

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX091521\
 Data File : VX024254.D
 Acq On : 15 Sep 2021 12:41
 Operator : JC/MD
 Sample : M3724-10 10X
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 TOTE-COMP

Quant Time: Sep 16 04:42:05 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091421W.M
 Quant Title : SW846 8260
 QLast Update : Tue Sep 14 12:18:04 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.562	168	121070	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	205971	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	198806	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	82353	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.964	65	90074	52.174	ug/l	0.00
Spiked Amount	50.000	Range 78 - 117	Recovery	= 104.340%		
35) Dibromofluoromethane	5.397	113	68505	49.594	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 99.180%		
50) Toluene-d8	8.653	98	257981	51.267	ug/l	0.00
Spiked Amount	50.000	Range 92 - 112	Recovery	= 102.540%		
62) 4-Bromofluorobenzene	11.085	95	92641	47.306	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 94.620%		
Target Compounds						
					Qvalue	
16) Acetone	2.398	43	453215	615.669	ug/l	91
20) Methylene Chloride	2.788	84	3684	2.606	ug/l	96
25) 2-Butanone	4.580	43	48531	44.761	ug/l	94
51) 4-Methyl-2-Pentanone	8.586	43	11893	5.313	ug/l	89
52) Toluene	8.720	92	4851	1.363	ug/l	98
86) p-Isopropyltoluene	12.012	119	33237	6.334	ug/l	96

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

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