

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF090925\
 Data File : BF143670.D
 Acq On : 08 Sep 2025 17:50
 Operator : RC/JU
 Sample : SSTDICC050
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC050

Manual Integrations
 APPROVED

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

Quant Time: Sep 09 04:49:12 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF090925.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 09 04:44:42 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.893	152	97020	20.000	ng	0.00
21) Naphthalene-d8	8.175	136	367625	20.000	ng	0.00
39) Acenaphthene-d10	9.928	164	206546	20.000	ng	0.00
64) Phenanthrene-d10	11.416	188	318098	20.000	ng	0.00
76) Chrysene-d12	14.051	240	191784	20.000	ng	0.00
86) Perylene-d12	15.527	264	205711	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.504	112	531717	94.737	ng	0.00
7) Phenol-d6	6.522	99	647805	94.580	ng	0.01
23) Nitrobenzene-d5	7.463	82	584502	98.991	ng	0.01
42) 2,4,6-Tribromophenol	10.722	330	203812	97.329	ng	0.00
45) 2-Fluorobiphenyl	9.251	172	1243890	93.909	ng	0.00
79) Terphenyl-d14	12.998	244	1068628	91.337	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.681	88	118620	47.168	ng	97
3) Pyridine	3.452	79	320416	49.190	ng	99
4) n-Nitrosodimethylamine	3.416	42	123949	49.137	ng	97
6) Aniline	6.557	93	445982	48.435	ng	100
8) 2-Chlorophenol	6.681	128	294266	48.255	ng	99
9) Benzaldehyde	6.446	77	205349	43.582	ng	98
10) Phenol	6.534	94	382886	49.348	ng	96
11) bis(2-Chloroethyl)ether	6.634	93	268243	47.205	ng	99
12) 1,3-Dichlorobenzene	6.834	146	336717	47.930	ng	99
13) 1,4-Dichlorobenzene	6.910	146	334375	47.494	ng	100
14) 1,2-Dichlorobenzene	7.063	146	316756	47.797	ng	99
15) Benzyl Alcohol	7.034	79	246788	48.469	ng	99
16) 2,2'-oxybis(1-Chloropr...	7.163	45	355662	46.601	ng	99
17) 2-Methylphenol	7.140	107	241054	48.094	ng	99
18) Hexachloroethane	7.404	117	111606	48.697	ng	97
19) n-Nitroso-di-n-propyla...	7.316	70	200895	46.958	ng	98
20) 3+4-Methylphenols	7.298	107	303487	47.125	ng	90
22) Acetophenone	7.304	105	388426	47.028	ng	# 93
24) Nitrobenzene	7.481	77	298631	49.141	ng	98
25) Isophorone	7.722	82	560951	49.016	ng	99
26) 2-Nitrophenol	7.793	139	155916	53.316	ng	98
27) 2,4-Dimethylphenol	7.822	122	259350	48.350	ng	98
28) bis(2-Chloroethoxy)met...	7.928	93	332295	47.530	ng	99
29) 2,4-Dichlorophenol	8.028	162	266324	49.264	ng	98
30) 1,2,4-Trichlorobenzene	8.116	180	271134	48.434	ng	99
31) Naphthalene	8.198	128	863873	47.783	ng	99
32) Benzoic acid	7.951	122	180969m	54.138	ng	
33) 4-Chloroaniline	8.245	127	303003	47.801	ng	99
34) Hexachlorobutadiene	8.316	225	178918	48.670	ng	98
35) Caprolactam	8.640	113	72486	48.831	ng	98
36) 4-Chloro-3-methylphenol	8.722	107	265019	48.866	ng	99
37) 2-Methylnaphthalene	8.887	142	589056	47.324	ng	99
38) 1-Methylnaphthalene	8.987	142	561441	47.102	ng	99
40) 1,2,4,5-Tetrachloroben...	9.051	216	283891	48.505	ng	99
41) Hexachlorocyclopentadiene	9.039	237	156085	51.110	ng	97
43) 2,4,6-Trichlorophenol	9.163	196	197790	50.926	ng	99

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF090925\
 Data File : BF143670.D
 Acq On : 08 Sep 2025 17:50
 Operator : RC/JU
 Sample : SSTDICC050
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC050

Manual Integrations
 APPROVED

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

Quant Time: Sep 09 04:49:12 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF090925.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 09 04:44:42 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.198	196	193931	48.673	ng	99
46) 1,1'-Biphenyl	9.357	154	700394	47.112	ng	99
47) 2-Chloronaphthalene	9.381	162	555244	48.064	ng	100
48) 2-Nitroaniline	9.469	65	159367	50.644	ng	99
49) Acenaphthylene	9.792	152	788615	47.740	ng	99
50) Dimethylphthalate	9.651	163	643992	47.627	ng	99
51) 2,6-Dinitrotoluene	9.716	165	133830	51.814	ng	97
52) Acenaphthene	9.963	154	550830	48.274	ng	99
53) 3-Nitroaniline	9.881	138	135153	49.002	ng	96
54) 2,4-Dinitrophenol	9.981	184	68136	50.218	ng	80
55) Dibenzofuran	10.134	168	769457	47.061	ng	98
56) 4-Nitrophenol	10.028	139	110520	49.100	ng	97
57) 2,4-Dinitrotoluene	10.116	165	175193	51.184	ng	98
58) Fluorene	10.481	166	586495	45.769	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.251	232	164452	50.350	ng	99
60) Diethylphthalate	10.351	149	601040	47.196	ng	100
61) 4-Chlorophenyl-phenyle...	10.469	204	301781	46.641	ng	97
62) 4-Nitroaniline	10.504	138	129918	48.939	ng	98
63) Azobenzene	10.634	77	533223	47.608	ng	99
65) 4,6-Dinitro-2-methylph...	10.528	198	88161	54.287	ng	98
66) n-Nitrosodiphenylamine	10.592	169	511118	49.331	ng	99
67) 4-Bromophenyl-phenylether	10.963	248	190903	50.174	ng	97
68) Hexachlorobenzene	11.028	284	197195	49.615	ng	98
69) Atrazine	11.116	200	165174	49.250	ng	100
70) Pentachlorophenol	11.216	266	129084	52.273	ng	99
71) Phenanthrene	11.439	178	822891	47.279	ng	99
72) Anthracene	11.492	178	834353	47.497	ng	99
73) Carbazole	11.645	167	681315	46.357	ng	99
74) Di-n-butylphthalate	11.975	149	831556	47.019	ng	100
75) Fluoranthene	12.628	202	713849	44.285	ng	99
77) Benzidine	12.745	184	144455	34.797	ng	99
78) Pyrene	12.857	202	715426	46.459	ng	100
80) Butylbenzylphthalate	13.469	149	255555	52.576	ng	100
81) Benzo(a)anthracene	14.039	228	587390	47.908	ng	99
82) 3,3'-Dichlorobenzidine	14.004	252	173688	50.175	ng	99
83) Chrysene	14.080	228	550796	49.275	ng	99
84) Bis(2-ethylhexyl)phtha...	14.027	149	355402	54.312	ng	100
85) Di-n-octyl phthalate	14.645	149	643390	56.430	ng	100
87) Indeno(1,2,3-cd)pyrene	17.021	276	721687	52.017	ng	100
88) Benzo(b)fluoranthene	15.092	252	614649	49.009	ng	99
89) Benzo(k)fluoranthene	15.127	252	519986	47.253	ng	99
90) Benzo(a)pyrene	15.463	252	528879	49.158	ng	99
91) Dibenzo(a,h)anthracene	17.045	278	596049	51.849	ng	98
92) Benzo(g,h,i)perylene	17.480	276	561319	51.622	ng	99

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

Manual Integrations APPROVED

Quant Time: Sep 09 04:49:12 2025
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF090925.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Sep 09 04:44:42 2025
Response via : Initial Calibration

