

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100621\
 Data File : BP007348.D
 Acq On : 06 Oct 2021 11:24
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

mohammad
 10/8/2021 8:34:18 AM

Quant Time: Oct 06 12:27:39 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP092121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 01 12:49:05 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.858	152	174028	20.00	ng	0.00
21) Naphthalene-d8	10.658	136	666060	20.00	ng	# 0.00
39) Acenaphthene-d10	14.498	164	430223	20.00	ng	0.00
64) Phenanthrene-d10	17.251	188	907150	20.00	ng	# 0.00
76) Chrysene-d12	21.345	240	920127	20.00	ng	0.00
86) Perylene-d12	23.745	264	982609	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.440	112	833209	80.14	ng	-0.01
7) Phenol-d6	7.040	99	1118616	78.72	ng	-0.01
23) Nitrobenzene-d5	9.022	82	942893	82.44	ng	0.00
42) 2,4,6-Tribromophenol	15.993	330	464885	83.35	ng	0.00
45) 2-Fluorobiphenyl	13.122	172	2341915	77.98	ng	0.00
79) Terphenyl-d14	19.810	244	3798341	79.10	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.328	88	169859	40.40	ng	# 88
3) Pyridine	3.740	79	463412	40.28	ng	# 86
4) n-Nitrosodimethylamine	3.652	42	180359	45.73	ng	# 89
6) Aniline	7.187	93	609668	38.93	ng	99
8) 2-Chlorophenol	7.428	128	437573	38.73	ng	98
9) Benzaldehyde	7.005	77	299432m	37.41	ng	
10) Phenol	7.069	94	532419	38.86	ng	93
11) bis(2-Chloroethyl)ether	7.287	93	411716	38.07	ng	95
12) 1,3-Dichlorobenzene	7.746	146	491430	38.56	ng	97
13) 1,4-Dichlorobenzene	7.893	146	499875	38.47	ng	98
14) 1,2-Dichlorobenzene	8.211	146	484784	38.50	ng	98
15) Benzyl Alcohol	8.111	79	365105	40.12	ng	95
16) 2,2'-oxybis(1-Chloropr...	8.393	45	473700	38.26	ng	90
17) 2-Methylphenol	8.316	107	362273	39.23	ng	94
18) Hexachloroethane	8.934	117	172601	39.57	ng	95
19) n-Nitroso-di-n-propyla...	8.675	70	298833	38.67	ng	99
20) 3+4-Methylphenols	8.646	107	500790	39.22	ng	94
22) Acetophenone	8.687	105	645729	40.23	ng	# 99
24) Nitrobenzene	9.069	77	448370	41.12	ng	96
25) Isophorone	9.599	82	822696	41.25	ng	98
26) 2-Nitrophenol	9.775	139	243737	42.87	ng	92
27) 2,4-Dimethylphenol	9.846	122	355088	39.66	ng	97
28) bis(2-Chloroethoxy)met...	10.075	93	539370	38.33	ng	99
29) 2,4-Dichlorophenol	10.316	162	422311	40.44	ng	98
30) 1,2,4-Trichlorobenzene	10.522	180	468239	39.36	ng	97
31) Naphthalene	10.710	128	1320672	38.85	ng	99
32) Benzoic acid	10.028	122	289270	42.04	ng	97
33) 4-Chloroaniline	10.828	127	565726	40.28	ng	99
34) Hexachlorobutadiene	10.987	225	286851	40.27	ng	98
35) Caprolactam	11.646	113	140911	43.86	ng	# 70
36) 4-Chloro-3-methylphenol	11.963	107	435936	41.07	ng	100
37) 2-Methylnaphthalene	12.322	142	925834	38.90	ng	97
38) 1-Methylnaphthalene	12.540	142	900915	38.74	ng	99
40) 1,2,4,5-Tetrachloroben...	12.693	216	517260	39.26	ng	98
41) Hexachlorocyclopentadiene	12.663	237	283431	38.82	ng	99
43) 2,4,6-Trichlorophenol	12.940	196	360748	40.13	ng	96

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.022	196	385174	39.79	ng	98
46) 1,1'-Biphenyl	13.334	154	1234533	38.58	ng	100
47) 2-Chloronaphthalene	13.375	162	968279	39.06	ng	98
48) 2-Nitroaniline	13.587	65	276304	45.72	ng	99
49) Acenaphthylene	14.222	152	1520875	39.83	ng	100
50) Dimethylphthalate	13.963	163	1231721	39.81	ng	100
51) 2,6-Dinitrotoluene	14.081	165	264244	41.81	ng	98
52) Acenaphthene	14.563	154	904154	39.09	ng	99
53) 3-Nitroaniline	14.416	138	291060	42.59	ng	99
54) 2,4-Dinitrophenol	14.640	184	143531	39.42	ng	94
55) Dibenzofuran	14.898	168	1511536	39.16	ng	96
56) 4-Nitrophenol	14.751	139	214292	38.62	ng	# 88
57) 2,4-Dinitrotoluene	14.875	165	364005	43.64	ng	98
58) Fluorene	15.545	166	1185938	39.78	ng	100
59) 2,3,4,6-Tetrachlorophenol	15.134	232	337126m	39.66	ng	
60) Diethylphthalate	15.322	149	1216184	40.56	ng	98
61) 4-Chlorophenyl-phenyle...	15.540	204	622040	39.48	ng	93
62) 4-Nitroaniline	15.581	138	309141	44.98	ng	# 86
63) Azobenzene	15.834	77	1077446	41.47	ng	96
65) 4,6-Dinitro-2-methylph...	15.645	198	226758	41.65	ng	96
66) n-Nitrosodiphenylamine	15.757	169	1047065	38.79	ng	99
67) 4-Bromophenyl-phenylether	16.440	248	388276	39.05	ng	95
68) Hexachlorobenzene	16.557	284	433188	39.42	ng	99
69) Atrazine	16.716	200	383719	40.33	ng	98
70) Pentachlorophenol	16.910	266	266460	39.05	ng	99
71) Phenanthrene	17.292	178	1884751	38.75	ng	99
72) Anthracene	17.387	178	1889066	39.70	ng	100
73) Carbazole	17.663	167	1881207	40.35	ng	98
74) Di-n-butylphthalate	18.204	149	2163655	41.87	ng	99
75) Fluoranthene	19.263	202	2272385	40.55	ng	98
77) Benzidine	19.445	184	1019462	36.05	ng	99
78) Pyrene	19.616	202	2352098	38.96	ng	99
80) Butylbenzylphthalate	20.486	149	1018532	42.13	ng	96
81) Benzo(a)anthracene	21.328	228	2351355	39.37	ng	99
82) 3,3'-Dichlorobenzidine	21.257	252	819291	39.94	ng	# 98
83) Chrysene	21.380	228	2219547	38.88	ng	98
84) Bis(2-ethylhexyl)phtha...	21.245	149	1444306	41.55	ng	98
85) Di-n-octyl phthalate	22.163	149	2503228	42.30	ng	100
87) Indeno(1,2,3-cd)pyrene	26.280	276	2822634	39.97	ng	# 96
88) Benzo(b)fluoranthene	23.016	252	2318353	39.48	ng	98
89) Benzo(k)fluoranthene	23.063	252	2254369	37.92	ng	99
90) Benzo(a)pyrene	23.639	252	1968342	39.45	ng	# 98
91) Dibenzo(a,h)anthracene	26.292	278	2347984	39.68	ng	# 97
92) Benzo(g,h,i)perylene	27.057	276	2265209	39.22	ng	# 95

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

