

Data Path : Z:\HPCHEM1\BNA F\DATA\BF061115\
 Data File : BF079892.D
 Acq On : 11 Jun 2015 13:38
 Operator : TP/IZ
 Sample : SSTDICCC040
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICCC040

Manual Integrations
 APPROVED

apatel
 6/13/2015 8:43:13 AM

Quant Time: Jun 12 01:24:15 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF061115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 12 00:59:58 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	39922	20.00	ng	0.00
21) Naphthalene-d8	8.68	136	168986	20.00	ng	0.00
38) Acenaphthene-d10	10.46	164	88661	20.00	ng	0.00
63) Phenanthrene-d10	11.98	188	194433	20.00	ng	0.00
75) Chrysene-d12	14.99	240	231855	20.00	ng	0.00
86) Perylene-d12	16.93	264	220048	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	6.00	112	177611	74.54	ng	0.00
7) Phenol-d6	6.99	99	254396	81.62	ng	0.00
23) Nitrobenzene-d5	7.95	82	210281	74.02	ng	0.00
41) 2,4,6-Tribromophenol	11.26	330	73290	92.62	ng	0.00
44) 2-Fluorobiphenyl	9.76	172	486395	77.32	ng	0.00
78) Terphenyl-d14	13.69	244	785904	77.02	ng	0.00

Target Compounds					Qvalue	
2) 1,4-Dioxane	3.34	88	40438	37.11	ng	100
3) Pyridine	4.12	79	94837	33.28	ng	100
4) n-Nitrosodimethylamine	4.03	42	45884	32.27	ng	100
6) Aniline	7.05	93	180962	40.31	ng	100
8) 2-Chlorophenol	7.18	128	110887	40.42	ng	100
9) Benzaldehyde	6.94	77	64145	35.81	ng	100
10) Phenol	7.00	94	140333	41.97	ng	100
11) bis(2-Chloroethyl)ether	7.11	93	113358	42.76	ng	100
12) 1,3-Dichlorobenzene	7.34	146	128276	40.91	ng	100
13) 1,4-Dichlorobenzene	7.42	146	131667	37.57	ng	100
14) 1,2-Dichlorobenzene	7.56	146	116998	38.47	ng	100
15) Benzyl Alcohol	7.52	79	100046	40.15	ng	100
16) 2,2'-oxybis(1-Chloropropan	7.66	45	162907	40.27	ng	100
17) 2-Methylphenol	7.62	107	97123	41.50	ng	100
18) Hexachloroethane	7.92	117	41278	37.72	ng	100
19) n-Nitroso-di-n-propylamine	7.78	70	89768	40.67	ng	100
20) 3+4-Methylphenols	7.77	107	126358	41.60	ng	100
22) Acetophenone	7.79	105	166010	39.80	ng	100
24) Nitrobenzene	7.96	77	112305	38.53	ng	100
25) Isophorone	8.20	82	248409	40.55	ng	100
26) 2-Nitrophenol	8.30	139	35561	31.87	ng	100
27) 2,4-Dimethylphenol	8.31	122	105631	38.52	ng	100
28) bis(2-Chloroethoxy)methane	8.41	93	145707	39.72	ng	100
29) 2,4-Dichlorophenol	8.52	162	89791	38.71	ng	100
30) 1,2,4-Trichlorobenzene	8.63	180	113841	38.24	ng	100
31) Naphthalene	8.71	128	376000	41.61	ng	100
32) Benzoic acid	8.36	122	38864	46.85	ng	100
33) 4-Chloroaniline	8.74	127	149048m	40.95	ng	
34) Hexachlorobutadiene	8.83	225	61832	37.91	ng	100
35) Caprolactam	9.08	113	32577	46.35	ng	100
36) 4-Chloro-3-methylphenol	9.21	107	109226	42.85	ng	100
37) 2-Methylnaphthalene	9.40	142	252696	39.52	ng	100
39) 1,2,4,5-Tetrachlorobenzene	9.58	216	117534	38.37	ng	100
40) Hexachlorocyclopentadiene	9.56	237	41311	29.28	ng	100

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.68	196	76852	41.09	nq	100
43) 2,4,5-Trichlorophenol	9.71	196	84962	43.33	nq	100
45) 1,1'-Biphenyl	9.86	154	283464	38.81	nq	100
46) 2-Chloronaphthalene	9.90	162	241390	40.08	nq	100
47) 2-Nitroaniline	9.98	65	49224	37.59	nq	100
48) Acenaphthylene	10.32	152	415891	40.44	nq	100
49) Dimethylphthalate	10.15	163	282994	40.50	nq	100
50) 2,6-Dinitrotoluene	10.22	165	51577	42.31	nq	100
51) Acenaphthene	10.49	154	232818	39.08	nq	100
52) 3-Nitroaniline	10.39	138	62230	41.64	nq	100
53) 2,4-Dinitrophenol	10.49	184	6929	26.50	nq	100
54) Dibenzofuran	10.66	168	345554	41.11	nq	100
55) 4-Nitrophenol	10.52	139	43398	40.33	nq	100
56) 2,4-Dinitrotoluene	10.63	165	66529	42.98	nq	100
57) Fluorene	11.02	166	312446	41.86	nq	100
58) 2,3,4,6-Tetrachlorophenol	10.78	232	67095	46.01	nq	100
59) Diethylphthalate	10.86	149	295954	43.07	nq	100
60) 4-Chlorophenyl-phenylether	10.99	204	133606	41.24	nq	100
61) 4-Nitroaniline	11.01	138	67533	45.59	nq	100
62) Azobenzene	11.15	77	280276	41.67	nq	100
64) 4,6-Dinitro-2-methylphenol	11.04	198	15640	29.84	nq	100
65) n-Nitrosodiphenylamine	11.11	169	258922	39.82	nq	100
66) 4-Bromophenyl-phenylether	11.50	248	91181	40.09	nq	100
67) Hexachlorobenzene	11.58	284	99759	42.14	nq	100
68) Atrazine	11.63	200	85231	45.58	nq	100
69) Pentachlorophenol	11.77	266	38563	43.77	nq	100
70) Phenanthrene	12.00	178	483942	41.23	nq	100
71) Anthracene	12.06	178	465349	41.04	nq	100
72) Carbazole	12.21	167	465462	43.48	nq	100
73) Di-n-butylphthalate	12.54	149	503683	45.40	nq	100
74) Fluoranthene	13.28	202	544462	44.13	nq	100
76) Benzidine	13.39	184	251207	37.27	nq	100
77) Pyrene	13.54	202	565157	39.37	nq	100
79) Butylbenzylphthalate	14.25	149	210052	42.90	nq	100
80) Benzo(a)anthracene	14.98	228	563015	40.88	nq	100
81) 3,3'-Dichlorobenzidine	14.93	252	194572	44.69	nq	100
82) Chrysene	15.03	228	532317	41.21	nq	100
83) Bis(2-ethylhexyl)phthalate	14.96	149	335424	44.00	nq	100
84) Di-n-octyl phthalate	15.78	149	561882	46.84	nq	100
85) Indeno(1,2,3-cd)pyrene	18.86	276	603625	43.81	nq	100
87) Benzo(b)fluoranthene	16.38	252	523267	40.00	nq	100
88) Benzo(k)fluoranthene	16.41	252	577316	42.40	nq	100
89) Benzo(a)pyrene	16.85	252	514346	41.61	nq	100
90) Dibenzo(a,h)anthracene	18.88	278	487048	42.62	nq	100
91) Benzo(g,h,i)perylene	19.43	276	503675	41.68	nq	100

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

