

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN103122\
 Data File : VN075083.D
 Acq On : 31 Oct 2022 13:46
 Operator : JC\MD
 Sample : N5316-09
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 COMPOSITE-PILE-2

Quant Time: Nov 01 06:15:53 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N102622W.M
 Quant Title : SW846 8260
 QLast Update : Thu Oct 27 09:48:49 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.230	168	95192	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	193700	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	173442	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	56218	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	86966	67.805	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 135.600%#	
35) Dibromofluoromethane	8.171	113	65835	57.896	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 115.800%	
50) Toluene-d8	10.571	98	236456	53.597	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 107.200%	
62) 4-Bromofluorobenzene	12.853	95	69456	44.818	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 89.640%	
Target Compounds						
					Qvalue	
16) Acetone	4.442	43	11347	20.894	ug/l	97
20) Methylene Chloride	5.289	84	49354	40.758	ug/l	94
37) Ethyl Acetate	7.571	43	16481	8.536	ug/l	99
43) Isopropyl Acetate	8.694	43	70332	23.485	ug/l #	91

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN103122\
Data File : VN075083.D
Acq On : 31 Oct 2022 13:46
Operator : JC\MD
Sample : N5316-09
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
COMPOSITE-PILE-2

Quant Time: Nov 01 06:15:53 2022
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N102622W.M
Quant Title : SW846 8260
QLast Update : Thu Oct 27 09:48:49 2022
Response via : Initial Calibration

