

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG020323\
 Data File : BG056536.D
 Acq On : 3 Feb 2023 12:58
 Operator : CG/JU
 Sample : PB150626BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB150626BS

Manual Integrations
 APPROVED

Reviewed By :Christian
 Giraldo

Quant Time: Feb 04 05:10:32 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG012723.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 01 01:15:52 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.271	152	30028	20.000	ng	0.00
21) Naphthalene-d8	11.109	136	117455	20.000	ng	# 0.00
39) Acenaphthene-d10	14.898	164	100480	20.000	ng	0.00
64) Phenanthrene-d10	17.630	188	272654	20.000	ng	# 0.00
76) Chrysene-d12	21.925	240	261179	20.000	ng	0.00
86) Perylene-d12	25.345	264	276112	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.791	112	197460	120.968	ng	0.00
7) Phenol-d6	7.419	99	272375	112.618	ng	0.00
23) Nitrobenzene-d5	9.452	82	140949	68.143	ng	0.00
42) 2,4,6-Tribromophenol	16.379	330	140033	104.829	ng	0.00
45) 2-Fluorobiphenyl	13.518	172	476364	65.729	ng	0.00
79) Terphenyl-d14	20.210	244	989947	66.571	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.582	88	25534	37.567	ng	# 93
3) Pyridine	4.005	79	68493	33.805	ng	95
4) n-Nitrosodimethylamine	3.917	42	17767	41.235	ng	# 100
6) Aniline	7.583	93	107842	34.056	ng	95
8) 2-Chlorophenol	7.836	128	86982	48.507	ng	99
9) Benzaldehyde	7.395	77	51733	51.345	ng	94
10) Phenol	7.448	94	123162	50.103	ng	97
11) bis(2-Chloroethyl)ether	7.677	93	72990	43.199	ng	98
12) 1,3-Dichlorobenzene	8.159	146	91836	44.481	ng	94
13) 1,4-Dichlorobenzene	8.306	146	94523	45.207	ng	95
14) 1,2-Dichlorobenzene	8.629	146	91756	45.231	ng	99
15) Benzyl Alcohol	8.506	79	74808	45.081	ng	99
16) 2,2'-oxybis(1-Chloropr...	8.788	45	14984	41.854	ng	90
17) 2-Methylphenol	8.711	107	83689	44.751	ng	95
18) Hexachloroethane	9.369	117	28715	45.406	ng	91
19) n-Nitroso-di-n-propyla...	9.076	70	56395	40.462	ng	96
20) 3+4-Methylphenols	9.040	107	117010	44.647	ng	99
22) Acetophenone	9.099	105	135674	43.474	ng	# 99
24) Nitrobenzene	9.493	77	89064	44.640	ng	94
25) Isophorone	10.010	82	176662	41.400	ng	99
26) 2-Nitrophenol	10.210	139	54003	44.612	ng	97
27) 2,4-Dimethylphenol	10.257	122	96773m	47.768	ng	
28) bis(2-Chloroethoxy)met...	10.486	93	100946	42.822	ng	98
29) 2,4-Dichlorophenol	10.750	162	109541	48.036	ng	96
30) 1,2,4-Trichlorobenzene	10.962	180	115481	45.571	ng	99
31) Naphthalene	11.156	128	268781	43.356	ng	99
32) Benzoic acid	10.409	122	47352m	41.777	ng	
33) 4-Chloroaniline	11.261	127	93521	32.318	ng	99
34) Hexachlorobutadiene	11.432	225	78899	44.356	ng	95
35) Caprolactam	12.031	113	35459m	46.106	ng	
36) 4-Chloro-3-methylphenol	12.366	107	104804	47.590	ng	96
37) 2-Methylnaphthalene	12.742	142	216092	42.281	ng	99
38) 1-Methylnaphthalene	12.959	142	198570	41.601	ng	95
40) 1,2,4,5-Tetrachloroben...	13.106	216	158530	43.186	ng	98
41) Hexachlorocyclopentadiene	13.077	237	155895	80.104	ng	97
43) 2,4,6-Trichlorophenol	13.341	196	106325	44.946	ng	98

02/06/2023
 Supervised By :Jagrut
 padhyay

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.418	196	115293	41.627	ng	99
46) 1,1'-Biphenyl	13.729	154	314271	43.122	ng	98
47) 2-Chloronaphthalene	13.782	162	252215	44.275	ng	98
48) 2-Nitroaniline	13.982	65	52900	44.921	ng	96
49) Acenaphthylene	14.622	152	400612	43.224	ng	99
50) Dimethylphthalate	14.334	163	349169	42.953	ng	99
51) 2,6-Dinitrotoluene	14.463	165	78438	44.246	ng	97
52) Acenaphthene	14.963	154	234532	42.672	ng	98
53) 3-Nitroaniline	14.798	138	56537	32.896	ng	99
54) 2,4-Dinitrophenol	15.010	184	104843	94.335	ng	96
55) Dibenzofuran	15.292	168	426101	43.229	ng	99
56) 4-Nitrophenol	15.104	139	108909	92.740	ng	97
57) 2,4-Dinitrotoluene	15.251	165	113642	45.510	ng	99
58) Fluorene	15.938	166	350870	44.736	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.515	232	116130	47.157	ng	98
60) Diethylphthalate	15.686	149	335856	43.998	ng	99
61) 4-Chlorophenyl-phenyle...	15.915	204	210791	44.045	ng	97
62) 4-Nitroaniline	15.956	138	74110	43.094	ng	97
63) Azobenzene	16.208	77	233817	44.516	ng	96
65) 4,6-Dinitro-2-methylph...	16.015	198	79187	41.310	ng	97
66) n-Nitrosodiphenylamine	16.132	169	324649	42.055	ng	99
67) 4-Bromophenyl-phenylether	16.814	248	148922	42.666	ng	99
68) Hexachlorobenzene	16.943	284	152341	44.562	ng	100
69) Atrazine	17.066	200	143056	48.035	ng	99
70) Pentachlorophenol	17.284	266	155891	75.362	ng	96
71) Phenanthrene	17.677	178	614704	43.076	ng	99
72) Anthracene	17.765	178	635820	43.404	ng	100
73) Carbazole	18.036	167	553834	44.391	ng	99
74) Di-n-butylphthalate	18.559	149	556089	42.450	ng	99
75) Fluoranthene	19.669	202	770278	42.768	ng	99
77) Benzidine	19.840	184	313849	44.405	ng	98
78) Pyrene	20.033	202	784146	42.214	ng	99
80) Butylbenzylphthalate	20.885	149	234901	41.406	ng	99
81) Benzo(a)anthracene	21.902	228	780022	42.719	ng	100
82) 3,3'-Dichlorobenzidine	21.802	252	191627	29.143	ng	100
83) Chrysene	21.978	228	722011	42.089	ng	99
84) Bis(2-ethylhexyl)phtha...	21.761	149	337675	41.516	ng	99
85) Di-n-octyl phthalate	23.036	149	567174	41.873	ng	100
87) Indeno(1,2,3-cd)pyrene	29.299	276	914003	45.221	ng	# 90
88) Benzo(b)fluoranthene	24.258	252	816790	47.659	ng	98
89) Benzo(k)fluoranthene	24.328	252	784610	45.345	ng	99
90) Benzo(a)pyrene	25.192	252	714778	42.340	ng	97
91) Dibenzo(a,h)anthracene	29.352	278	770273	45.402	ng	99
92) Benzo(g,h,i)perylene	30.545	276	744395	45.473	ng	99

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 Upadhyay

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

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