

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM051324\
 Data File : BM045718.D
 Acq On : 13 May 2024 17:58
 Operator : MA/JU
 Sample : SSTDICC010
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC010

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 05/14/2024
 Supervised By :mohammad ahmed 05/17/2024

Quant Time: May 14 02:25:46 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM051324.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 14 02:21:04 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.587	152	100363	20.000	ng	0.00
21) Naphthalene-d8	10.357	136	420425	20.000	ng	0.00
39) Acenaphthene-d10	14.216	164	264577	20.000	ng	0.00
64) Phenanthrene-d10	16.962	188	674257	20.000	ng	0.00
76) Chrysene-d12	21.180	240	385634	20.000	ng	0.00
86) Perylene-d12	23.985	264	423896	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.275	112	113068	19.333	ng	0.00
7) Phenol-d6	6.875	99	128562	18.027	ng	0.00
23) Nitrobenzene-d5	8.757	82	179138	19.570	ng	0.00
42) 2,4,6-Tribromophenol	15.721	330	62598	17.328	ng	0.00
45) 2-Fluorobiphenyl	12.845	172	353401	18.410	ng	0.00
79) Terphenyl-d14	19.598	244	664361	16.813	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.222	88	15811m	9.918	ng	Qvalue
3) Pyridine	3.681	79	70080m	11.171	ng	
4) n-Nitrosodimethylamine	3.534	42	31910	9.654	ng	# 95
6) Aniline	6.975	93	58568	9.959	ng	96
8) 2-Chlorophenol	7.192	128	49339	9.138	ng	99
9) Benzaldehyde	6.775	77	57689	11.303	ng	96
10) Phenol	6.898	94	61302	9.222	ng	100
11) bis(2-Chloroethyl)ether	7.051	93	59134	9.990	ng	98
12) 1,3-Dichlorobenzene	7.475	146	69960	9.560	ng	98
13) 1,4-Dichlorobenzene	7.622	146	71290	9.469	ng	99
14) 1,2-Dichlorobenzene	7.939	146	64319	9.573	ng	98
15) Benzyl Alcohol	7.898	79	61337	9.321	ng	93
16) 2,2'-oxybis(1-Chloropr...	8.145	45	58441	10.013	ng	87
17) 2-Methylphenol	8.110	107	43547	9.167	ng	95
18) Hexachloroethane	8.657	117	29570	9.127	ng	95
19) n-Nitroso-di-n-propyla...	8.445	70	47100	8.792	ng	94
20) 3+4-Methylphenols	8.445	107	59022	8.793	ng	# 87
22) Acetophenone	8.451	105	89263	9.424	ng	# 99
24) Nitrobenzene	8.798	77	90016	10.162	ng	99
25) Isophorone	9.369	82	125657	9.495	ng	98
26) 2-Nitrophenol	9.504	139	26878	8.936	ng	98
27) 2,4-Dimethylphenol	9.616	122	35547	9.408	ng	97
28) bis(2-Chloroethoxy)met...	9.828	93	70419	9.195	ng	98
29) 2,4-Dichlorophenol	10.057	162	48705	8.760	ng	97
30) 1,2,4-Trichlorobenzene	10.216	180	63656	9.461	ng	99
31) Naphthalene	10.404	128	197736	9.506	ng	98
32) Benzoic acid	9.822	122	29722m	8.015	ng	
33) 4-Chloroaniline	10.586	127	69147	9.261	ng	98
34) Hexachlorobutadiene	10.692	225	40345	9.408	ng	98
35) Caprolactam	11.557	113	14196	7.829	ng	89
36) 4-Chloro-3-methylphenol	11.739	107	57290	8.393	ng	95
37) 2-Methylnaphthalene	12.022	142	138300	9.091	ng	96
38) 1-Methylnaphthalene	12.245	142	134357	8.969	ng	96
40) 1,2,4,5-Tetrachloroben...	12.386	216	67342	9.589	ng	99
41) Hexachlorocyclopentadiene	12.374	237	29070	9.441	ng	98
43) 2,4,6-Trichlorophenol	12.669	196	41767	9.067	ng	99

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44) 2,4,5-Trichlorophenol	12.763	196	48723	9.420	ng	98
46) 1,1'-Biphenyl	13.051	154	186962	9.838	ng	98
47) 2-Chloronaphthalene	13.086	162	141254	9.609	ng	99
48) 2-Nitroaniline	13.339	65	45071	9.340	ng	96
49) Acenaphthylene	13.939	152	215230	9.419	ng	99
50) Dimethylphthalate	13.727	163	188268	9.764	ng	100
51) 2,6-Dinitrotoluene	13.821	165	39834	9.552	ng	94
52) Acenaphthene	14.280	154	129407	9.132	ng	97
53) 3-Nitroaniline	14.180	138	38914	9.494	ng	97
54) 2,4-Dinitrophenol	14.357	184	16667	11.441	ng	89
55) Dibenzofuran	14.621	168	219694	9.347	ng	98
56) 4-Nitrophenol	14.563	139	31662	8.975	ng	96
57) 2,4-Dinitrotoluene	14.610	165	56956	9.015	ng	92
58) Fluorene	15.268	166	190609	9.022	ng	100
59) 2,3,4,6-Tetrachlorophenol	14.874	232	44076	9.283	ng	96
60) Diethylphthalate	15.086	149	213264	9.728	ng	98
61) 4-Chlorophenyl-phenyle...	15.274	204	93254	9.102	ng	98
62) 4-Nitroaniline	15.357	138	40058	8.971	ng	88
63) Azobenzene	15.568	77	218622	10.165	ng	98
65) 4,6-Dinitro-2-methylph...	15.368	198	26930	10.938	ng	91
66) n-Nitrosodiphenylamine	15.498	169	149683	8.556	ng	98
67) 4-Bromophenyl-phenylether	16.162	248	60931	9.124	ng	95
68) Hexachlorobenzene	16.262	284	73901	9.206	ng	96
69) Atrazine	16.480	200	54459	9.879	ng	96
70) Pentachlorophenol	16.639	266	42584	8.374	ng	95
71) Phenanthrene	17.004	178	307801	9.054	ng	100
72) Anthracene	17.092	178	288534	8.810	ng	99
73) Carbazole	17.386	167	283056	8.911	ng	98
74) Di-n-butylphthalate	17.956	149	365756	8.857	ng	99
75) Fluoranthene	19.021	202	375220	9.070	ng	99
77) Benzidine	19.245	184	68340	8.127	ng	99
78) Pyrene	19.386	202	385789	9.807	ng	100
80) Butylbenzylphthalate	20.309	149	144198	9.905	ng	100
81) Benzo(a)anthracene	21.162	228	231954	8.789	ng	99
82) 3,3'-Dichlorobenzidine	21.109	252	82655	8.405	ng	98
83) Chrysene	21.215	228	208197	8.603	ng	99
84) Bis(2-ethylhexyl)phtha...	21.109	149	149808	9.362	ng	98
85) Di-n-octyl phthalate	22.174	149	243386	9.150	ng	# 99
87) Indeno(1,2,3-cd)pyrene	27.091	276	251341	9.422	ng	# 100
88) Benzo(b)fluoranthene	23.103	252	236591	8.635	ng	99
89) Benzo(k)fluoranthene	23.162	252	221065	8.808	ng	100
90) Benzo(a)pyrene	23.850	252	202649	8.526	ng	98
91) Dibenzo(a,h)anthracene	27.144	278	210075	9.347	ng	98
92) Benzo(g,h,i)perylene	28.050	276	215587	9.326	ng	99

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

