

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF032018\
 Data File : BF103826.D
 Acq On : 21 Mar 2018 2:58
 Operator : JU/SJ
 Sample : PB107522BS
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB107522BS

Manual Integrations
 APPROVED

Sohil
 3/22/2018 12:58:00 PM

Quant Time: Mar 21 04:46:32 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF031518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 15 18:31:53 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	151996	20.00	ng	0.00
21) Naphthalene-d8	8.26	136	640312	20.00	ng	0.00
38) Acenaphthene-d10	10.02	164	271516	20.00	ng	0.00
63) Phenanthrene-d10	11.51	188	474465	20.00	ng	0.01
75) Chrysene-d12	14.17	240	305171	20.00	ng	0.02
86) Perylene-d12	15.72	264	288505	20.00	ng	0.04

System Monitoring Compounds

5) 2-Fluorophenol	5.60	112	1136970	113.77	ng	0.02
7) Phenol-d6	6.60	99	1449639	118.57	ng	0.01
23) Nitrobenzene-d5	7.54	82	823671	89.71	ng	0.00
41) 2,4,6-Tribromophenol	10.81	330	327713	140.38	ng	0.01
44) 2-Fluorobiphenyl	9.33	172	1482176	87.08	ng	0.00
78) Terphenyl-d14	13.10	244	1203259	91.52	ng	0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.90	88	171086	34.769	ng	# 89
3) Pyridine	3.66	79	485452	35.868	ng	# 87
4) n-Nitrosodimethylamine	3.62	42	178578	35.785	ng	# 73
6) Aniline	6.63	93	263570	15.098	ng	97
8) 2-Chlorophenol	6.76	128	437063	41.750	ng	97
9) Benzaldehyde	6.52	77	138842	17.286	ng	98
10) Phenol	6.61	94	525321	40.436	ng	96
11) bis(2-Chloroethyl)ether	6.70	93	415107	38.706	ng	97
12) 1,3-Dichlorobenzene	6.92	146	438621	36.788	ng	98
13) 1,4-Dichlorobenzene	6.99	146	443468	37.014	ng	97
14) 1,2-Dichlorobenzene	7.15	146	399579	35.941	ng	97
15) Benzyl Alcohol	7.11	79	344606	36.379	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.24	45	501750	32.893	ng	93
17) 2-Methylphenol	7.22	107	344015	36.902	ng	98
18) Hexachloroethane	7.48	117	149052	35.368	ng	95
19) n-Nitroso-di-n-propylamine	7.38	70	266848	34.057	ng	92
20) 3+4-Methylphenols	7.37	107	422945	35.749	ng	90
22) Acetophenone	7.38	105	536642	32.610	ng	# 93
24) Nitrobenzene	7.56	77	409014	37.943	ng	95
25) Isophorone	7.79	82	766842	36.077	ng	98
26) 2-Nitrophenol	7.87	139	216548	51.855	ng	93
27) 2,4-Dimethylphenol	7.90	122	411117	45.148	ng	97
28) bis(2-Chloroethoxy)methane	8.00	93	524567	40.755	ng	99
29) 2,4-Dichlorophenol	8.11	162	355021	42.854	ng	98
30) 1,2,4-Trichlorobenzene	8.19	180	360300	37.497	ng	99
31) Naphthalene	8.28	128	1215282	37.572	ng	99
32) Benzoic acid	8.00	122	203297	35.017	ng	95
33) 4-Chloroaniline	8.32	127	110712	8.159	ng	96
34) Hexachlorobutadiene	8.39	225	192342	36.510	ng	99
35) Caprolactam	8.70	113	118060m	41.386	ng	
36) 4-Chloro-3-methylphenol	8.80	107	385638	42.225	ng	98
37) 2-Methylnaphthalene	8.97	142	801616	39.080	ng	99
39) 1,2,4,5-Tetrachlorobenzene	9.13	216	331830	42.839	ng	99
40) Hexachlorocyclopentadiene	9.12	237	163544	40.917	ng	96

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.25	196	236457	47.725	ng	100
43) 2,4,5-Trichlorophenol	9.29	196	248755	44.483	ng	98
45) 1,1'-Biphenyl	9.43	154	884901	39.085	ng	100
46) 2-Chloronaphthalene	9.46	162	700451	38.952	ng	98
47) 2-Nitroaniline	9.56	65	210412	45.193	ng	93
48) Acenaphthylene	9.88	152	1069006	38.336	ng	100
49) Dimethylphthalate	9.73	163	834773	40.298	ng	99
50) 2,6-Dinitrotoluene	9.80	165	186278	46.611	ng	90
51) Acenaphthene	10.05	154	666235	40.238	ng	99
52) 3-Nitroaniline	9.96	138	92100	21.467	ng	99
53) 2,4-Dinitrophenol	10.07	184	78449	70.034	ng	# 16
54) Dibenzofuran	10.23	168	945402	39.979	ng	99
55) 4-Nitrophenol	10.13	139	317556	83.499	ng	88
56) 2,4-Dinitrotoluene	10.20	165	242698	47.831	ng	# 91
57) Fluorene	10.57	166	753066	41.039	ng	100
58) 2,3,4,6-Tetrachlorophenol	10.34	232	190352	48.038	ng	95
59) Diethylphthalate	10.43	149	839700	40.841	ng	98
60) 4-Chlorophenyl-phenylether	10.56	204	351539	41.919	ng	96
61) 4-Nitroaniline	10.59	138	182055	40.059	ng	93
62) Azobenzene	10.72	77	793947	37.982	ng	95
64) 4,6-Dinitro-2-methylphenol	10.61	198	66576	41.243	ng	86
65) n-Nitrosodiphenylamine	10.67	169	691568	42.822	ng	100
66) 4-Bromophenyl-phenylether	11.05	248	212442	43.609	ng	# 87
67) Hexachlorobenzene	11.12	284	222220	39.967	ng	95
68) Atrazine	11.20	200	214938	43.023	ng	99
69) Pentachlorophenol	11.31	266	219052	68.528	ng	97
70) Phenanthrene	11.54	178	994555	38.319	ng	99
71) Anthracene	11.59	178	1044409	39.620	ng	100
72) Carbazole	11.74	167	968616	37.805	ng	99
73) Di-n-butylphthalate	12.06	149	1249144	40.632	ng	100
74) Fluoranthene	12.73	202	908202	33.966	ng	99
76) Benzidine	12.84	184	462316	35.520	ng	99
77) Pyrene	12.96	202	889356	39.683	ng	98
79) Butylbenzylphthalate	13.57	149	451881	45.509	ng	94
80) Benzo(a)anthracene	14.16	228	784679	40.621	ng	99
81) 3,3'-Dichlorobenzidine	14.12	252	237351	36.733	ng	97
82) Chrysene	14.20	228	724103	39.323	ng	98
83) Bis(2-ethylhexyl)phthalate	14.13	149	641622	45.232	ng	99
84) Di-n-octyl phthalate	14.76	149	1030791	49.949	ng	95
85) Indeno(1,2,3-cd)pyrene	17.30	276	657886	40.911	ng	98
87) Benzo(b)fluoranthene	15.26	252	728445	39.965	ng	97
88) Benzo(k)fluoranthene	15.29	252	651385	37.177	ng	98
89) Benzo(a)pyrene	15.65	252	684201	41.386	ng	98
90) Dibenzo(a,h)anthracene	17.32	278	556825	38.898	ng	98
91) Benzo(g,h,i)perylene	17.79	276	533746	38.326	ng	95

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

