

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080918\
 Data File : BG036256.D
 Acq On : 10 Aug 2018 6:43
 Operator : JU/SJ
 Sample : J4393-03
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 JTY-SB-01-1-62

Quant Time: Aug 10 08:10:29 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG071218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 31 12:30:06 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.01	152	2022	20.00	ng	-0.02
21) Naphthalene-d8	10.84	136	1534	20.00	ng	0.01
38) Acenaphthene-d10	14.64	164	1166	20.00	ng	-0.01
63) Phenanthrene-d10	17.39	188	3574	20.00	ng	0.00
75) Chrysene-d12	21.65	240	3993	20.00	ng	0.00
86) Perylene-d12	24.86	264	3200	20.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.60	112	237984	1954.05	ng	0.00
7) Phenol-d6	7.18	99	338126	1860.81	ng	-0.01
23) Nitrobenzene-d5	9.17	82	226657	6172.45	ng	-0.02
41) 2,4,6-Tribromophenol	16.13	330	165480	10767.39	ng	-0.02
44) 2-Fluorobiphenyl	13.25	172	494770	5684.94	ng	-0.02
78) Terphenyl-d14	19.99	244	877386	4843.53	ng	-0.02

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	3.52	88	536	8.647	ng	# 23
6) Aniline	7.33	93	526	2.233	ng	# 81
8) 2-Chlorophenol	7.57	128	907	7.322	ng	# 1
9) Benzaldehyde	7.17	77	306	2.745	ng	# 34
10) Phenol	7.20	94	4763	25.945	ng	# 1
11) bis(2-Chloroethyl)ether	7.43	93	795	5.530	ng	# 45
12) 1,3-Dichlorobenzene	7.90	146	869	5.810	ng	# 64
13) 1,4-Dichlorobenzene	8.04	146	837	5.507	ng	# 71
14) 1,2-Dichlorobenzene	8.37	146	853	5.842	ng	# 86
16) 2,2'-oxybis(1-Chloropropan	8.54	45	846	7.019	ng	# 75
17) 2-Methylphenol	8.47	107	395	3.165	ng	# 59
18) Hexachloroethane	9.09	117	470	8.088	ng	# 66
19) n-Nitroso-di-n-propylamine	9.17	70	31522	206.007	ng	# 67
24) Nitrobenzene	9.21	77	985	26.672	ng	# 75
25) Isophorone	9.77	82	1102	16.166	ng	# 69
26) 2-Nitrophenol	9.99	139	55	4.193	ng	# 29
27) 2,4-Dimethylphenol	10.07	122	64	2.883	ng	# 62
29) 2,4-Dichlorophenol	10.58	162	81	3.196	ng	# 23
30) 1,2,4-Trichlorobenzene	10.70	180	522	16.797	ng	# 12
31) Naphthalene	10.89	128	2620	32.788	ng	# 68
32) Benzoic acid	10.15	122	61	4.081	ng	# 1
33) 4-Chloroaniline	11.07	127	309	8.872	ng	# 46
34) Hexachlorobutadiene	11.13	225	732	30.207	ng	# 57
36) 4-Chloro-3-methylphenol	12.16	107	72	2.120	ng	# 22
37) 2-Methylnaphthalene	12.48	142	1679	28.203	ng	# 74
39) 1,2,4,5-Tetrachlorobenzene	12.84	216	1348	30.630	ng	# 83
42) 2,4,6-Trichlorophenol	13.13	196	315	12.239	ng	# 59
43) 2,4,5-Trichlorophenol	13.13	196	315	10.941	ng	# 63
45) 1,1'-Biphenyl	13.48	154	3481	38.507	ng	# 71
46) 2-Chloronaphthalene	13.52	162	1651	23.480	ng	# 76
47) 2-Nitroaniline	13.79	65	78	3.138	ng	# 60
48) Acenaphthylene	14.36	152	3544	30.088	ng	# 82
49) Dimethylphthalate	14.09	163	54318	531.702	ng	# 96
50) 2,6-Dinitrotoluene	14.24	165	346	16.632	ng	# 18

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
51) Acenaphthene	14.70	154	1813	24.709	nq	# 77
52) 3-Nitroaniline	14.93	138	53	2.493	nq	# 1
54) Dibenzofuran	15.06	168	3586	30.730	nq	90
55) 4-Nitrophenol	15.06	139	1292	85.324	nq	# 1
56) 2,4-Dinitrotoluene	15.02	165	286	9.583	nq	# 71
57) Fluorene	15.69	166	3047	31.154	nq	95
58) 2,3,4,6-Tetrachlorophenol	15.29	232	609	22.061	nq	# 64
59) Diethylphthalate	15.44	149	2879	27.407	nq	# 85
60) 4-Chlorophenyl-phenylether	15.68	204	1442	25.085	nq	# 83
61) 4-Nitroaniline	15.84	138	53	2.383	nq	# 12
62) Azobenzene	15.97	77	2628	25.363	nq	# 57
64) 4,6-Dinitro-2-methylphenol	15.43	198	60	3.016	nq	# 29
65) n-Nitrosodiphenylamine	15.90	169	1838	18.068	nq	89
66) 4-Bromophenyl-phenylether	16.57	248	958	23.104	nq	# 79
67) Hexachlorobenzene	16.70	284	952	22.295	nq	# 63
68) Atrazine	16.85	200	797	19.028	nq	# 58
69) Pentachlorophenol	17.13	266	53	2.038	nq	# 19
70) Phenanthrene	17.43	178	7095	39.784	nq	98
71) Anthracene	17.54	178	4862	27.380	nq	97
72) Carbazole	17.82	167	1809	10.727	nq	# 59
73) Di-n-butylphthalate	18.33	149	5361	26.901	nq	# 94
74) Fluoranthene	19.44	202	8160	33.969	nq	96
76) Benzydine	19.99	184	15496	147.613	nq	# 90
77) Pyrene	19.80	202	9006	35.670	nq	96
79) Butylbenzylphthalate	20.68	149	1929	21.365	nq	# 75
80) Benzo(a)anthracene	21.64	228	7141	28.698	nq	# 94
82) Chrysene	21.64	228	7141	30.969	nq	93
83) Bis(2-ethylhexyl)phthalate	21.52	149	4604	37.508	nq	# 92
84) Di-n-octyl phthalate	22.71	149	5540	26.955	nq	# 92
85) Indeno(1,2,3-cd)pyrene	28.50	276	1723	6.749	nq	# 85
87) Benzo(b)fluoranthene	23.83	252	5646	29.029	nq	# 75
88) Benzo(k)fluoranthene	23.90	252	5675	29.845	nq	# 88
89) Benzo(a)pyrene	24.71	252	5611	31.010	nq	# 82
90) Dibenzo(a,h)anthracene	28.57	278	2030	11.428	nq	# 88
91) Benzo(a,h,i)perylene	29.64	276	2799	16.102	nq	96

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

