

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061323\
 Data File : BF133725.D
 Acq On : 13 Jun 2023 12:05
 Operator : CG\JU
 Sample : 03047-02MS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB32BMS

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 06/14/2023
 Supervised By :mohammad ahmed 06/16/2023

Quant Time: Jun 13 12:39:27 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF060923.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 13 11:26:55 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.834	152	87761	20.000	ng	0.00
21) Naphthalene-d8	8.116	136	333257	20.000	ng	0.00
39) Acenaphthene-d10	9.875	164	149569	20.000	ng	0.00
64) Phenanthrene-d10	11.363	188	233053	20.000	ng	0.00
76) Chrysene-d12	14.010	240	161968	20.000	ng	0.00
86) Perylene-d12	15.474	264	197092	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.452	112	407316	80.799	ng	0.00
7) Phenol-d6	6.469	99	491871	74.959	ng	0.00
23) Nitrobenzene-d5	7.398	82	338898	55.389	ng	0.00
42) 2,4,6-Tribromophenol	10.663	330	122648	64.781	ng	0.00
45) 2-Fluorobiphenyl	9.192	172	590400	56.112	ng	0.00
79) Terphenyl-d14	12.951	244	423139	40.424	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.616	88	81394	36.726	ng	95
3) Pyridine	3.363	79	209003	32.355	ng	95
4) n-Nitrosodimethylamine	3.322	42	156935	43.950	ng	92
6) Aniline	6.493	93	287334	34.766	ng	98
8) 2-Chlorophenol	6.616	128	235474	44.324	ng	99
9) Benzaldehyde	6.381	77	141831	32.579	ng	94
10) Phenol	6.481	94	285250	41.631	ng	93
11) bis(2-Chloroethyl)ether	6.569	93	219336	43.119	ng	95
12) 1,3-Dichlorobenzene	6.775	146	257146	45.303	ng	97
13) 1,4-Dichlorobenzene	6.851	146	261382	45.561	ng	97
14) 1,2-Dichlorobenzene	7.004	146	250667	48.355	ng	98
15) Benzyl Alcohol	6.975	79	219963	42.919	ng	97
16) 2,2'-oxybis(1-Chloropr...	7.110	45	388932	45.105	ng	99
17) 2-Methylphenol	7.087	107	189843	40.317	ng	98
18) Hexachloroethane	7.345	117	100033	50.932	ng	97
19) n-Nitroso-di-n-propyla...	7.251	70	169951	40.707	ng	99
20) 3+4-Methylphenols	7.240	107	232870	38.016	ng	93
22) Acetophenone	7.245	105	331694	43.396	ng	# 95
24) Nitrobenzene	7.416	77	263825	42.822	ng	96
25) Isophorone	7.657	82	451136	40.162	ng	100
26) 2-Nitrophenol	7.734	139	123842	44.789	ng	95
27) 2,4-Dimethylphenol	7.769	122	199313	41.822	ng	95
28) bis(2-Chloroethoxy)met...	7.869	93	262379	40.414	ng	99
29) 2,4-Dichlorophenol	7.975	162	186304	39.887	ng	97
30) 1,2,4-Trichlorobenzene	8.057	180	210254	43.145	ng	96
31) Naphthalene	8.140	128	664611	40.934	ng	100
32) Benzoic acid	7.887	122	132128	43.960	ng	94
33) 4-Chloroaniline	8.187	127	175690	25.307	ng	99
34) Hexachlorobutadiene	8.251	225	133544	41.855	ng	97
35) Caprolactam	8.563	113	53029m	33.740	ng	
36) 4-Chloro-3-methylphenol	8.669	107	204720	38.025	ng	97
37) 2-Methylnaphthalene	8.834	142	432815	38.366	ng	100
38) 1-Methylnaphthalene	8.934	142	402020	38.727	ng	99
40) 1,2,4,5-Tetrachloroben...	8.998	216	205173	45.123	ng	98
41) Hexachlorocyclopentadiene	8.981	237	225553	111.313	ng	97
43) 2,4,6-Trichlorophenol	9.110	196	127999	42.126	ng	96

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.151	196	134428	38.028	ng	98
46) 1,1'-Biphenyl	9.298	154	536514	44.476	ng	99
47) 2-Chloronaphthalene	9.322	162	389813	45.050	ng	98
48) 2-Nitroaniline	9.416	65	126207	39.881	ng	95
49) Acenaphthylene	9.739	152	620951	41.230	ng	99
50) Dimethylphthalate	9.598	163	422822	37.873	ng	99
51) 2,6-Dinitrotoluene	9.657	165	92430	40.111	ng	# 71
52) Acenaphthene	9.910	154	359207m	40.747	ng	
53) 3-Nitroaniline	9.828	138	79697	28.420	ng	91
54) 2,4-Dinitrophenol	9.939	184	87445	77.823	ng	96
55) Dibenzofuran	10.081	168	539733	40.124	ng	95
56) 4-Nitrophenol	9.986	139	133137	70.347	ng	# 70
57) 2,4-Dinitrotoluene	10.063	165	111783	38.142	ng	97
58) Fluorene	10.428	166	406242	39.492	ng	98
59) 2,3,4,6-Tetrachlorophenol	10.198	232	102341	38.669	ng	95
60) Diethylphthalate	10.298	149	412150	36.273	ng	99
61) 4-Chlorophenyl-phenyle...	10.416	204	196785	38.750	ng	99
62) 4-Nitroaniline	10.439	138	83611	30.409	ng	89
63) Azobenzene	10.575	77	426373	39.639	ng	97
65) 4,6-Dinitro-2-methylph...	10.469	198	52583	47.217	ng	# 73
66) n-Nitrosodiphenylamine	10.534	169	343408	43.868	ng	99
67) 4-Bromophenyl-phenylether	10.904	248	111917	43.026	ng	95
68) Hexachlorobenzene	10.969	284	121811	44.532	ng	95
69) Atrazine	11.063	200	97648	42.991	ng	96
70) Pentachlorophenol	11.169	266	114112	69.668	ng	96
71) Phenanthrene	11.386	178	495424	41.900	ng	99
72) Anthracene	11.439	178	497801	40.377	ng	99
73) Carbazole	11.592	167	426017	37.101	ng	98
74) Di-n-butylphthalate	11.922	149	523220	35.544	ng	99
75) Fluoranthene	12.575	202	462846	36.862	ng	98
77) Benzidine	12.698	184	191914	35.311	ng	99
78) Pyrene	12.810	202	466047	30.889	ng	100
80) Butylbenzylphthalate	13.427	149	196743	30.888	ng	99
81) Benzo(a)anthracene	13.998	228	455102	40.622	ng	99
82) 3,3'-Dichlorobenzidine	13.963	252	147334	39.556	ng	99
83) Chrysene	14.033	228	429517	40.534	ng	99
84) Bis(2-ethylhexyl)phtha...	13.986	149	274969	35.060	ng	98
85) Di-n-octyl phthalate	14.604	149	534382	47.948	ng	98
87) Indeno(1,2,3-cd)pyrene	16.945	276	546135	40.277	ng	99
88) Benzo(b)fluoranthene	15.051	252	514824	43.001	ng	99
89) Benzo(k)fluoranthene	15.080	252	466775	40.023	ng	97
90) Benzo(a)pyrene	15.416	252	444344	39.990	ng	98
91) Dibenzo(a,h)anthracene	16.963	278	459033	40.891	ng	99
92) Benzo(g,h,i)perylene	17.392	276	412378	37.051	ng	98

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

