

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042723\
 Data File : BF133094.D
 Acq On : 27 Apr 2023 12:15
 Operator : CG\JU
 Sample : PB152410BS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB152410BS

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 04/28/2023
 Supervised By : Jagrut Upadhyay 04/28/2023

Quant Time: Apr 28 02:05:28 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042123.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 28 02:00:53 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.740	152	197994	20.000	ng	0.00
21) Naphthalene-d8	8.022	136	795223	20.000	ng	0.00
39) Acenaphthene-d10	9.775	164	414621	20.000	ng	0.00
64) Phenanthrene-d10	11.257	188	743732	20.000	ng	0.00
76) Chrysene-d12	13.898	240	522731	20.000	ng	0.00
86) Perylene-d12	15.310	264	466023	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.363	112	1610777	122.895	ng	0.02
7) Phenol-d6	6.387	99	1935068	118.947	ng	0.00
23) Nitrobenzene-d5	7.304	82	1201270	81.538	ng	0.00
42) 2,4,6-Tribromophenol	10.569	330	530094	127.127	ng	0.00
45) 2-Fluorobiphenyl	9.104	172	2027725	81.819	ng	0.00
79) Terphenyl-d14	12.845	244	2430867	80.203	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.481	88	203447	33.465	ng	97
3) Pyridine	3.175	79	535564	33.168	ng	98
4) n-Nitrosodimethylamine	3.134	42	305307	36.504	ng	95
6) Aniline	6.398	93	709472	33.860	ng	# 26
8) 2-Chlorophenol	6.522	128	601979	45.235	ng	97
9) Benzaldehyde	6.281	77	321788	44.626	ng	99
10) Phenol	6.398	94	698472	38.919	ng	# 70
11) bis(2-Chloroethyl)ether	6.475	93	564658	41.927	ng	99
12) 1,3-Dichlorobenzene	6.681	146	624949	44.170	ng	98
13) 1,4-Dichlorobenzene	6.757	146	628556	44.327	ng	98
14) 1,2-Dichlorobenzene	6.910	146	610170	46.674	ng	99
15) Benzyl Alcohol	6.887	79	465116	41.389	ng	99
16) 2,2'-oxybis(1-Chloropr...	7.022	45	918498	39.266	ng	95
17) 2-Methylphenol	7.004	107	505256	41.463	ng	98
18) Hexachloroethane	7.251	117	216616	43.939	ng	98
19) n-Nitroso-di-n-propyla...	7.163	70	383058	37.822	ng	94
20) 3+4-Methylphenols	7.157	107	582417	40.585	ng	# 89
22) Acetophenone	7.151	105	791973	42.993	ng	# 98
24) Nitrobenzene	7.328	77	607510	40.693	ng	95
25) Isophorone	7.563	82	1100750	39.351	ng	99
26) 2-Nitrophenol	7.640	139	334178	46.409	ng	96
27) 2,4-Dimethylphenol	7.687	122	536127	43.421	ng	98
28) bis(2-Chloroethoxy)met...	7.775	93	658007	39.763	ng	100
29) 2,4-Dichlorophenol	7.887	162	476855	42.424	ng	98
30) 1,2,4-Trichlorobenzene	7.969	180	503449	43.234	ng	99
31) Naphthalene	8.045	128	1613131	40.573	ng	99
32) Benzoic acid	7.816	122	383320	50.260	ng	96
33) 4-Chloroaniline	8.092	127	371860	23.149	ng	99
34) Hexachlorobutadiene	8.163	225	272242	42.182	ng	100
35) Caprolactam	8.469	113	175705m	46.532	ng	
36) 4-Chloro-3-methylphenol	8.581	107	516965	42.650	ng	98
37) 2-Methylnaphthalene	8.734	142	1087233	39.998	ng	99
38) 1-Methylnaphthalene	8.834	142	1016155	39.962	ng	100
40) 1,2,4,5-Tetrachloroben...	8.904	216	451187	40.473	ng	100
41) Hexachlorocyclopentadiene	8.892	237	589473	90.667	ng	98
43) 2,4,6-Trichlorophenol	9.016	196	322534	41.835	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.057	196	350065	39.261	ng	97
46) 1,1'-Biphenyl	9.198	154	1337776	40.690	ng	99
47) 2-Chloronaphthalene	9.222	162	997494	41.451	ng	98
48) 2-Nitroaniline	9.322	65	322894	40.600	ng	95
49) Acenaphthylene	9.639	152	1549742	40.304	ng	99
50) Dimethylphthalate	9.504	163	1170570	40.489	ng	99
51) 2,6-Dinitrotoluene	9.563	165	277464	42.641	ng	87
52) Acenaphthene	9.810	154	1156970m	44.926	ng	
53) 3-Nitroaniline	9.728	138	246716	31.888	ng	99
54) 2,4-Dinitrophenol	9.839	184	337658	103.235	ng	92
55) Dibenzofuran	9.981	168	1400804	39.876	ng	98
56) 4-Nitrophenol	9.904	139	539228	89.984	ng	97
57) 2,4-Dinitrotoluene	9.969	165	375972	45.307	ng	# 82
58) Fluorene	10.328	166	1044631	40.589	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.098	232	300324	43.446	ng	97
60) Diethylphthalate	10.204	149	1103604	40.413	ng	99
61) 4-Chlorophenyl-phenyle...	10.316	204	498846	39.774	ng	97
62) 4-Nitroaniline	10.345	138	330849	46.526	ng	97
63) Azobenzene	10.475	77	1095599	37.873	ng	98
65) 4,6-Dinitro-2-methylph...	10.375	198	217275	48.198	ng	91
66) n-Nitrosodiphenylamine	10.439	169	955070	39.589	ng	98
67) 4-Bromophenyl-phenylether	10.804	248	319755	39.641	ng	97
68) Hexachlorobenzene	10.875	284	361815	43.775	ng	96
69) Atrazine	10.969	200	318165	45.770	ng	96
70) Pentachlorophenol	11.069	266	453063	76.966	ng	97
71) Phenanthrene	11.280	178	1627290	40.709	ng	100
72) Anthracene	11.333	178	1653594	40.422	ng	99
73) Carbazole	11.492	167	1553560	42.650	ng	100
74) Di-n-butylphthalate	11.822	149	1754879	42.572	ng	100
75) Fluoranthene	12.469	202	1687026	42.052	ng	98
77) Benzidine	12.592	184	846513	95.763	ng	# 93
78) Pyrene	12.698	202	1752191	39.103	ng	99
80) Butylbenzylphthalate	13.321	149	743649	46.870	ng	98
81) Benzo(a)anthracene	13.886	228	1459280	40.596	ng	99
82) 3,3'-Dichlorobenzidine	13.845	252	413190	41.609	ng	99
83) Chrysene	13.921	228	1436785	39.980	ng	99
84) Bis(2-ethylhexyl)phtha...	13.880	149	851096	42.737	ng	99
85) Di-n-octyl phthalate	14.492	149	1591777	49.020	ng	97
87) Indeno(1,2,3-cd)pyrene	16.698	276	1194653	45.067	ng	99
88) Benzo(b)fluoranthene	14.910	252	1346760	45.425	ng	99
89) Benzo(k)fluoranthene	14.939	252	1269089	43.182	ng	99
90) Benzo(a)pyrene	15.257	252	1080979	40.058	ng	99
91) Dibenzo(a,h)anthracene	16.715	278	994816	44.995	ng	98
92) Benzo(g,h,i)perylene	17.115	276	999480	45.790	ng	98

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

