

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX091721\
 Data File : VX024306.D
 Acq On : 17 Sep 2021 13:59
 Operator : JC/MD
 Sample : M3791-01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 7 Sample Multiplier: 1

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

Quant Time: Sep 17 14:29:17 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	142414	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	244994	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.061	117	231502	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	103039	50.000	ug/l	0.00

System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.964	65	104998	51.703	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	103.400%
35) Dibromofluoromethane	5.397	113	82217	50.040	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	100.080%
50) Toluene-d8	8.653	98	301321	50.342	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	100.680%
62) 4-Bromofluorobenzene	11.085	95	112550	48.318	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	96.640%

Target Compounds						Qvalue
16) Acetone	2.386	43	67074	77.461	ug/l	90
25) 2-Butanone	4.587	43	8400	6.586	ug/l #	82
40) Benzene	6.050	78	9141	1.415	ug/l	95
52) Toluene	8.726	92	12081	2.853	ug/l	100
65) Chlorobenzene	10.086	112	8271	1.737	ug/l	96
67) Ethyl Benzene	10.201	91	1749	0.210	ug/l	90
86) p-Isopropyltoluene	12.012	119	14706	2.240	ug/l	95
91) 1,2-Dichlorobenzene	12.335	146	2201	0.649	ug/l	94
95) Naphthalene	13.780	128	58369	9.009	ug/l	99

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

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