

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF041822\
 Data File : BF127821.D
 Acq On : 18 Apr 2022 11:28
 Operator : CG\JU
 Sample : PB144160BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB144160BS

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 04/19/2022
 Supervised By : Jagrut Upadhyay 04/19/2022

Quant Time: Apr 19 05:14:44 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF040622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 13 16:40:15 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.951	152	176788	20.000	ng	-0.01
21) Naphthalene-d8	8.240	136	638262	20.000	ng	# 0.00
39) Acenaphthene-d10	9.998	164	357915	20.000	ng	0.00
64) Phenanthrene-d10	11.492	188	645376	20.000	ng	0.00
76) Chrysene-d12	14.139	240	433483	20.000	ng	0.00
86) Perylene-d12	15.657	264	353025	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.593	112	1259173	115.734	ng	0.01
7) Phenol-d6	6.587	99	1611705	113.235	ng	0.00
23) Nitrobenzene-d5	7.522	82	1168855	81.642	ng	0.00
42) 2,4,6-Tribromophenol	10.792	330	489159	128.430	ng	0.00
45) 2-Fluorobiphenyl	9.316	172	1956713	77.339	ng	0.00
79) Terphenyl-d14	13.080	244	2187501	83.295	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.963	88	184511	35.754	ng	83
3) Pyridine	3.693	79	508144	37.386	ng	# 81
4) n-Nitrosodimethylamine	3.669	42	292610	40.031	ng	92
6) Aniline	6.616	93	593654	34.971	ng	96
8) 2-Chlorophenol	6.740	128	528336	47.871	ng	94
9) Benzaldehyde	6.504	77	281836	39.214	ng	95
10) Phenol	6.598	94	632819m	44.464	ng	
11) bis(2-Chloroethyl)ether	6.687	93	539836	45.930	ng	93
12) 1,3-Dichlorobenzene	6.893	146	570850	45.267	ng	98
13) 1,4-Dichlorobenzene	6.969	146	579786	45.481	ng	99
14) 1,2-Dichlorobenzene	7.122	146	551070	45.224	ng	99
15) Benzyl Alcohol	7.098	79	488858m	39.287	ng	
16) 2,2'-oxybis(1-Chloropr...	7.222	45	773537	41.694	ng	95
17) 2-Methylphenol	7.204	107	428530	44.637	ng	98
18) Hexachloroethane	7.463	117	222288	45.697	ng	99
19) n-Nitroso-di-n-propyla...	7.369	70	393473	42.742	ng	93
20) 3+4-Methylphenols	7.357	107	504726	41.256	ng	# 79
22) Acetophenone	7.363	105	713725	41.614	ng	# 87
24) Nitrobenzene	7.540	77	622813	44.877	ng	100
25) Isophorone	7.775	82	1144377	45.808	ng	98
26) 2-Nitrophenol	7.851	139	270183	54.172	ng	94
27) 2,4-Dimethylphenol	7.887	122	476393	48.786	ng	94
28) bis(2-Chloroethoxy)met...	7.981	93	659218	42.795	ng	99
29) 2,4-Dichlorophenol	8.092	162	462212	43.940	ng	97
30) 1,2,4-Trichlorobenzene	8.175	180	537340	44.385	ng	99
31) Naphthalene	8.263	128	1457348	43.574	ng	99
32) Benzoic acid	8.016	122	315712	44.563	ng	93
33) 4-Chloroaniline	8.304	127	213697	16.224	ng	96
34) Hexachlorobutadiene	8.369	225	357662	42.948	ng	99
35) Caprolactam	8.687	113	133435m	42.919	ng	
36) 4-Chloro-3-methylphenol	8.787	107	486126	44.079	ng	97
37) 2-Methylnaphthalene	8.951	142	972596	43.480	ng	99
38) 1-Methylnaphthalene	9.051	142	930627	43.103	ng	98
40) 1,2,4,5-Tetrachloroben...	9.116	216	519960	44.292	ng	98
41) Hexachlorocyclopentadiene	9.098	237	702587	99.355	ng	98
43) 2,4,6-Trichlorophenol	9.228	196	345723	43.378	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.269	196	374720	46.594	ng	98
46) 1,1'-Biphenyl	9.416	154	1188876	43.433	ng	100
47) 2-Chloronaphthalene	9.445	162	943833	43.389	ng	100
48) 2-Nitroaniline	9.539	65	325129	48.125	ng	93
49) Acenaphthylene	9.863	152	1510672	45.658	ng	99
50) Dimethylphthalate	9.722	163	1133165	43.678	ng	100
51) 2,6-Dinitrotoluene	9.781	165	252591	51.652	ng	94
52) Acenaphthene	10.034	154	1013892m	48.639	ng	
53) 3-Nitroaniline	9.951	138	160773	31.120	ng	96
54) 2,4-Dinitrophenol	10.057	184	276356	111.761	ng	93
55) Dibenzofuran	10.204	168	1291622	42.461	ng	97
56) 4-Nitrophenol	10.116	139	384354	94.719	ng	88
57) 2,4-Dinitrotoluene	10.186	165	335660	55.363	ng	# 91
58) Fluorene	10.551	166	975808	43.416	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.322	232	314809	44.923	ng	100
60) Diethylphthalate	10.416	149	1065898	42.934	ng	98
61) 4-Chlorophenyl-phenyle...	10.539	204	523554	42.395	ng	99
62) 4-Nitroaniline	10.575	138	244198	50.189	ng	95
63) Azobenzene	10.698	77	1107613	39.927	ng	97
65) 4,6-Dinitro-2-methylph...	10.598	198	176704	56.687	ng	# 81
66) n-Nitrosodiphenylamine	10.663	169	887874	43.682	ng	99
67) 4-Bromophenyl-phenylether	11.033	248	348785	44.530	ng	97
68) Hexachlorobenzene	11.092	284	392943	45.893	ng	94
69) Atrazine	11.186	200	342204	50.673	ng	97
70) Pentachlorophenol	11.292	266	436183	85.296	ng	97
71) Phenanthrene	11.516	178	1474713	43.128	ng	99
72) Anthracene	11.569	178	1495083	44.202	ng	99
73) Carbazole	11.722	167	1348974	42.563	ng	99
74) Di-n-butylphthalate	12.045	149	1618898	42.729	ng	100
75) Fluoranthene	12.710	202	1648932	44.191	ng	99
77) Benzidine	12.827	184	527966	46.664	ng	99
78) Pyrene	12.939	202	1639040	44.777	ng	99
80) Butylbenzylphthalate	13.551	149	623752	45.766	ng	88
81) Benzo(a)anthracene	14.127	228	1352478	46.715	ng	98
82) 3,3'-Dichlorobenzidine	14.086	252	290872	32.226	ng	97
83) Chrysene	14.169	228	1223564	44.027	ng	99
84) Bis(2-ethylhexyl)phtha...	14.104	149	836240	43.529	ng	# 98
85) Di-n-octyl phthalate	14.721	149	1272223	45.291	ng	# 93
87) Indeno(1,2,3-cd)pyrene	17.215	276	1064578	48.265	ng	99
88) Benzo(b)fluoranthene	15.204	252	1152666	50.756	ng	99
89) Benzo(k)fluoranthene	15.239	252	1012714	44.962	ng	99
90) Benzo(a)pyrene	15.592	252	1025927	55.580	ng	98
91) Dibenzo(a,h)anthracene	17.239	278	920782	51.524	ng	98
92) Benzo(g,h,i)perylene	17.692	276	939013	51.253	ng	98

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

