

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF052025\
 Data File : BF142468.D
 Acq On : 20 May 2025 12:38
 Operator : RC/JU
 Sample : SSTDICC005
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC005

Quant Time: May 20 16:54:53 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF052025.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 20 16:26:47 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.898	152	87254	20.000	ng	0.00
21) Naphthalene-d8	8.181	136	345154	20.000	ng	0.00
39) Acenaphthene-d10	9.939	164	197100	20.000	ng	0.00
64) Phenanthrene-d10	11.427	188	349474	20.000	ng	0.00
76) Chrysene-d12	14.074	240	218524	20.000	ng	0.00
86) Perylene-d12	15.568	264	154887	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.504	112	55140	10.647	ng	-0.01
7) Phenol-d6	6.516	99	66768	10.710	ng	-0.02
23) Nitrobenzene-d5	7.457	82	64876	10.249	ng	-0.02
42) 2,4,6-Tribromophenol	10.727	330	22702	10.378	ng	-0.01
45) 2-Fluorobiphenyl	9.257	172	170142	11.583	ng	-0.01
79) Terphenyl-d14	13.016	244	177430	11.096	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.669	88	10748	5.186	ng	97
3) Pyridine	3.451	79	26993	5.115	ng	98
4) n-Nitrosodimethylamine	3.369	42	13346	4.870	ng	# 97
6) Aniline	6.557	93	44576	5.318	ng	98
8) 2-Chlorophenol	6.681	128	29341	5.201	ng	98
9) Benzaldehyde	6.445	77	22572	6.020	ng	98
10) Phenol	6.528	94	37431	5.336	ng	97
11) bis(2-Chloroethyl)ether	6.628	93	26215	5.192	ng	98
12) 1,3-Dichlorobenzene	6.839	146	34073	5.413	ng	97
13) 1,4-Dichlorobenzene	6.916	146	33593	5.298	ng	99
14) 1,2-Dichlorobenzene	7.069	146	32604	5.374	ng	99
15) Benzyl Alcohol	7.034	79	23103	5.015	ng	97
16) 2,2'-oxybis(1-Chloropr...	7.169	45	45414	5.355	ng	99
17) 2-Methylphenol	7.139	107	22693	5.153	ng	98
18) Hexachloroethane	7.416	117	11635	5.255	ng	99
19) n-Nitroso-di-n-propyla...	7.298	70	20527	5.313	ng	96
20) 3+4-Methylphenols	7.292	107	30797	5.460	ng	# 88
22) Acetophenone	7.304	105	41438	5.423	ng	99
24) Nitrobenzene	7.475	77	29193	5.114	ng	96
25) Isophorone	7.716	82	54838	5.182	ng	98
26) 2-Nitrophenol	7.798	139	14387	4.716	ng	99
27) 2,4-Dimethylphenol	7.828	122	27170	5.051	ng	99
28) bis(2-Chloroethoxy)met...	7.928	93	35010	5.296	ng	97
29) 2,4-Dichlorophenol	8.034	162	24812	5.072	ng	98
30) 1,2,4-Trichlorobenzene	8.122	180	28037	5.281	ng	97
31) Naphthalene	8.204	128	91563	5.388	ng	100
33) 4-Chloroaniline	8.251	127	36597	5.314	ng	98
34) Hexachlorobutadiene	8.322	225	17505	5.280	ng	99
35) Caprolactam	8.581	113	7020	5.109	ng	97
36) 4-Chloro-3-methylphenol	8.722	107	26197	5.225	ng	97
37) 2-Methylnaphthalene	8.898	142	58574	5.478	ng	99
38) 1-Methylnaphthalene	8.992	142	60674	5.486	ng	100
40) 1,2,4,5-Tetrachloroben...	9.063	216	29150	5.152	ng	99
41) Hexachlorocyclopentadiene	9.051	237	15667	4.105	ng	99
43) 2,4,6-Trichlorophenol	9.169	196	19079	5.001	ng	99
44) 2,4,5-Trichlorophenol	9.210	196	20200	5.020	ng	98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
46) 1,1'-Biphenyl	9.357	154	82376	5.387	ng	99
47) 2-Chloronaphthalene	9.386	162	60002	5.311	ng	97
48) 2-Nitroaniline	9.475	65	16399	5.022	ng	95
49) Acenaphthylene	9.798	152	101686	5.330	ng	99
50) Dimethylphthalate	9.651	163	71020	5.461	ng	99
51) 2,6-Dinitrotoluene	9.716	165	14295	5.093	ng	97
52) Acenaphthene	9.975	154	62026	5.325	ng	99
53) 3-Nitroaniline	9.886	138	15749	5.137	ng	99
55) Dibenzofuran	10.145	168	92955	5.545	ng	98
56) 4-Nitrophenol	10.039	139	10876	4.808	ng	93
57) 2,4-Dinitrotoluene	10.116	165	18322	4.998	ng	91
58) Fluorene	10.486	166	72775	5.620	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.263	232	17096	5.090	ng	98
60) Diethylphthalate	10.357	149	70046	5.515	ng	99
61) 4-Chlorophenyl-phenyle...	10.480	204	35694	5.611	ng	97
62) 4-Nitroaniline	10.492	138	14920	5.366	ng	95
63) Azobenzene	10.639	77	61041	5.351	ng	99
65) 4,6-Dinitro-2-methylph...	10.522	198	7475	3.833	ng	90
66) n-Nitrosodiphenylamine	10.592	169	62215	5.204	ng	98
67) 4-Bromophenyl-phenylether	10.975	248	20974	5.076	ng	98
68) Hexachlorobenzene	11.039	284	23159	5.067	ng	100
69) Atrazine	11.122	200	16045	5.053	ng	98
70) Pentachlorophenol	11.233	266	11357	4.401	ng	96
71) Phenanthrene	11.451	178	104484	5.594	ng	100
72) Anthracene	11.504	178	104017	5.457	ng	100
73) Carbazole	11.657	167	89967	5.521	ng	99
74) Di-n-butylphthalate	11.986	149	96777	5.426	ng	99
75) Fluoranthene	12.645	202	102797	5.890	ng	99
77) Benzidine	12.763	184	36585	4.499	ng	98
78) Pyrene	12.874	202	105394	5.170	ng	99
80) Butylbenzylphthalate	13.492	149	25251	4.483	ng	96
81) Benzo(a)anthracene	14.063	228	77800	5.332	ng	100
82) 3,3'-Dichlorobenzidine	14.021	252	19673	4.445	ng	98
83) Chrysene	14.098	228	66761	5.092	ng	99
84) Bis(2-ethylhexyl)phtha...	14.051	149	27864	3.864	ng	98
87) Indeno(1,2,3-cd)pyrene	17.068	276	55034	4.733	ng	99
88) Benzo(b)fluoranthene	15.121	252	51014	5.565	ng	98
89) Benzo(k)fluoranthene	15.151	252	46114	5.371	ng	98
90) Benzo(a)pyrene	15.498	252	44694	5.172	ng	99
91) Dibenzo(a,h)anthracene	17.092	278	44618	4.731	ng	98
92) Benzo(g,h,i)perylene	17.527	276	44684	4.737	ng	97

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

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