

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF071525\
 Data File : BF143095.D
 Acq On : 15 Jul 2025 14:05
 Operator : RC/JU
 Sample : SSTDICC010
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC010

Manual Integrations
 APPROVED

Reviewed By :Rahul Chavli 07/16/2025
 Supervised By :Jagrut Upadhyay 07/16/2025

Quant Time: Jul 15 17:38:34 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF071525.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 15 17:08:07 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.969	152	129158	20.000	ng	0.00
21) Naphthalene-d8	8.245	136	488635	20.000	ng	0.00
39) Acenaphthene-d10	10.004	164	248013	20.000	ng	0.00
64) Phenanthrene-d10	11.486	188	412977	20.000	ng	0.00
76) Chrysene-d12	14.127	240	231826	20.000	ng	0.00
86) Perylene-d12	15.633	264	195100	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.581	112	164234	20.090	ng	0.00
7) Phenol-d6	6.575	99	208774	20.317	ng	-0.01
23) Nitrobenzene-d5	7.516	82	156007m	17.910	ng	-0.01
42) 2,4,6-Tribromophenol	10.786	330	39055	17.661	ng	0.00
45) 2-Fluorobiphenyl	9.322	172	407861	21.860	ng	0.00
79) Terphenyl-d14	13.069	244	340992	21.502	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.810	88	40980	10.070	ng	98
3) Pyridine	3.581	79	99499	9.656	ng	98
4) n-Nitrosodimethylamine	3.504	42	50912	9.536	ng	# 96
6) Aniline	6.622	93	145577	10.060	ng	99
8) 2-Chlorophenol	6.745	128	78009	9.498	ng	98
9) Benzaldehyde	6.510	77	77218	9.862	ng	97
10) Phenol	6.587	94	112508	10.114	ng	98
11) bis(2-Chloroethyl)ether	6.693	93	88047	10.224	ng	99
12) 1,3-Dichlorobenzene	6.910	146	95695	10.313	ng	99
13) 1,4-Dichlorobenzene	6.987	146	95134	10.191	ng	99
14) 1,2-Dichlorobenzene	7.140	146	90960	10.218	ng	100
15) Benzyl Alcohol	7.092	79	72743	9.554	ng	99
16) 2,2'-oxybis(1-Chloropr...	7.240	45	163494	10.299	ng	99
17) 2-Methylphenol	7.204	107	69808	9.849	ng	99
18) Hexachloroethane	7.487	117	28400	9.345	ng	98
19) n-Nitroso-di-n-propyla...	7.363	70	63831	10.139	ng	95
20) 3+4-Methylphenols	7.351	107	90318	10.391	ng	# 83
22) Acetophenone	7.363	105	126127	10.708	ng	95
24) Nitrobenzene	7.540	77	77369	9.201	ng	97
25) Isophorone	7.775	82	167986	9.978	ng	99
26) 2-Nitrophenol	7.857	139	18432	8.966	ng	97
27) 2,4-Dimethylphenol	7.887	122	78706	10.072	ng	98
28) bis(2-Chloroethoxy)met...	7.987	93	106045	10.387	ng	98
29) 2,4-Dichlorophenol	8.092	162	62164	9.724	ng	99
30) 1,2,4-Trichlorobenzene	8.187	180	74581	10.280	ng	98
31) Naphthalene	8.269	128	256031	10.551	ng	99
32) Benzoic acid	7.939	122	19661m	10.880	ng	
33) 4-Chloroaniline	8.310	127	102550	10.510	ng	96
34) Hexachlorobutadiene	8.392	225	42959	9.901	ng	97
35) Caprolactam	8.634	113	17855	8.944	ng	99
36) 4-Chloro-3-methylphenol	8.775	107	70747	9.935	ng	97
37) 2-Methylnaphthalene	8.957	142	152822	10.557	ng	99
38) 1-Methylnaphthalene	9.057	142	161228	10.739	ng	99
40) 1,2,4,5-Tetrachloroben...	9.128	216	74348	10.287	ng	99
41) Hexachlorocyclopentadiene	9.122	237	31472	7.854	ng	100
43) 2,4,6-Trichlorophenol	9.234	196	38748	8.628	ng	98

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44) 2,4,5-Trichlorophenol	9.263	196	45455	9.653	ng	98
46) 1,1'-Biphenyl	9.422	154	209259	10.580	ng	99
47) 2-Chloronaphthalene	9.445	162	149175	10.267	ng	99
48) 2-Nitroaniline	9.534	65	27816	9.163	ng	98
49) Acenaphthylene	9.863	152	249435	10.302	ng	100
50) Dimethylphthalate	9.710	163	156072	10.003	ng	100
51) 2,6-Dinitrotoluene	9.769	165	21789	9.451	ng	96
52) Acenaphthene	10.033	154	149318	10.451	ng	100
53) 3-Nitroaniline	9.939	138	28105	8.336	ng	98
54) 2,4-Dinitrophenol	10.039	184	4729	10.239	ng #	1
55) Dibenzofuran	10.204	168	222555	10.451	ng	98
56) 4-Nitrophenol	10.081	139	20334	7.589	ng	97
57) 2,4-Dinitrotoluene	10.175	165	24420	9.293	ng #	95
58) Fluorene	10.551	166	171975	10.771	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.322	232	33995	9.084	ng	99
60) Diethylphthalate	10.416	149	150934	10.027	ng	100
61) 4-Chlorophenyl-phenyle...	10.545	204	79583	10.400	ng	99
62) 4-Nitroaniline	10.545	138	25757	8.785	ng	96
63) Azobenzene	10.698	77	173437	10.418	ng	98
65) 4,6-Dinitro-2-methylph...	10.581	198	7158	10.290	ng	96
66) n-Nitrosodiphenylamine	10.651	169	146433	10.094	ng	98
67) 4-Bromophenyl-phenylether	11.033	248	45279	9.641	ng	97
68) Hexachlorobenzene	11.104	284	49035	9.840	ng	96
69) Atrazine	11.175	200	35230	9.492	ng	99
70) Pentachlorophenol	11.286	266	20245	7.405	ng	97
71) Phenanthrene	11.510	178	234601	10.526	ng	100
72) Anthracene	11.563	178	232459	10.372	ng	99
73) Carbazole	11.710	167	211853	10.551	ng	99
74) Di-n-butylphthalate	12.051	149	174639	9.475	ng	100
75) Fluoranthene	12.704	202	217346	10.725	ng	99
77) Benzidine	12.816	184	76979	9.747	ng	99
78) Pyrene	12.927	202	226752	10.542	ng	99
80) Butylbenzylphthalate	13.545	149	26977	9.210	ng	94
81) Benzo(a)anthracene	14.116	228	151099	9.781	ng	99
82) 3,3'-Dichlorobenzidine	14.074	252	36435	8.122	ng	99
83) Chrysene	14.151	228	142990	10.050	ng	99
84) Bis(2-ethylhexyl)phtha...	14.110	149	39724	9.073	ng	97
85) Di-n-octyl phthalate	14.727	149	54952	10.526	ng	96
87) Indeno(1,2,3-cd)pyrene	17.168	276	137620	9.484	ng	100
88) Benzo(b)fluoranthene	15.186	252	116154	10.321	ng	100
89) Benzo(k)fluoranthene	15.216	252	103994	9.686	ng	98
90) Benzo(a)pyrene	15.568	252	99366	9.336	ng	99
91) Dibenzo(a,h)anthracene	17.186	278	113316	9.553	ng	99
92) Benzo(g,h,i)perylene	17.627	276	115816	9.577	ng	98

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

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