

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF100121\
 Data File : BF125743.D
 Acq On : 1 Oct 2021 9:53
 Operator : JU/CG
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

mohammad
 10/4/2021 11:53:21 AM

Quant Time: Oct 01 10:59:12 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF092421.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 24 15:38:32 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.893	152	251096	20.00	ng	0.00
21) Naphthalene-d8	8.175	136	987735	20.00	ng	# 0.00
39) Acenaphthene-d10	9.928	164	498321	20.00	ng	0.00
64) Phenanthrene-d10	11.410	188	776279	20.00	ng	# 0.00
76) Chrysene-d12	14.045	240	392693	20.00	ng	0.00
86) Perylene-d12	15.521	264	445028	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.504	112	1180280	77.44	ng	0.00
7) Phenol-d6	6.516	99	1563399	76.30	ng	0.00
23) Nitrobenzene-d5	7.451	82	1302609	91.99	ng	0.00
42) 2,4,6-Tribromophenol	10.716	330	400093	86.11	ng	0.00
45) 2-Fluorobiphenyl	9.251	172	2335285	84.06	ng	0.00
79) Terphenyl-d14	12.992	244	1817468	80.15	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.675	88	249750	38.51	ng	# 100
3) Pyridine	3.434	79	673723	39.27	ng	# 70
4) n-Nitrosodimethylamine	3.393	42	240006	37.70	ng	# 1
6) Aniline	6.551	93	907390	38.35	ng	94
8) 2-Chlorophenol	6.675	128	698631	40.41	ng	99
9) Benzaldehyde	6.440	77	402375	43.91	ng	94
10) Phenol	6.528	94	800558	38.65	ng	90
11) bis(2-Chloroethyl)ether	6.628	93	617583	38.37	ng	94
12) 1,3-Dichlorobenzene	6.834	146	725532m	38.01	ng	
13) 1,4-Dichlorobenzene	6.910	146	744304	38.32	ng	98
14) 1,2-Dichlorobenzene	7.063	146	693888	37.87	ng	99
15) Benzyl Alcohol	7.028	79	544826	38.33	ng	93
16) 2,2'-oxybis(1-Chloropr...	7.163	45	769921	37.58	ng	87
17) 2-Methylphenol	7.140	107	549077	38.73	ng	97
18) Hexachloroethane	7.404	117	257069	40.78	ng	92
19) n-Nitroso-di-n-propyla...	7.304	70	426340	36.55	ng	# 67
20) 3+4-Methylphenols	7.292	107	659446	37.00	ng	96
22) Acetophenone	7.298	105	844898	37.70	ng	# 91
24) Nitrobenzene	7.475	77	668450	46.40	ng	95
25) Isophorone	7.710	82	1210105	38.16	ng	94
26) 2-Nitrophenol	7.787	139	359881	47.67	ng	93
27) 2,4-Dimethylphenol	7.822	122	566835	40.19	ng	97
28) bis(2-Chloroethoxy)met...	7.916	93	706167	39.48	ng	98
29) 2,4-Dichlorophenol	8.028	162	588500	42.00	ng	99
30) 1,2,4-Trichlorobenzene	8.116	180	600507	39.37	ng	97
31) Naphthalene	8.198	128	1930415	38.06	ng	99
32) Benzoic acid	7.934	122	409750	41.14	ng	99
33) 4-Chloroaniline	8.239	127	806046	37.89	ng	98
34) Hexachlorobutadiene	8.316	225	334465	39.77	ng	100
35) Caprolactam	8.616	113	167480	36.63	ng	# 75
36) 4-Chloro-3-methylphenol	8.716	107	569509	38.79	ng	99
37) 2-Methylnaphthalene	8.886	142	1299287	37.77	ng	100
38) 1-Methylnaphthalene	8.986	142	1221028	37.56	ng	98
40) 1,2,4,5-Tetrachloroben...	9.051	216	534425	41.24	ng	98
41) Hexachlorocyclopentadiene	9.039	237	248801	45.37	ng	99
43) 2,4,6-Trichlorophenol	9.157	196	403467	44.39	ng	98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.198	196	436044	44.44	ng	92
46) 1,1'-Biphenyl	9.351	154	1456698	39.64	ng	98
47) 2-Chloronaphthalene	9.375	162	1212636	39.86	ng	98
48) 2-Nitroaniline	9.463	65	317123	44.52	ng	95
49) Acenaphthylene	9.792	152	1797456	39.25	ng	98
50) Dimethylphthalate	9.651	163	1331640	39.67	ng	98
51) 2,6-Dinitrotoluene	9.710	165	313181	45.98	ng	# 85
52) Acenaphthene	9.963	154	1136846	39.59	ng	98
53) 3-Nitroaniline	9.875	138	327785	42.39	ng	89
54) 2,4-Dinitrophenol	9.981	184	118844	46.86	ng	# 68
55) Dibenzofuran	10.133	168	1650798	38.32	ng	95
56) 4-Nitrophenol	10.028	139	244935	39.61	ng	# 83
57) 2,4-Dinitrotoluene	10.110	165	392245	46.19	ng	# 82
58) Fluorene	10.475	166	1212707	36.17	ng	98
59) 2,3,4,6-Tetrachlorophenol	10.251	232	323638	42.86	ng	# 87
60) Diethylphthalate	10.345	149	1223977	37.82	ng	99
61) 4-Chlorophenyl-phenyle...	10.469	204	549248	36.85	ng	87
62) 4-Nitroaniline	10.486	138	291004	38.44	ng	# 76
63) Azobenzene	10.628	77	1178320	37.71	ng	# 83
65) 4,6-Dinitro-2-methylph...	10.516	198	183011	48.35	ng	# 80
66) n-Nitrosodiphenylamine	10.586	169	1100068	41.28	ng	96
67) 4-Bromophenyl-phenylether	10.957	248	346615	42.55	ng	96
68) Hexachlorobenzene	11.028	284	396872	41.51	ng	# 80
69) Atrazine	11.110	200	302039	42.75	ng	92
70) Pentachlorophenol	11.216	266	226780	44.88	ng	95
71) Phenanthrene	11.439	178	1670252	38.37	ng	97
72) Anthracene	11.486	178	1666948	38.75	ng	99
73) Carbazole	11.639	167	1551418	36.65	ng	99
74) Di-n-butylphthalate	11.969	149	1592047	37.30	ng	99
75) Fluoranthene	12.622	202	1494808	32.13	ng	96
77) Benzidine	12.739	184	522212	50.62	ng	97
78) Pyrene	12.851	202	1485650	41.86	ng	96
80) Butylbenzylphthalate	13.469	149	479433	40.80	ng	95
81) Benzo(a)anthracene	14.039	228	1056019	39.59	ng	100
82) 3,3'-Dichlorobenzidine	13.998	252	334232	43.42	ng	98
83) Chrysene	14.074	228	1049459	39.20	ng	96
84) Bis(2-ethylhexyl)phtha...	14.027	149	549718	40.86	ng	# 98
85) Di-n-octyl phthalate	14.639	149	1007528	51.06	ng	# 88
87) Indeno(1,2,3-cd)pyrene	17.015	276	1301537	43.08	ng	97
88) Benzo(b)fluoranthene	15.092	252	1122727	36.01	ng	98
89) Benzo(k)fluoranthene	15.121	252	1099752m	36.81	ng	
90) Benzo(a)pyrene	15.463	252	1071383	38.60	ng	97
91) Dibenzo(a,h)anthracene	17.033	278	1080199	42.78	ng	98
92) Benzo(g,h,i)perylene	17.462	276	1121951	43.51	ng	93

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

