

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032222\
 Data File : BP009483.D
 Acq On : 22 Mar 2022 12:20
 Operator : CG/JU
 Sample : PB143421BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB143421BS

Quant Time: Mar 23 01:20:28 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP031722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 17 18:26:53 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.793	152	409221	20.000	ng	0.00
21) Naphthalene-d8	10.587	136	1540071	20.000	ng	0.00
39) Acenaphthene-d10	14.440	164	862084	20.000	ng	0.00
64) Phenanthrene-d10	17.198	188	1601892	20.000	ng	0.00
76) Chrysene-d12	21.310	240	1367524	20.000	ng	0.00
86) Perylene-d12	23.716	264	1324954	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.399	112	3237090	138.110	ng	0.00
7) Phenol-d6	6.987	99	3917256	123.583	ng	0.00
23) Nitrobenzene-d5	8.952	82	2423584	83.626	ng	0.00
42) 2,4,6-Tribromophenol	15.934	330	1352233	95.054	ng	-0.01
45) 2-Fluorobiphenyl	13.057	172	5216966	83.892	ng	0.00
79) Terphenyl-d14	19.775	244	6553739	86.019	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.334	88	369499	39.801	ng	99
3) Pyridine	3.722	79	976047	39.036	ng	99
4) n-Nitrosodimethylamine	3.634	42	444938	48.319	ng	98
6) Aniline	7.128	93	1390796	38.498	ng	98
8) 2-Chlorophenol	7.369	128	1167853	43.696	ng	98
9) Benzaldehyde	6.940	77	618864	40.976	ng	96
10) Phenol	7.011	94	1429656	44.958	ng	99
11) bis(2-Chloroethyl)ether	7.222	93	1108774	44.057	ng	97
12) 1,3-Dichlorobenzene	7.687	146	1286989	42.664	ng	98
13) 1,4-Dichlorobenzene	7.828	146	1306813	42.371	ng	99
14) 1,2-Dichlorobenzene	8.146	146	1252222	42.006	ng	99
15) Benzyl Alcohol	8.040	79	947136	37.478	ng	99
16) 2,2'-oxybis(1-Chloropr...	8.328	45	1256466	46.057	ng	97
17) 2-Methylphenol	8.252	107	917873	41.886	ng	99
18) Hexachloroethane	8.869	117	457342	42.312	ng	97
19) n-Nitroso-di-n-propyla...	8.605	70	803318	40.068	ng	99
20) 3+4-Methylphenols	8.581	107	1216861	40.385	ng	99
22) Acetophenone	8.616	105	1609830	42.233	ng	99
24) Nitrobenzene	8.993	77	1217254	44.462	ng	98
25) Isophorone	9.522	82	2164754	43.790	ng	99
26) 2-Nitrophenol	9.704	139	585968	41.062	ng	97
27) 2,4-Dimethylphenol	9.781	122	992616	49.295	ng	99
28) bis(2-Chloroethoxy)met...	10.004	93	1348615	41.899	ng	98
29) 2,4-Dichlorophenol	10.252	162	1003849	40.870	ng	99
30) 1,2,4-Trichlorobenzene	10.451	180	1147099	40.310	ng	99
31) Naphthalene	10.640	128	3323358	41.893	ng	100
32) Benzoic acid	9.957	122	553189	35.138	ng	95
33) 4-Chloroaniline	10.757	127	522120	16.129	ng	98
34) Hexachlorobutadiene	10.928	225	729737	37.846	ng	99
35) Caprolactam	11.546	113	279461	33.713	ng	99
36) 4-Chloro-3-methylphenol	11.898	107	981270	37.179	ng	99
37) 2-Methylnaphthalene	12.251	142	2232292	40.106	ng	99
38) 1-Methylnaphthalene	12.475	142	2133932	39.355	ng	99
40) 1,2,4,5-Tetrachloroben...	12.628	216	1234861	43.172	ng	99
41) Hexachlorocyclopentadiene	12.610	237	1309331	102.868	ng	100
43) 2,4,6-Trichlorophenol	12.875	196	764461	40.322	ng	98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.951	196	817157	39.691	ng	98
46) 1,1'-Biphenyl	13.269	154	2815822	43.784	ng	99
47) 2-Chloronaphthalene	13.310	162	2184522	43.144	ng	99
48) 2-Nitroaniline	13.516	65	575097	41.703	ng	99
49) Acenaphthylene	14.157	152	3524378	45.090	ng	100
50) Dimethylphthalate	13.898	163	2497494	38.233	ng	100
51) 2,6-Dinitrotoluene	14.016	165	556539	40.663	ng	99
52) Acenaphthene	14.498	154	1967332	42.134	ng	100
53) 3-Nitroaniline	14.345	138	295823	20.463	ng	99
54) 2,4-Dinitrophenol	14.563	184	562746	66.953	ng	98
55) Dibenzofuran	14.839	168	3143953	40.080	ng	100
56) 4-Nitrophenol	14.681	139	830065	77.031	ng	97
57) 2,4-Dinitrotoluene	14.810	165	726096	38.583	ng	99
58) Fluorene	15.492	166	2428276	39.361	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.075	232	680851	36.479	ng	95
60) Diethylphthalate	15.269	149	2376059	36.343	ng	100
61) 4-Chlorophenyl-phenyle...	15.487	204	1276212	37.553	ng	97
62) 4-Nitroaniline	15.516	138	507072	33.400	ng	97
63) Azobenzene	15.781	77	2309387	40.607	ng	98
65) 4,6-Dinitro-2-methylph...	15.581	198	376769	36.577	ng	98
66) n-Nitrosodiphenylamine	15.704	169	2105219	44.410	ng	98
67) 4-Bromophenyl-phenylether	16.386	248	786146	40.106	ng	96
68) Hexachlorobenzene	16.504	284	882886	39.136	ng	99
69) Atrazine	16.669	200	715938	42.956	ng	98
70) Pentachlorophenol	16.857	266	978742	81.160	ng	99
71) Phenanthrene	17.239	178	3630507	42.338	ng	100
72) Anthracene	17.334	178	3658122	43.208	ng	100
73) Carbazole	17.610	167	3236821	39.093	ng	99
74) Di-n-butylphthalate	18.169	149	3714041	38.279	ng	99
75) Fluoranthene	19.222	202	3980916	38.347	ng	99
77) Benzidine	19.404	184	1434299	42.790	ng	100
78) Pyrene	19.575	202	4081264	45.290	ng	100
80) Butylbenzylphthalate	20.457	149	1519627	39.675	ng	99
81) Benzo(a)anthracene	21.298	228	3776446	42.033	ng	100
82) 3,3'-Dichlorobenzidine	21.233	252	1007877	29.038	ng	99
83) Chrysene	21.345	228	3516020	41.328	ng	99
84) Bis(2-ethylhexyl)phtha...	21.233	149	2161822	40.493	ng	100
85) Di-n-octyl phthalate	22.157	149	3567139	38.764	ng	99
87) Indeno(1,2,3-cd)pyrene	26.209	276	4050225	40.292	ng	97
88) Benzo(b)fluoranthene	22.980	252	3635508	45.963	ng	98
89) Benzo(k)fluoranthene	23.027	252	3597440	46.453	ng	98
90) Benzo(a)pyrene	23.610	252	3491211	51.589	ng	98
91) Dibenzo(a,h)anthracene	26.233	278	3578486	42.714	ng	98
92) Benzo(g,h,i)perylene	26.980	276	3533937	42.879	ng	97

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

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