

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN091419\
 Data File : VN058013.D
 Acq On : 13 Sep 2019 13:08
 Operator : JC/SP
 Sample : K4859-06
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 IDW-GW200A-B4-091119

Manual Integrations
 APPROVED

MMDadoda
 9/16/2019 2:49:51 PM

Quant Time: Sep 14 03:11:17 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	265161	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.58	114	449089	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	407952	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	210483	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.01	65	321200	59.81	ug/l	0.00
Spiked Amount 50.000			Recovery	=	119.62%	
35) Dibromofluoromethane	7.58	113	226681	54.69	ug/l	0.00
Spiked Amount 50.000			Recovery	=	109.38%	
50) Toluene-d8	10.09	98	825640	51.30	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.60%	
62) 4-Bromofluorobenzene	12.40	95	279311	45.54	ug/l	0.00
Spiked Amount 50.000			Recovery	=	91.08%	
Target Compounds						
16) Acetone	3.80	43	24126	13.703	ug/l	94
25) 2-Butanone	6.84	43	7126m	2.796	ug/l	
30) Chloroform	7.36	83	89827	12.131	ug/l	96
40) Benzene	8.03	78	4879	0.311	ug/l #	88
44) Trichloroethene	8.83	130	1243	0.323	ug/l	96
51) 4-Methyl-2-Pentanone	9.98	43	15864	2.962	ug/l	98
52) Toluene	10.15	92	7283	0.737	ug/l	98
64) Tetrachloroethene	10.63	164	15346	5.095	ug/l	95
67) Ethyl Benzene	11.51	91	7933	0.424	ug/l	96
68) m/p-Xylenes	11.62	106	4792	0.699	ug/l	96
69) o-Xylene	11.95	106	6160	0.904	ug/l	92
95) Naphthalene	15.13	128	46969	2.313	ug/l	98

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

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