

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN061920\
 Data File : VN061950.D
 Acq On : 19 Jun 2020 14:25
 Operator : JC/MD
 Sample : L2984-04
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 S49-0564

Quant Time: Jun 20 08:19:10 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N061620W.M
 Quant Title : SW846 8260
 QLast Update : Wed Jun 17 01:19:44 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.63	168	152082	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	276099	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	266200	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	105007	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	100955	49.50	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.00%	
35) Dibromofluoromethane	7.55	113	87881	49.75	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.50%	
50) Toluene-d8	10.06	98	344283	49.58	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.16%	
62) 4-Bromofluorobenzene	12.38	95	114230	47.96	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.92%	
Target Compounds						
4) Vinyl Chloride	2.16	62	18806	7.473	ug/l	95
16) Acetone	3.78	43	17365	28.823	ug/l	95
25) 2-Butanone	6.79	43	112529	128.018	ug/l	96
27) cis-1,2-Dichloroethene	6.78	96	6589	3.217	ug/l	92
52) Toluene	10.12	92	3124	0.652	ug/l	90

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

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN061920\
Data File : VN061950.D
Acq On : 19 Jun 2020 14:25
Operator : JC/MD
Sample : L2984-04
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
S49-0564

Quant Time: Jun 20 08:19:10 2020
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N061620W.M
Quant Title : SW846 8260
QLast Update : Wed Jun 17 01:19:44 2020
Response via : Initial Calibration

