

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU100121\
 Data File : VU045072.D
 Acq On : 01 Oct 2021 14:52
 Operator : SY/MD
 Sample : PB139508ZHE#03
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 PB139508ZHE#03

Quant Time: Oct 02 03:11:34 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U092821W.M
 Quant Title : SW846 8260
 QLast Update : Tue Sep 28 12:32:13 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.375	168	201523	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.253	114	377457	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.420	117	360933	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.816	152	155033	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.703	65	169689	50.584	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	101.160%
35) Dibromofluoromethane	5.292	113	116151	48.780	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	97.560%
50) Toluene-d8	7.903	98	476308	50.627	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	101.260%
62) 4-Bromofluorobenzene	10.639	95	177919	45.806	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	91.620%
Target Compounds						
16) Acetone	2.642	43	33688	16.918	ug/l	99
18) Methyl Acetate	2.961	43	8597	1.477	ug/l	98
20) Methylene Chloride	3.047	84	8667	2.890	ug/l	93
43) Isopropyl Acetate	5.919	43	26446	3.366	ug/l	100

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

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