

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM071118\
 Data File : BM015890.D
 Acq On : 11 Jul 2018 15:53
 Operator : SJ/JU
 Sample : J3906-07
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 CB-7-8-6-ST-EW-3@2.5

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-BM070518.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	5.316	337	345	358	rBV	237411	372844	7.76%	1.141%
2	5.799	420	427	444	rBV	1880257	2903527	60.45%	8.888%
3	7.446	700	707	727	rBV	1925984	3162901	65.85%	9.682%
4	7.816	763	770	785	rBV	2035914	3294989	68.60%	10.087%
5	8.293	845	851	864	rBV	329011	546155	11.37%	1.672%
6	8.616	898	906	919	rBV	1474207	2369699	49.34%	7.254%
7	9.481	1045	1053	1066	rBV	1126382	1905825	39.68%	5.834%
8	11.122	1325	1332	1341	rBV	382091	672482	14.00%	2.059%
9	13.539	1736	1743	1755	rBV	2622429	3683733	76.70%	11.277%
10	14.363	1877	1883	1890	rBV	295075	413887	8.62%	1.267%
11	14.916	1971	1977	1987	rVB2	503688	784406	16.33%	2.401%
12	16.392	2222	2228	2238	rBV	2158527	2967486	61.78%	9.084%
13	17.639	2434	2440	2452	rBV2	605719	896627	18.67%	2.745%
14	18.474	2574	2582	2593	rBV	184681	294814	6.14%	0.902%
15	19.309	2719	2724	2731	rBV	39389	59818	1.25%	0.183%
16	19.727	2791	2795	2805	rBV	128106	183902	3.83%	0.563%
17	20.204	2870	2876	2882	rBV	4146522	4803001	100.00%	14.703%
18	20.951	3000	3003	3007	rBV	54120	55915	1.16%	0.171%
19	21.409	3074	3081	3084	rBV5	107335	192867	4.02%	0.590%
20	21.503	3093	3097	3104	rBV	199683	257197	5.35%	0.787%
21	21.750	3133	3139	3144	rBV	666317	921580	19.19%	2.821%
22	21.845	3151	3155	3158	rBV	95595	112193	2.34%	0.343%
23	22.297	3228	3232	3236	rBV	109885	136104	2.83%	0.417%
24	22.786	3311	3315	3321	rVB2	131183	180285	3.75%	0.552%
25	23.321	3402	3406	3413	rVB	142703	211044	4.39%	0.646%
26	23.927	3505	3509	3517	rVB2	151708	259849	5.41%	0.795%
27	24.144	3540	3546	3554	rVB	422811	822979	17.13%	2.519%
28	24.615	3623	3626	3635	rVB2	111972	200679	4.18%	0.614%

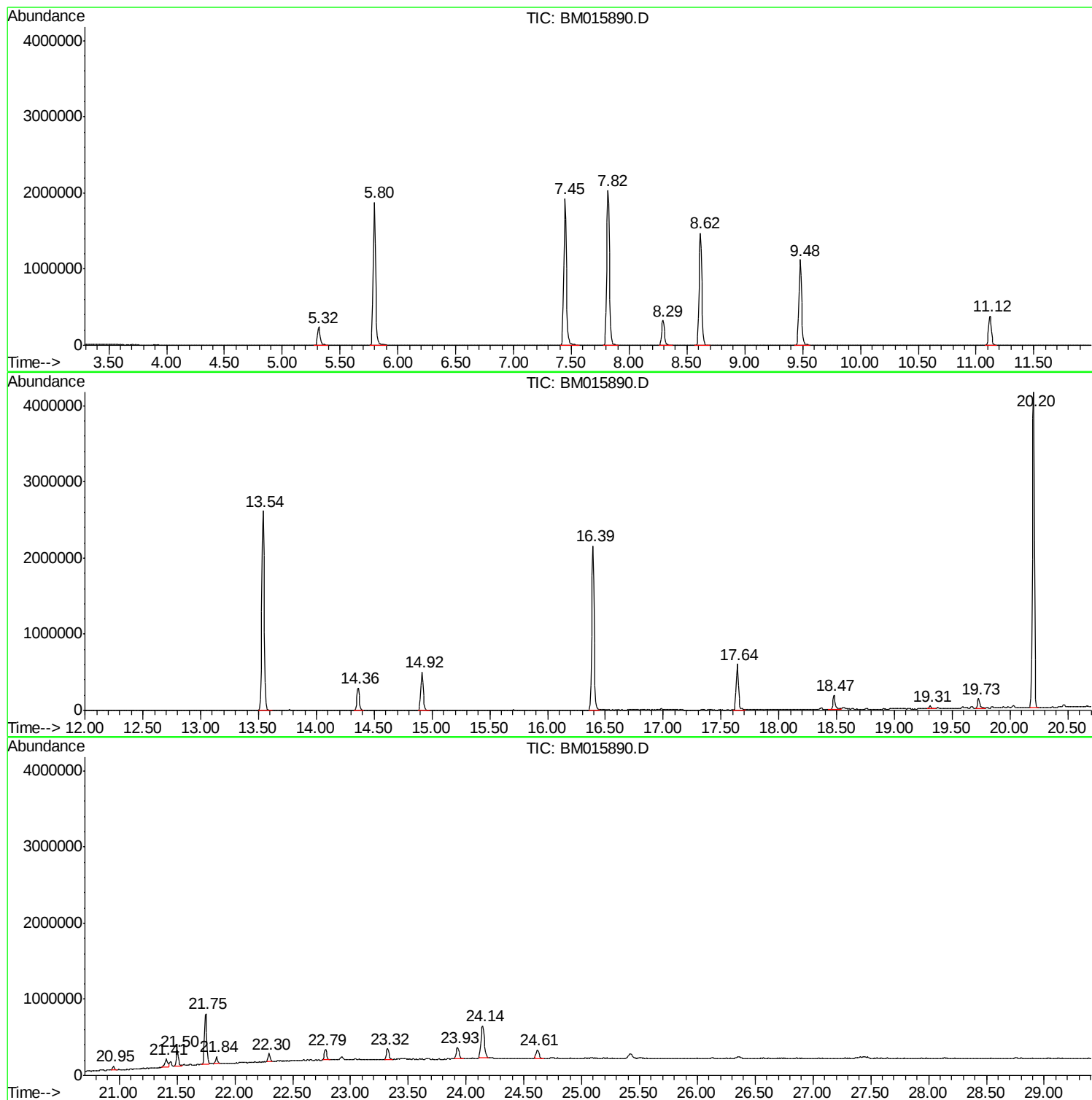
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM070518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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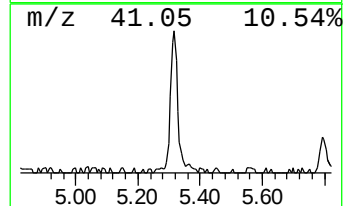
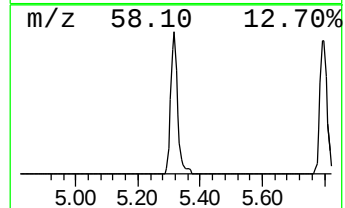
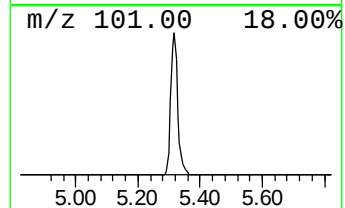
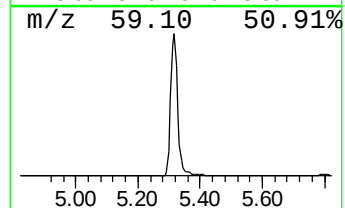
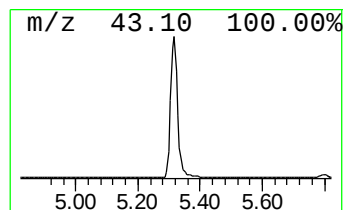
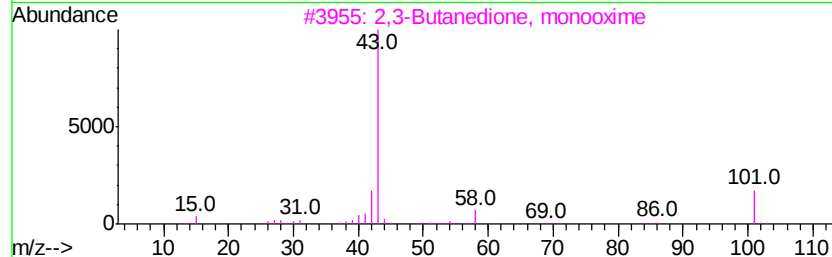
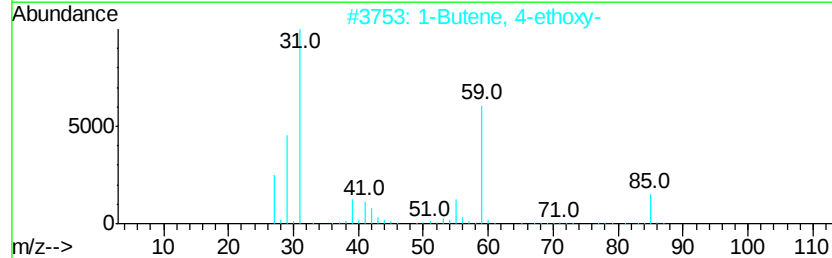
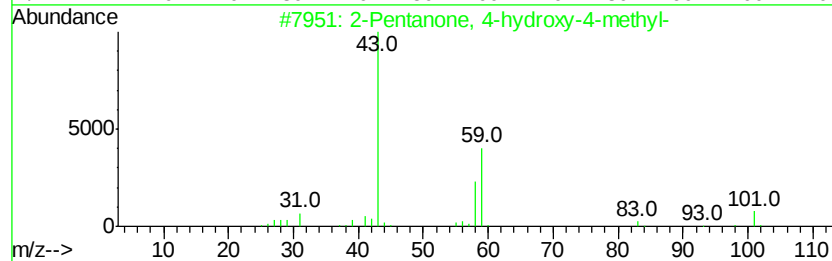
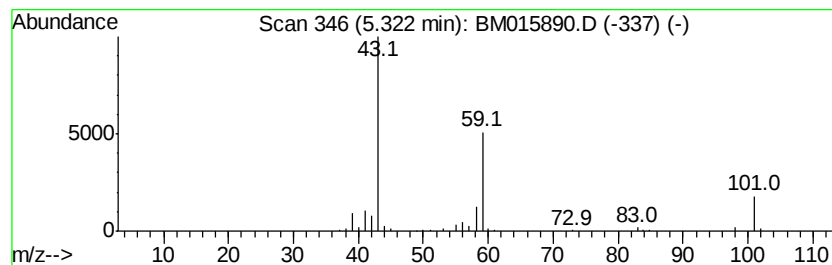
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.32	13.65 ng	372844	1,4-Dichlorobenzene-d4	8.29

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			1-Butene, 4-ethoxy-	100	C6H12O	044611-46-3	23
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16
4			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5			Butane	58	C4H10	000106-97-8	9



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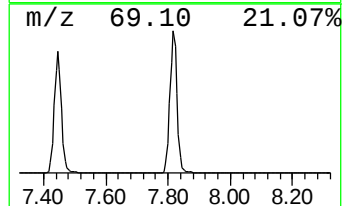
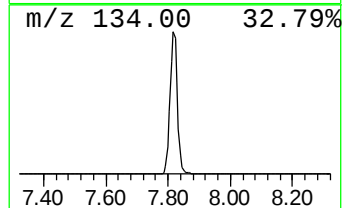
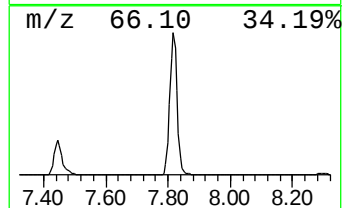
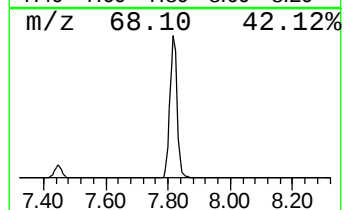
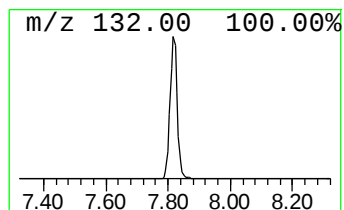
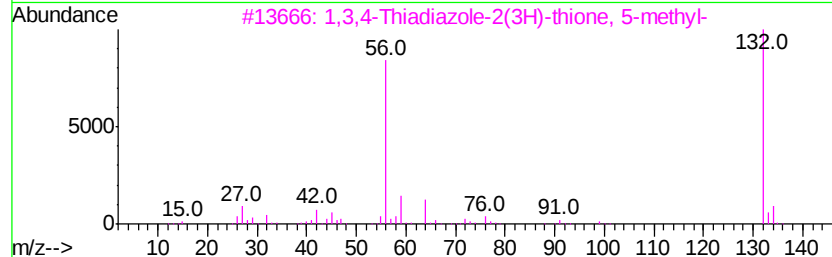
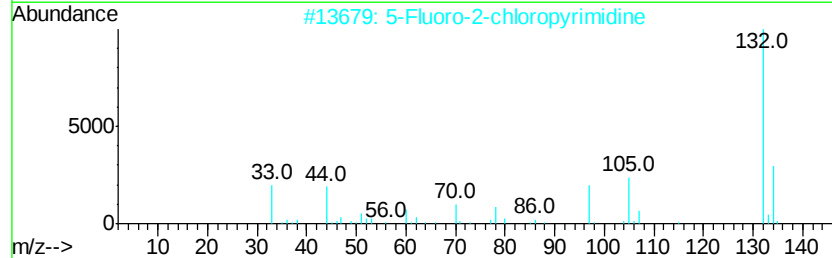
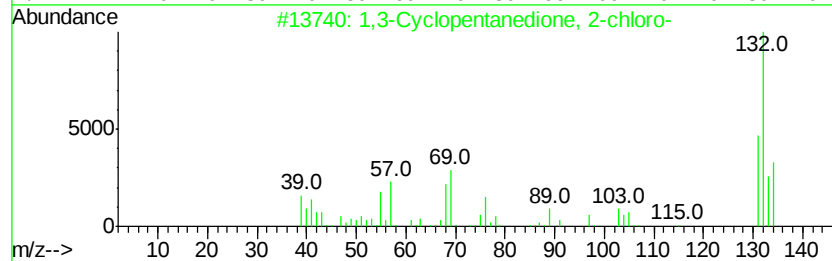
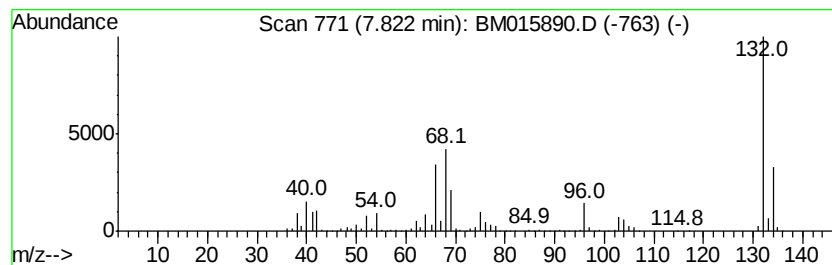
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Peak Number 2 unknown7.82 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.82	120.66 ng	3294990	1,4-Dichlorobenzene-d4	8.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	35
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	27
3		1,3,4-Thiadiazole-2(3H)-thione, ...	132	C3H4N2S2	029490-19-5	27
4		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	27
5		Imidazole, 4,5-dicyano-1-methyl-	132	C6H4N4	019485-35-9	22



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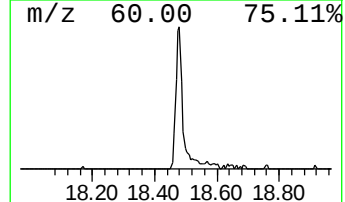
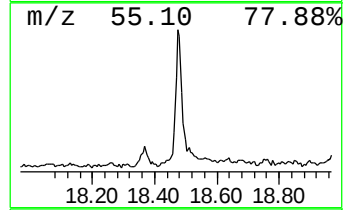
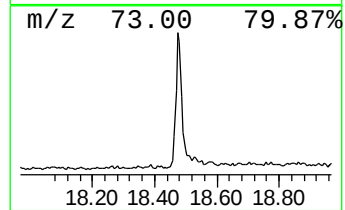
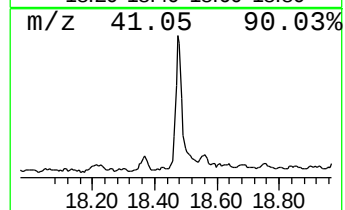
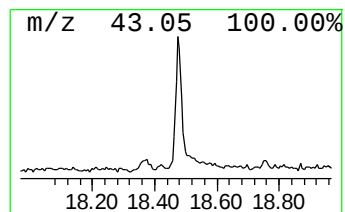
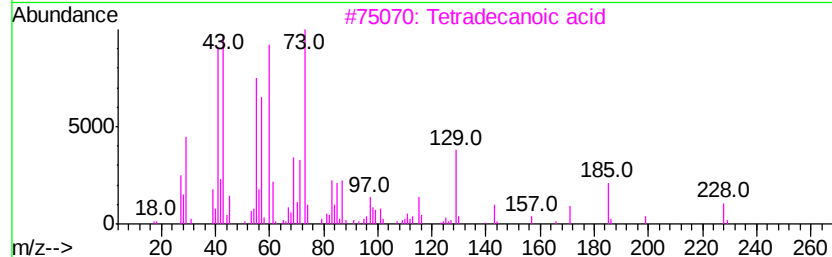
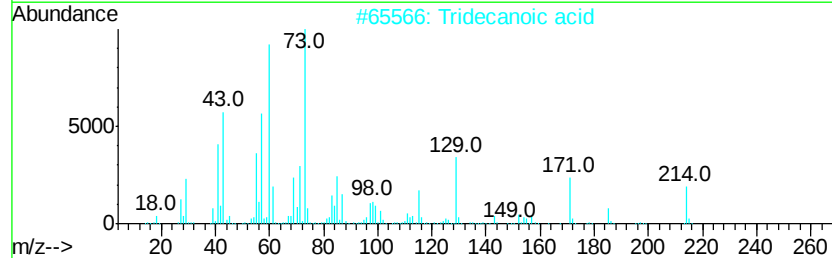
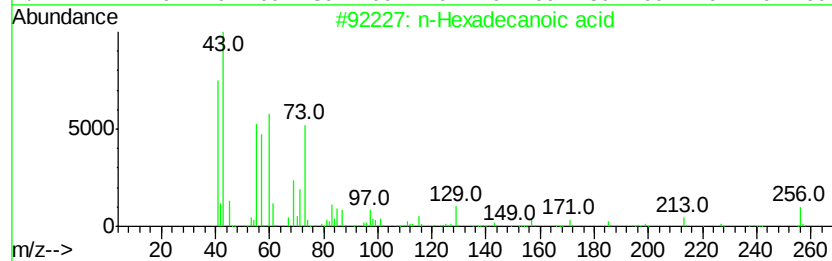
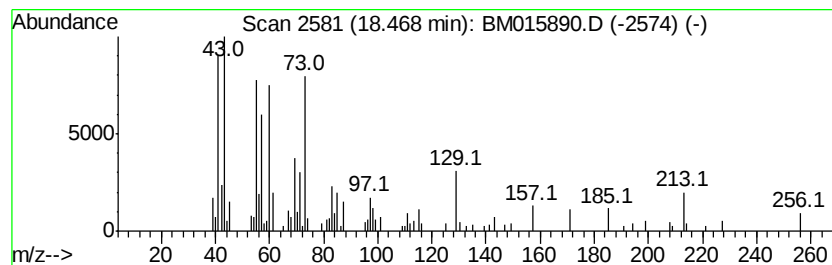
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Peak Number 4 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.47	6.58 ng	294814	Phenanthrene-d10	17.64

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	90
2		Tridecanoic acid	214	C13H26O2	000638-53-9	62
3		Tetradecanoic acid	228	C14H28O2	000544-63-8	53
4		Pentadecanoic acid	242	C15H30O2	001002-84-2	47
5		Undecanoic acid	186	C11H22O2	000112-37-8	47



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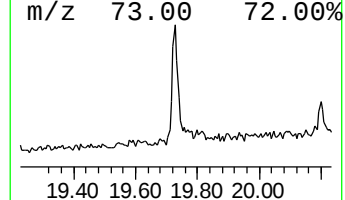
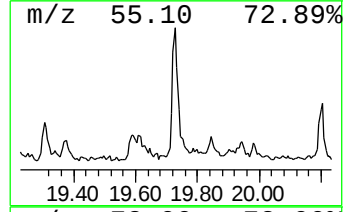
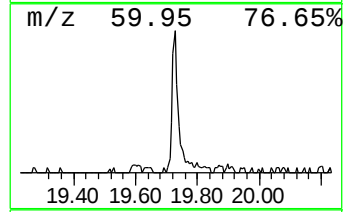
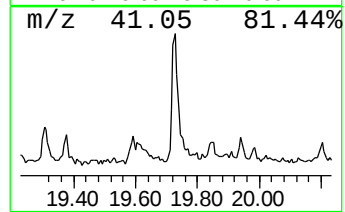
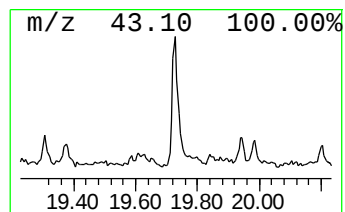
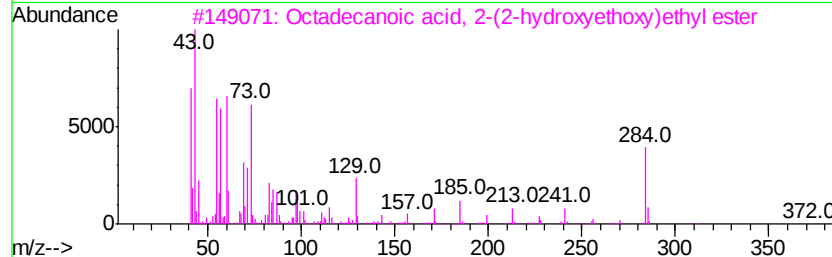
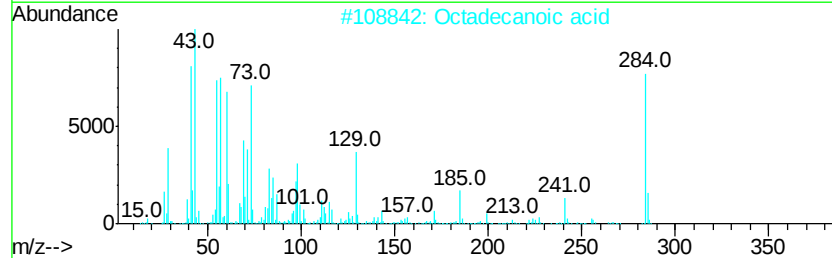
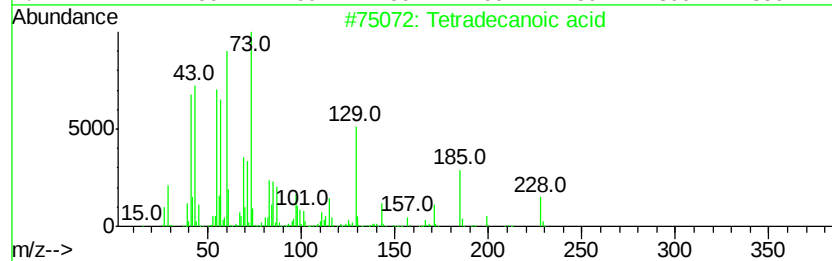
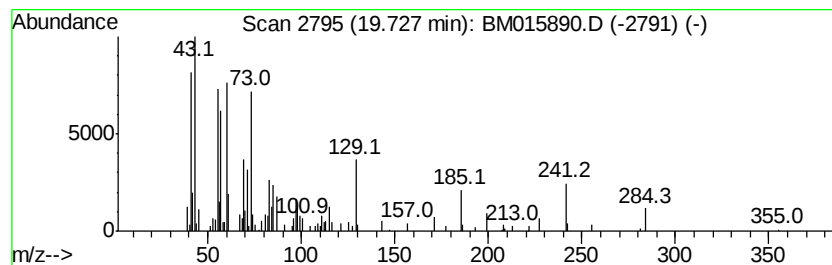
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Peak Number 5 Tetradecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.73	3.99 ng	183902	Chrysene-d12	21.75

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetradecanoic acid	228	C14H28O2	000544-63-8	83
2		Octadecanoic acid	284	C18H36O2	000057-11-4	78
3		Octadecanoic acid, 2-(2-hydroxyve...	372	C22H44O4	000106-11-6	72
4		Pentadecanoic acid	242	C15H30O2	001002-84-2	64
5		Undecanoic acid	186	C11H22O2	000112-37-8	53



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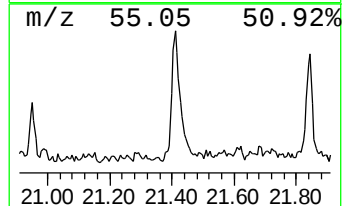
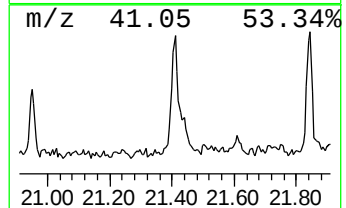
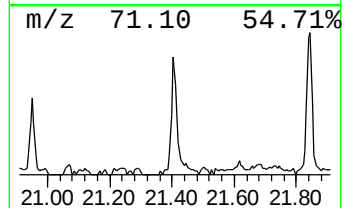
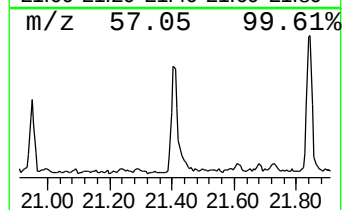
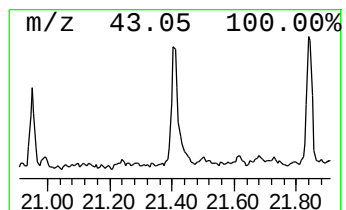
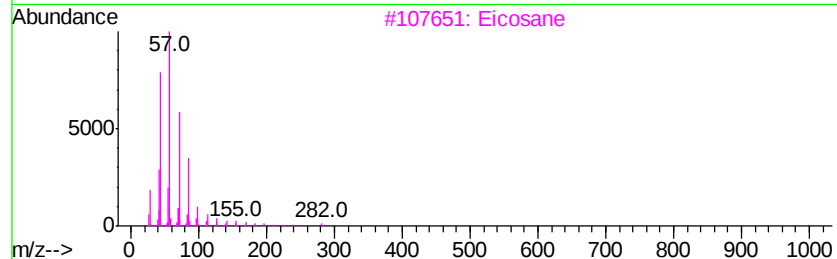
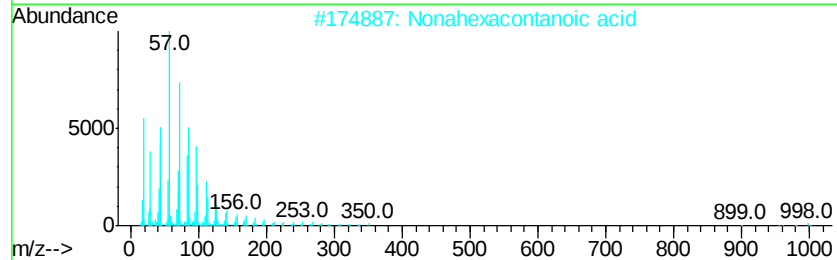
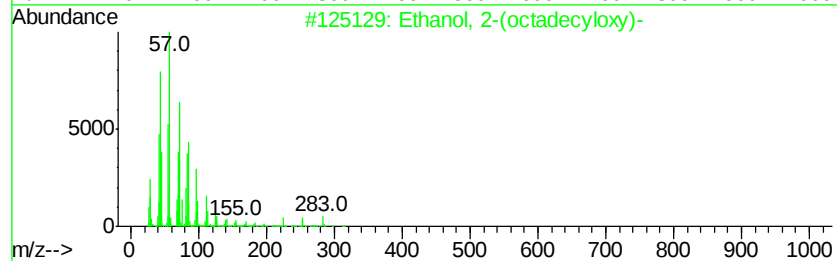
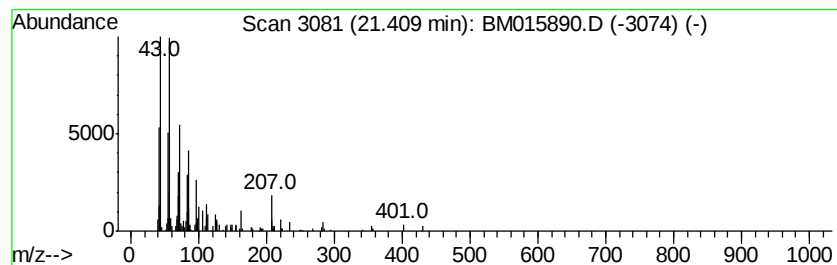
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 Peak Number 6 Ethanol, 2-(octadecyloxy)- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.41	4.19 ng	192867	Chrysene-d12	21.75

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethanol, 2-(octadecyloxy)-	314	C20H42O2	002136-72-3	68
2		Nonaheptacontanoic acid	999	C69H138O2	040710-32-5	68
3		Eicosane	282	C20H42	000112-95-8	55
4		1-Pentadecene	210	C15H30	013360-61-7	53
5		Nonadecane	268	C19H40	000629-92-5	49



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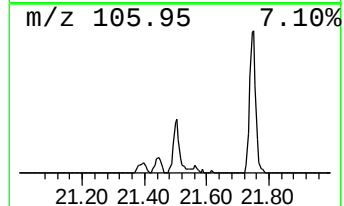
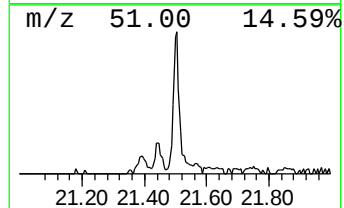
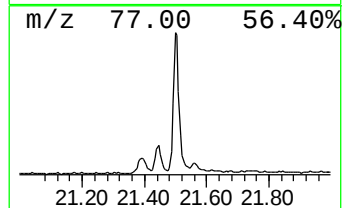
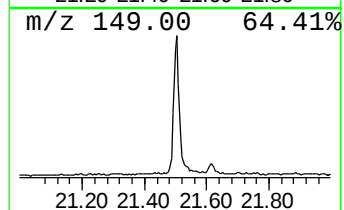
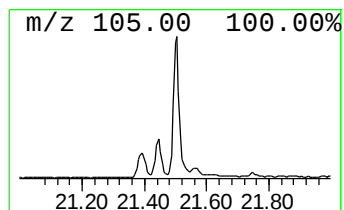
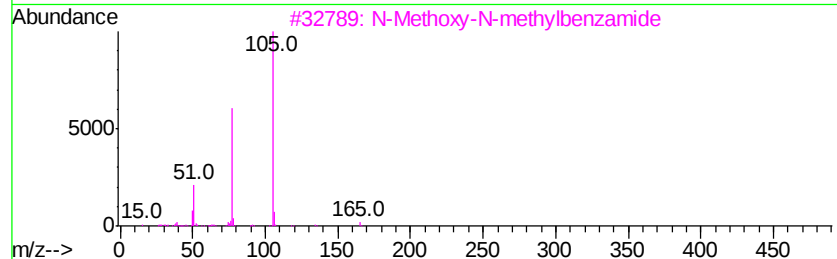
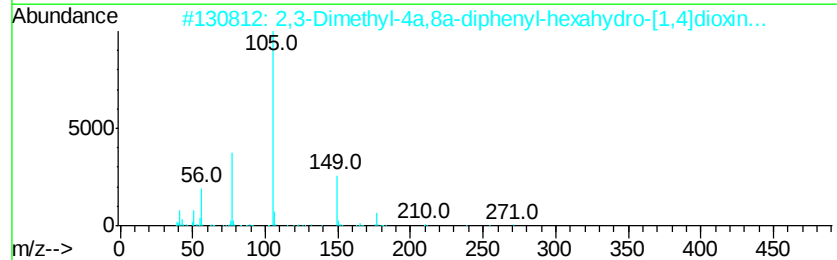
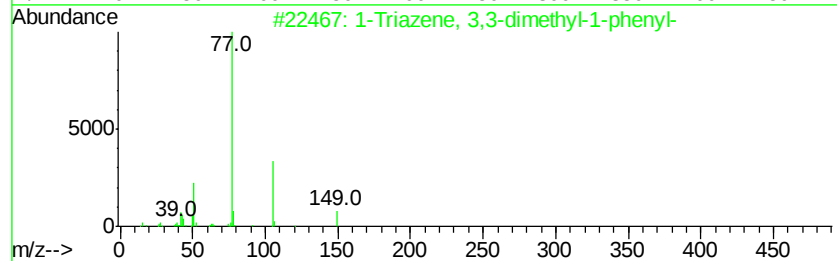
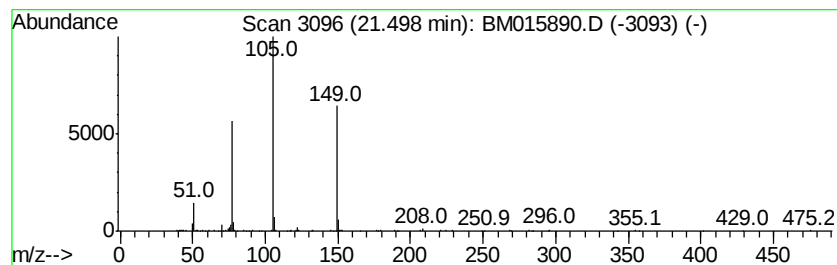
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 1-Triazene, 3,3-dimethyl-1-... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.50	5.58 ng	257197	Chrysene-d12	21.75

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Triazene, 3,3-dimethyl-1-phenyl-	149	C8H11N3	007227-91-0	64
2			2,3-Dimethyl-4a,8a-diphenyl-hexa...	326	C20H22O4	1000297-07-0	59
3			N-Methoxy-N-methylbenzamide	165	C9H11NO2	006919-61-5	43
4			Acetophenone, 2-chloro-	154	C8H7ClO	000532-27-4	43
5			Benzenepropanenitrile, .beta.-oxo-	145	C9H7NO	000614-16-4	43



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM071118\
Data File : BM015890.D
Acq On : 11 Jul 2018 15:53
Operator : SJ/JU
Sample : J3906-07
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
CB-7-8-6-ST-EW-3@2.5

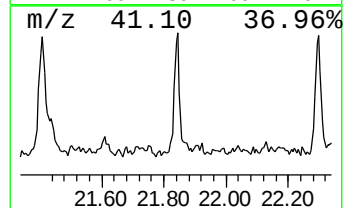
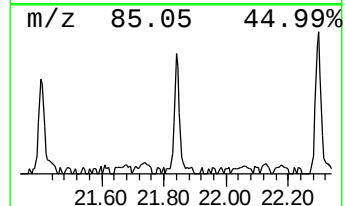
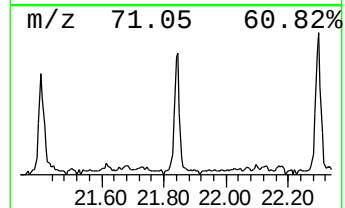
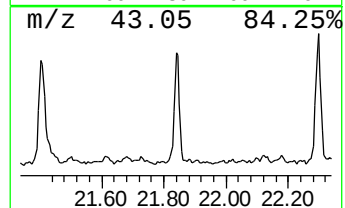
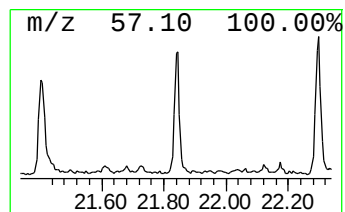
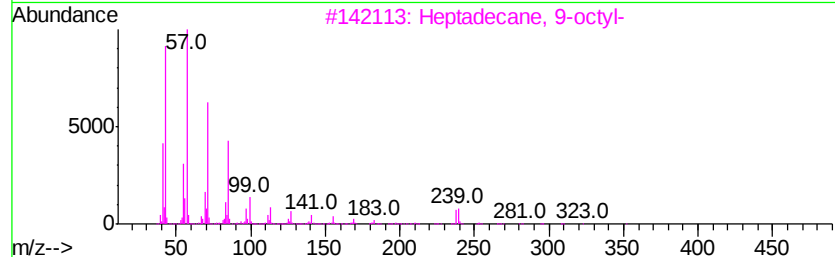
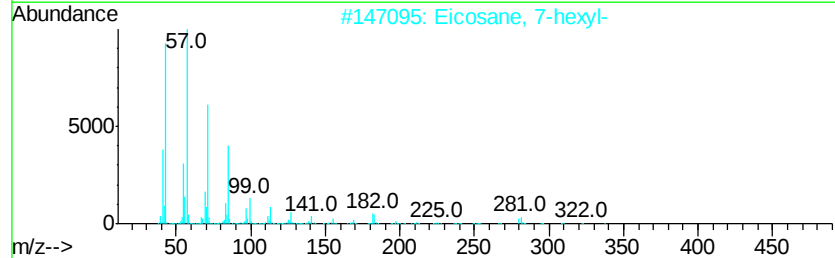
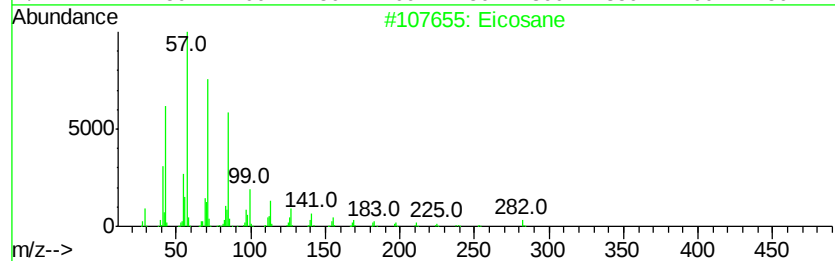
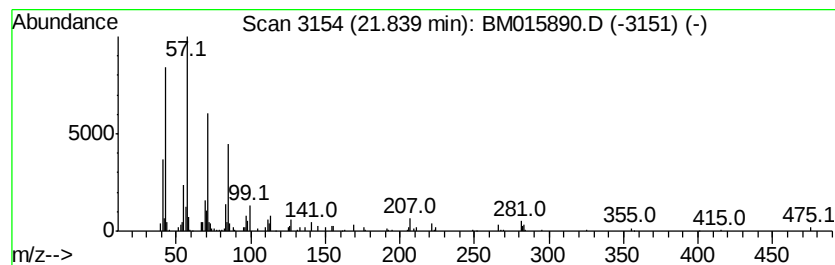
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Eicosane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.84	2.43 ng	112193	Chrysene-d12	21.75

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	98
2		Eicosane, 7-hexyl-	366	C26H54	055333-99-8	91
3		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	90
4		Heneicosane	296	C21H44	000629-94-7	90
5		Hexatriacontane	507	C36H74	000630-06-8	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM071118\
Data File : BM015890.D
Acq On : 11 Jul 2018 15:53
Operator : SJ/JU
Sample : J3906-07
Misc :
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Instrument :
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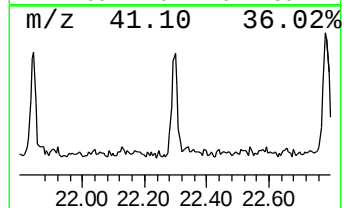
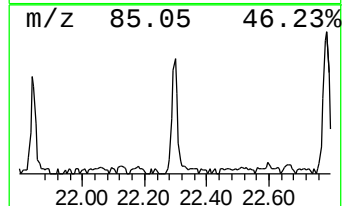
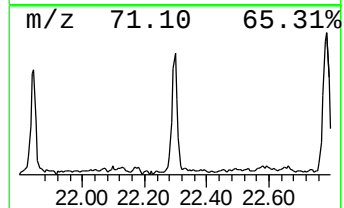
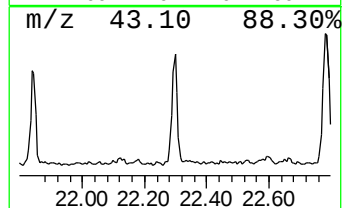
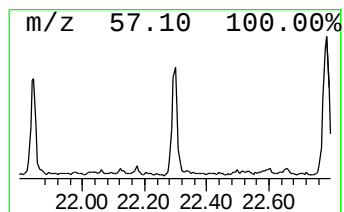
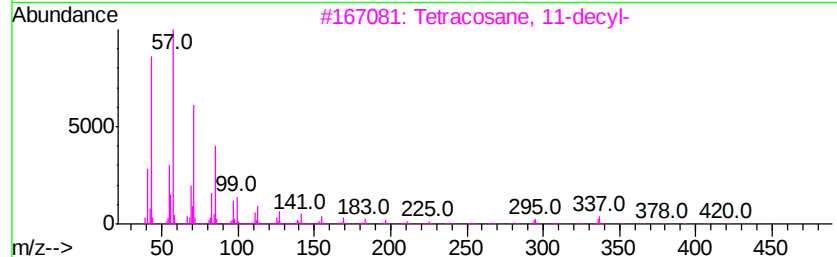
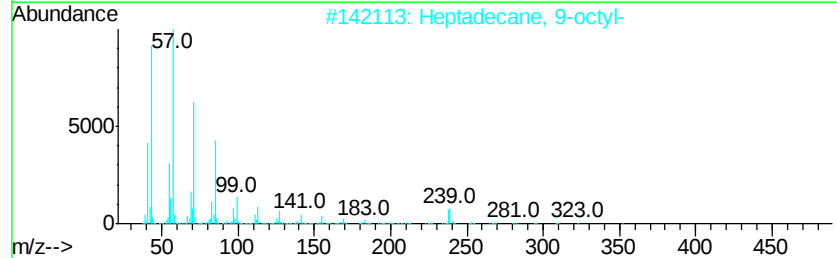
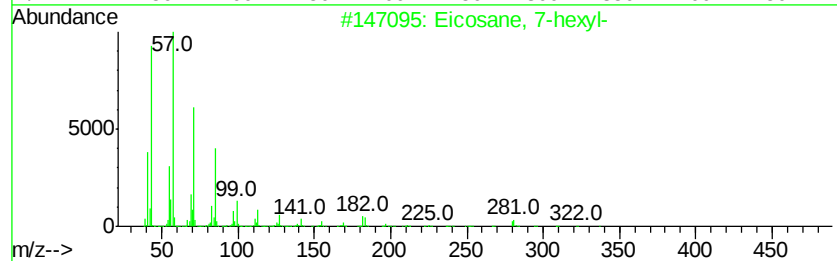
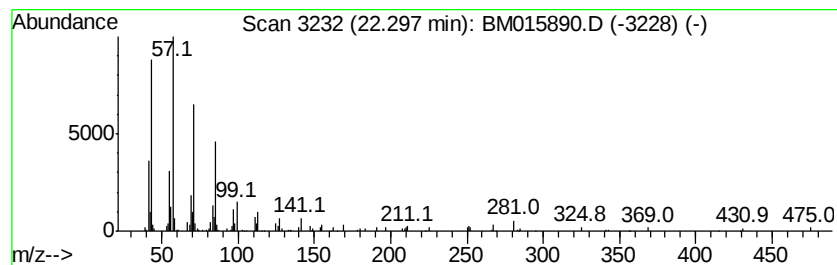
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Eicosane, 7-hexyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.30	2.95 ng	136104	Chrysene-d12	21.75

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane, 7-hexyl-	366	C26H54	055333-99-8	91
2		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	90
3		Tetracosane, 11-decyl-	479	C34H70	055429-84-0	90
4		Heptacosane	380	C27H56	000593-49-7	90
5		Hexatriacontane	507	C36H74	000630-06-8	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM071118\
Data File : BM015890.D
Acq On : 11 Jul 2018 15:53
Operator : SJ/JU
Sample : J3906-07
Misc :
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Instrument :
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ClientSampleId :
CB-7-8-6-ST-EW-3@2.5

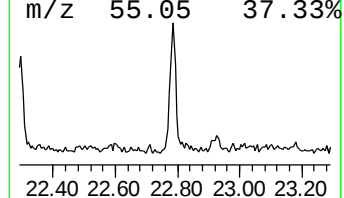
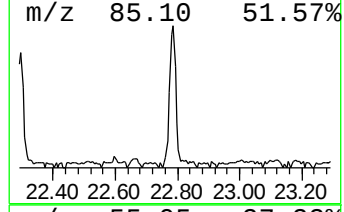
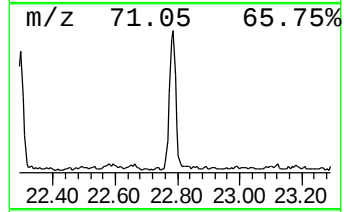
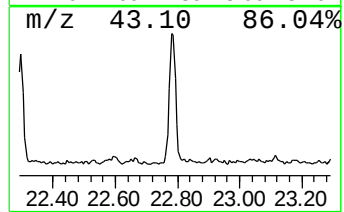
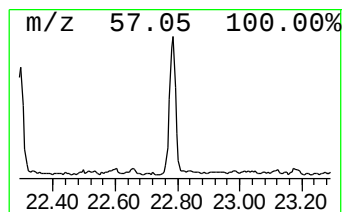
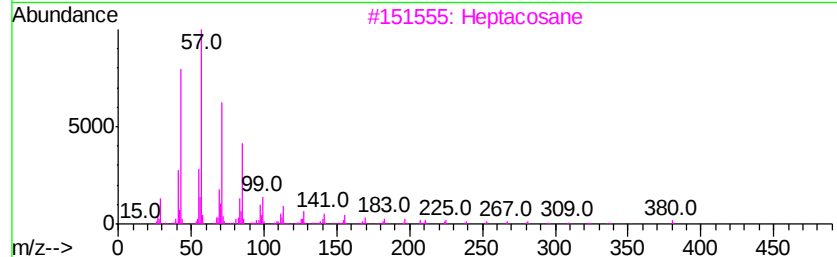
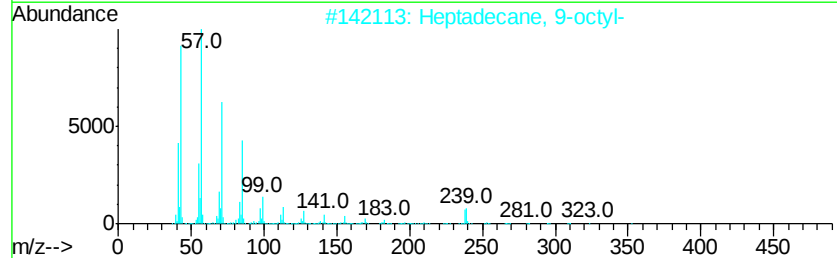
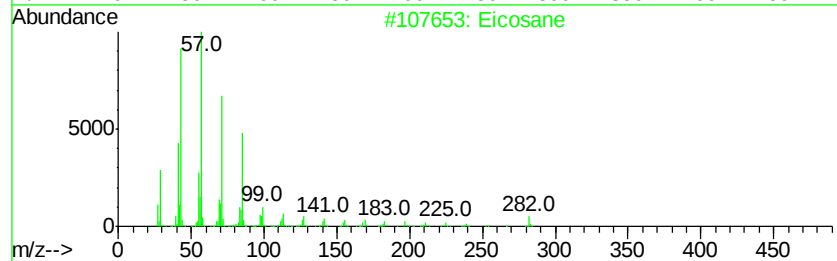
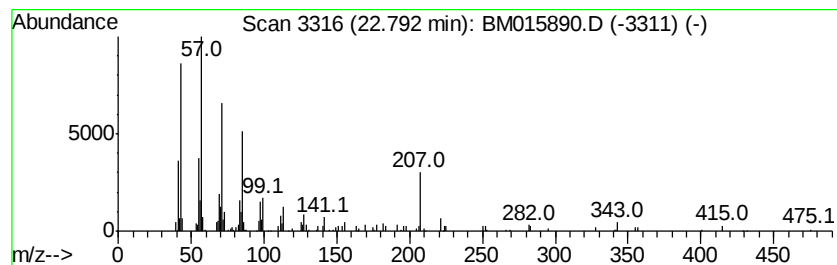
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Heptacosane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.79	3.91 ng	180285	Chrysene-d12	21.75

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane	282	C20H42	000112-95-8	95
2			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	87
3			Heptacosane	380	C27H56	000593-49-7	87
4			Tetracosane	338	C24H50	000646-31-1	87
5			Hexatriacontane	507	C36H74	000630-06-8	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM071118\
Data File : BM015890.D
Acq On : 11 Jul 2018 15:53
Operator : SJ/JU
Sample : J3906-07
Misc :
ALS Vial : 7 Sample Multiplier: 1

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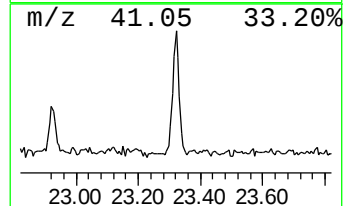
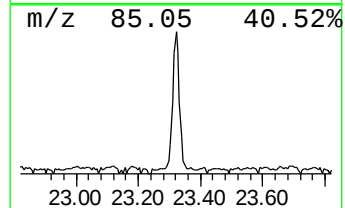
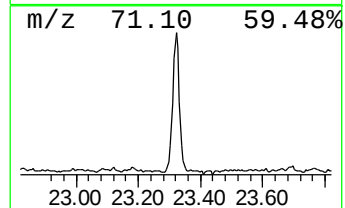
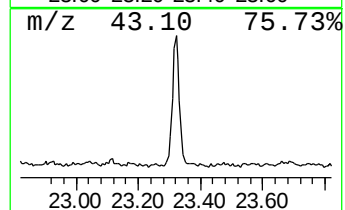
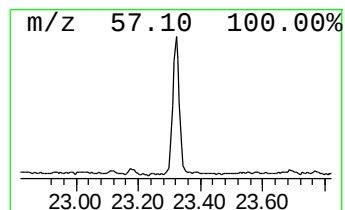
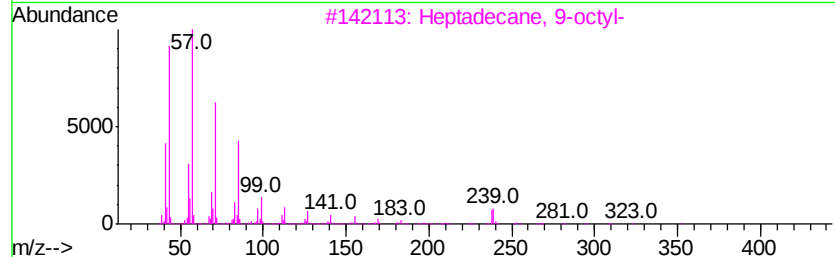
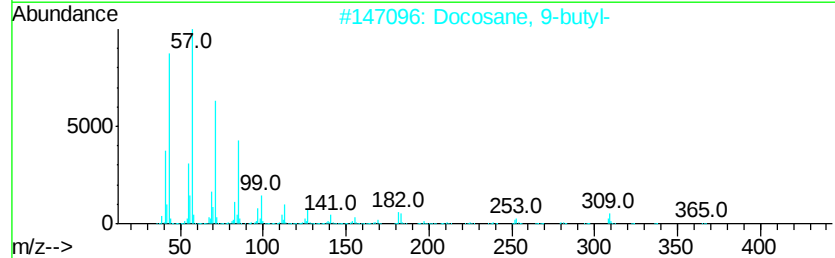
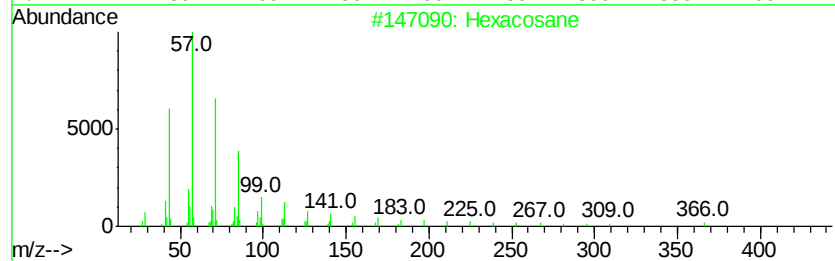
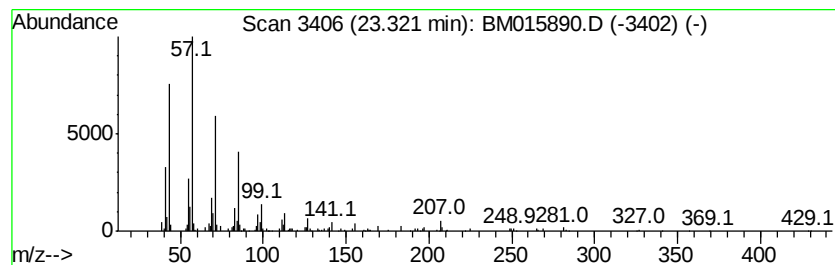
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Hexacosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.32	5.13 ng	211044	Perylene-d12	24.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexacosane	366	C26H54	000630-01-3	90
2		Docosane, 9-butyl-	366	C26H54	055282-14-9	90
3		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	90
4		Docosane, 11-butyl-	366	C26H54	013475-76-8	90
5		Heptacosane	380	C27H56	000593-49-7	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM071118\
Data File : BM015890.D
Acq On : 11 Jul 2018 15:53
Operator : SJ/JU
Sample : J3906-07
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
CB-7-8-6-ST-EW-3@2.5

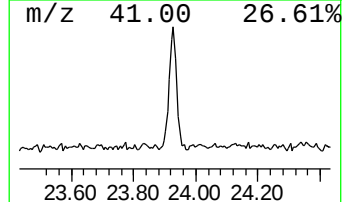
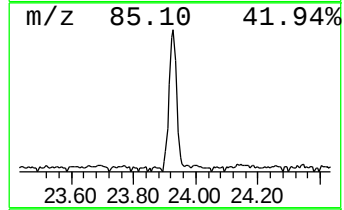
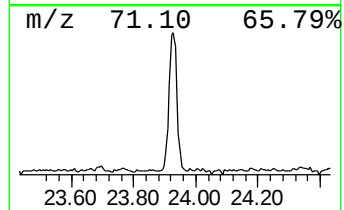
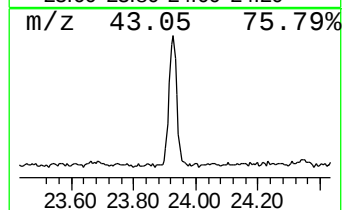
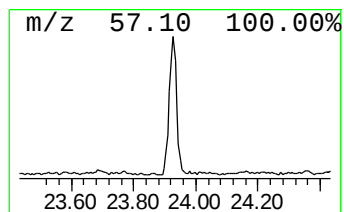
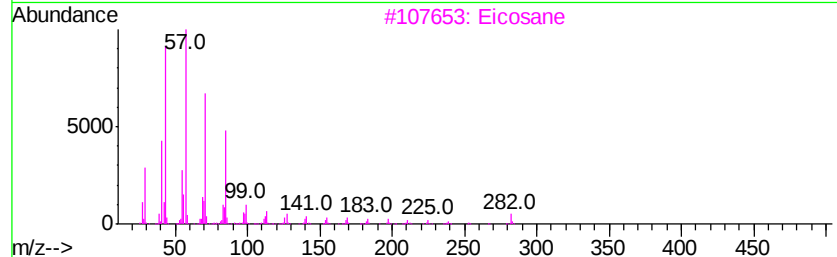
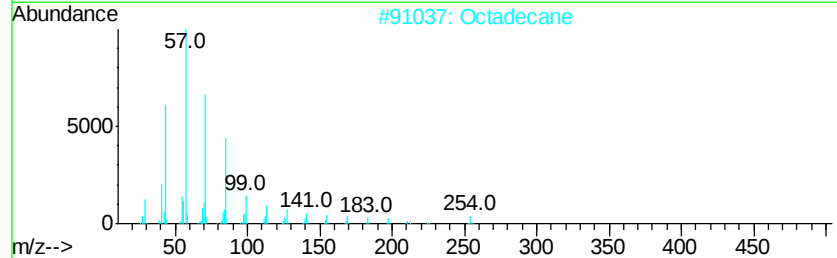
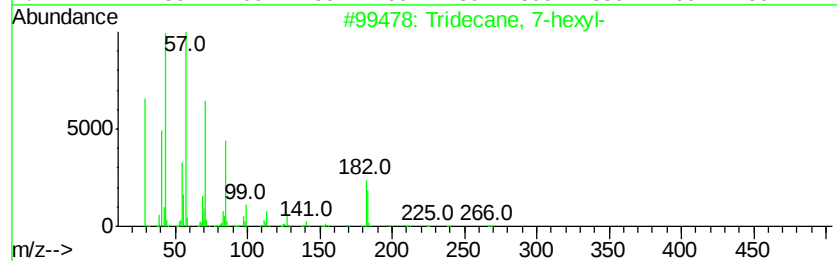
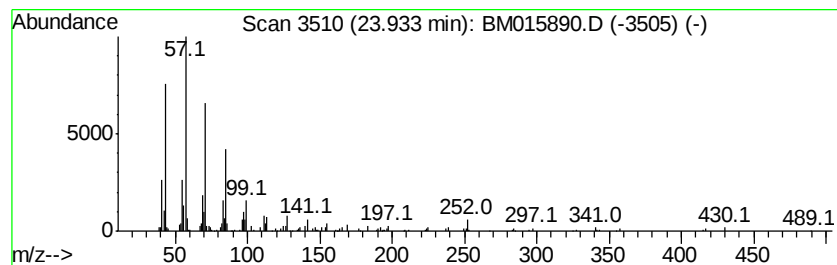
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Tridecane, 7-hexyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.93	6.31 ng	259849	Perylene-d12	24.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tridecane, 7-hexyl-	268	C19H40	007225-66-3	94
2		Octadecane	254	C18H38	000593-45-3	92
3		Eicosane	282	C20H42	000112-95-8	90
4		Tetracosane	338	C24H50	000646-31-1	90
5		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM071118\
Data File : BM015890.D
Acq On : 11 Jul 2018 15:53
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Sample : J3906-07
Misc :
ALS Vial : 7 Sample Multiplier: 1

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ClientSampleId :
CB-7-8-6-ST-EW-3@2.5

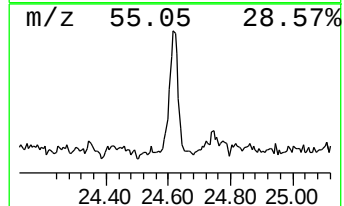
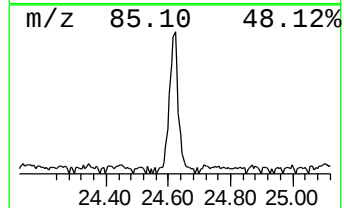
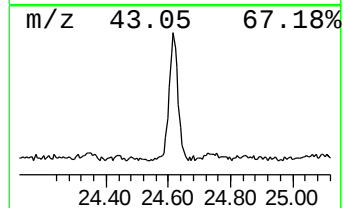
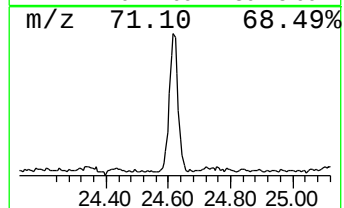
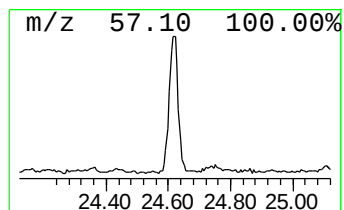
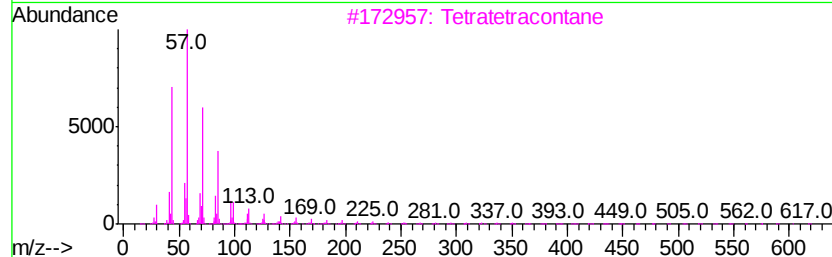
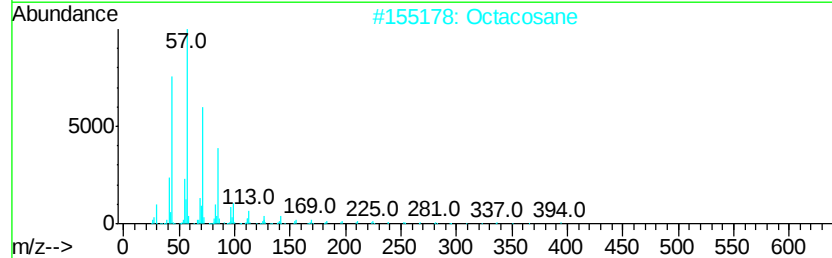
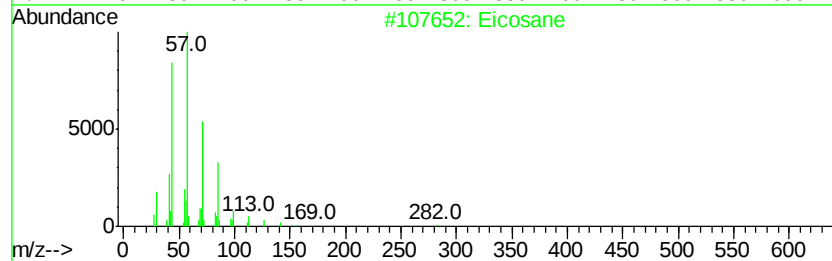
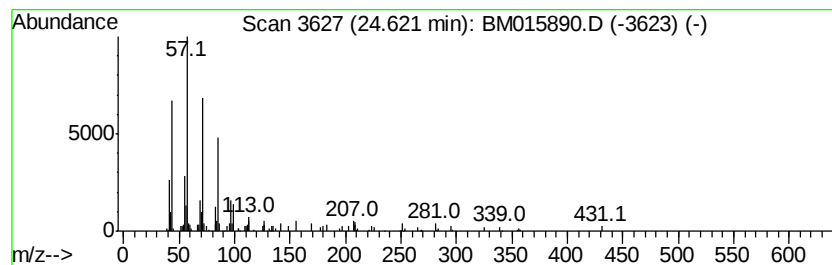
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Octacosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.62	4.88 ng	200679	Perylene-d12	24.14

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane	282	C20H42	000112-95-8	96
2			Octacosane	394	C28H58	000630-02-4	90
3			Tetratetracontane	619	C44H90	007098-22-8	90
4			Hexatriacontane	507	C36H74	000630-06-8	90
5			Tetracosane	338	C24H50	000646-31-1	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM071118\
Data File : BM015890.D
Acq On : 11 Jul 2018 15:53
Operator : SJ/JU
Sample : J3906-07
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
CB-7-8-6-ST-EW-3@2.5

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\8270-BM070518.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-Pentanone, 4-hy...	5.32	13.7	ng	372844	1	8.29	546155	20.0
unknown7.82	7.82	120.7	ng	3294990	1	8.29	546155	20.0
n-Hexadecanoic acid	18.47	6.6	ng	294814	4	17.64	896627	20.0
Tetradecanoic acid	19.73	4.0	ng	183902	5	21.75	921580	20.0
Ethanol, 2-(octad...	21.41	4.2	ng	192867	5	21.75	921580	20.0
1-Triazene, 3,3-d...	21.50	5.6	ng	257197	5	21.75	921580	20.0
Eicosane	21.84	2.4	ng	112193	5	21.75	921580	20.0
Eicosane, 7-hexyl-	22.30	3.0	ng	136104	5	21.75	921580	20.0
Heptacosane	22.79	3.9	ng	180285	5	21.75	921580	20.0
Hexacosane	23.32	5.1	ng	211044	6	24.14	822979	20.0
Tridecane, 7-hexyl-	23.93	6.3	ng	259849	6	24.14	822979	20.0
Octacosane	24.62	4.9	ng	200679	6	24.14	822979	20.0