

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN102120\
 Data File : VN064245.D
 Acq On : 22 Oct 2020 4:13
 Operator : JC/MD
 Sample : L4221-14
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 OK-03-102020

Manual Integrations
 APPROVED

MMDadoda
 10/23/2020 2:10:56 PM

Quant Time: Oct 22 06:03:11 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N102120W.M
 Quant Title : SW846 8260
 QLast Update : Wed Oct 21 12:48:40 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.63	168	260370	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	455263	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	466581	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	211643	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	203219	52.56	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.12%	
35) Dibromofluoromethane	7.55	113	149368	49.35	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.70%	
50) Toluene-d8	10.06	98	579854	52.81	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.62%	
62) 4-Bromofluorobenzene	12.38	95	260453	58.51	ug/l	0.00
Spiked Amount 50.000			Recovery	=	117.02%	
Target Compounds						
16) Acetone	3.78	43	29493	26.658	ug/l	93
18) Methyl Acetate	4.28	43	3961m	1.405	ug/l	
20) Methylene Chloride	4.51	84	6451	2.187	ug/l	95
43) Isopropyl Acetate	8.13	43	90368	11.767	ug/l #	87
52) Toluene	10.13	92	1328520	165.718	ug/l	100

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

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