

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN102721\
 Data File : VN069195.D
 Acq On : 26 Oct 2021 19:16
 Operator : JC/MD
 Sample : M4337-13
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 DDC-8-PS

Manual Integrations
 APPROVED

MMDadoda
 10/27/2021 5:05:45 PM

Quant Time: Oct 27 06:58:22 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	331650	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.971	114	577414	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.746	117	558379	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.677	152	203413	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	216095	49.419	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery = 98.840%			
35) Dibromofluoromethane	8.027	113	170252	50.679	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery = 101.360%			
50) Toluene-d8	10.443	98	716570	53.310	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery = 106.620%			
62) 4-Bromofluorobenzene	12.733	95	254960	50.274	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery = 100.540%			
Target Compounds					Qvalue	
16) Acetone	4.309	43	3009m	1.844	ug/l	
30) Chloroform	7.823	83	4027	0.589	ug/l	92

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN102721\
Data File : VN069195.D
Acq On : 26 Oct 2021 19:16
Operator : JC/MD
Sample : M4337-13
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
DDC-8-PS

Manual Integrations
APPROVED

MMDadoda
10/27/2021 5:05:45 PM

Quant Time: Oct 27 06:58:22 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

