

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN062218\
 Data File : VN049494.D
 Acq On : 22 Jun 2018 16:23
 Operator : MD\SY
 Sample : J3649-01
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 F-6-20-18

Manual Integrations
 APPROVED

MMDadoda
 6/25/2018 4:23:25 PM

Quant Time: Jun 22 23:02:05 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N061618W.M
 Quant Title : SW846 8260
 QLast Update : Sat Jun 16 01:02:01 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.67	168	1530828	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.59	114	2401929	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	2047151	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	630326	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.03	65	1066398	52.52	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.04%	
35) Dibromofluoromethane	7.59	113	908259	46.74	ug/l	0.00
Spiked Amount 50.000			Recovery	=	93.48%	
50) Toluene-d8	10.09	98	3386059	45.76	ug/l	0.00
Spiked Amount 50.000			Recovery	=	91.52%	
62) 4-Bromofluorobenzene	12.40	95	887329	36.27	ug/l	0.00
Spiked Amount 50.000			Recovery	=	72.54%	
Target Compounds						
16) Acetone	3.82	43	94959	19.26	ug/l	95
18) Methyl Acetate	4.33	43	14417m	0.96	ug/l	
25) 2-Butanone	6.85	43	44689	6.75	ug/l	94
27) cis-1,2-Dichloroethene	6.83	96	15035	0.73	ug/l	77
64) Tetrachloroethene	10.63	164	10072	0.48	ug/l	89

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

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