

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF060523\
 Data File : BF133564.D
 Acq On : 05 Jun 2023 12:53
 Operator : CG\JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 06/06/2023
 Supervised By :mohammad ahmed 06/08/2023

Quant Time: Jun 05 17:50:20 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF052523.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 05 17:49:53 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.834	152	133955	20.000	ng	0.00
21) Naphthalene-d8	8.116	136	489545	20.000	ng	0.00
39) Acenaphthene-d10	9.874	164	244121	20.000	ng	0.00
64) Phenanthrene-d10	11.357	188	471183	20.000	ng	0.00
76) Chrysene-d12	14.004	240	247744	20.000	ng	0.00
86) Perylene-d12	15.456	264	307678	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.434	112	638588	79.539	ng	0.00
7) Phenol-d6	6.457	99	762766	77.177	ng	0.00
23) Nitrobenzene-d5	7.398	82	707592	96.508	ng	0.00
42) 2,4,6-Tribromophenol	10.663	330	228767	94.010	ng	0.00
45) 2-Fluorobiphenyl	9.192	172	1298909	82.123	ng	0.00
79) Terphenyl-d14	12.945	244	1256072	90.145	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.534	88	142845	39.395	ng	100
3) Pyridine	3.275	79	396157	38.966	ng	97
4) n-Nitrosodimethylamine	3.228	42	204742	36.926	ng	95
6) Aniline	6.498	93	511102	38.087	ng	98
8) 2-Chlorophenol	6.616	128	321810	40.679	ng	95
9) Benzaldehyde	6.381	77	183711	36.748	ng	99
10) Phenol	6.469	94	392631	39.601	ng	98
11) bis(2-Chloroethyl)ether	6.569	93	314226	38.332	ng	98
12) 1,3-Dichlorobenzene	6.775	146	346874	39.495	ng	98
13) 1,4-Dichlorobenzene	6.851	146	348267	39.553	ng	98
14) 1,2-Dichlorobenzene	7.004	146	322315	39.668	ng	99
15) Benzyl Alcohol	6.975	79	294532	39.253	ng	97
16) 2,2'-oxybis(1-Chloropr...	7.110	45	567400	36.495	ng	98
17) 2-Methylphenol	7.081	107	289414	38.632	ng	99
18) Hexachloroethane	7.345	117	125055	41.073	ng	93
19) n-Nitroso-di-n-propyla...	7.251	70	238082	37.231	ng	92
20) 3+4-Methylphenols	7.234	107	355273	41.946	ng	# 88
22) Acetophenone	7.245	105	432965	38.405	ng	# 91
24) Nitrobenzene	7.416	77	363003	46.466	ng	100
25) Isophorone	7.657	82	659333	38.585	ng	98
26) 2-Nitrophenol	7.733	139	164432	46.137	ng	95
27) 2,4-Dimethylphenol	7.769	122	286605	39.420	ng	99
28) bis(2-Chloroethoxy)met...	7.869	93	385668	38.923	ng	99
29) 2,4-Dichlorophenol	7.969	162	271708	41.538	ng	98
30) 1,2,4-Trichlorobenzene	8.057	180	281373	39.768	ng	98
31) Naphthalene	8.139	128	944452	39.794	ng	99
32) Benzoic acid	7.863	122	175694m	38.418	ng	
33) 4-Chloroaniline	8.186	127	407959	40.105	ng	97
34) Hexachlorobutadiene	8.257	225	176311	40.966	ng	99
35) Caprolactam	8.557	113	94664	44.127	ng	99
36) 4-Chloro-3-methylphenol	8.663	107	302442	41.276	ng	97
37) 2-Methylnaphthalene	8.828	142	650290	40.130	ng	98
38) 1-Methylnaphthalene	8.928	142	613538	40.617	ng	97
40) 1,2,4,5-Tetrachloroben...	8.992	216	290496	40.149	ng	99
41) Hexachlorocyclopentadiene	8.980	237	133691	36.056	ng	98
43) 2,4,6-Trichlorophenol	9.104	196	195711	42.272	ng	100

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.139	196	227380	42.312	ng	97
46) 1,1'-Biphenyl	9.292	154	780324	39.343	ng	99
47) 2-Chloronaphthalene	9.316	162	569411	39.613	ng	98
48) 2-Nitroaniline	9.410	65	201137	43.950	ng	96
49) Acenaphthylene	9.733	152	973775	39.697	ng	99
50) Dimethylphthalate	9.598	163	703554	41.005	ng	100
51) 2,6-Dinitrotoluene	9.657	165	153573	47.147	ng	88
52) Acenaphthene	9.904	154	592444	39.871	ng	98
53) 3-Nitroaniline	9.827	138	181887	44.708	ng	# 89
54) 2,4-Dinitrophenol	9.922	184	49843	40.064	ng	89
55) Dibenzofuran	10.080	168	877303	39.685	ng	99
56) 4-Nitrophenol	9.975	139	134200	45.716	ng	93
57) 2,4-Dinitrotoluene	10.057	165	194630	48.063	ng	89
58) Fluorene	10.422	166	639757	39.499	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.192	232	179855	44.439	ng	98
60) Diethylphthalate	10.298	149	730500	41.013	ng	99
61) 4-Chlorophenyl-phenyle...	10.416	204	315645	40.587	ng	97
62) 4-Nitroaniline	10.439	138	173742	42.522	ng	96
63) Azobenzene	10.574	77	696641	39.073	ng	99
65) 4,6-Dinitro-2-methylph...	10.463	198	80180	42.772	ng	81
66) n-Nitrosodiphenylamine	10.533	169	612033	40.010	ng	99
67) 4-Bromophenyl-phenylether	10.904	248	207246	41.416	ng	94
68) Hexachlorobenzene	10.963	284	212157	40.544	ng	96
69) Atrazine	11.057	200	175464	41.805	ng	98
70) Pentachlorophenol	11.157	266	133626	42.707	ng	98
71) Phenanthrene	11.386	178	928466	38.905	ng	99
72) Anthracene	11.433	178	955600	38.906	ng	100
73) Carbazole	11.592	167	895746	38.563	ng	99
74) Di-n-butylphthalate	11.927	149	1127207	41.039	ng	100
75) Fluoranthene	12.574	202	920135	34.905	ng	97
77) Benzidine	12.692	184	207440m	28.629	ng	
78) Pyrene	12.804	202	911953	43.258	ng	99
80) Butylbenzylphthalate	13.427	149	382663	45.876	ng	95
81) Benzo(a)anthracene	13.992	228	661983	40.111	ng	99
82) 3,3'-Dichlorobenzidine	13.957	252	219885	42.620	ng	# 97
83) Chrysene	14.027	228	676067	40.578	ng	100
84) Bis(2-ethylhexyl)phtha...	13.986	149	473787	45.441	ng	99
85) Di-n-octyl phthalate	14.604	149	778332	44.813	ng	98
87) Indeno(1,2,3-cd)pyrene	16.909	276	738743	37.694	ng	99
88) Benzo(b)fluoranthene	15.033	252	752844	37.576	ng	99
89) Benzo(k)fluoranthene	15.062	252	681128	34.866	ng	99
90) Benzo(a)pyrene	15.398	252	709524	38.919	ng	99
91) Dibenzo(a,h)anthracene	16.927	278	629306	39.604	ng	98
92) Benzo(g,h,i)perylene	17.345	276	560172	34.962	ng	99

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

