

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU010319\
 Data File : VU029004.D
 Acq On : 03 Jan 2019 18:01
 Operator : MD/SY
 Sample : K1043-01
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 HB-1319-WS

Manual Integrations
 APPROVED

MMDadoda
 1/4/2019 9:51:24 AM

Quant Time: Jan 04 08:03:12 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U010319W.M
 Quant Title : SW846 8260
 QLast Update : Fri Jan 04 02:22:53 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	4.98	168	131977	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	5.88	114	219040	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.09	117	195294	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.48	152	79818	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.31	65	81782	50.13	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.26%	
35) Dibromofluoromethane	4.88	113	70366	49.56	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.12%	
50) Toluene-d8	7.56	98	270796	48.31	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.62%	
62) 4-Bromofluorobenzene	10.30	95	86742	43.06	ug/l	0.00
Spiked Amount 50.000			Recovery	=	86.12%	
Target Compounds						
					Qvalue	
16) Acetone	2.32	43	4362	4.244	ug/l	96
39) Methylcyclohexane	6.42	83	1337	0.497	ug/l #	87
40) Benzene	5.39	78	13727	2.010	ug/l	96
52) Toluene	7.64	92	69672	16.758	ug/l	98
67) Ethyl Benzene	9.25	91	35665	4.880	ug/l	98
68) m/p-Xylenes	9.37	106	59879	21.086	ug/l	99
69) o-Xylene	9.78	106	31161	11.388	ug/l	98
73) Isopropylbenzene	10.17	105	4496	0.761	ug/l	99
78) n-propylbenzene	10.59	91	15574	2.237	ug/l	99
80) 1,3,5-Trimethylbenzene	10.77	105	28031	5.682	ug/l	100
84) 1,2,4-Trimethylbenzene	11.15	105	113262	22.994	ug/l	99
85) sec-Butylbenzene	11.32	105	2932m	0.499	ug/l	
86) p-Isopropyltoluene	11.47	119	10514	2.098	ug/l	97
95) Naphthalene	13.74	128	32564	5.272	ug/l	98

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

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