

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU083019\
 Data File : VU034079.D
 Acq On : 29 Aug 2019 21:52
 Operator : JC/SP
 Sample : K4609-01
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 FERMENTATION-WASTE

Quant Time: Aug 30 05:36:35 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U082919W.M
 Quant Title : SW846 8260
 QLast Update : Thu Aug 29 04:11:12 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	4.96	168	469297	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.87	114	729968	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.07	117	700811	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.47	152	345186	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.29	65	292506	48.64	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.28%	
35) Dibromofluoromethane	4.86	113	242510	47.90	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.80%	
50) Toluene-d8	7.55	98	922317	48.65	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.30%	
62) 4-Bromofluorobenzene	10.29	95	324421	43.99	ug/l	0.00
Spiked Amount 50.000			Recovery	=	87.98%	
Target Compounds						
16) Acetone	2.31	43	42480	13.845	ug/l	99
17) Carbon Disulfide	2.47	76	34288	2.306	ug/l	97
25) 2-Butanone	4.26	43	25016	5.860	ug/l	100
29) Tetrahydrofuran	4.62	42	200348	76.105	ug/l	97
52) Toluene	7.62	92	44929	3.491	ug/l	100

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

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