

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN092520\
 Data File : VN063680.D
 Acq On : 25 Sep 2020 15:55
 Operator : JC/MD
 Sample : L4119-19
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 DUP06-20200922

Manual Integrations
 APPROVED

MMDadoda
 9/29/2020 10:40:43 AM

Quant Time: Sep 28 02:47:55 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	202061	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	374510	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	386286	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	158795	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	172788	50.23	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.46%	
35) Dibromofluoromethane	7.55	113	124815	48.14	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.28%	
50) Toluene-d8	10.06	98	486674	49.92	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.84%	
62) 4-Bromofluorobenzene	12.38	95	188297	52.69	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.38%	
Target Compounds						
3) Chloromethane	2.04	50	3122	1.031	ug/l	98
16) Acetone	3.78	43	5840	5.283	ug/l #	89
20) Methylene Chloride	4.50	84	2573m	0.919	ug/l	
44) Trichloroethene	8.80	130	7076	2.561	ug/l	98
64) Tetrachloroethene	10.60	164	4981	1.894	ug/l #	84

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

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