

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX102422\
 Data File : VX032348.D
 Acq On : 24 Oct 2022 18:17
 Operator : JC/MD
 Sample : N5260-03
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 EB-2022-10-21

Quant Time: Oct 25 03:58:50 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	133112	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	210962	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	182822	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	84914	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	83801	48.394	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery =	96.780%		
35) Dibromofluoromethane	5.385	113	66682	45.423	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	90.840%		
50) Toluene-d8	8.653	98	246771	46.698	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery =	93.400%		
62) 4-Bromofluorobenzene	11.079	95	85170	44.014	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	88.020%		
Target Compounds						Qvalue
16) Acetone	2.386	43	2329	4.144	ug/l	98
67) Ethyl Benzene	10.195	91	5949	0.896	ug/l	99
68) m/p-Xylenes	10.305	106	2911	1.111	ug/l	95
78) n-propylbenzene	11.305	91	2061	0.297	ug/l	93
84) 1,2,4-Trimethylbenzene	11.750	105	6148	1.208	ug/l	99

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

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