

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080618\
 Data File : BM016201.D
 Acq On : 06 Aug 2018 21:10
 Operator : SJ/JU
 Sample : J4333-13
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SOIL-4

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-BM072318.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	5.216	323	328	343	rVB	114102	173301	6.86%	1.158%
2	5.693	403	409	420	rBV	820963	1247383	49.40%	8.334%
3	7.334	682	688	707	rBV	833950	1338967	53.03%	8.946%
4	7.698	743	750	765	rBV	909629	1458890	57.77%	9.747%
5	8.169	825	830	839	rBV	177623	291222	11.53%	1.946%
6	8.493	878	885	897	rBB	657363	1064885	42.17%	7.115%
7	9.357	1026	1032	1053	rBV	486674	816057	32.32%	5.452%
8	10.987	1303	1309	1317	rBV	210202	354226	14.03%	2.367%
9	13.422	1717	1723	1735	rVB	1149572	1599642	63.35%	10.688%
10	14.251	1859	1864	1872	rBB	83787	116388	4.61%	0.778%
11	14.804	1952	1958	1965	rBB2	269207	404876	16.03%	2.705%
12	16.286	2204	2210	2225	rBV	1079002	1422123	56.32%	9.502%
13	17.533	2417	2422	2428	rBV	346456	487563	19.31%	3.258%
14	18.380	2561	2566	2577	rBV	109166	174932	6.93%	1.169%
15	19.568	2765	2768	2774	rVB	27580	33421	1.32%	0.223%
16	19.639	2776	2780	2785	rBV3	46031	59053	2.34%	0.395%
17	19.933	2827	2830	2834	rVB	24491	29048	1.15%	0.194%
18	20.109	2855	2860	2866	rBV	2074269	2525160	100.00%	16.872%
19	21.351	3066	3071	3079	rVB3	73790	138394	5.48%	0.925%
20	21.662	3119	3124	3129	rBV	490463	633037	25.07%	4.230%
21	24.015	3518	3524	3532	rVB	313251	598299	23.69%	3.997%

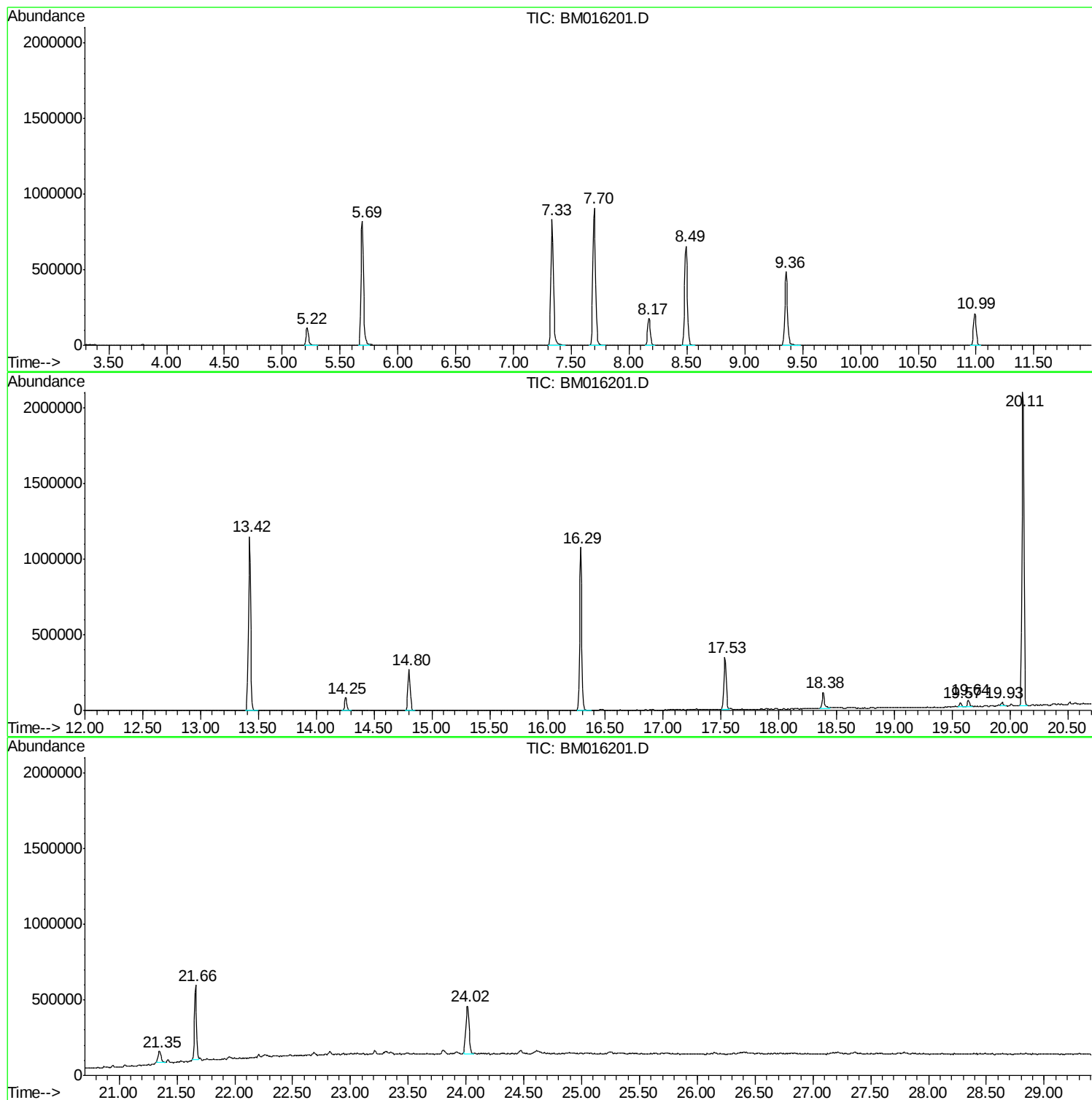
Sum of corrected areas: 14966867

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM072318.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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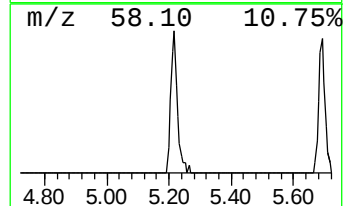
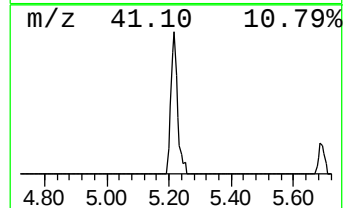
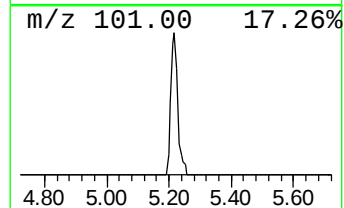
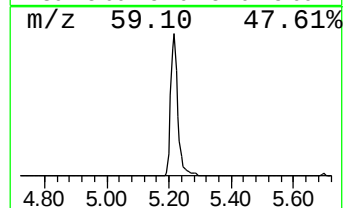
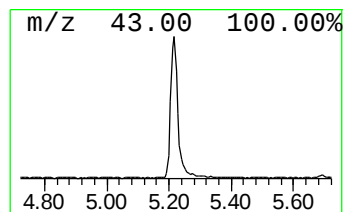
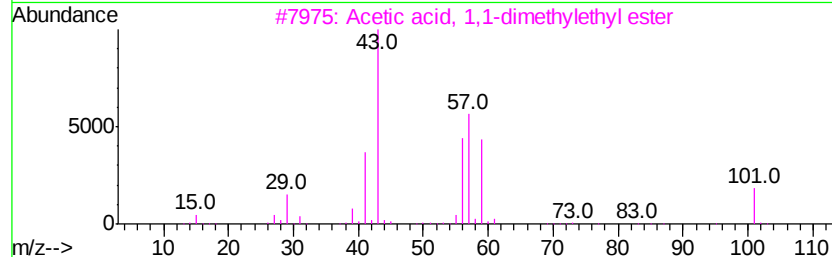
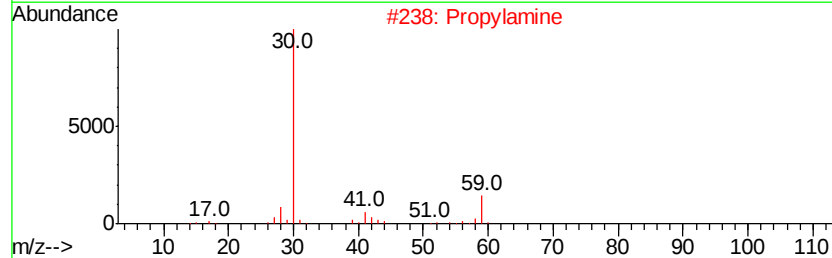
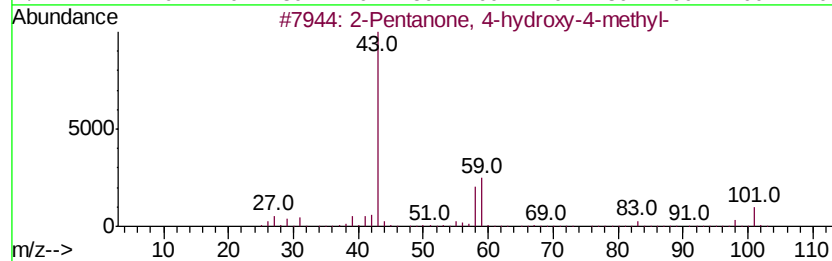
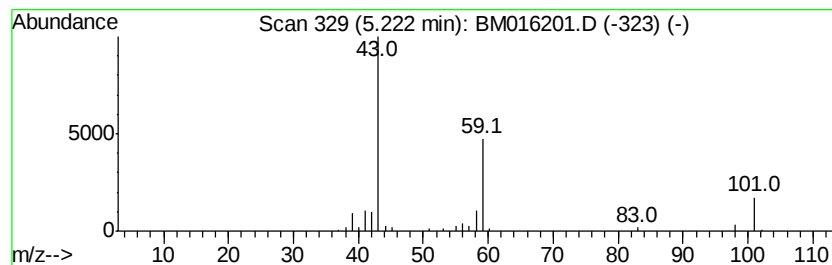
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM072318.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.22	11.90 ng	173301	1,4-Dichlorobenzene-d4	8.17

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			Propylamine	59	C3H9N	000107-10-8	40
3			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38
4			2-Nonanone, 9-hydroxy-	158	C9H18O2	025368-56-3	36
5			Butane	58	C4H10	000106-97-8	22



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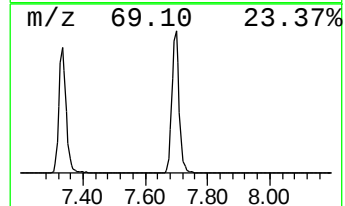
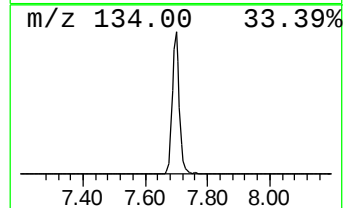
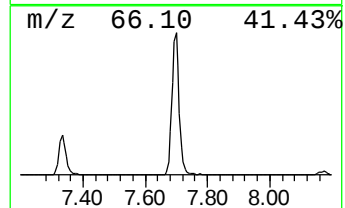
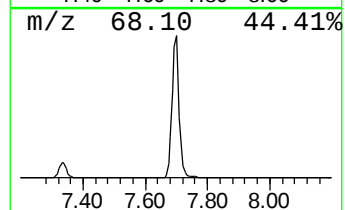
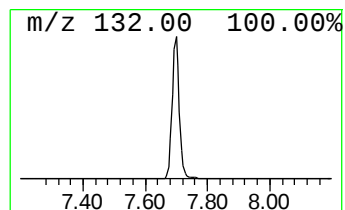
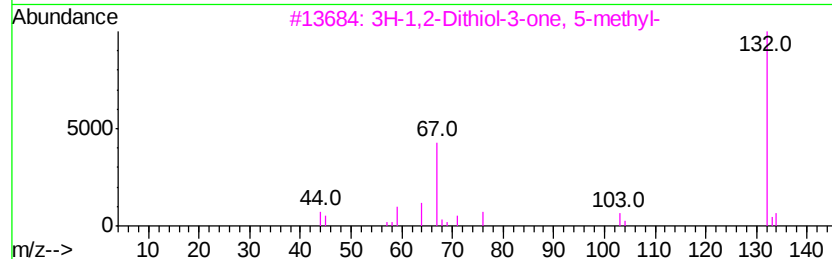
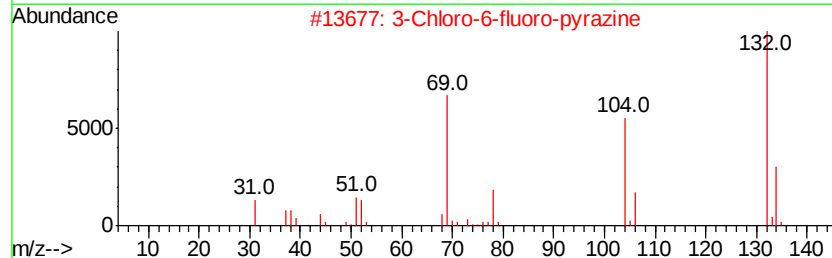
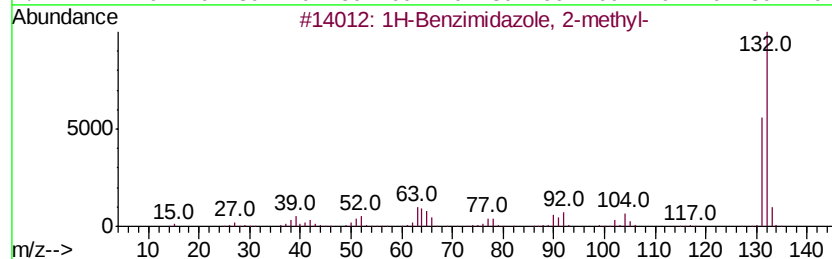
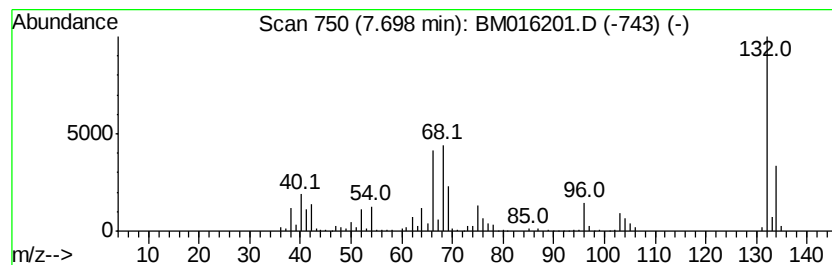
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown7.70 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.70	100.19 ng	1458890	1,4-Dichlorobenzene-d4	8.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	14
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	14
3		3H-1,2-Dithiol-3-one, 5-methyl-	132	C4H4OS2	003620-08-4	14
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	14
5		1,3,4-Thiadiazole-2(3H)-thione, ...	132	C3H4N2S2	029490-19-5	14



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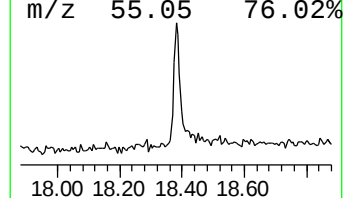
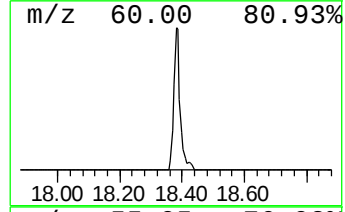
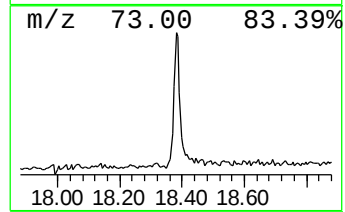
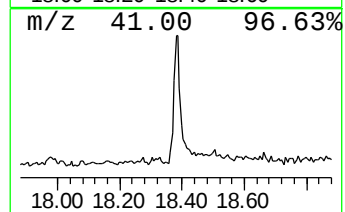
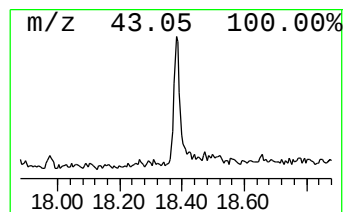
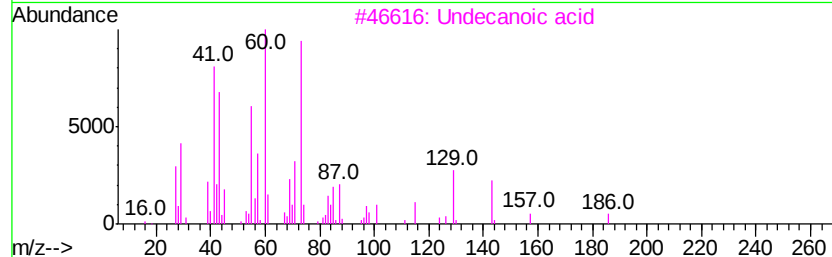
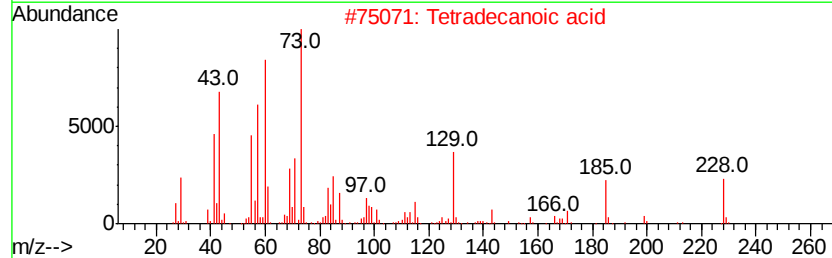
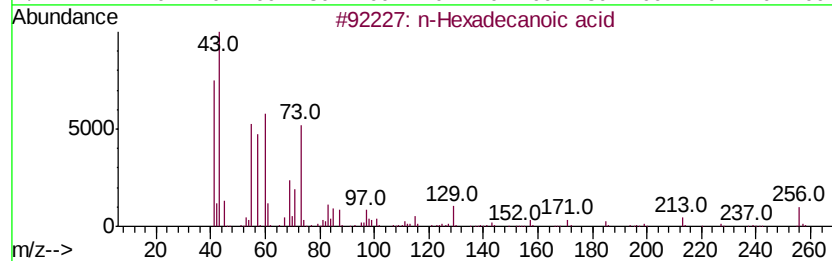
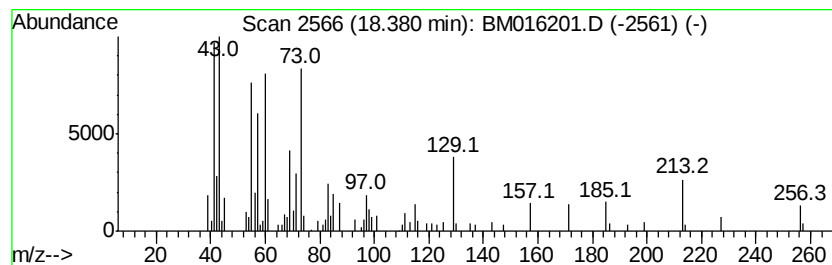
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Peak Number 4 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.38	7.18 ng	174932	Phenanthrene-d10	17.53

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	83
2		Tetradecanoic acid	228	C14H28O2	000544-63-8	74
3		Undecanoic acid	186	C11H22O2	000112-37-8	70
4		Tridecanoic acid	214	C13H26O2	000638-53-9	64
5		Dodecanoic acid	200	C12H24O2	000143-07-7	58



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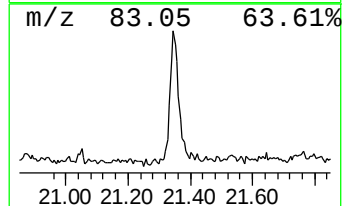
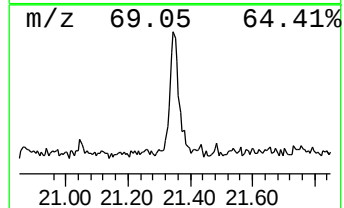
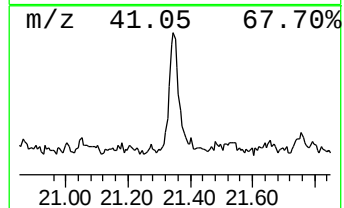
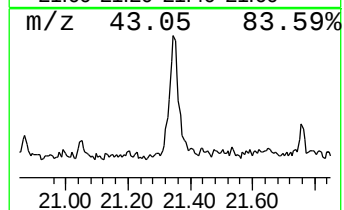
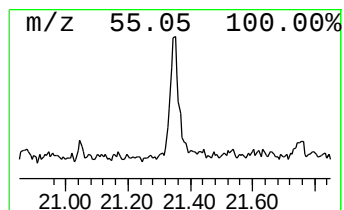
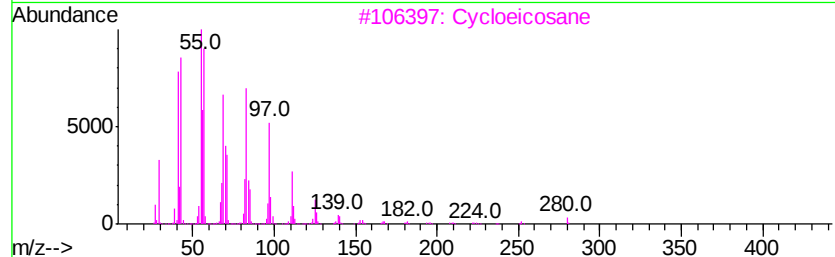
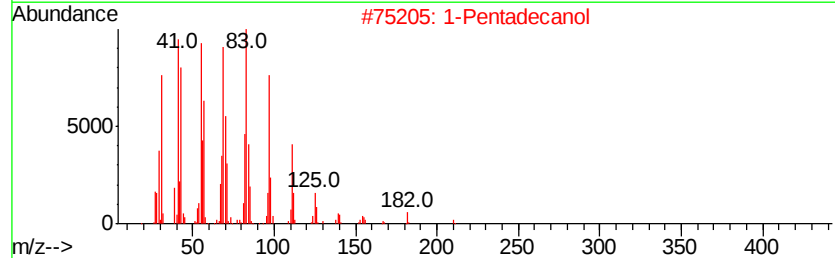
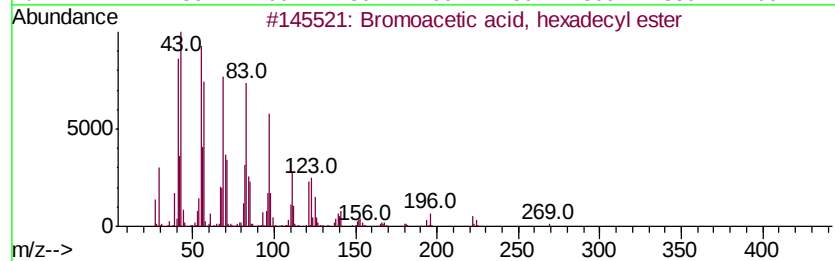
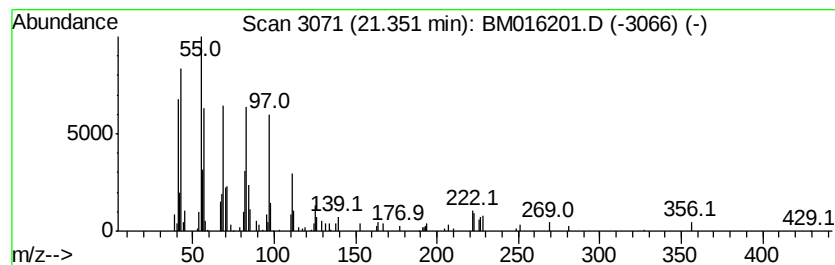
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Peak Number 5 Bromoacetic acid, hexadecyl... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.35	4.37 ng	138394	Chrysene-d12	21.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Bromoacetic acid, hexadecyl ester	362	C18H35BrO2	005454-48-8	80
2		1-Pentadecanol	228	C15H32O	000629-76-5	74
3		Cycloeicosane	280	C20H40	000296-56-0	74
4		Chloroacetic acid, pentadecyl ester	304	C17H33ClO2	070301-47-2	72
5		5-Eicosene, (E)-	280	C20H40	074685-30-6	72



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.22	11.9	ng	173301	1	8.17	291222	20.0
unknown7.70	7.70	100.2	ng	1458890	1	8.17	291222	20.0
n-Hexadecanoic acid	18.38	7.2	ng	174932	4	17.53	487563	20.0
Bromoacetic acid,...	21.35	4.4	ng	138394	5	21.66	633037	20.0