

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF091622\
 Data File : BF130285.D
 Acq On : 16 Sep 2022 11:14
 Operator : CG\JU
 Sample : PB147653BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB147653BS

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 09/19/2022
 Supervised By : Jagrut Upadhyay 09/19/2022

Quant Time: Sep 17 02:29:25 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 14 14:55:51 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.675	152	104233	20.000	ng	0.00
21) Naphthalene-d8	7.963	136	422197	20.000	ng	# 0.00
39) Acenaphthene-d10	9.710	164	219944	20.000	ng	0.00
64) Phenanthrene-d10	11.192	188	369324	20.000	ng	0.00
76) Chrysene-d12	13.821	240	205732	20.000	ng	0.00
86) Perylene-d12	15.210	264	172959	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.293	112	805884	134.101	ng	0.02
7) Phenol-d6	6.322	99	989991	131.660	ng	0.00
23) Nitrobenzene-d5	7.245	82	643802	77.904	ng	0.00
42) 2,4,6-Tribromophenol	10.504	330	324775	154.106	ng	0.00
45) 2-Fluorobiphenyl	9.039	172	1231233	81.517	ng	0.00
79) Terphenyl-d14	12.774	244	1227624	98.845	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.393	88	92434	33.491	ng	97
3) Pyridine	3.081	79	244410	33.948	ng	97
4) n-Nitrosodimethylamine	3.040	42	151478	38.684	ng	94
6) Aniline	6.340	93	394584	42.590	ng	96
8) 2-Chlorophenol	6.463	128	314684	45.969	ng	93
9) Benzaldehyde	6.228	77	198731	39.456	ng	95
10) Phenol	6.334	94	385172	43.711	ng	87
11) bis(2-Chloroethyl)ether	6.422	93	278355	43.795	ng	96
12) 1,3-Dichlorobenzene	6.616	146	330662	43.898	ng	99
13) 1,4-Dichlorobenzene	6.693	146	339587	44.209	ng	99
14) 1,2-Dichlorobenzene	6.845	146	319815	44.570	ng	98
15) Benzyl Alcohol	6.828	79	244766	41.056	ng	94
16) 2,2'-oxybis(1-Chloropr...	6.957	45	378016	40.778	ng	97
17) 2-Methylphenol	6.940	107	241021	43.311	ng	98
18) Hexachloroethane	7.187	117	130107	42.970	ng	100
19) n-Nitroso-di-n-propyla...	7.104	70	214901	42.355	ng	94
20) 3+4-Methylphenols	7.092	107	305058	42.199	ng	# 86
22) Acetophenone	7.098	105	441196	42.743	ng	# 95
24) Nitrobenzene	7.263	77	334300	39.839	ng	99
25) Isophorone	7.504	82	580385	43.574	ng	99
26) 2-Nitrophenol	7.581	139	163006	47.387	ng	91
27) 2,4-Dimethylphenol	7.622	122	269400	50.864	ng	98
28) bis(2-Chloroethoxy)met...	7.722	93	347269	46.302	ng	99
29) 2,4-Dichlorophenol	7.822	162	252603	43.522	ng	97
30) 1,2,4-Trichlorobenzene	7.904	180	263632	42.012	ng	94
31) Naphthalene	7.981	128	896474	41.215	ng	100
32) Benzoic acid	7.757	122	185494	45.288	ng	94
33) 4-Chloroaniline	8.034	127	293917	35.529	ng	99
34) Hexachlorobutadiene	8.098	225	170305	42.466	ng	99
35) Caprolactam	8.416	113	79402m	47.720	ng	
36) 4-Chloro-3-methylphenol	8.522	107	284780	44.266	ng	97
37) 2-Methylnaphthalene	8.675	142	584915	42.177	ng	99
38) 1-Methylnaphthalene	8.769	142	555333	41.684	ng	100
40) 1,2,4,5-Tetrachloroben...	8.839	216	271843	45.289	ng	99
41) Hexachlorocyclopentadiene	8.828	237	367520	114.363	ng	98
43) 2,4,6-Trichlorophenol	8.951	196	181909	43.424	ng	96

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	8.992	196	193861	42.885	ng	99
46) 1,1'-Biphenyl	9.139	154	716688	43.630	ng	99
47) 2-Chloronaphthalene	9.163	162	559103	42.076	ng	98
48) 2-Nitroaniline	9.263	65	178737	40.329	ng	90
49) Acenaphthylene	9.575	152	848073	42.567	ng	99
50) Dimethylphthalate	9.445	163	658119	43.318	ng	100
51) 2,6-Dinitrotoluene	9.504	165	143136	45.048	ng	93
52) Acenaphthene	9.745	154	625294m	47.768	ng	
53) 3-Nitroaniline	9.675	138	127474	36.004	ng	92
54) 2,4-Dinitrophenol	9.781	184	155820	89.589	ng	# 87
55) Dibenzofuran	9.922	168	775279	42.580	ng	98
56) 4-Nitrophenol	9.839	139	232133	90.018	ng	92
57) 2,4-Dinitrotoluene	9.904	165	190807	46.988	ng	98
58) Fluorene	10.257	166	636655	42.756	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.033	232	152722	43.163	ng	97
60) Diethylphthalate	10.145	149	649697	42.644	ng	99
61) 4-Chlorophenyl-phenylether	10.257	204	292113	44.719	ng	95
62) 4-Nitroaniline	10.286	138	150242	42.242	ng	98
63) Azobenzene	10.416	77	621339	39.606	ng	97
65) 4,6-Dinitro-2-methylph...	10.310	198	91448	44.481	ng	98
66) n-Nitrosodiphenylamine	10.375	169	536865	43.500	ng	100
67) 4-Bromophenyl-phenylether	10.745	248	187139	45.932	ng	92
68) Hexachlorobenzene	10.804	284	210317	45.539	ng	94
69) Atrazine	10.904	200	174495	48.381	ng	97
70) Pentachlorophenol	10.998	266	237145	95.675	ng	99
71) Phenanthrene	11.216	178	875303	42.213	ng	100
72) Anthracene	11.269	178	887270	44.341	ng	99
73) Carbazole	11.428	167	795238	44.036	ng	99
74) Di-n-butylphthalate	11.763	149	952367	43.733	ng	99
75) Fluoranthene	12.398	202	852244	42.535	ng	99
77) Benzidine	12.527	184	423778	70.958	ng	99
78) Pyrene	12.627	202	848151	47.126	ng	99
80) Butylbenzylphthalate	13.257	149	313307	46.991	ng	91
81) Benzo(a)anthracene	13.810	228	593633	43.374	ng	99
82) 3,3'-Dichlorobenzidine	13.780	252	178083	47.800	ng	98
83) Chrysene	13.845	228	554936	42.703	ng	99
84) Bis(2-ethylhexyl)phtha...	13.810	149	358425	46.933	ng	100
85) Di-n-octyl phthalate	14.421	149	521879	44.678	ng	97
87) Indeno(1,2,3-cd)pyrene	16.545	276	610770	49.797	ng	99
88) Benzo(b)fluoranthene	14.821	252	498646	44.171	ng	98
89) Benzo(k)fluoranthene	14.851	252	489092	44.043	ng	98
90) Benzo(a)pyrene	15.157	252	435017	48.640	ng	97
91) Dibenzo(a,h)anthracene	16.568	278	526469	52.323	ng	99
92) Benzo(g,h,i)perylene	16.951	276	519756	51.279	ng	98

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

