

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111122\
 Data File : BF131157.D
 Acq On : 11 Nov 2022 17:01
 Operator : CG\JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 11/12/2022
 Supervised By : Jagrut Upadhyay 11/12/2022

Quant Time: Nov 11 18:00:17 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF110922.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 11 18:00:01 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.875	152	64026	20.000	ng	0.00
21) Naphthalene-d8	8.157	136	231656	20.000	ng	0.00
39) Acenaphthene-d10	9.922	164	148801	20.000	ng	0.00
64) Phenanthrene-d10	11.410	188	242292	20.000	ng	0.00
76) Chrysene-d12	14.057	240	186742	20.000	ng	0.00
86) Perylene-d12	15.545	264	125056	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.487	112	341458	90.696	ng	0.00
7) Phenol-d6	6.504	99	485376	98.612	ng	0.00
23) Nitrobenzene-d5	7.445	82	474113	92.341	ng	0.00
42) 2,4,6-Tribromophenol	10.710	330	107725	70.051	ng	0.00
45) 2-Fluorobiphenyl	9.239	172	824721	77.371	ng	0.00
79) Terphenyl-d14	12.998	244	794056	74.554	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.634	88	74614	40.263	ng	96
3) Pyridine	3.405	79	219180	43.074	ng	96
4) n-Nitrosodimethylamine	3.369	42	129186	44.011	ng	91
6) Aniline	6.540	93	302473	50.059	ng	98
8) 2-Chlorophenol	6.657	128	194192	44.836	ng	96
9) Benzaldehyde	6.428	77	142613	44.525	ng	95
10) Phenol	6.516	94	287266	49.926	ng	99
11) bis(2-Chloroethyl)ether	6.610	93	184854	44.905	ng	96
12) 1,3-Dichlorobenzene	6.816	146	191589	39.777	ng	96
13) 1,4-Dichlorobenzene	6.893	146	193812	39.613	ng	98
14) 1,2-Dichlorobenzene	7.045	146	217155	47.915	ng	97
15) Benzyl Alcohol	7.016	79	194646	51.705	ng	98
16) 2,2'-oxybis(1-Chloropr...	7.145	45	354092	51.300	ng	99
17) 2-Methylphenol	7.128	107	175981	48.580	ng	97
18) Hexachloroethane	7.387	117	89685	49.160	ng	94
19) n-Nitroso-di-n-propyla...	7.293	70	154649	47.853	ng	100
20) 3+4-Methylphenols	7.281	107	215377	47.587	ng	# 89
22) Acetophenone	7.287	105	277062	47.792	ng	# 87
24) Nitrobenzene	7.463	77	239442	48.143	ng	98
25) Isophorone	7.698	82	326880	41.175	ng	100
26) 2-Nitrophenol	7.775	139	88214	39.973	ng	# 86
27) 2,4-Dimethylphenol	7.810	122	139092m	41.099	ng	
28) bis(2-Chloroethoxy)met...	7.904	93	183478	40.782	ng	98
29) 2,4-Dichlorophenol	8.016	162	164416	45.442	ng	97
30) 1,2,4-Trichlorobenzene	8.098	180	158639	40.659	ng	99
31) Naphthalene	8.181	128	498583	39.390	ng	100
32) Benzoic acid	7.940	122	85234m	37.341	ng	
33) 4-Chloroaniline	8.234	127	196518	38.719	ng	98
34) Hexachlorobutadiene	8.292	225	108852	40.830	ng	99
35) Caprolactam	8.616	113	45596m	41.108	ng	
36) 4-Chloro-3-methylphenol	8.716	107	161873	40.429	ng	99
37) 2-Methylnaphthalene	8.875	142	318117	38.988	ng	100
38) 1-Methylnaphthalene	8.975	142	351185	43.949	ng	99
40) 1,2,4,5-Tetrachloroben...	9.039	216	162039	33.348	ng	99
41) Hexachlorocyclopentadiene	9.022	237	65344	34.466	ng	100
43) 2,4,6-Trichlorophenol	9.151	196	106938	33.535	ng	97

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44) 2,4,5-Trichlorophenol	9.192	196	123339	34.571	ng	97
46) 1,1'-Biphenyl	9.339	154	447936	38.934	ng	99
47) 2-Chloronaphthalene	9.363	162	354489	37.970	ng	98
48) 2-Nitroaniline	9.463	65	123395	36.315	ng	98
49) Acenaphthylene	9.781	152	567608	39.788	ng	99
50) Dimethylphthalate	9.639	163	373402	35.278	ng	99
51) 2,6-Dinitrotoluene	9.704	165	81784	35.158	ng	97
52) Acenaphthene	9.951	154	331912	39.004	ng	99
53) 3-Nitroaniline	9.881	138	105408	39.635	ng	97
54) 2,4-Dinitrophenol	9.986	184	46193	37.833	ng	# 90
55) Dibenzofuran	10.128	168	425105	33.733	ng	99
56) 4-Nitrophenol	10.039	139	64802	36.610	ng	94
57) 2,4-Dinitrotoluene	10.110	165	111513	36.063	ng	94
58) Fluorene	10.469	166	355572	34.739	ng	98
59) 2,3,4,6-Tetrachlorophenol	10.245	232	94150	34.319	ng	96
60) Diethylphthalate	10.339	149	381494	37.424	ng	99
61) 4-Chlorophenyl-phenyle...	10.457	204	171291	35.960	ng	93
62) 4-Nitroaniline	10.498	138	90065	33.512	ng	95
63) Azobenzene	10.622	77	393287	36.860	ng	97
65) 4,6-Dinitro-2-methylph...	10.522	198	64277	41.261	ng	81
66) n-Nitrosodiphenylamine	10.581	169	313101	38.438	ng	99
67) 4-Bromophenyl-phenylether	10.951	248	106014	38.695	ng	96
68) Hexachlorobenzene	11.016	284	116249	38.072	ng	98
69) Atrazine	11.110	200	101055	40.077	ng	98
70) Pentachlorophenol	11.216	266	57898	40.144	ng	99
71) Phenanthrene	11.433	178	531137	39.858	ng	100
72) Anthracene	11.486	178	510616	39.480	ng	98
73) Carbazole	11.645	167	534324	46.457	ng	99
74) Di-n-butylphthalate	11.969	149	590375	45.511	ng	99
75) Fluoranthene	12.627	202	567189	42.521	ng	99
77) Benzidine	12.745	184	203598	36.956	ng	99
78) Pyrene	12.857	202	578437	36.104	ng	99
80) Butylbenzylphthalate	13.469	149	227882	39.727	ng	91
81) Benzo(a)anthracene	14.045	228	524735	39.313	ng	99
82) 3,3'-Dichlorobenzidine	14.010	252	152471	40.022	ng	99
83) Chrysene	14.086	228	494047	39.193	ng	99
84) Bis(2-ethylhexyl)phtha...	14.027	149	303067	48.944	ng	98
85) Di-n-octyl phthalate	14.639	149	378263	48.172	ng	98
87) Indeno(1,2,3-cd)pyrene	17.057	276	335472	34.011	ng	99
88) Benzo(b)fluoranthene	15.104	252	298206	33.728	ng	100
89) Benzo(k)fluoranthene	15.139	252	279018	32.833	ng	98
90) Benzo(a)pyrene	15.480	252	235160	33.175	ng	98
91) Dibenzo(a,h)anthracene	17.074	278	275879	34.450	ng	99
92) Benzo(g,h,i)perylene	17.515	276	285850	33.912	ng	99

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

