

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092425\
 Data File : BP025845.D
 Acq On : 24 Sep 2025 14:41
 Operator : CG/JU
 Sample : Q3159-01MSD
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 VNJ-245MSD

Manual Integrations
 APPROVED

Reviewed By :Rahul Chavli 09/25/2025
 Supervised By :Jagrut Upadhyay 09/25/2025

Quant Time: Sep 24 15:03:04 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP091225.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 11 16:45:04 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.743	152	152433	20.000	ng	-0.02
21) Naphthalene-d8	10.513	136	642516	20.000	ng	-0.01
39) Acenaphthene-d10	14.372	164	449880	20.000	ng	0.00
64) Phenanthrene-d10	17.178	188	981325	20.000	ng	0.00
76) Chrysene-d12	21.619	240	971996	20.000	ng	-0.02
86) Perylene-d12	24.977	264	1089001	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.384	112	789261	83.546	ng	0.00
7) Phenol-d6	6.955	99	1071951	85.368	ng	0.00
23) Nitrobenzene-d5	8.890	82	638880	43.861	ng	-0.01
42) 2,4,6-Tribromophenol	15.889	330	505896	90.156	ng	-0.02
45) 2-Fluorobiphenyl	12.984	172	1486713	44.002	ng	0.00
79) Terphenyl-d14	19.901	244	2643241	48.922	ng	-0.01
Target Compounds						Qvalue
2) 1,4-Dioxane	3.302	88	151194	38.132	ng	97
3) Pyridine	3.702	79	421056	38.566	ng	93
4) n-Nitrosodimethylamine	3.608	42	206365	40.057	ng	88
6) Aniline	7.084	93	426583	24.748	ng	99
8) 2-Chlorophenol	7.331	128	513522	49.550	ng	98
9) Benzaldehyde	6.902	77	290760	38.632	ng	99
10) Phenol	6.978	94	654088	49.401	ng	97
11) bis(2-Chloroethyl)ether	7.178	93	477622	44.899	ng	96
12) 1,3-Dichlorobenzene	7.637	146	528794	44.831	ng	97
13) 1,4-Dichlorobenzene	7.784	146	549703	45.751	ng	99
14) 1,2-Dichlorobenzene	8.096	146	527885	45.225	ng	98
15) Benzyl Alcohol	7.990	79	481357	46.391	ng	97
16) 2,2'-oxybis(1-Chloropr...	8.260	45	734247	42.207	ng	98
17) 2-Methylphenol	8.202	107	439962	49.286	ng	99
18) Hexachloroethane	8.813	117	205716	44.824	ng	98
19) n-Nitroso-di-n-propyla...	8.549	70	387626	44.259	ng	96
20) 3+4-Methylphenols	8.531	107	595930	47.668	ng	98
22) Acetophenone	8.560	105	842940	49.093	ng	# 99
24) Nitrobenzene	8.937	77	575680	44.056	ng	97
25) Isophorone	9.460	82	1121036	43.787	ng	99
26) 2-Nitrophenol	9.637	139	289604	49.918	ng	94
27) 2,4-Dimethylphenol	9.707	122	496294	48.762	ng	99
28) bis(2-Chloroethoxy)met...	9.925	93	678294	45.833	ng	99
29) 2,4-Dichlorophenol	10.190	162	511824	50.437	ng	99
30) 1,2,4-Trichlorobenzene	10.384	180	525655	45.962	ng	99
31) Naphthalene	10.566	128	1558447	44.983	ng	99
32) Benzoic acid	9.907	122	449198	59.362	ng	97
33) 4-Chloroaniline	10.684	127	256373	17.754	ng	99
34) Hexachlorobutadiene	10.843	225	334592	45.447	ng	100
35) Caprolactam	11.490	113	189259	48.381	ng	95
36) 4-Chloro-3-methylphenol	11.831	107	595223	49.198	ng	97
37) 2-Methylnaphthalene	12.178	142	1041680	46.813	ng	99
38) 1-Methylnaphthalene	12.395	142	1090767	45.968	ng	100
40) 1,2,4,5-Tetrachloroben...	12.548	216	702193	51.509	ng	99
41) Hexachlorocyclopentadiene	12.525	237	522588	70.384	ng	98
43) 2,4,6-Trichlorophenol	12.801	196	450018	50.214	ng	98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.895	196	495149	49.643	ng	95
46) 1,1'-Biphenyl	13.195	154	1677730	50.807	ng	99
47) 2-Chloronaphthalene	13.242	162	1197961	46.074	ng	98
48) 2-Nitroaniline	13.448	65	403742	46.582	ng	97
49) Acenaphthylene	14.089	152	2034141	47.217	ng	100
50) Dimethylphthalate	13.825	163	1560647	45.438	ng	100
51) 2,6-Dinitrotoluene	13.948	165	357875	49.180	ng	94
52) Acenaphthene	14.437	154	1128044	45.404	ng	100
53) 3-Nitroaniline	14.289	138	218551	28.560	ng	99
54) 2,4-Dinitrophenol	14.501	184	435998	91.419	ng	99
55) Dibenzofuran	14.778	168	1852056	45.739	ng	98
56) 4-Nitrophenol	14.654	139	548930	83.114	ng	97
57) 2,4-Dinitrotoluene	14.748	165	517775	49.997	ng	94
58) Fluorene	15.436	166	1509318	46.876	ng	100
59) 2,3,4,6-Tetrachlorophenol	15.025	232	452737	50.389	ng	96
60) Diethylphthalate	15.201	149	1657757	47.582	ng	100
61) 4-Chlorophenyl-phenyle...	15.431	204	767272	47.284	ng	96
62) 4-Nitroaniline	15.472	138	395195	50.411	ng	96
63) Azobenzene	15.731	77	1660111	49.280	ng	98
65) 4,6-Dinitro-2-methylph...	15.536	198	316584	48.510	ng	87
66) n-Nitrosodiphenylamine	15.654	169	1381799	46.013	ng	100
67) 4-Bromophenyl-phenylether	16.348	248	526507	48.838	ng	96
68) Hexachlorobenzene	16.472	284	583506	48.535	ng	94
69) Atrazine	16.631	200	579205	50.457	ng	97
70) Pentachlorophenol	16.831	266	732463	92.311	ng	99
71) Phenanthrene	17.225	178	2565970	46.095	ng	100
72) Anthracene	17.313	178	2601847	46.066	ng	100
73) Carbazole	17.595	167	2446842	46.598	ng	99
74) Di-n-butylphthalate	18.183	149	3090974	47.967	ng	100
75) Fluoranthene	19.307	202	3226908	48.335	ng	99
77) Benzidine	19.507	184	1606932	57.393	ng	100
78) Pyrene	19.683	202	3376227	52.057	ng	100
80) Butylbenzylphthalate	20.630	149	1457356	51.248	ng	97
81) Benzo(a)anthracene	21.601	228	3179373	47.778	ng	100
82) 3,3'-Dichlorobenzidine	21.524	252	806783	35.064	ng	100
83) Chrysene	21.671	228	3027196	47.722	ng	99
84) Bis(2-ethylhexyl)phtha...	21.536	149	2097741	50.421	ng	100
85) Di-n-octyl phthalate	22.807	149	3628821	49.034	ng	100
87) Indeno(1,2,3-cd)pyrene	28.818	276	3757549	47.447	ng	# 85
88) Benzo(b)fluoranthene	23.930	252	3204787	48.124	ng	98
89) Benzo(k)fluoranthene	23.995	252	3196181	47.034	ng	99
90) Benzo(a)pyrene	24.818	252	3053629	46.823	ng	99
91) Dibenzo(a,h)anthracene	28.895	278	3043235	46.958	ng	99
92) Benzo(g,h,i)perylene	30.012	276	3057868m	47.979	ng	

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

