

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN102120\
 Data File : VN064246.D
 Acq On : 22 Oct 2020 4:37
 Operator : JC/MD
 Sample : L4467-02
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 VB-15242

Manual Integrations
 APPROVED

MMDadoda
 10/23/2020 2:10:59 PM

Quant Time: Oct 22 06:06:17 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	7.62	168	258705	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	453410	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	464051	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	216749	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	202473	52.70	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.40%	
35) Dibromofluoromethane	7.55	113	149149	49.48	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.96%	
50) Toluene-d8	10.06	98	578436	52.90	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.80%	
62) 4-Bromofluorobenzene	12.38	95	259643	58.57	ug/l	0.00
Spiked Amount 50.000			Recovery	=	117.14%	
Target Compounds						
16) Acetone	3.78	43	42876	39.004	ug/l	93
18) Methyl Acetate	4.28	43	3565	1.273	ug/l #	78
20) Methylene Chloride	4.51	84	7319	2.497	ug/l #	84
25) 2-Butanone	6.79	43	10208m	5.845	ug/l	
43) Isopropyl Acetate	8.13	43	61208	8.002	ug/l #	88

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

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