

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA M\DATA\BM030718\
 Data File : BM014503.D
 Acq On : 06 Mar 2018 18:11
 Operator : SJ/JU
 Sample : J1699-13
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 B5

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0
 Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\HPCHEM1\BNA_M\METHODS\8270-BM020718.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	4.687	287	289	296	rVB2	32992	47349	1.50%	0.170%
2	5.140	360	366	380	rBV	1545236	2391306	75.73%	8.610%
3	6.716	627	634	651	rBV	1326102	2395373	75.86%	8.625%
4	7.045	682	690	702	rBV	1688679	2640566	83.62%	9.508%
5	7.504	762	768	775	rBV	448503	703630	22.28%	2.534%
6	7.816	814	821	835	rBV	1325493	2113850	66.94%	7.611%
7	8.669	958	966	981	rBV	829578	1650543	52.27%	5.943%
8	10.280	1231	1240	1253	rBV2	415012	865177	27.40%	3.115%
9	12.492	1610	1616	1624	rBV2	50189	88405	2.80%	0.318%
10	12.774	1655	1664	1683	rBV	1852602	3157783	100.00%	11.370%
11	13.692	1815	1820	1826	rVB6	15406	32540	1.03%	0.117%
12	14.168	1892	1901	1914	rBV3	587191	986123	31.23%	3.551%
13	14.780	1993	2005	2007	rBV6	25191	76963	2.44%	0.277%
14	14.968	2032	2037	2042	rBV4	20219	32071	1.02%	0.115%
15	15.021	2042	2046	2052	rVB5	31140	58473	1.85%	0.211%
16	15.680	2152	2158	2174	rBV	1260344	2218104	70.24%	7.987%
17	16.939	2366	2372	2381	rBV2	590047	995266	31.52%	3.584%
18	17.221	2415	2420	2424	rBV	51106	65130	2.06%	0.235%
19	17.603	2481	2485	2491	rVB2	127388	155662	4.93%	0.560%
20	17.839	2519	2525	2533	rBV2	290134	505405	16.01%	1.820%
21	17.903	2533	2536	2548	rVB	193266	268531	8.50%	0.967%
22	18.756	2676	2681	2684	rBV2	266771	337557	10.69%	1.215%
23	18.792	2684	2687	2691	rVB	281189	342303	10.84%	1.233%
24	18.927	2705	2710	2713	rBV4	72386	100545	3.18%	0.362%
25	19.133	2740	2745	2753	rBV2	248051	436543	13.82%	1.572%
26	19.350	2779	2782	2785	rBV4	39382	42710	1.35%	0.154%
27	19.580	2813	2821	2827	rBV	2498563	2993259	94.79%	10.778%
28	19.886	2870	2873	2876	rVB5	40953	44672	1.41%	0.161%
29	20.386	2956	2958	2962	rVB5	47505	49748	1.58%	0.179%
30	20.850	3034	3037	3043	rBV5	117207	160336	5.08%	0.577%
31	21.150	3084	3088	3100	rVB	698408	977124	30.94%	3.518%
32	23.333	3453	3459	3465	rBV	437516	839173	26.57%	3.022%

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Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
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Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

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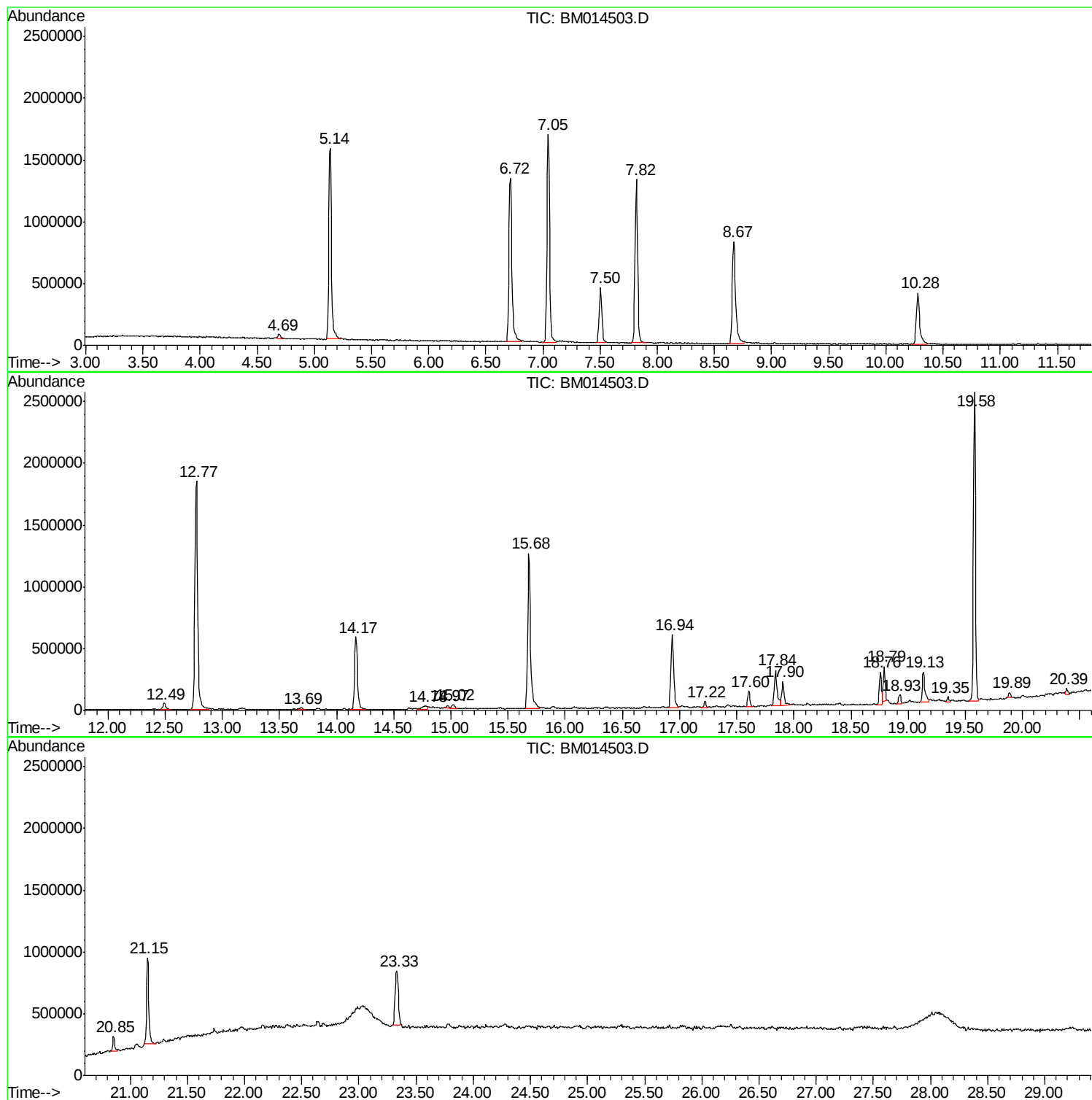
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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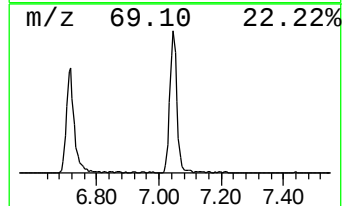
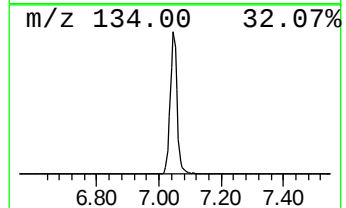
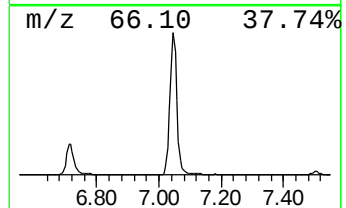
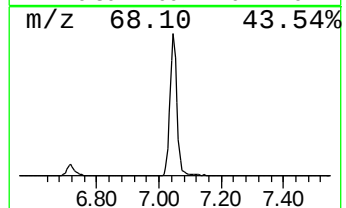
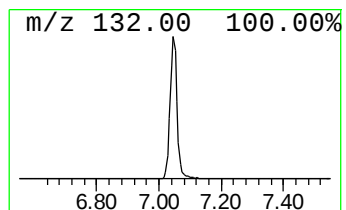
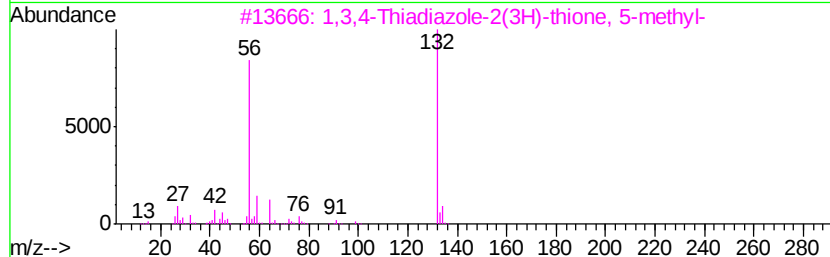
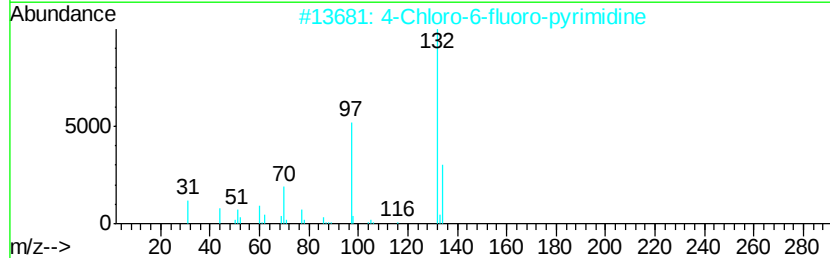
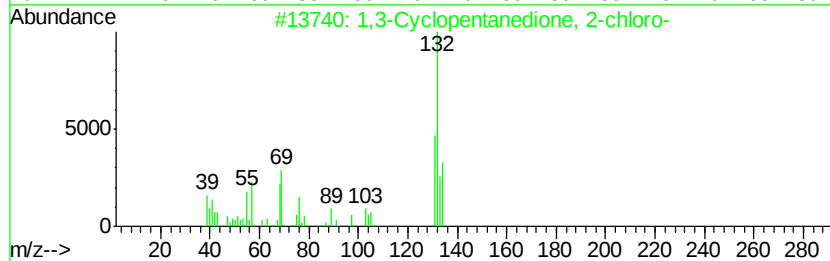
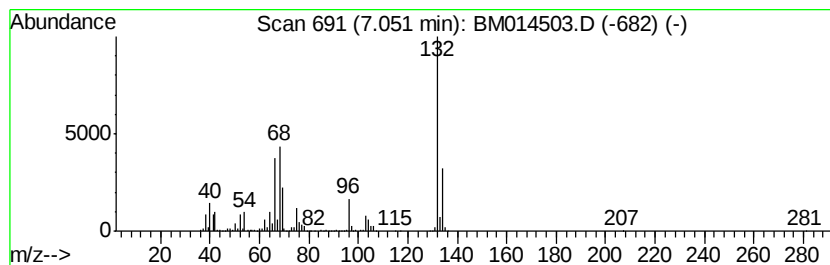
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown7.05 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.05	75.06 ng	2640570	1,4-Dichlorobenzene-d4	7.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	16
2		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	14
3		1,3,4-Thiadiazole-2(3H)-thione, ...	132	C3H4N2S2	029490-19-5	14
4		3H-1,2-Dithiol-3-one, 5-methyl-	132	C4H4OS2	003620-08-4	14
5		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10



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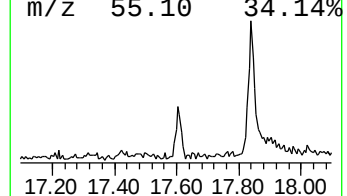
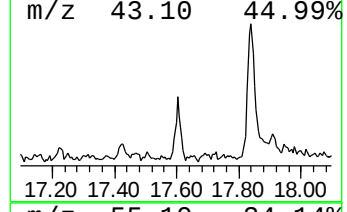
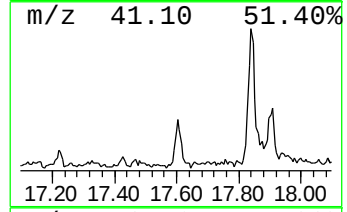
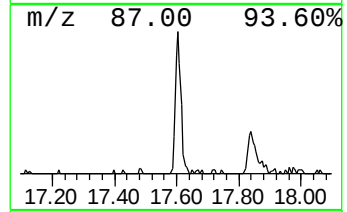
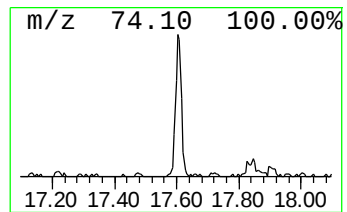
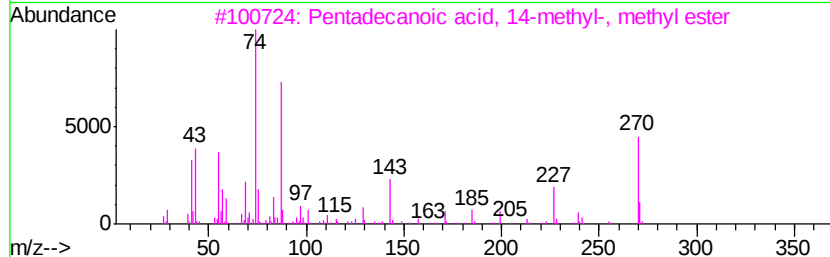
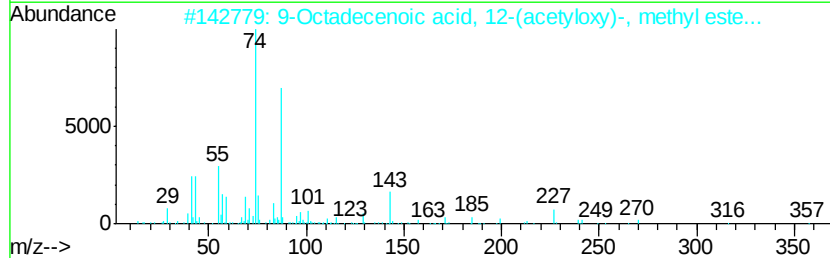
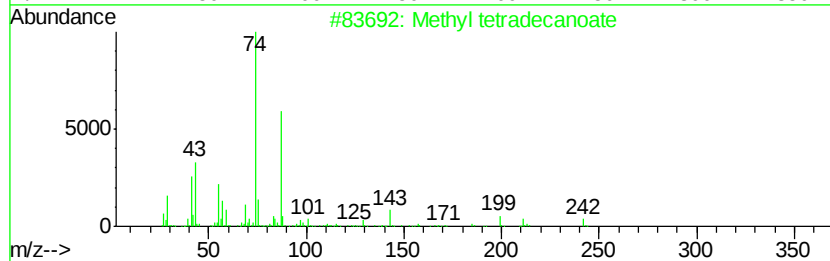
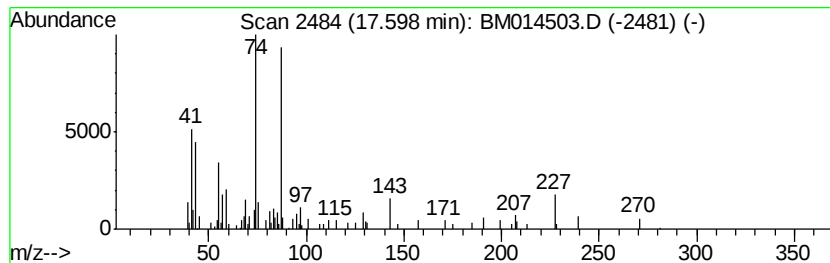
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 3 Methyl tetradecanoate Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.60	3.13 ng	155662	Phenanthrene-d10	16.94

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Methyl tetradecanoate	242	C15H30O2	000124-10-7	93
2		9-Octadecenoic acid, 12-(acetylo...	354	C21H38O4	000140-03-4	90
3		Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	83
4		Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	80
5		Hexadecanoic acid, 2-methyl-	270	C17H34O2	027147-71-3	74



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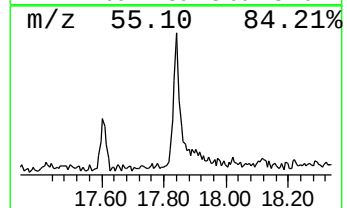
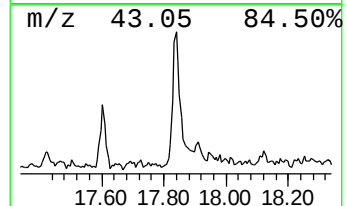
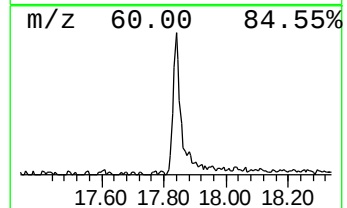
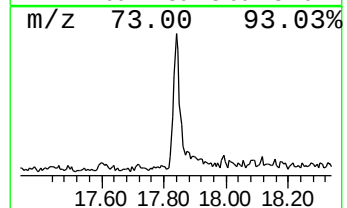
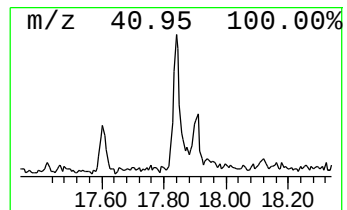
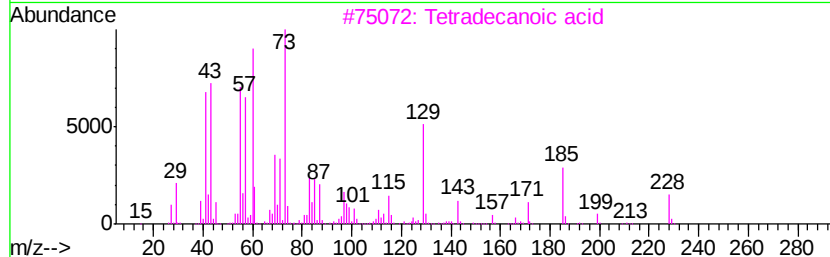
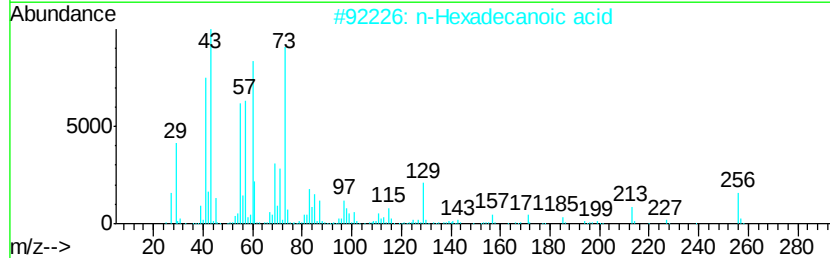
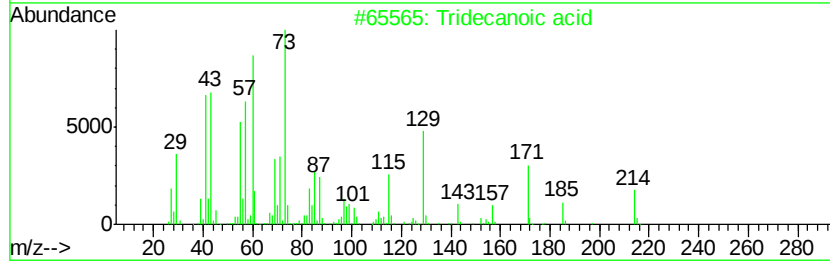
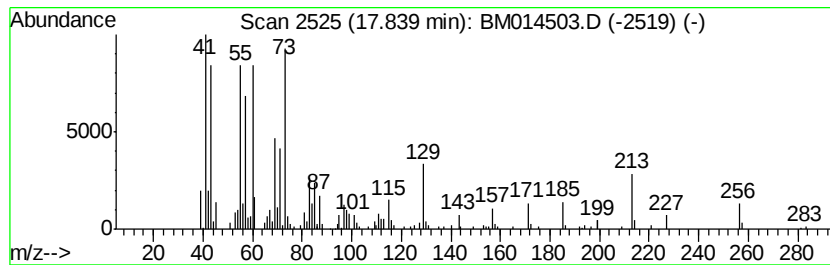
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Tridecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD		R.T.
17.84	10.16 ng	505405	Phenanthrene-d10		16.94
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecanoic acid	214	C13H26O2	000638-53-9	95
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	74
4	Pentadecanoic acid	242	C15H30O2	001002-84-2	53
5	n-Decanoic acid	172	C10H20O2	000334-48-5	43



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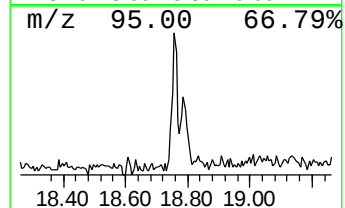
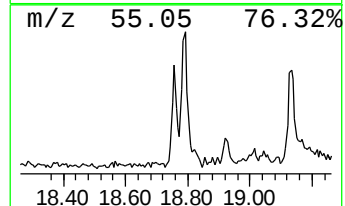
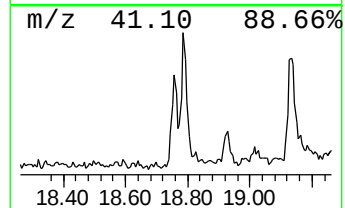
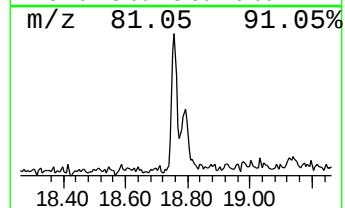
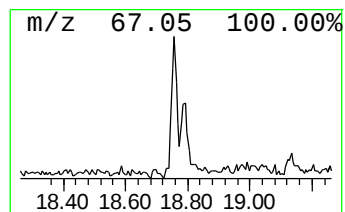
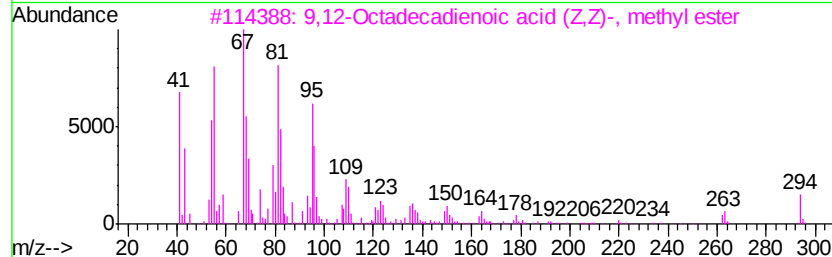
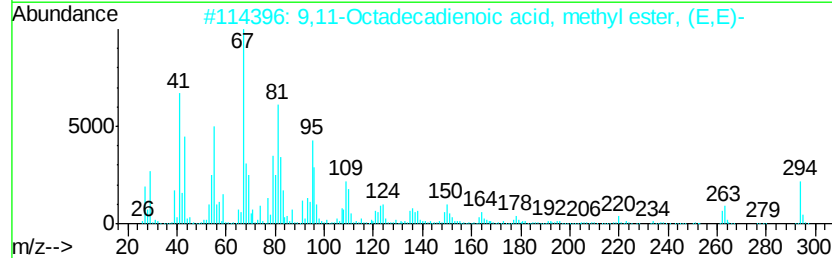
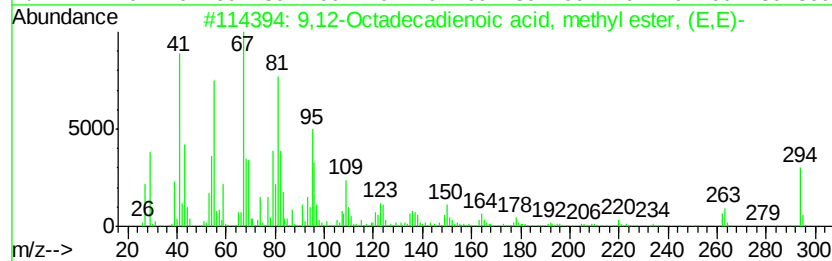
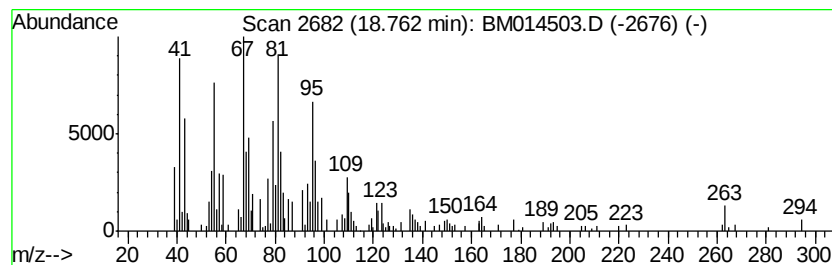
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 9,12-Octadecadienoic acid, ... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.76	6.78 ng	337557	Phenanthrene-d10	16.94	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,12-Octadecadienoic acid, methy...	294	C19H34O2	002566-97-4	97
2	9,11-Octadecadienoic acid, methy...	294	C19H34O2	013038-47-6	91
3	9,12-Octadecadienoic acid (Z,Z)-...	294	C19H34O2	000112-63-0	91
4	9,12-Octadecadienoic acid, methy...	294	C19H34O2	002462-85-3	89
5	9,15-Octadecadienoic acid, methy...	294	C19H34O2	017309-05-6	81



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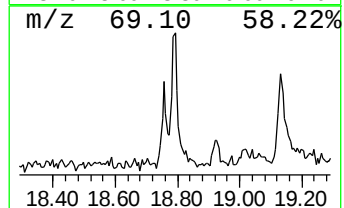
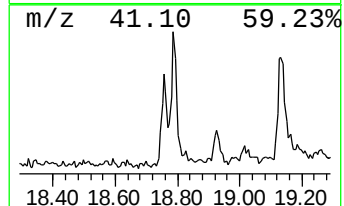
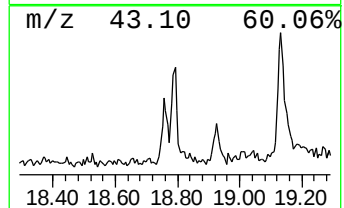
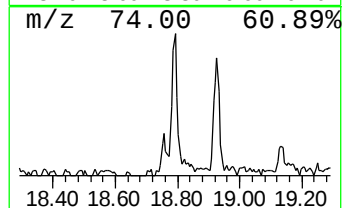
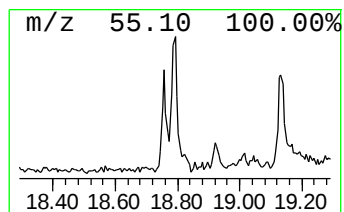
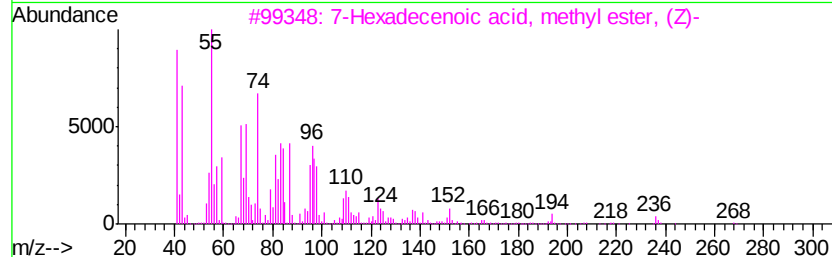
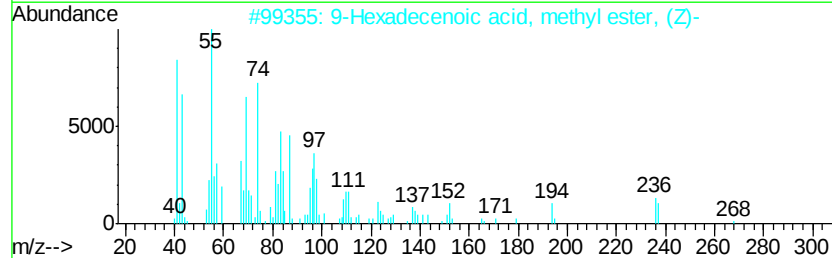
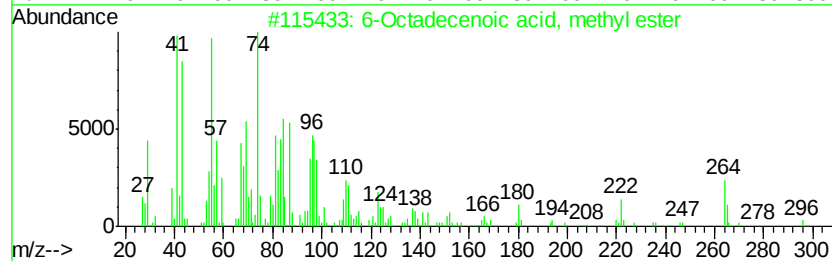
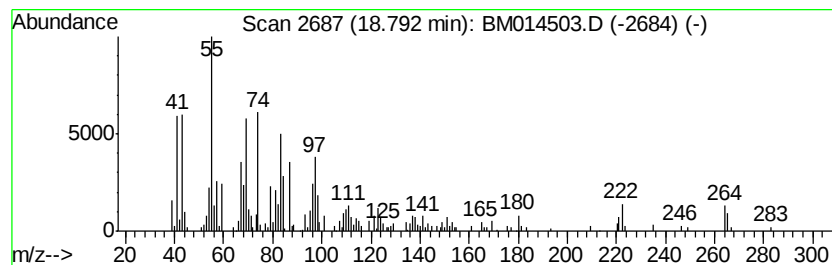
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Peak Number 6 6-Octadecenoic acid, methyl... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.79	6.88 ng	342303	Phenanthrene-d10	16.94

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		6-Octadecenoic acid, methyl ester	296	C19H36O2	052355-31-4	80
2		9-Hexadecenoic acid, methyl ester...	268	C17H32O2	001120-25-8	72
3		7-Hexadecenoic acid, methyl ester...	268	C17H32O2	056875-67-3	64
4		5-Octadecenoic acid, methyl ester	296	C19H36O2	056554-45-1	58
5		14-Octadecenoic acid, methyl ester	296	C19H36O2	056554-48-4	53



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Misc :
ALS Vial : 5 Sample Multiplier: 1

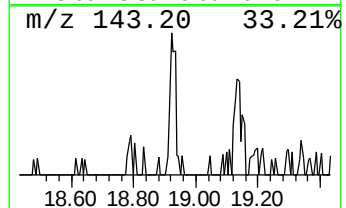
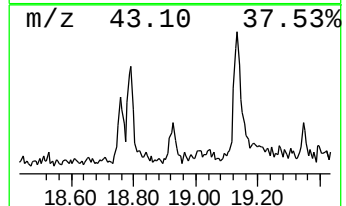
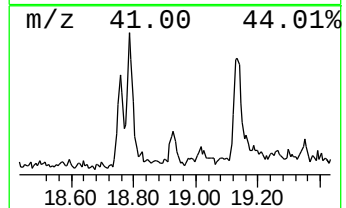
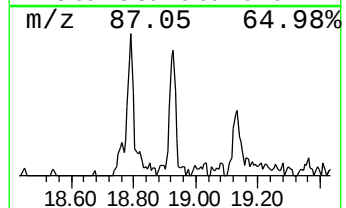
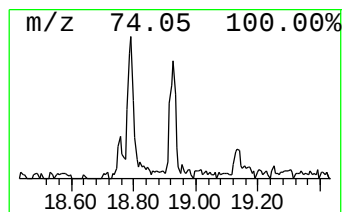
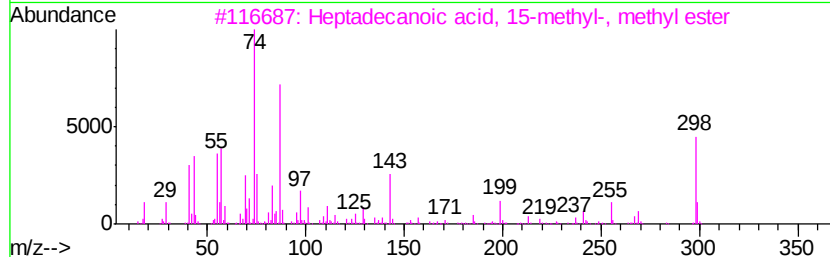
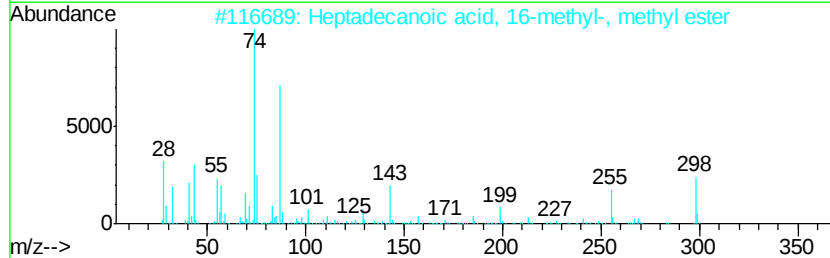
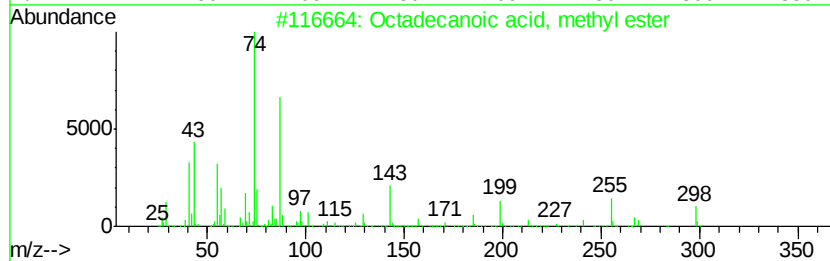
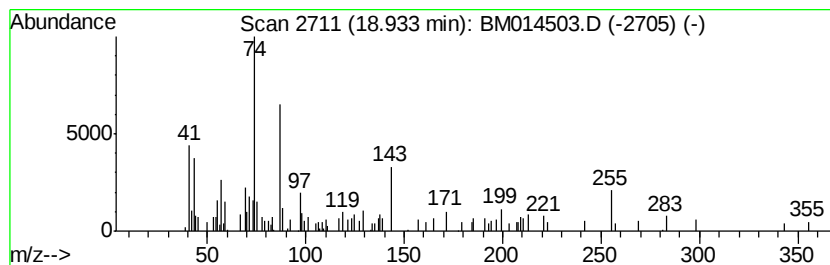
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BNA_M
ClientSampled :
B5

Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-BM020718.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Octadecanoic acid, methyl e... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.		
18.93	2.02 ng	100545	Phenanthrene-d10	16.94		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octadecanoic acid, methyl ester	298	C19H38O2	000112-61-8	93
2		Heptadecanoic acid, 16-methyl-, ...	298	C19H38O2	005129-61-3	90
3		Heptadecanoic acid, 15-methyl-, ...	298	C19H38O2	054833-55-5	72
4		Dodecanoic acid, 10-methyl-, met...	228	C14H28O2	005129-65-7	72
5		Tridecanoic acid, 12-methyl-, me...	242	C15H30O2	005129-58-8	72



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA M\DATA\BM030718\
Data File : BM014503.D
Acq On : 06 Mar 2018 18:11
Operator : SJ/JU
Sample : J1699-13
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
B5

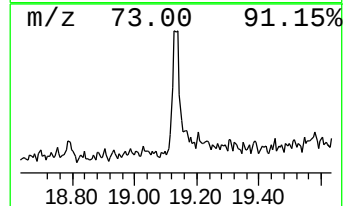
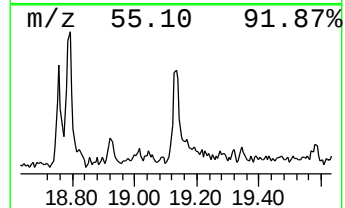
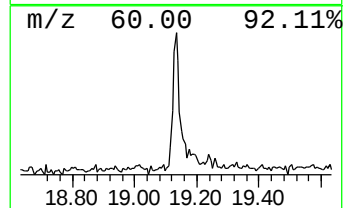
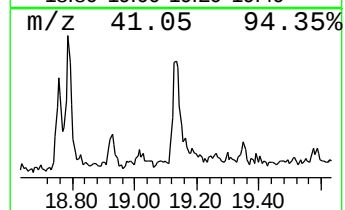
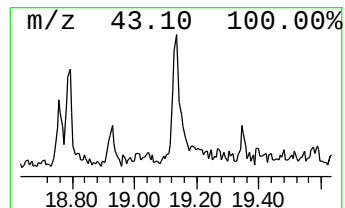
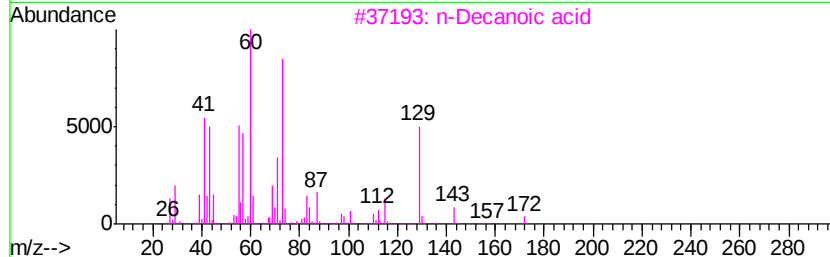
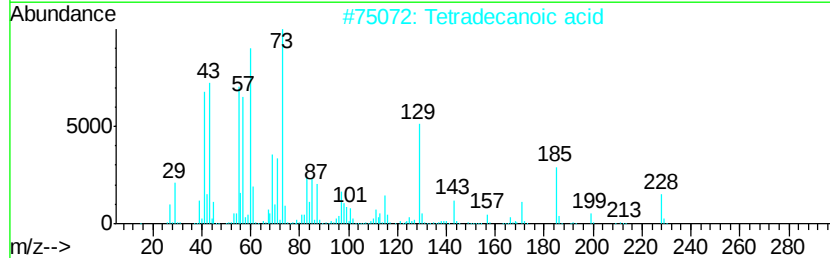
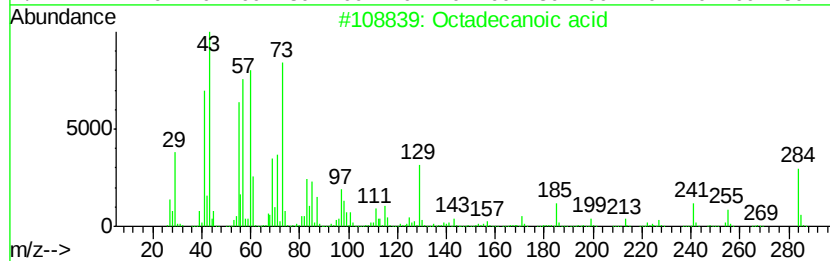
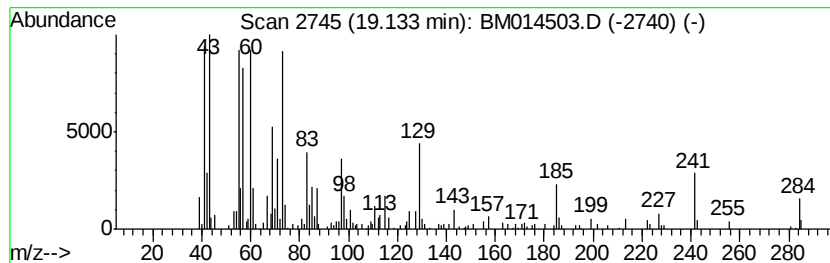
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Octadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.13	8.94 ng	436543	Chrysene-d12	21.15

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanoic acid	284	C18H36O2	000057-11-4	93
2		Tetradecanoic acid	228	C14H28O2	000544-63-8	76
3		n-Decanoic acid	172	C10H20O2	000334-48-5	70
4		Undecanoic acid	186	C11H22O2	000112-37-8	70
5		Pentadecanoic acid	242	C15H30O2	001002-84-2	60



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA M\DATA\BM030718\
Data File : BM014503.D
Acq On : 06 Mar 2018 18:11
Operator : SJ/JU
Sample : J1699-13
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
B5

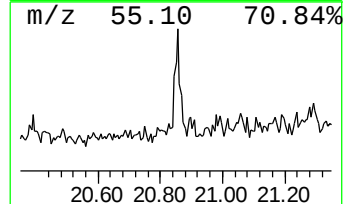
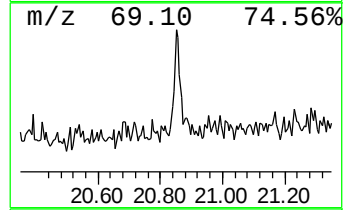
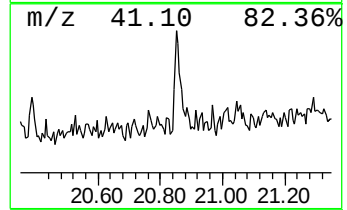
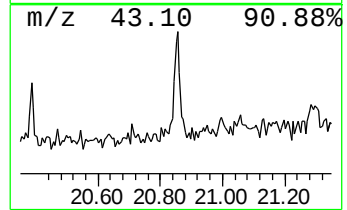
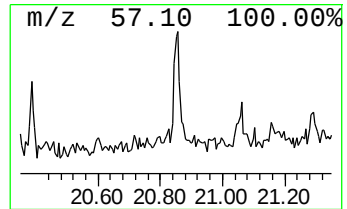
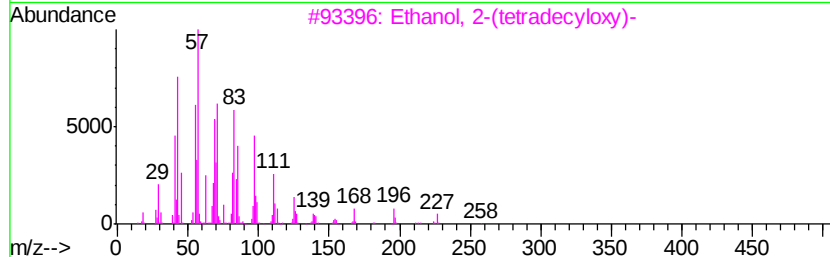
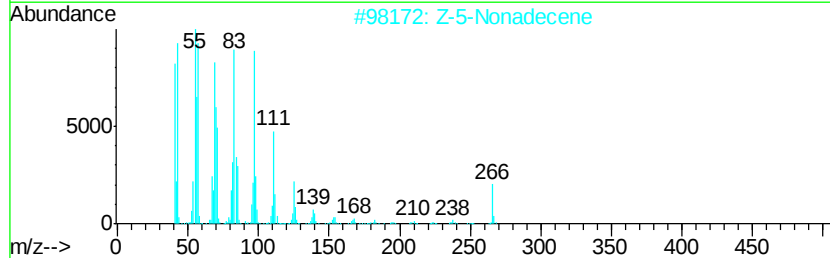
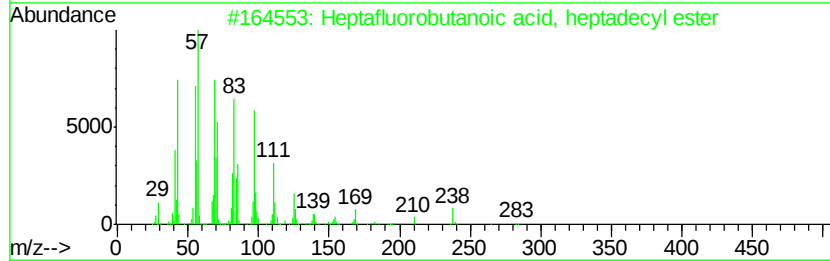
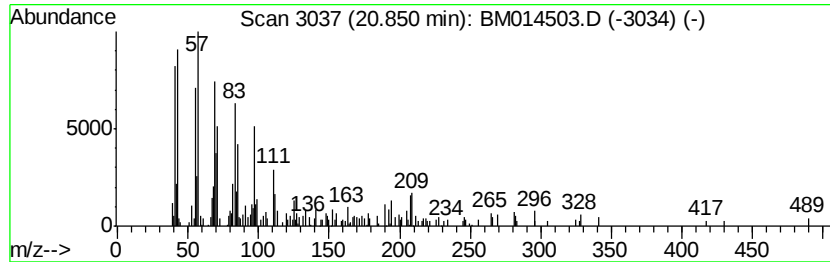
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Heptafluorobutanoic acid, h... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.85	3.28 ng	160336	Chrysene-d12	21.15

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptafluorobutanoic acid, heptad...	452	C21H35F7O2	1000282-97-3	81
2			Z-5-Nonadecene	266	C19H38	1000131-11-8	62
3			Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	60
4			Cyclopentane, (4-octyldodecyl)-	350	C25H50	005638-09-5	53
5			1-Hexadecanesulfonyl chloride	324	C16H33ClO2S	038775-38-1	49



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_M\DATA\BM030718\
Data File : BM014503.D
Acq On : 06 Mar 2018 18:11
Operator : SJ/JU
Sample : J1699-13
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
B5

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-BM020718.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
unknown7.05	7.05	75.1	ng	2640570	1	7.50	703630	20.0
Methyl tetradecan...	17.60	3.1	ng	155662	4	16.94	995266	20.0
Tridecanoic acid	17.84	10.2	ng	505405	4	16.94	995266	20.0
9,12-Octadecadien...	18.76	6.8	ng	337557	4	16.94	995266	20.0
6-Octadecenoic ac...	18.79	6.9	ng	342303	4	16.94	995266	20.0
Octadecanoic acid...	18.93	2.0	ng	100545	4	16.94	995266	20.0
Octadecanoic acid	19.13	8.9	ng	436543	5	21.15	977124	20.0
Heptafluorobutano...	20.85	3.3	ng	160336	5	21.15	977124	20.0