Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060625\
Data File : BP024868.D
Acq On : 06 Jun 2025 17:09
Operator : RC/JU
Sample : SSTDICV040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
ICVBP060625

Quant Time: Jun 06 17:28:57 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 16:20:27 2025
Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By :Rahul Chavli 06/09/2025
Supervised By :Jagrut Upadhyay 06/09/2025

Compound	R.T.	QIon	Response	Conc	Units		Dev(Min)
Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.596	152	259470	20.000	ng		-0.02
21) Naphthalene-d8	10.366	136	1129880	20.000	ng		-0.01
39) Acenaphthene-d10	14.248	164	734874	20.000	ng		0.00
64) Phenanthrene-d10	17.054	188	1466437	20.000	ng		0.00
76) Chrysene-d12	21.489	240	1557824	20.000	ng		0.00
86) Perylene-d12	24.736	264	1860226	20.000	ng		0.02
System Monitoring Compounds							
5) 2-Fluorophenol	5.231	112	1138471	73.239	ng		-0.01
7) Phenol-d6	6.802	99	1573129	76.488	ng		-0.01
23) Nitrobenzene-d5	8.749	82	1561363	67.150	ng		-0.01
42) 2,4,6-Tribromophenol	15.778	330	743311	73.160	ng		0.00
45) 2-Fluorobiphenyl	12.848	172	3405857	62.434	ng		-0.01
79) Terphenyl-d14	19.789	244	5436960	62.551	ng		0.00
Target Compounds							Qvalue
2) 1,4-Dioxane	3.161	88	241705	35.338	ng		98
3) Pyridine	3.555	79	659477	40.092	ng		98
4) n-Nitrosodimethylamine	3.467	42	278482	41.411	ng		94
6) Aniline	6.937	93	1054913	40.192	ng		100
8) 2-Chlorophenol	7.172	128	687858	39.034	ng		97
9) Benzaldehyde	6.755	77	576121m	48.599	ng		
10) Phenol	6.825	94	845877	39.895	ng		98
11) bis(2-Chloroethyl)ether	7.031	93	656501	39.368	ng		99
12) 1,3-Dichlorobenzene	7.490	146	735256	37.402	ng		99
13) 1,4-Dichlorobenzene	7.631	146	742188	37.417	ng		99
14) 1,2-Dichlorobenzene	7.943	146	722064	37.071	ng		99
15) Benzyl Alcohol	7.843	79	644939	40.859	ng		98
16) 2,2'-oxybis(1-Chloropr...	8.125	45	849644	38.927	ng		98
17) 2-Methylphenol	8.055	107	591959	39.981	ng		98
18) Hexachloroethane	8.660	117	277433	37.151	ng		98
19) n-Nitroso-di-n-propyla...	8.402	70	559389	40.009	ng		99
20) 3+4-Methylphenols	8.384	107	797217	39.466	ng		99
22) Acetophenone	8.419	105	1100455	38.549	ng	#	99
24) Nitrobenzene	8.790	77	810932	39.258	ng		99
25) Isophorone	9.313	82	1576033	39.109	ng		99
26) 2-Nitrophenol	9.496	139	403179	39.558	ng		99
27) 2,4-Dimethylphenol	9.566	122	676106	38.761	ng		99
28) bis(2-Chloroethoxy)met...	9.784	93	921897	38.351	ng		99
29) 2,4-Dichlorophenol	10.043	162	653496	39.046	ng		99
30) 1,2,4-Trichlorobenzene	10.231	180	689288	36.514	ng		100
31) Naphthalene	10.413	128	2147539	37.089	ng		100
32) Benzoic acid	9.749	122	476446	40.420	ng		99
33) 4-Chloroaniline	10.543	127	960185	39.581	ng		99
34) Hexachlorobutadiene	10.696	225	413168	36.314	ng		100
35) Caprolactam	11.337	113	254008	41.229	ng		98
36) 4-Chloro-3-methylphenol	11.690	107	781340	40.474	ng		100
37) 2-Methylnaphthalene	12.037	142	1411483	38.430	ng		99
38) 1-Methylnaphthalene	12.254	142	1479329	37.674	ng		97
40) 1,2,4,5-Tetrachloroben...	12.413	216	774379	37.136	ng		100
41) Hexachlorocyclopentadiene	12.384	237	485560	38.177	ng		97
43) 2,4,6-Trichlorophenol	12.666	196	543521	38.812	ng		98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.760	196	585859	39.064	ng	99
46) 1,1'-Biphenyl	13.066	154	2002791	37.502	ng	99
47) 2-Chloronaphthalene	13.107	162	1514031	36.886	ng	100
48) 2-Nitroaniline	13.331	65	515024	40.813	ng	99
49) Acenaphthylene	13.966	152	2551023	37.274	ng	100
50) Dimethylphthalate	13.707	163	2015820	37.185	ng	99
51) 2,6-Dinitrotoluene	13.831	165	445393	38.111	ng	95
52) Acenaphthene	14.313	154	1469566	37.472	ng	98
53) 3-Nitroaniline	14.172	138	492714	40.626	ng	100
54) 2,4-Dinitrophenol	14.401	184	266900	38.197	ng	98
55) Dibenzofuran	14.660	168	2304340	36.605	ng	98
56) 4-Nitrophenol	14.531	139	362887	38.363	ng	94
57) 2,4-Dinitrotoluene	14.637	165	636513	38.936	ng	99
58) Fluorene	15.319	166	1883765	37.043	ng	99
59) 2,3,4,6-Tetrachlorophenol	14.901	232	514561	38.401	ng	100
60) Diethylphthalate	15.101	149	2033664	37.642	ng	99
61) 4-Chlorophenyl-phenyle...	15.313	204	915384	36.814	ng	98
62) 4-Nitroaniline	15.366	138	464418	42.233	ng	99
63) Azobenzene	15.619	77	1893664	38.217	ng	99
65) 4,6-Dinitro-2-methylph...	15.425	198	375376	38.956	ng	97
66) n-Nitrosodiphenylamine	15.542	169	1674745	36.846	ng	99
67) 4-Bromophenyl-phenylether	16.237	248	593842	35.890	ng	98
68) Hexachlorobenzene	16.348	284	725771	36.182	ng	100
69) Atrazine	16.531	200	611440	37.108	ng	99
70) Pentachlorophenol	16.719	266	407441	39.247	ng	99
71) Phenanthrene	17.107	178	2965530	36.595	ng	99
72) Anthracene	17.207	178	3028697	36.903	ng	99
73) Carbazole	17.489	167	2872120	37.755	ng	100
74) Di-n-butylphthalate	18.072	149	3440228	36.500	ng	100
75) Fluoranthene	19.195	202	3413021	36.337	ng	100
77) Benzidine	19.395	184	2156506	44.696	ng	100
78) Pyrene	19.572	202	3527120	36.241	ng	99
80) Butylbenzylphthalate	20.519	149	1664895	37.347	ng	99
81) Benzo(a)anthracene	21.472	228	3660732	36.741	ng	100
82) 3,3'-Dichlorobenzidine	21.395	252	1487801	37.613	ng	98
83) Chrysene	21.536	228	3438391	36.421	ng	99
84) Bis(2-ethylhexyl)phtha...	21.407	149	2365859	37.041	ng	100
85) Di-n-octyl phthalate	22.642	149	4126479	36.654	ng	100
87) Indeno(1,2,3-cd)pyrene	28.436	276	5100289	37.578	ng	100
88) Benzo(b)fluoranthene	23.718	252	3910900	36.735	ng	99
89) Benzo(k)fluoranthene	23.795	252	3904662	36.032	ng	100
90) Benzo(a)pyrene	24.595	252	3839776	36.932	ng	100
91) Dibenzo(a,h)anthracene	28.501	278	4123111	37.321	ng	99
92) Benzo(g,h,i)perylene	29.577	276	4097986	37.385	ng	99

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

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