

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN082223\
 Data File : VN078998.D
 Acq On : 22 Aug 2023 16:48
 Operator : JC\MD
 Sample : 04090-05
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TP-K

Quant Time: Aug 23 03:05:15 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N081523W.M
 Quant Title : SW846 8260
 QLast Update : Wed Aug 16 02:42:02 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.230	168	176958	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	354877	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	355139	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	108636	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	148230	53.831	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 107.660%	
35) Dibromofluoromethane	8.171	113	121651	50.925	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 101.840%	
50) Toluene-d8	10.571	98	444831	50.150	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 100.300%	
62) 4-Bromofluorobenzene	12.853	95	142327	49.091	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 98.180%	
Target Compounds						
					Qvalue	
16) Acetone	4.442	43	26591	23.404	ug/l	92
20) Methylene Chloride	5.283	84	13308	5.496	ug/l	90
25) 2-Butanone	7.500	43	9437	5.181	ug/l #	66
43) Isopropyl Acetate	8.694	43	223717	31.482	ug/l #	69
51) 4-Methyl-2-Pentanone	10.453	43	37959	9.006	ug/l #	78

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

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