

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU011619\
 Data File : VU029201.D
 Acq On : 16 Jan 2019 11:52
 Operator : JC/SP
 Sample : K1147-09
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 TP-3

Quant Time: Jan 17 01:16:59 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	109296	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.88	114	193634	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	184022	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.48	152	73856	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	76559	56.66	ug/l	0.00
Spiked Amount	50.000		Recovery	=	113.32%	
35) Dibromofluoromethane	4.88	113	66710	53.15	ug/l	0.00
Spiked Amount	50.000		Recovery	=	106.30%	
50) Toluene-d8	7.56	98	253942	51.25	ug/l	0.00
Spiked Amount	50.000		Recovery	=	102.50%	
62) 4-Bromofluorobenzene	10.30	95	79719	44.76	ug/l	0.00
Spiked Amount	50.000		Recovery	=	89.52%	
Target Compounds						
16) Acetone	2.32	43	5267	6.188	ug/l	95
20) Methylene Chloride	2.70	84	59290	41.028	ug/l	98
43) Isopropyl Acetate	5.55	43	6669	2.046	ug/l	96
67) Ethyl Benzene	9.25	91	9652	1.402	ug/l	91
68) m/p-Xylenes	9.37	106	18892	7.060	ug/l	95
69) o-Xylene	9.78	106	8648	3.354	ug/l	99
84) 1,2,4-Trimethylbenzene	11.15	105	13366	2.933	ug/l	98

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

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