

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN061219\
 Data File : VN056254.D
 Acq On : 12 Jun 2019 17:25
 Operator : JC/SP
 Sample : K3327-01
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-3

Manual Integrations
 APPROVED

MMDadoda
 6/13/2019 11:08:46 AM

Quant Time: Jun 13 05:30:08 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N060419W.M
 Quant Title : SW846 8260
 QLast Update : Tue Jun 04 11:02:09 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.66	168	312441	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.59	114	468216	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	465097	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.34	152	154934	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.03	65	153310	49.42	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.84%	
35) Dibromofluoromethane	7.59	113	145289	50.58	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.16%	
50) Toluene-d8	10.09	98	575094	52.03	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.06%	
62) 4-Bromofluorobenzene	12.40	95	188387	50.11	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.22%	
Target Compounds						
16) Acetone	3.83	43	5971	4.024	ug/l	94
44) Trichloroethene	8.84	130	1123	0.320	ug/l	95
87) 1,3-Dichlorobenzene	13.28	146	5457	1.063	ug/l	97
88) 1,4-Dichlorobenzene	13.36	146	5602m	1.134	ug/l	
91) 1,2-Dichlorobenzene	13.65	146	14132	2.726	ug/l	97

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

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