

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE011120\
 Data File : BE101031.D
 Acq On : 10 Jan 2020 20:05
 Operator : JU
 Sample : K3591-07
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 LOD-MDL-WATER-01-QT3-2019

Quant Time: Jan 11 01:45:30 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE010820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jan 08 18:36:25 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	3124	0.40	ng	0.00
7) Naphthalene-d8	10.51	136	13796	0.40	ng	0.01
13) Acenaphthene-d10	14.36	164	8833	0.40	ng	0.00
19) Phenanthrene-d10	17.09	188	19093	0.40	ng	0.00
27) Chrysene-d12	21.28	240	14126	0.40	ng	0.00
34) Perylene-d12	23.70	264	18277	0.40	ng	0.00

System Monitoring Compounds						
4) 2-Fluorophenol	5.34	112	2329	0.33	ng	0.00
5) Phenol-d6	6.92	99	2301	0.26	ng	0.00
8) Nitrobenzene-d5	8.90	82	2356	0.24	ng	0.03
11) 2-Methylnaphthalene-d10	12.12	152	8482	0.32	ng	0.01
14) 2,4,6-Tribromophenol	15.85	330	534	0.21	ng	0.00
15) 2-Fluorobiphenyl	13.00	172	12597	0.38	ng	0.00
25) Fluoranthene-d10	19.13	212	76114	0.29	ng	0.00
29) Terphenyl-d14	19.74	244	14845	0.43	ng	0.00

Target Compounds				Qvalue		
2) 1,4-Dioxane	3.29	88	1393m	0.088	ng	
3) n-Nitrosodimethylamine	3.67	42	422m	0.097	ng	
6) bis(2-Chloroethyl)ether	7.17	93	659	0.062	ng	# 88
9) Naphthalene	10.56	128	13523	0.088	ng	# 96
10) Hexachlorobutadiene	10.85	225	610	0.094	ng	96
12) 2-Methylnaphthalene	12.19	142	2035	0.088	ng	95
16) Acenaphthylene	14.08	152	2501	0.070	ng	99
17) Acenaphthene	14.43	154	1955	0.076	ng	97
18) Fluorene	15.41	166	2480	0.076	ng	98
20) 4-Bromophenyl-phenylether	16.31	248	647	0.069	ng	# 84
21) Hexachlorobenzene	16.41	284	869	0.085	ng	98
22) Pentachlorophenol	16.76	266	259	0.322	ng	93
23) Phenanthrene	17.15	178	4013	0.077	ng	99
24) Anthracene	17.24	178	4002	0.081	ng	96
26) Fluoranthene	19.16	202	4989	0.076	ng	98
28) Pyrene	19.52	202	4865	0.093	ng	98
30) Benzo(a)anthracene	21.26	228	2292	0.057	ng	93
31) Chrysene	21.32	228	6400	0.107	ng	97
32) Bis(2-ethylhexyl)phthalate	21.20	149	5138	0.076	ng	# 99
33) Indeno(1,2,3-cd)pyrene	26.24	276	3654m	0.077	ng	
35) Benzo(b)fluoranthene	22.96	252	3118	0.061	ng	# 89
36) Benzo(k)fluoranthene	23.01	252	5747	0.077	ng	# 79
37) Benzo(a)pyrene	23.59	252	3507	0.071	ng	# 77
38) Dibenzo(a,h)anthracene	26.27	278	2965m	0.065	ng	
39) Benzo(g,h,i)perylene	27.02	276	3996	0.071	ng	94

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

