

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN092220\
 Data File : VN063563.D
 Acq On : 23 Sep 2020 4:27
 Operator : JC/MD
 Sample : PB131772ZHE#01
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB131772ZHE#01

Manual Integrations
 APPROVED

MMDadoda
 9/23/2020 1:06:47 PM

Quant Time: Sep 23 07:17:59 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	219384	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	422712	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	439081	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	190559	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	188633	50.51	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.02%	
35) Dibromofluoromethane	7.55	113	138863	47.45	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.90%	
50) Toluene-d8	10.06	98	549707	49.96	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.92%	
62) 4-Bromofluorobenzene	12.38	95	219968	54.53	ug/l	0.00
Spiked Amount 50.000			Recovery	=	109.06%	
Target Compounds						
16) Acetone	3.78	43	32816	27.342	ug/l	97
18) Methyl Acetate	4.30	43	4183m	1.373	ug/l	
20) Methylene Chloride	4.51	84	19565	6.434	ug/l	86
43) Isopropyl Acetate	8.13	43	80339	10.165	ug/l	99

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

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