

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN100120\
 Data File : VN063778.D
 Acq On : 1 Oct 2020 17:46
 Operator : JC/MD
 Sample : L4200-02
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB-2020-WW-000-01

Quant Time: Oct 02 03:07:11 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N093020W.M
 Quant Title : SW846 8260
 QLast Update : Thu Oct 01 01:46:38 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.63	168	215132	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	405253	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	407203	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	170332	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	187521	52.76	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.52%	
35) Dibromofluoromethane	7.55	113	135191	48.08	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.16%	
50) Toluene-d8	10.06	98	512073	48.04	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.08%	
62) 4-Bromofluorobenzene	12.38	95	200806	50.11	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.22%	
Target Compounds						
3) Chloromethane	2.04	50	1946	0.731	ug/l	95
16) Acetone	3.78	43	14569	13.316	ug/l	100
40) Benzene	8.01	78	5167	0.441	ug/l #	88
44) Trichloroethene	8.80	130	34528	11.238	ug/l	97
64) Tetrachloroethene	10.60	164	89469	28.399	ug/l	98

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

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