

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102825\
 Data File : BF144097.D
 Acq On : 28 Oct 2025 13:17
 Operator : RC/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Quant Time: Oct 28 14:10:36 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF102425.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 24 17:30:25 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.881	152	66521	20.000	ng	0.00
21) Naphthalene-d8	8.163	136	250146	20.000	ng	0.00
39) Acenaphthene-d10	9.916	164	147935	20.000	ng	0.00
64) Phenanthrene-d10	11.410	188	264042	20.000	ng	0.00
76) Chrysene-d12	14.057	240	128248	20.000	ng	0.00
86) Perylene-d12	15.539	264	169914	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.498	112	346602	82.434	ng	0.00
7) Phenol-d6	6.510	99	390159	84.022	ng	0.00
23) Nitrobenzene-d5	7.445	82	371128	81.285	ng	0.00
42) 2,4,6-Tribromophenol	10.710	330	169601	90.002	ng	0.00
45) 2-Fluorobiphenyl	9.239	172	816145	77.635	ng	0.00
79) Terphenyl-d14	12.992	244	806709	109.967	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.628	88	69149	38.059	ng	96
3) Pyridine	3.404	79	202882	41.710	ng	96
4) n-Nitrosodimethylamine	3.351	42	112807	37.743	ng	91
6) Aniline	6.545	93	275642	41.492	ng	99
8) 2-Chlorophenol	6.663	128	172921	41.762	ng	100
9) Benzaldehyde	6.434	77	147003	40.774	ng	98
10) Phenol	6.528	94	236104	41.214	ng	92
11) bis(2-Chloroethyl)ether	6.616	93	176755	42.699	ng	96
12) 1,3-Dichlorobenzene	6.822	146	212619	41.627	ng	98
13) 1,4-Dichlorobenzene	6.898	146	210959	41.909	ng	98
14) 1,2-Dichlorobenzene	7.051	146	203316	41.744	ng	98
15) Benzyl Alcohol	7.022	79	164325	43.521	ng	97
16) 2,2'-oxybis(1-Chloropr...	7.151	45	334275	42.966	ng	96
17) 2-Methylphenol	7.134	107	136788	42.952	ng	97
18) Hexachloroethane	7.392	117	74260	41.263	ng	98
19) n-Nitroso-di-n-propyla...	7.292	70	133622	44.499	ng	92
20) 3+4-Methylphenols	7.287	107	161750	42.964	ng	# 86
22) Acetophenone	7.292	105	224937	41.609	ng	95
24) Nitrobenzene	7.463	77	204243	42.287	ng	99
25) Isophorone	7.704	82	357344	44.766	ng	99
26) 2-Nitrophenol	7.781	139	101537	44.191	ng	95
27) 2,4-Dimethylphenol	7.816	122	151833	44.467	ng	98
28) bis(2-Chloroethoxy)met...	7.910	93	222369	44.670	ng	97
29) 2,4-Dichlorophenol	8.022	162	169254	42.625	ng	97
30) 1,2,4-Trichlorobenzene	8.104	180	172640	42.262	ng	99
31) Naphthalene	8.186	128	493628	41.811	ng	98
32) Benzoic acid	7.934	122	115219	53.198	ng	# 84
33) 4-Chloroaniline	8.234	127	182107	44.308	ng	98
34) Hexachlorobutadiene	8.298	225	134816	40.361	ng	99
35) Caprolactam	8.610	113	48594	48.802	ng	91
36) 4-Chloro-3-methylphenol	8.716	107	161239	46.014	ng	96
37) 2-Methylnaphthalene	8.875	142	333422	42.835	ng	96
38) 1-Methylnaphthalene	8.975	142	320892	43.513	ng	100
40) 1,2,4,5-Tetrachloroben...	9.039	216	197569	38.859	ng	98
41) Hexachlorocyclopentadiene	9.028	237	70289	44.760	ng	97
43) 2,4,6-Trichlorophenol	9.151	196	140181	42.418	ng	98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.198	196	154404	44.519	ng	96
46) 1,1'-Biphenyl	9.339	154	411658	38.873	ng	100
47) 2-Chloronaphthalene	9.363	162	360439	40.627	ng	97
48) 2-Nitroaniline	9.463	65	114922	43.622	ng	98
49) Acenaphthylene	9.780	152	498378	41.661	ng	99
50) Dimethylphthalate	9.633	163	423610	43.685	ng	100
51) 2,6-Dinitrotoluene	9.704	165	87885	45.846	ng	92
52) Acenaphthene	9.951	154	328569	41.127	ng	98
53) 3-Nitroaniline	9.875	138	96397	45.229	ng	98
54) 2,4-Dinitrophenol	9.980	184	49169	47.423	ng	98
55) Dibenzofuran	10.128	168	463130	41.757	ng	97
56) 4-Nitrophenol	10.033	139	65421	45.329	ng	92
57) 2,4-Dinitrotoluene	10.110	165	120319	46.505	ng	95
58) Fluorene	10.469	166	357025	42.247	ng	97
59) 2,3,4,6-Tetrachlorophenol	10.245	232	109986	45.002	ng	98
60) Diethylphthalate	10.339	149	401002	43.974	ng	100
61) 4-Chlorophenyl-phenyle...	10.457	204	205056	43.147	ng	97
62) 4-Nitroaniline	10.486	138	89874	44.504	ng	96
63) Azobenzene	10.622	77	376152	43.581	ng	98
65) 4,6-Dinitro-2-methylph...	10.522	198	71598	42.264	ng	98
66) n-Nitrosodiphenylamine	10.580	169	344740	40.864	ng	98
67) 4-Bromophenyl-phenylether	10.951	248	140573	42.521	ng	97
68) Hexachlorobenzene	11.022	284	169947	41.916	ng	96
69) Atrazine	11.104	200	109902	39.700	ng	96
70) Pentachlorophenol	11.216	266	89552	46.401	ng	96
71) Phenanthrene	11.433	178	534720	40.767	ng	99
72) Anthracene	11.486	178	546494	40.333	ng	98
73) Carbazole	11.639	167	458674	39.771	ng	98
74) Di-n-butylphthalate	11.969	149	571521	40.420	ng	100
75) Fluoranthene	12.621	202	528275	36.455	ng	99
77) Benzidine	12.751	184	201207	50.411	ng	99
78) Pyrene	12.857	202	530531	57.540	ng	99
80) Butylbenzylphthalate	13.468	149	167106	47.907	ng	98
81) Benzo(a)anthracene	14.045	228	379691	43.808	ng	99
82) 3,3'-Dichlorobenzidine	14.004	252	132543	43.932	ng	100
83) Chrysene	14.080	228	342254	42.081	ng	99
84) Bis(2-ethylhexyl)phtha...	14.027	149	198759	41.739	ng	99
85) Di-n-octyl phthalate	14.639	149	339385	41.443	ng	100
87) Indeno(1,2,3-cd)pyrene	17.045	276	525344	45.034	ng	98
88) Benzo(b)fluoranthene	15.104	252	382543	39.824	ng	100
89) Benzo(k)fluoranthene	15.133	252	355845	40.151	ng	100
90) Benzo(a)pyrene	15.474	252	371712	43.134	ng	99
91) Dibenzo(a,h)anthracene	17.062	278	413954	44.829	ng	99
92) Benzo(g,h,i)perylene	17.503	276	429795	46.294	ng	99

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

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