

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN081919\
 Data File : VN057388.D
 Acq On : 19 Aug 2019 10:16
 Operator : JC/SP
 Sample : K4231-03
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 REACTOR-3-F-(2)

Quant Time: Aug 20 07:37:32 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N081519W.M
 Quant Title : SW846 8260
 QLast Update : Thu Aug 15 11:33:14 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.66	168	1046634	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.58	114	1435536	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.40	117	1195801	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	285263	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.02	65	473814	41.64	ug/l	0.00
Spiked Amount 50.000			Recovery	=	83.28%	
35) Dibromofluoromethane	7.58	113	443624	47.14	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.28%	
50) Toluene-d8	10.09	98	1668032	48.46	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.92%	
62) 4-Bromofluorobenzene	12.40	95	424152	34.87	ug/l	0.00
Spiked Amount 50.000			Recovery	=	69.74%	
Target Compounds						
16) Acetone	3.81	43	22353	5.663	ug/l	97
20) Methylene Chloride	4.55	84	17969	1.627	ug/l	89
43) Isopropyl Acetate	8.16	43	78122	3.056	ug/l	93
93) 1,2,4-Trichlorobenzene	14.91	180	2795	6.026	ug/l	90

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

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