

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN102618\
 Data File : VN052030.D
 Acq On : 26 Oct 2018 16:07
 Operator : MD\SY
 Sample : PB114195ZHE#07
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 13 Sample Multiplier: 28

Instrument :
 MSVOA_N
 ClientSampleId :
 PB114195ZHE#07

Manual Integrations
 APPROVED

MMDadoda
 10/29/2018 3:46:39 PM

Quant Time: Oct 27 00:34:42 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N102318W.M
 Quant Title : SW846 8260
 QLast Update : Wed Oct 24 10:12:55 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.67	168	1149272	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.59	114	1676025	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	1606805	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.34	152	749136	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.03	65	590590	57.70	ug/l	0.00
Spiked Amount 50.000			Recovery	=	115.40%	
35) Dibromofluoromethane	7.59	113	615058	57.51	ug/l	0.00
Spiked Amount 50.000			Recovery	=	115.02%	
50) Toluene-d8	10.09	98	2290682	59.27	ug/l	0.00
Spiked Amount 50.000			Recovery	=	118.54%	
62) 4-Bromofluorobenzene	12.40	95	726944	54.72	ug/l	0.00
Spiked Amount 50.000			Recovery	=	109.44%	
Target Compounds						
16) Acetone	3.82	43	13397m	7.944	ug/l	
20) Methylene Chloride	4.55	84	22476	2.272	ug/l	92
43) Isopropyl Acetate	8.17	43	327123	30.382	ug/l	98

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN102618\
Data File : VN052030.D
Acq On : 26 Oct 2018 16:07
Operator : MD\SY
Sample : PB114195ZHE#07
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 13 Sample Multiplier: 28

Instrument :
MSVOA_N
ClientSampleId :
PB114195ZHE#07

Manual Integrations
APPROVED

MMDadoda
10/29/2018 3:46:39 PM

Quant Time: Oct 27 00:34:42 2018
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N102318W.M
Quant Title : SW846 8260
QLast Update : Wed Oct 24 10:12:55 2018
Response via : Initial Calibration

