

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU010919\
 Data File : VU029097.D
 Acq On : 09 Jan 2019 18:15
 Operator : MD/SY
 Sample : PB116248ZHE#12
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 PB116248ZHE#12

Quant Time: Jan 10 05:42:06 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U010319W.M
 Quant Title : SW846 8260
 QLast Update : Fri Jan 04 02:22:53 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	4.98	168	110232	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.89	114	189701	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	179692	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.48	152	66249	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	77386	56.79	ug/l	0.00
Spiked Amount 50.000			Recovery	=	113.58%	
35) Dibromofluoromethane	4.88	113	66282	53.90	ug/l	0.00
Spiked Amount 50.000			Recovery	=	107.80%	
50) Toluene-d8	7.56	98	250103	51.52	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.04%	
62) 4-Bromofluorobenzene	10.30	95	74620	42.77	ug/l	0.00
Spiked Amount 50.000			Recovery	=	85.54%	
Target Compounds						
16) Acetone	2.32	43	5391	6.280	ug/l	97
18) Methyl Acetate	2.63	43	2036	1.020	ug/l #	78
20) Methylene Chloride	2.70	84	1873	1.285	ug/l	93
43) Isopropyl Acetate	5.55	43	8227	2.577	ug/l	93

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

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