

Data Path : Z:\HPCHEM1\BNA F\DATA\BF020116\
 Data File : BF084712.D
 Acq On : 1 Feb 2016 16:35
 Operator : SJ/IZ
 Sample : PB88012BS
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB88012BS

Manual Integrations
 APPROVED

mohammad
 2/2/2016 4:17:58 PM

Quant Time: Feb 02 00:46:23 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF020116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 01 14:55:11 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.72	152	179367	20.00	ng	-0.01
21) Naphthalene-d8	8.00	136	709297	20.00	ng	-0.02
38) Acenaphthene-d10	9.76	164	359559	20.00	ng	-0.01
63) Phenanthrene-d10	11.23	188	653877	20.00	ng	-0.01
75) Chrysene-d12	13.87	240	499782	20.00	ng	0.00
86) Perylene-d12	15.24	264	431063	20.00	ng	-0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.32	112	1434608	124.58	ng	0.01
7) Phenol-d6	6.36	99	1744298	122.18	ng	-0.01
23) Nitrobenzene-d5	7.29	82	1120849	89.02	ng	-0.01
41) 2,4,6-Tribromophenol	10.54	330	459123	122.42	ng	-0.01
44) 2-Fluorobiphenyl	9.08	172	1867500	88.79	ng	-0.01
78) Terphenyl-d14	12.82	244	1818694	82.46	ng	-0.01

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	2.30	88	176033	34.90	ng	# 72
3) Pyridine	2.97	79	541514	41.20	ng	# 74
4) n-Nitrosodimethylamine	2.92	42	271276	39.91	ng	# 81
6) Aniline	6.37	93	558550	31.97	ng	# 75
8) 2-Chlorophenol	6.50	128	520694	39.95	ng	# 92
9) Benzaldehyde	6.26	77	23284	2.76	ng	# 85
10) Phenol	6.38	94	590381	36.92	ng	# 90
11) bis(2-Chloroethyl)ether	6.45	93	479675	38.46	ng	# 83
12) 1,3-Dichlorobenzene	6.65	146	500928m	36.37	ng	# 97
13) 1,4-Dichlorobenzene	6.73	146	513452	37.30	ng	# 95
14) 1,2-Dichlorobenzene	6.89	146	489146	38.15	ng	# 91
15) Benzyl Alcohol	6.86	79	417098	42.31	ng	# 87
16) 2,2'-oxybis(1-Chloropropan	7.00	45	768698	36.64	ng	# 92
17) 2-Methylphenol	6.99	107	437131	42.40	ng	# 92
18) Hexachloroethane	7.22	117	204908	38.65	ng	# 92
19) n-Nitroso-di-n-propylamine	7.14	70	361942	40.45	ng	# 76
20) 3+4-Methylphenols	7.14	107	523050	40.88	ng	# 75
22) Acetophenone	7.13	105	721780	38.61	ng	# 99
24) Nitrobenzene	7.30	77	494644	36.69	ng	# 95
25) Isophorone	7.55	82	993798	40.59	ng	# 96
26) 2-Nitrophenol	7.62	139	278609	41.90	ng	# 87
27) 2,4-Dimethylphenol	7.66	122	470808	40.70	ng	# 96
28) bis(2-Chloroethoxy)methane	7.76	93	595880	41.02	ng	# 99
29) 2,4-Dichlorophenol	7.87	162	432662	43.72	ng	# 93
30) 1,2,4-Trichlorobenzene	7.94	180	440701	39.42	ng	# 96
31) Naphthalene	8.02	128	1378602	38.75	ng	# 99
32) Benzoic acid	7.80	122	305358	41.27	ng	# 93
33) 4-Chloroaniline	8.08	127	483473	32.86	ng	# 98
34) Hexachlorobutadiene	8.14	225	264141	40.98	ng	# 97
35) Caprolactam	8.45	113	139498m	45.73	ng	# 92
36) 4-Chloro-3-methylphenol	8.57	107	490280	43.42	ng	# 98
37) 2-Methylnaphthalene	8.72	142	1052150	47.25	ng	# 98
39) 1,2,4,5-Tetrachlorobenzene	8.88	216	396463	34.72	ng	# 98
40) Hexachlorocyclopentadiene	8.86	237	406795	80.49	ng	# 97

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42) 2,4,6-Trichlorophenol	9.00	196	303644	41.27	nq	94
43) 2,4,5-Trichlorophenol	9.05	196	294397	39.38	nq #	91
45) 1,1'-Biphenyl	9.18	154	1209938	37.67	nq	95
46) 2-Chloronaphthalene	9.21	162	889998	37.63	nq #	85
47) 2-Nitroaniline	9.30	65	285151	38.08	nq	97
48) Acenaphthylene	9.62	152	1427572	39.78	nq	98
49) Dimethylphthalate	9.49	163	1137412	41.53	nq #	91
50) 2,6-Dinitrotoluene	9.55	165	258853	43.27	nq #	84
51) Acenaphthene	9.79	154	930216	39.84	nq	96
52) 3-Nitroaniline	9.71	138	200884	29.88	nq	97
53) 2,4-Dinitrophenol	9.82	184	253517	90.96	nq #	84
54) Dibenzofuran	9.96	168	1306270	45.77	nq	97
55) 4-Nitrophenol	9.89	139	350220	70.89	nq #	80
56) 2,4-Dinitrotoluene	9.95	165	306843	40.21	nq #	84
57) Fluorene	10.30	166	1017423	42.69	nq	99
58) 2,3,4,6-Tetrachlorophenol	10.08	232	237179	41.68	nq	92
59) Diethylphthalate	10.19	149	1137559	42.54	nq	97
60) 4-Chlorophenyl-phenylether	10.29	204	428272	42.35	nq	94
61) 4-Nitroaniline	10.33	138	248423	37.92	nq	90
62) Azobenzene	10.45	77	1113604	43.31	nq	90
64) 4,6-Dinitro-2-methylphenol	10.36	198	194368	48.16	nq	86
65) n-Nitrosodiphenylamine	10.42	169	929385	44.76	nq	99
66) 4-Bromophenyl-phenylether	10.78	248	308995	42.78	nq #	93
67) Hexachlorobenzene	10.84	284	311870	39.63	nq #	83
68) Atrazine	10.94	200	264882	40.40	nq	97
69) Pentachlorophenol	11.05	266	347203	93.88	nq	98
70) Phenanthrene	11.25	178	1430679	39.45	nq	97
71) Anthracene	11.31	178	1605968	43.95	nq	99
72) Carbazole	11.47	167	1488939	46.73	nq	99
73) Di-n-butylphthalate	11.80	149	1569291	38.68	nq #	96
74) Fluoranthene	12.44	202	1619827	44.13	nq	96
76) Benzidine	12.57	184	565142	31.93	nq	98
77) Pyrene	12.67	202	1697384	44.36	nq	98
79) Butylbenzylphthalate	13.30	149	786972	45.34	nq	96
80) Benzo(a)anthracene	13.85	228	1330452	42.89	nq	99
81) 3,3'-Dichlorobenzidine	13.83	252	413332	37.63	nq #	99
82) Chrysene	13.89	228	1319015	43.85	nq	99
83) Bis(2-ethylhexyl)phthalate	13.86	149	935505	43.31	nq	97
84) Di-n-octyl phthalate	14.47	149	1590511	44.63	nq #	90
85) Indeno(1,2,3-cd)pyrene	16.55	276	996566	42.09	nq	99
87) Benzo(b)fluoranthene	14.87	252	1271452m	43.90	nq	
88) Benzo(k)fluoranthene	14.89	252	944002m	39.42	nq	
89) Benzo(a)pyrene	15.20	252	1073637	43.51	nq #	98
90) Dibenzo(a,h)anthracene	16.57	278	833519	40.12	nq #	95
91) Benzo(g,h,i)perylene	16.95	276	837448	40.02	nq	98

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

