

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN021224\
 Data File : VN081011.D
 Acq On : 12 Feb 2024 18:21
 Operator : JC\MD
 Sample : P1449-02
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 ETGI-362

Quant Time: Feb 13 03:24:16 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N020624W.M
 Quant Title : SW846 8260
 QLast Update : Wed Feb 07 06:22:23 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.224	168	588477	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	9.100	114	1020764	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	1011655	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	371241	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	291196	39.225	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	78.460%
35) Dibromofluoromethane	8.171	113	276847	50.543	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	101.080%
50) Toluene-d8	10.571	98	920693	46.471	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	92.940%
62) 4-Bromofluorobenzene	12.853	95	330037	45.855	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	91.720%
Target Compounds						
					Qvalue	
3) Chloromethane	2.365	50	20176	1.585	ug/l	100
13) Acrolein	4.177	56	8895	6.900	ug/l	# 81
16) Acetone	4.430	43	111044	45.466	ug/l	100
20) Methylene Chloride	5.271	84	22665	3.063	ug/l	# 82
37) Ethyl Acetate	7.553	43	12296	1.994	ug/l	# 83
43) Isopropyl Acetate	8.694	43	455929	40.212	ug/l	# 87

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

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