

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF122821\
 Data File : BF126378.D
 Acq On : 28 Dec 2021 13:07
 Operator : JU/CG
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Quant Time: Dec 28 14:59:05 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF121521.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 22 13:00:11 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.810	152	147822	20.000	ng	0.00
21) Naphthalene-d8	8.092	136	528888	20.000	ng	#-0.01
39) Acenaphthene-d10	9.851	164	213848	20.000	ng	0.00
64) Phenanthrene-d10	11.333	188	355594	20.000	ng	# 0.00
76) Chrysene-d12	13.974	240	173430	20.000	ng	0.00
86) Perylene-d12	15.421	264	220185	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.416	112	645443	64.301	ng	-0.01
7) Phenol-d6	6.440	99	981498	77.007	ng	0.00
23) Nitrobenzene-d5	7.375	82	828357	84.825	ng	0.00
42) 2,4,6-Tribromophenol	10.639	330	169452	98.600	ng	0.00
45) 2-Fluorobiphenyl	9.175	172	1085277	89.009	ng	0.00
79) Terphenyl-d14	12.922	244	873185	87.511	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.528	88	210815	43.025	ng	# 100
3) Pyridine	3.257	79	547704	42.092	ng	# 80
4) n-Nitrosodimethylamine	3.205	42	195143	39.021	ng	# 1
6) Aniline	6.475	93	630056	38.917	ng	96
8) 2-Chlorophenol	6.598	128	424180	41.671	ng	92
9) Benzaldehyde	6.357	77	259542	33.746	ng	94
10) Phenol	6.457	94	541230	37.830	ng	93
11) bis(2-Chloroethyl)ether	6.545	93	453825	40.866	ng	96
12) 1,3-Dichlorobenzene	6.751	146	424534	41.345	ng	97
13) 1,4-Dichlorobenzene	6.828	146	425897	41.073	ng	95
14) 1,2-Dichlorobenzene	6.987	146	341486	35.618	ng	97
15) Benzyl Alcohol	6.951	79	314952	33.849	ng	95
16) 2,2'-oxybis(1-Chloropr...	7.093	45	650639	35.802	ng	93
17) 2-Methylphenol	7.063	107	320043	35.890	ng	97
18) Hexachloroethane	7.328	117	151884	37.238	ng	98
19) n-Nitroso-di-n-propyla...	7.228	70	263768	33.711	ng	# 73
20) 3+4-Methylphenols	7.216	107	376342	35.699	ng	# 85
22) Acetophenone	7.222	105	493552	40.548	ng	93
24) Nitrobenzene	7.393	77	415816	42.654	ng	89
25) Isophorone	7.634	82	721998	39.152	ng	# 87
26) 2-Nitrophenol	7.710	139	196960	44.818	ng	# 82
27) 2,4-Dimethylphenol	7.751	122	299820	40.173	ng	97
28) bis(2-Chloroethoxy)met...	7.845	93	486282	40.108	ng	# 97
29) 2,4-Dichlorophenol	7.951	162	285236	42.996	ng	92
30) 1,2,4-Trichlorobenzene	8.040	180	287850	41.949	ng	99
31) Naphthalene	8.116	128	1046484	42.310	ng	98
32) Benzoic acid	7.881	122	216591	33.687	ng	91
33) 4-Chloroaniline	8.163	127	434663	42.089	ng	# 95
34) Hexachlorobutadiene	8.240	225	148051	43.930	ng	98
35) Caprolactam	8.528	113	67828	41.140	ng	86
36) 4-Chloro-3-methylphenol	8.645	107	326967	41.782	ng	91
37) 2-Methylnaphthalene	8.810	142	628923	41.290	ng	99
38) 1-Methylnaphthalene	8.910	142	609288	41.218	ng	95
40) 1,2,4,5-Tetrachloroben...	8.975	216	225283	48.811	ng	# 96
41) Hexachlorocyclopentadiene	8.963	237	135996	51.455	ng	95
43) 2,4,6-Trichlorophenol	9.087	196	170298	46.905	ng	92

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.128	196	182928	48.653	ng	99
46) 1,1'-Biphenyl	9.275	154	746721	47.225	ng	97
47) 2-Chloronaphthalene	9.298	162	560450	47.109	ng	97
48) 2-Nitroaniline	9.392	65	206284	47.505	ng	91
49) Acenaphthylene	9.710	152	798385	43.919	ng	98
50) Dimethylphthalate	9.575	163	612944	45.248	ng	98
51) 2,6-Dinitrotoluene	9.634	165	129839	44.283	ng	89
52) Acenaphthene	9.886	154	534016	45.298	ng	98
53) 3-Nitroaniline	9.804	138	153315	41.140	ng #	79
54) 2,4-Dinitrophenol	9.910	184	66719	55.265	ng #	77
55) Dibenzofuran	10.057	168	716160	44.619	ng	99
56) 4-Nitrophenol	9.963	139	129500	48.108	ng #	82
57) 2,4-Dinitrotoluene	10.039	165	168508	45.054	ng #	95
58) Fluorene	10.398	166	515346	44.351	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.175	232	128495	46.092	ng #	98
60) Diethylphthalate	10.275	149	574053	42.169	ng	98
61) 4-Chlorophenyl-phenyle...	10.392	204	235185	44.953	ng	94
62) 4-Nitroaniline	10.416	138	157181	50.279	ng #	75
63) Azobenzene	10.551	77	646654	39.218	ng	85
65) 4,6-Dinitro-2-methylph...	10.445	198	95387	49.124	ng	93
66) n-Nitrosodiphenylamine	10.510	169	476876	41.786	ng	98
67) 4-Bromophenyl-phenylether	10.881	248	134251	40.227	ng	93
68) Hexachlorobenzene	10.951	284	152706	39.785	ng #	82
69) Atrazine	11.033	200	82169	39.289	ng	96
70) Pentachlorophenol	11.139	266	82472	38.730	ng	93
71) Phenanthrene	11.363	178	770380	41.975	ng	98
72) Anthracene	11.410	178	781770	42.438	ng	99
73) Carbazole	11.563	167	725981	41.847	ng	100
74) Di-n-butylphthalate	11.898	149	899354	40.894	ng	100
75) Fluoranthene	12.545	202	717037	43.362	ng	91
77) Benzidine	12.669	184	152252	55.684	ng	98
78) Pyrene	12.775	202	713469	44.778	ng	96
80) Butylbenzylphthalate	13.398	149	256237	34.174	ng #	83
81) Benzo(a)anthracene	13.963	228	414766	40.374	ng	99
82) 3,3'-Dichlorobenzidine	13.927	252	198573	42.310	ng	99
83) Chrysene	14.004	228	425061	39.823	ng	98
84) Bis(2-ethylhexyl)phtha...	13.963	149	288265	34.405	ng	100
85) Di-n-octyl phthalate	14.574	149	556931	37.287	ng	97
87) Indeno(1,2,3-cd)pyrene	16.857	276	671198	45.701	ng	94
88) Benzo(b)fluoranthene	15.004	252	476325	36.004	ng	98
89) Benzo(k)fluoranthene	15.033	252	492976	38.952	ng #	95
90) Benzo(a)pyrene	15.363	252	459355	41.776	ng #	96
91) Dibenzo(a,h)anthracene	16.874	278	548730	45.962	ng	95
92) Benzo(g,h,i)perylene	17.292	276	578401	46.825	ng	93

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

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