

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042121\
 Data File : BM029567.D
 Acq On : 21 Apr 2021 12:41
 Operator : CG/JU
 Sample : SSTDICC050
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC050

Manual Integrations
 APPROVED

mohammad
 4/22/2021 12:11:14 PM

Quant Time: Apr 21 14:32:53 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM042121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 21 13:05:45 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.651	152	83761	20.00	ng	0.00
21) Naphthalene-d8	10.428	136	337254	20.00	ng	0.00
39) Acenaphthene-d10	14.292	164	255238	20.00	ng	0.00
64) Phenanthrene-d10	17.033	188	554168	20.00	ng	0.00
76) Chrysene-d12	21.221	240	646384	20.00	ng	0.00
86) Perylene-d12	23.415	264	671553	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.246	112	404664	98.07	ng	0.00
7) Phenol-d6	6.828	99	626731	101.39	ng	0.00
23) Nitrobenzene-d5	8.804	82	657212	101.31	ng	0.00
42) 2,4,6-Tribromophenol	15.786	330	354756	99.61	ng	0.00
45) 2-Fluorobiphenyl	12.910	172	1839838	97.91	ng	0.00
79) Terphenyl-d14	19.668	244	3413719	99.68	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.175	88	78828	48.46	ng	100
3) Pyridine	3.581	79	217458m	52.18	ng	
4) n-Nitrosodimethylamine	3.493	42	105700	51.49	ng	99
6) Aniline	6.987	93	341350	50.24	ng	99
8) 2-Chlorophenol	7.216	128	249664	48.69	ng	99
9) Benzaldehyde	6.798	77	144295	41.06	ng	99
10) Phenol	6.857	94	302227	50.88	ng	99
11) bis(2-Chloroethyl)ether	7.087	93	225906	48.95	ng	99
12) 1,3-Dichlorobenzene	7.539	146	285418	47.79	ng	99
13) 1,4-Dichlorobenzene	7.687	146	293720	47.93	ng	98
14) 1,2-Dichlorobenzene	7.998	146	292692	48.21	ng	99
15) Benzyl Alcohol	7.892	79	232884	51.37	ng	99
16) 2,2'-oxybis(1-Chloropr...	8.175	45	289880	48.49	ng	98
17) 2-Methylphenol	8.092	107	210738	50.28	ng	98
18) Hexachloroethane	8.722	117	106267	47.48	ng	97
19) n-Nitroso-di-n-propyla...	8.463	70	225544	49.69	ng	97
20) 3+4-Methylphenols	8.428	107	289634	49.75	ng	99
22) Acetophenone	8.469	105	404729	50.08	ng	# 97
24) Nitrobenzene	8.845	77	326282	49.58	ng	100
25) Isophorone	9.375	82	608274	48.97	ng	99
26) 2-Nitrophenol	9.551	139	149781	51.57	ng	99
27) 2,4-Dimethylphenol	9.616	122	219905	49.47	ng	98
28) bis(2-Chloroethoxy)met...	9.851	93	300375	48.76	ng	99
29) 2,4-Dichlorophenol	10.081	162	293258	49.89	ng	98
30) 1,2,4-Trichlorobenzene	10.298	180	326683	48.56	ng	97
31) Naphthalene	10.480	128	835371	48.09	ng	100
32) Benzoic acid	9.792	122	186852	53.56	ng	96
33) 4-Chloroaniline	10.592	127	348140	50.22	ng	98
34) Hexachlorobutadiene	10.763	225	207476	47.93	ng	96
35) Caprolactam	11.398	113	87354	49.42	ng	96
36) 4-Chloro-3-methylphenol	11.727	107	286731	49.69	ng	100
37) 2-Methylnaphthalene	12.098	142	674675	48.77	ng	98
38) 1-Methylnaphthalene	12.322	142	637205	47.89	ng	100
40) 1,2,4,5-Tetrachloroben...	12.474	216	376635	49.36	ng	99
41) Hexachlorocyclopentadiene	12.451	237	213786	51.35	ng	99
43) 2,4,6-Trichlorophenol	12.716	196	266063	50.19	ng	99

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44) 2,4,5-Trichlorophenol	12.786	196	289756	49.58	ng	99
46) 1,1'-Biphenyl	13.121	154	910927	48.61	ng	100
47) 2-Chloronaphthalene	13.163	162	723623	48.61	ng	99
48) 2-Nitroaniline	13.374	65	198415	53.74	ng	98
49) Acenaphthylene	14.016	152	1119036	48.86	ng	99
50) Dimethylphthalate	13.763	163	926577	48.02	ng	100
51) 2,6-Dinitrotoluene	13.874	165	210929	49.37	ng	98
52) Acenaphthene	14.357	154	696616	48.78	ng	98
53) 3-Nitroaniline	14.204	138	187215	52.89	ng	100
54) 2,4-Dinitrophenol	14.416	184	127281	54.81	ng	90
55) Dibenzofuran	14.692	168	1169620	48.60	ng	98
56) 4-Nitrophenol	14.521	139	157963	51.11	ng	98
57) 2,4-Dinitrotoluene	14.668	165	298159	51.25	ng	96
58) Fluorene	15.345	166	1000912	49.30	ng	100
59) 2,3,4,6-Tetrachlorophenol	14.921	232	259019	49.31	ng	98
60) Diethylphthalate	15.127	149	929829	48.48	ng	99
61) 4-Chlorophenyl-phenyle...	15.339	204	498279	48.50	ng	94
62) 4-Nitroaniline	15.374	138	202531	54.92	ng	99
63) Azobenzene	15.633	77	885359	49.01	ng	99
65) 4,6-Dinitro-2-methylph...	15.427	198	186602	50.05	ng	98
66) n-Nitrosodiphenylamine	15.557	169	849841	49.65	ng	100
67) 4-Bromophenyl-phenylether	16.233	248	285615	48.26	ng	98
68) Hexachlorobenzene	16.345	284	322539	48.77	ng	100
69) Atrazine	16.515	200	301770	48.67	ng	98
70) Pentachlorophenol	16.686	266	217163	50.52	ng	97
71) Phenanthrene	17.080	178	1556925	49.13	ng	100
72) Anthracene	17.168	178	1569589	50.11	ng	100
73) Carbazole	17.439	167	1478899	50.27	ng	99
74) Di-n-butylphthalate	18.009	149	1657656	50.63	ng	99
75) Fluoranthene	19.092	202	1922873	49.52	ng	100
77) Benzidine	19.286	184	684660	49.95	ng	99
78) Pyrene	19.456	202	2036645	49.22	ng	99
80) Butylbenzylphthalate	20.368	149	797064	50.18	ng	94
81) Benzo(a)anthracene	21.203	228	2094056	48.93	ng	99
82) 3,3'-Dichlorobenzidine	21.139	252	684268	49.42	ng	99
83) Chrysene	21.256	228	2049414	48.73	ng	99
84) Bis(2-ethylhexyl)phtha...	21.144	149	1299751	50.07	ng	99
85) Di-n-octyl phthalate	22.009	149	2138356	48.71	ng	100
87) Indeno(1,2,3-cd)pyrene	25.632	276	2600189	49.54	ng	100
88) Benzo(b)fluoranthene	22.762	252	2169007	49.27	ng	99
89) Benzo(k)fluoranthene	22.803	252	2169354	50.35	ng	99
90) Benzo(a)pyrene	23.321	252	2020931	49.38	ng	99
91) Dibenzo(a,h)anthracene	25.644	278	2256755	49.87	ng	99
92) Benzo(g,h,i)perylene	26.309	276	2045297	48.75	ng	99

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

