

Data Path : Z:\HPCHEM1\BNA F\DATA\BF092517\
 Data File : BF098814.D
 Acq On : 26 Sep 2017 6:38
 Operator : SJ/JU
 Sample : PB102566BS
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB102566BS

Manual Integrations
 APPROVED

Sohil
 9/26/2017 4:58:59 PM

Quant Time: Sep 26 09:01:08 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 23 13:50:58 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.92	152	137351	20.00	ng	0.00
21) Naphthalene-d8	8.20	136	585966	20.00	ng	0.00
38) Acenaphthene-d10	9.96	164	263717	20.00	ng	0.00
63) Phenanthrene-d10	11.44	188	443862	20.00	ng	0.00
75) Chrysene-d12	14.08	240	251004	20.00	ng	0.00
86) Perylene-d12	15.57	264	239884	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.54	112	800365	92.76	ng	0.00
7) Phenol-d6	6.54	99	967079	90.86	ng	0.00
23) Nitrobenzene-d5	7.48	82	846961	92.69	ng	0.00
41) 2,4,6-Tribromophenol	10.74	330	196697	91.15	ng	0.00
44) 2-Fluorobiphenyl	9.27	172	1536679	97.28	ng	0.00
78) Terphenyl-d14	13.02	244	1104748	100.09	ng	-0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.79	88	144152	34.975	ng	97
3) Pyridine	3.56	79	413448	37.082	ng	98
4) n-Nitrosodimethylamine	3.51	42	190373	40.265	ng	96
6) Aniline	6.57	93	487990	33.970	ng	99
8) 2-Chlorophenol	6.70	128	414391	42.410	ng	99
9) Benzaldehyde	6.46	77	98876	16.015	ng	99
10) Phenol	6.55	94	503555	42.302	ng	98
11) bis(2-Chloroethyl)ether	6.65	93	393071	42.475	ng	98
12) 1,3-Dichlorobenzene	6.86	146	432843	42.299	ng	98
13) 1,4-Dichlorobenzene	6.93	146	435418	41.821	ng	98
14) 1,2-Dichlorobenzene	7.09	146	402750	41.866	ng	98
15) Benzyl Alcohol	7.05	79	353717	45.300	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.19	45	614930	42.650	ng	98
17) 2-Methylphenol	7.16	107	354440	44.333	ng	97
18) Hexachloroethane	7.43	117	147752	39.482	ng	97
19) n-Nitroso-di-n-propylamine	7.33	70	281273	41.391	ng	96
20) 3+4-Methylphenols	7.32	107	431138	42.627	ng	# 82
22) Acetophenone	7.32	105	559836	39.434	ng	# 99
24) Nitrobenzene	7.50	77	427542	41.249	ng	97
25) Isophorone	7.73	82	784672	41.537	ng	99
26) 2-Nitrophenol	7.81	139	210818	42.297	ng	96
27) 2,4-Dimethylphenol	7.85	122	393712	41.827	ng	98
28) bis(2-Chloroethoxy)methane	7.94	93	504298	42.836	ng	100
29) 2,4-Dichlorophenol	8.05	162	348760	43.155	ng	98
30) 1,2,4-Trichlorobenzene	8.14	180	341417	42.239	ng	99
31) Naphthalene	8.22	128	1167529	42.424	ng	100
32) Benzoic acid	7.96	122	209263	27.065	ng	97
33) 4-Chloroaniline	8.26	127	344480	29.974	ng	99
34) Hexachlorobutadiene	8.33	225	172938	41.539	ng	100
35) Caprolactam	8.64	113	116916m	45.100	ng	
36) 4-Chloro-3-methylphenol	8.74	107	374630	41.242	ng	99
37) 2-Methylnaphthalene	8.91	142	807085	45.112	ng	99
39) 1,2,4,5-Tetrachlorobenzene	9.07	216	313773	39.896	ng	99
40) Hexachlorocyclopentadiene	9.06	237	160189	44.569	ng	97

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.19	196	238876	42.873	nq	100
43) 2,4,5-Trichlorophenol	9.23	196	240204	43.187	nq	95
45) 1,1'-Biphenyl	9.37	154	921071	40.639	nq	100
46) 2-Chloronaphthalene	9.40	162	725148	42.612	nq	97
47) 2-Nitroaniline	9.49	65	234035	44.914	nq	98
48) Acenaphthylene	9.82	152	1100259	42.235	nq	100
49) Dimethylphthalate	9.67	163	870031	43.712	nq	100
50) 2,6-Dinitrotoluene	9.73	165	193431	44.639	nq	99
51) Acenaphthene	9.99	154	674992	40.904	nq	97
52) 3-Nitroaniline	9.90	138	175678	34.770	nq	96
53) 2,4-Dinitrophenol	10.00	184	37958	21.428	nq	95
54) Dibenzofuran	10.16	168	1001948	45.602	nq	99
55) 4-Nitrophenol	10.06	139	346754	81.164	nq	91
56) 2,4-Dinitrotoluene	10.14	165	252846	44.961	nq	# 95
57) Fluorene	10.50	166	750760	42.512	nq	100
58) 2,3,4,6-Tetrachlorophenol	10.27	232	176115	45.838	nq	98
59) Diethylphthalate	10.37	149	858873	44.160	nq	97
60) 4-Chlorophenyl-phenylether	10.49	204	343061	43.246	nq	98
61) 4-Nitroaniline	10.52	138	197547	40.527	nq	90
62) Azobenzene	10.65	77	814337	45.537	nq	98
64) 4,6-Dinitro-2-methylphenol	10.55	198	36857	13.648	nq	83
65) n-Nitrosodiphenylamine	10.61	169	697866	45.385	nq	99
66) 4-Bromophenyl-phenylether	10.99	248	199468	45.911	nq	# 90
67) Hexachlorobenzene	11.05	284	194684	41.955	nq	95
68) Atrazine	11.14	200	218889	47.313	nq	99
69) Pentachlorophenol	11.24	266	189390	65.580	nq	98
70) Phenanthrene	11.47	178	1043328	43.913	nq	99
71) Anthracene	11.52	178	1080000	44.427	nq	99
72) Carbazole	11.67	167	1014051	42.597	nq	99
73) Di-n-butylphthalate	11.99	149	1325413	46.255	nq	99
74) Fluoranthene	12.65	202	951093	39.533	nq	99
76) Benzidine	12.77	184	618618	66.634	nq	98
77) Pyrene	12.89	202	938287	49.022	nq	99
79) Butylbenzylphthalate	13.49	149	486279	54.059	nq	98
80) Benzo(a)anthracene	14.07	228	678166	44.619	nq	100
81) 3,3'-Dichlorobenzidine	14.03	252	238863	42.870	nq	98
82) Chrysene	14.10	228	635990	43.952	nq	100
83) Bis(2-ethylhexyl)phthalate	14.05	149	672624	54.305	nq	99
84) Di-n-octyl phthalate	14.66	149	1011712	50.727	nq	98
85) Indeno(1,2,3-cd)pyrene	17.08	276	571833	49.402	nq	96
87) Benzo(b)fluoranthene	15.13	252	658745m	44.664	nq	
88) Benzo(k)fluoranthene	15.16	252	590170	40.512	nq	98
89) Benzo(a)pyrene	15.50	252	597487	44.132	nq	# 96
90) Dibenzo(a,h)anthracene	17.09	278	491060	45.322	nq	97
91) Benzo(g,h,i)perylene	17.53	276	448542	41.881	nq	97

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

