

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032523\
 Data File : BP014300.D
 Acq On : 24 Mar 2023 21:49
 Operator : CG/JU
 Sample : PB151516BS
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB151516BS

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 03/27/2023
 Supervised By : Jagrut Upadhyay 03/27/2023

Quant Time: Mar 25 03:22:41 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP030623.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Mar 25 03:19:01 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.028	152	186669	20.000	ng	0.00
21) Naphthalene-d8	10.851	136	676742	20.000	ng	# 0.00
39) Acenaphthene-d10	14.675	164	381644	20.000	ng	0.00
64) Phenanthrene-d10	17.433	188	736976	20.000	ng	# 0.00
76) Chrysene-d12	21.515	240	610059	20.000	ng	0.00
86) Perylene-d12	24.033	264	620464	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.581	112	1699241	147.627	ng	0.00
7) Phenol-d6	7.199	99	2158376	142.823	ng	0.00
23) Nitrobenzene-d5	9.210	82	1462671	98.786	ng	0.00
42) 2,4,6-Tribromophenol	16.175	330	806542	130.344	ng	0.00
45) 2-Fluorobiphenyl	13.298	172	2871318	97.474	ng	0.00
79) Terphenyl-d14	19.974	244	3740273	101.097	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.440	88	200887	40.174	ng	94
3) Pyridine	3.852	79	530969	38.236	ng	94
4) n-Nitrosodimethylamine	3.763	42	286980	46.557	ng	96
6) Aniline	7.357	93	772995	38.939	ng	97
8) 2-Chlorophenol	7.593	128	586516	47.576	ng	95
9) Benzaldehyde	7.169	77	456640m	60.382	ng	
10) Phenol	7.222	94	765230	48.653	ng	99
11) bis(2-Chloroethyl)ether	7.451	93	600531	48.222	ng	96
12) 1,3-Dichlorobenzene	7.916	146	664971	49.086	ng	99
13) 1,4-Dichlorobenzene	8.069	146	677994	49.470	ng	100
14) 1,2-Dichlorobenzene	8.387	146	643953	49.056	ng	99
15) Benzyl Alcohol	8.281	79	583827m	47.384	ng	
16) 2,2'-oxybis(1-Chloropr...	8.563	45	787742m	58.102	ng	
17) 2-Methylphenol	8.481	107	501226	44.683	ng	99
18) Hexachloroethane	9.116	117	261809	50.851	ng	94
19) n-Nitroso-di-n-propyla...	8.845	70	484939	48.752	ng	97
20) 3+4-Methylphenols	8.810	107	663272	43.905	ng	99
22) Acetophenone	8.869	105	930721	49.031	ng	# 98
24) Nitrobenzene	9.251	77	735352	48.538	ng	97
25) Isophorone	9.781	82	1245965	45.669	ng	98
26) 2-Nitrophenol	9.963	139	311585	47.283	ng	96
27) 2,4-Dimethylphenol	10.022	122	513433	47.918	ng	96
28) bis(2-Chloroethoxy)met...	10.257	93	764047	48.386	ng	99
29) 2,4-Dichlorophenol	10.498	162	531147	45.285	ng	98
30) 1,2,4-Trichlorobenzene	10.716	180	617426	46.367	ng	98
31) Naphthalene	10.904	128	1723222	46.341	ng	99
32) Benzoic acid	10.187	122	328200	37.583	ng	92
33) 4-Chloroaniline	11.028	127	192802	12.095	ng	98
34) Hexachlorobutadiene	11.181	225	427303	44.137	ng	97
35) Caprolactam	11.822	113	142171m	42.346	ng	
36) 4-Chloro-3-methylphenol	12.139	107	583160	44.580	ng	96
37) 2-Methylnaphthalene	12.504	142	1162770	42.384	ng	99
38) 1-Methylnaphthalene	12.728	142	1081500	42.291	ng	99
40) 1,2,4,5-Tetrachloroben...	12.875	216	710948	48.872	ng	# 99
41) Hexachlorocyclopentadiene	12.845	237	1018793	111.594	ng	100
43) 2,4,6-Trichlorophenol	13.116	196	424221	46.135	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.186	196	450651	42.294	ng	99
46) 1,1'-Biphenyl	13.510	154	1533754	49.928	ng	98
47) 2-Chloronaphthalene	13.557	162	1171293	49.530	ng	99
48) 2-Nitroaniline	13.763	65	369683	51.511	ng	92
49) Acenaphthylene	14.398	152	1782947	47.410	ng	100
50) Dimethylphthalate	14.133	163	1399839	44.460	ng	99
51) 2,6-Dinitrotoluene	14.257	165	302584	45.636	ng	94
52) Acenaphthene	14.739	154	1058808	46.770	ng	99
53) 3-Nitroaniline	14.592	138	175725	27.043	ng	92
54) 2,4-Dinitrophenol	14.798	184	383768	78.137	ng	93
55) Dibenzofuran	15.075	168	1709439	46.318	ng	99
56) 4-Nitrophenol	14.898	139	468069	88.460	ng	92
57) 2,4-Dinitrotoluene	15.045	165	403352	45.510	ng	# 94
58) Fluorene	15.727	166	1332054	45.454	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.304	232	390537	42.013	ng	95
60) Diethylphthalate	15.492	149	1374270	44.208	ng	99
61) 4-Chlorophenyl-phenyle...	15.716	204	712307	42.877	ng	96
62) 4-Nitroaniline	15.757	138	281506	43.508	ng	91
63) Azobenzene	16.010	77	1454881	50.578	ng	96
65) 4,6-Dinitro-2-methylph...	15.810	198	242199	44.369	ng	91
66) n-Nitrosodiphenylamine	15.933	169	1138543	48.890	ng	99
67) 4-Bromophenyl-phenylether	16.616	248	439251	45.059	ng	95
68) Hexachlorobenzene	16.733	284	502633	48.103	ng	99
69) Atrazine	16.886	200	408384	49.948	ng	98
70) Pentachlorophenol	17.080	266	608528	80.751	ng	99
71) Phenanthrene	17.480	178	1985419	48.224	ng	99
72) Anthracene	17.569	178	1999451	47.529	ng	100
73) Carbazole	17.839	167	1733032	47.758	ng	100
74) Di-n-butylphthalate	18.369	149	2164715	47.712	ng	99
75) Fluoranthene	19.433	202	2164369	43.982	ng	97
77) Benzidine	19.616	184	1069385	93.441	ng	100
78) Pyrene	19.786	202	2247065	47.989	ng	100
80) Butylbenzylphthalate	20.645	149	894068	50.365	ng	97
81) Benzo(a)anthracene	21.498	228	2095644	46.817	ng	99
82) 3,3'-Dichlorobenzidine	21.427	252	516131	36.036	ng	99
83) Chrysene	21.557	228	1951817	46.044	ng	99
84) Bis(2-ethylhexyl)phtha...	21.398	149	1292466	49.629	ng	100
85) Di-n-octyl phthalate	22.357	149	2122533	50.209	ng	97
87) Indeno(1,2,3-cd)pyrene	26.680	276	2333622	48.676	ng	98
88) Benzo(b)fluoranthene	23.262	252	2051511	50.495	ng	99
89) Benzo(k)fluoranthene	23.309	252	1997015	49.576	ng	98
90) Benzo(a)pyrene	23.921	252	1764279	45.527	ng	99
91) Dibenzo(a,h)anthracene	26.698	278	1946682	48.843	ng	98
92) Benzo(g,h,i)perylene	27.497	276	1915535	48.497	ng	98

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

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