

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF120822\
 Data File : BF131554.D
 Acq On : 08 Dec 2022 13:15
 Operator : CG\JU
 Sample : SSTDICC020
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC020

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 12/14/2022
 Supervised By :mohammad ahmed 12/14/2022

Quant Time: Dec 09 05:18:11 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF120822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 09 05:15:18 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.910	152	51997	20.000	ng	0.00
21) Naphthalene-d8	8.198	136	181787	20.000	ng	# 0.00
39) Acenaphthene-d10	9.963	164	109679	20.000	ng	0.00
64) Phenanthrene-d10	11.457	188	207587	20.000	ng	0.00
76) Chrysene-d12	14.104	240	114188	20.000	ng	# 0.00
86) Perylene-d12	15.615	264	91690	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.510	112	130574	42.414	ng	0.00
7) Phenol-d6	6.528	99	169711	41.629	ng	0.00
23) Nitrobenzene-d5	7.475	82	197395	42.493	ng	0.00
42) 2,4,6-Tribromophenol	10.751	330	49509	38.704	ng	0.00
45) 2-Fluorobiphenyl	9.280	172	350478	40.545	ng	0.00
79) Terphenyl-d14	13.045	244	342319	39.937	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.657	88	28998	21.660	ng	99
3) Pyridine	3.422	79	82417	22.125	ng	99
4) n-Nitrosodimethylamine	3.375	42	44992	22.425	ng	# 98
6) Aniline	6.569	93	100631	20.573	ng	98
8) 2-Chlorophenol	6.692	128	67816	20.714	ng	93
9) Benzaldehyde	6.457	77	60226	21.274	ng	94
10) Phenol	6.539	94	93782	20.524	ng	100
11) bis(2-Chloroethyl)ether	6.639	93	71113	20.463	ng	97
12) 1,3-Dichlorobenzene	6.851	146	78656	20.426	ng	97
13) 1,4-Dichlorobenzene	6.928	146	79087	20.360	ng	99
14) 1,2-Dichlorobenzene	7.081	146	75955	20.794	ng	97
15) Benzyl Alcohol	7.045	79	78785	20.724	ng	95
16) 2,2'-oxybis(1-Chloropr...	7.181	45	92971	21.353	ng	99
17) 2-Methylphenol	7.157	107	61096	20.338	ng	97
18) Hexachloroethane	7.422	117	32535	20.714	ng	88
19) n-Nitroso-di-n-propyla...	7.316	70	63043	20.107	ng	99
20) 3+4-Methylphenols	7.310	107	83911	20.297	ng	92
22) Acetophenone	7.322	105	116669	20.738	ng	94
24) Nitrobenzene	7.492	77	99994	21.372	ng	99
25) Isophorone	7.733	82	160523	20.646	ng	99
26) 2-Nitrophenol	7.810	139	32594	19.683	ng	94
27) 2,4-Dimethylphenol	7.845	122	51419	19.353	ng	96
28) bis(2-Chloroethoxy)met...	7.945	93	81632	20.217	ng	99
29) 2,4-Dichlorophenol	8.051	162	63804	20.407	ng	97
30) 1,2,4-Trichlorobenzene	8.139	180	78101	20.342	ng	98
31) Naphthalene	8.222	128	207654	20.223	ng	100
32) Benzoic acid	7.939	122	33652m	17.376	ng	
33) 4-Chloroaniline	8.269	127	80609	21.026	ng	98
34) Hexachlorobutadiene	8.333	225	55109	19.995	ng	97
35) Caprolactam	8.628	113	16387	19.596	ng	92
36) 4-Chloro-3-methylphenol	8.745	107	71704	20.344	ng	100
37) 2-Methylnaphthalene	8.916	142	142832	20.116	ng	100
38) 1-Methylnaphthalene	9.016	142	137615	19.993	ng	99
40) 1,2,4,5-Tetrachloroben...	9.080	216	82980	20.343	ng	99
41) Hexachlorocyclopentadiene	9.063	237	39179	18.606	ng	98
43) 2,4,6-Trichlorophenol	9.192	196	51370	20.098	ng	99

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44) 2,4,5-Trichlorophenol	9.227	196	54000	20.125	ng	95
46) 1,1'-Biphenyl	9.380	154	175872	20.363	ng	97
47) 2-Chloronaphthalene	9.404	162	146645	20.853	ng	98
48) 2-Nitroaniline	9.504	65	51719	21.319	ng	98
49) Acenaphthylene	9.822	152	213236	20.637	ng	99
50) Dimethylphthalate	9.680	163	171839	20.274	ng	99
51) 2,6-Dinitrotoluene	9.745	165	35066	20.111	ng	84
52) Acenaphthene	9.998	154	140181	20.146	ng	97
53) 3-Nitroaniline	9.916	138	33509	20.296	ng	99
54) 2,4-Dinitrophenol	10.016	184	16897	18.109	ng	# 90
55) Dibenzofuran	10.169	168	201339	20.394	ng	97
56) 4-Nitrophenol	10.063	139	23178	19.809	ng	97
57) 2,4-Dinitrotoluene	10.145	165	46487	20.102	ng	96
58) Fluorene	10.510	166	161267	19.888	ng	97
59) 2,3,4,6-Tetrachlorophenol	10.280	232	45566m	19.655	ng	
60) Diethylphthalate	10.380	149	166679	20.033	ng	99
61) 4-Chlorophenyl-phenyle...	10.504	204	89998	19.864	ng	97
62) 4-Nitroaniline	10.527	138	33074	19.975	ng	98
63) Azobenzene	10.663	77	190052	20.891	ng	96
65) 4,6-Dinitro-2-methylph...	10.551	198	23548	17.022	ng	89
66) n-Nitrosodiphenylamine	10.622	169	132011	19.528	ng	99
67) 4-Bromophenyl-phenylether	10.998	248	53856	19.745	ng	97
68) Hexachlorobenzene	11.057	284	56611	19.269	ng	# 88
69) Atrazine	11.145	200	43605	19.754	ng	100
70) Pentachlorophenol	11.251	266	28070	17.995	ng	99
71) Phenanthrene	11.480	178	229204	19.760	ng	100
72) Anthracene	11.533	178	224521	19.914	ng	99
73) Carbazole	11.686	167	183183	19.784	ng	99
74) Di-n-butylphthalate	12.016	149	220305	17.760	ng	99
75) Fluoranthene	12.674	202	237829	19.127	ng	99
77) Benzidine	12.798	184	40211	18.602	ng	98
78) Pyrene	12.904	202	233767	21.000	ng	100
80) Butylbenzylphthalate	13.521	149	64479	17.836	ng	98
81) Benzo(a)anthracene	14.098	228	167561	20.271	ng	98
82) 3,3'-Dichlorobenzidine	14.057	252	47543	20.665	ng	99
83) Chrysene	14.133	228	156141	19.863	ng	96
84) Bis(2-ethylhexyl)phtha...	14.080	149	78233	17.964	ng	97
85) Di-n-octyl phthalate	14.698	149	122059	18.263	ng	99
87) Indeno(1,2,3-cd)pyrene	17.151	276	155992	23.233	ng	99
88) Benzo(b)fluoranthene	15.168	252	120647	19.242	ng	96
89) Benzo(k)fluoranthene	15.198	252	134341	20.948	ng	99
90) Benzo(a)pyrene	15.551	252	102547	19.987	ng	98
91) Dibenzo(a,h)anthracene	17.168	278	128478	22.842	ng	97
92) Benzo(g,h,i)perylene	17.615	276	131934	24.403	ng	99

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

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