

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF020325\
 Data File : BF141364.D
 Acq On : 03 Feb 2025 12:44
 Operator : RC/JU
 Sample : PB166468BSD
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB166468BSD

Manual Integrations APPROVED

Reviewed By :Jagrut Upadhyay 02/04/2025
 Supervised By :mohammad ahmed 02/06/2025

Quant Time: Feb 03 13:04:39 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF012925.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jan 29 17:41:25 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.798	152	92246	20.000	ng	0.00
21) Naphthalene-d8	8.081	136	377694	20.000	ng	0.00
39) Acenaphthene-d10	9.833	164	203609	20.000	ng	0.00
64) Phenanthrene-d10	11.321	188	343734	20.000	ng	0.00
76) Chrysene-d12	13.963	240	243486	20.000	ng	-0.01
86) Perylene-d12	15.421	264	234414	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.422	112	546522	90.955	ng	0.01
7) Phenol-d6	6.439	99	661544	86.546	ng	0.00
23) Nitrobenzene-d5	7.357	82	642716	89.707	ng	-0.01
42) 2,4,6-Tribromophenol	10.627	330	179421	96.121	ng	0.00
45) 2-Fluorobiphenyl	9.157	172	1187345	89.758	ng	0.00
79) Terphenyl-d14	12.910	244	1315661	83.331	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.563	88	110356	41.760	ng	99
3) Pyridine	3.293	79	255116	40.068	ng	96
4) n-Nitrosodimethylamine	3.240	42	167602	46.079	ng	94
6) Aniline	6.457	93	267476	41.338	ng	# 32
8) 2-Chlorophenol	6.587	128	279168	43.347	ng	97
9) Benzaldehyde	6.345	77	225965	51.036	ng	99
10) Phenol	6.451	94	340789	42.058	ng	# 69
11) bis(2-Chloroethyl)ether	6.534	93	259872	41.828	ng	97
12) 1,3-Dichlorobenzene	6.739	146	286427	40.336	ng	99
13) 1,4-Dichlorobenzene	6.816	146	290586	40.734	ng	99
14) 1,2-Dichlorobenzene	6.963	146	280032	41.921	ng	100
15) Benzyl Alcohol	6.939	79	245815	47.057	ng	99
16) 2,2'-oxybis(1-Chloropr...	7.075	45	442997	41.005	ng	75
17) 2-Methylphenol	7.063	107	218408	41.727	ng	# 90
18) Hexachloroethane	7.310	117	109768	41.710	ng	96
19) n-Nitroso-di-n-propyla...	7.210	70	195251	42.988	ng	100
20) 3+4-Methylphenols	7.216	107	286970	44.652	ng	# 83
22) Acetophenone	7.204	105	425016	47.341	ng	# 96
24) Nitrobenzene	7.381	77	294998	39.732	ng	99
25) Isophorone	7.616	82	526883	42.543	ng	98
26) 2-Nitrophenol	7.692	139	146508	41.840	ng	96
27) 2,4-Dimethylphenol	7.739	122	235515m	52.842	ng	
28) bis(2-Chloroethoxy)met...	7.828	93	324070	41.022	ng	99
29) 2,4-Dichlorophenol	7.945	162	229279	42.020	ng	100
30) 1,2,4-Trichlorobenzene	8.022	180	244875	39.298	ng	100
31) Naphthalene	8.098	128	800918	40.141	ng	100
32) Benzoic acid	7.863	122	191082	46.637	ng	98
33) 4-Chloroaniline	8.151	127	131775	19.476	ng	98
34) Hexachlorobutadiene	8.216	225	147922	40.479	ng	99
35) Caprolactam	8.516	113	80262m	45.041	ng	
36) 4-Chloro-3-methylphenol	8.645	107	253967	43.393	ng	97
37) 2-Methylnaphthalene	8.792	142	507995	40.058	ng	100
38) 1-Methylnaphthalene	8.892	142	500486	40.177	ng	99
40) 1,2,4,5-Tetrachloroben...	8.957	216	271467	46.866	ng	99
41) Hexachlorocyclopentadiene	8.945	237	320846	187.555	ng	100
43) 2,4,6-Trichlorophenol	9.075	196	164100	43.283	ng	99

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44) 2,4,5-Trichlorophenol	9.122	196	178649	43.293	ng	99
46) 1,1'-Biphenyl	9.257	154	735017	46.289	ng	99
47) 2-Chloronaphthalene	9.280	162	501873	40.549	ng	99
48) 2-Nitroaniline	9.375	65	170006	43.669	ng	96
49) Acenaphthylene	9.698	152	788250	45.497	ng	100
50) Dimethylphthalate	9.557	163	589543	42.871	ng	99
51) 2,6-Dinitrotoluene	9.622	165	130295	40.807	ng	99
52) Acenaphthene	9.869	154	588646m	47.559	ng	
53) 3-Nitroaniline	9.786	138	74473	23.797	ng	96
54) 2,4-Dinitrophenol	9.898	184	155968	105.080	ng	94
55) Dibenzofuran	10.039	168	709857	42.846	ng	98
56) 4-Nitrophenol	9.969	139	226393	98.918	ng	93
57) 2,4-Dinitrotoluene	10.022	165	179950	43.529	ng	95
58) Fluorene	10.386	166	553666	43.083	ng	98
59) 2,3,4,6-Tetrachlorophenol	10.163	232	150622	46.397	ng	98
60) Diethylphthalate	10.257	149	580043	43.502	ng	99
61) 4-Chlorophenyl-phenyle...	10.374	204	269308	42.901	ng	98
62) 4-Nitroaniline	10.404	138	134490	43.712	ng	96
63) Azobenzene	10.533	77	570513	42.745	ng	98
65) 4,6-Dinitro-2-methylph...	10.433	198	101114	49.732	ng	100
66) n-Nitrosodiphenylamine	10.498	169	493450	44.385	ng	100
67) 4-Bromophenyl-phenylether	10.869	248	166220	43.194	ng	98
68) Hexachlorobenzene	10.933	284	175203	42.921	ng	97
69) Atrazine	11.027	200	184369	49.599	ng	100
70) Pentachlorophenol	11.133	266	222097	92.914	ng	98
71) Phenanthrene	11.345	178	828491	44.839	ng	99
72) Anthracene	11.398	178	851738	47.046	ng	100
73) Carbazole	11.551	167	742030	46.223	ng	99
74) Di-n-butylphthalate	11.886	149	882495	45.361	ng	99
75) Fluoranthene	12.533	202	864959	47.410	ng	99
77) Benzidine	12.657	184	365599	46.837	ng	99
78) Pyrene	12.763	202	878288	37.578	ng	99
80) Butylbenzylphthalate	13.386	149	351233	43.015	ng	99
81) Benzo(a)anthracene	13.951	228	679307	40.430	ng	99
82) 3,3'-Dichlorobenzidine	13.915	252	131566	24.620	ng	100
83) Chrysene	13.992	228	663360	43.080	ng	99
84) Bis(2-ethylhexyl)phtha...	13.945	149	464794	44.873	ng	99
85) Di-n-octyl phthalate	14.568	149	757337	47.075	ng	99
87) Indeno(1,2,3-cd)pyrene	16.874	276	709714	46.909	ng	98
88) Benzo(b)fluoranthene	15.004	252	683657	46.365	ng	99
89) Benzo(k)fluoranthene	15.033	252	543607	41.616	ng	99
90) Benzo(a)pyrene	15.362	252	588326	47.216	ng	99
91) Dibenzo(a,h)anthracene	16.892	278	577441	47.439	ng	99
92) Benzo(g,h,i)perylene	17.309	276	556775	43.990	ng	98

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

