

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN100120\
 Data File : VN063802.D
 Acq On : 2 Oct 2020 4:32
 Operator : JC/MD
 Sample : L4190-16
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 45 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PAV-B18-092920-19-20

Quant Time: Oct 02 05:50:15 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N093020W.M
 Quant Title : SW846 8260
 QLast Update : Thu Oct 01 01:46:38 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.63	168	204899	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	393647	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	411643	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	167613	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	185469	54.79	ug/l	0.00
Spiked Amount 50.000			Recovery	=	109.58%	
35) Dibromofluoromethane	7.55	113	132485	48.50	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.00%	
50) Toluene-d8	10.06	98	515898	49.82	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.64%	
62) 4-Bromofluorobenzene	12.38	95	199784	51.33	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.66%	
Target Compounds						
16) Acetone	3.78	43	27505	26.395	ug/l	96
18) Methyl Acetate	4.30	43	5444	2.012	ug/l #	70
20) Methylene Chloride	4.50	84	9121	3.912	ug/l	96
43) Isopropyl Acetate	8.13	43	75507	11.554	ug/l	99

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

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