

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN041224\
 Data File : VN081721.D
 Acq On : 12 Apr 2024 16:31
 Operator : JC\MD
 Sample : P2070-02
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TP-1

Quant Time: Apr 13 06:58:30 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N031824W.M
 Quant Title : SW846 8260
 QLast Update : Tue Mar 19 06:00:00 2024
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By : John Carlone 04/15/2024
 Supervised By : Mahesh Dadoda 04/15/2024

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

Internal Standards						
1) Pentafluorobenzene	8.224	168	258518	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	496924	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	510425	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	193074	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	198112	60.092	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 120.180%			
35) Dibromofluoromethane	8.165	113	160283	51.109	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 102.220%			
50) Toluene-d8	10.571	98	643719	56.782	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 113.560%#			
62) 4-Bromofluorobenzene	12.853	95	202091	50.591	ug/l	0.00
Spiked Amount 50.000	Range 64 - 133		Recovery = 101.180%			
Target Compounds						
					Qvalue	
16) Acetone	4.430	43	40188	37.720	ug/l	89
20) Methylene Chloride	5.283	84	8697	2.492	ug/l #	71
37) Ethyl Acetate	7.565	43	4690	1.038	ug/l #	87
43) Isopropyl Acetate	8.688	43	282877m	37.762	ug/l	

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

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