

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092021\
 Data File : VX024356.D
 Acq On : 20 Sep 2021 17:16
 Operator : JC/MD
 Sample : M3817-01RE
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 SW-RR-02-(1.0)RE

Quant Time: Sep 21 08:42:20 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091421W.M
 Quant Title : SW846 8260
 QLast Update : Tue Sep 14 12:18:04 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.562	168	144964	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	245811	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.061	117	205729	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	94213	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.964	65	103998	50.310	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery = 100.620%			
35) Dibromofluoromethane	5.397	113	83654	50.745	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 101.500%			
50) Toluene-d8	8.653	98	258589	43.059	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery = 86.120%#			
62) 4-Bromofluorobenzene	11.085	95	97142	41.565	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery = 83.120%			
Target Compounds					Qvalue	
16) Acetone	2.386	43	79642	90.357	ug/l #	88
17) Carbon Disulfide	2.520	76	1530	0.457	ug/l #	87
25) 2-Butanone	4.580	43	12445	9.586	ug/l	98
40) Benzene	6.050	78	17076	2.634	ug/l	92
52) Toluene	8.726	92	60657	14.278	ug/l	99
65) Chlorobenzene	10.085	112	7520	1.777	ug/l	99
86) p-Isopropyltoluene	12.012	119	10567	1.760	ug/l	90
91) 1,2-Dichlorobenzene	12.341	146	1554	0.502	ug/l	97
95) Naphthalene	13.780	128	38141	6.806	ug/l	99

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

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