

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF031618\
 Data File : BF103700.D
 Acq On : 16 Mar 2018 17:05
 Operator : JU/SJ
 Sample : PB107409BS
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB107409BS

Manual Integrations
 APPROVED

Sohil
 3/19/2018 2:00:53 PM

Quant Time: Mar 19 10:35:16 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	166834	20.00	ng	0.00
21) Naphthalene-d8	8.25	136	684416	20.00	ng	0.00
38) Acenaphthene-d10	10.01	164	322422	20.00	ng	0.00
63) Phenanthrene-d10	11.50	188	594726	20.00	ng	0.00
75) Chrysene-d12	14.15	240	516888	20.00	ng	0.00
86) Perylene-d12	15.67	264	478593	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.60	112	1306519	119.11	ng	0.02
7) Phenol-d6	6.59	99	1619098	120.65	ng	0.00
23) Nitrobenzene-d5	7.53	82	997091	101.59	ng	0.00
41) 2,4,6-Tribromophenol	10.80	330	377384	136.13	ng	0.00
44) 2-Fluorobiphenyl	9.33	172	1652382	81.75	ng	0.00
78) Terphenyl-d14	13.09	244	1718051	77.15	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.90	88	197064	36.486	ng	93
3) Pyridine	3.66	79	558102	37.568	ng	90
4) n-Nitrosodimethylamine	3.62	42	240551	43.916	ng	90
6) Aniline	6.63	93	306943	16.018	ng	98
8) 2-Chlorophenol	6.75	128	487114	42.393	ng	98
9) Benzaldehyde	6.52	77	192785	21.867	ng	98
10) Phenol	6.60	94	594555	41.695	ng	96
11) bis(2-Chloroethyl)ether	6.70	93	473751	40.245	ng	95
12) 1,3-Dichlorobenzene	6.91	146	500682	38.258	ng	99
13) 1,4-Dichlorobenzene	6.99	146	505642	38.450	ng	99
14) 1,2-Dichlorobenzene	7.14	146	465243	38.125	ng	100
15) Benzyl Alcohol	7.10	79	432158	41.564	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.23	45	652914	38.996	ng	96
17) 2-Methylphenol	7.21	107	417041	40.757	ng	98
18) Hexachloroethane	7.48	117	188820	40.819	ng	98
19) n-Nitroso-di-n-propylamine	7.37	70	333686	38.800	ng	96
20) 3+4-Methylphenols	7.36	107	511720	39.406	ng	90
22) Acetophenone	7.37	105	655317	37.255	ng	# 95
24) Nitrobenzene	7.55	77	519196	45.061	ng	97
25) Isophorone	7.79	82	963899	42.426	ng	99
26) 2-Nitrophenol	7.86	139	254374	56.053	ng	98
27) 2,4-Dimethylphenol	7.89	122	465697	47.847	ng	98
28) bis(2-Chloroethoxy)methane	7.99	93	603488	43.866	ng	99
29) 2,4-Dichlorophenol	8.10	162	396851	44.816	ng	98
30) 1,2,4-Trichlorobenzene	8.19	180	399751	38.922	ng	99
31) Naphthalene	8.27	128	1357577	39.267	ng	99
32) Benzoic acid	7.96	122	75082m	18.582	ng	
33) 4-Chloroaniline	8.31	127	116596	8.039	ng	99
34) Hexachlorobutadiene	8.38	225	214125	38.026	ng	98
35) Caprolactam	8.69	113	123689m	40.565	ng	
36) 4-Chloro-3-methylphenol	8.79	107	436992	44.765	ng	99
37) 2-Methylnaphthalene	8.96	142	901598	41.122	ng	99
39) 1,2,4,5-Tetrachlorobenzene	9.13	216	354163	38.503	ng	98
40) Hexachlorocyclopentadiene	9.12	237	449568	94.718	ng	98

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42) 2,4,6-Trichlorophenol	9.23	196	273869	46.549	nq	99
43) 2,4,5-Trichlorophenol	9.27	196	289012	43.522	nq	97
45) 1,1'-Biphenyl	9.43	154	1037062	38.574	nq	99
46) 2-Chloronaphthalene	9.45	162	833648	39.040	nq	100
47) 2-Nitroaniline	9.55	65	270253	48.325	nq	98
48) Acenaphthylene	9.87	152	1280341	38.666	nq	99
49) Dimethylphthalate	9.73	163	979494	39.818	nq	100
50) 2,6-Dinitrotoluene	9.79	165	219934	46.366	nq	96
51) Acenaphthene	10.05	154	809191	41.156	nq	99
52) 3-Nitroaniline	9.96	138	132726	25.238	nq	98
53) 2,4-Dinitrophenol	10.06	184	105925	75.328	nq #	23
54) Dibenzofuran	10.22	168	1147421	40.861	nq	98
55) 4-Nitrophenol	10.12	139	421464	92.739	nq	95
56) 2,4-Dinitrotoluene	10.19	165	301658	49.774	nq	97
57) Fluorene	10.56	166	866550	39.768	nq	98
58) 2,3,4,6-Tetrachlorophenol	10.33	232	228850	48.636	nq	97
59) Diethylphthalate	10.43	149	988068	40.469	nq	98
60) 4-Chlorophenyl-phenylether	10.55	204	394700	39.635	nq	99
61) 4-Nitroaniline	10.58	138	252649	46.183	nq	99
62) Azobenzene	10.71	77	995336	40.098	nq	96
64) 4,6-Dinitro-2-methylphenol	10.60	198	101422	46.925	nq	89
65) n-Nitrosodiphenylamine	10.67	169	794890	39.267	nq	99
66) 4-Bromophenyl-phenylether	11.04	248	244193	39.991	nq	96
67) Hexachlorobenzene	11.10	284	265656	38.118	nq	97
68) Atrazine	11.19	200	251575	40.174	nq	99
69) Pentachlorophenol	11.30	266	272742	68.070	nq	97
70) Phenanthrene	11.53	178	1269545	39.023	nq	99
71) Anthracene	11.57	178	1300531	39.360	nq	100
72) Carbazole	11.73	167	1254689	39.068	nq	98
73) Di-n-butylphthalate	12.05	149	1549665	40.214	nq	99
74) Fluoranthene	12.72	202	1339178	39.956	nq	100
76) Benzidine	12.83	184	577488	26.195	nq	99
77) Pyrene	12.95	202	1337783	35.242	nq	100
79) Butylbenzylphthalate	13.56	149	698864	41.554	nq	97
80) Benzo(a)anthracene	14.14	228	1260933	38.538	nq	100
81) 3,3'-Dichlorobenzidine	14.10	252	167308	15.287	nq	98
82) Chrysene	14.18	228	1165359	37.364	nq	99
83) Bis(2-ethylhexyl)phthalate	14.12	149	957553	39.854	nq	100
84) Di-n-octyl phthalate	14.74	149	1609648	46.051	nq	98
85) Indeno(1,2,3-cd)pyrene	17.24	276	1052128	38.628	nq	98
87) Benzo(b)fluoranthene	15.23	252	1179803	39.019	nq	98
88) Benzo(k)fluoranthene	15.26	252	1115514	38.380	nq	100
89) Benzo(a)pyrene	15.62	252	1105469	40.309	nq	98
90) Dibenzo(a,h)anthracene	17.26	278	879334	37.029	nq	98
91) Benzo(g,h,i)perylene	17.72	276	845478	36.597	nq	95

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

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