

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX103024\
 Data File : VX043628.D
 Acq On : 30 Oct 2024 16:44
 Operator : JC/MD
 Sample : PB164501ZHE#06
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 PB164501ZHE#06

Quant Time: Oct 31 23:01:30 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	140990	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.757	114	261231	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	228877	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.018	152	108679	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	93709	41.683	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery =	83.360%		
35) Dibromofluoromethane	5.385	113	79264	44.747	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	89.500%		
50) Toluene-d8	8.647	98	302084	49.264	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery =	98.520%		
62) 4-Bromofluorobenzene	11.079	95	106496	46.869	ug/l	0.00
Spiked Amount 50.000	Range 77 - 121		Recovery =	93.740%		
Target Compounds						
16) Acetone	2.386	43	29816	33.398	ug/l	97
18) Methyl Acetate	2.709	43	3374	1.293	ug/l	97
20) Methylene Chloride	2.788	84	2840	1.476	ug/l	89
43) Isopropyl Acetate	6.342	43	121433	28.795	ug/l	98

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX103024\
Data File : VX043628.D
Acq On : 30 Oct 2024 16:44
Operator : JC/MD
Sample : PB164501ZHE#06
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
PB164501ZHE#06

Quant Time: Oct 31 23:01:30 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

