

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN102721\
 Data File : VN069206.D
 Acq On : 27 Oct 2021 00:44
 Operator : JC/MD
 Sample : M4337-18
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-1D(OFFSITE)

Quant Time: Oct 27 09:26:00 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	329072	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.971	114	574765	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.746	117	554376	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.677	152	206105	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	217299	50.084	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	100.160%
35) Dibromofluoromethane	8.024	113	168843	50.491	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	100.980%
50) Toluene-d8	10.443	98	718851	53.726	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	107.460%
62) 4-Bromofluorobenzene	12.736	95	257498	51.009	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	102.020%
Target Compounds						
					Qvalue	
16) Acetone	4.309	43	3991	2.465	ug/l #	84
27) cis-1,2-Dichloroethene	7.329	96	52554	12.656	ug/l	88
30) Chloroform	7.820	83	1767	0.260	ug/l #	83
44) Trichloroethene	9.217	130	14954	3.997	ug/l	98
64) Tetrachloroethene	10.979	164	39019	10.926	ug/l	98

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

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