

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX060220\
 Data File : VX016548.D
 Acq On : 02 Jun 2020 11:56
 Operator : JC/SP
 Sample : L2801-08
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 SS050MW641-200528

Quant Time: Jun 03 04:28:40 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X052720W.M
 Quant Title : SW846 8260
 QLast Update : Wed May 27 16:31:05 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.63	168	352951	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.83	114	583295	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.10	117	569038	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.06	152	291867	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.03	65	202969	46.31	ug/l	0.00
Spiked Amount 50.000			Recovery	=	92.62%	
35) Dibromofluoromethane	5.46	113	170969	47.60	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.20%	
50) Toluene-d8	8.70	98	727622	52.05	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.10%	
62) 4-Bromofluorobenzene	11.12	95	289316	53.56	ug/l	0.00
Spiked Amount 50.000			Recovery	=	107.12%	
Target Compounds						
16) Acetone	2.42	43	1555501	809.997	ug/l	98
17) Carbon Disulfide	2.55	76	5832	0.578	ug/l #	87
25) 2-Butanone	4.63	43	2874839	883.477	ug/l	97
27) cis-1,2-Dichloroethene	4.56	96	7098	1.633	ug/l	82
51) 4-Methyl-2-Pentanone	8.62	43	56175	8.723	ug/l	98
52) Toluene	8.76	92	2998	0.289	ug/l	99
59) 2-Hexanone	9.47	43	14650	2.925	ug/l	94

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

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