

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX080723\
 Data File : VX036846.D
 Acq On : 07 Aug 2023 18:53
 Operator : JC/MD
 Sample : 03887-11
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MCLEAN-TP2

Quant Time: Aug 08 01:40:24 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072123W.M
 Quant Title : SW846 8260
 QLast Update : Fri Jul 21 12:14:12 2023
 Response via : Initial Calibration

Manual Integrations
 APPROVED

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

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	138228	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	254963	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	241138	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	113942	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	96853	47.354	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery =	94.700%		
35) Dibromofluoromethane	5.385	113	81573	46.547	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	93.100%		
50) Toluene-d8	8.647	98	305852	49.215	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery =	98.420%		
62) 4-Bromofluorobenzene	11.079	95	120789	51.516	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	103.040%		
Target Compounds						
16) Acetone	2.380	43	33251	39.178	ug/l #	89
20) Methylene Chloride	2.794	84	5627	2.981	ug/l	91
25) 2-Butanone	4.568	43	15925	11.565	ug/l	95
37) Ethyl Acetate	4.721	43	5100	1.725	ug/l	98
43) Isopropyl Acetate	6.342	43	110433	21.700	ug/l #	94
51) 4-Methyl-2-Pentanone	8.568	43	32158m	11.419	ug/l	

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

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