Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060625\
Data File : BP024865.D
Acq On : 06 Jun 2025 13:56
Operator : RC/JU
Sample : SSTDICC050
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SSTDICC050

Quant Time: Jun 06 16:03:11 2025
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP060625.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 06 15:53:04 2025
Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.613	152	235518	20.000	ng	0.00
21) Naphthalene-d8	10.384	136	975092	20.000	ng	0.00
39) Acenaphthene-d10	14.260	164	631564	20.000	ng	0.01
64) Phenanthrene-d10	17.072	188	1197709	20.000	ng	0.01
76) Chrysene-d12	21.501	240	1276071	20.000	ng	0.02
86) Perylene-d12	24.742	264	1475251	20.000	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.243	112	1510671	107.067	ng	0.00
7) Phenol-d6	6.813	99	1973616	105.719	ng	0.00
23) Nitrobenzene-d5	8.760	82	2165794	107.931	ng	0.00
42) 2,4,6-Tribromophenol	15.784	330	940592	107.721	ng	0.00
45) 2-Fluorobiphenyl	12.860	172	4934980	105.263	ng	0.00
79) Terphenyl-d14	19.801	244	7428863	104.339	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.167	88	321356	51.761	ng	100
3) Pyridine	3.567	79	806405	54.010	ng	98
4) n-Nitrosodimethylamine	3.478	42	319345	52.317	ng	98
6) Aniline	6.955	93	1263250	53.024	ng	100
8) 2-Chlorophenol	7.190	128	855311	53.473	ng	99
9) Benzaldehyde	6.766	77	579996	53.748	ng	99
10) Phenol	6.843	94	1015401	52.761	ng	99
11) bis(2-Chloroethyl)ether	7.049	93	802393	53.010	ng	99
12) 1,3-Dichlorobenzene	7.502	146	925262	51.854	ng	99
13) 1,4-Dichlorobenzene	7.643	146	944146	52.440	ng	100
14) 1,2-Dichlorobenzene	7.960	146	906741	51.287	ng	99
15) Benzyl Alcohol	7.860	79	758909	52.968	ng	100
16) 2,2'-oxybis(1-Chloropr...	8.131	45	1030911	52.035	ng	99
17) 2-Methylphenol	8.066	107	712455	53.013	ng	100
18) Hexachloroethane	8.672	117	359979	53.107	ng	99
19) n-Nitroso-di-n-propyla...	8.413	70	671592	52.919	ng	99
20) 3+4-Methylphenols	8.396	107	960504	52.385	ng	97
22) Acetophenone	8.431	105	1304864	52.966	ng	99
24) Nitrobenzene	8.802	77	955872	53.620	ng	99
25) Isophorone	9.325	82	1862532	53.556	ng	100
26) 2-Nitrophenol	9.513	139	489550	55.657	ng	99
27) 2,4-Dimethylphenol	9.578	122	806896	53.602	ng	99
28) bis(2-Chloroethoxy)met...	9.801	93	1133590	54.643	ng	99
29) 2,4-Dichlorophenol	10.054	162	797316	55.201	ng	99
30) 1,2,4-Trichlorobenzene	10.243	180	857570	52.640	ng	98
31) Naphthalene	10.437	128	2630251	52.637	ng	100
32) Benzoic acid	9.772	122	561400	55.936	ng	98
33) 4-Chloroaniline	10.548	127	1148401	54.855	ng	99
34) Hexachlorobutadiene	10.713	225	532031	54.184	ng	98
35) Caprolactam	11.354	113	288081	54.183	ng	98
36) 4-Chloro-3-methylphenol	11.701	107	912092	54.747	ng	99
37) 2-Methylnaphthalene	12.048	142	1697018	53.539	ng	99
38) 1-Methylnaphthalene	12.272	142	1807332	53.333	ng	98
40) 1,2,4,5-Tetrachloroben...	12.419	216	949667	52.992	ng	100
41) Hexachlorocyclopentadiene	12.395	237	638568	58.420	ng	100
43) 2,4,6-Trichlorophenol	12.678	196	649483	53.966	ng	99

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Compound	R.T.	QIon	Response	Conc	Units		Dev(Min)
44) 2,4,5-Trichlorophenol	12.760	196	707702	54.908	ng		99
46) 1,1'-Biphenyl	13.072	154	2382821	51.916	ng		99
47) 2-Chloronaphthalene	13.113	162	1849683	52.435	ng		100
48) 2-Nitroaniline	13.337	65	586484	54.079	ng		99
49) Acenaphthylene	13.978	152	3062130	52.061	ng		100
50) Dimethylphthalate	13.719	163	2447212	52.527	ng		100
51) 2,6-Dinitrotoluene	13.837	165	535050	53.272	ng		98
52) Acenaphthene	14.325	154	1716357	50.924	ng		99
53) 3-Nitroaniline	14.178	138	580183	55.664	ng		99
54) 2,4-Dinitrophenol	14.401	184	321110	51.248	ng		95
55) Dibenzofuran	14.666	168	2773911	51.272	ng		99
56) 4-Nitrophenol	14.531	139	435101	51.384	ng		96
57) 2,4-Dinitrotoluene	14.648	165	768586	54.706	ng		98
58) Fluorene	15.331	166	2264731	51.819	ng		100
59) 2,3,4,6-Tetrachlorophenol	14.907	232	623685	54.063	ng		100
60) Diethylphthalate	15.101	149	2438733	52.523	ng		99
61) 4-Chlorophenyl-phenyle...	15.319	204	1118999	52.365	ng		99
62) 4-Nitroaniline	15.372	138	530012	56.050	ng		96
63) Azobenzene	15.619	77	2249353	52.821	ng		100
65) 4,6-Dinitro-2-methylph...	15.431	198	439647	55.863	ng		98
66) n-Nitrosodiphenylamine	15.542	169	1974147	53.178	ng		100
67) 4-Bromophenyl-phenylether	16.236	248	737798	54.595	ng		97
68) Hexachlorobenzene	16.354	284	867444	52.948	ng		99
69) Atrazine	16.531	200	732007	54.392	ng		98
70) Pentachlorophenol	16.725	266	486264	57.349	ng		99
71) Phenanthrene	17.113	178	3477677	52.543	ng		100
72) Anthracene	17.201	178	3558691	53.089	ng		99
73) Carbazole	17.495	167	3330211	53.599	ng		100
74) Di-n-butylphthalate	18.077	149	4253447	55.253	ng		100
75) Fluoranthene	19.201	202	4024039	52.455	ng		100
77) Benzidine	19.413	184	2202092	55.718	ng		100
78) Pyrene	19.577	202	4217295	52.900	ng		100
80) Butylbenzylphthalate	20.530	149	2045307	56.010	ng		99
81) Benzo(a)anthracene	21.477	228	4296840	52.648	ng		100
82) 3,3'-Dichlorobenzidine	21.401	252	1730292	53.402	ng		99
83) Chrysene	21.548	228	4080671	52.768	ng		99
84) Bis(2-ethylhexyl)phtha...	21.407	149	2938577	56.166	ng		99
85) Di-n-octyl phthalate	22.642	149	5061827	54.889	ng		100
87) Indeno(1,2,3-cd)pyrene	28.424	276	5749111	53.412	ng	#	93
88) Benzo(b)fluoranthene	23.701	252	4545535	53.840	ng		100
89) Benzo(k)fluoranthene	23.771	252	4641829	54.012	ng		100
90) Benzo(a)pyrene	24.583	252	4478222	54.313	ng		99
91) Dibenzo(a,h)anthracene	28.500	278	4715596	53.822	ng		99
92) Benzo(g,h,i)perylene	29.547	276	4649887	53.490	ng		100

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

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