

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081324\
 Data File : BF138965.D
 Acq On : 13 Aug 2024 10:08
 Operator : RC/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Quant Time: Aug 13 11:45:55 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF073024.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 30 17:50:01 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.839	152	46987	20.000	ng	0.00
21) Naphthalene-d8	8.122	136	189694	20.000	ng	0.00
39) Acenaphthene-d10	9.875	164	108534	20.000	ng	0.00
64) Phenanthrene-d10	11.363	188	193718	20.000	ng	0.00
76) Chrysene-d12	13.998	240	93804	20.000	ng	0.00
86) Perylene-d12	15.462	264	93639	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.469	112	241732	79.416	ng	0.00
7) Phenol-d6	6.492	99	325552	79.661	ng	0.00
23) Nitrobenzene-d5	7.410	82	329140	84.832	ng	0.00
42) 2,4,6-Tribromophenol	10.669	330	74564	83.870	ng	0.00
45) 2-Fluorobiphenyl	9.198	172	596252	82.543	ng	0.00
79) Terphenyl-d14	12.939	244	504053	89.966	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.587	88	47721	35.810	ng	97
3) Pyridine	3.346	79	115174	35.677	ng	96
4) n-Nitrosodimethylamine	3.316	42	85416	44.426	ng	# 80
6) Aniline	6.510	93	137097	37.616	ng	# 65
8) 2-Chlorophenol	6.634	128	127368	39.771	ng	100
9) Benzaldehyde	6.398	77	90277	36.851	ng	99
10) Phenol	6.504	94	168509	39.162	ng	# 76
11) bis(2-Chloroethyl)ether	6.581	93	126683	38.259	ng	99
12) 1,3-Dichlorobenzene	6.781	146	142552	39.765	ng	98
13) 1,4-Dichlorobenzene	6.857	146	144375	39.907	ng	98
14) 1,2-Dichlorobenzene	7.010	146	139453	41.246	ng	98
15) Benzyl Alcohol	6.992	79	126510	42.950	ng	98
16) 2,2'-oxybis(1-Chloropr...	7.116	45	210772	36.988	ng	66
17) 2-Methylphenol	7.104	107	104957	39.689	ng	# 91
18) Hexachloroethane	7.345	117	55102	40.463	ng	96
19) n-Nitroso-di-n-propyla...	7.257	70	103190	41.806	ng	97
20) 3+4-Methylphenols	7.263	107	143621	42.329	ng	# 81
22) Acetophenone	7.257	105	196063	42.213	ng	96
24) Nitrobenzene	7.428	77	161626	40.938	ng	99
25) Isophorone	7.663	82	271632	41.000	ng	98
26) 2-Nitrophenol	7.745	139	69439	40.880	ng	95
27) 2,4-Dimethylphenol	7.781	122	81769	40.235	ng	98
28) bis(2-Chloroethoxy)met...	7.869	93	157086	38.936	ng	99
29) 2,4-Dichlorophenol	7.992	162	107612	41.207	ng	99
30) 1,2,4-Trichlorobenzene	8.063	180	123977	41.137	ng	99
31) Naphthalene	8.145	128	402596	40.320	ng	100
32) Benzoic acid	7.934	122	41753	26.136	ng	98
33) 4-Chloroaniline	8.198	127	128082	38.214	ng	98
34) Hexachlorobutadiene	8.251	225	80583	44.145	ng	97
35) Caprolactam	8.581	113	33299	42.733	ng	91
36) 4-Chloro-3-methylphenol	8.686	107	126700	42.452	ng	98
37) 2-Methylnaphthalene	8.833	142	263939	41.855	ng	100
38) 1-Methylnaphthalene	8.933	142	258191	41.783	ng	99
40) 1,2,4,5-Tetrachloroben...	8.998	216	120374	39.926	ng	99
41) Hexachlorocyclopentadiene	8.980	237	22558	32.772	ng	98
43) 2,4,6-Trichlorophenol	9.122	196	70499	38.351	ng	100

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.169	196	77343	38.487	ng	98
46) 1,1'-Biphenyl	9.298	154	339171	39.901	ng	99
47) 2-Chloronaphthalene	9.322	162	250981	39.700	ng	99
48) 2-Nitroaniline	9.428	65	91268	42.585	ng	95
49) Acenaphthylene	9.739	152	357198	39.838	ng	100
50) Dimethylphthalate	9.598	163	291241	41.967	ng	99
51) 2,6-Dinitrotoluene	9.669	165	67281	42.959	ng	87
52) Acenaphthene	9.910	154	238425	39.557	ng	99
53) 3-Nitroaniline	9.845	138	64317	39.724	ng	97
54) 2,4-Dinitrophenol	9.957	184	26906	37.319	ng	# 78
55) Dibenzofuran	10.080	168	349823	41.116	ng	99
56) 4-Nitrophenol	10.022	139	41207	42.323	ng	87
57) 2,4-Dinitrotoluene	10.075	165	89344	44.712	ng	# 80
58) Fluorene	10.422	166	291898	43.082	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.204	232	63795	41.523	ng	97
60) Diethylphthalate	10.298	149	299444	45.507	ng	100
61) 4-Chlorophenyl-phenyle...	10.410	204	144826	43.462	ng	95
62) 4-Nitroaniline	10.457	138	63424	41.221	ng	84
63) Azobenzene	10.575	77	306678	42.022	ng	96
65) 4,6-Dinitro-2-methylph...	10.486	198	45307	38.336	ng	91
66) n-Nitrosodiphenylamine	10.533	169	236601	39.074	ng	99
67) 4-Bromophenyl-phenylether	10.904	248	81521	38.868	ng	97
68) Hexachlorobenzene	10.969	284	85468	39.467	ng	97
69) Atrazine	11.063	200	64679	41.401	ng	96
70) Pentachlorophenol	11.174	266	34309	35.149	ng	98
71) Phenanthrene	11.386	178	399518	40.052	ng	99
72) Anthracene	11.439	178	393434	40.037	ng	100
73) Carbazole	11.598	167	338647	39.944	ng	99
74) Di-n-butylphthalate	11.916	149	448632	47.073	ng	100
75) Fluoranthene	12.574	202	381687	40.988	ng	98
77) Benzidine	12.698	184	65303	29.106	ng	98
78) Pyrene	12.804	202	372704	42.200	ng	99
80) Butylbenzylphthalate	13.410	149	132619	46.891	ng	99
81) Benzo(a)anthracene	13.992	228	265759	41.142	ng	99
82) 3,3'-Dichlorobenzidine	13.951	252	67501	40.835	ng	98
83) Chrysene	14.027	228	230220	39.504	ng	99
84) Bis(2-ethylhexyl)phtha...	13.963	149	167488	40.442	ng	98
85) Di-n-octyl phthalate	14.574	149	266460	34.775	ng	100
87) Indeno(1,2,3-cd)pyrene	16.939	276	234418	34.933	ng	96
88) Benzo(b)fluoranthene	15.039	252	239407	41.244	ng	97
89) Benzo(k)fluoranthene	15.062	252	207707	41.328	ng	98
90) Benzo(a)pyrene	15.404	252	200130	40.989	ng	97
91) Dibenzo(a,h)anthracene	16.945	278	191802	34.819	ng	98
92) Benzo(g,h,i)perylene	17.380	276	187222	32.753	ng	97

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

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