

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091624\
 Data File : VN083913.D
 Acq On : 16 Sep 2024 13:27
 Operator : JC\MD
 Sample : P3874-03RE
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 BPOW6-8-HYD-20240904RE

Quant Time: Sep 17 02:31:20 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N091024W.M
 Quant Title : SW846 8260
 QLast Update : Wed Sep 11 15:18:56 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) Pentafluorobenzene	8.224	168	147918	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.100	114	254251	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	210527	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	79246	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.582	65	115695	51.872	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 103.740%	
35) Dibromofluoromethane	8.171	113	82527	50.702	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 101.400%	
50) Toluene-d8	10.565	98	303518	53.940	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 107.880%	
62) 4-Bromofluorobenzene	12.847	95	96906	41.742	ug/l	0.00
Spiked Amount	50.000	Range	77 - 121	Recovery	= 83.480%	
Target Compounds						
					Qvalue	
44) Trichloroethene	9.353	130	1846	1.081	ug/l	89
47) Bromodichloromethane	9.882	83	4666	1.727	ug/l #	89
60) Dibromochloromethane	11.353	129	4964	2.634	ug/l	99
71) Bromoform	12.570	173	2173	1.988	ug/l #	88

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

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