

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX100224\
 Data File : VX043234.D
 Acq On : 02 Oct 2024 16:18
 Operator : JC/MD
 Sample : P4267-01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 BPOW6-11-HYD-20241001

Quant Time: Oct 03 02:34:59 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100124W.M
 Quant Title : SW846 8260
 QLast Update : Wed Oct 02 16:50:57 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	81727	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.757	114	156737	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	139076	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.018	152	57983	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	71278	53.403	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 106.800%			
35) Dibromofluoromethane	5.385	113	52658	48.716	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 97.440%			
50) Toluene-d8	8.647	98	187525	50.062	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 100.120%			
62) 4-Bromofluorobenzene	11.079	95	67198	49.379	ug/l	0.00
Spiked Amount 50.000	Range 77 - 121		Recovery = 98.760%			
Target Compounds						
16) Acetone	2.386	43	1468	2.634	ug/l	92
18) Methyl Acetate	2.715	43	705	0.504	ug/l #	72
30) Chloroform	5.093	83	1490	0.776	ug/l #	83
47) Bromodichloromethane	7.824	83	2191	1.369	ug/l #	98
60) Dibromochloromethane	9.525	129	2342	1.958	ug/l	94
71) Bromoform	10.799	173	814	1.005	ug/l #	90

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

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