

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF051024\
 Data File : BF137881.D
 Acq On : 10 May 2024 15:17
 Operator : CG\JU
 Sample : PB160847BS
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB160847BS

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 05/13/2024
 Supervised By :mohammad ahmed 05/13/2024

Quant Time: May 11 00:35:17 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF043024.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 01 03:25:08 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.051	152	164207	20.000	ng	0.00
21) Naphthalene-d8	8.333	136	662141	20.000	ng	-0.01
39) Acenaphthene-d10	10.098	164	382223	20.000	ng	-0.01
64) Phenanthrene-d10	11.592	188	708935	20.000	ng	0.00
76) Chrysene-d12	14.245	240	444978	20.000	ng	0.00
86) Perylene-d12	15.809	264	360487	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.686	112	1206166	122.983	ng	0.01
7) Phenol-d6	6.669	99	1494682	120.152	ng	0.00
23) Nitrobenzene-d5	7.616	82	919480	80.678	ng	0.00
42) 2,4,6-Tribromophenol	10.892	330	523351	134.103	ng	0.00
45) 2-Fluorobiphenyl	9.410	172	1955341	82.995	ng	-0.01
79) Terphenyl-d14	13.180	244	2440208	91.902	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.081	88	154232	34.922	ng	97
3) Pyridine	3.810	79	386199	36.131	ng	95
4) n-Nitrosodimethylamine	3.781	42	187336	38.921	ng	94
6) Aniline	6.710	93	391212m	31.729	ng	
8) 2-Chlorophenol	6.833	128	473933	44.342	ng	99
9) Benzaldehyde	6.604	77	198392	29.614	ng	98
10) Phenol	6.681	94	531053	41.662	ng	98
11) bis(2-Chloroethyl)ether	6.781	93	414830	41.562	ng	99
12) 1,3-Dichlorobenzene	6.992	146	503825	41.802	ng	98
13) 1,4-Dichlorobenzene	7.069	146	511469	42.292	ng	98
14) 1,2-Dichlorobenzene	7.222	146	487109	42.764	ng	98
15) Benzyl Alcohol	7.186	79	375183	43.612	ng	99
16) 2,2'-oxybis(1-Chloropr...	7.310	45	535243	38.152	ng	98
17) 2-Methylphenol	7.292	107	380112	44.074	ng	100
18) Hexachloroethane	7.563	117	176396	43.145	ng	97
19) n-Nitroso-di-n-propyla...	7.457	70	303688	40.850	ng	96
20) 3+4-Methylphenols	7.439	107	482997	42.568	ng	95
22) Acetophenone	7.457	105	633741	42.556	ng	99
24) Nitrobenzene	7.633	77	463276	39.440	ng	95
25) Isophorone	7.869	82	844915	41.118	ng	98
26) 2-Nitrophenol	7.945	139	255303	47.996	ng	93
27) 2,4-Dimethylphenol	7.975	122	345001	45.052	ng	97
28) bis(2-Chloroethoxy)met...	8.075	93	511983	41.151	ng	99
29) 2,4-Dichlorophenol	8.186	162	405382	43.797	ng	98
30) 1,2,4-Trichlorobenzene	8.275	180	426261	42.244	ng	99
31) Naphthalene	8.357	128	1391359	42.283	ng	100
32) Benzoic acid	8.086	122	284636	43.085	ng	93
33) 4-Chloroaniline	8.398	127	188031	15.410	ng	98
34) Hexachlorobutadiene	8.469	225	259651	41.913	ng	99
35) Caprolactam	8.769	113	134126m	45.735	ng	
36) 4-Chloro-3-methylphenol	8.875	107	426796	44.216	ng	95
37) 2-Methylnaphthalene	9.051	142	908479	42.652	ng	99
38) 1-Methylnaphthalene	9.151	142	853456	40.727	ng	100
40) 1,2,4,5-Tetrachloroben...	9.216	216	432711	42.744	ng	99
41) Hexachlorocyclopentadiene	9.198	237	470204	114.353	ng	99
43) 2,4,6-Trichlorophenol	9.322	196	295537	42.887	ng	100

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.363	196	325274	42.748	ng	97
46) 1,1'-Biphenyl	9.516	154	1122298	42.785	ng	98
47) 2-Chloronaphthalene	9.545	162	873080	41.226	ng	99
48) 2-Nitroaniline	9.633	65	252368	43.939	ng	99
49) Acenaphthylene	9.963	152	1400623	45.719	ng	100
50) Dimethylphthalate	9.810	163	1041130	43.148	ng	100
51) 2,6-Dinitrotoluene	9.874	165	234692	42.560	ng	99
52) Acenaphthene	10.133	154	937510	46.168	ng	99
53) 3-Nitroaniline	10.045	138	161874	28.032	ng	96
54) 2,4-Dinitrophenol	10.151	184	256901	94.815	ng	89
55) Dibenzofuran	10.304	168	1265525	42.826	ng	98
56) 4-Nitrophenol	10.204	139	397017	82.953	ng	93
57) 2,4-Dinitrotoluene	10.280	165	330241	45.841	ng	99
58) Fluorene	10.651	166	1012457	43.050	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.416	232	274367	44.414	ng	99
60) Diethylphthalate	10.516	149	1000297	43.649	ng	98
61) 4-Chlorophenyl-phenyle...	10.639	204	490338	42.315	ng	99
62) 4-Nitroaniline	10.669	138	241829	41.287	ng	95
63) Azobenzene	10.798	77	929130	41.949	ng	97
65) 4,6-Dinitro-2-methylph...	10.692	198	182663	44.891	ng	80
66) n-Nitrosodiphenylamine	10.757	169	886256	41.763	ng	100
67) 4-Bromophenyl-phenylether	11.133	248	322966	42.567	ng	95
68) Hexachlorobenzene	11.198	284	354730	42.783	ng	96
69) Atrazine	11.280	200	309458	47.826	ng	98
70) Pentachlorophenol	11.392	266	445550	78.272	ng	99
71) Phenanthrene	11.621	178	1495439	43.277	ng	99
72) Anthracene	11.674	178	1527840	44.438	ng	100
73) Carbazole	11.821	167	1402166	42.317	ng	100
74) Di-n-butylphthalate	12.139	149	1609118	44.518	ng	99
75) Fluoranthene	12.815	202	1565026	41.433	ng	98
77) Benzidine	12.927	184	339917	33.029	ng	99
78) Pyrene	13.045	202	1577530	44.713	ng	100
80) Butylbenzylphthalate	13.651	149	582028	51.994	ng	96
81) Benzo(a)anthracene	14.233	228	1259404	44.251	ng	99
82) 3,3'-Dichlorobenzidine	14.192	252	266341	31.260	ng	98
83) Chrysene	14.274	228	1170176	43.912	ng	100
84) Bis(2-ethylhexyl)phtha...	14.204	149	740624	52.497	ng	99
85) Di-n-octyl phthalate	14.833	149	983815	51.817	ng	98
87) Indeno(1,2,3-cd)pyrene	17.450	276	1030999	50.305	ng	99
88) Benzo(b)fluoranthene	15.339	252	971544	43.317	ng	99
89) Benzo(k)fluoranthene	15.374	252	944942	43.590	ng	99
90) Benzo(a)pyrene	15.745	252	871220	47.613	ng	99
91) Dibenzo(a,h)anthracene	17.474	278	831740	49.858	ng	100
92) Benzo(g,h,i)perylene	17.956	276	801400	45.825	ng	99

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

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