

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN092120\
 Data File : VN063499.D
 Acq On : 21 Sep 2020 20:30
 Operator : JC/MD
 Sample : L4072-02
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-8

Manual Integrations
 APPROVED

MMDadoda
 9/22/2020 2:07:24 PM

Quant Time: Sep 22 04:32:12 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	262927	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	482500	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	481602	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	212262	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	208202	46.52	ug/l	0.00
Spiked Amount 50.000			Recovery	=	93.04%	
35) Dibromofluoromethane	7.55	113	152596	45.68	ug/l	0.00
Spiked Amount 50.000			Recovery	=	91.36%	
50) Toluene-d8	10.06	98	622704	49.58	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.16%	
62) 4-Bromofluorobenzene	12.38	95	250169	54.33	ug/l	0.00
Spiked Amount 50.000			Recovery	=	108.66%	
Target Compounds						
27) cis-1,2-Dichloroethene	6.79	96	11728	3.152	ug/l	82
30) Chloroform	7.34	83	3530m	0.538	ug/l	
44) Trichloroethene	8.80	130	3144	0.883	ug/l	79
64) Tetrachloroethene	10.60	164	7020	2.141	ug/l	95

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN092120\
Data File : VN063499.D
Acq On : 21 Sep 2020 20:30
Operator : JC/MD
Sample : L4072-02
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-8

Manual Integrations
APPROVED

MMDadoda
9/22/2020 2:07:24 PM

Quant Time: Sep 22 04:32:12 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

