

Data Path : Z:\HPCHEM1\BNA F\DATA\BF082616\
 Data File : BF089915.D
 Acq On : 26 Aug 2016 18:44
 Operator : UM/SJ
 Sample : PB93122BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB93122BS

Quant Time: Aug 26 19:24:32 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 26 18:51:25 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.65	152	332167	20.00	ng	0.00
21) Naphthalene-d8	7.94	136	1321874	20.00	ng	0.00
38) Acenaphthene-d10	9.68	164	607334	20.00	ng	-0.01
63) Phenanthrene-d10	11.15	188	1106585	20.00	ng	0.00
75) Chrysene-d12	13.78	240	800543	20.00	ng	0.00
86) Perylene-d12	15.13	264	734703	20.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.24	112	2145741	110.78	ng	0.02
7) Phenol-d6	6.30	99	2445949	102.87	ng	0.00
23) Nitrobenzene-d5	7.22	82	1448279	68.07	ng	0.00
41) 2,4,6-Tribromophenol	10.47	330	580914	100.65	ng	-0.01
44) 2-Fluorobiphenyl	9.02	172	2437009	70.52	ng	0.00
78) Terphenyl-d14	12.74	244	2106499	59.53	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	2.19	88	266778	27.83	ng	97
3) Pyridine	2.82	79	787486	31.32	ng	92
4) n-Nitrosodimethylamine	2.79	42	331686	35.44	ng	# 80
6) Aniline	6.31	93	884702	27.74	ng	# 36
8) 2-Chlorophenol	6.43	128	775508	33.48	ng	91
9) Benzaldehyde	6.18	77	153340	9.90	ng	95
10) Phenol	6.31	94	788389	28.29	ng	85
11) bis(2-Chloroethyl)ether	6.40	93	797344	36.90	ng	92
12) 1,3-Dichlorobenzene	6.59	146	850258m	35.11	ng	
13) 1,4-Dichlorobenzene	6.67	146	845718m	33.98	ng	
14) 1,2-Dichlorobenzene	6.82	146	819966m	35.29	ng	
15) Benzyl Alcohol	6.80	79	539645	34.22	ng	95
16) 2,2'-oxybis(1-Chloropropan	6.95	45	930865	33.66	ng	91
17) 2-Methylphenol	6.92	107	703342	38.77	ng	95
18) Hexachloroethane	7.16	117	307611	34.65	ng	94
19) n-Nitroso-di-n-propylamine	7.08	70	559306	36.78	ng	# 86
20) 3+4-Methylphenols	7.07	107	781330	35.13	ng	# 82
22) Acetophenone	7.07	105	992962	32.71	ng	# 88
24) Nitrobenzene	7.24	77	786084	32.93	ng	# 74
25) Isophorone	7.48	82	1498096	34.91	ng	96
26) 2-Nitrophenol	7.55	139	406014	34.09	ng	# 83
27) 2,4-Dimethylphenol	7.61	122	794253	37.00	ng	97
28) bis(2-Chloroethoxy)methane	7.70	93	985200	37.11	ng	97
29) 2,4-Dichlorophenol	7.80	162	675802	37.52	ng	95
30) 1,2,4-Trichlorobenzene	7.88	180	710354	35.65	ng	96
31) Naphthalene	7.95	128	2193671	35.51	ng	99
32) Benzoic acid	7.74	122	397067	23.74	ng	95
33) 4-Chloroaniline	8.01	127	502773	18.77	ng	# 94
34) Hexachlorobutadiene	8.08	225	346616	33.22	ng	100
35) Caprolactam	8.39	113	213180m	38.86	ng	
36) 4-Chloro-3-methylphenol	8.50	107	679175	34.82	ng	96
37) 2-Methylnaphthalene	8.65	142	1550823	38.80	ng	99
39) 1,2,4,5-Tetrachlorobenzene	8.82	216	543688	27.65	ng	96
40) Hexachlorocyclopentadiene	8.81	237	680171	99.61	ng	98

Data Path : Z:\HPCHEM1\BNA F\DATA\BF082616\
 Data File : BF089915.D
 Acq On : 26 Aug 2016 18:44
 Operator : UM/SJ
 Sample : PB93122BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB93122BS

Quant Time: Aug 26 19:24:32 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 26 18:51:25 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	8.92	196	462662	37.39	nq	99
43) 2,4,5-Trichlorophenol	8.97	196	462154	37.56	nq	99
45) 1,1'-Biphenyl	9.12	154	1673956	35.62	nq	94
46) 2-Chloronaphthalene	9.13	162	1343194	36.15	nq	98
47) 2-Nitroaniline	9.23	65	449762	42.44	nq	# 83
48) Acenaphthylene	9.54	152	2078997	37.24	nq	99
49) Dimethylphthalate	9.43	163	1740118	40.69	nq	100
50) 2,6-Dinitrotoluene	9.48	165	364035	36.87	nq	# 58
51) Acenaphthene	9.71	154	1495869	40.49	nq	99
52) 3-Nitroaniline	9.64	138	331420	30.30	nq	# 76
53) 2,4-Dinitrophenol	9.75	184	344465	64.73	nq	95
54) Dibenzofuran	9.88	168	1907383	41.10	nq	98
55) 4-Nitrophenol	9.82	139	714366	80.01	nq	98
56) 2,4-Dinitrotoluene	9.88	165	497889	38.87	nq	91
57) Fluorene	10.23	166	1228413	33.22	nq	98
58) 2,3,4,6-Tetrachlorophenol	10.01	232	397530	44.39	nq	97
59) Diethylphthalate	10.12	149	1701855	39.18	nq	98
60) 4-Chlorophenyl-phenylether	10.23	204	572159	33.91	nq	99
61) 4-Nitroaniline	10.26	138	444610	43.36	nq	83
62) Azobenzene	10.39	77	1712305	41.93	nq	94
64) 4,6-Dinitro-2-methylphenol	10.28	198	273166	42.96	nq	# 78
65) n-Nitrosodiphenylamine	10.35	169	1414317	40.65	nq	100
66) 4-Bromophenyl-phenylether	10.72	248	417444	35.98	nq	98
67) Hexachlorobenzene	10.78	284	500456	37.81	nq	97
68) Atrazine	10.88	200	371975	34.91	nq	97
69) Pentachlorophenol	10.97	266	550534	73.39	nq	98
70) Phenanthrene	11.18	178	2134564	38.19	nq	99
71) Anthracene	11.23	178	2177891	38.33	nq	99
72) Carbazole	11.39	167	2209364	43.36	nq	96
73) Di-n-butylphthalate	11.74	149	2412864	37.85	nq	98
74) Fluoranthene	12.35	202	2112326	39.73	nq	94
76) Benzidine	12.49	184	1078184	41.62	nq	96
77) Pyrene	12.58	202	1997511	30.64	nq	98
79) Butylbenzylphthalate	13.22	149	1114727	37.86	nq	98
80) Benzo(a)anthracene	13.77	228	1843862	40.22	nq	100
81) 3,3'-Dichlorobenzidine	13.74	252	405635	26.99	nq	98
82) Chrysene	13.81	228	1621052	41.24	nq	99
83) Bis(2-ethylhexyl)phthalate	13.78	149	1388523	34.24	nq	# 97
84) Di-n-octyl phthalate	14.40	149	2491256	46.23	nq	# 92
85) Indeno(1,2,3-cd)pyrene	16.37	276	1440575	28.73	nq	99
87) Benzo(b)fluoranthene	14.77	252	1768536m	43.76	nq	
88) Benzo(k)fluoranthene	14.79	252	1219253m	33.25	nq	
89) Benzo(a)pyrene	15.07	252	1407757	36.80	nq	99
90) Dibenzo(a,h)anthracene	16.39	278	1212988	29.21	nq	99
91) Benzo(g,h,i)perylene	16.74	276	1230011	28.13	nq	97

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

