

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN101223\
 Data File : VN079552.D
 Acq On : 12 Oct 2023 17:59
 Operator : JC\MD
 Sample : PB156319ZHE#03
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB156319ZHE#03

Quant Time: Oct 13 01:23:36 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N100523W.M
 Quant Title : SW846 8260
 QLast Update : Fri Oct 06 03:34:49 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.230	168	175020	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	372641	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	379013	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	103856	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	158585	55.883	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	=	111.760%	
35) Dibromofluoromethane	8.171	113	131020	48.926	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	=	97.860%	
50) Toluene-d8	10.571	98	502944	51.906	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	=	103.820%	
62) 4-Bromofluorobenzene	12.853	95	144591	46.499	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	=	93.000%	
Target Compounds						
					Qvalue	
16) Acetone	4.442	43	30349	26.963	ug/l	88
20) Methylene Chloride	5.277	84	19804	7.523	ug/l #	81
43) Isopropyl Acetate	8.694	43	316853	45.223	ug/l #	69
51) 4-Methyl-2-Pentanone	10.447	43	21257	5.374	ug/l #	73

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

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