

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN082318\
 Data File : VN050849.D
 Acq On : 23 Aug 2018 13:32
 Operator : MD\SY
 Sample : PB112301ZHE#03
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB112301ZHE#03

Manual Integrations
 APPROVED

MMDadoda
 8/24/2018 10:40:53 AM

Quant Time: Aug 24 03:51:39 2018
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N082118W.M
 Quant Title : SW846 8260
 QLast Update : Wed Aug 22 05:38:12 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.67	168	615215	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.59	114	1031600	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	859741	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	286495	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.03	65	409102	55.24	ug/l	0.00
Spiked Amount 50.000			Recovery	=	110.48%	
35) Dibromofluoromethane	7.59	113	392886	47.87	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.74%	
50) Toluene-d8	10.09	98	1392787	45.39	ug/l	0.00
Spiked Amount 50.000			Recovery	=	90.78%	
62) 4-Bromofluorobenzene	12.40	95	366278	35.92	ug/l	0.00
Spiked Amount 50.000			Recovery	=	71.84%	
Target Compounds						
16) Acetone	3.82	43	15400	1.77	ug/l	95
18) Methyl Acetate	4.34	43	13259	1.90	ug/l	96
20) Methylene Chloride	4.55	84	21967	2.63	ug/l	91
29) Tetrahydrofuran	7.23	42	2050m	1.20	ug/l	
43) Isopropyl Acetate	8.17	43	332532	27.66	ug/l #	90

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

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