

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN042419\
 Data File : VN055228.D
 Acq On : 24 Apr 2019 15:19
 Operator : JC/SP
 Sample : K2200-12
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 15 Sample Multiplier: 1

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

Quant Time: Apr 25 06:42:48 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	366025	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.59	114	641496	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	585762	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.34	152	181652	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.03	65	256860	51.85	ug/l	0.00
Spiked Amount	50.000		Recovery	=	103.70%	
35) Dibromofluoromethane	7.59	113	209652	48.11	ug/l	0.00
Spiked Amount	50.000		Recovery	=	96.22%	
50) Toluene-d8	10.09	98	791108	48.99	ug/l	0.00
Spiked Amount	50.000		Recovery	=	97.98%	
62) 4-Bromofluorobenzene	12.40	95	246926	48.29	ug/l	0.00
Spiked Amount	50.000		Recovery	=	96.58%	
Target Compounds						
16) Acetone	3.83	43	12958	7.912	ug/l	97
17) Carbon Disulfide	4.06	76	4760	0.452	ug/l	99
18) Methyl Acetate	4.34	43	9208	1.523	ug/l #	84
20) Methylene Chloride	4.55	84	12858	2.731	ug/l	95
30) Chloroform	7.37	83	9702	1.278	ug/l	96
43) Isopropyl Acetate	8.17	43	21008	2.700	ug/l #	73

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

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