

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\U101018\
 Data File : VU027564.D
 Acq On : 10 Oct 2018 12:53
 Operator : MD/SY
 Sample : J5324-03
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 OILY-WATER

Quant Time: Oct 11 05:06:10 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U100118W.M
 Quant Title : SW846 8260
 QLast Update : Tue Oct 02 02:04:34 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	4.99	168	129508	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.89	114	212184	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	185061	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.48	152	85643	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	98865	45.72	ug/l	0.00
Spiked Amount 50.000			Recovery	=	91.44%	
35) Dibromofluoromethane	4.89	113	71206	48.97	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.94%	
50) Toluene-d8	7.57	98	267860	49.22	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.44%	
62) 4-Bromofluorobenzene	10.31	95	89838	44.42	ug/l	0.00
Spiked Amount 50.000			Recovery	=	88.84%	
Target Compounds						
					Qvalue	
11) Tert butyl alcohol	2.83	59	2975	6.533	ug/l #	80
16) Acetone	2.32	43	217558	170.292	ug/l	98
25) 2-Butanone	4.27	43	27281	15.659	ug/l	100
30) Chloroform	4.68	83	4836	1.604	ug/l	94
40) Benzene	5.40	78	1475	0.220	ug/l #	79
51) 4-Methyl-2-Pentanone	7.47	43	4455	1.483	ug/l	91
52) Toluene	7.64	92	993	0.250	ug/l	100
59) 2-Hexanone	8.37	43	5804	2.490	ug/l	98
95) Naphthalene	13.75	128	2612	2.050	ug/l #	92

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

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