

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051225\
 Data File : BP024608.D
 Acq On : 12 May 2025 18:42
 Operator : RC/JU
 Sample : SSTDICC050
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SSTDICC050

Quant Time: May 13 02:55:42 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP051225.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 13 02:51:00 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc Units	Dev(Min)
Internal Standards					
1) 1,4-Dichlorobenzene-d4	7.669	152	148140	20.00 ng	0.00
21) Naphthalene-d8	10.434	136	636656	20.00 ng	0.00
39) Acenaphthene-d10	14.298	164	403739	20.00 ng	0.00
64) Phenanthrene-d10	17.092	188	772415	20.00 ng	0.00
76) Chrysene-d12	21.510	240	834598	20.00 ng	0.00
86) Perylene-d12	24.786	264	958174	20.00 ng	0.00
System Monitoring Compounds					
5) 2-Fluorophenol	5.287	112	851301	101.54 ng	0.00
7) Phenol-d6	6.858	99	1142036	103.57 ng	0.00
23) Nitrobenzene-d5	8.816	82	1100904	99.39 ng	0.00
42) 2,4,6-Tribromophenol	15.810	330	660002	99.75 ng	0.00
45) 2-Fluorobiphenyl	12.910	172	2672191	99.00 ng	0.00
79) Terphenyl-d14	19.810	244	4275676	100.29 ng	0.00
Target Compounds					Qvalue
2) 1,4-Dioxane	3.228	88	178202	47.91 ng	99
3) Pyridine	3.628	79	448907m	51.58 ng	
4) n-Nitrosodimethylamine	3.534	42	190784	49.84 ng	98
6) Aniline	7.005	93	512884	51.85 ng	99
8) 2-Chlorophenol	7.240	128	486614	50.42 ng	100
9) Benzaldehyde	6.822	77	289704	47.52 ng	96
10) Phenol	6.881	94	566334	51.57 ng	98
11) bis(2-Chloroethyl)ether	7.099	93	447378	50.16 ng	100
12) 1,3-Dichlorobenzene	7.557	146	516215	48.61 ng	100
13) 1,4-Dichlorobenzene	7.699	146	525637	48.61 ng	99
14) 1,2-Dichlorobenzene	8.010	146	504489	48.20 ng	98
15) Benzyl Alcohol	7.910	79	416637	54.81 ng	99
16) 2,2'-oxybis(1-Chloropr...	8.193	45	502203	49.11 ng	100
17) 2-Methylphenol	8.116	107	394481	52.24 ng	99
18) Hexachloroethane	8.728	117	198910	48.48 ng	99
19) n-Nitroso-di-n-propyla...	8.463	70	369088	50.19 ng	98
20) 3+4-Methylphenols	8.440	107	546753	52.60 ng	99
22) Acetophenone	8.481	105	748961	49.13 ng	100
24) Nitrobenzene	8.857	77	534613	49.74 ng	99
25) Isophorone	9.375	82	965795	49.75 ng	99
26) 2-Nitrophenol	9.563	139	289691	52.70 ng	99
27) 2,4-Dimethylphenol	9.628	122	338207	50.98 ng	99
28) bis(2-Chloroethoxy)met...	9.851	93	627150	49.72 ng	99
29) 2,4-Dichlorophenol	10.099	162	477747	52.01 ng	98
30) 1,2,4-Trichlorobenzene	10.299	180	494183	48.35 ng	99
31) Naphthalene	10.487	128	1604352	48.32 ng	100
32) Benzoic acid	9.804	122	355171	48.99 ng	99
33) 4-Chloroaniline	10.604	127	582128	52.04 ng	99
34) Hexachlorobutadiene	10.769	225	315698	47.90 ng	98
35) Caprolactam	11.398	113	165189	51.48 ng	96
36) 4-Chloro-3-methylphenol	11.740	107	561391	49.93 ng	98
37) 2-Methylnaphthalene	12.098	142	1138927	48.94 ng	99
38) 1-Methylnaphthalene	12.322	142	1109251	48.23 ng	100
40) 1,2,4,5-Tetrachloroben...	12.475	216	566781	50.31 ng	100
41) Hexachlorocyclopentadiene	12.451	237	210496	56.98 ng	98
43) 2,4,6-Trichlorophenol	12.722	196	396736	52.38 ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.798	196	433223	51.39	ng	99
46) 1,1'-Biphenyl	13.116	154	1444812	49.79	ng	100
47) 2-Chloronaphthalene	13.157	162	1089486	49.44	ng	99
48) 2-Nitroaniline	13.369	65	335112	51.90	ng	98
49) Acenaphthylene	14.010	152	1713049	49.75	ng	99
50) Dimethylphthalate	13.757	163	1469533	48.80	ng	100
51) 2,6-Dinitrotoluene	13.875	165	322875	50.25	ng	99
52) Acenaphthene	14.357	154	1079801	49.23	ng	99
53) 3-Nitroaniline	14.210	138	328887	53.07	ng	99
54) 2,4-Dinitrophenol	14.422	184	190669	48.53	ng	97
55) Dibenzofuran	14.698	168	1733801	48.53	ng	99
56) 4-Nitrophenol	14.545	139	273025	53.57	ng	97
57) 2,4-Dinitrotoluene	14.675	165	455759	51.47	ng	98
58) Fluorene	15.357	166	1364198	49.16	ng	99
59) 2,3,4,6-Tetrachlorophenol	14.934	232	400390	50.36	ng	99
60) Diethylphthalate	15.134	149	1509202	48.39	ng	99
61) 4-Chlorophenyl-phenyle...	15.357	204	682495	48.75	ng	96
62) 4-Nitroaniline	15.392	138	328268	53.32	ng	99
63) Azobenzene	15.651	77	1307644	49.16	ng	98
65) 4,6-Dinitro-2-methylph...	15.457	198	261750	55.21	ng	100
66) n-Nitrosodiphenylamine	15.575	169	1157921	50.13	ng	99
67) 4-Bromophenyl-phenylether	16.263	248	455525	49.90	ng	99
68) Hexachlorobenzene	16.381	284	557991	49.09	ng	98
69) Atrazine	16.551	200	377067	48.41	ng	99
70) Pentachlorophenol	16.739	266	391082	53.52	ng	98
71) Phenanthrene	17.133	178	2039584	48.96	ng	100
72) Anthracene	17.228	178	2011752	49.43	ng	100
73) Carbazole	17.510	167	1891671	49.56	ng	99
74) Di-n-butylphthalate	18.098	149	2569452	50.03	ng	100
75) Fluoranthene	19.216	202	2486056	48.96	ng	99
77) Benzidine	19.422	184	604467	62.20	ng	99
78) Pyrene	19.592	202	2552368	50.54	ng	99
80) Butylbenzylphthalate	20.539	149	1173770	52.12	ng	98
81) Benzo(a)anthracene	21.492	228	2499172	49.18	ng	100
82) 3,3'-Dichlorobenzidine	21.416	252	627836	51.68	ng	99
83) Chrysene	21.557	228	2394884	48.64	ng	100
84) Bis(2-ethylhexyl)phtha...	21.427	149	1725117	52.19	ng	99
85) Di-n-octyl phthalate	22.680	149	2878138	52.51	ng	100
87) Indeno(1,2,3-cd)pyrene	28.503	276	3286670	50.19	ng	# 93
88) Benzo(b)fluoranthene	23.757	252	2778244	50.24	ng	99
89) Benzo(k)fluoranthene	23.821	252	2737321	50.16	ng	100
90) Benzo(a)pyrene	24.633	252	2419904	50.55	ng	99
91) Dibenzo(a,h)anthracene	28.580	278	2738395	50.97	ng	99
92) Benzo(g,h,i)perylene	29.668	276	2785599	50.19	ng	99

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

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