

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF050421\
 Data File : BF123807.D
 Acq On : 4 May 2021 14:23
 Operator : JU/CG
 Sample : PB136158BS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB136158BS

Manual Integrations
 APPROVED

mohammad
 5/5/2021 11:25:04 AM

Quant Time: May 04 14:57:38 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF042721.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 04 12:58:27 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.804	152	279829	20.00	ng	0.00
21) Naphthalene-d8	8.086	136	1070301	20.00	ng	0.00
39) Acenaphthene-d10	9.845	164	543826	20.00	ng	0.00
64) Phenanthrene-d10	11.333	188	1059607	20.00	ng	0.00
76) Chrysene-d12	13.980	240	756998	20.00	ng	0.00
86) Perylene-d12	15.427	264	588035	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.440	112	2138487	121.07	ng	0.02
7) Phenol-d6	6.451	99	2742394	112.71	ng	0.00
23) Nitrobenzene-d5	7.369	82	1995494	84.16	ng	0.00
42) 2,4,6-Tribromophenol	10.639	330	837823	123.75	ng	0.00
45) 2-Fluorobiphenyl	9.169	172	3027855	81.08	ng	0.00
79) Terphenyl-d14	12.921	244	3433539	87.34	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.616	88	338757	35.60	ng	96
3) Pyridine	3.351	79	832666	35.81	ng	91
4) n-Nitrosodimethylamine	3.304	42	567595	44.31	ng	97
6) Aniline	6.463	93	835043	29.67	ng	# 1
8) 2-Chlorophenol	6.592	128	862772	45.41	ng	99
9) Benzaldehyde	6.351	77	727913	61.81	ng	99
10) Phenol	6.463	94	1091654	41.77	ng	80
11) bis(2-Chloroethyl)ether	6.539	93	863349	45.24	ng	99
12) 1,3-Dichlorobenzene	6.745	146	830815	43.28	ng	96
13) 1,4-Dichlorobenzene	6.822	146	842029	43.35	ng	95
14) 1,2-Dichlorobenzene	6.975	146	808905	42.89	ng	98
15) Benzyl Alcohol	6.951	79	846211	44.37	ng	98
16) 2,2'-oxybis(1-Chloropr...	7.081	45	1843896	45.43	ng	96
17) 2-Methylphenol	7.069	107	712960	44.03	ng	95
18) Hexachloroethane	7.316	117	353526	45.98	ng	97
19) n-Nitroso-di-n-propyla...	7.222	70	666596	44.37	ng	100
20) 3+4-Methylphenols	7.222	107	875607	42.50	ng	93
22) Acetophenone	7.216	105	1283861	44.13	ng	94
24) Nitrobenzene	7.386	77	1018606	41.24	ng	95
25) Isophorone	7.628	82	1841715	41.43	ng	99
26) 2-Nitrophenol	7.704	139	444943	44.15	ng	99
27) 2,4-Dimethylphenol	7.745	122	760134	48.01	ng	98
28) bis(2-Chloroethoxy)met...	7.839	93	1078609	47.61	ng	99
29) 2,4-Dichlorophenol	7.951	162	669942	42.94	ng	97
30) 1,2,4-Trichlorobenzene	8.028	180	705929	43.17	ng	98
31) Naphthalene	8.110	128	2336720	42.40	ng	99
32) Benzoic acid	7.892	122	449033	42.05	ng	96
33) 4-Chloroaniline	8.157	127	254041	11.05	ng	96
34) Hexachlorobutadiene	8.228	225	424480	42.59	ng	99
35) Caprolactam	8.539	113	232935m	42.52	ng	
36) 4-Chloro-3-methylphenol	8.651	107	877415	45.07	ng	96
37) 2-Methylnaphthalene	8.804	142	1552943	43.24	ng	98
38) 1-Methylnaphthalene	8.904	142	1436474	42.55	ng	99
40) 1,2,4,5-Tetrachloroben...	8.969	216	670343	41.88	ng	99
41) Hexachlorocyclopentadiene	8.957	237	793840	113.40	ng	99
43) 2,4,6-Trichlorophenol	9.086	196	477348	43.30	ng	98

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44) 2,4,5-Trichlorophenol	9.128	196	501912	42.44	ng	99
46) 1,1'-Biphenyl	9.269	154	1864505	41.85	ng	99
47) 2-Chloronaphthalene	9.292	162	1400267	41.68	ng	98
48) 2-Nitroaniline	9.392	65	640331	43.27	ng	97
49) Acenaphthylene	9.710	152	2311648	42.63	ng	99
50) Dimethylphthalate	9.575	163	1674510	43.72	ng	99
51) 2,6-Dinitrotoluene	9.633	165	377352	42.61	ng	98
52) Acenaphthene	9.880	154	1376542	41.67	ng	99
53) 3-Nitroaniline	9.798	138	200597	18.97	ng	97
54) 2,4-Dinitrophenol	9.910	184	426115	90.60	ng	96
55) Dibenzofuran	10.051	168	2032687	40.68	ng	100
56) 4-Nitrophenol	9.980	139	643372	83.52	ng	93
57) 2,4-Dinitrotoluene	10.039	165	510018	42.24	ng	# 84
58) Fluorene	10.398	166	1612143	41.99	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.175	232	417477	44.58	ng	99
60) Diethylphthalate	10.275	149	1816320	44.55	ng	98
61) 4-Chlorophenyl-phenyle...	10.386	204	764434	42.72	ng	98
62) 4-Nitroaniline	10.422	138	429959	41.02	ng	96
63) Azobenzene	10.545	77	1963095	41.89	ng	97
65) 4,6-Dinitro-2-methylph...	10.451	198	268118	40.84	ng	88
66) n-Nitrosodiphenylamine	10.510	169	1436563	41.45	ng	99
67) 4-Bromophenyl-phenylether	10.874	248	496709	44.51	ng	94
68) Hexachlorobenzene	10.945	284	559193	44.03	ng	97
69) Atrazine	11.039	200	483129	46.11	ng	98
70) Pentachlorophenol	11.145	266	586526	90.61	ng	99
71) Phenanthrene	11.357	178	2311745	41.59	ng	99
72) Anthracene	11.410	178	2399221	43.42	ng	99
73) Carbazole	11.563	167	2268377	40.61	ng	99
74) Di-n-butylphthalate	11.898	149	2856616	45.56	ng	99
75) Fluoranthene	12.545	202	2538609	41.88	ng	99
77) Benzidine	12.668	184	1401284	71.48	ng	99
78) Pyrene	12.774	202	2572579	43.78	ng	99
80) Butylbenzylphthalate	13.398	149	1216572	48.46	ng	96
81) Benzo(a)anthracene	13.968	228	1957483	42.67	ng	99
82) 3,3'-Dichlorobenzidine	13.927	252	246763	17.52	ng	98
83) Chrysene	14.004	228	1945327	42.15	ng	100
84) Bis(2-ethylhexyl)phtha...	13.957	149	1553704	49.82	ng	99
85) Di-n-octyl phthalate	14.568	149	2617828	52.50	ng	99
87) Indeno(1,2,3-cd)pyrene	16.880	276	1590216	46.43	ng	99
88) Benzo(b)fluoranthene	15.009	252	1699506m	42.89	ng	
89) Benzo(k)fluoranthene	15.039	252	1659089	43.87	ng	99
90) Benzo(a)pyrene	15.374	252	1542240	45.51	ng	96
91) Dibenzo(a,h)anthracene	16.892	278	1307288	45.21	ng	99
92) Benzo(g,h,i)perylene	17.315	276	1330736	45.95	ng	100

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

