

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN081219\
 Data File : VN057236.D
 Acq On : 12 Aug 2019 23:44
 Operator : JC/SP
 Sample : K4231-01
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 REACTOR-3-E-(3)

Quant Time: Aug 13 03:45:58 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	1293977	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.01	114	1799381	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.91	117	1476825	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.88	152	373444	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.44	65	608213	45.70	ug/l	0.00
Spiked Amount 50.000			Recovery	=	91.40%	
35) Dibromofluoromethane	6.99	113	546133	48.37	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.74%	
50) Toluene-d8	9.56	98	2091462	48.41	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.82%	
62) 4-Bromofluorobenzene	11.92	95	538931	38.72	ug/l	0.00
Spiked Amount 50.000			Recovery	=	77.44%	
Target Compounds						
16) Acetone	3.48	43	21753	5.413	ug/l	99
18) Methyl Acetate	3.90	43	29529	1.833	ug/l	100
20) Methylene Chloride	4.10	84	14494	1.093	ug/l	97
43) Isopropyl Acetate	7.57	43	167944	8.058	ug/l #	75

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

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