

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VO51619\
 Data File : VX009614.D
 Acq On : 16 May 2019 17:55
 Operator : JC/SP
 Sample : K2866-08
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 TP-22-S

Quant Time: May 17 07:15:04 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	5.67	168	169103	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.86	114	273328	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.11	117	241763	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	103000	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.06	65	114099	49.43	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.86%	
35) Dibromofluoromethane	5.50	113	83481	48.58	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.16%	
50) Toluene-d8	8.71	98	343915	49.58	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.16%	
62) 4-Bromofluorobenzene	11.13	95	112787	44.75	ug/l	0.00
Spiked Amount 50.000			Recovery	=	89.50%	
Target Compounds						
16) Acetone	2.45	43	30681	24.752	ug/l	99
18) Methyl Acetate	2.78	43	4536	1.539	ug/l	95
20) Methylene Chloride	2.85	84	198836	99.071	ug/l	96
43) Isopropyl Acetate	6.45	43	12227	2.160	ug/l	98

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX051619\
Data File : VX009614.D
Acq On : 16 May 2019 17:55
Operator : JC/SP
Sample : K2866-08
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
TP-22-S

Quant Time: May 17 07:15:04 2019
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X042919W.M
Quant Title : SW846 8260
QLast Update : Thu May 09 07:31:15 2019
Response via : Initial Calibration

