

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX102122\
 Data File : VX032312.D
 Acq On : 22 Oct 2022 01:42
 Operator : JC/MD
 Sample : N5248-11
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 FSND-MW-5-20221018

Quant Time: Oct 22 06:04:28 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X101922W.M
 Quant Title : SW846 8260
 QLast Update : Sat Oct 22 04:56:40 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	149344	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	232460	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	201161	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	93526	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	94503	48.642	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery =	97.280%		
35) Dibromofluoromethane	5.391	113	73632	45.519	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	91.040%		
50) Toluene-d8	8.647	98	271932	46.700	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery =	93.400%		
62) 4-Bromofluorobenzene	11.079	95	96373	45.198	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	90.400%		
Target Compounds						Qvalue
30) Chloroform	5.099	83	8499	2.948	ug/l	100
47) Bromodichloromethane	7.824	83	7544	3.599	ug/l	92
60) Dibromochloromethane	9.525	129	6245	3.839	ug/l	99
71) Bromoform	10.799	173	1694	3.436	ug/l #	100

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

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