

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX100224\
 Data File : VX043240.D
 Acq On : 02 Oct 2024 18:37
 Operator : JC/MD
 Sample : P4242-04
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 TP-Q

Quant Time: Oct 03 04:58:46 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	76288	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.757	114	151070	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	136426	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	59897	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	68487	54.970	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	109.940%
35) Dibromofluoromethane	5.385	113	50292	48.272	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	96.540%
50) Toluene-d8	8.647	98	182780	50.625	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	101.260%
62) 4-Bromofluorobenzene	11.079	95	67581	51.524	ug/l	0.00
Spiked Amount	50.000	Range	77 - 121	Recovery	=	103.040%
Target Compounds						
					Qvalue	
16) Acetone	2.386	43	14610	28.080	ug/l	93
18) Methyl Acetate	2.709	43	2405	1.843	ug/l	93
20) Methylene Chloride	2.782	84	3120	3.252	ug/l	91
25) 2-Butanone	4.593	43	4213	5.906	ug/l	93
43) Isopropyl Acetate	6.342	43	89670	34.348	ug/l	100

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

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