

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX091620\
 Data File : VX018443.D
 Acq On : 16 Sep 2020 16:27
 Operator : JC/SP
 Sample : L3988-26
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 SA-DUP05-20200910

Manual Integrations
 APPROVED

MMDadoda
 9/17/2020 11:24:17 AM

Quant Time: Sep 21 09:03:08 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X082120W.M
 Quant Title : SW846 8260
 QLast Update : Fri Aug 21 03:03:56 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.63	168	430278	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.84	114	693667	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.10	117	676918	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.06	152	339416	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.04	65	314193	52.60	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.20%	
35) Dibromofluoromethane	5.47	113	240676	50.95	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.90%	
50) Toluene-d8	8.70	98	844686	48.01	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.02%	
62) 4-Bromofluorobenzene	11.12	95	325274	46.99	ug/l	0.00
Spiked Amount 50.000			Recovery	=	93.98%	
Target Compounds						
12) 1,1-Dichloroethene	2.36	96	2790	0.664	ug/l #	90
24) 1,1-Dichloroethane	3.67	63	30303	3.570	ug/l	97
64) Tetrachloroethene	9.32	164	1992	0.400	ug/l #	51
87) 1,3-Dichlorobenzene	12.01	146	2593	0.239	ug/l	94
88) 1,4-Dichlorobenzene	12.08	146	6687m	0.615	ug/l	

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

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