

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN110722\
 Data File : VN075195.D
 Acq On : 07 Nov 2022 18:10
 Operator : JC\MD
 Sample : N5491-02
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TP-1

Quant Time: Nov 08 01:12:36 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N102622W.M
 Quant Title : SW846 8260
 QLast Update : Thu Oct 27 09:48:49 2022
 Response via : Initial Calibration

Manual Integrations
 APPROVED

Reviewed By : John Carlone 11/08/2022
 Supervised By : Mahesh Dadoda 11/08/2022

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

Internal Standards						
1) Pentafluorobenzene	8.230	168	97818	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	199481	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	185335	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	72126	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.582	65	77989	59.173	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 118.340%			
35) Dibromofluoromethane	8.171	113	62333	53.228	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 106.460%			
50) Toluene-d8	10.571	98	241604	53.177	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 106.360%			
62) 4-Bromofluorobenzene	12.853	95	76917	48.194	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery = 96.380%			
Target Compounds						
16) Acetone	4.453	43	9786	17.536	ug/l	95
18) Methyl Acetate	5.047	43	4479	2.457	ug/l	95
20) Methylene Chloride	5.277	84	4678	3.760	ug/l #	89
37) Ethyl Acetate	7.577	43	17655	8.880	ug/l	99
43) Isopropyl Acetate	8.694	43	110517m	35.835	ug/l	

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

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