

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN100623\
 Data File : VN079446.D
 Acq On : 06 Oct 2023 14:42
 Operator : JC\MD
 Sample : PB156147ZHE#01
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB156147ZHE#01

Quant Time: Oct 06 23:03:50 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	151143	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	309120	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	309370	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	77006	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	133327	54.404	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	108.800%
35) Dibromofluoromethane	8.171	113	110291	49.648	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	99.300%
50) Toluene-d8	10.571	98	414757	51.600	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	103.200%
62) 4-Bromofluorobenzene	12.853	95	109257	42.356	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	84.720%
Target Compounds						
					Qvalue	
16) Acetone	4.442	43	32942	33.890	ug/l	94
20) Methylene Chloride	5.283	84	44598	20.598	ug/l #	82
43) Isopropyl Acetate	8.694	43	204659	35.212	ug/l #	75
51) 4-Methyl-2-Pentanone	10.453	43	16829	5.129	ug/l #	60

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

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