

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN102021\
 Data File : VN069057.D
 Acq On : 20 Oct 2021 13:15
 Operator : JC/MD
 Sample : M4286-01
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 SW-RR-01-(2.0)

Quant Time: Oct 20 16:40:55 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N101221W.M
 Quant Title : SW846 8260
 QLast Update : Wed Oct 13 05:44:40 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.088	168	321036	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.971	114	558628	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.747	117	512010	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.678	152	181264	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	210223	49.666	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	99.340%
35) Dibromofluoromethane	8.024	113	163976	50.452	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	100.900%
50) Toluene-d8	10.443	98	693262	53.310	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	106.620%
62) 4-Bromofluorobenzene	12.734	95	232175	47.321	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	94.640%
Target Compounds						
					Qvalue	
59) 2-Hexanone	11.092	43	4507	1.247	ug/l	90
74) N-amyl acetate	12.390	43	2714	0.475	ug/l	95

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN102021\
Data File : VN069057.D
Acq On : 20 Oct 2021 13:15
Operator : JC/MD
Sample : M4286-01
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
SW-RR-01-(2.0)

Quant Time: Oct 20 16:40:55 2021
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N101221W.M
Quant Title : SW846 8260
QLast Update : Wed Oct 13 05:44:40 2021
Response via : Initial Calibration

