

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU010919\
 Data File : VU029098.D
 Acq On : 09 Jan 2019 18:39
 Operator : MD/SY
 Sample : PB116248ZHE#13
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 PB116248ZHE#13

Quant Time: Jan 10 05:43:44 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	107447	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.88	114	187921	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	178145	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.48	152	69499	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	73958	55.68	ug/l	0.00
Spiked Amount 50.000			Recovery	=	111.36%	
35) Dibromofluoromethane	4.88	113	64426	52.89	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.78%	
50) Toluene-d8	7.56	98	244935	50.93	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.86%	
62) 4-Bromofluorobenzene	10.30	95	75783	43.85	ug/l	0.00
Spiked Amount 50.000			Recovery	=	87.70%	
Target Compounds						
16) Acetone	2.32	43	5483	6.553	ug/l	97
18) Methyl Acetate	2.62	43	2050	1.054	ug/l #	88
20) Methylene Chloride	2.70	84	1915	1.348	ug/l	88
43) Isopropyl Acetate	5.55	43	8170	2.583	ug/l #	89

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

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