Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF031325\
Data File : BF141939.D
Acq On : 13 Mar 2025 12:17
Operator : RC/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SSTDCCC040

Quant Time: Mar 13 13:05:23 2025
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF031025.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Mar 10 15:46:22 2025
Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.875	152	198004	20.000	ng	0.00
21) Naphthalene-d8	8.157	136	767232	20.000	ng	0.00
39) Acenaphthene-d10	9.910	164	426931	20.000	ng	0.00
64) Phenanthrene-d10	11.398	188	726402	20.000	ng	0.00
76) Chrysene-d12	14.033	240	451214	20.000	ng	0.00
86) Perylene-d12	15.504	264	394917	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.493	112	914193	77.057	ng	0.00
7) Phenol-d6	6.504	99	1145375	75.826	ng	0.00
23) Nitrobenzene-d5	7.439	82	1055101	77.391	ng	0.00
42) 2,4,6-Tribromophenol	10.698	330	432513	79.856	ng	0.00
45) 2-Fluorobiphenyl	9.233	172	2159613	76.930	ng	0.00
79) Terphenyl-d14	12.980	244	2456339	80.485	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.681	88	195530	39.334	ng	99
3) Pyridine	3.446	79	472100	38.717	ng	99
4) n-Nitrosodimethylamine	3.399	42	226886	39.034	ng	98
6) Aniline	6.540	93	558500	37.480	ng	100
8) 2-Chlorophenol	6.663	128	505200	38.395	ng	100
9) Benzaldehyde	6.428	77	297311	35.193	ng	99
10) Phenol	6.516	94	605095	38.077	ng	100
11) bis(2-Chloroethyl)ether	6.616	93	446408	37.411	ng	99
12) 1,3-Dichlorobenzene	6.816	146	542765	38.208	ng	99
13) 1,4-Dichlorobenzene	6.892	146	551564	38.375	ng	99
14) 1,2-Dichlorobenzene	7.051	146	522830	38.620	ng	99
15) Benzyl Alcohol	7.016	79	463386	37.946	ng	99
16) 2,2'-oxybis(1-Chloropr...	7.151	45	508939	35.329	ng	99
17) 2-Methylphenol	7.122	107	393438	37.607	ng	99
18) Hexachloroethane	7.392	117	205179	38.069	ng	95
19) n-Nitroso-di-n-propyla...	7.292	70	347590	35.812	ng	100
20) 3+4-Methylphenols	7.275	107	502037	37.464	ng	98
22) Acetophenone	7.287	105	704297	38.332	ng	99
24) Nitrobenzene	7.457	77	523612	38.634	ng	100
25) Isophorone	7.698	82	889235	36.899	ng	100
26) 2-Nitrophenol	7.775	139	263074	39.549	ng	100
27) 2,4-Dimethylphenol	7.804	122	352630	38.499	ng	100
28) bis(2-Chloroethoxy)met...	7.904	93	559932	37.044	ng	99
29) 2,4-Dichlorophenol	8.010	162	435680	38.322	ng	99
30) 1,2,4-Trichlorobenzene	8.098	180	482699	38.527	ng	99
31) Naphthalene	8.181	128	1512818	38.385	ng	100
32) Benzoic acid	7.922	122	338815	42.081	ng	99
33) 4-Chloroaniline	8.228	127	523526	37.629	ng	99
34) Hexachlorobutadiene	8.298	225	319154	39.060	ng	99
35) Caprolactam	8.598	113	131311	37.884	ng	94
36) 4-Chloro-3-methylphenol	8.698	107	479000	37.769	ng	100
37) 2-Methylnaphthalene	8.869	142	1002209	38.045	ng	100
38) 1-Methylnaphthalene	8.969	142	966389	38.016	ng	99
40) 1,2,4,5-Tetrachloroben...	9.034	216	506575	38.972	ng	98
41) Hexachlorocyclopentadiene	9.022	237	213556	41.981	ng	99
43) 2,4,6-Trichlorophenol	9.145	196	335717	39.520	ng	100

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.181	196	335620	39.009	ng	100
46) 1,1'-Biphenyl	9.333	154	1256074	38.721	ng	100
47) 2-Chloronaphthalene	9.357	162	946087	39.130	ng	99
48) 2-Nitroaniline	9.451	65	273545	39.074	ng	100
49) Acenaphthylene	9.775	152	1399495	39.005	ng	100
50) Dimethylphthalate	9.633	163	1139250	38.106	ng	100
51) 2,6-Dinitrotoluene	9.692	165	243366	39.096	ng	97
52) Acenaphthene	9.945	154	991587	39.326	ng	99
53) 3-Nitroaniline	9.863	138	243288	38.530	ng	99
54) 2,4-Dinitrophenol	9.963	184	132695	41.009	ng	88
55) Dibenzofuran	10.116	168	1389049	38.554	ng	99
56) 4-Nitrophenol	10.010	139	196664	41.301	ng	100
57) 2,4-Dinitrotoluene	10.098	165	318648	39.193	ng	98
58) Fluorene	10.463	166	1061221	38.195	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.233	232	310873	39.708	ng	98
60) Diethylphthalate	10.333	149	1134119	38.044	ng	100
61) 4-Chlorophenyl-phenyle...	10.451	204	559798	38.646	ng	98
62) 4-Nitroaniline	10.475	138	242993	39.528	ng	99
63) Azobenzene	10.610	77	1050098	37.888	ng	99
65) 4,6-Dinitro-2-methylph...	10.504	198	181277	41.862	ng	99
66) n-Nitrosodiphenylamine	10.569	169	906795	38.576	ng	100
67) 4-Bromophenyl-phenylether	10.939	248	355291	38.946	ng	100
68) Hexachlorobenzene	11.004	284	395519	39.530	ng	99
69) Atrazine	11.092	200	238990	33.164	ng	99
70) Pentachlorophenol	11.198	266	261878	42.480	ng	100
71) Phenanthrene	11.422	178	1501673	38.274	ng	99
72) Anthracene	11.475	178	1507154	38.288	ng	100
73) Carbazole	11.627	167	1302310	38.363	ng	100
74) Di-n-butylphthalate	11.957	149	1710894	37.983	ng	99
75) Fluoranthene	12.610	202	1559838	37.478	ng	99
77) Benzidine	12.727	184	397420	63.130	ng	99
78) Pyrene	12.839	202	1536542	39.368	ng	99
80) Butylbenzylphthalate	13.451	149	585688	38.339	ng	99
81) Benzo(a)anthracene	14.021	228	1129185	38.137	ng	100
82) 3,3'-Dichlorobenzidine	13.986	252	353908	41.361	ng	99
83) Chrysene	14.063	228	1052493	39.214	ng	100
84) Bis(2-ethylhexyl)phtha...	14.010	149	770893	36.599	ng	100
85) Di-n-octyl phthalate	14.621	149	1091714	37.251	ng	99
87) Indeno(1,2,3-cd)pyrene	16.986	276	1036676	40.580	ng	98
88) Benzo(b)fluoranthene	15.074	252	995633	36.786	ng	99
89) Benzo(k)fluoranthene	15.098	252	862161	37.223	ng	100
90) Benzo(a)pyrene	15.439	252	809840	38.240	ng	100
91) Dibenzo(a,h)anthracene	17.004	278	863026	40.945	ng	99
92) Benzo(g,h,i)perylene	17.439	276	863842	41.473	ng	98

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

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