

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN061918\
 Data File : VN049422.D
 Acq On : 20 Jun 2018 7:54
 Operator : MD\SY
 Sample : J3498-04
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TP-G-S

Quant Time: Jun 20 11:19:40 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N061618W.M
 Quant Title : SW846 8260
 QLast Update : Sat Jun 16 01:02:01 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.67	168	1154736	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.59	114	1853316	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	1505100	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	486085	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.03	65	768262	50.16	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.32%	
35) Dibromofluoromethane	7.59	113	687619	45.86	ug/l	0.00
Spiked Amount 50.000			Recovery	=	91.72%	
50) Toluene-d8	10.09	98	2530533	44.32	ug/l	0.00
Spiked Amount 50.000			Recovery	=	88.64%	
62) 4-Bromofluorobenzene	12.40	95	664376	35.19	ug/l	0.00
Spiked Amount 50.000			Recovery	=	70.38%	
Target Compounds						
16) Acetone	3.82	43	33575	9.029	ug/l	99
20) Methylene Chloride	4.55	84	3021833	194.565	ug/l	92
84) 1,2,4-Trimethylbenzene	13.04	105	47116	1.414	ug/l	98
95) Naphthalene	15.14	128	283338	21.934	ug/l	99

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN061918\
Data File : VN049422.D
Acq On : 20 Jun 2018 7:54
Operator : MD\SY
Sample : J3498-04
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 43 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
TP-G-S

Quant Time: Jun 20 11:19:40 2018
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N061618W.M
Quant Title : SW846 8260
QLast Update : Sat Jun 16 01:02:01 2018
Response via : Initial Calibration

