

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081721\
 Data File : BM031776.D
 Acq On : 17 Aug 2021 10:31
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDCCC040

Quant Time: Aug 17 12:17:43 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM081621.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 16 16:42:38 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.552	152	43792	20.000	ng	0.00
21) Naphthalene-d8	10.310	136	193668	20.000	ng	0.00
39) Acenaphthene-d10	14.187	164	139957	20.000	ng	0.00
64) Phenanthrene-d10	16.939	188	303148	20.000	ng	0.00
76) Chrysene-d12	21.139	240	259597	20.000	ng	0.00
86) Perylene-d12	23.286	264	208601	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.181	112	229952	84.278	ng	0.00
7) Phenol-d6	6.746	99	332142	84.009	ng	0.00
23) Nitrobenzene-d5	8.705	82	385138	83.267	ng	0.00
42) 2,4,6-Tribromophenol	15.686	330	146853	85.194	ng	0.00
45) 2-Fluorobiphenyl	12.798	172	875098	83.223	ng	0.00
79) Terphenyl-d14	19.586	244	1526927	99.102	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.193	88	46325	40.610	ng	92
3) Pyridine	3.570	79	126465	40.116	ng	# 84
4) n-Nitrosodimethylamine	3.493	42	76560	41.616	ng	87
6) Aniline	6.899	93	174858	40.448	ng	98
8) 2-Chlorophenol	7.122	128	134684	42.156	ng	97
9) Benzaldehyde	6.716	77	116225	41.418	ng	98
10) Phenol	6.769	94	165122	42.041	ng	88
11) bis(2-Chloroethyl)ether	6.999	93	120657	40.895	ng	94
12) 1,3-Dichlorobenzene	7.434	146	142458	41.686	ng	# 93
13) 1,4-Dichlorobenzene	7.587	146	145689	41.375	ng	97
14) 1,2-Dichlorobenzene	7.893	146	141812	41.001	ng	96
15) Benzyl Alcohol	7.799	79	141893	41.676	ng	95
16) 2,2'-oxybis(1-Chloropr...	8.081	45	168137	39.883	ng	97
17) 2-Methylphenol	7.999	107	116505	41.948	ng	95
18) Hexachloroethane	8.605	117	61368	42.045	ng	90
19) n-Nitroso-di-n-propyla...	8.363	70	128060	41.020	ng	95
20) 3+4-Methylphenols	8.328	107	163183	41.799	ng	95
22) Acetophenone	8.369	105	231273	42.056	ng	# 92
24) Nitrobenzene	8.746	77	194054	40.993	ng	98
25) Isophorone	9.269	82	340383	41.029	ng	97
26) 2-Nitrophenol	9.440	139	82234	44.130	ng	# 84
27) 2,4-Dimethylphenol	9.510	122	119649	41.990	ng	94
28) bis(2-Chloroethoxy)met...	9.746	93	166537	41.720	ng	98
29) 2,4-Dichlorophenol	9.969	162	140772	43.482	ng	95
30) 1,2,4-Trichlorobenzene	10.175	180	149078	43.123	ng	98
31) Naphthalene	10.357	128	469105	42.307	ng	99
32) Benzoic acid	9.681	122	95069	37.264	ng	94
33) 4-Chloroaniline	10.481	127	195734	42.469	ng	94
34) Hexachlorobutadiene	10.634	225	95607	43.313	ng	95
35) Caprolactam	11.304	113	48072	42.764	ng	87
36) 4-Chloro-3-methylphenol	11.616	107	168997	42.098	ng	91
37) 2-Methylnaphthalene	11.975	142	351007	42.302	ng	# 95
38) 1-Methylnaphthalene	12.198	142	334444	42.202	ng	# 100
40) 1,2,4,5-Tetrachloroben...	12.351	216	166688	41.929	ng	# 97
41) Hexachlorocyclopentadiene	12.322	237	78299	39.369	ng	98
43) 2,4,6-Trichlorophenol	12.604	196	124463	42.472	ng	98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.675	196	133710	41.751	ng	# 90
46) 1,1'-Biphenyl	13.010	154	454934	41.306	ng	99
47) 2-Chloronaphthalene	13.051	162	358841	41.501	ng	94
48) 2-Nitroaniline	13.275	65	134211	41.825	ng	88
49) Acenaphthylene	13.904	152	588737	41.590	ng	98
50) Dimethylphthalate	13.663	163	469059	40.221	ng	99
51) 2,6-Dinitrotoluene	13.787	165	106249	41.025	ng	# 77
52) Acenaphthene	14.251	154	356486	41.338	ng	99
53) 3-Nitroaniline	14.116	138	110464	41.540	ng	94
54) 2,4-Dinitrophenol	14.328	184	55505	36.355	ng	# 79
55) Dibenzofuran	14.592	168	597345	41.452	ng	92
56) 4-Nitrophenol	14.434	139	88694	39.101	ng	# 71
57) 2,4-Dinitrotoluene	14.581	165	156669	42.003	ng	# 60
58) Fluorene	15.245	166	505512	42.257	ng	99
59) 2,3,4,6-Tetrachlorophenol	14.822	232	120428	41.469	ng	# 86
60) Diethylphthalate	15.034	149	529687	41.322	ng	96
61) 4-Chlorophenyl-phenyle...	15.245	204	242433	41.983	ng	# 85
62) 4-Nitroaniline	15.286	138	113694	41.613	ng	# 79
63) Azobenzene	15.539	77	587767	41.422	ng	92
65) 4,6-Dinitro-2-methylph...	15.345	198	77157	37.231	ng	# 72
66) n-Nitrosodiphenylamine	15.469	169	424882	42.631	ng	99
67) 4-Bromophenyl-phenylether	16.139	248	137825	42.713	ng	# 89
68) Hexachlorobenzene	16.245	284	151325	42.871	ng	# 92
69) Atrazine	16.433	200	152325	42.947	ng	97
70) Pentachlorophenol	16.592	266	94723	39.952	ng	98
71) Phenanthrene	16.980	178	772817	41.800	ng	99
72) Anthracene	17.075	178	768089	42.377	ng	99
73) Carbazole	17.351	167	747587	41.172	ng	97
74) Di-n-butylphthalate	17.922	149	952702	42.124	ng	# 98
75) Fluoranthene	19.004	202	897212	40.515	ng	98
77) Benzidine	19.204	184	376252	48.772	ng	99
78) Pyrene	19.369	202	913475	48.573	ng	99
80) Butylbenzylphthalate	20.292	149	414868	46.028	ng	92
81) Benzo(a)anthracene	21.127	228	798704	42.004	ng	99
82) 3,3'-Dichlorobenzidine	21.068	252	236526	40.252	ng	99
83) Chrysene	21.174	228	769091	41.495	ng	98
84) Bis(2-ethylhexyl)phtha...	21.068	149	588223	42.231	ng	# 94
85) Di-n-octyl phthalate	21.927	149	875915	38.060	ng	# 94
87) Indeno(1,2,3-cd)pyrene	25.421	276	625382	42.542	ng	98
88) Benzo(b)fluoranthene	22.651	252	687538	43.898	ng	98
89) Benzo(k)fluoranthene	22.698	252	632723	41.855	ng	98
90) Benzo(a)pyrene	23.198	252	585968	42.256	ng	99
91) Dibenzo(a,h)anthracene	25.427	278	545420	42.187	ng	98
92) Benzo(g,h,i)perylene	26.062	276	511945	43.485	ng	98

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

