

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX042020\
 Data File : VX015776.D
 Acq On : 20 Apr 2020 19:13
 Operator : JC/SP
 Sample : L2342-06
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 TP-18-COMP

Manual Integrations
 APPROVED

MMDadoda
 4/21/2020 12:16:31 PM

Quant Time: Apr 21 08:47:00 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	1029777	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.83	114	1632290	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.10	117	1557907	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.06	152	881993	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.03	65	624576	43.00	ug/l	0.00
Spiked Amount 50.000			Recovery	=	86.00%	
35) Dibromofluoromethane	5.47	113	500802	48.00	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.00%	
50) Toluene-d8	8.70	98	1923172	50.06	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.12%	
62) 4-Bromofluorobenzene	11.12	95	823926	52.77	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.54%	
Target Compounds						
16) Acetone	2.42	43	122639	19.654	ug/l	99
17) Carbon Disulfide	2.55	76	151342	5.103	ug/l	99
20) Methylene Chloride	2.84	84	33445	2.870	ug/l	91
25) 2-Butanone	4.65	43	53740m	5.355	ug/l	
43) Isopropyl Acetate	6.42	43	375349	11.354	ug/l	99

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

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