

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX101923\
 Data File : VX038387.D
 Acq On : 19 Oct 2023 16:10
 Operator : JC/MD
 Sample : 04966-01DL 10X
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 231013059-VOADL

Quant Time: Oct 20 03:42:00 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100523W.M
 Quant Title : SW846 8260
 QLast Update : Thu Oct 05 15:27:13 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) Pentafluorobenzene	5.550	168	129389	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	229632	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	236831	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	125343	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	90257	43.369	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	86.740%
35) Dibromofluoromethane	5.391	113	73234	44.260	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	88.520%
50) Toluene-d8	8.647	98	287767	45.513	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	91.020%#
62) 4-Bromofluorobenzene	11.079	95	110475	46.088	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	92.180%
Target Compounds						
					Qvalue	
11) Tert butyl alcohol	2.953	59	148540	326.737	ug/l #	91
16) Acetone	2.373	43	142880	164.502	ug/l	95
25) 2-Butanone	4.550	43	213407	152.281	ug/l	93
29) Tetrahydrofuran	5.007	42	41951	45.979	ug/l	89
51) 4-Methyl-2-Pentanone	8.574	43	12834	4.599	ug/l	99
52) Toluene	8.720	92	1741	0.419	ug/l	87
67) Ethyl Benzene	10.195	91	4196	0.474	ug/l	91
68) m/p-Xylenes	10.299	106	2034	0.595	ug/l	95
69) o-Xylene	10.646	106	1039	0.311	ug/l	95
95) Naphthalene	13.774	128	17906	1.935	ug/l	99

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

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