

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX100924\
 Data File : VX043337.D
 Acq On : 09 Oct 2024 14:22
 Operator : JC/MD
 Sample : PB163972ZHE#06
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 PB163972ZHE#06

Quant Time: Oct 10 01:31:47 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	83387	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.757	114	155882	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	132559	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	58033	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	71747	52.684	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	105.360%
35) Dibromofluoromethane	5.385	113	50988	47.430	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	94.860%
50) Toluene-d8	8.647	98	182812	49.071	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	98.140%
62) 4-Bromofluorobenzene	11.079	95	65407	48.327	ug/l	0.00
Spiked Amount	50.000	Range	77 - 121	Recovery	=	96.660%
Target Compounds						
					Qvalue	
16) Acetone	2.392	43	14877	26.159	ug/l	99
18) Methyl Acetate	2.709	43	2959	2.074	ug/l	90
20) Methylene Chloride	2.782	84	3596	3.429	ug/l #	92
25) 2-Butanone	4.587	43	5880	7.541	ug/l	94
43) Isopropyl Acetate	6.342	43	86466	32.098	ug/l	99

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

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