

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011116\
 Data File : BF084383.D
 Acq On : 11 Jan 2016 16:53
 Operator : SJ/UM
 Sample : PB87733BS
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB87733BS

Quant Time: Jan 12 03:27:38 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF123115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 04 12:22:40 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	254949	20.00	ng	-0.05
21) Naphthalene-d8	8.14	136	1055674	20.00	ng	-0.05
38) Acenaphthene-d10	9.89	164	450907	20.00	ng	-0.06
63) Phenanthrene-d10	11.38	188	765979	20.00	ng	-0.05
75) Chrysene-d12	14.02	240	511788	20.00	ng	-0.05
86) Perylene-d12	15.45	264	426239	20.00	ng	-0.06

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.48	112	1919692	121.79	ng	-0.01
7) Phenol-d6	6.51	99	2315661	116.98	ng	-0.02
23) Nitrobenzene-d5	7.42	82	1387610	73.15	ng	-0.05
41) 2,4,6-Tribromophenol	10.69	330	536837	117.93	ng	-0.05
44) 2-Fluorobiphenyl	9.22	172	2207214	86.20	ng	-0.04
78) Terphenyl-d14	12.97	244	1715549	74.82	ng	-0.05

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	2.54	88	220244	29.50	ng	# 71
3) Pyridine	3.26	79	619653	29.77	ng	# 71
4) n-Nitrosodimethylamine	3.22	42	361206	33.80	ng	79
6) Aniline	6.52	93	286541	9.89	ng	# 1
8) 2-Chlorophenol	6.65	128	698699	39.07	ng	98
9) Benzaldehyde	6.40	77	154299	11.97	ng	99
10) Phenol	6.52	94	925119	41.10	ng	94
11) bis(2-Chloroethyl)ether	6.59	93	620150	35.57	ng	# 80
12) 1,3-Dichlorobenzene	6.80	146	702097m	38.06	ng	
13) 1,4-Dichlorobenzene	6.88	146	703840m	38.05	ng	
14) 1,2-Dichlorobenzene	7.02	146	618625	34.92	ng	96
15) Benzyl Alcohol	7.00	79	524502	31.50	ng	91
16) 2,2'-oxybis(1-Chloropropan	7.14	45	1036857	32.66	ng	88
17) 2-Methylphenol	7.13	107	556281	37.11	ng	# 91
18) Hexachloroethane	7.37	117	260562	35.22	ng	97
19) n-Nitroso-di-n-propylamine	7.28	70	454624	33.93	ng	# 72
20) 3+4-Methylphenols	7.28	107	669457	38.00	ng	# 67
22) Acetophenone	7.26	105	946011	37.27	ng	95
24) Nitrobenzene	7.45	77	747872	38.87	ng	90
25) Isophorone	7.69	82	1313217	36.20	ng	95
26) 2-Nitrophenol	7.76	139	317741	36.29	ng	# 76
27) 2,4-Dimethylphenol	7.80	122	600366	33.88	ng	98
28) bis(2-Chloroethoxy)methane	7.89	93	758176	34.22	ng	99
29) 2,4-Dichlorophenol	8.01	162	558555	37.83	ng	99
30) 1,2,4-Trichlorobenzene	8.09	180	598061	39.24	ng	# 95
31) Naphthalene	8.17	128	1907906	39.83	ng	99
32) Benzoic acid	7.93	122	273929	27.15	ng	92
33) 4-Chloroaniline	8.21	127	128222	5.84	ng	99
34) Hexachlorobutadiene	8.28	225	322787	36.84	ng	97
35) Caprolactam	8.59	113	170157	34.19	ng	# 42
36) 4-Chloro-3-methylphenol	8.72	107	663427	40.12	ng	99
37) 2-Methylnaphthalene	8.85	142	1181323	34.36	ng	97
39) 1,2,4,5-Tetrachlorobenzene	9.02	216	550715	42.48	ng	99
40) Hexachlorocyclopentadiene	9.01	237	279830	35.76	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.14	196	353499	36.45	ng	93
43) 2,4,5-Trichlorophenol	9.18	196	347549	37.38	ng #	81
45) 1,1'-Biphenyl	9.32	154	1447329	40.39	ng	99
46) 2-Chloronaphthalene	9.34	162	1052737	37.40	ng	96
47) 2-Nitroaniline	9.45	65	420498	45.26	ng #	85
48) Acenaphthylene	9.76	152	1690719	40.65	ng	97
49) Dimethylphthalate	9.63	163	1441338	44.71	ng #	90
50) 2,6-Dinitrotoluene	9.69	165	297390	42.06	ng #	62
51) Acenaphthene	9.93	154	1128496	40.45	ng	97
52) 3-Nitroaniline	9.86	138	142620	16.28	ng #	70
53) 2,4-Dinitrophenol	9.96	184	131008	45.12	ng	93
54) Dibenzofuran	10.11	168	1467397	40.31	ng	96
55) 4-Nitrophenol	10.03	139	367648	57.81	ng #	76
56) 2,4-Dinitrotoluene	10.09	165	348545	38.95	ng #	86
57) Fluorene	10.44	166	1097639	38.97	ng	99
58) 2,3,4,6-Tetrachlorophenol	10.22	232	289399	36.46	ng	95
59) Diethylphthalate	10.33	149	1322206	41.60	ng	98
60) 4-Chlorophenyl-phenylether	10.44	204	574048	49.49	ng	97
61) 4-Nitroaniline	10.46	138	261493	33.48	ng	93
62) Azobenzene	10.60	77	1419881	40.73	ng	95
64) 4,6-Dinitro-2-methylphenol	10.50	198	116823	26.04	ng	83
65) n-Nitrosodiphenylamine	10.56	169	1005587	41.06	ng	98
66) 4-Bromophenyl-phenylether	10.93	248	360435	43.72	ng	98
67) Hexachlorobenzene	10.99	284	351581	37.89	ng #	83
68) Atrazine	11.09	200	381024	47.54	ng	97
69) Pentachlorophenol	11.20	266	322112	66.95	ng	99
70) Phenanthrene	11.41	178	1642625	41.00	ng	99
71) Anthracene	11.46	178	1606562	40.90	ng	98
72) Carbazole	11.62	167	1490646	38.82	ng	98
73) Di-n-butylphthalate	11.95	149	1971981	49.57	ng #	97
74) Fluoranthene	12.59	202	1337112	31.95	ng	99
76) Benzidine	12.72	184	629281	34.29	ng	99
77) Pyrene	12.82	202	1330609	33.13	ng	99
79) Butylbenzylphthalate	13.45	149	679061	39.07	ng #	83
80) Benzo(a)anthracene	14.01	228	1075246	35.78	ng	99
81) 3,3'-Dichlorobenzidine	13.97	252	376852	35.93	ng #	96
82) Chrysene	14.04	228	909425	31.49	ng	99
83) Bis(2-ethylhexyl)phthalate	14.01	149	877406	39.96	ng #	95
84) Di-n-octyl phthalate	14.61	149	1382374	37.29	ng #	87
85) Indeno(1,2,3-cd)pyrene	16.84	276	858853	37.52	ng	100
87) Benzo(h)fluoranthene	15.04	252	992442	35.59	ng	98
88) Benzo(k)fluoranthene	15.07	252	988533m	43.35	ng	
89) Benzo(a)pyrene	15.39	252	944999	39.96	ng	100
90) Dibenzo(a,h)anthracene	16.85	278	735301	36.13	ng	98
91) Benzo(g,h,i)perylene	17.27	276	681315	32.33	ng	98

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

