

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN022522\
 Data File : VN071076.D
 Acq On : 25 Feb 2022 17:46
 Operator : JC\MD
 Sample : N1664-11
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TP-4

Quant Time: Feb 26 03:43:59 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N021522W.M
 Quant Title : SW846 8260
 QLast Update : Mon Feb 21 07:57:15 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.086	168	850267	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.963	114	1374088	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.739	117	1201686	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.674	152	421695	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.433	65	508889	52.394	ug/l	0.00
Spiked Amount	50.000	Range 61 - 141	Recovery	= 104.780%		
35) Dibromofluoromethane	8.016	113	391699	47.926	ug/l	0.00
Spiked Amount	50.000	Range 69 - 133	Recovery	= 95.860%		
50) Toluene-d8	10.439	98	1664196	50.703	ug/l	0.00
Spiked Amount	50.000	Range 65 - 126	Recovery	= 101.400%		
62) 4-Bromofluorobenzene	12.727	95	564812	48.142	ug/l	0.00
Spiked Amount	50.000	Range 58 - 135	Recovery	= 96.280%		
Target Compounds						
					Qvalue	
16) Acetone	4.293	43	51519	9.631	ug/l	94
18) Methyl Acetate	4.863	43	19207	1.050	ug/l	96
20) Methylene Chloride	5.098	84	110871	11.405	ug/l	93
43) Isopropyl Acetate	8.557	43	129195	5.708	ug/l #	82
93) 1,2,4-Trichlorobenzene	15.268	180	1367	2.168	ug/l	97
96) 1,2,3-Trichlorobenzene	15.709	180	1454	2.709	ug/l	95

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

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