

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX011419\
 Data File : VX007046.D
 Acq On : 14 Jan 2019 18:16
 Operator : JC/SP
 Sample : K1128-04
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 RBR-200036

Quant Time: Jan 15 00:40:38 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X010819W.M
 Quant Title : SW846 8260
 QLast Update : Tue Jan 08 14:06:59 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.67	168	603101	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.86	114	923361	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.12	117	880947	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	431943	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.07	65	322826	51.22	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.44%	
35) Dibromofluoromethane	5.51	113	286592	48.47	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.94%	
50) Toluene-d8	8.71	98	1129006	52.16	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.32%	
62) 4-Bromofluorobenzene	11.13	95	403619	53.49	ug/l	0.00
Spiked Amount 50.000			Recovery	=	106.98%	
Target Compounds						
16) Acetone	2.45	43	22504	7.222	ug/l	97
18) Methyl Acetate	2.78	43	20088	2.226	ug/l	94
20) Methylene Chloride	2.86	84	1960391	333.994	ug/l	96
43) Isopropyl Acetate	6.46	43	24197	1.782	ug/l	96

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

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