Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF043025\
Data File : BF142244.D
Acq On : 30 Apr 2025 13:46
Operator : RC/JU
Sample : SSTDICC050
Misc :
ALS Vial : 7 Sample Multiplier: 1

Quant Time: Apr 30 15:43:55 2025
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF043025.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Apr 30 15:28:40 2025
Response via : Initial Calibration

Instrument :
BNA_F
ClientSampleId :
SSTDICC050

Manual Integrations APPROVED

Reviewed By :Rahul Chavli 05/01/2025
Supervised By :Jagrut Upadhyay 05/01/2025

Compound	R.T.	QIon	Response	Conc	Units		Dev(Min)
Internal Standards							
1) 1,4-Dichlorobenzene-d4	6.910	152	179073	20.000	ng		0.00
21) Naphthalene-d8	8.192	136	691236	20.000	ng		0.00
39) Acenaphthene-d10	9.951	164	360144	20.000	ng		0.00
64) Phenanthrene-d10	11.433	188	604171	20.000	ng		0.00
76) Chrysene-d12	14.074	240	322957	20.000	ng		0.00
86) Perylene-d12	15.568	264	337726	20.000	ng		0.00
System Monitoring Compounds							
5) 2-Fluorophenol	5.516	112	1150166	102.577	ng		0.00
7) Phenol-d6	6.528	99	1387418	102.468	ng		0.00
23) Nitrobenzene-d5	7.475	82	1213856	115.369	ng		0.00
42) 2,4,6-Tribromophenol	10.739	330	369823	112.641	ng		0.00
45) 2-Fluorobiphenyl	9.269	172	2425708	94.966	ng		0.00
79) Terphenyl-d14	13.021	244	2247982	101.236	ng		0.00
Target Compounds							Qvalue
2) 1,4-Dioxane	2.693	88	259869	51.894	ng		99
3) Pyridine	3.463	79	672711	51.688	ng		100
4) n-Nitrosodimethylamine	3.416	42	353028	52.416	ng		96
6) Aniline	6.575	93	986118	51.454	ng		99
8) 2-Chlorophenol	6.692	128	606252	52.325	ng		99
9) Benzaldehyde	6.463	77	482619	50.998	ng		98
10) Phenol	6.539	94	755326m	51.820	ng		
11) bis(2-Chloroethyl)ether	6.645	93	575347	50.934	ng		100
12) 1,3-Dichlorobenzene	6.851	146	661612	50.476	ng		99
13) 1,4-Dichlorobenzene	6.928	146	666083	50.439	ng		99
14) 1,2-Dichlorobenzene	7.081	146	640983	50.843	ng		100
15) Benzyl Alcohol	7.045	79	504297	52.514	ng		99
16) 2,2'-oxybis(1-Chloropr...	7.186	45	1049319	51.444	ng		100
17) 2-Methylphenol	7.151	107	485145	51.113	ng		100
18) Hexachloroethane	7.428	117	229911	53.193	ng		97
19) n-Nitroso-di-n-propyla...	7.328	70	420535	50.261	ng		99
20) 3+4-Methylphenols	7.310	107	611404	51.510	ng	#	89
22) Acetophenone	7.322	105	796460	49.805	ng		96
24) Nitrobenzene	7.492	77	584670	56.634	ng		99
25) Isophorone	7.734	82	1129933	51.170	ng		99
26) 2-Nitrophenol	7.810	139	228021	51.526	ng		97
27) 2,4-Dimethylphenol	7.839	122	576054	51.741	ng		100
28) bis(2-Chloroethoxy)met...	7.939	93	708439	50.870	ng		99
29) 2,4-Dichlorophenol	8.045	162	494488	53.125	ng		99
30) 1,2,4-Trichlorobenzene	8.133	180	524508	50.497	ng		100
31) Naphthalene	8.216	128	1725818	50.064	ng		100
32) Benzoic acid	7.939	122	252184	47.667	ng		98
33) 4-Chloroaniline	8.257	127	730860	50.914	ng		99
34) Hexachlorobutadiene	8.333	225	307118	50.679	ng		99
35) Caprolactam	8.628	113	154579	55.251	ng		97
36) 4-Chloro-3-methylphenol	8.733	107	505527	52.891	ng		100
37) 2-Methylnaphthalene	8.904	142	1057924	49.915	ng		100
38) 1-Methylnaphthalene	9.004	142	1107217	50.427	ng		98
40) 1,2,4,5-Tetrachloroben...	9.069	216	509426	50.261	ng		99
41) Hexachlorocyclopentadiene	9.063	237	331287	57.591	ng		99
43) 2,4,6-Trichlorophenol	9.180	196	337910	53.801	ng		100

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.216	196	370489	56.106	ng	98
46) 1,1'-Biphenyl	9.369	154	1392513	49.736	ng	100
47) 2-Chloronaphthalene	9.398	162	1041842	50.015	ng	100
48) 2-Nitroaniline	9.486	65	284370	51.394	ng	100
49) Acenaphthylene	9.810	152	1769441	50.339	ng	100
50) Dimethylphthalate	9.669	163	1178638	51.156	ng	99
51) 2,6-Dinitrotoluene	9.727	165	232906	52.309	ng	99
52) Acenaphthene	9.986	154	1024409	49.912	ng	100
53) 3-Nitroaniline	9.898	138	279608	51.745	ng	97
54) 2,4-Dinitrophenol	9.998	184	62086	50.725	ng	98
55) Dibenzofuran	10.157	168	1542097	50.123	ng	100
56) 4-Nitrophenol	10.039	139	210153	58.029	ng	99
57) 2,4-Dinitrotoluene	10.127	165	280193	51.918	ng	98
58) Fluorene	10.498	166	1146269	49.000	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.269	232	300782	54.989	ng	99
60) Diethylphthalate	10.369	149	1170793	51.459	ng	99
61) 4-Chlorophenyl-phenyle...	10.492	204	553271	49.802	ng	98
62) 4-Nitroaniline	10.510	138	258117	52.547	ng	100
63) Azobenzene	10.651	77	1150899	51.226	ng	99
65) 4,6-Dinitro-2-methylph...	10.539	198	98114	49.012	ng	95
66) n-Nitrosodiphenylamine	10.610	169	1041277	50.442	ng	100
67) 4-Bromophenyl-phenylether	10.980	248	340298	50.576	ng	97
68) Hexachlorobenzene	11.045	284	384807	51.064	ng	99
69) Atrazine	11.133	200	287405	52.819	ng	100
70) Pentachlorophenol	11.233	266	219414	55.114	ng	98
71) Phenanthrene	11.463	178	1607091	49.244	ng	99
72) Anthracene	11.510	178	1659788	49.744	ng	100
73) Carbazole	11.663	167	1504715	49.436	ng	100
74) Di-n-butylphthalate	11.998	149	1578769	51.573	ng	100
75) Fluoranthene	12.651	202	1563998	48.481	ng	99
77) Benzidine	12.768	184	762858	55.590	ng	99
78) Pyrene	12.880	202	1563568	52.484	ng	100
80) Butylbenzylphthalate	13.498	149	423529	51.685	ng	100
81) Benzo(a)anthracene	14.068	228	1120699	51.691	ng	100
82) 3,3'-Dichlorobenzidine	14.027	252	363124	57.021	ng	98
83) Chrysene	14.104	228	990991	50.214	ng	100
84) Bis(2-ethylhexyl)phtha...	14.057	149	585800	51.638	ng	100
85) Di-n-octyl phthalate	14.680	149	1065626	51.078	ng	99
87) Indeno(1,2,3-cd)pyrene	17.080	276	1313336	53.091	ng	99
88) Benzo(b)fluoranthene	15.127	252	1073096	51.153	ng	100
89) Benzo(k)fluoranthene	15.162	252	967248	50.408	ng	98
90) Benzo(a)pyrene	15.504	252	996599	51.890	ng	100
91) Dibenzo(a,h)anthracene	17.103	278	1073246	53.457	ng	99
92) Benzo(g,h,i)perylene	17.539	276	1069402	52.984	ng	98

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

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