

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN110620\
 Data File : VN064493.D
 Acq On : 6 Nov 2020 19:15
 Operator : JC/MD
 Sample : L4656-03
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 16 Sample Multiplier: 1

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

Manual Integrations
 APPROVED

MMDadoda
 11/9/2020 9:56:43 AM

Quant Time: Nov 07 04:37:06 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	250273	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	523192	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	558729	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	198916	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	244334	56.05	ug/l	0.00
Spiked Amount 50.000			Recovery	=	112.10%	
35) Dibromofluoromethane	7.55	113	178828	50.20	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.40%	
50) Toluene-d8	10.06	98	686270	49.85	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.70%	
62) 4-Bromofluorobenzene	12.38	95	258041	49.56	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.12%	
Target Compounds						
16) Acetone	3.78	43	29378	17.692	ug/l	95
18) Methyl Acetate	4.29	43	5639m	1.727	ug/l	
20) Methylene Chloride	4.51	84	14146	4.294	ug/l	96
43) Isopropyl Acetate	8.13	43	93208	10.406	ug/l	99

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

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