

Data Path : Z:\HPCHEM1\BNA F\DATA\BF013017\
 Data File : BF092657.D
 Acq On : 31 Jan 2017 2:06
 Operator : SJ/MA
 Sample : I1552-01
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FK-EF-01-003-ABCD

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-BF013017.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	5.150	265	268	274	rVB	1503816	1902359	40.95%	4.881%
2	5.573	302	305	307	rBV	3627473	3638644	78.32%	9.335%
3	6.601	392	395	397	rBV	3569380	4202593	90.46%	10.782%
4	6.727	403	406	409	rBV	3543534	3564615	76.73%	9.145%
5	6.956	424	426	429	rBV	1236151	1037302	22.33%	2.661%
6	7.116	437	440	442	rBV	3270588	2979682	64.14%	7.645%
7	7.527	473	476	478	rBV	2417860	2641574	56.86%	6.777%
8	8.247	536	539	542	rBV	1382975	1333871	28.71%	3.422%
9	9.333	630	634	636	rBV	4032158	4645689	100.00%	11.919%
10	9.722	665	668	670	rBV	366916	315041	6.78%	0.808%
11	9.916	682	685	689	rVB2	77367	153375	3.30%	0.393%
12	10.008	690	693	695	rBV	1841356	1632155	35.13%	4.187%
13	10.430	727	730	732	rVB	92502	97080	2.09%	0.249%
14	10.648	746	749	753	rBV	29450	46856	1.01%	0.120%
15	10.796	759	762	765	rBV	2259866	2405841	51.79%	6.172%
16	10.910	770	772	775	rBV	94088	152981	3.29%	0.392%
17	11.356	808	811	812	rBV	88219	98132	2.11%	0.252%
18	11.493	820	823	825	rBV	1774105	1565546	33.70%	4.016%
19	12.705	927	929	931	rBV	86790	77290	1.66%	0.198%
20	12.934	946	949	951	rBV	65883	71399	1.54%	0.183%
21	13.082	959	962	964	rBV	3574302	3880729	83.53%	9.956%
22	13.962	1037	1039	1042	rBV2	88889	120404	2.59%	0.309%
23	14.134	1051	1054	1056	rBV	860142	1133662	24.40%	2.908%
24	14.979	1125	1128	1132	rVB	43359	56502	1.22%	0.145%
25	15.174	1143	1145	1149	rVB	26155	53551	1.15%	0.137%
26	15.597	1179	1182	1190	rVB	662095	991245	21.34%	2.543%
27	16.065	1220	1223	1225	rBV	32791	55601	1.20%	0.143%
28	16.568	1264	1267	1274	rVB2	31734	73141	1.57%	0.188%
29	17.163	1316	1319	1323	rVB3	23410	51233	1.10%	0.131%

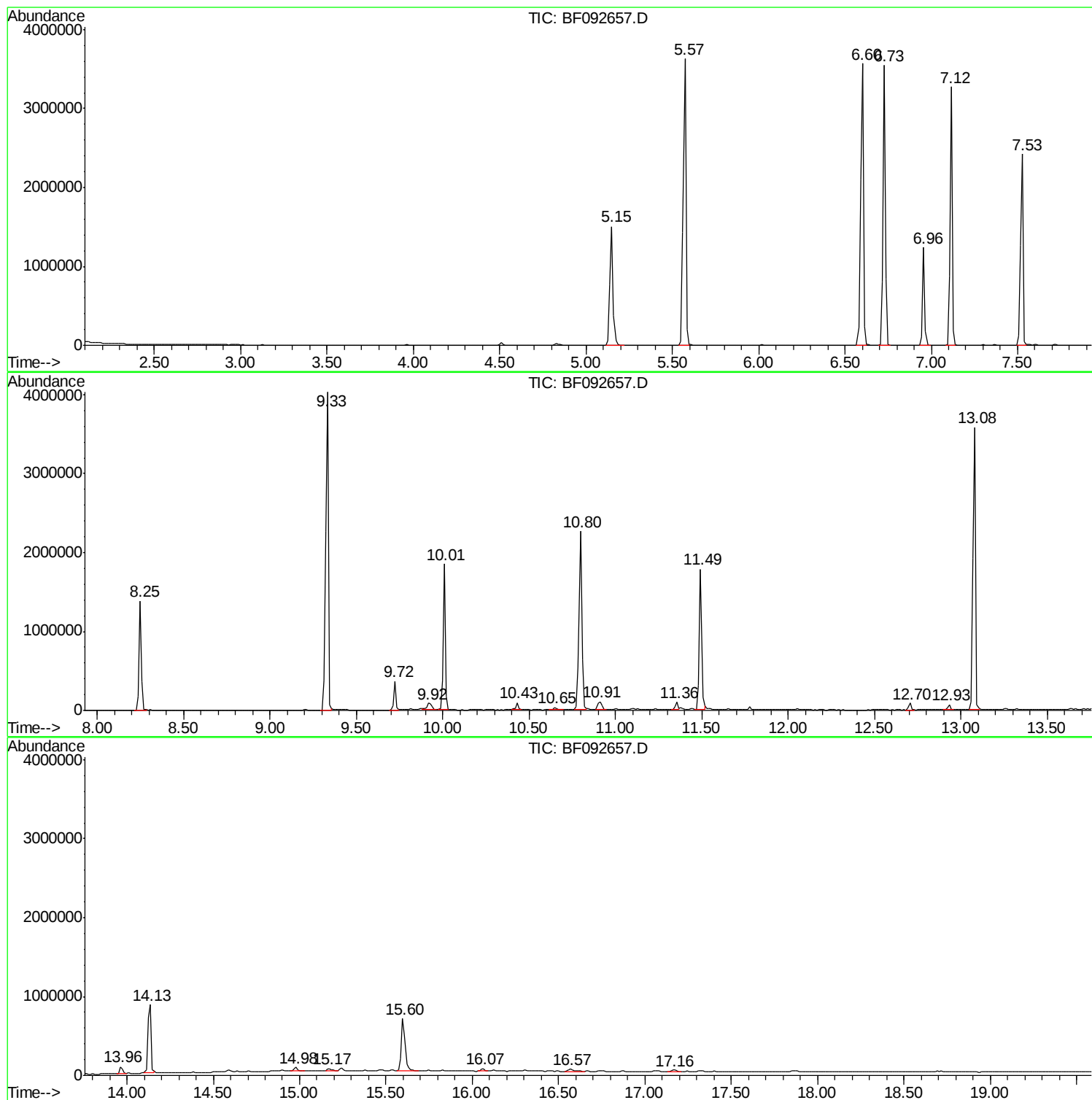
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FK-EF-01-003-ABCD

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF013017.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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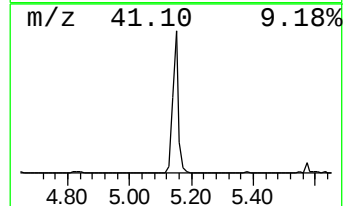
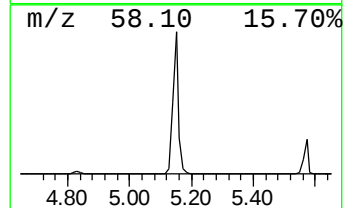
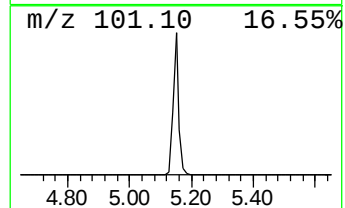
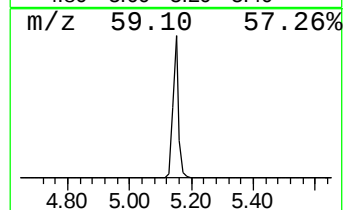
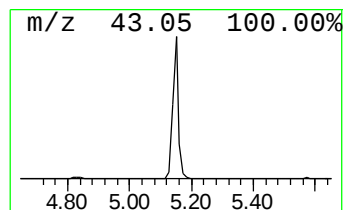
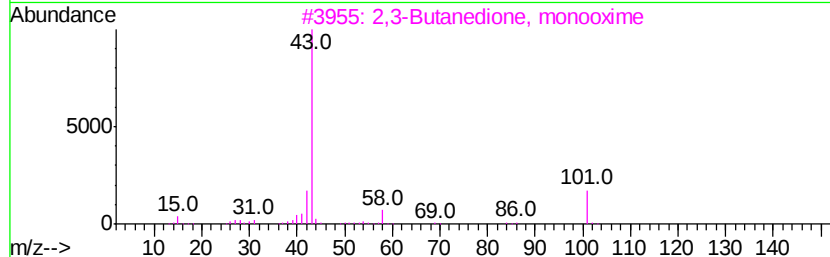
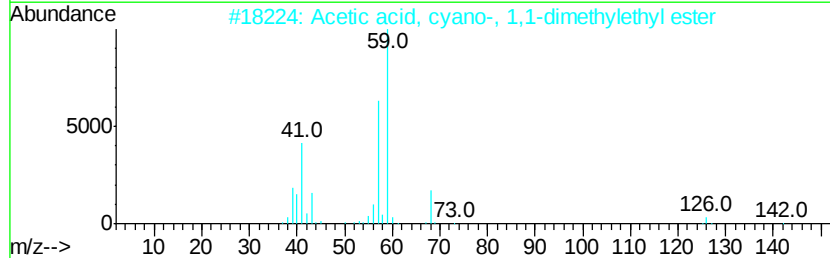
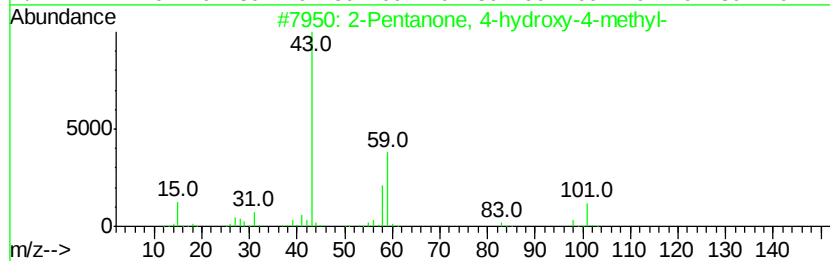
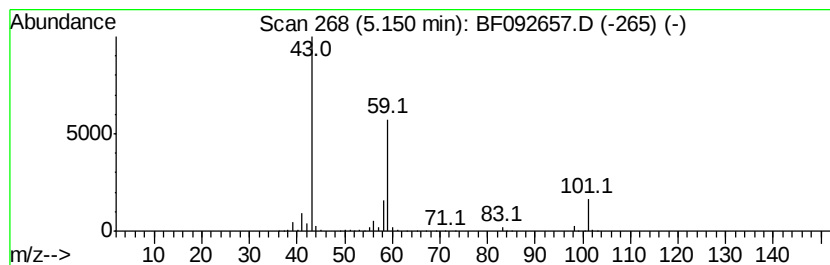
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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.
5.15	36.68 ng	1902360	1,4-Dichlorobenzene-d4	6.96

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	25
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
4			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	9
5			3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	9



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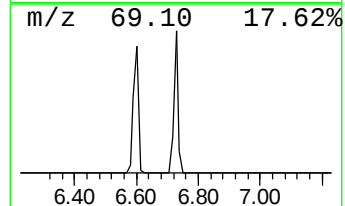
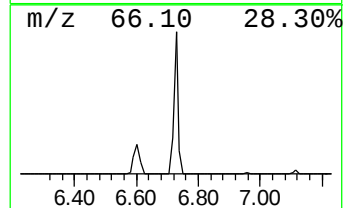
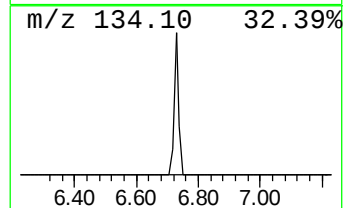
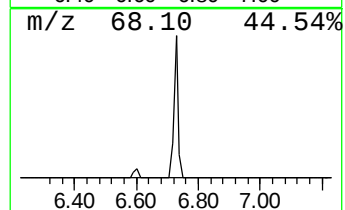
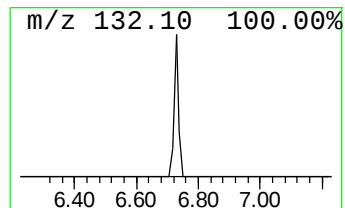
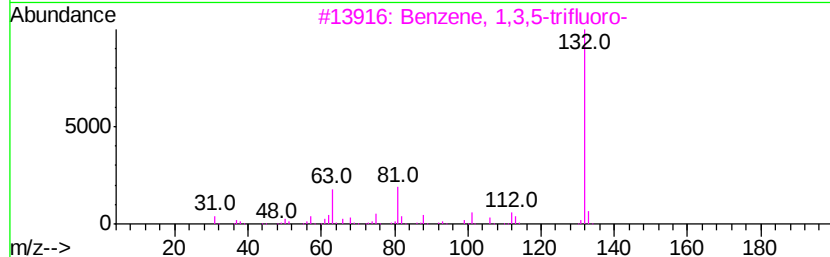
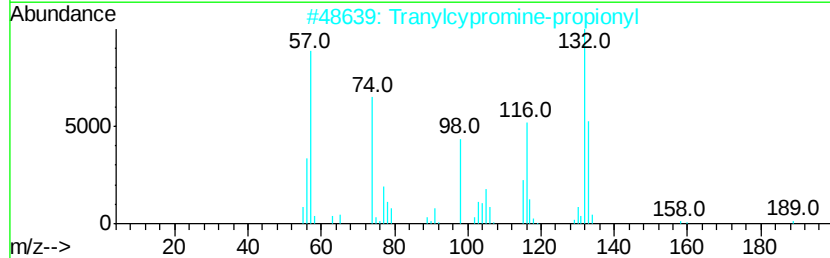
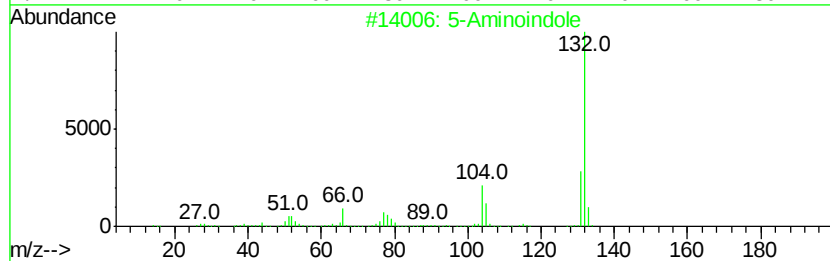
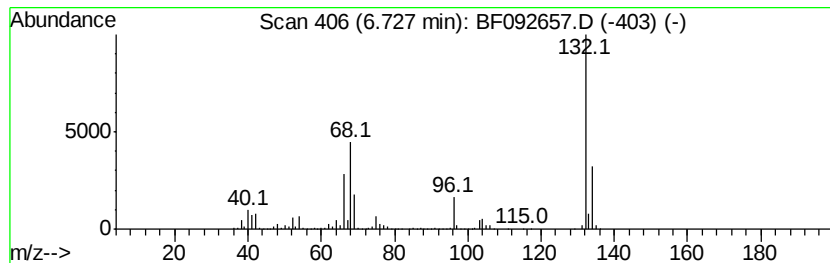
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Peak Number 2 unknown6.73 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.73	68.73 ng	3564620	1,4-Dichlorobenzene-d4	6.96

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Aminoindole	132	C8H8N2	005192-03-0	12
2		Tranvlcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9
4		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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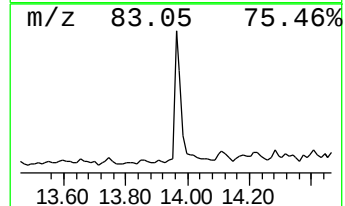
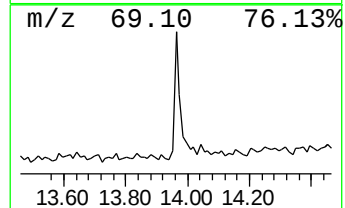
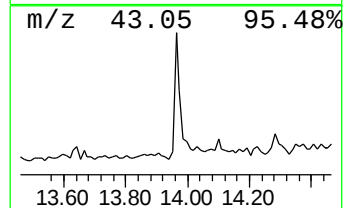
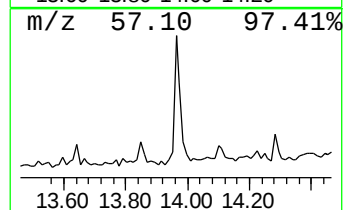
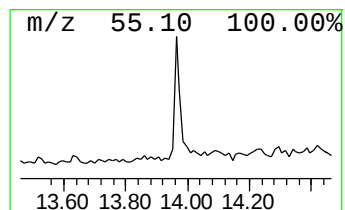
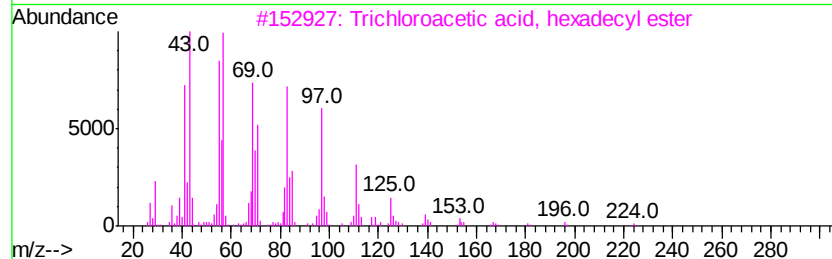
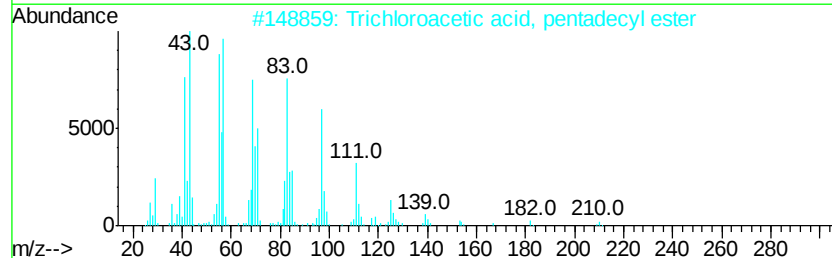
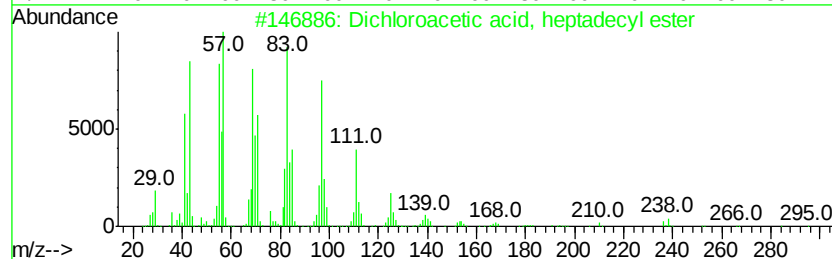
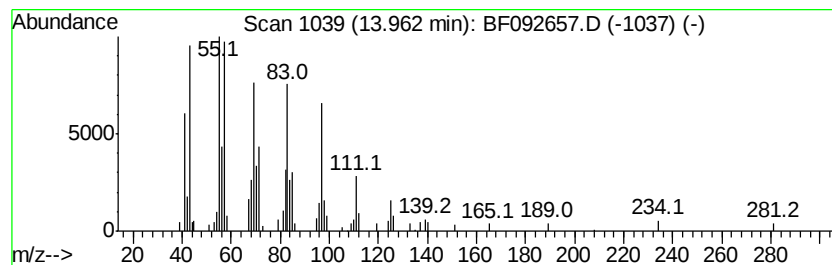
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Peak Number 4 Dichloroacetic acid, heptad... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.96	2.12 ng	120404	Chrysene-d12	14.13

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	95
2		Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	95
3		Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	95
4		1-Docosene	308	C22H44	001599-67-3	92
5		1-Nonadecanol	284	C19H40O	001454-84-8	91



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.15	36.7	ng	1902360	1	6.96	1037300	20.0
unknown6.73	6.73	68.7	ng	3564620	1	6.96	1037300	20.0
Dichloroacetic ac...	13.96	2.1	ng	120404	5	14.13	1133660	20.0