

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN092820\
 Data File : VN063709.D
 Acq On : 28 Sep 2020 16:45
 Operator : JC/MD
 Sample : PB131914ZHE#04
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB131914ZHE#04

Manual Integrations
 APPROVED

MMDadoda
 9/29/2020 4:38:05 PM

Quant Time: Sep 29 04:26:39 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	218774	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	406310	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	434405	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	182034	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	189705	50.94	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.88%	
35) Dibromofluoromethane	7.55	113	138275	49.16	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.32%	
50) Toluene-d8	10.06	98	541059	51.16	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.32%	
62) 4-Bromofluorobenzene	12.38	95	214628	55.36	ug/l	0.00
Spiked Amount 50.000			Recovery	=	110.72%	
Target Compounds						
16) Acetone	3.78	43	17104	14.291	ug/l	94
18) Methyl Acetate	4.28	43	3443m	1.133	ug/l	
20) Methylene Chloride	4.51	84	13829	4.560	ug/l #	90
25) 2-Butanone	6.80	43	10098	5.486	ug/l	94
43) Isopropyl Acetate	8.13	43	70864	9.328	ug/l	99

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

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