

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX102124\
 Data File : VX043453.D
 Acq On : 21 Oct 2024 14:39
 Operator : JC/MD
 Sample : P4425-06
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :

Quant Time: Oct 21 17:41:51 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	5.556	168	88841	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	165761	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	147262	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	63074	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	76113	52.459	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	104.920%
35) Dibromofluoromethane	5.385	113	56280	49.232	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	98.460%
50) Toluene-d8	8.647	98	198044	49.992	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	99.980%
62) 4-Bromofluorobenzene	11.079	95	72094	50.093	ug/l	0.00
Spiked Amount	50.000	Range	77 - 121	Recovery	=	100.180%
Target Compounds						
						Qvalue
16) Acetone	2.386	43	33351	55.043	ug/l	98
18) Methyl Acetate	2.715	43	2401	1.580	ug/l	98
20) Methylene Chloride	2.788	84	2978	2.666	ug/l	88
43) Isopropyl Acetate	6.342	43	89380	31.202	ug/l	100

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX102124\
Data File : VX043453.D
Acq On : 21 Oct 2024 14:39
Operator : JC/MD
Sample : P4425-06
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :

Quant Time: Oct 21 17:41:51 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

