

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF022723\
 Data File : BF132545.D
 Acq On : 27 Feb 2023 11:00
 Operator : CG\JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 02/28/2023
 Supervised By :Jagrut Upadhyay 02/28/2023

Quant Time: Feb 27 13:18:38 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF022223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 24 23:17:30 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.007	152	177623	20.000	ng	0.00
21) Naphthalene-d8	8.289	136	619750	20.000	ng	0.00
39) Acenaphthene-d10	10.054	164	335185	20.000	ng	0.00
64) Phenanthrene-d10	11.548	188	653363	20.000	ng	0.00
76) Chrysene-d12	14.207	240	421616	20.000	ng	0.00
86) Perylene-d12	15.759	264	357603	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.625	112	883079	80.688	ng	0.00
7) Phenol-d6	6.637	99	1134103	79.401	ng	0.00
23) Nitrobenzene-d5	7.572	82	1080806	82.740	ng	0.00
42) 2,4,6-Tribromophenol	10.848	330	325824	90.010	ng	0.00
45) 2-Fluorobiphenyl	9.372	172	1714216	77.873	ng	0.00
79) Terphenyl-d14	13.142	244	2045570	82.031	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.872	88	259720m	42.901	ng	
3) Pyridine	3.648	79	639705	39.808	ng	99
4) n-Nitrosodimethylamine	3.613	42	257879	42.485	ng	98
6) Aniline	6.672	93	798876	41.561	ng	99
8) 2-Chlorophenol	6.795	128	462777	40.742	ng	94
9) Benzaldehyde	6.560	77	299104	38.808	ng	98
10) Phenol	6.648	94	654923	39.994	ng	100
11) bis(2-Chloroethyl)ether	6.742	93	505245	39.967	ng	95
12) 1,3-Dichlorobenzene	6.948	146	495939	40.686	ng	98
13) 1,4-Dichlorobenzene	7.025	146	497904	40.908	ng	99
14) 1,2-Dichlorobenzene	7.178	146	452699	38.755	ng	99
15) Benzyl Alcohol	7.148	79	478997	43.544	ng	99
16) 2,2'-oxybis(1-Chloropr...	7.278	45	656267	40.710	ng	99
17) 2-Methylphenol	7.254	107	431914	40.743	ng	99
18) Hexachloroethane	7.519	117	200344	42.132	ng	99
19) n-Nitroso-di-n-propyla...	7.419	70	342347	38.945	ng	97
20) 3+4-Methylphenols	7.407	107	518140	37.832	ng	# 77
22) Acetophenone	7.419	105	639139	40.899	ng	# 86
24) Nitrobenzene	7.595	77	592653	42.422	ng	98
25) Isophorone	7.825	82	1040326	41.540	ng	98
26) 2-Nitrophenol	7.907	139	258852	41.994	ng	99
27) 2,4-Dimethylphenol	7.936	122	398762	40.989	ng	97
28) bis(2-Chloroethoxy)met...	8.031	93	605369	41.321	ng	99
29) 2,4-Dichlorophenol	8.148	162	405937	41.947	ng	97
30) 1,2,4-Trichlorobenzene	8.231	180	441037	41.970	ng	97
31) Naphthalene	8.313	128	1299654	40.963	ng	99
32) Benzoic acid	8.060	122	330542m	43.507	ng	
33) 4-Chloroaniline	8.360	127	575981	42.364	ng	99
34) Hexachlorobutadiene	8.425	225	296214	44.360	ng	99
35) Caprolactam	8.742	113	137635	43.140	ng	95
36) 4-Chloro-3-methylphenol	8.836	107	450473	42.386	ng	99
37) 2-Methylnaphthalene	9.007	142	894063	41.350	ng	99
38) 1-Methylnaphthalene	9.107	142	834917	41.319	ng	99
40) 1,2,4,5-Tetrachloroben...	9.172	216	470370	42.892	ng	100
41) Hexachlorocyclopentadiene	9.154	237	273514	45.983	ng	100
43) 2,4,6-Trichlorophenol	9.283	196	312041	41.987	ng	99

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44) 2,4,5-Trichlorophenol	9.325	196	368552	42.489	ng	95
46) 1,1'-Biphenyl	9.472	154	1040136	40.999	ng	98
47) 2-Chloronaphthalene	9.501	162	832982	41.413	ng	99
48) 2-Nitroaniline	9.595	65	308340	41.622	ng	98
49) Acenaphthylene	9.919	152	1315179	41.084	ng	99
50) Dimethylphthalate	9.772	163	1017352	41.685	ng	99
51) 2,6-Dinitrotoluene	9.836	165	235764	42.418	ng	98
52) Acenaphthene	10.089	154	856127	42.117	ng	100
53) 3-Nitroaniline	10.007	138	263215	42.335	ng	96
54) 2,4-Dinitrophenol	10.113	184	147773	43.062	ng	97
55) Dibenzofuran	10.260	168	1215753	41.026	ng	95
56) 4-Nitrophenol	10.166	139	194792	42.010	ng	95
57) 2,4-Dinitrotoluene	10.242	165	319842	43.255	ng	98
58) Fluorene	10.607	166	938547	41.406	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.377	232	283893	44.059	ng	98
60) Diethylphthalate	10.472	149	1021338	42.849	ng	99
61) 4-Chlorophenyl-phenyle...	10.595	204	495312	42.122	ng	97
62) 4-Nitroaniline	10.630	138	259461	42.513	ng	97
63) Azobenzene	10.754	77	1079703	42.771	ng	95
65) 4,6-Dinitro-2-methylph...	10.654	198	197220	44.590	ng	99
66) n-Nitrosodiphenylamine	10.713	169	833934	40.941	ng	99
67) 4-Bromophenyl-phenylether	11.089	248	317154	42.913	ng	98
68) Hexachlorobenzene	11.154	284	329858	42.417	ng	# 90
69) Atrazine	11.242	200	275129	43.266	ng	97
70) Pentachlorophenol	11.354	266	221971	47.975	ng	98
71) Phenanthrene	11.577	178	1386940	40.976	ng	99
72) Anthracene	11.630	178	1422084	40.800	ng	100
73) Carbazole	11.783	167	1279339	40.710	ng	99
74) Di-n-butylphthalate	12.101	149	1563100	42.404	ng	99
75) Fluoranthene	12.771	202	1557840	42.101	ng	98
77) Benzidine	12.889	184	440347	42.905	ng	100
78) Pyrene	13.007	202	1583040	41.298	ng	100
80) Butylbenzylphthalate	13.613	149	650847	43.026	ng	99
81) Benzo(a)anthracene	14.201	228	1321677	41.896	ng	99
82) 3,3'-Dichlorobenzidine	14.154	252	379707	40.827	ng	98
83) Chrysene	14.236	228	1211494	41.068	ng	100
84) Bis(2-ethylhexyl)phtha...	14.165	149	802254	43.173	ng	100
85) Di-n-octyl phthalate	14.795	149	1164937	42.900	ng	99
87) Indeno(1,2,3-cd)pyrene	17.365	276	1076036	45.922	ng	98
88) Benzo(b)fluoranthene	15.295	252	1061754	44.305	ng	100
89) Benzo(k)fluoranthene	15.330	252	944747	40.281	ng	99
90) Benzo(a)pyrene	15.695	252	950386	42.373	ng	98
91) Dibenzo(a,h)anthracene	17.383	278	876837	45.928	ng	98
92) Benzo(g,h,i)perylene	17.853	276	924001	43.001	ng	98

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

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