

Data Path : Z:\VOASRV\HPCHEM1\MSVOA F\DATA\VF052019\
 Data File : VF062663.D
 Acq On : 20 May 2019 16:25
 Operator : FY/SY
 Sample : K2894-10
 Misc : 5.00µ/5mL/MSVOA F/SOIL
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_F
 ClientSampleId :
 SB-07-2.5-3.5

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:\VOASRV\HPCHEM1\MSVOA_F\METHODS\82F051619S.M
 Title : SW846 8260

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.125	57	60	63	rBV2	24806	54051	3.20%	0.932%
2	1.460	92	94	99	rVV5	8230	21480	1.27%	0.370%
3	1.707	115	119	125	rBV9	12971	46412	2.75%	0.800%
4	1.865	132	135	137	rBV2	13226	28346	1.68%	0.489%
5	2.101	156	159	161	rBV4	9599	18130	1.07%	0.313%
6	2.298	176	179	181	rBV	51291	101344	6.00%	1.747%
7	2.348	181	184	188	rVB5	42751	110384	6.53%	1.903%
8	2.505	197	200	204	rBV2	29336	62227	3.68%	1.073%
9	3.077	256	258	262	rVB	31003	54064	3.20%	0.932%
10	4.319	378	384	395	rVV	188905	581193	34.39%	10.021%
11	5.068	451	460	470	rBV2	531293	1690045	100.00%	29.141%
12	5.788	527	533	542	rBV	473543	1235560	73.11%	21.305%
13	7.444	698	701	710	rVB10	7273	22738	1.35%	0.392%
14	7.720	724	729	735	rBV	299142	731137	43.26%	12.607%
15	8.361	790	794	798	rBV7	8650	27979	1.66%	0.482%
16	9.918	948	952	959	rVV2	150548	353497	20.92%	6.095%
17	10.263	982	987	989	rBV4	6007	17675	1.05%	0.305%
18	11.505	1109	1113	1117	rBV	61868	103110	6.10%	1.778%
19	11.633	1123	1126	1133	rBV3	33863	72857	4.31%	1.256%
20	11.732	1133	1136	1139	rVB4	10475	18995	1.12%	0.328%
21	12.619	1223	1226	1230	rVB2	55105	87092	5.15%	1.502%
22	13.329	1294	1298	1303	rVB2	55882	98947	5.85%	1.706%
23	13.398	1303	1305	1310	rBV6	7372	17535	1.04%	0.302%
24	13.723	1334	1338	1339	rVV3	11262	21956	1.30%	0.379%
25	13.752	1339	1341	1346	rVV5	20810	43514	2.57%	0.750%
26	14.058	1367	1372	1376	rBV6	7468	18468	1.09%	0.318%
27	14.186	1382	1385	1389	rVB6	9732	19142	1.13%	0.330%
28	14.502	1411	1417	1424	rBV10	18081	66454	3.93%	1.146%
29	14.630	1427	1430	1434	rBV5	13307	29210	1.73%	0.504%
30	14.906	1455	1458	1464	rVB6	8196	17844	1.06%	0.308%
31	15.201	1485	1488	1491	rVB5	13254	28115	1.66%	0.485%

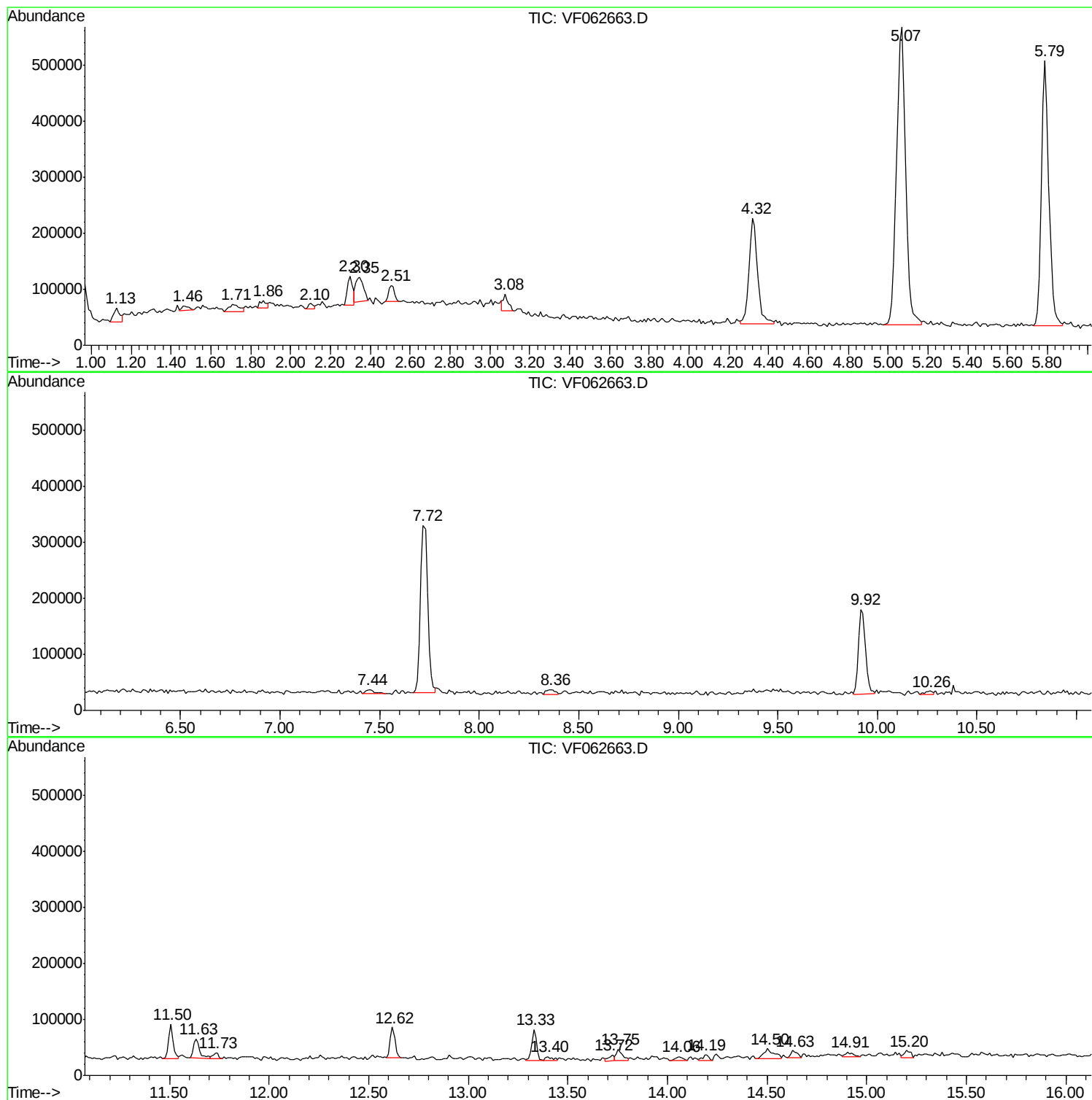
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F051619S.M
Quant Title : SW846 8260

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



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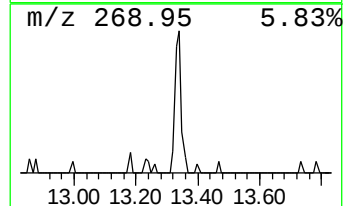
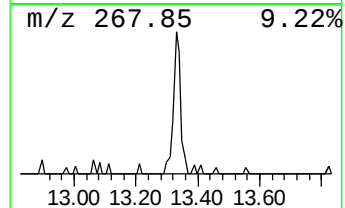
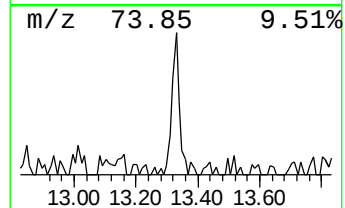
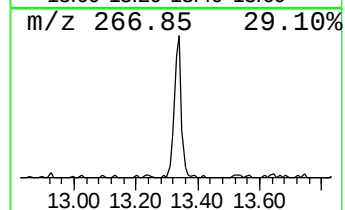
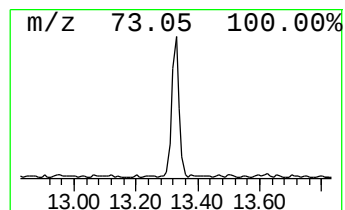
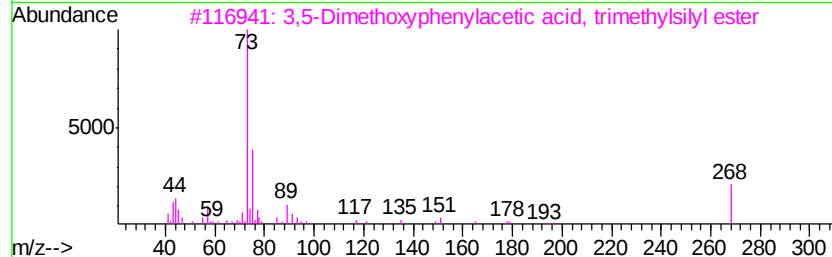
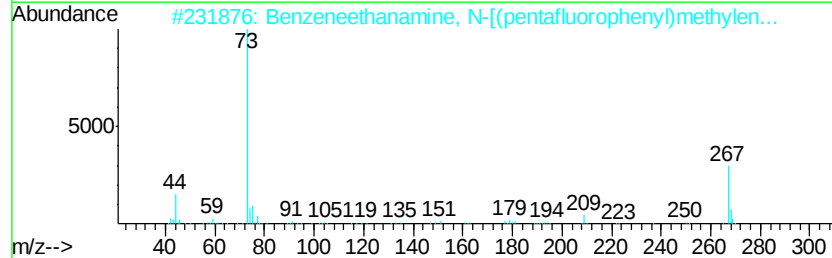
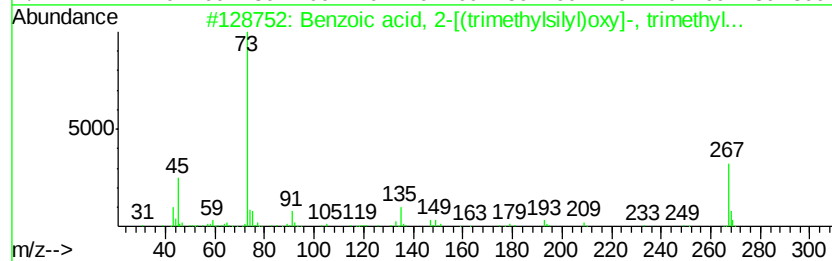
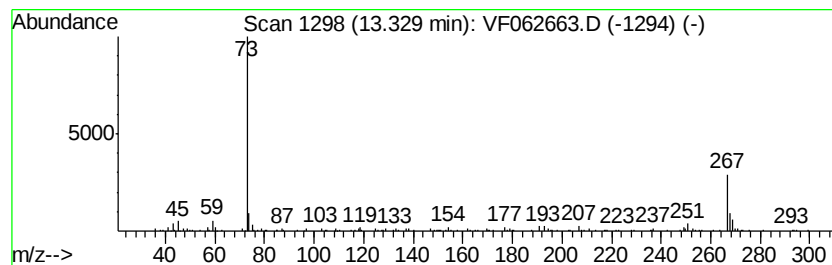
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TIC Integration Parameters: LSCINT.P

Peak Number 3 Benzoic acid, 2-[(trimethyl... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.		
13.33	56.81 ug/l	98947	1,4-Dichlorobenzene-d4	12.62		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzoic acid, 2-[(trimethylsilyl)oxy]-, trimethylsilyl ester	282	C13H22O3Si2	003789-85-3	59
2		Benzenethanamine, N-[(pentafluorophenyl)methyl]-, trimethylsilyl ester	475	C21H26F5N2Si2	055429-85-1	50
3		3,5-Dimethoxyphenylacetic acid, trimethylsilyl ester	268	C13H20O4Si	1000071-82-4	16
4		1-(3-Chlorophenyl)piperazine, 4-(trimethylsilyl)oxy-	268	C13H21ClN2Si	1000373-99-3	12
5		Benzenepropanoic acid, .alpha.-(trimethylsilyl)oxy-	265	C13H19NO3Si	055494-13-8	10



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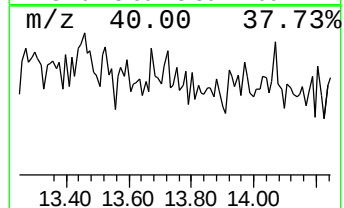
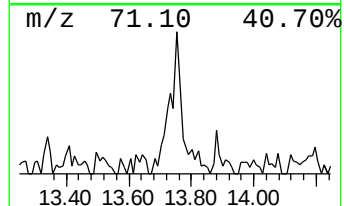
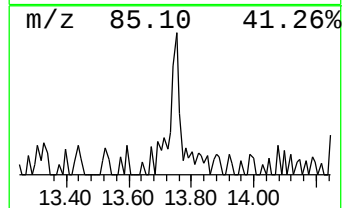
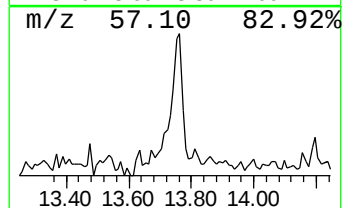
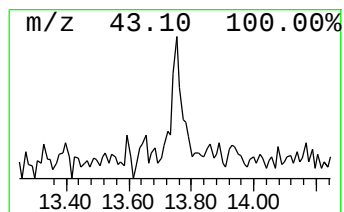
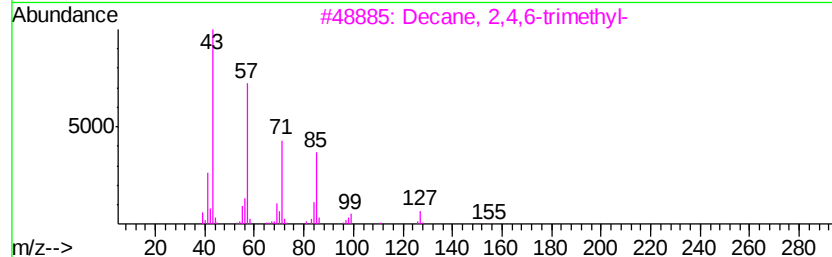
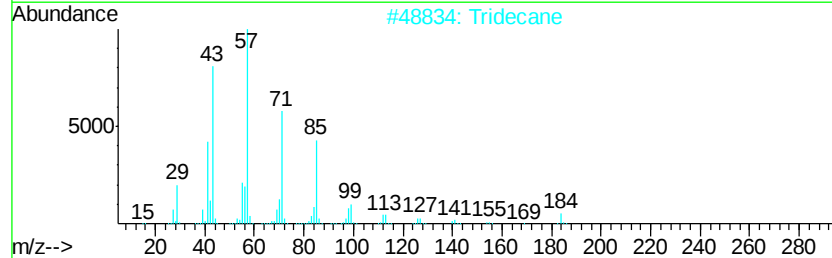
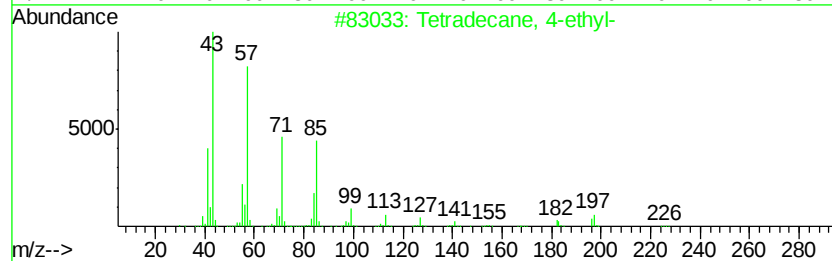
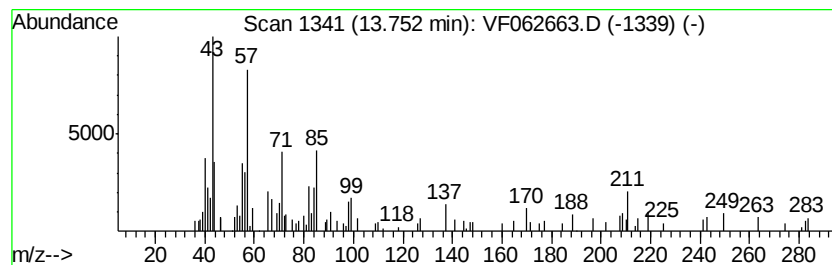
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Peak Number 5 unknown13.75 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.75	24.98 ug/l	43514	1,4-Dichlorobenzene-d4	12.62

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetradecane, 4-ethyl-	226	C16H34	055045-14-2	43
2		Tridecane	184	C13H28	000629-50-5	43
3		Decane, 2,4,6-trimethyl-	184	C13H28	062108-27-4	38
4		Undecane	156	C11H24	001120-21-4	38
5		Octacosane	394	C28H58	000630-02-4	38



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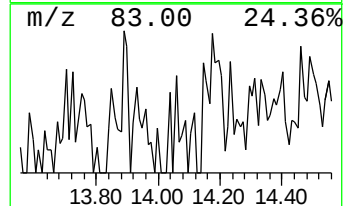
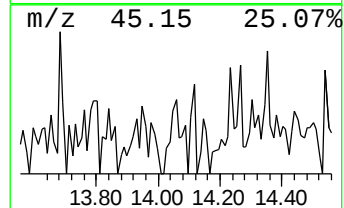
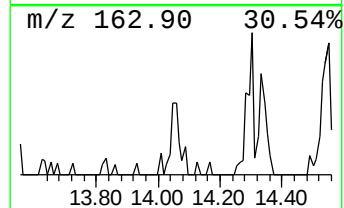
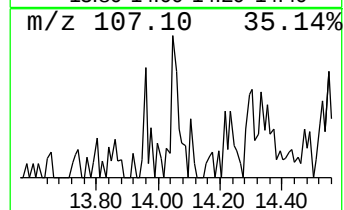
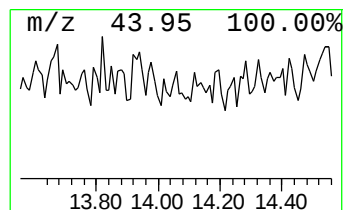
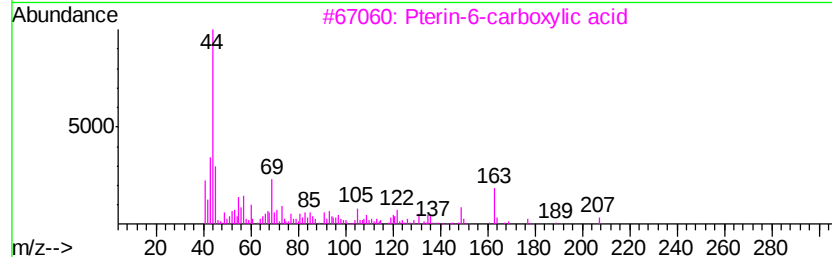
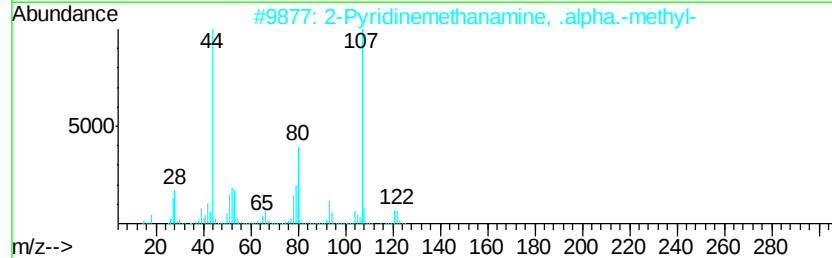
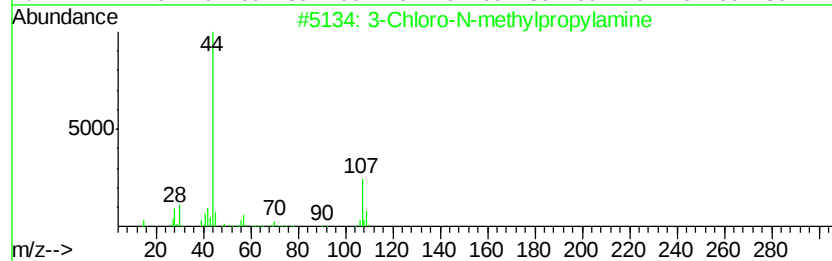
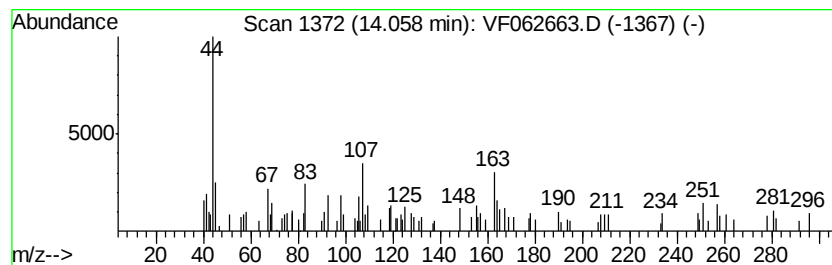
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 unknown14.06 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.06	10.60 ug/l	18468	1,4-Dichlorobenzene-d4	12.62

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-N-methylpropylamine	107	C4H10ClN	065232-62-4	43
2		2-Pyridinemethanamine, .alpha.-m...	122	C7H10N2	042088-91-5	43
3		Pterin-6-carboxylic acid	207	C7H5N5O3	000948-60-7	35
4		5,6-Dihydro-2,4,6-trimethyl-4H-1...	163	C6H13NS2	1000360-36-4	32
5		D-Ribo-Hexitol, 2-(acetylamino)-...	273	C12H19NO6	043180-03-6	30



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzoic acid, 2-[...	13.33	56.8	ug/l	98947	4	12.62	87092	50.0
unknown13.75	13.75	25.0	ug/l	43514	4	12.62	87092	50.0
unknown14.06	14.06	10.6	ug/l	18468	4	12.62	87092	50.0