

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN061820\
 Data File : VN061913.D
 Acq On : 18 Jun 2020 20:28
 Operator : JC/MD
 Sample : L3024-11
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 RO-7

Quant Time: Jun 19 06:33:05 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N061620W.M
 Quant Title : SW846 8260
 QLast Update : Wed Jun 17 01:19:44 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.63	168	138243	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	259537	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	260511	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	98961	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	95316	51.41	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.82%	
35) Dibromofluoromethane	7.55	113	83863	50.51	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.02%	
50) Toluene-d8	10.06	98	333792	51.14	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.28%	
62) 4-Bromofluorobenzene	12.38	95	108844	48.61	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.22%	
Target Compounds						
16) Acetone	3.78	43	12307	22.472	ug/l	99
18) Methyl Acetate	4.28	43	2580	1.636	ug/l #	80
20) Methylene Chloride	4.50	84	9044	5.069	ug/l	87
43) Isopropyl Acetate	8.13	43	29377	9.430	ug/l #	91

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

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