

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX082923\
 Data File : VX037371.D
 Acq On : 30 Aug 2023 05:35
 Operator : JC/MD
 Sample : 04198-10
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 49 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 OR-309-OK-370-VOC-25

Quant Time: Aug 31 01:26:39 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X081723W.M
 Quant Title : SW846 8260
 QLast Update : Fri Aug 18 09:30:38 2023
 Response via : Initial Calibration

Manual Integrations
 APPROVED

Reviewed By : John Carlone 08/31/2023
 Supervised By : Mahesh Dadoda 08/31/2023

Supervised By : Mahesh
 Dadoda
 08/31/2023

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	145256	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	244567	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	220673	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	98855	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	91287	43.592	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery =	87.180%		
35) Dibromofluoromethane	5.385	113	75157	46.402	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	92.800%		
50) Toluene-d8	8.647	98	288556	49.113	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery =	98.220%		
62) 4-Bromofluorobenzene	11.079	95	105445	45.669	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	91.340%		
Target Compounds					Qvalue	
8) Diethyl Ether	2.136	74	1085	1.065	ug/l	79
16) Acetone	2.380	43	21408	26.131	ug/l #	84
20) Methylene Chloride	2.788	84	15571	8.499	ug/l #	84
25) 2-Butanone	4.568	43	7619	6.177	ug/l #	82
37) Ethyl Acetate	4.733	43	2717	1.156	ug/l #	91
43) Isopropyl Acetate	6.342	43	84081	22.530	ug/l #	91
51) 4-Methyl-2-Pentanone	8.567	43	15405m	6.798	ug/l	

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

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