

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU080118\
 Data File : VU025783.D
 Acq On : 01 Aug 2018 17:53
 Operator : MD/SY
 Sample : J4284-02
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 48 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 OILY-WATER

Quant Time: Aug 02 05:07:53 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U072718W.M
 Quant Title : SW846 8260
 QLast Update : Tue Jul 31 16:07:47 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	4.99	168	202811	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.89	114	318914	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	313988	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.49	152	193770	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	153646	48.62	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.24%	
35) Dibromofluoromethane	4.89	113	132065	47.61	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.22%	
50) Toluene-d8	7.57	98	467672	48.18	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.36%	
62) 4-Bromofluorobenzene	10.31	95	181623	48.21	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.42%	
Target Compounds						
						Qvalue
11) Tert butyl alcohol	2.83	59	20515	29.652	ug/l	98
16) Acetone	2.33	43	50324	32.664	ug/l	97
20) Methylene Chloride	2.71	84	1299	0.476	ug/l #	85
25) 2-Butanone	4.27	43	11725	5.591	ug/l #	89
29) Tetrahydrofuran	4.65	42	16729	13.314	ug/l	96
30) Chloroform	4.68	83	3250	0.665	ug/l	87
51) 4-Methyl-2-Pentanone	7.47	43	2625	0.657	ug/l	98
95) Naphthalene	13.75	128	56453	5.076	ug/l	99

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

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