

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF010324\
 Data File : BF136908.D
 Acq On : 03 Jan 2024 15:23
 Operator : CG\JU
 Sample : 06038-02MS
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 T-402MS

Quant Time: Jan 03 16:10:11 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122023.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 21 01:54:25 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.781	152	82442	20.000	ng	-0.01
21) Naphthalene-d8	8.057	136	327893	20.000	ng	-0.01
39) Acenaphthene-d10	9.816	164	166394	20.000	ng	-0.01
64) Phenanthrene-d10	11.310	188	300770	20.000	ng	0.00
76) Chrysene-d12	13.969	240	157812	20.000	ng	# 0.00
86) Perylene-d12	15.445	264	164013	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.446	112	577571	109.310	ng	0.02
7) Phenol-d6	6.440	99	688915	100.623	ng	0.00
23) Nitrobenzene-d5	7.351	82	489711	76.424	ng	0.00
42) 2,4,6-Tribromophenol	10.616	330	234058	119.438	ng	0.00
45) 2-Fluorobiphenyl	9.139	172	987621	79.831	ng	-0.01
79) Terphenyl-d14	12.904	244	975318	86.367	ng	-0.01
Target Compounds						Qvalue
2) 1,4-Dioxane	2.746	88	72779	33.058	ng	# 82
3) Pyridine	3.487	79	154902	28.838	ng	93
4) n-Nitrosodimethylamine	3.416	42	106531	37.138	ng	98
6) Aniline	6.451	93	89793	13.118	ng	# 1
8) 2-Chlorophenol	6.575	128	235622	40.749	ng	95
9) Benzaldehyde	6.340	77	11134	2.859	ng	92
10) Phenol	6.451	94	256616	35.111	ng	92
11) bis(2-Chloroethyl)ether	6.528	93	225582	40.586	ng	96
12) 1,3-Dichlorobenzene	6.722	146	228497	36.225	ng	97
13) 1,4-Dichlorobenzene	6.798	146	227577	35.310	ng	98
14) 1,2-Dichlorobenzene	6.951	146	215936	36.006	ng	98
15) Benzyl Alcohol	6.934	79	199930	43.285	ng	98
16) 2,2'-oxybis(1-Chloropr...	7.057	45	342677	37.710	ng	63
17) 2-Methylphenol	7.051	107	188978	39.451	ng	94
18) Hexachloroethane	7.287	117	80793	34.517	ng	95
19) n-Nitroso-di-n-propyla...	7.204	70	164678	39.679	ng	94
20) 3+4-Methylphenols	7.204	107	239512	40.023	ng	# 62
22) Acetophenone	7.192	105	339712	41.334	ng	# 88
24) Nitrobenzene	7.369	77	237313	36.971	ng	96
25) Isophorone	7.604	82	465464	39.921	ng	99
26) 2-Nitrophenol	7.681	139	127294	41.449	ng	97
27) 2,4-Dimethylphenol	7.722	122	190450	50.938	ng	98
28) bis(2-Chloroethoxy)met...	7.816	93	287345	40.544	ng	98
29) 2,4-Dichlorophenol	7.928	162	198660	41.271	ng	96
30) 1,2,4-Trichlorobenzene	8.004	180	201900	37.646	ng	98
31) Naphthalene	8.087	128	1051380	58.366	ng	100
32) Benzoic acid	7.881	122	154423	42.686	ng	99
33) 4-Chloroaniline	8.145	127	16298	2.450	ng	98
34) Hexachlorobutadiene	8.192	225	115165	35.344	ng	98
35) Caprolactam	8.522	113	58617	36.914	ng	96
36) 4-Chloro-3-methylphenol	8.639	107	219508	40.508	ng	99
37) 2-Methylnaphthalene	8.775	142	498383	42.990	ng	97
38) 1-Methylnaphthalene	8.875	142	483437	42.505	ng	99
40) 1,2,4,5-Tetrachloroben...	8.939	216	222527	43.183	ng	98
41) Hexachlorocyclopentadiene	8.922	237	148736	89.880	ng	100
43) 2,4,6-Trichlorophenol	9.063	196	147381	44.596	ng	98

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44) 2,4,5-Trichlorophenol	9.110	196	156469	42.487	ng	98
46) 1,1'-Biphenyl	9.239	154	613731	43.954	ng	99
47) 2-Chloronaphthalene	9.263	162	437308	41.198	ng	99
48) 2-Nitroaniline	9.369	65	133572	40.184	ng	97
49) Acenaphthylene	9.681	152	723720	44.606	ng	99
50) Dimethylphthalate	9.551	163	528455	43.481	ng	100
51) 2,6-Dinitrotoluene	9.616	165	115094	41.726	ng	97
52) Acenaphthene	9.851	154	438243	43.574	ng	100
53) 3-Nitroaniline	9.781	138	26796	9.171	ng	85
54) 2,4-Dinitrophenol	9.898	184	109601	72.082	ng	92
55) Dibenzofuran	10.028	168	628574	41.230	ng	98
56) 4-Nitrophenol	9.969	139	133733	67.799	ng	95
57) 2,4-Dinitrotoluene	10.022	165	143037	39.945	ng	93
58) Fluorene	10.369	166	523910	43.310	ng	98
59) 2,3,4,6-Tetrachlorophenol	10.151	232	128186	45.217	ng	99
60) Diethylphthalate	10.245	149	539247	42.762	ng	99
61) 4-Chlorophenyl-phenyle...	10.363	204	239237	40.687	ng	97
62) 4-Nitroaniline	10.404	138	90008	31.954	ng	93
63) Azobenzene	10.522	77	480027	40.756	ng	98
65) 4,6-Dinitro-2-methylph...	10.433	198	71298	41.363	ng	87
66) n-Nitrosodiphenylamine	10.486	169	448140	45.867	ng	99
67) 4-Bromophenyl-phenylether	10.851	248	150916	45.181	ng	98
68) Hexachlorobenzene	10.916	284	164627	44.226	ng	95
69) Atrazine	11.022	200	106244	42.515	ng	99
70) Pentachlorophenol	11.122	266	159391	91.864	ng	99
71) Phenanthrene	11.339	178	723935	46.909	ng	100
72) Anthracene	11.392	178	710419	45.455	ng	99
73) Carbazole	11.551	167	608784	41.314	ng	99
74) Di-n-butylphthalate	11.875	149	847029	47.107	ng	100
75) Fluoranthene	12.527	202	668776	40.513	ng	98
77) Benzidine	12.657	184	31161	6.697	ng	99
78) Pyrene	12.757	202	654138	42.225	ng	99
80) Butylbenzylphthalate	13.380	149	275570	48.885	ng	98
81) Benzo(a)anthracene	13.957	228	493669	45.050	ng	99
82) 3,3'-Dichlorobenzidine	13.916	252	73659	23.669	ng	# 98
83) Chrysene	13.992	228	463175	43.222	ng	98
84) Bis(2-ethylhexyl)phtha...	13.945	149	397385	56.029	ng	97
85) Di-n-octyl phthalate	14.563	149	643476	58.653	ng	97
87) Indeno(1,2,3-cd)pyrene	16.962	276	530255	44.778	ng	99
88) Benzo(b)fluoranthene	15.010	252	478925	43.719	ng	99
89) Benzo(k)fluoranthene	15.045	252	432042	41.739	ng	99
90) Benzo(a)pyrene	15.386	252	414047	44.777	ng	98
91) Dibenzo(a,h)anthracene	16.974	278	444689	45.773	ng	98
92) Benzo(g,h,i)perylene	17.415	276	407332	40.936	ng	99

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

