

Data Path : Z:\VOASRV\HPCHEM1\MSVOA W\DATA\VW030219\
 Data File : VW008958.D
 Acq On : 02 Mar 2019 09:07
 Operator : SY/VA
 Sample : K1671-02
 Misc : 4.88G/5ML/MSVOA W/SOIL
 ALS Vial : 51 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 FK-BPM-04-002-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:\VOASRV\HPCHEM1\MSVOA_W\METHOD\82W022119S.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	1.965	6	10	15	rVB4	5429	7097	2.58%	0.361%
2	2.019	16	19	23	rVB3	4456	6163	2.24%	0.313%
3	2.154	36	41	48	rBV3	3928	6533	2.37%	0.332%
4	2.257	48	58	60	rBV8	6827	21786	7.91%	1.107%
5	2.629	112	119	126	rBV10	7886	19582	7.11%	0.995%
6	2.739	135	137	141	rBV4	2541	4199	1.52%	0.213%
7	2.836	149	153	154	rBV4	2584	3405	1.24%	0.173%
8	3.031	183	185	186	rBV2	4746	2842	1.03%	0.144%
9	4.135	363	366	372	rVB5	3819	6639	2.41%	0.337%
10	4.915	486	494	506	rBV9	8649	27763	10.08%	1.411%
11	5.946	657	663	671	rBV9	2988	7269	2.64%	0.369%
12	7.147	848	860	869	rBV	24582	75302	27.33%	3.826%
13	7.884	972	981	986	rBV2	39155	93752	34.02%	4.764%
14	7.951	986	992	1002	rVB	78900	177107	64.27%	9.000%
15	8.256	1034	1042	1046	rBV	36740	91611	33.25%	4.655%
16	8.305	1047	1050	1060	rVV2	42822	88809	32.23%	4.513%
17	8.683	1106	1112	1115	rBV5	2277	4957	1.80%	0.252%
18	8.841	1131	1138	1148	rBV	101520	202937	73.65%	10.312%
19	9.494	1239	1245	1248	rVB5	1481	2922	1.06%	0.148%
20	10.323	1373	1381	1390	rBV	147561	275546	100.00%	14.002%
21	10.512	1408	1412	1416	rBV	7139	11998	4.35%	0.610%
22	10.713	1437	1445	1448	rBV3	16930	37919	13.76%	1.927%
23	11.067	1497	1503	1504	rBV6	3871	6302	2.29%	0.320%
24	11.445	1560	1565	1572	rVB4	8109	16169	5.87%	0.822%
25	11.634	1587	1596	1607	rBV	148507	257349	93.40%	13.077%
26	11.902	1637	1640	1643	rBV5	2012	3072	1.11%	0.156%
27	12.621	1750	1758	1765	rBV	136693	242796	88.11%	12.337%
28	13.304	1865	1870	1871	rBV4	3565	5288	1.92%	0.269%
29	13.560	1906	1912	1919	rBV	139203	222130	80.61%	11.287%
30	13.651	1923	1927	1930	rBV5	2131	2909	1.06%	0.148%
31	13.962	1974	1978	1982	rBV7	2513	4811	1.75%	0.244%
32	14.078	1993	1997	2003	rVB3	8247	14035	5.09%	0.713%
33	14.535	2065	2072	2073	rBV7	3784	7034	2.55%	0.357%
34	14.938	2136	2138	2142	rVV4	2295	3973	1.44%	0.202%

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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 >
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35	15.096	2158	2164	2166	rVV6	1436	2889	1.05%	0.147%
36	16.724	2427	2431	2435	rBV6	1705	3061	1.11%	0.156%

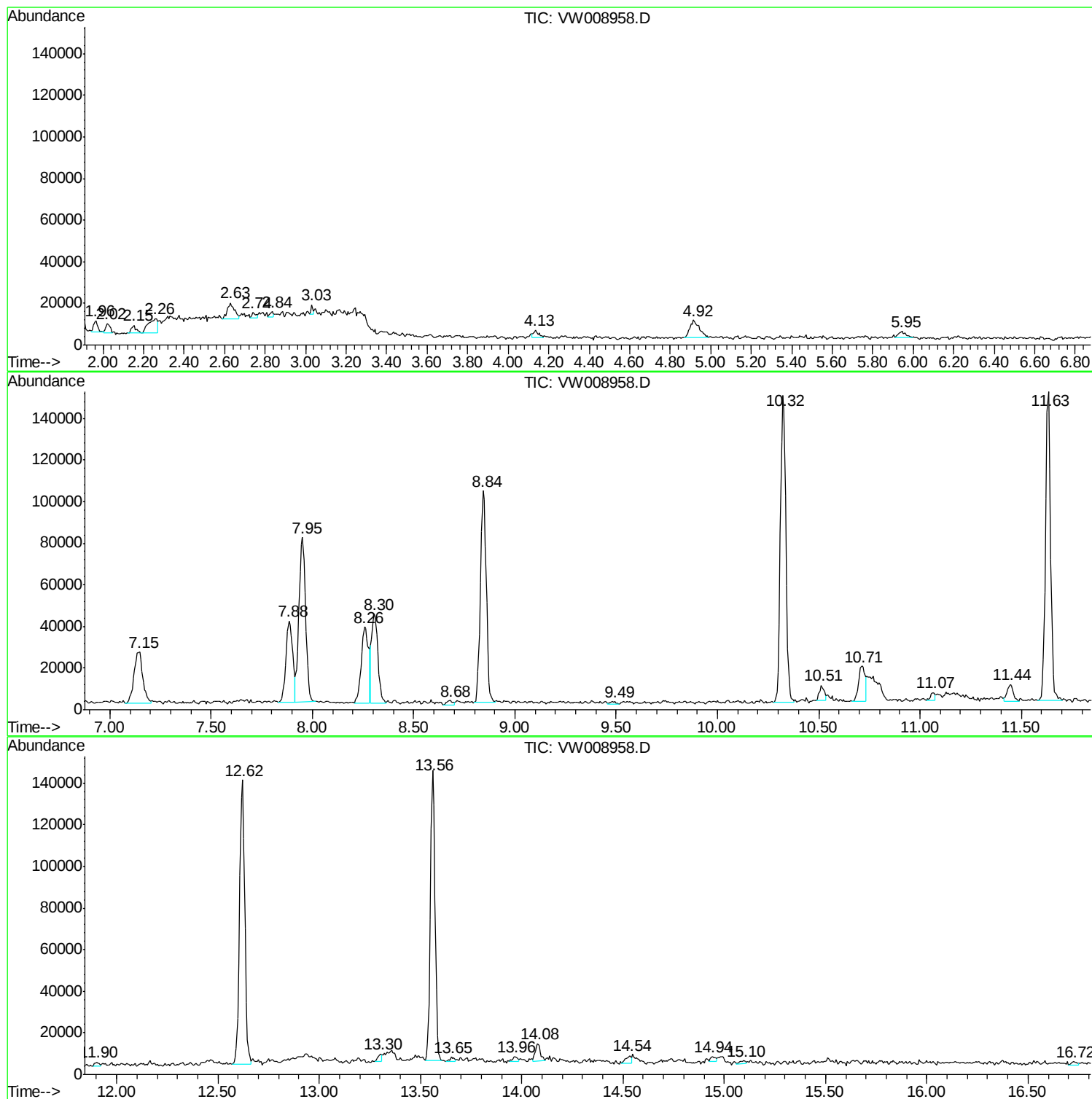
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W022119S.M
Quant Title : SW846 8260

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



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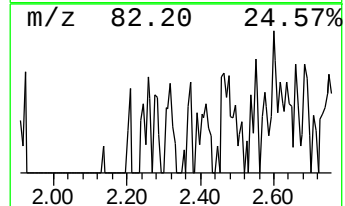
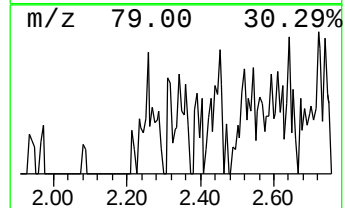
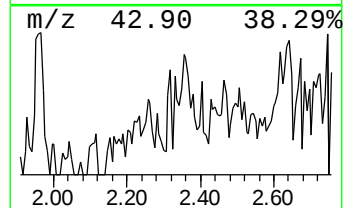
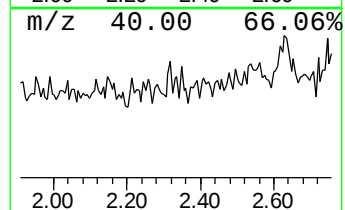
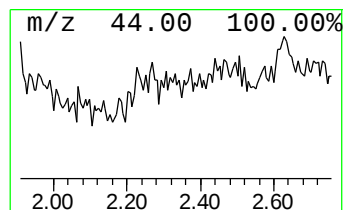
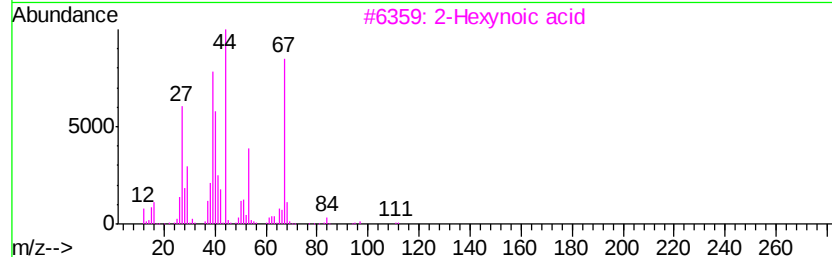
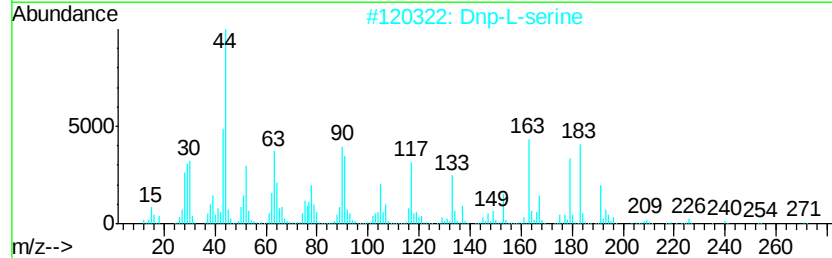
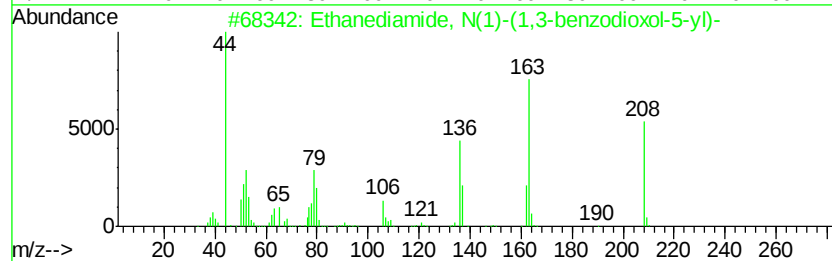
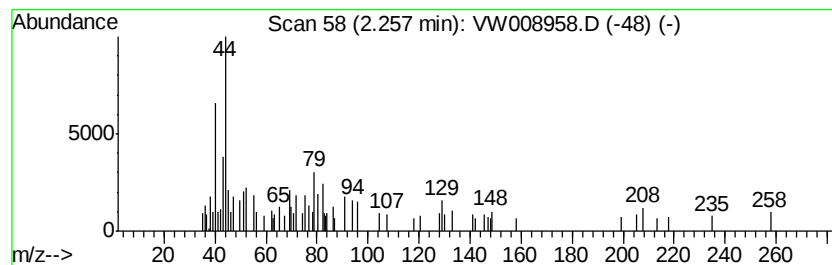
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W022119S.M
Quant Title : SW846 8260

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

Peak Number 1 unknown-01 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.26	6.15 ug/l	21786	Pentafluorobenzene	7.95

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanediamide, N(1)-(1,3-benzodi...	208	C9H8N2O4	1000351-43-9	43
2			Dnp-L-serine	271	C9H9N3O7	001655-64-7	33
3			2-Hexynoic acid	112	C6H8O2	000764-33-0	27
4			Propanenitrile, 3-chloro-2,3-dih...	147	C3H2ClN3O2	139912-33-7	23
5			1-Nitrobiuret	148	C2H4N4O4	016326-62-8	17



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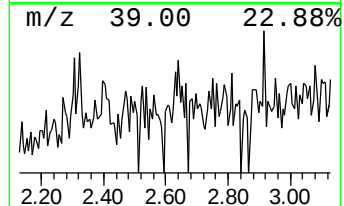
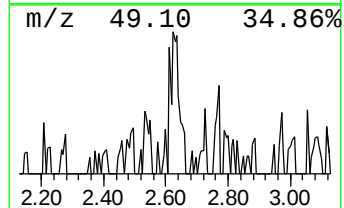
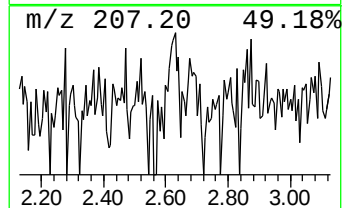
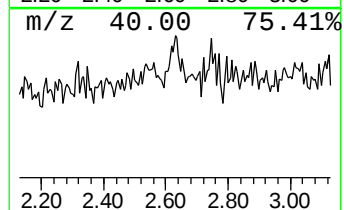
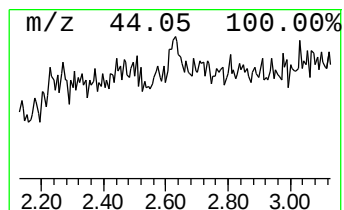
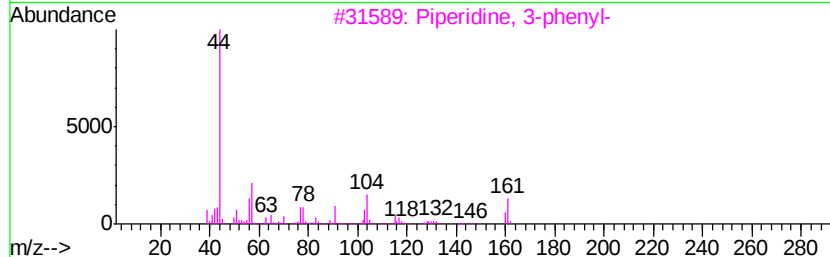
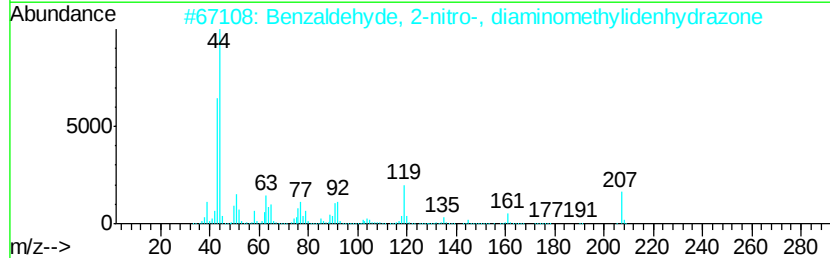
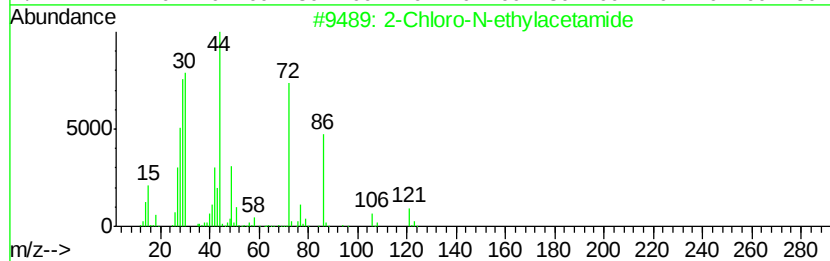
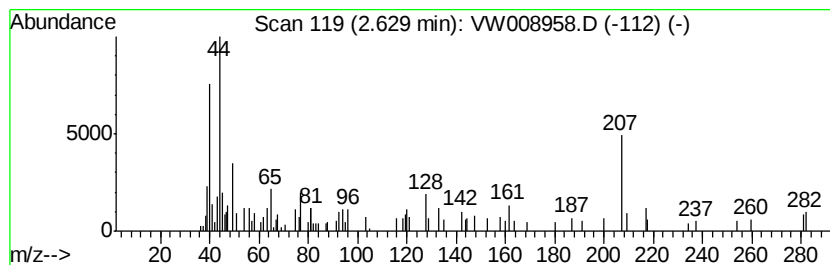
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Peak Number 2 unknown-02 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.63	5.53 ug/l	19582	Pentafluorobenzene	7.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Chloro-N-ethylacetamide	121	C4H8ClNO	000105-35-1	28
2		Benzaldehyde, 2-nitro-, diaminom...	207	C8H9N5O2	102632-31-5	17
3		Piperidine, 3-phenyl-	161	C11H15N	003973-62-4	14
4		Acetic acid, 2-(1-methyl-2-oxohy...	252	C10H12N4O4	1000327-59-6	14
5		Acetamide, 2-cyano-	84	C3H4N2O	000107-91-5	9



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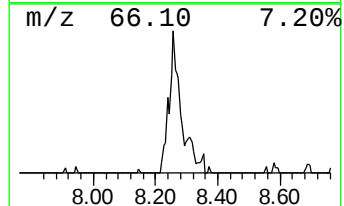
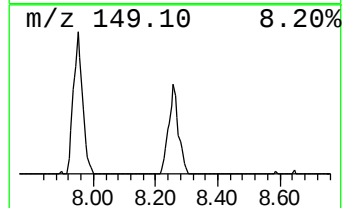
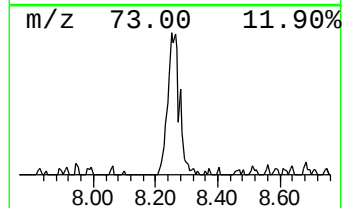
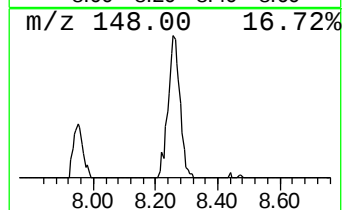
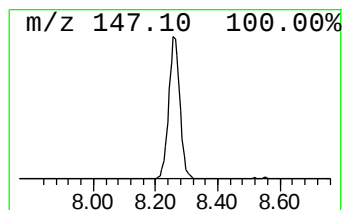
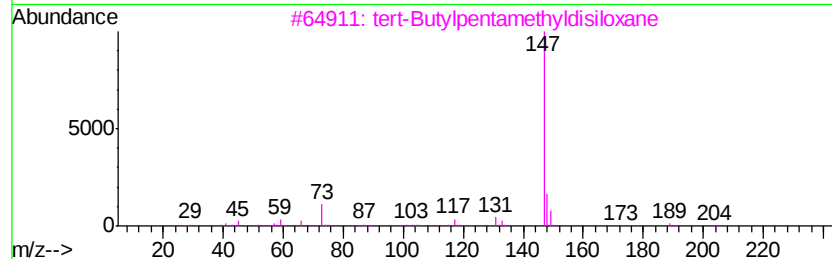
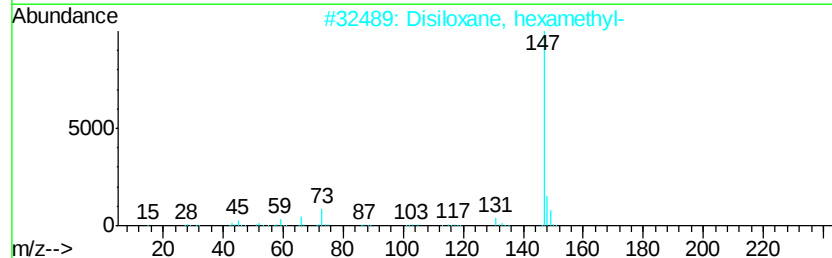
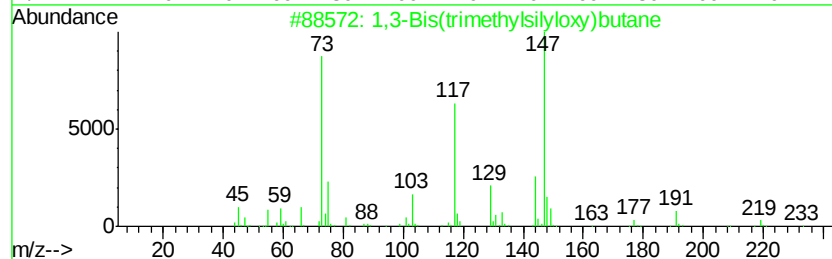
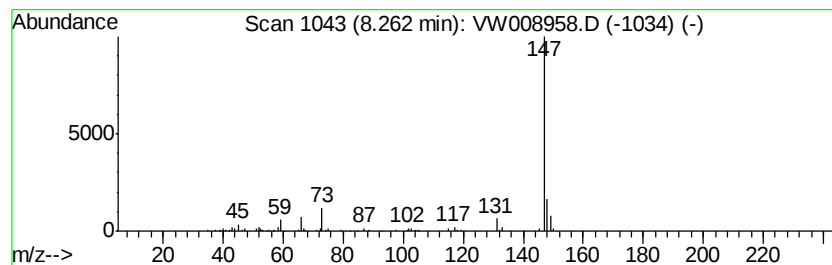
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TIC Integration Parameters: LSCINT.P

Peak Number 4 1,3-Bis(trimethylsilyloxy)b... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.26	25.86 ug/l	91611	Pentafluorobenzene	7.95

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,3-Bis(trimethylsilyloxy)butane	234	C10H26O2Si2	056771-47-2	83
2			Disiloxane, hexamethyl-	162	C6H18OSi2	000107-46-0	83
3			tert-Butylpentamethyldisiloxane	204	C9H24OSi2	067875-54-1	83
4			1-Pentamethyldisilyloxy-3-phenyl...	264	C14H24OSi2	1000216-94-9	78
5			3,7-Dioxa-2,8-disilanonane, 2,2,...	220	C9H24O2Si2	017887-80-8	74



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown-01	2.26	6.2	ug/l	21786	1	7.95	177107	50.0
unknown-02	2.63	5.5	ug/l	19582	1	7.95	177107	50.0
1,3-Bis(trimethyl...	8.26	25.9	ug/l	91611	1	7.95	177107	50.0