

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX101922\
 Data File : VX032222.D
 Acq On : 19 Oct 2022 20:51
 Operator : JC/MD
 Sample : N5170-09 50X
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 TRE-1667

Quant Time: Oct 20 04:32:08 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X101922W.M
 Quant Title : SW846 8260
 QLast Update : Wed Oct 19 16:51:09 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	137353	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	220290	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	193466	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	93553	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	98373	55.055	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 110.100%	
35) Dibromofluoromethane	5.385	113	75706	49.387	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 98.780%	
50) Toluene-d8	8.647	98	276801	50.163	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	= 100.320%	
62) 4-Bromofluorobenzene	11.079	95	100706	49.839	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 99.680%	
Target Compounds						
					Qvalue	
16) Acetone	2.380	43	8907	15.361	ug/l	98
18) Methyl Acetate	2.709	43	1628	0.827	ug/l	97
19) Methyl tert-butyl Ether	3.111	73	110871	26.244	ug/l	98
52) Toluene	8.720	92	2634	0.711	ug/l	93
67) Ethyl Benzene	10.195	91	3530	0.503	ug/l	96
68) m/p-Xylenes	10.305	106	4513	1.627	ug/l	95
69) o-Xylene	10.646	106	2747	1.034	ug/l	90
78) n-propylbenzene	11.305	91	2145	0.281	ug/l	99
80) 1,3,5-Trimethylbenzene	11.451	105	2473	0.429	ug/l	93
84) 1,2,4-Trimethylbenzene	11.750	105	12355	2.204	ug/l	98

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

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