

Data Path : Z:\HPCHEM1\BNA F\DATA\BF120815\
 Data File : BF083605.D
 Acq On : 8 Dec 2015 20:21
 Operator : UM/IZ
 Sample : PB87006BS
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB87006BS

Manual Integrations
 APPROVED

Mmdadoda
 12/9/2015 6:13:23 PM

Quant Time: Dec 17 10:17:57 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF120815.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 17 09:56:43 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	5.70	152	134070	20.00	ng	0.00
21) Naphthalene-d8	7.60	136	537524	20.00	ng	-0.01
38) Acenaphthene-d10	10.38	164	266682	20.00	ng	0.00
63) Phenanthrene-d10	12.79	188	509454	20.00	ng	0.00
75) Chrysene-d12	16.99	240	423257	20.00	ng	0.00
86) Perylene-d12	18.64	264	324311	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	3.98	112	802016	90.73	ng	0.01
7) Phenol-d6	5.28	99	1034336	87.57	ng	0.00
23) Nitrobenzene-d5	6.53	82	964215	83.70	ng	0.00
41) 2,4,6-Tribromophenol	11.68	330	212704	89.53	ng	-0.01
44) 2-Fluorobiphenyl	9.34	172	1568131	82.76	ng	0.00
78) Terphenyl-d14	15.41	244	1563621	86.92	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	1.86	88	141694	32.45	ng	# 90
3) Pyridine	2.30	79	355476	33.50	ng	98
4) n-Nitrosodimethylamine	2.26	42	230575	41.28	ng	83
6) Aniline	5.26	93	411410	28.27	ng	95
8) 2-Chlorophenol	5.42	128	407515	42.15	ng	95
9) Benzaldehyde	5.13	77	85437	11.31	ng	97
10) Phenol	5.29	94	611469	42.74	ng	82
11) bis(2-Chloroethyl)ether	5.36	93	405520	38.53	ng	98
12) 1,3-Dichlorobenzene	5.61	146	432927m	40.18	ng	
13) 1,4-Dichlorobenzene	5.72	146	445182	40.27	ng	99
14) 1,2-Dichlorobenzene	5.94	146	430986	42.01	ng	98
15) Benzyl Alcohol	5.95	79	406713	48.40	ng	98
16) 2,2'-oxybis(1-Chloropropan	6.13	45	787076	40.44	ng	78
17) 2-Methylphenol	6.14	107	342001	43.48	ng	93
18) Hexachloroethane	6.41	117	159011	40.74	ng	# 86
19) n-Nitroso-di-n-propylamine	6.34	70	353010	42.82	ng	98
20) 3+4-Methylphenols	6.38	107	444506	42.92	ng	100
22) Acetophenone	6.32	105	583248	38.68	ng	# 98
24) Nitrobenzene	6.56	77	484794	38.08	ng	98
25) Isophorone	6.93	82	895269	40.05	ng	99
26) 2-Nitrophenol	7.04	139	209827	41.50	ng	92
27) 2,4-Dimethylphenol	7.16	122	381434	43.13	ng	97
28) bis(2-Chloroethoxy)methane	7.29	93	536014	43.18	ng	97
29) 2,4-Dichlorophenol	7.45	162	343009	42.65	ng	97
30) 1,2,4-Trichlorobenzene	7.53	180	353878	39.83	ng	97
31) Naphthalene	7.63	128	1164595	40.91	ng	99
32) Benzoic acid	7.48	122	205623	37.02	ng	96
33) 4-Chloroaniline	7.78	127	201476	19.41	ng	93
34) Hexachlorobutadiene	7.84	225	208030	40.08	ng	98
35) Caprolactam	8.37	113	111925m	46.37	ng	
36) 4-Chloro-3-methylphenol	8.62	107	392941	43.49	ng	93
37) 2-Methylnaphthalene	8.73	142	777296	43.47	ng	99
39) 1,2,4,5-Tetrachlorobenzene	9.00	216	345931	39.51	ng	# 100
40) Hexachlorocyclopentadiene	8.98	237	141093	72.13	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.23	196	212071	40.71	nq	99
43) 2,4,5-Trichlorophenol	9.32	196	222814	43.20	nq	98
45) 1,1'-Biphenyl	9.49	154	916971	39.56	nq	100
46) 2-Chloronaphthalene	9.50	162	719324	41.10	nq	96
47) 2-Nitroaniline	9.72	65	276001	43.73	nq	97
48) Acenaphthylene	10.16	152	1172841	40.64	nq	100
49) Dimethylphthalate	10.04	163	896347	42.22	nq	99
50) 2,6-Dinitrotoluene	10.13	165	191502	43.69	nq	87
51) Acenaphthene	10.44	154	680190	41.17	nq	99
52) 3-Nitroaniline	10.41	138	109433	23.49	nq	85
53) 2,4-Dinitrophenol	10.60	184	168637	77.90	nq	100
54) Dibenzofuran	10.73	168	1042557	45.51	nq	97
55) 4-Nitrophenol	10.81	139	190828	93.05	nq	93
56) 2,4-Dinitrotoluene	10.80	165	262303	45.12	nq	# 82
57) Fluorene	11.28	166	842763	42.15	nq	99
58) 2,3,4,6-Tetrachlorophenol	10.97	232	201405	48.46	nq	# 100
59) Diethylphthalate	11.18	149	858189	41.58	nq	98
60) 4-Chlorophenyl-phenylether	11.30	204	404356	42.48	nq	99
61) 4-Nitroaniline	11.40	138	178527	42.19	nq	87
62) Azobenzene	11.56	77	1084076	47.72	nq	97
64) 4,6-Dinitro-2-methylphenol	11.46	198	135057	39.28	nq	88
65) n-Nitrosodiphenylamine	11.52	169	716065	42.59	nq	99
66) 4-Bromophenyl-phenylether	12.09	248	227336	40.01	nq	97
67) Hexachlorobenzene	12.17	284	256031	42.47	nq	# 91
68) Atrazine	12.43	200	212177	42.41	nq	100
69) Pentachlorophenol	12.53	266	152369	91.34	nq	98
70) Phenanthrene	12.82	178	1241567	42.42	nq	99
71) Anthracene	12.91	178	1292461	42.62	nq	100
72) Carbazole	13.21	167	1136043	43.69	nq	99
73) Di-n-butylphthalate	13.84	149	1425578	43.59	nq	99
74) Fluoranthene	14.75	202	1278439	42.69	nq	99
76) Benzidine	15.04	184	208182	15.81	nq	98
77) Pyrene	15.11	202	1320721	43.34	nq	99
79) Butylbenzylphthalate	16.24	149	576440	42.62	nq	91
80) Benzo(a)anthracene	16.98	228	1060785	42.82	nq	99
81) 3,3'-Dichlorobenzidine	16.97	252	277991	33.22	nq	97
82) Chrysene	17.03	228	1038543	41.82	nq	98
83) Bis(2-ethylhexyl)phthalate	17.09	149	818374	43.59	nq	99
84) Di-n-octyl phthalate	17.86	149	1283321	44.74	nq	# 100
85) Indeno(1,2,3-cd)pyrene	19.86	276	684574	40.84	nq	# 100
87) Benzo(b)fluoranthene	18.25	252	804068m	43.54	nq	
88) Benzo(k)fluoranthene	18.28	252	908351m	44.64	nq	
89) Benzo(a)pyrene	18.58	252	787666	43.71	nq	# 95
90) Dibenzo(a,h)anthracene	19.86	278	585245	42.48	nq	99
91) Benzo(g,h,i)perylene	20.23	276	543724	40.47	nq	96

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

