

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN081419\
 Data File : VN057286.D
 Acq On : 14 Aug 2019 14:33
 Operator : JC/SP
 Sample : K4231-07
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 12 Sample Multiplier: 1

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

Quant Time: Aug 14 15:03:22 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N081319W.M
 Quant Title : SW846 8260
 QLast Update : Wed Aug 14 11:11:44 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.66	168	1094747	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.58	114	1530386	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.40	117	1269140	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	311021	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.02	65	529381	47.94	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.88%	
35) Dibromofluoromethane	7.58	113	467495	47.84	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.68%	
50) Toluene-d8	10.09	98	1783125	50.03	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.06%	
62) 4-Bromofluorobenzene	12.40	95	467768	38.27	ug/l	0.00
Spiked Amount 50.000			Recovery	=	76.54%	
Target Compounds						
16) Acetone	3.81	43	52616	13.897	ug/l	97
18) Methyl Acetate	4.32	43	28108	1.757	ug/l	98
20) Methylene Chloride	4.54	84	68148	6.162	ug/l	96
43) Isopropyl Acetate	8.16	43	150640	8.505	ug/l #	82

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

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