

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN100820\
 Data File : VN064026.D
 Acq On : 9 Oct 2020 7:16
 Operator : JC/MD
 Sample : L4311-02
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 53 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PAV-B10-100720-0-1

Manual Integrations
 APPROVED

MMDadoda
 10/9/2020 4:53:27 PM

Quant Time: Oct 09 12:24:53 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	237428	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	458034	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	480744	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	182254	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	201664	51.41	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.82%	
35) Dibromofluoromethane	7.55	113	153562	48.32	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.64%	
50) Toluene-d8	10.06	98	582698	48.36	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.72%	
62) 4-Bromofluorobenzene	12.38	95	221948	49.01	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.02%	
Target Compounds						
16) Acetone	3.78	43	15171	12.564	ug/l	95
20) Methylene Chloride	4.51	84	11773m	4.346	ug/l	
43) Isopropyl Acetate	8.13	43	88296	11.611	ug/l	99
51) 4-Methyl-2-Pentanone	9.95	43	8196	1.837	ug/l	97

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

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