

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102524\
 Data File : BF140033.D
 Acq On : 25 Oct 2024 13:05
 Operator : RC/JU
 Sample : PB164211BSD
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB164211BSD

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 10/25/2024
 Supervised By :mohammad ahmed 10/28/2024

Quant Time: Oct 25 13:46:48 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF101824.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 18 15:07:50 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.887	152	90907	20.000	ng	0.00
21) Naphthalene-d8	8.169	136	349292	20.000	ng	0.00
39) Acenaphthene-d10	9.928	164	193554	20.000	ng	0.00
64) Phenanthrene-d10	11.416	188	342208	20.000	ng	0.00
76) Chrysene-d12	14.051	240	182421	20.000	ng	0.00
86) Perylene-d12	15.527	264	173566	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.516	112	817434	140.768	ng	0.02
7) Phenol-d6	6.516	99	1035519	137.685	ng	0.00
23) Nitrobenzene-d5	7.451	82	682878	108.355	ng	0.00
42) 2,4,6-Tribromophenol	10.722	330	302175	166.907	ng	0.00
45) 2-Fluorobiphenyl	9.245	172	1210570	103.367	ng	0.00
79) Terphenyl-d14	12.998	244	1275444	113.929	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.769	88	108647	40.129	ng	97
3) Pyridine	3.522	79	291337	43.412	ng	98
4) n-Nitrosodimethylamine	3.475	42	176609	48.981	ng	94
6) Aniline	6.551	93	304089	43.383	ng	100
8) 2-Chlorophenol	6.675	128	315172	53.091	ng	97
9) Benzaldehyde	6.440	77	75497	17.844	ng	98
10) Phenol	6.528	94	387548m	49.982	ng	
11) bis(2-Chloroethyl)ether	6.622	93	300215	50.094	ng	100
12) 1,3-Dichlorobenzene	6.834	146	339253	49.784	ng	99
13) 1,4-Dichlorobenzene	6.910	146	344093	50.541	ng	100
14) 1,2-Dichlorobenzene	7.057	146	331495	52.170	ng	98
15) Benzyl Alcohol	7.028	79	292437	53.008	ng	98
16) 2,2'-oxybis(1-Chloropr...	7.163	45	472633	46.251	ng	97
17) 2-Methylphenol	7.134	107	257565	50.882	ng	97
18) Hexachloroethane	7.404	117	123876	51.024	ng	95
19) n-Nitroso-di-n-propyla...	7.298	70	221948	48.658	ng	98
20) 3+4-Methylphenols	7.287	107	317999	49.114	ng	91
22) Acetophenone	7.298	105	434118	50.743	ng	99
24) Nitrobenzene	7.469	77	344294	49.828	ng	99
25) Isophorone	7.710	82	609300	50.938	ng	99
26) 2-Nitrophenol	7.787	139	161263	61.864	ng	97
27) 2,4-Dimethylphenol	7.816	122	255086	58.687	ng	98
28) bis(2-Chloroethoxy)met...	7.916	93	354871	48.967	ng	100
29) 2,4-Dichlorophenol	8.028	162	252813	51.065	ng	99
30) 1,2,4-Trichlorobenzene	8.110	180	274975	50.194	ng	99
31) Naphthalene	8.192	128	916220	50.860	ng	100
32) Benzoic acid	7.934	122	190989	50.379	ng	99
33) 4-Chloroaniline	8.239	127	135526	22.063	ng	98
34) Hexachlorobutadiene	8.310	225	183797	53.398	ng	98
35) Caprolactam	8.610	113	81804m	51.965	ng	
36) 4-Chloro-3-methylphenol	8.716	107	283139	51.649	ng	99
37) 2-Methylnaphthalene	8.887	142	576248	52.231	ng	99
38) 1-Methylnaphthalene	8.987	142	530285	48.990	ng	99
40) 1,2,4,5-Tetrachloroben...	9.051	216	272619	52.208	ng	99
41) Hexachlorocyclopentadiene	9.039	237	363352	199.981	ng	99
43) 2,4,6-Trichlorophenol	9.157	196	198341	53.650	ng	99

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44) 2,4,5-Trichlorophenol	9.198	196	198905	52.148	ng	99
46) 1,1'-Biphenyl	9.351	154	681920	50.609	ng	98
47) 2-Chloronaphthalene	9.375	162	544289	50.154	ng	99
48) 2-Nitroaniline	9.469	65	189567	57.114	ng	94
49) Acenaphthylene	9.792	152	854895	54.489	ng	99
50) Dimethylphthalate	9.651	163	628143	52.102	ng	100
51) 2,6-Dinitrotoluene	9.710	165	139423	53.411	ng	98
52) Acenaphthene	9.963	154	605657	59.675	ng	99
53) 3-Nitroaniline	9.875	138	93535	34.746	ng	99
54) 2,4-Dinitrophenol	9.986	184	157105	139.564	ng	# 1
55) Dibenzofuran	10.134	168	757662	52.031	ng	99
56) 4-Nitrophenol	10.034	139	236405	114.147	ng	96
57) 2,4-Dinitrotoluene	10.116	165	193572	60.301	ng	97
58) Fluorene	10.481	166	578953	52.200	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.251	232	166114	55.554	ng	100
60) Diethylphthalate	10.351	149	601279	50.795	ng	99
61) 4-Chlorophenyl-phenyle...	10.469	204	289782	51.737	ng	100
62) 4-Nitroaniline	10.492	138	139851	55.007	ng	95
63) Azobenzene	10.628	77	618361	49.274	ng	99
65) 4,6-Dinitro-2-methylph...	10.522	198	104067	74.554	ng	92
66) n-Nitrosodiphenylamine	10.586	169	535013	52.236	ng	99
67) 4-Bromophenyl-phenylether	10.957	248	185841	52.715	ng	99
68) Hexachlorobenzene	11.022	284	207805	52.411	ng	99
69) Atrazine	11.110	200	173450	62.517	ng	99
70) Pentachlorophenol	11.216	266	261617	108.721	ng	99
71) Phenanthrene	11.439	178	863098	53.395	ng	100
72) Anthracene	11.492	178	875234	55.495	ng	100
73) Carbazole	11.645	167	768644	52.404	ng	99
74) Di-n-butylphthalate	11.975	149	870114	51.346	ng	99
75) Fluoranthene	12.627	202	878759	53.776	ng	99
77) Benzidine	12.745	184	155302	51.774	ng	99
78) Pyrene	12.857	202	895412	56.076	ng	100
80) Butylbenzylphthalate	13.474	149	286198	59.784	ng	98
81) Benzo(a)anthracene	14.045	228	661522	55.707	ng	100
82) 3,3'-Dichlorobenzidine	14.004	252	136010	39.283	ng	98
83) Chrysene	14.080	228	588384	54.040	ng	99
84) Bis(2-ethylhexyl)phtha...	14.027	149	316231	58.877	ng	99
85) Di-n-octyl phthalate	14.645	149	455924	46.669	ng	97
87) Indeno(1,2,3-cd)pyrene	17.033	276	652789	58.462	ng	97
88) Benzo(b)fluoranthene	15.098	252	530580	50.183	ng	99
89) Benzo(k)fluoranthene	15.127	252	525217	57.601	ng	100
90) Benzo(a)pyrene	15.468	252	514414	59.130	ng	99
91) Dibenzo(a,h)anthracene	17.051	278	536640	57.611	ng	97
92) Benzo(g,h,i)perylene	17.486	276	500426	53.784	ng	97

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

