

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX042020\
 Data File : VX015773.D
 Acq On : 20 Apr 2020 18:05
 Operator : JC/SP
 Sample : L2341-02
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 CHASSIS-WASH-2

Manual Integrations
 APPROVED

MMDadoda
 4/21/2020 12:16:22 PM

Quant Time: Apr 21 08:37:52 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X032720W.M
 Quant Title : SW846 8260
 QLast Update : Fri Mar 27 09:35:51 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.63	168	1068963	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.83	114	1675754	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.10	117	1621366	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.06	152	937316	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.03	65	630422	41.81	ug/l	0.00
Spiked Amount 50.000			Recovery	=	83.62%	
35) Dibromofluoromethane	5.47	113	507959	47.42	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.84%	
50) Toluene-d8	8.70	98	2007308	50.90	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.80%	
62) 4-Bromofluorobenzene	11.12	95	847284	52.86	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.72%	
Target Compounds						
16) Acetone	2.42	43	900408	139.011	ug/l	98
17) Carbon Disulfide	2.55	76	207584	6.743	ug/l	98
20) Methylene Chloride	2.83	84	54219	4.481	ug/l	89
25) 2-Butanone	4.64	43	139790	13.420	ug/l	94
51) 4-Methyl-2-Pentanone	8.62	43	251006m	13.051	ug/l	

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

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