

Data Path : Z:\HPCHEM1\BNA F\DATA\BF042015\
 Data File : BF078650.D
 Acq On : 20 Apr 2015 18:09
 Operator : TP/IZ
 Sample : PB82843BS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB82843BS

Manual Integrations
 APPROVED

apatel
 4/21/2015 5:03:33 PM

Quant Time: Apr 21 01:19:17 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 21 00:52:21 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	5.86	152	34816	20.00	ng	0.00
21) Naphthalene-d8	8.02	136	145055	20.00	ng	0.00
38) Acenaphthene-d10	11.21	164	73132	20.00	ng	0.00
63) Phenanthrene-d10	13.94	188	141188	20.00	ng	0.00
75) Chrysene-d12	17.81	240	145837	20.00	ng	0.00
86) Perylene-d12	19.53	264	137628	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	3.87	112	253545	120.07	ng	0.00
7) Phenol-d6	5.36	99	323667	122.31	ng	0.00
23) Nitrobenzene-d5	6.80	82	189785	79.30	ng	0.00
41) 2,4,6-Tribromophenol	12.69	330	82147	135.44	ng	0.00
44) 2-Fluorobiphenyl	10.03	172	414037	80.93	ng	0.00
78) Terphenyl-d14	16.47	244	487453	75.53	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	1.45	88	29228m	30.61	ng	
3) Pyridine	1.94	79	72950	29.17	ng	98
4) n-Nitrosodimethylamine	1.90	42	37811	39.27	ng	93
6) Aniline	5.35	93	87142	23.22	ng	91
8) 2-Chlorophenol	5.52	128	97020	38.86	ng	100
9) Benzaldehyde	5.16	77	14216	8.11	ng	95
10) Phenol	5.38	94	123453	38.93	ng	98
11) bis(2-Chloroethyl)ether	5.48	93	85553	37.52	ng	96
12) 1,3-Dichlorobenzene	5.76	146	100956	36.81	ng	# 93
13) 1,4-Dichlorobenzene	5.88	146	101617	36.81	ng	98
14) 1,2-Dichlorobenzene	6.12	146	97146	37.53	ng	94
15) Benzyl Alcohol	6.14	79	76806	41.30	ng	89
16) 2,2'-oxybis(1-Chloropropan	6.38	45	116251	38.22	ng	89
17) 2-Methylphenol	6.36	107	76738	39.58	ng	97
18) Hexachloroethane	6.67	117	35656	38.30	ng	93
19) n-Nitroso-di-n-propylamine	6.59	70	67756	38.81	ng	95
20) 3+4-Methylphenols	6.64	107	102716	39.72	ng	91
22) Acetophenone	6.56	105	135759	40.17	ng	# 93
24) Nitrobenzene	6.83	77	96429	38.54	ng	89
25) Isophorone	7.26	82	177325	39.04	ng	94
26) 2-Nitrophenol	7.38	139	48752	40.89	ng	95
27) 2,4-Dimethylphenol	7.54	122	86007	38.80	ng	98
28) bis(2-Chloroethoxy)methane	7.70	93	111086	38.14	ng	99
29) 2,4-Dichlorophenol	7.83	162	83469	41.39	ng	96
30) 1,2,4-Trichlorobenzene	7.94	180	84265	36.44	ng	97
31) Naphthalene	8.06	128	281047	37.23	ng	99
32) Benzoic acid	7.78	122	49706m	47.72	ng	
33) 4-Chloroaniline	8.22	127	63853	20.38	ng	97
34) Hexachlorobutadiene	8.31	225	47618	37.50	ng	98
35) Caprolactam	8.84	113	24739	41.09	ng	96
36) 4-Chloro-3-methylphenol	9.18	107	85506	42.02	ng	89
37) 2-Methylnaphthalene	9.32	142	194838	38.01	ng	100
39) 1,2,4,5-Tetrachlorobenzene	9.63	216	84323	37.21	ng	99
40) Hexachlorocyclopentadiene	9.61	237	68911	71.44	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.88	196	57593	40.30	nq	98
43) 2,4,5-Trichlorophenol	9.95	196	58883	39.44	nq	# 92
45) 1,1'-Biphenyl	10.19	154	226499	37.86	nq	96
46) 2-Chloronaphthalene	10.19	162	176545	37.51	nq	96
47) 2-Nitroaniline	10.44	65	51020	43.81	nq	# 67
48) Acenaphthylene	10.95	152	294569	38.76	nq	99
49) Dimethylphthalate	10.84	163	217855	39.65	nq	# 98
50) 2,6-Dinitrotoluene	10.94	165	46344	41.00	nq	# 56
51) Acenaphthene	11.27	154	168285	39.33	nq	100
52) 3-Nitroaniline	11.21	138	33273	25.89	nq	# 97
53) 2,4-Dinitrophenol	11.44	184	35143	79.96	nq	92
54) Dibenzofuran	11.60	168	267346	40.91	nq	95
55) 4-Nitrophenol	11.65	139	57807	61.73	nq	90
56) 2,4-Dinitrotoluene	11.67	165	64307	43.92	nq	# 81
57) Fluorene	12.23	166	217508	41.06	nq	100
58) 2,3,4,6-Tetrachlorophenol	11.86	232	41847	37.81	nq	98
59) Diethylphthalate	12.17	149	217300	41.02	nq	99
60) 4-Chlorophenyl-phenylether	12.29	204	99732	40.92	nq	# 83
61) 4-Nitroaniline	12.34	138	47123	36.27	nq	99
62) Azobenzene	12.57	77	206228	42.65	nq	97
64) 4,6-Dinitro-2-methylphenol	12.41	198	28041	41.43	nq	87
65) n-Nitrosodiphenylamine	12.53	169	186571	38.20	nq	97
66) 4-Bromophenyl-phenylether	13.18	248	61240	40.96	nq	99
67) Hexachlorobenzene	13.22	284	63118	38.52	nq	# 88
68) Atrazine	13.60	200	70287	49.69	nq	98
69) Pentachlorophenol	13.63	266	46301	65.30	nq	99
70) Phenanthrene	13.99	178	319329	39.10	nq	99
71) Anthracene	14.08	178	320797	39.86	nq	99
72) Carbazole	14.41	167	310379	41.15	nq	99
73) Di-n-butylphthalate	15.09	149	364614	42.68	nq	100
74) Fluoranthene	15.86	202	347925	40.40	nq	97
76) Benzidine	16.13	184	110246	19.63	nq	97
77) Pyrene	16.17	202	355639	38.85	nq	98
79) Butylbenzylphthalate	17.17	149	164446	47.13	nq	# 82
80) Benzo(a)anthracene	17.79	228	338731	39.04	nq	99
81) 3,3'-Dichlorobenzidine	17.79	252	75629	26.02	nq	# 97
82) Chrysene	17.84	228	309233	37.93	nq	99
83) Bis(2-ethylhexyl)phthalate	17.93	149	234301	42.81	nq	99
84) Di-n-octyl phthalate	18.72	149	405180	45.09	nq	# 100
85) Indeno(1,2,3-cd)pyrene	20.73	276	350209	37.86	nq	# 87
87) Benzo(b)fluoranthene	19.10	252	369163	43.40	nq	# 97
88) Benzo(k)fluoranthene	19.13	252	259033m	34.27	nq	
89) Benzo(a)pyrene	19.47	252	310057	41.77	nq	97
90) Dibenzo(a,h)anthracene	20.75	278	286772	38.59	nq	99
91) Benzo(g,h,i)perylene	21.05	276	285190	38.27	nq	99

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

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