

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN111124\
 Data File : VN084773.D
 Acq On : 11 Nov 2024 15:42
 Operator : JC\MD
 Sample : P4722-18
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 WC-2(2)

Quant Time: Nov 12 00:33:17 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N103024W.M
 Quant Title : SW846 8260
 QLast Update : Thu Oct 31 18:45:38 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.218	168	171996	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.100	114	301059	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	280468	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	127677	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.576	65	121045	48.726	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery =	97.460%		
35) Dibromofluoromethane	8.165	113	100566	49.350	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	98.700%		
50) Toluene-d8	10.565	98	364159	48.512	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery =	97.020%		
62) 4-Bromofluorobenzene	12.847	95	137389	48.972	ug/l	0.00
Spiked Amount 50.000	Range 77 - 121		Recovery =	97.940%		
Target Compounds						
16) Acetone	4.436	43	13805	16.575	ug/l	96
20) Methylene Chloride	5.271	84	3239	1.482	ug/l #	84
25) 2-Butanone	7.482	43	8580	7.403	ug/l	97
43) Isopropyl Acetate	8.682	43	95568	19.587	ug/l #	89

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN111124\
Data File : VN084773.D
Acq On : 11 Nov 2024 15:42
Operator : JC\MD
Sample : P4722-18
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
WC-2(2)

Quant Time: Nov 12 00:33:17 2024
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N103024W.M
Quant Title : SW846 8260
QLast Update : Thu Oct 31 18:45:38 2024
Response via : Initial Calibration

