

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN092520\
 Data File : VN063689.D
 Acq On : 25 Sep 2020 19:30
 Operator : JC/MD
 Sample : L4133-06
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 FIELD-BLANK

Quant Time: Sep 28 03:11:07 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	198824	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	374585	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	380894	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	157610	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	179090	52.91	ug/l	0.00
Spiked Amount	50.000		Recovery	=	105.82%	
35) Dibromofluoromethane	7.55	113	124566	48.04	ug/l	0.00
Spiked Amount	50.000		Recovery	=	96.08%	
50) Toluene-d8	10.06	98	480393	49.27	ug/l	0.00
Spiked Amount	50.000		Recovery	=	98.54%	
62) 4-Bromofluorobenzene	12.38	95	188481	52.73	ug/l	0.00
Spiked Amount	50.000		Recovery	=	105.46%	
Target Compounds						
					Qvalue	
3) Chloromethane	2.04	50	3308	1.110	ug/l	96
16) Acetone	3.78	43	8339	7.666	ug/l	100
20) Methylene Chloride	4.51	84	5501	1.996	ug/l #	84
44) Trichloroethene	8.80	130	938	0.339	ug/l	98
64) Tetrachloroethene	10.61	164	1078	0.416	ug/l #	72
65) Chlorobenzene	11.40	112	2591	0.320	ug/l #	76

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

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