

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN031621\
 Data File : VN066212.D
 Acq On : 16 Mar 2021 13:18
 Operator : JC/MD
 Sample : IBLK
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 IBLK

Quant Time: Mar 17 05:21:36 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N031621W.M
 Quant Title : SW846 8260
 QLast Update : Wed Mar 17 05:05:15 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) Pentafluorobenzene	7.590	168	118699	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.515	114	202013	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.347	117	180210	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.286	152	59042	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.949	65	85991	50.28	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery = 100.56%			
35) Dibromofluoromethane	7.512	113	64328	49.92	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery = 99.84%			
50) Toluene-d8	10.028	98	251119	50.39	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery = 100.78%			
62) 4-Bromofluorobenzene	12.345	95	75257	41.90	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery = 83.80%			
Target Compounds						Qvalue
2) Dichlorodifluoromethane	1.821	85	566	0.48	ug/l	84
3) Chloromethane	2.017	50	5056	2.39	ug/l	99
4) Vinyl Chloride	2.145	62	1210	0.72	ug/l	91
7) Trichlorofluoromethane	2.969	101	861	0.44	ug/l #	77
17) Carbon Disulfide	3.969	76	12058	3.83	ug/l	98
25) 2-Butanone	6.753	43	4268	3.34	ug/l #	74
30) Chloroform	7.292	83	936	0.38	ug/l #	82
39) Methylcyclohexane	9.008	83	1152	0.52	ug/l #	81
51) 4-Methyl-2-Pentanone	9.920	43	10322	3.78	ug/l	95
53) t-1,3-Dichloropropene	10.320	75	1954	0.80	ug/l #	83
59) 2-Hexanone	10.695	43	15121	8.38	ug/l	88
64) Tetrachloroethene	10.567	164	1352	0.90	ug/l #	86
65) Chlorobenzene	11.377	112	3090	0.85	ug/l	93
67) Ethyl Benzene	11.454	91	4146	0.65	ug/l #	86
68) m/p-Xylenes	11.564	106	3034	1.28	ug/l	92
69) o-Xylene	11.886	106	1130	0.49	ug/l	89
70) Styrene	11.910	104	2551	2.72	ug/l	98
73) Isopropylbenzene	12.192	105	4069	0.98	ug/l	99
78) n-propylbenzene	12.538	91	7343	1.54	ug/l	99
79) 2-Chlorotoluene	12.616	91	3593	1.24	ug/l	74
80) 1,3,5-Trimethylbenzene	12.677	105	4484	1.31	ug/l	99
82) 4-Chlorotoluene	12.718	91	4050	1.74	ug/l	95
83) tert-Butylbenzene	12.935	119	3905	1.30	ug/l	97
84) 1,2,4-Trimethylbenzene	12.986	105	5467	1.63	ug/l	94
85) sec-Butylbenzene	13.120	105	7372	1.92	ug/l	100
86) p-Isopropyltoluene	13.235	119	6997	2.08	ug/l	88
87) 1,3-Dichlorobenzene	13.227	146	4355	2.37	ug/l	92
88) 1,4-Dichlorobenzene	13.305	146	5281	2.74	ug/l	85
89) n-Butylbenzene	13.562	91	13856	6.16	ug/l	98
91) 1,2-Dichlorobenzene	13.592	146	5091	2.81	ug/l	92
93) 1,2,4-Trichlorobenzene	14.852	180	14116	11.92	ug/l	98
94) Hexachlorobutadiene	14.949	225	5342	9.28	ug/l	100
95) Naphthalene	15.070	128	108363	31.30	ug/l	100
96) 1,2,3-Trichlorobenzene	15.249	180	17437	14.79	ug/l	97

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

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