

Data Path : Z:\HPCHEM1\BNA F\DATA\BF092817\
 Data File : BF098914.D
 Acq On : 28 Sep 2017 18:42
 Operator : SJ/JU
 Sample : PB102707BS
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB102707BS

Manual Integrations
 APPROVED

Sohil
 9/29/2017 6:04:23 PM

Quant Time: Sep 29 01:43:13 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 23 13:50:58 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.92	152	109915	20.00	ng	0.00
21) Naphthalene-d8	8.20	136	469498	20.00	ng	0.00
38) Acenaphthene-d10	9.96	164	217700	20.00	ng	0.00
63) Phenanthrene-d10	11.44	188	391367	20.00	ng	0.00
75) Chrysene-d12	14.09	240	282790	20.00	ng	0.00
86) Perylene-d12	15.57	264	237252	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.55	112	798158	115.60	ng	0.00
7) Phenol-d6	6.55	99	964649	113.25	ng	0.00
23) Nitrobenzene-d5	7.48	82	585659	80.00	ng	0.00
41) 2,4,6-Tribromophenol	10.74	330	203641	114.32	ng	0.00
44) 2-Fluorobiphenyl	9.27	172	1074129	82.37	ng	0.00
78) Terphenyl-d14	13.03	244	996075	80.10	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.76	88	101024	30.629	ng	99
3) Pyridine	3.55	79	289643	32.462	ng	99
4) n-Nitrosodimethylamine	3.49	42	142670	37.708	ng	96
6) Aniline	6.58	93	338614	29.455	ng	99
8) 2-Chlorophenol	6.70	128	292744	37.439	ng	97
9) Benzaldehyde	6.47	77	66536	13.466	ng	96
10) Phenol	6.56	94	374394	39.302	ng	99
11) bis(2-Chloroethyl)ether	6.65	93	280995	37.944	ng	99
12) 1,3-Dichlorobenzene	6.86	146	299262	36.545	ng	99
13) 1,4-Dichlorobenzene	6.93	146	307480	36.905	ng	99
14) 1,2-Dichlorobenzene	7.09	146	286827	37.258	ng	100
15) Benzyl Alcohol	7.06	79	253891	40.632	ng	97
16) 2,2'-oxybis(1-Chloropropan	7.19	45	442484	38.350	ng	99
17) 2-Methylphenol	7.16	107	246742	38.566	ng	99
18) Hexachloroethane	7.43	117	112337	37.511	ng	93
19) n-Nitroso-di-n-propylamine	7.32	70	198771	36.551	ng	99
20) 3+4-Methylphenols	7.32	107	308281	38.089	ng	# 85
22) Acetophenone	7.32	105	395222	34.744	ng	# 97
24) Nitrobenzene	7.50	77	309656	37.287	ng	99
25) Isophorone	7.73	82	553724	36.583	ng	98
26) 2-Nitrophenol	7.82	139	160251	40.127	ng	90
27) 2,4-Dimethylphenol	7.85	122	271000	35.933	ng	100
28) bis(2-Chloroethoxy)methane	7.94	93	351672	37.282	ng	99
29) 2,4-Dichlorophenol	8.05	162	241591	37.310	ng	98
30) 1,2,4-Trichlorobenzene	8.14	180	239246	36.942	ng	98
31) Naphthalene	8.22	128	825476	37.436	ng	100
32) Benzoic acid	7.96	122	180889	29.199	ng	99
33) 4-Chloroaniline	8.26	127	245571	26.668	ng	99
34) Hexachlorobutadiene	8.33	225	122162	36.622	ng	98
35) Caprolactam	8.63	113	83377m	40.141	ng	
36) 4-Chloro-3-methylphenol	8.75	107	270811	37.209	ng	99
37) 2-Methylnaphthalene	8.91	142	570723	39.814	ng	99
39) 1,2,4,5-Tetrachlorobenzene	9.07	216	216060	33.279	ng	98
40) Hexachlorocyclopentadiene	9.06	237	236297	79.641	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.19	196	167767	36.476	nq	98
43) 2,4,5-Trichlorophenol	9.23	196	175157	38.149	nq	97
45) 1,1'-Biphenyl	9.37	154	658477	35.194	nq	99
46) 2-Chloronaphthalene	9.40	162	517999	36.874	nq	97
47) 2-Nitroaniline	9.50	65	172086	40.006	nq	97
48) Acenaphthylene	9.82	152	798065	37.111	nq	99
49) Dimethylphthalate	9.67	163	623047	37.919	nq	99
50) 2,6-Dinitrotoluene	9.74	165	138805	38.804	nq	89
51) Acenaphthene	9.99	154	550860	40.438	nq	99
52) 3-Nitroaniline	9.90	138	126549	30.341	nq	99
53) 2,4-Dinitrophenol	10.02	184	118740	64.941	nq	93
54) Dibenzofuran	10.16	168	727130	40.090	nq	99
55) 4-Nitrophenol	10.07	139	262873	74.536	nq	90
56) 2,4-Dinitrotoluene	10.14	165	189768	40.877	nq	99
57) Fluorene	10.50	166	563785	38.673	nq	100
58) 2,3,4,6-Tetrachlorophenol	10.27	232	133287	42.024	nq	97
59) Diethylphthalate	10.37	149	629801	39.227	nq	98
60) 4-Chlorophenyl-phenylether	10.49	204	251489	38.404	nq	98
61) 4-Nitroaniline	10.52	138	158781	39.459	nq	95
62) Azobenzene	10.66	77	612184	41.469	nq	96
64) 4,6-Dinitro-2-methylphenol	10.55	198	86671	36.398	nq	83
65) n-Nitrosodiphenylamine	10.62	169	506087	37.327	nq	100
66) 4-Bromophenyl-phenylether	10.99	248	143060	37.344	nq	94
67) Hexachlorobenzene	11.05	284	146942	35.914	nq	96
68) Atrazine	11.14	200	160549	39.358	nq	99
69) Pentachlorophenol	11.24	266	164522	64.610	nq	99
70) Phenanthrene	11.47	178	820850	39.184	nq	99
71) Anthracene	11.52	178	850562	39.682	nq	99
72) Carbazole	11.67	167	798340	38.034	nq	99
73) Di-n-butylphthalate	12.00	149	1012988	40.094	nq	99
74) Fluoranthene	12.66	202	857606	40.428	nq	99
76) Benzidine	12.77	184	308896	29.533	nq	97
77) Pyrene	12.89	202	878121	40.722	nq	100
79) Butylbenzylphthalate	13.50	149	434367	42.860	nq	94
80) Benzo(a)anthracene	14.07	228	688081	40.183	nq	100
81) 3,3'-Dichlorobenzidine	14.03	252	204498	32.577	nq	99
82) Chrysene	14.12	228	640725	39.302	nq	98
83) Bis(2-ethylhexyl)phthalate	14.05	149	580813	41.622	nq	100
84) Di-n-octyl phthalate	14.66	149	918292	40.868	nq	99
85) Indeno(1,2,3-cd)pyrene	17.09	276	551106	42.259	nq	96
87) Benzo(b)fluoranthene	15.13	252	576827m	39.543	nq	
88) Benzo(k)fluoranthene	15.17	252	542614	37.661	nq	99
89) Benzo(a)pyrene	15.52	252	526122	39.292	nq	98
90) Dibenzo(a,h)anthracene	17.11	278	462079	43.121	nq	98
91) Benzo(g,h,i)perylene	17.55	276	464261	43.829	nq	96

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

