

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN112420\
 Data File : VN064806.D
 Acq On : 24 Nov 2020 13:26
 Operator : JC/MD
 Sample : L4829-04
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-4

Manual Integrations
 APPROVED

MMDadoda
 11/25/2020 3:57:15 PM

Quant Time: Nov 25 04:44:56 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N112320W.M
 Quant Title : SW846 8260
 QLast Update : Mon Nov 23 13:54:12 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.63	168	173480	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	315088	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	324568	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	149225	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	104565	51.85	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery	=	103.70%	
35) Dibromofluoromethane	7.55	113	97739	51.84	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery	=	103.68%	
50) Toluene-d8	10.06	98	390012	50.63	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery	=	101.26%	
62) 4-Bromofluorobenzene	12.37	95	156627	56.81	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery	=	113.62%	
Target Compounds						
16) Acetone	3.78	43	2372	3.578	ug/l #	81
30) Chloroform	7.33	83	1699m	0.496	ug/l	
64) Tetrachloroethene	10.60	164	1180	0.374	ug/l #	91
67) Ethyl Benzene	11.48	91	3321	0.285	ug/l	99
84) 1,2,4-Trimethylbenzene	13.01	105	4976	0.570	ug/l	99
95) Naphthalene	15.10	128	24028	2.656	ug/l	99

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

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