

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM121518\
 Data File : BM018211.D
 Acq On : 15 Dec 2018 19:31
 Operator : JU/SJ
 Sample : J6340-02
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SB-02B

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

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	4.687	302	306	314	rBV	88094	134509	1.55%	0.289%
2	5.134	375	382	403	rBV	2579287	3746831	43.16%	8.061%
3	6.704	641	649	668	rBV	2589360	4349825	50.11%	9.358%
4	7.040	699	706	722	rBV	2790609	4302145	49.56%	9.256%
5	7.498	778	784	792	rBV	423690	671305	7.73%	1.444%
6	7.810	830	837	855	rBV	1896507	3080322	35.48%	6.627%
7	8.657	971	981	994	rBV	1450854	2459962	28.34%	5.292%
8	10.269	1248	1255	1272	rBV	475254	927870	10.69%	1.996%
9	12.769	1672	1680	1700	rBV	3663029	5683282	65.47%	12.227%
10	13.633	1822	1827	1839	rBV	161944	308864	3.56%	0.664%
11	14.157	1910	1916	1932	rBV2	799057	1231912	14.19%	2.650%
12	15.663	2166	2172	2191	rBV	3239182	4861642	56.00%	10.459%
13	16.921	2380	2386	2401	rBV	935684	1433516	16.51%	3.084%
14	17.833	2535	2541	2558	rBV	436163	654580	7.54%	1.408%
15	19.127	2756	2761	2772	rBV	243228	391915	4.51%	0.843%
16	19.574	2831	2837	2847	rBV	7058253	8680813	100.00%	18.676%
17	20.862	3052	3056	3065	rBV	233869	335891	3.87%	0.723%
18	21.139	3098	3103	3111	rBV	1247818	1733956	19.97%	3.730%
19	23.303	3465	3471	3479	rVB2	812210	1492889	17.20%	3.212%

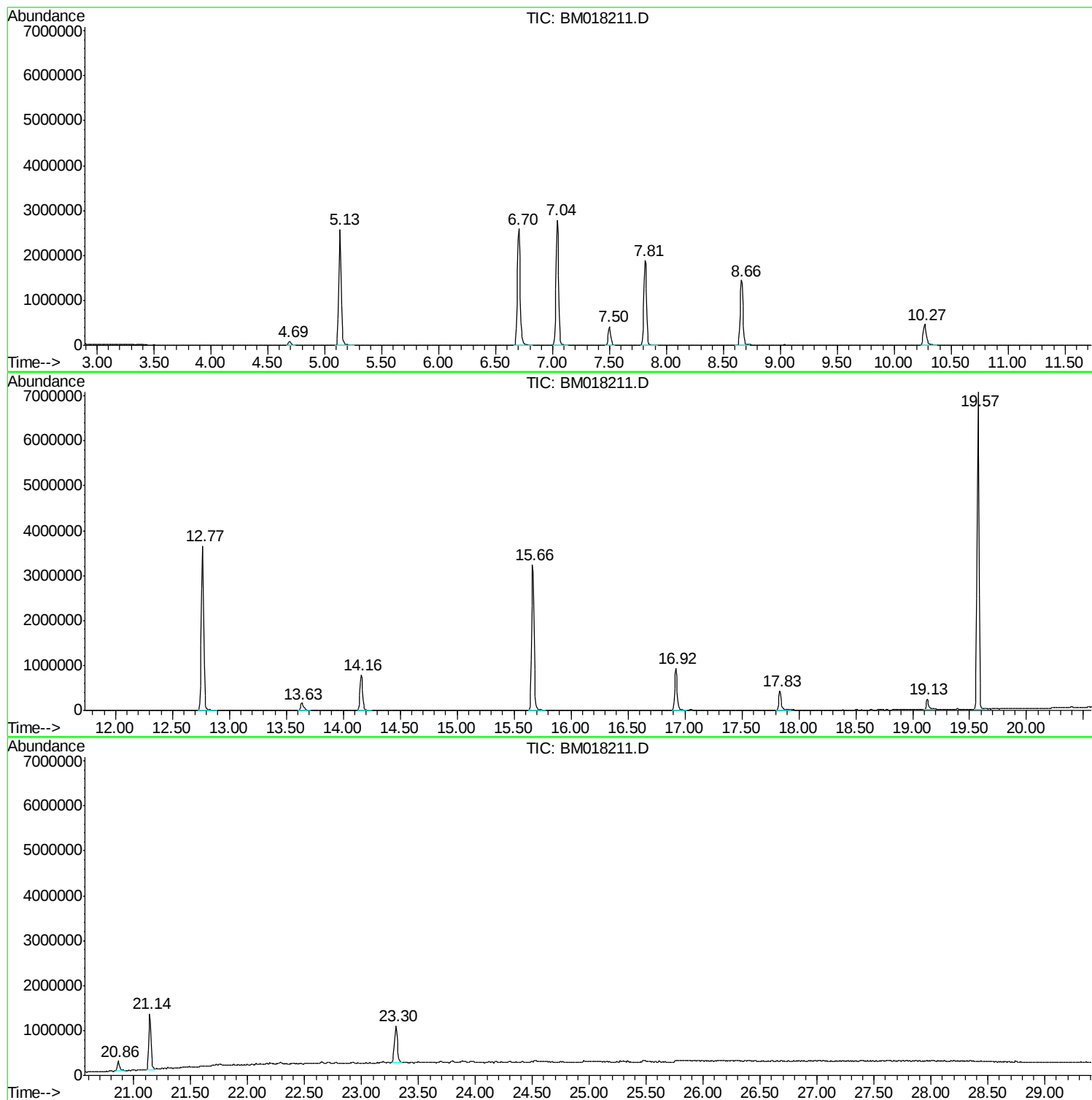
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM121518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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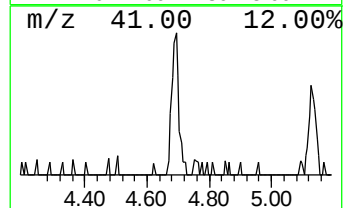
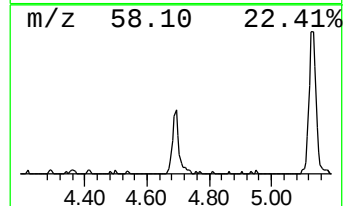
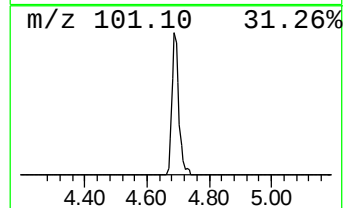
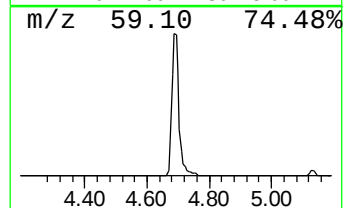
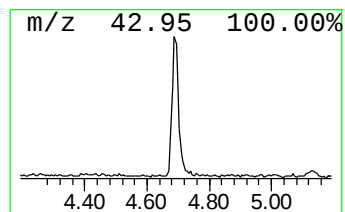
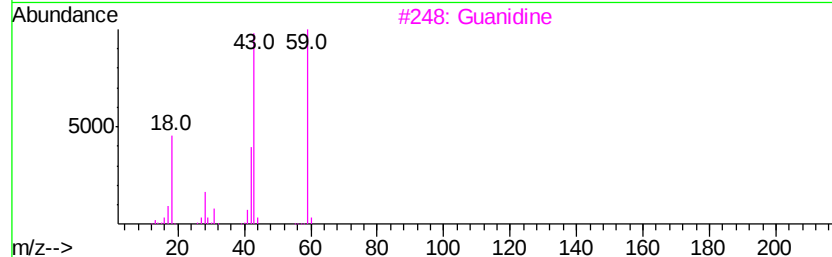
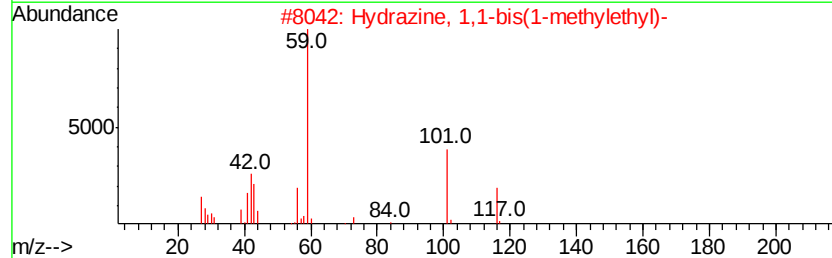
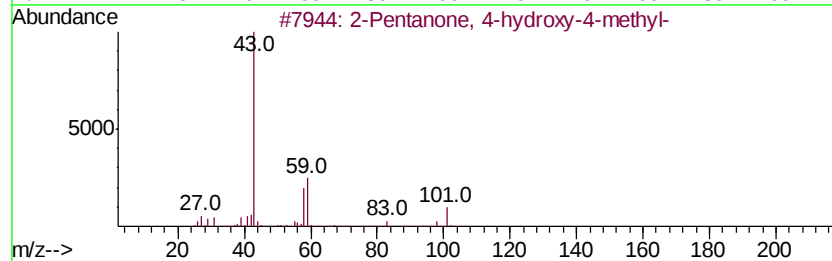
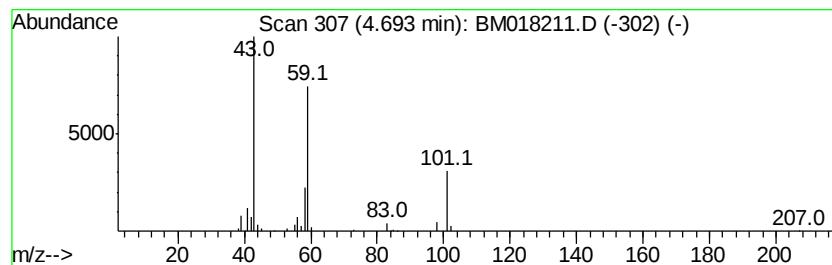
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM121518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.69	4.01 ng	134509	1,4-Dichlorobenzene-d4	7.50

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			Hydrazine, 1,1-bis(1-methylethyl)-	116	C6H16N2	000921-14-2	47
3			Guanidine	59	CH5N3	000113-00-8	43
4			(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7N02	016504-58-8	38
5			Propane, 1,1-dipropoxy-	160	C9H20O2	004744-11-0	36



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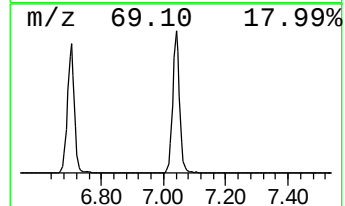
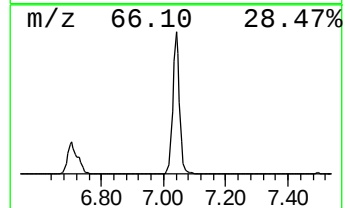
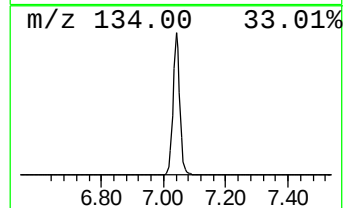
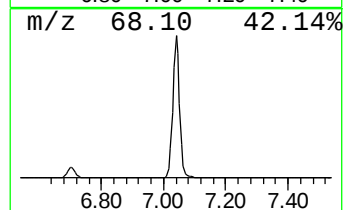
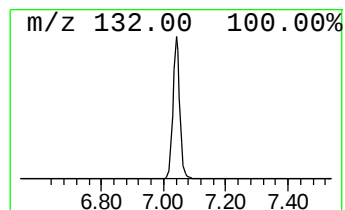
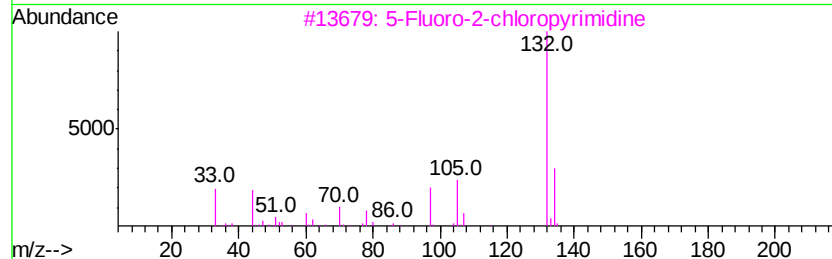
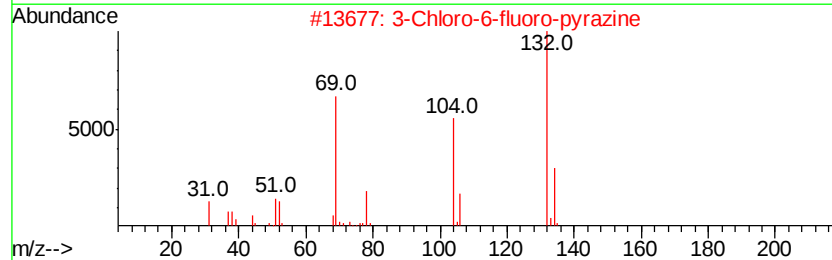
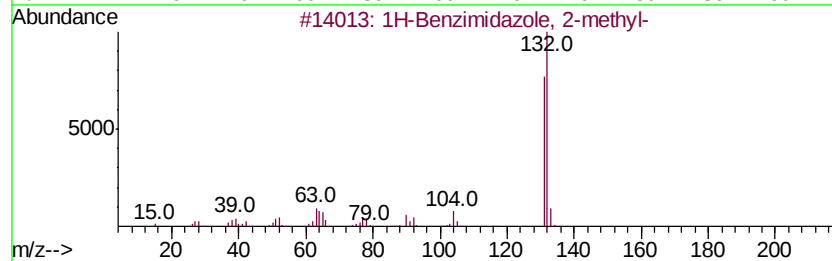
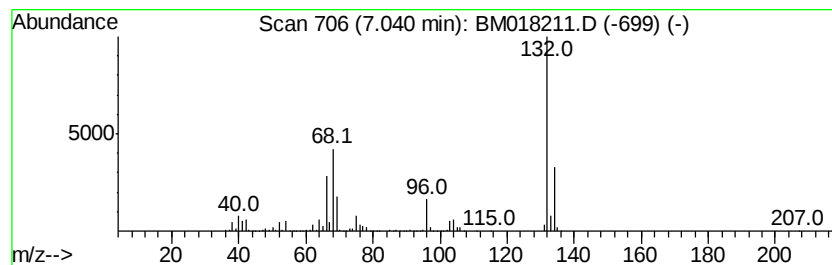
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 2 unknown7.04 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.04	128.17 ng	4302150	1,4-Dichlorobenzene-d4	7.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		(5-METHYL-2-PYRIDYL)ACETONITRILE	132	C8H8N2	1000241-93-9	10



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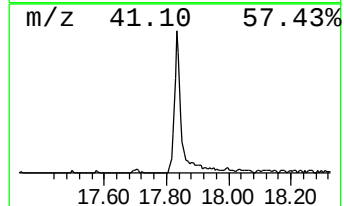
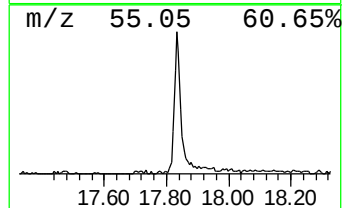
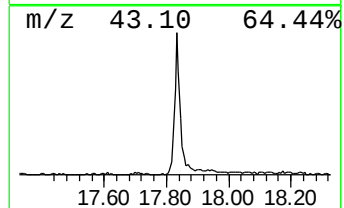
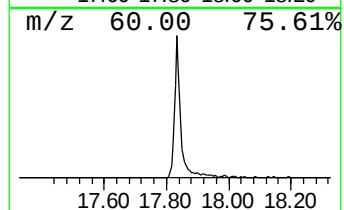
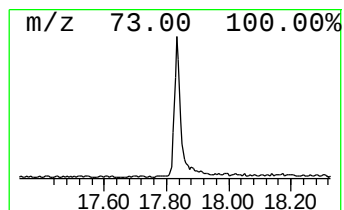
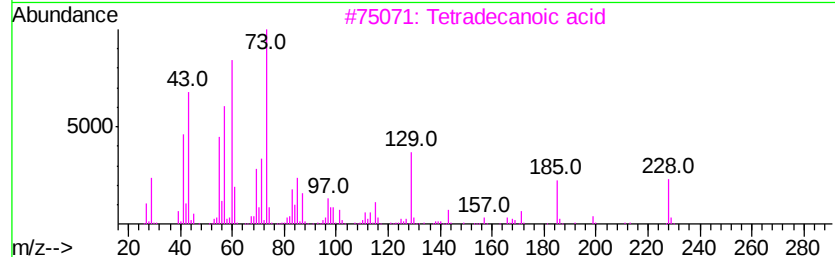
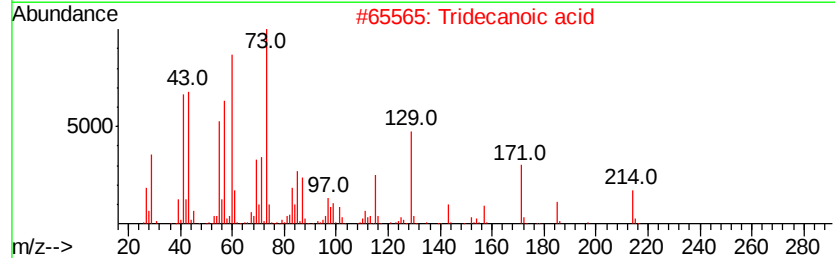
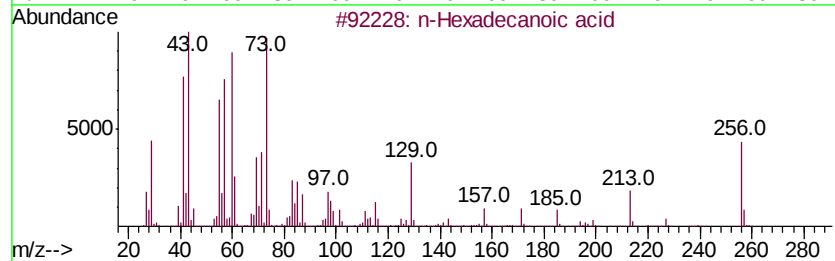
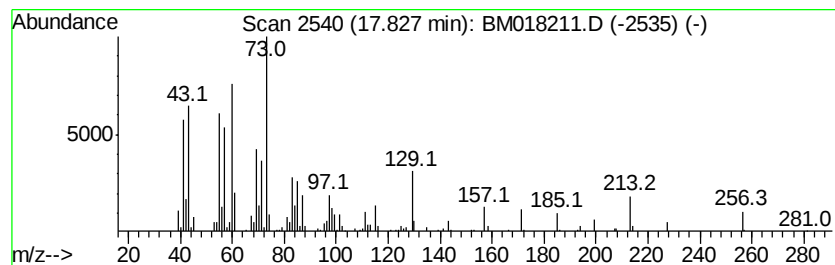
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Peak Number 4 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.83	9.13 ng	654580	Phenanthrene-d10	16.92

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
2		Tridecanoic acid	214	C13H26O2	000638-53-9	93
3		Tetradecanoic acid	228	C14H28O2	000544-63-8	90
4		n-Decanoic acid	172	C10H20O2	000334-48-5	70
5		Dodecanoic acid	200	C12H24O2	000143-07-7	68



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Instrument :
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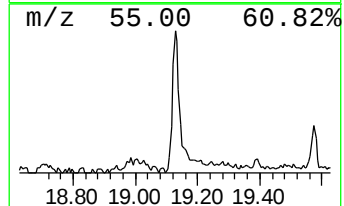
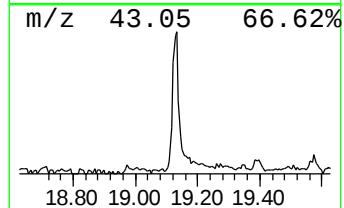
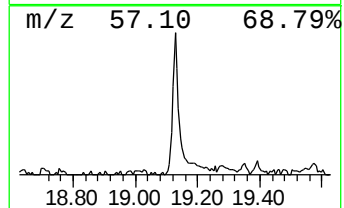
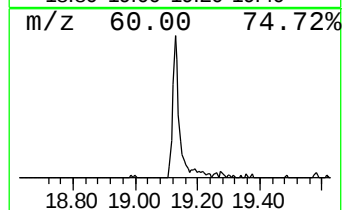
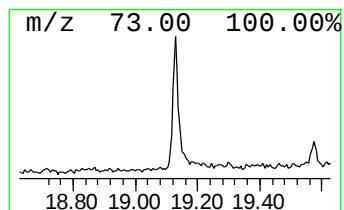
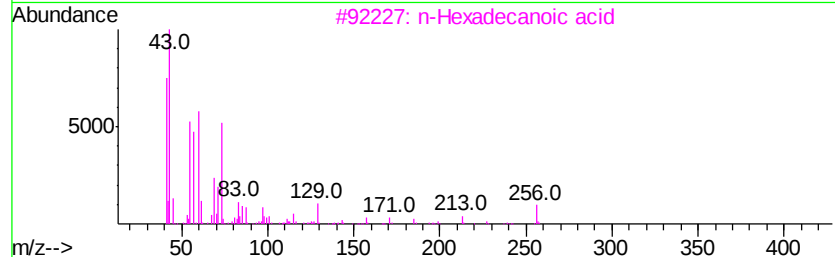
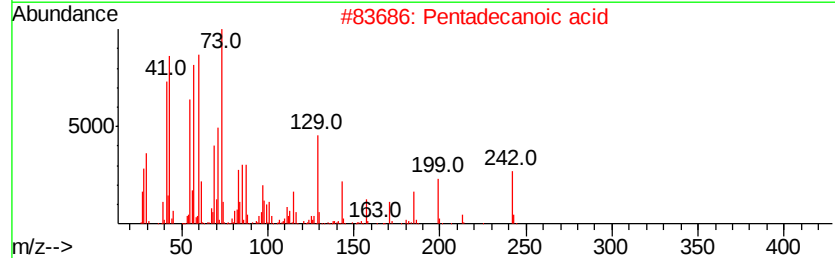
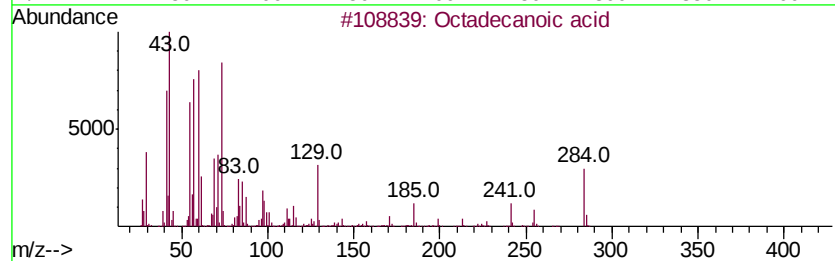
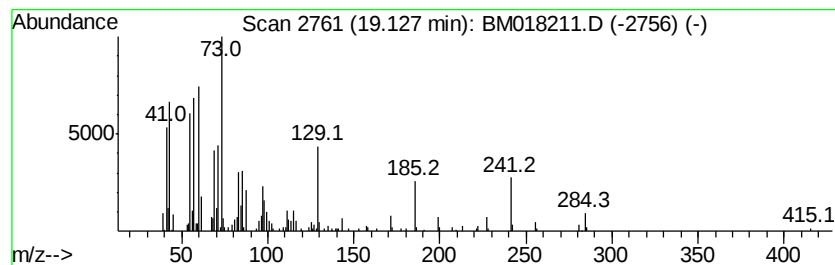
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Peak Number 5 Octadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.13	4.52 ng	391915	Chrysene-d12	21.14

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	90
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	70
5		n-Decanoic acid	172	C10H20O2	000334-48-5	62



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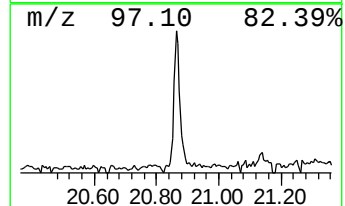
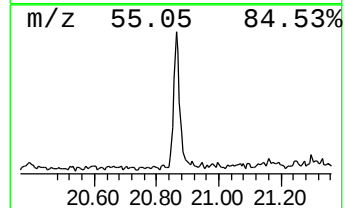
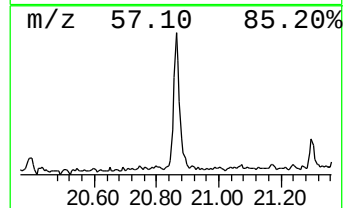
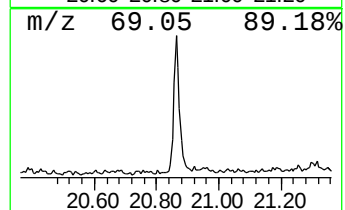
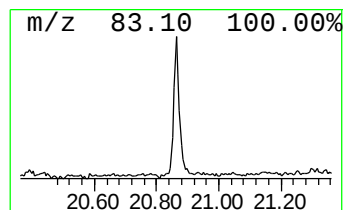
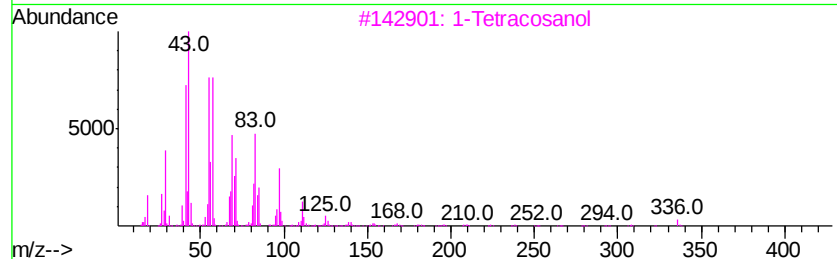
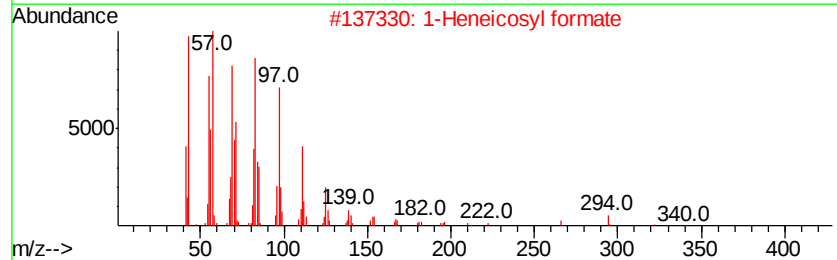
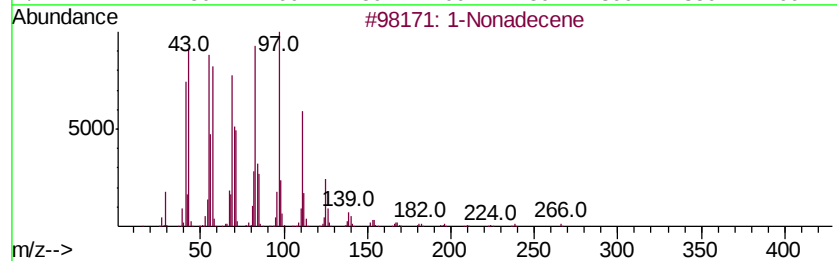
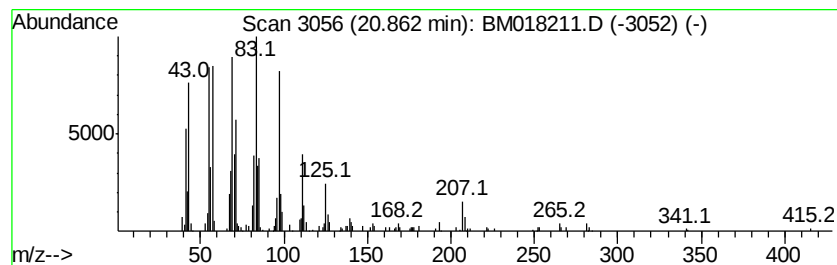
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 Peak Number 6 1-Nonadecene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD		R.T.
20.86	3.87 ng	335891	Chrysene-d12		21.14
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Nonadecene	266	C19H38	018435-45-5	92
2	1-Heneicosyl formate	340	C22H44O2	077899-03-7	91
3	1-Tetracosanol	354	C24H50O	000506-51-4	91
4	1-Nonadecanol	284	C19H40O	001454-84-8	90
5	Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	90



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.69	4.0	ng	134509	1	7.50	671305	20.0
unknown7.04	7.04	128.2	ng	4302150	1	7.50	671305	20.0
n-Hexadecanoic acid	17.83	9.1	ng	654580	4	16.92	1433520	20.0
Octadecanoic acid	19.13	4.5	ng	391915	5	21.14	1733960	20.0
1-Nonadecene	20.86	3.9	ng	335891	5	21.14	1733960	20.0