

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN022420\
 Data File : VN060216.D
 Acq On : 24 Feb 2020 17:52
 Operator : JC/MD
 Sample : L1671-04
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 E156-B4-022420-GW-BD

Manual Integrations
 APPROVED

MMDadoda
 2/25/2020 9:42:48 AM

Quant Time: Feb 25 00:31:04 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N021220W.M
 Quant Title : SW846 8260
 QLast Update : Wed Feb 12 14:12:21 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.64	168	176332	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.57	114	276682	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.40	117	257418	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.34	152	116603	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.00	65	102841	47.93	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.86%	
35) Dibromofluoromethane	7.57	113	83827	49.29	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.58%	
50) Toluene-d8	10.08	98	331700	49.53	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.06%	
62) 4-Bromofluorobenzene	12.40	95	119406	48.52	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.04%	
Target Compounds						
16) Acetone	3.80	43	2169m	2.317	ug/l	Qvalue
30) Chloroform	7.35	83	3446	1.010	ug/l	95
64) Tetrachloroethene	10.62	164	15954	8.153	ug/l	97

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

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