

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN100120\
 Data File : VN063783.D
 Acq On : 1 Oct 2020 19:45
 Operator : JC/MD
 Sample : PB132009ZHE#05
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB132009ZHE#05

Manual Integrations
 APPROVED

MMDadoda
 10/2/2020 3:11:54 PM

Quant Time: Oct 02 03:17:50 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	211643	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	401095	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	418881	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	170400	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	185363	53.01	ug/l	0.00
Spiked Amount 50.000			Recovery	=	106.02%	
35) Dibromofluoromethane	7.55	113	135266	48.60	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.20%	
50) Toluene-d8	10.06	98	520791	49.36	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.72%	
62) 4-Bromofluorobenzene	12.38	95	208977	52.69	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.38%	
Target Compounds						
16) Acetone	3.78	43	20587	19.127	ug/l	99
18) Methyl Acetate	4.30	43	4361m	1.561	ug/l	
20) Methylene Chloride	4.51	84	7099	2.972	ug/l #	91
43) Isopropyl Acetate	8.13	43	79202	11.894	ug/l	99

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

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