

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN110620\
 Data File : VN064494.D
 Acq On : 6 Nov 2020 19:40
 Operator : JC/MD
 Sample : L4656-10
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 SR-L14-L13-SOUTH

Manual Integrations
 APPROVED

MMDadoda
 11/9/2020 9:56:38 AM

Quant Time: Nov 07 04:42:13 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N110420W.M
 Quant Title : SW846 8260
 QLast Update : Wed Nov 04 12:09:03 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.63	168	242842	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	523148	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	551534	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	206689	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	237974	56.26	ug/l	0.00
Spiked Amount 50.000			Recovery	=	112.52%	
35) Dibromofluoromethane	7.55	113	176306	49.50	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.00%	
50) Toluene-d8	10.06	98	675117	49.04	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.08%	
62) 4-Bromofluorobenzene	12.38	95	257559	49.47	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.94%	
Target Compounds						
16) Acetone	3.78	43	42810	26.569	ug/l	100
18) Methyl Acetate	4.29	43	6243m	1.971	ug/l	
20) Methylene Chloride	4.50	84	14158	4.429	ug/l	93
25) 2-Butanone	6.80	43	11101	5.407	ug/l	91
43) Isopropyl Acetate	8.13	43	98810	11.032	ug/l	99

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

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