

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN042419\
 Data File : VN055230.D
 Acq On : 24 Apr 2019 16:10
 Operator : JC/SP
 Sample : K2200-10
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 EO-02-042319-B

Quant Time: Apr 25 06:55:57 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N042219W.M
 Quant Title : SW846 8260
 QLast Update : Tue Apr 23 05:52:53 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.67	168	382544	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.59	114	653165	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	608558	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.34	152	191742	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.03	65	262504	50.70	ug/l	0.00
Spiked Amount	50.000		Recovery	=	101.40%	
35) Dibromofluoromethane	7.59	113	214928	48.44	ug/l	0.00
Spiked Amount	50.000		Recovery	=	96.88%	
50) Toluene-d8	10.09	98	804133	48.91	ug/l	0.00
Spiked Amount	50.000		Recovery	=	97.82%	
62) 4-Bromofluorobenzene	12.40	95	261341	50.20	ug/l	0.00
Spiked Amount	50.000		Recovery	=	100.40%	
Target Compounds						
					Qvalue	
16) Acetone	3.82	43	13590	7.940	ug/l	100
17) Carbon Disulfide	4.06	76	4513	0.410	ug/l #	94
18) Methyl Acetate	4.34	43	9998	1.627	ug/l #	77
20) Methylene Chloride	4.55	84	152936	31.079	ug/l	98
30) Chloroform	7.38	83	5183	0.653	ug/l #	82
43) Isopropyl Acetate	8.18	43	21373	2.698	ug/l #	83

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

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