

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN102120\
 Data File : VN064249.D
 Acq On : 22 Oct 2020 5:48
 Operator : JC/MD
 Sample : L4468-02
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 72-11977

Manual Integrations
 APPROVED

MMDadoda
 10/23/2020 2:11:04 PM

Quant Time: Oct 22 06:14:24 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	254968	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	451153	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	459665	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	211656	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	199775	52.76	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.52%	
35) Dibromofluoromethane	7.55	113	152950	50.99	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.98%	
50) Toluene-d8	10.06	98	578551	53.17	ug/l	0.00
Spiked Amount 50.000			Recovery	=	106.34%	
62) 4-Bromofluorobenzene	12.38	95	255450	57.91	ug/l	0.00
Spiked Amount 50.000			Recovery	=	115.82%	
Target Compounds						
16) Acetone	3.78	43	21482	19.828	ug/l	94
18) Methyl Acetate	4.30	43	4971m	1.801	ug/l	
20) Methylene Chloride	4.50	84	6553m	2.268	ug/l	
43) Isopropyl Acetate	8.13	43	87186	11.456	ug/l #	89

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

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