

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX102522\
 Data File : VX032376.D
 Acq On : 25 Oct 2022 19:53
 Operator : JC/MD
 Sample : N5249-14
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 FSND-RB-20221020

Quant Time: Oct 27 14:18:53 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	124452	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	197006	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	171734	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	77293	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	79582	49.155	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	98.320%
35) Dibromofluoromethane	5.385	113	63997	46.682	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	93.360%
50) Toluene-d8	8.653	98	233564	47.330	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	94.660%
62) 4-Bromofluorobenzene	11.079	95	78459	43.418	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	86.840%
Target Compounds						
					Qvalue	
16) Acetone	2.385	43	2357	4.486	ug/l	89
40) Benzene	6.043	78	1743	0.341	ug/l	95
52) Toluene	8.720	92	8045	2.430	ug/l	96
68) m/p-Xylenes	10.305	106	2158	0.877	ug/l	100
69) o-Xylene	10.646	106	674	0.286	ug/l	80

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

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