

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN092220\
 Data File : VN063573.D
 Acq On : 23 Sep 2020 8:28
 Operator : JC/MD
 Sample : L4101-02
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 52 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 0-6FT

Quant Time: Sep 23 09:29:17 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N091620W.M
 Quant Title : SW846 8260
 QLast Update : Wed Sep 16 13:05:20 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.63	168	201310	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	388058	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	394445	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	159161	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	173827	50.73	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.46%	
35) Dibromofluoromethane	7.55	113	125260	46.63	ug/l	0.00
Spiked Amount 50.000			Recovery	=	93.26%	
50) Toluene-d8	10.06	98	500181	49.52	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.04%	
62) 4-Bromofluorobenzene	12.38	95	197844	53.43	ug/l	0.00
Spiked Amount 50.000			Recovery	=	106.86%	
Target Compounds						
16) Acetone	3.78	43	13079	11.876	ug/l	95
18) Methyl Acetate	4.30	43	4622	1.653	ug/l #	54
20) Methylene Chloride	4.50	84	13087	4.690	ug/l	85
43) Isopropyl Acetate	8.13	43	79660	10.980	ug/l	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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