

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX040920\
 Data File : VX015602.D
 Acq On : 09 Apr 2020 17:30
 Operator : JC/SP
 Sample : L2190-05
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 KY062MW071-200406

Quant Time: Apr 10 08:55:47 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	1224185	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.83	114	1928198	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.10	117	1868071	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.06	152	1048198	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.03	65	806254	46.70	ug/l	0.00
Spiked Amount 50.000			Recovery	=	93.40%	
35) Dibromofluoromethane	5.47	113	597906	48.51	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.02%	
50) Toluene-d8	8.70	98	2337209	51.50	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.00%	
62) 4-Bromofluorobenzene	11.12	95	1017468	55.17	ug/l	0.00
Spiked Amount 50.000			Recovery	=	110.34%	
Target Compounds						
16) Acetone	2.42	43	28524	3.845	ug/l	95
27) cis-1,2-Dichloroethene	4.56	96	257897	17.379	ug/l	91
30) Chloroform	5.17	83	18720	0.772	ug/l	98
44) Trichloroethene	7.19	130	714405	50.334	ug/l	95
64) Tetrachloroethene	9.32	164	4460973	316.429	ug/l	95

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

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