

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN092220\
 Data File : VN063564.D
 Acq On : 23 Sep 2020 4:51
 Operator : JC/MD
 Sample : PB131772ZHE#02
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB131772ZHE#02

Manual Integrations
 APPROVED

MMDadoda
 9/23/2020 1:06:48 PM

Quant Time: Sep 23 07:19:40 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	206228	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	394010	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	397928	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	163541	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	179962	51.26	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.52%	
35) Dibromofluoromethane	7.55	113	128767	47.21	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.42%	
50) Toluene-d8	10.06	98	509242	49.65	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.30%	
62) 4-Bromofluorobenzene	12.38	95	196263	52.20	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.40%	
Target Compounds						
16) Acetone	3.78	43	32873	29.137	ug/l	94
18) Methyl Acetate	4.30	43	4424m	1.545	ug/l	
20) Methylene Chloride	4.51	84	18671	6.532	ug/l	93
43) Isopropyl Acetate	8.13	43	81322	11.039	ug/l	100

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

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