

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN091620\
 Data File : VN063402.D
 Acq On : 17 Sep 2020 7:03
 Operator : JC/MD
 Sample : L4015-11
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 52 Sample Multiplier: 1

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

Quant Time: Sep 17 07:32:16 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	224890	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	417968	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	414363	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	179266	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	186845	48.81	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.62%	
35) Dibromofluoromethane	7.55	113	135977	46.99	ug/l	0.00
Spiked Amount 50.000			Recovery	=	93.98%	
50) Toluene-d8	10.06	98	539022	49.54	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.08%	
62) 4-Bromofluorobenzene	12.38	95	213547	53.54	ug/l	0.00
Spiked Amount 50.000			Recovery	=	107.08%	
Target Compounds						
16) Acetone	3.78	43	15737	12.791	ug/l	96
18) Methyl Acetate	4.30	43	5874	1.881	ug/l #	70
20) Methylene Chloride	4.50	84	13381	4.293	ug/l	96
43) Isopropyl Acetate	8.13	43	84461	10.808	ug/l	99

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN091620\
Data File : VN063402.D
Acq On : 17 Sep 2020 7:03
Operator : JC/MD
Sample : L4015-11
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 52 Sample Multiplier: 1

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

Quant Time: Sep 17 07:32:16 2020
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N091620W.M
Quant Title : SW846 8260
QLast Update : Wed Sep 16 13:05:20 2020
Response via : Initial Calibration

