

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN022421\
 Data File : VN066004.D
 Acq On : 24 Feb 2021 15:18
 Operator : JC/MD
 Sample : M1473-05
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 GWBR1-132-149-022321

Manual Integrations
 APPROVED

MMDadoda
 2/25/2021 12:15:57 PM

Quant Time: Feb 25 02:14:18 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N021721W.M
 Quant Title : SW846 8260
 QLast Update : Wed Feb 17 15:48:13 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	7.631	168	125493	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.550	114	207386	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.380	117	187066	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.315	152	75224	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.991	65	82761	47.16	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery =	94.32%		
35) Dibromofluoromethane	7.553	113	61795	48.15	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery =	96.30%		
50) Toluene-d8	10.058	98	250376	47.52	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery =	95.04%		
62) 4-Bromofluorobenzene	12.373	95	89993	45.50	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery =	91.00%		
Target Compounds					Qvalue	
12) 1,1-Dichloroethene	3.714	96	1441m	1.02	ug/l	
16) Acetone	3.795	43	3137	4.11	ug/l	99
27) cis-1,2-Dichloroethene	6.788	96	6672	3.95	ug/l	84
30) Chloroform	7.335	83	10383	3.78	ug/l	93
44) Trichloroethene	8.801	130	19093	12.30	ug/l	99
52) Toluene	10.122	92	33416	8.67	ug/l	94

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

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