

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN091019\
 Data File : VN057854.D
 Acq On : 10 Sep 2019 12:02
 Operator : JC/SP
 Sample : K4622-34
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 60 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PAR-11-B4-(0)

Manual Integrations
 APPROVED

MMDadoda
 9/13/2019 8:16:44 AM

Quant Time: Sep 13 05:01:23 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N091019W.M
 Quant Title : SW846 8260
 QLast Update : Tue Sep 10 03:02:13 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.65	168	292053	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.58	114	516574	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	484340	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	244404	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.01	65	335974	56.80	ug/l	0.00
Spiked Amount 50.000			Recovery	=	113.60%	
35) Dibromofluoromethane	7.58	113	225549	47.31	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.62%	
50) Toluene-d8	10.09	98	950636	51.36	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.72%	
62) 4-Bromofluorobenzene	12.40	95	299239	42.42	ug/l	0.00
Spiked Amount 50.000			Recovery	=	84.84%	
Target Compounds						
16) Acetone	3.81	43	13634	7.031	ug/l	96
20) Methylene Chloride	4.53	84	8208m	1.018	ug/l	
43) Isopropyl Acetate	8.15	43	96593m	9.280	ug/l	

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

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