

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN122118\
 Data File : VN053112.D
 Acq On : 21 Dec 2018 19:49
 Operator : MD\SY
 Sample : J6377-14ME 10X
 Misc : 4.66g/10mL/100uL/5.00mL/MSVOA_N/MEOH
 ALS Vial : 22 Sample Multiplier: 28

Instrument :
 MSVOA_N
 ClientSampleId :
 P001-SD005-0612-01ME

Quant Time: Dec 22 02:37:39 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N121818W.M
 Quant Title : SW846 8260
 QLast Update : Tue Dec 18 13:03:37 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.67	168	1216147	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.59	114	1847174	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	1645923	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	678935	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.03	65	596224	48.31	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.62%	
35) Dibromofluoromethane	7.59	113	478435	56.70	ug/l	0.00
Spiked Amount 50.000			Recovery	=	113.40%	
50) Toluene-d8	10.09	98	2250386	49.68	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.36%	
62) 4-Bromofluorobenzene	12.40	95	751707	47.27	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.54%	
Target Compounds						
16) Acetone	3.82	43	2580	1.066	ug/l	89
68) m/p-Xylenes	11.62	106	42006	1.761	ug/l	97
95) Naphthalene	15.14	128	3548	1.276	ug/l #	91

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_N\DATA\VN122118\
Data File : VN053112.D
Acq On : 21 Dec 2018 19:49
Operator : MD\SY
Sample : J6377-14ME 10X
Misc : 4.66g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 22 Sample Multiplier: 28

Instrument :
MSVOA_N
ClientSampleId :
P001-SD005-0612-01ME

Quant Time: Dec 22 02:37:39 2018
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N121818W.M
Quant Title : SW846 8260
QLast Update : Tue Dec 18 13:03:37 2018
Response via : Initial Calibration

