

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX061121\
 Data File : VX022671.D
 Acq On : 11 Jun 2021 16:43
 Operator : JC/MD
 Sample : M2645-05
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 BUR-1048

Quant Time: Jun 12 02:30:19 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061021W.M
 Quant Title : SW846 8260
 QLast Update : Thu Jun 10 12:20:00 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.562	168	55548	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	6.769	114	112741	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	96717	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	34971	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.964	65	47868	47.915	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery =	95.840%		
35) Dibromofluoromethane	5.397	113	33006	45.885	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	91.780%		
50) Toluene-d8	8.653	98	136494	48.136	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery =	96.280%		
62) 4-Bromofluorobenzene	11.085	95	46124	43.009	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	86.020%		
Target Compounds						
					Qvalue	
16) Acetone	2.392	43	2319	5.184	ug/l	# 84
30) Chloroform	5.105	83	5052	3.385	ug/l	97
52) Toluene	8.720	92	7845	3.772	ug/l	96
67) Ethyl Benzene	10.195	91	4695	1.265	ug/l	# 85
68) m/p-Xylenes	10.305	106	8965	6.670	ug/l	99
69) o-Xylene	10.646	106	3138	2.410	ug/l	91
84) 1,2,4-Trimethylbenzene	11.756	105	7864	3.241	ug/l	98
95) Naphthalene	13.780	128	2332	2.236	ug/l	96

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

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