

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN060524\
 Data File : VN082474.D
 Acq On : 05 Jun 2024 16:38
 Operator : JC\MD
 Sample : P2715-08
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 WCS-TPM

Manual Integrations
 APPROVED

Reviewed By : John Carlone 06/06/2024
 Supervised By : Mahesh Dadoda 06/06/2024

Quant Time: Jun 06 00:59:09 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N051524W.M
 Quant Title : SW846 8260
 QLast Update : Thu May 16 05:21:06 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.230	168	152061	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	287878	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.870	117	287926	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	119784	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.582	65	134551	56.038	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 112.080%			
35) Dibromofluoromethane	8.171	113	98427	55.580	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 111.160%			
50) Toluene-d8	10.571	98	380142	55.990	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 111.980%			
62) 4-Bromofluorobenzene	12.853	95	125746	51.890	ug/l	0.00
Spiked Amount 50.000	Range 64 - 133		Recovery = 103.780%			
Target Compounds						
16) Acetone	4.430	43	100050	101.577	ug/l #	88
20) Methylene Chloride	5.277	84	5566	2.801	ug/l #	66
25) 2-Butanone	7.494	43	24703	16.542	ug/l #	87
43) Isopropyl Acetate	8.694	43	281178	47.741	ug/l #	71
51) 4-Methyl-2-Pentanone	10.447	43	23630m	7.061	ug/l	

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

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