

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081122\
 Data File : BM036210.D
 Acq On : 11 Aug 2022 15:35
 Operator : CG/JU
 Sample : PB146776BS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB146776BS

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 08/12/2022
 Supervised By : Jagrut Upadhyay 08/12/2022

Quant Time: Aug 12 02:14:53 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM080122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 10 22:19:26 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.828	152	262194	20.000	ng	0.00
21) Naphthalene-d8	10.628	136	1067468	20.000	ng	-0.01
39) Acenaphthene-d10	14.469	164	555674	20.000	ng	0.00
64) Phenanthrene-d10	17.215	188	958218	20.000	ng	0.00
76) Chrysene-d12	21.398	240	747486	20.000	ng	0.00
86) Perylene-d12	23.739	264	721905	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.410	112	2059225	127.331	ng	0.00
7) Phenol-d6	7.028	99	2571633	124.970	ng	0.00
23) Nitrobenzene-d5	8.998	82	1472602	77.222	ng	0.00
42) 2,4,6-Tribromophenol	15.963	330	699379	107.287	ng	0.00
45) 2-Fluorobiphenyl	13.098	172	2932092	77.908	ng	0.00
79) Terphenyl-d14	19.839	244	3194365	84.292	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.263	88	204893	31.441	ng	# 93
3) Pyridine	3.675	79	584714	33.460	ng	# 89
4) n-Nitrosodimethylamine	3.587	42	311335	43.355	ng	# 68
6) Aniline	7.157	93	1012331	44.424	ng	95
8) 2-Chlorophenol	7.398	128	761349	44.154	ng	95
9) Benzaldehyde	6.969	77	346477	31.804	ng	94
10) Phenol	7.051	94	963667	46.509	ng	91
11) bis(2-Chloroethyl)ether	7.257	93	696951	40.833	ng	96
12) 1,3-Dichlorobenzene	7.710	146	781884	40.920	ng	99
13) 1,4-Dichlorobenzene	7.863	146	801656	41.140	ng	98
14) 1,2-Dichlorobenzene	8.175	146	780736	41.525	ng	98
15) Benzyl Alcohol	8.081	79	598734m	43.358	ng	
16) 2,2'-oxybis(1-Chloropr...	8.357	45	904658	39.860	ng	97
17) 2-Methylphenol	8.293	107	607194	44.270	ng	96
18) Hexachloroethane	8.898	117	286607	40.859	ng	95
19) n-Nitroso-di-n-propyla...	8.645	70	504270	40.343	ng	88
20) 3+4-Methylphenols	8.628	107	809558	43.444	ng	98
22) Acetophenone	8.663	105	1066670	42.213	ng	# 92
24) Nitrobenzene	9.040	77	787430	43.923	ng	97
25) Isophorone	9.569	82	1335002	43.033	ng	98
26) 2-Nitrophenol	9.751	139	383634	45.370	ng	93
27) 2,4-Dimethylphenol	9.822	122	641670	48.988	ng	98
28) bis(2-Chloroethoxy)met...	10.051	93	840179	38.014	ng	100
29) 2,4-Dichlorophenol	10.292	162	625608	42.471	ng	99
30) 1,2,4-Trichlorobenzene	10.492	180	650550	38.972	ng	99
31) Naphthalene	10.681	128	2203862	41.526	ng	100
32) Benzoic acid	9.992	122	431435	41.551	ng	92
33) 4-Chloroaniline	10.804	127	649573	30.370	ng	96
34) Hexachlorobutadiene	10.957	225	371893	37.913	ng	98
35) Caprolactam	11.616	113	181427	39.987	ng	90
36) 4-Chloro-3-methylphenol	11.945	107	636135	41.086	ng	99
37) 2-Methylnaphthalene	12.292	142	1410920	39.951	ng	99
38) 1-Methylnaphthalene	12.516	142	1353956	39.148	ng	97
40) 1,2,4,5-Tetrachloroben...	12.663	216	653512	43.380	ng	98
41) Hexachlorocyclopentadiene	12.633	237	823402	103.199	ng	98
43) 2,4,6-Trichlorophenol	12.910	196	439433	42.867	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.992	196	475424	43.935	ng	96
46) 1,1'-Biphenyl	13.310	154	1747924	43.163	ng	99
47) 2-Chloronaphthalene	13.351	162	1350435	43.571	ng	99
48) 2-Nitroaniline	13.569	65	390881	46.277	ng	89
49) Acenaphthylene	14.192	152	2113105	44.105	ng	99
50) Dimethylphthalate	13.939	163	1453507	38.465	ng	99
51) 2,6-Dinitrotoluene	14.063	165	328333	43.305	ng	91
52) Acenaphthene	14.533	154	1477050m	47.129	ng	
53) 3-Nitroaniline	14.392	138	288525	32.617	ng	98
54) 2,4-Dinitrophenol	14.598	184	365785	91.170	ng	92
55) Dibenzofuran	14.869	168	1910959	40.717	ng	97
56) 4-Nitrophenol	14.727	139	614365	86.789	ng	86
57) 2,4-Dinitrotoluene	14.845	165	436276	43.901	ng	98
58) Fluorene	15.516	166	1511118	40.818	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.104	232	360103	39.284	ng	97
60) Diethylphthalate	15.298	149	1394950	35.922	ng	99
61) 4-Chlorophenyl-phenyle...	15.516	204	695075	37.653	ng	95
62) 4-Nitroaniline	15.557	138	358279m	39.371	ng	
63) Azobenzene	15.804	77	1414973	38.329	ng	94
65) 4,6-Dinitro-2-methylph...	15.610	198	208624	42.581	ng	95
66) n-Nitrosodiphenylamine	15.733	169	1228155	45.403	ng	99
67) 4-Bromophenyl-phenylether	16.404	248	395442	41.258	ng	97
68) Hexachlorobenzene	16.516	284	478239	43.487	ng	91
69) Atrazine	16.686	200	376504	49.743	ng	100
70) Pentachlorophenol	16.868	266	770404	119.614	ng	100
71) Phenanthrene	17.257	178	2139752	43.831	ng	99
72) Anthracene	17.345	178	2145883	45.404	ng	99
73) Carbazole	17.627	167	2007431	41.920	ng	98
74) Di-n-butylphthalate	18.186	149	2033816	35.578	ng	99
75) Fluoranthene	19.274	202	2192145	40.116	ng	99
77) Benzidine	19.468	184	1237542	110.395	ng	99
78) Pyrene	19.633	202	2267132	48.912	ng	99
80) Butylbenzylphthalate	20.539	149	798332	39.906	ng	96
81) Benzo(a)anthracene	21.380	228	1988921	43.199	ng	99
82) 3,3'-Dichlorobenzidine	21.321	252	666227	59.594	ng	99
83) Chrysene	21.433	228	1879737	42.950	ng	99
84) Bis(2-ethylhexyl)phtha...	21.309	149	1053514	34.906	ng	99
85) Di-n-octyl phthalate	22.215	149	1700807	34.530	ng	95
87) Indeno(1,2,3-cd)pyrene	26.174	276	2361484	46.549	ng	98
88) Benzo(b)fluoranthene	23.027	252	1972570	46.638	ng	98
89) Benzo(k)fluoranthene	23.074	252	1892538	44.302	ng	99
90) Benzo(a)pyrene	23.639	252	1739550	49.019	ng	98
91) Dibenzo(a,h)anthracene	26.191	278	2072317	48.800	ng	98
92) Benzo(g,h,i)perylene	26.927	276	2121724	51.611	ng	98

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

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