

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX051420\
 Data File : VX016264.D
 Acq On : 15 May 2020 04:51
 Operator : JC/SP
 Sample : L2641-04
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 50 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 OWS-BUS-WASH

Manual Integrations
 APPROVED

MMDadoda
 5/15/2020 10:10:45 AM

Quant Time: May 15 07:07:09 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X050420W.M
 Quant Title : SW846 8260
 QLast Update : Wed May 06 06:43:32 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.63	168	256933	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.83	114	432190	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.10	117	429227	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.06	152	226185	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.04	65	160099	47.35	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.70%	
35) Dibromofluoromethane	5.47	113	128610	46.97	ug/l	0.00
Spiked Amount 50.000			Recovery	=	93.94%	
50) Toluene-d8	8.70	98	540489	51.23	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.46%	
62) 4-Bromofluorobenzene	11.12	95	217414	54.18	ug/l	0.00
Spiked Amount 50.000			Recovery	=	108.36%	
Target Compounds						
16) Acetone	2.42	43	73070	48.437	ug/l	99
18) Methyl Acetate	2.75	43	4014	1.082	ug/l	100
20) Methylene Chloride	2.84	84	40958	13.170	ug/l	95
25) 2-Butanone	4.64	43	30983	12.745	ug/l	95
43) Isopropyl Acetate	6.42	43	70795	8.999	ug/l	99
51) 4-Methyl-2-Pentanone	8.62	43	35347m	7.342	ug/l	
52) Toluene	8.76	92	15965	2.058	ug/l	97

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

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