

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN031822\
 Data File : VN071372.D
 Acq On : 18 Mar 2022 15:06
 Operator : JC\MD
 Sample : N1960-18
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TB-20220316

Quant Time: Mar 19 07:07:36 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N030722W.M
 Quant Title : SW846 8260
 QLast Update : Tue Mar 08 06:52:46 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.086	168	771668	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.963	114	1295437	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.739	117	1235505	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.674	152	427447	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.434	65	501906	47.457	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	94.920%
35) Dibromofluoromethane	8.016	113	383354	46.895	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	93.800%
50) Toluene-d8	10.439	98	1590871	48.335	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	96.660%
62) 4-Bromofluorobenzene	12.727	95	578169	51.875	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	103.760%
Target Compounds						
					Qvalue	
3) Chloromethane	2.299	50	3768	0.297	ug/l #	79
16) Acetone	4.304	43	8867	1.723	ug/l #	83
18) Methyl Acetate	4.863	43	9344	1.140	ug/l	92

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

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