

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX080723\
 Data File : VX036847.D
 Acq On : 07 Aug 2023 19:17
 Operator : JC/MD
 Sample : 03887-14
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MCLEAN-TP9

Quant Time: Aug 08 01:40:45 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	136128	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	248579	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	237241	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	110455	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	94571	46.952	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery =	93.900%		
35) Dibromofluoromethane	5.385	113	79958	46.798	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	93.600%		
50) Toluene-d8	8.647	98	302756	49.968	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery =	99.940%		
62) 4-Bromofluorobenzene	11.079	95	117846	51.552	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	103.100%		
Target Compounds						
					Qvalue	
16) Acetone	2.386	43	32085	38.387	ug/l	93
18) Methyl Acetate	2.709	43	2192	0.623	ug/l	95
20) Methylene Chloride	2.788	84	5040	2.711	ug/l	89
25) 2-Butanone	4.568	43	14730	10.862	ug/l	93
37) Ethyl Acetate	4.733	43	5781	2.006	ug/l	96
43) Isopropyl Acetate	6.342	43	100106	20.176	ug/l	95
51) 4-Methyl-2-Pentanone	8.568	43	40242m	14.656	ug/l	

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

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