

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN121218\
 Data File : VN052849.D
 Acq On : 13 Dec 2018 1:07
 Operator : MD\SY
 Sample : J6373-10
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 34 Sample Multiplier: 28

Instrument :
 MSVOA_N
 ClientSampleId :
 OILY-WATER

Manual Integrations
 APPROVED

MMDadoda
 12/13/2018 1:36:44 PM

Quant Time: Dec 13 13:39:30 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N121218W.M
 Quant Title : SW846 8260
 QLast Update : Thu Dec 13 01:08:35 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.67	168	1161193	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.59	114	1805919	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	1614860	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	674774	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.03	65	596987	48.35	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.70%	
35) Dibromofluoromethane	7.59	113	508080	50.40	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.80%	
50) Toluene-d8	10.09	98	2222414	50.85	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.70%	
62) 4-Bromofluorobenzene	12.40	95	720817	48.67	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.34%	
Target Compounds						
						Qvalue
11) Tert butyl alcohol	4.78	59	8528m	12.362	ug/l	
16) Acetone	3.79	43	41925791m	15119.279	ug/l	
17) Carbon Disulfide	4.06	76	51398	1.567	ug/l	98
25) 2-Butanone	6.84	43	66581	15.474	ug/l	98
39) Methylcyclohexane	9.08	83	22614	1.108	ug/l	92
51) 4-Methyl-2-Pentanone	9.99	43	490923	52.269	ug/l	100
52) Toluene	10.14	92	25759204m	852.095	ug/l	

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

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