

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN111020\
 Data File : VN064540.D
 Acq On : 10 Nov 2020 21:02
 Operator : JC/MD
 Sample : L4680-08
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PAV-B7-110920-19-20

Manual Integrations
 APPROVED

MMDadoda
 11/11/2020 1:02:06 PM

Quant Time: Nov 11 03:35:17 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N110420W.M
 Quant Title : SW846 8260
 QLast Update : Wed Nov 04 12:09:03 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.63	168	240430	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	514439	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	566592	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	204634	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	237039	56.60	ug/l	0.00
Spiked Amount 50.000			Recovery	=	113.20%	
35) Dibromofluoromethane	7.55	113	177414	50.65	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.30%	
50) Toluene-d8	10.06	98	676058	49.94	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.88%	
62) 4-Bromofluorobenzene	12.38	95	255983	50.00	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.00%	
Target Compounds						
16) Acetone	3.79	43	9956	6.241	ug/l	99
20) Methylene Chloride	4.51	84	13857m	4.378	ug/l	
43) Isopropyl Acetate	8.13	43	44707	5.076	ug/l	97

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN111020\
Data File : VN064540.D
Acq On : 10 Nov 2020 21:02
Operator : JC/MD
Sample : L4680-08
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
PAV-B7-110920-19-20

Manual Integrations
APPROVED

MMDadoda
11/11/2020 1:02:06 PM

Quant Time: Nov 11 03:35:17 2020
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N110420W.M
Quant Title : SW846 8260
QLast Update : Wed Nov 04 12:09:03 2020
Response via : Initial Calibration

