

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042523\
 Data File : BF133027.D
 Acq On : 25 Apr 2023 12:36
 Operator : CG\JU
 Sample : PB152211BS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB152211BS

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 04/26/2023
 Supervised By : Jagrut Upadhyay 04/26/2023

Quant Time: Apr 25 13:43:39 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042123.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Apr 22 03:30:56 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.745	152	250618	20.000	ng	0.00
21) Naphthalene-d8	8.033	136	951591	20.000	ng	0.00
39) Acenaphthene-d10	9.786	164	519715	20.000	ng	0.00
64) Phenanthrene-d10	11.269	188	910422	20.000	ng	0.00
76) Chrysene-d12	13.904	240	586442	20.000	ng	0.00
86) Perylene-d12	15.321	264	487957	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.375	112	1638165	98.741	ng	0.02
7) Phenol-d6	6.392	99	2018099	98.003	ng	0.00
23) Nitrobenzene-d5	7.316	82	1203450	68.263	ng	0.00
42) 2,4,6-Tribromophenol	10.574	330	551807	105.574	ng	0.00
45) 2-Fluorobiphenyl	9.110	172	2122337	68.320	ng	0.00
79) Terphenyl-d14	12.851	244	2351164	69.145	ng	-0.01
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.528	88	306866	39.877	ng	98
3) Pyridine	3.222	79	766419	37.499	ng	99
4) n-Nitrosodimethylamine	3.187	42	415088	39.209	ng	96
6) Aniline	6.410	93	878993	33.142	ng	# 13
8) 2-Chlorophenol	6.534	128	727831	43.208	ng	99
9) Benzaldehyde	6.292	77	409630	45.036	ng	99
10) Phenol	6.410	94	910959	40.100	ng	78
11) bis(2-Chloroethyl)ether	6.487	93	718534	42.150	ng	98
12) 1,3-Dichlorobenzene	6.686	146	777767	43.429	ng	99
13) 1,4-Dichlorobenzene	6.763	146	783969	43.678	ng	99
14) 1,2-Dichlorobenzene	6.916	146	715394	43.232	ng	96
15) Benzyl Alcohol	6.892	79	573548	40.321	ng	99
16) 2,2'-oxybis(1-Chloropr...	7.028	45	1166448	39.395	ng	96
17) 2-Methylphenol	7.010	107	628783	40.765	ng	97
18) Hexachloroethane	7.257	117	263564	42.236	ng	94
19) n-Nitroso-di-n-propyla...	7.169	70	493157	38.468	ng	98
20) 3+4-Methylphenols	7.169	107	720621	39.671	ng	# 75
22) Acetophenone	7.157	105	991914	44.999	ng	# 96
24) Nitrobenzene	7.334	77	741787	41.523	ng	99
25) Isophorone	7.575	82	1424164	42.547	ng	100
26) 2-Nitrophenol	7.651	139	414106	48.059	ng	# 89
27) 2,4-Dimethylphenol	7.692	122	671998	45.482	ng	100
28) bis(2-Chloroethoxy)met...	7.786	93	835562	42.196	ng	99
29) 2,4-Dichlorophenol	7.898	162	599633	44.581	ng	96
30) 1,2,4-Trichlorobenzene	7.975	180	620856	44.555	ng	100
31) Naphthalene	8.057	128	1998971	42.016	ng	99
32) Benzoic acid	7.833	122	508848	55.755	ng	97
33) 4-Chloroaniline	8.098	127	322062	16.754	ng	98
34) Hexachlorobutadiene	8.175	225	342490	44.346	ng	99
35) Caprolactam	8.486	113	216878m	47.998	ng	
36) 4-Chloro-3-methylphenol	8.592	107	669447	46.154	ng	98
37) 2-Methylnaphthalene	8.745	142	1344838	41.345	ng	98
38) 1-Methylnaphthalene	8.845	142	1345591	44.222	ng	99
40) 1,2,4,5-Tetrachloroben...	8.916	216	565436	40.464	ng	99
41) Hexachlorocyclopentadiene	8.904	237	807805	99.123	ng	98
43) 2,4,6-Trichlorophenol	9.022	196	415801	43.027	ng	98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.063	196	447964	40.081	ng	99
46) 1,1'-Biphenyl	9.210	154	1678087	40.719	ng	99
47) 2-Chloronaphthalene	9.233	162	1258396	41.719	ng	99
48) 2-Nitroaniline	9.328	65	412202	41.348	ng	95
49) Acenaphthylene	9.645	152	1975295	40.984	ng	100
50) Dimethylphthalate	9.516	163	1501580	41.436	ng	99
51) 2,6-Dinitrotoluene	9.575	165	354681	43.485	ng	# 83
52) Acenaphthene	9.822	154	1447733m	44.849	ng	
53) 3-Nitroaniline	9.739	138	279708	28.841	ng	99
54) 2,4-Dinitrophenol	9.845	184	427076	104.170	ng	95
55) Dibenzofuran	9.992	168	1780428	40.434	ng	99
56) 4-Nitrophenol	9.910	139	675753	89.964	ng	92
57) 2,4-Dinitrotoluene	9.975	165	470394	45.223	ng	97
58) Fluorene	10.333	166	1297042	40.205	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.110	232	381981	44.085	ng	97
60) Diethylphthalate	10.216	149	1417926	41.424	ng	98
61) 4-Chlorophenyl-phenyle...	10.327	204	635069	40.396	ng	100
62) 4-Nitroaniline	10.357	138	404409	45.370	ng	98
63) Azobenzene	10.486	77	1445790	39.872	ng	99
65) 4,6-Dinitro-2-methylph...	10.386	198	269276	48.797	ng	98
66) n-Nitrosodiphenylamine	10.445	169	1205245	40.812	ng	99
67) 4-Bromophenyl-phenylether	10.816	248	417871	42.320	ng	98
68) Hexachlorobenzene	10.880	284	456975	45.165	ng	99
69) Atrazine	10.980	200	408869	48.049	ng	98
70) Pentachlorophenol	11.074	266	561249	77.888	ng	98
71) Phenanthrene	11.292	178	1968950	40.238	ng	99
72) Anthracene	11.345	178	2037223	40.682	ng	99
73) Carbazole	11.498	167	1882830	42.225	ng	99
74) Di-n-butylphthalate	11.833	149	2235332	44.298	ng	100
75) Fluoranthene	12.474	202	2039561	41.531	ng	99
77) Benzidine	12.598	184	1108571	111.784	ng	# 93
78) Pyrene	12.704	202	2076636	41.308	ng	98
80) Butylbenzylphthalate	13.327	149	906614	50.934	ng	95
81) Benzo(a)anthracene	13.892	228	1631646	40.460	ng	99
82) 3,3'-Dichlorobenzidine	13.857	252	363870	32.662	ng	100
83) Chrysene	13.933	228	1657070	41.100	ng	99
84) Bis(2-ethylhexyl)phtha...	13.892	149	969760	43.405	ng	100
85) Di-n-octyl phthalate	14.504	149	1615645	44.349	ng	98
87) Indeno(1,2,3-cd)pyrene	16.709	276	1401774	50.503	ng	100
88) Benzo(b)fluoranthene	14.915	252	1455485	46.886	ng	98
89) Benzo(k)fluoranthene	14.951	252	1255152	40.788	ng	99
90) Benzo(a)pyrene	15.262	252	1145018	40.523	ng	99
91) Dibenzo(a,h)anthracene	16.727	278	1166786	50.401	ng	99
92) Benzo(g,h,i)perylene	17.133	276	1184478	51.826	ng	99

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

