

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN092220\
 Data File : VN063548.D
 Acq On : 22 Sep 2020 21:18
 Operator : JC/MD
 Sample : L4082-15
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 RE131D1-20200916

Manual Integrations
 APPROVED

MMDadoda
 9/23/2020 1:06:29 PM

Quant Time: Sep 23 05:22:06 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N091620W.M
 Quant Title : SW846 8260
 QLast Update : Wed Sep 16 13:05:20 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.63	168	217070	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	401677	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	408517	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	169346	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	181083	49.01	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.02%	
35) Dibromofluoromethane	7.55	113	130663	46.99	ug/l	0.00
Spiked Amount 50.000			Recovery	=	93.98%	
50) Toluene-d8	10.06	98	514483	49.20	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.40%	
62) 4-Bromofluorobenzene	12.38	95	206472	53.87	ug/l	0.00
Spiked Amount 50.000			Recovery	=	107.74%	
Target Compounds						
					Qvalue	
3) Chloromethane	2.04	50	2113	0.649	ug/l	98
9) 1,1,2-Trichlorotrifluoroet	3.73	101	10688	4.419	ug/l #	76
12) 1,1-Dichloroethene	3.70	96	2643	1.079	ug/l	97
19) Methyl tert-butyl Ether	4.99	73	7083m	0.776	ug/l	
27) cis-1,2-Dichloroethene	6.78	96	14285	4.650	ug/l	89
30) Chloroform	7.33	83	8739	1.614	ug/l	91
44) Trichloroethene	8.80	130	679391	229.300	ug/l	99
49) 1,4-Dioxane	9.17	88	1365	25.779	ug/l #	36
64) Tetrachloroethene	10.60	164	38886	13.979	ug/l	96

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

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