

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN102120\
 Data File : VN064244.D
 Acq On : 22 Oct 2020 3:50
 Operator : JC/MD
 Sample : PB132513ZHE#09
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB132513ZHE#09

Manual Integrations
 APPROVED

MMDadoda
 10/23/2020 2:10:53 PM

Quant Time: Oct 22 05:58:39 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N102120W.M
 Quant Title : SW846 8260
 QLast Update : Wed Oct 21 12:48:40 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.63	168	262549	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	458607	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	459722	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	210763	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	204010	52.32	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.64%	
35) Dibromofluoromethane	7.55	113	149009	48.87	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.74%	
50) Toluene-d8	10.06	98	583697	52.77	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.54%	
62) 4-Bromofluorobenzene	12.38	95	262057	58.45	ug/l	0.00
Spiked Amount 50.000			Recovery	=	116.90%	
Target Compounds						
16) Acetone	3.79	43	10271	9.207	ug/l	97
18) Methyl Acetate	4.29	43	4402m	1.548	ug/l	
20) Methylene Chloride	4.50	84	7461	2.508	ug/l	94
43) Isopropyl Acetate	8.13	43	92391	11.942	ug/l #	90

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

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