

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU010919\
 Data File : VU029101.D
 Acq On : 09 Jan 2019 19:51
 Operator : MD/SY
 Sample : PB116248ZHE#16
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 PB116248ZHE#16

Quant Time: Jan 10 05:48:35 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	107085	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.88	114	185440	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	180174	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.48	152	70634	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	74601	56.35	ug/l	0.00
Spiked Amount 50.000			Recovery	=	112.70%	
35) Dibromofluoromethane	4.88	113	64396	53.57	ug/l	0.00
Spiked Amount 50.000			Recovery	=	107.14%	
50) Toluene-d8	7.56	98	246944	52.04	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.08%	
62) 4-Bromofluorobenzene	10.30	95	76178	44.67	ug/l	0.00
Spiked Amount 50.000			Recovery	=	89.34%	
Target Compounds						
16) Acetone	2.32	43	5561	6.668	ug/l	98
18) Methyl Acetate	2.63	43	2106	1.086	ug/l	93
20) Methylene Chloride	2.70	84	2213	1.563	ug/l	99
43) Isopropyl Acetate	5.55	43	7937	2.543	ug/l	95

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

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