

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN081823\
 Data File : VN078953.D
 Acq On : 18 Aug 2023 19:36
 Operator : JC\MD
 Sample : PB154902ZHE#10
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB154902ZHE#10

Quant Time: Aug 18 23:48:04 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	8.224	168	149778	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	293103	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.870	117	289928	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	87755	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.582	65	119835	51.416	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 102.840%	
35) Dibromofluoromethane	8.177	113	101138	51.261	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 102.520%	
50) Toluene-d8	10.570	98	382946	52.272	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 104.540%	
62) 4-Bromofluorobenzene	12.853	95	113151	47.253	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 94.500%	
Target Compounds						
					Qvalue	
16) Acetone	4.441	43	27465	28.560	ug/l	100
20) Methylene Chloride	5.283	84	11796	5.756	ug/l #	83
25) 2-Butanone	7.494	43	11031	7.155	ug/l #	68
43) Isopropyl Acetate	8.694	43	266196	45.354	ug/l #	69
51) 4-Methyl-2-Pentanone	10.453	43	47482	13.640	ug/l #	78

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

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