

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061021\
 Data File : BF124225.D
 Acq On : 10 Jun 2021 9:48
 Operator : JU/CG
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Quant Time: Jun 10 11:31:29 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF060321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 10 01:53:56 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.851	152	36999	20.000	ng	0.00
21) Naphthalene-d8	8.134	136	133442	20.000	ng	0.00
39) Acenaphthene-d10	9.892	164	78062	20.000	ng	0.00
64) Phenanthrene-d10	11.380	188	146599	20.000	ng	0.00
76) Chrysene-d12	14.021	240	90193	20.000	ng	0.00
86) Perylene-d12	15.498	264	64326	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.481	112	184992	75.266	ng	0.00
7) Phenol-d6	6.498	99	245377	74.273	ng	0.00
23) Nitrobenzene-d5	7.422	82	261459	77.710	ng	0.00
42) 2,4,6-Tribromophenol	10.686	330	88856	80.344	ng	0.00
45) 2-Fluorobiphenyl	9.210	172	493250	79.512	ng	0.00
79) Terphenyl-d14	12.963	244	527598	80.313	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.599	88	38500	35.793	ng	95
3) Pyridine	3.363	79	97978	35.567	ng	86
4) n-Nitrosodimethylamine	3.316	42	56126	34.289	ng	# 95
6) Aniline	6.516	93	134064	35.528	ng	# 44
8) 2-Chlorophenol	6.645	128	99066	36.223	ng	92
9) Benzaldehyde	6.404	77	61913	42.161	ng	87
10) Phenol	6.510	94	127109	35.601	ng	# 64
11) bis(2-Chloroethyl)ether	6.587	93	95135	36.551	ng	97
12) 1,3-Dichlorobenzene	6.792	146	120349	37.642	ng	97
13) 1,4-Dichlorobenzene	6.869	146	119622	37.301	ng	98
14) 1,2-Dichlorobenzene	7.022	146	115758	37.732	ng	96
15) Benzyl Alcohol	6.998	79	112410	37.743	ng	94
16) 2,2'-oxybis(1-Chloropr...	7.128	45	135159	33.287	ng	71
17) 2-Methylphenol	7.116	107	84052	37.806	ng	# 87
18) Hexachloroethane	7.363	117	47898	37.926	ng	# 87
19) n-Nitroso-di-n-propyla...	7.269	70	86273	36.993	ng	# 85
20) 3+4-Methylphenols	7.269	107	107968	36.651	ng	90
22) Acetophenone	7.263	105	152723	40.147	ng	92
24) Nitrobenzene	7.439	77	128471	38.062	ng	95
25) Isophorone	7.675	82	222565	36.709	ng	97
26) 2-Nitrophenol	7.751	139	59211	39.271	ng	94
27) 2,4-Dimethylphenol	7.792	122	78662	36.999	ng	97
28) bis(2-Chloroethoxy)met...	7.881	93	112032	37.272	ng	99
29) 2,4-Dichlorophenol	7.998	162	97189	39.506	ng	93
30) 1,2,4-Trichlorobenzene	8.075	180	107015	38.816	ng	98
31) Naphthalene	8.157	128	294871	36.960	ng	98
32) Benzoic acid	7.928	122	47164	36.152	ng	97
33) 4-Chloroaniline	8.210	127	120099	40.472	ng	# 92
34) Hexachlorobutadiene	8.275	225	74772	38.396	ng	96
35) Caprolactam	8.586	113	26316	38.938	ng	# 73
36) 4-Chloro-3-methylphenol	8.698	107	99425	36.940	ng	# 87
37) 2-Methylnaphthalene	8.851	142	212403	37.965	ng	97
38) 1-Methylnaphthalene	8.951	142	202037	38.643	ng	97
40) 1,2,4,5-Tetrachloroben...	9.016	216	112832	40.796	ng	97
41) Hexachlorocyclopentadiene	9.004	237	44911	34.760	ng	99
43) 2,4,6-Trichlorophenol	9.133	196	69957	38.782	ng	97

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.181	196	74293	38.366	ng	# 90
46) 1,1'-Biphenyl	9.316	154	262477	39.832	ng	99
47) 2-Chloronaphthalene	9.339	162	204326	38.094	ng	95
48) 2-Nitroaniline	9.439	65	77716	40.150	ng	99
49) Acenaphthylene	9.757	152	303028	38.867	ng	99
50) Dimethylphthalate	9.616	163	245461	38.006	ng	98
51) 2,6-Dinitrotoluene	9.681	165	55943	37.778	ng	86
52) Acenaphthene	9.928	154	196889	38.665	ng	95
53) 3-Nitroaniline	9.851	138	51570	37.209	ng	95
54) 2,4-Dinitrophenol	9.963	184	30255	41.718	ng	# 84
55) Dibenzofuran	10.098	168	309590	38.845	ng	97
56) 4-Nitrophenol	10.028	139	47143	47.596	ng	86
57) 2,4-Dinitrotoluene	10.086	165	75981	38.905	ng	# 93
58) Fluorene	10.445	166	241167	39.421	ng	94
59) 2,3,4,6-Tetrachlorophenol	10.222	232	61050	38.905	ng	96
60) Diethylphthalate	10.310	149	255799	39.246	ng	99
61) 4-Chlorophenyl-phenyle...	10.433	204	119785	38.293	ng	97
62) 4-Nitroaniline	10.469	138	54895	39.393	ng	87
63) Azobenzene	10.592	77	251890	37.074	ng	97
65) 4,6-Dinitro-2-methylph...	10.498	198	46825	39.746	ng	# 74
66) n-Nitrosodiphenylamine	10.551	169	212217	38.099	ng	98
67) 4-Bromophenyl-phenylether	10.922	248	77455	38.493	ng	93
68) Hexachlorobenzene	10.992	284	83952	36.836	ng	96
69) Atrazine	11.080	200	69291	43.945	ng	96
70) Pentachlorophenol	11.192	266	40912	36.743	ng	92
71) Phenanthrene	11.404	178	350321	38.254	ng	98
72) Anthracene	11.457	178	343062	37.197	ng	100
73) Carbazole	11.616	167	301103	37.464	ng	95
74) Di-n-butylphthalate	11.939	149	399520	38.680	ng	99
75) Fluoranthene	12.598	202	350913	36.505	ng	95
77) Benzidine	12.710	184	85586	41.332	ng	# 93
78) Pyrene	12.827	202	346647	40.025	ng	98
80) Butylbenzylphthalate	13.433	149	137858	39.238	ng	99
81) Benzo(a)anthracene	14.010	228	245545	37.371	ng	98
82) 3,3'-Dichlorobenzidine	13.968	252	73045	37.895	ng	98
83) Chrysene	14.051	228	238969	37.696	ng	98
84) Bis(2-ethylhexyl)phtha...	13.992	149	156745	37.505	ng	97
85) Di-n-octyl phthalate	14.604	149	211353	37.893	ng	# 92
87) Indeno(1,2,3-cd)pyrene	16.992	276	206131	39.360	ng	98
88) Benzo(b)fluoranthene	15.068	252	187251	37.473	ng	98
89) Benzo(k)fluoranthene	15.098	252	182348	38.548	ng	97
90) Benzo(a)pyrene	15.439	252	169132	38.969	ng	98
91) Dibenzo(a,h)anthracene	17.004	278	170635	38.274	ng	98
92) Benzo(g,h,i)perylene	17.445	276	166682	37.842	ng	98

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

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