

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN070120\
 Data File : VN062321.D
 Acq On : 2 Jul 2020 4:45
 Operator : JC/MD
 Sample : PB129873ZHE#07
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB129873ZHE#07

Manual Integrations
 APPROVED

MMDadoda
 7/2/2020 12:41:01 PM

Quant Time: Jul 02 08:21:28 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N062420W.M
 Quant Title : SW846 8260
 QLast Update : Wed Jun 24 15:33:00 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.63	168	192430	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	415149	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	400745	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	137788	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	161773	59.91	ug/l	0.00
Spiked Amount 50.000			Recovery	=	119.82%	
35) Dibromofluoromethane	7.55	113	130417	48.97	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.94%	
50) Toluene-d8	10.06	98	538713	50.60	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.20%	
62) 4-Bromofluorobenzene	12.38	95	169566	47.14	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.28%	
Target Compounds						
16) Acetone	3.78	43	11998	10.270	ug/l	98
18) Methyl Acetate	4.28	43	6699m	3.121	ug/l	
20) Methylene Chloride	4.50	84	7443	2.798	ug/l #	94
43) Isopropyl Acetate	8.13	43	74298	13.190	ug/l	95

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

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