

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN123120\
 Data File : VN065357.D
 Acq On : 31 Dec 2020 20:03
 Operator : JC/MD
 Sample : L5241-07
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 JG-BUS-WASH

Quant Time: Jan 01 01:54:03 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122220W.M
 Quant Title : SW846 8260
 QLast Update : Tue Dec 22 15:42:18 2020
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	7.630	168	257276	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.553	114	428363	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.379	117	458116	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.318	152	201532	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.994	65	148013	50.96	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	101.92%
35) Dibromofluoromethane	7.553	113	132045	51.09	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	102.18%
50) Toluene-d8	10.061	98	533567	55.09	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	110.18%
62) 4-Bromofluorobenzene	12.376	95	210076	54.35	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	108.70%
Target Compounds						
					Qvalue	
16) Acetone	3.795	43	69378	48.24	ug/l	92
18) Methyl Acetate	4.303	43	4739	1.82	ug/l #	76
20) Methylene Chloride	4.515	84	14124	4.98	ug/l	92
25) 2-Butanone	6.811	43	17284	10.60	ug/l	98
51) 4-Methyl-2-Pentanone	9.955	43	91499	24.46	ug/l	97
52) Toluene	10.125	92	60222	8.25	ug/l	98

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

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