

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF071422\
 Data File : BF129220.D
 Acq On : 14 Jul 2022 14:58
 Operator : CG\JU
 Sample : SSTDICC060
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC060

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 07/15/2022
 Supervised By : Jagrut Upadhyay 07/15/2022

Quant Time: Jul 14 15:47:03 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF071422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 14 14:17:14 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.904	152	330932	20.000	ng	0.00
21) Naphthalene-d8	8.187	136	1255669	20.000	ng	0.00
39) Acenaphthene-d10	9.945	164	628303	20.000	ng	0.00
64) Phenanthrene-d10	11.433	188	1027350	20.000	ng	# 0.00
76) Chrysene-d12	14.080	240	634146	20.000	ng	0.00
86) Perylene-d12	15.568	264	531685	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.528	112	2274889	116.217	ng	0.00
7) Phenol-d6	6.540	99	3022966	124.346	ng	0.00
23) Nitrobenzene-d5	7.475	82	2760663	131.136	ng	0.00
42) 2,4,6-Tribromophenol	10.739	330	648285	94.891	ng	0.00
45) 2-Fluorobiphenyl	0.000	172	0d	0.000	ng	
79) Terphenyl-d14	13.021	244	3644711	119.353	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.740	88	566587	64.924	ng	93
3) Pyridine	3.510	79	1459212	68.468	ng	95
4) n-Nitrosodimethylamine	3.487	42	725901	75.177	ng	95
6) Aniline	6.563	93	1834513m	67.473	ng	
8) 2-Chlorophenol	6.692	128	1217161	59.185	ng	99
9) Benzaldehyde	6.457	77	667055	52.066	ng	99
10) Phenol	6.551	94	1574924	63.940	ng	97
11) bis(2-Chloroethyl)ether	6.645	93	1253063	64.274	ng	95
12) 1,3-Dichlorobenzene	6.845	146	1297067	57.294	ng	98
13) 1,4-Dichlorobenzene	6.922	146	1294628	56.696	ng	98
14) 1,2-Dichlorobenzene	7.075	146	1174218	54.703	ng	97
15) Benzyl Alcohol	7.045	79	1087255	63.580	ng	99
16) 2,2'-oxybis(1-Chloropr...	7.175	45	1983895	71.248	ng	95
17) 2-Methylphenol	7.151	107	1021091	61.016	ng	99
18) Hexachloroethane	7.416	117	506035	63.123	ng	90
19) n-Nitroso-di-n-propyla...	7.328	70	874345	62.172	ng	94
20) 3+4-Methylphenols	7.310	107	1200212	56.398	ng	97
22) Acetophenone	7.322	105	1605428	55.734	ng	# 97
24) Nitrobenzene	7.492	77	1297901	63.719	ng	97
25) Isophorone	7.728	82	2314264	65.047	ng	100
26) 2-Nitrophenol	7.804	139	668170	70.582	ng	94
27) 2,4-Dimethylphenol	7.834	122	1030298	60.284	ng	99
28) bis(2-Chloroethoxy)met...	7.934	93	1505246	61.604	ng	99
29) 2,4-Dichlorophenol	8.045	162	998490	57.278	ng	99
30) 1,2,4-Trichlorobenzene	8.128	180	1007950	52.544	ng	97
31) Naphthalene	8.210	128	3294542	57.795	ng	98
32) Benzoic acid	7.981	122	883820	64.455	ng	98
33) 4-Chloroaniline	8.257	127	1402202	61.046	ng	99
34) Hexachlorobutadiene	8.322	225	569440	49.714	ng	99
35) Caprolactam	8.657	113	349329m	63.195	ng	
36) 4-Chloro-3-methylphenol	8.739	107	1098778	62.037	ng	97
37) 2-Methylnaphthalene	8.898	142	2141882	55.840	ng	99
38) 1-Methylnaphthalene	8.998	142	2060232	55.308	ng	99
40) 1,2,4,5-Tetrachloroben...	9.063	216	884184	50.066	ng	98
41) Hexachlorocyclopentadiene	9.051	237	498917	53.519	ng	99
43) 2,4,6-Trichlorophenol	9.175	196	681645	56.665	ng	98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.216	196	700839	54.773	ng	# 91
46) 1,1'-Biphenyl	9.369	154	2471778	54.721	ng	99
47) 2-Chloronaphthalene	9.392	162	1991755	57.099	ng	98
48) 2-Nitroaniline	9.486	65	731823	79.656	ng	97
49) Acenaphthylene	9.810	152	2930765	55.294	ng	98
50) Dimethylphthalate	9.669	163	2305746	58.660	ng	99
51) 2,6-Dinitrotoluene	9.733	165	515698	64.205	ng	97
52) Acenaphthene	9.980	154	1936455m	56.088	ng	
53) 3-Nitroaniline	9.904	138	626460	65.404	ng	# 96
54) 2,4-Dinitrophenol	10.004	184	311629	86.619	ng	97
55) Dibenzofuran	10.151	168	2677649	53.262	ng	93
56) 4-Nitrophenol	10.057	139	501409	64.069	ng	94
57) 2,4-Dinitrotoluene	10.133	165	679932	68.570	ng	93
58) Fluorene	10.498	166	1983018	50.994	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.269	232	533841m	50.360	ng	
60) Diethylphthalate	10.363	149	2197083	55.783	ng	99
61) 4-Chlorophenyl-phenyle...	10.486	204	900161	46.459	ng	90
62) 4-Nitroaniline	10.528	138	611477	64.379	ng	93
63) Azobenzene	10.645	77	2389432	61.920	ng	96
65) 4,6-Dinitro-2-methylph...	10.545	198	405952	89.316	ng	94
66) n-Nitrosodiphenylamine	10.610	169	1826683	58.545	ng	95
67) 4-Bromophenyl-phenylether	10.975	248	596660	54.598	ng	90
68) Hexachlorobenzene	11.045	284	632253	51.957	ng	97
69) Atrazine	11.133	200	472257	53.789	ng	97
70) Pentachlorophenol	11.233	266	389435	53.730	ng	99
71) Phenanthrene	11.463	178	2794266	56.925	ng	99
72) Anthracene	11.516	178	2780208	57.447	ng	99
73) Carbazole	11.669	167	2839689	58.557	ng	99
74) Di-n-butylphthalate	11.986	149	3134875	61.311	ng	97
75) Fluoranthene	12.651	202	2828442	55.097	ng	96
77) Benzidine	12.769	184	657224	58.204	ng	99
78) Pyrene	12.880	202	2884338	65.618	ng	99
80) Butylbenzylphthalate	13.492	149	1341699	74.926	ng	97
81) Benzo(a)anthracene	14.068	228	2286425	59.050	ng	98
82) 3,3'-Dichlorobenzidine	14.027	252	518357	61.957	ng	# 96
83) Chrysene	14.110	228	2164716	60.221	ng	98
84) Bis(2-ethylhexyl)phtha...	14.045	149	1690368	71.029	ng	99
85) Di-n-octyl phthalate	14.657	149	2556046	72.757	ng	99
87) Indeno(1,2,3-cd)pyrene	17.092	276	2070755	63.502	ng	97
88) Benzo(b)fluoranthene	15.133	252	1931402	60.403	ng	98
89) Benzo(k)fluoranthene	15.163	252	1782207	57.465	ng	98
90) Benzo(a)pyrene	15.504	252	1583207	60.737	ng	# 96
91) Dibenzo(a,h)anthracene	17.104	278	1639614	61.679	ng	# 96
92) Benzo(g,h,i)perylene	17.551	276	1764219	66.150	ng	96

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

