

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN080521\
 Data File : VN068016.D
 Acq On : 5 Aug 2021 15:18
 Operator : JC/MD
 Sample : M3261-03DL 10X
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 41936DL

Manual Integrations
 APPROVED

MMDadoda
 8/6/2021 5:43:00 PM

Quant Time: Aug 05 15:57:29 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N072921W.M
 Quant Title : SW846 8260
 QLast Update : Thu Jul 29 15:02:35 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) Pentafluorobenzene	8.088	168	290965	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	8.971	114	557398	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.747	117	545816	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.678	152	203273	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.437	65	223962	54.371	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery = 108.740%			
35) Dibromofluoromethane	8.024	113	169839	52.283	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery = 104.560%			
50) Toluene-d8	10.443	98	706523	47.349	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery = 94.700%			
62) 4-Bromofluorobenzene	12.734	95	236914	43.854	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery = 87.700%			
Target Compounds						
16) Acetone	4.291	43	7524	5.006	ug/l	# 84
25) 2-Butanone	7.340	43	3076m	1.411	ug/l	
52) Toluene	10.507	92	360248	36.322	ug/l	98
68) m/p-Xylenes	11.956	106	2387	4.327	ug/l	88
79) 2-Chlorotoluene	13.010	91	83123	7.288	ug/l	91
82) 4-Chlorotoluene	13.109	91	84016	7.336	ug/l	94
84) 1,2,4-Trimethylbenzene	13.372	105	7200	0.553	ug/l	98
86) p-Isopropyltoluene	13.619	119	9918	0.797	ug/l	92

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

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