

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU010919\
 Data File : VU029091.D
 Acq On : 09 Jan 2019 15:49
 Operator : MD/SY
 Sample : PB116248ZHE#06
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 PB116248ZHE#06

Quant Time: Jan 10 05:33:02 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	109986	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.88	114	189644	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	177951	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.48	152	66674	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	75174	55.29	ug/l	0.00
Spiked Amount 50.000			Recovery	=	110.58%	
35) Dibromofluoromethane	4.88	113	64258	52.27	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.54%	
50) Toluene-d8	7.57	98	247763	51.05	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.10%	
62) 4-Bromofluorobenzene	10.31	95	76546	43.89	ug/l	0.00
Spiked Amount 50.000			Recovery	=	87.78%	
Target Compounds						
16) Acetone	2.32	43	5857	6.838	ug/l	96
18) Methyl Acetate	2.63	43	2279	1.144	ug/l #	81
20) Methylene Chloride	2.70	84	2145	1.475	ug/l	96
43) Isopropyl Acetate	5.55	43	9127	2.860	ug/l	97

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

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