

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070721\
 Data File : BP006128.D
 Acq On : 07 Jul 2021 11:54
 Operator : CG/JU
 Sample : SSTDICC010
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SSTDICC010

Manual Integrations
 APPROVED

mohammad
 7/8/2021 12:38:53 PM

Quant Time: Jul 07 13:50:11 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP070721.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 07 13:48:57 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.875	152	167698	20.000	ng	0.00
21) Naphthalene-d8	10.681	136	679606	20.000	ng	# 0.00
39) Acenaphthene-d10	14.510	164	397479	20.000	ng	0.00
64) Phenanthrene-d10	17.263	188	751129	20.000	ng	# 0.00
76) Chrysene-d12	21.345	240	719031	20.000	ng	# 0.00
86) Perylene-d12	23.739	264	815835	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.458	112	208217	19.923	ng	0.00
7) Phenol-d6	7.069	99	266025	19.639	ng	0.00
23) Nitrobenzene-d5	9.046	82	217439	15.686	ng	0.00
42) 2,4,6-Tribromophenol	16.010	330	85532	12.539	ng	0.00
45) 2-Fluorobiphenyl	13.140	172	554037	18.303	ng	0.00
79) Terphenyl-d14	19.816	244	729267	18.991	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.323	88	43765	9.269	ng	90
3) Pyridine	3.758	79	96766	8.720	ng	# 85
4) n-Nitrosodimethylamine	3.658	42	40563	7.864	ng	# 90
6) Aniline	7.211	93	140752	9.307	ng	99
8) 2-Chlorophenol	7.446	128	121427	10.134	ng	99
9) Benzaldehyde	7.028	77	80439	13.930	ng	94
10) Phenol	7.093	94	128768	9.516	ng	94
11) bis(2-Chloroethyl)ether	7.305	93	102346	9.981	ng	98
12) 1,3-Dichlorobenzene	7.763	146	129947	9.698	ng	97
13) 1,4-Dichlorobenzene	7.916	146	132219	9.676	ng	98
14) 1,2-Dichlorobenzene	8.228	146	127702	9.777	ng	99
15) Benzyl Alcohol	8.140	79	73768m	7.230	ng	
16) 2,2'-oxybis(1-Chloropr...	8.411	45	120020	12.662	ng	91
17) 2-Methylphenol	8.346	107	87767	9.519	ng	95
18) Hexachloroethane	8.952	117	43695	8.840	ng	91
19) n-Nitroso-di-n-propyla...	8.693	70	74746	8.970	ng	97
20) 3+4-Methylphenols	8.681	107	116875	9.384	ng	94
22) Acetophenone	8.711	105	158844	8.880	ng	98
24) Nitrobenzene	9.093	77	112473	7.814	ng	95
25) Isophorone	9.622	82	207120	8.091	ng	97
26) 2-Nitrophenol	9.805	139	55787	8.123	ng	90
27) 2,4-Dimethylphenol	9.875	122	87118	8.514	ng	98
28) bis(2-Chloroethoxy)met...	10.099	93	115003	8.608	ng	98
29) 2,4-Dichlorophenol	10.369	162	94248	7.665	ng	99
30) 1,2,4-Trichlorobenzene	10.546	180	105150	7.611	ng	94
31) Naphthalene	10.734	128	366357	9.317	ng	99
32) Benzoic acid	10.034	122	35675	5.435	ng	91
33) 4-Chloroaniline	10.863	127	133676	8.415	ng	98
34) Hexachlorobutadiene	11.010	225	56390	6.034	ng	97
35) Caprolactam	11.675	113	29082	8.211	ng	# 74
36) 4-Chloro-3-methylphenol	11.999	107	106960	8.567	ng	97
37) 2-Methylnaphthalene	12.346	142	249174	8.900	ng	97
38) 1-Methylnaphthalene	12.563	142	240519	9.121	ng	99
40) 1,2,4,5-Tetrachloroben...	12.716	216	97519	6.970	ng	97
43) 2,4,6-Trichlorophenol	12.969	196	66625	6.943	ng	93
44) 2,4,5-Trichlorophenol	13.057	196	72788	7.043	ng	97

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
46) 1,1'-Biphenyl	13.351	154	307792	9.604	ng	99
47) 2-Chloronaphthalene	13.393	162	234476	8.960	ng	95
48) 2-Nitroaniline	13.616	65	56214	8.170	ng	94
49) Acenaphthylene	14.240	152	378100	9.417	ng	99
50) Dimethylphthalate	13.975	163	291003	8.904	ng	99
51) 2,6-Dinitrotoluene	14.098	165	63010	8.483	ng	99
52) Acenaphthene	14.575	154	232808	9.548	ng	98
53) 3-Nitroaniline	14.445	138	61476	8.539	ng	95
54) 2,4-Dinitrophenol	14.693	184	12343	3.258	ng	94
55) Dibenzofuran	14.916	168	356432	8.889	ng	98
56) 4-Nitrophenol	14.822	139	23745	4.069	ng	# 81
57) 2,4-Dinitrotoluene	14.904	165	83365	8.391	ng	# 89
58) Fluorene	15.563	166	288295	9.228	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.151	232	55630	6.107	ng	# 72
60) Diethylphthalate	15.334	149	301193	9.186	ng	98
61) 4-Chlorophenyl-phenyle...	15.551	204	125708	7.760	ng	# 89
62) 4-Nitroaniline	15.610	138	49735	6.712	ng	# 81
63) Azobenzene	15.851	77	252305	8.505	ng	96
65) 4,6-Dinitro-2-methylph...	15.675	198	31610	5.727	ng	84
66) n-Nitrosodiphenylamine	15.775	169	250367	10.051	ng	99
67) 4-Bromophenyl-phenylether	16.451	248	72648	7.535	ng	# 90
68) Hexachlorobenzene	16.569	284	84552	7.317	ng	# 90
69) Atrazine	16.734	200	76064	9.815	ng	90
70) Pentachlorophenol	16.934	266	29592	4.605	ng	99
71) Phenanthrene	17.304	178	437074	9.689	ng	99
72) Anthracene	17.398	178	423472	9.539	ng	99
73) Carbazole	17.681	167	417739	9.801	ng	99
74) Di-n-butylphthalate	18.216	149	497627	10.077	ng	# 99
75) Fluoranthene	19.275	202	467148	8.528	ng	95
77) Benzidine	19.463	184	196282	13.119	ng	98
78) Pyrene	19.628	202	489138	9.744	ng	100
80) Butylbenzylphthalate	20.486	149	244536	11.841	ng	98
81) Benzo(a)anthracene	21.328	228	467552	9.008	ng	99
82) 3,3'-Dichlorobenzidine	21.269	252	130434	7.269	ng	# 97
83) Chrysene	21.380	228	463746	9.225	ng	98
84) Bis(2-ethylhexyl)phtha...	21.245	149	366653	12.106	ng	# 98
85) Di-n-octyl phthalate	22.157	149	660045	12.182	ng	99
87) Indeno(1,2,3-cd)pyrene	26.257	276	597224	8.560	ng	# 87
88) Benzo(b)fluoranthene	23.004	252	487772	8.729	ng	# 94
89) Benzo(k)fluoranthene	23.057	252	511272	9.436	ng	# 95
90) Benzo(a)pyrene	23.633	252	469771	8.942	ng	# 94
91) Dibenzo(a,h)anthracene	26.274	278	510822	8.507	ng	# 92
92) Benzo(g,h,i)perylene	27.039	276	514253	8.669	ng	# 87

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

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