

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX051119\
 Data File : VX009522.D
 Acq On : 11 May 2019 13:57
 Operator : JC/SP
 Sample : K2799-01
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-1

Manual Integrations
 APPROVED

MMDadoda
 5/13/2019 4:27:15 PM

Quant Time: May 12 02:24:32 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X042919W.M
 Quant Title : SW846 8260
 QLast Update : Thu May 09 07:31:15 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.66	168	191346	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.86	114	306093	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.11	117	268202	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	114941	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.06	65	125580	48.08	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.16%	
35) Dibromofluoromethane	5.50	113	93463	48.57	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.14%	
50) Toluene-d8	8.71	98	382783	49.27	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.54%	
62) 4-Bromofluorobenzene	11.13	95	127154	45.05	ug/l	0.00
Spiked Amount 50.000			Recovery	=	90.10%	
Target Compounds						
16) Acetone	2.44	43	4578	3.264	ug/l	98
19) Methyl tert-butyl Ether	3.20	73	2263	0.325	ug/l	99
87) 1,3-Dichlorobenzene	12.02	146	4291	1.160	ug/l	96
88) 1,4-Dichlorobenzene	12.10	146	4500m	1.207	ug/l	
91) 1,2-Dichlorobenzene	12.39	146	11264	2.990	ug/l	97

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

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