

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061623\
 Data File : VN078265.D
 Acq On : 16 Jun 2023 17:17
 Operator : JC\MD
 Sample : 03262-08
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 WS-SOIL3-0615

Quant Time: Jun 17 02:26:33 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N060623W.M
 Quant Title : SW846 8260
 QLast Update : Wed Jun 07 05:05:00 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.235	168	445838	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	850077	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.870	117	803693	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	258184	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	326731	54.965	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	= 109.920%		
35) Dibromofluoromethane	8.177	113	292150	54.354	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 108.700%		
50) Toluene-d8	10.570	98	1040197	50.200	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	= 100.400%		
62) 4-Bromofluorobenzene	12.853	95	329660	51.177	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 102.360%		
Target Compounds						
					Qvalue	
16) Acetone	4.441	43	181338	105.983	ug/l	96
20) Methylene Chloride	5.288	84	21727	3.843	ug/l #	86
25) 2-Butanone	7.494	43	65562	24.917	ug/l #	83
37) Ethyl Acetate	7.577	43	68981	11.453	ug/l #	86
43) Isopropyl Acetate	8.700	43	353733	30.710	ug/l #	79
51) 4-Methyl-2-Pentanone	10.453	43	433496	71.470	ug/l	90

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

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