

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX031622\
 Data File : VX027473.D
 Acq On : 16 Mar 2022 16:03
 Operator : JC/MD
 Sample : N1950-10RE
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 3222RE

Quant Time: Mar 17 00:51:55 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X030822W.M
 Quant Title : SW846 8260
 QLast Update : Tue Mar 08 15:54:34 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	360592	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	588159	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	541890	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	254259	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	217665	51.763	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	= 103.520%	
35) Dibromofluoromethane	5.391	113	191925	50.571	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	= 101.140%	
50) Toluene-d8	8.653	98	684857	47.799	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	= 95.600%	
62) 4-Bromofluorobenzene	11.085	95	261108	47.815	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	= 95.640%	
Target Compounds						
						Qvalue
11) Tert butyl alcohol	2.953	59	5769	6.998	ug/l	# 87
16) Acetone	2.386	43	10571	3.640	ug/l	94
52) Toluene	8.720	92	2607	0.255	ug/l	99
67) Ethyl Benzene	10.195	91	5303	0.270	ug/l	93
68) m/p-Xylenes	10.305	106	2071	0.269	ug/l	90
69) o-Xylene	10.647	106	7343	0.964	ug/l	96

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

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