

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF101916\
 Data File : BF090665.D
 Acq On : 19 Oct 2016 23:18
 Operator : UM/SJ
 Sample : PB94148BS
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB94148BS

Manual Integrations
 APPROVED

Sohil
 10/20/2016 10:36:59 AM

Quant Time: Oct 20 02:17:15 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF101316.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 19 15:13:23 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	209171	20.00	ng	0.00
21) Naphthalene-d8	8.18	136	903105	20.00	ng	0.00
38) Acenaphthene-d10	9.94	164	442145	20.00	ng	0.00
63) Phenanthrene-d10	11.42	188	659575	20.00	ng	0.00
75) Chrysene-d12	14.06	240	453752	20.00	ng	0.00
86) Perylene-d12	15.50	264	381370	20.00	ng	-0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.50	112	1031026	90.36	ng	0.01
7) Phenol-d6	6.53	99	1358445	80.13	ng	0.00
23) Nitrobenzene-d5	7.46	82	1316986	81.01	ng	0.00
41) 2,4,6-Tribromophenol	10.73	330	288156	73.10	ng	0.00
44) 2-Fluorobiphenyl	9.25	172	2016654	75.15	ng	-0.01
78) Terphenyl-d14	13.00	244	1639852	84.15	ng	-0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.58	88	207827	31.96	ng	96
3) Pyridine	3.31	79	504702	34.63	ng	94
4) n-Nitrosodimethylamine	3.26	42	230810	44.80	ng	91
6) Aniline	6.55	93	446538	21.80	ng	96
8) 2-Chlorophenol	6.67	128	597595	40.39	ng	99
9) Benzaldehyde	6.43	77	147776	13.71	ng	92
10) Phenol	6.54	94	792629	40.88	ng	99
11) bis(2-Chloroethyl)ether	6.62	93	642946	40.19	ng	99
12) 1,3-Dichlorobenzene	6.83	146	574297	38.92	ng	98
13) 1,4-Dichlorobenzene	6.91	146	586776	36.57	ng	99
14) 1,2-Dichlorobenzene	7.06	146	606009	40.40	ng	99
15) Benzyl Alcohol	7.03	79	532715	46.13	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.17	45	952643	42.31	ng	91
17) 2-Methylphenol	7.15	107	496828	43.09	ng	98
18) Hexachloroethane	7.40	117	238943	37.68	ng	99
19) n-Nitroso-di-n-propylamine	7.31	70	506728	43.59	ng	# 88
20) 3+4-Methylphenols	7.30	107	616037	44.07	ng	93
22) Acetophenone	7.30	105	815396	40.87	ng	# 93
24) Nitrobenzene	7.47	77	630658	37.80	ng	98
25) Isophorone	7.71	82	1259905	42.58	ng	100
26) 2-Nitrophenol	7.79	139	274861	36.31	ng	99
27) 2,4-Dimethylphenol	7.83	122	573452	45.65	ng	99
28) bis(2-Chloroethoxy)methane	7.93	93	771898	44.18	ng	99
29) 2,4-Dichlorophenol	8.04	162	438516	39.86	ng	98
30) 1,2,4-Trichlorobenzene	8.12	180	487309	37.82	ng	100
31) Naphthalene	8.20	128	1754691	38.14	ng	99
32) Benzoic acid	7.95	122	214585	27.81	ng	98
33) 4-Chloroaniline	8.25	127	254166	14.46	ng	98
34) Hexachlorobutadiene	8.31	225	270602	37.59	ng	100
35) Caprolactam	8.62	113	153317	49.93	ng	95
36) 4-Chloro-3-methylphenol	8.74	107	526329	41.46	ng	# 86
37) 2-Methylnaphthalene	8.89	142	1031200	38.17	ng	99
39) 1,2,4,5-Tetrachlorobenzene	9.06	216	511702	43.65	ng	99
40) Hexachlorocyclopentadiene	9.05	237	342084	59.73	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.17	196	321424	48.61	ng	98
43) 2,4,5-Trichlorophenol	9.22	196	318889	40.43	ng	# 85
45) 1,1'-Biphenyl	9.35	154	1326964	39.44	ng	100
46) 2-Chloronaphthalene	9.38	162	1003342	40.00	ng	100
47) 2-Nitroaniline	9.48	65	366200	40.26	ng	85
48) Acenaphthylene	9.80	152	1553233	38.88	ng	98
49) Dimethylphthalate	9.65	163	1146718	39.97	ng	100
50) 2,6-Dinitrotoluene	9.72	165	282313	44.97	ng	97
51) Acenaphthene	9.97	154	992408	37.38	ng	100
52) 3-Nitroaniline	9.89	138	199110	25.28	ng	89
53) 2,4-Dinitrophenol	9.99	184	47355	18.58	ng	90
54) Dibenzofuran	10.14	168	1399451	40.07	ng	98
55) 4-Nitrophenol	10.06	139	313631	66.16	ng	# 80
56) 2,4-Dinitrotoluene	10.12	165	332234	38.97	ng	91
57) Fluorene	10.49	166	1140745	39.59	ng	100
58) 2,3,4,6-Tetrachlorophenol	10.26	232	250592	38.14	ng	99
59) Diethylphthalate	10.36	149	1268675	43.61	ng	# 94
60) 4-Chlorophenyl-phenylether	10.47	204	562341	41.41	ng	98
61) 4-Nitroaniline	10.51	138	258522	32.73	ng	93
62) Azobenzene	10.63	77	1422473	42.27	ng	97
64) 4,6-Dinitro-2-methylphenol	10.53	198	41973	10.51	ng	96
65) n-Nitrosodiphenylamine	10.59	169	898855	41.26	ng	99
66) 4-Bromophenyl-phenylether	10.97	248	307404	43.43	ng	97
67) Hexachlorobenzene	11.03	284	329571	42.89	ng	# 63
68) Atrazine	11.13	200	292964	48.62	ng	91
69) Pentachlorophenol	11.23	266	269335	70.72	ng	98
70) Phenanthrene	11.45	178	1481943	42.09	ng	100
71) Anthracene	11.49	178	1467639	38.73	ng	99
72) Carbazole	11.65	167	1368326	40.34	ng	100
73) Di-n-butylphthalate	11.98	149	1844680	46.42	ng	99
74) Fluoranthene	12.63	202	1304724	36.84	ng	97
76) Benzidine	12.76	184	282142	24.99	ng	96
77) Pyrene	12.86	202	1289916	42.92	ng	99
79) Butylbenzylphthalate	13.48	149	692511	52.01	ng	97
80) Benzo(a)anthracene	14.05	228	1056821	40.91	ng	98
81) 3,3'-Dichlorobenzidine	14.01	252	289398	32.79	ng	# 95
82) Chrysene	14.09	228	1104168	44.36	ng	100
83) Bis(2-ethylhexyl)phthalate	14.04	149	945443	52.63	ng	100
84) Di-n-octyl phthalate	14.65	149	1339700	55.62	ng	96
85) Indeno(1,2,3-cd)pyrene	16.94	276	916936	63.72	ng	99
87) Benzo(b)fluoranthene	15.09	252	1046820	41.95	ng	# 94
88) Benzo(k)fluoranthene	15.12	252	901936m	33.59	ng	
89) Benzo(a)pyrene	15.45	252	927383	40.59	ng	99
90) Dibenzo(a,h)anthracene	16.96	278	786620	45.15	ng	99
91) Benzo(g,h,i)perylene	17.37	276	679690	40.91	ng	97

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

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