

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN100120\
 Data File : VN063777.D
 Acq On : 1 Oct 2020 17:22
 Operator : JC/MD
 Sample : L4200-01
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB-2020-WS-000-01

Quant Time: Oct 02 03:02:56 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N093020W.M
 Quant Title : SW846 8260
 QLast Update : Thu Oct 01 01:46:38 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.63	168	212199	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	399911	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	408816	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	171537	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	180756	51.56	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.12%	
35) Dibromofluoromethane	7.55	113	134454	48.45	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.90%	
50) Toluene-d8	10.06	98	511525	48.63	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.26%	
62) 4-Bromofluorobenzene	12.38	95	202749	51.27	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.54%	
Target Compounds						
16) Acetone	3.78	43	31844	29.508	ug/l	98
18) Methyl Acetate	4.29	43	4321	1.542	ug/l #	69
20) Methylene Chloride	4.51	84	9143	3.790	ug/l	91
43) Isopropyl Acetate	8.13	43	72121	10.863	ug/l	100

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

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