

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN091620\
 Data File : VN063397.D
 Acq On : 17 Sep 2020 5:03
 Operator : JC/MD
 Sample : L4015-06
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 47 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 REACTOR-6E-(C-0)

Manual Integrations
 APPROVED

MMDadoda
 9/17/2020 1:44:17 PM

Quant Time: Sep 17 06:42:37 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	231866	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	439518	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	426015	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	182475	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	198054	50.18	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.36%	
35) Dibromofluoromethane	7.55	113	143764	47.25	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.50%	
50) Toluene-d8	10.06	98	563734	49.27	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.54%	
62) 4-Bromofluorobenzene	12.38	95	215473	51.38	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.76%	
Target Compounds						
16) Acetone	3.79	43	16238	12.801	ug/l	98
18) Methyl Acetate	4.28	43	5759m	1.788	ug/l	
20) Methylene Chloride	4.51	84	12325	3.835	ug/l	90
43) Isopropyl Acetate	8.13	43	75137	9.144	ug/l	100

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

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