

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112723\
 Data File : BF136219.D
 Acq On : 27 Nov 2023 12:51
 Operator : CG\JU
 Sample : SSTDICC010
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC010

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 11/28/2023
 Supervised By :mohammad ahmed 11/29/2023

Quant Time: Nov 28 02:58:10 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF112723.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 28 02:38:48 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.804	152	164747	20.000	ng	0.00
21) Naphthalene-d8	8.086	136	733417	20.000	ng	0.00
39) Acenaphthene-d10	9.839	164	378219	20.000	ng	0.00
64) Phenanthrene-d10	11.333	188	725372	20.000	ng	0.00
76) Chrysene-d12	13.992	240	440598	20.000	ng	0.00
86) Perylene-d12	15.474	264	298146	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.434	112	229062	19.352	ng	0.00
7) Phenol-d6	6.434	99	287703	19.439	ng	-0.01
23) Nitrobenzene-d5	7.363	82	259737	18.714	ng	-0.01
42) 2,4,6-Tribromophenol	10.633	330	85793	18.554	ng	0.00
45) 2-Fluorobiphenyl	9.163	172	588956	19.890	ng	0.00
79) Terphenyl-d14	12.933	244	669922	17.539	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.610	88	43010	9.603	ng	95
3) Pyridine	3.363	79	106999	9.678	ng	97
4) n-Nitrosodimethylamine	3.310	42	52587	9.410	ng	95
6) Aniline	6.469	93	141326	10.103	ng	95
8) 2-Chlorophenol	6.586	128	122522	9.507	ng	97
9) Benzaldehyde	6.357	77	89625	10.865	ng	99
10) Phenol	6.445	94	154758	9.747	ng	93
11) bis(2-Chloroethyl)ether	6.539	93	110423	9.480	ng	97
12) 1,3-Dichlorobenzene	6.745	146	138010	9.436	ng	98
13) 1,4-Dichlorobenzene	6.822	146	140769	9.468	ng	96
14) 1,2-Dichlorobenzene	6.975	146	135912	9.656	ng	96
15) Benzyl Alcohol	6.945	79	98783	9.520	ng	97
16) 2,2'-oxybis(1-Chloropr...	7.081	45	158772	9.813	ng	99
17) 2-Methylphenol	7.051	107	95750	9.450	ng	99
18) Hexachloroethane	7.316	117	50217	9.629	ng	95
19) n-Nitroso-di-n-propyla...	7.210	70	87375	9.871	ng	96
20) 3+4-Methylphenols	7.204	107	128459	9.755	ng	92
22) Acetophenone	7.210	105	183214	9.512	ng	# 92
24) Nitrobenzene	7.381	77	131697	9.364	ng	99
25) Isophorone	7.622	82	236659	9.457	ng	96
26) 2-Nitrophenol	7.698	139	53453	8.920	ng	98
27) 2,4-Dimethylphenol	7.733	122	76502	9.699	ng	98
28) bis(2-Chloroethoxy)met...	7.833	93	142936	9.442	ng	99
29) 2,4-Dichlorophenol	7.939	162	101369	9.528	ng	99
30) 1,2,4-Trichlorobenzene	8.022	180	122078	9.372	ng	98
31) Naphthalene	8.104	128	386590	9.378	ng	98
32) Benzoic acid	7.810	122	58673	8.162	ng	92
33) 4-Chloroaniline	8.151	127	137252	9.795	ng	98
34) Hexachlorobutadiene	8.222	225	71265	9.352	ng	99
35) Caprolactam	8.492	113	28123	9.456	ng	# 84
36) 4-Chloro-3-methylphenol	8.627	107	109704	9.507	ng	100
37) 2-Methylnaphthalene	8.798	142	249112	9.633	ng	95
38) 1-Methylnaphthalene	8.892	142	239527	9.495	ng	94
40) 1,2,4,5-Tetrachloroben...	8.963	216	114985	9.553	ng	98
41) Hexachlorocyclopentadiene	8.951	237	39487	8.526	ng	93
43) 2,4,6-Trichlorophenol	9.075	196	74210	9.499	ng	94

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44) 2,4,5-Trichlorophenol	9.110	196	82220	9.643	ng	98
46) 1,1'-Biphenyl	9.263	154	321934	9.652	ng	97
47) 2-Chloronaphthalene	9.286	162	244858	9.453	ng	97
48) 2-Nitroaniline	9.380	65	62106	9.184	ng	90
49) Acenaphthylene	9.698	152	353562	9.565	ng	98
50) Dimethylphthalate	9.563	163	273990	9.585	ng	99
51) 2,6-Dinitrotoluene	9.622	165	54329	9.303	ng	85
52) Acenaphthene	9.874	154	225818m	9.667	ng	
53) 3-Nitroaniline	9.792	138	54730	9.445	ng	98
54) 2,4-Dinitrophenol	9.898	184	19975	7.354	ng	88
55) Dibenzofuran	10.045	168	337967	9.723	ng	99
56) 4-Nitrophenol	9.945	139	40987	9.313	ng	94
57) 2,4-Dinitrotoluene	10.027	165	71961	9.225	ng	98
58) Fluorene	10.392	166	272027	9.977	ng	96
59) 2,3,4,6-Tetrachlorophenol	10.163	232	65303	9.632	ng	98
60) Diethylphthalate	10.269	149	273194	9.554	ng	99
61) 4-Chlorophenyl-phenyle...	10.386	204	134328	10.145	ng	95
62) 4-Nitroaniline	10.404	138	53971	9.515	ng	97
63) Azobenzene	10.545	77	254126	9.586	ng	99
65) 4,6-Dinitro-2-methylph...	10.433	198	30291	7.636	ng	92
66) n-Nitrosodiphenylamine	10.504	169	224545	9.660	ng	94
67) 4-Bromophenyl-phenylether	10.880	248	78781	9.309	ng	91
68) Hexachlorobenzene	10.939	284	93964	9.594	ng	96
69) Atrazine	11.033	200	67446	9.781	ng	95
70) Pentachlorophenol	11.133	266	48846	8.823	ng	99
71) Phenanthrene	11.357	178	395092	9.691	ng	99
72) Anthracene	11.410	178	388925	9.513	ng	99
73) Carbazole	11.563	167	331688	9.539	ng	99
74) Di-n-butylphthalate	11.904	149	403173	9.267	ng	98
75) Fluoranthene	12.551	202	399171	9.645	ng	96
77) Benzidine	12.680	184	43480	6.003	ng	97
78) Pyrene	12.780	202	408515	8.466	ng	98
80) Butylbenzylphthalate	13.415	149	131842	8.779	ng	97
81) Benzo(a)anthracene	13.980	228	318136	9.721	ng	99
82) 3,3'-Dichlorobenzidine	13.945	252	76258	9.680	ng	95
83) Chrysene	14.015	228	292890	9.269	ng	98
84) Bis(2-ethylhexyl)phtha...	13.980	149	160897	8.927	ng	# 99
85) Di-n-octyl phthalate	14.598	149	200461	11.274	ng	# 97
87) Indeno(1,2,3-cd)pyrene	16.968	276	155457	9.265	ng	96
88) Benzo(b)fluoranthene	15.033	252	211059	10.090	ng	98
89) Benzo(k)fluoranthene	15.062	252	196855	9.909	ng	98
90) Benzo(a)pyrene	15.404	252	159813	9.346	ng	99
91) Dibenzo(a,h)anthracene	16.998	278	128251	9.478	ng	98
92) Benzo(g,h,i)perylene	17.427	276	135825	9.169	ng	99

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

