

Data Path : Z:\HPCHEM1\BNA F\DATA\BF101917\
 Data File : BF099669.D
 Acq On : 19 Oct 2017 20:55
 Operator : SJ/JU
 Sample : PB103438BS
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB103438BS

Quant Time: Oct 20 13:33:25 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 20 13:11:43 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.82	152	89824	20.00	ng	0.00
21) Naphthalene-d8	8.10	136	376134	20.00	ng	0.00
38) Acenaphthene-d10	9.86	164	167352	20.00	ng	0.00
63) Phenanthrene-d10	11.35	188	253247	20.00	ng	0.00
75) Chrysene-d12	13.99	240	157127	20.00	ng	0.00
86) Perylene-d12	15.44	264	172624	20.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.46	112	427944	75.84	ng	0.02
7) Phenol-d6	6.46	99	508437	73.04	ng	0.00
23) Nitrobenzene-d5	7.39	82	433104	73.84	ng	0.00
41) 2,4,6-Tribromophenol	10.65	330	96025	70.12	ng	0.00
44) 2-Fluorobiphenyl	9.17	172	863459	86.13	ng	0.00
78) Terphenyl-d14	12.92	244	565736	81.88	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	2.65	88	75756	28.106	ng	95
3) Pyridine	3.42	79	197531	27.090	ng	91
4) n-Nitrosodimethylamine	3.37	42	79764	25.797	ng	# 74
6) Aniline	6.49	93	174495	18.574	ng	# 84
8) 2-Chlorophenol	6.61	128	221392	34.647	ng	93
9) Benzaldehyde	6.37	77	55926	13.851	ng	97
10) Phenol	6.48	94	265599	34.118	ng	95
11) bis(2-Chloroethyl)ether	6.56	93	202649	33.485	ng	94
12) 1,3-Dichlorobenzene	6.76	146	232662	34.767	ng	97
13) 1,4-Dichlorobenzene	6.84	146	238361	35.008	ng	97
14) 1,2-Dichlorobenzene	6.99	146	217099	34.508	ng	97
15) Benzyl Alcohol	6.97	79	156533	30.654	ng	97
16) 2,2'-oxybis(1-Chloropropan	7.09	45	264816	28.085	ng	74
17) 2-Methylphenol	7.08	107	178324	34.107	ng	97
18) Hexachloroethane	7.33	117	79286	32.397	ng	96
19) n-Nitroso-di-n-propylamine	7.23	70	137547	30.950	ng	93
20) 3+4-Methylphenols	7.24	107	224221	33.899	ng	# 85
22) Acetophenone	7.23	105	287212	31.516	ng	# 98
24) Nitrobenzene	7.41	77	211151	31.736	ng	90
25) Isophorone	7.65	82	393736	32.470	ng	98
26) 2-Nitrophenol	7.72	139	119921	37.482	ng	89
27) 2,4-Dimethylphenol	7.76	122	194200	32.141	ng	96
28) bis(2-Chloroethoxy)methane	7.85	93	260628	34.489	ng	99
29) 2,4-Dichlorophenol	7.97	162	182632	35.205	ng	97
30) 1,2,4-Trichlorobenzene	8.05	180	188931	36.414	ng	99
31) Naphthalene	8.12	128	625284	35.396	ng	99
32) Benzoic acid	7.89	122	102963	20.745	ng	95
33) 4-Chloroaniline	8.17	127	109147	14.795	ng	97
34) Hexachlorobutadiene	8.23	225	93533	34.999	ng	100
35) Caprolactam	8.55	113	56023	33.667	ng	# 85
36) 4-Chloro-3-methylphenol	8.67	107	183810	31.524	ng	93
37) 2-Methylnaphthalene	8.82	142	425419	37.044	ng	97
39) 1,2,4,5-Tetrachlorobenzene	8.98	216	173649	34.793	ng	98
40) Hexachlorocyclopentadiene	8.96	237	40815	17.895	ng	97

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.10	196	114436	32.366	nq	99
43) 2,4,5-Trichlorophenol	9.15	196	120339	34.095	nq	95
45) 1,1'-Biphenyl	9.27	154	503427	35.002	nq	99
46) 2-Chloronaphthalene	9.30	162	389176	36.038	nq	96
47) 2-Nitroaniline	9.41	65	100680	30.448	nq	84
48) Acenaphthylene	9.72	152	572149	34.610	nq	99
49) Dimethylphthalate	9.57	163	449947	35.623	nq	99
50) 2,6-Dinitrotoluene	9.65	165	96529	35.104	nq	91
51) Acenaphthene	9.89	154	334574	31.950	nq	98
52) 3-Nitroaniline	9.82	138	69287	21.610	nq	84
53) 2,4-Dinitrophenol	9.94	184	28825	24.495	nq	# 87
54) Dibenzofuran	10.06	168	491987	35.286	nq	99
55) 4-Nitrophenol	9.99	139	107196	39.539	nq	88
56) 2,4-Dinitrotoluene	10.06	165	112497	31.523	nq	97
57) Fluorene	10.40	166	392886	35.058	nq	100
58) 2,3,4,6-Tetrachlorophenol	10.19	232	80045	32.830	nq	94
59) Diethylphthalate	10.27	149	420980	34.109	nq	98
60) 4-Chlorophenyl-phenylether	10.39	204	181529	36.060	nq	97
61) 4-Nitroaniline	10.43	138	80494	26.022	nq	90
62) Azobenzene	10.56	77	365705	32.226	nq	89
64) 4,6-Dinitro-2-methylphenol	10.47	198	28849	18.723	nq	# 78
65) n-Nitrosodiphenylamine	10.52	169	346201	39.461	nq	99
66) 4-Bromophenyl-phenylether	10.88	248	99680	40.212	nq	96
67) Hexachlorobenzene	10.96	284	95705	36.149	nq	# 88
68) Atrazine	11.04	200	102965	39.008	nq	99
69) Pentachlorophenol	11.15	266	64833	39.347	nq	99
70) Phenanthrene	11.37	178	487702	35.978	nq	98
71) Anthracene	11.42	178	504130	36.347	nq	98
72) Carbazole	11.58	167	435198	32.041	nq	98
73) Di-n-butylphthalate	11.89	149	610132	37.320	nq	99
74) Fluoranthene	12.56	202	418745	30.506	nq	100
76) Benzidine	12.68	184	216020	37.171	nq	97
77) Pyrene	12.79	202	412854	34.458	nq	99
79) Butylbenzylphthalate	13.39	149	198905	35.323	nq	91
80) Benzo(a)anthracene	13.97	228	340811	35.820	nq	100
81) 3,3'-Dichlorobenzidine	13.93	252	109619	31.429	nq	99
82) Chrysene	14.02	228	325317	35.914	nq	99
83) Bis(2-ethylhexyl)phthalate	13.95	149	276015	35.598	nq	# 99
84) Di-n-octyl phthalate	14.56	149	468283	37.508	nq	# 94
85) Indeno(1,2,3-cd)pyrene	16.92	276	355940	49.122	nq	96
87) Benzo(b)fluoranthene	15.02	252	369165	34.782	nq	# 96
88) Benzo(k)fluoranthene	15.05	252	323985	30.905	nq	99
89) Benzo(a)pyrene	15.39	252	345797	35.493	nq	# 93
90) Dibenzo(a,h)anthracene	16.93	278	301589	38.681	nq	98
91) Benzo(g,h,i)perylene	17.36	276	287889	37.354	nq	96

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

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