

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF100424\
 Data File : BF139652.D
 Acq On : 04 Oct 2024 12:43
 Operator : RC/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Quant Time: Oct 04 13:05:03 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF100124.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 02 04:58:47 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.863	152	102462	20.000	ng	0.00
21) Naphthalene-d8	8.151	136	379259	20.000	ng	0.00
39) Acenaphthene-d10	9.904	164	208025	20.000	ng	0.00
64) Phenanthrene-d10	11.386	188	382146	20.000	ng	0.00
76) Chrysene-d12	14.027	240	210514	20.000	ng	0.00
86) Perylene-d12	15.498	264	230633	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.481	112	470342	79.563	ng	0.00
7) Phenol-d6	6.498	99	619401	77.914	ng	0.00
23) Nitrobenzene-d5	7.434	82	579590	77.374	ng	0.00
42) 2,4,6-Tribromophenol	10.692	330	150960	75.174	ng	0.00
45) 2-Fluorobiphenyl	9.222	172	1040674	76.800	ng	0.00
79) Terphenyl-d14	12.975	244	959279	74.770	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.628	88	106022	41.483	ng	100
3) Pyridine	3.375	79	260893	41.972	ng	99
4) n-Nitrosodimethylamine	3.340	42	159200	39.128	ng	95
6) Aniline	6.528	93	274760	38.949	ng	99
8) 2-Chlorophenol	6.651	128	251229	39.666	ng	99
9) Benzaldehyde	6.416	77	181334	43.060	ng	99
10) Phenol	6.516	94	328411	39.287	ng	94
11) bis(2-Chloroethyl)ether	6.604	93	246076	36.155	ng	98
12) 1,3-Dichlorobenzene	6.804	146	278590	39.100	ng	99
13) 1,4-Dichlorobenzene	6.881	146	279643	38.752	ng	100
14) 1,2-Dichlorobenzene	7.040	146	261177	38.617	ng	99
15) Benzyl Alcohol	7.010	79	226288	37.254	ng	98
16) 2,2'-oxybis(1-Chloropr...	7.140	45	437957	37.170	ng	98
17) 2-Methylphenol	7.122	107	204976	38.767	ng	97
18) Hexachloroethane	7.375	117	103094	38.302	ng	98
19) n-Nitroso-di-n-propyla...	7.281	70	185437	36.282	ng	98
20) 3+4-Methylphenols	7.275	107	252378	38.002	ng	98
22) Acetophenone	7.275	105	343673	38.176	ng	98
24) Nitrobenzene	7.451	77	302558	39.006	ng	100
25) Isophorone	7.687	82	498991	37.506	ng	99
26) 2-Nitrophenol	7.763	139	133875	40.148	ng	100
27) 2,4-Dimethylphenol	7.804	122	157086	38.481	ng	97
28) bis(2-Chloroethoxy)met...	7.898	93	299551	37.705	ng	99
29) 2,4-Dichlorophenol	8.010	162	207293	38.931	ng	98
30) 1,2,4-Trichlorobenzene	8.087	180	232441	38.601	ng	99
31) Naphthalene	8.169	128	752920	38.828	ng	100
32) Benzoic acid	7.922	122	159788	39.365	ng	99
33) 4-Chloroaniline	8.222	127	248727	37.894	ng	98
34) Hexachlorobutadiene	8.287	225	150300	38.570	ng	100
35) Caprolactam	8.592	113	66657	38.168	ng	99
36) 4-Chloro-3-methylphenol	8.704	107	229034	37.999	ng	99
37) 2-Methylnaphthalene	8.863	142	476538	37.733	ng	99
38) 1-Methylnaphthalene	8.957	142	464905	37.660	ng	100
40) 1,2,4,5-Tetrachloroben...	9.028	216	226915	39.282	ng	98
41) Hexachlorocyclopentadiene	9.010	237	63461	35.828	ng	99
43) 2,4,6-Trichlorophenol	9.139	196	150287	38.593	ng	99

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44) 2,4,5-Trichlorophenol	9.181	196	165622	39.438	ng	99
46) 1,1'-Biphenyl	9.322	154	596198	38.533	ng	99
47) 2-Chloronaphthalene	9.351	162	453208	39.073	ng	98
48) 2-Nitroaniline	9.445	65	168647	40.504	ng	98
49) Acenaphthylene	9.763	152	685469	38.860	ng	100
50) Dimethylphthalate	9.628	163	525252	38.572	ng	100
51) 2,6-Dinitrotoluene	9.686	165	120472	39.163	ng	99
52) Acenaphthene	9.939	154	467215	38.582	ng	99
53) 3-Nitroaniline	9.857	138	122775	39.056	ng	97
54) 2,4-Dinitrophenol	9.963	184	61761	37.328	ng	98
55) Dibenzofuran	10.110	168	653933	38.568	ng	100
56) 4-Nitrophenol	10.022	139	89676	37.408	ng	92
57) 2,4-Dinitrotoluene	10.092	165	162138	40.324	ng	97
58) Fluorene	10.451	166	513669	38.037	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.228	232	129875	38.290	ng	98
60) Diethylphthalate	10.322	149	536166	38.120	ng	99
61) 4-Chlorophenyl-phenyle...	10.445	204	254896	37.728	ng	97
62) 4-Nitroaniline	10.469	138	125846	39.138	ng	99
63) Azobenzene	10.604	77	536029	37.895	ng	99
65) 4,6-Dinitro-2-methylph...	10.498	198	87719	40.643	ng	99
66) n-Nitrosodiphenylamine	10.563	169	442746	39.305	ng	99
67) 4-Bromophenyl-phenylether	10.933	248	150009	38.286	ng	98
68) Hexachlorobenzene	10.998	284	166667	38.334	ng	99
69) Atrazine	11.086	200	126041	41.705	ng	99
70) Pentachlorophenol	11.192	266	96694	37.350	ng	99
71) Phenanthrene	11.416	178	692249	38.419	ng	100
72) Anthracene	11.463	178	683891	38.365	ng	100
73) Carbazole	11.622	167	653732	38.184	ng	100
74) Di-n-butylphthalate	11.945	149	813276	39.063	ng	100
75) Fluoranthene	12.598	202	721008	36.605	ng	100
77) Benzidine	12.727	184	228756	61.606	ng	99
78) Pyrene	12.827	202	715774	38.016	ng	100
80) Butylbenzylphthalate	13.445	149	281407	41.785	ng	98
81) Benzo(a)anthracene	14.016	228	541176	39.101	ng	100
82) 3,3'-Dichlorobenzidine	13.980	252	184505	43.558	ng	98
83) Chrysene	14.057	228	473814	37.445	ng	99
84) Bis(2-ethylhexyl)phtha...	14.004	149	374783	44.566	ng	100
85) Di-n-octyl phthalate	14.616	149	566236	45.832	ng	100
87) Indeno(1,2,3-cd)pyrene	16.968	276	558741	41.161	ng	99
88) Benzo(b)fluoranthene	15.068	252	509730	34.182	ng	100
89) Benzo(k)fluoranthene	15.098	252	495647	39.183	ng	100
90) Benzo(a)pyrene	15.433	252	452282	38.465	ng	99
91) Dibenzo(a,h)anthracene	16.986	278	467106	41.488	ng	99
92) Benzo(g,h,i)perylene	17.415	276	458425	40.599	ng	99

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

