

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX102022\
 Data File : VX032257.D
 Acq On : 20 Oct 2022 13:00
 Operator : JC/MD
 Sample : N5149-04
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 BP-VPB-RW5-GW-918-920

Quant Time: Oct 21 01:23:20 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X101922W.M
 Quant Title : SW846 8260
 QLast Update : Wed Oct 19 16:51:09 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	159003	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	243853	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	206906	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	93485	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	100752	48.709	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	97.420%
35) Dibromofluoromethane	5.391	113	76249	44.934	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	89.860%
50) Toluene-d8	8.647	98	293912	48.117	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	96.240%
62) 4-Bromofluorobenzene	11.079	95	103341	46.201	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	92.400%
Target Compounds						
					Qvalue	
16) Acetone	2.380	43	2407	3.586	ug/l	90
17) Carbon Disulfide	2.514	76	1284	0.386	ug/l	97

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

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