

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF110124\
 Data File : BF140203.D
 Acq On : 01 Nov 2024 10:41
 Operator : RC/JU
 Sample : PB164531BS
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB164531BS

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 11/04/2024
 Supervised By :mohammad ahmed 11/04/2024

Quant Time: Nov 01 11:48:10 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	6.886	152	129622	20.000	ng	0.00
21) Naphthalene-d8	8.169	136	521877	20.000	ng	0.00
39) Acenaphthene-d10	9.927	164	295286	20.000	ng	0.00
64) Phenanthrene-d10	11.421	188	550886	20.000	ng	0.00
76) Chrysene-d12	14.080	240	342085	20.000	ng	0.03
86) Perylene-d12	15.580	264	274353	20.000	ng	0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.516	112	1009520	121.923	ng	0.02
7) Phenol-d6	6.516	99	1276406	119.024	ng	0.00
23) Nitrobenzene-d5	7.451	82	879938	93.450	ng	0.00
42) 2,4,6-Tribromophenol	10.727	330	394547	142.848	ng	0.01
45) 2-Fluorobiphenyl	9.245	172	1553470	86.947	ng	0.00
79) Terphenyl-d14	13.015	244	1849152	88.082	ng	0.02
Target Compounds						Qvalue
2) 1,4-Dioxane	2.775	88	125492	32.507	ng	99
3) Pyridine	3.522	79	330816	34.571	ng	98
4) n-Nitrosodimethylamine	3.481	42	219658	42.725	ng	# 91
6) Aniline	6.545	93	472772	47.303	ng	92
8) 2-Chlorophenol	6.669	128	410734	48.523	ng	99
9) Benzaldehyde	6.434	77	74764	12.393	ng	100
10) Phenol	6.528	94	527398m	47.703	ng	
11) bis(2-Chloroethyl)ether	6.622	93	408546	47.809	ng	99
12) 1,3-Dichlorobenzene	6.828	146	435029	44.771	ng	99
13) 1,4-Dichlorobenzene	6.904	146	437670	45.085	ng	99
14) 1,2-Dichlorobenzene	7.057	146	417872	46.122	ng	99
15) Benzyl Alcohol	7.022	79	391305	49.744	ng	97
16) 2,2'-oxybis(1-Chloropr...	7.157	45	710084	48.733	ng	99
17) 2-Methylphenol	7.133	107	347643	48.165	ng	98
18) Hexachloroethane	7.398	117	162590	46.968	ng	97
19) n-Nitroso-di-n-propyla...	7.298	70	307367	47.258	ng	99
20) 3+4-Methylphenols	7.286	107	419923	45.485	ng	# 83
22) Acetophenone	7.292	105	565334	44.227	ng	95
24) Nitrobenzene	7.469	77	469489	45.476	ng	99
25) Isophorone	7.704	82	849700	47.544	ng	99
26) 2-Nitrophenol	7.781	139	217132	55.751	ng	99
27) 2,4-Dimethylphenol	7.816	122	344208	53.002	ng	98
28) bis(2-Chloroethoxy)met...	7.910	93	497630	45.958	ng	99
29) 2,4-Dichlorophenol	8.022	162	343723	46.467	ng	99
30) 1,2,4-Trichlorobenzene	8.104	180	358513	43.801	ng	98
31) Naphthalene	8.186	128	1192622	44.310	ng	100
32) Benzoic acid	7.939	122	263842	46.581	ng	99
33) 4-Chloroaniline	8.233	127	300591	32.752	ng	100
34) Hexachlorobutadiene	8.304	225	232640	45.237	ng	98
35) Caprolactam	8.616	113	107742m	45.808	ng	
36) 4-Chloro-3-methylphenol	8.716	107	396151	48.366	ng	99
37) 2-Methylnaphthalene	8.880	142	755708	45.845	ng	99
38) 1-Methylnaphthalene	8.980	142	708867	43.831	ng	99
40) 1,2,4,5-Tetrachloroben...	9.045	216	363492	45.628	ng	99
41) Hexachlorocyclopentadiene	9.033	237	455655	164.383	ng	99
43) 2,4,6-Trichlorophenol	9.157	196	274874	48.736	ng	99

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44) 2,4,5-Trichlorophenol	9.204	196	271693	46.690	ng	98
46) 1,1'-Biphenyl	9.345	154	928665	45.177	ng	99
47) 2-Chloronaphthalene	9.375	162	732952	44.270	ng	99
48) 2-Nitroaniline	9.469	65	280152	55.326	ng	98
49) Acenaphthylene	9.792	152	1164380	48.647	ng	100
50) Dimethylphthalate	9.651	163	897277	48.784	ng	100
51) 2,6-Dinitrotoluene	9.716	165	197741	49.654	ng	99
52) Acenaphthene	9.963	154	835850	53.983	ng	99
53) 3-Nitroaniline	9.880	138	164214	39.986	ng	99
54) 2,4-Dinitrophenol	9.992	184	242202	140.964	ng	# 1
55) Dibenzofuran	10.139	168	1024518	46.117	ng	99
56) 4-Nitrophenol	10.045	139	335799	106.279	ng	93
57) 2,4-Dinitrotoluene	10.122	165	273892	55.927	ng	96
58) Fluorene	10.480	166	798373	47.184	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.257	232	230068	50.435	ng	99
60) Diethylphthalate	10.357	149	880306	48.746	ng	99
61) 4-Chlorophenyl-phenyle...	10.469	204	404336	47.319	ng	98
62) 4-Nitroaniline	10.504	138	212570	54.804	ng	96
63) Azobenzene	10.633	77	911991	47.635	ng	99
65) 4,6-Dinitro-2-methylph...	10.533	198	159299	70.893	ng	96
66) n-Nitrosodiphenylamine	10.592	169	741018	44.943	ng	99
67) 4-Bromophenyl-phenylether	10.963	248	262959	46.335	ng	98
68) Hexachlorobenzene	11.033	284	285481	44.728	ng	98
69) Atrazine	11.121	200	254138	56.901	ng	98
70) Pentachlorophenol	11.227	266	357120	92.191	ng	99
71) Phenanthrene	11.451	178	1186870	45.611	ng	100
72) Anthracene	11.504	178	1209649	47.645	ng	100
73) Carbazole	11.657	167	1084663	45.937	ng	99
74) Di-n-butylphthalate	11.986	149	1347316	49.389	ng	100
75) Fluoranthene	12.639	202	1279951	48.657	ng	99
77) Benzidine	12.763	184	387457	68.881	ng	99
78) Pyrene	12.868	202	1303458	43.530	ng	100
80) Butylbenzylphthalate	13.492	149	537469	59.870	ng	100
81) Benzo(a)anthracene	14.068	228	1068499	47.982	ng	100
82) 3,3'-Dichlorobenzidine	14.027	252	278367	42.873	ng	99
83) Chrysene	14.104	228	984160	48.201	ng	100
84) Bis(2-ethylhexyl)phtha...	14.051	149	701582	69.657	ng	100
85) Di-n-octyl phthalate	14.680	149	1105441	60.342	ng	99
87) Indeno(1,2,3-cd)pyrene	17.109	276	828153	46.921	ng	98
88) Benzo(b)fluoranthene	15.139	252	810103	48.473	ng	99
89) Benzo(k)fluoranthene	15.168	252	790884	54.873	ng	100
90) Benzo(a)pyrene	15.515	252	716832	52.128	ng	99
91) Dibenzo(a,h)anthracene	17.127	278	677815	46.035	ng	99
92) Benzo(g,h,i)perylene	17.574	276	624704	42.476	ng	99

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

