

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080825\
 Data File : BP025326.D
 Acq On : 08 Aug 2025 10:15
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Reviewed By :Rahul
 Chavli

Quant Time: Aug 08 11:12:11 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP080525.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 06 06:41:44 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.802	152	240709	20.000	ng	0.00
21) Naphthalene-d8	10.572	136	976682	20.000	ng	0.00
39) Acenaphthene-d10	14.419	164	575605	20.000	ng	0.00
64) Phenanthrene-d10	17.225	188	1048506	20.000	ng	0.00
76) Chrysene-d12	21.666	240	989888	20.000	ng	0.00
86) Perylene-d12	25.059	264	1196041	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.408	112	1139134	76.114	ng	0.00
7) Phenol-d6	6.960	99	1423958	72.295	ng	0.00
23) Nitrobenzene-d5	8.943	82	1455838	82.581	ng	0.00
42) 2,4,6-Tribromophenol	15.936	330	491175	85.536	ng	0.01
45) 2-Fluorobiphenyl	13.031	172	3287807	78.893	ng	0.00
79) Terphenyl-d14	19.936	244	4177606	79.712	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.331	88	222915	37.744	ng	97
3) Pyridine	3.725	79	586352	35.734	ng	97
4) n-Nitrosodimethylamine	3.631	42	260317	34.172	ng	97
6) Aniline	7.131	93	967841	35.795	ng	98
8) 2-Chlorophenol	7.366	128	633395	38.664	ng	98
9) Benzaldehyde	6.943	77	668759	47.805	ng	97
10) Phenol	6.990	94	747065	35.449	ng	99
11) bis(2-Chloroethyl)ether	7.219	93	607130	36.768	ng	98
12) 1,3-Dichlorobenzene	7.690	146	706677	38.391	ng	99
13) 1,4-Dichlorobenzene	7.831	146	717825	38.546	ng	99
14) 1,2-Dichlorobenzene	8.149	146	684747	37.957	ng	99
15) Benzyl Alcohol	8.025	79	543783	35.616	ng	98
16) 2,2'-oxybis(1-Chloropr...	8.319	45	878026	33.660	ng	98
17) 2-Methylphenol	8.225	107	513129	36.388	ng	99
18) Hexachloroethane	8.872	117	263322	39.538	ng	97
19) n-Nitroso-di-n-propyla...	8.596	70	452293	34.950	ng	97
20) 3+4-Methylphenols	8.549	107	690868	35.579	ng	98
22) Acetophenone	8.607	105	909064	37.077	ng	# 97
24) Nitrobenzene	8.978	77	652814	39.118	ng	97
25) Isophorone	9.501	82	1272966	36.275	ng	100
26) 2-Nitrophenol	9.684	139	311077	43.187	ng	98
27) 2,4-Dimethylphenol	9.743	122	591009	36.530	ng	96
28) bis(2-Chloroethoxy)met...	9.984	93	809318	37.912	ng	100
29) 2,4-Dichlorophenol	10.219	162	567733	40.082	ng	97
30) 1,2,4-Trichlorobenzene	10.437	180	625060	39.999	ng	100
31) Naphthalene	10.619	128	1934670	38.158	ng	100
32) Benzoic acid	9.854	122	376644	42.801	ng	100
33) 4-Chloroaniline	10.719	127	839289	37.219	ng	98
34) Hexachlorobutadiene	10.913	225	374078	41.702	ng	98
35) Caprolactam	11.478	113	182502	33.213	ng	96
36) 4-Chloro-3-methylphenol	11.837	107	597779	36.473	ng	96
37) 2-Methylnaphthalene	12.231	142	1200862	36.592	ng	99
38) 1-Methylnaphthalene	12.448	142	1258723	36.759	ng	99
40) 1,2,4,5-Tetrachloroben...	12.595	216	640167	40.736	ng	99
41) Hexachlorocyclopentadiene	12.584	237	447894	46.179	ng	98
43) 2,4,6-Trichlorophenol	12.842	196	435004	42.850	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.907	196	463206	40.675	ng	99
46) 1,1'-Biphenyl	13.248	154	1681680	39.644	ng	100
47) 2-Chloronaphthalene	13.284	162	1269125	39.216	ng	99
48) 2-Nitroaniline	13.478	65	351935	37.795	ng	93
49) Acenaphthylene	14.137	152	2093588	38.783	ng	99
50) Dimethylphthalate	13.872	163	1561021	37.202	ng	99
51) 2,6-Dinitrotoluene	13.978	165	321759	41.572	ng	91
52) Acenaphthene	14.489	154	1274700	38.713	ng	99
53) 3-Nitroaniline	14.307	138	370794	40.941	ng	96
54) 2,4-Dinitrophenol	14.525	184	141775	45.329	ng	# 46
55) Dibenzofuran	14.831	168	1876057	38.127	ng	99
56) 4-Nitrophenol	14.619	139	302516	37.847	ng	96
57) 2,4-Dinitrotoluene	14.784	165	439835	37.428	ng	97
58) Fluorene	15.489	166	1442226	37.318	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.054	232	383749	40.341	ng	98
60) Diethylphthalate	15.266	149	1525689	36.065	ng	99
61) 4-Chlorophenyl-phenyle...	15.484	204	704373	38.445	ng	100
62) 4-Nitroaniline	15.489	138	361326	40.264	ng	95
63) Azobenzene	15.784	77	1427610	36.326	ng	99
65) 4,6-Dinitro-2-methylph...	15.560	198	207397	45.566	ng	98
66) n-Nitrosodiphenylamine	15.701	169	1260046	39.469	ng	100
67) 4-Bromophenyl-phenylether	16.401	248	435742	41.303	ng	96
68) Hexachlorobenzene	16.519	284	483113	40.812	ng	97
69) Atrazine	16.678	200	448699	39.965	ng	99
70) Pentachlorophenol	16.872	266	323077	42.838	ng	100
71) Phenanthrene	17.272	178	2241309	39.028	ng	100
72) Anthracene	17.360	178	2253461	38.917	ng	100
73) Carbazole	17.636	167	2132429	38.364	ng	100
74) Di-n-butylphthalate	18.242	149	2726922	40.553	ng	99
75) Fluoranthene	19.354	202	2566801	38.610	ng	99
77) Benzidine	19.536	184	2224681	62.093	ng	100
78) Pyrene	19.724	202	2597289	39.457	ng	100
80) Butylbenzylphthalate	20.671	149	1249709	46.812	ng	99
81) Benzo(a)anthracene	21.648	228	2561922	38.946	ng	100
82) 3,3'-Dichlorobenzidine	21.554	252	1005276	43.314	ng	98
83) Chrysene	21.713	228	2387480	39.131	ng	99
84) Bis(2-ethylhexyl)phtha...	21.595	149	1891251	44.366	ng	100
85) Di-n-octyl phthalate	22.901	149	3396499	47.602	ng	98
87) Indeno(1,2,3-cd)pyrene	28.900	276	3393119m	39.580	ng	
88) Benzo(b)fluoranthene	23.971	252	2704358	37.772	ng	100
89) Benzo(k)fluoranthene	24.048	252	2648921	37.321	ng	99
90) Benzo(a)pyrene	24.895	252	2674274	38.786	ng	100
91) Dibenzo(a,h)anthracene	28.995	278	2769517	39.446	ng	99
92) Benzo(g,h,i)perylene	30.089	276	2697139	39.142	ng	99

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

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