

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120821\
 Data File : VN069895.D
 Acq On : 9 Dec 2021 6:28
 Operator : JC/MD
 Sample : M4959-14
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 47 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TP-5S

Quant Time: Dec 09 12:00:36 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N120221W.M
 Quant Title : SW846 8260
 QLast Update : Fri Dec 03 08:46:47 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.088	168	1320440	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.968	114	2496899	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.744	117	2781200	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.675	152	1094476	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.437	65	1036884	50.859	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	101.720%
35) Dibromofluoromethane	8.021	113	735692	47.039	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	94.080%
50) Toluene-d8	10.441	98	3292285	52.425	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	104.860%
62) 4-Bromofluorobenzene	12.731	95	1316912	57.884	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	115.760%
Target Compounds						
						Qvalue
16) Acetone	4.291	43	111122	9.061	ug/l	97
18) Methyl Acetate	4.862	43	55853	1.509	ug/l	93
20) Methylene Chloride	5.103	84	33894	2.079	ug/l	92
43) Isopropyl Acetate	8.563	43	172825	3.301	ug/l #	79

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120821\
Data File : VN069895.D
Acq On : 9 Dec 2021 6:28
Operator : JC/MD
Sample : M4959-14
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 47 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
TP-5S

Quant Time: Dec 09 12:00:36 2021
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N120221W.M
Quant Title : SW846 8260
QLast Update : Fri Dec 03 08:46:47 2021
Response via : Initial Calibration

