

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN091620\
 Data File : VN063399.D
 Acq On : 17 Sep 2020 5:51
 Operator : JC/MD
 Sample : L4015-08
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 49 Sample Multiplier: 1

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

Quant Time: Sep 17 06:45:14 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	225711	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	427101	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	422051	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	180893	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	191453	49.83	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.66%	
35) Dibromofluoromethane	7.55	113	140471	47.51	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.02%	
50) Toluene-d8	10.06	98	560688	50.43	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.86%	
62) 4-Bromofluorobenzene	12.38	95	215099	52.78	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.56%	
Target Compounds						
16) Acetone	3.78	43	16815	13.617	ug/l	99
18) Methyl Acetate	4.29	43	5917	1.888	ug/l #	65
20) Methylene Chloride	4.51	84	13403	4.284	ug/l	87
43) Isopropyl Acetate	8.13	43	88203	11.046	ug/l	99

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

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