

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX083118\
 Data File : VX004333.D
 Acq On : 31 Aug 2018 18:20
 Operator : JC/MD
 Sample : J4721-01
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 CB-0818-S-TCLP-GRID42(4)

Quant Time: Sep 01 05:17:56 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X082218W.M
 Quant Title : SW846 8260
 QLast Update : Thu Aug 23 05:39:44 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.67	168	266024	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.86	114	433001	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.12	117	419869	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	225230	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.07	65	190355	54.90	ug/l	0.00
Spiked Amount	50.000		Recovery	=	109.80%	
35) Dibromofluoromethane	5.50	113	165195	51.28	ug/l	0.00
Spiked Amount	50.000		Recovery	=	102.56%	
50) Toluene-d8	8.71	98	622274	51.18	ug/l	0.00
Spiked Amount	50.000		Recovery	=	102.36%	
62) 4-Bromofluorobenzene	11.14	95	213405	47.04	ug/l	0.00
Spiked Amount	50.000		Recovery	=	94.08%	
Target Compounds						
16) Acetone	2.45	43	17133	4.17	ug/l	97
17) Carbon Disulfide	2.57	76	2221	1.08	ug/l	99
18) Methyl Acetate	2.78	43	6100	1.19	ug/l	93
20) Methylene Chloride	2.86	84	1464268	438.71	ug/l	97
30) Chloroform	5.21	83	8488	1.50	ug/l	95
43) Isopropyl Acetate	6.46	43	164631	22.38	ug/l	100

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

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