

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN092220\
 Data File : VN063576.D
 Acq On : 23 Sep 2020 9:40
 Operator : JC/MD
 Sample : L4104-12
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 55 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PAV-B29-092120-16-17

Quant Time: Sep 23 10:27:16 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N091620W.M
 Quant Title : SW846 8260
 QLast Update : Wed Sep 16 13:05:20 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.63	168	212532	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	407676	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	415006	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	169681	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	181353	50.13	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.26%	
35) Dibromofluoromethane	7.55	113	130629	46.29	ug/l	0.00
Spiked Amount 50.000			Recovery	=	92.58%	
50) Toluene-d8	10.06	98	521873	49.18	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.36%	
62) 4-Bromofluorobenzene	12.38	95	200957	51.66	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.32%	
Target Compounds						
16) Acetone	3.78	43	22169	19.067	ug/l	93
20) Methylene Chloride	4.50	84	12176	4.133	ug/l	94
25) 2-Butanone	6.80	43	6163	3.446	ug/l	99
43) Isopropyl Acetate	8.13	43	79138	10.383	ug/l	99
51) 4-Methyl-2-Pentanone	9.96	43	8507	1.994	ug/l	97
59) 2-Hexanone	10.74	43	87007	28.145	ug/l #	54

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN092220\
Data File : VN063576.D
Acq On : 23 Sep 2020 9:40
Operator : JC/MD
Sample : L4104-12
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 55 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
PAV-B29-092120-16-17

Quant Time: Sep 23 10:27:16 2020
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N091620W.M
Quant Title : SW846 8260
QLast Update : Wed Sep 16 13:05:20 2020
Response via : Initial Calibration

