

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011223\
 Data File : BP013461.D
 Acq On : 12 Jan 2023 16:17
 Operator : CG/JU
 Sample : SSTDICC010
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SSTDICC010

Quant Time: Jan 13 02:32:58 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP011223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 13 02:28:25 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.910	152	475052	20.000	ng	0.00
21) Naphthalene-d8	10.722	136	2080241	20.000	ng	0.00
39) Acenaphthene-d10	14.557	164	1431538	20.000	ng	0.00
64) Phenanthrene-d10	17.316	188	3300517	20.000	ng	0.00
76) Chrysene-d12	21.404	240	3349376	20.000	ng	0.00
86) Perylene-d12	23.868	264	4036170	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.463	112	546281	21.011	ng	0.00
7) Phenol-d6	7.075	99	756890	22.292	ng	0.00
23) Nitrobenzene-d5	9.075	82	762791	20.102	ng	0.00
42) 2,4,6-Tribromophenol	16.051	330	428059	24.801	ng	0.00
45) 2-Fluorobiphenyl	13.175	172	2037417	19.324	ng	0.00
79) Terphenyl-d14	19.869	244	3505347	20.647	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.334	88	109907	9.541	ng	96
3) Pyridine	3.752	79	325250	9.942	ng	99
4) n-Nitrosodimethylamine	3.658	42	156354	10.664	ng	# 95
6) Aniline	7.234	93	509713	12.397	ng	100
8) 2-Chlorophenol	7.469	128	319225	10.415	ng	98
9) Benzaldehyde	7.046	77	258953	12.711	ng	99
10) Phenol	7.099	94	390697	10.280	ng	98
11) bis(2-Chloroethyl)ether	7.328	93	325734	10.646	ng	99
12) 1,3-Dichlorobenzene	7.799	146	338564	9.494	ng	98
13) 1,4-Dichlorobenzene	7.946	146	344394	9.483	ng	98
14) 1,2-Dichlorobenzene	8.263	146	336018	9.748	ng	99
15) Benzyl Alcohol	8.151	79	291002	11.814	ng	97
16) 2,2'-oxybis(1-Chloropr...	8.440	45	511474	11.541	ng	99
17) 2-Methylphenol	8.357	107	295115	11.748	ng	99
18) Hexachloroethane	8.993	117	120523	9.799	ng	97
19) n-Nitroso-di-n-propyla...	8.716	70	284683	12.910	ng	99
20) 3+4-Methylphenols	8.687	107	403778	11.823	ng	98
22) Acetophenone	8.740	105	522667	10.097	ng	# 99
24) Nitrobenzene	9.122	77	388771	9.831	ng	98
25) Isophorone	9.646	82	799168	12.400	ng	98
26) 2-Nitrophenol	9.834	139	184758	10.206	ng	95
27) 2,4-Dimethylphenol	9.893	122	319644	11.722	ng	99
28) bis(2-Chloroethoxy)met...	10.128	93	471781	11.416	ng	99
29) 2,4-Dichlorophenol	10.369	162	329114	9.779	ng	99
30) 1,2,4-Trichlorobenzene	10.587	180	358832	9.033	ng	99
31) Naphthalene	10.775	128	1084069	9.517	ng	99
32) Benzoic acid	9.993	122	239507	11.132	ng	97
33) 4-Chloroaniline	10.887	127	480811	11.318	ng	100
34) Hexachlorobutadiene	11.051	225	226902	8.954	ng	99
35) Caprolactam	11.669	113	118968	13.691	ng	98
36) 4-Chloro-3-methylphenol	12.010	107	366846	11.670	ng	98
37) 2-Methylnaphthalene	12.381	142	816680	10.661	ng	99
38) 1-Methylnaphthalene	12.598	142	778750	10.440	ng	98
40) 1,2,4,5-Tetrachloroben...	12.745	216	452939	8.869	ng	99
41) Hexachlorocyclopentadiene	12.722	237	205191	7.968	ng	97
43) 2,4,6-Trichlorophenol	12.992	196	311399	9.669	ng	99

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44) 2,4,5-Trichlorophenol	13.063	196	358941	10.316	ng	99
46) 1,1'-Biphenyl	13.387	154	1076921	9.514	ng	99
47) 2-Chloronaphthalene	13.434	162	821704	9.134	ng	99
48) 2-Nitroaniline	13.639	65	256422	11.521	ng	94
49) Acenaphthylene	14.281	152	1366608	10.327	ng	99
50) Dimethylphthalate	14.010	163	1137252	10.861	ng	99
51) 2,6-Dinitrotoluene	14.134	165	236001	11.014	ng	96
52) Acenaphthene	14.622	154	805476	9.791	ng	99
53) 3-Nitroaniline	14.469	138	249967	11.441	ng	100
54) 2,4-Dinitrophenol	14.681	184	136993	10.731	ng	98
55) Dibenzofuran	14.957	168	1351890	10.268	ng	99
56) 4-Nitrophenol	14.781	139	194522	10.863	ng	98
57) 2,4-Dinitrotoluene	14.928	165	314812	11.754	ng	96
58) Fluorene	15.604	166	1083718	10.635	ng	100
59) 2,3,4,6-Tetrachlorophenol	15.186	232	317441	10.673	ng	99
60) Diethylphthalate	15.375	149	1114388	11.577	ng	99
61) 4-Chlorophenyl-phenyle...	15.598	204	582644	11.048	ng	99
62) 4-Nitroaniline	15.633	138	250350	12.277	ng	97
63) Azobenzene	15.892	77	1014642	11.342	ng	100
65) 4,6-Dinitro-2-methylph...	15.692	198	198392	9.357	ng	94
66) n-Nitrosodiphenylamine	15.816	169	958489	9.301	ng	99
67) 4-Bromophenyl-phenylether	16.498	248	367207	9.409	ng	99
68) Hexachlorobenzene	16.616	284	403366	8.827	ng	99
69) Atrazine	16.769	200	350291	12.115	ng	98
70) Pentachlorophenol	16.963	266	270223	10.041	ng	99
71) Phenanthrene	17.357	178	1753703	9.338	ng	99
72) Anthracene	17.451	178	1775962	10.073	ng	98
73) Carbazole	17.722	167	1609872	10.415	ng	100
74) Di-n-butylphthalate	18.263	149	1919822	11.292	ng	99
75) Fluoranthene	19.327	202	2176645	10.768	ng	99
77) Benzidine	19.504	184	1010769	21.246	ng	100
78) Pyrene	19.680	202	2272419	9.442	ng	98
80) Butylbenzylphthalate	20.545	149	925262	11.858	ng	100
81) Benzo(a)anthracene	21.392	228	2310067	10.090	ng	98
82) 3,3'-Dichlorobenzidine	21.321	252	852123	12.499	ng	99
83) Chrysene	21.445	228	2195495	9.831	ng	98
84) Bis(2-ethylhexyl)phtha...	21.298	149	1384191	13.042	ng	99
85) Di-n-octyl phthalate	22.239	149	2419433	13.574	ng	99
87) Indeno(1,2,3-cd)pyrene	26.439	276	3001015	9.100	ng	100
88) Benzo(b)fluoranthene	23.115	252	2427500	9.341	ng	100
89) Benzo(k)fluoranthene	23.162	252	2333731	8.877	ng	99
90) Benzo(a)pyrene	23.757	252	2347692	10.765	ng	99
91) Dibenzo(a,h)anthracene	26.451	278	2493596	9.098	ng	99
92) Benzo(g,h,i)perylene	27.227	276	2479666	9.131	ng	99

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

