

Data Path : Z:\HPCHEM1\BNA F\DATA\BF102915\
 Data File : BF082577.D
 Acq On : 29 Oct 2015 18:31
 Operator : UM/IZ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040EC

Quant Time: Oct 30 06:04:49 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF102615.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 29 19:06:23 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.63	152	105145	20.00	ng	0.00
21) Naphthalene-d8	7.92	136	432249	20.00	ng	0.00
38) Acenaphthene-d10	9.67	164	215602	20.00	ng	0.00
63) Phenanthrene-d10	11.15	188	421564	20.00	ng	0.00
75) Chrysene-d12	13.79	240	305977	20.00	ng	0.00
86) Perylene-d12	15.21	264	225924	20.00	ng	0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.20	112	472653	88.74	ng	0.00
7) Phenol-d6	6.28	99	580734	83.66	ng	0.00
23) Nitrobenzene-d5	7.21	82	559006	83.43	ng	0.00
41) 2,4,6-Tribromophenol	10.47	330	128988	79.89	ng	0.00
44) 2-Fluorobiphenyl	8.99	172	1027100	75.75	ng	0.00
78) Terphenyl-d14	12.74	244	991844	93.39	ng	0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.06	88	112970	42.26	ng	99
3) Pyridine	2.71	79	252400	40.50	ng	99
4) n-Nitrosodimethylamine	2.70	42	127380	46.48	ng	94
6) Aniline	6.30	93	379472	43.38	ng	91
8) 2-Chlorophenol	6.41	128	275877	44.16	ng	94
9) Benzaldehyde	6.18	77	141048	41.08	ng	93
10) Phenol	6.31	94	328455	41.68	ng	89
11) bis(2-Chloroethyl)ether	6.38	93	280759	44.89	ng	95
12) 1,3-Dichlorobenzene	6.56	146	336064	45.10	ng	98
13) 1,4-Dichlorobenzene	6.64	146	337191	43.98	ng	99
14) 1,2-Dichlorobenzene	6.80	146	293839	41.29	ng	96
15) Benzyl Alcohol	6.80	79	224470	46.45	ng	# 88
16) 2,2'-oxybis(1-Chloropropan	6.91	45	349599	41.28	ng	# 33
17) 2-Methylphenol	6.91	107	216184	43.15	ng	93
18) Hexachloroethane	7.13	117	109042	41.78	ng	# 82
19) n-Nitroso-di-n-propylamine	7.06	70	203513	45.00	ng	99
20) 3+4-Methylphenols	7.07	107	298514	46.64	ng	91
22) Acetophenone	7.06	105	372776	39.75	ng	# 93
24) Nitrobenzene	7.23	77	292039	39.03	ng	94
25) Isophorone	7.47	82	548788	41.23	ng	97
26) 2-Nitrophenol	7.54	139	166017	45.93	ng	# 71
27) 2,4-Dimethylphenol	7.60	122	222625	37.90	ng	97
28) bis(2-Chloroethoxy)methane	7.68	93	320317	41.04	ng	98
29) 2,4-Dichlorophenol	7.79	162	274203	46.80	ng	98
30) 1,2,4-Trichlorobenzene	7.86	180	266434	40.22	ng	99
31) Naphthalene	7.94	128	758889	38.42	ng	99
32) Benzoic acid	7.78	122	127488	38.74	ng	97
33) 4-Chloroaniline	8.01	127	311683	40.70	ng	99
34) Hexachlorobutadiene	8.04	225	151895	39.64	ng	98
35) Caprolactam	8.41	113	69363	42.59	ng	# 84
36) 4-Chloro-3-methylphenol	8.50	107	251000	44.06	ng	90
37) 2-Methylnaphthalene	8.63	142	537396	40.27	ng	98
39) 1,2,4,5-Tetrachlorobenzene	8.80	216	266365	40.93	ng	100
40) Hexachlorocyclopentadiene	8.78	237	25800	26.64	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	8.92	196	162268	40.00	nq	97
43) 2,4,5-Trichlorophenol	8.97	196	186512	43.85	nq	97
45) 1,1'-Biphenyl	9.10	154	648207	40.40	nq	98
46) 2-Chloronaphthalene	9.12	162	510948	40.40	nq	96
47) 2-Nitroaniline	9.23	65	166707	45.39	nq	99
48) Acenaphthylene	9.53	152	798316	39.41	nq	100
49) Dimethylphthalate	9.40	163	602426	40.16	nq	99
50) 2,6-Dinitrotoluene	9.48	165	126140	38.35	nq	94
51) Acenaphthene	9.70	154	460195	39.62	nq	99
52) 3-Nitroaniline	9.66	138	143577	40.23	nq	100
53) 2,4-Dinitrophenol	9.78	184	55244	43.85	nq	91
54) Dibenzofuran	9.87	168	678552	40.03	nq	97
55) 4-Nitrophenol	9.85	139	32407m	29.38	nq	
56) 2,4-Dinitrotoluene	9.88	165	181442	45.48	nq	# 84
57) Fluorene	10.22	166	540547	38.85	nq	99
58) 2,3,4,6-Tetrachlorophenol	10.01	232	129125	39.65	nq	93
59) Diethylphthalate	10.10	149	565947	38.76	nq	97
60) 4-Chlorophenyl-phenylether	10.22	204	248032	37.90	nq	97
61) 4-Nitroaniline	10.27	138	137230	38.39	nq	92
62) Azobenzene	10.38	77	574202	40.72	nq	93
64) 4,6-Dinitro-2-methylphenol	10.30	198	94850	38.35	nq	# 76
65) n-Nitrosodiphenylamine	10.34	169	506381	39.70	nq	99
66) 4-Bromophenyl-phenylether	10.70	248	154633	39.09	nq	95
67) Hexachlorobenzene	10.76	284	163026	38.64	nq	98
68) Atrazine	10.87	200	133254	33.75	nq	98
69) Pentachlorophenol	10.97	266	57829	37.84	nq	98
70) Phenanthrene	11.18	178	815696	37.60	nq	100
71) Anthracene	11.23	178	800769	36.65	nq	99
72) Carbazole	11.39	167	757976	39.34	nq	98
73) Di-n-butylphthalate	11.71	149	852307	37.09	nq	100
74) Fluoranthene	12.36	202	845554	37.97	nq	98
76) Benzidine	12.50	184	319029	37.99	nq	98
77) Pyrene	12.59	202	840707	44.67	nq	100
79) Butylbenzylphthalate	13.21	149	350625	44.31	nq	94
80) Benzo(a)anthracene	13.78	228	672522	41.00	nq	100
81) 3,3'-Dichlorobenzidine	13.75	252	227646	40.02	nq	# 97
82) Chrysene	13.83	228	638563	41.21	nq	99
83) Bis(2-ethylhexyl)phthalate	13.77	149	451817	42.64	nq	98
84) Di-n-octyl phthalate	14.39	149	635810	36.36	nq	99
85) Indeno(1,2,3-cd)pyrene	16.51	276	505285	35.49	nq	100
87) Benzo(b)fluoranthene	14.82	252	534798m	39.86	nq	
88) Benzo(k)fluoranthene	14.85	252	461316m	42.18	nq	
89) Benzo(a)pyrene	15.14	252	478706	41.18	nq	# 97
90) Dibenzo(a,h)anthracene	16.50	278	422581	43.45	nq	97
91) Benzo(g,h,i)perylene	16.90	276	411754	42.79	nq	98

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

