

Data Path : Z:\voasrv\HPCHEM1\MSVOA N\Data\VN112420\
 Data File : VN064810.D
 Acq On : 24 Nov 2020 15:03
 Operator : JC/MD
 Sample : L4829-08
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-13D

Quant Time: Dec 05 22:49:24 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N112320W.M
 Quant Title : SW846 8260
 QLast Update : Mon Nov 23 13:54:12 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.63	168	172737	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	310156	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	319819	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	149048	50.00	ug/l	0.00

System Monitoring Compounds

33) 1,2-Dichloroethane-d4	7.99	65	106460	53.01	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	106.02%
35) Dibromofluoromethane	7.55	113	95496	51.46	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	102.92%
50) Toluene-d8	10.06	98	385491	50.84	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	101.68%
62) 4-Bromofluorobenzene	12.38	95	155677	57.37	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	114.74%

Target Compounds

					Qvalue
3) Chloromethane	2.04	50	1007	0.518 ug/l #	74
16) Acetone	3.78	43	3591	5.440 ug/l	98
30) Chloroform	7.33	83	1777	0.521 ug/l	98
64) Tetrachloroethene	10.61	164	2044	0.657 ug/l #	84
67) Ethyl Benzene	11.48	91	4344	0.379 ug/l	94
84) 1,2,4-Trimethylbenzene	13.02	105	3957	0.454 ug/l	93
95) Naphthalene	15.10	128	39942	4.420 ug/l	99

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

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