

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU121118\
 Data File : VU028582.D
 Acq On : 11 Dec 2018 04:28
 Operator : MD/SY
 Sample : J6305-07
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 P001-SW004-02

Quant Time: Dec 11 08:19:08 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U120618W.M
 Quant Title : SW846 8260
 QLast Update : Fri Dec 07 00:39:42 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	4.98	168	79178	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.88	114	148108	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	134265	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.48	152	54870	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	69272	57.33	ug/l	0.00
Spiked Amount			Recovery	=	114.66%	
35) Dibromofluoromethane	4.88	113	49418	53.02	ug/l	0.00
Spiked Amount			Recovery	=	106.04%	
50) Toluene-d8	7.56	98	190352	51.63	ug/l	0.00
Spiked Amount			Recovery	=	103.26%	
62) 4-Bromofluorobenzene	10.31	95	64718	46.96	ug/l	0.00
Spiked Amount			Recovery	=	93.92%	

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
7) Trichlorofluoromethane	1.88	101	9908	6.688	ug/l	99
16) Acetone	2.32	43	134649	170.771	ug/l	99
17) Carbon Disulfide	2.48	76	2797	0.881	ug/l #	91
18) Methyl Acetate	2.62	43	3468	1.697	ug/l #	88
25) 2-Butanone	4.27	43	23370	21.688	ug/l	96
40) Benzene	5.39	78	8239	1.726	ug/l	99
51) 4-Methyl-2-Pentanone	7.46	43	4701	2.249	ug/l	93
52) Toluene	7.64	92	2932	1.044	ug/l	82
67) Ethyl Benzene	9.25	91	3457	0.657	ug/l	94
68) m/p-Xylenes	9.38	106	2005	1.048	ug/l	100
69) o-Xylene	9.78	106	779	0.419	ug/l	99
70) Styrene	9.79	104	2085	0.677	ug/l	93
84) 1,2,4-Trimethylbenzene	11.15	105	2485	0.692	ug/l	88

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

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