

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF071625\
 Data File : BF143117.D
 Acq On : 16 Jul 2025 19:39
 Operator : RC/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Quant Time: Jul 17 03:44:09 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 17:53:25 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.969	152	135149	20.000	ng	0.00
21) Naphthalene-d8	8.251	136	510164	20.000	ng	0.00
39) Acenaphthene-d10	10.004	164	254065	20.000	ng	0.00
64) Phenanthrene-d10	11.492	188	397173	20.000	ng	0.00
76) Chrysene-d12	14.127	240	210714	20.000	ng	0.00
86) Perylene-d12	15.633	264	255524	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.587	112	675027	78.913	ng	0.00
7) Phenol-d6	6.587	99	845178	78.603	ng	0.00
23) Nitrobenzene-d5	7.528	82	820016	90.170	ng	0.00
42) 2,4,6-Tribromophenol	10.792	330	193992	85.634	ng	0.00
45) 2-Fluorobiphenyl	9.322	172	1444367	75.570	ng	0.00
79) Terphenyl-d14	13.075	244	1097674	76.150	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.811	88	169746	39.864	ng	98
3) Pyridine	3.581	79	425792	39.489	ng	98
4) n-Nitrosodimethylamine	3.522	42	217941	39.012	ng	99
6) Aniline	6.628	93	595971	39.360	ng	99
8) 2-Chlorophenol	6.751	128	355815	41.401	ng	98
9) Benzaldehyde	6.516	77	335257	40.919	ng	99
10) Phenol	6.598	94	490852	42.169	ng	100
11) bis(2-Chloroethyl)ether	6.698	93	350864	38.936	ng	99
12) 1,3-Dichlorobenzene	6.910	146	386579	39.815	ng	99
13) 1,4-Dichlorobenzene	6.987	146	383281	39.238	ng	99
14) 1,2-Dichlorobenzene	7.140	146	363768	39.051	ng	98
15) Benzyl Alcohol	7.098	79	319688	40.125	ng	99
16) 2,2'-oxybis(1-Chloropr...	7.240	45	645668	38.870	ng	100
17) 2-Methylphenol	7.204	107	295753	39.876	ng	99
18) Hexachloroethane	7.487	117	131364	41.310	ng	98
19) n-Nitroso-di-n-propyla...	7.375	70	260963	39.615	ng	99
20) 3+4-Methylphenols	7.363	107	369651	40.641	ng	99
22) Acetophenone	7.369	105	476586	38.753	ng	# 93
24) Nitrobenzene	7.545	77	376400	42.874	ng	99
25) Isophorone	7.787	82	688462	39.167	ng	100
26) 2-Nitrophenol	7.863	139	148006	42.663	ng	100
27) 2,4-Dimethylphenol	7.893	122	326049	39.965	ng	99
28) bis(2-Chloroethoxy)met...	7.993	93	420098	39.412	ng	100
29) 2,4-Dichlorophenol	8.098	162	274747	41.164	ng	99
30) 1,2,4-Trichlorobenzene	8.187	180	297978	39.339	ng	98
31) Naphthalene	8.269	128	991571	39.140	ng	100
32) Benzoic acid	7.987	122	150277	38.466	ng	97
33) 4-Chloroaniline	8.310	127	401159	39.377	ng	100
34) Hexachlorobutadiene	8.393	225	181527	40.072	ng	99
35) Caprolactam	8.675	113	85847	41.190	ng	95
36) 4-Chloro-3-methylphenol	8.787	107	298817	40.192	ng	100
37) 2-Methylnaphthalene	8.963	142	589636	39.014	ng	100
38) 1-Methylnaphthalene	9.063	142	610968	38.976	ng	100
40) 1,2,4,5-Tetrachloroben...	9.128	216	287903	38.885	ng	99
41) Hexachlorocyclopentadiene	9.122	237	173333	42.225	ng	99
43) 2,4,6-Trichlorophenol	9.234	196	195982	42.600	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.269	196	196557	40.747	ng	99
46) 1,1'-Biphenyl	9.422	154	779366	38.466	ng	100
47) 2-Chloronaphthalene	9.451	162	573731	38.546	ng	100
48) 2-Nitroaniline	9.534	65	178618	40.710	ng	99
49) Acenaphthylene	9.869	152	971647	39.176	ng	99
50) Dimethylphthalate	9.722	163	628010	39.291	ng	100
51) 2,6-Dinitrotoluene	9.775	165	130906	41.829	ng	95
52) Acenaphthene	10.039	154	566434	38.701	ng	99
53) 3-Nitroaniline	9.945	138	151765	43.942	ng	97
54) 2,4-Dinitrophenol	10.045	184	41904	42.980	ng	# 44
55) Dibenzofuran	10.210	168	838011	38.413	ng	100
56) 4-Nitrophenol	10.086	139	117068	42.650	ng	99
57) 2,4-Dinitrotoluene	10.181	165	155813	40.106	ng	96
58) Fluorene	10.551	166	619482	37.875	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.322	232	164509	42.914	ng	98
60) Diethylphthalate	10.422	149	615586	39.920	ng	100
61) 4-Chlorophenyl-phenyle...	10.545	204	306293	39.072	ng	100
62) 4-Nitroaniline	10.557	138	129761	43.204	ng	99
63) Azobenzene	10.704	77	652575	38.265	ng	99
65) 4,6-Dinitro-2-methylph...	10.586	198	62192	43.295	ng	97
66) n-Nitrosodiphenylamine	10.657	169	545898	39.129	ng	99
67) 4-Bromophenyl-phenylether	11.034	248	180440	39.949	ng	98
68) Hexachlorobenzene	11.104	284	191457	39.950	ng	97
69) Atrazine	11.181	200	145300	40.708	ng	99
70) Pentachlorophenol	11.292	266	114049	43.376	ng	99
71) Phenanthrene	11.516	178	831246	38.780	ng	100
72) Anthracene	11.569	178	844296	39.169	ng	99
73) Carbazole	11.716	167	742234	38.436	ng	100
74) Di-n-butylphthalate	12.051	149	743944	41.970	ng	100
75) Fluoranthene	12.704	202	752396	38.604	ng	99
77) Benzidine	12.816	184	274069	38.177	ng	99
78) Pyrene	12.933	202	747120	38.214	ng	100
80) Butylbenzylphthalate	13.545	149	153672	39.754	ng	95
81) Benzo(a)anthracene	14.116	228	570413	40.624	ng	100
82) 3,3'-Dichlorobenzidine	14.074	252	164768	40.409	ng	99
83) Chrysene	14.157	228	486986	37.657	ng	100
84) Bis(2-ethylhexyl)phtha...	14.110	149	241107	41.219	ng	100
85) Di-n-octyl phthalate	14.727	149	439529	40.440	ng	99
87) Indeno(1,2,3-cd)pyrene	17.180	276	781535	41.122	ng	100
88) Benzo(b)fluoranthene	15.192	252	580612	39.393	ng	99
89) Benzo(k)fluoranthene	15.221	252	558211	39.698	ng	99
90) Benzo(a)pyrene	15.574	252	570084	40.896	ng	99
91) Dibenzo(a,h)anthracene	17.198	278	635338	40.895	ng	99
92) Benzo(g,h,i)perylene	17.645	276	642024	40.534	ng	98

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

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