

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081925\
 Data File : BP025473.D
 Acq On : 19 Aug 2025 12:47
 Operator : CG/JU
 Sample : PB169244BSD
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB169244BSD

Quant Time: Aug 19 13:10:51 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP081425.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 14 18:14:31 2025
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By :Rahul Chavli 08/20/2025
 Supervised By :Jagrut Upadhyay 08/20/2025

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.790	152	172374	20.000	ng	0.00
21) Naphthalene-d8	10.555	136	704478	20.000	ng	0.00
39) Acenaphthene-d10	14.401	164	488785	20.000	ng	0.00
64) Phenanthrene-d10	17.207	188	1043571	20.000	ng	0.00
76) Chrysene-d12	21.654	240	1070063	20.000	ng	0.00
86) Perylene-d12	25.030	264	1176738	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.408	112	1466573	141.294	ng	0.00
7) Phenol-d6	6.972	99	1870146	140.533	ng	0.00
23) Nitrobenzene-d5	8.931	82	1177998	84.587	ng	0.00
42) 2,4,6-Tribromophenol	15.919	330	948405	144.154	ng	0.00
45) 2-Fluorobiphenyl	13.013	172	2831786	79.179	ng	0.00
79) Terphenyl-d14	19.925	244	4922368	84.162	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.331	88	158108	38.889	ng	99
3) Pyridine	3.726	79	436550	39.230	ng	97
4) n-Nitrosodimethylamine	3.631	42	221517	48.284	ng	97
6) Aniline	7.119	93	672688	37.518	ng	99
8) 2-Chlorophenol	7.367	128	584248	49.881	ng	100
9) Benzaldehyde	6.937	77	263740	28.288	ng	98
10) Phenol	6.996	94	701844	50.262	ng	99
11) bis(2-Chloroethyl)ether	7.214	93	510577	46.912	ng	99
12) 1,3-Dichlorobenzene	7.684	146	601240	46.552	ng	99
13) 1,4-Dichlorobenzene	7.825	146	610323	46.233	ng	100
14) 1,2-Dichlorobenzene	8.137	146	589005	46.093	ng	99
15) Benzyl Alcohol	8.025	79	504404	47.704	ng	99
16) 2,2'-oxybis(1-Chloropr...	8.308	45	689083	47.263	ng	98
17) 2-Methylphenol	8.225	107	482348	49.736	ng	99
18) Hexachloroethane	8.861	117	227208	46.811	ng	97
19) n-Nitroso-di-n-propyla...	8.584	70	393822	44.678	ng	98
20) 3+4-Methylphenols	8.549	107	660509	49.134	ng	97
22) Acetophenone	8.596	105	819556	46.389	ng	# 99
24) Nitrobenzene	8.972	77	591064	47.489	ng	99
25) Isophorone	9.496	82	1134513	46.032	ng	98
26) 2-Nitrophenol	9.678	139	314959	49.309	ng	97
27) 2,4-Dimethylphenol	9.737	122	558293	50.825	ng	100
28) bis(2-Chloroethoxy)met...	9.966	93	698931	46.914	ng	98
29) 2,4-Dichlorophenol	10.208	162	564908	51.769	ng	99
30) 1,2,4-Trichlorobenzene	10.419	180	553924	46.307	ng	99
31) Naphthalene	10.608	128	1710559	46.717	ng	99
32) Benzoic acid	9.896	122	439462	54.190	ng	98
33) 4-Chloroaniline	10.708	127	440457	27.233	ng	99
34) Hexachlorobutadiene	10.896	225	342743	46.008	ng	98
35) Caprolactam	11.502	113	210088m	52.506	ng	
36) 4-Chloro-3-methylphenol	11.837	107	646361	51.621	ng	99
37) 2-Methylnaphthalene	12.213	142	1107270	46.469	ng	100
38) 1-Methylnaphthalene	12.431	142	1168249	46.102	ng	100
40) 1,2,4,5-Tetrachloroben...	12.584	216	626241	44.362	ng	99
41) Hexachlorocyclopentadiene	12.566	237	771660	90.496	ng	98
43) 2,4,6-Trichlorophenol	12.825	196	479514	50.315	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.896	196	536382	51.353	ng	99
46) 1,1'-Biphenyl	13.225	154	1638603	46.160	ng	99
47) 2-Chloronaphthalene	13.266	162	1259783	45.586	ng	98
48) 2-Nitroaniline	13.466	65	415059	51.546	ng	98
49) Acenaphthylene	14.125	152	2177367	47.615	ng	100
50) Dimethylphthalate	13.854	163	1738994	48.585	ng	99
51) 2,6-Dinitrotoluene	13.966	165	395341	52.404	ng	98
52) Acenaphthene	14.472	154	1220137	45.457	ng	99
53) 3-Nitroaniline	14.296	138	300405	35.392	ng	97
54) 2,4-Dinitrophenol	14.513	184	495694	123.038	ng	99
55) Dibenzofuran	14.813	168	2007042	47.295	ng	100
56) 4-Nitrophenol	14.625	139	746663	107.051	ng	98
57) 2,4-Dinitrotoluene	14.766	165	583246	54.187	ng	99
58) Fluorene	15.466	166	1589403	47.465	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.037	232	472386	51.220	ng	100
60) Diethylphthalate	15.243	149	1844493	50.715	ng	99
61) 4-Chlorophenyl-phenyle...	15.460	204	789982	47.436	ng	98
62) 4-Nitroaniline	15.484	138	442261	50.825	ng	99
63) Azobenzene	15.754	77	1593888	51.181	ng	99
65) 4,6-Dinitro-2-methylph...	15.548	198	340236	52.303	ng	96
66) n-Nitrosodiphenylamine	15.678	169	1483025	46.603	ng	100
67) 4-Bromophenyl-phenylether	16.372	248	531300	46.291	ng	96
68) Hexachlorobenzene	16.495	284	616466	46.335	ng	98
69) Atrazine	16.660	200	590070	49.512	ng	97
70) Pentachlorophenol	16.848	266	843937	100.486	ng	97
71) Phenanthrene	17.254	178	2687053	46.873	ng	99
72) Anthracene	17.348	178	2780840	47.501	ng	100
73) Carbazole	17.619	167	2645585	47.978	ng	99
74) Di-n-butylphthalate	18.213	149	3375219	49.751	ng	100
75) Fluoranthene	19.336	202	3282061	47.997	ng	99
77) Benzidine	19.519	184	1590772	43.560	ng	100
78) Pyrene	19.713	202	3315062	47.664	ng	100
80) Butylbenzylphthalate	20.654	149	1618023	51.331	ng	98
81) Benzo(a)anthracene	21.630	228	3439380	48.736	ng	99
82) 3,3'-Dichlorobenzidine	21.548	252	816647	30.701	ng	99
83) Chrysene	21.701	228	3127598	47.324	ng	99
84) Bis(2-ethylhexyl)phtha...	21.566	149	2323944	50.084	ng	98
85) Di-n-octyl phthalate	22.860	149	4076449	51.076	ng	99
87) Indeno(1,2,3-cd)pyrene	28.859	276	3998661m	46.336	ng	
88) Benzo(b)fluoranthene	23.954	252	3354449	48.343	ng	100
89) Benzo(k)fluoranthene	24.019	252	3390357	48.827	ng	99
90) Benzo(a)pyrene	24.854	252	3251341	48.136	ng	100
91) Dibenzo(a,h)anthracene	28.971	278	3297613	46.760	ng	97
92) Benzo(g,h,i)perylene	30.054	276	3230088	46.595	ng	98

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

