

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN091119\
 Data File : VN057906.D
 Acq On : 11 Sep 2019 8:05
 Operator : JC/SP
 Sample : K4624-19
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 49 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PAR-11-E4-(4)

Quant Time: Sep 13 05:43:33 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N091019W.M
 Quant Title : SW846 8260
 QLast Update : Tue Sep 10 03:02:13 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.65	168	259989	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.58	114	461783	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	420922	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	207484	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.01	65	334535	63.53	ug/l	0.00
Spiked Amount 50.000			Recovery	=	127.06%	
35) Dibromofluoromethane	7.58	113	222960	52.31	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.62%	
50) Toluene-d8	10.09	98	858610	51.89	ug/l	0.00
Spiked Amount 50.000			Recovery	=	103.78%	
62) 4-Bromofluorobenzene	12.41	95	281364	44.62	ug/l	0.00
Spiked Amount 50.000			Recovery	=	89.24%	
Target Compounds						
16) Acetone	3.80	43	18231	10.561	ug/l	96
18) Methyl Acetate	4.31	43	5917	1.286	ug/l #	72
20) Methylene Chloride	4.53	84	5563	0.473	ug/l #	80
43) Isopropyl Acetate	8.16	43	75568	8.121	ug/l #	90

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

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN091119\
Data File : VN057906.D
Acq On : 11 Sep 2019 8:05
Operator : JC/SP
Sample : K4624-19
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 49 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
PAR-11-E4-(4)

Quant Time: Sep 13 05:43:33 2019
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N091019W.M
Quant Title : SW846 8260
QLast Update : Tue Sep 10 03:02:13 2019
Response via : Initial Calibration

