

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF071421\
 Data File : BF124567.D
 Acq On : 14 Jul 2021 16:59
 Operator : JU/CG
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040EC

Quant Time: Jul 14 17:25:41 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF070721.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 14 13:25:43 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.928	152	7132	20.000	ng	0.00
21) Naphthalene-d8	8.210	136	27553	20.000	ng	0.00
39) Acenaphthene-d10	9.969	164	15208	20.000	ng	0.00
64) Phenanthrene-d10	11.457	188	27426	20.000	ng	0.00
76) Chrysene-d12	14.098	240	18168	20.000	ng	0.00
86) Perylene-d12	15.592	264	14270	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.545	112	34179	83.031	ng	0.00
7) Phenol-d6	6.551	99	42502	84.874	ng	0.00
23) Nitrobenzene-d5	7.492	82	35803	82.029	ng	0.00
42) 2,4,6-Tribromophenol	10.763	330	9919	83.813	ng	0.00
45) 2-Fluorobiphenyl	9.292	172	83924	80.472	ng	0.00
79) Terphenyl-d14	13.045	244	90554	83.696	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.799	88	6047	44.911	ng	# 100
3) Pyridine	3.546	79	14945	43.277	ng	94
4) n-Nitrosodimethylamine	3.516	42	12557	44.737	ng	# 95
6) Aniline	6.598	93	21685	41.099	ng	# 96
8) 2-Chlorophenol	6.716	128	19482	42.215	ng	97
9) Benzaldehyde	6.481	77	10570	35.979	ng	# 87
10) Phenol	6.563	94	20073	41.057	ng	97
11) bis(2-Chloroethyl)ether	6.663	93	15378	40.571	ng	97
12) 1,3-Dichlorobenzene	6.869	146	21920	43.397	ng	97
13) 1,4-Dichlorobenzene	6.945	146	22715	44.029	ng	97
14) 1,2-Dichlorobenzene	7.098	146	20654	42.126	ng	97
15) Benzyl Alcohol	7.069	79	14137	40.800	ng	88
16) 2,2'-oxybis(1-Chloropr...	7.204	45	28088	42.994	ng	97
17) 2-Methylphenol	7.169	107	13982	41.392	ng	97
18) Hexachloroethane	7.445	117	8076	41.569	ng	87
19) n-Nitroso-di-n-propyla...	7.345	70	11786	41.862	ng	98
20) 3+4-Methylphenols	7.328	107	18530	41.665	ng	92
22) Acetophenone	7.339	105	25542	40.847	ng	# 96
24) Nitrobenzene	7.510	77	17841	41.223	ng	96
25) Isophorone	7.751	82	31873	39.779	ng	98
26) 2-Nitrophenol	7.828	139	10133	43.403	ng	97
27) 2,4-Dimethylphenol	7.857	122	16115	41.656	ng	97
28) bis(2-Chloroethoxy)met...	7.957	93	18492	39.565	ng	98
29) 2,4-Dichlorophenol	8.063	162	16560	41.501	ng	99
30) 1,2,4-Trichlorobenzene	8.151	180	18063	40.905	ng	99
31) Naphthalene	8.233	128	58769	41.239	ng	99
32) Benzoic acid	7.981	122	12396	45.973	ng	97
33) 4-Chloroaniline	8.281	127	21961	40.780	ng	98
34) Hexachlorobutadiene	8.351	225	10546	42.311	ng	95
35) Caprolactam	8.657	113	3807	37.719	ng	85
36) 4-Chloro-3-methylphenol	8.751	107	16044	40.660	ng	99
37) 2-Methylnaphthalene	8.928	142	39111	40.671	ng	96
38) 1-Methylnaphthalene	9.028	142	36456	40.366	ng	100
40) 1,2,4,5-Tetrachloroben...	9.092	216	16686	41.137	ng	97
41) Hexachlorocyclopentadiene	9.080	237	10437	41.192	ng	99
43) 2,4,6-Trichlorophenol	9.198	196	12368	42.236	ng	95

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.233	196	12315	40.776	ng	95
46) 1,1'-Biphenyl	9.392	154	47136	40.809	ng	99
47) 2-Chloronaphthalene	9.416	162	35995	39.747	ng	97
48) 2-Nitroaniline	9.510	65	9682	44.459	ng	92
49) Acenaphthylene	9.833	152	57221	40.490	ng	99
50) Dimethylphthalate	9.692	163	43090	40.724	ng	98
51) 2,6-Dinitrotoluene	9.751	165	9404	45.084	ng	96
52) Acenaphthene	10.004	154	37461	41.456	ng	98
53) 3-Nitroaniline	9.922	138	10075	43.285	ng	99
54) 2,4-Dinitrophenol	10.022	184	4939	45.378	ng	87
55) Dibenzofuran	10.175	168	53882	40.555	ng	99
56) 4-Nitrophenol	10.069	139	8561	44.191	ng	97
57) 2,4-Dinitrotoluene	10.151	165	13204	46.238	ng #	97
58) Fluorene	10.522	166	40780	39.682	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.286	232	9756	41.827	ng #	98
60) Diethylphthalate	10.392	149	44207	39.959	ng	99
61) 4-Chlorophenyl-phenyle...	10.510	204	19165	41.086	ng	98
62) 4-Nitroaniline	10.539	138	10297	43.894	ng	90
63) Azobenzene	10.669	77	38441	39.751	ng	97
65) 4,6-Dinitro-2-methylph...	10.563	198	6565	41.903	ng	92
66) n-Nitrosodiphenylamine	10.627	169	35994	40.305	ng	99
67) 4-Bromophenyl-phenylether	11.004	248	11301	39.849	ng	97
68) Hexachlorobenzene	11.063	284	11767	40.498	ng	90
69) Atrazine	11.157	200	11205	40.968	ng	94
70) Pentachlorophenol	11.257	266	7819	42.544	ng	92
71) Phenanthrene	11.486	178	59944	40.062	ng	97
72) Anthracene	11.533	178	59350	39.538	ng	99
73) Carbazole	11.686	167	55732	40.181	ng	99
74) Di-n-butylphthalate	12.016	149	75428	40.369	ng	99
75) Fluoranthene	12.674	202	61481	40.726	ng	96
77) Benzidine	12.792	184	27136	39.953	ng	99
78) Pyrene	12.904	202	62693	41.534	ng	98
80) Butylbenzylphthalate	13.515	149	29897	40.964	ng	94
81) Benzo(a)anthracene	14.086	228	49838	41.363	ng	99
82) 3,3'-Dichlorobenzidine	14.051	252	13418	42.514	ng	99
83) Chrysene	14.127	228	48172	41.734	ng	98
84) Bis(2-ethylhexyl)phtha...	14.074	149	40939	41.780	ng	98
85) Di-n-octyl phthalate	14.692	149	60979	43.007	ng	100
87) Indeno(1,2,3-cd)pyrene	17.109	276	35451	42.836	ng	97
88) Benzo(b)fluoranthene	15.151	252	41472	40.490	ng	100
89) Benzo(k)fluoranthene	15.186	252	37370	39.765	ng #	97
90) Benzo(a)pyrene	15.533	252	34581	40.901	ng	97
91) Dibenzo(a,h)anthracene	17.133	278	30898	43.862	ng	97
92) Benzo(g,h,i)perylene	17.574	276	29433	42.795	ng	93

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

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