

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN101220\
 Data File : VN064090.D
 Acq On : 12 Oct 2020 20:14
 Operator : JC/MD
 Sample : L4333-09
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 OH-WC-C-G2

Quant Time: Oct 13 02:41:59 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	226418	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	410953	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	428366	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	171573	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	186739	49.92	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.84%	
35) Dibromofluoromethane	7.55	113	138441	48.55	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.10%	
50) Toluene-d8	10.06	98	531214	49.14	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.28%	
62) 4-Bromofluorobenzene	12.38	95	206494	50.82	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.64%	
Target Compounds						
16) Acetone	3.78	43	102487	89.005	ug/l	99
18) Methyl Acetate	4.29	43	4168	1.394	ug/l #	63
20) Methylene Chloride	4.50	84	12535	4.841	ug/l #	93
25) 2-Butanone	6.80	43	9452	5.790	ug/l	91

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

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