

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF031618\
 Data File : BF103687.D
 Acq On : 16 Mar 2018 11:03
 Operator : JU/SJ
 Sample : PB107245BS
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB107245BS

Manual Integrations
 APPROVED

Sohil
 3/19/2018 2:00:36 PM

Quant Time: Mar 16 14:17:29 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF031518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 16 13:24:19 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	160266	20.00	ng	0.00
21) Naphthalene-d8	8.25	136	659103	20.00	ng	0.00
38) Acenaphthene-d10	10.01	164	311120	20.00	ng	0.00
63) Phenanthrene-d10	11.50	188	566445	20.00	ng	0.00
75) Chrysene-d12	14.15	240	452925	20.00	ng	0.00
86) Perylene-d12	15.67	264	363021	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	1153388	109.46	ng	0.01
7) Phenol-d6	6.59	99	1422658	110.36	ng	0.00
23) Nitrobenzene-d5	7.53	82	866929	91.72	ng	0.00
41) 2,4,6-Tribromophenol	10.80	330	334320	124.98	ng	0.00
44) 2-Fluorobiphenyl	9.33	172	1488321	76.31	ng	0.00
78) Terphenyl-d14	13.09	244	1532576	78.54	ng	0.00

Target Compounds					Qvalue	
2) 1,4-Dioxane	2.90	88	173868	33.511	ng	93
3) Pyridine	3.66	79	481973	33.773	ng	90
4) n-Nitrosodimethylamine	3.62	42	209865	39.884	ng	87
6) Aniline	6.63	93	307479	16.704	ng	98
8) 2-Chlorophenol	6.75	128	439713	39.836	ng	98
9) Benzaldehyde	6.52	77	199889	23.602	ng	99
10) Phenol	6.60	94	530540	38.730	ng	99
11) bis(2-Chloroethyl)ether	6.70	93	422067	37.324	ng	99
12) 1,3-Dichlorobenzene	6.91	146	448761	35.696	ng	100
13) 1,4-Dichlorobenzene	6.99	146	455050	36.021	ng	99
14) 1,2-Dichlorobenzene	7.14	146	420629	35.882	ng	98
15) Benzyl Alcohol	7.10	79	388031	38.849	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.23	45	588864	36.612	ng	96
17) 2-Methylphenol	7.21	107	370224	37.664	ng	99
18) Hexachloroethane	7.48	117	169861	38.225	ng	97
19) n-Nitroso-di-n-propylamine	7.37	70	302848	36.658	ng	96
20) 3+4-Methylphenols	7.36	107	465161	37.289	ng	99
22) Acetophenone	7.37	105	599158	35.371	ng	# 95
24) Nitrobenzene	7.55	77	467532	42.135	ng	96
25) Isophorone	7.79	82	865522	39.559	ng	99
26) 2-Nitrophenol	7.86	139	209055	49.221	ng	99
27) 2,4-Dimethylphenol	7.89	122	418877	44.689	ng	99
28) bis(2-Chloroethoxy)methane	7.99	93	546461	41.246	ng	99
29) 2,4-Dichlorophenol	8.10	162	359479	42.155	ng	99
30) 1,2,4-Trichlorobenzene	8.19	180	359264	36.324	ng	99
31) Naphthalene	8.27	128	1232217	37.010	ng	100
32) Benzoic acid	7.92	122	13046m	11.471	ng	
33) 4-Chloroaniline	8.31	127	157816	11.298	ng	99
34) Hexachlorobutadiene	8.39	225	194072	35.788	ng	99
35) Caprolactam	8.68	113	109540m	37.304	ng	
36) 4-Chloro-3-methylphenol	8.79	107	392355	41.736	ng	99
37) 2-Methylnaphthalene	8.96	142	811639	38.441	ng	98
39) 1,2,4,5-Tetrachlorobenzene	9.13	216	323317	36.426	ng	98
40) Hexachlorocyclopentadiene	9.12	237	392404	85.678	ng	97

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF031618\
 Data File : BF103687.D
 Acq On : 16 Mar 2018 11:03
 Operator : JU/SJ
 Sample : PB107245BS
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB107245BS

Manual Integrations
 APPROVED

Sohil
 3/19/2018 2:00:36 PM

Quant Time: Mar 16 14:17:29 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF031518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 16 13:24:19 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.23	196	248651	43.798	ng	99
43) 2,4,5-Trichlorophenol	9.27	196	259485	40.495	ng	97
45) 1,1'-Biphenyl	9.43	154	944819	36.419	ng	99
46) 2-Chloronaphthalene	9.45	162	766029	37.176	ng	100
47) 2-Nitroaniline	9.55	65	241297	45.224	ng	95
48) Acenaphthylene	9.87	152	1173988	36.742	ng	99
49) Dimethylphthalate	9.73	163	906638	38.195	ng	99
50) 2,6-Dinitrotoluene	9.79	165	197935	43.512	ng	95
51) Acenaphthene	10.05	154	702373	37.021	ng	98
52) 3-Nitroaniline	9.96	138	119771	23.849	ng	96
53) 2,4-Dinitrophenol	10.06	184	25802	33.986	ng	# 33
54) Dibenzofuran	10.22	168	1053667	38.885	ng	98
55) 4-Nitrophenol	10.11	139	362041	83.103	ng	100
56) 2,4-Dinitrotoluene	10.19	165	269753	46.582	ng	98
57) Fluorene	10.56	166	794803	37.800	ng	99
58) 2,3,4,6-Tetrachlorophenol	10.33	232	197251	43.443	ng	98
59) Diethylphthalate	10.43	149	900875	38.238	ng	98
60) 4-Chlorophenyl-phenylether	10.55	204	364127	37.893	ng	99
61) 4-Nitroaniline	10.57	138	225427	42.986	ng	97
62) Azobenzene	10.71	77	925786	38.651	ng	96
64) 4,6-Dinitro-2-methylphenol	10.60	198	39612	26.412	ng	83
65) n-Nitrosodiphenylamine	10.67	169	738622	38.309	ng	100
66) 4-Bromophenyl-phenylether	11.04	248	225928	38.847	ng	95
67) Hexachlorobenzene	11.10	284	249258	37.551	ng	97
68) Atrazine	11.19	200	239009	40.073	ng	98
69) Pentachlorophenol	11.30	266	187898	49.236	ng	98
70) Phenanthrene	11.53	178	1167796	37.688	ng	99
71) Anthracene	11.57	178	1218945	38.732	ng	99
72) Carbazole	11.73	167	1145859	37.461	ng	99
73) Di-n-butylphthalate	12.05	149	1461626	39.824	ng	100
74) Fluoranthene	12.72	202	1248881	39.122	ng	100
76) Benzidine	12.83	184	451957	23.396	ng	99
77) Pyrene	12.95	202	1272395	38.253	ng	100
79) Butylbenzylphthalate	13.56	149	628827	42.670	ng	96
80) Benzo(a)anthracene	14.14	228	1112112	38.790	ng	99
81) 3,3'-Dichlorobenzidine	14.10	252	119279	12.438	ng	97
82) Chrysene	14.17	228	1041998	38.127	ng	100
83) Bis(2-ethylhexyl)phthalate	14.12	149	860658	40.880	ng	99
84) Di-n-octyl phthalate	14.74	149	1376486	44.941	ng	98
85) Indeno(1,2,3-cd)pyrene	17.23	276	800512	33.541	ng	97
87) Benzo(b)fluoranthene	15.22	252	978256	42.654	ng	98
88) Benzo(k)fluoranthene	15.26	252	872027	39.554	ng	99
89) Benzo(a)pyrene	15.61	252	865324	41.598	ng	98
90) Dibenzo(a,h)anthracene	17.26	278	673254	37.377	ng	99
91) Benzo(g,h,i)perylene	17.72	276	659461	37.633	ng	96

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

