

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU070318\
 Data File : VU025198.D
 Acq On : 03 Jul 2018 18:50
 Operator : MD/SY
 Sample : J3826-02
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 OK-03-070218

Quant Time: Jul 04 06:01:56 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U061318W.M
 Quant Title : SW846 8260
 QLast Update : Wed Jun 13 13:55:26 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	4.99	168	202014	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.89	114	312142	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	281949	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.49	152	150005	50.00	ug/l	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
33) 1,2-Dichloroethane-d4	5.31	65	154207	46.76	ug/l	0.00
Spiked Amount			Recovery	=	93.52%	
35) Dibromofluoromethane	4.89	113	120078	46.35	ug/l	0.00
Spiked Amount			Recovery	=	92.70%	
50) Toluene-d8	7.57	98	436517	46.20	ug/l	0.00
Spiked Amount			Recovery	=	92.40%	
62) 4-Bromofluorobenzene	10.31	95	161152	42.95	ug/l	0.00
Spiked Amount			Recovery	=	85.90%	

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
16) Acetone	2.32	43	20964	11.831	ug/l	100
18) Methyl Acetate	2.62	43	19615	5.164	ug/l	97
20) Methylene Chloride	2.71	84	257609	92.654	ug/l	98
51) 4-Methyl-2-Pentanone	7.46	43	2028	0.467	ug/l #	74
67) Ethyl Benzene	9.26	91	4217	0.334	ug/l	100
68) m/p-Xylenes	9.38	106	1440	0.297	ug/l	93
69) o-Xylene	9.79	106	1141	0.238	ug/l	93
84) 1,2,4-Trimethylbenzene	11.16	105	4898	0.481	ug/l	99
95) Naphthalene	13.75	128	93630	9.053	ug/l	98

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

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