

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX111422\
 Data File : VX032797.D
 Acq On : 14 Nov 2022 16:13
 Operator : JC/MD
 Sample : N5572-06
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 31911

Manual Integrations
 APPROVED

Reviewed By : John Carlone 11/15/2022
 Supervised By : Mahesh Dadoda 11/15/2022

Quant Time : Nov 15 05:08:53 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102622W.M
 Quant Title : SW846 8260
 QLast Update : Thu Oct 27 08:28:06 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.550	168	101910	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	164098	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	146592	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	69590	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	69605	49.156	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery =	98.320%		
35) Dibromofluoromethane	5.391	113	56827	48.548	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	97.100%		
50) Toluene-d8	8.647	98	202806	49.827	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery =	99.660%		
62) 4-Bromofluorobenzene	11.079	95	76611	48.186	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	96.380%		
Target Compounds					Qvalue	
11) Tert butyl alcohol	2.995	59	6625m	32.145	ug/l	
16) Acetone	2.392	43	94214	183.062	ug/l	100
18) Methyl Acetate	2.709	43	8554	5.111	ug/l	99
25) 2-Butanone	4.562	43	168528	211.361	ug/l	98
51) 4-Methyl-2-Pentanone	8.574	43	1939	1.182	ug/l	94
52) Toluene	8.720	92	1131	0.376	ug/l	98

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

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