

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF010423\
 Data File : BF131894.D
 Acq On : 04 Jan 2023 12:15
 Operator : CG\JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 01/05/2023
 Supervised By : Jagrut Upadhyay 01/05/2023

Quant Time: Jan 05 05:33:04 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 30 02:06:40 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.816	152	70957	20.000	ng	0.00
21) Naphthalene-d8	8.104	136	251515	20.000	ng	0.00
39) Acenaphthene-d10	9.869	164	148792	20.000	ng	0.00
64) Phenanthrene-d10	11.363	188	298626	20.000	ng	0.00
76) Chrysene-d12	14.015	240	196599	20.000	ng	0.00
86) Perylene-d12	15.492	264	135672	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.416	112	357811	83.568	ng	0.00
7) Phenol-d6	6.451	99	464271	82.532	ng	0.00
23) Nitrobenzene-d5	7.392	82	509406	80.863	ng	0.00
42) 2,4,6-Tribromophenol	10.663	330	136006	84.129	ng	0.00
45) 2-Fluorobiphenyl	9.186	172	894572	82.351	ng	0.00
79) Terphenyl-d14	12.957	244	1082447	93.133	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.493	88	79582	40.081	ng	98
3) Pyridine	3.228	79	232846	41.715	ng	97
4) n-Nitrosodimethylamine	3.204	42	108433	41.314	ng	99
6) Aniline	6.481	93	284692	41.730	ng	99
8) 2-Chlorophenol	6.598	128	188079	41.693	ng	93
9) Benzaldehyde	6.363	77	142088	39.365	ng	95
10) Phenol	6.463	94	276424	41.263	ng	95
11) bis(2-Chloroethyl)ether	6.557	93	197189	41.163	ng	99
12) 1,3-Dichlorobenzene	6.757	146	210075	40.751	ng	98
13) 1,4-Dichlorobenzene	6.833	146	215856	41.549	ng	97
14) 1,2-Dichlorobenzene	6.986	146	202006	41.381	ng	98
15) Benzyl Alcohol	6.963	79	209258	42.420	ng	96
16) 2,2'-oxybis(1-Chloropr...	7.092	45	252776	41.348	ng	97
17) 2-Methylphenol	7.075	107	172315	41.784	ng	100
18) Hexachloroethane	7.328	117	87205	41.569	ng	96
19) n-Nitroso-di-n-propyla...	7.239	70	163000	41.028	ng	98
20) 3+4-Methylphenols	7.233	107	224043	41.637	ng	98
22) Acetophenone	7.233	105	297579	40.357	ng	98
24) Nitrobenzene	7.410	77	258398	40.929	ng	99
25) Isophorone	7.651	82	424482	41.261	ng	98
26) 2-Nitrophenol	7.722	139	100972	41.135	ng	93
27) 2,4-Dimethylphenol	7.763	122	145492	41.553	ng	97
28) bis(2-Chloroethoxy)met...	7.857	93	226865	41.105	ng	99
29) 2,4-Dichlorophenol	7.969	162	174990	41.482	ng	98
30) 1,2,4-Trichlorobenzene	8.045	180	204067	40.801	ng	99
31) Naphthalene	8.128	128	565608	41.063	ng	99
32) Benzoic acid	7.892	122	99385m	38.835	ng	
33) 4-Chloroaniline	8.180	127	219440	41.520	ng	98
34) Hexachlorobutadiene	8.239	225	139368	40.729	ng	99
35) Caprolactam	8.569	113	52605	42.119	ng	95
36) 4-Chloro-3-methylphenol	8.669	107	204224	42.836	ng	97
37) 2-Methylnaphthalene	8.822	142	383436	41.477	ng	99
38) 1-Methylnaphthalene	8.922	142	370768	41.290	ng	100
40) 1,2,4,5-Tetrachloroben...	8.986	216	211547	40.492	ng	99
41) Hexachlorocyclopentadiene	8.969	237	66184	32.134	ng	99
43) 2,4,6-Trichlorophenol	9.104	196	137966	41.232	ng	98

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44) 2,4,5-Trichlorophenol	9.145	196	148805	41.140	ng	98
46) 1,1'-Biphenyl	9.292	154	468704	41.145	ng	98
47) 2-Chloronaphthalene	9.316	162	387013	41.065	ng	99
48) 2-Nitroaniline	9.416	65	145297	42.061	ng	95
49) Acenaphthylene	9.733	152	585942	41.754	ng	99
50) Dimethylphthalate	9.598	163	476706	42.349	ng	99
51) 2,6-Dinitrotoluene	9.663	165	103298	42.391	ng	91
52) Acenaphthene	9.904	154	352266	41.593	ng	99
53) 3-Nitroaniline	9.833	138	103712	41.770	ng	97
54) 2,4-Dinitrophenol	9.939	184	56456	40.259	ng	96
55) Dibenzofuran	10.074	168	550416	42.182	ng	98
56) 4-Nitrophenol	9.998	139	68023	39.109	ng	94
57) 2,4-Dinitrotoluene	10.069	165	140983	42.305	ng	97
58) Fluorene	10.421	166	439231	41.232	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.198	232	116833m	39.361	ng	
60) Diethylphthalate	10.292	149	453239	41.835	ng	98
61) 4-Chlorophenyl-phenyle...	10.410	204	234850	42.409	ng	98
62) 4-Nitroaniline	10.451	138	108652	41.989	ng	98
63) Azobenzene	10.574	77	499491	41.542	ng	99
65) 4,6-Dinitro-2-methylph...	10.480	198	83321	40.980	ng	98
66) n-Nitrosodiphenylamine	10.533	169	372003	40.837	ng	96
67) 4-Bromophenyl-phenylether	10.904	248	140700	40.652	ng	94
68) Hexachlorobenzene	10.969	284	156818	41.245	ng	92
69) Atrazine	11.063	200	129748	43.264	ng	97
70) Pentachlorophenol	11.168	266	74519	39.609	ng	98
71) Phenanthrene	11.386	178	657700	40.050	ng	99
72) Anthracene	11.439	178	650473	40.491	ng	99
73) Carbazole	11.598	167	565990	39.909	ng	99
74) Di-n-butylphthalate	11.927	149	714054	41.655	ng	99
75) Fluoranthene	12.580	202	758298	39.643	ng	99
77) Benzidine	12.704	184	138729	42.939	ng	99
78) Pyrene	12.810	202	758537	46.033	ng	99
80) Butylbenzylphthalate	13.433	149	289310	47.335	ng	99
81) Benzo(a)anthracene	14.004	228	612557	43.034	ng	98
82) 3,3'-Dichlorobenzidine	13.968	252	170845	41.526	ng	98
83) Chrysene	14.045	228	564050	42.169	ng	99
84) Bis(2-ethylhexyl)phtha...	13.992	149	336548	45.224	ng	100
85) Di-n-octyl phthalate	14.604	149	441102	44.031	ng	100
87) Indeno(1,2,3-cd)pyrene	16.980	276	382774	44.409	ng	99
88) Benzo(b)fluoranthene	15.062	252	391564	39.361	ng	99
89) Benzo(k)fluoranthene	15.092	252	401722	41.794	ng	99
90) Benzo(a)pyrene	15.433	252	310780	40.351	ng	98
91) Dibenzo(a,h)anthracene	16.998	278	321993	45.606	ng	98
92) Benzo(g,h,i)perylene	17.427	276	320662	39.449	ng	98

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

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