

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061722\
 Data File : BF128865.D
 Acq On : 17 Jun 2022 20:41
 Operator : CG\JU
 Sample : N3383-02MS 2X
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RB-03-(0-3.5)MS

Quant Time: Jun 18 05:07:50 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF060222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 14 14:53:53 2022
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By : Christian Giraldo 06/20/2022
 Supervised By : Jagrut Upadhyay 06/20/2022

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.781	152	96012	20.000	ng	0.00
21) Naphthalene-d8	8.069	136	320804	20.000	ng	# 0.00
39) Acenaphthene-d10	9.822	164	128411	20.000	ng	0.00
64) Phenanthrene-d10	11.316	188	239956	20.000	ng	# 0.00
76) Chrysene-d12	13.962	240	160540	20.000	ng	0.00
86) Perylene-d12	15.421	264	120886	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.422	112	218676	36.278	ng	0.00
7) Phenol-d6	6.439	99	266039	36.157	ng	0.00
23) Nitrobenzene-d5	7.345	82	182328	26.461	ng	0.00
42) 2,4,6-Tribromophenol	10.616	330	45350	29.890	ng	0.00
45) 2-Fluorobiphenyl	9.139	172	304841	33.254	ng	0.00
79) Terphenyl-d14	12.898	244	285364	30.887	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.499	88	40116	17.927	ng	# 95
3) Pyridine	3.234	79	112202	18.875	ng	95
4) n-Nitrosodimethylamine	3.181	42	75808	18.389	ng	# 80
6) Aniline	6.445	93	141322	17.809	ng	# 66
8) 2-Chlorophenol	6.581	128	126524	20.513	ng	99
9) Benzaldehyde	6.334	77	94923	21.176	ng	97
10) Phenol	6.451	94	166765	21.441	ng	# 67
11) bis(2-Chloroethyl)ether	6.516	93	120518	21.146	ng	99
12) 1,3-Dichlorobenzene	6.722	146	147086	20.628	ng	98
13) 1,4-Dichlorobenzene	6.798	146	149120	20.707	ng	97
14) 1,2-Dichlorobenzene	6.951	146	138748	20.597	ng	98
15) Benzyl Alcohol	6.928	79	118324	19.220	ng	91
16) 2,2'-oxybis(1-Chloropr...	7.063	45	194400	20.702	ng	# 45
17) 2-Methylphenol	7.057	107	98871	20.282	ng	# 83
18) Hexachloroethane	7.292	117	34381	12.832	ng	98
19) n-Nitroso-di-n-propyla...	7.192	70	95713	20.922	ng	97
20) 3+4-Methylphenols	7.210	107	131155	20.153	ng	# 79
22) Acetophenone	7.192	105	188462	22.461	ng	# 96
24) Nitrobenzene	7.369	77	138187	21.845	ng	96
25) Isophorone	7.604	82	234490	22.407	ng	95
26) 2-Nitrophenol	7.686	139	56192	20.161	ng	# 95
27) 2,4-Dimethylphenol	7.733	122	102709	24.192	ng	94
28) bis(2-Chloroethoxy)met...	7.816	93	139549	21.094	ng	98
29) 2,4-Dichlorophenol	7.939	162	95832	19.503	ng	99
30) 1,2,4-Trichlorobenzene	8.010	180	106550	19.258	ng	98
31) Naphthalene	8.092	128	525471	31.917	ng	98
32) Benzoic acid	7.851	122	19214m	6.379	ng	
33) 4-Chloroaniline	8.139	127	97539	15.714	ng	97
34) Hexachlorobutadiene	8.204	225	75491	19.239	ng	98
35) Caprolactam	8.504	113	27079	19.929	ng	95
36) 4-Chloro-3-methylphenol	8.645	107	96279	17.954	ng	97
37) 2-Methylnaphthalene	8.780	142	326641	31.192	ng	99
38) 1-Methylnaphthalene	8.880	142	287611	28.427	ng	99
40) 1,2,4,5-Tetrachloroben...	8.945	216	103792	26.026	ng	98
43) 2,4,6-Trichlorophenol	9.069	196	60033	22.260	ng	100
44) 2,4,5-Trichlorophenol	9.122	196	62074	22.272	ng	# 95

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46) 1,1'-Biphenyl	9.245	154	228642	23.731	ng	97
47) 2-Chloronaphthalene	9.269	162	159317	21.670	ng	99
48) 2-Nitroaniline	9.369	65	57131	22.725	ng	92
49) Acenaphthylene	9.686	152	299734	26.631	ng	100
50) Dimethylphthalate	9.545	163	197098	22.424	ng	100
51) 2,6-Dinitrotoluene	9.610	165	37285	21.493	ng	# 75
52) Acenaphthene	9.857	154	169244	23.091	ng	98
53) 3-Nitroaniline	9.780	138	40980	21.766	ng	99
55) Dibenzofuran	10.027	168	237867	22.581	ng	92
56) 4-Nitrophenol	9.980	139	40376	29.584	ng	# 74
57) 2,4-Dinitrotoluene	10.022	165	48737	21.051	ng	# 92
58) Fluorene	10.374	166	251018	30.727	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.157	232	36229	15.738	ng	# 96
60) Diethylphthalate	10.245	149	182825	20.696	ng	98
61) 4-Chlorophenyl-phenyle...	10.363	204	85215	20.111	ng	93
62) 4-Nitroaniline	10.392	138	44353	23.266	ng	# 84
63) Azobenzene	10.522	77	163546	18.806	ng	98
66) n-Nitrosodiphenylamine	10.480	169	158346	21.469	ng	99
67) 4-Bromophenyl-phenylether	10.857	248	54131	20.229	ng	96
68) Hexachlorobenzene	10.922	284	59391	19.985	ng	97
69) Atrazine	11.016	200	53961	22.751	ng	97
70) Pentachlorophenol	11.133	266	25381	17.109	ng	95
71) Phenanthrene	11.339	178	750248	62.335	ng	99
72) Anthracene	11.392	178	366607	30.774	ng	99
73) Carbazole	11.551	167	249927	22.588	ng	98
74) Di-n-butylphthalate	11.874	149	286883	21.295	ng	99
75) Fluoranthene	12.533	202	881778	70.089	ng	98
77) Benzidine	12.657	184	186184	58.177	ng	96
78) Pyrene	12.768	202	914885	76.597	ng	98
80) Butylbenzylphthalate	13.380	149	122089	24.340	ng	85
81) Benzo(a)anthracene	13.957	228	462555	46.223	ng	97
82) 3,3'-Dichlorobenzidine	13.915	252	74456	34.778	ng	96
83) Chrysene	13.992	228	416664	43.673	ng	99
84) Bis(2-ethylhexyl)phtha...	13.939	149	157486	24.151	ng	# 97
85) Di-n-octyl phthalate	14.551	149	220660	22.467	ng	# 99
87) Indeno(1,2,3-cd)pyrene	16.880	276	215581	29.588	ng	# 90
88) Benzo(b)fluoranthene	15.004	252	361806	46.839	ng	97
89) Benzo(k)fluoranthene	15.027	252	229035m	31.520	ng	
90) Benzo(a)pyrene	15.362	252	281899	46.433	ng	97
91) Dibenzo(a,h)anthracene	16.886	278	143441	23.734	ng	# 93
92) Benzo(g,h,i)perylene	17.321	276	184973	30.881	ng	# 92

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

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