

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080422\
 Data File : BM036095.D
 Acq On : 04 Aug 2022 12:48
 Operator : CG/JU
 Sample : PB146679BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB146679BS

Quant Time: Aug 05 03:38:45 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM080122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 05 03:36:25 2022
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By : Christian Giraldo 08/05/2022
 Supervised By : Jagrut Upadhyay 08/05/2022

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.857	152	348043	20.000	ng	0.00
21) Naphthalene-d8	10.663	136	1463601	20.000	ng	0.00
39) Acenaphthene-d10	14.498	164	807587	20.000	ng	0.00
64) Phenanthrene-d10	17.239	188	1442062	20.000	ng	0.00
76) Chrysene-d12	21.427	240	1108692	20.000	ng	0.00
86) Perylene-d12	23.786	264	1014257	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.428	112	2819910	131.358	ng	0.00
7) Phenol-d6	7.040	99	3478741	127.353	ng	0.00
23) Nitrobenzene-d5	9.034	82	2159600	82.596	ng	0.00
42) 2,4,6-Tribromophenol	15.992	330	1182155	124.778	ng	0.00
45) 2-Fluorobiphenyl	13.127	172	4753908	86.913	ng	0.00
79) Terphenyl-d14	19.868	244	5819551	103.534	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.293	88	234509	27.110	ng	93
3) Pyridine	3.704	79	692079	29.835	ng	89
4) n-Nitrosodimethylamine	3.610	42	377654	39.619	ng	# 65
6) Aniline	7.187	93	1353782	44.755	ng	96
8) 2-Chlorophenol	7.422	128	1016996	44.432	ng	96
9) Benzaldehyde	6.998	77	481794	33.316	ng	96
10) Phenol	7.063	94	1238300	45.022	ng	90
11) bis(2-Chloroethyl)ether	7.293	93	950530	41.953	ng	94
12) 1,3-Dichlorobenzene	7.745	146	1015286	40.029	ng	98
13) 1,4-Dichlorobenzene	7.898	146	1041455	40.263	ng	99
14) 1,2-Dichlorobenzene	8.210	146	1013704	40.617	ng	99
15) Benzyl Alcohol	8.110	79	808649m	44.115	ng	
16) 2,2'-oxybis(1-Chloropr...	8.398	45	1291323	42.862	ng	96
17) 2-Methylphenol	8.310	107	821307	45.110	ng	95
18) Hexachloroethane	8.939	117	388882	41.764	ng	96
19) n-Nitroso-di-n-propyla...	8.681	70	718499	43.304	ng	89
20) 3+4-Methylphenols	8.645	107	1093498	44.207	ng	96
22) Acetophenone	8.692	105	1448233	41.801	ng	# 93
24) Nitrobenzene	9.075	77	1058227	43.052	ng	96
25) Isophorone	9.604	82	1976518	46.468	ng	98
26) 2-Nitrophenol	9.781	139	522702	45.086	ng	92
27) 2,4-Dimethylphenol	9.845	122	910125	50.677	ng	98
28) bis(2-Chloroethoxy)met...	10.086	93	1256994	41.480	ng	99
29) 2,4-Dichlorophenol	10.316	162	896200	44.374	ng	100
30) 1,2,4-Trichlorobenzene	10.528	180	911746	39.836	ng	98
31) Naphthalene	10.716	128	3020555	41.510	ng	100
32) Benzoic acid	10.004	122	563262	39.565	ng	93
33) 4-Chloroaniline	10.833	127	963143	32.843	ng	96
34) Hexachlorobutadiene	10.992	225	537367	39.956	ng	99
35) Caprolactam	11.639	113	255722m	41.107	ng	
36) 4-Chloro-3-methylphenol	11.963	107	909244	42.831	ng	100
37) 2-Methylnaphthalene	12.327	142	2045537	42.243	ng	98
38) 1-Methylnaphthalene	12.545	142	1956979	41.269	ng	98
40) 1,2,4,5-Tetrachloroben...	12.692	216	973753	44.475	ng	98
41) Hexachlorocyclopentadiene	12.669	237	1410185	121.610	ng	98
43) 2,4,6-Trichlorophenol	12.939	196	675360	45.331	ng	98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.010	196	715094	45.469	ng	97
46) 1,1'-Biphenyl	13.339	154	2606278	44.283	ng	99
47) 2-Chloronaphthalene	13.380	162	1985258	44.073	ng	99
48) 2-Nitroaniline	13.592	65	557753	45.435	ng	90
49) Acenaphthylene	14.227	152	3117558	44.773	ng	100
50) Dimethylphthalate	13.969	163	2328308	42.395	ng	99
51) 2,6-Dinitrotoluene	14.086	165	504440	45.779	ng	93
52) Acenaphthene	14.563	154	2209229m	48.503	ng	
53) 3-Nitroaniline	14.416	138	437437	34.026	ng	97
54) 2,4-Dinitrophenol	14.621	184	573282	98.316	ng	91
55) Dibenzofuran	14.898	168	2870615	42.086	ng	99
56) 4-Nitrophenol	14.727	139	919240	89.351	ng	88
57) 2,4-Dinitrotoluene	14.874	165	668520	46.287	ng	96
58) Fluorene	15.545	166	2288813	42.540	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.127	232	569812	42.772	ng	97
60) Diethylphthalate	15.327	149	2365639	41.916	ng	99
61) 4-Chlorophenyl-phenyle...	15.545	204	1123430	41.874	ng	96
62) 4-Nitroaniline	15.580	138	533716	40.355	ng	90
63) Azobenzene	15.833	77	2274019	42.384	ng	93
65) 4,6-Dinitro-2-methylph...	15.627	198	329565	44.696	ng	95
66) n-Nitrosodiphenylamine	15.763	169	1911474	46.955	ng	99
67) 4-Bromophenyl-phenylether	16.433	248	665736	46.154	ng	97
68) Hexachlorobenzene	16.545	284	772086	46.651	ng	92
69) Atrazine	16.715	200	637483	55.964	ng	98
70) Pentachlorophenol	16.892	266	921243	95.043	ng	99
71) Phenanthrene	17.286	178	3258557	44.354	ng	99
72) Anthracene	17.374	178	3328368	46.795	ng	99
73) Carbazole	17.651	167	2996913	41.585	ng	99
74) Di-n-butylphthalate	18.215	149	3873406	45.024	ng	99
75) Fluoranthene	19.298	202	3429017	41.696	ng	98
77) Benzidine	19.492	184	2102108	126.427	ng	99
78) Pyrene	19.662	202	3491210	50.781	ng	99
80) Butylbenzylphthalate	20.568	149	1528476	51.511	ng	97
81) Benzo(a)anthracene	21.409	228	3113578	45.594	ng	99
82) 3,3'-Dichlorobenzidine	21.345	252	998211	60.200	ng	99
83) Chrysene	21.462	228	2912642	44.869	ng	99
84) Bis(2-ethylhexyl)phtha...	21.345	149	2297751	51.328	ng	97
85) Di-n-octyl phthalate	22.256	149	3562840	48.767	ng	95
87) Indeno(1,2,3-cd)pyrene	26.244	276	3139368	44.045	ng	98
88) Benzo(b)fluoranthene	23.068	252	2958475	49.785	ng	99
89) Benzo(k)fluoranthene	23.115	252	2891224	48.172	ng	98
90) Benzo(a)pyrene	23.686	252	2566471	51.475	ng	99
91) Dibenzo(a,h)anthracene	26.262	278	2800780	46.944	ng	99
92) Benzo(g,h,i)perylene	27.003	276	2737192	47.390	ng	99

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

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