

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX100224\
 Data File : VX043239.D
 Acq On : 02 Oct 2024 18:14
 Operator : JC/MD
 Sample : P4236-16
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 TP2

Quant Time: Oct 03 04:59:20 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100124W.M
 Quant Title : SW846 8260
 QLast Update : Wed Oct 02 16:50:57 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.550	168	77427	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.757	114	153567	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	134445	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	59046	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	69367	54.857	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 109.720%			
35) Dibromofluoromethane	5.385	113	51092	48.243	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 96.480%			
50) Toluene-d8	8.647	98	184452	50.258	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 100.520%			
62) 4-Bromofluorobenzene	11.079	95	66664	49.998	ug/l	0.00
Spiked Amount 50.000	Range 77 - 121		Recovery = 100.000%			
Target Compounds						
16) Acetone	2.386	43	29970	56.754	ug/l	99
18) Methyl Acetate	2.709	43	2321	1.752	ug/l #	89
20) Methylene Chloride	2.788	84	3501	3.596	ug/l	90
25) 2-Butanone	4.593	43	4289	5.924	ug/l	93
43) Isopropyl Acetate	6.342	43	83044	31.292	ug/l	100

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

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