

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN080924\
 Data File : VN083184.D
 Acq On : 09 Aug 2024 16:00
 Operator : JC\MD
 Sample : P3394-02ME
 Misc : 4.95g/10mL/100uL/5.00mL/MSVOA_N/MEOH
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 B-105-SB02ME

Quant Time: Aug 10 01:00:02 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N080724W.M
 Quant Title : SW846 8260
 QLast Update : Thu Aug 08 06:30:41 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.218	168	132260	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.100	114	245759	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	245304	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	116637	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.577	65	106957	56.814	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	= 113.620%		
35) Dibromofluoromethane	8.165	113	83384	54.358	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 108.720%		
50) Toluene-d8	10.565	98	320087	55.939	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	= 111.880%		
62) 4-Bromofluorobenzene	12.853	95	134045	60.088	ug/l	0.00
Spiked Amount	50.000	Range 64 - 133	Recovery	= 120.180%		
Target Compounds						
					Qvalue	
18) Methyl Acetate	5.036	43	7148	3.257	ug/l #	91
84) 1,2,4-Trimethylbenzene	13.482	105	10106	1.228	ug/l	99

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN080924\
Data File : VN083184.D
Acq On : 09 Aug 2024 16:00
Operator : JC\MD
Sample : P3394-02ME
Misc : 4.95g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
B-105-SB02ME

Quant Time: Aug 10 01:00:02 2024
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N080724W.M
Quant Title : SW846 8260
QLast Update : Thu Aug 08 06:30:41 2024
Response via : Initial Calibration

