

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN031921\
 Data File : VN066279.D
 Acq On : 19 Mar 2021 14:43
 Operator : JC/MD
 Sample : M1693-01
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TP-5A

Manual Integrations
 APPROVED

MMDadoda
 3/22/2021 11:50:31 AM

Quant Time: Mar 20 04:58:18 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N031621W.M
 Quant Title : SW846 8260
 QLast Update : Wed Mar 17 05:05:15 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	7.587	168	321908	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.515	114	481200	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.350	117	451645	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.286	152	146456	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.949	65	207168	44.67	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery =	89.34%		
35) Dibromofluoromethane	7.509	113	169705	55.28	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	110.56%		
50) Toluene-d8	10.028	98	666015	56.11	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery =	112.22%#		
62) 4-Bromofluorobenzene	12.345	95	179241	41.89	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	83.78%		
Target Compounds					Qvalue	
16) Acetone	3.738	43	78491	35.23	ug/l	95
20) Methylene Chloride	4.449	84	16682m	4.79	ug/l	
43) Isopropyl Acetate	8.097	43	28563	2.65	ug/l	97
65) Chlorobenzene	11.371	112	11672	1.29	ug/l	99

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

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