

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN080620\
 Data File : VN062769.D
 Acq On : 6 Aug 2020 16:18
 Operator : JC/MD
 Sample : PB130737ZHE#11
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB130737ZHE#11

Manual Integrations
 APPROVED

MMDadoda
 8/7/2020 12:50:48 PM

Quant Time: Aug 07 04:30:28 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N080520W.M
 Quant Title : SW846 8260
 QLast Update : Wed Aug 05 13:26:42 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.63	168	135054	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	238511	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	248892	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	86311	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	115271	51.61	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.22%	
35) Dibromofluoromethane	7.55	113	81543	49.98	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.96%	
50) Toluene-d8	10.06	98	321705	50.51	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.02%	
62) 4-Bromofluorobenzene	12.38	95	116442	48.37	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.74%	
Target Compounds						
16) Acetone	3.78	43	21929	27.173	ug/l	94
18) Methyl Acetate	4.28	43	4062m	1.022	ug/l	
20) Methylene Chloride	4.50	84	8991	5.410	ug/l #	91
43) Isopropyl Acetate	8.13	43	51935	12.835	ug/l	98

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

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