

Data Path : Z:\HPCHEM1\BNA F\DATA\BF122716\
 Data File : BF091969.D
 Acq On : 27 Dec 2016 15:54
 Operator : UM/SJ
 Sample : PB95629BS
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB95629BS

Quant Time: Dec 28 00:39:26 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF122116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 21 16:17:51 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.73	152	213713	20.00	ng	-0.02
21) Naphthalene-d8	8.02	136	908228	20.00	ng	-0.02
38) Acenaphthene-d10	9.78	164	477659	20.00	ng	-0.01
63) Phenanthrene-d10	11.25	188	760275	20.00	ng	-0.01
75) Chrysene-d12	13.89	240	566130	20.00	ng	-0.01
86) Perylene-d12	15.29	264	471897	20.00	ng	-0.01

System Monitoring Compounds						
5) 2-Fluorophenol	5.36	112	1601502	125.12	ng	0.01
7) Phenol-d6	6.41	99	1687992	108.02	ng	0.00
23) Nitrobenzene-d5	7.30	82	1066520	65.30	ng	-0.02
41) 2,4,6-Tribromophenol	10.57	330	374620	94.77	ng	-0.01
44) 2-Fluorobiphenyl	9.10	172	1890189	80.38	ng	-0.01
78) Terphenyl-d14	12.84	244	1622171	69.49	ng	-0.01

Target Compounds						Qvalue
2) 1,4-Dioxane	2.32	88	252222	40.03	ng	95
3) Pyridine	2.99	79	684791	31.14	ng	98
4) n-Nitrosodimethylamine	2.94	42	338791	40.84	ng	97
6) Aniline	6.41	93	439924	21.67	ng	# 64
8) 2-Chlorophenol	6.53	128	607800	41.75	ng	87
9) Benzaldehyde	6.27	77	63725	5.29	ng	99
10) Phenol	6.43	94	712023	39.99	ng	# 64
11) bis(2-Chloroethyl)ether	6.48	93	718746	50.73	ng	88
12) 1,3-Dichlorobenzene	6.67	146	627832m	40.39	ng	
13) 1,4-Dichlorobenzene	6.75	146	636052	40.21	ng	94
14) 1,2-Dichlorobenzene	6.90	146	547267	39.87	ng	97
15) Benzyl Alcohol	6.89	79	405649	33.41	ng	95
16) 2,2'-oxybis(1-Chloropropan	7.01	45	822610	38.99	ng	83
17) 2-Methylphenol	7.02	107	491309	41.96	ng	# 89
18) Hexachloroethane	7.24	117	238179	40.61	ng	96
19) n-Nitroso-di-n-propylamine	7.16	70	499938	48.09	ng	94
20) 3+4-Methylphenols	7.18	107	708765	50.75	ng	# 73
22) Acetophenone	7.15	105	923924	41.24	ng	# 90
24) Nitrobenzene	7.32	77	708678	40.74	ng	95
25) Isophorone	7.56	82	1181608	39.21	ng	96
26) 2-Nitrophenol	7.64	139	354743	41.35	ng	# 88
27) 2,4-Dimethylphenol	7.70	122	509748	35.82	ng	98
28) bis(2-Chloroethoxy)methane	7.78	93	780139	42.69	ng	99
29) 2,4-Dichlorophenol	7.90	162	506425	42.03	ng	97
30) 1,2,4-Trichlorobenzene	7.96	180	508619	38.52	ng	# 95
31) Naphthalene	8.04	128	1695519	40.79	ng	99
32) Benzoic acid	7.85	122	327582	31.04	ng	95
33) 4-Chloroaniline	8.14	127	312299	16.66	ng	96
34) Hexachlorobutadiene	8.16	225	282818	40.58	ng	99
35) Caprolactam	8.48	113	150292m	37.96	ng	
36) 4-Chloro-3-methylphenol	8.61	107	541794	39.08	ng	99
37) 2-Methylnaphthalene	8.74	142	1108422	43.46	ng	98
39) 1,2,4,5-Tetrachlorobenzene	8.90	216	447987	37.12	ng	# 98
40) Hexachlorocyclopentadiene	8.89	237	423097	78.78	ng	95

Data Path : Z:\HPCHEM1\BNA F\DATA\BF122716\
 Data File : BF091969.D
 Acq On : 27 Dec 2016 15:54
 Operator : UM/SJ
 Sample : PB95629BS
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB95629BS

Quant Time: Dec 28 00:39:26 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF122116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 21 16:17:51 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.02	196	328185	38.89	nq	97
43) 2,4,5-Trichlorophenol	9.08	196	337267	38.83	nq	# 89
45) 1,1'-Biphenyl	9.20	154	1253454	38.33	nq	97
46) 2-Chloronaphthalene	9.22	162	983638	37.98	nq	94
47) 2-Nitroaniline	9.33	65	375431	40.71	nq	# 77
48) Acenaphthylene	9.64	152	1596790	40.09	nq	98
49) Dimethylphthalate	9.50	163	1256790	41.14	nq	99
50) 2,6-Dinitrotoluene	9.57	165	306307	45.81	nq	93
51) Acenaphthene	9.81	154	988598	36.61	nq	98
52) 3-Nitroaniline	9.74	138	202946	24.52	nq	84
53) 2,4-Dinitrophenol	9.85	184	226720	60.94	nq	# 52
54) Dibenzofuran	9.98	168	1441092	39.51	nq	94
55) 4-Nitrophenol	9.94	139	371901	66.19	nq	96
56) 2,4-Dinitrotoluene	9.97	165	340238	40.54	nq	# 84
57) Fluorene	10.33	166	1064801	40.13	nq	99
58) 2,3,4,6-Tetrachlorophenol	10.11	232	294801	42.91	nq	98
59) Diethylphthalate	10.20	149	1204223	40.59	nq	100
60) 4-Chlorophenyl-phenylether	10.32	204	488008	42.00	nq	95
61) 4-Nitroaniline	10.36	138	302585	35.29	nq	96
62) Azobenzene	10.48	77	1440267	44.09	nq	96
64) 4,6-Dinitro-2-methylphenol	10.38	198	203449	44.25	nq	81
65) n-Nitrosodiphenylamine	10.44	169	1011565	44.84	nq	100
66) 4-Bromophenyl-phenylether	10.81	248	339783	42.96	nq	# 82
67) Hexachlorobenzene	10.86	284	326498	38.62	nq	# 78
68) Atrazine	10.98	200	278136	38.22	nq	93
69) Pentachlorophenol	11.07	266	305163	73.57	nq	99
70) Phenanthrene	11.28	178	1552090	37.65	nq	99
71) Anthracene	11.33	178	1756722	49.93	nq	100
72) Carbazole	11.49	167	1391013	38.49	nq	100
73) Di-n-butylphthalate	11.83	149	1767777	45.28	nq	99
74) Fluoranthene	12.47	202	1628494	45.03	nq	97
76) Benzidine	12.59	184	691945	49.01	nq	97
77) Pyrene	12.69	202	1693253	40.76	nq	100
79) Butylbenzylphthalate	13.32	149	805105	42.62	nq	89
80) Benzo(a)anthracene	13.88	228	1320585	41.32	nq	100
81) 3,3'-Dichlorobenzidine	13.85	252	398109	32.85	nq	98
82) Chrysene	13.92	228	1104049	34.44	nq	99
83) Bis(2-ethylhexyl)phthalate	13.88	149	926546	41.65	nq	99
84) Di-n-octyl phthalate	14.49	149	1682313	40.82	nq	98
85) Indeno(1,2,3-cd)pyrene	16.63	276	1151428	41.46	nq	99
87) Benzo(b)fluoranthene	14.90	252	1231615m	41.92	nq	
88) Benzo(k)fluoranthene	14.92	252	991198m	39.56	nq	
89) Benzo(a)pyrene	15.23	252	1058794	40.60	nq	99
90) Dibenzo(a,h)anthracene	16.64	278	950868	44.36	nq	99
91) Benzo(g,h,i)perylene	17.03	276	984031	44.15	nq	96

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

