

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN012819\
 Data File : VN053537.D
 Acq On : 28 Jan 2019 11:27
 Operator : JC/SP
 Sample : K1256-01
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 6 Sample Multiplier: 28

Instrument :
 MSVOA_N
 ClientSampleId :
 30093

Manual Integrations
 APPROVED

MMDadoda
 1/29/2019 9:52:24 AM

Quant Time: Jan 29 07:05:21 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N011519W.M
 Quant Title : SW846 8260
 QLast Update : Wed Jan 16 03:10:53 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.67	168	823991	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.59	114	1312167	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	1206301	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	556110	50.00	ug/l	0.00

System Monitoring Compounds

33) 1,2-Dichloroethane-d4	8.03	65	431019	52.92	ug/l	0.00
Spiked Amount	50.000		Recovery	=	105.84%	
35) Dibromofluoromethane	7.59	113	415328	55.46	ug/l	0.00
Spiked Amount	50.000		Recovery	=	110.92%	
50) Toluene-d8	10.09	98	1591507	50.25	ug/l	0.00
Spiked Amount	50.000		Recovery	=	100.50%	
62) 4-Bromofluorobenzene	12.40	95	552184	49.72	ug/l	0.00
Spiked Amount	50.000		Recovery	=	99.44%	

Target Compounds

					Qvalue
16) Acetone	3.82	43	21544	11.275	ug/l 96
20) Methylene Chloride	4.55	84	9196	1.054	ug/l # 92
25) 2-Butanone	6.84	43	20001	6.792	ug/l 99
40) Benzene	8.04	78	37923	1.044	ug/l 96
51) 4-Methyl-2-Pentanone	9.99	43	70967	10.516	ug/l 100
52) Toluene	10.16	92	131540	5.835	ug/l 99
68) m/p-Xylenes	11.62	106	147257	8.963	ug/l 100
69) o-Xylene	11.95	106	184742	11.572	ug/l 100
80) 1,3,5-Trimethylbenzene	12.73	105	203383	6.138	ug/l 99
84) 1,2,4-Trimethylbenzene	13.04	105	230566	6.901	ug/l 99
86) p-Isopropyltoluene	13.29	119	52516	1.507	ug/l 98
88) 1,4-Dichlorobenzene	13.36	146	86947	4.592	ug/l 96

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

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