

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX040220\
 Data File : VX015498.D
 Acq On : 02 Apr 2020 21:24
 Operator : JC/SP
 Sample : L2107-03
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 S32-MULTI-IDW01

Quant Time: Apr 03 11:22:54 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	5.64	168	1090920	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.83	114	1723595	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.10	117	1664126	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.07	152	891871	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.03	65	729099	47.39	ug/l	0.00
Spiked Amount	50.000		Recovery	=	94.78%	
35) Dibromofluoromethane	5.47	113	547046	49.65	ug/l	0.00
Spiked Amount	50.000		Recovery	=	99.30%	
50) Toluene-d8	8.70	98	2089864	51.52	ug/l	0.00
Spiked Amount	50.000		Recovery	=	103.04%	
62) 4-Bromofluorobenzene	11.12	95	915619	55.54	ug/l	0.00
Spiked Amount	50.000		Recovery	=	111.08%	
Target Compounds						
16) Acetone	2.42	43	112768	17.059	ug/l	98
21) trans-1,2-Dichloroethene	3.15	96	12525	1.096	ug/l	98
27) cis-1,2-Dichloroethene	4.58	96	385706	29.166	ug/l	93
29) Tetrahydrofuran	5.11	42	16487	2.412	ug/l	94
43) Isopropyl Acetate	6.42	43	249602	7.151	ug/l	98
44) Trichloroethene	7.19	130	127118	10.019	ug/l	96

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

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