

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111125\
 Data File : BP026093.D
 Acq On : 11 Nov 2025 16:52
 Operator : RC/JU
 Sample : Q3586-03MS
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SED-3MS

Manual Integrations
 APPROVED

Reviewed By :Rahul Chavli 11/12/2025
 Supervised By :Jagrut Upadhyay 11/12/2025

Quant Time: Nov 11 17:12:09 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP102925.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 30 11:51:34 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.678	152	296222	20.000	ng	-0.01
21) Naphthalene-d8	10.449	136	1166343	20.000	ng	0.00
39) Acenaphthene-d10	14.313	164	745428	20.000	ng	0.00
64) Phenanthrene-d10	17.125	188	1549759	20.000	ng	0.00
76) Chrysene-d12	21.577	240	1762341	20.000	ng	0.01
86) Perylene-d12	24.924	264	2068063	20.000	ng	0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.296	112	1778196	99.054	ng	0.00
7) Phenol-d6	6.849	99	2260243	98.920	ng	0.00
23) Nitrobenzene-d5	8.813	82	1249195	66.768	ng	-0.01
42) 2,4,6-Tribromophenol	15.825	330	939233	116.547	ng	0.00
45) 2-Fluorobiphenyl	12.919	172	3047663	58.748	ng	-0.01
79) Terphenyl-d14	19.866	244	5237660	60.381	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.225	88	284762	37.628	ng	98
3) Pyridine	3.614	79	786988	36.581	ng	97
4) n-Nitrosodimethylamine	3.525	42	362450	42.035	ng	90
6) Aniline	7.008	93	934451	30.643	ng	98
8) 2-Chlorophenol	7.249	128	946013	47.589	ng	97
9) Benzaldehyde	6.819	77	463796	39.047	ng	99
10) Phenol	6.878	94	1148006	45.257	ng	100
11) bis(2-Chloroethyl)ether	7.102	93	833295	41.626	ng	99
12) 1,3-Dichlorobenzene	7.566	146	1009516	44.220	ng	99
13) 1,4-Dichlorobenzene	7.713	146	1027189	44.502	ng	100
14) 1,2-Dichlorobenzene	8.025	146	980092	44.498	ng	100
15) Benzyl Alcohol	7.908	79	748510	45.441	ng	98
16) 2,2'-oxybis(1-Chloropr...	8.202	45	1165540	38.523	ng	98
17) 2-Methylphenol	8.113	107	766752	46.161	ng	99
18) Hexachloroethane	8.749	117	355877	45.057	ng	97
19) n-Nitroso-di-n-propyla...	8.472	70	609970	43.005	ng	97
20) 3+4-Methylphenols	8.431	107	1042051	45.095	ng	98
22) Acetophenone	8.484	105	1300398	44.085	ng	# 97
24) Nitrobenzene	8.855	77	906907	45.910	ng	98
25) Isophorone	9.384	82	1731602	43.240	ng	99
26) 2-Nitrophenol	9.560	139	459763	62.382	ng	96
27) 2,4-Dimethylphenol	9.625	122	854401	50.736	ng	98
28) bis(2-Chloroethoxy)met...	9.860	93	1086520	43.051	ng	100
29) 2,4-Dichlorophenol	10.096	162	868157	48.825	ng	98
30) 1,2,4-Trichlorobenzene	10.307	180	905744	46.866	ng	97
31) Naphthalene	10.496	128	2743320	43.132	ng	100
32) Benzoic acid	9.755	122	691403	61.547	ng	96
33) 4-Chloroaniline	10.596	127	470005	19.230	ng	99
34) Hexachlorobutadiene	10.790	225	548423	44.654	ng	98
35) Caprolactam	11.366	113	308604	49.424	ng	# 88
36) 4-Chloro-3-methylphenol	11.725	107	912347	48.947	ng	99
37) 2-Methylnaphthalene	12.113	142	1770111	39.764	ng	99
38) 1-Methylnaphthalene	12.331	142	1816726	42.008	ng	98
40) 1,2,4,5-Tetrachloroben...	12.484	216	1008627	43.548	ng	99
41) Hexachlorocyclopentadiene	12.472	237	1120849	132.265	ng	97
43) 2,4,6-Trichlorophenol	12.725	196	708087	51.722	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.790	196	788989	50.507	ng	99
46) 1,1'-Biphenyl	13.131	154	2485365	43.543	ng	99
47) 2-Chloronaphthalene	13.172	162	1920130	42.757	ng	99
48) 2-Nitroaniline	13.372	65	556442	48.138	ng	95
49) Acenaphthylene	14.031	152	3279242	50.100	ng	100
50) Dimethylphthalate	13.766	163	2430796	44.831	ng	100
51) 2,6-Dinitrotoluene	13.878	165	530255	52.105	ng	93
52) Acenaphthene	14.378	154	1822943	40.569	ng	100
53) 3-Nitroaniline	14.201	138	344027	29.253	ng	# 96
54) 2,4-Dinitrophenol	14.419	184	617165	110.722	ng	92
55) Dibenzofuran	14.719	168	2957373	43.561	ng	98
56) 4-Nitrophenol	14.525	139	1117576	103.365	ng	94
57) 2,4-Dinitrotoluene	14.678	165	777166	51.982	ng	91
58) Fluorene	15.384	166	2341092	44.015	ng	100
59) 2,3,4,6-Tetrachlorophenol	14.948	232	699686	56.966	ng	98
60) Diethylphthalate	15.160	149	2469705	45.307	ng	99
61) 4-Chlorophenyl-phenyle...	15.384	204	1137548	42.828	ng	98
62) 4-Nitroaniline	15.395	138	605617	52.270	ng	94
63) Azobenzene	15.684	77	2303527	46.628	ng	96
65) 4,6-Dinitro-2-methylph...	15.454	198	416417	61.165	ng	91
66) n-Nitrosodiphenylamine	15.595	169	2155394	44.482	ng	100
67) 4-Bromophenyl-phenylether	16.301	248	754222	43.874	ng	97
68) Hexachlorobenzene	16.419	284	877752	44.852	ng	96
69) Atrazine	16.578	200	819183	50.115	ng	98
70) Pentachlorophenol	16.772	266	1254838	104.976	ng	99
71) Phenanthrene	17.172	178	3856008	42.748	ng	100
72) Anthracene	17.260	178	3964031	44.333	ng	99
73) Carbazole	17.542	167	3883757	45.885	ng	100
74) Di-n-butylphthalate	18.154	149	4567374	48.225	ng	99
75) Fluoranthene	19.266	202	4810970	45.721	ng	98
77) Benzidine	19.460	184	2760993	61.722	ng	99
78) Pyrene	19.642	202	5019988	42.116	ng	99
80) Butylbenzylphthalate	20.607	149	2238758	50.144	ng	99
81) Benzo(a)anthracene	21.560	228	5361718	44.274	ng	100
82) 3,3'-Dichlorobenzidine	21.489	252	1330753	34.105	ng	99
83) Chrysene	21.625	228	5030796	43.854	ng	100
84) Bis(2-ethylhexyl)phtha...	21.525	149	3357041	47.305	ng	99
85) Di-n-octyl phthalate	22.819	149	5992300	57.119	ng	97
87) Indeno(1,2,3-cd)pyrene	28.718	276	6960277m	45.588	ng	
88) Benzo(b)fluoranthene	23.877	252	5541720	43.165	ng	99
89) Benzo(k)fluoranthene	23.948	252	5709385	43.549	ng	99
90) Benzo(a)pyrene	24.771	252	5501616	46.645	ng	99
91) Dibenzo(a,h)anthracene	28.818	278	5718465	46.024	ng	99
92) Benzo(g,h,i)perylene	29.895	276	5625013	47.046	ng	97

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

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