

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN071918\
 Data File : VN049911.D
 Acq On : 19 Jul 2018 12:58
 Operator : MD\SY
 Sample : J4043-03
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 SD-BASELINE-WATER

Manual Integrations
 APPROVED

MMDadoda
 7/20/2018 11:08:03 AM

Quant Time: Jul 20 01:43:46 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N071818W.M
 Quant Title : SW846 8260
 QLast Update : Thu Jul 19 01:58:39 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.67	168	601893	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.59	114	970595	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	821471	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	349717	50.00	ug/l	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
33) 1,2-Dichloroethane-d4	8.03	65	398152	48.13	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.26%	
35) Dibromofluoromethane	7.59	113	358739	44.03	ug/l	0.00
Spiked Amount 50.000			Recovery	=	88.06%	
50) Toluene-d8	10.09	98	1316378	44.34	ug/l	0.00
Spiked Amount 50.000			Recovery	=	88.68%	
62) 4-Bromofluorobenzene	12.40	95	395980	38.73	ug/l	0.00
Spiked Amount 50.000			Recovery	=	77.46%	

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
11) Tert butyl alcohol	4.79	59	11323m	19.21	ug/l	
16) Acetone	3.82	43	120238	56.14	ug/l	99
17) Carbon Disulfide	4.05	76	160488	8.22	ug/l	100
18) Methyl Acetate	4.33	43	7274	0.26	ug/l #	85
25) 2-Butanone	6.84	43	36593	11.95	ug/l	96
51) 4-Methyl-2-Pentanone	9.99	43	13042	1.80	ug/l	91
59) 2-Hexanone	10.75	43	7024	1.44	ug/l	85
95) Naphthalene	15.13	128	6077	4.83	ug/l	99

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

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