

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX101524\
 Data File : VX043403.D
 Acq On : 15 Oct 2024 14:19
 Operator : JC/MD
 Sample : P3932-03
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 FM1

Quant Time: Oct 16 01:56:14 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	84635	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	155406	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	129351	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	55975	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	69939	50.599	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	=	101.200%	
35) Dibromofluoromethane	5.385	113	50015	46.667	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	=	93.340%	
50) Toluene-d8	8.647	98	181721	48.928	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	=	97.860%	
62) 4-Bromofluorobenzene	11.079	95	62367	46.222	ug/l	0.00
Spiked Amount	50.000	Range 77 - 121	Recovery	=	92.440%	
Target Compounds						
					Qvalue	
16) Acetone	2.392	43	25820	44.731	ug/l	96
18) Methyl Acetate	2.709	43	3126	2.159	ug/l	93
20) Methylene Chloride	2.788	84	2164	2.033	ug/l	92
43) Isopropyl Acetate	6.342	43	84822	31.584	ug/l	99

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX101524\
Data File : VX043403.D
Acq On : 15 Oct 2024 14:19
Operator : JC/MD
Sample : P3932-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
FM1

Quant Time: Oct 16 01:56:14 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

