

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112724\
 Data File : VX044031.D
 Acq On : 27 Nov 2024 11:36
 Operator : JC/MD
 Sample : P5001-01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 SPLP-C1-1

Quant Time: Nov 28 00:43:02 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	113027	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.757	114	217889	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.049	117	202444	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.018	152	89439	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.946	65	100098	51.634	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	= 103.260%		
35) Dibromofluoromethane	5.385	113	54173	34.795	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 69.580%#		
50) Toluene-d8	8.647	98	276760	51.568	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	= 103.140%		
62) 4-Bromofluorobenzene	11.079	95	94376	51.670	ug/l	0.00
Spiked Amount	50.000	Range 77 - 121	Recovery	= 103.340%		
Target Compounds						
					Qvalue	
16) Acetone	2.386	43	7740	8.693	ug/l	98
18) Methyl Acetate	2.703	43	4615	2.226	ug/l	93
40) Benzene	6.038	78	25802	4.269	ug/l	99
52) Toluene	8.714	92	64884	17.938	ug/l	99
67) Ethyl Benzene	10.189	91	461277	61.225	ug/l	100
68) m/p-Xylenes	10.299	106	84417	30.045	ug/l	98
69) o-Xylene	10.640	106	56854	20.725	ug/l	94
73) Isopropylbenzene	10.963	105	23457	3.412	ug/l	99
80) 1,3,5-Trimethylbenzene	11.451	105	23909	4.227	ug/l	91
84) 1,2,4-Trimethylbenzene	11.750	105	81289	14.435	ug/l	99
95) Naphthalene	13.780	128	8232638	1286.861	ug/l	96

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

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