

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX031623\
 Data File : VX034547.D
 Acq On : 16 Mar 2023 15:48
 Operator : JC/MD
 Sample : 01945-02
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 230315024-01-TRIP-BLANK

Quant Time: Mar 17 06:26:48 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X031023W.M
 Quant Title : SW846 8260
 QLast Update : Mon Mar 13 10:29:32 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.550	168	245721	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	422566	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	404988	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	176024	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	170329	47.648	ug/l	0.01
Spiked Amount	50.000	Range	78 - 117	Recovery	=	95.300%
35) Dibromofluoromethane	5.385	113	132756	47.583	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	95.160%
50) Toluene-d8	8.647	98	504645	48.453	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	96.900%
62) 4-Bromofluorobenzene	11.079	95	222448	52.877	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	105.760%
Target Compounds						
					Qvalue	
16) Acetone	2.374	43	78330	46.517	ug/l	97
18) Methyl Acetate	2.703	43	3589	0.652	ug/l	95
30) Chloroform	5.111	83	3772	0.688	ug/l	88

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX031623\
Data File : VX034547.D
Acq On : 16 Mar 2023 15:48
Operator : JC/MD
Sample : 01945-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
230315024-01-TRIP-BLANK

Quant Time: Mar 17 06:26:48 2023
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X031023W.M
Quant Title : SW846 8260
QLast Update : Mon Mar 13 10:29:32 2023
Response via : Initial Calibration

