

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF012318\
 Data File : BF102370.D
 Acq On : 24 Jan 2018 4:03
 Operator : SJ/JU
 Sample : J1249-13
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 AB-7(0-2)

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_F\METHODS\8270-BF012218.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	4.916	477	481	492	rVB	61360	85529	3.18%	0.380%
2	5.328	546	551	554	rBV	2591251	2466533	91.65%	10.953%
3	6.357	721	726	743	rBV	2252450	2691244	100.00%	11.951%
4	6.487	743	748	751	rBV	2860242	2573177	95.61%	11.427%
5	6.716	783	787	795	rBV	887671	751734	27.93%	3.338%
6	6.869	809	813	817	rBV	2174889	1885623	70.07%	8.373%
7	7.281	878	883	886	rBV	1663504	1540499	57.24%	6.841%
8	7.998	1001	1005	1020	rBV	1129069	1022792	38.00%	4.542%
9	8.557	1097	1100	1104	rBV	37124	31830	1.18%	0.141%
10	9.075	1183	1188	1191	rBV	2728899	2401514	89.23%	10.664%
11	9.469	1251	1255	1262	rVB	289902	244158	9.07%	1.084%
12	9.680	1287	1291	1295	rVB	56283	43148	1.60%	0.192%
13	9.751	1299	1303	1312	rVB	1189475	1073647	39.89%	4.768%
14	10.180	1373	1376	1379	rVB	63372	49573	1.84%	0.220%
15	10.463	1419	1424	1427	rBV	75277	64272	2.39%	0.285%
16	10.539	1433	1437	1447	rVB	1335888	1211523	45.02%	5.380%
17	10.651	1453	1456	1460	rBV	45403	43281	1.61%	0.192%
18	11.233	1551	1555	1563	rBV	1076952	937630	34.84%	4.164%
19	12.639	1791	1794	1800	rBV2	46981	39326	1.46%	0.175%
20	12.716	1805	1807	1811	rVB	65554	53474	1.99%	0.237%
21	12.821	1820	1825	1828	rBV	1786414	1569249	58.31%	6.969%
22	13.286	1902	1904	1911	rVB3	24733	35190	1.31%	0.156%
23	13.710	1973	1976	1985	rBV	193023	229886	8.54%	1.021%
24	13.874	1997	2004	2013	rBV	712133	723156	26.87%	3.211%
25	14.339	2080	2083	2089	rBV7	19179	28595	1.06%	0.127%
26	15.298	2241	2246	2256	rBV	544819	656990	24.41%	2.917%
27	16.186	2393	2397	2403	rBV5	34087	65404	2.43%	0.290%

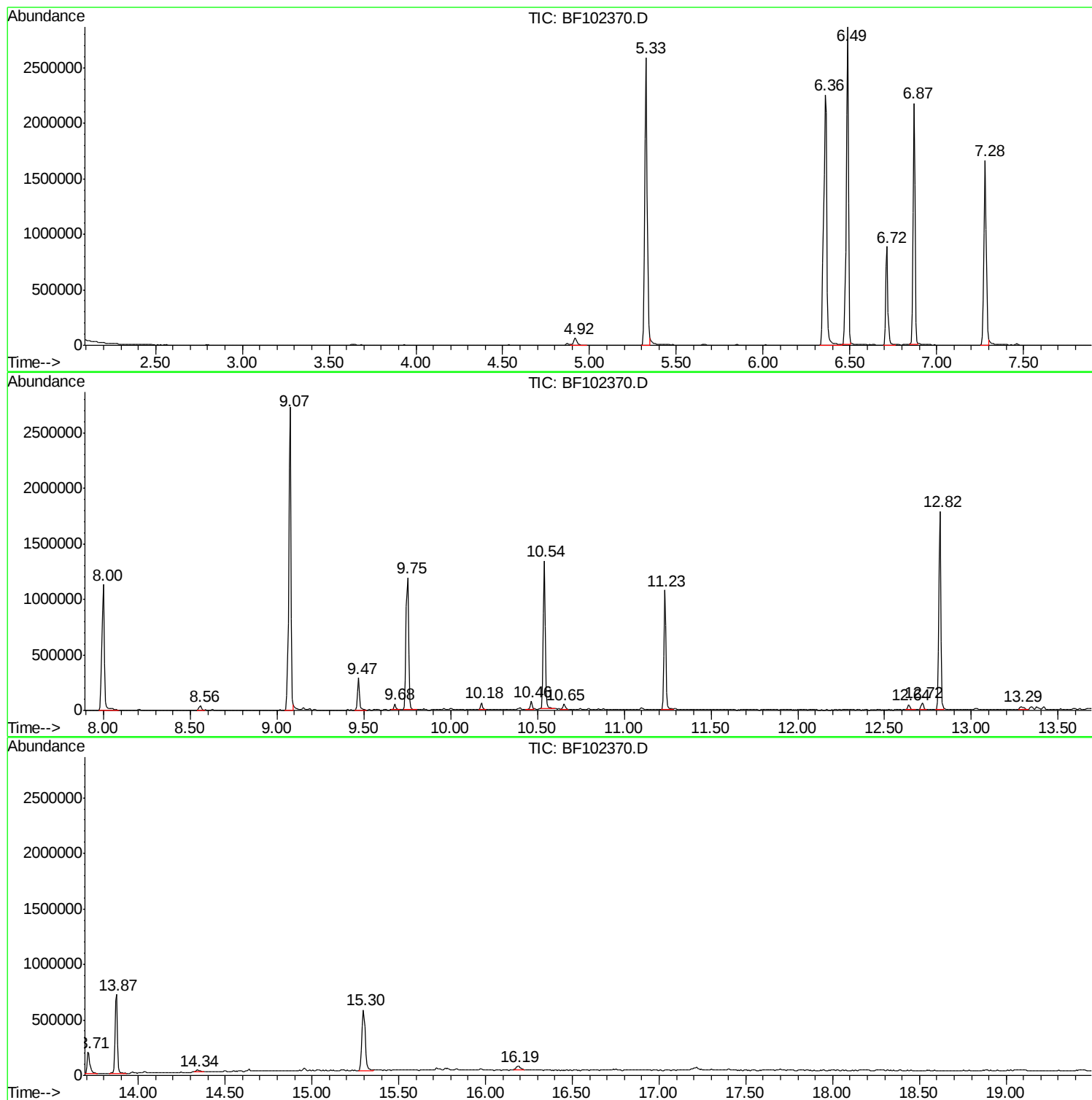
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF012218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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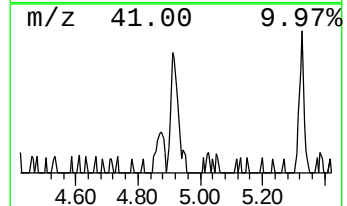
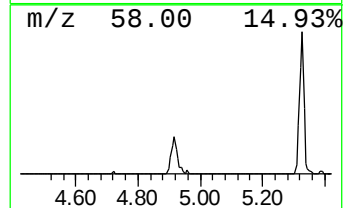
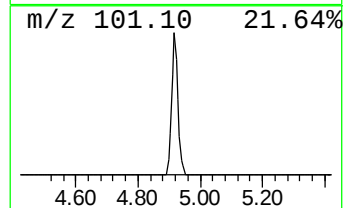
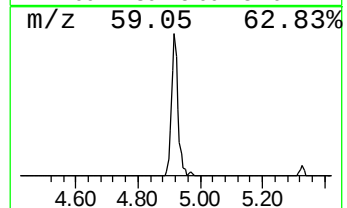
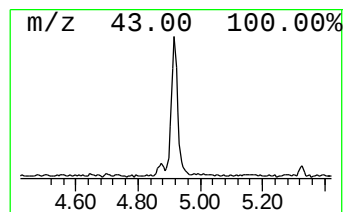
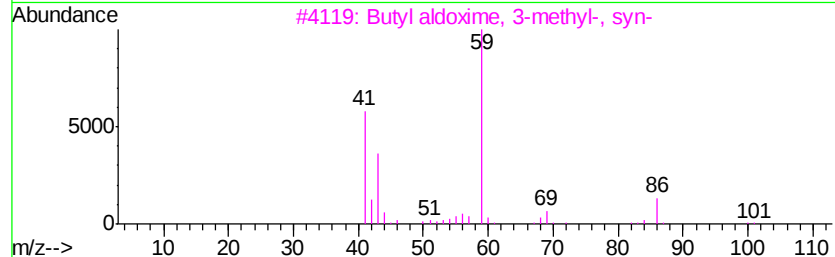
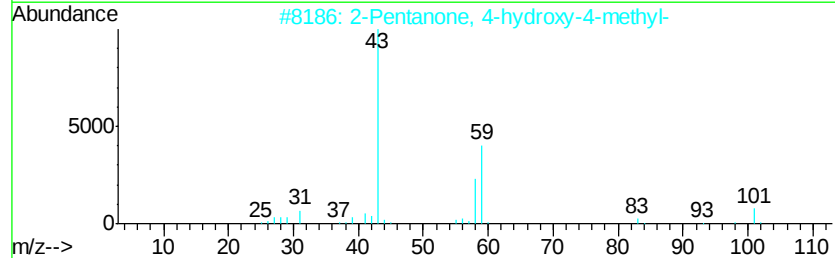
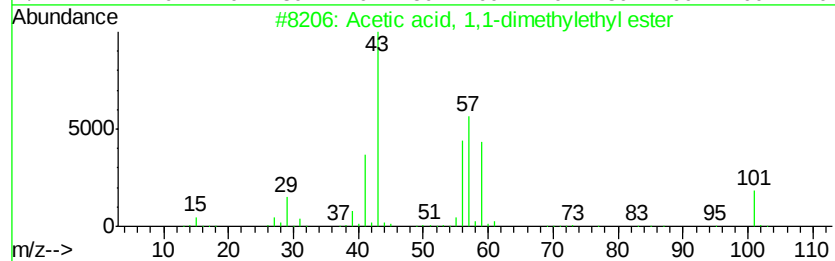
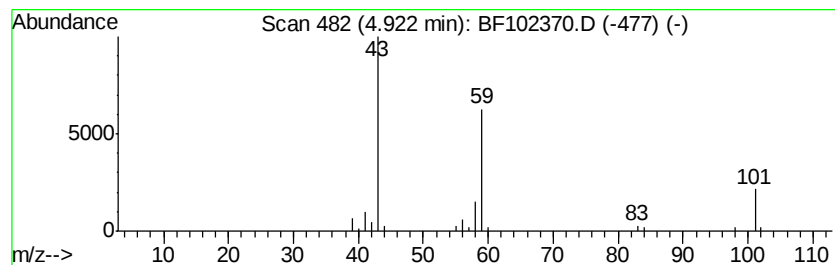
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF012218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown4.92 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.92	2.28 ng	85529	1,4-Dichlorobenzene-d4	6.72

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
2			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	36
3			Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	32
4			Ethanone, 1-(2-methylcyclopropyl)-	98	C6H10O	000930-56-3	9
5			Hexanamide	115	C6H13NO	000628-02-4	9



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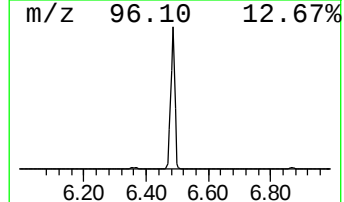
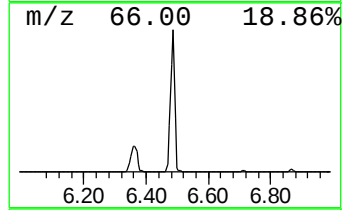
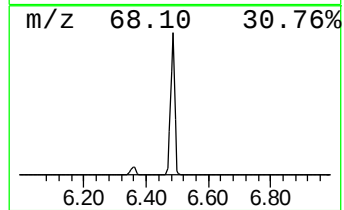
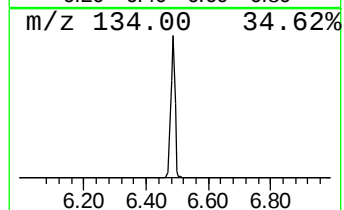
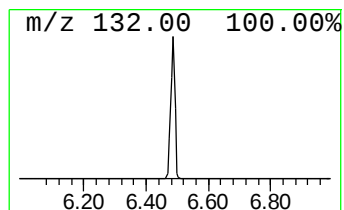
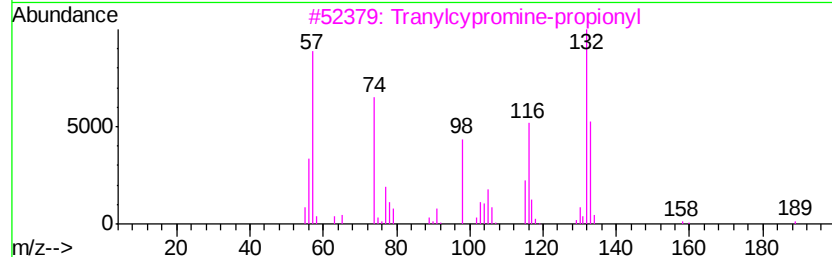
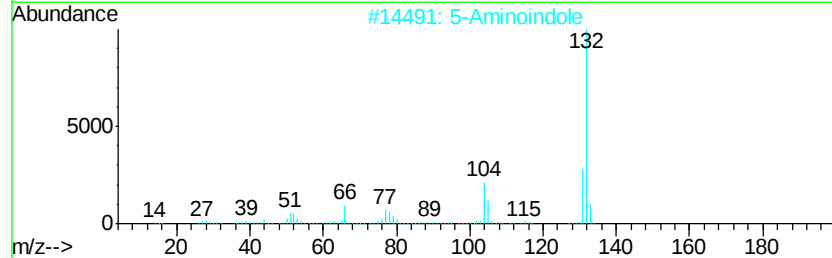
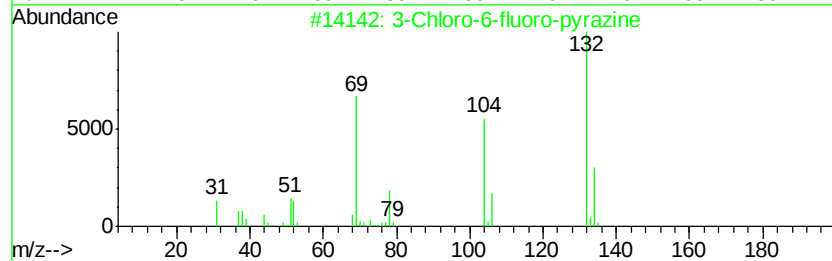
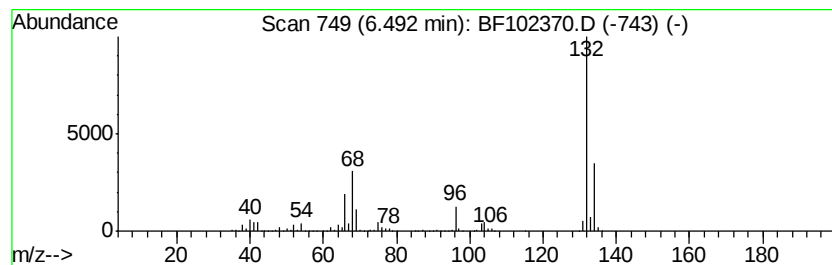
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 Peak Number 2 unknown6.49 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.49	68.46 ng	2573180	1,4-Dichlorobenzene-d4	6.72

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Aminoindole	132	C8H8N2	005192-03-0	12
3		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9
5		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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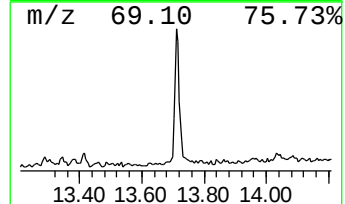
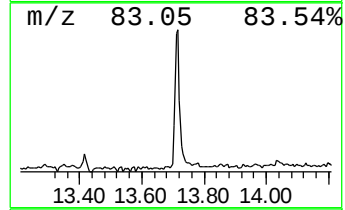
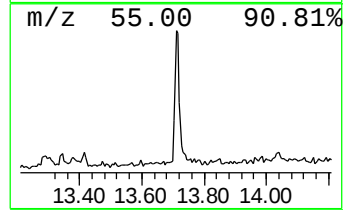
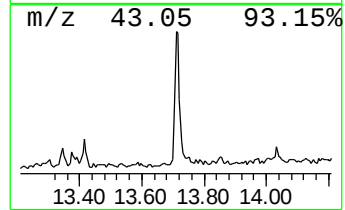
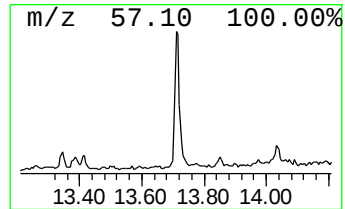
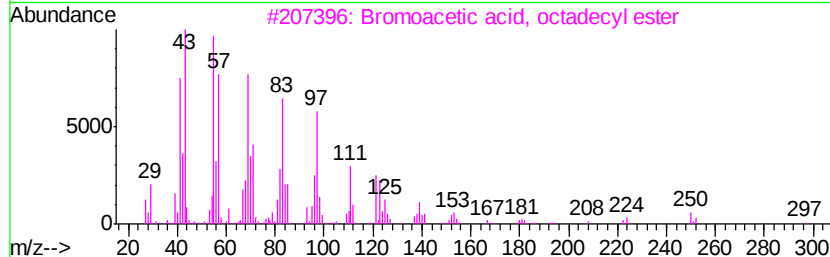
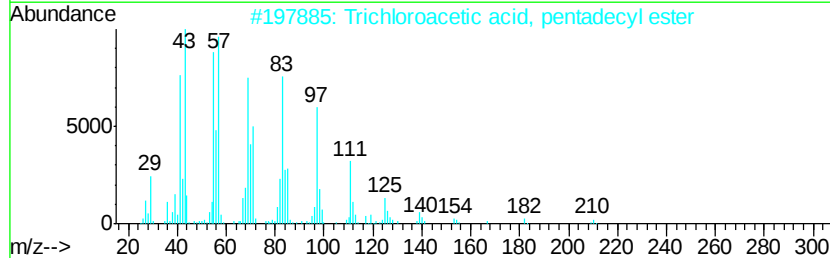
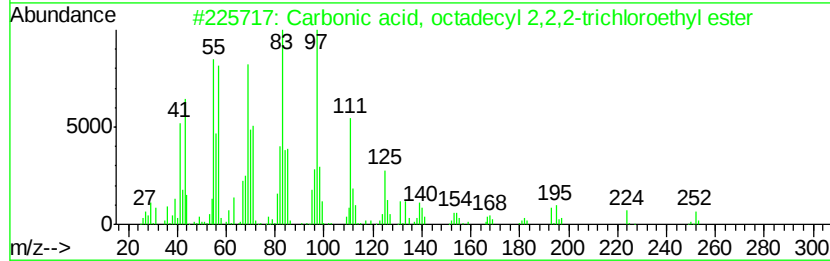
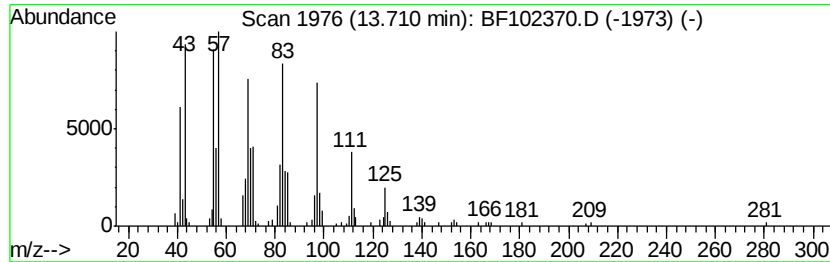
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Peak Number 4 Carbonic acid, octadecyl 2,... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.71	6.36 ng	229886	Chrysene-d12	13.87

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Carbonic acid, octadecyl 2,2,2-t...	444	C21H39Cl3O3	1000314-56-3	91
2		Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	91
3		Bromoacetic acid, octadecyl ester	390	C20H39BrO2	018992-03-5	91
4		Behenic alcohol	326	C22H46O	000661-19-8	91
5		Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	91



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown4.92	4.92	2.3	ng	85529	1	6.72	751734	20.0
unknown6.49	6.49	68.5	ng	2573180	1	6.72	751734	20.0
Carbonic acid, oc...	13.71	6.4	ng	229886	5	13.87	723156	20.0