

Data Path : W:\HPCHEM1\MSVOA X\DATA\VX040418\
 Data File : VX000683.D
 Acq On : 05 Apr 2018 02:48
 Operator : JC/MD
 Sample : J2209-02 5X
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 OILY-DEBRIS

Quant Time: Apr 05 06:00:48 2018
 Quant Method : W:\HPCHEM1\MSVOA_X\METHOD\82X040418W.M
 Quant Title : SW846 8260
 QLast Update : Thu Apr 05 03:34:50 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.68	168	85929	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.87	114	148062	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.12	117	144008	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.09	152	75153	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.07	65	65825	45.09	ug/l	0.00
Spiked Amount 50.000			Recovery	=	90.18%	
35) Dibromofluoromethane	5.51	113	47347	46.84	ug/l	0.00
Spiked Amount 50.000			Recovery	=	93.68%	
50) Toluene-d8	8.72	98	189470	44.09	ug/l	0.00
Spiked Amount 50.000			Recovery	=	88.18%	
62) 4-Bromofluorobenzene	11.15	95	64937	38.17	ug/l	0.00
Spiked Amount 50.000			Recovery	=	76.34%	
Target Compounds						
16) Acetone	2.45	43	9255	7.71	ug/l	96
20) Methylene Chloride	2.86	84	31860	23.26	ug/l	93
51) 4-Methyl-2-Pentanone	8.66	43	4515	1.83	ug/l	97
95) Naphthalene	13.84	128	20881	3.69	ug/l	98

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

Data Path : W:\HPCHEM1\MSVOA X\DATA\VX040418\
Data File : VX000683.D
Acq On : 05 Apr 2018 02:48
Operator : JC/MD
Sample : J2209-02 5X
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
OILY-DEBRIS

Quant Time: Apr 05 06:00:48 2018
Quant Method : W:\HPCHEM1\MSVOA_X\METHOD\82X040418W.M
Quant Title : SW846 8260
QLast Update : Thu Apr 05 03:34:50 2018
Response via : Initial Calibration

