

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN103118\
 Data File : VN052088.D
 Acq On : 31 Oct 2018 21:40
 Operator : MD\SY
 Sample : J5728-02
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 24 Sample Multiplier: 28

Instrument :
 MSVOA_N
 ClientSampleId :
 EG-DRUM

Manual Integrations
 APPROVED

MMDadoda
 11/2/2018 1:11:45 PM

Quant Time: Nov 01 03:10:30 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N103018W.M
 Quant Title : SW846 8260
 QLast Update : Wed Oct 31 02:45:22 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.67	168	286557	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.59	114	536565	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	512999	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	208386	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.03	65	250656	54.51	ug/l	0.00
Spiked Amount 50.000			Recovery	=	109.02%	
35) Dibromofluoromethane	7.59	113	205743	61.29	ug/l	0.00
Spiked Amount 50.000			Recovery	=	122.58%	
50) Toluene-d8	10.09	98	812095	60.53	ug/l	0.00
Spiked Amount 50.000			Recovery	=	121.06%	
62) 4-Bromofluorobenzene	12.40	95	275890	56.02	ug/l	0.00
Spiked Amount 50.000			Recovery	=	112.04%	
Target Compounds						
16) Acetone	3.82	43	8680m	8.84	ug/l	Qvalue
18) Methyl Acetate	4.33	43	3457	1.11	ug/l #	79
20) Methylene Chloride	4.55	84	19020	5.09	ug/l	88
43) Isopropyl Acetate	8.17	43	148372	22.84	ug/l	98

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_N\DATA\VN103118\
Data File : VN052088.D
Acq On : 31 Oct 2018 21:40
Operator : MD\SY
Sample : J5728-02
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 24 Sample Multiplier: 28

Instrument :
MSVOA_N
ClientSampleId :
EG-DRUM

Manual Integrations
APPROVED

MMDadoda
11/2/2018 1:11:45 PM

Quant Time: Nov 01 03:10:30 2018
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N103018W.M
Quant Title : SW846 8260
QLast Update : Wed Oct 31 02:45:22 2018
Response via : Initial Calibration

