

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU061021\
 Data File : VU044163.D
 Acq On : 10 Jun 2021 18:08
 Operator : SY/MD
 Sample : PB136963TB
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 PB136963TB

Quant Time: Jun 11 01:21:38 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U060921W.M
 Quant Title : SW846 8260
 QLast Update : Thu Jun 10 05:11:37 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.382	168	145634	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.256	114	296045	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.423	117	312696	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.819	152	148734	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.710	65	126143	54.467	ug/l	0.00
Spiked Amount 50.000			Recovery	=	108.940%	
35) Dibromofluoromethane	5.298	113	100673	48.885	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.780%	
50) Toluene-d8	7.906	98	401966	50.824	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.640%	
62) 4-Bromofluorobenzene	10.639	95	151121	47.680	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.360%	
Target Compounds						
					Qvalue	
16) Acetone	2.633	43	17513	12.612	ug/l	100
20) Methylene Chloride	3.054	84	15515	6.388	ug/l	99
43) Isopropyl Acetate	5.919	43	16993	3.037	ug/l	98
51) 4-Methyl-2-Pentanone	7.800	43	5093	1.257	ug/l	93

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

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