

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN091620\
 Data File : VN063401.D
 Acq On : 17 Sep 2020 6:39
 Operator : JC/MD
 Sample : L4015-10
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 51 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 REACTOR-6E-(C-4)

Quant Time: Sep 17 07:30:51 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N091620W.M
 Quant Title : SW846 8260
 QLast Update : Wed Sep 16 13:05:20 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.63	168	229787	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	428251	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	426555	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	180971	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	193629	49.50	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.00%	
35) Dibromofluoromethane	7.55	113	142727	48.14	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.28%	
50) Toluene-d8	10.06	98	560735	50.30	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.60%	
62) 4-Bromofluorobenzene	12.38	95	219361	53.68	ug/l	0.00
Spiked Amount 50.000			Recovery	=	107.36%	
Target Compounds						
16) Acetone	3.79	43	16475	13.105	ug/l	100
18) Methyl Acetate	4.29	43	5363	1.680	ug/l #	85
20) Methylene Chloride	4.51	84	12858	4.037	ug/l	94
43) Isopropyl Acetate	8.13	43	87594	10.940	ug/l	100

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

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