

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042623\
 Data File : BF133074.D
 Acq On : 26 Apr 2023 20:02
 Operator : CG\JU
 Sample : 02491-01MSD
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :

BNA_F

ClientSampleId :

S2S-TP3MSD

Manual Integrations

APPROVED

Reviewed By : Christian Giraldo 04/27/2023
 Supervised By : Jagrut Upadhyay 04/27/2023

Quant Time: Apr 27 02:00:50 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042123.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 27 01:56:03 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.734	152	168004	20.000	ng	0.00
21) Naphthalene-d8	8.022	136	691110	20.000	ng	0.00
39) Acenaphthene-d10	9.774	164	356311	20.000	ng	0.00
64) Phenanthrene-d10	11.257	188	645956	20.000	ng	0.00
76) Chrysene-d12	13.892	240	371338	20.000	ng	0.00
86) Perylene-d12	15.304	264	353101	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.351	112	1193003	107.269	ng	0.01
7) Phenol-d6	6.381	99	1509565	109.355	ng	0.00
23) Nitrobenzene-d5	7.304	82	912627	71.278	ng	0.00
42) 2,4,6-Tribromophenol	10.563	330	389596	108.723	ng	0.00
45) 2-Fluorobiphenyl	9.098	172	1562805	73.379	ng	0.00
79) Terphenyl-d14	12.839	244	1799978	83.599	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.463	88	230187	44.622	ng	96
3) Pyridine	3.157	79	582522	42.516	ng	95
4) n-Nitrosodimethylamine	3.116	42	302924	42.685	ng	92
6) Aniline	6.398	93	630396	35.456	ng	# 89
8) 2-Chlorophenol	6.522	128	574385	50.866	ng	95
9) Benzaldehyde	6.281	77	332482	63.491	ng	99
10) Phenol	6.392	94	699736	45.949	ng	80
11) bis(2-Chloroethyl)ether	6.475	93	540561	47.303	ng	97
12) 1,3-Dichlorobenzene	6.675	146	599306	49.919	ng	99
13) 1,4-Dichlorobenzene	6.751	146	606816	50.433	ng	99
14) 1,2-Dichlorobenzene	6.904	146	577516	52.062	ng	97
15) Benzyl Alcohol	6.881	79	468672	49.150	ng	98
16) 2,2'-oxybis(1-Chloropr...	7.016	45	885908	44.634	ng	97
17) 2-Methylphenol	6.998	107	463755	44.851	ng	99
18) Hexachloroethane	7.251	117	205778	49.192	ng	98
19) n-Nitroso-di-n-propyla...	7.157	70	376396	43.798	ng	97
20) 3+4-Methylphenols	7.151	107	568463	46.684	ng	98
22) Acetophenone	7.145	105	750590	46.885	ng	# 97
24) Nitrobenzene	7.322	77	575773	44.377	ng	100
25) Isophorone	7.563	82	1081589	44.491	ng	99
26) 2-Nitrophenol	7.639	139	319696	51.086	ng	92
27) 2,4-Dimethylphenol	7.681	122	521787	48.626	ng	100
28) bis(2-Chloroethoxy)met...	7.775	93	642388	44.668	ng	99
29) 2,4-Dichlorophenol	7.880	162	473034	48.424	ng	100
30) 1,2,4-Trichlorobenzene	7.963	180	488838	48.303	ng	100
31) Naphthalene	8.045	128	1590417	46.028	ng	99
32) Benzoic acid	7.804	122	349075m	52.665	ng	
33) 4-Chloroaniline	8.092	127	233430	16.720	ng	98
34) Hexachlorobutadiene	8.163	225	266505	47.513	ng	99
35) Caprolactam	8.469	113	172356m	52.521	ng	
36) 4-Chloro-3-methylphenol	8.580	107	523690	49.713	ng	100
37) 2-Methylnaphthalene	8.733	142	1060658	44.899	ng	98
38) 1-Methylnaphthalene	8.833	142	999470	45.227	ng	99
40) 1,2,4,5-Tetrachloroben...	8.904	216	443336	46.276	ng	99
41) Hexachlorocyclopentadiene	8.892	237	509044	91.109	ng	98
43) 2,4,6-Trichlorophenol	9.016	196	326173	49.231	ng	99

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Thu Apr 27 01:56:03 2023

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.051	196	346255	45.189	ng	99
46) 1,1'-Biphenyl	9.198	154	1304779	46.181	ng	99
47) 2-Chloronaphthalene	9.222	162	988224	47.787	ng	100
48) 2-Nitroaniline	9.316	65	322344	47.163	ng	95
49) Acenaphthylene	9.633	152	1550059	46.910	ng	100
50) Dimethylphthalate	9.504	163	1169879	47.087	ng	98
51) 2,6-Dinitrotoluene	9.563	165	276995	49.535	ng	# 81
52) Acenaphthene	9.810	154	1112348m	50.262	ng	
53) 3-Nitroaniline	9.727	138	207970	31.279	ng	95
54) 2,4-Dinitrophenol	9.833	184	284303	101.147	ng	98
55) Dibenzofuran	9.980	168	1395897	46.239	ng	99
56) 4-Nitrophenol	9.898	139	505519	98.164	ng	98
57) 2,4-Dinitrotoluene	9.963	165	367261	51.500	ng	91
58) Fluorene	10.322	166	1031387	46.632	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.098	232	292674	49.269	ng	98
60) Diethylphthalate	10.204	149	1099935	46.871	ng	98
61) 4-Chlorophenyl-phenyle...	10.316	204	499105	46.307	ng	100
62) 4-Nitroaniline	10.345	138	315288	51.593	ng	95
63) Azobenzene	10.474	77	1138687	45.804	ng	98
65) 4,6-Dinitro-2-methylph...	10.369	198	207206	52.922	ng	85
66) n-Nitrosodiphenylamine	10.433	169	961597	45.893	ng	99
67) 4-Bromophenyl-phenylether	10.804	248	320249	45.712	ng	98
68) Hexachlorobenzene	10.869	284	350849	48.873	ng	98
69) Atrazine	10.963	200	333864	55.298	ng	98
70) Pentachlorophenol	11.063	266	402138	78.656	ng	97
71) Phenanthrene	11.280	178	1586059	45.684	ng	99
72) Anthracene	11.333	178	1644887	46.295	ng	100
73) Carbazole	11.486	167	1489466	47.080	ng	100
74) Di-n-butylphthalate	11.821	149	1740941	48.626	ng	100
75) Fluoranthene	12.463	202	1648320	47.306	ng	100
77) Benzidine	12.586	184	707700	112.700	ng	# 93
78) Pyrene	12.692	202	1680999	52.808	ng	99
80) Butylbenzylphthalate	13.315	149	691409	61.344	ng	95
81) Benzo(a)anthracene	13.880	228	1212114	47.467	ng	99
82) 3,3'-Dichlorobenzidine	13.845	252	227758	32.286	ng	99
83) Chrysene	13.915	228	1168957	45.789	ng	98
84) Bis(2-ethylhexyl)phtha...	13.880	149	790487	55.876	ng	99
85) Di-n-octyl phthalate	14.492	149	1265031	54.840	ng	97
87) Indeno(1,2,3-cd)pyrene	16.686	276	1321137	65.777	ng	100
88) Benzo(b)fluoranthene	14.904	252	1063974	47.364	ng	98
89) Benzo(k)fluoranthene	14.933	252	1018500	45.738	ng	99
90) Benzo(a)pyrene	15.251	252	946075	46.270	ng	99
91) Dibenzo(a,h)anthracene	16.709	278	1101588	65.759	ng	99
92) Benzo(g,h,i)perylene	17.103	276	1103462	66.721	ng	98

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

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