

Data Path : Z:\HPCHEM1\BNA G\DATA\BG092517\
 Data File : BG028910.D
 Acq On : 25 Sep 2017 16:34
 Operator : SJ/JU
 Sample : PB102294BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB102294BS

Quant Time: Sep 26 01:23:49 2017
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG091217.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 12 16:57:11 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.28	152	28931	20.00	ng	-0.05
21) Naphthalene-d8	11.10	136	120223	20.00	ng	-0.05
38) Acenaphthene-d10	14.90	164	89944	20.00	ng	-0.05
63) Phenanthrene-d10	17.65	188	248757	20.00	ng	-0.05
75) Chrysene-d12	21.97	240	295939	20.00	ng	-0.06
86) Perylene-d12	25.42	264	296400	20.00	ng	-0.11

System Monitoring Compounds

5) 2-Fluorophenol	5.85	112	248154	141.53	ng	-0.05
7) Phenol-d6	7.44	99	343990	130.31	ng	-0.05
23) Nitrobenzene-d5	9.45	82	209587	83.40	ng	-0.05
41) 2,4,6-Tribromophenol	16.39	330	213360	143.42	ng	-0.05
44) 2-Fluorobiphenyl	13.52	172	560821	83.14	ng	-0.05
78) Terphenyl-d14	20.23	244	1102186	79.42	ng	-0.05

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.77	88	21765	30.654	ng	96
3) Pyridine	4.16	79	60836	30.036	ng	# 86
4) n-Nitrosodimethylamine	4.08	42	32887	35.695	ng	# 61
6) Aniline	7.61	93	93503	30.434	ng	97
8) 2-Chlorophenol	7.85	128	73129	37.414	ng	93
9) Benzaldehyde	7.42	77	38452	23.478	ng	86
10) Phenol	7.47	94	107086	38.437	ng	88
11) bis(2-Chloroethyl)ether	7.69	93	69479	34.736	ng	91
12) 1,3-Dichlorobenzene	8.17	146	73676	35.006	ng	96
13) 1,4-Dichlorobenzene	8.31	146	75462	34.868	ng	99
14) 1,2-Dichlorobenzene	8.64	146	72700	34.002	ng	97
15) Benzyl Alcohol	8.52	79	66759	32.395	ng	86
16) 2,2'-oxybis(1-Chloropropan	8.80	45	121389	33.392	ng	99
17) 2-Methylphenol	8.72	107	65825	36.157	ng	94
18) Hexachloroethane	9.36	117	27125	34.196	ng	99
19) n-Nitroso-di-n-propylamine	9.08	70	59002	31.529	ng	# 86
20) 3+4-Methylphenols	9.05	107	92279	36.239	ng	94
22) Acetophenone	9.11	105	114257	32.503	ng	# 87
24) Nitrobenzene	9.49	77	92865	33.370	ng	# 89
25) Isophorone	10.01	82	165098	32.467	ng	99
26) 2-Nitrophenol	10.21	139	38635	32.620	ng	# 92
27) 2,4-Dimethylphenol	10.25	122	74628	34.471	ng	95
28) bis(2-Chloroethoxy)methane	10.48	93	97508	35.310	ng	96
29) 2,4-Dichlorophenol	10.75	162	78773	36.569	ng	94
30) 1,2,4-Trichlorobenzene	10.96	180	85237	34.687	ng	98
31) Naphthalene	11.15	128	218194	34.750	ng	98
32) Benzoic acid	10.39	122	30827	23.594	ng	90
33) 4-Chloroaniline	11.27	127	37712	14.337	ng	97
34) Hexachlorobutadiene	11.42	225	60581	33.471	ng	95
35) Caprolactam	12.04	113	30265	39.180	ng	# 85
36) 4-Chloro-3-methylphenol	12.37	107	98270	35.054	ng	94
37) 2-Methylnaphthalene	12.74	142	171063	36.490	ng	94
39) 1,2,4,5-Tetrachlorobenzene	13.10	216	112523	30.426	ng	# 94
40) Hexachlorocyclopentadiene	13.07	237	103300	72.672	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.34	196	74313	31.660	nq	94
43) 2,4,5-Trichlorophenol	13.42	196	83483	35.396	nq	96
45) 1,1'-Biphenyl	13.73	154	230643	30.971	nq	97
46) 2-Chloronaphthalene	13.78	162	180921	33.308	nq	99
47) 2-Nitroaniline	13.99	65	73530	36.425	nq	88
48) Acenaphthylene	14.62	152	296264	34.140	nq	96
49) Dimethylphthalate	14.34	163	264224	35.032	nq	100
50) 2,6-Dinitrotoluene	14.47	165	60925	36.303	nq	# 80
51) Acenaphthene	14.96	154	175834	33.356	nq	95
52) 3-Nitroaniline	14.81	138	42876	28.890	nq	92
53) 2,4-Dinitrophenol	15.02	184	62721	71.834	nq	98
54) Dibenzofuran	15.29	168	313900	37.340	nq	93
55) 4-Nitrophenol	15.12	139	79251	72.885	nq	# 71
56) 2,4-Dinitrotoluene	15.26	165	91975	38.976	nq	# 88
57) Fluorene	15.94	166	271366	36.158	nq	97
58) 2,3,4,6-Tetrachlorophenol	15.52	232	84288	40.548	nq	# 95
59) Diethylphthalate	15.69	149	284037	36.646	nq	99
60) 4-Chlorophenyl-phenylether	15.92	204	152862	35.769	nq	92
61) 4-Nitroaniline	15.97	138	62673	39.861	nq	89
62) Azobenzene	16.22	77	254442	39.033	nq	92
64) 4,6-Dinitro-2-methylphenol	16.03	198	57296	35.695	nq	96
65) n-Nitrosodiphenylamine	16.14	169	242406	33.993	nq	99
66) 4-Bromophenyl-phenylether	16.82	248	110834	34.627	nq	96
67) Hexachlorobenzene	16.96	284	118140	32.429	nq	# 83
68) Atrazine	17.08	200	106452	34.767	nq	95
69) Pentachlorophenol	17.30	266	113062	68.689	nq	97
70) Phenanthrene	17.69	178	468317	36.815	nq	94
71) Anthracene	17.78	178	470019	36.577	nq	98
72) Carbazole	18.05	167	432704	37.388	nq	96
73) Di-n-butylphthalate	18.58	149	511526	36.046	nq	# 94
74) Fluoranthene	19.69	202	614833	39.584	nq	92
76) Benzidine	19.86	184	319344	41.612	nq	94
77) Pyrene	20.05	202	630664	36.946	nq	97
79) Butylbenzylphthalate	20.92	149	242863	35.228	nq	90
80) Benzo(a)anthracene	21.95	228	652675	36.639	nq	96
81) 3,3'-Dichlorobenzidine	21.85	252	165007	22.847	nq	# 96
82) Chrysene	22.02	228	608086	35.977	nq	97
83) Bis(2-ethylhexyl)phthalate	21.80	149	339207	34.610	nq	97
84) Di-n-octyl phthalate	23.09	149	592460	34.598	nq	# 89
85) Indeno(1,2,3-cd)pyrene	29.39	276	774564	35.667	nq	# 55
87) Benzo(b)fluoranthene	24.32	252	655653	36.922	nq	97
88) Benzo(k)fluoranthene	24.39	252	662849	37.369	nq	95
89) Benzo(a)pyrene	25.26	252	643915	38.202	nq	95
90) Dibenzo(a,h)anthracene	29.45	278	654784	37.260	nq	99
91) Benzo(g,h,i)perylene	30.65	276	637702	37.191	nq	97

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

