

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN100620\
 Data File : VN063930.D
 Acq On : 7 Oct 2020 7:06
 Operator : JC/MD
 Sample : L4260-16
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 49 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PAV-B23-100220-18-19

Quant Time: Oct 07 07:35:05 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	226746	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	427844	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	448792	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	174701	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	202417	54.04	ug/l	0.00
Spiked Amount	50.000		Recovery	=	108.08%	
35) Dibromofluoromethane	7.55	113	145913	49.15	ug/l	0.00
Spiked Amount	50.000		Recovery	=	98.30%	
50) Toluene-d8	10.06	98	553751	49.20	ug/l	0.00
Spiked Amount	50.000		Recovery	=	98.40%	
62) 4-Bromofluorobenzene	12.38	95	213510	50.47	ug/l	0.00
Spiked Amount	50.000		Recovery	=	100.94%	
Target Compounds						
16) Acetone	3.78	43	17031	14.769	ug/l	92
17) Carbon Disulfide	4.01	76	6049	1.667	ug/l #	94
18) Methyl Acetate	4.29	43	3185	1.064	ug/l #	75
20) Methylene Chloride	4.51	84	9735	3.777	ug/l	88
27) cis-1,2-Dichloroethene	6.78	96	6738	2.291	ug/l	94
43) Isopropyl Acetate	8.13	43	85447	12.029	ug/l	99

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

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