

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX010322\
 Data File : VX026223.D
 Acq On : 03 Jan 2022 16:23
 Operator : JC/MD
 Sample : M5335-01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 20968

Manual Integrations
 APPROVED

Reviewed By : John Carlone 01/04/2022
 Supervised By : Mahesh Dadoda 01/04/2022

Quant Time: Jan 04 03:51:49 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X123021W.M
 Quant Title : SW846 8260
 QLast Update : Thu Dec 30 08:02:21 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	180375	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	315939	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	283403	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	126948	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	141225	54.120	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery = 108.240%			
35) Dibromofluoromethane	5.391	113	103234	50.259	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 100.520%			
50) Toluene-d8	8.659	98	394988	51.699	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery = 103.400%			
62) 4-Bromofluorobenzene	11.079	95	136910	50.338	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery = 100.680%			
Target Compounds					Qvalue	
16) Acetone	2.373	43	1834176	1297.280	ug/l	95
25) 2-Butanone	4.562	43	6344	3.221	ug/l #	84
51) 4-Methyl-2-Pentanone	8.586	43	7738184	1986.205	ug/l	97
52) Toluene	8.726	92	2060	0.354	ug/l	91
68) m/p-Xylenes	10.305	106	1437	0.346	ug/l	77
69) o-Xylene	10.646	106	5535	1.361	ug/l	93
80) 1,3,5-Trimethylbenzene	11.451	105	7980	0.974	ug/l	97
85) sec-Butylbenzene	11.890	105	2023	0.204	ug/l	98
86) p-Isopropyltoluene	12.012	119	3632m	0.439	ug/l	

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

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