

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX102124\
 Data File : VX043452.D
 Acq On : 21 Oct 2024 14:16
 Operator : JC/MD
 Sample : P4425-04
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :

Quant Time: Oct 21 17:41:42 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	98579	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	180350	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	153403	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	65906	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	81525	50.638	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	101.280%
35) Dibromofluoromethane	5.385	113	60782	48.869	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	97.740%
50) Toluene-d8	8.647	98	212093	49.207	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	98.420%
62) 4-Bromofluorobenzene	11.079	95	75931	48.491	ug/l	0.00
Spiked Amount	50.000	Range	77 - 121	Recovery	=	96.980%
Target Compounds						
						Qvalue
16) Acetone	2.386	43	33058	49.170	ug/l	97
18) Methyl Acetate	2.709	43	1748	1.036	ug/l	99
20) Methylene Chloride	2.794	84	2712	2.188	ug/l	87
43) Isopropyl Acetate	6.342	43	90716	29.107	ug/l	99

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX102124\
Data File : VX043452.D
Acq On : 21 Oct 2024 14:16
Operator : JC/MD
Sample : P4425-04
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :

Quant Time: Oct 21 17:41:42 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

