

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN011222\
 Data File : VN070594.D
 Acq On : 12 Jan 2022 18:33
 Operator : JC/MD
 Sample : PB141975ZHE#01
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB141975ZHE#01

Quant Time: Jan 13 04:28:06 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
 Quant Title : SW846 8260
 QLast Update : Mon Dec 27 18:47:12 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.086	168	791602	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.965	114	1391412	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.744	117	1238608	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.675	152	429855	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.437	65	660040	57.991	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	115.980%
35) Dibromofluoromethane	8.021	113	466704	55.436	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	110.880%
50) Toluene-d8	10.440	98	1916734	55.733	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	111.460%
62) 4-Bromofluorobenzene	12.731	95	692335	55.737	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	111.480%
Target Compounds						
						Qvalue
16) Acetone	4.291	43	91098	12.330	ug/l	96
18) Methyl Acetate	4.862	43	49316	2.307	ug/l	91
20) Methylene Chloride	5.090	84	57427	5.618	ug/l	87
43) Isopropyl Acetate	8.563	43	191651	6.734	ug/l #	80
80) 1,3,5-Trimethylbenzene	13.061	105	9984	0.312	ug/l	99

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

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