

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060722\
 Data File : BP010709.D
 Acq On : 07 Jun 2022 15:16
 Operator : CG/JU
 Sample : SSTDICC010
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SSTDICC010

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 06/08/2022
 Supervised By : Jagrut Upadhyay 06/09/2022

Quant Time: Jun 07 17:16:34 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP060722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 07 17:14:59 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.840	152	84532	20.000	ng	0.00
21) Naphthalene-d8	10.640	136	339498	20.000	ng	0.00
39) Acenaphthene-d10	14.481	164	222967	20.000	ng	0.00
64) Phenanthrene-d10	17.234	188	488378	20.000	ng	0.00
76) Chrysene-d12	21.327	240	492284	20.000	ng	0.00
86) Perylene-d12	23.733	264	496129	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.428	112	95844	20.684	ng	0.00
7) Phenol-d6	7.016	99	126762	19.270	ng	0.00
23) Nitrobenzene-d5	8.999	82	136650	19.030	ng	0.00
42) 2,4,6-Tribromophenol	15.975	330	55962	22.173	ng	0.00
45) 2-Fluorobiphenyl	13.098	172	298064	18.973	ng	0.00
79) Terphenyl-d14	19.792	244	499772	19.069	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.334	88	19685	10.004	ng	94
3) Pyridine	3.746	79	54486	9.887	ng	97
4) n-Nitrosodimethylamine	3.646	42	28760	8.559	ng	# 98
6) Aniline	7.169	93	73623	9.513	ng	99
8) 2-Chlorophenol	7.411	128	51922	9.928	ng	98
9) Benzaldehyde	6.987	77	44354	10.246	ng	99
10) Phenol	7.046	94	65436	9.619	ng	96
11) bis(2-Chloroethyl)ether	7.263	93	49716	9.221	ng	98
12) 1,3-Dichlorobenzene	7.728	146	61101	10.081	ng	99
13) 1,4-Dichlorobenzene	7.875	146	61667	9.930	ng	96
14) 1,2-Dichlorobenzene	8.193	146	58886	9.905	ng	97
15) Benzyl Alcohol	8.093	79	49313	9.351	ng	98
16) 2,2'-oxybis(1-Chloropr...	8.375	45	87484	8.679	ng	95
17) 2-Methylphenol	8.293	107	44778	9.808	ng	98
18) Hexachloroethane	8.916	117	24366	10.143	ng	98
19) n-Nitroso-di-n-propyla...	8.652	70	45745	9.534	ng	93
20) 3+4-Methylphenols	8.628	107	60707	9.673	ng	98
22) Acetophenone	8.669	105	85997	9.713	ng	# 96
24) Nitrobenzene	9.046	77	68694	9.469	ng	97
25) Isophorone	9.575	82	114051	9.507	ng	99
26) 2-Nitrophenol	9.757	139	27087	9.580	ng	96
27) 2,4-Dimethylphenol	9.828	122	40801	9.709	ng	96
28) bis(2-Chloroethoxy)met...	10.052	93	64052	9.666	ng	97
29) 2,4-Dichlorophenol	10.305	162	48062	9.513	ng	96
30) 1,2,4-Trichlorobenzene	10.505	180	57745	9.717	ng	97
31) Naphthalene	10.687	128	174165	9.738	ng	99
32) Benzoic acid	9.934	122	19436	6.577	ng	92
33) 4-Chloroaniline	10.810	127	65656	9.503	ng	96
34) Hexachlorobutadiene	10.975	225	38720	9.651	ng	96
35) Caprolactam	11.593	113	15719	10.779	ng	# 87
36) 4-Chloro-3-methylphenol	11.946	107	58117	10.097	ng	98
37) 2-Methylnaphthalene	12.304	142	119700	9.592	ng	98
38) 1-Methylnaphthalene	12.522	142	118964	9.761	ng	99
40) 1,2,4,5-Tetrachloroben...	12.675	216	65371	9.425	ng	99
41) Hexachlorocyclopentadiene	12.651	237	17403	4.713	ng	96
43) 2,4,6-Trichlorophenol	12.922	196	40528	9.360	ng	93

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44) 2,4,5-Trichlorophenol	12.998	196	45293	9.248	ng	97
46) 1,1'-Biphenyl	13.310	154	159277	9.577	ng	99
47) 2-Chloronaphthalene	13.351	162	121446	9.291	ng	99
48) 2-Nitroaniline	13.563	65	42887	9.466	ng	96
49) Acenaphthylene	14.198	152	193634	9.644	ng	99
50) Dimethylphthalate	13.940	163	162580	10.130	ng	98
51) 2,6-Dinitrotoluene	14.057	165	34624	10.144	ng	93
52) Acenaphthene	14.545	154	117998	9.572	ng	98
53) 3-Nitroaniline	14.393	138	32579	9.746	ng	94
54) 2,4-Dinitrophenol	14.622	184	13668	7.772	ng	92
55) Dibenzofuran	14.881	168	191269	9.812	ng	99
56) 4-Nitrophenol	14.740	139	21064	8.346	ng	96
57) 2,4-Dinitrotoluene	14.857	165	45789	10.306	ng	# 94
58) Fluorene	15.528	166	158002	9.914	ng	96
59) 2,3,4,6-Tetrachlorophenol	15.116	232	40735	9.702	ng	98
60) Diethylphthalate	15.304	149	167298	10.471	ng	100
61) 4-Chlorophenyl-phenyle...	15.522	204	82220	10.265	ng	96
62) 4-Nitroaniline	15.557	138	32681m	10.081	ng	
63) Azobenzene	15.816	77	165455	9.805	ng	97
65) 4,6-Dinitro-2-methylph...	15.628	198	22767	7.952	ng	95
66) n-Nitrosodiphenylamine	15.740	169	133212	9.059	ng	98
67) 4-Bromophenyl-phenylether	16.422	248	49077	9.237	ng	97
68) Hexachlorobenzene	16.545	284	59626	10.160	ng	# 89
69) Atrazine	16.698	200	49366	10.350	ng	99
70) Pentachlorophenol	16.898	266	24944	7.405	ng	92
71) Phenanthrene	17.275	178	257837	9.627	ng	99
72) Anthracene	17.369	178	242616	9.460	ng	99
73) Carbazole	17.645	167	214644	9.848	ng	99
74) Di-n-butylphthalate	18.192	149	280144	10.110	ng	98
75) Fluoranthene	19.245	202	300767	10.339	ng	98
77) Benzidine	19.433	184	97240	10.283	ng	98
78) Pyrene	19.604	202	315864	9.207	ng	98
80) Butylbenzylphthalate	20.469	149	130371	9.649	ng	97
81) Benzo(a)anthracene	21.310	228	308906	9.564	ng	98
82) 3,3'-Dichlorobenzidine	21.245	252	92090	10.396	ng	97
83) Chrysene	21.363	228	301689	9.523	ng	98
84) Bis(2-ethylhexyl)phtha...	21.233	149	187841	9.862	ng	100
85) Di-n-octyl phthalate	22.157	149	315531	9.941	ng	100
87) Indeno(1,2,3-cd)pyrene	26.233	276	332574	9.195	ng	98
88) Benzo(b)fluoranthene	22.998	252	282388	9.040	ng	99
89) Benzo(k)fluoranthene	23.045	252	301387	9.518	ng	99
90) Benzo(a)pyrene	23.627	252	236912	9.169	ng	99
91) Dibenzo(a,h)anthracene	26.251	278	278622	9.211	ng	98
92) Benzo(g,h,i)perylene	27.004	276	278654	9.266	ng	96

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

