

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU010919\
 Data File : VU029090.D
 Acq On : 09 Jan 2019 15:25
 Operator : MD/SY
 Sample : PB116248ZHE#05
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 PB116248ZHE#05

Quant Time: Jan 10 05:01:39 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	112012	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.88	114	191011	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	181377	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.48	152	69918	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	75484	54.51	ug/l	0.00
Spiked Amount 50.000			Recovery	=	109.02%	
35) Dibromofluoromethane	4.88	113	66369	53.60	ug/l	0.00
Spiked Amount 50.000			Recovery	=	107.20%	
50) Toluene-d8	7.57	98	246137	50.35	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.70%	
62) 4-Bromofluorobenzene	10.31	95	75872	43.19	ug/l	0.00
Spiked Amount 50.000			Recovery	=	86.38%	
Target Compounds						
16) Acetone	2.32	43	5762	6.606	ug/l	99
18) Methyl Acetate	2.62	43	2148	1.059	ug/l #	75
20) Methylene Chloride	2.70	84	2284	1.542	ug/l	94
43) Isopropyl Acetate	5.55	43	9212	2.866	ug/l	97

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

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