

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN013124\
 Data File : VN080843.D
 Acq On : 31 Jan 2024 16:20
 Operator : JC\MD
 Sample : P1276-04
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TP-37

Manual Integrations
 APPROVED

Reviewed By : John Carlone 02/01/2024
 Supervised By : Mahesh Dadoda 02/01/2024

Quant Time: Jan 31 23:18:52 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N012924W.M
 Quant Title : SW846 8260
 QLast Update : Mon Jan 29 12:54:23 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.224	168	94914	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.100	114	221298	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	231745	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	91595	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.577	65	82116	50.010	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 100.020%			
35) Dibromofluoromethane	8.165	113	60119	44.404	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 88.800%			
50) Toluene-d8	10.571	98	275035	51.902	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 103.800%			
62) 4-Bromofluorobenzene	12.853	95	98506	53.345	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery = 106.680%			
Target Compounds						
					Qvalue	
3) Chloromethane	2.354	50	2161	1.020	ug/l #	85
16) Acetone	4.430	43	26761	44.938	ug/l #	78
20) Methylene Chloride	5.277	84	3763m	2.407	ug/l	
25) 2-Butanone	7.483	43	3933	3.477	ug/l	91
37) Ethyl Acetate	7.565	43	3192	1.232	ug/l #	77
43) Isopropyl Acetate	8.694	43	98648	20.916	ug/l #	90

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

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