

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN092520\
 Data File : VN063695.D
 Acq On : 25 Sep 2020 21:54
 Operator : JC/MD
 Sample : L4133-05
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW4S

Quant Time: Sep 28 09:01:13 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	198501	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	372671	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	379750	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	159928	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	176160	52.13	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.26%	
35) Dibromofluoromethane	7.55	113	125315	48.57	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.14%	
50) Toluene-d8	10.06	98	481278	49.61	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.22%	
62) 4-Bromofluorobenzene	12.38	95	186246	52.37	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.74%	
Target Compounds						
					Qvalue	
3) Chloromethane	2.04	50	919	0.309	ug/l	93
16) Acetone	3.78	43	30604	28.182	ug/l	94
27) cis-1,2-Dichloroethene	6.78	96	12380	4.407	ug/l	84
44) Trichloroethene	8.81	130	5593	2.035	ug/l	87
64) Tetrachloroethene	10.60	164	106890	41.337	ug/l	95

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

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