

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF092821\
 Data File : BF125688.D
 Acq On : 28 Sep 2021 14:30
 Operator : JU/CG
 Sample : PB139432BSD
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB139432BSD

Manual Integrations
 APPROVED

mohammad
 9/29/2021 12:55:01 PM

Quant Time: Sep 28 15:16:57 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF092421.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 24 15:38:32 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.892	152	339868	20.000	ng	0.00
21) Naphthalene-d8	8.175	136	1452185	20.000	ng	# 0.00
39) Acenaphthene-d10	9.927	164	775508	20.000	ng	0.00
64) Phenanthrene-d10	11.410	188	1184530	20.000	ng	# 0.00
76) Chrysene-d12	14.051	240	620586	20.000	ng	0.00
86) Perylene-d12	15.521	264	696569	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.522	112	2432537	117.921	ng	0.02
7) Phenol-d6	6.522	99	3064251	110.490	ng	0.01
23) Nitrobenzene-d5	7.457	82	1868650	89.754	ng	0.00
42) 2,4,6-Tribromophenol	10.721	330	825210	114.128	ng	0.00
45) 2-Fluorobiphenyl	9.251	172	3255271	73.635	ng	0.00
79) Terphenyl-d14	12.998	244	2579489	71.981	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.763	88	306486	34.911	ng	# 100
3) Pyridine	3.510	79	719402	30.981	ng	# 71
4) n-Nitrosodimethylamine	3.469	42	373092	43.302	ng	# 1
6) Aniline	6.551	93	1333877	41.655	ng	94
8) 2-Chlorophenol	6.675	128	1061831	45.378	ng	98
9) Benzaldehyde	6.439	77	575918	47.092	ng	94
10) Phenol	6.534	94	1186325m	42.317	ng	
11) bis(2-Chloroethyl)ether	6.628	93	963321	44.217	ng	95
12) 1,3-Dichlorobenzene	6.833	146	1042722	40.356	ng	95
13) 1,4-Dichlorobenzene	6.910	146	1063376	40.452	ng	98
14) 1,2-Dichlorobenzene	7.063	146	1039099	41.896	ng	99
15) Benzyl Alcohol	7.033	79	804755	41.834	ng	92
16) 2,2'-oxybis(1-Chloropr...	7.163	45	1194329	43.068	ng	87
17) 2-Methylphenol	7.139	107	862848	44.967	ng	96
18) Hexachloroethane	7.404	117	380520	44.596	ng	93
19) n-Nitroso-di-n-propyla...	7.310	70	643367	40.744	ng	# 67
20) 3+4-Methylphenols	7.292	107	999642	41.438	ng	94
22) Acetophenone	7.304	105	1338115	40.608	ng	# 89
24) Nitrobenzene	7.475	77	988299	46.666	ng	98
25) Isophorone	7.716	82	1847075	39.614	ng	94
26) 2-Nitrophenol	7.786	139	550100	49.330	ng	# 92
27) 2,4-Dimethylphenol	7.822	122	975373	47.037	ng	95
28) bis(2-Chloroethoxy)met...	7.922	93	1185919	45.093	ng	98
29) 2,4-Dichlorophenol	8.028	162	882148	42.826	ng	98
30) 1,2,4-Trichlorobenzene	8.116	180	892863	39.818	ng	96
31) Naphthalene	8.198	128	2795321	37.489	ng	99
32) Benzoic acid	7.957	122	667483	44.656	ng	97
33) 4-Chloroaniline	8.239	127	934416	29.876	ng	98
34) Hexachlorobutadiene	8.316	225	485433	39.262	ng	99
35) Caprolactam	8.616	113	278419m	41.419	ng	
36) 4-Chloro-3-methylphenol	8.722	107	885036	41.003	ng	98
37) 2-Methylnaphthalene	8.886	142	1930939	38.183	ng	96
38) 1-Methylnaphthalene	8.986	142	1798983	37.641	ng	100
40) 1,2,4,5-Tetrachloroben...	9.051	216	788488	39.094	ng	98
41) Hexachlorocyclopentadiene	9.045	237	844300	98.938	ng	97
43) 2,4,6-Trichlorophenol	9.163	196	608685	43.036	ng	97

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.198	196	652364	42.721	ng	93
46) 1,1'-Biphenyl	9.351	154	2145144	37.513	ng	94
47) 2-Chloronaphthalene	9.374	162	1829484	38.642	ng	99
48) 2-Nitroaniline	9.469	65	506531	45.588	ng	89
49) Acenaphthylene	9.792	152	2645465	37.118	ng	94
50) Dimethylphthalate	9.651	163	2048864	39.217	ng	97
51) 2,6-Dinitrotoluene	9.710	165	480263	45.367	ng	93
52) Acenaphthene	9.963	154	1819203	40.711	ng	96
53) 3-Nitroaniline	9.880	138	427989	36.070	ng	# 97
54) 2,4-Dinitrophenol	9.986	184	400565	80.423	ng	# 72
55) Dibenzofuran	10.133	168	2427597	36.210	ng	98
56) 4-Nitrophenol	10.039	139	763217	79.306	ng	# 86
57) 2,4-Dinitrotoluene	10.116	165	615080	46.502	ng	# 80
58) Fluorene	10.480	166	1804057	34.572	ng	97
59) 2,3,4,6-Tetrachlorophenol	10.251	232	487686	41.496	ng	# 89
60) Diethylphthalate	10.351	149	1935310	38.424	ng	98
61) 4-Chlorophenyl-phenyle...	10.469	204	838783	36.162	ng	88
62) 4-Nitroaniline	10.498	138	477767	40.449	ng	# 73
63) Azobenzene	10.627	77	1807068	37.158	ng	86
65) 4,6-Dinitro-2-methylph...	10.522	198	264662	46.298	ng	# 75
66) n-Nitrosodiphenylamine	10.586	169	1692652	41.626	ng	94
67) 4-Bromophenyl-phenylether	10.957	248	527747	42.453	ng	97
68) Hexachlorobenzene	11.027	284	617984	42.357	ng	# 82
69) Atrazine	11.110	200	487767	45.239	ng	95
70) Pentachlorophenol	11.216	266	638984	82.866	ng	96
71) Phenanthrene	11.439	178	2498905	37.620	ng	94
72) Anthracene	11.492	178	2526633	38.494	ng	95
73) Carbazole	11.639	167	2280746	35.308	ng	98
74) Di-n-butylphthalate	11.968	149	2466641	37.877	ng	99
75) Fluoranthene	12.621	202	2226253	31.360	ng	94
77) Benzidine	12.739	184	1031811	63.290	ng	98
78) Pyrene	12.851	202	2208125	39.372	ng	95
80) Butylbenzylphthalate	13.468	149	792920	42.693	ng	98
81) Benzo(a)anthracene	14.039	228	1658128	39.334	ng	99
82) 3,3'-Dichlorobenzidine	13.998	252	599574	49.292	ng	99
83) Chrysene	14.074	228	1672730	39.540	ng	96
84) Bis(2-ethylhexyl)phtha...	14.027	149	968078	45.533	ng	98
85) Di-n-octyl phthalate	14.639	149	1846454	59.213	ng	# 89
87) Indeno(1,2,3-cd)pyrene	17.015	276	2184248	46.195	ng	97
88) Benzo(b)fluoranthene	15.092	252	1997655	40.933	ng	98
89) Benzo(k)fluoranthene	15.127	252	1735725	37.121	ng	99
90) Benzo(a)pyrene	15.462	252	1888499	43.472	ng	97
91) Dibenzo(a,h)anthracene	17.039	278	1839860	46.554	ng	98
92) Benzo(g,h,i)perylene	17.474	276	1868001	46.281	ng	91

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

