

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN041519\
 Data File : VN055060.D
 Acq On : 15 Apr 2019 23:05
 Operator : JC/SP
 Sample : K2390-04
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 5124

Quant Time: Apr 16 07:21:19 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N041519W.M
 Quant Title : SW846 8260
 QLast Update : Mon Apr 15 12:28:28 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.66	168	522337	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.58	114	845297	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	715130	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.34	152	231080	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.02	65	335185	48.28	ug/l	0.00
Spiked Amount	50.000		Recovery	=	96.56%	
35) Dibromofluoromethane	7.59	113	272734	48.28	ug/l	0.00
Spiked Amount	50.000		Recovery	=	96.56%	
50) Toluene-d8	10.09	98	1012631	50.08	ug/l	0.00
Spiked Amount	50.000		Recovery	=	100.16%	
62) 4-Bromofluorobenzene	12.40	95	298436	45.40	ug/l	0.00
Spiked Amount	50.000		Recovery	=	90.80%	
Target Compounds						
16) Acetone	3.82	43	11358	6.027	ug/l	97
20) Methylene Chloride	4.54	84	458128	69.155	ug/l	98
43) Isopropyl Acetate	8.17	43	12467	1.247	ug/l #	89
95) Naphthalene	15.13	128	4487	3.588	ug/l #	94

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN041519\
Data File : VN055060.D
Acq On : 15 Apr 2019 23:05
Operator : JC/SP
Sample : K2390-04
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 29 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
5124

Quant Time: Apr 16 07:21:19 2019
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N041519W.M
Quant Title : SW846 8260
QLast Update : Mon Apr 15 12:28:28 2019
Response via : Initial Calibration

