

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN103020\
 Data File : VN064407.D
 Acq On : 30 Oct 2020 18:03
 Operator : JC/MD
 Sample : L4568-01
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 CHASSIS-WASH

Manual Integrations
 APPROVED

MMDadoda
 11/2/2020 4:12:09 PM

Quant Time: Nov 02 03:17:13 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N102720W.M
 Quant Title : SW846 8260
 QLast Update : Tue Oct 27 12:23:17 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.63	168	201978	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	365197	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	363771	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	154371	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	169790	52.31	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.62%	
35) Dibromofluoromethane	7.55	113	120070	49.30	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.60%	
50) Toluene-d8	10.06	98	457117	50.74	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.48%	
62) 4-Bromofluorobenzene	12.37	95	194058	55.25	ug/l	0.00
Spiked Amount 50.000			Recovery	=	110.50%	
Target Compounds						
16) Acetone	3.78	43	43859	33.267	ug/l	98
18) Methyl Acetate	4.28	43	3442m	1.412	ug/l	
20) Methylene Chloride	4.51	84	17640	7.663	ug/l	94
25) 2-Butanone	6.79	43	15018	9.215	ug/l #	83
43) Isopropyl Acetate	8.13	43	28413	4.465	ug/l #	87
51) 4-Methyl-2-Pentanone	9.95	43	76962	20.804	ug/l	95

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

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