

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121225\
 Data File : BP026307.D
 Acq On : 12 Dec 2025 10:44
 Operator : RC/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SSTDCCC040

Quant Time: Dec 12 11:08:23 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP111825.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 05 14:52:44 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.648	152	311819	20.000	ng	0.00
21) Naphthalene-d8	10.413	136	1120151	20.000	ng	0.00
39) Acenaphthene-d10	14.277	164	664855	20.000	ng	-0.01
64) Phenanthrene-d10	17.089	188	1293189	20.000	ng	0.00
76) Chrysene-d12	21.536	240	1396095	20.000	ng	0.00
86) Perylene-d12	24.830	264	1677512	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.272	112	1571033	81.000	ng	0.00
7) Phenol-d6	6.837	99	1811356	73.986	ng	0.00
23) Nitrobenzene-d5	8.790	82	1654197	84.058	ng	0.00
42) 2,4,6-Tribromophenol	15.807	330	781169	90.348	ng	0.00
45) 2-Fluorobiphenyl	12.883	172	3969551	86.805	ng	-0.02
79) Terphenyl-d14	19.830	244	6011490	88.010	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.213	88	333156	42.241	ng	97
3) Pyridine	3.602	79	878709	38.794	ng	100
4) n-Nitrosodimethylamine	3.513	42	333876	39.526	ng	95
6) Aniline	6.984	93	1132086	35.239	ng	99
8) 2-Chlorophenol	7.225	128	866919	39.334	ng	99
9) Benzaldehyde	6.807	77	659810	49.992	ng	99
10) Phenol	6.860	94	960449	36.470	ng	95
11) bis(2-Chloroethyl)ether	7.084	93	777995	37.679	ng	97
12) 1,3-Dichlorobenzene	7.537	146	992776	41.627	ng	100
13) 1,4-Dichlorobenzene	7.684	146	1000480	41.629	ng	99
14) 1,2-Dichlorobenzene	7.995	146	946983	41.049	ng	100
15) Benzyl Alcohol	7.884	79	652501	36.663	ng	98
16) 2,2'-oxybis(1-Chloropr...	8.166	45	1041180	36.129	ng	98
17) 2-Methylphenol	8.090	107	659243	36.883	ng	98
18) Hexachloroethane	8.719	117	370465	42.436	ng	96
19) n-Nitroso-di-n-propyla...	8.448	70	528043	34.989	ng	98
20) 3+4-Methylphenols	8.413	107	871289	35.301	ng	99
22) Acetophenone	8.460	105	1154320	40.626	ng	100
24) Nitrobenzene	8.831	77	835747	42.014	ng	99
25) Isophorone	9.354	82	1537304	39.817	ng	99
26) 2-Nitrophenol	9.537	139	451442	44.887	ng	96
27) 2,4-Dimethylphenol	9.601	122	672512	37.780	ng	99
28) bis(2-Chloroethoxy)met...	9.831	93	951456	39.212	ng	100
29) 2,4-Dichlorophenol	10.066	162	766225	42.237	ng	99
30) 1,2,4-Trichlorobenzene	10.278	180	828268	44.416	ng	99
31) Naphthalene	10.466	128	2520150	41.615	ng	99
32) Benzoic acid	9.760	122	660362	46.310	ng	99
33) 4-Chloroaniline	10.566	127	981425	38.685	ng	99
34) Hexachlorobutadiene	10.754	225	573826	47.129	ng	99
35) Caprolactam	11.360	113	229353	35.802	ng	95
36) 4-Chloro-3-methylphenol	11.701	107	756361	39.626	ng	99
37) 2-Methylnaphthalene	12.072	142	1708079	40.020	ng	98
38) 1-Methylnaphthalene	12.301	142	1675180	40.175	ng	99
40) 1,2,4,5-Tetrachloroben...	12.442	216	924968	45.575	ng	100
41) Hexachlorocyclopentadiene	12.431	237	500232	39.513	ng	98
43) 2,4,6-Trichlorophenol	12.695	196	610189	44.238	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.772	196	666526	43.388	ng	98
46) 1,1'-Biphenyl	13.101	154	2111084	42.051	ng	99
47) 2-Chloronaphthalene	13.130	162	1687798	42.671	ng	99
48) 2-Nitroaniline	13.342	65	450173	41.554	ng	96
49) Acenaphthylene	13.995	152	2716872	41.857	ng	100
50) Dimethylphthalate	13.736	163	2093854	42.266	ng	99
51) 2,6-Dinitrotoluene	13.842	165	434492	44.381	ng	99
52) Acenaphthene	14.348	154	1556802	41.019	ng	100
53) 3-Nitroaniline	14.177	138	451172	39.424	ng	97
54) 2,4-Dinitrophenol	14.395	184	337105	64.586	ng	96
55) Dibenzofuran	14.689	168	2480135	42.020	ng	99
56) 4-Nitrophenol	14.507	139	366854	35.901	ng	95
57) 2,4-Dinitrotoluene	14.654	165	637345	43.763	ng	95
58) Fluorene	15.348	166	1939398	41.535	ng	99
59) 2,3,4,6-Tetrachlorophenol	14.919	232	565967	43.527	ng	98
60) Diethylphthalate	15.130	149	2085357	41.969	ng	100
61) 4-Chlorophenyl-phenyle...	15.348	204	999931	42.949	ng	97
62) 4-Nitroaniline	15.366	138	464972	39.148	ng	95
63) Azobenzene	15.642	77	1686773	40.372	ng	99
65) 4,6-Dinitro-2-methylph...	15.436	198	387473	50.730	ng	95
66) n-Nitrosodiphenylamine	15.566	169	1673759	41.861	ng	99
67) 4-Bromophenyl-phenylether	16.266	248	650388	45.366	ng	98
68) Hexachlorobenzene	16.383	284	771132	46.281	ng	95
69) Atrazine	16.554	200	622891	41.434	ng	98
70) Pentachlorophenol	16.736	266	510022	44.124	ng	99
71) Phenanthrene	17.130	178	3041991	41.727	ng	100
72) Anthracene	17.224	178	3140388	42.293	ng	100
73) Carbazole	17.507	167	2789021	39.675	ng	99
74) Di-n-butylphthalate	18.107	149	3588402	43.208	ng	99
75) Fluoranthene	19.230	202	3711845	41.836	ng	98
77) Benzidine	19.418	184	902695	20.594	ng	98
78) Pyrene	19.607	202	3843478	42.247	ng	100
80) Butylbenzylphthalate	20.565	149	1702391	42.637	ng	98
81) Benzo(a)anthracene	21.512	228	3933483	41.073	ng	100
82) 3,3'-Dichlorobenzidine	21.436	252	1410708	40.486	ng	99
83) Chrysene	21.583	228	3745037	41.300	ng	99
84) Bis(2-ethylhexyl)phtha...	21.465	149	2535340	42.214	ng	99
85) Di-n-octyl phthalate	22.730	149	4480918	42.319	ng	99
87) Indeno(1,2,3-cd)pyrene	28.577	276	5514018	43.493	ng	# 92
88) Benzo(b)fluoranthene	23.801	252	4220400	41.396	ng	99
89) Benzo(k)fluoranthene	23.865	252	4132090	40.518	ng	99
90) Benzo(a)pyrene	24.689	252	4014382	41.525	ng	99
91) Dibenzo(a,h)anthracene	28.653	278	4484023	43.242	ng	98
92) Benzo(g,h,i)perylene	29.747	276	4489584	43.430	ng	97

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

