

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120324\
 Data File : VX044089.D
 Acq On : 03 Dec 2024 13:47
 Operator : JC/MD
 Sample : P5006-02
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 SPLP-C4-3

Manual Integrations
 APPROVED

Reviewed By : John Carlone 12/04/2024
 Supervised By : Mahesh Dadoda 12/04/2024

Quant Time: Dec 03 23:44:00 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112124W.M
 Quant Title : SW846 8260
 QLast Update : Fri Nov 22 03:45:52 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.544	168	112757	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.751	114	220233	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.049	117	195286	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.018	152	89527	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.946	65	103087	53.303	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 106.600%			
35) Dibromofluoromethane	5.379	113	60552	38.478	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 76.960%			
50) Toluene-d8	8.647	98	271904	50.124	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 100.240%			
62) 4-Bromofluorobenzene	11.079	95	97762	52.955	ug/l	0.00
Spiked Amount 50.000	Range 77 - 121		Recovery = 105.900%			
Target Compounds						
16) Acetone	2.380	43	17325	19.504	ug/l	96
27) cis-1,2-Dichloroethene	4.483	96	2630m	1.586	ug/l	
67) Ethyl Benzene	10.195	91	22563	3.105	ug/l	99
84) 1,2,4-Trimethylbenzene	11.750	105	12797	2.270	ug/l	93
95) Naphthalene	13.774	128	1092260	170.566	ug/l	99

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

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