

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM052423\
 Data File : BM039979.D
 Acq On : 24 May 2023 11:58
 Operator : CG/JU
 Sample : PB152902BSD
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB152902BSD

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 05/25/2023
 Supervised By : Jagrut Upadhyay 05/25/2023

Quant Time: May 24 13:18:04 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM052323.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 24 01:06:13 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.016	152	254454	20.000	ng	0.00
21) Naphthalene-d8	10.833	136	985060	20.000	ng	0.00
39) Acenaphthene-d10	14.651	164	470358	20.000	ng	0.00
64) Phenanthrene-d10	17.398	188	816635	20.000	ng	0.00
76) Chrysene-d12	21.574	240	675316	20.000	ng	0.00
86) Perylene-d12	24.027	264	746680	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.563	112	1623722	97.884	ng	0.00
7) Phenol-d6	7.169	99	1874015	86.603	ng	0.00
23) Nitrobenzene-d5	9.186	82	1652350	90.513	ng	0.00
42) 2,4,6-Tribromophenol	16.139	330	534440	87.543	ng	0.00
45) 2-Fluorobiphenyl	13.286	172	3721842	93.330	ng	0.00
79) Terphenyl-d14	20.015	244	4482695	103.067	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.422	88	288593	43.079	ng	98
3) Pyridine	3.828	79	749421	40.746	ng	97
4) n-Nitrosodimethylamine	3.740	42	356431	45.615	ng	95
6) Aniline	7.339	93	812942	30.362	ng	99
8) 2-Chlorophenol	7.575	128	847532	48.538	ng	97
9) Benzaldehyde	7.145	77	499506	47.751	ng	99
10) Phenol	7.198	94	1038107	47.860	ng	99
11) bis(2-Chloroethyl)ether	7.434	93	828896	46.396	ng	99
12) 1,3-Dichlorobenzene	7.904	146	947751	51.256	ng	99
13) 1,4-Dichlorobenzene	8.051	146	969061	51.509	ng	99
14) 1,2-Dichlorobenzene	8.375	146	919645	49.851	ng	99
15) Benzyl Alcohol	8.257	79	636914	45.661	ng	98
16) 2,2'-oxybis(1-Chloropr...	8.557	45	1034038	45.120	ng	100
17) 2-Methylphenol	8.457	107	666458	42.946	ng	98
18) Hexachloroethane	9.110	117	352099	51.623	ng	97
19) n-Nitroso-di-n-propyla...	8.828	70	561135	44.772	ng	97
20) 3+4-Methylphenols	8.786	107	865266	42.253	ng	98
22) Acetophenone	8.845	105	1253135	47.132	ng	# 98
24) Nitrobenzene	9.228	77	849431	45.872	ng	98
25) Isophorone	9.751	82	1496918	44.964	ng	99
26) 2-Nitrophenol	9.939	139	363280	55.442	ng	96
27) 2,4-Dimethylphenol	9.998	122	723396	48.471	ng	99
28) bis(2-Chloroethoxy)met...	10.239	93	1001757	45.896	ng	99
29) 2,4-Dichlorophenol	10.475	162	673697	47.645	ng	98
30) 1,2,4-Trichlorobenzene	10.692	180	766315	49.503	ng	99
31) Naphthalene	10.886	128	2516124	46.294	ng	100
32) Benzoic acid	10.116	122	365775	38.259	ng	93
33) 4-Chloroaniline	10.986	127	283905	12.396	ng	99
34) Hexachlorobutadiene	11.174	225	420547	48.401	ng	98
35) Caprolactam	11.763	113	176857	44.294	ng	88
36) 4-Chloro-3-methylphenol	12.104	107	659431	45.120	ng	99
37) 2-Methylnaphthalene	12.486	142	1601091	44.195	ng	99
38) 1-Methylnaphthalene	12.704	142	1484881	43.786	ng	99
40) 1,2,4,5-Tetrachloroben...	12.851	216	780691	52.477	ng	99
41) Hexachlorocyclopentadiene	12.839	237	1096877	142.527	ng	99
43) 2,4,6-Trichlorophenol	13.086	196	451136	49.962	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.157	196	490041	45.595	ng	99
46) 1,1'-Biphenyl	13.492	154	2012779	49.983	ng	99
47) 2-Chloronaphthalene	13.533	162	1535549	51.084	ng	100
48) 2-Nitroaniline	13.733	65	357738	44.200	ng	95
49) Acenaphthylene	14.374	152	2285399	48.566	ng	99
50) Dimethylphthalate	14.110	163	1676508	47.282	ng	100
51) 2,6-Dinitrotoluene	14.227	165	337915	49.818	ng	97
52) Acenaphthene	14.715	154	1593761m	51.835	ng	05/25/2023
53) 3-Nitroaniline	14.551	138	169056	21.069	ng	100
54) 2,4-Dinitrophenol	14.757	184	312622	82.534	ng	98
55) Dibenzofuran	15.051	168	2136458	47.070	ng	99
56) 4-Nitrophenol	14.851	139	661891	106.635	ng	94
57) 2,4-Dinitrotoluene	15.010	165	439734	45.257	ng	94
58) Fluorene	15.698	166	1761203	46.677	ng	98
59) 2,3,4,6-Tetrachlorophenol	15.274	232	387731	48.511	ng	100
60) Diethylphthalate	15.468	149	1639964	46.793	ng	99
61) 4-Chlorophenyl-phenyle...	15.692	204	856028	46.372	ng	100
62) 4-Nitroaniline	15.715	138	375985	44.992	ng	96
63) Azobenzene	15.986	77	1627174	46.568	ng	98
65) 4,6-Dinitro-2-methylph...	15.768	198	195040	44.094	ng	96
66) n-Nitrosodiphenylamine	15.904	169	1396956	49.743	ng	99
67) 4-Bromophenyl-phenylether	16.586	248	439458	48.702	ng	98
68) Hexachlorobenzene	16.698	284	530400	51.436	ng	98
69) Atrazine	16.851	200	420269	60.163	ng	99
70) Pentachlorophenol	17.039	266	630206	93.098	ng	99
71) Phenanthrene	17.439	178	2340170	47.908	ng	100
72) Anthracene	17.527	178	2354749	48.655	ng	100
73) Carbazole	17.798	167	2166895	49.558	ng	99
74) Di-n-butylphthalate	18.362	149	2396370	43.260	ng	100
75) Fluoranthene	19.450	202	2266083	46.546	ng	100
77) Benzidine	19.633	184	714690	68.151	ng	99
78) Pyrene	19.809	202	2417369	48.782	ng	100
80) Butylbenzylphthalate	20.709	149	918223	45.355	ng	92
81) Benzo(a)anthracene	21.556	228	2289064	47.882	ng	99
82) 3,3'-Dichlorobenzidine	21.491	252	527034	36.905	ng	100
83) Chrysene	21.615	228	2158095	47.048	ng	100
84) Bis(2-ethylhexyl)phtha...	21.486	149	1437286	44.543	ng	98
85) Di-n-octyl phthalate	22.433	149	2365078	47.512	ng	98
87) Indeno(1,2,3-cd)pyrene	26.591	276	3315672	55.839	ng	98
88) Benzo(b)fluoranthene	23.280	252	2375688	49.695	ng	99
89) Benzo(k)fluoranthene	23.327	252	2444812	46.769	ng	100
90) Benzo(a)pyrene	23.915	252	2082806	45.430	ng	98
91) Dibenzo(a,h)anthracene	26.603	278	2834853	55.299	ng	99
92) Benzo(g,h,i)perylene	27.373	276	2527834	55.792	ng	99

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

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