

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092721\
 Data File : VU044923.D
 Acq On : 28 Sep 2021 05:48
 Operator : SY/MD
 Sample : M3935-04
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 TP-1-S

Quant Time: Sep 28 06:32:38 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U092421W.M
 Quant Title : SW846 8260
 QLast Update : Sat Sep 25 09:30:46 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.375	168	193200	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.253	114	344093	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.420	117	334351	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.816	152	161563	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.703	65	157108	46.594	ug/l	0.00
Spiked Amount	50.000		Recovery	=	93.180%	
35) Dibromofluoromethane	5.292	113	107375	45.488	ug/l	0.00
Spiked Amount	50.000		Recovery	=	90.980%	
50) Toluene-d8	7.899	98	444337	49.840	ug/l	0.00
Spiked Amount	50.000		Recovery	=	99.680%	
62) 4-Bromofluorobenzene	10.639	95	178721	53.973	ug/l	0.00
Spiked Amount	50.000		Recovery	=	107.940%	
Target Compounds						
					Qvalue	
16) Acetone	2.629	43	100072	52.615	ug/l	99
20) Methylene Chloride	3.044	84	16225	5.270	ug/l	97
25) 2-Butanone	4.713	43	15152	5.250	ug/l	97
43) Isopropyl Acetate	5.916	43	22936	3.060	ug/l	99

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

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