

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU011519\
 Data File : VU029192.D
 Acq On : 15 Jan 2019 17:57
 Operator : JC/SP
 Sample : K1139-08
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 S-4

Manual Integrations
 APPROVED

MMDadoda
 1/16/2019 9:42:09 AM

Quant Time: Jan 16 06:50:29 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U010319W.M
 Quant Title : SW846 8260
 QLast Update : Sat Jan 12 01:33:59 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	4.98	168	107682	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.88	114	190793	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	183559	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.48	152	82711	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	75362	56.61	ug/l	0.00
Spiked Amount 50.000			Recovery	=	113.22%	
35) Dibromofluoromethane	4.88	113	64815	52.41	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.82%	
50) Toluene-d8	7.56	98	254724	52.17	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.34%	
62) 4-Bromofluorobenzene	10.30	95	82589	47.07	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.14%	
Target Compounds						
					Qvalue	
16) Acetone	2.32	43	10481	12.499	ug/l	98
20) Methylene Chloride	2.70	84	276100	193.921	ug/l	97
43) Isopropyl Acetate	5.55	43	7477	2.328	ug/l	95
67) Ethyl Benzene	9.25	91	682091	99.293	ug/l	100
68) m/p-Xylenes	9.37	106	1337474	501.092	ug/l	100
69) o-Xylene	9.78	106	703088	273.384	ug/l	98
73) Isopropylbenzene	10.16	105	20021	3.270	ug/l	100
78) n-propylbenzene	10.58	91	111105	15.400	ug/l	99
80) 1,3,5-Trimethylbenzene	10.77	105	489044	95.658	ug/l	99
84) 1,2,4-Trimethylbenzene	11.15	105	2787503	546.118	ug/l	99
85) sec-Butylbenzene	11.32	105	12148m	1.993	ug/l	
86) p-Isopropyltoluene	11.48	119	6562m	1.264	ug/l	
95) Naphthalene	13.74	128	32573	5.089	ug/l	99

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

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