

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN051623\
 Data File : VN077691.D
 Acq On : 16 May 2023 14:39
 Operator : JC\MD
 Sample : 02804-02
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 OWBR-03-128-148-051523

Quant Time: May 17 04:01:04 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N051523W.M
 Quant Title : SW846 8260
 QLast Update : Tue May 16 04:07:42 2023
 Response via : Initial Calibration

Manual Integrations
 APPROVED

Reviewed By : John Carlone 05/17/2023
 Supervised By : Mahesh Dadoda 05/17/2023

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

Internal Standards						
1) Pentafluorobenzene	8.230	168	366692	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	663829	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	593952	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	236696	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	258732	47.157	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery =	94.320%		
35) Dibromofluoromethane	8.171	113	209851	48.478	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	96.960%		
50) Toluene-d8	10.571	98	828319	49.714	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery =	99.420%		
62) 4-Bromofluorobenzene	12.853	95	272814	46.361	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	92.720%		
Target Compounds						
16) Acetone	4.448	43	8213m	4.758	ug/l	
25) 2-Butanone	7.500	43	9881	3.659	ug/l	99
30) Chloroform	7.965	83	3265	0.365	ug/l #	74
40) Benzene	8.618	78	47142	2.523	ug/l	97
67) Ethyl Benzene	11.965	91	6490	0.281	ug/l #	86

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

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