

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX100924\
 Data File : VX043341.D
 Acq On : 09 Oct 2024 15:54
 Operator : JC/MD
 Sample : PB163972ZHE#10
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 PB163972ZHE#10

Quant Time: Oct 10 01:33:36 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	85176	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.757	114	158460	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	134221	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.018	152	59608	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	72017	51.772	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	103.540%
35) Dibromofluoromethane	5.385	113	51577	47.197	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	94.400%
50) Toluene-d8	8.647	98	185746	49.047	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	98.100%
62) 4-Bromofluorobenzene	11.079	95	67586	49.124	ug/l	0.00
Spiked Amount	50.000	Range	77 - 121	Recovery	=	98.240%
Target Compounds						
					Qvalue	
16) Acetone	2.392	43	14358	24.716	ug/l	98
18) Methyl Acetate	2.715	43	4021	2.759	ug/l	96
20) Methylene Chloride	2.788	84	3418	3.191	ug/l	93
25) 2-Butanone	4.587	43	6002	7.535	ug/l	96
43) Isopropyl Acetate	6.342	43	85896	31.368	ug/l	100

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

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