

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX102122\
 Data File : VX032299.D
 Acq On : 21 Oct 2022 19:16
 Operator : JC/MD
 Sample : N5248-06
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 FSND-MW11R-20221018

Quant Time: Oct 22 05:14:18 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	145879	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	223583	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	194217	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	89624	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	90724	47.806	ug/l	0.00
Spiked Amount	50.000	Range 78 - 117	Recovery	=	95.620%	
35) Dibromofluoromethane	5.391	113	70825	45.522	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	=	91.040%	
50) Toluene-d8	8.647	98	261213	46.641	ug/l	0.00
Spiked Amount	50.000	Range 92 - 112	Recovery	=	93.280%	
62) 4-Bromofluorobenzene	11.079	95	92262	44.988	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	=	89.980%	
Target Compounds						
					Qvalue	
16) Acetone	2.386	43	1238	2.010	ug/l #	86
30) Chloroform	5.099	83	2688	0.954	ug/l	97
32) 1,1,1-Trichloroethane	5.373	97	762	0.317	ug/l #	48
38) Carbon Tetrachloride	5.684	117	10336	4.892	ug/l	94
44) Trichloroethene	7.129	130	654	0.409	ug/l	96

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

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