

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX103124\
 Data File : VX043644.D
 Acq On : 31 Oct 2024 12:23
 Operator : JC/MD
 Sample : P4616-04
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 BP-F10

Quant Time: Nov 01 00:51:23 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102824W.M
 Quant Title : SW846 8260
 QLast Update : Mon Oct 28 13:41:34 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) Pentafluorobenzene	5.550	168	135972	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.757	114	248880	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	220188	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.018	152	100651	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	89925	41.476	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	82.960%
35) Dibromofluoromethane	5.385	113	76855	45.540	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	91.080%
50) Toluene-d8	8.647	98	291084	49.826	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	99.660%
62) 4-Bromofluorobenzene	11.079	95	103430	47.779	ug/l	0.00
Spiked Amount	50.000	Range	77 - 121	Recovery	=	95.560%
Target Compounds						
					Qvalue	
16) Acetone	2.392	43	36147	41.984	ug/l	97
20) Methylene Chloride	2.782	84	2718	1.465	ug/l	87
43) Isopropyl Acetate	6.342	43	97827	24.348	ug/l	98

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX103124\
Data File : VX043644.D
Acq On : 31 Oct 2024 12:23
Operator : JC/MD
Sample : P4616-04
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BP-F10

Quant Time: Nov 01 00:51:23 2024
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102824W.M
Quant Title : SW846 8260
QLast Update : Mon Oct 28 13:41:34 2024
Response via : Initial Calibration

