

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN091620\
 Data File : VN063400.D
 Acq On : 17 Sep 2020 6:15
 Operator : JC/MD
 Sample : L4015-09
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 50 Sample Multiplier: 1

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

Quant Time: Sep 17 06:46:26 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	227080	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	428800	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	425877	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	184189	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	191763	49.61	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.22%	
35) Dibromofluoromethane	7.55	113	140261	47.25	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.50%	
50) Toluene-d8	10.06	98	557803	49.97	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.94%	
62) 4-Bromofluorobenzene	12.38	95	217079	53.05	ug/l	0.00
Spiked Amount 50.000			Recovery	=	106.10%	
Target Compounds						
16) Acetone	3.78	43	16284	13.108	ug/l	96
18) Methyl Acetate	4.29	43	4385	1.390	ug/l #	85
20) Methylene Chloride	4.51	84	13648	4.336	ug/l #	84
43) Isopropyl Acetate	8.13	43	89120	11.116	ug/l	99

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

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