

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN091119\
 Data File : VN057868.D
 Acq On : 10 Sep 2019 17:24
 Operator : JC/SP
 Sample : K4623-23
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PAR-11-C4-(0)

Manual Integrations
 APPROVED

MMDadoda
 9/13/2019 11:46:29 PM

Quant Time: Sep 13 05:59:24 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N091019W.M
 Quant Title : SW846 8260
 QLast Update : Tue Sep 10 03:02:13 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.65	168	294151	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.58	114	493597	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	454448	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	228927	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.01	65	294419	49.42	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.84%	
35) Dibromofluoromethane	7.58	113	199941	43.89	ug/l	0.00
Spiked Amount 50.000			Recovery	=	87.78%	
50) Toluene-d8	10.09	98	815236	46.09	ug/l	0.00
Spiked Amount 50.000			Recovery	=	92.18%	
62) 4-Bromofluorobenzene	12.40	95	262746	38.98	ug/l	0.00
Spiked Amount 50.000			Recovery	=	77.96%	
Target Compounds						
16) Acetone	3.81	43	10981m	5.622	ug/l	
18) Methyl Acetate	4.31	43	7107	1.365	ug/l #	86
43) Isopropyl Acetate	8.16	43	74963	7.537	ug/l #	90

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_N\DATA\VN091119\
Data File : VN057868.D
Acq On : 10 Sep 2019 17:24
Operator : JC/SP
Sample : K4623-23
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
PAR-11-C4-(0)

Manual Integrations
APPROVED

MMDadoda
9/13/2019 11:46:29 PM

Quant Time: Sep 13 05:59:24 2019
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N091019W.M
Quant Title : SW846 8260
QLast Update : Tue Sep 10 03:02:13 2019
Response via : Initial Calibration

