

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN091919\
 Data File : VN058229.D
 Acq On : 20 Sep 2019 00:47
 Operator : JC/SP
 Sample : K4974-08
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 RE134D3-20190916

Manual Integrations
 APPROVED

MMDadoda
 9/20/2019 3:36:51 PM

Quant Time: Sep 20 09:08:21 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N091819W.M
 Quant Title : SW846 8260
 QLast Update : Thu Sep 19 09:27:53 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.65	168	516502	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.57	114	838992	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	718918	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	296625	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.01	65	366181	50.90	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.80%	
35) Dibromofluoromethane	7.58	113	253004	49.60	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.20%	
50) Toluene-d8	10.08	98	1056176	53.15	ug/l	0.00
Spiked Amount 50.000			Recovery	=	106.30%	
62) 4-Bromofluorobenzene	12.40	95	339031	45.93	ug/l	0.00
Spiked Amount 50.000			Recovery	=	91.86%	
Target Compounds						
3) Chloromethane	2.04	50	4489	0.346	ug/l	90
9) 1,1,2-Trichlorotrifluoroet	3.74	101	1723206	412.690	ug/l	99
12) 1,1-Dichloroethene	3.74	96	7864	2.141	ug/l #	38
16) Acetone	3.80	43	13855m	5.674	ug/l	
27) cis-1,2-Dichloroethene	6.82	96	12825	2.449	ug/l	89
44) Trichloroethene	8.82	130	577014	127.594	ug/l	97
49) 1,4-Dioxane	9.19	88	2028	23.199	ug/l #	79
64) Tetrachloroethene	10.63	164	140016	39.953	ug/l	99

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

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