

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN091019\
 Data File : VN057816.D
 Acq On : 9 Sep 2019 21:19
 Operator : JC/SP
 Sample : K4781-03
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 GW801-20.0-25.0-090719

Quant Time: Sep 13 04:37:30 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N091019W.M
 Quant Title : SW846 8260
 QLast Update : Tue Sep 10 03:02:13 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.65	168	305526	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.58	114	511668	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	453946	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	228002	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.01	65	328431	53.07	ug/l	0.00
Spiked Amount 50.000			Recovery	=	106.14%	
35) Dibromofluoromethane	7.58	113	230365	48.78	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.56%	
50) Toluene-d8	10.09	98	929514	50.70	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.40%	
62) 4-Bromofluorobenzene	12.40	95	294889	42.20	ug/l	0.00
Spiked Amount 50.000			Recovery	=	84.40%	
Target Compounds						
16) Acetone	3.80	43	39793	19.616	ug/l	99
17) Carbon Disulfide	4.03	76	3671	0.881	ug/l #	73
30) Chloroform	7.36	83	190456	22.323	ug/l	98
47) Bromodichloromethane	9.40	83	22096	3.522	ug/l	93
60) Dibromochloromethane	10.90	129	2958	0.668	ug/l	93
64) Tetrachloroethene	10.63	164	2369	0.707	ug/l	86

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN091019\
Data File : VN057816.D
Acq On : 9 Sep 2019 21:19
Operator : JC/SP
Sample : K4781-03
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
GW801-20.0-25.0-090719

Quant Time: Sep 13 04:37:30 2019
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N091019W.M
Quant Title : SW846 8260
QLast Update : Tue Sep 10 03:02:13 2019
Response via : Initial Calibration

