

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX032619\
 Data File : VX008462.D
 Acq On : 26 Mar 2019 18:21
 Operator : JC/SP
 Sample : K2088-01
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-104

Quant Time: Mar 27 03:18:41 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X032519W.M
 Quant Title : SW846 8260
 QLast Update : Tue Mar 26 07:57:20 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.66	168	242384	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.86	114	407886	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.11	117	351523	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	148564	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.06	65	171473	46.66	ug/l	0.00
Spiked Amount 50.000			Recovery	=	93.32%	
35) Dibromofluoromethane	5.50	113	124345	46.75	ug/l	0.00
Spiked Amount 50.000			Recovery	=	93.50%	
50) Toluene-d8	8.71	98	509646	47.92	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.84%	
62) 4-Bromofluorobenzene	11.13	95	176486	44.78	ug/l	0.00
Spiked Amount 50.000			Recovery	=	89.56%	
Target Compounds						
16) Acetone	2.44	43	18140	9.445	ug/l	94
51) 4-Methyl-2-Pentanone	8.64	43	8175	1.397	ug/l	99
52) Toluene	8.78	92	16197	2.081	ug/l	97
68) m/p-Xylenes	10.35	106	6009	1.145	ug/l	96
69) o-Xylene	10.69	106	2951	0.579	ug/l	91
84) 1,2,4-Trimethylbenzene	11.80	105	5158	0.499	ug/l	99

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX032619\
Data File : VX008462.D
Acq On : 26 Mar 2019 18:21
Operator : JC/SP
Sample : K2088-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-104

Quant Time: Mar 27 03:18:41 2019
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X032519W.M
Quant Title : SW846 8260
QLast Update : Tue Mar 26 07:57:20 2019
Response via : Initial Calibration

