

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN111418\
 Data File : VN052316.D
 Acq On : 14 Nov 2018 16:42
 Operator : MD\SY
 Sample : J5164-09 0.5ppb
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 14 Sample Multiplier: 28

Instrument :
 MSVOA_N
 ClientSampleId :
 MDL-WATER-03-QT4-2018

Manual Integrations
 APPROVED

Quant Time: Nov 15 04:39:33 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N111418W.M
 Quant Title : SW846 8260
 QLast Update : Thu Nov 15 03:02:37 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.67	168	254198	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.59	114	450017	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	389875	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.34	152	136921	50.00	ug/l	0.00

System Monitoring Compounds

33) 1,2-Dichloroethane-d4	8.03	65	190875	53.17	ug/l	0.00
Spiked Amount	50.000		Recovery	=	106.34%	
35) Dibromofluoromethane	7.59	113	145373	50.08	ug/l	0.00
Spiked Amount	50.000		Recovery	=	100.16%	
50) Toluene-d8	10.09	98	551259	48.28	ug/l	0.00
Spiked Amount	50.000		Recovery	=	96.56%	
62) 4-Bromofluorobenzene	12.40	95	163191	43.44	ug/l	0.00
Spiked Amount	50.000		Recovery	=	86.88%	

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Target Compounds

						Qvalue
6) Chloroethane	2.71	64	1241	0.678	ug/l	97
8) Diethyl Ether	3.41	74	869m	0.509	ug/l	
12) 1,1-Dichloroethene	3.74	96	1489	0.569	ug/l #	76
15) Acrylonitrile	4.99	53	3209m	2.964	ug/l	
37) Ethyl Acetate	6.94	43	1790m	0.547	ug/l	
81) trans-1,4-Dichloro-2-buten	12.29	75	287m	0.401	ug/l	

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

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