

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN091118\
 Data File : VN051123.D
 Acq On : 11 Sep 2018 13:33
 Operator : MD\SY
 Sample : J4849-09
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 EP1-MW-G

Quant Time: Sep 12 02:31:22 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N091018W.M
 Quant Title : SW846 8260
 QLast Update : Tue Sep 11 02:27:09 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.67	168	660855	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.59	114	1016701	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	909078	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.34	152	417295	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	385950	52.48	ug/l	0.00
Spiked Amount			Recovery	=	104.96%	
35) Dibromofluoromethane	7.59	113	377734	49.44	ug/l	0.00
Spiked Amount			Recovery	=	98.88%	
50) Toluene-d8	10.09	98	1405625	48.49	ug/l	0.00
Spiked Amount			Recovery	=	96.98%	
62) 4-Bromofluorobenzene	12.40	95	475853	49.62	ug/l	0.00
Spiked Amount			Recovery	=	99.24%	

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Chloromethane	2.06	50	35490	3.81	ug/l	99
17) Carbon Disulfide	4.05	76	13535	0.76	ug/l	99
18) Methyl Acetate	4.34	43	8062	1.37	ug/l	96
27) cis-1,2-Dichloroethene	6.83	96	16561	1.97	ug/l	92
39) Methylcyclohexane	9.08	83	5692	0.46	ug/l #	90
67) Ethyl Benzene	11.51	91	21230	0.57	ug/l	99
69) o-Xylene	11.95	106	8910	0.65	ug/l	57
73) Isopropylbenzene	12.25	105	3226302	92.19	ug/l	100
78) n-propylbenzene	12.59	91	7008664	182.20	ug/l	99
83) tert-Butylbenzene	12.99	119	595585	23.84	ug/l	96
84) 1,2,4-Trimethylbenzene	13.04	105	1249931	44.93	ug/l	99
85) sec-Butylbenzene	13.17	105	2882957	90.00	ug/l	81
86) p-Isopropyltoluene	13.29	119	1146884	42.22	ug/l	100
89) n-Butylbenzene	13.61	91	1316276	62.27	ug/l #	75

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

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