

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN080819\
 Data File : VN057132.D
 Acq On : 8 Aug 2019 15:09
 Operator : JC/SP
 Sample : K4204-08
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 13967

Quant Time: Aug 09 04:05:03 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N080719W.M
 Quant Title : SW846 8260
 QLast Update : Thu Aug 08 06:49:25 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.05	168	1402493	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.01	114	1987175	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.90	117	1668952	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.87	152	480076	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.44	65	657487	45.58	ug/l	0.00
Spiked Amount 50.000			Recovery	=	91.16%	
35) Dibromofluoromethane	6.99	113	598586	48.00	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.00%	
50) Toluene-d8	9.56	98	2308072	48.38	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.76%	
62) 4-Bromofluorobenzene	11.92	95	666654	43.37	ug/l	0.00
Spiked Amount 50.000			Recovery	=	86.74%	
Target Compounds						
16) Acetone	3.47	43	970870	222.884	ug/l	99
18) Methyl Acetate	3.90	43	34440	2.023	ug/l	100
20) Methylene Chloride	4.10	84	35736	2.486	ug/l	96
25) 2-Butanone	6.20	43	339001	50.938	ug/l	100
30) Chloroform	6.76	83	232287	9.651	ug/l	98
51) 4-Methyl-2-Pentanone	9.45	43	131807	9.840	ug/l	99

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

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