

Data Path : Z:\HPCHEM1\BNA F\DATA\BF081916\
 Data File : BF089798.D
 Acq On : 19 Aug 2016 16:29
 Operator : UM/SJ
 Sample : PB92975BSD
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB92975BSD

Manual Integrations
 APPROVED

sohil
 8/22/2016 3:47:35 PM

Quant Time: Aug 22 01:35:28 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 19 14:02:57 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.68	152	568868	20.00	ng	0.00
21) Naphthalene-d8	7.98	136	2212914	20.00	ng	0.00
38) Acenaphthene-d10	9.72	164	1026038	20.00	ng	0.01
63) Phenanthrene-d10	11.20	188	1860031	20.00	ng	0.01
75) Chrysene-d12	13.85	240	1184954	20.00	ng	-0.02
86) Perylene-d12	15.33	264	985235	20.00	ng	-0.05

System Monitoring Compounds

5) 2-Fluorophenol	5.28	112	2741523	82.65	ng	0.02
7) Phenol-d6	6.32	99	3271697	80.34	ng	0.00
23) Nitrobenzene-d5	7.26	82	3012774	84.58	ng	0.00
41) 2,4,6-Tribromophenol	10.50	330	761464	78.09	ng	0.00
44) 2-Fluorobiphenyl	9.05	172	4492379	76.95	ng	0.00
78) Terphenyl-d14	12.79	244	4189078	79.98	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.27	88	546659	33.30	ng	97
3) Pyridine	2.91	79	1519322	35.29	ng	95
4) n-Nitrosodimethylamine	2.88	42	617771	38.54	ng	# 78
6) Aniline	6.34	93	1101767	20.17	ng	# 38
8) 2-Chlorophenol	6.47	128	1605613	40.48	ng	87
9) Benzaldehyde	6.22	77	391791	14.77	ng	98
10) Phenol	6.34	94	1935463	40.55	ng	95
11) bis(2-Chloroethyl)ether	6.42	93	1349203	36.46	ng	96
12) 1,3-Dichlorobenzene	6.63	146	1660553m	40.04	ng	
13) 1,4-Dichlorobenzene	6.71	146	1613938m	37.86	ng	
14) 1,2-Dichlorobenzene	6.86	146	1568391	39.41	ng	96
15) Benzyl Alcohol	6.83	79	1100946	40.77	ng	95
16) 2,2'-oxybis(1-Chloropropan	6.98	45	1830269	38.64	ng	91
17) 2-Methylphenol	6.95	107	1243480	40.02	ng	98
18) Hexachloroethane	7.20	117	590675	38.85	ng	90
19) n-Nitroso-di-n-propylamine	7.12	70	1109506	42.60	ng	# 84
20) 3+4-Methylphenols	7.11	107	1483178	38.93	ng	# 66
22) Acetophenone	7.10	105	1865152	36.71	ng	# 90
24) Nitrobenzene	7.27	77	1467831	36.73	ng	99
25) Isophorone	7.52	82	2894054	40.28	ng	# 92
26) 2-Nitrophenol	7.59	139	798985	40.07	ng	# 83
27) 2,4-Dimethylphenol	7.63	122	1352825	37.64	ng	95
28) bis(2-Chloroethoxy)methane	7.74	93	1988188	44.73	ng	96
29) 2,4-Dichlorophenol	7.83	162	1179694	39.12	ng	93
30) 1,2,4-Trichlorobenzene	7.92	180	1353367	40.58	ng	95
31) Naphthalene	8.00	128	4199331	40.60	ng	100
32) Benzoic acid	7.76	122	824647	29.45	ng	98
33) 4-Chloroaniline	8.04	127	607964	13.56	ng	# 95
34) Hexachlorobutadiene	8.12	225	674093	38.59	ng	96
35) Caprolactam	8.42	113	413146	44.99	ng	# 85
36) 4-Chloro-3-methylphenol	8.54	107	1362202	41.71	ng	99
37) 2-Methylnaphthalene	8.68	142	2851452	42.62	ng	98
39) 1,2,4,5-Tetrachlorobenzene	8.86	216	1006278	30.29	ng	97
40) Hexachlorocyclopentadiene	8.84	237	1131086	98.05	ng	97

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	8.96	196	891928	42.67	nq	99
43) 2,4,5-Trichlorophenol	9.00	196	880088	42.33	nq	97
45) 1,1'-Biphenyl	9.15	154	3049834	38.41	nq	95
46) 2-Chloronaphthalene	9.16	162	2486822	39.62	nq	99
47) 2-Nitroaniline	9.27	65	893446	49.91	nq	# 80
48) Acenaphthylene	9.58	152	4072325	43.18	nq	99
49) Dimethylphthalate	9.46	163	3178302	43.99	nq	# 99
50) 2,6-Dinitrotoluene	9.51	165	637511	38.22	nq	# 85
51) Acenaphthene	9.76	154	3024462	48.45	nq	99
52) 3-Nitroaniline	9.67	138	365170	19.76	nq	97
53) 2,4-Dinitrophenol	9.78	184	761583	78.24	nq	# 82
54) Dibenzofuran	9.93	168	3756940	47.92	nq	91
55) 4-Nitrophenol	9.84	139	1136984	75.38	nq	91
56) 2,4-Dinitrotoluene	9.91	165	839267	38.78	nq	# 61
57) Fluorene	10.26	166	2664165	42.64	nq	97
58) 2,3,4,6-Tetrachlorophenol	10.04	232	755766	49.95	nq	98
59) Diethylphthalate	10.16	149	3071188	41.85	nq	97
60) 4-Chlorophenyl-phenylether	10.26	204	1104721	38.76	nq	92
61) 4-Nitroaniline	10.28	138	659914	38.09	nq	97
62) Azobenzene	10.42	77	3094728	44.85	nq	95
64) 4,6-Dinitro-2-methylphenol	10.32	198	458348	42.88	nq	# 58
65) n-Nitrosodiphenylamine	10.39	169	2671875	45.69	nq	98
66) 4-Bromophenyl-phenylether	10.75	248	831334	42.63	nq	# 86
67) Hexachlorobenzene	10.81	284	868001	39.02	nq	# 85
68) Atrazine	10.91	200	687715	38.40	nq	96
69) Pentachlorophenol	11.00	266	1015874	80.56	nq	97
70) Phenanthrene	11.22	178	3993614	42.51	nq	95
71) Anthracene	11.27	178	3754470	39.32	nq	98
72) Carbazole	11.43	167	3754180	43.83	nq	97
73) Di-n-butylphthalate	11.78	149	4555471	42.51	nq	# 98
74) Fluoranthene	12.40	202	3573278	39.99	nq	93
76) Benzidine	12.52	184	1186372	30.94	nq	98
77) Pyrene	12.63	202	3920376	40.63	nq	95
79) Butylbenzylphthalate	13.28	149	2053139	47.11	nq	91
80) Benzo(a)anthracene	13.84	228	3067284	45.20	nq	98
81) 3,3'-Dichlorobenzidine	13.82	252	611011	27.46	nq	# 94
82) Chrysene	13.89	228	2701462	46.43	nq	98
83) Bis(2-ethylhexyl)phthalate	13.87	149	2635932	43.91	nq	98
84) Di-n-octyl phthalate	14.55	149	4056464	50.85	nq	# 95
85) Indeno(1,2,3-cd)pyrene	16.62	276	2690780	36.25	nq	99
87) Benzo(b)fluoranthene	14.95	252	2527982	46.64	nq	# 95
88) Benzo(k)fluoranthene	14.97	252	2110938m	42.93	nq	
89) Benzo(a)pyrene	15.28	252	2254803	43.95	nq	# 94
90) Dibenzo(a,h)anthracene	16.64	278	2256158	40.52	nq	97
91) Benzo(g,h,i)perylene	17.01	276	2317636	39.53	nq	98

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

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