

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN092520\
 Data File : VN063691.D
 Acq On : 25 Sep 2020 20:18
 Operator : JC/MD
 Sample : L4133-01
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW1S

Manual Integrations
 APPROVED

MMDadoda
 9/29/2020 10:41:02 AM

Quant Time: Sep 28 08:53:35 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	205194	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	388004	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	394152	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	162691	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	180591	51.70	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.40%	
35) Dibromofluoromethane	7.55	113	130037	48.41	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.82%	
50) Toluene-d8	10.06	98	495949	49.10	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.20%	
62) 4-Bromofluorobenzene	12.38	95	193633	52.30	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.60%	
Target Compounds						
16) Acetone	3.78	43	3652	3.253	ug/l	93
27) cis-1,2-Dichloroethene	6.78	96	41842	14.409	ug/l	89
44) Trichloroethene	8.80	130	17576	6.141	ug/l	81
52) Toluene	10.13	92	2808m	0.396	ug/l	
64) Tetrachloroethene	10.60	164	239706	89.313	ug/l	98

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

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