

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN081219\
 Data File : VN057237.D
 Acq On : 13 Aug 2019 00:08
 Operator : JC/SP
 Sample : K4231-02
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 REACTOR-3-F-(0-3)

Quant Time: Aug 13 03:47:48 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	1268032	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.01	114	1775422	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.91	117	1441808	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.88	152	359626	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.44	65	595821	45.69	ug/l	0.00
Spiked Amount 50.000			Recovery	=	91.38%	
35) Dibromofluoromethane	6.99	113	535140	48.03	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.06%	
50) Toluene-d8	9.56	98	2038571	47.83	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.66%	
62) 4-Bromofluorobenzene	11.92	95	515014	37.50	ug/l	0.00
Spiked Amount 50.000			Recovery	=	75.00%	
Target Compounds						
8) Diethyl Ether	3.15	74	7985	1.142	ug/l	99
16) Acetone	3.48	43	27574	7.001	ug/l	91
18) Methyl Acetate	3.90	43	29464	1.879	ug/l	98
20) Methylene Chloride	4.10	84	24068	1.852	ug/l #	95
43) Isopropyl Acetate	7.57	43	145060	7.054	ug/l #	79

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

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