

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101823\
 Data File : BF135838.D
 Acq On : 18 Oct 2023 10:27
 Operator : CG\JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Quant Time: Oct 19 01:13:33 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF101323.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 17 01:22:00 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.740	152	95918	20.000	ng	0.00
21) Naphthalene-d8	8.022	136	334120	20.000	ng	0.00
39) Acenaphthene-d10	9.781	164	174848	20.000	ng	0.00
64) Phenanthrene-d10	11.275	188	340988	20.000	ng	0.00
76) Chrysene-d12	13.945	240	227882	20.000	ng	0.00
86) Perylene-d12	15.421	264	206413	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.375	112	529503	88.754	ng	0.00
7) Phenol-d6	6.404	99	572187	76.867	ng	0.00
23) Nitrobenzene-d5	7.316	82	544645	89.654	ng	0.00
42) 2,4,6-Tribromophenol	10.581	330	142575	84.977	ng	0.00
45) 2-Fluorobiphenyl	9.104	172	949959	84.946	ng	0.00
79) Terphenyl-d14	12.874	244	1031526	83.828	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.416	88	139666	48.535	ng	99
3) Pyridine	3.175	79	285833	40.483	ng	99
4) n-Nitrosodimethylamine	3.152	42	161825	42.619	ng	93
6) Aniline	6.410	93	352664	37.646	ng	89
8) 2-Chlorophenol	6.534	128	253039	39.331	ng	99
9) Benzaldehyde	6.304	77	164238	38.765	ng	98
10) Phenol	6.416	94	309076	39.626	ng	96
11) bis(2-Chloroethyl)ether	6.487	93	236514	37.633	ng	99
12) 1,3-Dichlorobenzene	6.681	146	257953	39.550	ng	99
13) 1,4-Dichlorobenzene	6.757	146	256180	38.773	ng	99
14) 1,2-Dichlorobenzene	6.910	146	236341	39.727	ng	99
15) Benzyl Alcohol	6.898	79	199863	40.380	ng	98
16) 2,2'-oxybis(1-Chloropr...	7.016	45	406039	35.723	ng	64
17) 2-Methylphenol	7.016	107	202485	36.357	ng	97
18) Hexachloroethane	7.240	117	93583	39.064	ng	94
19) n-Nitroso-di-n-propyla...	7.163	70	175393	38.107	ng	99
20) 3+4-Methylphenols	7.169	107	276619	39.841	ng	95
22) Acetophenone	7.157	105	342826	44.739	ng	98
24) Nitrobenzene	7.334	77	266947	43.783	ng	99
25) Isophorone	7.569	82	482287	42.081	ng	99
26) 2-Nitrophenol	7.645	139	124032	41.685	ng	98
27) 2,4-Dimethylphenol	7.692	122	217179	44.338	ng	99
28) bis(2-Chloroethoxy)met...	7.775	93	296931	43.504	ng	99
29) 2,4-Dichlorophenol	7.892	162	205918	44.859	ng	98
30) 1,2,4-Trichlorobenzene	7.963	180	208477	44.405	ng	99
31) Naphthalene	8.045	128	660857	40.551	ng	99
32) Benzoic acid	7.845	122	143921	43.407	ng	96
33) 4-Chloroaniline	8.104	127	316261	44.425	ng	97
34) Hexachlorobutadiene	8.151	225	128447	46.221	ng	97
35) Caprolactam	8.498	113	67404	42.954	ng	# 85
36) 4-Chloro-3-methylphenol	8.598	107	227628	44.716	ng	95
37) 2-Methylnaphthalene	8.734	142	444392	41.388	ng	100
38) 1-Methylnaphthalene	8.834	142	417903	42.198	ng	100
40) 1,2,4,5-Tetrachloroben...	8.904	216	212190	42.188	ng	100
41) Hexachlorocyclopentadiene	8.881	237	18863	36.877	ng	98
43) 2,4,6-Trichlorophenol	9.028	196	131612	40.607	ng	97

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44) 2,4,5-Trichlorophenol	9.081	196	159757	42.379	ng	96
46) 1,1'-Biphenyl	9.204	154	575424	42.962	ng	98
47) 2-Chloronaphthalene	9.228	162	411182	42.597	ng	99
48) 2-Nitroaniline	9.339	65	163109	42.809	ng	96
49) Acenaphthylene	9.645	152	652475	40.846	ng	99
50) Dimethylphthalate	9.510	163	477439	40.449	ng	99
51) 2,6-Dinitrotoluene	9.586	165	108928	42.914	ng	# 80
52) Acenaphthene	9.816	154	400215	41.800	ng	98
53) 3-Nitroaniline	9.757	138	133441	42.886	ng	99
54) 2,4-Dinitrophenol	9.886	184	45470	42.915	ng	95
55) Dibenzofuran	9.986	168	616071	42.644	ng	99
56) 4-Nitrophenol	9.951	139	44410	34.504	ng	99
57) 2,4-Dinitrotoluene	9.998	165	140693	43.569	ng	95
58) Fluorene	10.334	166	480179	42.456	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.122	232	114808	40.300	ng	98
60) Diethylphthalate	10.210	149	499385	41.443	ng	99
61) 4-Chlorophenyl-phenyle...	10.322	204	232724	42.575	ng	97
62) 4-Nitroaniline	10.381	138	124747	43.574	ng	97
63) Azobenzene	10.486	77	482082	41.178	ng	98
65) 4,6-Dinitro-2-methylph...	10.416	198	71680	41.770	ng	88
66) n-Nitrosodiphenylamine	10.451	169	430983	41.468	ng	98
67) 4-Bromophenyl-phenylether	10.816	248	143846	42.024	ng	97
68) Hexachlorobenzene	10.886	284	143216	41.096	ng	97
69) Atrazine	10.981	200	115767	40.951	ng	99
70) Pentachlorophenol	11.092	266	52916	38.499	ng	96
71) Phenanthrene	11.304	178	689379	42.203	ng	98
72) Anthracene	11.357	178	709902	41.940	ng	99
73) Carbazole	11.522	167	663998	42.061	ng	99
74) Di-n-butylphthalate	11.839	149	788556	40.684	ng	100
75) Fluoranthene	12.504	202	755569	41.676	ng	97
77) Benzidine	12.627	184	188951	37.998	ng	99
78) Pyrene	12.733	202	765704	42.284	ng	99
80) Butylbenzylphthalate	13.351	149	318269	39.996	ng	97
81) Benzo(a)anthracene	13.933	228	637527	42.592	ng	99
82) 3,3'-Dichlorobenzidine	13.898	252	203831	41.060	ng	# 97
83) Chrysene	13.974	228	607912	41.931	ng	99
84) Bis(2-ethylhexyl)phtha...	13.910	149	420804	39.017	ng	98
85) Di-n-octyl phthalate	14.533	149	635137	37.138	ng	# 99
87) Indeno(1,2,3-cd)pyrene	16.939	276	470615	41.844	ng	98
88) Benzo(b)fluoranthene	14.992	252	469751	40.663	ng	99
89) Benzo(k)fluoranthene	15.021	252	494418	44.083	ng	99
90) Benzo(a)pyrene	15.363	252	445079	40.972	ng	99
91) Dibenzo(a,h)anthracene	16.939	278	383521	41.597	ng	98
92) Benzo(g,h,i)perylene	17.392	276	426652	45.049	ng	97

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

