

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081425\
 Data File : BF143377.D
 Acq On : 14 Aug 2025 12:55
 Operator : RC/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :

BNA_F

ClientSampleId :

SSTDCCC040

Manual Integrations

APPROVED

Reviewed By :Rahul Chavli 08/15/2025

Supervised By :Jagrut Upadhyay 08/15/2025

Quant Time: Aug 14 13:16:25 2025

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF081225.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Tue Aug 12 13:25:39 2025

Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.934	152	143032	20.000	ng	0.00
21) Naphthalene-d8	8.216	136	542236	20.000	ng	0.00
39) Acenaphthene-d10	9.969	164	283411	20.000	ng	0.00
64) Phenanthrene-d10	11.457	188	404824	20.000	ng	0.00
76) Chrysene-d12	14.098	240	239234	20.000	ng	0.00
86) Perylene-d12	15.592	264	289673	20.000	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.563	112	619218	74.863	ng	0.00
7) Phenol-d6	6.563	99	763071	71.941	ng	0.00
23) Nitrobenzene-d5	7.492	82	688096	79.025	ng	0.00
42) 2,4,6-Tribromophenol	10.763	330	172272	83.174	ng	0.00
45) 2-Fluorobiphenyl	9.286	172	1221958	73.114	ng	0.00
79) Terphenyl-d14	13.039	244	982722	73.338	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.740	88	152653	35.666	ng	98
3) Pyridine	3.522	79	412638	36.179	ng	99
4) n-Nitrosodimethylamine	3.451	42	187367	35.267	ng	98
6) Aniline	6.598	93	530102	35.450	ng	97
8) 2-Chlorophenol	6.722	128	337460	38.182	ng	98
9) Benzaldehyde	6.487	77	233521	33.310	ng	98
10) Phenol	6.581	94	456448	36.326	ng	98
11) bis(2-Chloroethyl)ether	6.663	93	333589	36.022	ng	99
12) 1,3-Dichlorobenzene	6.881	146	368645	36.162	ng	98
13) 1,4-Dichlorobenzene	6.951	146	368057	35.973	ng	99
14) 1,2-Dichlorobenzene	7.104	146	353914	36.440	ng	99
15) Benzyl Alcohol	7.075	79	283480	37.284	ng	100
16) 2,2'-oxybis(1-Chloropr...	7.204	45	598078	36.174	ng	98
17) 2-Methylphenol	7.187	107	271579	36.403	ng	100
18) Hexachloroethane	7.451	117	125788	39.214	ng	95
19) n-Nitroso-di-n-propyla...	7.339	70	228789	34.908	ng	98
20) 3+4-Methylphenols	7.339	107	331798	36.775	ng	97
22) Acetophenone	7.339	105	425087	35.580	ng	98
24) Nitrobenzene	7.510	77	359814	38.049	ng	98
25) Isophorone	7.751	82	628488	35.357	ng	99
26) 2-Nitrophenol	7.828	139	165535	44.009	ng	100
27) 2,4-Dimethylphenol	7.863	122	282954m	38.155	ng	
28) bis(2-Chloroethoxy)met...	7.957	93	395236	36.418	ng	100
29) 2,4-Dichlorophenol	8.075	162	280574	38.736	ng	99
30) 1,2,4-Trichlorobenzene	8.157	180	286115	37.448	ng	99
31) Naphthalene	8.239	128	950950	36.447	ng	100
32) Benzoic acid	7.975	122	133895	43.135	ng	99
33) 4-Chloroaniline	8.281	127	341490	36.287	ng	99
34) Hexachlorobutadiene	8.357	225	181385	38.164	ng	99
35) Caprolactam	8.639	113	76059	37.353	ng	96
36) 4-Chloro-3-methylphenol	8.763	107	282403	37.191	ng	99
37) 2-Methylnaphthalene	8.928	142	622239	36.276	ng	99
38) 1-Methylnaphthalene	9.028	142	599818	36.308	ng	99
40) 1,2,4,5-Tetrachloroben...	9.098	216	293706	37.871	ng	99
41) Hexachlorocyclopentadiene	9.086	237	122851	47.813	ng	99
43) 2,4,6-Trichlorophenol	9.210	196	179680	40.834	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.251	196	195288	39.301	ng	98
46) 1,1'-Biphenyl	9.392	154	727640	36.312	ng	100
47) 2-Chloronaphthalene	9.416	162	575800	36.485	ng	99
48) 2-Nitroaniline	9.510	65	174216	38.835	ng	98
49) Acenaphthylene	9.833	152	822757	36.100	ng	99
50) Dimethylphthalate	9.681	163	605945	36.590	ng	100
51) 2,6-Dinitrotoluene	9.745	165	122638	39.683	ng	97
52) Acenaphthene	10.004	154	543801	35.890	ng	99
53) 3-Nitroaniline	9.916	138	138234	40.760	ng	98
54) 2,4-Dinitrophenol	10.028	184	37998	39.881	ng	90
55) Dibenzofuran	10.175	168	787502	36.113	ng	99
56) 4-Nitrophenol	10.080	139	80105	33.008	ng	96
57) 2,4-Dinitrotoluene	10.151	165	163031	39.758	ng	92
58) Fluorene	10.522	166	593649	35.641	ng	98
59) 2,3,4,6-Tetrachlorophenol	10.298	232	141719	42.309	ng	97
60) Diethylphthalate	10.386	149	549288	37.031	ng	99
61) 4-Chlorophenyl-phenyle...	10.510	204	293085	36.416	ng	99
62) 4-Nitroaniline	10.528	138	124208	37.855	ng	99
63) Azobenzene	10.669	77	600012	35.414	ng	99
65) 4,6-Dinitro-2-methylph...	10.563	198	69944	43.506	ng	95
66) n-Nitrosodiphenylamine	10.627	169	498392	37.389	ng	99
67) 4-Bromophenyl-phenylether	10.998	248	178131	38.474	ng	99
68) Hexachlorobenzene	11.075	284	186624	37.395	ng	99
69) Atrazine	11.151	200	135519	38.979	ng	100
70) Pentachlorophenol	11.269	266	75453	34.523	ng	94
71) Phenanthrene	11.480	178	808775	36.618	ng	100
72) Anthracene	11.533	178	823989	36.786	ng	99
73) Carbazole	11.686	167	672887	35.273	ng	99
74) Di-n-butylphthalate	12.016	149	677330	41.959	ng	100
75) Fluoranthene	12.674	202	695831	35.965	ng	98
77) Benzidine	12.786	184	257151	39.317	ng	98
78) Pyrene	12.904	202	703812	35.069	ng	100
80) Butylbenzylphthalate	13.510	149	198206	49.838	ng	98
81) Benzo(a)anthracene	14.086	228	575793	37.867	ng	99
82) 3,3'-Dichlorobenzidine	14.045	252	165667	40.349	ng	99
83) Chrysene	14.127	228	496975	35.052	ng	100
84) Bis(2-ethylhexyl)phtha...	14.074	149	293859	50.177	ng	98
85) Di-n-octyl phthalate	14.686	149	552169	47.599	ng	97
87) Indeno(1,2,3-cd)pyrene	17.121	276	820270	39.525	ng	99
88) Benzo(b)fluoranthene	15.151	252	585035	36.925	ng	99
89) Benzo(k)fluoranthene	15.180	252	538537	34.980	ng	100
90) Benzo(a)pyrene	15.533	252	550144	36.867	ng	99
91) Dibenzo(a,h)anthracene	17.133	278	657670	39.667	ng	99
92) Benzo(g,h,i)perylene	17.586	276	662349	39.657	ng	100

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

