

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF091516\
 Data File : BF090086.D
 Acq On : 15 Sep 2016 12:25
 Operator : UM/SJ
 Sample : PB93381BS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB93381BS

Manual Integrations
 APPROVED

sohil
 9/16/2016 11:34:53 AM

Quant Time: Sep 15 13:01:47 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF091516.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 15 09:41:41 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.55	152	142239	20.00	ng	0.00
21) Naphthalene-d8	7.84	136	583740	20.00	ng	0.00
38) Acenaphthene-d10	9.59	164	306273	20.00	ng	0.00
63) Phenanthrene-d10	11.05	188	485002	20.00	ng	0.00
75) Chrysene-d12	13.67	240	311819	20.00	ng	0.00
86) Perylene-d12	14.99	264	260029	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.13	112	1118619	126.48	ng	0.01
7) Phenol-d6	6.20	99	1324407	114.38	ng	0.00
23) Nitrobenzene-d5	7.12	82	805070	77.54	ng	-0.01
41) 2,4,6-Tribromophenol	10.36	330	298513	97.03	ng	-0.01
44) 2-Fluorobiphenyl	8.92	172	1433840	73.52	ng	0.00
78) Terphenyl-d14	12.63	244	1099677	83.75	ng	-0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.06	88	162575	35.21	ng	99
3) Pyridine	2.65	79	454458	38.93	ng	98
4) n-Nitrosodimethylamine	2.62	42	138838	37.78	ng	98
6) Aniline	6.22	93	301183	20.96	ng	# 26
8) 2-Chlorophenol	6.33	128	400428	39.30	ng	95
9) Benzaldehyde	6.09	77	41932	6.79	ng	97
10) Phenol	6.22	94	550916	42.39	ng	# 74
11) bis(2-Chloroethyl)ether	6.30	93	416586	39.97	ng	97
12) 1,3-Dichlorobenzene	6.49	146	439616	41.94	ng	97
13) 1,4-Dichlorobenzene	6.57	146	443978	41.03	ng	97
14) 1,2-Dichlorobenzene	6.72	146	403302m	41.02	ng	
15) Benzyl Alcohol	6.71	79	305343	46.83	ng	96
16) 2,2'-oxybis(1-Chloropropan	6.84	45	476128	38.55	ng	97
17) 2-Methylphenol	6.83	107	360988	43.31	ng	97
18) Hexachloroethane	7.06	117	159266	40.56	ng	94
19) n-Nitroso-di-n-propylamine	6.98	70	276407	38.42	ng	98
20) 3+4-Methylphenols	6.98	107	438067	44.91	ng	# 86
22) Acetophenone	6.97	105	517786	36.64	ng	# 98
24) Nitrobenzene	7.14	77	449217	41.08	ng	90
25) Isophorone	7.38	82	824877	37.94	ng	98
26) 2-Nitrophenol	7.46	139	231283	43.62	ng	# 89
27) 2,4-Dimethylphenol	7.52	122	379250	40.65	ng	94
28) bis(2-Chloroethoxy)methane	7.61	93	548172	42.99	ng	99
29) 2,4-Dichlorophenol	7.70	162	336172	40.79	ng	92
30) 1,2,4-Trichlorobenzene	7.78	180	322421	39.66	ng	98
31) Naphthalene	7.86	128	1265066	44.09	ng	99
32) Benzoic acid	7.63	122	119974	16.78	ng	97
33) 4-Chloroaniline	7.92	127	134444	11.01	ng	96
34) Hexachlorobutadiene	7.99	225	172053	41.33	ng	99
35) Caprolactam	8.28	113	127632m	46.37	ng	
36) 4-Chloro-3-methylphenol	8.41	107	370272	38.70	ng	98
37) 2-Methylnaphthalene	8.55	142	707105	39.73	ng	100
39) 1,2,4,5-Tetrachlorobenzene	8.72	216	267267	33.94	ng	99
40) Hexachlorocyclopentadiene	8.71	237	290589	71.00	ng	97

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42) 2,4,6-Trichlorophenol	8.83	196	220179	38.57	ng	99
43) 2,4,5-Trichlorophenol	8.88	196	230805	37.59	ng #	95
45) 1,1'-Biphenyl	9.02	154	836697	33.39	ng	98
46) 2-Chloronaphthalene	9.04	162	720178	38.78	ng	95
47) 2-Nitroaniline	9.13	65	223641	37.72	ng #	75
48) Acenaphthylene	9.44	152	1091840	35.01	ng	99
49) Dimethylphthalate	9.32	163	798737	35.11	ng	98
50) 2,6-Dinitrotoluene	9.38	165	182530	35.91	ng	92
51) Acenaphthene	9.61	154	637796	36.62	ng	100
52) 3-Nitroaniline	9.54	138	107646	17.40	ng	87
53) 2,4-Dinitrophenol	9.66	184	141020	51.02	ng	91
54) Dibenzofuran	9.79	168	985136	40.89	ng	97
55) 4-Nitrophenol	9.72	139	291353	66.94	ng	87
56) 2,4-Dinitrotoluene	9.78	165	233562	37.69	ng	92
57) Fluorene	10.12	166	661978	35.17	ng	99
58) 2,3,4,6-Tetrachlorophenol	9.91	232	186273	39.78	ng	100
59) Diethylphthalate	10.02	149	774316	35.16	ng	100
60) 4-Chlorophenyl-phenylether	10.12	204	281942	37.94	ng	96
61) 4-Nitroaniline	10.16	138	215909	34.49	ng	88
62) Azobenzene	10.28	77	941720	40.75	ng	97
64) 4,6-Dinitro-2-methylphenol	10.18	198	102230	31.62	ng #	43
65) n-Nitrosodiphenylamine	10.25	169	703967	45.81	ng	100
66) 4-Bromophenyl-phenylether	10.62	248	209530	44.82	ng #	93
67) Hexachlorobenzene	10.67	284	237009	43.02	ng #	88
68) Atrazine	10.78	200	196140	42.95	ng	95
69) Pentachlorophenol	10.87	266	242902	75.26	ng	99
70) Phenanthrene	11.07	178	973005	41.51	ng	99
71) Anthracene	11.13	178	1162544	45.43	ng	100
72) Carbazole	11.29	167	1155285	44.01	ng	99
73) Di-n-butylphthalate	11.63	149	1224638	41.60	ng	100
74) Fluoranthene	12.25	202	1067498	43.47	ng	99
76) Benzidine	12.39	184	414859	32.62	ng	96
77) Pyrene	12.47	202	1007236	44.26	ng	99
79) Butylbenzylphthalate	13.11	149	488427	43.10	ng #	72
80) Benzo(a)anthracene	13.66	228	788577	43.32	ng	99
81) 3,3'-Dichlorobenzidine	13.62	252	115214	15.10	ng #	95
82) Chrysene	13.69	228	797673	45.01	ng	99
83) Bis(2-ethylhexyl)phthalate	13.68	149	593705	42.36	ng	99
84) Di-n-octyl phthalate	14.29	149	940908	37.55	ng	98
85) Indeno(1,2,3-cd)pyrene	16.17	276	739792	49.87	ng	99
87) Benzo(b)fluoranthene	14.64	252	594219m	38.98	ng	
88) Benzo(k)fluoranthene	14.66	252	621790m	45.31	ng	
89) Benzo(a)pyrene	14.94	252	600569	43.93	ng #	94
90) Dibenzo(a,h)anthracene	16.19	278	622686	58.16	ng	99
91) Benzo(g,h,i)perylene	16.53	276	639681	62.61	ng	100

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

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