

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091724\
 Data File : VN083943.D
 Acq On : 17 Sep 2024 14:17
 Operator : JC\MD
 Sample : P3989-04
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TP-2

Quant Time: Sep 18 01:48:52 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N091024W.M
 Quant Title : SW846 8260
 QLast Update : Wed Sep 11 15:18:56 2024
 Response via : Initial Calibration

Manual Integrations
 APPROVED

Reviewed By : John Carlone 09/18/2024
 Supervised By : Mahesh Dadoda 09/18/2024

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

Internal Standards						
1) Pentafluorobenzene	8.224	168	155639	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.100	114	274380	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	238290	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	96984	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.577	65	123028	52.423	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 104.840%			
35) Dibromofluoromethane	8.171	113	88101	50.156	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 100.320%			
50) Toluene-d8	10.565	98	331842	54.647	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 109.300%			
62) 4-Bromofluorobenzene	12.847	95	117589	46.935	ug/l	0.00
Spiked Amount 50.000	Range 77 - 121		Recovery = 93.880%			
Target Compounds						
16) Acetone	4.424	43	15757	16.749	ug/l	95
20) Methylene Chloride	5.277	84	6970	3.620	ug/l	97
43) Isopropyl Acetate	8.688	43	183364m	42.164	ug/l	

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

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