

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN092818\
 Data File : VN051622.D
 Acq On : 28 Sep 2018 18:23
 Operator : MD\SY
 Sample : J4870-02DL 2X
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 18 Sample Multiplier: 28

Instrument :
 MSVOA_N
 ClientSampleId :
 PT-VOA-WPDL

Manual Integrations
 APPROVED

MMDadoda
 10/3/2018 5:27:46 PM

Quant Time: Sep 29 05:21:32 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA N\METHODS\624N092818W.M
 Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS
 QLast Update : Sat Sep 29 01:07:29 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	7.20	128	71419	30.00	ug/l	0.00
28) 1,4-Difluorobenzene	8.59	114	381039	30.00	ug/l	0.00
57) Chlorobenzene-d5	11.41	117	340812	30.00	ug/l	0.00

System Monitoring Compounds

27) 1,2-Dichloroethane-d4	8.03	65	148021	29.77	ug/l	0.00
Spiked Amount	30.000	Range	50 - 169	Recovery	=	99.23%
60) 4-Bromofluorobenzene	12.40	95	142415	27.17	ug/l	0.00
Spiked Amount	30.000	Range	56 - 143	Recovery	=	90.57%
63) Toluene-d8	10.09	98	460272	28.62	ug/l	0.00
Spiked Amount	30.000	Range	66 - 137	Recovery	=	95.40%

Target Compounds

					Ovalue
3) Chloromethane	2.06	50	4003	0.571	ug/l 90
4) Vinyl Chloride	2.19	62	184053	26.750	ug/l 95
5) Bromomethane	2.57	94	156837	43.883	ug/l 98
6) Chloroethane	2.71	64	74294	21.524	ug/l 97
11) Methyl Iodide	3.95	142	5773m	1.040	ug/l
15) Acetone	3.82	58	19343	48.402	ug/l 95
18) Methylene Chloride	4.55	84	63875	12.884	ug/l 94
19) trans-1,2-Dichloroethene	5.05	96	47098	10.289	ug/l 89
21) 1,1-Dichloroethane	5.85	63	203497	22.436	ug/l 99
24) Methyl tert-Butyl Ether	5.06	73	519458	45.833	ug/l 100
25) Chloroform	7.37	83	210513	24.274	ug/l 99
30) 2-Butanone	6.84	43	117006	66.166	ug/l # 95
32) 1,1,1-Trichloroethane	7.57	97	60478	8.284	ug/l 98
34) Benzene	8.04	78	680075	35.279	ug/l # 95
37) Trichloroethene	8.84	130	174540	35.266	ug/l 93
39) 1,2-Dichloropropane	9.12	63	190635	36.091	ug/l 98
40) Dibromomethane	9.21	93	161304	54.974	ug/l 91
49) cis-1,3-Dichloropropene	9.84	75	334488	44.755	ug/l 100
53) Dibromochloromethane	10.90	129	169324	35.853	ug/l 99
54) 1,2-Dibromoethane	11.01	107	190468	47.871	ug/l 99
56) Bromoform	12.13	173	150777	48.114	ug/l 98
58) 4-Methyl-2-Pentanone	9.99	43	204538	52.860	ug/l 97
59) 2-Hexanone	10.75	43	127860	50.228	ug/l 99
61) Tetrachloroethene	10.63	164	129694	26.878	ug/l 98
62) Toluene	10.16	91	485552	23.235	ug/l 99
64) Chlorobenzene	11.44	112	259114	20.507	ug/l 99
65) 1,1,1,2-Tetrachloroethane	11.51	131	337996	70.055	ug/l 98
66) Ethyl Benzene	11.51	91	159513	7.458	ug/l 96
67) m/p-Xylenes	11.62	106	148736	18.044	ug/l 97
68) o-Xylene	11.95	106	291450	37.039	ug/l 97
69) Styrene	11.97	104	685190	55.478	ug/l 95
71) 1,1,2,2-Tetrachloroethane	12.50	83	295379	58.167	ug/l 99
72) 1,2,3-Trichloropropane	12.55	75	252827	58.233	ug/l 1
76) 1,3,5-Trimethylbenzene	12.73	105	423405	24.141	ug/l 100
80) 1,2,4-Trimethylbenzene	13.04	105	705207	40.257	ug/l 100
83) 1,3-Dichlorobenzene	13.28	146	504219	55.398	ug/l 99
84) 1,4-Dichlorobenzene	13.36	146	362691	42.392	ug/l 99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
85) n-Butylbenzene	13.62	91	3166	0.215	ug/l	89
86) Hexachloroethane	13.87	117	3211	0.895	ug/l	96
87) 1,2-Dichlorobenzene	13.65	146	267647	30.060	ug/l	99
88) 1,2-Dibromo-3-Chloropropan	14.27	75	27844	37.620	ug/l	93
89) 1,2,4-Trichlorobenzene	14.91	180	258823	62.107	ug/l	97
90) Hexachlorobutadiene	15.01	225	336155	98.943	ug/l	97
91) Naphthalene	15.13	128	522454	50.937	ug/l	99
92) 1,2,3-Trichlorobenzene	15.31	180	6601	1.575	ug/l	96

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

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