

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF091025\
 Data File : BF143683.D
 Acq On : 09 Sep 2025 13:54
 Operator : RC/JU
 Sample : SSTDICC080
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC080

Manual Integrations
 APPROVED

Reviewed By :Rahul Chavli 09/10/2025
 Supervised By :Jagrut Upadhyay 09/10/2025

Quant Time: Sep 09 15:31:16 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091025.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 09 15:20:16 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.892	152	56382	20.000	ng	0.00
21) Naphthalene-d8	8.175	136	213611	20.000	ng	0.00
39) Acenaphthene-d10	9.927	164	124115	20.000	ng	0.00
64) Phenanthrene-d10	11.416	188	187688	20.000	ng	0.00
76) Chrysene-d12	14.057	240	116895	20.000	ng	0.01
86) Perylene-d12	15.527	264	129691	20.000	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.504	112	518124	156.786	ng	0.00
7) Phenol-d6	6.516	99	624813	155.282	ng	0.01
23) Nitrobenzene-d5	7.463	82	559223	162.838	ng	0.01
42) 2,4,6-Tribromophenol	10.722	330	193329	154.564	ng	0.01
45) 2-Fluorobiphenyl	9.251	172	1179990	143.618	ng	0.00
79) Terphenyl-d14	12.998	244	1031371	142.979	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.675	88	113617	77.075	ng	99
3) Pyridine	3.440	79	302753	79.506	ng	98
4) n-Nitrosodimethylamine	3.410	42	116844	80.779	ng	94
6) Aniline	6.557	93	412609	77.423	ng	99
8) 2-Chlorophenol	6.675	128	283491	79.429	ng	100
9) Benzaldehyde	6.445	77	255985	78.850	ng	99
10) Phenol	6.534	94	346126m	77.819	ng	
11) bis(2-Chloroethyl)ether	6.634	93	257746	77.590	ng	97
12) 1,3-Dichlorobenzene	6.834	146	315159	77.209	ng	99
13) 1,4-Dichlorobenzene	6.910	146	314684	76.379	ng	100
14) 1,2-Dichlorobenzene	7.063	146	298298	76.670	ng	99
15) Benzyl Alcohol	7.034	79	230026	78.590	ng	99
16) 2,2'-oxybis(1-Chloropr...	7.163	45	337215	77.591	ng	97
17) 2-Methylphenol	7.139	107	230006	80.110	ng	98
18) Hexachloroethane	7.404	117	105200	79.044	ng	98
19) n-Nitroso-di-n-propyla...	7.316	70	192814	78.023	ng	99
20) 3+4-Methylphenols	7.298	107	288678	76.353	ng	91
22) Acetophenone	7.304	105	371310	76.965	ng	# 93
24) Nitrobenzene	7.481	77	284138	79.794	ng	98
25) Isophorone	7.722	82	528121	80.197	ng	99
26) 2-Nitrophenol	7.792	139	150727	86.090	ng	99
27) 2,4-Dimethylphenol	7.822	122	245459	78.096	ng	99
28) bis(2-Chloroethoxy)met...	7.928	93	323002	79.814	ng	100
29) 2,4-Dichlorophenol	8.028	162	253340	80.128	ng	99
30) 1,2,4-Trichlorobenzene	8.116	180	256397	78.277	ng	99
31) Naphthalene	8.198	128	818642	77.417	ng	99
32) Benzoic acid	7.957	122	186144m	97.074	ng	
33) 4-Chloroaniline	8.245	127	284322	77.235	ng	99
34) Hexachlorobutadiene	8.316	225	167644	78.469	ng	99
35) Caprolactam	8.639	113	69698	82.827	ng	99
36) 4-Chloro-3-methylphenol	8.722	107	248979	79.866	ng	98
37) 2-Methylnaphthalene	8.886	142	552024	76.261	ng	99
38) 1-Methylnaphthalene	8.986	142	527838	75.984	ng	98
40) 1,2,4,5-Tetrachloroben...	9.051	216	271043	76.281	ng	100
41) Hexachlorocyclopentadiene	9.039	237	146134	80.108	ng	99
43) 2,4,6-Trichlorophenol	9.163	196	189455	79.850	ng	98

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF091025\
 Data File : BF143683.D
 Acq On : 09 Sep 2025 13:54
 Operator : RC/JU
 Sample : SSTDICC080
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC080

Manual Integrations
 APPROVED

Reviewed By :Rahul Chavli 09/10/2025
 Supervised By :Jagrut Upadhyay 09/10/2025

Quant Time: Sep 09 15:31:16 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091025.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 09 15:20:16 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.198	196	189631	78.619	ng	98
46) 1,1'-Biphenyl	9.357	154	666722	74.617	ng	100
47) 2-Chloronaphthalene	9.380	162	530685	75.759	ng	100
48) 2-Nitroaniline	9.469	65	152123	82.062	ng	96
49) Acenaphthylene	9.792	152	759223	75.664	ng	100
50) Dimethylphthalate	9.651	163	617247	76.436	ng	100
51) 2,6-Dinitrotoluene	9.716	165	128331	83.117	ng	97
52) Acenaphthene	9.969	154	526045	76.080	ng	99
53) 3-Nitroaniline	9.886	138	132242	79.589	ng	99
54) 2,4-Dinitrophenol	9.986	184	68283	82.053	ng	# 53
55) Dibenzofuran	10.139	168	732979	73.721	ng	99
56) 4-Nitrophenol	10.033	139	105333	79.683	ng	99
57) 2,4-Dinitrotoluene	10.116	165	174053	82.828	ng	98
58) Fluorene	10.480	166	567673	72.939	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.251	232	155147	79.326	ng	99
60) Diethylphthalate	10.357	149	575874	75.960	ng	99
61) 4-Chlorophenyl-phenyle...	10.474	204	291237	73.784	ng	99
62) 4-Nitroaniline	10.504	138	126541	80.410	ng	100
63) Azobenzene	10.633	77	504293	75.241	ng	98
65) 4,6-Dinitro-2-methylph...	10.527	198	87568	91.637	ng	99
66) n-Nitrosodiphenylamine	10.592	169	484542	78.804	ng	99
67) 4-Bromophenyl-phenylether	10.963	248	179680	79.387	ng	97
68) Hexachlorobenzene	11.027	284	191089	79.958	ng	98
69) Atrazine	11.116	200	149614	77.606	ng	99
70) Pentachlorophenol	11.216	266	123112	85.217	ng	99
71) Phenanthrene	11.445	178	783137	75.729	ng	99
72) Anthracene	11.498	178	788152	75.796	ng	99
73) Carbazole	11.645	167	649839	76.309	ng	100
74) Di-n-butylphthalate	11.974	149	798541	77.888	ng	99
75) Fluoranthene	12.627	202	695885	73.302	ng	99
77) Benzidine	12.745	184	242284	63.559	ng	100
78) Pyrene	12.857	202	697866	74.668	ng	99
80) Butylbenzylphthalate	13.474	149	249215	85.104	ng	99
81) Benzo(a)anthracene	14.045	228	595222	79.148	ng	99
82) 3,3'-Dichlorobenzidine	14.004	252	185097	80.004	ng	99
83) Chrysene	14.080	228	531262	77.133	ng	98
84) Bis(2-ethylhexyl)phtha...	14.033	149	361768	86.584	ng	99
85) Di-n-octyl phthalate	14.645	149	647392	86.465	ng	100
87) Indeno(1,2,3-cd)pyrene	17.033	276	709706	80.182	ng	100
88) Benzo(b)fluoranthene	15.098	252	621698	80.528	ng	99
89) Benzo(k)fluoranthene	15.127	252	507280	72.806	ng	99
90) Benzo(a)pyrene	15.468	252	528783	77.963	ng	100
91) Dibenzo(a,h)anthracene	17.056	278	585296	79.703	ng	99
92) Benzo(g,h,i)perylene	17.492	276	547026	79.639	ng	99

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

Manual Integrations APPROVED

Quant Time: Sep 09 15:31:16 2025
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091025.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Sep 09 15:20:16 2025
Response via : Initial Calibration

