

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN101421\
 Data File : VN068930.D
 Acq On : 14 Oct 2021 17:31
 Operator : JC/MD
 Sample : M4192-18
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 HAR-18

Quant Time: Oct 14 22:48:05 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N101221W.M
 Quant Title : SW846 8260
 QLast Update : Wed Oct 13 05:44:40 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.088	168	378602	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.971	114	617345	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.746	117	570594	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.677	152	228753	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	224471	44.969	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery = 89.940%			
35) Dibromofluoromethane	8.024	113	185439	51.629	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery = 103.260%			
50) Toluene-d8	10.446	98	755212	52.550	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery = 105.100%			
62) 4-Bromofluorobenzene	12.733	95	259672	47.891	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery = 95.780%			
Target Compounds						Qvalue
16) Acetone	4.299	43	20564	11.039	ug/l	90
18) Methyl Acetate	4.864	43	11877	1.681	ug/l	92
20) Methylene Chloride	5.103	84	23615	5.508	ug/l	95
25) 2-Butanone	7.348	43	5168	1.835	ug/l	97
43) Isopropyl Acetate	8.571	43	73575	7.657	ug/l #	76

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

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