

Data Path : Z:\voasrv\HPCHEM1\MSVOA X\Data\VX101819\
 Data File : VX013128.D
 Acq On : 18 Oct 2019 15:34
 Operator : JC/SP
 Sample : K5414-06
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 DUP-1-SRI4

Quant Time: Oct 22 07:51:57 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	167991	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.85	114	265261	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.11	117	215885	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.07	152	80268	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.05	65	101302	48.25	ug/l	0.00
Spiked Amount	50.000		Recovery	=	96.50%	
35) Dibromofluoromethane	5.49	113	79913	52.49	ug/l	0.00
Spiked Amount	50.000		Recovery	=	104.98%	
50) Toluene-d8	8.71	98	306234	51.57	ug/l	0.00
Spiked Amount	50.000		Recovery	=	103.14%	
62) 4-Bromofluorobenzene	11.14	95	97536	42.27	ug/l	0.00
Spiked Amount	50.000		Recovery	=	84.54%	
Target Compounds						
					Qvalue	
16) Acetone	2.44	43	2232	2.017	ug/l #	81
17) Carbon Disulfide	2.56	76	3575	0.739	ug/l	98
30) Chloroform	5.20	83	3382	1.011	ug/l	86
44) Trichloroethene	7.21	130	3628	1.899	ug/l	98
60) Dibromochloromethane	9.58	129	570	2.140	ug/l	79
64) Tetrachloroethene	9.33	164	9515	6.354	ug/l	95

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

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