

Data Path : Z:\HPCHEM1\BNA F\DATA\BF022317\
 Data File : BF093108.D
 Acq On : 23 Feb 2017 17:01
 Operator : SJ/MA
 Sample : PB97124BS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB97124BS

Manual Integrations
 APPROVED

mohammad
 2/24/2017 2:11:04 PM

Quant Time: Feb 24 00:05:18 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF013017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 17 17:31:55 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.84	152	164493	20.00	ng	-0.01
21) Naphthalene-d8	8.13	136	622852	20.00	ng	-0.01
38) Acenaphthene-d10	9.88	164	308558	20.00	ng	-0.01
63) Phenanthrene-d10	11.36	188	542824	20.00	ng	-0.02
75) Chrysene-d12	14.00	240	433846	20.00	ng	-0.02
86) Perylene-d12	15.43	264	350323	20.00	ng	-0.02

System Monitoring Compounds						
5) 2-Fluorophenol	5.44	112	1126711	117.51	ng	-0.01
7) Phenol-d6	6.49	99	1477349	119.75	ng	0.00
23) Nitrobenzene-d5	7.41	82	929249	83.08	ng	-0.01
41) 2,4,6-Tribromophenol	10.67	330	292432	131.04	ng	-0.01
44) 2-Fluorobiphenyl	9.21	172	1730989	101.63	ng	-0.01
78) Terphenyl-d14	12.94	244	1373839	74.41	ng	-0.02

Target Compounds						Qvalue
2) 1,4-Dioxane	2.43	88	146641	28.30	ng	# 86
3) Pyridine	3.15	79	433298	32.89	ng	88
4) n-Nitrosodimethylamine	3.10	42	251624	40.68	ng	86
6) Aniline	6.51	93	564633	33.55	ng	# 46
8) 2-Chlorophenol	6.62	128	441037	41.70	ng	93
9) Benzaldehyde	6.38	77	191837	21.71	ng	94
10) Phenol	6.50	94	621610	43.07	ng	91
11) bis(2-Chloroethyl)ether	6.58	93	432592	37.55	ng	91
12) 1,3-Dichlorobenzene	6.78	146	490463	39.59	ng	99
13) 1,4-Dichlorobenzene	6.85	146	489488	39.04	ng	99
14) 1,2-Dichlorobenzene	7.01	146	490264	40.89	ng	96
15) Benzyl Alcohol	6.99	79	352498	38.60	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.13	45	799519	39.95	ng	94
17) 2-Methylphenol	7.10	107	382314	42.13	ng	97
18) Hexachloroethane	7.36	117	170366	39.80	ng	89
19) n-Nitroso-di-n-propylamine	7.26	70	313231	39.89	ng	99
20) 3+4-Methylphenols	7.26	107	450540	41.53	ng	97
22) Acetophenone	7.25	105	604473	42.51	ng	# 95
24) Nitrobenzene	7.44	77	527910	45.97	ng	# 87
25) Isophorone	7.68	82	938897	45.05	ng	99
26) 2-Nitrophenol	7.74	139	240025	43.97	ng	94
27) 2,4-Dimethylphenol	7.79	122	435231	45.39	ng	97
28) bis(2-Chloroethoxy)methane	7.88	93	579384	41.97	ng	97
29) 2,4-Dichlorophenol	8.00	162	406400	45.45	ng	97
30) 1,2,4-Trichlorobenzene	8.08	180	404689	42.00	ng	98
31) Naphthalene	8.14	128	1246270	39.47	ng	99
32) Benzoic acid	7.94	122	250355	46.37	ng	94
33) 4-Chloroaniline	8.20	127	331525	26.77	ng	98
34) Hexachlorobutadiene	8.27	225	214494	40.46	ng	98
35) Caprolactam	8.58	113	130477m	46.39	ng	
36) 4-Chloro-3-methylphenol	8.69	107	431202	46.25	ng	99
37) 2-Methylnaphthalene	8.84	142	906557	44.72	ng	97
39) 1,2,4,5-Tetrachlorobenzene	9.00	216	361168	41.70	ng	99
40) Hexachlorocyclopentadiene	8.99	237	320702	82.07	ng	95

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42) 2,4,6-Trichlorophenol	9.13	196	275725	48.45	nq	95
43) 2,4,5-Trichlorophenol	9.16	196	286964	46.02	nq	96
45) 1,1'-Biphenyl	9.30	154	1012159	45.30	nq	96
46) 2-Chloronaphthalene	9.33	162	848177	48.53	nq	96
47) 2-Nitroaniline	9.42	65	265067	45.11	nq	94
48) Acenaphthylene	9.74	152	1355939	44.91	nq	98
49) Dimethylphthalate	9.61	163	949241	45.67	nq	99
50) 2,6-Dinitrotoluene	9.68	165	245615	54.72	nq	88
51) Acenaphthene	9.92	154	809924	44.87	nq	96
52) 3-Nitroaniline	9.84	138	151906	29.57	nq	90
53) 2,4-Dinitrophenol	9.95	184	219827	95.31	nq	# 56
54) Dibenzofuran	10.09	168	1142552	47.32	nq	95
55) 4-Nitrophenol	10.02	139	355036	95.97	nq	89
56) 2,4-Dinitrotoluene	10.08	165	259684	43.46	nq	# 92
57) Fluorene	10.43	166	920682	49.56	nq	98
58) 2,3,4,6-Tetrachlorophenol	10.21	232	218713	52.58	nq	96
59) Diethylphthalate	10.30	149	926555	47.31	nq	98
60) 4-Chlorophenyl-phenylether	10.42	204	407806	45.70	nq	87
61) 4-Nitroaniline	10.46	138	263552	50.10	nq	# 76
62) Azobenzene	10.58	77	1081515	53.45	nq	96
64) 4,6-Dinitro-2-methylphenol	10.49	198	166598	50.29	nq	# 67
65) n-Nitrosodiphenylamine	10.54	169	831678	47.96	nq	99
66) 4-Bromophenyl-phenylether	10.91	248	260206	45.88	nq	# 86
67) Hexachlorobenzene	10.98	284	247473	44.16	nq	# 89
68) Atrazine	11.07	200	259732	46.67	nq	98
69) Pentachlorophenol	11.17	266	266094	90.90	nq	98
70) Phenanthrene	11.39	178	1370301	44.06	nq	98
71) Anthracene	11.44	178	1264745	40.50	nq	99
72) Carbazole	11.60	167	1189741	39.85	nq	99
73) Di-n-butylphthalate	11.93	149	1601838	49.61	nq	# 96
74) Fluoranthene	12.58	202	1362218	42.46	nq	90
76) Benzidine	12.69	184	558495	30.21	nq	97
77) Pyrene	12.80	202	1380643	42.52	nq	99
79) Butylbenzylphthalate	13.41	149	643544	48.81	nq	99
80) Benzo(a)anthracene	13.99	228	1104489	41.62	nq	99
81) 3,3'-Dichlorobenzidine	13.95	252	293155	30.99	nq	98
82) Chrysene	14.03	228	1145009	46.09	nq	99
83) Bis(2-ethylhexyl)phthalate	13.97	149	829328	45.99	nq	99
84) Di-n-octyl phthalate	14.59	149	1512916	51.53	nq	99
85) Indeno(1,2,3-cd)pyrene	16.82	276	825402	42.89	nq	# 94
87) Benzo(b)fluoranthene	15.01	252	1239570m	53.76	nq	
88) Benzo(k)fluoranthene	15.05	252	705397m	35.43	nq	
89) Benzo(a)pyrene	15.37	252	905051	45.38	nq	99
90) Dibenzo(a,h)anthracene	16.83	278	684295	42.45	nq	96
91) Benzo(g,h,i)perylene	17.24	276	702133	43.12	nq	# 90

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

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