

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN061820\
 Data File : VN061938.D
 Acq On : 19 Jun 2020 7:42
 Operator : JC/MD
 Sample : L2809-04
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 54 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 HD-01-061720

Quant Time: Jun 19 08:06:57 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	143835	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	272650	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	270336	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	99776	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	98989	51.32	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.64%	
35) Dibromofluoromethane	7.55	113	86052	49.33	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.66%	
50) Toluene-d8	10.06	98	349914	51.03	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.06%	
62) 4-Bromofluorobenzene	12.38	95	113028	48.05	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.10%	
Target Compounds						
16) Acetone	3.77	43	9591	16.832	ug/l	98
18) Methyl Acetate	4.28	43	3746	2.283	ug/l #	89
20) Methylene Chloride	4.50	84	6161	3.319	ug/l #	92
43) Isopropyl Acetate	8.13	43	32217	9.844	ug/l #	90

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

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