

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF091823\
 Data File : BF135298.D
 Acq On : 18 Sep 2023 10:41
 Operator : CG\JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 09/19/2023
 Supervised By :mohammad ahmed 09/19/2023

Quant Time: Sep 18 11:53:24 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091123.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 12 00:19:07 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.763	152	95683	20.000	ng	0.00
21) Naphthalene-d8	8.039	136	380844	20.000	ng	0.00
39) Acenaphthene-d10	9.798	164	202362	20.000	ng	0.00
64) Phenanthrene-d10	11.292	188	397254	20.000	ng	0.00
76) Chrysene-d12	13.951	240	180433	20.000	ng	# 0.00
86) Perylene-d12	15.410	264	187310	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.387	112	461175	78.392	ng	0.00
7) Phenol-d6	6.410	99	574081	76.744	ng	0.00
23) Nitrobenzene-d5	7.328	82	543239	77.496	ng	0.00
42) 2,4,6-Tribromophenol	10.592	330	160046	79.586	ng	0.00
45) 2-Fluorobiphenyl	9.122	172	1013857	75.589	ng	0.00
79) Terphenyl-d14	12.886	244	1030726	88.555	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.463	88	105112	40.757	ng	98
3) Pyridine	3.210	79	289580	41.720	ng	96
4) n-Nitrosodimethylamine	3.175	42	162084	44.005	ng	92
6) Aniline	6.428	93	366373	37.349	ng	# 87
8) 2-Chlorophenol	6.551	128	244568	38.943	ng	95
9) Benzaldehyde	6.316	77	161316	37.854	ng	99
10) Phenol	6.422	94	310005	37.793	ng	84
11) bis(2-Chloroethyl)ether	6.504	93	233099	37.884	ng	100
12) 1,3-Dichlorobenzene	6.698	146	250304	39.218	ng	98
13) 1,4-Dichlorobenzene	6.781	146	249222	38.381	ng	99
14) 1,2-Dichlorobenzene	6.934	146	232880	38.619	ng	96
15) Benzyl Alcohol	6.910	79	213105	39.700	ng	99
16) 2,2'-oxybis(1-Chloropr...	7.040	45	431320	37.190	ng	97
17) 2-Methylphenol	7.022	107	208557	37.633	ng	96
18) Hexachloroethane	7.269	117	94679	39.462	ng	97
19) n-Nitroso-di-n-propyla...	7.181	70	175951	37.974	ng	99
20) 3+4-Methylphenols	7.175	107	252623	37.909	ng	90
22) Acetophenone	7.175	105	331743	38.100	ng	99
24) Nitrobenzene	7.345	77	269078	38.836	ng	99
25) Isophorone	7.587	82	488407	37.943	ng	99
26) 2-Nitrophenol	7.663	139	131888	39.661	ng	95
27) 2,4-Dimethylphenol	7.704	122	213298	38.160	ng	98
28) bis(2-Chloroethoxy)met...	7.798	93	294221	38.189	ng	99
29) 2,4-Dichlorophenol	7.904	162	200997	38.808	ng	98
30) 1,2,4-Trichlorobenzene	7.981	180	210282	38.844	ng	98
31) Naphthalene	8.063	128	714813	38.630	ng	99
32) Benzoic acid	7.834	122	160844m	38.215	ng	
33) 4-Chloroaniline	8.116	127	309295	38.097	ng	98
34) Hexachlorobutadiene	8.175	225	129933	39.805	ng	98
35) Caprolactam	8.498	113	68811m	39.586	ng	
36) 4-Chloro-3-methylphenol	8.604	107	226772	39.395	ng	100
37) 2-Methylnaphthalene	8.757	142	484489	38.951	ng	99
38) 1-Methylnaphthalene	8.851	142	445145	38.225	ng	100
40) 1,2,4,5-Tetrachloroben...	8.922	216	224476	38.289	ng	99
41) Hexachlorocyclopentadiene	8.904	237	96357	41.242	ng	100
43) 2,4,6-Trichlorophenol	9.039	196	147263	38.365	ng	99

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44) 2,4,5-Trichlorophenol	9.081	196	171764	38.857	ng	97
46) 1,1'-Biphenyl	9.222	154	590053	37.943	ng	99
47) 2-Chloronaphthalene	9.245	162	423329	37.518	ng	99
48) 2-Nitroaniline	9.345	65	169429	38.651	ng	96
49) Acenaphthylene	9.663	152	716510	38.210	ng	100
50) Dimethylphthalate	9.528	163	529157	38.676	ng	99
51) 2,6-Dinitrotoluene	9.592	165	114783	38.297	ng	93
52) Acenaphthene	9.833	154	423711	38.249	ng	98
53) 3-Nitroaniline	9.763	138	133796	38.029	ng	97
54) 2,4-Dinitrophenol	9.875	184	54349	35.695	ng	97
55) Dibenzofuran	10.004	168	639526	37.832	ng	97
56) 4-Nitrophenol	9.933	139	79044	35.350	ng	97
57) 2,4-Dinitrotoluene	9.998	165	143714	38.301	ng	# 86
58) Fluorene	10.351	166	501857	38.618	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.128	232	132381	38.931	ng	99
60) Diethylphthalate	10.228	149	533075	38.823	ng	99
61) 4-Chlorophenyl-phenyle...	10.345	204	244677	39.195	ng	93
62) 4-Nitroaniline	10.380	138	130587	38.614	ng	98
63) Azobenzene	10.504	77	514756	38.305	ng	98
65) 4,6-Dinitro-2-methylph...	10.410	198	82917	39.298	ng	99
66) n-Nitrosodiphenylamine	10.463	169	457816	38.067	ng	98
67) 4-Bromophenyl-phenylether	10.833	248	150907	38.195	ng	97
68) Hexachlorobenzene	10.898	284	157321	39.085	ng	93
69) Atrazine	10.998	200	119697	36.989	ng	98
70) Pentachlorophenol	11.098	266	85243	35.396	ng	100
71) Phenanthrene	11.316	178	721497	38.401	ng	99
72) Anthracene	11.369	178	737747	38.201	ng	99
73) Carbazole	11.527	167	659520	37.514	ng	100
74) Di-n-butylphthalate	11.863	149	807978	39.003	ng	100
75) Fluoranthene	12.510	202	744516	38.029	ng	99
77) Benzidine	12.639	184	127839	34.584	ng	99
78) Pyrene	12.739	202	735888	42.837	ng	99
80) Butylbenzylphthalate	13.369	149	263846	43.623	ng	98
81) Benzo(a)anthracene	13.939	228	450443	38.661	ng	100
82) 3,3'-Dichlorobenzidine	13.904	252	137655	37.825	ng	99
83) Chrysene	13.974	228	439191	37.914	ng	98
84) Bis(2-ethylhexyl)phtha...	13.927	149	287337	46.295	ng	99
85) Di-n-octyl phthalate	14.545	149	397915	40.342	ng	99
87) Indeno(1,2,3-cd)pyrene	16.898	276	491395	40.012	ng	98
88) Benzo(b)fluoranthene	14.986	252	350621	34.579	ng	99
89) Benzo(k)fluoranthene	15.015	252	364580	36.399	ng	99
90) Benzo(a)pyrene	15.351	252	360391	37.270	ng	98
91) Dibenzo(a,h)anthracene	16.909	278	403680	40.172	ng	98
92) Benzo(g,h,i)perylene	17.345	276	406518	38.736	ng	96

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

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