

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN090619\
 Data File : VN057728.D
 Acq On : 6 Sep 2019 21:11
 Operator : JC/SP
 Sample : K4580-27
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 T12-B4-(0)

Manual Integrations
 APPROVED

MMDadoda

9/10/2019 9:56:47 AM

Quant Time: Sep 07 00:01:29 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N090619W.M
 Quant Title : SW846 8260
 QLast Update : Fri Sep 06 03:46:54 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.65	168	191791	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.58	114	322273	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	299840	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	152499	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.01	65	260599	55.85	ug/l	0.00
Spiked Amount 50.000			Recovery	=	111.70%	
35) Dibromofluoromethane	7.58	113	154655	52.69	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.38%	
50) Toluene-d8	10.09	98	572387	52.71	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.42%	
62) 4-Bromofluorobenzene	12.40	95	211398	46.87	ug/l	0.00
Spiked Amount 50.000			Recovery	=	93.74%	
Target Compounds						
16) Acetone	3.80	43	10819	7.535	ug/l	97
20) Methylene Chloride	4.53	84	4033m	1.299	ug/l	
43) Isopropyl Acetate	8.16	43	53094	6.499	ug/l	100

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN090619\
Data File : VN057728.D
Acq On : 6 Sep 2019 21:11
Operator : JC/SP
Sample : K4580-27
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
T12-B4-(0)

Manual Integrations
APPROVED

MMDadoda

9/10/2019 9:56:47 AM

Quant Time: Sep 07 00:01:29 2019
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N090619W.M
Quant Title : SW846 8260
QLast Update : Fri Sep 06 03:46:54 2019
Response via : Initial Calibration

