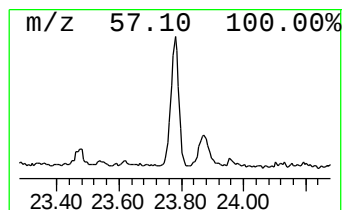
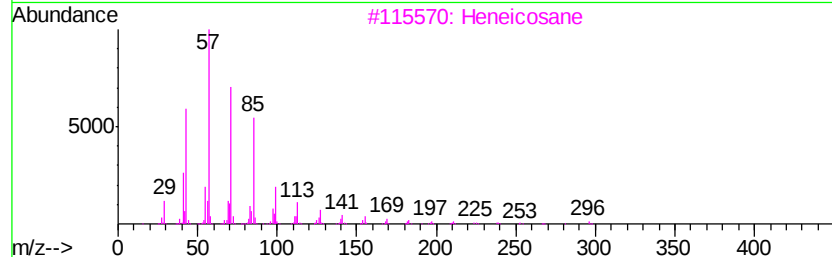
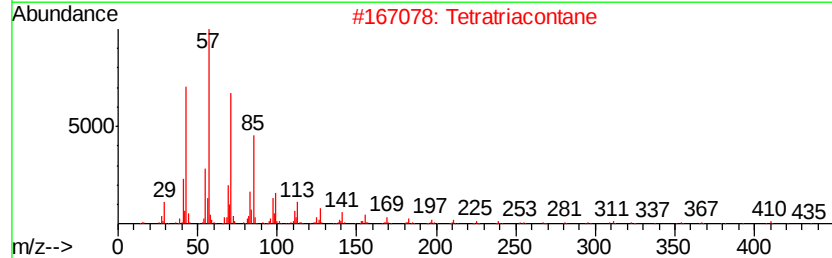
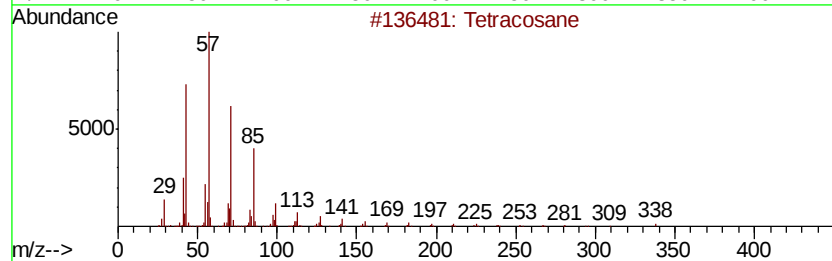
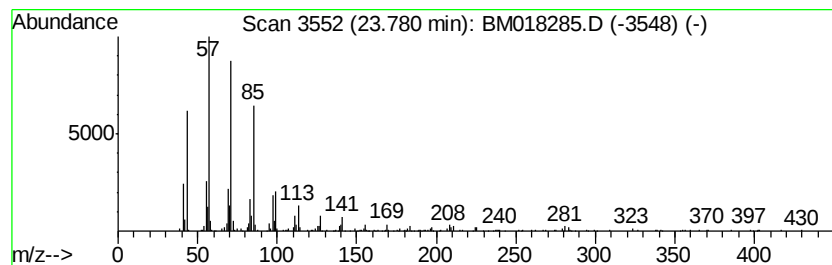
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Tetracosane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.78	9.50 ng	695024	Perylene-d12	23.29

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetracosane	338	C24H50	000646-31-1	91
2			Tetratriacontane	479	C34H70	014167-59-0	91
3			Heneicosane	296	C21H44	000629-94-7	90
4			Tricosane, 2-methyl-	338	C24H50	001928-30-9	89
5			Triacontane	422	C30H62	000638-68-6	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM121818\
 Data File : BM018285.D
 Acq On : 18 Dec 2018 23:36
 Operator : JU/SJ
 Sample : J6389-04
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 P001-SS012-1824-01

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
unknown4.68	4.68	2.7	ng	133221	1	7.49	977657	20.0
unknown7.03	7.03	88.6	ng	4329900	1	7.49	977657	20.0
n-Hexadecanoic acid	17.83	15.2	ng	1123640	4	16.91	1479270	20.0
Octadecanoic acid	19.12	6.0	ng	509582	5	21.13	1689560	20.0
1,1'-Biphenyl, 2,...	19.84	2.6	ng	222699	5	21.13	1689560	20.0
1,1'-Biphenyl, 2,...	20.13	2.0	ng	169002	5	21.13	1689560	20.0
Heptafluorobutano...	20.85	14.1	ng	1193020	5	21.13	1689560	20.0
Decane, 3,8-dimet...	21.29	3.2	ng	266808	5	21.13	1689560	20.0
Heptacosane	21.71	7.2	ng	606453	5	21.13	1689560	20.0
Hexadecanedinitrile	21.74	6.6	ng	559111	5	21.13	1689560	20.0
Heneicosane, 11-(...	22.16	6.4	ng	542989	5	21.13	1689560	20.0
Triacontane	22.64	10.9	ng	800704	6	23.29	1462540	20.0
Trichloroacetic a...	22.69	4.8	ng	352923	6	23.29	1462540	20.0
Tetracosane	23.78	9.5	ng	695024	6	23.29	1462540	20.0