

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042723\
 Data File : BF133092.D
 Acq On : 27 Apr 2023 11:14
 Operator : CG\JU
 Sample : PB152384BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB152384BS

Quant Time: Apr 28 02:04:12 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042123.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 28 02:00:53 2023
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By : Christian Giraldo 04/28/2023
 Supervised By : Jagrut Upadhyay 04/28/2023

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.740	152	210908	20.000	ng	0.00
21) Naphthalene-d8	8.022	136	850691	20.000	ng	0.00
39) Acenaphthene-d10	9.775	164	443617	20.000	ng	0.00
64) Phenanthrene-d10	11.257	188	808689	20.000	ng	0.00
76) Chrysene-d12	13.898	240	565734	20.000	ng	0.00
86) Perylene-d12	15.316	264	518473	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.363	112	1435940	102.848	ng	0.02
7) Phenol-d6	6.387	99	1765195	101.861	ng	0.00
23) Nitrobenzene-d5	7.304	82	1066587	67.676	ng	0.00
42) 2,4,6-Tribromophenol	10.569	330	475202	106.514	ng	0.00
45) 2-Fluorobiphenyl	9.104	172	1873680	70.662	ng	0.00
79) Terphenyl-d14	12.845	244	2139665	65.229	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.499	88	257795	39.808	ng	96
3) Pyridine	3.193	79	633305	36.820	ng	96
4) n-Nitrosodimethylamine	3.152	42	333121	37.391	ng	94
6) Aniline	6.398	93	761321	34.110	ng	# 18
8) 2-Chlorophenol	6.528	128	622175	43.890	ng	96
9) Benzaldehyde	6.281	77	341589	44.378	ng	98
10) Phenol	6.398	94	772939	40.431	ng	# 72
11) bis(2-Chloroethyl)ether	6.475	93	596727	41.595	ng	99
12) 1,3-Dichlorobenzene	6.681	146	668649	44.365	ng	97
13) 1,4-Dichlorobenzene	6.757	146	672578	44.527	ng	98
14) 1,2-Dichlorobenzene	6.910	146	620801	44.580	ng	97
15) Benzyl Alcohol	6.887	79	498207	41.619	ng	99
16) 2,2'-oxybis(1-Chloropr...	7.022	45	964395	38.704	ng	94
17) 2-Methylphenol	7.004	107	559252	43.084	ng	97
18) Hexachloroethane	7.251	117	231396	44.063	ng	99
19) n-Nitroso-di-n-propyla...	7.163	70	416793	38.633	ng	94
20) 3+4-Methylphenols	7.157	107	633513	41.442	ng	# 88
22) Acetophenone	7.151	105	861719	43.729	ng	# 97
24) Nitrobenzene	7.328	77	650856	40.754	ng	96
25) Isophorone	7.563	82	1210790	40.463	ng	99
26) 2-Nitrophenol	7.640	139	367090	47.656	ng	95
27) 2,4-Dimethylphenol	7.687	122	592832	44.883	ng	99
28) bis(2-Chloroethoxy)met...	7.781	93	716154	40.455	ng	100
29) 2,4-Dichlorophenol	7.887	162	528470	43.950	ng	99
30) 1,2,4-Trichlorobenzene	7.969	180	551372	44.262	ng	98
31) Naphthalene	8.045	128	1797835	42.270	ng	99
32) Benzoic acid	7.822	122	435460	53.373	ng	96
33) 4-Chloroaniline	8.092	127	361657	21.045	ng	99
34) Hexachlorobutadiene	8.163	225	295994	42.871	ng	100
35) Caprolactam	8.475	113	194355m	48.115	ng	
36) 4-Chloro-3-methylphenol	8.581	107	578194	44.591	ng	98
37) 2-Methylnaphthalene	8.739	142	1184084	40.721	ng	98
38) 1-Methylnaphthalene	8.834	142	1176244	43.241	ng	99
40) 1,2,4,5-Tetrachloroben...	8.904	216	496064	41.590	ng	100
41) Hexachlorocyclopentadiene	8.892	237	672270	96.643	ng	99
43) 2,4,6-Trichlorophenol	9.016	196	352429	42.725	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.057	196	390781	40.963	ng	99
46) 1,1'-Biphenyl	9.204	154	1470476	41.802	ng	98
47) 2-Chloronaphthalene	9.222	162	1104337	42.892	ng	99
48) 2-Nitroaniline	9.322	65	347691	40.860	ng	95
49) Acenaphthylene	9.639	152	1708354	41.525	ng	100
50) Dimethylphthalate	9.504	163	1281480	41.428	ng	99
51) 2,6-Dinitrotoluene	9.563	165	308715	44.342	ng	93
52) Acenaphthene	9.810	154	1264266m	45.884	ng	
53) 3-Nitroaniline	9.734	138	252155	30.460	ng	94
54) 2,4-Dinitrophenol	9.839	184	370370	105.835	ng	97
55) Dibenzofuran	9.986	168	1534243	40.820	ng	98
56) 4-Nitrophenol	9.904	139	579557	90.393	ng	95
57) 2,4-Dinitrotoluene	9.969	165	405280	45.647	ng	90
58) Fluorene	10.328	166	1140967	41.434	ng	98
59) 2,3,4,6-Tetrachlorophenol	10.104	232	323672	43.764	ng	98
60) Diethylphthalate	10.204	149	1197139	40.973	ng	99
61) 4-Chlorophenyl-phenyle...	10.316	204	550744	41.041	ng	96
62) 4-Nitroaniline	10.351	138	354721	46.622	ng	95
63) Azobenzene	10.481	77	1205997	38.964	ng	95
65) 4,6-Dinitro-2-methylph...	10.375	198	241897	49.350	ng	84
66) n-Nitrosodiphenylamine	10.439	169	1059065	40.373	ng	99
67) 4-Bromophenyl-phenylether	10.804	248	355948	40.584	ng	96
68) Hexachlorobenzene	10.875	284	396834	44.155	ng	98
69) Atrazine	10.969	200	348260	46.075	ng	97
70) Pentachlorophenol	11.069	266	492840	76.998	ng	98
71) Phenanthrene	11.286	178	1768693	40.693	ng	100
72) Anthracene	11.339	178	1818665	40.886	ng	100
73) Carbazole	11.492	167	1665681	42.055	ng	99
74) Di-n-butylphthalate	11.828	149	1884611	42.046	ng	99
75) Fluoranthene	12.469	202	1860656	42.654	ng	98
77) Benzidine	12.592	184	1028369	107.493	ng	# 93
78) Pyrene	12.698	202	1913657	39.460	ng	99
80) Butylbenzylphthalate	13.322	149	814856	47.454	ng	100
81) Benzo(a)anthracene	13.886	228	1567028	40.280	ng	99
82) 3,3'-Dichlorobenzidine	13.851	252	371851	34.600	ng	97
83) Chrysene	13.922	228	1573810	40.464	ng	100
84) Bis(2-ethylhexyl)phtha...	13.886	149	923213	42.834	ng	98
85) Di-n-octyl phthalate	14.492	149	1727080	49.143	ng	96
87) Indeno(1,2,3-cd)pyrene	16.698	276	1357661	46.035	ng	100
88) Benzo(b)fluoranthene	14.910	252	1561245	47.332	ng	98
89) Benzo(k)fluoranthene	14.939	252	1299101	39.731	ng	99
90) Benzo(a)pyrene	15.257	252	1212286	40.379	ng	99
91) Dibenzo(a,h)anthracene	16.715	278	1108569	45.068	ng	99
92) Benzo(g,h,i)perylene	17.115	276	1137014	46.821	ng	99

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

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