

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN031920\
 Data File : BN010302.D
 Acq On : 19 Mar 2020 23:27
 Operator : CG/JU
 Sample : L1969-03
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 OUTFALL001-200317

Quant Time: Mar 20 01:49:52 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN031920.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 19 15:26:45 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	4415	0.40	ng	0.00
7) Naphthalene-d8	10.38	136	15532	0.40	ng	0.00
13) Acenaphthene-d10	14.24	164	9355	0.40	ng	-0.01
19) Phenanthrene-d10	16.99	188	22166	0.40	ng	0.00
27) Chrysene-d12	21.19	240	24434	0.40	ng	0.00
34) Perylene-d12	23.36	264	25806	0.40	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Fluorophenol	5.25	112	1277	0.15	ng	0.00
5) Phenol-d6	6.80	99	976	0.09	ng	0.00
8) Nitrobenzene-d5	8.74	82	5744	0.58	ng	0.00
11) 2-Methylnaphthalene-d10	11.97	152	7508	0.30	ng	0.00
14) 2,4,6-Tribromophenol	15.73	330	1119	0.31	ng	0.00
15) 2-Fluorobiphenyl	12.86	172	24319	0.68	ng	0.00
25) Fluoranthene-d10	19.02	212	20259	0.30	ng	0.00
29) Terphenyl-d14	19.63	244	30662	0.58	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) n-Nitrosodimethylamine	3.54	42	51	0.015	ng	# 1
6) bis(2-Chloroethyl)ether	7.07	93	15	0.001	ng	# 51
9) Naphthalene	10.43	128	156	0.004	ng	# 23
10) Hexachlorobutadiene	10.75	225	3	0.000	ng	# 98
12) 2-Methylnaphthalene	12.05	142	63	0.002	ng	# 60
16) Acenaphthylene	13.96	152	84	0.002	ng	# 36
17) Acenaphthene	14.30	154	37	0.001	ng	# 62
18) Fluorene	15.29	166	53	0.002	ng	# 91
20) 4-Bromophenyl-phenylether	16.18	248	3	0.000	ng	# 48
21) Hexachlorobenzene	16.31	284	14	0.001	ng	# 19
22) Pentachlorophenol	16.65	266	11	0.002	ng	# 57
23) Phenanthrene	17.03	178	286	0.005	ng	# 80
24) Anthracene	17.11	178	84	0.002	ng	# 82
26) Fluoranthene	19.05	202	465	0.006	ng	# 89
28) Pyrene	19.41	202	446	0.006	ng	# 90
30) Benzo(a)anthracene	21.16	228	622	0.008	ng	# 71
31) Chrysene	21.22	228	610	0.007	ng	# 83
32) Bis(2-ethylhexyl)phthalate	21.13	149	1464	0.056	ng	# 97
33) Indeno(1,2,3-cd)pyrene	25.50	276	258	0.003	ng	# 65
35) Benzo(b)fluoranthene	22.71	252	655	0.007	ng	# 1
36) Benzo(k)fluoranthene	22.76	252	649	0.007	ng	# 1
37) Benzo(a)pyrene	23.26	252	319	0.004	ng	# 1
38) Dibenzo(a,h)anthracene	25.52	278	117	0.001	ng	# 1
39) Benzo(g,h,i)perylene	26.15	276	108	0.001	ng	# 1

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

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