

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF090424\
 Data File : BF139238.D
 Acq On : 04 Sep 2024 13:36
 Operator : RC/JU
 Sample : SSTDICCC040
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICCC040

Quant Time: Sep 04 15:47:19 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF090424.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 04 15:37:35 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.893	152	130545	20.000	ng	0.00
21) Naphthalene-d8	8.175	136	471691	20.000	ng	0.00
39) Acenaphthene-d10	9.928	164	256219	20.000	ng	0.00
64) Phenanthrene-d10	11.416	188	463513	20.000	ng	0.00
76) Chrysene-d12	14.051	240	215446	20.000	ng	0.00
86) Perylene-d12	15.527	264	225523	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.504	112	597448	76.560	ng	0.00
7) Phenol-d6	6.516	99	805982	77.366	ng	0.00
23) Nitrobenzene-d5	7.451	82	711055	78.036	ng	0.00
42) 2,4,6-Tribromophenol	10.716	330	173686	78.502	ng	0.00
45) 2-Fluorobiphenyl	9.251	172	1244011	74.948	ng	0.00
79) Terphenyl-d14	12.998	244	1057635	79.261	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.675	88	142095	39.101	ng	100
3) Pyridine	3.428	79	339786	38.735	ng	100
4) n-Nitrosodimethylamine	3.387	42	224145	39.348	ng	100
6) Aniline	6.551	93	370740	38.700	ng	100
8) 2-Chlorophenol	6.675	128	308088	38.094	ng	100
9) Benzaldehyde	6.440	77	223693	35.473	ng	100
10) Phenol	6.528	94	425350	38.708	ng	100
11) bis(2-Chloroethyl)ether	6.628	93	301241	38.265	ng	100
12) 1,3-Dichlorobenzene	6.834	146	350915	38.334	ng	100
13) 1,4-Dichlorobenzene	6.910	146	352882	38.514	ng	100
14) 1,2-Dichlorobenzene	7.063	146	330177	38.638	ng	100
15) Benzyl Alcohol	7.028	79	297767	39.193	ng	100
16) 2,2'-oxybis(1-Chloropr...	7.163	45	584261	38.986	ng	100
17) 2-Methylphenol	7.140	107	254485	38.521	ng	100
18) Hexachloroethane	7.404	117	125952	38.489	ng	100
19) n-Nitroso-di-n-propyla...	7.304	70	229707	37.956	ng	100
20) 3+4-Methylphenols	7.293	107	323601	38.866	ng	100
22) Acetophenone	7.298	105	433365	38.565	ng	100
24) Nitrobenzene	7.475	77	374579	38.923	ng	100
25) Isophorone	7.710	82	604293	38.983	ng	100
26) 2-Nitrophenol	7.787	139	154758	40.944	ng	100
27) 2,4-Dimethylphenol	7.822	122	212822	39.480	ng	100
28) bis(2-Chloroethoxy)met...	7.922	93	356208	38.569	ng	100
29) 2,4-Dichlorophenol	8.028	162	254754	39.470	ng	100
30) 1,2,4-Trichlorobenzene	8.110	180	286730	38.897	ng	100
31) Naphthalene	8.198	128	926389	38.782	ng	100
32) Benzoic acid	7.928	122	192798	39.342	ng	100
33) 4-Chloroaniline	8.240	127	315728	38.173	ng	100
34) Hexachlorobutadiene	8.310	225	179696	38.470	ng	100
35) Caprolactam	8.610	113	80380	40.227	ng	100
36) 4-Chloro-3-methylphenol	8.716	107	277907	39.512	ng	100
37) 2-Methylnaphthalene	8.887	142	574761	38.459	ng	100
38) 1-Methylnaphthalene	8.987	142	568379	38.670	ng	100
40) 1,2,4,5-Tetrachloroben...	9.051	216	276336	38.120	ng	100
41) Hexachlorocyclopentadiene	9.040	237	98261	42.178	ng	100
43) 2,4,6-Trichlorophenol	9.163	196	191881	40.455	ng	100

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF090424\
 Data File : BF139238.D
 Acq On : 04 Sep 2024 13:36
 Operator : RC/JU
 Sample : SSTDICCC040
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICCC040

Quant Time: Sep 04 15:47:19 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF090424.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 04 15:37:35 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.198	196	194088	39.050	ng	100
46) 1,1'-Biphenyl	9.351	154	729306	38.418	ng	100
47) 2-Chloronaphthalene	9.375	162	547186	38.174	ng	100
48) 2-Nitroaniline	9.469	65	197640	40.031	ng	100
49) Acenaphthylene	9.792	152	831934	38.535	ng	100
50) Dimethylphthalate	9.651	163	595907	38.373	ng	100
51) 2,6-Dinitrotoluene	9.710	165	142139	40.346	ng	100
52) Acenaphthene	9.963	154	556791	38.884	ng	100
53) 3-Nitroaniline	9.881	138	149592	39.853	ng	100
54) 2,4-Dinitrophenol	9.981	184	62228	38.067	ng	100
55) Dibenzofuran	10.134	168	787591	38.364	ng	100
56) 4-Nitrophenol	10.028	139	115336	41.630	ng	100
57) 2,4-Dinitrotoluene	10.110	165	184677	41.185	ng	100
58) Fluorene	10.475	166	610821	38.375	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.251	232	153801	39.311	ng	100
60) Diethylphthalate	10.351	149	595972	38.997	ng	100
61) 4-Chlorophenyl-phenyle...	10.469	204	297041	38.697	ng	100
62) 4-Nitroaniline	10.492	138	145015	40.586	ng	100
63) Azobenzene	10.628	77	621547	38.588	ng	100
65) 4,6-Dinitro-2-methylph...	10.516	198	88814	41.469	ng	100
66) n-Nitrosodiphenylamine	10.586	169	516265	38.264	ng	100
67) 4-Bromophenyl-phenylether	10.957	248	175568	39.040	ng	100
68) Hexachlorobenzene	11.022	284	197042	38.274	ng	100
69) Atrazine	11.110	200	115561	37.124	ng	100
70) Pentachlorophenol	11.216	266	126074	42.394	ng	100
71) Phenanthrene	11.439	178	824396	38.072	ng	100
72) Anthracene	11.492	178	822280	38.874	ng	100
73) Carbazole	11.645	167	765295	38.710	ng	100
74) Di-n-butylphthalate	11.975	149	818910	40.455	ng	100
75) Fluoranthene	12.628	202	834957	38.462	ng	100
77) Benzidine	12.751	184	187373	43.193	ng	100
78) Pyrene	12.857	202	836549	40.195	ng	100
80) Butylbenzylphthalate	13.475	149	209437	42.747	ng	100
81) Benzo(a)anthracene	14.039	228	525882	38.226	ng	100
82) 3,3'-Dichlorobenzidine	14.004	252	149497	40.919	ng	100
83) Chrysene	14.080	228	514546	40.074	ng	100
84) Bis(2-ethylhexyl)phtha...	14.033	149	248996	43.299	ng	100
85) Di-n-octyl phthalate	14.651	149	461517	41.919	ng	100
87) Indeno(1,2,3-cd)pyrene	17.010	276	566838	40.233	ng	100
88) Benzo(b)fluoranthene	15.098	252	516473	39.766	ng	100
89) Benzo(k)fluoranthene	15.127	252	460455	39.540	ng	100
90) Benzo(a)pyrene	15.463	252	444040	40.263	ng	100
91) Dibenzo(a,h)anthracene	17.033	278	472304	40.679	ng	100
92) Benzo(g,h,i)perylene	17.462	276	479258	39.990	ng	100

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

