

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN091019\
 Data File : VN057840.D
 Acq On : 10 Sep 2019 6:52
 Operator : JC/SP
 Sample : K4622-25
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 46 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PAR-11-A4-(0)

Manual Integrations
 APPROVED

MMDadoda
 9/13/2019 5:16:28 AM

Quant Time: Sep 10 08:55:59 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	304422	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.58	114	523056	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	485584	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	246873	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.01	65	339294	55.03	ug/l	0.00
Spiked Amount 50.000			Recovery	=	110.06%	
35) Dibromofluoromethane	7.58	113	228117	47.25	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.50%	
50) Toluene-d8	10.09	98	948569	50.61	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.22%	
62) 4-Bromofluorobenzene	12.40	95	303856	42.54	ug/l	0.00
Spiked Amount 50.000			Recovery	=	85.08%	
Target Compounds						
16) Acetone	3.80	43	9718	4.808	ug/l	98
18) Methyl Acetate	4.32	43	8568m	1.590	ug/l	
43) Isopropyl Acetate	8.16	43	100453	9.531	ug/l #	87

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

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