

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010422\
 Data File : BP008654.D
 Acq On : 04 Jan 2022 10:31
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SSTDCCC040

Quant Time: Jan 04 13:13:13 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP122821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 30 08:44:33 2021
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By :Christian
 Giraldo
 01/05/2022
 Supervised By :Jagrut
 Upadhyay
 01/05/2022

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.887	152	503685	20.000	ng	-0.01
21) Naphthalene-d8	10.687	136	2013505	20.000	ng	-0.01
39) Acenaphthene-d10	14.522	164	1446002	20.000	ng	0.00
64) Phenanthrene-d10	17.269	188	3267783	20.000	ng	0.00
76) Chrysene-d12	21.351	240	3433142	20.000	ng	-0.01
86) Perylene-d12	23.774	264	3724535	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.463	112	2297978	83.275	ng	0.00
7) Phenol-d6	7.058	99	3094256	80.559	ng	0.00
23) Nitrobenzene-d5	9.046	82	2791117	82.463	ng	-0.01
42) 2,4,6-Tribromophenol	16.010	330	2051732	104.874	ng	0.00
45) 2-Fluorobiphenyl	13.151	172	8168266	81.842	ng	0.00
79) Terphenyl-d14	19.822	244	12726305	75.687	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.369	88	437711	39.264	ng	100
3) Pyridine	3.769	79	1241933	40.840	ng	100
4) n-Nitrosodimethylamine	3.675	42	408740	41.627	ng	99
6) Aniline	7.210	93	1761034	39.448	ng	99
8) 2-Chlorophenol	7.452	128	1331300	40.956	ng	99
9) Benzaldehyde	7.022	77	888732	40.654	ng	98
10) Phenol	7.087	94	1556837	40.149	ng	100
11) bis(2-Chloroethyl)ether	7.310	93	1216962	38.913	ng	99
12) 1,3-Dichlorobenzene	7.775	146	1510219	39.691	ng	100
13) 1,4-Dichlorobenzene	7.922	146	1543413	39.879	ng	99
14) 1,2-Dichlorobenzene	8.240	146	1501103	39.701	ng	98
15) Benzyl Alcohol	8.128	79	1127195	40.971	ng	99
16) 2,2'-oxybis(1-Chloropr...	8.422	45	1266462	35.588	ng	99
17) 2-Methylphenol	8.334	107	1091786	40.461	ng	99
18) Hexachloroethane	8.975	117	496504	39.349	ng	94
19) n-Nitroso-di-n-propyla...	8.699	70	923934	37.668	ng	100
20) 3+4-Methylphenols	8.663	107	1529457	40.346	ng	100
22) Acetophenone	8.710	105	1937698	40.132	ng	# 99
24) Nitrobenzene	9.087	77	1311616	40.689	ng	97
25) Isophorone	9.622	82	2518373	39.829	ng	98
26) 2-Nitrophenol	9.799	139	779104	50.887	ng	99
27) 2,4-Dimethylphenol	9.869	122	1131508	40.956	ng	98
28) bis(2-Chloroethoxy)met...	10.099	93	1692438	38.661	ng	100
29) 2,4-Dichlorophenol	10.340	162	1446257	43.253	ng	99
30) 1,2,4-Trichlorobenzene	10.551	180	1620658	41.103	ng	99
31) Naphthalene	10.740	128	4217986	40.360	ng	99
32) Benzoic acid	10.022	122	1012455	58.021	ng	100
33) 4-Chloroaniline	10.846	127	1828942	40.812	ng	99
34) Hexachlorobutadiene	11.034	225	1003232	41.912	ng	100
35) Caprolactam	11.646	113	484968	44.126	ng	98
36) 4-Chloro-3-methylphenol	11.975	107	1419766	42.478	ng	99
37) 2-Methylnaphthalene	12.345	142	3147618	40.013	ng	99
38) 1-Methylnaphthalene	12.569	142	3072631	39.950	ng	100
40) 1,2,4,5-Tetrachloroben...	12.716	216	1876509	41.192	ng	100
41) Hexachlorocyclopentadiene	12.704	237	964799	39.827	ng	100
43) 2,4,6-Trichlorophenol	12.957	196	1353901	45.680	ng	100

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.028	196	1463505	44.720	ng	99
46) 1,1'-Biphenyl	13.357	154	4261110	39.841	ng	99
47) 2-Chloronaphthalene	13.398	162	3405382	40.627	ng	99
48) 2-Nitroaniline	13.592	65	817210	44.474	ng	97
49) Acenaphthylene	14.240	152	5313617	41.270	ng	99
50) Dimethylphthalate	13.981	163	4476706	40.196	ng	100
51) 2,6-Dinitrotoluene	14.092	165	983267	43.025	ng	96
52) Acenaphthene	14.587	154	3468107m	41.169	ng	
53) 3-Nitroaniline	14.422	138	1034142	43.477	ng	98
54) 2,4-Dinitrophenol	14.622	184	538699	52.923	ng	99
55) Dibenzofuran	14.922	168	5445124	40.319	ng	99
56) 4-Nitrophenol	14.734	139	882620	47.122	ng	96
57) 2,4-Dinitrotoluene	14.881	165	1363800	44.725	ng	98
58) Fluorene	15.569	166	4262730	40.505	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.145	232	1361919m	46.452	ng	
60) Diethylphthalate	15.345	149	4354809	40.002	ng	100
61) 4-Chlorophenyl-phenyle...	15.563	204	2365011	40.342	ng	100
62) 4-Nitroaniline	15.586	138	1096692	45.399	ng	99
63) Azobenzene	15.857	77	3388167	38.561	ng	98
65) 4,6-Dinitro-2-methylph...	15.645	198	848222	51.620	ng	97
66) n-Nitrosodiphenylamine	15.775	169	3890687	39.451	ng	100
67) 4-Bromophenyl-phenylether	16.457	248	1612054	45.084	ng	99
68) Hexachlorobenzene	16.581	284	1865588	42.504	ng	95
69) Atrazine	16.733	200	1551019	42.044	ng	98
70) Pentachlorophenol	16.922	266	1223553	51.181	ng	99
71) Phenanthrene	17.310	178	7099662	40.298	ng	100
72) Anthracene	17.404	178	7015750	40.869	ng	100
73) Carbazole	17.675	167	6928015	41.119	ng	100
74) Di-n-butylphthalate	18.228	149	7668155	41.599	ng	100
75) Fluoranthene	19.275	202	8710457	42.983	ng	98
77) Benzidine	19.451	184	4151643	40.251	ng	99
78) Pyrene	19.627	202	8814517	39.466	ng	99
80) Butylbenzylphthalate	20.504	149	3717586	41.731	ng	95
81) Benzo(a)anthracene	21.339	228	8975892	40.301	ng	99
82) 3,3'-Dichlorobenzidine	21.269	252	3660496	44.699	ng	99
83) Chrysene	21.392	228	8475134	40.695	ng	99
84) Bis(2-ethylhexyl)phtha...	21.269	149	5101102	37.907	ng	100
85) Di-n-octyl phthalate	22.198	149	9120647	43.786	ng	99
87) Indeno(1,2,3-cd)pyrene	26.298	276	10353979	41.386	ng	98
88) Benzo(b)fluoranthene	23.033	252	9623608	40.178	ng	99
89) Benzo(k)fluoranthene	23.086	252	8888488	37.329	ng	99
90) Benzo(a)pyrene	23.668	252	7896457	40.894	ng	99
91) Dibenzo(a,h)anthracene	26.321	278	8581014	40.658	ng	99
92) Benzo(g,h,i)perylene	27.074	276	8110129	41.095	ng	99

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

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