

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU100121\
 Data File : VU045073.D
 Acq On : 01 Oct 2021 15:15
 Operator : SY/MD
 Sample : PB139508ZHE#04
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 PB139508ZHE#04

Quant Time: Oct 02 03:11:48 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	202254	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.250	114	376633	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.423	117	360080	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.816	152	156026	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.706	65	167684	49.806	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery =	99.620%		
35) Dibromofluoromethane	5.292	113	113605	47.815	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	95.640%		
50) Toluene-d8	7.902	98	474052	50.497	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery =	101.000%		
62) 4-Bromofluorobenzene	10.639	95	175954	45.399	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	90.800%		
Target Compounds					Qvalue	
16) Acetone	2.642	43	34451	17.239	ug/l	99
18) Methyl Acetate	2.961	43	9067	1.552	ug/l	100
20) Methylene Chloride	3.044	84	9264	3.078	ug/l	96
43) Isopropyl Acetate	5.922	43	26647	3.399	ug/l	99

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

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