

Data Path : Z:\HPCHEM1\BNA N\DATA\BN020718\
 Data File : BN000037.D
 Acq On : 07 Feb 2018 14:23
 Operator : JU/SJ
 Sample : SSTDICC060
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTDICC060

Manual Integrations
 APPROVED

Sohil
 2/8/2018 3:37:29 PM

Quant Time: Feb 07 18:20:12 2018
 Quant Method : Z:\HPCHEM1\BNA N\METHODS\8270-BN020718.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 07 14:04:10 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.47	152	409307	20.00	ng	0.00
21) Naphthalene-d8	11.32	136	2033829	20.00	ng	0.00
38) Acenaphthene-d10	15.08	164	1193440	20.00	ng	0.00
63) Phenanthrene-d10	17.80	188	2531370	20.00	ng	0.00
75) Chrysene-d12	21.93	240	2607216	20.00	ng	0.00
86) Perylene-d12	24.42	264	2836226	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.95	112	3265614	117.64	ng	0.00
7) Phenol-d6	7.59	99	4942723	117.22	ng	0.00
23) Nitrobenzene-d5	9.65	82	4643964	138.07	ng	0.00
41) 2,4,6-Tribromophenol	16.55	330	1564609	130.42	ng	0.00
44) 2-Fluorobiphenyl	13.72	172	10191771	114.37	ng	0.00
78) Terphenyl-d14	20.37	244	11549694	116.92	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.66	88	615719	56.066	ng	94
3) Pyridine	4.10	79	2075796	62.324	ng	98
4) n-Nitrosodimethylamine	4.01	42	570106	58.179	ng	# 88
6) Aniline	7.78	93	3373309	58.685	ng	91
8) 2-Chlorophenol	8.02	128	1880997	61.250	ng	87
9) Benzaldehyde	7.59	77	1420185	65.962	ng	87
10) Phenol	7.62	94	2590059	60.093	ng	85
11) bis(2-Chloroethyl)ether	7.88	93	2063357	60.008	ng	# 81
12) 1,3-Dichlorobenzene	8.36	146	2065701	60.589	ng	99
13) 1,4-Dichlorobenzene	8.51	146	2093865	59.799	ng	97
14) 1,2-Dichlorobenzene	8.83	146	2077572	60.053	ng	96
15) Benzyl Alcohol	8.70	79	1836021	59.292	ng	90
16) 2,2'-oxybis(1-Chloropropan	9.00	45	2162040	58.134	ng	78
17) 2-Methylphenol	8.90	107	1874647	60.318	ng	95
18) Hexachloroethane	9.58	117	761263	63.399	ng	# 81
19) n-Nitroso-di-n-propylamine	9.28	70	1680971	59.158	ng	# 82
20) 3+4-Methylphenols	9.23	107	2665678	61.238	ng	92
22) Acetophenone	9.30	105	3625229	60.350	ng	# 86
24) Nitrobenzene	9.69	77	2482663	68.819	ng	# 92
25) Isophorone	10.22	82	4742338	59.524	ng	96
26) 2-Nitrophenol	10.41	139	971154	87.896	ng	89
27) 2,4-Dimethylphenol	10.45	122	1990585	60.936	ng	95
28) bis(2-Chloroethoxy)methane	10.70	93	2900604	63.120	ng	96
29) 2,4-Dichlorophenol	10.95	162	1861424	66.410	ng	96
30) 1,2,4-Trichlorobenzene	11.17	180	2051547	62.505	ng	97
31) Naphthalene	11.36	128	6909489	60.575	ng	98
32) Benzoic acid	10.59	122	1453341	85.916	ng	86
33) 4-Chloroaniline	11.46	127	3189246	60.309	ng	94
34) Hexachlorobutadiene	11.64	225	1060403	63.415	ng	96
35) Caprolactam	12.22	113	782419	62.799	ng	95
36) 4-Chloro-3-methylphenol	12.55	107	2336578	61.309	ng	96
37) 2-Methylnaphthalene	12.94	142	4863650	60.342	ng	95
39) 1,2,4,5-Tetrachlorobenzene	13.30	216	2218639	63.823	ng	# 96
40) Hexachlorocyclopentadiene	13.27	237	993484	85.567	ng	93

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42) 2,4,6-Trichlorophenol	13.52	196	1410559	68.651	nq	98
43) 2,4,5-Trichlorophenol	13.59	196	1669238	68.986	nq #	94
45) 1,1'-Biphenyl	13.93	154	6498247	60.941	nq	97
46) 2-Chloronaphthalene	13.97	162	4943871	61.758	nq	97
47) 2-Nitroaniline	14.16	65	1464682	77.470	nq #	78
48) Acenaphthylene	14.81	152	8146737	60.604	nq	98
49) Dimethylphthalate	14.53	163	6348727	59.522	nq	99
50) 2,6-Dinitrotoluene	14.65	165	1363887	71.661	nq	92
51) Acenaphthene	15.15	154	5176489	61.026	nq	98
52) 3-Nitroaniline	14.97	138	1646347	70.564	nq #	89
53) 2,4-Dinitrophenol	15.17	184	421112	96.227	nq #	10
54) Dibenzofuran	15.47	168	7178171	58.970	nq	95
55) 4-Nitrophenol	15.25	139	1243557	71.843	nq #	81
56) 2,4-Dinitrotoluene	15.42	165	1862525	75.024	nq #	89
57) Fluorene	16.12	166	5852074	57.096	nq	96
58) 2,3,4,6-Tetrachlorophenol	15.69	232	1311957m	66.579	nq	
59) Diethylphthalate	15.87	149	6510372	58.418	nq	99
60) 4-Chlorophenyl-phenylether	16.10	204	2676173	59.076	nq	93
61) 4-Nitroaniline	16.13	138	1685093	67.764	nq	90
62) Azobenzene	16.39	77	6456048	57.971	nq	90
64) 4,6-Dinitro-2-methylphenol	16.17	198	754712	96.582	nq	87
65) n-Nitrosodiphenylamine	16.32	169	5274120	63.758	nq	97
66) 4-Bromophenyl-phenylether	16.99	248	1539962	65.012	nq #	87
67) Hexachlorobenzene	17.11	284	1755745	64.337	nq #	89
68) Atrazine	17.25	200	1729278	68.580	nq	95
69) Pentachlorophenol	17.45	266	1153526	70.577	nq	98
70) Phenanthrene	17.85	178	9080922	59.620	nq	98
71) Anthracene	17.94	178	9142329	60.591	nq	99
72) Carbazole	18.20	167	9173232	58.761	nq #	97
73) Di-n-butylphthalate	18.73	149	10464984	60.363	nq #	97
74) Fluoranthene	19.83	202	9857870	57.000	nq	91
76) Benzidine	19.99	184	5542547m	83.983	nq	
77) Pyrene	20.19	202	10393799	63.190	nq	95
79) Butylbenzylphthalate	21.05	149	5011922	76.442	nq	97
80) Benzo(a)anthracene	21.92	228	9962429	61.631	nq	98
81) 3,3'-Dichlorobenzidine	21.83	252	3827609	70.660	nq #	95
82) Chrysene	21.97	228	9749607	60.730	nq	95
83) Bis(2-ethylhexyl)phthalate	21.81	149	7416284	68.359	nq #	97
84) Di-n-octyl phthalate	22.77	149	11731069	66.807	nq #	93
85) Indeno(1,2,3-cd)pyrene	27.01	276	12928306	62.430	nq #	86
87) Benzo(b)fluoranthene	23.67	252	10735438	63.660	nq #	96
88) Benzo(k)fluoranthene	23.72	252	10414270	61.593	nq #	93
89) Benzo(a)pyrene	24.32	252	10333057	63.074	nq #	95
90) Dibenzo(a,h)anthracene	27.02	278	10787585	62.912	nq #	89
91) Benzo(g,h,i)perylene	27.80	276	10913121	62.622	nq #	87

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

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