

Data Path : Z:\HPCHEM1\BNA F\DATA\BF012617\
 Data File : BF092535.D
 Acq On : 26 Jan 2017 20:04
 Operator : SJ/MA
 Sample : PB96342BS
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB96342BS

Manual Integrations
 APPROVED

mohammad
 1/27/2017 3:03:04 PM

Quant Time: Jan 27 01:21:30 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF012417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 24 13:48:07 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	158959	20.00	ng	0.00
21) Naphthalene-d8	8.27	136	655850	20.00	ng	0.00
38) Acenaphthene-d10	10.03	164	358994	20.00	ng	0.00
63) Phenanthrene-d10	11.52	188	601635	20.00	ng	-0.01
75) Chrysene-d12	14.17	240	493763	20.00	ng	0.00
86) Perylene-d12	15.64	264	401177	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.58	112	1396785	136.71	ng	0.03
7) Phenol-d6	6.60	99	1616485	124.21	ng	0.01
23) Nitrobenzene-d5	7.54	82	1059450	88.24	ng	0.00
41) 2,4,6-Tribromophenol	10.83	330	384329	157.10	ng	0.01
44) 2-Fluorobiphenyl	9.36	172	1932654	99.55	ng	0.00
78) Terphenyl-d14	13.11	244	1680725	90.64	ng	-0.01

Target Compounds						Qvalue
2) 1,4-Dioxane	2.72	88	199980	36.80	ng	# 83
3) Pyridine	3.47	79	588705	38.06	ng	89
4) n-Nitrosodimethylamine	3.42	42	304515	40.13	ng	86
6) Aniline	6.64	93	393698	20.28	ng	# 43
8) 2-Chlorophenol	6.76	128	489628	45.62	ng	89
9) Benzaldehyde	6.51	77	236293	25.59	ng	87
10) Phenol	6.61	94	691830	44.04	ng	# 73
11) bis(2-Chloroethyl)ether	6.70	93	466570	39.70	ng	92
12) 1,3-Dichlorobenzene	6.91	146	525960m	41.44	ng	
13) 1,4-Dichlorobenzene	6.99	146	554078	43.38	ng	96
14) 1,2-Dichlorobenzene	7.15	146	503265m	41.53	ng	
15) Benzyl Alcohol	7.12	79	490461	50.01	ng	96
16) 2,2'-oxybis(1-Chloropropan	7.25	45	915868	39.39	ng	92
17) 2-Methylphenol	7.23	107	456632	49.80	ng	94
18) Hexachloroethane	7.49	117	193982	44.08	ng	92
19) n-Nitroso-di-n-propylamine	7.39	70	355210	40.08	ng	94
20) 3+4-Methylphenols	7.38	107	481709	43.30	ng	93
22) Acetophenone	7.39	105	710379	45.48	ng	# 94
24) Nitrobenzene	7.56	77	596607	47.43	ng	# 90
25) Isophorone	7.80	82	1039511	43.94	ng	97
26) 2-Nitrophenol	7.88	139	289523	54.93	ng	# 84
27) 2,4-Dimethylphenol	7.92	122	514863	46.97	ng	96
28) bis(2-Chloroethoxy)methane	8.01	93	634898	40.88	ng	99
29) 2,4-Dichlorophenol	8.12	162	463773	48.80	ng	94
30) 1,2,4-Trichlorobenzene	8.20	180	455015	44.30	ng	95
31) Naphthalene	8.29	128	1425454	42.47	ng	99
32) Benzoic acid	8.04	122	368313	45.25	ng	97
33) 4-Chloroaniline	8.33	127	151695	10.68	ng	94
34) Hexachlorobutadiene	8.41	225	245611	43.72	ng	99
35) Caprolactam	8.72	113	153040m	48.22	ng	
36) 4-Chloro-3-methylphenol	8.82	107	467451	45.61	ng	99
37) 2-Methylnaphthalene	8.99	142	991329	46.65	ng	96
39) 1,2,4,5-Tetrachlorobenzene	9.15	216	414837	45.81	ng	99
40) Hexachlorocyclopentadiene	9.14	237	427046	94.17	ng	96

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.26	196	357016	51.46	nq	93
43) 2,4,5-Trichlorophenol	9.31	196	298389	47.00	nq #	83
45) 1,1'-Biphenyl	9.45	154	1165640	44.92	nq	96
46) 2-Chloronaphthalene	9.48	162	937817	44.25	nq	96
47) 2-Nitroaniline	9.57	65	348752	48.86	nq	89
48) Acenaphthylene	9.89	152	1417495	45.60	nq	99
49) Dimethylphthalate	9.76	163	1136848	49.80	nq	99
50) 2,6-Dinitrotoluene	9.81	165	235282	50.42	nq #	82
51) Acenaphthene	10.06	154	928475	48.06	nq	97
52) 3-Nitroaniline	9.98	138	123748	21.18	nq	80
53) 2,4-Dinitrophenol	10.09	184	224615	75.25	nq #	63
54) Dibenzofuran	10.24	168	1239155	47.24	nq	97
55) 4-Nitrophenol	10.14	139	398010	85.76	nq	95
56) 2,4-Dinitrotoluene	10.22	165	364876	58.91	nq #	72
57) Fluorene	10.58	166	911240	46.85	nq	99
58) 2,3,4,6-Tetrachlorophenol	10.36	232	257355m	54.11	nq	
59) Diethylphthalate	10.45	149	1007580	46.03	nq	99
60) 4-Chlorophenyl-phenylether	10.58	204	450489	47.76	nq #	72
61) 4-Nitroaniline	10.60	138	246416	42.16	nq	93
62) Azobenzene	10.74	77	1241926	49.86	nq	89
64) 4,6-Dinitro-2-methylphenol	10.64	198	179045	54.48	nq	80
65) n-Nitrosodiphenylamine	10.69	169	853818	45.44	nq	100
66) 4-Bromophenyl-phenylether	11.07	248	273266	44.99	nq #	75
67) Hexachlorobenzene	11.13	284	260612	42.44	nq #	81
68) Atrazine	11.22	200	261935	40.73	nq	99
69) Pentachlorophenol	11.32	266	316177	80.54	nq	97
70) Phenanthrene	11.55	178	1543070	46.57	nq	99
71) Anthracene	11.60	178	1385236	42.78	nq	99
72) Carbazole	11.76	167	1486499	48.08	nq	97
73) Di-n-butylphthalate	12.09	149	1716127	48.55	nq #	95
74) Fluoranthene	12.74	202	1575666	48.62	nq	95
76) Benzidine	12.85	184	422211	23.03	nq	97
77) Pyrene	12.97	202	1562229	43.67	nq	99
79) Butylbenzylphthalate	13.59	149	768330	53.01	nq #	82
80) Benzo(a)anthracene	14.16	228	1211842	45.67	nq	100
81) 3,3'-Dichlorobenzidine	14.11	252	310082	30.78	nq	96
82) Chrysene	14.19	228	1116835	39.41	nq	99
83) Bis(2-ethylhexyl)phthalate	14.15	149	934492	51.20	nq #	95
84) Di-n-octyl phthalate	14.75	149	1660047	50.10	nq	99
85) Indeno(1,2,3-cd)pyrene	17.14	276	976256	46.56	nq	96
87) Benzo(b)fluoranthene	15.21	252	1354609m	52.60	nq	
88) Benzo(k)fluoranthene	15.24	252	781315m	36.48	nq	
89) Benzo(a)pyrene	15.59	252	1014629	46.57	nq	99
90) Dibenzo(a,h)anthracene	17.16	278	817183	46.38	nq	97
91) Benzo(g,h,i)perylene	17.59	276	796930	44.73	nq #	91

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

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