

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX051420\
 Data File : VX016223.D
 Acq On : 14 May 2020 12:14
 Operator : JC/SP
 Sample : L2596-02 20X
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 E-1U

Manual Integrations
 APPROVED

MMDadoda
 5/15/2020 12:29:02 PM

Quant Time: May 15 04:37:44 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	233118	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.83	114	392070	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.10	117	390997	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.07	152	216735	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.03	65	145485	47.43	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.86%	
35) Dibromofluoromethane	5.47	113	120434	48.49	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.98%	
50) Toluene-d8	8.70	98	487325	50.92	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.84%	
62) 4-Bromofluorobenzene	11.12	95	204482	56.17	ug/l	0.00
Spiked Amount 50.000			Recovery	=	112.34%	
Target Compounds						
16) Acetone	2.43	43	9701	7.088	ug/l #	86
17) Carbon Disulfide	2.54	76	69639	9.401	ug/l	99
18) Methyl Acetate	2.75	43	1184	0.352	ug/l	98
25) 2-Butanone	4.65	43	3893	1.765	ug/l	92
29) Tetrahydrofuran	5.10	42	7495	5.129	ug/l	99
30) Chloroform	5.19	83	2707m	0.580	ug/l	
95) Naphthalene	13.82	128	19288	1.144	ug/l	99

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

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