

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112524\
 Data File : VX043995.D
 Acq On : 25 Nov 2024 18:31
 Operator : JC/MD
 Sample : P4846-05
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 FSND-MW-EVAL-04D-20241113

Quant Time: Nov 26 00:52:33 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112124W.M
 Quant Title : SW846 8260
 QLast Update : Fri Nov 22 03:45:52 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.550	168	113117	50.000	ug/l	0.01
34) 1,4-Difluorobenzene	6.757	114	215357	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	184073	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.018	152	75257	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	96160	49.563	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	99.120%
35) Dibromofluoromethane	5.379	113	70790	46.002	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	92.000%
50) Toluene-d8	8.647	98	262547	49.495	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	99.000%
62) 4-Bromofluorobenzene	11.079	95	84222	46.653	ug/l	0.00
Spiked Amount	50.000	Range	77 - 121	Recovery	=	93.300%
Target Compounds						
					Qvalue	
3) Chloromethane	1.295	50	488	0.324	ug/l	90
16) Acetone	2.386	43	2079	2.333	ug/l	94
44) Trichloroethene	7.123	130	96519	57.081	ug/l	95

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112524\
Data File : VX043995.D
Acq On : 25 Nov 2024 18:31
Operator : JC/MD
Sample : P4846-05
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
FSND-MW-EVAL-04D-20241113

Quant Time: Nov 26 00:52:33 2024
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112124W.M
Quant Title : SW846 8260
QLast Update : Fri Nov 22 03:45:52 2024
Response via : Initial Calibration

