

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX092718\
 Data File : VX004772.D
 Acq On : 27 Sep 2018 14:23
 Operator : JC/MD
 Sample : J5048-05
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 DMW-35A-SEP18

Quant Time: Sep 28 12:50:50 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.M
 Quant Title : SW846 8260
 QLast Update : Wed Sep 26 14:17:02 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.67	168	260628	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.87	114	422308	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.12	117	404658	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	225863	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.07	65	191708	51.49	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.98%	
35) Dibromofluoromethane	5.50	113	162954	45.53	ug/l	0.00
Spiked Amount 50.000			Recovery	=	91.06%	
50) Toluene-d8	8.72	98	599982	50.42	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.84%	
62) 4-Bromofluorobenzene	11.14	95	219403	51.18	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.36%	
Target Compounds						
16) Acetone	2.45	43	288472	132.716	ug/l	96
27) cis-1,2-Dichloroethene	4.60	96	20881	6.046	ug/l	96
44) Trichloroethene	7.21	130	85394	22.722	ug/l	93
64) Tetrachloroethene	9.34	164	106857	26.830	ug/l	97

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX092718\
Data File : VX004772.D
Acq On : 27 Sep 2018 14:23
Operator : JC/MD
Sample : J5048-05
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
DMW-35A-SEP18

Quant Time: Sep 28 12:50:50 2018
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.M
Quant Title : SW846 8260
QLast Update : Wed Sep 26 14:17:02 2018
Response via : Initial Calibration

