

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN092820\
 Data File : VN063706.D
 Acq On : 28 Sep 2020 15:33
 Operator : JC/MD
 Sample : L4158-04
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 BLUE-FRAC-TANK

Quant Time: Sep 29 16:23:00 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	221755	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	407532	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	425331	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	176298	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	188217	49.86	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.72%	
35) Dibromofluoromethane	7.55	113	137777	48.84	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.68%	
50) Toluene-d8	10.06	98	531379	50.09	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.18%	
62) 4-Bromofluorobenzene	12.38	95	208657	53.66	ug/l	0.00
Spiked Amount 50.000			Recovery	=	107.32%	
Target Compounds						
16) Acetone	3.78	43	12052	9.934	ug/l	94
20) Methylene Chloride	4.51	84	2303	0.749	ug/l #	69
30) Chloroform	7.33	83	141265	25.547	ug/l	99
47) Bromodichloromethane	9.37	83	21837	4.738	ug/l #	98
60) Dibromochloromethane	10.87	129	3355	1.003	ug/l	88

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

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