

Data Path : Z:\voasrv\HPCHEM1\MSVOA X\Data\VX101819\
 Data File : VX013129.D
 Acq On : 18 Oct 2019 15:58
 Operator : JC/SP
 Sample : K5414-02
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-309-SRI4

Quant Time: Oct 22 07:57:32 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X100819W.M
 Quant Title : SW846 8260
 QLast Update : Wed Oct 16 07:26:22 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.65	168	168115	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.85	114	266345	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.11	117	223090	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.07	152	88478	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.05	65	102632	48.84	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.68%	
35) Dibromofluoromethane	5.49	113	81044	53.02	ug/l	0.00
Spiked Amount 50.000			Recovery	=	106.04%	
50) Toluene-d8	8.71	98	313776	52.63	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.26%	
62) 4-Bromofluorobenzene	11.13	95	103913	44.85	ug/l	0.00
Spiked Amount 50.000			Recovery	=	89.70%	
Target Compounds						
16) Acetone	2.43	43	2245	2.027	ug/l	97
17) Carbon Disulfide	2.56	76	3215	0.664	ug/l	98
30) Chloroform	5.20	83	3527	1.054	ug/l	90
44) Trichloroethene	7.21	130	3701	1.930	ug/l	92
60) Dibromochloromethane	9.58	129	722	2.214	ug/l	91
64) Tetrachloroethene	9.34	164	9764	6.310	ug/l	94

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

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