LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA F\DATA\VF080218\
Data File : VF059835.D
Acq On : 3 Aug 2018 12:52
Operator : VA/AP
Sample : J4285-09
Misc : 6.01g/5mL/MSVOA F/SOIL
ALS Vial : 100 Sample Multiplier: 1

Instrument : MSVOA_F
ClientSampleId : KG-PIT-3

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:\VOASRV\HPCHEM1\MSVOA_F\METHODS\82F073018S.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	1.042	50	52	54	rVB2	2201	3208	2.05%	0.480%
2	1.101	54	58	60	rBV2	7181	12890	8.23%	1.930%
3	1.190	60	67	68	rBV5	3962	12480	7.97%	1.869%
4	1.634	110	112	115	rVB2	2088	3665	2.34%	0.549%
5	1.673	115	116	117	rBV	1899	1914	1.22%	0.287%
6	1.752	120	124	125	rVB2	1253	2473	1.58%	0.370%
7	1.772	125	126	128	rBV2	1219	1773	1.13%	0.265%
8	1.831	130	132	137	rVV4	1859	3998	2.55%	0.599%
9	1.900	137	139	140	rVV	1535	1896	1.21%	0.284%
10	1.929	140	142	146	rVV3	1762	3156	2.02%	0.473%
11	2.156	162	165	166	rVB2	1908	2564	1.64%	0.384%
12	2.186	166	168	171	rBV	2033	3855	2.46%	0.577%
13	2.245	173	174	175	rBV	1689	1730	1.11%	0.259%
14	2.274	175	177	179	rVV2	1831	2485	1.59%	0.372%
15	2.314	179	181	185	rVB3	1464	3026	1.93%	0.453%
16	2.481	196	198	201	rVB2	2004	3913	2.50%	0.586%
17	2.550	203	205	206	rVV	1297	2026	1.29%	0.303%
18	2.580	206	208	211	rVV	2127	3428	2.19%	0.513%
19	2.659	213	216	218	rVV	1272	2973	1.90%	0.445%
20	2.748	221	225	228	rVV2	1124	3827	2.44%	0.573%
21	2.797	228	230	235	rVV2	1857	3937	2.52%	0.589%
22	2.994	246	250	251	rVV3	1493	2211	1.41%	0.331%
23	3.033	251	254	255	rVV	1083	1966	1.26%	0.294%
24	3.083	258	259	264	rVV2	1299	2136	1.36%	0.320%
25	3.181	266	269	273	rVB2	1081	2133	1.36%	0.319%
26	3.300	279	281	284	rVV	1388	2502	1.60%	0.375%
27	3.398	288	291	294	rVB	1723	3608	2.30%	0.540%
28	3.497	298	301	302	rVB2	1441	1654	1.06%	0.248%
29	3.615	310	313	316	rVB2	1475	3187	2.04%	0.477%
30	3.674	316	319	320	rBV	1215	1750	1.12%	0.262%
31	3.704	320	322	324	rBV2	1441	1691	1.08%	0.253%
32	3.822	332	334	338	rVV	1133	2298	1.47%	0.344%
33	4.078	357	360	362	rBV2	1622	2691	1.72%	0.403%
34	4.128	364	365	368	rBV2	1579	2464	1.57%	0.369%

LSC Area Percent Report

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ALS Vial : 100 Sample Multiplier: 1

Instrument : MSVOA_F
ClientSampleId : KG-PIT-3

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

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Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_F\METHODS\82F073018S.M
Title : SW846 8260

35	4.285	374	381	385	rBV3	6305	17455	11.15%	2.614%
36	4.344	385	387	389	rVB2	1652	1935	1.24%	0.290%
37	4.482	397	401	403	rBV2	1658	4230	2.70%	0.633%
38	4.768	428	430	433	rBV2	1033	2236	1.43%	0.335%
39	5.034	449	457	466	rBV3	48912	156531	100.00%	23.438%
40	5.212	473	475	479	rBV3	1238	2661	1.70%	0.398%
41	5.370	488	491	492	rBV	1480	1936	1.24%	0.290%
42	5.439	494	498	499	rVV3	1223	2166	1.38%	0.324%
43	5.596	509	514	519	rVB2	1425	2941	1.88%	0.440%
44	5.754	525	530	537	rBV	28958	78150	49.93%	11.702%
45	5.862	539	541	544	rBV	974	1652	1.06%	0.247%
46	5.981	552	553	555	rBV	1670	1741	1.11%	0.261%
47	6.060	558	561	564	rBV2	671	1673	1.07%	0.251%
48	6.405	591	596	597	rBV2	784	1751	1.12%	0.262%
49	6.513	603	607	608	rVB2	1712	2209	1.41%	0.331%
50	6.789	632	635	636	rBV2	1693	2124	1.36%	0.318%
51	6.976	653	654	658	rVB	1194	1907	1.22%	0.286%
52	7.144	668	671	673	rBV2	1302	2571	1.64%	0.385%
53	7.706	723	728	734	rBV	25268	63138	40.34%	9.454%
54	8.051	761	763	766	rBV3	1149	1711	1.09%	0.256%
55	8.268	782	785	786	rBV	1197	2093	1.34%	0.313%
56	8.337	788	792	795	rVV3	1505	4619	2.95%	0.692%
57	8.425	798	801	803	rVV	9368	9867	6.30%	1.477%
58	8.859	842	845	848	rVV2	1002	2669	1.71%	0.400%
59	9.027	860	862	866	rBV4	1200	2525	1.61%	0.378%
60	9.293	887	889	893	rVB4	1022	1843	1.18%	0.276%
61	9.460	904	906	908	rVB	1550	1676	1.07%	0.251%
62	9.510	908	911	914	rVB2	1728	2941	1.88%	0.440%
63	9.638	921	924	927	rBV2	1677	3388	2.16%	0.507%
64	9.914	947	952	959	rVB2	17263	40965	26.17%	6.134%
65	10.022	962	963	966	rVV2	1227	1698	1.08%	0.254%
66	10.111	970	972	974	rVB	1355	1758	1.12%	0.263%
67	10.564	1016	1018	1021	rVB2	1285	2150	1.37%	0.322%
68	10.614	1021	1023	1027	rBV2	1026	2082	1.33%	0.312%
69	10.722	1030	1034	1036	rVB2	1377	2602	1.66%	0.390%
70	10.752	1036	1037	1039	rBV	1361	1910	1.22%	0.286%
71	11.156	1074	1078	1079	rBV2	1078	2280	1.46%	0.341%

LSC Area Percent Report

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Misc : 6.01g/5mL/MSVOA F/SOIL
ALS Vial : 100 Sample Multiplier: 1

Instrument : MSVOA_F
ClientSampleId : KG-PIT-3

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

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Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_F\METHODS\82F073018S.M
Title : SW846 8260

72	11.323	1092	1095	1098	rVB3	1390	2452	1.57%	0.367%
73	11.392	1100	1102	1104	rBV2	1662	2977	1.90%	0.446%
74	11.501	1110	1113	1115	rBV2	10582	15243	9.74%	2.282%
75	11.737	1134	1137	1139	rVB3	2964	4204	2.69%	0.629%
76	11.856	1147	1149	1152	rVB2	1239	2242	1.43%	0.336%
77	11.905	1152	1154	1155	rBV2	1201	1889	1.21%	0.283%
78	12.092	1172	1173	1175	rBV2	1662	1937	1.24%	0.290%
79	12.132	1175	1177	1179	rBV2	1880	3090	1.97%	0.463%
80	12.368	1199	1201	1202	rVB	2268	2274	1.45%	0.340%
81	12.408	1202	1205	1206	rBV3	1541	2963	1.89%	0.444%
82	12.516	1215	1216	1218	rBV	1302	1823	1.16%	0.273%
83	12.625	1224	1227	1231	rVB2	8489	13656	8.72%	2.045%
84	12.674	1231	1232	1235	rBV2	1391	2138	1.37%	0.320%
85	12.802	1243	1245	1246	rVB	1657	1918	1.23%	0.287%
86	12.920	1256	1257	1260	rVB2	2136	2022	1.29%	0.303%
87	13.186	1282	1284	1285	rBV2	1633	2291	1.46%	0.343%
88	13.344	1299	1300	1303	rBV2	2054	3231	2.06%	0.484%
89	13.443	1308	1310	1312	rBV2	1922	2100	1.34%	0.314%
90	13.522	1316	1318	1319	rVB2	2132	1997	1.28%	0.299%
91	13.679	1331	1334	1335	rBV3	1916	2598	1.66%	0.389%
92	14.212	1384	1388	1390	rBV4	2828	5994	3.83%	0.897%
93	14.251	1390	1392	1394	rBV3	1775	2806	1.79%	0.420%
94	14.389	1403	1406	1408	rBV3	3328	5614	3.59%	0.841%
95	14.724	1438	1440	1443	rBV4	2761	5363	3.43%	0.803%
96	14.892	1455	1457	1461	rBV4	2850	5268	3.37%	0.789%
97	15.168	1483	1485	1487	rBV2	4285	5087	3.25%	0.762%
98	15.582	1525	1527	1528	rVB	3654	3576	2.28%	0.535%
99	15.661	1533	1535	1537	rVV	3505	4119	2.63%	0.617%
100	15.769	1544	1546	1549	rBV3	3602	7694	4.92%	1.152%

Sum of corrected areas: 667858

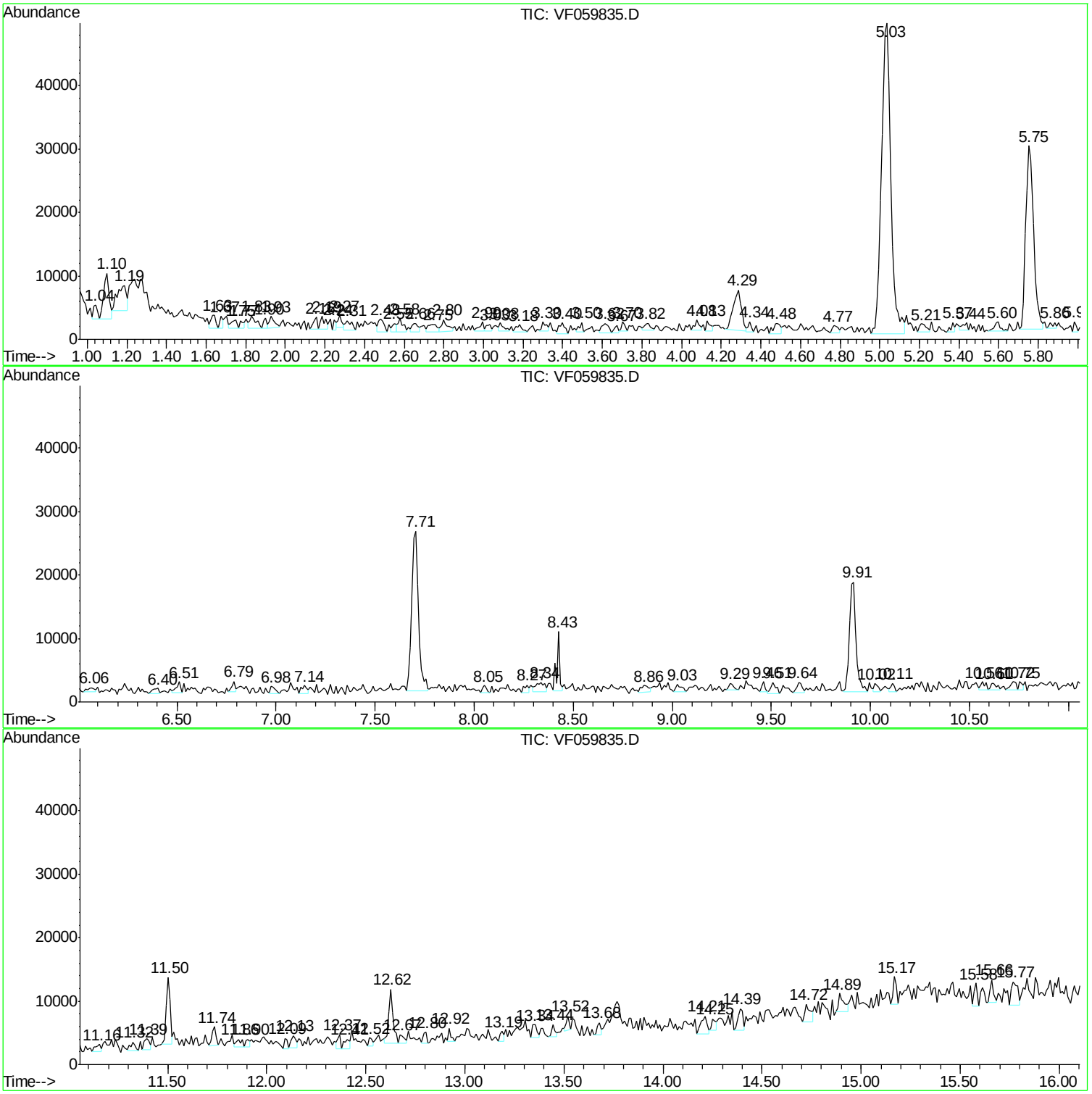
LSC Report - Integrated Chromatogram

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_F\DATA\VF080218\
Data File : VF059835.D
Acq On : 3 Aug 2018 12:52
Operator : VA/AP
Sample : J4285-09
Misc : 6.01g/5mL/MSVOA_F/SOIL
ALS Vial : 100 Sample Multiplier: 1

Instrument : MSVOA_F
ClientSampleId : KG-PIT-3

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_F\METHODS\82F073018S.M
Quant Title : SW846 8260

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



Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA F\DATA\VF080218\
Data File : VF059835.D
Acq On : 3 Aug 2018 12:52
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ALS Vial : 100 Sample Multiplier: 1

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Quant Title : SW846 8260

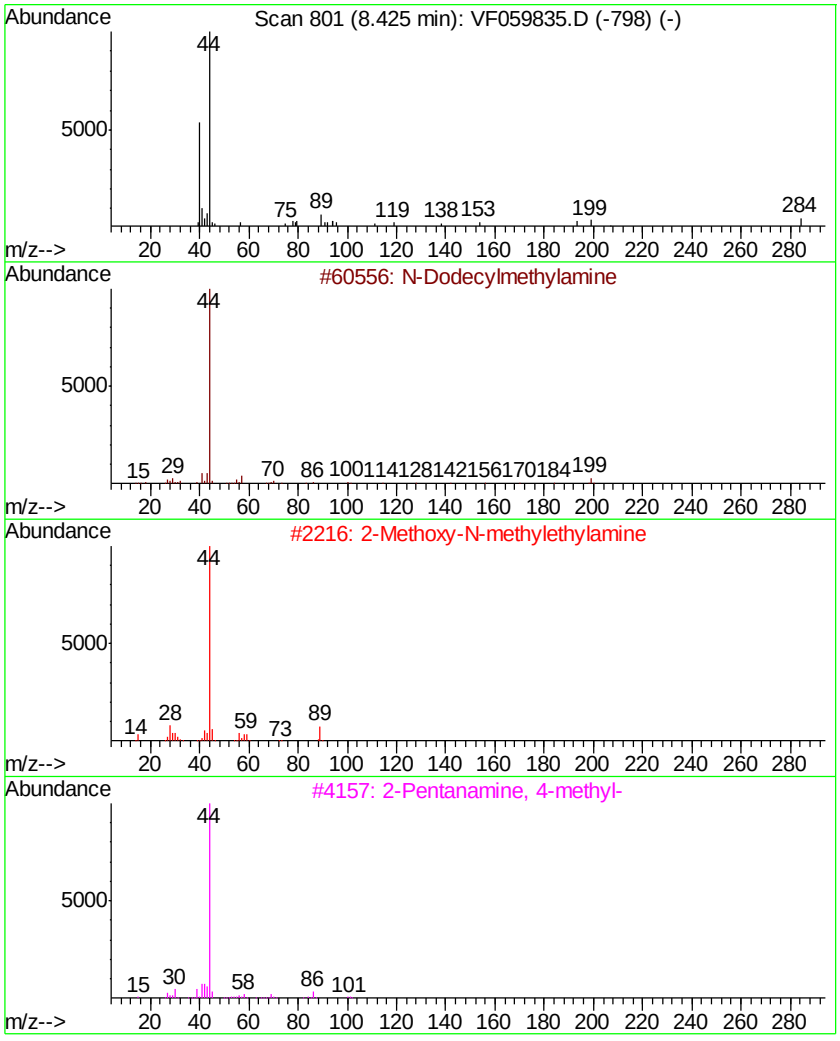
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

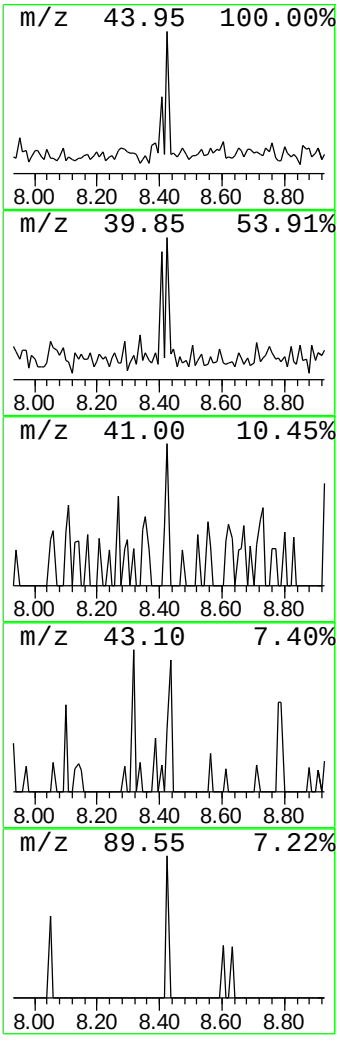
****

Peak Number 1 unknown8.43 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.43	12.04 ug/l	9867	Chlorobenzene-d5	9.91

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	N-Dodecylmethylamine	199	C13H29N	1000342-26-1	9
2	2-Methoxy-N-methylethylamine	89	C4H11NO	038256-93-8	9
3	2-Pentanamine, 4-methyl-	101	C6H15N	000108-09-8	9
4	2-Hexanamine, 5-methyl-	115	C7H17N	028292-43-5	9
5	Acetamide, 2,2,2-trifluoro-	113	C2H2F3NO	000354-38-1	9





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA F\DATA\VF080218\
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Acq On : 3 Aug 2018 12:52
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Sample : J4285-09
Misc : 6.01g/5mL/MSVOA F/SOIL
ALS Vial : 100 Sample Multiplier: 1

Instrument : MSVOA_F
ClientSampleId : KG-PIT-3

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_F\METHODS\82F073018S.M
Quant Title : SW846 8260

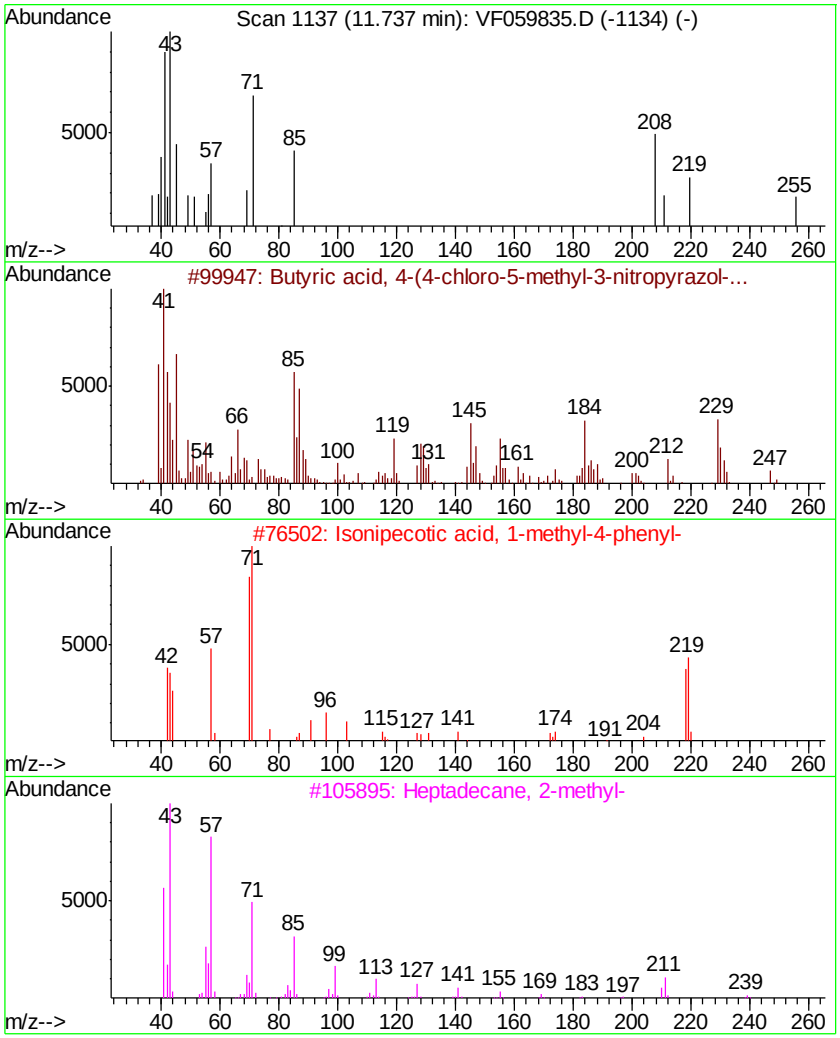
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

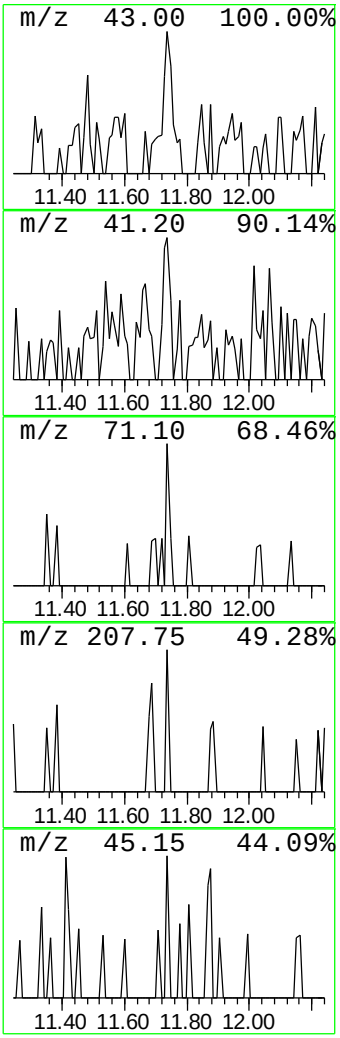
****

Peak Number 2 unknown11.74 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.74	15.39 ug/l	4204	1,4-Dichlorobenzene-d4	12.62

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butyric acid, 4-(4-chloro-5-meth...	247	C8H10ClN3O4	1000304-30-9	12
2	Isonipecotic acid, 1-methyl-4-ph...	219	C13H17NO2	003627-48-3	10
3	Heptadecane, 2-methyl-	254	C18H38	001560-89-0	10
4	2-Bromononane	206	C9H19Br	002216-35-5	10
5	Heptadecane, 9-hexyl-	324	C23H48	055124-79-3	10





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Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_F\METHODS\82F073018S.M
Quant Title : SW846 8260

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

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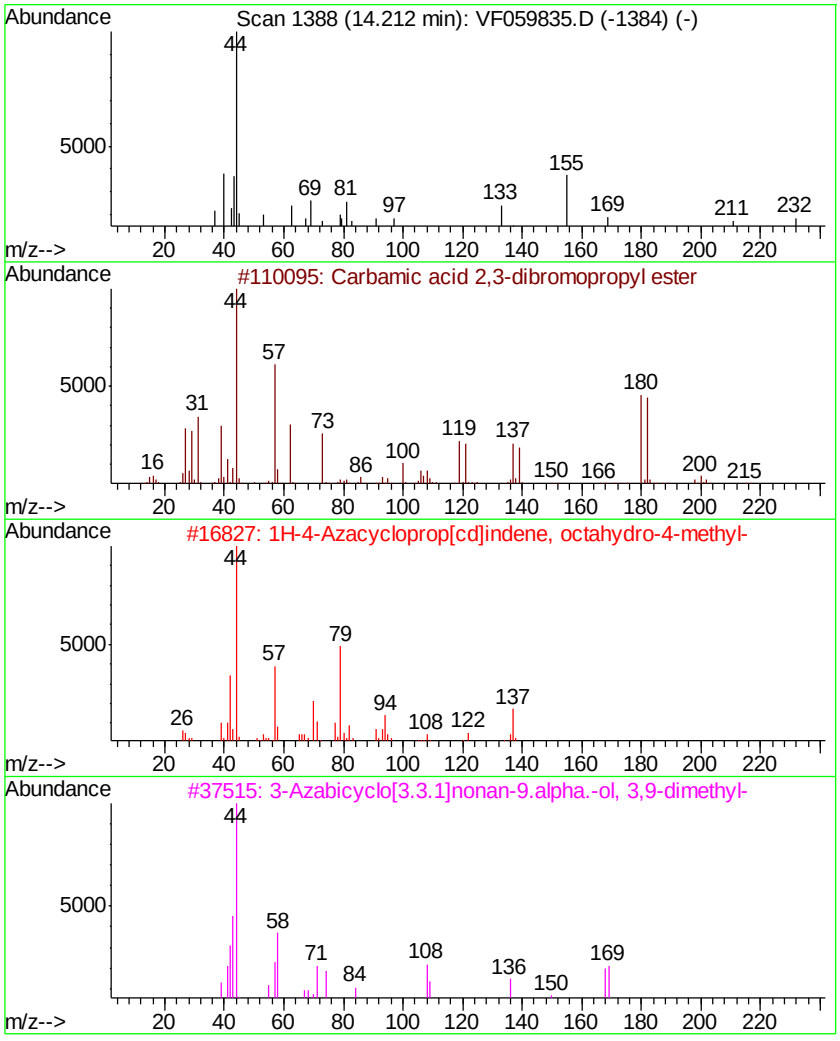
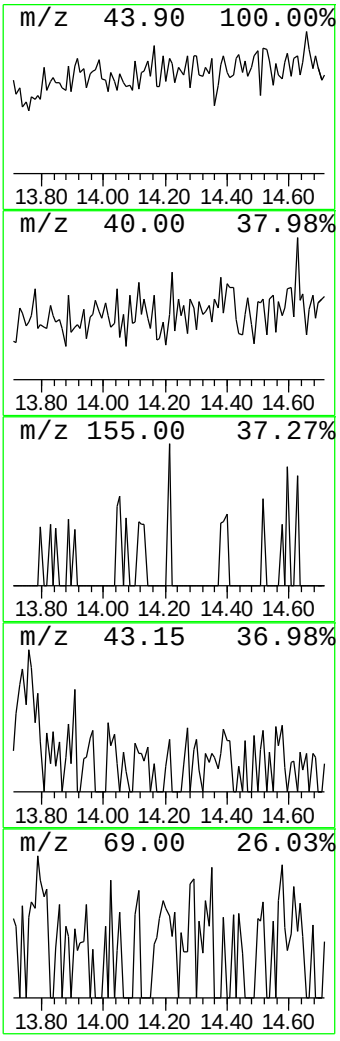
Peak Number 3 unknown14.21 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.21	21.95 ug/l	5994	1,4-Dichlorobenzene-d4	12.62

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Carbamic acid 2,3-dibromopropyl ...	259	C4H7Br2NO2	055190-46-0	17
2	1H-4-Azacycloprop[cd]indene, oct...	137	C9H15N	016967-50-3	12
3	3-Azabicyclo[3.3.1]nonan-9.alpha...	169	C10H19NO	014948-72-2	9
4	Pteridine-8-oxide, 6-aldoximino-...	222	C7H6N6O3	023663-23-2	9
5	1,1-Dichloro-2-methyl-3-(4,4-dif...	232	C10H10Cl2O2	132607-16-0	9





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Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_F\METHODS\82F073018S.M
Quant Title : SW846 8260

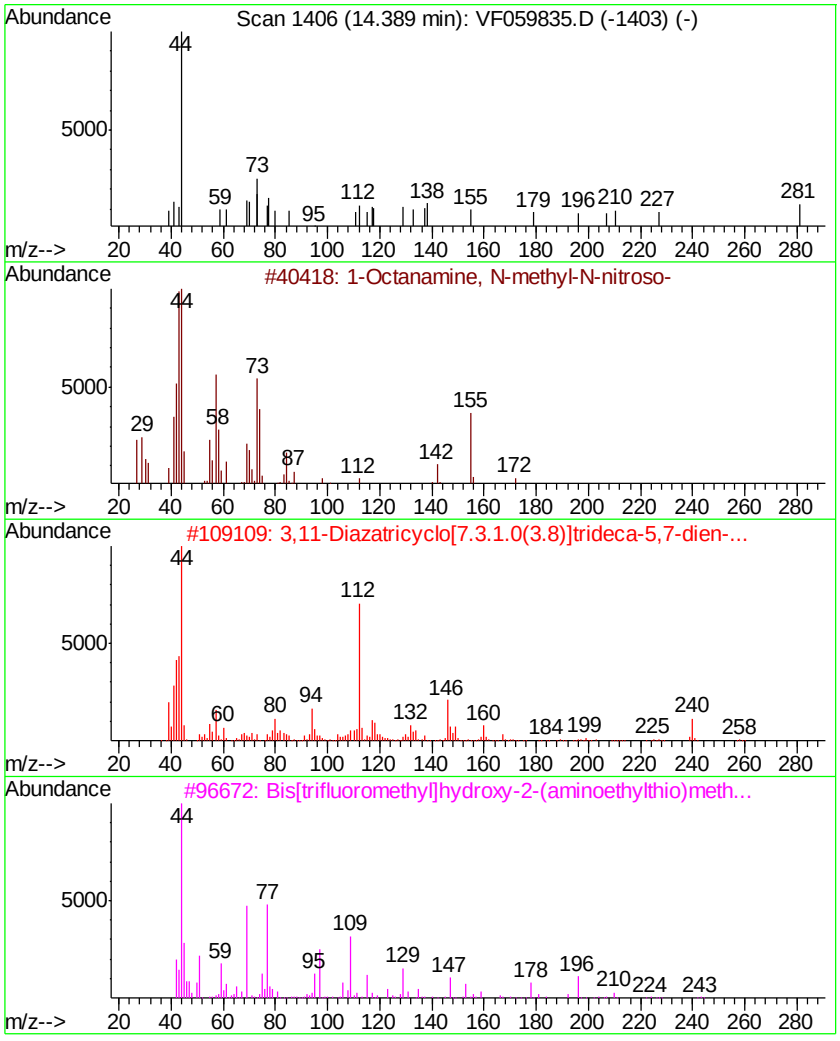
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

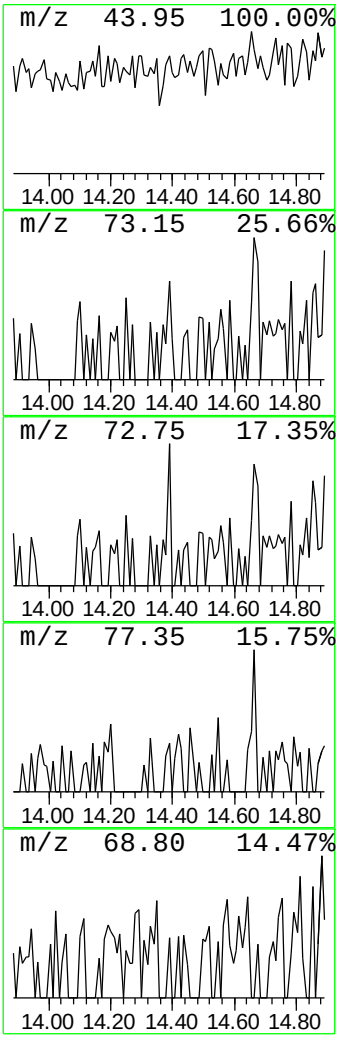
****

Peak Number 4 unknown14.39 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.39	20.56 ug/l	5614	1,4-Dichlorobenzene-d4	12.62

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Octanamine, N-methyl-N-nitroso-	172	C9H20N2O	034423-54-6	9
2	3,11-Diazatricyclo[7.3.1.0(3.8)]...	258	C15H18N2O2	1000276-73-9	9
3	Bis[trifluoromethyl]hydroxy-2-(a...	243	C5H7F6NOS	021918-92-3	9
4	8-Azabicyclo[4.3.1]decan-10-one,...	167	C10H17NO	004146-36-5	9
5	2-(2-Benzylphenoxy)ethylamine	227	C15H17NO	1000303-73-3	7





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Data File : VF059835.D
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Operator : VA/AP
Sample : J4285-09
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ALS Vial : 100 Sample Multiplier: 1

Instrument : MSVOA_F
ClientSampleId : KG-PIT-3

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_F\METHODS\82F073018S.M
Quant Title : SW846 8260

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

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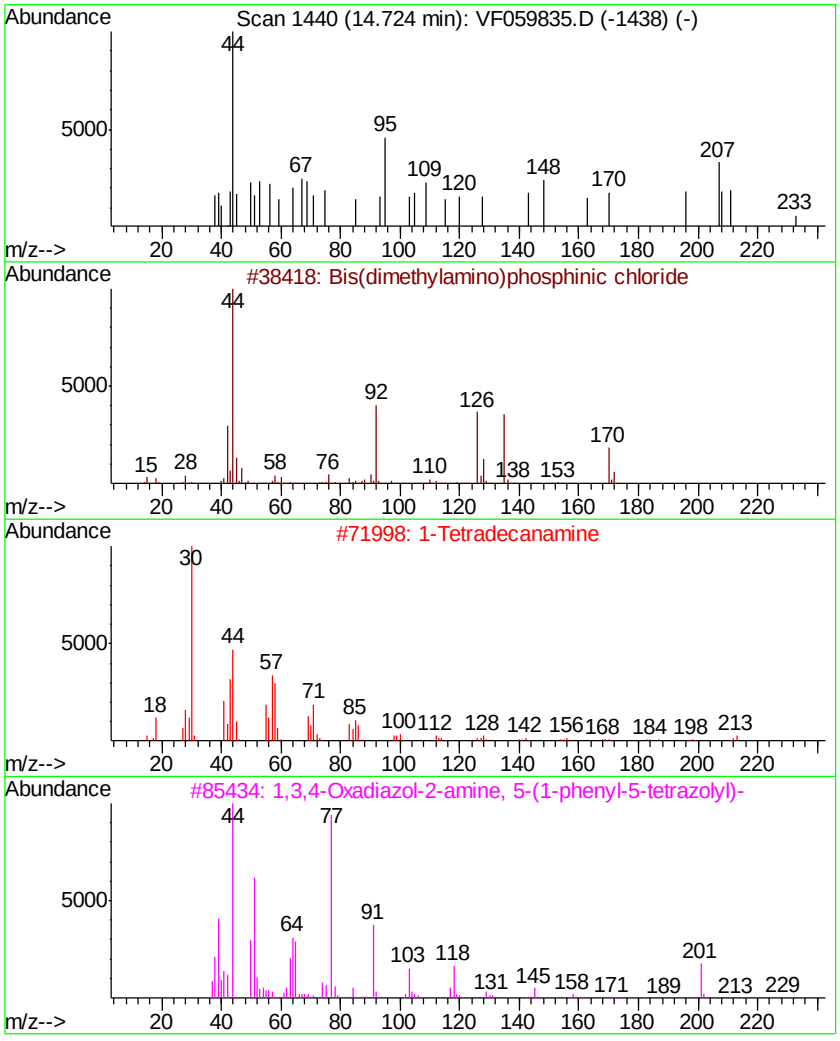
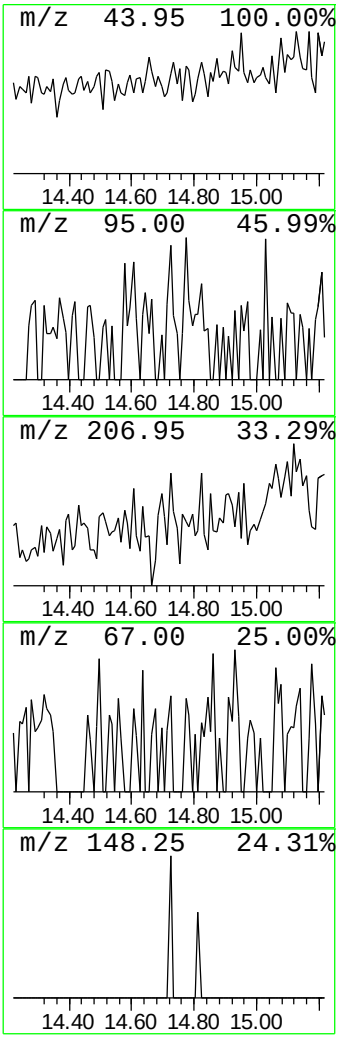
Peak Number 5 unknown14.72 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.72	19.64 ug/l	5363	1,4-Dichlorobenzene-d4	12.62

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bis(dimethylamino)phosphinic chl...	170	C4H12ClN2OP	001605-65-8	12
2	1-Tetradecanamine	213	C14H31N	002016-42-4	10
3	1,3,4-Oxadiazol-2-amine, 5-(1-ph...	229	C9H7N7O	339244-78-9	10
4	Benzaldehyde, 2,4,5-trimethoxy-	196	C10H12O4	004460-86-0	9
5	Bis[trifluoromethyl]hydroxy-2-(a...	243	C5H7F6NOS	021918-92-3	9





Data Path : Z:\VOASRV\HPCHEM1\MSVOA F\DATA\VF080218\
Data File : VF059835.D
Acq On : 3 Aug 2018 12:52
Operator : VA/AP
Sample : J4285-09
Misc : 6.01g/5mL/MSVOA F/SOIL
ALS Vial : 100 Sample Multiplier: 1

Instrument : MSVOA_F
ClientSampleId : KG-PIT-3

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_F\METHODS\82F073018S.M
Quant Title : SW846 8260

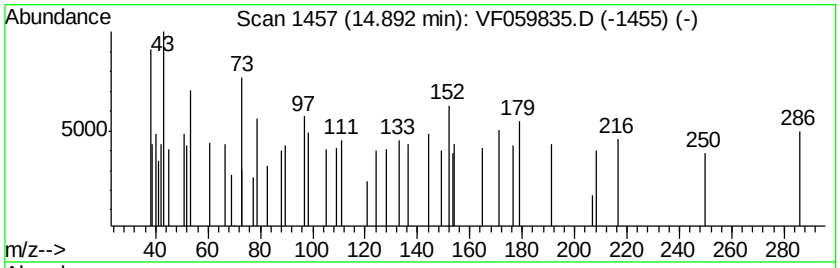
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

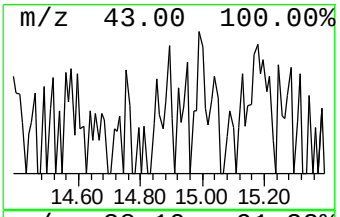
****

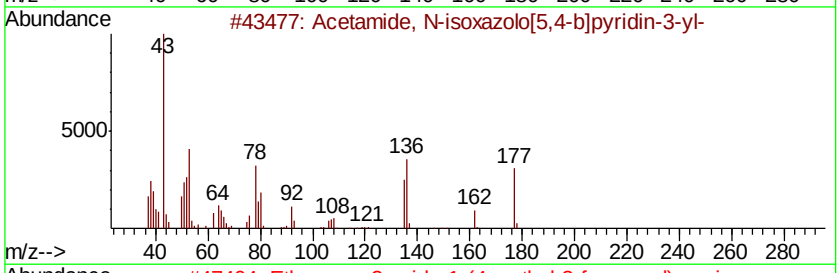
Peak Number 6 unknown14.89 Concentration Rank 5

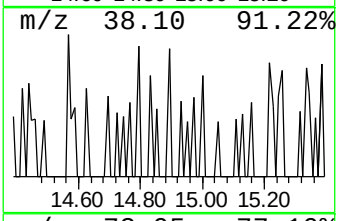
R.T.	EstConc	Area	Relative to ISTD	R.T.
14.89	19.29 ug/l	5268	1,4-Dichlorobenzene-d4	12.62

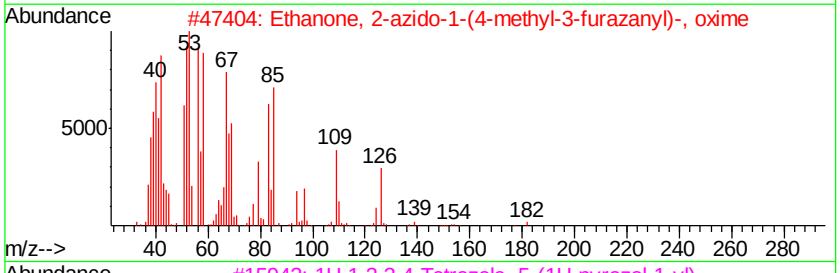
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetamide, N-isoxazolo[5,4-b]pyr...	177	C8H7N3O2	1000362-63-0	22
2	Ethanone, 2-azido-1-(4-methyl-3-...	182	C5H6N6O2	1000270-71-4	22
3	1H-1,2,3,4-Tetrazole, 5-(1H-pyra...	136	C4H4N6	1000351-28-9	12
4	3-Formylhept-3-enedinitrile	148	C8H8N2O	1000192-13-5	10
5	2-Cyclobutene-1-carboxamide	97	C5H7NO	053778-54-4	10

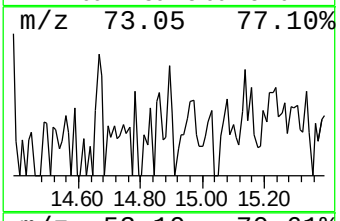


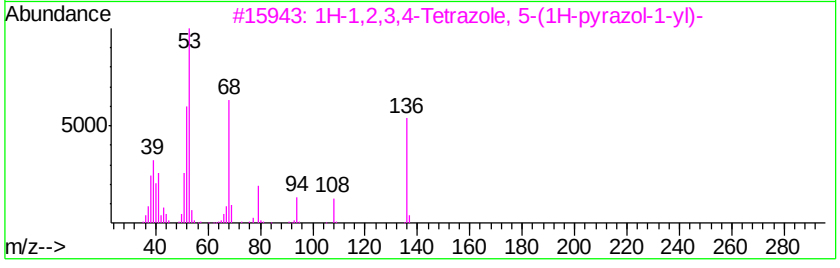


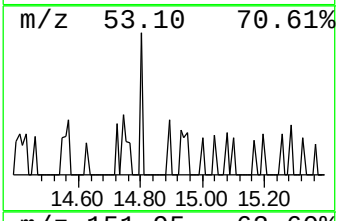


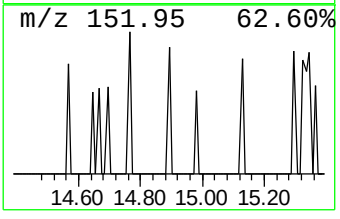












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Data File : VF059835.D
Acq On : 3 Aug 2018 12:52
Operator : VA/AP
Sample : J4285-09
Misc : 6.01g/5mL/MSVOA F/SOIL
ALS Vial : 100 Sample Multiplier: 1

Instrument : MSVOA_F
ClientSampleId : KG-PIT-3

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_F\METHODS\82F073018S.M
Quant Title : SW846 8260

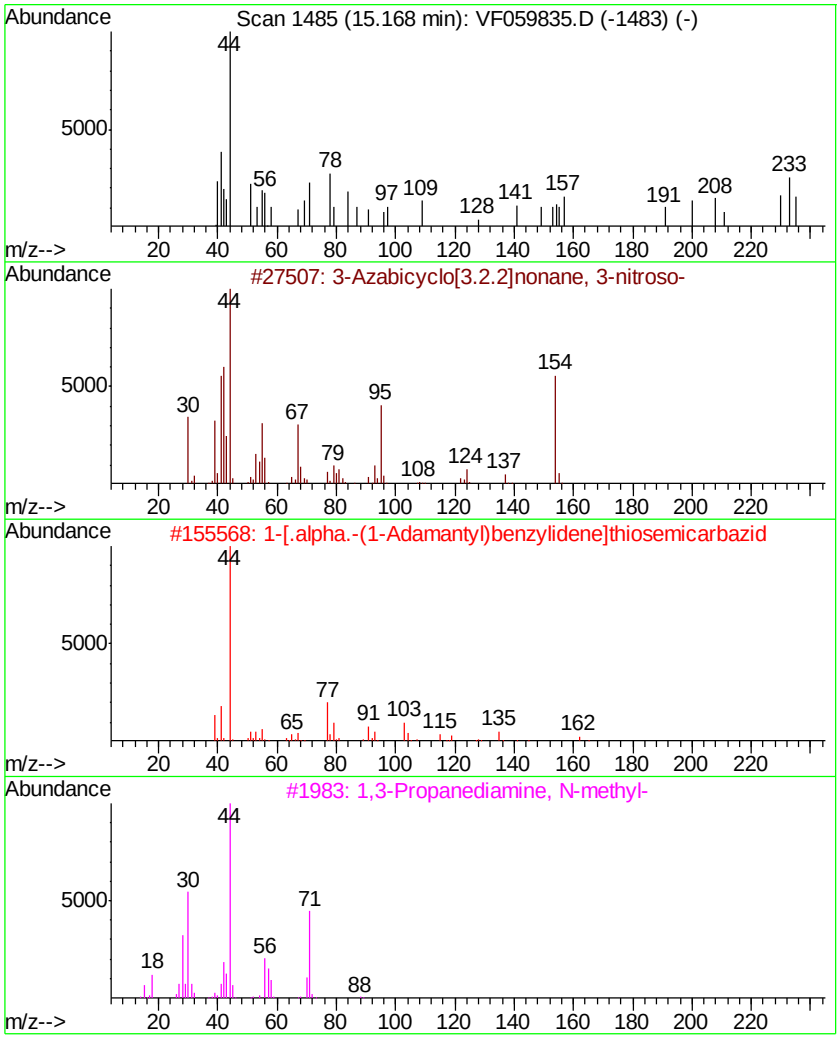
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

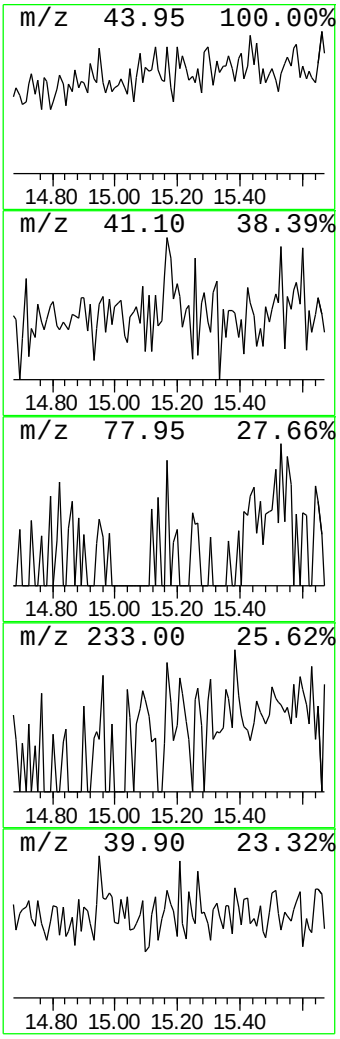
****

Peak Number 7 unknown15.17 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.17	18.63 ug/l	5087	1,4-Dichlorobenzene-d4	12.62

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Azabicyclo[3.2.2]nonane, 3-nit...	154	C8H14N2O	001522-09-4	23
2	1-[.alpha.-(1-Adamantyl)benzylid...	313	C18H23N3S	1000222-83-3	9
3	1,3-Propanediamine, N-methyl-	88	C4H12N2	006291-84-5	9
4	3,3'-Iminobispropylamine	131	C6H17N3	000056-18-8	9
5	3-Azahexan-1-ol, 6-cyclohexyl-	185	C11H23NO	1000251-64-6	9





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Data File : VF059835.D
Acq On : 3 Aug 2018 12:52
Operator : VA/AP
Sample : J4285-09
Misc : 6.01g/5mL/MSVOA F/SOIL
ALS Vial : 100 Sample Multiplier: 1

Instrument : MSVOA_F
ClientSampleId : KG-PIT-3

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_F\METHODS\82F073018S.M
Quant Title : SW846 8260

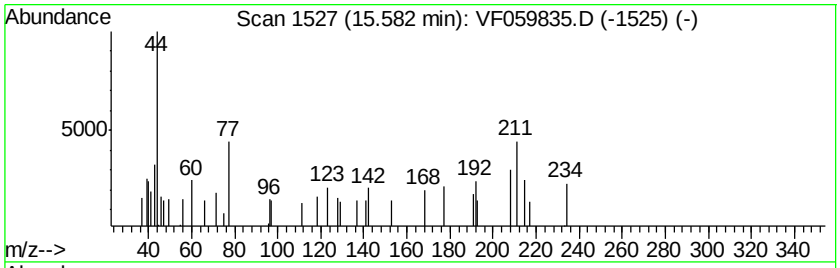
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

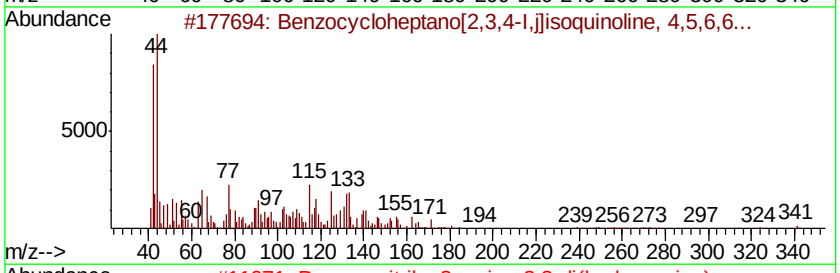
****

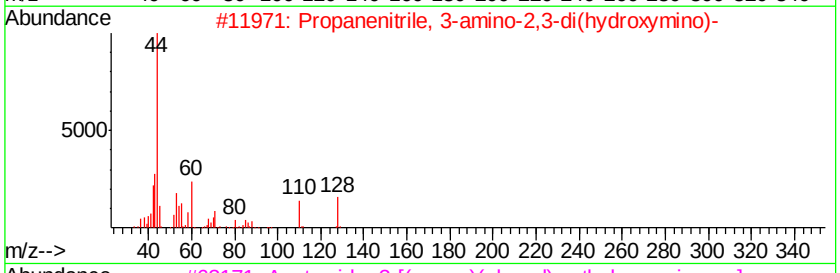
Peak Number 8 unknown15.58 Concentration Rank 9

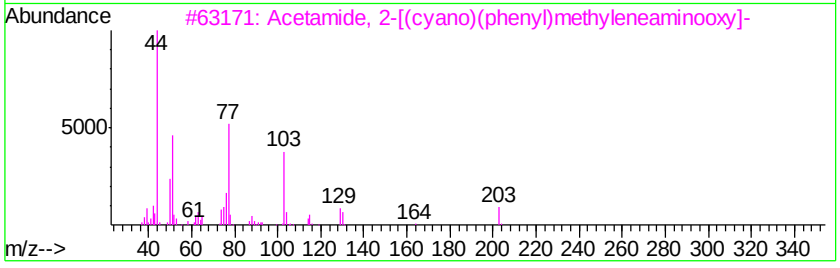
R.T.	EstConc	Area	Relative to ISTD	R.T.
15.58	13.09 ug/l	3576	1,4-Dichlorobenzene-d4	12.62

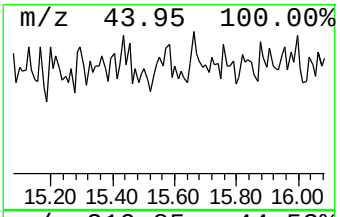
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzocycloheptano[2,3,4-I,j]isoq...	341	C20H23NO4	1000127-82-2	10
2	Propanenitrile, 3-amino-2,3-di(h...	128	C3H4N4O2	206363-14-6	9
3	Acetamide, 2-[(cyano)(phenyl)met...	203	C10H9N3O2	1000317-12-4	9
4	Acetamide, 2-fluoro-	77	C2H4FNO	000640-19-7	7
5	N,N-Dimethylethanesulfonamide	137	C4H11NO2S	006338-68-7	5

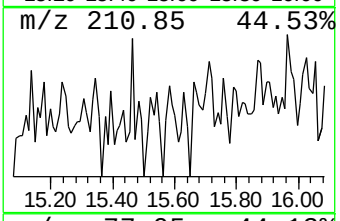


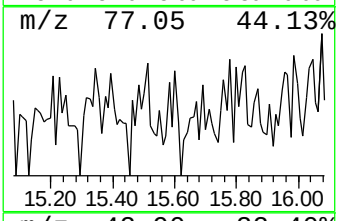


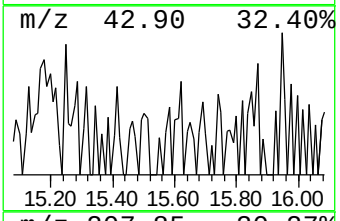


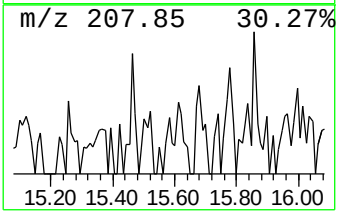












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Data File : VF059835.D
Acq On : 3 Aug 2018 12:52
Operator : VA/AP
Sample : J4285-09
Misc : 6.01g/5mL/MSVOA F/SOIL
ALS Vial : 100 Sample Multiplier: 1

Instrument : MSVOA_F
ClientSampleId : KG-PIT-3

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_F\METHODS\82F073018S.M
Quant Title : SW846 8260

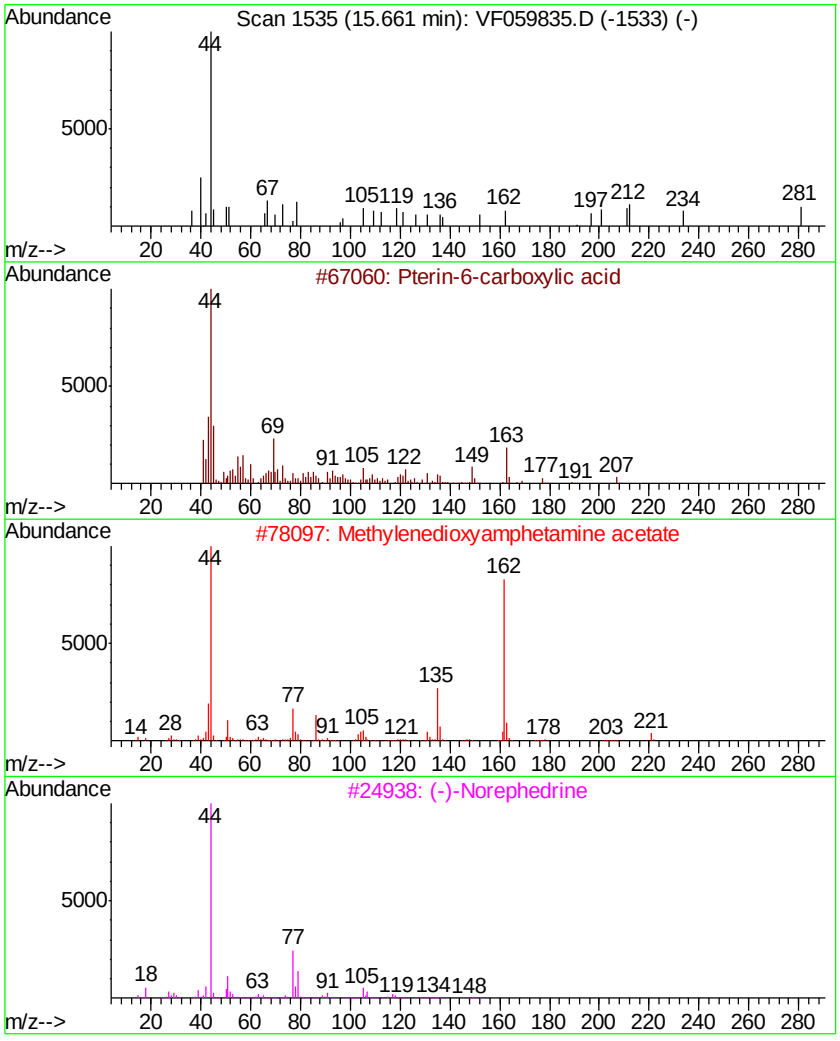
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

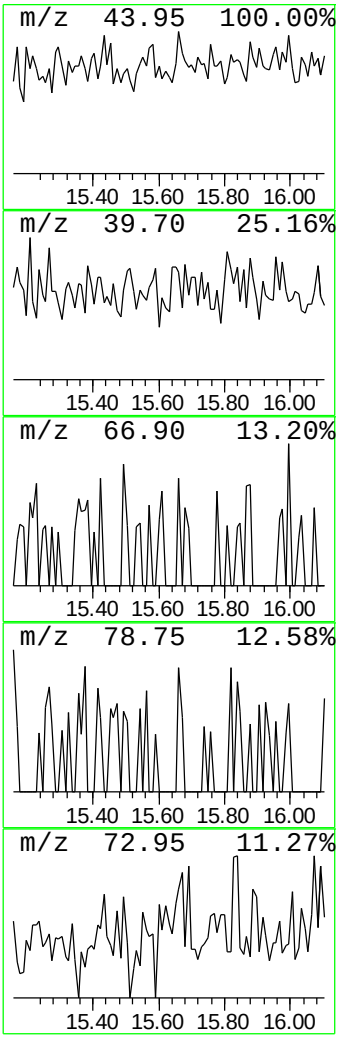
****

Peak Number 9 unknown15.66 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.66	15.08 ug/l	4119	1,4-Dichlorobenzene-d4	12.62

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pterin-6-carboxylic acid	207	C7H5N5O3	000948-60-7	28
2	Methylenedioxyamphetamine acetate	221	C12H15NO3	036209-71-9	9
3	(-)-Norephedrine	151	C9H13NO	000492-41-1	9
4	Metaraminol	167	C9H13NO2	000054-49-9	9
5	Pteridine-8-oxide, 6-aldoximino-...	222	C7H6N6O3	023663-23-2	9





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Operator : VA/AP
Sample : J4285-09
Misc : 6.01g/5mL/MSVOA F/SOIL
ALS Vial : 100 Sample Multiplier: 1

Instrument : MSVOA_F
ClientSampleId : KG-PIT-3

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_F\METHODS\82F073018S.M
Quant Title : SW846 8260

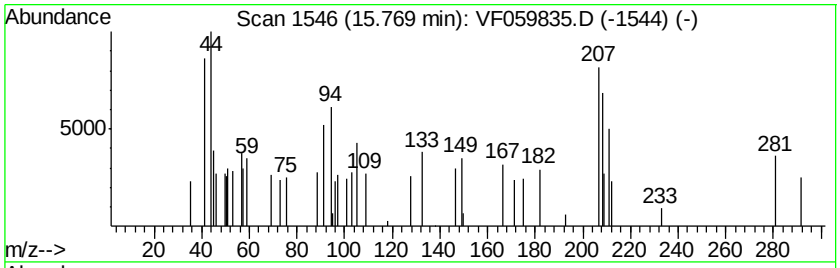
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

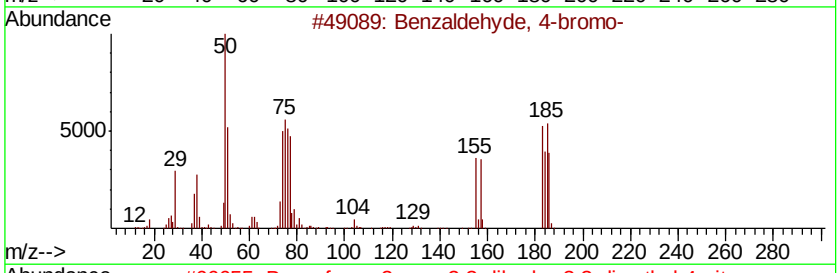
****

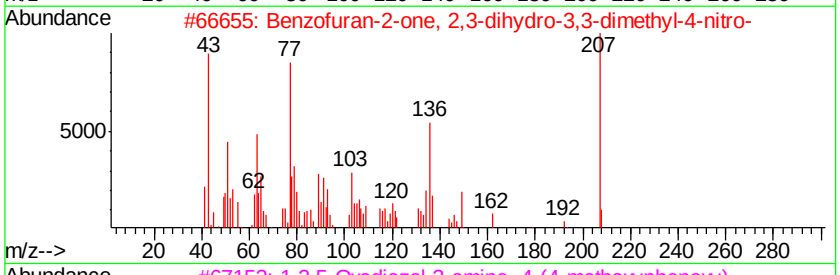
Peak Number 10 unknown15.77 Concentration Rank 1

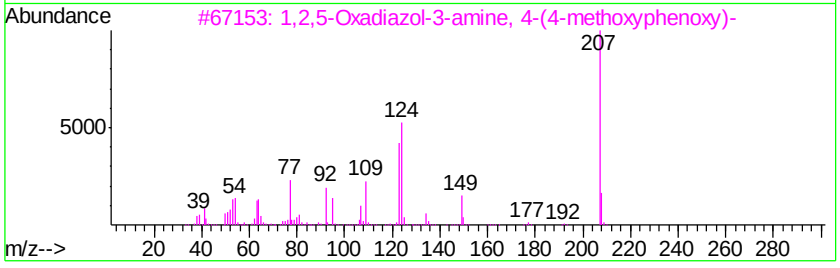
R.T.	EstConc	Area	Relative to ISTD	R.T.
15.77	28.17 ug/l	7694	1,4-Dichlorobenzene-d4	12.62

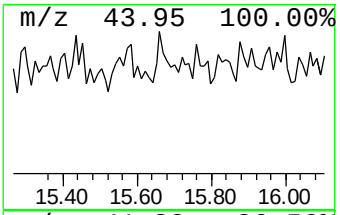
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzaldehyde, 4-bromo-	184	C7H5BrO	001122-91-4	14
2	Benzofuran-2-one, 2,3-dihydro-3,...	207	C10H9NO4	1000129-53-4	10
3	1,2,5-Oxadiazol-3-amine, 4-(4-me...	207	C9H9N3O3	1000351-72-7	9
4	1,3,5-Triazine, 2-chloro-4,6-bis...	207	C5H6ClN3S2	004407-40-3	9
5	2-Amino-7,7-dimethyl-5,6,7,8-tet...	211	C9H13N3OS	123973-48-8	9

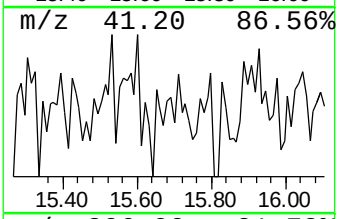


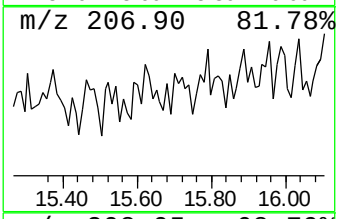


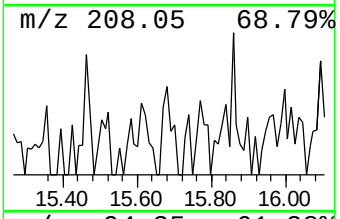


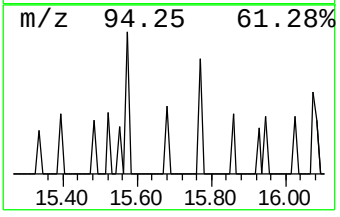












Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_F\DATA\VF080218\
Data File : VF059835.D
Acq On : 3 Aug 2018 12:52
Operator : VA/AP
Sample : J4285-09
Misc : 6.01g/5mL/MSVOA_F/SOIL
ALS Vial : 100 Sample Multiplier: 1

Instrument : MSVOA_F
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Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_F\METHODS\82F073018S.M
Quant Title : SW846 8260

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

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
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unknown14.21	14.21	21.9	ug/l	5994	4	12.62	13656	50.0
unknown14.39	14.39	20.6	ug/l	5614	4	12.62	13656	50.0
unknown14.72	14.72	19.6	ug/l	5363	4	12.62	13656	50.0
unknown14.89	14.89	19.3	ug/l	5268	4	12.62	13656	50.0
unknown15.17	15.17	18.6	ug/l	5087	4	12.62	13656	50.0
unknown15.58	15.58	13.1	ug/l	3576	4	12.62	13656	50.0
unknown15.66	15.66	15.1	ug/l	4119	4	12.62	13656	50.0
unknown15.77	15.77	28.2	ug/l	7694	4	12.62	13656	50.0