

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN100520\
 Data File : VN063869.D
 Acq On : 5 Oct 2020 14:33
 Operator : JC/MD
 Sample : PB132063ZHE#01
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB132063ZHE#01

Manual Integrations
 APPROVED

MMDadoda
 10/6/2020 12:56:48 PM

Quant Time: Oct 06 02:12:19 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	227368	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	425351	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	445504	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	175933	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	199685	53.16	ug/l	0.00
Spiked Amount 50.000			Recovery	=	106.32%	
35) Dibromofluoromethane	7.55	113	142012	48.12	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.24%	
50) Toluene-d8	10.06	98	546565	48.85	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.70%	
62) 4-Bromofluorobenzene	12.38	95	215833	51.32	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.64%	
Target Compounds						
16) Acetone	3.78	43	11425	9.881	ug/l	100
18) Methyl Acetate	4.29	43	3556m	1.184	ug/l	
20) Methylene Chloride	4.50	84	11945	4.599	ug/l	97
43) Isopropyl Acetate	8.13	43	79257	11.223	ug/l	100

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

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