

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF020625\
 Data File : BF141480.D
 Acq On : 06 Feb 2025 15:03
 Operator : RC/JU
 Sample : SSTDICV040
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ICVBF020625

Quant Time: Feb 06 16:01:10 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF020625.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 06 15:49:10 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.787	152	90009	20.000	ng	0.00
21) Naphthalene-d8	8.069	136	362467	20.000	ng	0.00
39) Acenaphthene-d10	9.828	164	197991	20.000	ng	0.00
64) Phenanthrene-d10	11.316	188	337289	20.000	ng	0.00
76) Chrysene-d12	13.963	240	252530	20.000	ng	0.00
86) Perylene-d12	15.427	264	199389	20.000	ng	0.01

System Monitoring Compounds						
5) 2-Fluorophenol	5.404	112	470176	81.894	ng	0.00
7) Phenol-d6	6.428	99	586892	80.878	ng	-0.01
23) Nitrobenzene-d5	7.357	82	544613	78.369	ng	0.00
42) 2,4,6-Tribromophenol	10.622	330	158098	79.490	ng	0.00
45) 2-Fluorobiphenyl	9.151	172	1012492	78.786	ng	0.00
79) Terphenyl-d14	12.904	244	1133126	75.750	ng	0.00

Target Compounds					Qvalue
2) 1,4-Dioxane	2.475	88	93282	40.369	ng 99
3) Pyridine	3.205	79	230604	41.938	ng 99
4) n-Nitrosodimethylamine	3.158	42	151656	41.653	ng 97
6) Aniline	6.451	93	284762	41.834	ng 92
8) 2-Chlorophenol	6.575	128	251415	40.527	ng 96
9) Benzaldehyde	6.340	77	154649	40.105	ng 99
10) Phenol	6.446	94	306753	40.615	ng 97
11) bis(2-Chloroethyl)ether	6.528	93	226152	39.866	ng 100
12) 1,3-Dichlorobenzene	6.728	146	264052	40.317	ng 98
13) 1,4-Dichlorobenzene	6.804	146	267068	39.911	ng 99
14) 1,2-Dichlorobenzene	6.957	146	250623	40.354	ng 99
15) Benzyl Alcohol	6.934	79	228379	41.041	ng 100
16) 2,2'-oxybis(1-Chloropr...	7.069	45	393626	40.535	ng 98
17) 2-Methylphenol	7.051	107	197905	40.768	ng 99
18) Hexachloroethane	7.298	117	101687	41.061	ng 99
19) n-Nitroso-di-n-propyla...	7.204	70	173832	40.048	ng 100
20) 3+4-Methylphenols	7.204	107	247254	40.073	ng 98
22) Acetophenone	7.198	105	353867	39.246	ng 100
24) Nitrobenzene	7.375	77	267193	39.429	ng 100
25) Isophorone	7.610	82	440640	39.767	ng 100
26) 2-Nitrophenol	7.687	139	136248	40.413	ng 98
27) 2,4-Dimethylphenol	7.728	122	157242	39.924	ng 99
28) bis(2-Chloroethoxy)met...	7.822	93	282744	39.822	ng 100
29) 2,4-Dichlorophenol	7.934	162	211891	39.818	ng 99
30) 1,2,4-Trichlorobenzene	8.016	180	227379	39.237	ng 100
31) Naphthalene	8.093	128	741255	39.287	ng 100
32) Benzoic acid	7.857	122	164966	40.324	ng 98
33) 4-Chloroaniline	8.145	127	267977	40.680	ng 99
34) Hexachlorobutadiene	8.210	225	139658	40.144	ng 99
35) Caprolactam	8.516	113	66599	41.104	ng 99
36) 4-Chloro-3-methylphenol	8.634	107	235824	40.006	ng 100
37) 2-Methylnaphthalene	8.787	142	475273	38.710	ng 100
38) 1-Methylnaphthalene	8.881	142	464346	39.049	ng 99
40) 1,2,4,5-Tetrachloroben...	8.951	216	230166	40.182	ng 99
41) Hexachlorocyclopentadiene	8.934	237	73834	38.754	ng 97
43) 2,4,6-Trichlorophenol	9.063	196	147770	39.914	ng 99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.110	196	160926	40.108	ng	99
46) 1,1'-Biphenyl	9.251	154	606611	39.519	ng	100
47) 2-Chloronaphthalene	9.275	162	447817	39.453	ng	100
48) 2-Nitroaniline	9.375	65	154092	40.451	ng	98
49) Acenaphthylene	9.687	152	670833	39.734	ng	99
50) Dimethylphthalate	9.551	163	530769	39.488	ng	99
51) 2,6-Dinitrotoluene	9.616	165	117431	39.965	ng	99
52) Acenaphthene	9.863	154	452794	39.893	ng	100
53) 3-Nitroaniline	9.787	138	127965	40.463	ng	99
54) 2,4-Dinitrophenol	9.892	184	74461	42.639	ng	97
55) Dibenzofuran	10.034	168	664069	39.860	ng	99
56) 4-Nitrophenol	9.957	139	98990	42.166	ng	99
57) 2,4-Dinitrotoluene	10.022	165	158941	40.633	ng	100
58) Fluorene	10.375	166	508833	39.501	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.157	232	139688	40.438	ng	97
60) Diethylphthalate	10.251	149	521933	39.467	ng	100
61) 4-Chlorophenyl-phenyle...	10.369	204	243659	39.318	ng	98
62) 4-Nitroaniline	10.398	138	132059	41.124	ng	99
63) Azobenzene	10.528	77	522705	39.759	ng	99
65) 4,6-Dinitro-2-methylph...	10.428	198	98944	42.230	ng	98
66) n-Nitrosodiphenylamine	10.492	169	432307	40.040	ng	99
67) 4-Bromophenyl-phenylether	10.857	248	154044	40.864	ng	95
68) Hexachlorobenzene	10.928	284	156075	40.208	ng	99
69) Atrazine	11.016	200	127486	39.358	ng	100
70) Pentachlorophenol	11.122	266	105427	42.476	ng	99
71) Phenanthrene	11.339	178	744761	40.114	ng	100
72) Anthracene	11.392	178	753295	40.362	ng	100
73) Carbazole	11.551	167	676218	40.267	ng	100
74) Di-n-butylphthalate	11.881	149	794101	40.953	ng	100
75) Fluoranthene	12.528	202	807256	41.011	ng	100
77) Benzidine	12.651	184	211734	38.017	ng	100
78) Pyrene	12.757	202	812250	37.784	ng	100
80) Butylbenzylphthalate	13.380	149	311103	41.781	ng	99
81) Benzo(a)anthracene	13.951	228	638036	38.009	ng	100
82) 3,3'-Dichlorobenzidine	13.916	252	185650	38.608	ng	99
83) Chrysene	13.992	228	610091	39.361	ng	100
84) Bis(2-ethylhexyl)phtha...	13.945	149	371434	44.229	ng	100
85) Di-n-octyl phthalate	14.574	149	534947	44.652	ng	100
87) Indeno(1,2,3-cd)pyrene	16.874	276	474949	38.551	ng	99
88) Benzo(b)fluoranthene	15.010	252	526847	39.352	ng	100
89) Benzo(k)fluoranthene	15.039	252	470794	41.182	ng	98
90) Benzo(a)pyrene	15.369	252	427462	39.459	ng	99
91) Dibenzo(a,h)anthracene	16.892	278	388028	38.870	ng	99
92) Benzo(g,h,i)perylene	17.310	276	402566	38.536	ng	99

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

