

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120424\
 Data File : VX044117.D
 Acq On : 04 Dec 2024 15:20
 Operator : JC/MD
 Sample : P5006-12DL 5X
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 SPLP-C11-1DL

Quant Time: Dec 05 00:28:08 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	101061	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	6.757	114	197189	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.049	117	174734	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.018	152	75611	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	91646	52.872	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 105.740%			
35) Dibromofluoromethane	5.379	113	62986	44.702	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 89.400%			
50) Toluene-d8	8.647	98	242319	49.891	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 99.780%			
62) 4-Bromofluorobenzene	11.079	95	84083	50.868	ug/l	0.00
Spiked Amount 50.000	Range 77 - 121		Recovery = 101.740%			
Target Compounds						
16) Acetone	2.386	43	3538	4.444	ug/l	99
40) Benzene	6.031	78	22219	4.063	ug/l	97
52) Toluene	8.714	92	11352	3.468	ug/l	96
67) Ethyl Benzene	10.189	91	232333	35.728	ug/l	98
68) m/p-Xylenes	10.299	106	17970	7.410	ug/l	96
69) o-Xylene	10.640	106	21278	8.987	ug/l	98
73) Isopropylbenzene	10.963	105	18497	3.182	ug/l	92
78) n-propylbenzene	11.305	91	9593	1.478	ug/l	97
80) 1,3,5-Trimethylbenzene	11.451	105	14799	3.095	ug/l	99
84) 1,2,4-Trimethylbenzene	11.750	105	53643	11.268	ug/l	98
95) Naphthalene	13.774	128	2198121	406.431	ug/l	99

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

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