

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX102522\
 Data File : VX032372.D
 Acq On : 25 Oct 2022 18:19
 Operator : JC/MD
 Sample : N5249-10
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 FSND-RB-20221019

Quant Time: Oct 27 14:18:21 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	124378	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	199164	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	169303	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	77249	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	79505	49.137	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	98.280%
35) Dibromofluoromethane	5.391	113	63985	46.168	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	92.340%
50) Toluene-d8	8.653	98	232543	46.612	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	93.220%
62) 4-Bromofluorobenzene	11.079	95	77626	42.492	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	84.980%
Target Compounds						
					Qvalue	
16) Acetone	2.380	43	2057	3.917	ug/l	97
40) Benzene	6.043	78	1850	0.358	ug/l	99
52) Toluene	8.720	92	8502	2.540	ug/l	98
67) Ethyl Benzene	10.195	91	1502	0.244	ug/l #	82
68) m/p-Xylenes	10.305	106	2104	0.867	ug/l	97
69) o-Xylene	10.646	106	827	0.356	ug/l	93

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

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