

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF100422\
 Data File : BF130517.D
 Acq On : 04 Oct 2022 14:40
 Operator : CG\JU
 Sample : N4914-02MS
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 S1MS

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 10/05/2022
 Supervised By :mohammad ahmed 10/06/2022

Quant Time: Oct 05 02:17:21 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 28 02:59:31 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.633	152	81495	20.000	ng	-0.01
21) Naphthalene-d8	7.922	136	305728	20.000	ng	0.00
39) Acenaphthene-d10	9.663	164	137174	20.000	ng	-0.02
64) Phenanthrene-d10	11.145	188	194309	20.000	ng	-0.01
76) Chrysene-d12	13.780	240	162067	20.000	ng	-0.01
86) Perylene-d12	15.168	264	209306	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.257	112	702770	149.571	ng	0.01
7) Phenol-d6	6.292	99	768547	130.728	ng	0.00
23) Nitrobenzene-d5	7.204	82	589108	98.443	ng	-0.01
42) 2,4,6-Tribromophenol	10.457	330	200216	152.326	ng	-0.01
45) 2-Fluorobiphenyl	8.998	172	994275	105.549	ng	-0.01
79) Terphenyl-d14	12.733	244	802700	82.044	ng	-0.01
Target Compounds						Qvalue
2) 1,4-Dioxane	2.404	88	131389	60.888	ng	# 84
3) Pyridine	3.069	79	294814	52.374	ng	93
4) n-Nitrosodimethylamine	2.993	42	136788	44.679	ng	92
6) Aniline	6.304	93	125954	17.388	ng	# 1
8) 2-Chlorophenol	6.422	128	279157	52.157	ng	98
9) Benzaldehyde	6.186	77	90513	22.984	ng	91
10) Phenol	6.304	94	852002	123.665	ng	99
11) bis(2-Chloroethyl)ether	6.381	93	274626	55.264	ng	97
12) 1,3-Dichlorobenzene	6.575	146	298371	50.663	ng	98
13) 1,4-Dichlorobenzene	6.651	146	304665	50.729	ng	100
14) 1,2-Dichlorobenzene	6.804	146	288110	51.354	ng	98
15) Benzyl Alcohol	6.792	79	190995	40.975	ng	95
16) 2,2'-oxybis(1-Chloropr...	6.916	45	345617	47.685	ng	# 35
17) 2-Methylphenol	6.910	107	347089	79.774	ng	95
18) Hexachloroethane	7.139	117	120677	50.976	ng	95
19) n-Nitroso-di-n-propyla...	7.063	70	192813	48.604	ng	93
20) 3+4-Methylphenols	7.069	107	601534	106.427	ng	# 55
22) Acetophenone	7.057	105	405894	54.303	ng	# 98
24) Nitrobenzene	7.222	77	300543	49.461	ng	99
25) Isophorone	7.469	82	528177	54.760	ng	99
26) 2-Nitrophenol	7.539	139	144806	58.134	ng	92
27) 2,4-Dimethylphenol	7.592	122	324507	84.609	ng	97
28) bis(2-Chloroethoxy)met...	7.680	93	350625	64.559	ng	98
29) 2,4-Dichlorophenol	7.792	162	206074	49.031	ng	96
30) 1,2,4-Trichlorobenzene	7.863	180	229859	50.585	ng	97
31) Naphthalene	7.951	128	5447693	345.869	ng	97
32) Benzoic acid	7.728	122	223044	75.201	ng	# 33
33) 4-Chloroaniline	7.998	127	78482	13.101	ng	99
34) Hexachlorobutadiene	8.057	225	146009	50.278	ng	99
35) Caprolactam	8.380	113	45840	38.045	ng	# 86
36) 4-Chloro-3-methylphenol	8.492	107	214865	46.122	ng	96
37) 2-Methylnaphthalene	8.633	142	748827	74.566	ng	99
38) 1-Methylnaphthalene	8.727	142	657256	68.128	ng	99
40) 1,2,4,5-Tetrachloroben...	8.798	216	219655	58.676	ng	97
41) Hexachlorocyclopentadiene	8.780	237	175722	87.674	ng	99
43) 2,4,6-Trichlorophenol	8.916	196	130177	49.826	ng	95

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	8.963	196	137172	48.654	ng	99
46) 1,1'-Biphenyl	9.092	154	630440	61.538	ng	99
47) 2-Chloronaphthalene	9.116	162	448544	54.124	ng	99
48) 2-Nitroaniline	9.222	65	125771	45.502	ng	95
49) Acenaphthylene	9.527	152	823637	66.285	ng	99
50) Dimethylphthalate	9.404	163	466840	49.269	ng	99
51) 2,6-Dinitrotoluene	9.463	165	98865	49.890	ng	90
52) Acenaphthene	9.698	154	409007	50.099	ng	98
53) 3-Nitroaniline	9.633	138	41718	18.893	ng	91
54) 2,4-Dinitrophenol	9.745	184	89853	83.305	ng	88
55) Dibenzofuran	9.874	168	637056	56.101	ng	97
56) 4-Nitrophenol	9.810	139	107796	67.025	ng	89
57) 2,4-Dinitrotoluene	9.869	165	123226	48.656	ng	# 93
58) Fluorene	10.216	166	512499	55.186	ng	99
59) 2,3,4,6-Tetrachlorophenol	9.998	232	99870	45.257	ng	95
60) Diethylphthalate	10.098	149	438365	46.134	ng	100
61) 4-Chlorophenyl-phenyle...	10.210	204	207659	50.973	ng	96
62) 4-Nitroaniline	10.245	138	84661m	38.166	ng	
63) Azobenzene	10.369	77	440783	45.050	ng	96
65) 4,6-Dinitro-2-methylph...	10.274	198	57408	53.074	ng	86
66) n-Nitrosodiphenylamine	10.327	169	362496	55.827	ng	98
67) 4-Bromophenyl-phenylether	10.698	248	125489	58.543	ng	94
68) Hexachlorobenzene	10.757	284	139480	57.403	ng	96
69) Atrazine	10.863	200	86804	45.745	ng	98
70) Pentachlorophenol	10.963	266	115575	88.626	ng	97
71) Phenanthrene	11.174	178	685265	62.815	ng	99
72) Anthracene	11.221	178	590578	56.097	ng	99
73) Carbazole	11.386	167	700889	73.769	ng	99
74) Di-n-butylphthalate	11.716	149	614223	53.610	ng	99
75) Fluoranthene	12.357	202	562081	53.320	ng	98
77) Benzidine	12.486	184	69111	14.690	ng	99
78) Pyrene	12.580	202	569931	40.199	ng	99
80) Butylbenzylphthalate	13.210	149	260288	49.557	ng	99
81) Benzo(a)anthracene	13.768	228	569679	52.839	ng	99
82) 3,3'-Dichlorobenzidine	13.739	252	138886	47.323	ng	99
83) Chrysene	13.804	228	546572	53.392	ng	99
84) Bis(2-ethylhexyl)phtha...	13.768	149	372294	61.883	ng	99
85) Di-n-octyl phthalate	14.380	149	701025	76.184	ng	97
87) Indeno(1,2,3-cd)pyrene	16.486	276	863420	58.171	ng	98
88) Benzo(b)fluoranthene	14.780	252	757210	55.427	ng	99
89) Benzo(k)fluoranthene	14.809	252	602169	44.809	ng	99
90) Benzo(a)pyrene	15.115	252	613740	56.707	ng	99
91) Dibenzo(a,h)anthracene	16.503	278	743130	61.031	ng	98
92) Benzo(g,h,i)perylene	16.886	276	732304	59.703	ng	# 95

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

