

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112522\
 Data File : BF131390.D
 Acq On : 25 Nov 2022 11:44
 Operator : CG\JU
 Sample : PB149192BS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB149192BS

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 12/02/2022
 Supervised By : mohammad ahmed 12/02/2022

Quant Time: Nov 25 23:15:44 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF112122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 25 23:13:01 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.957	152	110871	20.000	ng	0.00
21) Naphthalene-d8	8.251	136	438902	20.000	ng	0.00
39) Acenaphthene-d10	10.016	164	253182	20.000	ng	0.00
64) Phenanthrene-d10	11.516	188	464515	20.000	ng	0.00
76) Chrysene-d12	14.174	240	298800	20.000	ng	0.00
86) Perylene-d12	15.710	264	247028	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.581	112	769601	132.110	ng	0.02
7) Phenol-d6	6.581	99	981414	132.042	ng	0.00
23) Nitrobenzene-d5	7.528	82	641856	73.718	ng	0.00
42) 2,4,6-Tribromophenol	10.816	330	298379	140.081	ng	0.00
45) 2-Fluorobiphenyl	9.333	172	1185195	68.103	ng	0.00
79) Terphenyl-d14	13.110	244	1389748	76.058	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.852	88	93188	33.554	ng	96
3) Pyridine	3.604	79	257169	33.348	ng	96
4) n-Nitrosodimethylamine	3.575	42	182572	40.146	ng	95
6) Aniline	6.622	93	393305m	40.377	ng	
8) 2-Chlorophenol	6.740	128	306305	46.411	ng	100
9) Benzaldehyde	6.510	77	182032	39.804	ng	97
10) Phenol	6.592	94	373287	45.298	ng	99
11) bis(2-Chloroethyl)ether	6.692	93	284915	41.266	ng	98
12) 1,3-Dichlorobenzene	6.898	146	313426	39.847	ng	98
13) 1,4-Dichlorobenzene	6.975	146	319402	39.963	ng	98
14) 1,2-Dichlorobenzene	7.128	146	310495	42.192	ng	98
15) Benzyl Alcohol	7.098	79	285432	46.836	ng	99
16) 2,2'-oxybis(1-Chloropr...	7.234	45	513615	38.898	ng	99
17) 2-Methylphenol	7.204	107	249015	44.269	ng	99
18) Hexachloroethane	7.475	117	121694	42.564	ng	98
19) n-Nitroso-di-n-propyla...	7.375	70	223092	41.001	ng	99
20) 3+4-Methylphenols	7.363	107	313338m	43.627	ng	
22) Acetophenone	7.375	105	425268	38.066	ng	# 95
24) Nitrobenzene	7.545	77	344445	37.970	ng	98
25) Isophorone	7.787	82	620662	39.955	ng	100
26) 2-Nitrophenol	7.863	139	156714	51.954	ng	91
27) 2,4-Dimethylphenol	7.892	122	277309	46.049	ng	98
28) bis(2-Chloroethoxy)met...	7.992	93	354518	40.182	ng	99
29) 2,4-Dichlorophenol	8.104	162	267001	40.875	ng	95
30) 1,2,4-Trichlorobenzene	8.186	180	290911	35.951	ng	99
31) Naphthalene	8.275	128	847699	36.244	ng	99
32) Benzoic acid	8.022	122	209057	53.339	ng	95
33) 4-Chloroaniline	8.316	127	253508	27.474	ng	99
34) Hexachlorobutadiene	8.386	225	181234	34.588	ng	97
35) Caprolactam	8.698	113	86898m	50.625	ng	
36) 4-Chloro-3-methylphenol	8.798	107	292404	43.087	ng	95
37) 2-Methylnaphthalene	8.969	142	578970	36.534	ng	100
38) 1-Methylnaphthalene	9.069	142	543683	35.379	ng	99
40) 1,2,4,5-Tetrachloroben...	9.133	216	297757	36.241	ng	99
41) Hexachlorocyclopentadiene	9.116	237	367948	99.621	ng	98
43) 2,4,6-Trichlorophenol	9.239	196	206755	44.360	ng	98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.281	196	217676	41.533	ng	98
46) 1,1'-Biphenyl	9.433	154	714266	36.969	ng	98
47) 2-Chloronaphthalene	9.463	162	568816	36.700	ng	98
48) 2-Nitroaniline	9.557	65	209676	44.556	ng	95
49) Acenaphthylene	9.880	152	885701	37.487	ng	100
50) Dimethylphthalate	9.739	163	675533	38.295	ng	99
51) 2,6-Dinitrotoluene	9.798	165	153720	43.574	ng	93
52) Acenaphthene	10.051	154	636465m	43.020	ng	
53) 3-Nitroaniline	9.969	138	121807	33.602	ng	91
54) 2,4-Dinitrophenol	10.075	184	169948	96.529	ng	98
55) Dibenzofuran	10.228	168	777373	36.523	ng	100
56) 4-Nitrophenol	10.128	139	246499	97.980	ng	99
57) 2,4-Dinitrotoluene	10.210	165	211421	47.631	ng	85
58) Fluorene	10.569	166	622338	37.464	ng	98
59) 2,3,4,6-Tetrachlorophenol	10.339	232	189667	43.454	ng	99
60) Diethylphthalate	10.439	149	644209	38.676	ng	99
61) 4-Chlorophenyl-phenyle...	10.563	204	318542	37.407	ng	95
62) 4-Nitroaniline	10.592	138	157355	43.617	ng	98
63) Azobenzene	10.722	77	648459	36.207	ng	97
65) 4,6-Dinitro-2-methylph...	10.616	198	104902	51.429	ng	85
66) n-Nitrosodiphenylamine	10.680	169	546591	36.928	ng	100
67) 4-Bromophenyl-phenylether	11.057	248	206862	36.335	ng	95
68) Hexachlorobenzene	11.116	284	224873	37.215	ng	98
69) Atrazine	11.210	200	203356	44.622	ng	100
70) Pentachlorophenol	11.316	266	272268	77.696	ng	99
71) Phenanthrene	11.539	178	924291	36.819	ng	99
72) Anthracene	11.592	178	943942	38.261	ng	99
73) Carbazole	11.745	167	849246	40.354	ng	100
74) Di-n-butylphthalate	12.074	149	970840	39.149	ng	99
75) Fluoranthene	12.733	202	1010963	38.217	ng	100
77) Benzidine	12.857	184	441003	79.903	ng	96
78) Pyrene	12.969	202	1017113	38.581	ng	100
80) Butylbenzylphthalate	13.586	149	359243	41.151	ng	93
81) Benzo(a)anthracene	14.163	228	792307	38.685	ng	99
82) 3,3'-Dichlorobenzidine	14.121	252	226073	41.597	ng	100
83) Chrysene	14.204	228	748124	37.231	ng	99
84) Bis(2-ethylhexyl)phtha...	14.145	149	432484	37.623	ng	# 99
85) Di-n-octyl phthalate	14.768	149	626783	39.658	ng	97
87) Indeno(1,2,3-cd)pyrene	17.292	276	661630	39.878	ng	100
88) Benzo(b)fluoranthene	15.257	252	684569	41.715	ng	99
89) Benzo(k)fluoranthene	15.286	252	657802	39.005	ng	99
90) Benzo(a)pyrene	15.645	252	583534	43.336	ng	# 98
91) Dibenzo(a,h)anthracene	17.315	278	559243	40.827	ng	99
92) Benzo(g,h,i)perylene	17.780	276	557417	40.772	ng	98

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

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