

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN081718\
 Data File : VN050731.D
 Acq On : 17 Aug 2018 15:37
 Operator : MD\SY
 Sample : J4548-01
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 001M-WILLETS-PT-BLVD(AUG)

Manual Integrations
 APPROVED

MMDadoda
 8/20/2018 10:53:10 AM

Quant Time: Aug 18 00:31:23 2018
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA N\methods\624N072518W.M
 Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS
 QLast Update : Thu Jul 26 15:21:50 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	7.20	128	112827	30.00	ug/l	0.00
28) 1,4-Difluorobenzene	8.59	114	615119	30.00	ug/l	0.00
57) Chlorobenzene-d5	11.41	117	503348	30.00	ug/l	0.00

System Monitoring Compounds

27) 1,2-Dichloroethane-d4	8.03	65	231940	30.24	ug/l	0.00
Spiked Amount	30.000	Range	50 - 169	Recovery	=	100.80%
60) 4-Bromofluorobenzene	12.40	95	204106	26.39	ug/l	0.00
Spiked Amount	30.000	Range	56 - 143	Recovery	=	87.97%
63) Toluene-d8	10.09	98	692552	28.81	ug/l	0.00
Spiked Amount	30.000	Range	66 - 137	Recovery	=	96.03%

Target Compounds

						Ovalue
15) Acetone	3.82	58	1518939	2485.63	ug/l	91
16) Carbon Disulfide	4.05	76	12126m	0.66	ug/l	
30) 2-Butanone	6.84	43	72048	24.05	ug/l	99
58) 4-Methyl-2-Pentanone	9.99	43	14929	2.55	ug/l	98
62) Toluene	10.16	91	2289466	79.72	ug/l	100
82) p-Isopropyltoluene	13.29	119	224501	9.64	ug/l	98

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

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