

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN081821\
 Data File : VN068187.D
 Acq On : 18 Aug 2021 12:42
 Operator : JC/MD
 Sample : M3408-09
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 33S

Quant Time: Aug 18 13:44:16 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N081721W.M
 Quant Title : SW846 8260
 QLast Update : Tue Aug 17 13:09:49 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.088	168	415084	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	8.971	114	679538	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.749	117	677712	50.000	ug/l	# 0.00
72) 1,4-Dichlorobenzene-d4	13.678	152	254650	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	232005	52.348	ug/l	0.00
Spiked Amount	50.000	Range 61 - 141	Recovery	= 104.700%		
35) Dibromofluoromethane	8.024	113	210582	49.359	ug/l	0.00
Spiked Amount	50.000	Range 69 - 133	Recovery	= 98.720%		
50) Toluene-d8	10.443	98	834981	51.364	ug/l	0.00
Spiked Amount	50.000	Range 65 - 126	Recovery	= 102.720%		
62) 4-Bromofluorobenzene	12.736	95	264325	48.844	ug/l	0.00
Spiked Amount	50.000	Range 58 - 135	Recovery	= 97.680%		
Target Compounds						
						Qvalue
74) N-amyl acetate	12.396	43	3306	0.701	ug/l	94
78) n-propylbenzene	12.921	91	4451	0.229	ug/l	89
82) 4-Chlorotoluene	13.112	91	3284	0.282	ug/l	82
84) 1,2,4-Trimethylbenzene	13.369	105	3267	0.228	ug/l	98
85) sec-Butylbenzene	13.506	105	3518	0.205	ug/l	98
86) p-Isopropyltoluene	13.613	119	3659	0.259	ug/l	91
87) 1,3-Dichlorobenzene	13.619	146	3166	0.372	ug/l	86
89) n-Butylbenzene	13.949	91	7572	0.718	ug/l	96
91) 1,2-Dichlorobenzene	13.991	146	3505	0.433	ug/l	92
93) 1,2,4-Trichlorobenzene	15.268	180	6044	4.031	ug/l	93
95) Naphthalene	15.509	128	24618	5.642	ug/l	# 94
96) 1,2,3-Trichlorobenzene	15.703	180	7552	4.598	ug/l	93

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

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