Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF071525\
Data File : BF143098.D
Acq On : 15 Jul 2025 15:35
Operator : RC/JU
Sample : SSTDICC050
Misc :
ALS Vial : 7 Sample Multiplier: 1

Quant Time: Jul 15 17:41:05 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

Instrument :
BNA_F
ClientSampleId :
SSTDICC050

Manual Integrations APPROVED

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

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.969	152	119035	20.000	ng	0.00
21) Naphthalene-d8	8.251	136	456501	20.000	ng	0.00
39) Acenaphthene-d10	10.004	164	223126	20.000	ng	0.00
64) Phenanthrene-d10	11.492	188	354075	20.000	ng	0.00
76) Chrysene-d12	14.127	240	183692	20.000	ng	0.00
86) Perylene-d12	15.633	264	221502	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.593	112	776953	103.124	ng	0.00
7) Phenol-d6	6.587	99	978105	103.280	ng	0.00
23) Nitrobenzene-d5	7.528	82	911721	112.035	ng	0.00
42) 2,4,6-Tribromophenol	10.792	330	227678	114.439	ng	0.00
45) 2-Fluorobiphenyl	9.328	172	1628659	97.028	ng	0.00
79) Terphenyl-d14	13.075	244	1271626	101.195	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.811	88	196652	52.435	ng	99
3) Pyridine	3.581	79	500208	52.670	ng	99
4) n-Nitrosodimethylamine	3.528	42	261539	53.153	ng	100
6) Aniline	6.628	93	699680	52.465	ng	100
8) 2-Chlorophenol	6.751	128	405741	53.601	ng	100
9) Benzaldehyde	6.516	77	431994	59.864	ng	98
10) Phenol	6.598	94	529842m	51.680	ng	
11) bis(2-Chloroethyl)ether	6.698	93	408091	51.417	ng	98
12) 1,3-Dichlorobenzene	6.910	146	437876	51.203	ng	99
13) 1,4-Dichlorobenzene	6.987	146	441012	51.260	ng	100
14) 1,2-Dichlorobenzene	7.140	146	421724	51.401	ng	98
15) Benzyl Alcohol	7.104	79	373560	53.234	ng	99
16) 2,2'-oxybis(1-Chloropr...	7.240	45	748892	51.187	ng	100
17) 2-Methylphenol	7.210	107	342357	52.409	ng	99
18) Hexachloroethane	7.487	117	148761	53.114	ng	98
19) n-Nitroso-di-n-propyla...	7.381	70	302685	52.169	ng	99
20) 3+4-Methylphenols	7.363	107	422794	52.776	ng	93
22) Acetophenone	7.375	105	549763	49.959	ng	97
24) Nitrobenzene	7.545	77	439273	55.918	ng	100
25) Isophorone	7.787	82	819274	52.088	ng	100
26) 2-Nitrophenol	7.863	139	160386	50.837	ng	99
27) 2,4-Dimethylphenol	7.893	122	379432	51.976	ng	100
28) bis(2-Chloroethoxy)met...	7.993	93	483608	50.704	ng	99
29) 2,4-Dichlorophenol	8.098	162	319968	53.575	ng	99
30) 1,2,4-Trichlorobenzene	8.193	180	345389	50.958	ng	99
31) Naphthalene	8.275	128	1126659	49.700	ng	100
32) Benzoic acid	7.998	122	180966	49.535	ng	98
33) 4-Chloroaniline	8.310	127	467074	51.237	ng	99
34) Hexachlorobutadiene	8.392	225	207797	51.263	ng	98
35) Caprolactam	8.681	113	99144	53.162	ng	96
36) 4-Chloro-3-methylphenol	8.787	107	350906	52.746	ng	98
37) 2-Methylnaphthalene	8.963	142	681726	50.410	ng	99
38) 1-Methylnaphthalene	9.063	142	701548	50.016	ng	100
40) 1,2,4,5-Tetrachloroben...	9.128	216	335206	51.551	ng	100
41) Hexachlorocyclopentadiene	9.122	237	198811	55.147	ng	98
43) 2,4,6-Trichlorophenol	9.234	196	223131	55.226	ng	99

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Compound	R.T.	QIon	Response	Conc	Units		Dev(Min)
44) 2,4,5-Trichlorophenol	9.269	196	232905	54.977	ng		99
46) 1,1'-Biphenyl	9.428	154	897834	50.458	ng		100
47) 2-Chloronaphthalene	9.451	162	666760	51.007	ng		100
48) 2-Nitroaniline	9.539	65	207291	52.775	ng		98
49) Acenaphthylene	9.869	152	1109586	50.940	ng		99
50) Dimethylphthalate	9.722	163	726121	51.728	ng		100
51) 2,6-Dinitrotoluene	9.781	165	146385	52.497	ng		92
52) Acenaphthene	10.039	154	651291	50.669	ng		100
53) 3-Nitroaniline	9.945	138	177602	58.553	ng		99
54) 2,4-Dinitrophenol	10.051	184	47800	51.931	ng	#	1
55) Dibenzofuran	10.210	168	964545	50.344	ng		100
56) 4-Nitrophenol	10.092	139	133073	55.203	ng		98
57) 2,4-Dinitrotoluene	10.181	165	181610	52.116	ng		99
58) Fluorene	10.557	166	716173	49.859	ng		100
59) 2,3,4,6-Tetrachlorophenol	10.322	232	189211	56.202	ng		98
60) Diethylphthalate	10.422	149	702995	51.909	ng		99
61) 4-Chlorophenyl-phenyle...	10.545	204	348876	50.675	ng		99
62) 4-Nitroaniline	10.557	138	150470	57.045	ng		99
63) Azobenzene	10.704	77	770603	51.451	ng		99
65) 4,6-Dinitro-2-methylph...	10.592	198	70324	51.847	ng		92
66) n-Nitrosodiphenylamine	10.663	169	634539	51.019	ng		100
67) 4-Bromophenyl-phenylether	11.034	248	211371	52.494	ng		99
68) Hexachlorobenzene	11.104	284	221245	51.785	ng		98
69) Atrazine	11.181	200	172972	54.359	ng		98
70) Pentachlorophenol	11.292	266	131545	56.120	ng		99
71) Phenanthrene	11.516	178	963568	50.425	ng		100
72) Anthracene	11.569	178	979118	50.953	ng		99
73) Carbazole	11.716	167	883827	51.340	ng		100
74) Di-n-butylphthalate	12.051	149	870362	55.079	ng		100
75) Fluoranthene	12.704	202	883558	50.852	ng		100
77) Benzidine	12.816	184	309313	49.425	ng		99
78) Pyrene	12.933	202	880826	51.681	ng		99
80) Butylbenzylphthalate	13.545	149	176615	51.324	ng		96
81) Benzo(a)anthracene	14.122	228	644353	52.641	ng		99
82) 3,3'-Dichlorobenzidine	14.080	252	196472	55.273	ng		100
83) Chrysene	14.157	228	583302	51.740	ng		99
84) Bis(2-ethylhexyl)phtha...	14.110	149	265225	51.119	ng		99
85) Di-n-octyl phthalate	14.727	149	492840	50.100	ng		99
87) Indeno(1,2,3-cd)pyrene	17.180	276	896872	54.440	ng		100
88) Benzo(b)fluoranthene	15.192	252	718106	56.204	ng		100
89) Benzo(k)fluoranthene	15.221	252	602261	49.410	ng		100
90) Benzo(a)pyrene	15.574	252	662193	54.800	ng		99
91) Dibenzo(a,h)anthracene	17.204	278	736672	54.700	ng		99
92) Benzo(g,h,i)perylene	17.651	276	746789	54.390	ng		98

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

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