

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052822\
 Data File : BP010590.D
 Acq On : 28 May 2022 02:55
 Operator : CG/JU
 Sample : PB145173BS
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB145173BS

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 05/28/2022
 Supervised By : Jagrut Upadhyay 05/31/2022

Quant Time: May 28 04:17:09 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP052822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat May 28 02:18:04 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.869	152	105939	20.000	ng	0.00
21) Naphthalene-d8	10.669	136	435429	20.000	ng	0.00
39) Acenaphthene-d10	14.504	164	290051	20.000	ng	0.00
64) Phenanthrene-d10	17.257	188	599777	20.000	ng	0.00
76) Chrysene-d12	21.339	240	534039	20.000	ng	0.00
86) Perylene-d12	23.751	264	492919	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.446	112	803898	138.434	ng	0.00
7) Phenol-d6	7.040	99	1020608	123.799	ng	0.00
23) Nitrobenzene-d5	9.028	82	750053	81.440	ng	0.00
42) 2,4,6-Tribromophenol	15.998	330	429886	130.933	ng	0.00
45) 2-Fluorobiphenyl	13.128	172	1598338	78.208	ng	0.00
79) Terphenyl-d14	19.816	244	2491962	87.647	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.352	88	76429	30.992	ng	96
3) Pyridine	3.758	79	196495	28.452	ng	94
4) n-Nitrosodimethylamine	3.664	42	184733	43.869	ng	94
6) Aniline	7.193	93	381454	39.328	ng	98
8) 2-Chlorophenol	7.434	128	281480	42.947	ng	97
9) Benzaldehyde	7.005	77	172014m	31.707	ng	
10) Phenol	7.064	94	360761	42.317	ng	94
11) bis(2-Chloroethyl)ether	7.293	93	268721	39.771	ng	98
12) 1,3-Dichlorobenzene	7.758	146	307206	40.444	ng	98
13) 1,4-Dichlorobenzene	7.905	146	314311	40.383	ng	100
14) 1,2-Dichlorobenzene	8.222	146	301302	40.439	ng	98
15) Benzyl Alcohol	8.111	79	284977m	43.121	ng	
16) 2,2'-oxybis(1-Chloropr...	8.399	45	502951	39.814	ng	99
17) 2-Methylphenol	8.316	107	239655	41.887	ng	97
18) Hexachloroethane	8.946	117	123985	41.182	ng	99
19) n-Nitroso-di-n-propyla...	8.681	70	234060	38.926	ng	95
20) 3+4-Methylphenols	8.646	107	325564	41.395	ng	97
22) Acetophenone	8.693	105	447700	39.426	ng	# 98
24) Nitrobenzene	9.069	77	369254	39.685	ng	98
25) Isophorone	9.599	82	623704	40.538	ng	99
26) 2-Nitrophenol	9.781	139	149057	41.105	ng	97
27) 2,4-Dimethylphenol	9.846	122	255671	47.436	ng	98
28) bis(2-Chloroethoxy)met...	10.081	93	366061	43.069	ng	98
29) 2,4-Dichlorophenol	10.316	162	273083	42.145	ng	97
30) 1,2,4-Trichlorobenzene	10.534	180	304235	39.916	ng	100
31) Naphthalene	10.716	128	890548	38.823	ng	98
32) Benzoic acid	9.993	122	156225	38.435	ng	98
33) 4-Chloroaniline	10.828	127	261783	29.542	ng	97
34) Hexachlorobutadiene	11.004	225	199419	38.753	ng	99
35) Caprolactam	11.622	113	83604	44.700	ng	96
36) 4-Chloro-3-methylphenol	11.957	107	318777	43.180	ng	99
37) 2-Methylnaphthalene	12.328	142	626919	39.168	ng	97
38) 1-Methylnaphthalene	12.546	142	604764	38.690	ng	98
40) 1,2,4,5-Tetrachloroben...	12.698	216	344762	38.209	ng	98
41) Hexachlorocyclopentadiene	12.681	237	420376	87.519	ng	99
43) 2,4,6-Trichlorophenol	12.940	196	226944	40.289	ng	98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.010	196	251086	39.409	ng	98
46) 1,1'-Biphenyl	13.340	154	835410	38.615	ng	100
47) 2-Chloronaphthalene	13.381	162	650605	38.260	ng	99
48) 2-Nitroaniline	13.581	65	240513	40.808	ng	96
49) Acenaphthylene	14.222	152	1075818	41.188	ng	100
50) Dimethylphthalate	13.963	163	815775	39.072	ng	100
51) 2,6-Dinitrotoluene	14.081	165	183943	41.428	ng	97
52) Acenaphthene	14.569	154	614184	38.300	ng	97
53) 3-Nitroaniline	14.410	138	145896	33.551	ng	97
54) 2,4-Dinitrophenol	14.622	184	209478	76.747	ng	98
55) Dibenzofuran	14.904	168	1007015	39.711	ng	98
56) 4-Nitrophenol	14.722	139	308218	84.057	ng	97
57) 2,4-Dinitrotoluene	14.869	165	251885	43.583	ng	99
58) Fluorene	15.551	166	815406	39.332	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.134	232	220246	40.326	ng	98
60) Diethylphthalate	15.328	149	849889	40.892	ng	99
61) 4-Chlorophenyl-phenyle...	15.545	204	420749	40.379	ng	99
62) 4-Nitroaniline	15.575	138	177909	38.865	ng	95
63) Azobenzene	15.840	77	890861	40.585	ng	98
65) 4,6-Dinitro-2-methylph...	15.640	198	136265	36.948	ng	99
66) n-Nitrosodiphenylamine	15.763	169	714484	39.562	ng	99
67) 4-Bromophenyl-phenylether	16.445	248	253776	38.894	ng	99
68) Hexachlorobenzene	16.563	284	279449	38.774	ng	99
69) Atrazine	16.722	200	257159	43.901	ng	98
70) Pentachlorophenol	16.910	266	348483	84.233	ng	99
71) Phenanthrene	17.298	178	1293058	39.313	ng	100
72) Anthracene	17.392	178	1272125	40.391	ng	99
73) Carbazole	17.663	167	1161428	43.389	ng	99
74) Di-n-butylphthalate	18.210	149	1397577	41.069	ng	100
75) Fluoranthene	19.263	202	1443197	40.397	ng	98
77) Benzidine	19.445	184	554164	54.020	ng	98
78) Pyrene	19.616	202	1494733	40.163	ng	100
80) Butylbenzylphthalate	20.486	149	598187	40.810	ng	99
81) Benzo(a)anthracene	21.327	228	1410763	40.263	ng	99
82) 3,3'-Dichlorobenzidine	21.257	252	460646	47.936	ng	99
83) Chrysene	21.380	228	1330571	38.715	ng	100
84) Bis(2-ethylhexyl)phtha...	21.251	149	850278	41.153	ng	100
85) Di-n-octyl phthalate	22.174	149	1392736	40.449	ng	100
87) Indeno(1,2,3-cd)pyrene	26.268	276	1565302	43.561	ng	100
88) Benzo(b)fluoranthene	23.016	252	1417431	45.672	ng	99
89) Benzo(k)fluoranthene	23.069	252	1326256	42.157	ng	98
90) Benzo(a)pyrene	23.645	252	1322813	51.529	ng	99
91) Dibenzo(a,h)anthracene	26.286	278	1388856	46.211	ng	98
92) Benzo(g,h,i)perylene	27.039	276	1420719	47.549	ng	99

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

