

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF071916\
 Data File : BF089116.D
 Acq On : 19 Jul 2016 14:13
 Operator : UM/SJ
 Sample : PB92230BS
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB92230BS

Quant Time: Jul 20 03:11:35 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF063016.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 19 15:28:17 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.63	152	50837	20.00	ng	0.00
21) Naphthalene-d8	7.91	136	196766	20.00	ng	0.00
38) Acenaphthene-d10	9.66	164	90162	20.00	ng	0.00
63) Phenanthrene-d10	11.13	188	155420	20.00	ng	0.00
75) Chrysene-d12	13.75	240	88679	20.00	ng	-0.01
86) Perylene-d12	15.11	264	89829	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.22	112	231149	68.14	ng	0.01
7) Phenol-d6	6.28	99	319557	72.91	ng	0.00
23) Nitrobenzene-d5	7.20	82	304190	81.01	ng	0.00
41) 2,4,6-Tribromophenol	10.44	330	50785	60.66	ng	0.00
44) 2-Fluorobiphenyl	8.99	172	543472	95.21	ng	0.00
78) Terphenyl-d14	12.71	244	327689	82.30	ng	-0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.17	88	45720	27.19	ng	# 17
3) Pyridine	2.79	79	137078	28.09	ng	96
4) n-Nitrosodimethylamine	2.75	42	60054	32.21	ng	99
6) Aniline	6.28	93	74408	11.65	ng	# 27
8) 2-Chlorophenol	6.41	128	134257	36.83	ng	97
9) Benzaldehyde	6.16	77	35098	11.82	ng	86
10) Phenol	6.30	94	184476	36.88	ng	87
11) bis(2-Chloroethyl)ether	6.36	93	130238	34.91	ng	89
12) 1,3-Dichlorobenzene	6.56	146	135845	33.86	ng	97
13) 1,4-Dichlorobenzene	6.64	146	142178	35.71	ng	97
14) 1,2-Dichlorobenzene	6.80	146	129747	34.81	ng	99
15) Benzyl Alcohol	6.78	79	114957	35.64	ng	90
16) 2,2'-oxybis(1-Chloropropan	6.91	45	143446	26.55	ng	# 45
17) 2-Methylphenol	6.90	107	103535	34.81	ng	# 91
18) Hexachloroethane	7.13	117	48088	31.22	ng	# 83
19) n-Nitroso-di-n-propylamine	7.05	70	92393	31.38	ng	94
20) 3+4-Methylphenols	7.06	107	141536	39.65	ng	# 78
22) Acetophenone	7.04	105	195076	40.85	ng	# 93
24) Nitrobenzene	7.21	77	132718	35.06	ng	# 81
25) Isophorone	7.46	82	262351	37.72	ng	97
26) 2-Nitrophenol	7.53	139	66548	42.87	ng	95
27) 2,4-Dimethylphenol	7.59	122	125325	38.76	ng	89
28) bis(2-Chloroethoxy)methane	7.68	93	168373	37.20	ng	# 97
29) 2,4-Dichlorophenol	7.78	162	112306	42.98	ng	98
30) 1,2,4-Trichlorobenzene	7.86	180	117033	40.81	ng	97
31) Naphthalene	7.93	128	361736	37.29	ng	99
32) Benzoic acid	7.72	122	66359	28.27	ng	93
33) 4-Chloroaniline	7.99	127	44147	10.23	ng	98
34) Hexachlorobutadiene	8.06	225	65032	42.35	ng	98
35) Caprolactam	8.35	113	31198	35.15	ng	94
36) 4-Chloro-3-methylphenol	8.49	107	123447	39.95	ng	94
37) 2-Methylnaphthalene	8.63	142	257200	41.18	ng	# 95
39) 1,2,4,5-Tetrachlorobenzene	8.79	216	99700	33.25	ng	# 1
40) Hexachlorocyclopentadiene	8.78	237	22900	15.56	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	8.91	196	73066	36.95	ng	99
43) 2,4,5-Trichlorophenol	8.96	196	71558	42.06	ng	97
45) 1,1'-Biphenyl	9.08	154	266966	35.76	ng	98
46) 2-Chloronaphthalene	9.11	162	210481	35.91	ng	95
47) 2-Nitroaniline	9.21	65	74709	35.92	ng	97
48) Acenaphthylene	9.52	152	332432	35.65	ng	99
49) Dimethylphthalate	9.39	163	245052	38.15	ng	100
50) 2,6-Dinitrotoluene	9.46	165	56975	40.02	ng	# 90
51) Acenaphthene	9.69	154	201822	35.31	ng	99
52) 3-Nitroaniline	9.62	138	32689	18.06	ng	84
53) 2,4-Dinitrophenol	9.74	184	20873	33.21	ng	# 84
54) Dibenzofuran	9.86	168	286715	38.32	ng	# 86
55) 4-Nitrophenol	9.80	139	73828	53.68	ng	92
56) 2,4-Dinitrotoluene	9.86	165	80235	43.11	ng	# 61
57) Fluorene	10.20	166	241189	41.83	ng	97
58) 2,3,4,6-Tetrachlorophenol	9.99	232	50962	38.72	ng	# 47
59) Diethylphthalate	10.09	149	253416	39.31	ng	98
60) 4-Chlorophenyl-phenylether	10.20	204	111613	41.66	ng	89
61) 4-Nitroaniline	10.23	138	47592	27.33	ng	79
62) Azobenzene	10.35	77	282987	38.49	ng	89
64) 4,6-Dinitro-2-methylphenol	10.26	198	16483	19.83	ng	86
65) n-Nitrosodiphenylamine	10.32	169	194082	36.90	ng	99
66) 4-Bromophenyl-phenylether	10.68	248	59621	38.07	ng	# 78
67) Hexachlorobenzene	10.75	284	64081	36.96	ng	# 83
68) Atrazine	10.86	200	68498	41.79	ng	94
69) Pentachlorophenol	10.95	266	48227	47.00	ng	98
70) Phenanthrene	11.15	178	295470	34.78	ng	99
71) Anthracene	11.21	178	347720	41.16	ng	98
72) Carbazole	11.37	167	290639	36.55	ng	98
73) Di-n-butylphthalate	11.71	149	385411	41.67	ng	97
74) Fluoranthene	12.33	202	269338	31.27	ng	97
76) Benzidine	12.47	184	129190	28.82	ng	97
77) Pyrene	12.56	202	272808	36.28	ng	99
79) Butylbenzylphthalate	13.19	149	127621	38.05	ng	# 75
80) Benzo(a)anthracene	13.75	228	220674	38.65	ng	98
81) 3,3'-Dichlorobenzidine	13.71	252	72113	33.59	ng	# 97
82) Chrysene	13.78	228	214240	37.92	ng	98
83) Bis(2-ethylhexyl)phthalate	13.76	149	180617	47.17	ng	# 99
84) Di-n-octyl phthalate	14.37	149	307645	47.29	ng	# 100
85) Indeno(1,2,3-cd)pyrene	16.35	276	177162	40.76	ng	# 100
87) Benzo(b)fluoranthene	14.74	252	226022m	40.11	ng	
88) Benzo(k)fluoranthene	14.77	252	174299m	35.53	ng	
89) Benzo(a)pyrene	15.05	252	186098	39.01	ng	98
90) Dibenzo(a,h)anthracene	16.35	278	151742	38.09	ng	97
91) Benzo(g,h,i)perylene	16.72	276	135502	32.87	ng	95

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

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