

Data Path : W:\HPCHEM1\MSVOA N\DATA\VN092115\
 Data File : VN027338.D
 Acq On : 21 Sep 2015 18:41
 Operator : MD\FY
 Sample : G3692-01
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PT-VOA-WP

Manual Integrations
 APPROVED

MMDadoda
 9/22/2015 2:28:23 PM

Quant Time: Sep 22 15:06:53 2015
 Quant Method : W:\HPCHEM1\MSVOA_N\METHODS\82N092115W.M
 Quant Title : SW846 8260
 QLast Update : Mon Sep 21 13:12:32 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.75	168	542165	50.00	ug/l	0.00
35) 1,4-Difluorobenzene	8.68	114	1013466	50.00	ug/l	0.00
64) Chlorobenzene-d5	11.52	117	870795	50.00	ug/l	0.00
73) 1,4-Dichlorobenzene-d4	13.47	152	337180	50.00	ug/l	0.00

System Monitoring Compounds

34) 1,2-Dichloroethane-d4	8.12	65	379102	35.72	ug/l	0.00
Spiked Amount	50.000		Recovery	=	71.44%	
36) Dibromofluoromethane	7.68	113	236353	35.74	ug/l	0.00
Spiked Amount	50.000		Recovery	=	71.48%	
51) Toluene-d8	10.19	98	915194	35.80	ug/l	0.00
Spiked Amount	50.000		Recovery	=	71.60%	
63) 4-Bromofluorobenzene	12.52	95	324830	34.97	ug/l	0.00
Spiked Amount	50.000		Recovery	=	69.94%	

Target Compounds

						Qvalue
6) Bromomethane	2.54	94	201619	68.84	ug/l	100
7) Chloroethane	2.69	64	284885	62.30	ug/l	100
8) Trichlorofluoromethane	3.02	101	1030905	98.71	ug/l	98
13) 1,1-Dichloroethene	3.77	96	661440	106.84	ug/l	95
17) Acetone	3.83	43	271552	94.05	ug/l	96
21) Methylene Chloride	4.60	84	531792	69.17	ug/l	97
22) trans-1,2-Dichloroethene	5.12	96	756515	111.14	ug/l	95
26) 2-Butanone	6.91	43	752305	198.96	ug/l	97
31) Chloroform	7.46	83	478562	33.30	ug/l	99
33) 1,1,1-Trichloroethane	7.67	97	198220	15.56	ug/l	98
39) Carbon Tetrachloride	7.88	117	569996	55.45	ug/l	100
41) Benzene	8.14	78	850815	26.84	ug/l	99
43) 1,2-Dichloroethane	8.22	62	1555049	120.32	ug/l	95
45) Trichloroethene	8.93	130	536573	84.49	ug/l	93
46) 1,2-Dichloropropane	9.22	63	732202	82.52	ug/l	98
47) Dibromomethane	9.31	93	178933	33.60	ug/l	92
48) Bromodichloromethane	9.50	83	496281	43.94	ug/l	99
52) 4-Methyl-2-Pentanone	10.07	43	211696	25.53	ug/l	99
54) t-1,3-Dichloropropene	10.47	75	360563	29.91	ug/l	99
55) cis-1,3-Dichloropropene	9.93	75	577455	43.60	ug/l	94
56) 1,1,2-Trichloroethane	10.66	97	254981	36.46	ug/l	98
60) 2-Hexanone	10.84	43	642851	114.08	ug/l	99
62) 1,2-Dibromoethane	11.11	107	257842	38.77	ug/l	99
65) Tetrachloroethene	10.74	164	364153	70.26	ug/l	95
66) Chlorobenzene	11.54	112	550460	29.48	ug/l	97
67) 1,1,1,2-Tetrachloroethane	11.62	131	121069	19.32	ug/l	98
68) Ethyl Benzene	11.62	91	1438658	40.85	ug/l	99
69) m/p-Xylenes	11.73	106	1645877	134.81	ug/l	91
71) Styrene	12.08	104	1327465	67.59	ug/l	94
72) Bromoform	12.24	173	137826	32.79	ug/l	99
76) 1,1,2,2-Tetrachloroethane	12.61	83	1096066	115.27	ug/l	99
77) 1,2,3-Trichloropropane	12.66	75	203695	24.06	ug/l	# 1
85) 1,2,4-Trimethylbenzene	13.16	105	2048281	75.43	ug/l	95
89) 1,4-Dichlorobenzene	13.49	146	273772	23.72	ug/l	93

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
92) 1,2-Dichlorobenzene	13.78	146	344277	30.22	ug/l	96
93) 1,2-Dibromo-3-Chloropropan	14.40	75	158711	85.17	ug/l	88
94) 1,2,4-Trichlorobenzene	15.06	180	660996	112.25	ug/l	98
95) Hexachlorobutadiene	15.16	225	475076	147.25	ug/l	98
96) Naphthalene	15.29	128	1302579	70.51	ug/l	100

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

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