

Data Path : Z:\HPCHEM1\BNA F\DATA\BF020316\
 Data File : BF084789.D
 Acq On : 3 Feb 2016 17:21
 Operator : SJ/IZ
 Sample : PB88159BS
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB88159BS

Manual Integrations
 APPROVED

mohammad
 2/4/2016 2:04:29 PM

Quant Time: Feb 04 00:41:05 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF020116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 03 14:16:01 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.68	152	184199	20.00	ng	0.00
21) Naphthalene-d8	7.97	136	772740	20.00	ng	0.00
38) Acenaphthene-d10	9.73	164	348545	20.00	ng	0.00
63) Phenanthrene-d10	11.21	188	662869	20.00	ng	0.00
75) Chrysene-d12	13.84	240	485596	20.00	ng	0.00
86) Perylene-d12	15.21	264	374399	20.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.29	112	1164896	98.50	ng	0.01
7) Phenol-d6	6.34	99	1517796	103.53	ng	0.00
23) Nitrobenzene-d5	7.26	82	976602	71.19	ng	0.00
41) 2,4,6-Tribromophenol	10.52	330	409818	112.72	ng	0.00
44) 2-Fluorobiphenyl	9.06	172	1695311	80.44	ng	0.00
78) Terphenyl-d14	12.80	244	1630950	76.11	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	2.28	88	161668	31.21	ng	# 74
3) Pyridine	2.93	79	475388	35.22	ng	# 74
4) n-Nitrosodimethylamine	2.89	42	240767	34.49	ng	80
6) Aniline	6.35	93	318710	17.77	ng	# 40
8) 2-Chlorophenol	6.48	128	485826	36.29	ng	99
9) Benzaldehyde	6.22	77	170693	19.72	ng	92
10) Phenol	6.35	94	486826	29.64	ng	# 63
11) bis(2-Chloroethyl)ether	6.43	93	444530m	34.71	ng	
12) 1,3-Dichlorobenzene	6.62	146	484137	34.23	ng	# 95
13) 1,4-Dichlorobenzene	6.70	146	500432	35.40	ng	96
14) 1,2-Dichlorobenzene	6.85	146	426575	32.39	ng	95
15) Benzyl Alcohol	6.84	79	379625	37.50	ng	91
16) 2,2'-oxybis(1-Chloropropan	6.98	45	696993	32.35	ng	87
17) 2-Methylphenol	6.97	107	395140	37.33	ng	# 91
18) Hexachloroethane	7.20	117	184467	33.88	ng	95
19) n-Nitroso-di-n-propylamine	7.12	70	327793	35.67	ng	# 77
20) 3+4-Methylphenols	7.12	107	475498	36.19	ng	# 79
22) Acetophenone	7.10	105	650378	31.93	ng	99
24) Nitrobenzene	7.28	77	475862	32.39	ng	97
25) Isophorone	7.52	82	884138	33.15	ng	98
26) 2-Nitrophenol	7.60	139	266650	36.81	ng	# 90
27) 2,4-Dimethylphenol	7.64	122	421715	33.46	ng	97
28) bis(2-Chloroethoxy)methane	7.73	93	527690	33.35	ng	99
29) 2,4-Dichlorophenol	7.85	162	373610	34.65	ng	92
30) 1,2,4-Trichlorobenzene	7.92	180	398640	32.73	ng	96
31) Naphthalene	8.00	128	1285086	33.15	ng	99
32) Benzoic acid	7.78	122	260559	32.33	ng	96
33) 4-Chloroaniline	8.05	127	141218	8.81	ng	95
34) Hexachlorobutadiene	8.12	225	235966	33.60	ng	99
35) Caprolactam	8.43	113	118996m	35.81	ng	
36) 4-Chloro-3-methylphenol	8.56	107	454643	36.96	ng	98
37) 2-Methylnaphthalene	8.69	142	931035	38.38	ng	98
39) 1,2,4,5-Tetrachlorobenzene	8.85	216	348517	31.48	ng	98
40) Hexachlorocyclopentadiene	8.84	237	349225	73.81	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	8.98	196	260004	36.46	nq	92
43) 2,4,5-Trichlorophenol	9.01	196	255773	35.30	nq	# 82
45) 1,1'-Biphenyl	9.15	154	1009132	32.41	nq	98
46) 2-Chloronaphthalene	9.17	162	754927	32.93	nq	95
47) 2-Nitroaniline	9.28	65	235488	32.44	nq	98
48) Acenaphthylene	9.58	152	1157421	33.27	nq	97
49) Dimethylphthalate	9.47	163	963143	36.28	nq	# 91
50) 2,6-Dinitrotoluene	9.53	165	224292	38.68	nq	# 91
51) Acenaphthene	9.76	154	748381	33.06	nq	97
52) 3-Nitroaniline	9.69	138	104757	16.08	nq	95
53) 2,4-Dinitrophenol	9.80	184	203757	76.41	nq	# 89
54) Dibenzofuran	9.94	168	1098352	39.70	nq	98
55) 4-Nitrophenol	9.87	139	275630	57.55	nq	# 81
56) 2,4-Dinitrotoluene	9.93	165	271690	36.73	nq	# 88
57) Fluorene	10.27	166	802814	34.75	nq	98
58) 2,3,4,6-Tetrachlorophenol	10.05	232	206185	37.38	nq	93
59) Diethylphthalate	10.17	149	940243	36.27	nq	96
60) 4-Chlorophenyl-phenylether	10.27	204	385729	38.73	nq	96
61) 4-Nitroaniline	10.30	138	200347	31.55	nq	92
62) Azobenzene	10.43	77	997842	40.03	nq	93
64) 4,6-Dinitro-2-methylphenol	10.34	198	152464	37.27	nq	73
65) n-Nitrosodiphenylamine	10.40	169	807890	38.38	nq	99
66) 4-Bromophenyl-phenylether	10.76	248	259635	35.46	nq	97
67) Hexachlorobenzene	10.82	284	281874	35.33	nq	# 94
68) Atrazine	10.92	200	226333	34.05	nq	99
69) Pentachlorophenol	11.02	266	259468	69.21	nq	99
70) Phenanthrene	11.23	178	1195787	32.52	nq	97
71) Anthracene	11.29	178	1290625	34.84	nq	99
72) Carbazole	11.45	167	1145914	35.47	nq	97
73) Di-n-butylphthalate	11.78	149	1283611	31.21	nq	# 96
74) Fluoranthene	12.42	202	1282453	34.46	nq	92
76) Benzidine	12.55	184	344132	21.46	nq	97
77) Pyrene	12.64	202	1216012	32.71	nq	99
79) Butylbenzylphthalate	13.28	149	591900	35.09	nq	# 84
80) Benzo(a)anthracene	13.83	228	1055002	35.00	nq	98
81) 3,3'-Dichlorobenzidine	13.79	252	158220	14.83	nq	# 94
82) Chrysene	13.86	228	918062	31.41	nq	99
83) Bis(2-ethylhexyl)phthalate	13.84	149	762190	36.31	nq	# 96
84) Di-n-octyl phthalate	14.44	149	1165772	33.67	nq	# 89
85) Indeno(1,2,3-cd)pyrene	16.50	276	778497	33.84	nq	99
87) Benzo(b)fluoranthene	14.83	252	830030m	33.00	nq	
88) Benzo(k)fluoranthene	14.85	252	847712m	40.76	nq	
89) Benzo(a)pyrene	15.15	252	771115	35.98	nq	# 96
90) Dibenzo(a,h)anthracene	16.51	278	642526	35.61	nq	97
91) Benzo(g,h,i)perylene	16.89	276	662663	36.46	nq	98

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

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