

Data Path : Z:\HPCHEM1\BNA F\DATA\BF122415\
 Data File : BF084087.D
 Acq On : 24 Dec 2015 19:03
 Operator : UM/SJ
 Sample : PB87523BSD
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB87523BSD

Manual Integrations
 APPROVED

Sohil
 12/28/2015 1:09:47 PM

Quant Time: Dec 25 04:15:53 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF122415.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 24 16:48:39 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.80	152	48580	20.00	ng	0.00
21) Naphthalene-d8	8.09	136	190347	20.00	ng	0.00
38) Acenaphthene-d10	9.84	164	89270	20.00	ng	0.00
63) Phenanthrene-d10	11.32	188	165494	20.00	ng	0.00
75) Chrysene-d12	13.96	240	122167	20.00	ng	0.00
86) Perylene-d12	15.38	264	93639	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.42	112	217626	75.95	ng	0.01
7) Phenol-d6	6.45	99	262218	73.96	ng	0.00
23) Nitrobenzene-d5	7.37	82	243106	74.76	ng	0.00
41) 2,4,6-Tribromophenol	10.62	330	70543	74.10	ng	0.00
44) 2-Fluorobiphenyl	9.16	172	485276	74.38	ng	0.00
78) Terphenyl-d14	12.90	244	448680	66.46	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.43	88	38533	32.81	ng	98
3) Pyridine	3.14	79	96206	33.54	ng	98
4) n-Nitrosodimethylamine	3.08	42	51137	40.46	ng	91
6) Aniline	6.46	93	105291	23.49	ng	# 22
8) 2-Chlorophenol	6.59	128	145393	40.86	ng	94
9) Benzaldehyde	6.35	77	34440	18.16	ng	88
10) Phenol	6.46	94	154646	38.14	ng	86
11) bis(2-Chloroethyl)ether	6.53	93	111369	38.08	ng	95
12) 1,3-Dichlorobenzene	6.74	146	145983	37.58	ng	98
13) 1,4-Dichlorobenzene	6.82	146	146226m	37.40	ng	
14) 1,2-Dichlorobenzene	6.97	146	137736	38.23	ng	99
15) Benzyl Alcohol	6.96	79	114777	42.75	ng	94
16) 2,2'-oxybis(1-Chloropropan	7.08	45	104203	36.48	ng	# 50
17) 2-Methylphenol	7.07	107	104853	38.38	ng	# 91
18) Hexachloroethane	7.31	117	54103	37.92	ng	97
19) n-Nitroso-di-n-propylamine	7.22	70	87259	39.95	ng	# 89
20) 3+4-Methylphenols	7.23	107	143931	41.30	ng	# 48
22) Acetophenone	7.21	105	184472	38.79	ng	93
24) Nitrobenzene	7.39	77	128439	37.20	ng	# 84
25) Isophorone	7.63	82	239652	39.84	ng	95
26) 2-Nitrophenol	7.70	139	71260	40.33	ng	# 82
27) 2,4-Dimethylphenol	7.76	122	127266	43.39	ng	96
28) bis(2-Chloroethoxy)methane	7.84	93	140964	40.20	ng	99
29) 2,4-Dichlorophenol	7.95	162	112518	41.31	ng	96
30) 1,2,4-Trichlorobenzene	8.03	180	117557	39.38	ng	98
31) Naphthalene	8.11	128	404931	39.98	ng	99
32) Benzoic acid	7.89	122	60858	43.28	ng	93
33) 4-Chloroaniline	8.16	127	56810	14.10	ng	97
34) Hexachlorobutadiene	8.22	225	63951	37.56	ng	98
35) Caprolactam	8.53	113	36373m	48.15	ng	
36) 4-Chloro-3-methylphenol	8.66	107	125573	41.00	ng	93
37) 2-Methylnaphthalene	8.80	142	258784	40.82	ng	97
39) 1,2,4,5-Tetrachlorobenzene	8.97	216	118286	42.03	ng	98
40) Hexachlorocyclopentadiene	8.96	237	69211	83.12	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.08	196	72590	39.56	nq	96
43) 2,4,5-Trichlorophenol	9.14	196	75840	41.29	nq	97
45) 1,1'-Biphenyl	9.26	154	348949	43.27	nq	98
46) 2-Chloronaphthalene	9.29	162	255317	41.64	nq	95
47) 2-Nitroaniline	9.39	65	76118	46.72	nq	# 67
48) Acenaphthylene	9.70	152	390549	38.79	nq	99
49) Dimethylphthalate	9.57	163	311403	42.12	nq	# 90
50) 2,6-Dinitrotoluene	9.63	165	70931	45.20	nq	# 82
51) Acenaphthene	9.87	154	227128	38.79	nq	99
52) 3-Nitroaniline	9.80	138	35720	21.21	nq	# 71
53) 2,4-Dinitrophenol	9.92	184	48504	79.70	nq	96
54) Dibenzofuran	10.04	168	332361	41.27	nq	# 90
55) 4-Nitrophenol	9.98	139	71215	82.84	nq	# 76
56) 2,4-Dinitrotoluene	10.04	165	91903	44.36	nq	89
57) Fluorene	10.38	166	268163	39.18	nq	98
58) 2,3,4,6-Tetrachlorophenol	10.17	232	59111	45.15	nq	97
59) Diethylphthalate	10.27	149	299205	39.84	nq	98
60) 4-Chlorophenyl-phenylether	10.37	204	116432	36.88	nq	# 81
61) 4-Nitroaniline	10.42	138	73582	47.27	nq	96
62) Azobenzene	10.53	77	272025	43.96	nq	88
64) 4,6-Dinitro-2-methylphenol	10.45	198	40232	43.80	nq	81
65) n-Nitrosodiphenylamine	10.50	169	258379	43.80	nq	99
66) 4-Bromophenyl-phenylether	10.86	248	80580	42.03	nq	# 79
67) Hexachlorobenzene	10.93	284	84646	40.33	nq	# 72
68) Atrazine	11.02	200	71162	40.05	nq	97
69) Pentachlorophenol	11.14	266	61328	87.98	nq	98
70) Phenanthrene	11.34	178	410263	41.82	nq	99
71) Anthracene	11.40	178	433065	42.21	nq	99
72) Carbazole	11.55	167	361031	41.37	nq	99
73) Di-n-butylphthalate	11.88	149	458987	40.82	nq	# 98
74) Fluoranthene	12.53	202	444127	43.25	nq	96
76) Benzidine	12.65	184	109789	26.44	nq	99
77) Pyrene	12.76	202	450502	42.74	nq	99
79) Butylbenzylphthalate	13.38	149	193543	44.36	nq	# 81
80) Benzo(a)anthracene	13.95	228	321584	42.79	nq	100
81) 3,3'-Dichlorobenzidine	13.91	252	68298	30.01	nq	# 96
82) Chrysene	13.99	228	293385m	42.34	nq	
83) Bis(2-ethylhexyl)phthalate	13.94	149	247217	43.85	nq	# 97
84) Di-n-octyl phthalate	14.55	149	356798	44.98	nq	100
85) Indeno(1,2,3-cd)pyrene	16.77	276	283302	40.83	nq	99
87) Benzo(b)fluoranthene	14.98	252	297146m	42.73	nq	
88) Benzo(k)fluoranthene	15.00	252	220983m	46.89	nq	
89) Benzo(a)pyrene	15.32	252	248702	43.41	nq	99
90) Dibenzo(a,h)anthracene	16.77	278	236561	42.31	nq	# 95
91) Benzo(g,h,i)perylene	17.19	276	239184	41.92	nq	98

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

