

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF100724\
 Data File : BF139685.D
 Acq On : 07 Oct 2024 14:35
 Operator : RC/JU
 Sample : PB163902BS
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB163902BS

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 10/09/2024
 Supervised By :mohammad ahmed 10/09/2024

Quant Time: Oct 07 14:59:01 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF100124.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 02 04:58:47 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.869	152	100095	20.000	ng	0.00
21) Naphthalene-d8	8.151	136	391689	20.000	ng	0.00
39) Acenaphthene-d10	9.910	164	230281	20.000	ng	0.00
64) Phenanthrene-d10	11.392	188	448928	20.000	ng	0.00
76) Chrysene-d12	14.039	240	273215	20.000	ng	0.00
86) Perylene-d12	15.504	264	208036	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.499	112	553335	95.816	ng	0.01
7) Phenol-d6	6.504	99	734302	94.552	ng	0.00
23) Nitrobenzene-d5	7.434	82	684647	88.498	ng	0.00
42) 2,4,6-Tribromophenol	10.698	330	207138	93.180	ng	0.00
45) 2-Fluorobiphenyl	9.228	172	1256347	83.755	ng	0.00
79) Terphenyl-d14	12.980	244	1509747	90.670	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.746	88	105750	42.355	ng	99
3) Pyridine	3.481	79	277261	45.660	ng	99
4) n-Nitrosodimethylamine	3.452	42	189741	47.737	ng	96
6) Aniline	6.534	93	356632	51.750	ng	99
8) 2-Chlorophenol	6.657	128	316635	51.175	ng	98
9) Benzaldehyde	6.416	77	100188	24.353	ng	99
10) Phenol	6.522	94	416124	50.957	ng	89
11) bis(2-Chloroethyl)ether	6.604	93	313128	47.094	ng	100
12) 1,3-Dichlorobenzene	6.810	146	321310	46.163	ng	99
13) 1,4-Dichlorobenzene	6.887	146	322832	45.795	ng	100
14) 1,2-Dichlorobenzene	7.040	146	308525	46.697	ng	100
15) Benzyl Alcohol	7.016	79	297384	50.117	ng	97
16) 2,2'-oxybis(1-Chloropr...	7.140	45	550758	47.848	ng	79
17) 2-Methylphenol	7.128	107	259212	50.183	ng	99
18) Hexachloroethane	7.381	117	122094	46.433	ng	99
19) n-Nitroso-di-n-propyla...	7.287	70	239296	47.927	ng	98
20) 3+4-Methylphenols	7.281	107	321888	49.614	ng	95
22) Acetophenone	7.281	105	431574	46.418	ng	98
24) Nitrobenzene	7.457	77	366843	45.793	ng	99
25) Isophorone	7.693	82	653687	47.574	ng	100
26) 2-Nitrophenol	7.769	139	171334	49.751	ng	99
27) 2,4-Dimethylphenol	7.804	122	248881m	59.033	ng	
28) bis(2-Chloroethoxy)met...	7.898	93	384475	46.859	ng	98
29) 2,4-Dichlorophenol	8.016	162	269825	49.067	ng	98
30) 1,2,4-Trichlorobenzene	8.092	180	272456	43.811	ng	100
31) Naphthalene	8.175	128	925954	46.236	ng	100
32) Benzoic acid	7.945	122	204268	48.726	ng	99
33) 4-Chloroaniline	8.222	127	117413	17.320	ng	99
34) Hexachlorobutadiene	8.287	225	175591	43.630	ng	99
35) Caprolactam	8.604	113	92938m	51.528	ng	
36) 4-Chloro-3-methylphenol	8.716	107	308180	49.507	ng	98
37) 2-Methylnaphthalene	8.863	142	610438	46.802	ng	99
38) 1-Methylnaphthalene	8.963	142	569133	44.640	ng	99
40) 1,2,4,5-Tetrachloroben...	9.034	216	283973	44.408	ng	99
41) Hexachlorocyclopentadiene	9.016	237	329647	168.123	ng	98
43) 2,4,6-Trichlorophenol	9.145	196	203240	47.147	ng	99

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44) 2,4,5-Trichlorophenol	9.192	196	216870	46.651	ng	99
46) 1,1'-Biphenyl	9.328	154	774495	45.219	ng	100
47) 2-Chloronaphthalene	9.357	162	575827	44.846	ng	99
48) 2-Nitroaniline	9.451	65	226385	49.116	ng	98
49) Acenaphthylene	9.769	152	951796	48.743	ng	99
50) Dimethylphthalate	9.634	163	727668	48.272	ng	100
51) 2,6-Dinitrotoluene	9.698	165	160416	47.108	ng	93
52) Acenaphthene	9.939	154	697764m	52.052	ng	
53) 3-Nitroaniline	9.863	138	117596	33.793	ng	97
54) 2,4-Dinitrophenol	9.975	184	199141	108.728	ng	98
55) Dibenzofuran	10.116	168	878524	46.807	ng	100
56) 4-Nitrophenol	10.034	139	262912	99.074	ng	99
57) 2,4-Dinitrotoluene	10.104	165	213859	48.047	ng	95
58) Fluorene	10.457	166	703720	47.074	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.234	232	190728	50.796	ng	99
60) Diethylphthalate	10.328	149	741788	47.642	ng	99
61) 4-Chlorophenyl-phenyle...	10.445	204	341535	45.666	ng	98
62) 4-Nitroaniline	10.486	138	177621	49.901	ng	98
63) Azobenzene	10.610	77	756440	48.309	ng	99
65) 4,6-Dinitro-2-methylph...	10.510	198	131001	51.668	ng	97
66) n-Nitrosodiphenylamine	10.569	169	628509	47.496	ng	99
67) 4-Bromophenyl-phenylether	10.939	248	206931	44.957	ng	99
68) Hexachlorobenzene	11.004	284	230531	45.135	ng	99
69) Atrazine	11.098	200	204756	57.671	ng	98
70) Pentachlorophenol	11.204	266	271923	89.410	ng	99
71) Phenanthrene	11.422	178	1011818	47.801	ng	100
72) Anthracene	11.475	178	1044748	49.890	ng	100
73) Carbazole	11.628	167	971439	48.300	ng	100
74) Di-n-butylphthalate	11.951	149	1191401	48.713	ng	99
75) Fluoranthene	12.610	202	1121318	48.460	ng	100
77) Benzidine	12.727	184	391839	81.308	ng	99
78) Pyrene	12.839	202	1145049	46.859	ng	99
80) Butylbenzylphthalate	13.451	149	460085	52.638	ng	99
81) Benzo(a)anthracene	14.027	228	891569	49.634	ng	100
82) 3,3'-Dichlorobenzidine	13.986	252	193599	35.216	ng	100
83) Chrysene	14.063	228	781572	47.591	ng	99
84) Bis(2-ethylhexyl)phtha...	14.010	149	603106	55.258	ng	99
85) Di-n-octyl phthalate	14.621	149	859040	53.575	ng	99
87) Indeno(1,2,3-cd)pyrene	16.986	276	619756	50.615	ng	100
88) Benzo(b)fluoranthene	15.074	252	672649	50.006	ng	99
89) Benzo(k)fluoranthene	15.104	252	559417	49.027	ng	99
90) Benzo(a)pyrene	15.445	252	558749	52.681	ng	99
91) Dibenzo(a,h)anthracene	17.004	278	507642	49.986	ng	99
92) Benzo(g,h,i)perylene	17.433	276	472795	46.420	ng	100

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

