

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF090523\
 Data File : BF135088.D
 Acq On : 05 Sep 2023 10:50
 Operator : CG\JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 09/06/2023
 Supervised By :mohammad ahmed 09/07/2023

Quant Time: Sep 05 11:19:04 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF083023.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 31 03:48:08 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.775	152	119294	20.000	ng	0.00
21) Naphthalene-d8	8.057	136	479258	20.000	ng	0.00
39) Acenaphthene-d10	9.816	164	256331	20.000	ng	0.00
64) Phenanthrene-d10	11.304	188	487998	20.000	ng	0.00
76) Chrysene-d12	13.963	240	310368	20.000	ng	0.00
86) Perylene-d12	15.427	264	257762	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.399	112	565742	73.252	ng	0.00
7) Phenol-d6	6.416	99	719169	74.707	ng	0.00
23) Nitrobenzene-d5	7.340	82	641293	75.175	ng	0.00
42) 2,4,6-Tribromophenol	10.610	330	214613	78.927	ng	0.00
45) 2-Fluorobiphenyl	9.134	172	1145333	72.747	ng	0.00
79) Terphenyl-d14	12.904	244	1280395	67.323	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.487	88	117466	35.906	ng	100
3) Pyridine	3.228	79	315701	35.795	ng	97
4) n-Nitrosodimethylamine	3.205	42	157271	36.463	ng	97
6) Aniline	6.446	93	441534	36.999	ng	99
8) 2-Chlorophenol	6.563	128	296497	37.789	ng	98
9) Benzaldehyde	6.328	77	192442	39.233	ng	98
10) Phenol	6.428	94	376419	37.247	ng	95
11) bis(2-Chloroethyl)ether	6.516	93	280374	36.842	ng	97
12) 1,3-Dichlorobenzene	6.716	146	303480	37.542	ng	98
13) 1,4-Dichlorobenzene	6.793	146	307122	37.960	ng	98
14) 1,2-Dichlorobenzene	6.945	146	284997	37.748	ng	99
15) Benzyl Alcohol	6.922	79	258295	39.164	ng	97
16) 2,2'-oxybis(1-Chloropr...	7.051	45	449505	35.272	ng	99
17) 2-Methylphenol	7.034	107	258288	37.777	ng	98
18) Hexachloroethane	7.281	117	112451	37.521	ng	96
19) n-Nitroso-di-n-propyla...	7.193	70	209633	36.661	ng	95
20) 3+4-Methylphenols	7.187	107	312430	38.733	ng	97
22) Acetophenone	7.187	105	390783	36.659	ng	# 99
24) Nitrobenzene	7.363	77	325672	37.217	ng	99
25) Isophorone	7.598	82	586954	36.290	ng	99
26) 2-Nitrophenol	7.675	139	163033	37.709	ng	98
27) 2,4-Dimethylphenol	7.716	122	253673	36.444	ng	97
28) bis(2-Chloroethoxy)met...	7.810	93	349294	35.801	ng	99
29) 2,4-Dichlorophenol	7.916	162	248493	37.563	ng	99
30) 1,2,4-Trichlorobenzene	7.998	180	265729	37.170	ng	100
31) Naphthalene	8.075	128	860636	36.761	ng	99
32) Benzoic acid	7.845	122	206634m	37.111	ng	
33) 4-Chloroaniline	8.128	127	373944	36.518	ng	96
34) Hexachlorobutadiene	8.192	225	149786	36.920	ng	97
35) Caprolactam	8.510	113	83043	38.339	ng	94
36) 4-Chloro-3-methylphenol	8.616	107	278749	38.795	ng	97
37) 2-Methylnaphthalene	8.769	142	584729	37.433	ng	99
38) 1-Methylnaphthalene	8.869	142	546863	37.415	ng	99
40) 1,2,4,5-Tetrachloroben...	8.934	216	260344	36.281	ng	99
41) Hexachlorocyclopentadiene	8.916	237	122355	40.062	ng	99
43) 2,4,6-Trichlorophenol	9.051	196	172390	35.772	ng	97

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.092	196	202353	36.269	ng	97
46) 1,1'-Biphenyl	9.234	154	705547	36.026	ng	99
47) 2-Chloronaphthalene	9.257	162	529205	35.934	ng	98
48) 2-Nitroaniline	9.363	65	192803	38.244	ng	99
49) Acenaphthylene	9.675	152	822187	36.931	ng	99
50) Dimethylphthalate	9.545	163	668044	37.220	ng	99
51) 2,6-Dinitrotoluene	9.604	165	150692	38.263	ng	98
52) Acenaphthene	9.851	154	515880	36.362	ng	99
53) 3-Nitroaniline	9.775	138	169157	36.872	ng	100
54) 2,4-Dinitrophenol	9.886	184	74692	39.682	ng	90
55) Dibenzofuran	10.022	168	749230	36.609	ng	99
56) 4-Nitrophenol	9.939	139	122304	37.925	ng	93
57) 2,4-Dinitrotoluene	10.016	165	191546	39.027	ng	96
58) Fluorene	10.369	166	588134	37.695	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.139	232	156485	36.424	ng	96
60) Diethylphthalate	10.245	149	631674	37.387	ng	100
61) 4-Chlorophenyl-phenyle...	10.357	204	284921	36.919	ng	100
62) 4-Nitroaniline	10.392	138	174210	38.809	ng	97
63) Azobenzene	10.522	77	624898	36.732	ng	99
65) 4,6-Dinitro-2-methylph...	10.428	198	112428	38.066	ng	90
66) n-Nitrosodiphenylamine	10.481	169	541199	35.667	ng	100
67) 4-Bromophenyl-phenylether	10.851	248	184956	35.734	ng	99
68) Hexachlorobenzene	10.916	284	193045	36.069	ng	99
69) Atrazine	11.010	200	141174	40.586	ng	97
70) Pentachlorophenol	11.110	266	119062	36.771	ng	100
71) Phenanthrene	11.333	178	909763	36.136	ng	100
72) Anthracene	11.386	178	939220	36.513	ng	99
73) Carbazole	11.545	167	832844	37.539	ng	99
74) Di-n-butylphthalate	11.875	149	992196	37.099	ng	99
75) Fluoranthene	12.527	202	973224	38.234	ng	99
77) Benzidine	12.651	184	248720	44.455	ng	99
78) Pyrene	12.757	202	975520	33.104	ng	99
80) Butylbenzylphthalate	13.386	149	395691	37.169	ng	92
81) Benzo(a)anthracene	13.957	228	737206	34.821	ng	99
82) 3,3'-Dichlorobenzidine	13.916	252	237599	37.192	ng	# 97
83) Chrysene	13.992	228	736402	35.393	ng	99
84) Bis(2-ethylhexyl)phtha...	13.945	149	444020	40.303	ng	100
85) Di-n-octyl phthalate	14.563	149	672660	40.001	ng	100
87) Indeno(1,2,3-cd)pyrene	16.915	276	625209	36.010	ng	98
88) Benzo(b)fluoranthene	15.004	252	584136	36.421	ng	98
89) Benzo(k)fluoranthene	15.033	252	573159	36.991	ng	99
90) Benzo(a)pyrene	15.368	252	532707	35.754	ng	99
91) Dibenzo(a,h)anthracene	16.933	278	506143	35.994	ng	99
92) Benzo(g,h,i)perylene	17.368	276	519053	35.490	ng	99

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

