

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX113022\
 Data File : VX033131.D
 Acq On : 30 Nov 2022 14:55
 Operator : JC/MD
 Sample : N5806-09
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 COMP-2

Quant Time: Dec 01 00:20:29 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112222W.M
 Quant Title : SW846 8260
 QLast Update : Tue Nov 22 13:58:46 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	121855	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	198895	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	185263	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	83423	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	67502	57.138	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	114.280%
35) Dibromofluoromethane	5.391	113	53774	47.957	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	95.920%
50) Toluene-d8	8.647	98	187067	55.693	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	111.380%
62) 4-Bromofluorobenzene	11.079	95	69013	46.854	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	93.700%
Target Compounds						
					Qvalue	
16) Acetone	2.392	43	12202	22.017	ug/l	100
20) Methylene Chloride	2.794	84	4310	2.505	ug/l	92
37) Ethyl Acetate	4.714	43	18332	11.720	ug/l	100
43) Isopropyl Acetate	6.342	43	85038	32.873	ug/l	100

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX113022\
Data File : VX033131.D
Acq On : 30 Nov 2022 14:55
Operator : JC/MD
Sample : N5806-09
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
COMP-2

Quant Time: Dec 01 00:20:29 2022
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112222W.M
Quant Title : SW846 8260
QLast Update : Tue Nov 22 13:58:46 2022
Response via : Initial Calibration

