

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX100924\
 Data File : VX043332.D
 Acq On : 09 Oct 2024 12:26
 Operator : JC/MD
 Sample : PB163972ZHE#01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 PB163972ZHE#01

Quant Time: Oct 10 01:29:28 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	5.544	168	78860	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.757	114	149834	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	125076	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	53239	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	66891	51.938	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	103.880%
35) Dibromofluoromethane	5.385	113	48249	46.693	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	93.380%
50) Toluene-d8	8.647	98	176913	49.404	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	98.800%
62) 4-Bromofluorobenzene	11.079	95	61097	46.964	ug/l	0.00
Spiked Amount	50.000	Range	77 - 121	Recovery	=	93.920%
Target Compounds						
						Qvalue
16) Acetone	2.392	43	15820	29.414	ug/l	95
18) Methyl Acetate	2.709	43	1909	1.415	ug/l	95
20) Methylene Chloride	2.788	84	6656	6.712	ug/l	98
25) 2-Butanone	4.587	43	4249	5.762	ug/l	92
43) Isopropyl Acetate	6.342	43	87639	33.847	ug/l	99

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

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